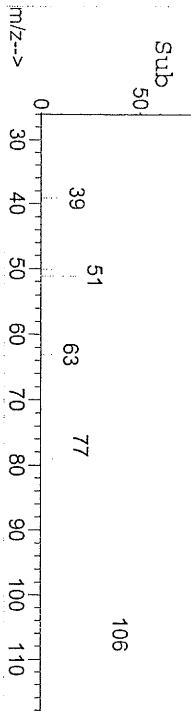
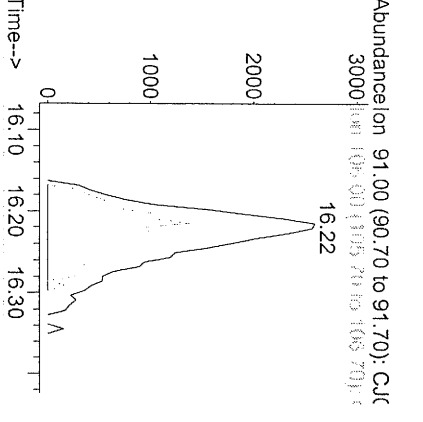
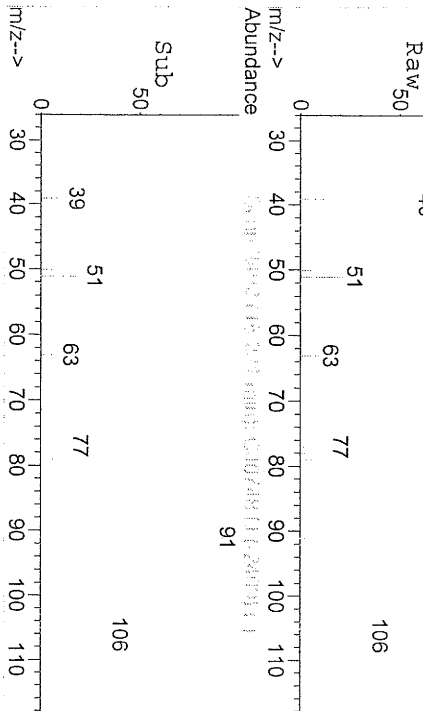


#65
 O-Xylene 0.20 PPB
 Concn: 16.22 min Scan# 2492
 Delta R.T. -0.01 min
 Lab File: CJ0748.D
 Acq: 29 Oct 2005 17:29

Tgt Ion	Ratio	Resp
91	100	Lower Upper
106	44.2	43.4 65.0



0000

LANCASTER LABORATORIES
VOLATILE ORGANICS IN AIR
SUMMA CANISTER SAMPLE
ANALYSIS DATA SHEET

Sample No.: 1FLDS
 Lab Sample ID: 4628493 Date Collected: 10/19/05 Date Received: 10/20/05
 Canister ID: SUMMA0016 Date Analyzed: 10/29/05 Time Analyzed: 18:45
 Injection Volume: 750 cc ²³² Pressure Rec'd: 7.6 psia Final Pressure: 22.9 psia
 Instrument ID: HP09464 ⁴¹²⁶⁵ Nominal Volume: 250 cc Dilution Factor: 1.0
 Lab File ID: C:\MSDCHEM\1\DATA\OCT29\CJ0750.D

CAS RN	COMPOUND NAME	CONCENTRATION	UNITS	MDL	ug/m3	Q
75-71-8	Dichlorodifluoromethane		4			J
76-14-2	Freon 114		1			U
74-87-3	Chloromethane		0.4			U
75-01-4	Vinyl Chloride		0.5			U
74-83-9	Bromomethane		0.8			U
75-00-3	Chloroethane		0.5			U
75-69-4	Trichlorofluoromethane		2			J
75-35-4	1,1-Dichloroethene		0.8			U
76-13-1	Freon 113		4			U
107-05-1	3-Chloropropene		2			U
75-09-2	Methylene Chloride		5			U
75-34-3	1,1-Dichloroethane		0.8			U
156-59-2	cis-1,2-Dichloroethene		3			J
67-66-3	Chloroform		1			U
71-55-6	1,1,1-Trichloroethane		1			U
56-23-5	Carbon Tetrachloride		1			U
107-06-2	1,2-Dichloroethane		1			J
71-43-2	Benzene		2			J
79-01-6	Trichloroethene		1			U
78-87-5	1,2-Dichloropropane		0.9			U
10061-01-5	cis-1,3-Dichloropropene		0.9			U
108-88-3	Toluene		24			U
10061-02-6	trans-1,3-Dichloropropene		0.9			U
79-00-5	1,1,2-Trichloroethane		1			U
127-18-4	Tetrachloroethene		32			U
106-93-4	1,2-Dibromoethane		2			U
108-90-7	Chlorobenzene		0.9			U
100-41-4	Ethylbenzene		5			U
1330-20-7	m/p-Xylene		13			J
95-47-6	o-Xylene		4			J
100-42-5	Styrene		3			J
79-34-5	1,1,2,2-Tetrachloroethane		1			U
622-96-8	4-Ethyltoluene		3			J
108-67-8	1,3,5-Trimethylbenzene		2			J

U = Compound was undetected at the specified limit of quantitation.
 B = Compound was found in method blank. D = analysis of diluted sample.
 J = Compound was detected, but below the limit of quantitation.

LANCASTER LABORATORIES
VOLATILE ORGANICS IN AIR
SUMMA CANISTER SAMPLE
ANALYSIS DATA SHEET

Sample No.: 1F1DS
 Lab Sample ID: 4628493 Date Collected: 10/19/05 Date Received: 10/20/05
 Canister ID: SUMMA0016 Date Analyzed: 10/29/05 Time Analyzed: 18:45
 Injection Volume: 750 cc ~~750~~ Pressure Rec'd: 7.6 psia Final Pressure: 22.9 psia
 Instrument ID: HP09464 ~~750~~ ^{4/4/05} Lab File ID: C:\MSDCHEM\1\DATA\OCT29\CJ0750.D Dilution Factor: 1.0

CAS RN	COMPOUND NAME	CONCENTRATION	UNITS	MDL	ug/m3	Q
95-63-6	1,2,4-Trimethylbenzene		3			J
541-73-1	1,3-Dichlorobenzene		3			U
106-46-7	1,4-Dichlorobenzene		3			U
95-50-1	1,2-Dichlorobenzene		3			U
120-82-1	1,2,4-Trichlorobenzene		7			U

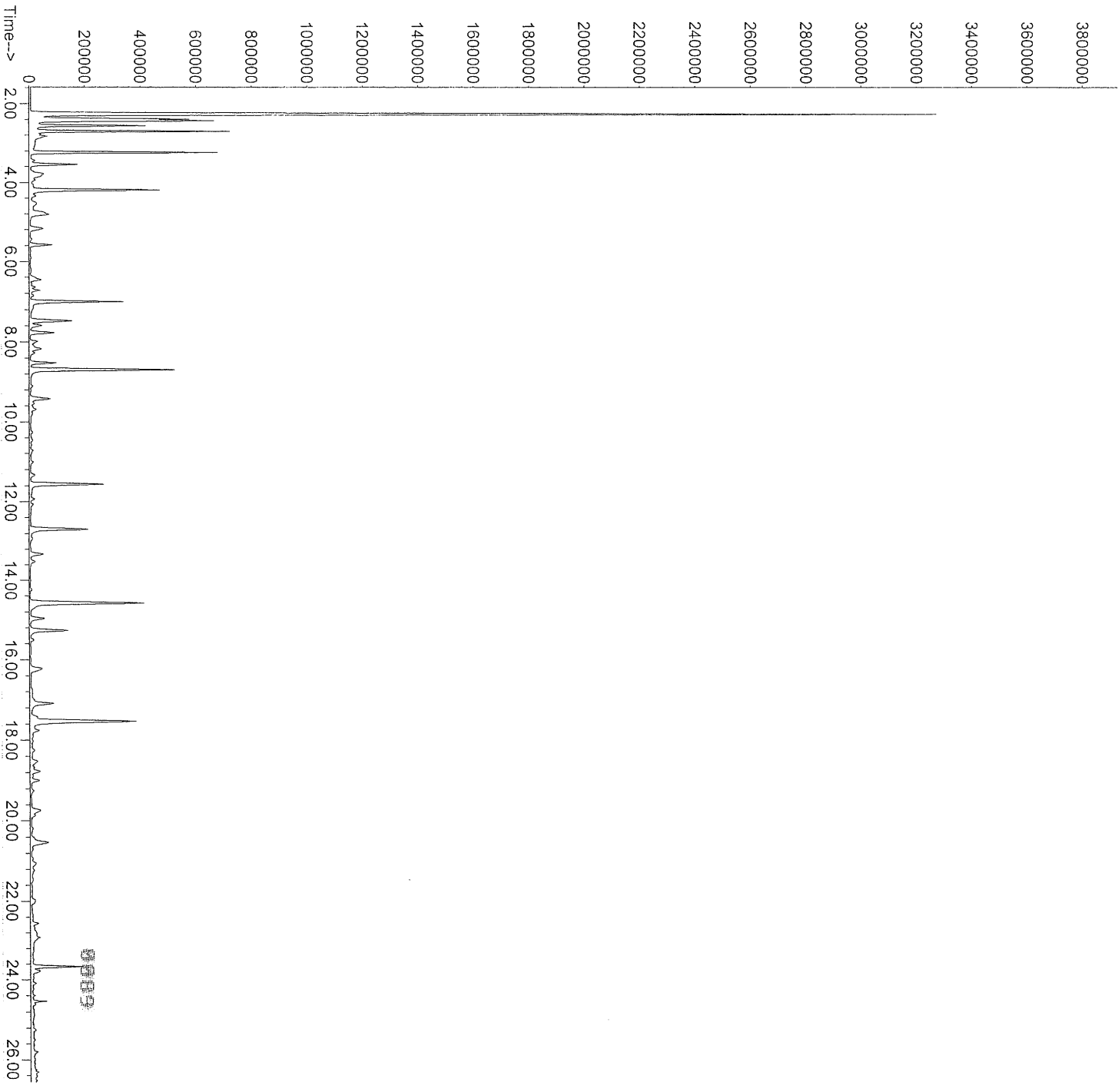
U = Compound was undetected at the specified limit of quantitation.
 B = Compound was found in method blank. D = analysis of diluted sample.
 J = Compound was detected, but below the limit of quantitation.

Data Path : C:\MSDCHEM\1\DATA\OCT29\
Data File : CJ0750.D
Acq On : 29 Oct 2005 18:45
Operator : JBS
Sample : 4628493 750CC
Misc :
ALS Vial : 45 Sample Multiplier: 1

Quant Time: Nov 07 23:08:40 2005
Quant Method : C:\MSDCHEM\1\METHODS\ALLTO15.M
Quant Title :
Quant Update : Sat Oct 29 13:19:44 2005
Response via : Initial Calibration

Abundance

TIC: CJ0750.D



Data Path : C:\MSDCHEM\1\DATA\OCT29\
 Data File : CJ0750.D
 Acq On : 29 Oct 2005 18:45
 Operator : JBS
 Sample : 4628493 750CC
 Misc :
 ALS Vial : 45 Sample Multiplier: 1

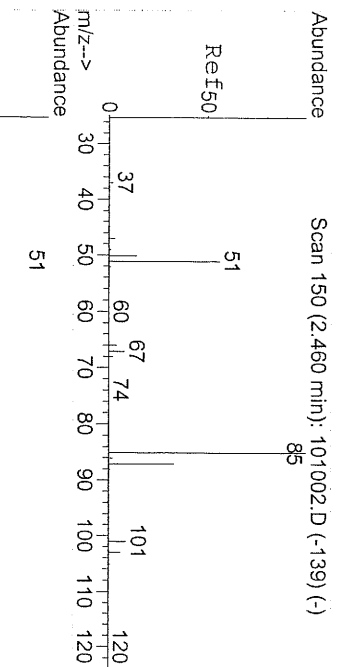
Quant Time: Nov 07 23:08:40 2005
 Quant Method : C:\MSDCHEM\1\METHODS\ALLTO15.M
 Quant Title :
 Qlast Update : Sat Oct 29 13:19:44 2005
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	7.01	130	1543513	10.000	PPB	0.00
37) 1,4-Difluorobenzene	8.69	114	6314405	10.000	PPB	-0.01
51) Chlorobenzene d5	14.57	117	4856798	10.000	PPB	-0.02

Target Compounds						Qvalue
3) Dichlorodifluoromethane	2.42	85	369520	0.818	PPB	96
12) Trichlorofluoromethane	3.44	101	161148	0.364	PPB	100
22) Methylene Chloride	4.76	84	196952	1.520	PPB	76
30) cis-1,2-Dichloroethene	6.65	61	193361	0.765	PPB	79
38) 1,2-Dichloroethane	8.04	62	92689	0.365	PPB	99
39) Benzene	7.99	78	305562	0.592	PPB	94
52) Toluene	11.56	91	3372699	6.468	PPB	99
57) Tetrachloroethene	12.69	166	1114097	4.648	PPB	96
63) Ethylbenzene	14.97	91	828121	1.134	PPB	90
64) m/p-Xylene	15.26	91	1722234	2.925	PPB	91
65) o-Xylene	16.22	91	553407	0.969	PPB	90
66) Styrene	16.27	104	257040	0.622	PPB	100
72) 4-Ethyltoluene	18.53	105	426879	0.554	PPB	98
73) 1,3,5-Trimethylbenzene	18.77	105	218010	0.329	PPB	87
75) 1,2,4-Trimethylbenzene	19.74	105	417122	0.638	PPB	82

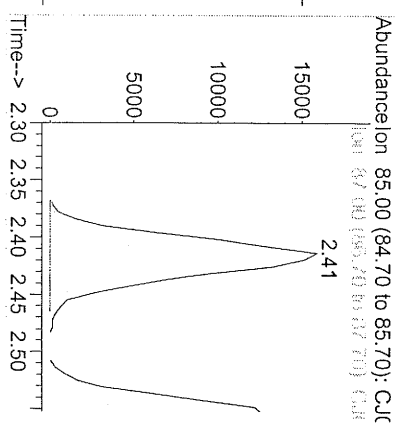
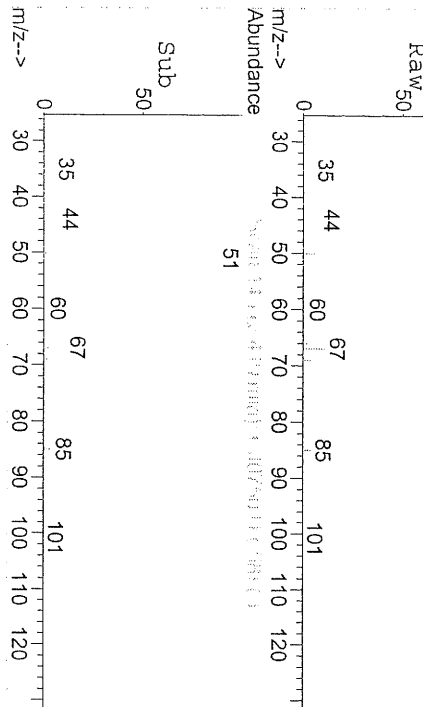
(#) = qualifier out of range (m) = manual integration (+) = signals summed

9999



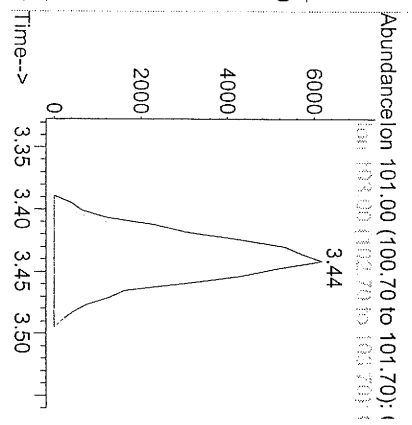
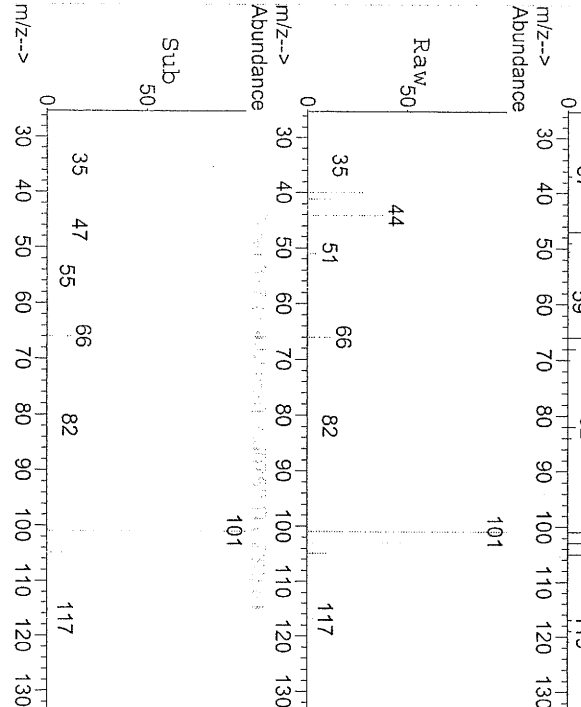
#3
 Dichlorodifluoromethane
 Concen: 0.82 PPB
 RT: 2.42 min Scan# 143
 Delta R.T. 0.00 min
 Lab File: CJ0750.D
 Acq: 29 Oct 2005 18:45

Tgt Ion:	85	Resp:	369520
Ion Ratio	100	Lower	Upper
	87	34.4	25.7 38.5

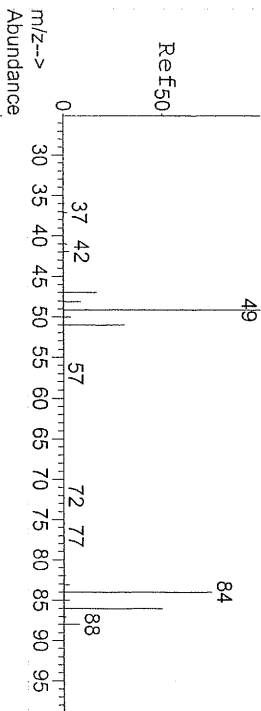


#12
 Trichlorofluoromethane
 Concen: 0.36 PPB
 RT: 3.44 min Scan# 317
 Delta R.T. -0.00 min
 Lab File: CJ0750.D
 Acq: 29 Oct 2005 18:45

Tgt Ion:	101	Resp:	161148
Ion Ratio	100	Lower	Upper
	103	65.2	52.5 78.7



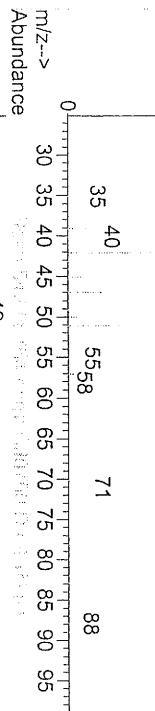
Abundance Scan 539 (4.746 min): 101002.D (-525) (-)



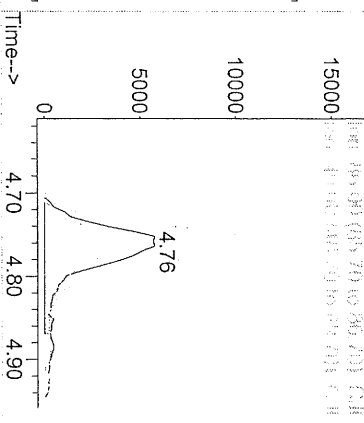
#22
Methylene Chloride
Concn: 1.52 PPB
RT: 4.76 min Scan# 542
Delta R.T. -0.01 min
Lab File: CJ0750.D
Acq: 29 Oct 2005 18:45

Tgt Ion: 84 Resp: 196952
Ion Ratio Lower Upper
84 100
86 70.3 50.7 76.1
49 192.7 123.5 185.3#

Raw 50

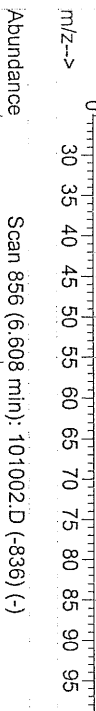


Abundance 84.00 (83.70 to 84.70): CJ

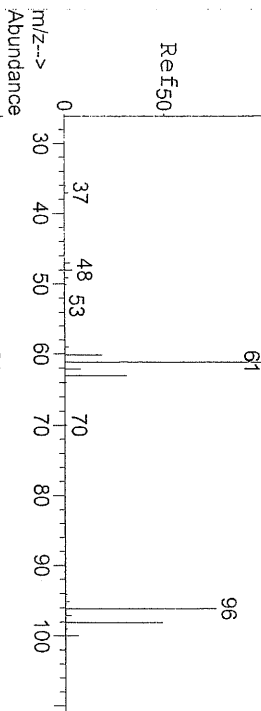


4.76

Sub 50



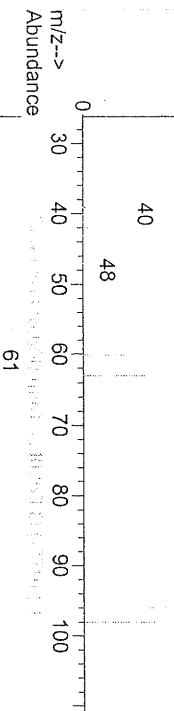
Scan 856 (6.608 min): 101002.D (-836) (-)



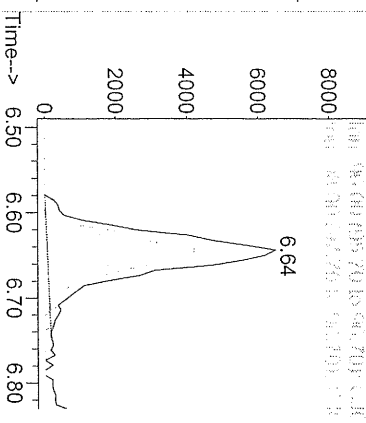
#30
cis-1,2-Dichloroethene
Concn: 0.77 PPB
RT: 6.65 min Scan# 863
Delta R.T. -0.01 min
Lab File: CJ0750.D
Acq: 29 Oct 2005 18:45

Tgt Ion: 61 Resp: 193361
Ion Ratio Lower Upper
61 100
96 63.2 66.6 99.8#
98 40.4 43.7 65.5#

Raw 50

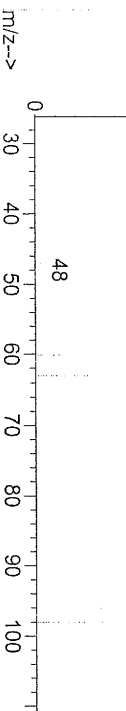


Abundance 61.00 (60.70 to 61.70): CJ

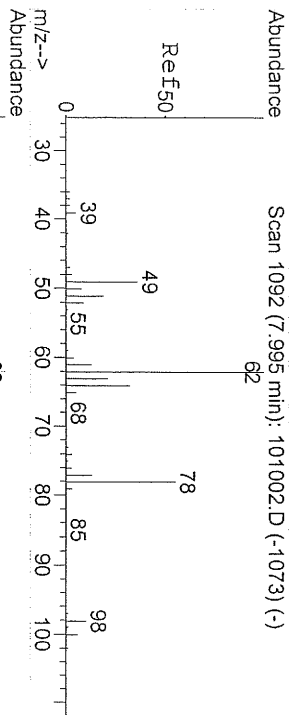


6.64

Sub 50

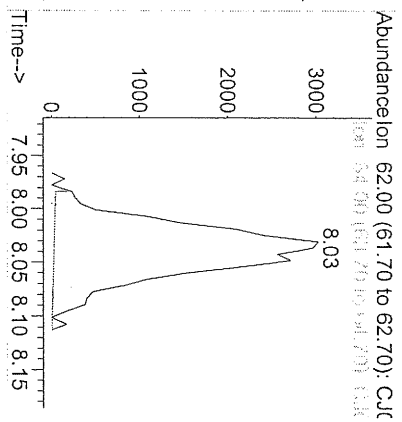
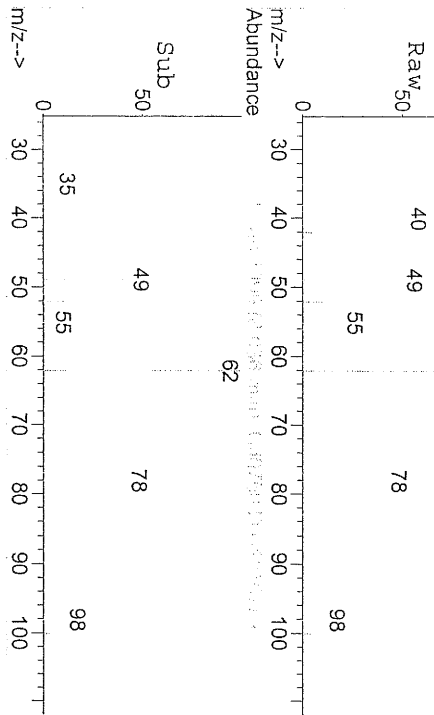


0092



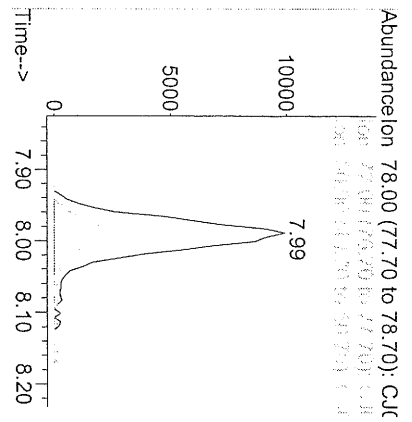
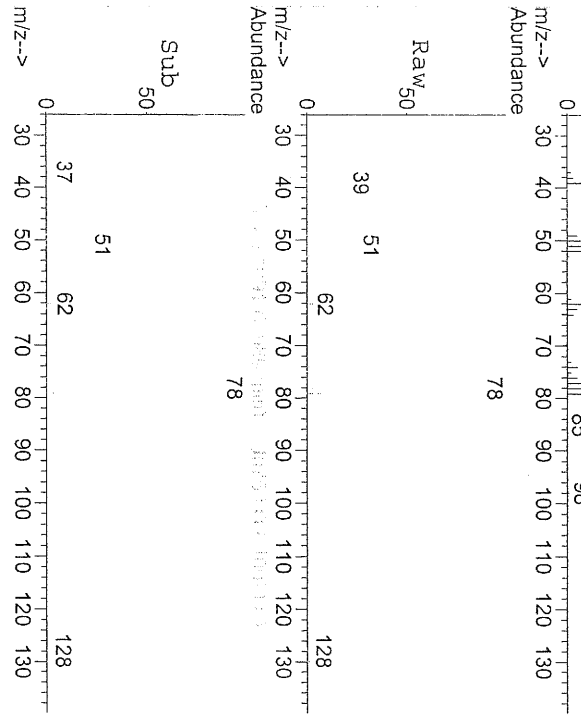
#38
 1,2-Dichloroethane
 Concen: 0.36 PPB
 RT: 8.04 min Scan# 1099
 Delta R.T. -0.02 min
 Lab File: CJ0750.D
 Acq: 29 Oct 2005 18:45

Tgt Ion:	Resp:	92689
Ion Ratio Lower	Upper	
62	100	
64	31.8	25.0
		37.4



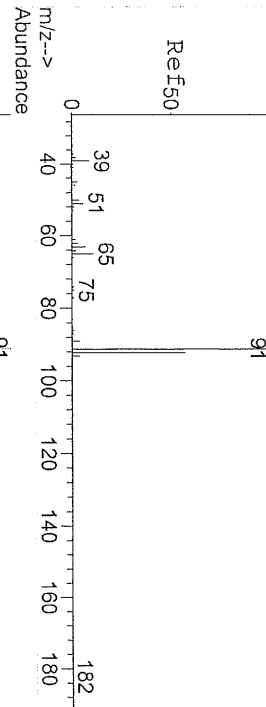
#39
 Benzene
 Concen: 0.59 PPB
 RT: 7.99 min Scan# 1091
 Delta R.T. -0.01 min
 Lab File: CJ0750.D
 Acq: 29 Oct 2005 18:45

Tgt Ion:	Resp:	305562
Ion Ratio Lower	Upper	
78	100	
77	23.0	19.4
50	23.8	15.7
		23.5#



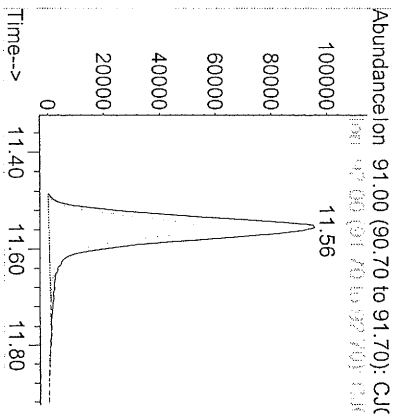
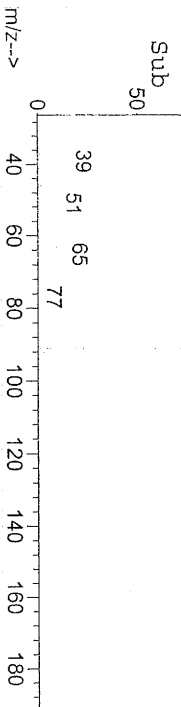
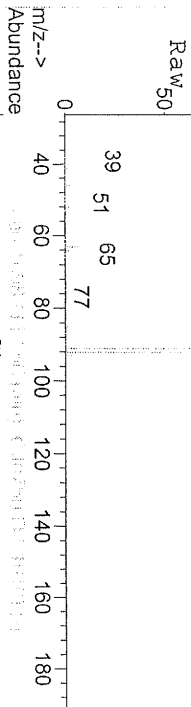
0093

Scan 1689 (11.503 min): 101002.D (-1671) (-)

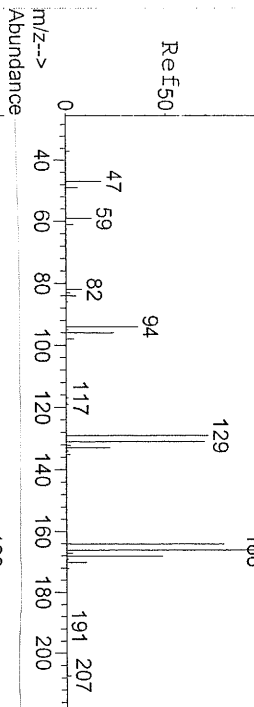


#52
Toluene
Concen: 6.47 PPB
RT: 11.56 min Scan# 1698
Delta R.T. -0.02 min
Lab File: CJ0750.D
Acq: 29 Oct 2005 18:45

Tgt Ion: 91 Resp: 3372699
Ion Ratio Lower Upper
91 100
92 57.4 45.4 68.0

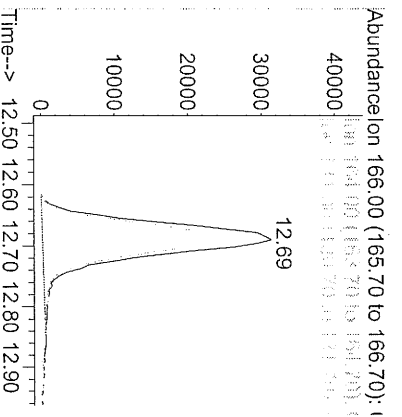
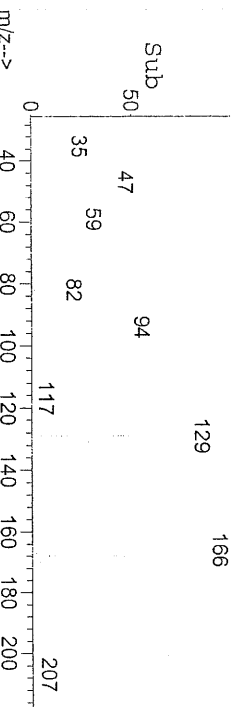
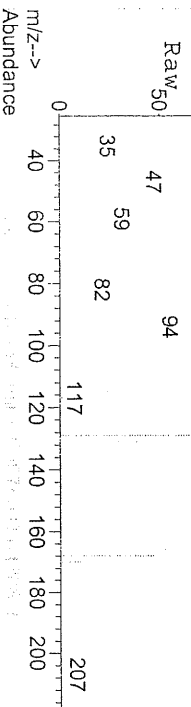


Scan 1883 (12.643 min): 101002.D (-1866) (-)

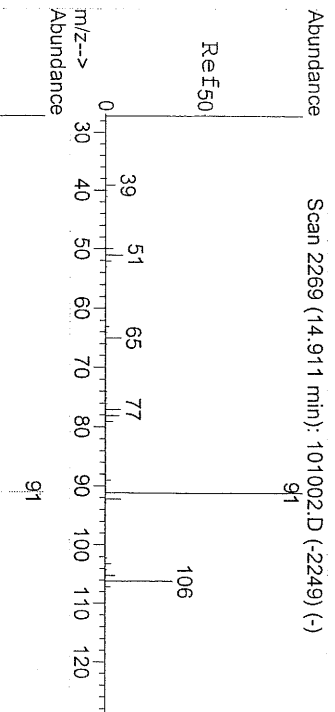


#57
Tetrachloroethene
Concen: 4.65 PPB
RT: 12.69 min Scan# 1891
Delta R.T. -0.01 min
Lab File: CJ0750.D
Acq: 29 Oct 2005 18:45

Tgt Ion: 166 Resp: 1114097
Ion Ratio Lower Upper
166 100
164 79.9 63.0 94.6
131 82.0 61.1 91.7

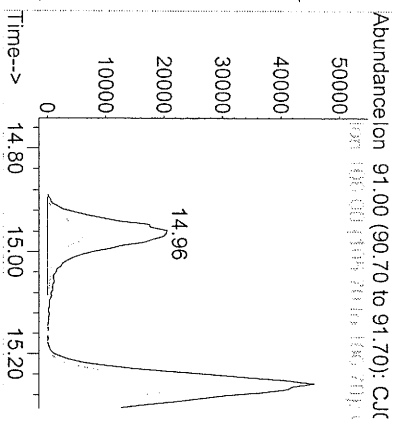
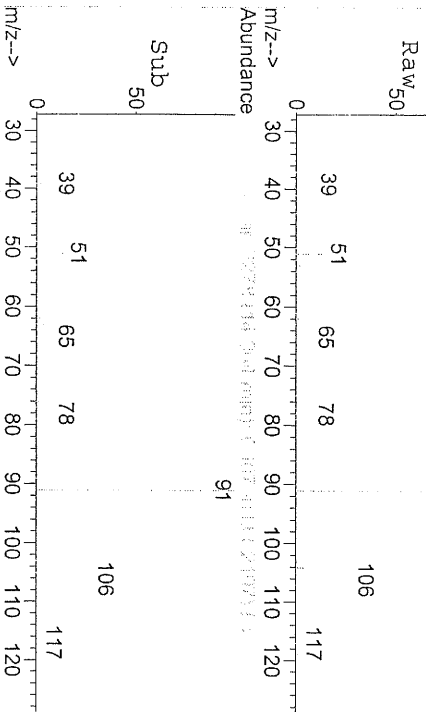


0094



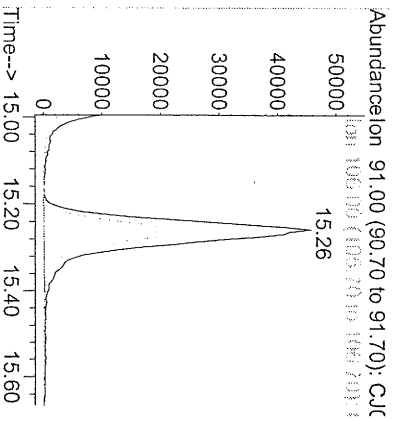
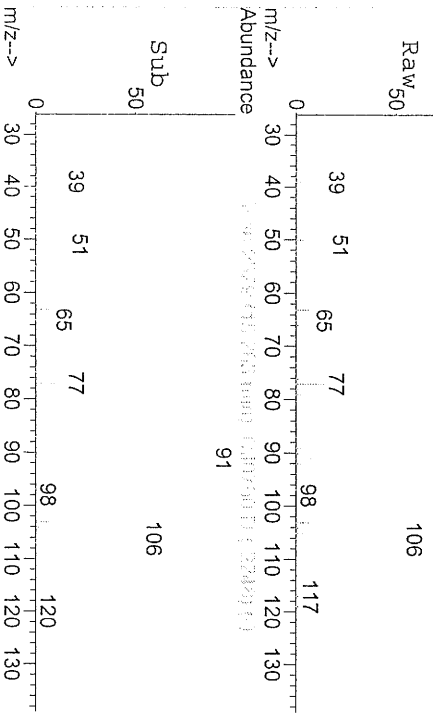
#63
Ethylbenzene
Concen: 1.13 PPB
RT: 14.97 min Scan# 2278
Delta R.T. -0.02 min
Lab File: CJ0750.D
Acq: 29 Oct 2005 18:45

Tgt Ion: 91 Resp: 828121
Ion Ratio Lower Upper
91 100
106 30.1 28.8 43.2

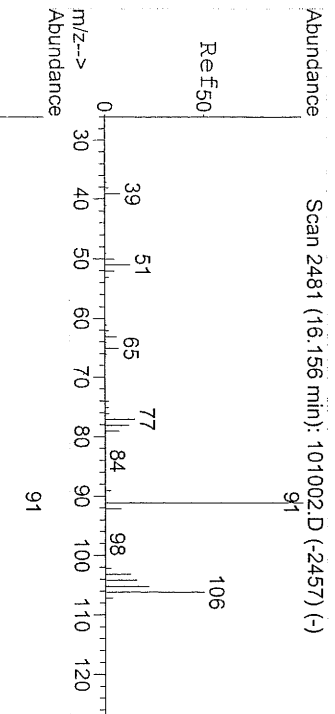


#64
m/p-Xylene
Concen: 2.92 PPB
RT: 15.26 min Scan# 2329
Delta R.T. -0.02 min
Lab File: CJ0750.D
Acq: 29 Oct 2005 18:45

Tgt Ion: 91 Resp: 1722234
Ion Ratio Lower Upper
91 100
106 49.7 45.0 67.4

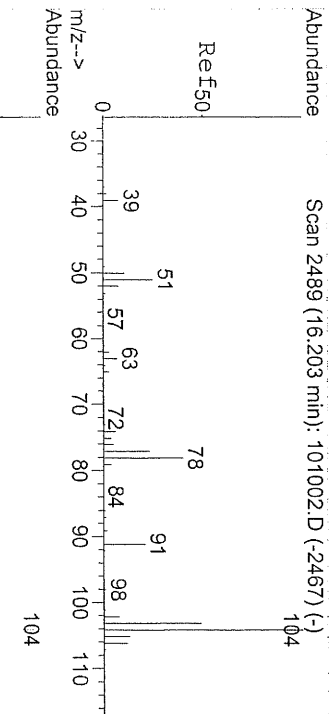
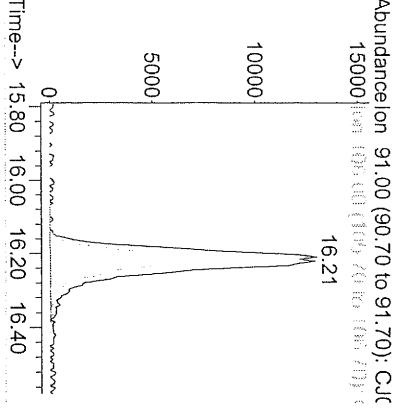
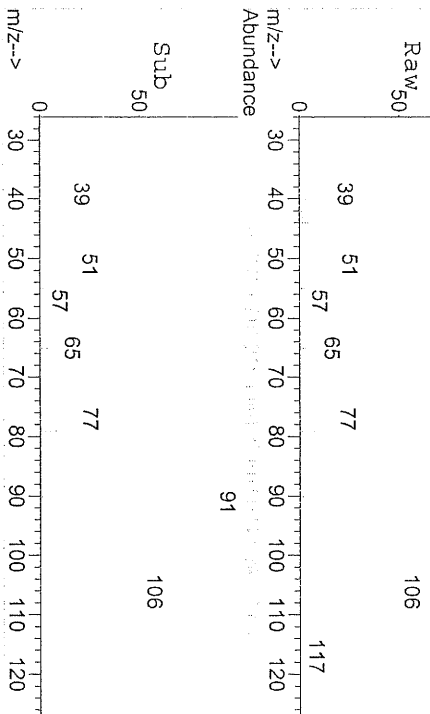


0095



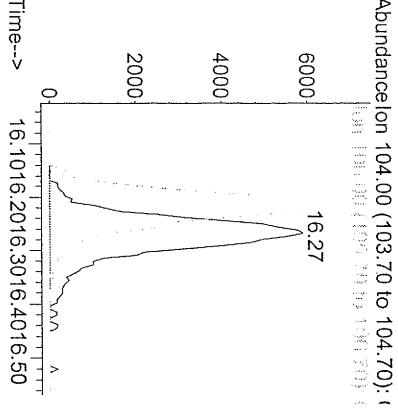
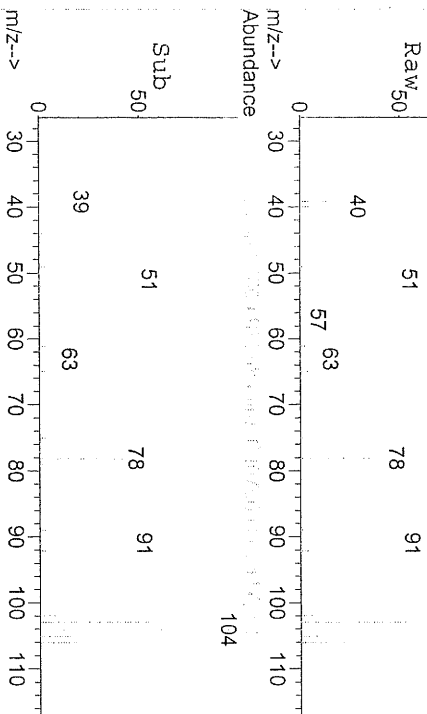
#65
o-Xylene 0.97 PPB
Concn: 16.22 min Scan# 2491
RT: 16.22 min Scan# 2491
Delta R.T. -0.02 min
Lab File: CJ0750.D
Acq: 29 Oct 2005 18:45

Tgt Ion: 91 Resp: 553407
Ion Ratio Lower Upper
91 100
106 47.2 43.4 65.0

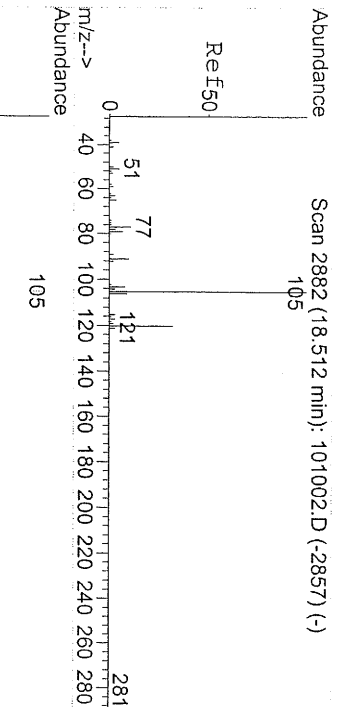


#66
Styrene 0.62 PPB
Concn: 16.27 min Scan# 2500
RT: 16.27 min Scan# 2500
Delta R.T. -0.01 min
Lab File: CJ0750.D
Acq: 29 Oct 2005 18:45

Tgt Ion: 104 Resp: 257040
Ion Ratio Lower Upper
104 100
106 0.0 0.0 0.0

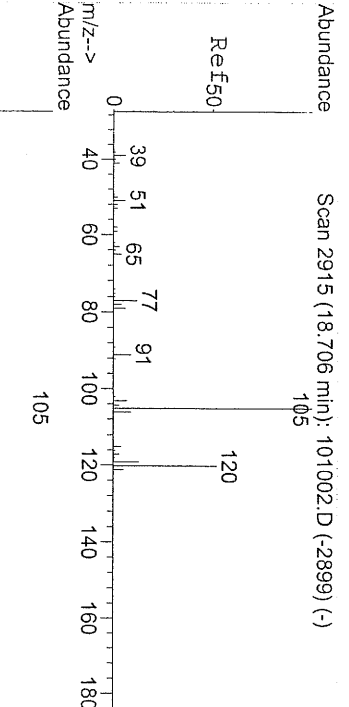
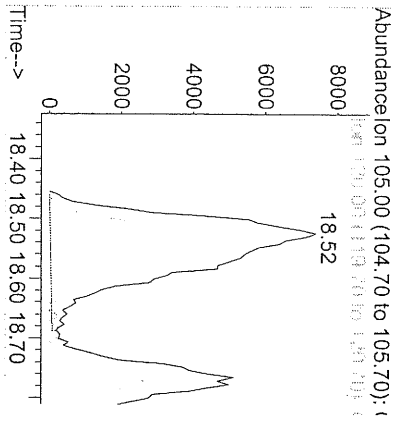
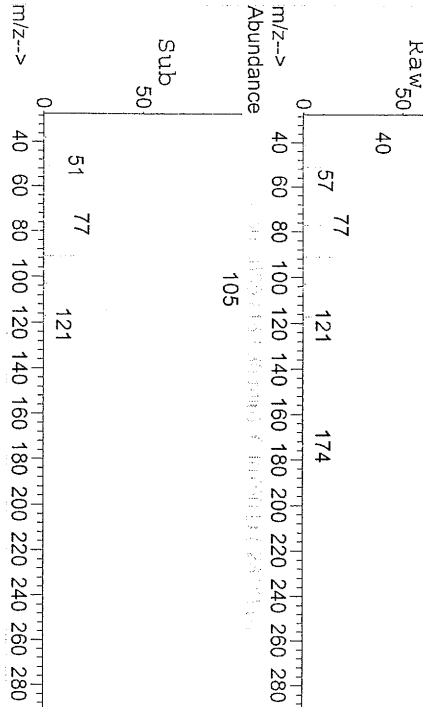


0096



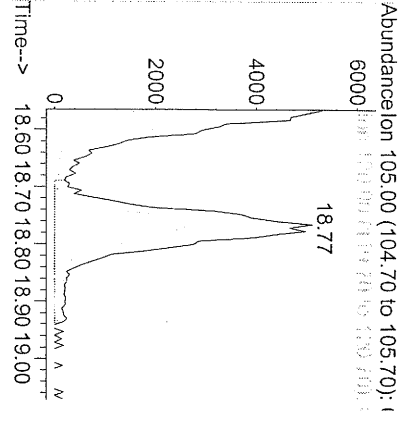
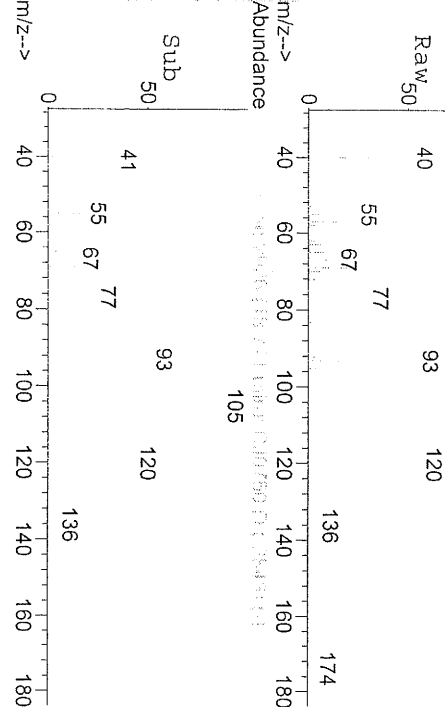
#72
 4-Ethyltoluene
 Concn: 0.55 PPB
 RT: 18.53 min Scan# 2885
 Delta R.T. -0.07 min
 Lab File: CJ0750.D
 Acq: 29 Oct 2005 18:45

Tgt Ion	Ratio	Resp
105	100	426879
120	32.5	27.0
		40.6

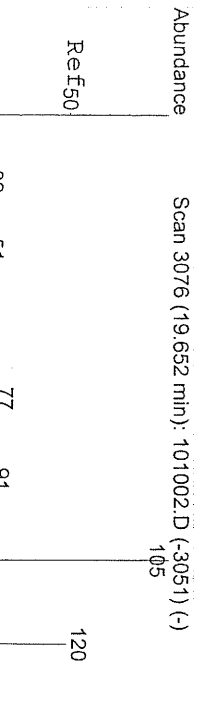


#73
 1,3,5-Trimethylbenzene
 Concn: 0.33 PPB
 RT: 18.77 min Scan# 2926
 Delta R.T. -0.01 min
 Lab File: CJ0750.D
 Acq: 29 Oct 2005 18:45

Tgt Ion	Ratio	Resp
105	100	218010
120	45.8	43.9
		65.9

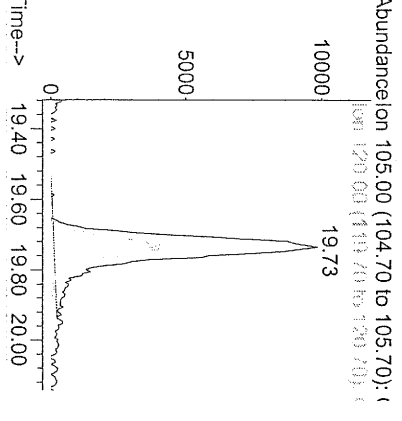
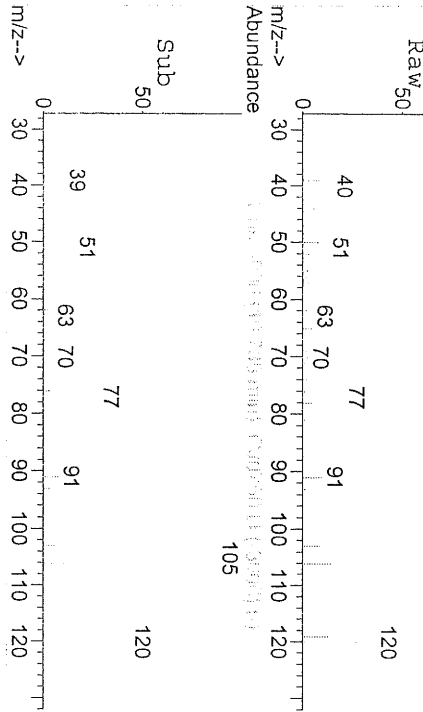


0097



#75
 1,2,4-Trimethylbenzene
 Concen: 0.64 PPB
 RT: 19.74 min Scan# 3090
 Delta R.T. -0.01 min
 Lab File: CJ0750.D
 Acq: 29 Oct 2005 18:45

Tgt Ion	Ratio	Resp
105	100	417122
120	40.2	63.5#



0000

LANCASTER LABORATORIES
VOLATILE ORGANICS IN AIR
SUMMA CANISTER SAMPLE
ANALYSIS DATA SHEET

Sample No.: 1FACC Date Collected: 10/19/05 Date Received: 10/20/05
 Lab Sample ID: 4628494 Date Analyzed: 10/29/05 Time Analyzed: 00:42
 Canister ID: SUMMA0232 Pressure Rec'd: 14.7 psia Final Pressure: 29.4 psia
 Injection Volume: 500 cc Nominal Volume: 250 cc Dilution Factor: 1.0
 Instrument ID: HP09464 ⁸⁷³² Lab File ID: C:\MSDCHEM\1\DATA\OCT28\CJ0722.D

CAS RN	COMPOUND NAME	CONCENTRATION	UNITS	MDL	ug/m3	Q
75-71-8	Dichlorodifluoromethane		6			U
76-14-2	Freon 114		1			U
74-87-3	Chloromethane		2			J
75-01-4	Vinyl Chloride		0.8			J
74-83-9	Bromomethane		0.8			J
75-00-3	Chloroethane		0.5			J
75-69-4	Trichlorofluoromethane		2			J
75-35-4	1,1-Dichloroethene		0.8			J
76-13-1	Freon 113		4			J
107-05-1	3-Chloropropene		2			J
75-09-2	Methylene Chloride		3			J
75-34-3	1,1-Dichloroethane		0.8			J
156-59-2	cis-1,2-Dichloroethene		6			J
67-66-3	Chloroform		1			J
71-55-6	1,1,1-Trichloroethane		1			J
56-23-5	Carbon Tetrachloride		1			J
107-06-2	1,2-Dichloroethane		4			J
71-43-2	Benzene		4			J
79-01-6	Trichloroethene		1			J
78-87-5	1,2-Dichloropropane		0.9			J
10061-01-5	cis-1,3-Dichloropropene		0.9			J
108-88-3	Toluene		45			J
10061-02-6	trans-1,3-Dichloropropene		0.9			J
79-00-5	1,1,2-Trichloroethane		1			J
127-18-4	Tetrachloroethene		74			J
106-93-4	1,2-Dibromoethane		2			J
108-90-7	Chlorobenzene		0.9			J
100-41-4	Ethylbenzene		10			J
1330-20-7	m/p-Xylene		23			J
95-47-6	o-Xylene		8			J
100-42-5	Styrene		6			J
79-34-5	1,1,2,2-Tetrachloroethane		1			J
622-96-8	4-Ethyltoluene		4			J
108-67-8	1,3,5-Trimethylbenzene		4			J

U = Compound was undetected at the specified limit of quantitation.
 B = Compound was found in method blank. D = analysis of diluted sample.
 J = Compound was detected, but below the limit of quantitation.

LANCASTER LABORATORIES
 VOLATILE ORGANICS IN AIR
 SUMMA CANISTER SAMPLE
 ANALYSIS DATA SHEET

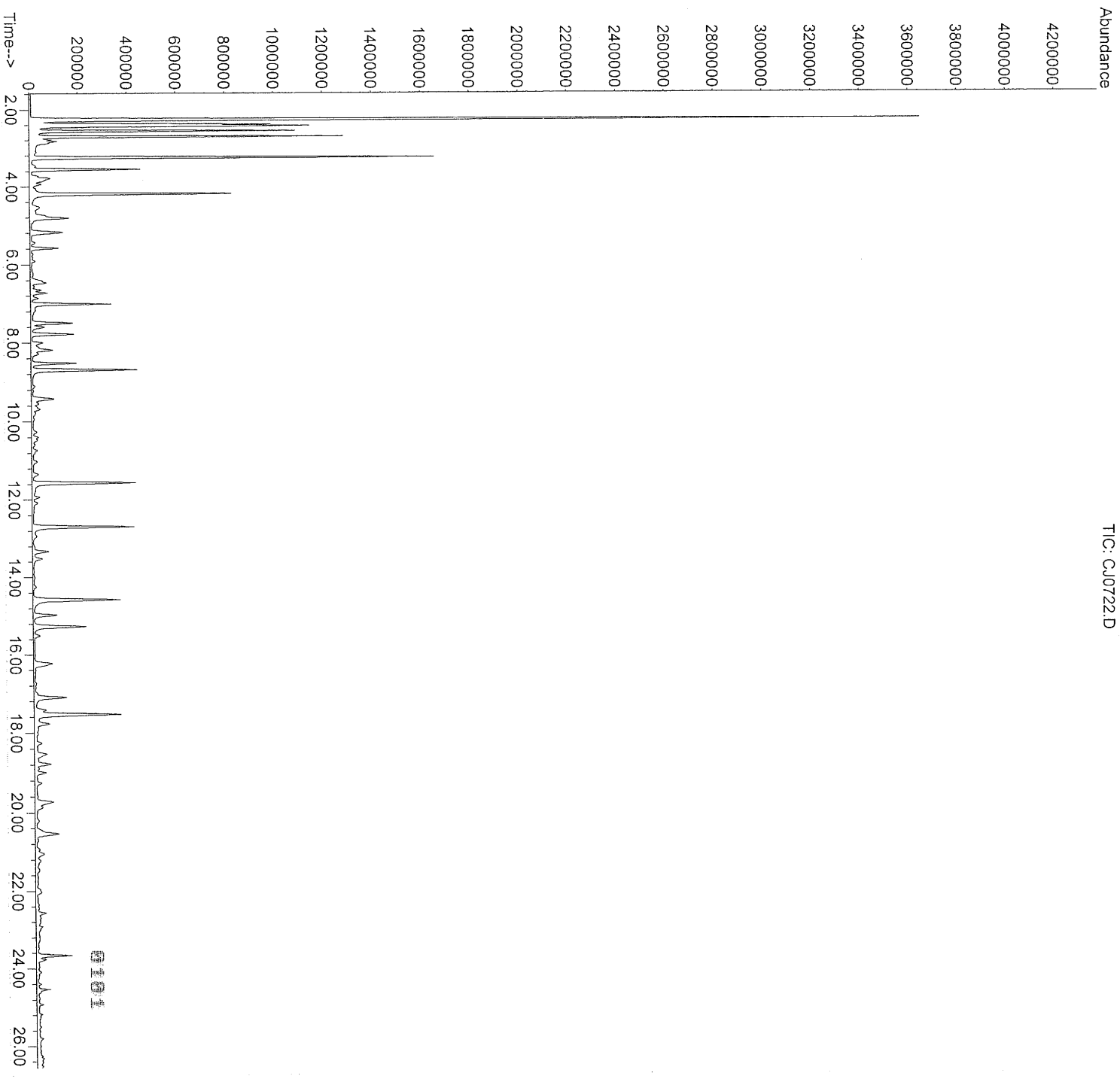
Sample No.: 1FACC
 Lab Sample ID: 4628494 Date Collected: 10/19/05 Date Received: 10/20/05
 Canister ID: SUMMA0232 Date Analyzed: 10/29/05 Time Analyzed: 00:42
 Injection Volume: 500 cc ~~796~~ Nominal Volume: 250 cc Final Pressure: 29.4 psia
 Instrument ID: HP09464 ~~11/14/05~~ Lab File ID: C:\MSDCHEM\1\DATA\OCT28\CJ0722.D Dilution Factor: 1.0

CAS RN	COMPOUND NAME	CONCENTRATION	UNITS: MDL	ug/m3	Q
95-63-6	1,2,4-Trimethylbenzene			7	U
541-73-1	1,3-Dichlorobenzene			3	U
106-46-7	1,4-Dichlorobenzene			3	U
95-50-1	1,2-Dichlorobenzene			3	U
120-82-1	1,2,4-Trichlorobenzene			7	U

U = Compound was undetected at the specified limit of quantitation.
 B = Compound was found in method blank. D = analysis of diluted sample.
 J = Compound was detected, but below the limit of quantitation.

Data Path : C:\MSDCHEM\1\DATA\OCT28\
Data File : CJ0722.D
Acq On : 29 Oct 2005 00:42
Operator : JBS
Sample : 4628494 500CC
Misc :
ALS Vial : 47 Sample Multiplier: 1

Quant Time: Nov 07 23:04:08 2005
Quant Method : C:\MSDCHEM\1\METHODS\ALLT015.M
Quant Title :
Quant Update : Fri Oct 28 08:49:53 2005
Response via : Initial Calibration



Data Path : C:\MSDCHEM\1\DATA\OCT28\
 Data File : CJ0722.D
 Acq On : 29 Oct 2005 00:42
 Operator : JBS
 Sample : 4628494 500CC
 Misc :
 ALS Vial : 47 Sample Multiplier: 1

Quant Time: Nov 07 23:04:08 2005
 Quant Method : C:\MSDCHEM\1\METHODS\ALLT015.M
 Quant Title :
 Qlast Update : Fri Oct 28 08:49:53 2005
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	7.01	130	1528717	10.000	PPB	0.00
37) 1,4-Difluorobenzene	8.70	114	5244632	10.000	PPB	0.00
51) Chlorobenzene d5	14.58	117	4301798	10.000	PPB	0.00

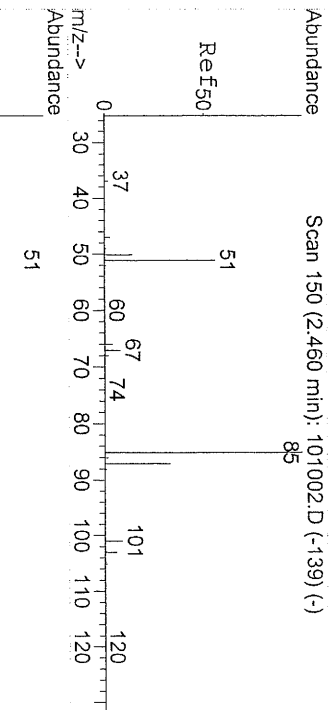
Target Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)	Qvalue
3) Dichlorodifluoromethane	2.42	85	527104	1.178	PPB	99	
6) Chloromethane	2.56	50	302777m	1.053	PPB		
7) Vinyl Chloride	2.70	62	64954	0.322	PPB	#	45
12) Trichlorofluoromethane	3.45	101	133440	0.304	PPB		98
22) Methylene Chloride	4.77	84	94661	0.738	PPB		80
26) Methyl t-Butyl Ether	5.20	73	1064305	2.486	PPB	#	48
30) cis-1,2-Dichloroethene	6.66	61	398124	1.591	PPB	#	78
38) 1,2-Dichloroethane	8.04	62	196238	0.930	PPB		100
39) Benzene	8.00	78	524374	1.224	PPB		95
42) Trichloroethene	9.11	130	49589	0.211	PPB	#	58
52) Toluene	11.56	91	5543988	12.003	PPB		100
57) Tetrachloroethene	12.69	166	2311722	10.889	PPB		97
63) Ethylbenzene	14.97	91	1468730	2.272	PPB		92
64) m/p-Xylene	15.27	91	2721982	5.219	PPB		91
65) o-Xylene	16.22	91	933865	1.847	PPB		90
66) Styrene	16.27	104	477423	1.303	PPB	#	100
72) 4-Ethyltoluene	18.53	105	551566m	0.808	PPB		
73) 1,3,5-Trimethylbenzene	18.78	105	420885	0.718	PPB		91
75) 1,2,4-Trimethylbenzene	19.74	105	880088	1.519	PPB	#	76
81) 1,2,4-Trichlorobenzene	24.92	180	86830m	0.621	PPB		

ADITYA
11/14/05
CS

(#) = qualifier out of range (m) = manual integration (+) = signals summed

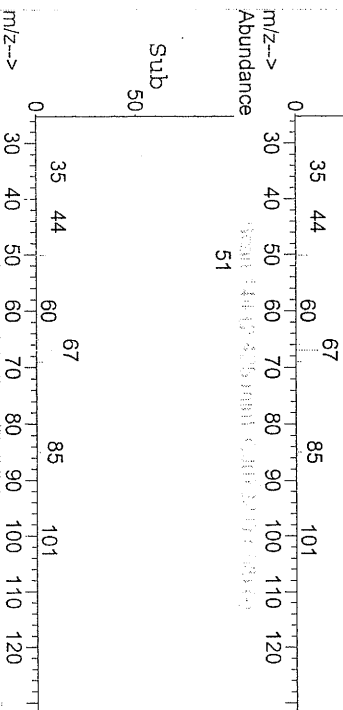
0102

#3
 Dichlorodifluoromethane
 Concn: 1.18 PPB
 RT: 2.42 min Scan# 144
 Delta R.T. 0.00 min
 Lab File: CJ0722.D
 Acq: 29 Oct 2005 00:42



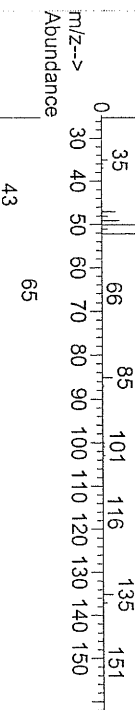
Tgt Ion: 85 Resp: 527104
 Ion Ratio Lower Upper
 85 100
 87 32.8 25.7 38.5

Abundance Ion 85.00 (84.70 to 85.70): CJ0722.D (82.00 (85.70 to 87.70) 4.00



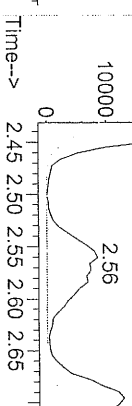
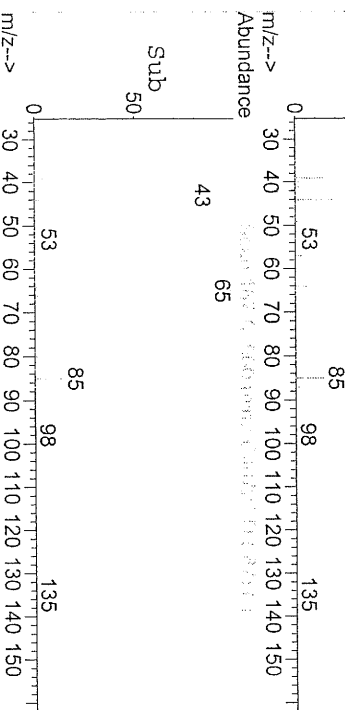
Time--> 2.30 2.35 2.40 2.45 2.50

#6
 Chloromethane
 Concn: 1.05 PPB m
 RT: 2.56 min Scan# 167
 Delta R.T. -0.03 min
 Lab File: CJ0722.D
 Acq: 29 Oct 2005 00:42



Tgt Ion: 50 Resp: 302777
 Ion Ratio Lower Upper
 50 100
 52 0.0 18.6 28.0#

Abundance Ion 50.00 (49.70 to 50.70): CJ0722.D (52.00 (51.70 to 53.70) 4.00

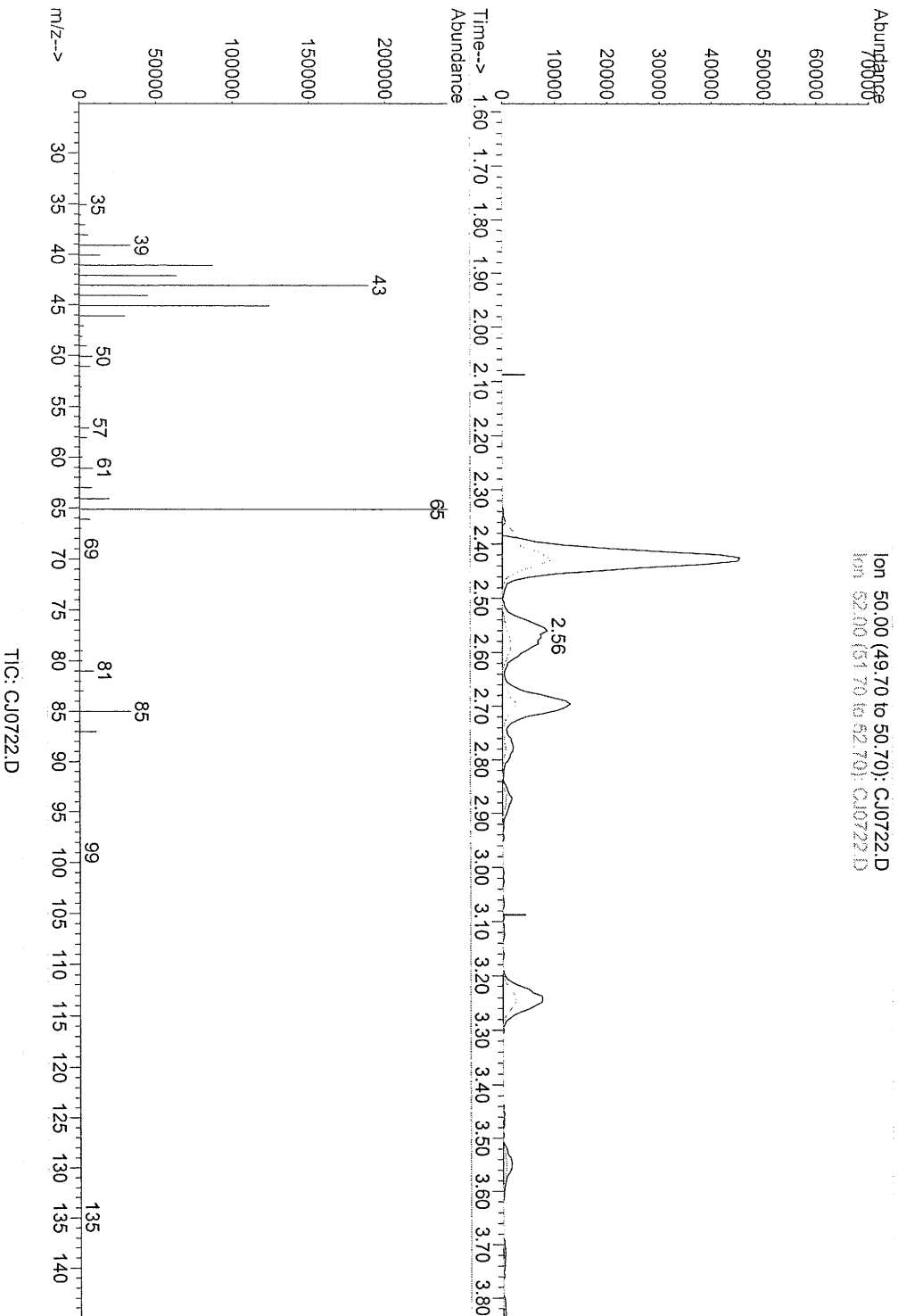


0103

Quantitation Report (Qedit)

Data Path : C:\MSDCHEM\1\DATA\OCT28\
 Data File : CJ0722.D
 Acq On : 29 Oct 2005 00:42
 Operator : JBS
 Sample : 4628494 500CC
 Misc :
 ALS Vial : 47 Sample Multiplier: 1
 Quant Time: Oct 29 00:09:34 2005
 Quant Method : C:\MSDCHEM\1\METHODS\ALLT015.M
 Quant Title :
 Qast Update : Fri Oct 28 08:49:53 2005
 Response via : Initial Calibration

Ion 50.00 (49.70 to 50.70): CJ0722.D
 Ion 52.00 (51.70 to 52.70): CJ0722.D



(6) Chloromethane

2.560min (-0.028) 1.05PPB m

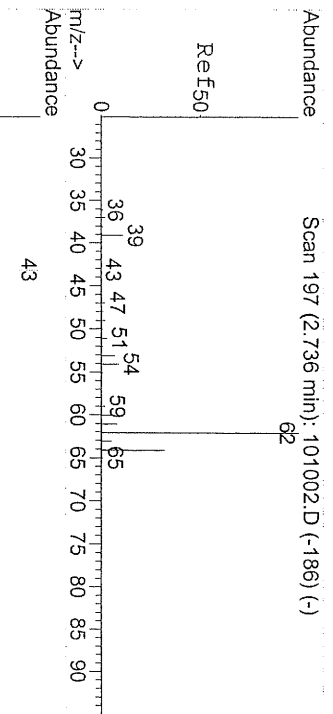
response 302777

Ion	Exp%	Act%
50.00	100	100
52.00	23.30	0.00#
0.00	0.00	0.00
0.00	0.00	0.00

*split peak
 present
 10/31/05*

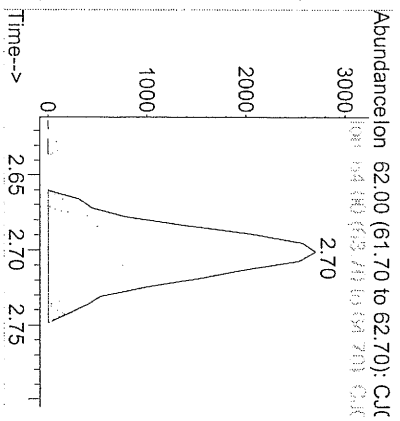
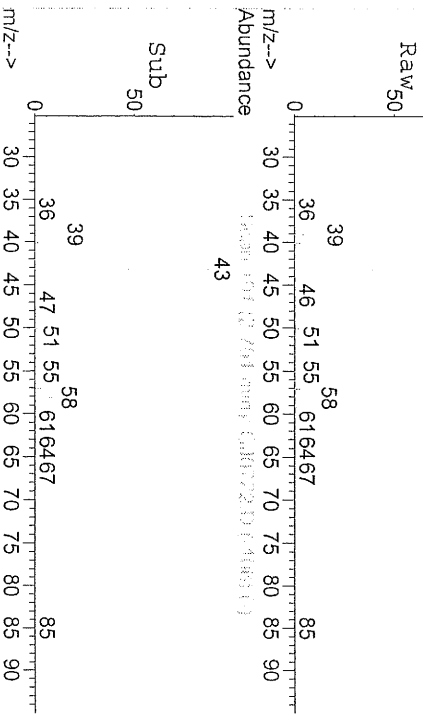
*OME/412
 11/12/05*

8184



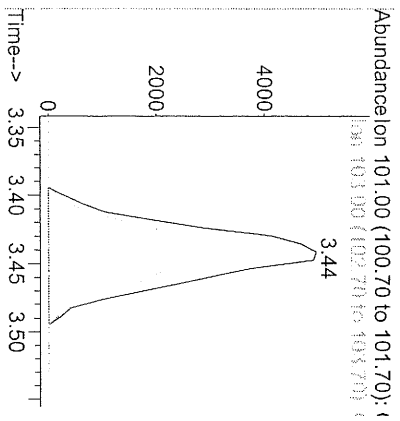
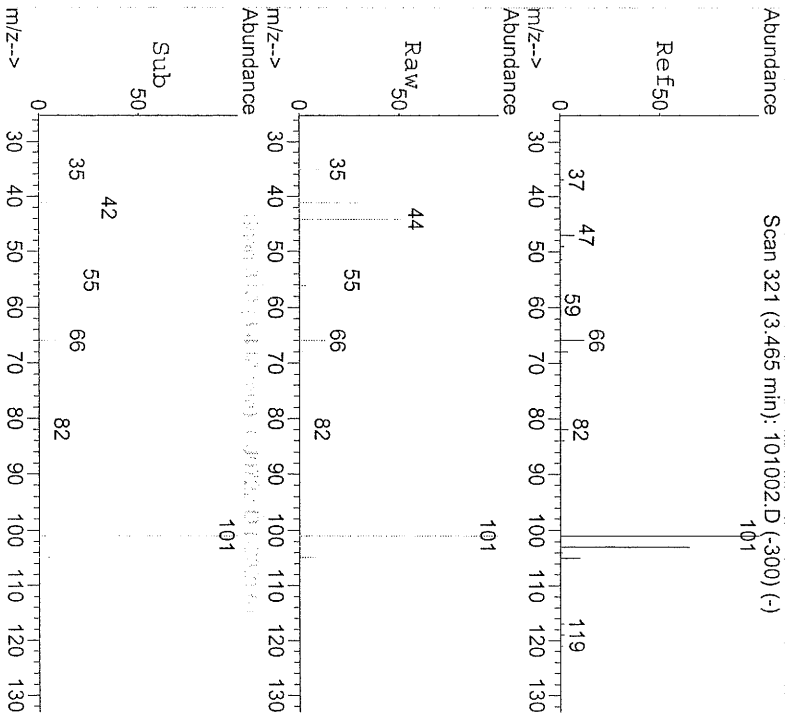
#7
 Vinyl Chloride
 Concen: 0.32 PPB
 RT: 2.70 min Scan# 191
 Delta R.T. 0.00 min
 Lab File: CJ0722.D
 Acq: 29 Oct 2005 00:42

Tgt Ion: 62 Resp: 64954
 Ion Ratio Lower Upper
 62 100
 64 0.0 23.8 35.6#

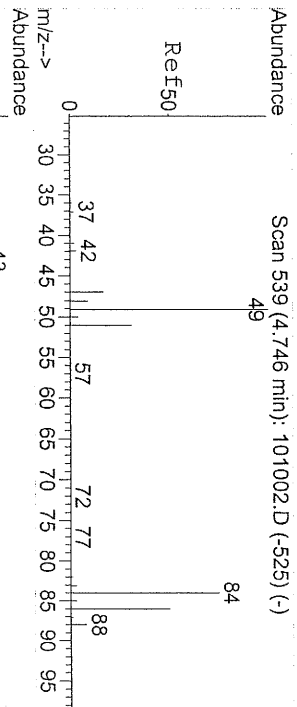


#12
 Trichlorofluoromethane
 Concen: 0.30 PPB
 RT: 3.45 min Scan# 318
 Delta R.T. 0.00 min
 Lab File: CJ0722.D
 Acq: 29 Oct 2005 00:42

Tgt Ion: 101 Resp: 133440
 Ion Ratio Lower Upper
 101 100
 103 63.9 52.5 78.7

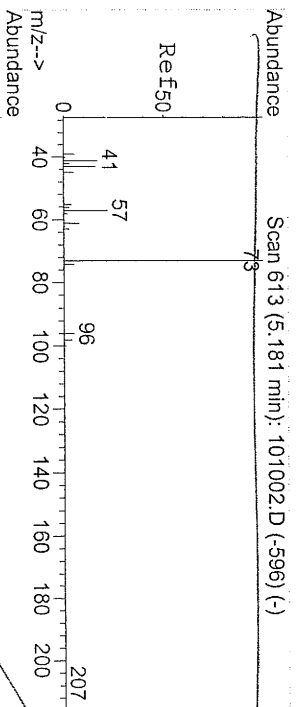
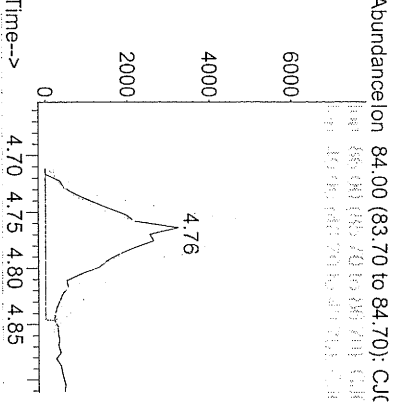
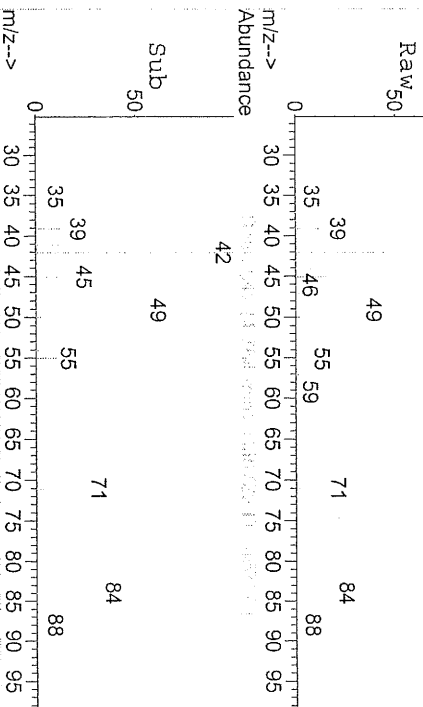


0105



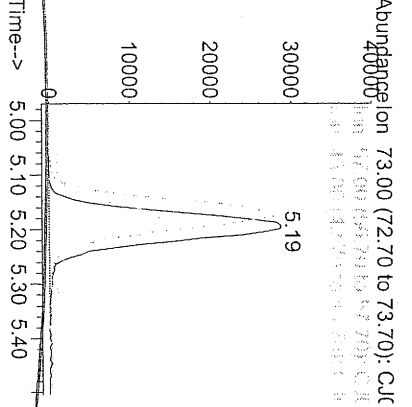
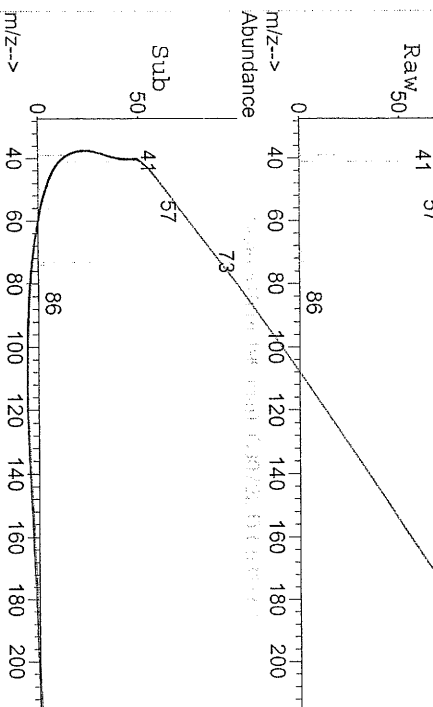
#22
 Methylene Chloride
 Concen: 0.74 PPB
 RT: 4.77 min Scan# 542
 Delta R.T. 0.00 min
 Lab File: CJ0722.D
 Acq: 29 Oct 2005 00:42

Tgt Ion:	84	Resp:	94661
Ion Ratio	100	Lower	Upper
86	75.2	50.7	76.1
49	182.8	123.5	185.3



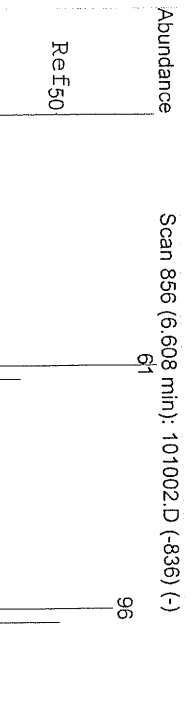
#26
 Methyl t-Butyl Ether
 Concen: 2.49 PPB
 RT: 5.20 min Scan# 616
 Delta R.T. -0.00 min
 Lab File: CJ0722.D
 Acq: 29 Oct 2005 00:42

Tgt Ion:	73	Resp:	1064305
Ion Ratio	100	Lower	Upper
57	0.0	20.6	31.0#
43	0.0	22.2	33.2#



Handwritten signature
 11/11/05

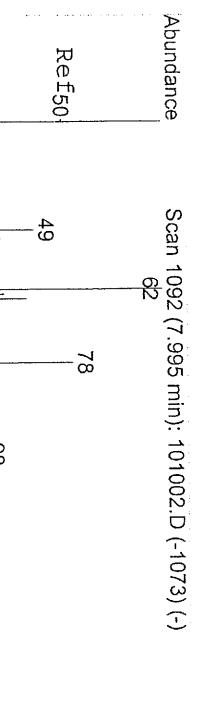
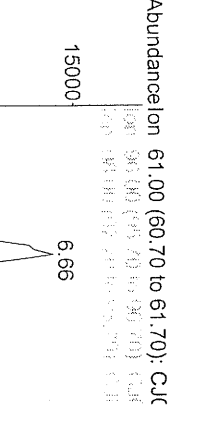
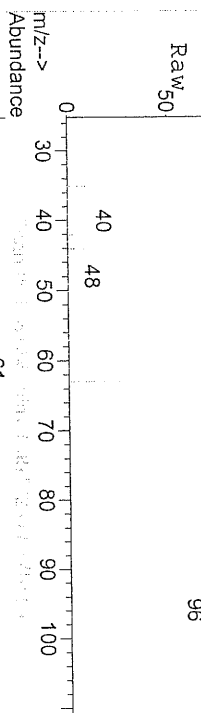
0106



Scan 856 (6.608 min): 101002.D (-836) (-)

#30
 cis-1,2-Dichloroethene
 Concen: 1.59 PPB
 RT: 6.66 min Scan# 864
 Delta R.T. 0.00 min
 Lab File: CJ0722.D
 Acq: 29 Oct 2005 00:42

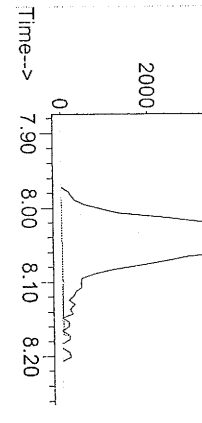
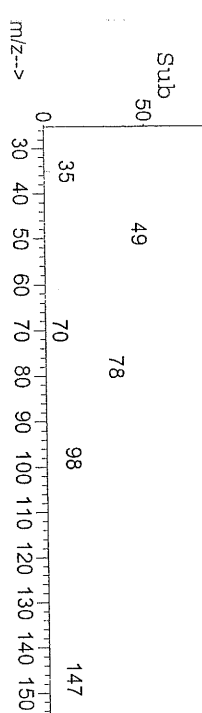
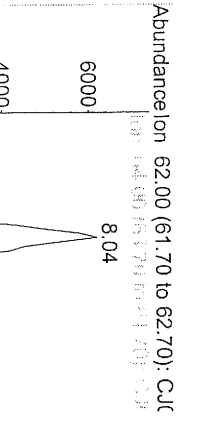
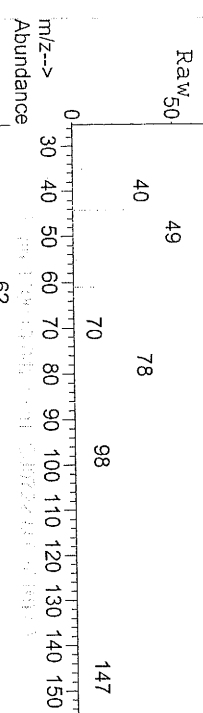
Tgt Ion	Ratio	Lower	Upper
61	100		
96	61.5	66.6	99.8#
98	40.8	43.7	65.5#



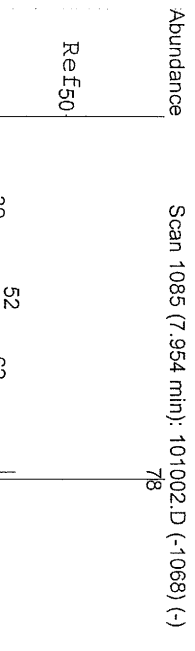
Scan 1092 (7.995 min): 101002.D (-1073) (-)

#38
 1,2-Dichloroethane
 Concen: 0.93 PPB
 RT: 8.04 min Scan# 1100
 Delta R.T. -0.00 min
 Lab File: CJ0722.D
 Acq: 29 Oct 2005 00:42

Tgt Ion	Ratio	Lower	Upper
62	100		
64	31.3	25.0	37.4

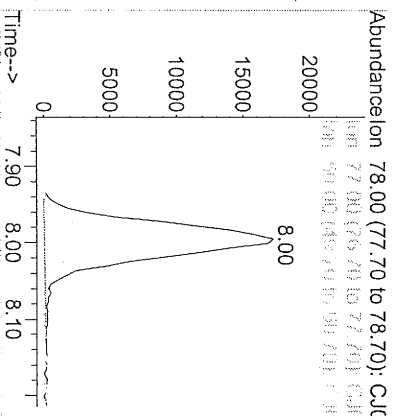
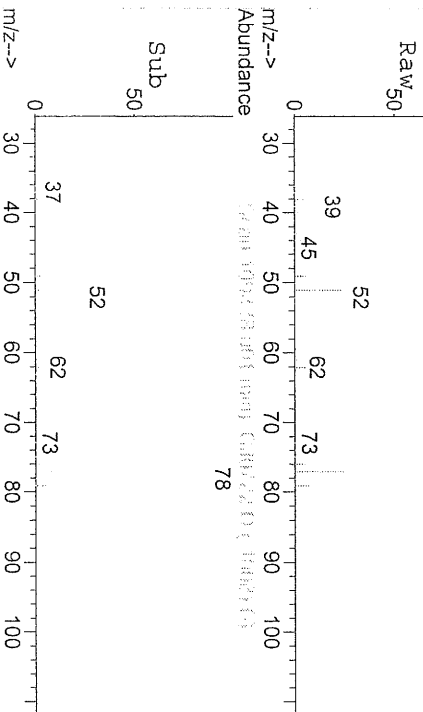


0107



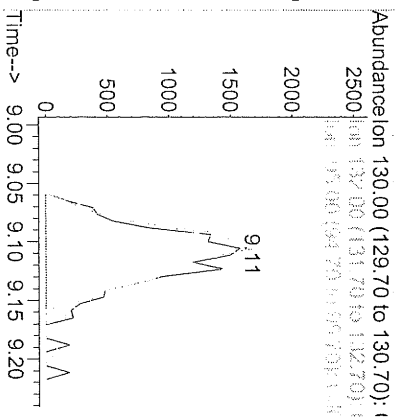
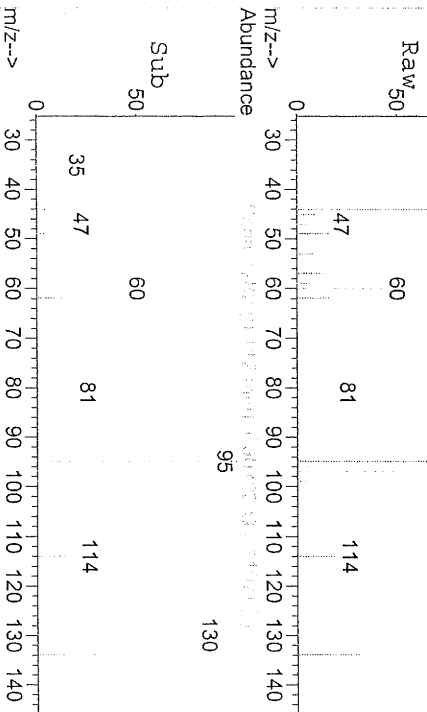
#39
Benzene
Concen: 1.22 PPB
RT: 8.00 min Scan# 1093
Delta R.T. -0.00 min
Lab File: CJ0722.D
Acq: 29 Oct 2005 00:42

Tgt Ion	Resp	524374
78	100	
77	23.0	19.4
50	23.3	15.7
		29.2
		23.5

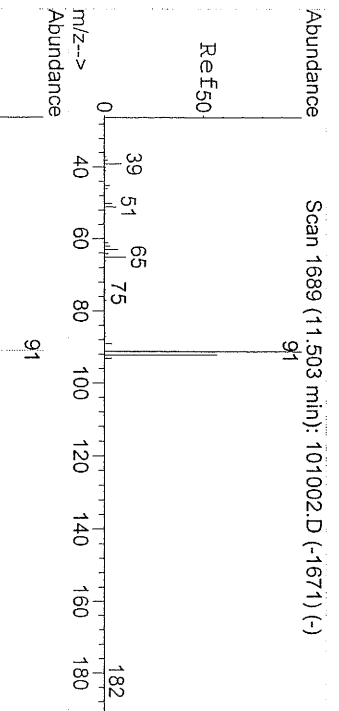


#42
Trichloroethene
Concen: 0.21 PPB
RT: 9.11 min Scan# 1282
Delta R.T. -0.00 min
Lab File: CJ0722.D
Acq: 29 Oct 2005 00:42

Tgt Ion	Resp	49589
130	100	
132	98.9	79.0
95	0.0	66.1
		118.6
		99.1#

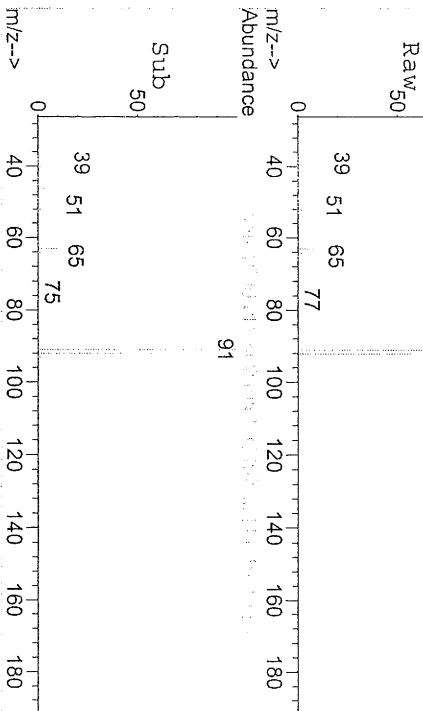


0100



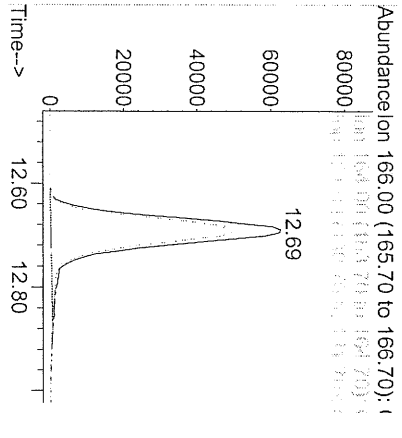
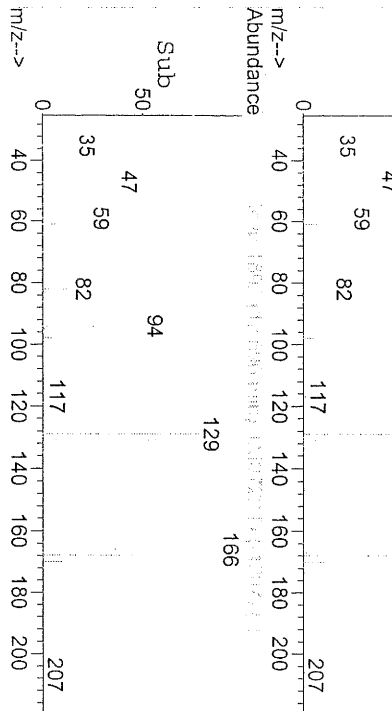
#52
Toluene
Concn: 12.00 PPB
RT: 11.56 min Scan# 1699
Delta R.T. -0.01 min
Lab File: CJ0722.D
Acq: 29 Oct 2005 00:42

Tgt Ion	Ratio	Lower	Upper
91	100		
92	57.0	45.4	68.0



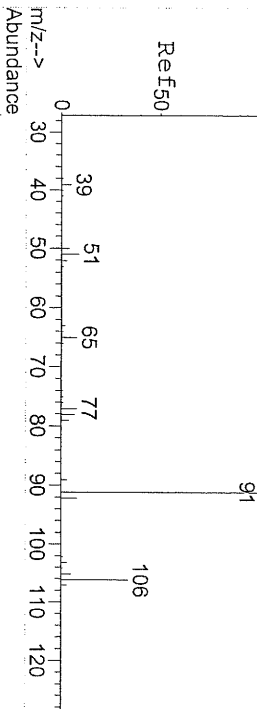
#57
Tetrachloroethene
Concn: 10.89 PPB
RT: 12.69 min Scan# 1892
Delta R.T. -0.00 min
Lab File: CJ0722.D
Acq: 29 Oct 2005 00:42

Tgt Ion	Ratio	Lower	Upper
166	100		
164	81.0	63.0	94.6
131	79.5	61.1	91.7



8189

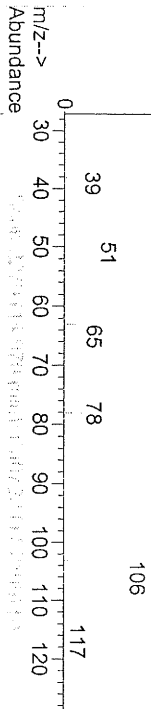
Abundance Scan 2269 (14.911 min): 101002.D (-2249) (-)



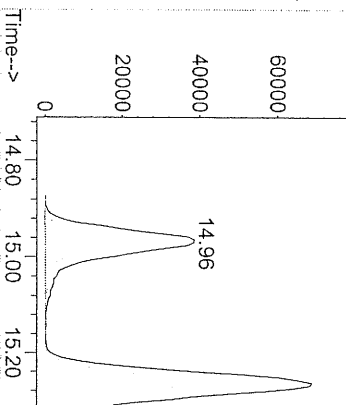
#63
Ethylbenzene
Concen: 2.27 PPB
RT: 14.97 min Scan# 2279
Delta R.T. -0.00 min
Lab File: CJ0722.D
Acq: 29 Oct 2005 00:42

Tgt Ion: 91 Resp: 1468730
Ion Ratio Lower Upper
91 100
106 31.1 28.8 43.2

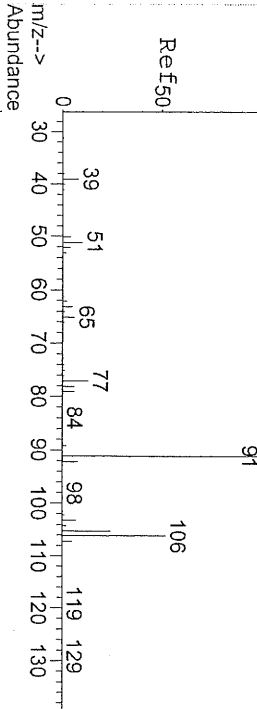
Raw 50



Abundance Ion 91.00 (90.70 to 91.70): CJ0722.D



Abundance Scan 2321 (15.216 min): 101002.D (-2301) (-)



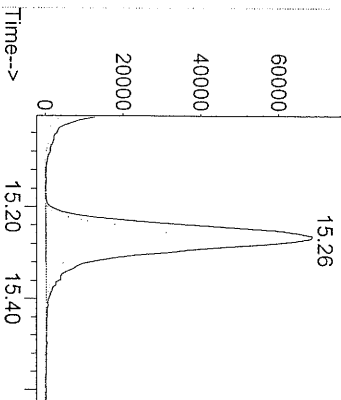
#64
m/p-Xylene
Concen: 5.22 PPB
RT: 15.27 min Scan# 2330
Delta R.T. -0.01 min
Lab File: CJ0722.D
Acq: 29 Oct 2005 00:42

Tgt Ion: 91 Resp: 2721982
Ion Ratio Lower Upper
91 100
106 49.8 45.0 67.4

Raw 50



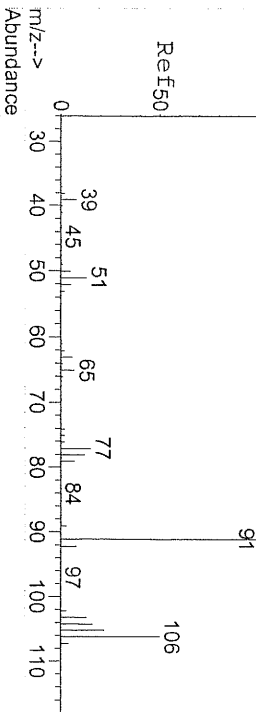
Abundance Ion 91.00 (90.70 to 91.70): CJ0722.D



Sub 50

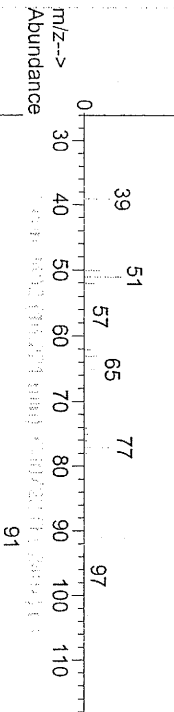


Abundance Scan 2481 (16.156 min): 101002.D (-2457) (-)

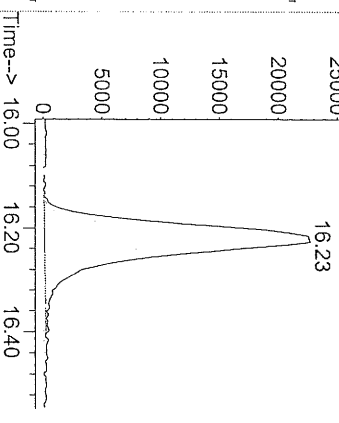


#65
o-Xylene 1.85 PPB
Concn: 1.85 PPB
RT: 16.22 min Scan# 2492
Delta R.T. 0.00 min
Lab File: CJ0722.D
Acq: 29 Oct 2005 00:42
Tgt Ion: 91 Resp: 933865
Ion Ratio Lower Upper
91 100
106 46.9 43.4 65.0

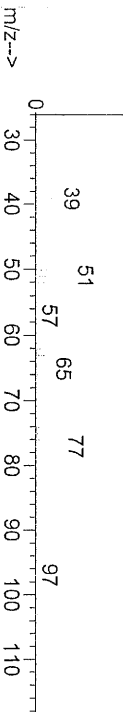
Raw 50



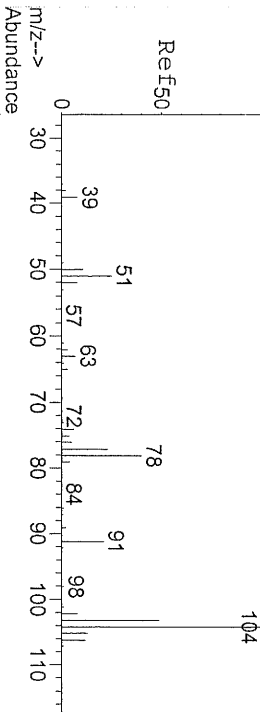
Abundance Ion 91.00 (90.70 to 91.70): CJ



Sub 50

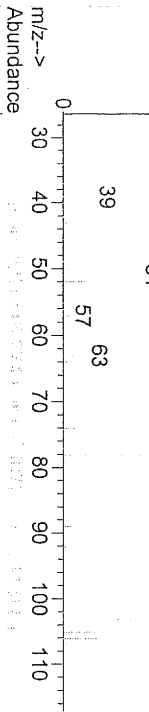


Abundance Scan 2489 (16.203 min): 101002.D (-2467) (-)

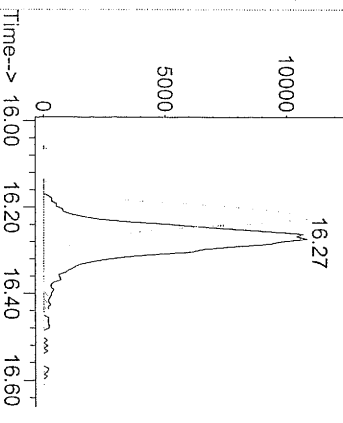


#66
Styrene 1.30 PPB
Concn: 1.30 PPB
RT: 16.27 min Scan# 2501
Delta R.T. -0.00 min
Lab File: CJ0722.D
Acq: 29 Oct 2005 00:42
Tgt Ion: 104 Resp: 477423
Ion Ratio Lower Upper
104 100
106 0.0 0.0 0.0

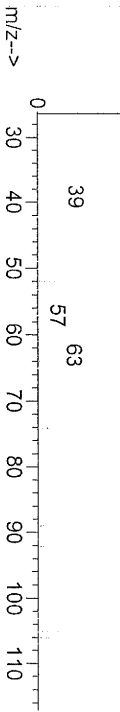
Raw 50



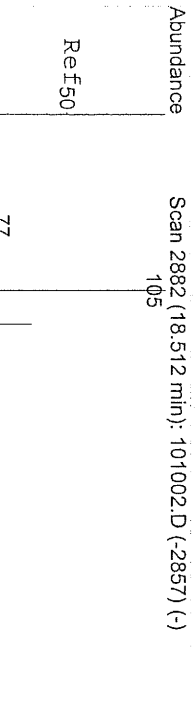
Abundance Ion 104.00 (103.70 to 104.70):



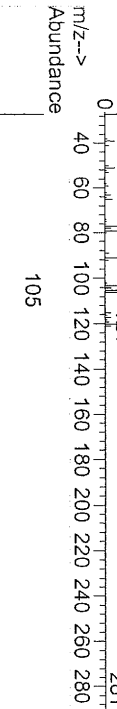
Sub 50



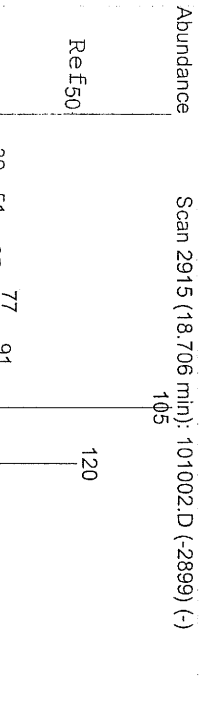
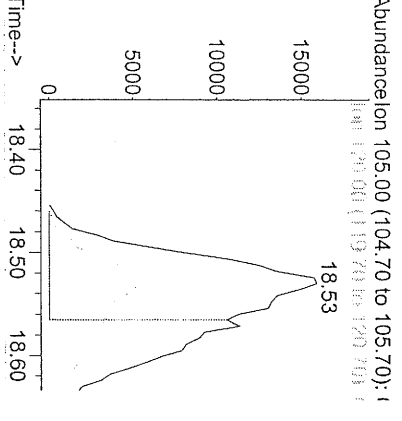
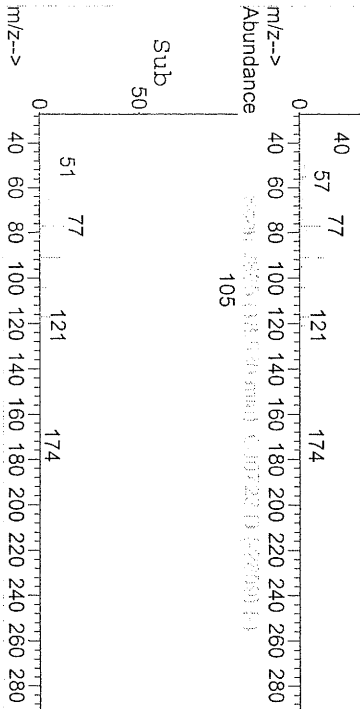
0000



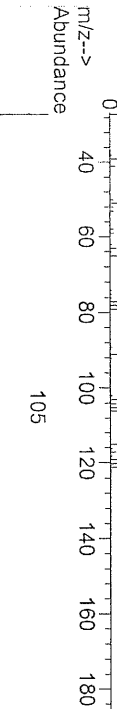
#72
 4-Ethyltoluene
 Concen: 0.81 PPB m
 RT: 18.53 min Scan# 2885
 Delta R.T. -0.05 min
 Lab File: CJ0722.D
 Acq: 29 Oct 2005 00:42



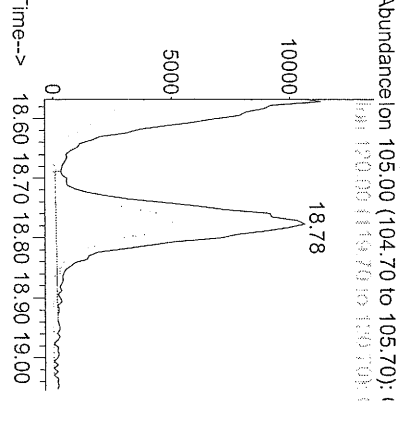
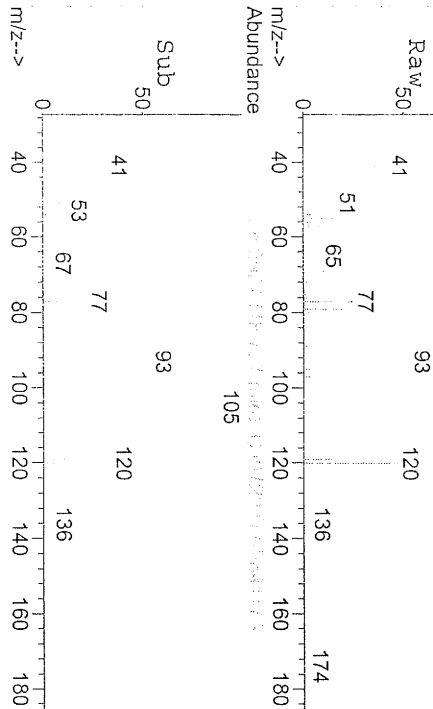
Tgt Ion:105 Resp: 551566
 Ion Ratio Lower Upper
 105 100
 120 50.4 27.0 40.6#



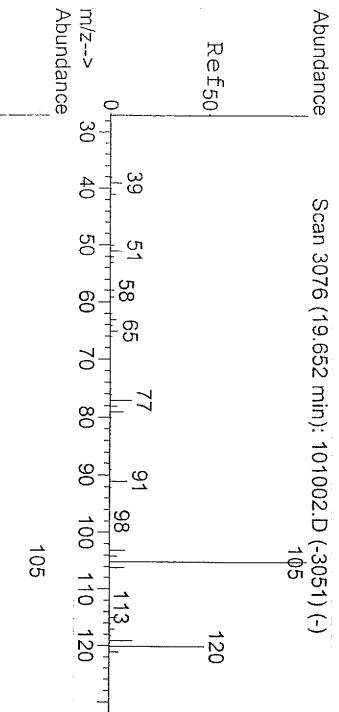
#73
 1,3,5-Trimethylbenzene
 Concen: 0.72 PPB
 RT: 18.78 min Scan# 2927
 Delta R.T. 0.01 min
 Lab File: CJ0722.D
 Acq: 29 Oct 2005 00:42



Tgt Ion:105 Resp: 420885
 Ion Ratio Lower Upper
 105 100
 120 48.6 43.9 65.9

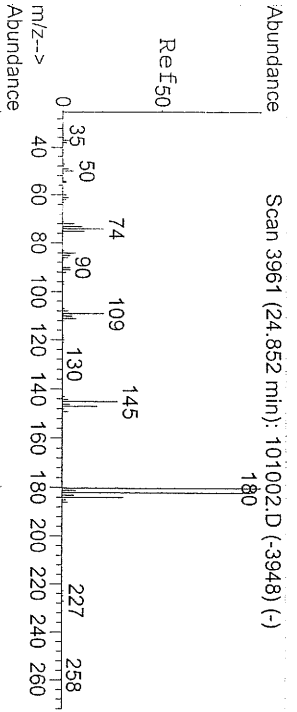
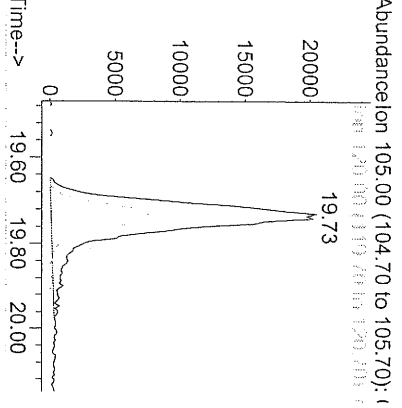
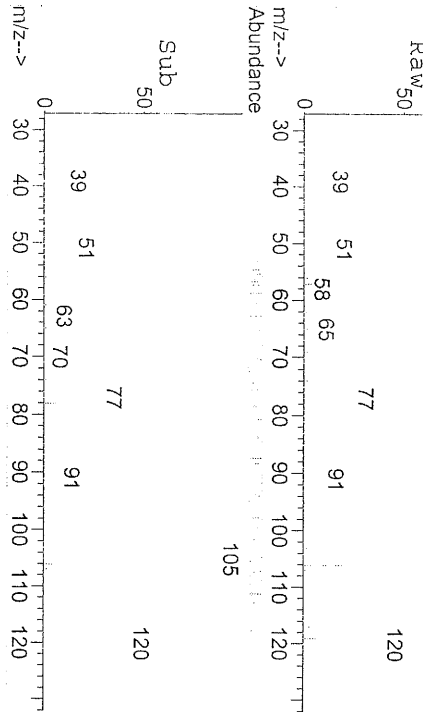


0112



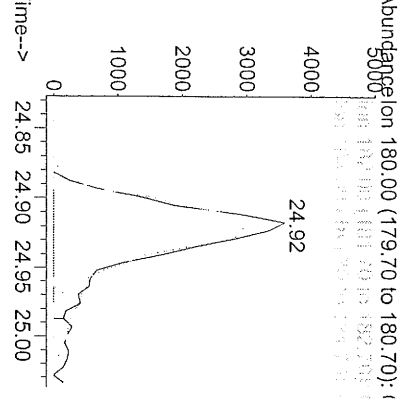
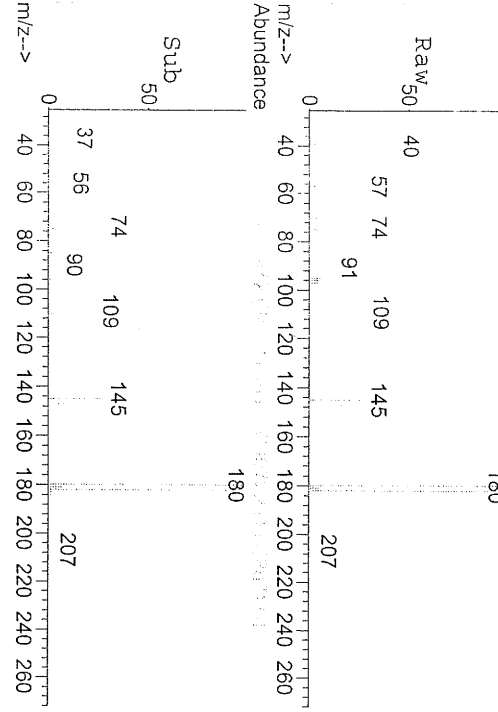
#75
 1,2,4-Trimethylbenzene
 Concn: 1.52 PPB
 RT: 19.74 min Scan# 3090
 Delta R.T. 0.00 min
 Lab File: CJ0722.D
 Acq: 29 Oct 2005 00:42

Tgt Ion: 105	Resp: 880088
Ion Ratio Lower	Upper
105	100
120	35.8
	42.3
	63.5#



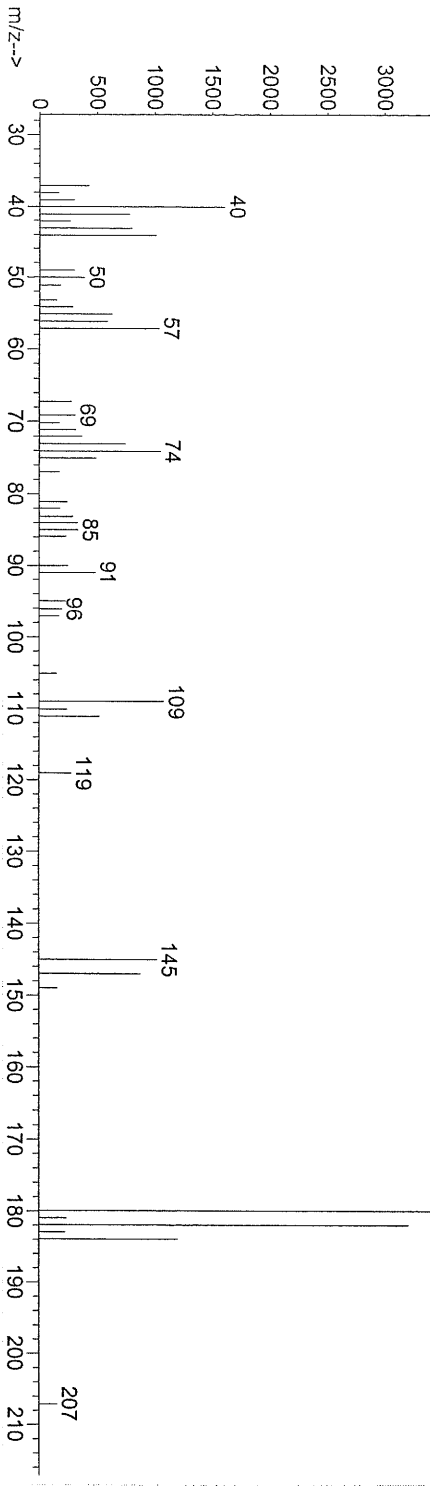
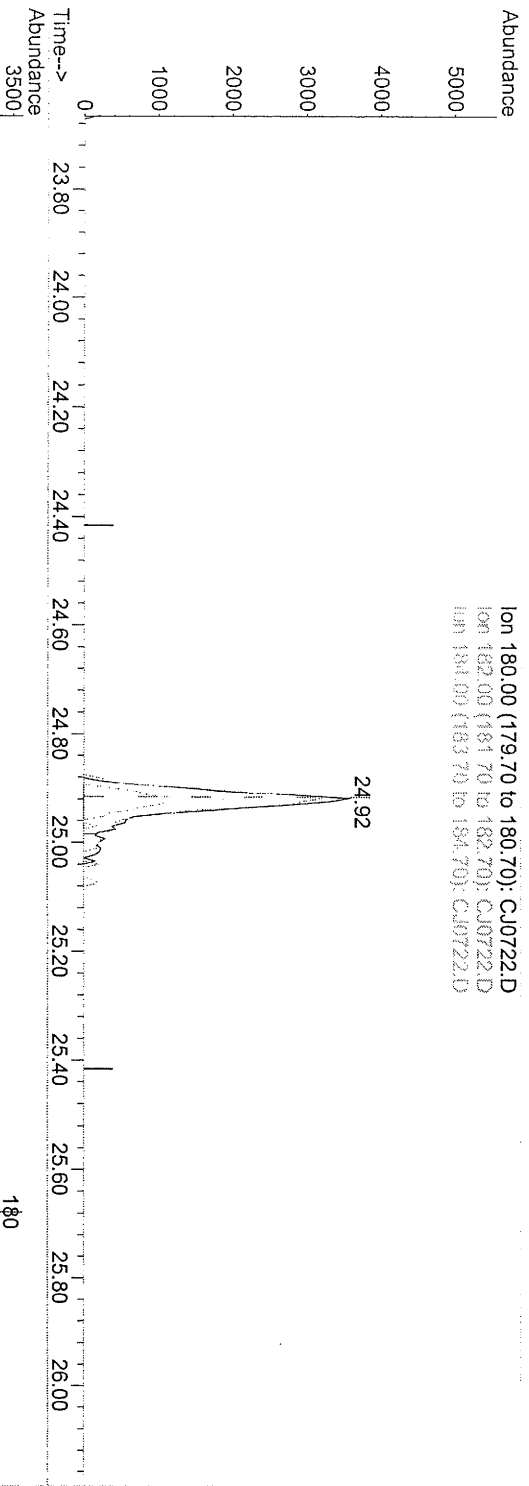
#81
 1,2,4-Trichlorobenzene
 Concn: 0.62 PPB m
 RT: 24.92 min Scan# 3972
 Delta R.T. 0.00 min
 Lab File: CJ0722.D
 Acq: 29 Oct 2005 00:42

Tgt Ion: 180	Resp: 86830
Ion Ratio Lower	Upper
180	100
182	101.6
184	0.0
	76.8
	25.4
	115.2
	38.2#



1111

Data Path : C:\MSDCHEM\1\DATA\OCT28\
Data File : CJ0722.D
Acq On : 29 Oct 2005 00:42
Operator : JBS
Sample : 4628494 500CC
Misc :
ALS Vial : 47 Sample Multiplier: 1
Quant Time: Oct 29 00:09:34 2005
Quant Method : C:\MSDCHEM\1\METHODS\ALLT015.M
Quant Title :
Quant Update : Fri Oct 28 08:49:53 2005
Response via : Initial Calibration



(81) 1,2,4-Trichlorobenzene
24.917min (+0.001) 0.62PPB m
response 86830

Ion	Exp%	Act%
180.00	100	100
182.00	96.00	101.63
184.00	31.80	0.00#
0.00	0.00	0.00

*split peak
182.00
184.00*

Omni 1/12/05

0115

Standards Data

Theoretical Standard Concentrations for EPA Method TO-15.

Compound Name	Cas No	VSTD001	VSTD002	VSTD005	VSTD010	VSTD025
Bromochloromethane	74-97-5	10.00	10.00	10.00	10.00	10.00
Propene	115-07-1	1.01	2.01	5.03	10.05	25.13
Dichlorodifluoromethane	75-71-8	1.01	2.01	5.03	10.05	25.13
Chlorodifluoromethane	75-45-6	1.00	2.00	5.00	10.00	25.00
Freon 114	76-14-2	1.00	2.00	5.00	10.00	25.00
Chloromethane	74-87-3	1.00	2.00	5.00	10.00	25.00
Vinyl Chloride	75-01-4	1.00	2.00	5.00	10.00	25.00
1,3-Butadiene	106-99-0	1.01	2.01	5.03	10.05	25.13
Bromomethane	74-83-9	1.00	2.00	5.00	10.00	25.00
Chloroethane	75-00-3	1.00	2.00	5.00	10.00	25.00
Dichlorofluoromethane	75-43-4	1.01	2.01	5.03	10.05	25.13
Trichlorofluoromethane	75-69-4	1.00	2.00	5.00	10.00	25.00
Pentane	109-66-0	1.05	2.10	5.25	10.50	26.25
Acrolein	107-02-8	1.07	2.14	5.35	10.70	26.75
1,1-Dichloroethene	75-35-4	1.00	2.00	5.00	10.00	25.00
Freon 113	76-13-1	1.00	2.00	5.00	10.00	25.00
Acetone	67-64-1	1.07	2.13	5.33	10.65	26.63
Methyl Iodide	74-88-4	1.05	2.10	5.25	10.50	26.25
Carbon Disulfide	75-15-0	1.09	2.17	5.43	10.85	27.13
Acetonitrile	75-05-8	1.07	2.13	5.33	10.65	26.63
3-Chloropropene	107-05-1	1.00	2.00	5.00	10.00	25.00
Methylene Chloride	75-09-2	1.00	2.00	5.00	10.00	25.00
tert-Butyl Alcohol	75-65-0	1.06	2.11	5.28	10.55	26.38
Acrylonitrile	107-13-1	1.07	2.13	5.33	10.65	26.63
trans-1,2-Dichloroethene	156-60-5	1.05	2.10	5.25	10.50	26.25
Methyl t-Butyl Ether	1634-04-4	1.05	2.10	5.25	10.50	26.25
Hexane	110-54-3	1.06	2.12	5.30	10.60	26.50
1,1-Dichloroethane	75-34-3	0.95	1.90	4.75	9.50	23.75
Vinyl Acetate	108-05-4	1.05	2.09	5.23	10.45	26.13
cis-1,2-Dichloroethene	156-59-2	0.95	1.90	4.75	9.50	23.75
2-Butanone	78-93-3	1.05	2.10	5.25	10.50	26.25
Ethyl Acetate	141-78-6	1.06	2.12	5.30	10.60	26.50
Methyl Acrylate	96-33-3	1.05	2.10	5.25	10.50	26.25
Chloroform	67-66-3	1.00	2.00	5.00	10.00	25.00
1,1,1-Trichloroethane	71-55-6	1.00	2.00	5.00	10.00	25.00
Carbon Tetrachloride	56-23-5	0.95	1.90	4.75	9.50	23.75
1,4-Difluorobenzene	540-36-3	10.00	10.00	10.00	10.00	10.00
1,2-Dichloroethane	107-06-2	1.00	2.00	5.00	10.00	25.00
Benzene	71-43-2	1.00	2.00	5.00	10.00	25.00
Isooctane	540-84-1	1.05	2.10	5.25	10.50	26.25
Heptane	142-82-5	1.06	2.12	5.30	10.60	26.50
Trichloroethene	79-01-6	1.00	2.00	5.00	10.00	25.00
Ethyl Acrylate	140-88-5	1.06	2.12	5.30	10.60	26.50
1,2-Dichloropropane	78-87-5	1.00	2.00	5.00	10.00	25.00
Methyl Methacrylate	80-62-6	1.07	2.13	5.33	10.65	26.63

0117

Theoretical Standard Concentrations for EPA Method TO-15.

Compound Name	Gas No	VSTD001	VSTD002	VSTD005	VSTD010	VSTD025
Dibromomethane	74-95-3	1.07	2.13	5.33	10.65	26.63
1,4-Dioxane	123-91-1	1.07	2.13	5.33	10.65	26.63
Bromodichloromethane	75-27-4	1.06	2.12	5.30	10.60	26.50
cis-1,3-Dichloropropene	10061-01-5	0.90	1.80	4.50	9.00	22.50
4-Methyl-2-Pentanone	108-10-1	1.05	2.10	5.25	10.50	26.25
Chlorobenzene d5	3114-55-4	10.00	10.00	10.00	10.00	10.00
Toluene	108-88-3	1.00	2.00	5.00	10.00	25.00
Octane	111-65-9	1.04	2.08	5.20	10.40	26.00
trans-1,3-Dichloropropene	10061-02-6	0.90	1.80	4.50	9.00	22.50
Ethyl Methacrylate	97-63-2	1.00	1.99	4.98	9.95	24.88
1,1,2-Trichloroethane	79-00-5	0.95	1.90	4.75	9.50	23.75
Tetrachloroethene	127-18-4	1.00	2.00	5.00	10.00	25.00
2-Hexanone	591-78-6	1.05	2.10	5.25	10.50	26.25
Dibromochloromethane	124-48-1	1.00	1.99	4.98	9.95	24.88
1,2-Dibromoethane	106-93-4	1.00	2.00	5.00	10.00	25.00
Chlorobenzene	108-90-7	1.00	2.00	5.00	10.00	25.00
1,1,1,2-Tetrachloroethane	630-20-6	0.95	1.90	4.75	9.50	23.75
Ethylbenzene	100-41-4	1.00	2.00	5.00	10.00	25.00
m/p-Xylene	1330-20-7	1.95	3.90	9.75	19.50	48.75
o-Xylene	95-47-6	1.00	2.00	5.00	10.00	25.00
Styrene	100-42-5	1.00	2.00	5.00	10.00	25.00
Bromoform	75-25-2	1.06	2.11	5.28	10.55	26.38
Cumene	98-82-8	1.05	2.10	5.25	10.50	26.25
1,1,2,2-Tetrachloroethane	79-34-5	1.05	2.10	5.25	10.50	26.25
1,2,3-Trichloropropane	96-18-4	1.05	2.10	5.25	10.50	26.25
Bromobenzene	108-86-1	1.06	2.12	5.30	10.60	26.50
4-Ethyltoluene	622-96-8	0.95	1.90	4.75	9.50	23.75
1,3,5-Trimethylbenzene	108-67-8	1.00	2.00	5.00	10.00	25.00
Alpha Methyl Styrene	611-15-1	1.05	2.10	5.25	10.50	26.25
1,2,4-Trimethylbenzene	95-63-6	0.95	1.90	4.75	9.50	23.75
1,3-Dichlorobenzene	541-73-1	1.00	2.00	5.00	10.00	25.00
1,4-Dichlorobenzene	106-46-7	1.00	2.00	5.00	10.00	25.00
Benzyl chloride	100-44-7	1.00	2.00	5.00	10.00	25.00
1,2-Dichlorobenzene	95-50-1	1.00	2.00	5.00	10.00	25.00
Hexachloroethane	67-72-1	1.06	2.12	5.30	10.60	26.50
1,2,4-Trichlorobenzene	120-82-1	1.00	2.00	5.00	10.00	25.00
Hexachlorobutadiene	87-68-3	0.95	1.90	4.75	9.50	23.75

FORM6
LANCASTER LABORATORIES
VOLATILE ORGANICS IN AIR

INITIAL CALIBRATION DATA SHEET

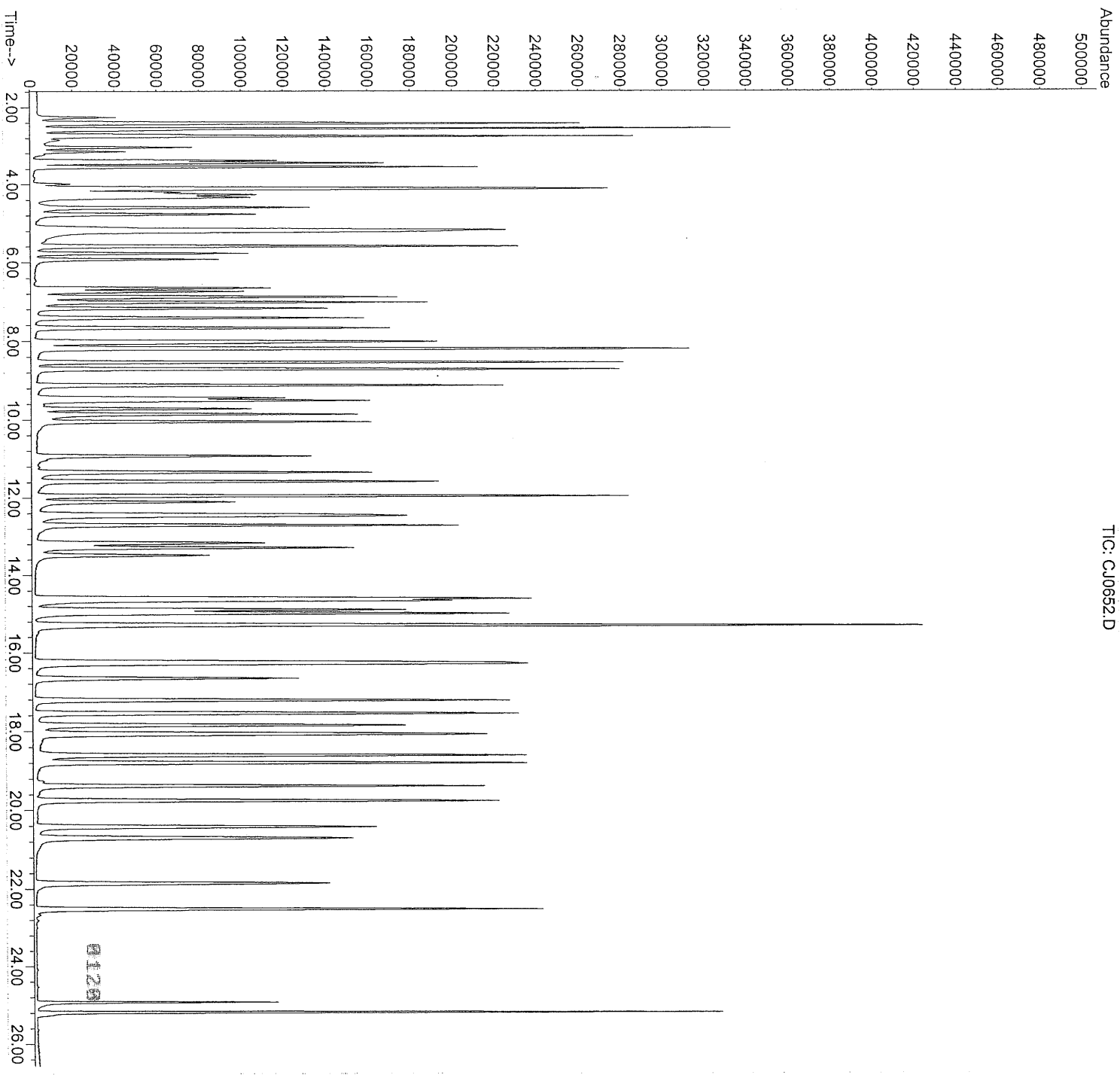
Instrument ID: HP09464

Data File Path: C:\MSDCHEM\1\DATA\OCT27

COMPOUND NAME	Relative Response Factor (RRF)					MEAN	% RSD
	CAL 1	CAL 2	CAL 3	CAL 4	CAL 5		
Dichlorodifluoromethane	3.797	2.920	2.677	2.308	2.936	2.928	19
Freon 114	4.303	3.322	3.038	2.643	3.317	3.324	18
Chloromethane	2.418	1.891	1.720	1.478	1.900	1.882	18
Vinyl Chloride	1.685	1.313	1.208	1.054	1.341	1.320	18
Bromomethane	1.232	1.015	0.919	0.833	1.007	1.001	15
Chloroethane	0.756	0.683	0.611	0.546	0.678	0.655	12
Trichlorofluoromethane	3.556	2.922	2.670	2.304	2.891	2.869	16
1,1-Dichloroethene	2.371	2.107	1.971	1.694	2.148	2.058	12
Freon 113	1.506	1.445	1.330	1.174	1.434	1.378	9
3-Chloropropene	0.612	0.595	0.469	0.414	0.506	0.519	16
Methylene Chloride	0.920	0.858	0.803	0.723	0.895	0.840	9
1,1-Dichloroethane	2.540	2.387	2.228	1.997	2.478	2.326	9
Cis-1,2-Dichloroethene	1.740	1.664	1.539	1.419	1.821	1.637	10
Chloroform	2.306	2.234	2.016	1.806	2.292	2.131	10
1,1,1-Trichloroethane	2.339	2.266	2.105	1.874	2.346	2.186	9
Carbon Tetrachloride	2.497	2.514	2.308	2.033	2.526	2.376	9
1,2-Dichloroethane	0.452	0.376	0.358	0.332	0.492	0.402	17
Benzene	0.900	0.802	0.747	0.676	0.961	0.817	14
Trichloroethene	0.558	0.451	0.410	0.339	0.478	0.447	18
1,2-Dichloropropane	0.410	0.359	0.340	0.297	0.426	0.366	14
Cis-1,3-Dichloropropene	0.489	0.447	0.431	0.390	0.580	0.468	15
Toluene	1.042	1.226	1.100	0.874	1.126	1.074	12
trans-1,3-Dichloropropene	0.407	0.395	0.418	0.361	0.506	0.417	13
1,1,2-Trichloroethane	0.454	0.434	0.400	0.299	0.382	0.394	15
Tetrachloroethene	0.566	0.565	0.484	0.373	0.480	0.494	16
1,2-Dibromoethane	0.581	0.550	0.508	0.463	0.616	0.543	11
Chlorobenzene	0.980	0.920	0.846	0.729	0.954	0.886	11
Ethylbenzene	1.500	1.568	1.497	1.265	1.684	1.503	10
m/p-Xylene	1.215	1.250	1.200	1.022	1.374	1.212	10
o-Xylene	1.159	1.225	1.159	0.990	1.346	1.176	11
Styrene	0.808	0.880	0.825	0.733	1.012	0.852	12
1,1,2,2-Tetrachloroethane	0.543	0.500	0.530	0.478	0.678	0.546	14
4-Ethyltoluene	1.537	1.573	1.574	1.331	1.924	1.588	13
1,3,5-Trimethylbenzene	1.476	1.330	1.328	1.091	1.587	1.362	14
1,2,4-Trimethylbenzene	1.409	1.285	1.325	1.109	1.606	1.347	13
1,3-Dichlorobenzene	0.865	0.695	0.746	0.627	0.922	0.771	16
1,4-Dichlorobenzene	0.815	0.680	0.715	0.592	0.890	0.738	16
1,2-Dichlorobenzene	0.944	0.712	0.681	0.554	0.788	0.736	20
1,2,4-Trichlorobenzene		0.225	0.251	0.200	0.322	0.250	21

Data Path : C:\MSDCHEM\1\DATA\OCT27\
Data File : CJ0652.D
Acq On : 27 Oct 2005 7:58
Operator : JBS
Sample : VSTD010
Misc :
ALS Vial : 27 Sample Multiplier: 1

Quant Time: Oct 27 09:30:56 2005
Quant Method : C:\MSDCHEM\1\METHODS\ALLT015.M
Quant Title :
Qlast Update : Thu Oct 27 09:25:01 2005
Response via : Initial Calibration



Data Path : C:\MSDCHEM\1\DATA\OCT27\
 Data File : CT0652.D
 Acq On : 27 Oct 2005 7:58
 Operator : JBS
 Sample : VSTD010
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Quant Time: Oct 27 09:30:56 2005
 Quant Method : C:\MSDCHEM\1\METHODS\ALLT015.M
 Quant Title :
 Quant Update : Thu Oct 27 09:25:01 2005
 Response via : Initial Calibration

Internal Standards	R.T.	Qion	Response	Conc Units	Dev(Min)
1) Bromochloromethane	7.01	130	838998	10.000	0.00
37) 1,4-Difluorobenzene	8.70	114	3345664	10.000	0.00
51) Chlorobenzene d5	14.59	117	2751751	10.000	0.00

Target Compounds	R.T.	Qion	Response	Conc Units	Dev(Min)	Qvalue
2) Propene	2.39	41	899344	13.852	PPB	96
3) Dichlorodifluoromethane	2.42	85	1945691	13.021	PPB	99
4) Chlorodifluoromethane	2.43	51	1628575	13.407	PPB	96
5) Freon 114	2.54	85	2217428	12.915	PPB	#
6) Chloromethane	2.59	50	1240161	13.606	PPB	96
7) Vinyl Chloride	2.70	62	884548	13.036	PPB	98
8) 1,3-Butadiene	2.74	54	1632019	13.962	PPB	96
9) Bromomethane	3.04	94	698493	13.216	PPB	95
10) Chloroethane	3.15	64	457935	12.856	PPB	97
11) Dichlorofluoromethane	3.39	67	1587167	13.013	PPB	98
12) Trichlorofluoromethane	3.44	101	1933239	12.800	PPB	100
13) Pentane	3.55	43	2118594	14.369	PPB	98
14) Acrolein	3.99	56	278327	15.424	PPB	94
15) 1,1-Dichloroethene	4.09	61	1421085	12.925	PPB	#
16) Freon 113	4.12	103	984952	12.735	PPB	93
17) Acetone	4.21	43	1561923	11.893	PPB	99
18) Methyl Iodide	4.27	142	1748017	13.428	PPB	100
19) Carbon Disulfide	4.35	76	2052125	14.068	PPB	91
20) Acetonitrile	4.60	41	1694232	14.481	PPB	95
21) 3-Chloropropene	4.60	76	347067	11.809	PPB	85
22) Methylene Chloride	4.77	84	606425	9.574	PPB	77
23) tert-Butyl Alcohol	5.12	59	1816260	13.943	PPB	100
24) Acrylonitrile	5.15	53	1053737	14.900	PPB	99
25) trans-1,2-Dichloroethene	5.16	61	1218909	14.064	PPB	77
26) Methyl t-Butyl Ether	5.21	73	2159096	13.813	PPB	85
27) Hexane	5.59	57	1693243	12.987	PPB	#
28) 1,1-Dichloroethane	5.77	63	1591337	12.220	PPB	99
29) Vinyl Acetate	5.92	43	2271652	17.351	PPB	94
30) cis-1,2-Dichloroethene	6.65	61	1131326	12.483	PPB	#
31) 2-Butanone	6.74	72	359256	13.183	PPB	100
32) Ethyl Acetate	6.87	70	171658	13.385	PPB	#
33) Methyl Acrylate	6.90	55	1718293	14.587	PPB	94
34) Chloroform	7.17	83	1515320	12.758	PPB	96
35) 1,1,1-Trichloroethane	7.41	97	1571954	13.002	PPB	94
36) Carbon Tetrachloride	7.67	117	1620188	12.311	PPB	100
38) 1,2-Dichloroethane	8.00	62	1112412	13.787	PPB	99
39) Benzene	8.00	78	2261134	12.828	PPB	96
40) Isooctane	8.18	57	5283084	14.153	PPB	95
41) Heptane	8.53	43	2335529	14.981	PPB	92
42) Trichloroethene	9.12	130	1134795	12.457	PPB	88
43) Ethyl Acrylate	9.44	55	2322802	14.270	PPB	91
44) 1,2-Dichloropropane	9.51	63	993304	12.553	PPB	97
45) Methyl Methacrylate	9.85	69	767128	13.126	PPB	80
46) Dibromomethane	9.72	174	661425	13.909	PPB	78
47) 1,4-Dioxane	9.88	88	240566	7.760	PPB	1
48) Bromodichloromethane	10.05	83	1626480	13.855	PPB	92

0121

Data Path : C:\MSDCHEM\1\DATA\OCT27\
 Data File : CJ0652.D
 Acq On : 27 Oct 2005 7:58
 Operator : JBS
 Sample : VSTD010
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Quant Time: Oct 27 09:30:56 2005

Quant Method : C:\MSDCHEM\1\METHODS\ALLT015.M

Quant Title :

Quant Update : Thu Oct 27 09:25:01 2005

Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
49) cis-1,3-Dichloropropene	10.93	75	1174874	11.729 PPB	# 71
50) 4-Methyl-2-Pentanone	11.33	43	2641430	12.204 PPB	# 96
52) Toluene	11.57	91	2404832	11.917 PPB	# 97
53) Octane	11.94	43	2983990	13.325 PPB	# 89
54) trans-1,3-Dichloropropene	12.10	75	893009	12.047 PPB	# 88
55) Ethyl Methacrylate	12.42	69	1172777	11.570 PPB	# 79
56) 1,1,2-Trichloroethane	12.46	97	780338	11.115 PPB	# 85
57) Tetrachloroethene	12.70	166	1026723	11.923 PPB	# 98
58) 2-Hexanone	13.16	43	2292001	11.333 PPB	# 74
59) Dibromochloromethane	13.28	127	1270854	13.243 PPB	# 99
60) 1,2-Dibromoethane	13.48	107	1274087	13.283 PPB	# 98
61) Chlorobenzene	14.65	112	2006277	12.867 PPB	# 85
62) 1,1,1,2-Tetrachloroethane	14.89	131	1146042	12.239 PPB	# 98
63) Ethylbenzene	14.98	91	3481941	13.168 PPB	# 93
64) m/p-Xylene	15.28	91	5484720	25.707 PPB	# 91
65) o-Xylene	16.23	91	2723080	13.325 PPB	# 90
66) Styrene	16.28	104	2017291	13.614 PPB	# 100
67) Bromoform	16.65	173	1339226	14.330 PPB	# 97
68) Cumene	17.21	105	3392516	14.812 PPB	# 94
69) 1,1,2,2-Tetrachloroethane	18.04	83	1380475	13.916 PPB	# 96
70) 1,2,3-Trichloropropane	18.07	110	486070	14.100 PPB	# 65
71) Bromobenzene	17.84	156	1041536	14.648 PPB	# 99
72) 4-Ethyltoluene	18.60	105	3479136	13.537 PPB	# 92
73) 1,3,5-Trimethylbenzene	18.79	105	3001578	13.718 PPB	# 90
74) Alpha Methyl Styrene	19.37	118	1454491	14.243 PPB	# 94
75) 1,2,4-Trimethylbenzene	19.75	105	2899380	12.963 PPB	# 89
76) 1,3-Dichlorobenzene	20.41	146	1726062	13.874 PPB	# 96
77) 1,4-Dichlorobenzene	20.70	146	1628337	13.372 PPB	# 96
78) Benzyl Chloride	16.23	91	2723080	13.325 PPB	# 100
79) 1,2-Dichlorobenzene	21.85	146	1524163	13.236 PPB	# 96
80) Hexachloroethane	22.51	117	1164521	13.990 PPB	# 97
81) 1,2,4-Trichlorobenzene	24.92	180	551388	11.036 PPB	# 97
82) Hexachlorobutadiene	25.18	225	830677	11.617 PPB	# 97

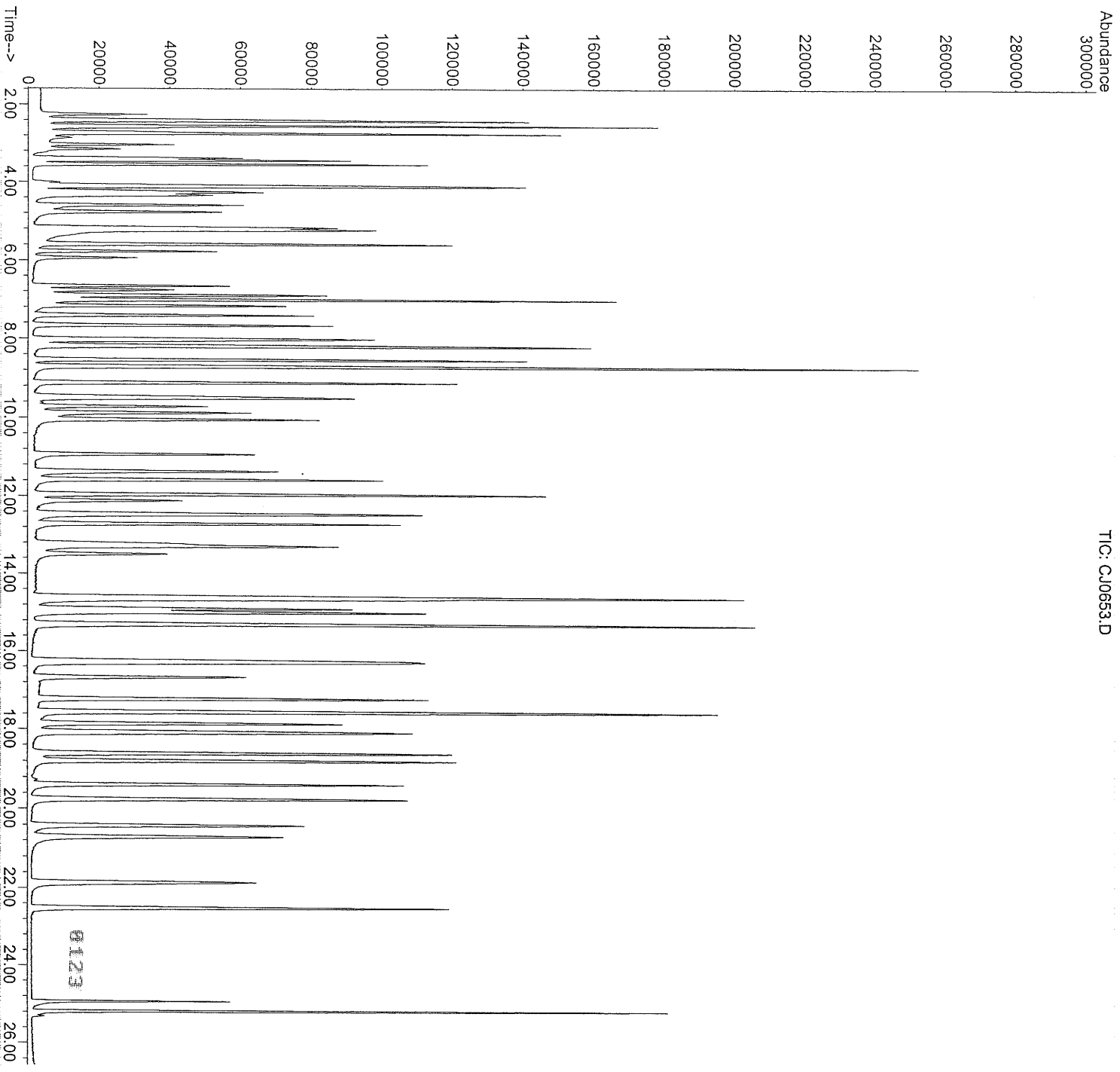
(#) = qualifier out of range (m) = manual integration (+) = signals summed

0122

Data Path : C:\MSDCHEM\1\DATA\OCT27\
Data File : CJ0653.D
Acq On : 27 Oct 2005 8:35
Operator : JBS
Sample : VSTD005
Misc :
ALS Vial : 28 Sample Multiplier: 1

Quant Time: Oct 27 13:02:40 2005
Quant Method : C:\MSDCHEM\1\METHODS\ALLT015.M
Quant Title :
Quant Update : Thu Oct 27 12:54:28 2005
Response via : Initial Calibration

TIC: CJ0653.D



Data Path : C:\MSDCHEM\1\DATA\OCT27\
Data File : CJ0653.D
Acq On : 27 Oct 2005 8:35
Operator : JBS
Sample : VSTD005
Misc :
ALS Vial : 28 Sample Multiplier: 1

Quant Time: Oct 27 13:02:40 2005

Quant Method : C:\MSDCHEM\1\METHODS\ALLT015.M

Quant Title :

Quant Update : Thu Oct 27 12:54:28 2005

Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
1) Bromochloromethane	7.02	130	772479	10.000	0.00
37) 1,4-Difluorobenzene	8.71	114	3124266	10.000	0.00
51) Chlorobenzene d5	14.60	117	2342017	10.000	0.01

Target Compounds	R.T.	QIon	Response	Conc Units	Dev(Min)	Qvalue
2) Propene	2.39	41	489592	4.599 PPB		96
3) Dichlorodifluoromethane	2.42	85	1039144	4.596 PPB		98
4) Chlorodifluoromethane	2.43	51	878350	4.590 PPB		94
5) Freon 114	2.54	85	1173292	4.570 PPB	#	81
6) Chloromethane	2.59	50	664385	4.577 PPB		97
7) Vinyl Chloride	2.70	62	466436	4.575 PPB		98
8) 1,3-Butadiene	2.75	54	847194	4.641 PPB		97
9) Bromomethane	3.04	94	354895	4.590 PPB		97
10) Chloroethane	3.16	64	236053	4.667 PPB		99
11) Dichlorofluoromethane	3.39	67	823157	4.757 PPB		98
12) Trichlorofluoromethane	3.44	101	1031326	4.655 PPB		99
13) Pentane	3.55	43	1122702	4.968 PPB		97
14) Acrolein	4.01	56	123515	4.979 PPB	#	95
15) 1,1-Dichloroethene	4.09	61	761333	4.789 PPB		87
16) Freon 113	4.12	103	513701	4.827 PPB		93
17) Acetone	4.23	43	789628	5.082 PPB		97
18) Methyl Iodide	4.27	142	917690	5.000 PPB		99
19) Carbon Disulfide	4.35	76	1055303	5.031 PPB	#	91
20) Acetonitrile	4.60	41	820519	5.163 PPB		97
21) 3-Chloropropene	4.60	76	181155	4.516 PPB		83
22) Methylene Chloride	4.77	84	310285	4.786 PPB	#	77
23) tert-Butyl Alcohol	5.24	59	739781	4.123 PPB		100
24) Acrylonitrile	5.17	53	528570	5.127 PPB		93
25) trans-1,2-Dichloroethene	5.17	61	609316	4.999 PPB	#	79
26) Methyl t-Butyl Ether	5.23	73	1091585	5.047 PPB	#	86
27) Hexane	5.59	57	872018	5.032 PPB	#	86
28) 1,1-Dichloroethane	5.78	63	817696	4.552 PPB	#	99
29) Vinyl Acetate	5.93	43	827182	4.525 PPB		94
30) cis-1,2-Dichloroethene	6.66	61	564698	4.477 PPB	#	80
31) 2-Butanone	6.77	72	176358	5.160 PPB	#	100
32) Ethyl Acetate	6.90	70	81100	5.288 PPB	#	100
33) Methyl Acrylate	6.91	55	813663	4.977 PPB	#	94
34) Chloroform	7.17	83	778636	4.732 PPB		97
35) 1,1,1-Trichloroethane	7.41	97	813112	4.816 PPB		93
36) Carbon Tetrachloride	7.67	117	846772	4.628 PPB		100
38) 1,2-Dichloroethane	8.05	62	559592	4.453 PPB		100
39) Benzene	8.01	78	1167404	4.573 PPB	#	95
40) Isooctane	8.19	57	2712511	4.800 PPB	#	96
41) Heptane	8.53	43	1185142	4.698 PPB	#	93
42) Trichloroethene	9.12	130	640146	4.581 PPB	#	89
43) Ethyl Acrylate	9.47	55	1080012	4.769 PPB	#	84
44) 1,2-Dichloropropane	9.51	63	530849	4.639 PPB	#	94
45) Methyl Methacrylate	9.88	69	388970	5.032 PPB	#	85
46) Dibromomethane	9.72	174	340757	4.897 PPB	#	80
47) 1,4-Dioxane	9.96	88	155284	4.547 PPB	#	21
48) Bromodichloromethane	10.05	83	843118	4.796 PPB	#	93

Data Path : C:\MSDCHEM\1\DATA\OCT27\
 Data File : CJ0653.D
 Acq On : 27 Oct 2005 8:35
 Operator : JBS
 Sample : VSTD005
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Quant Time: Oct 27 13:02:40 2005
 Quant Method : C:\MSDCHEM\1\METHODS\ALLT015.M
 Quant Title :
 QIast Update : Thu Oct 27 12:54:28 2005
 Response via : Initial Calibration

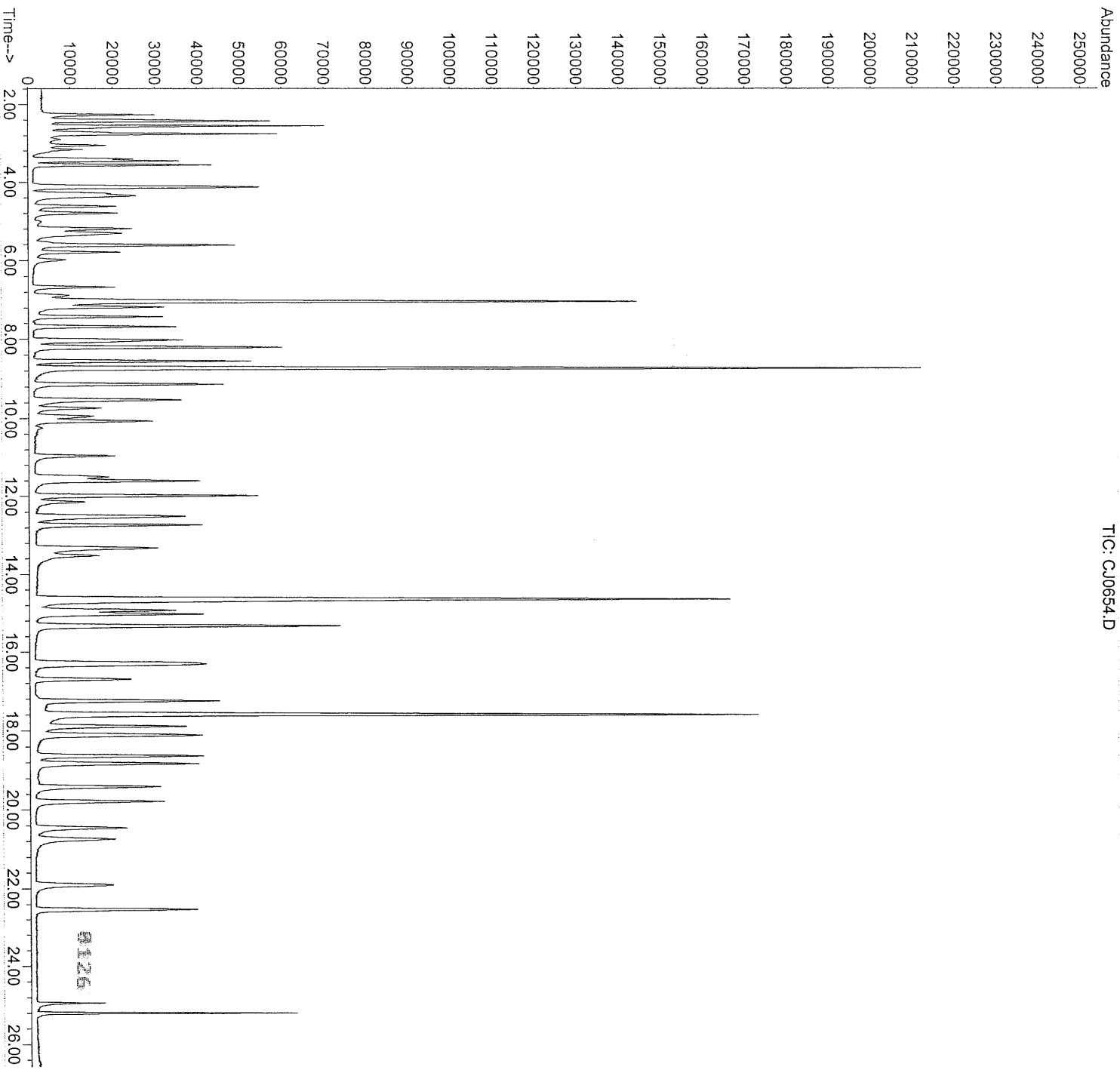
Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
49) cis-1,3-Dichloropropene	10.93	75	605872	4.148	PPB	# 69
50) 4-Methyl-2-Pentanone	11.37	43	1307857	4.326	PPB	# 95
52) Toluene	11.58	91	1288102	5.120	PPB	# 100
53) Octane	11.95	43	1601281	5.338	PPB	# 90
54) trans-1,3-Dichloropropene	12.12	75	440670	4.505	PPB	# 93
55) Ethyl Methacrylate	12.45	69	589088	4.977	PPB	# 80
56) 1,1,2-Trichloroethane	12.48	97	444593	4.818	PPB	# 87
57) Tetrachloroethene	12.71	166	566474	4.899	PPB	# 98
58) 2-Hexanone	13.23	43	1111009	4.927	PPB	# 80
59) Dibromochloromethane	13.29	127	611626	4.714	PPB	# 99
60) 1,2-Dibromoethane	13.50	107	594994	4.672	PPB	# 99
61) Chlorobenzene	14.66	112	990206	4.770	PPB	# 82
62) 1,1,1,2-Tetrachloroethane	14.90	131	592309	4.697	PPB	# 98
63) Ethylbenzene	14.99	91	1753256	4.978	PPB	# 92
64) m/p-Xylene	15.30	91	2739470	9.643	PPB	# 91
65) o-Xylene	16.25	91	1357639	4.931	PPB	# 91
66) Styrene	16.30	104	966113	4.836	PPB	# 100
67) Bromoform	16.67	173	654320	4.886	PPB	# 97
68) Cumene	17.22	105	1622301	5.084	PPB	# 93
69) 1,1,2,2-Tetrachloroethane	18.06	83	651725	5.096	PPB	# 97
70) 1,2,3-Trichloropropane	18.09	110	246163	4.971	PPB	# 61
71) Bromobenzene	17.86	156	507668	4.877	PPB	# 98
72) 4-Ethyltoluene	18.61	105	1750621	4.706	PPB	# 92
73) 1,3,5-Trimethylbenzene	18.80	105	1554617	4.870	PPB	# 90
74) Alpha Methyl Styrene	19.38	118	755902	5.319	PPB	# 94
75) 1,2,4-Trimethylbenzene	19.76	105	1473628	4.670	PPB	# 89
76) 1,3-Dichlorobenzene	20.42	146	873894	4.841	PPB	# 96
77) 1,4-Dichlorobenzene	20.71	146	836793	4.830	PPB	# 95
78) Benzyl Chloride	16.25	91	1357639	4.931	PPB	# 100
79) 1,2-Dichlorobenzene	21.87	146	797342	4.624	PPB	# 96
80) Hexachloroethane	22.51	117	598632	5.081	PPB	# 97
81) 1,2,4-Trichlorobenzene	24.93	180	294041	3.915	PPB	# 98
82) Hexachlorobutadiene	25.19	225	434013	4.159	PPB	# 97

(#) = qualifier out of range (m) = manual integration (+) = signals summed

8125

Data Path : C:\MSDCHEM\1\DATA\OCT27\
Data File : CJ0654.D
Acq On : 27 Oct 2005 9:13
Operator : JBS
Sample : VSTD002
Misc :
ALS Vial : 29 Sample Multiplier: 1

Quant Time: Oct 27 09:41:12 2005
Quant Method : C:\MSDCHEM\1\METHODS\ALLT015.M
Quant Title :
Quant Update : Thu Oct 27 09:25:01 2005
Response via : Initial Calibration



Data Path : C:\MSDCHEM\1\DATA\OCT27\
 Data File : C:\MSDCHEM\1\DATA\OCT27\
 Acq On : 27 Oct 2005 9:13
 Operator : JBS
 Sample : VSTD002
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Quant Time: Oct 27 09:41:12 2005
 Quant Method : C:\MSDCHEM\1\METHODS\ALLT015.M
 Quant Title :
 QLast Update : Thu Oct 27 09:25:01 2005
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) Bromochloromethane	7.03	130	701657	10.000	PPB	0.02
37) 1,4-Difluorobenzene	8.72	114	2892749	10.000	PPB	0.02
51) Chlorobenzene d5	14.62	117	2099790	10.000	PPB	0.03

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Dev (Min)	Qvalue
2) Propene	2.39	41	195347	3.598	PPB		96
3) Dichlorodifluoromethane	2.42	85	411878	3.296	PPB		100
4) Chlorodifluoromethane	2.44	51	348484	3.430	PPB		95
5) Freon 114	2.55	85	466221	3.247	PPB	#	81
6) Chloromethane	2.60	50	265386	3.481	PPB		97
7) Vinyl Chloride	2.71	62	184248	3.247	PPB		100
8) 1,3-Butadiene	2.75	54	338409	3.462	PPB		95
9) Bromomethane	3.05	94	142403	3.222	PPB		95
10) Chloroethane	3.17	64	95900	3.219	PPB	#	43
11) Dichlorofluoromethane	3.40	67	335851	3.293	PPB	#	89
12) Trichlorofluoromethane	3.45	101	410047	3.246	PPB		97
13) Pentane	3.56	43	440177	3.570	PPB	#	97
14) Acrolein	4.07	56	41232	2.732	PPB		19
15) 1,1-Dichloroethene	4.10	61	295701	3.216	PPB	#	86
16) Freon 113	4.14	103	202832	3.136	PPB		93
17) Acetone	4.33	43	296480m	2.699	PPB		
18) Methyl Iodide	4.28	142	362414	3.329	PPB		99
19) Carbon Disulfide	4.36	76	398470	3.266	PPB		97
20) Acetonitrile	4.62	41	249485	2.550	PPB	#	85
21) 3-Chloropropene	4.79	84	83549m	3.399	PPB		
22) Methylene Chloride	4.62	76	120335	2.272	PPB	#	66
23) tert-Butyl Alcohol	5.55	59	315808m	2.899	PPB		
24) Acrylonitrile	5.21	53	166602	2.817	PPB	#	71
25) trans-1,2-Dichloroethene	5.19	61	220937	3.048	PPB	#	80
26) Methyl t-Butyl Ether	5.31	73	402461	3.079	PPB	#	48
27) Hexane	5.61	57	364279	3.341	PPB	#	90
28) 1,1-Dichloroethane	5.79	63	318287	2.923	PPB		98
29) Vinyl Acetate	5.98	43	242255	2.212	PPB	#	78
30) cis-1,2-Dichloroethene	6.68	61	221805	2.926	PPB	#	77
31) 2-Butanone	6.88	72	54960	2.412	PPB	#	100
32) Ethyl Acetate	6.99	70	24305	2.266	PPB	#	76
33) Methyl Acrylate	6.99	55	257940	2.618	PPB	#	
34) Chloroform	7.18	83	313535	3.157	PPB		96
35) 1,1,1-Trichloroethane	7.43	97	317921	3.144	PPB		94
36) Carbon Tetrachloride	7.68	117	335210	3.046	PPB		99
38) 1,2-Dichloroethane	8.07	62	217816	3.122	PPB		96
39) Benzene	8.02	78	464135	3.045	PPB		96
40) Isooctane	8.20	57	1040761	3.225	PPB	#	96
41) Heptane	8.55	43	455865	3.382	PPB	#	93
42) Trichloroethene	9.14	130	261013	3.314	PPB	#	90
43) Ethyl Acrylate	9.55	55	347987	2.473	PPB	#	77
44) 1,2-Dichloropropane	9.53	63	207679	3.035	PPB		95
45) Methyl Methacrylate	9.95	69	133326	2.639	PPB	#	73
46) Dibromomethane	9.74	174	137085	3.334	PPB	#	79
47) 1,4-Dioxane	10.26	88	46700m	1.742	PPB		
48) Bromodichloromethane	10.07	83	334451	3.295	PPB	#	94

0127

Data Path : C:\MSDCHEM\1\DATA\OCT27\
 Data File : CJ0654.D
 Acq On : 27 Oct 2005 9:13
 Operator : JBS
 Sample : VSTD002
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

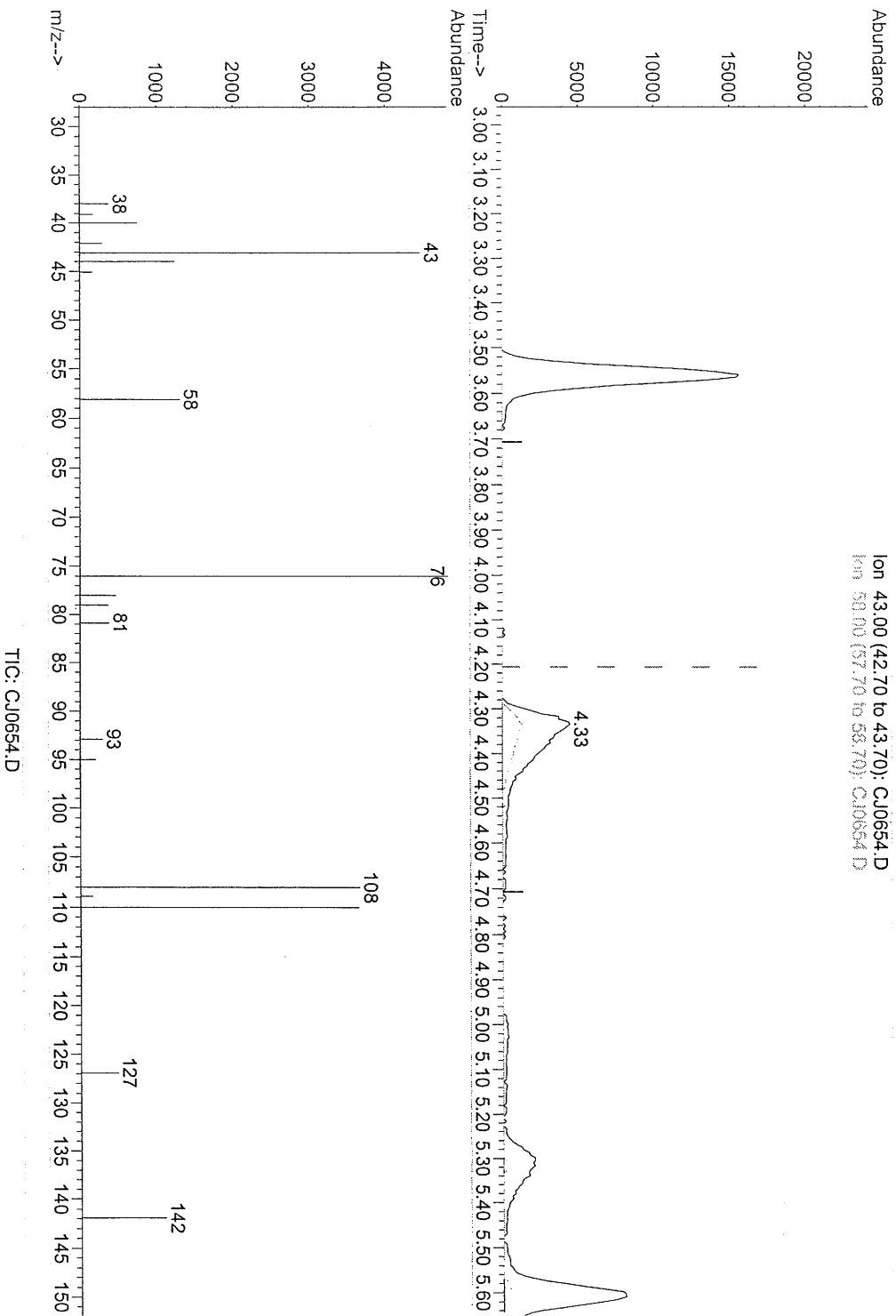
Quant Time: Oct 27 09:41:12 2005
 Quant Method : C:\MSDCHEM\1\METHODS\ALLT015.M
 Quant Title :
 Quant Update : Thu Oct 27 09:25:01 2005
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
49) cis-1,3-Dichloropropene	10.96	75	232765	2.688 PPB	75
50) 4-Methyl-2-Pentanone	11.50	43	519191	2.774 PPB	90
52) Toluene	11.60	91	514682	3.342 PPB	97
53) Octane	11.98	43	613642	3.591 PPB	90
54) trans-1,3-Dichloropropene	12.15	75	149391	2.641 PPB	84
55) Ethyl Methacrylate	12.53	69	214565	2.774 PPB	80
56) 1,1,2-Trichloroethane	12.50	97	173298	3.235 PPB	84
57) Tetrachloroethene	12.73	166	237244	3.610 PPB	99
58) 2-Hexanone	13.35	43	305195	1.978 PPB	97
59) Dibromochloromethane	13.32	127	243654	3.327 PPB	98
60) 1,2-Dibromoethane	13.52	107	230909	3.155 PPB	98
61) Chlorobenzene	14.69	112	386539	3.249 PPB	79
62) 1,1,1,2-Tetrachloroethane	14.92	131	228995	3.205 PPB	74
63) Ethylbenzene	15.02	91	658644	3.264 PPB	92
64) m/p-Xylene	15.32	91	1023607	6.287 PPB	89
65) o-Xylene	16.27	91	514313	3.298 PPB	90
66) Styrene	16.33	104	369519	3.268 PPB	100
67) Bromoform	16.70	173	257217	3.607 PPB	96
68) Cumene	17.25	105	625092	3.577 PPB	92
69) 1,1,2,2-Tetrachloroethane	18.08	83	220356	2.911 PPB	94
70) 1,2,3-Trichloropropane	18.12	110	92317	3.509 PPB	50
71) Bromobenzene	17.88	156	194554	3.586 PPB	99
72) 4-Ethyltoluene	18.63	105	627447	3.199 PPB	93
73) 1,3,5-Trimethylbenzene	18.82	105	558392	3.344 PPB	94
74) Alpha Methyl Styrene	19.40	118	253161	3.249 PPB	91
75) 1,2,4-Trimethylbenzene	19.78	105	512492	3.003 PPB	90
76) 1,3-Dichlorobenzene	20.45	146	292009	3.076 PPB	93
77) 1,4-Dichlorobenzene	20.74	146	285466	3.072 PPB	86
78) Benzyl Chloride	16.27	91	514313	3.298 PPB	#
79) 1,2-Dichlorobenzene	21.90	146	299161	3.405 PPB	100
80) Hexachloroethane	22.53	117	213633	3.363 PPB	94
81) 1,2,4-Trichlorobenzene	24.94	180	94332	2.474 PPB	97
82) Hexachlorobutadiene	25.19	225	160309	2.938 PPB	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

0128

Data Path : C:\MSDCHEM\1\DATA\OCT27\
 Data File : CJ0654.D
 Acq On : 27 Oct 2005 9:13
 Operator : JBS
 Sample : VSTD002
 Misc :
 ALS Vial : 29 Sample Multiplier: 1
 Quant Time: Oct 27 09:41:12 2005
 Quant Method : C:\MSDCHEM\1\METHODS\ALLTO15.M
 Quant Title :
 Qlast Update : Thu Oct 27 09:25:01 2005
 Response via : Initial Calibration



Ion 43.00 (42.70 to 43.70): CJ0654.D
 Ion 58.00 (57.70 to 58.70): CJ0654.D

(17) Acetone

4.335min (+0.128) 2.70PPB m

response 296480

Ion	Exp%	Act%
43.00	100	100
58.00	28.20	0.00#
0.00	0.00	0.00
0.00	0.00	0.00

Silent peak

AKASH
10/27/05

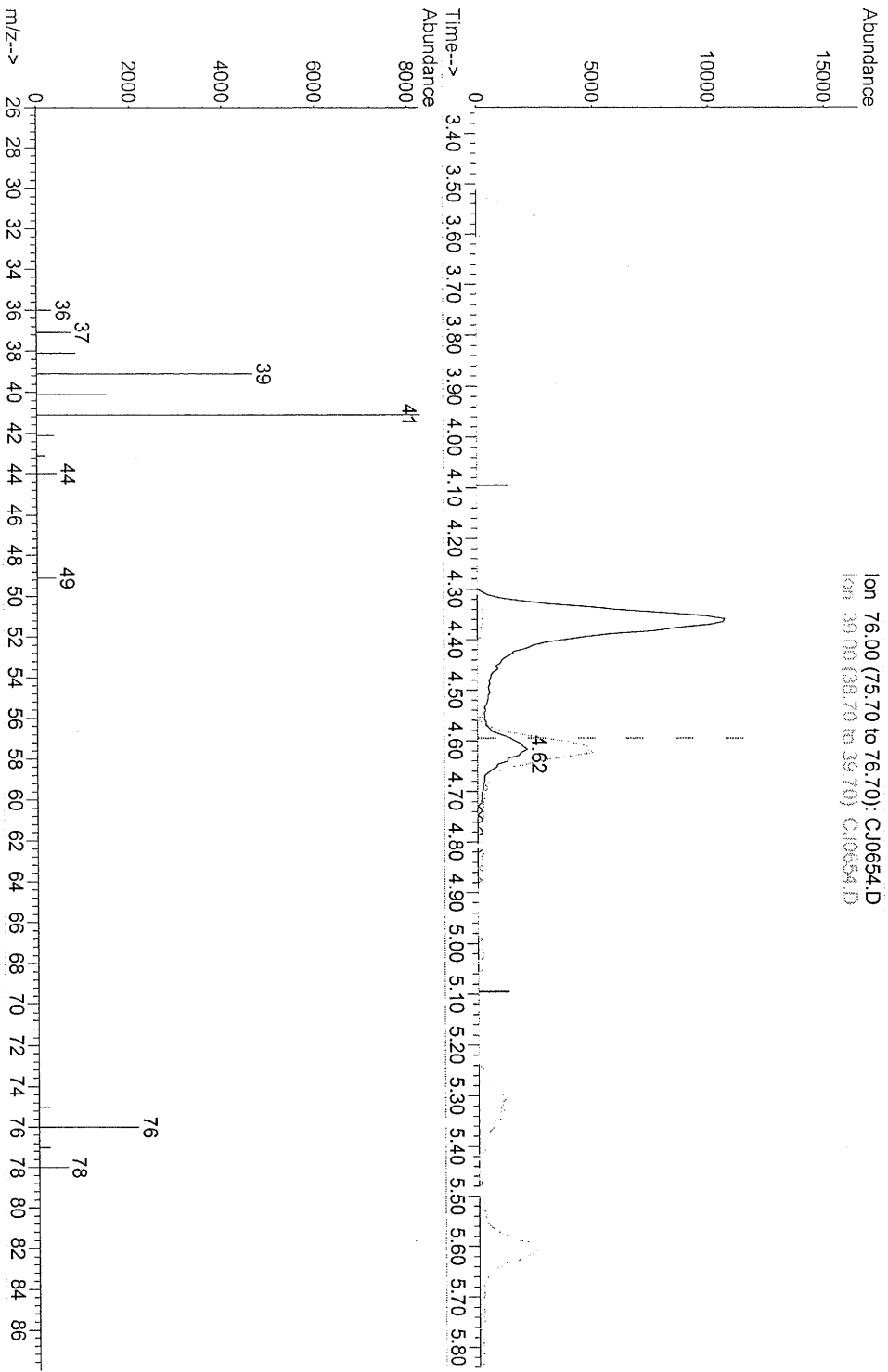
CMR/412
11/1/05

0129

Quantitation Report (Qedit)

Data Path : C:\MSDCHEM\1\DATA\OCT27\
Data File : CJ0654.D
Acq On : 27 Oct 2005 9:13
Operator : JBS
Sample : VSTD002
Misc :
ALS Vial : 29 Sample Multiplier: 1

Quant Time: Oct 27 09:41:12 2005
Quant Method : C:\MSDCHEM\1\METHODS\ALLT015.M
Quant Title :
Quant Update : Thu Oct 27 09:25:01 2005
Response via : Initial Calibration



(21) 3-Chloropropene

4.617min (+0.021) 3.40PPB m

response 83549

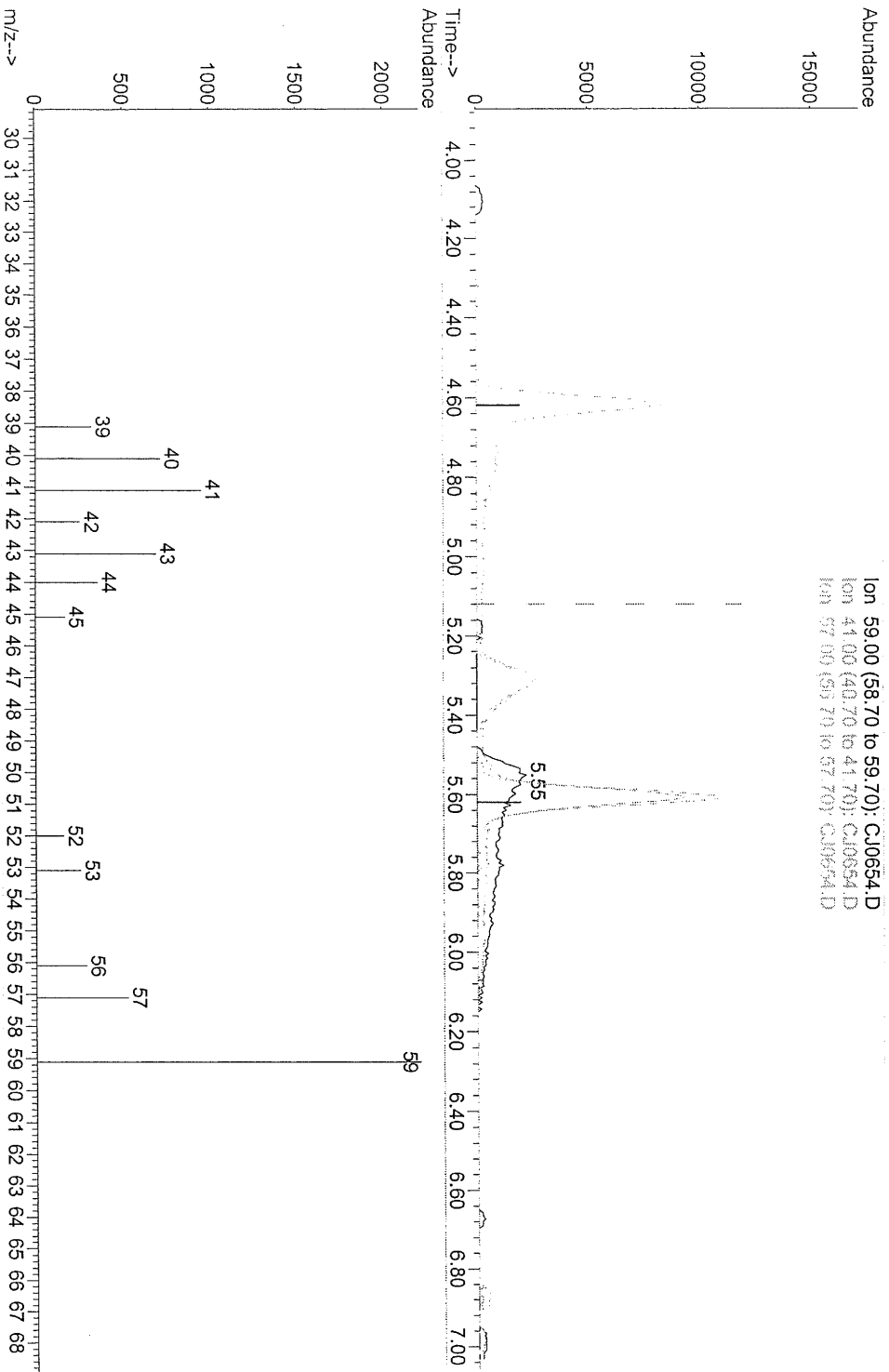
Ion	Exp%	Act%
76.00	100	100
39.00	289.60	0.00#
0.00	0.00	0.00
0.00	0.00	0.00

Spun-punk
11/11/05

Om2/412
11/11/05

Data Path : C:\MSDCHEM\1\DATA\OCT27\
Data File : C:\0654.D
Acq On : 27 Oct 2005 9:13
Operator : JBS
Sample : VSTD002
Misc :
ALS Vial : 29 Sample Multiplier: 1

Quant Time: Oct 27 09:41:12 2005
Quant Method : C:\MSDCHEM\1\METHODS\ALLT015.M
Quant Title :
Quant Update : Thu Oct 27 09:25:01 2005
Response via : Initial Calibration



(23) tert-Butyl Alcohol
5.551min (+0.432) 2.90PPB m
response 315808

Ion	Exp%	Act%
59.00	100	100
41.00	0.00	0.00
57.00	0.00	0.00
0.00	0.00	0.00

Sample
One/412
11/1/05
Known
0131

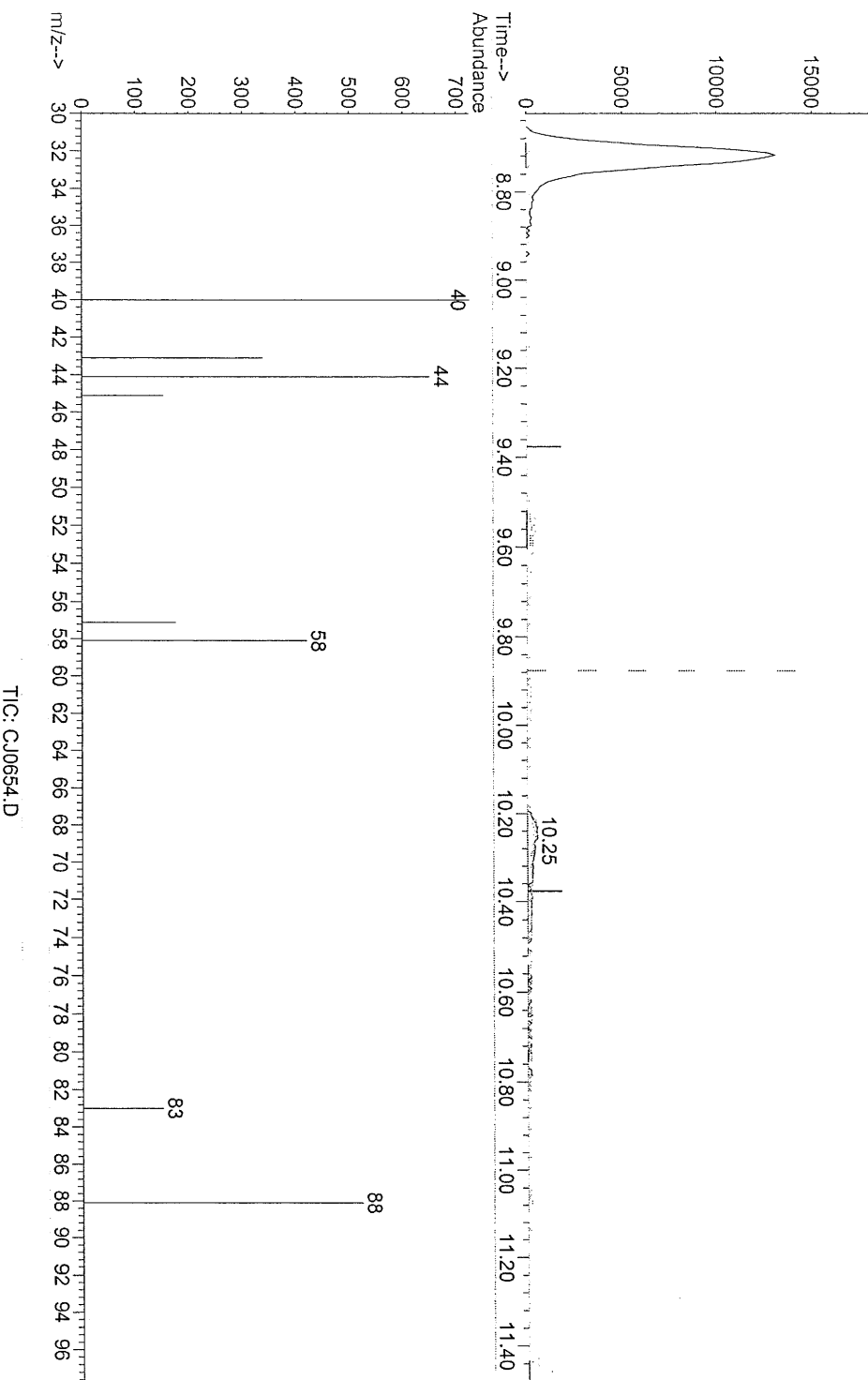
Quantitation Report (Qedit)

Data Path : C:\MSDCHEM\1\DATA\OCT27\
Data File : CJ0654.D
Acq On : 27 Oct 2005 9:13
Operator : JBS
Sample : VSTD002
Misc :
ALS Vial : 29 Sample Multiplier: 1

Quant Time: Oct 27 09:41:12 2005
Quant Method : C:\MSDCHEM\1\METHODS\ALLTO15.M
Quant Title :
Quant Update : Thu Oct 27 09:25:01 2005
Response via : Initial Calibration

Abundance
20000
15000
10000
5000
0

Ion 88.00 (87.70 to 88.70): CJ0654.D
Ion 58.00 (57.70 to 58.70): CJ0654.D
Ion 43.00 (42.70 to 43.70): CJ0654.D



(47) 1,4-Dioxane
10.257min (+0.381) 1.74PPB m
response 46700

Ion	Exp%	Act%
88.00	100	100
58.00	63.20	0.00#
43.00	2.30	0.00#
0.00	0.00	0.00

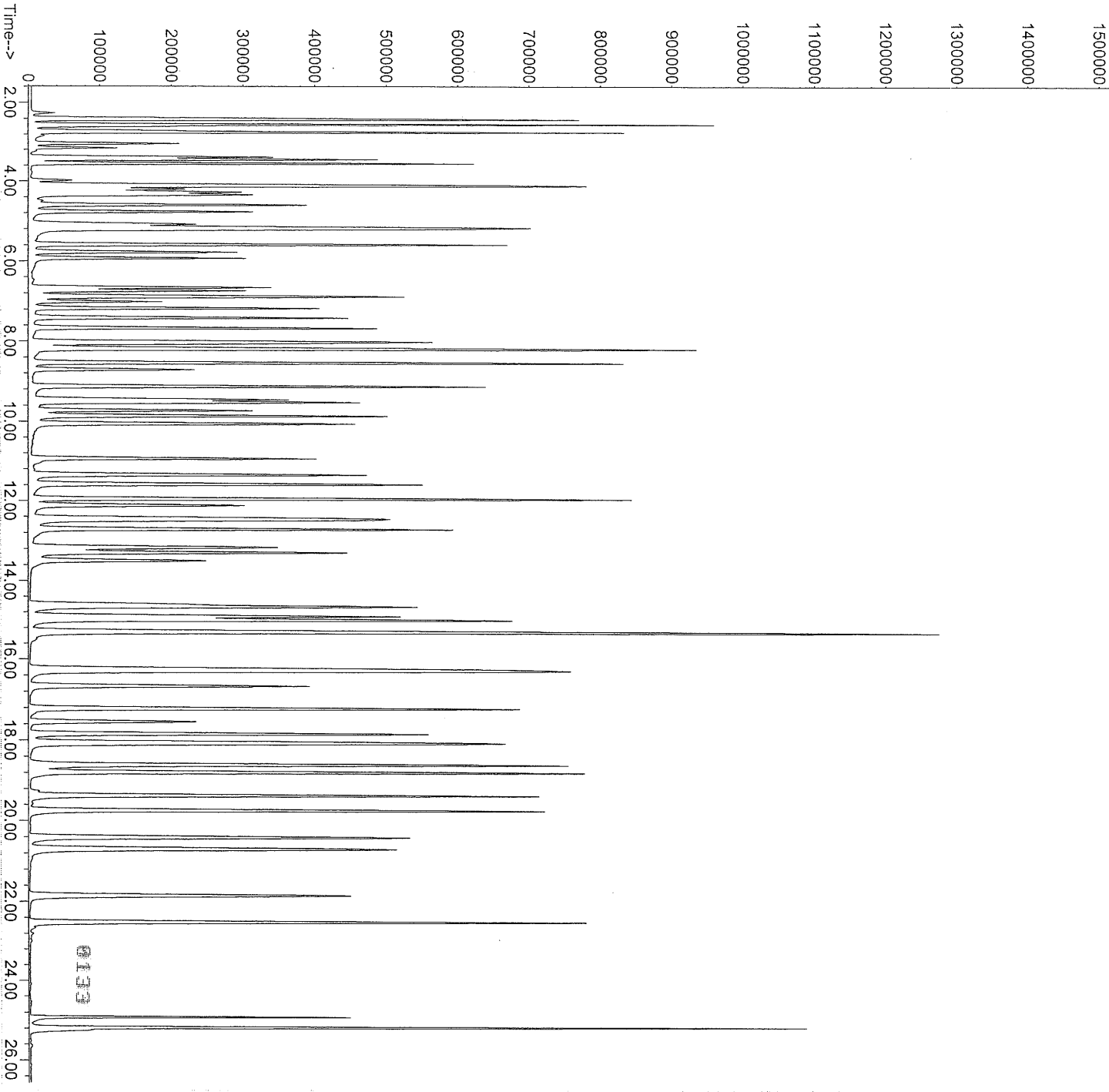
Spent sample
Am 27 2/23/05
CJ0654.D
11/1/05
0132

Data Path : C:\MSDCHEM\1\DATA\OCT27\
Data File : CJ0657.D
Acq On : 27 Oct 2005 11:06
Operator : JBS
Sample : VSTD025
Misc :
ALS Vial : 32 Sample Multiplier: 1

Quant Time: Oct 27 11:33:12 2005
Quant Method : C:\MSDCHEM\1\METHODS\ALLT015.M
Quant Title :
Quant Update : Thu Oct 27 09:25:01 2005
Response via : Initial Calibration

Abundance

TIC: CJ0657.D



Data Path : C:\MSDCHEM\1\DATA\OCT27\
 Data File : CJ0657.D
 Acq On : 27 Oct 2005 11:06
 Operator : JBS
 Sample : VSTD025
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Quant Time: Oct 27 11:33:12 2005
 Quant Method : C:\MSDCHEM\1\METHODS\ALLT015.M
 Quant Title :
 QIast Update : Thu Oct 27 09:25:01 2005
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
1) Bromochloromethane	7.02	130	773250	10.000	0.00
37) 1,4-Difluorobenzene	8.71	114	2754443	10.000	0.00
51) Chlorobenzene d5	14.59	117	2555380	10.000	0.00

Target Compounds	R.T.	QIon	Response	Conc Units	Dev(Min)	Qvalue
2) Propene	2.38	41	2732643	45.666 PPB		96
3) Dichlorodifluoromethane	2.42	85	5703620	41.415 PPB		99
4) Chlorodifluoromethane	2.43	51	4888965	43.671 PPB		95
5) Freon 114	2.54	85	6411369	40.517 PPB	#	80
6) Chloromethane	2.59	50	3672676	43.719 PPB		97
7) Vinyl Chloride	2.70	62	2591975	41.446 PPB		99
8) 1,3-Butadiene	2.74	54	4806190	44.614 PPB		97
9) Bromomethane	3.04	94	1947458	39.979 PPB		95
10) Chloroethane	3.15	64	1310849	39.931 PPB		99
11) Dichlorofluoromethane	3.38	67	4565985	40.619 PPB		98
12) Trichlorofluoromethane	3.44	101	5587750	40.143 PPB		99
13) Pentane	3.55	43	6209792	45.697 PPB		97
14) Acrolein	3.98	56	856236	51.484 PPB		87
15) 1,1-Dichloroethene	4.09	61	4152712	40.981 PPB	#	86
16) Freon 113	4.12	103	27722943	38.901 PPB		93
17) Acetone	4.20	43	4693807	38.780 PPB		99
18) Methyl Iodide	4.27	142	4959373	41.336 PPB		98
19) Carbon Disulfide	4.35	76	5927685	44.090 PPB	#	93
20) Acetonitrile	4.60	41	5267886	48.854 PPB		93
21) 3-Chloropropene	4.60	76	978710	36.133 PPB		92
22) Methylene Chloride	4.77	84	1730212	29.637 PPB		73
23) tert-Butyl Alcohol	5.08	59	5545082	46.188 PPB	#	100
24) Acrylonitrile	5.15	53	3470338	53.244 PPB		98
25) trans-1,2-Dichloroethene	5.16	61	3654712	45.756 PPB	#	76
26) Methyl t-Butyl Ether	5.20	73	6168905	42.821 PPB	#	93
27) Hexane	5.59	57	4831617	40.208 PPB		88
28) 1,1-Dichloroethane	5.78	63	4550579	37.915 PPB		99
29) Vinyl Acetate	5.92	43	7734689	64.100 PPB	#	94
30) cis-1,2-Dichloroethene	6.66	61	3344758	40.043 PPB	#	77
31) 2-Butanone	6.74	72	1069403	42.578 PPB	#	100
32) Ethyl Acetate	6.87	70	519350	43.940 PPB	#	100
33) Methyl Acrylate	6.90	55	5429483	50.011 PPB	#	94
34) Chloroform	7.17	83	4429768	40.468 PPB		96
35) 1,1,1-Trichloroethane	7.42	97	4535474	40.705 PPB	#	93
36) Carbon Tetrachloride	7.67	117	4638944	38.245 PPB		100
38) 1,2-Dichloroethane	8.06	62	3390970	51.047 PPB	#	94
39) Benzene	8.01	78	6614694	45.582 PPB	#	94
40) Isooctane	8.19	57	15648319	50.919 PPB	#	96
41) Heptane	8.53	43	7114589	55.430 PPB	#	90
42) Trichloroethene	9.12	130	3292151	43.895 PPB	#	87
43) Ethyl Acrylate	9.44	55	7366483	54.969 PPB	#	92
44) 1,2-Dichloropropane	9.52	63	2932676	45.017 PPB	#	97
45) Methyl Methacrylate	9.86	69	2352352	48.890 PPB	#	74
46) Dibromomethane	9.72	174	1955375	49.944 PPB	#	76
47) 1,4-Dioxane	9.85	88	992483	38.886 PPB	#	40
48) Bromodichloromethane	10.06	83	4867382	50.363 PPB	#	92

01134

Data Path : C:\MSDCHEM\1\DATA\OCT27\
 Data File : CJ0657.D
 Acq On : 27 Oct 2005 11:06
 Operator : JBS
 Sample : VSTD025
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

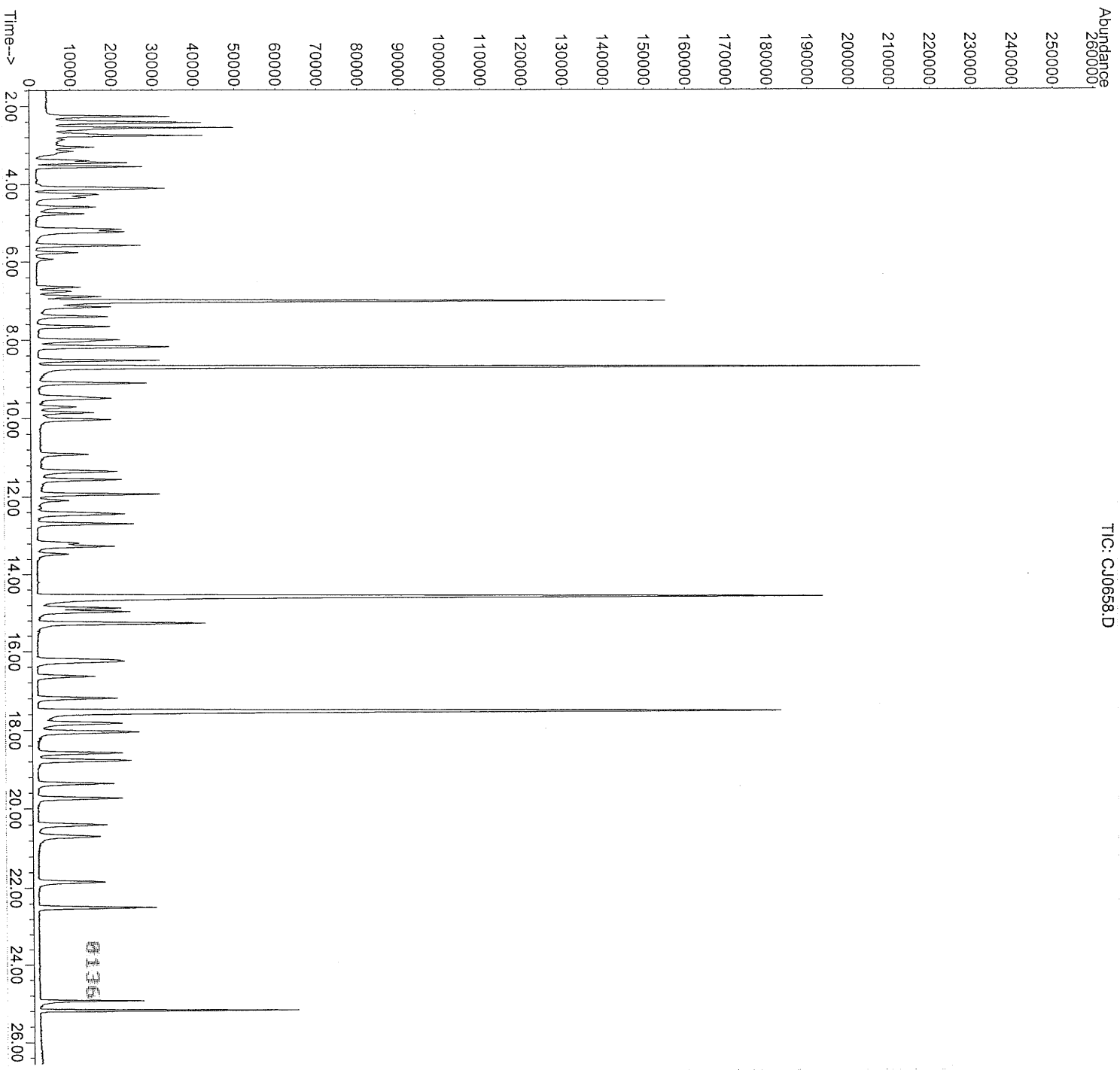
Quant Time: Oct 27 11:33:12 2005
 Quant Method : C:\MSDCHEM\1\METHODS\ALLT015.M
 Quant Title :
 QLast Update : Thu Oct 27 09:25:01 2005
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev	(Min)
49) cis-1,3-Dichloropropene	10.93	75	3597573	43.625	PPB	#	70
50) 4-Methyl-2-Pentanone	11.34	43	8302189	46.592	PPB	#	95
52) Toluene	11.58	91	7196153	38.399	PPB	#	99
53) Octane	11.95	43	9048032	43.509	PPB	#	89
54) trans-1,3-Dichloropropene	12.11	75	2907823	42.244	PPB	#	88
55) Ethyl Methacrylate	12.43	69	3698882	39.294	PPB	#	77
56) 1,1,2-Trichloroethane	12.48	97	2319792	35.583	PPB	#	86
57) Tetrachloroethene	12.71	166	3067347	38.357	PPB	#	97
58) 2-Hexanone	13.16	43	7570856	40.313	PPB	#	76
59) Dibromochloromethane	13.29	127	3810758	42.763	PPB	#	98
60) 1,2-Dibromoethane	13.49	107	3933018	44.155	PPB	#	98
61) Chlorobenzene	14.66	112	6096907	42.105	PPB	#	86
62) 1,1,1,2-Tetrachloroethane	14.90	131	3446450	39.633	PPB	#	99
63) Ethylbenzene	14.99	91	10761082	43.823	PPB	#	91
64) m/p-Xylene	15.29	91	17117942	86.396	PPB	#	90
65) o-Xylene	16.24	91	8595783	45.295	PPB	#	90
66) Styrene	16.29	104	6463519	46.971	PPB	#	100
67) Bromoform	16.66	173	4250283	48.974	PPB	#	97
68) Cumene	17.21	105	10911193	51.302	PPB	#	93
69) 1,1,2,2-Tetrachloroethane	18.04	83	4546195	49.352	PPB	#	97
70) 1,2,3-Trichloropropane	18.08	110	1583361	49.458	PPB	#	65
71) Bromobenzene	17.84	156	3458319	52.375	PPB	#	98
72) 4-Ethyltoluene	18.60	105	11679205	48.936	PPB	#	93
73) 1,3,5-Trimethylbenzene	18.79	105	10140606	49.907	PPB	#	91
74) Alpha Methyl Styrene	19.37	118	5007483	52.803	PPB	#	95
75) 1,2,4-Trimethylbenzene	19.74	105	9749454	46.939	PPB	#	90
76) 1,3-Dichlorobenzene	20.41	146	5888971	50.971	PPB	#	97
77) 1,4-Dichlorobenzene	20.69	146	5686485	50.287	PPB	#	97
78) Benzyl Chloride	16.24	91	8595783	45.295	PPB	#	100
79) 1,2-Dichlorobenzene	21.85	146	5036112	47.095	PPB	#	96
80) Hexachloroethane	22.51	117	3918641	50.696	PPB	#	96
81) 1,2,4-Trichlorobenzene	24.92	180	2056813	44.331	PPB	#	99
82) Hexachlorobutadiene	25.19	225	2400911	36.156	PPB	#	77

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : C:\MSDCHEM\1\DATA\OCT27\
Data File : C:\MSDCHEM\1\DATA\OCT27\
Acq On : 27 Oct 2005 11:47
Operator : JBS
Sample : VSTD001
Misc :
ALS Vial : 33 Sample Multiplier: 1

Quant Time: Oct 27 12:51:29 2005
Quant Method : C:\MSDCHEM\1\METHODS\ALLT015.M
Quant Title :
Quant Update : Thu Oct 27 09:25:01 2005
Response via : Initial Calibration



Data Path : C:\MSDCHEM\1\DATA\OCT27\
 Data File : CJ0658.D
 Acq On : 27 Oct 2005 11:47
 Operator : JBS
 Sample : VSTD001
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Quant Time: Oct 27 12:51:29 2005
 Quant Method : C:\MSDCHEM\1\METHODS\ALLT015.M
 Quant Title :
 Qlast Update : Thu Oct 27 09:25:01 2005
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
1) Bromochloromethane	7.01	130	726821	10.000	0.00
37) 1,4-Difluorobenzene	8.69	114	2643010	10.000	-0.01
51) Chlorobenzene d5	14.57	117	2324876	10.000	-0.01

Target Compounds	R.T.	QIon	Response	Conc Units	Dev(Min)	Qvalue
2) Propene	2.39	41	129506	2.302 PPB		95
3) Dichlorodifluoromethane	2.42	85	277385	2.143 PPB		99
4) Chlorodifluoromethane	2.43	51	229731	2.183 PPB		94
5) Freon 114	2.54	85	312722	2.103 PPB		79
6) Chloromethane	2.59	50	175775	2.226 PPB		97
7) Vinyl Chloride	2.70	62	122482	2.084 PPB		98
8) 1,3-Butadiene	2.75	54	206370	2.038 PPB		98
9) Bromomethane	3.04	94	89558	1.956 PPB		97
10) Chloroethane	3.15	64	54954	1.781 PPB		43
11) Dichlorofluoromethane	3.39	67	180262	1.706 PPB		88
12) Trichlorofluoromethane	3.44	101	258448	1.975 PPB		98
13) Pentane	3.55	43	260268	2.038 PPB		97
14) Acrolein	4.01	56	22774m	1.457 PPB		
15) 1,1-Dichloroethene	4.09	61	172331	1.809 PPB		87
16) Freon 113	4.11	103	109443	1.633 PPB		95
17) Acetone	4.24	43	167719	1.474 PPB		98
18) Methyl Iodide	4.27	142	208497	1.849 PPB		99
19) Carbon Disulfide	4.34	76	265270	2.099 PPB		95
20) Acetonitrile	4.59	41	167873	1.656 PPB		82
21) 3-Chloropropene	4.59	76	44504m	1.748 PPB		
22) Methylene Chloride	4.77	84	66832	1.218 PPB		66
23) tert-Butyl Alcohol	5.15	59	218450	1.936 PPB		100
24) Acrylonitrile	5.17	53	107301	1.751 PPB		65
25) trans-1,2-Dichloroethene	5.16	61	130196	1.734 PPB		75
26) Methyl t-Butyl Ether	5.23	73	235492	1.739 PPB		58
27) Hexane	5.58	57	183259	1.622 PPB		91
28) 1,1-Dichloroethane	5.77	63	175377	1.555 PPB		99
29) Vinyl Acetate	5.93	43	130229	1.148 PPB		99
30) cis-1,2-Dichloroethene	6.65	61	120140	1.530 PPB		74
31) 2-Butanone	6.76	72	35877	1.520 PPB		100
32) Ethyl Acetate	6.87	70	14253m	1.283 PPB		
33) Methyl Acrylate	6.90	55	167005	1.637 PPB		69
34) Chloroform	7.16	83	167576	1.629 PPB		77
35) 1,1,1-Trichloroethane	7.40	97	169989	1.623 PPB		91
36) Carbon Tetrachloride	7.66	117	172436	1.512 PPB		99
38) 1,2-Dichloroethane	8.04	62	119364	1.873 PPB		99
39) Benzene	7.99	78	237783	1.708 PPB		95
40) Isooctane	8.17	57	557464	1.890 PPB		97
41) Heptane	8.52	43	264694	2.149 PPB		89
42) Trichloroethene	9.10	130	147523	2.050 PPB		88
43) Ethyl Acrylate	9.44	55	208265	1.620 PPB		77
44) 1,2-Dichloropropane	9.50	63	108350	1.733 PPB		91
45) Methyl Methacrylate	9.85	69	70733	1.532 PPB		67
46) Dibromomethane	9.71	174	65883	1.754 PPB		71
47) 1,4-Dioxane	9.98	88	41364m	1.689 PPB		
48) Bromodichloromethane	10.03	83	177705	1.916 PPB		87

0137

Data Path : C:\MSDCHEM\1\DATA\OCT27\
 Data File : C:\MSDCHEM\1\DATA\OCT27\
 Acq On : 27 Oct 2005 11:47
 Operator : JBS
 Sample : VSTD001
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Quant Time: Oct 27 12:51:29 2005
 Quant Method : C:\MSDCHEM\1\METHODS\ALLT015.M
 Quant Title :
 QLast Update : Thu Oct 27 09:25:01 2005
 Response via : Initial Calibration

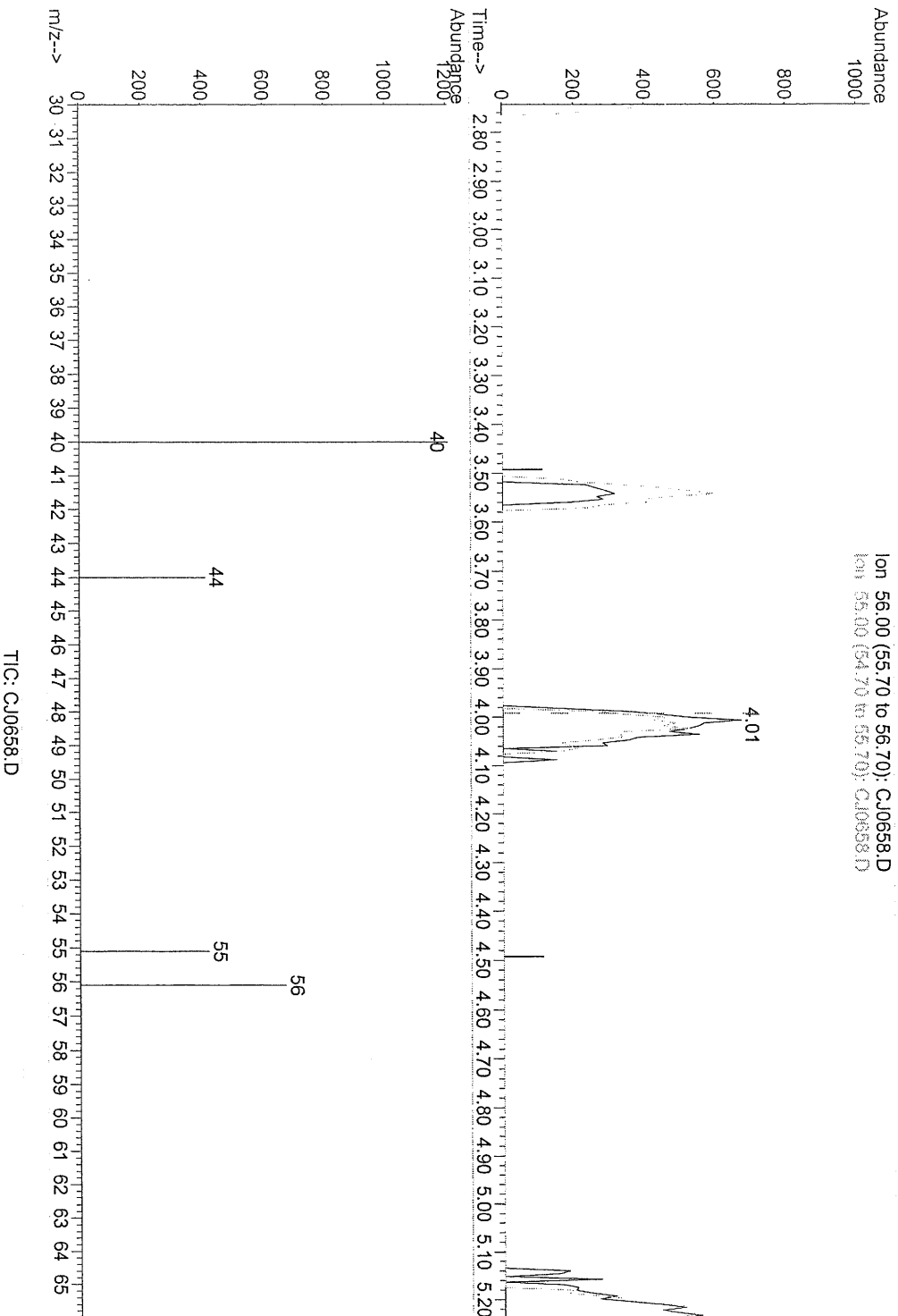
Internal Standards	R.T.	QIon	Response	Conc Units	Dev (Min)
49) cis-1,3-Dichloropropene	10.92	75	116363	1.471 PPB	# 68
50) 4-Methyl-2-Pentanone	11.34	43	356479	2.085 PPB	# 90
52) Toluene	11.56	91	242352	1.421 PPB	# 95
53) Octane	11.93	43	309331	1.635 PPB	# 90
54) trans-1,3-Dichloropropene	12.10	75	85223	1.361 PPB	# 70
55) Ethyl Methacrylate	12.42	69	120044	1.402 PPB	# 86
56) 1,1,2-Trichloroethane	12.45	97	100273	1.691 PPB	# 88
57) Tetrachloroethene	12.69	166	131492	1.807 PPB	# 99
58) 2-Hexanone	13.18	43	316302	1.851 PPB	# 78
59) Dibromochloromethane	13.27	127	138161	1.704 PPB	# 98
60) 1,2-Dibromoethane	13.47	107	135031	1.666 PPB	# 80
61) Chlorobenzene	14.64	112	227726	1.729 PPB	# 75
62) 1,1,1,2-Tetrachloroethane	14.87	131	127627	1.613 PPB	# 51
63) Ethylbenzene	14.97	91	348677	1.561 PPB	# 89
64) m/p-Xylene	15.27	91	550700	3.055 PPB	# 90
65) o-Xylene	16.21	91	269471	1.561 PPB	# 90
66) Styrene	16.26	104	187934	1.501 PPB	# 100
67) Bromoform	16.63	173	160667	2.035 PPB	# 95
68) Cumene	17.19	105	310608	1.605 PPB	# 93
69) 1,1,2,2-Tetrachloroethane	18.02	83	132661	1.583 PPB	# 94
70) 1,2,3-Trichloropropane	18.05	110	59258	2.035 PPB	# 48
71) Bromobenzene	17.82	156	125029	2.081 PPB	# 99
72) 4-Ethyltoluene	18.58	105	339440	1.563 PPB	# 93
73) 1,3,5-Trimethylbenzene	18.77	105	343252	1.857 PPB	# 89
74) Alpha Methyl Styrene	19.36	118	147905	1.714 PPB	# 95
75) 1,2,4-Trimethylbenzene	19.73	105	311252	1.647 PPB	# 84
76) 1,3-Dichlorobenzene	20.40	146	201187	1.914 PPB	# 89
77) 1,4-Dichlorobenzene	20.68	146	189507	1.842 PPB	# 82
78) Benzyl Chloride	16.21	91	269471	1.561 PPB	# 100
79) 1,2-Dichlorobenzene	21.84	146	219460	2.256 PPB	# 95
80) Hexachloroethane	22.50	117	141378	2.010 PPB	# 96
81) 1,2,4-Trichlorobenzene	24.92	180	145433	3.445 PPB	# 97
82) Hexachlorobutadiene	25.18	225	159312	2.637 PPB	# 98

(#) = qualifier out of range (m) = manual integration (+) = signals summed

0133

Data Path : C:\MSDCHEM\1\DATA\OCT27\
 Data File : CJ0658.D
 Acq On : 27 Oct 2005 11:47
 Operator : JBS
 Sample : VSTD001
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Quant Time: Oct 27 12:51:29 2005
 Quant Method : C:\MSDCHEM\1\METHODS\ALLT015.M
 Quant Title :
 Quant Update : Thu Oct 27 09:25:01 2005
 Response via : Initial Calibration



Ion 56.00 (55.70 to 56.70): CJ0658.D
 Ion 55.00 (54.70 to 55.70): CJ0658.D

(14) Acrolein

4.006min (+0.013) 1.46PPB m

response 22774

Ion	Exp%	Act%
56.00	100	100
55.00	62.80	0.00#
0.00	0.00	0.00
0.00	0.00	0.00

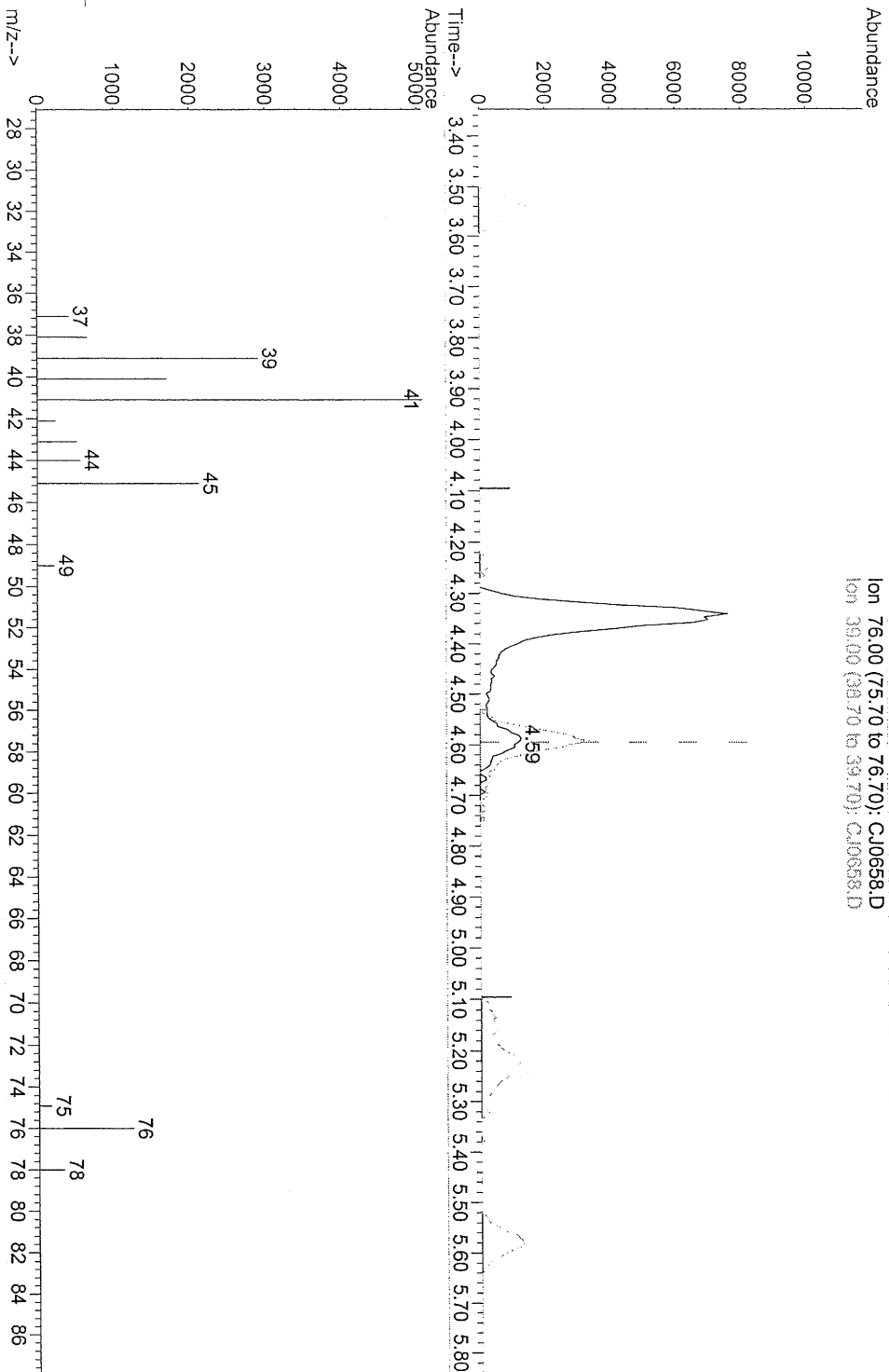
*Crack/HLZ
11/1/05*

*Synt peak
Narrow
1/12/05*

0139

Data Path : C:\MSDCHEM\1\DATA\OCT27\
Data File : C:\MSDCHEM\1\DATA\OCT27\
Acq On : 27 Oct 2005 11:47
Operator : JBS
Sample : VSTD001
Misc :
ALS Vial : 33 Sample Multiplier: 1

Quant Time: Oct 27 12:51:29 2005
Quant Method : C:\MSDCHEM\1\METHODS\ALLT015.M
Quant Title :
Quant Update : Thu Oct 27 09:25:01 2005
Response via : Initial Calibration



Ion 76.00 (75.70 to 76.70): CJ0658.D
Ion 39.00 (38.70 to 39.70): CJ0658.D

TIC: C:\MSDCHEM\1\DATA\OCT27\CJ0658.D

(21) 3-Chloropropene

response	44504	
4.587min (-0.008) 1.75PPB m		
Ion	Exp%	Act%
76.00	100	100
39.00	289.60	0.00#
0.00	0.00	0.00
0.00	0.00	0.00

AMR/ylz
11/1/05

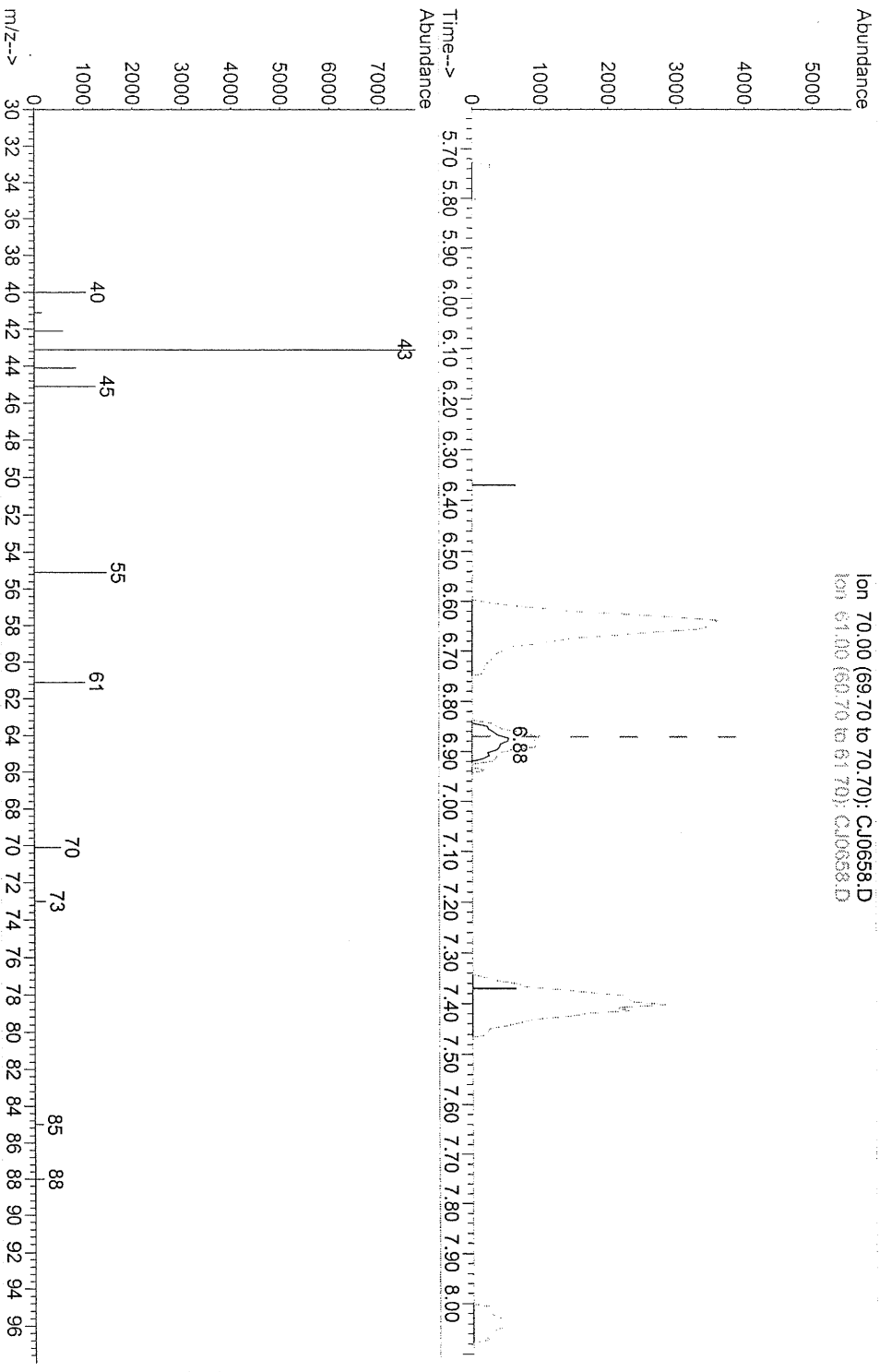
3-Chloropropene
AMR/ylz
11/1/05



Quantitation Report (Qedit)

Data Path : C:\MSDCHEM\1\DATA\OCT27\
Data File : CJ0658.D
Acq On : 27 Oct 2005 11:47
Operator : JBS
Sample : VSTD001
Misc :
ALS Vial : 33 Sample Multiplier: 1

Quant Time: Oct 27 12:51:29 2005
Quant Method : C:\MSDCHEM\1\METHODS\ALLT015.M
Quant Title :
Quant Update : Thu Oct 27 09:25:01 2005
Response via : Initial Calibration



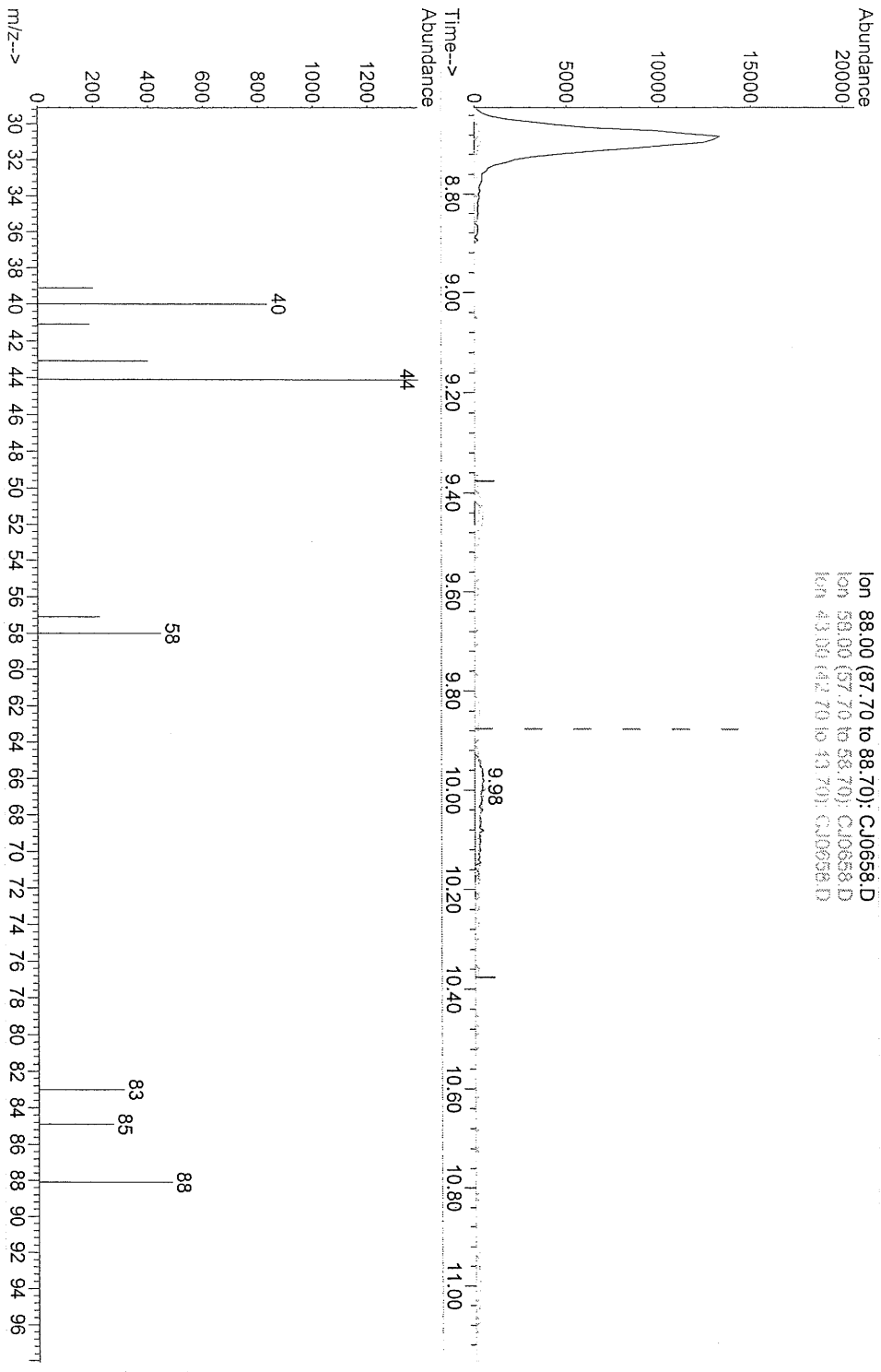
(32) Ethyl Acetate
6.873min (+0.003) 1.28PPB m
response 14253

Ion	Exp%	Act%
70.00	100	100
61.00	0.00	0.00
0.00	0.00	0.00
0.00	0.00	0.00

Omni/912
11/1/05
Sample
11/1/05

0141

Data Path : C:\MSDCHEM\1\DATA\OCT27\
 Data File : C:\MSDCHEM\1\DATA\OCT27\
 Acq On : 27 Oct 2005 11:47
 Operator : JBS
 Sample : VSTD001
 Misc :
 ALS Vial : 33 Sample Multiplier: 1
 Quant Time: Oct 27 12:51:29 2005
 Quant Method : C:\MSDCHEM\1\METHODS\ALLT015.M
 Quant Title :
 Quant Update : Thu Oct 27 09:25:01 2005
 Response via : Initial Calibration



Ion 88.00 (87.70 to 88.70): C:\MSDCHEM\1\DATA\OCT27\CJ10658.D
 Ion 58.00 (57.70 to 58.70): C:\MSDCHEM\1\DATA\OCT27\CJ10658.D
 Ion 43.00 (42.70 to 43.70): C:\MSDCHEM\1\DATA\OCT27\CJ10658.D

(47) 1,4-Dioxane
 9.981min (+0.105) 1.69PPB m
 response 41364

Ion	Exp%	Act%
88.00	100	100
58.00	63.20	0.00#
43.00	2.30	0.00#
0.00	0.00	0.00

*CONF/MLD
11/1/05*

*CONF/MLD
11/1/05*

0142

FORM7
LANCASTER LABORATORIES
VOLATILE ORGANICS IN AIR

CONTINUING CALIBRATION DATA SHEET

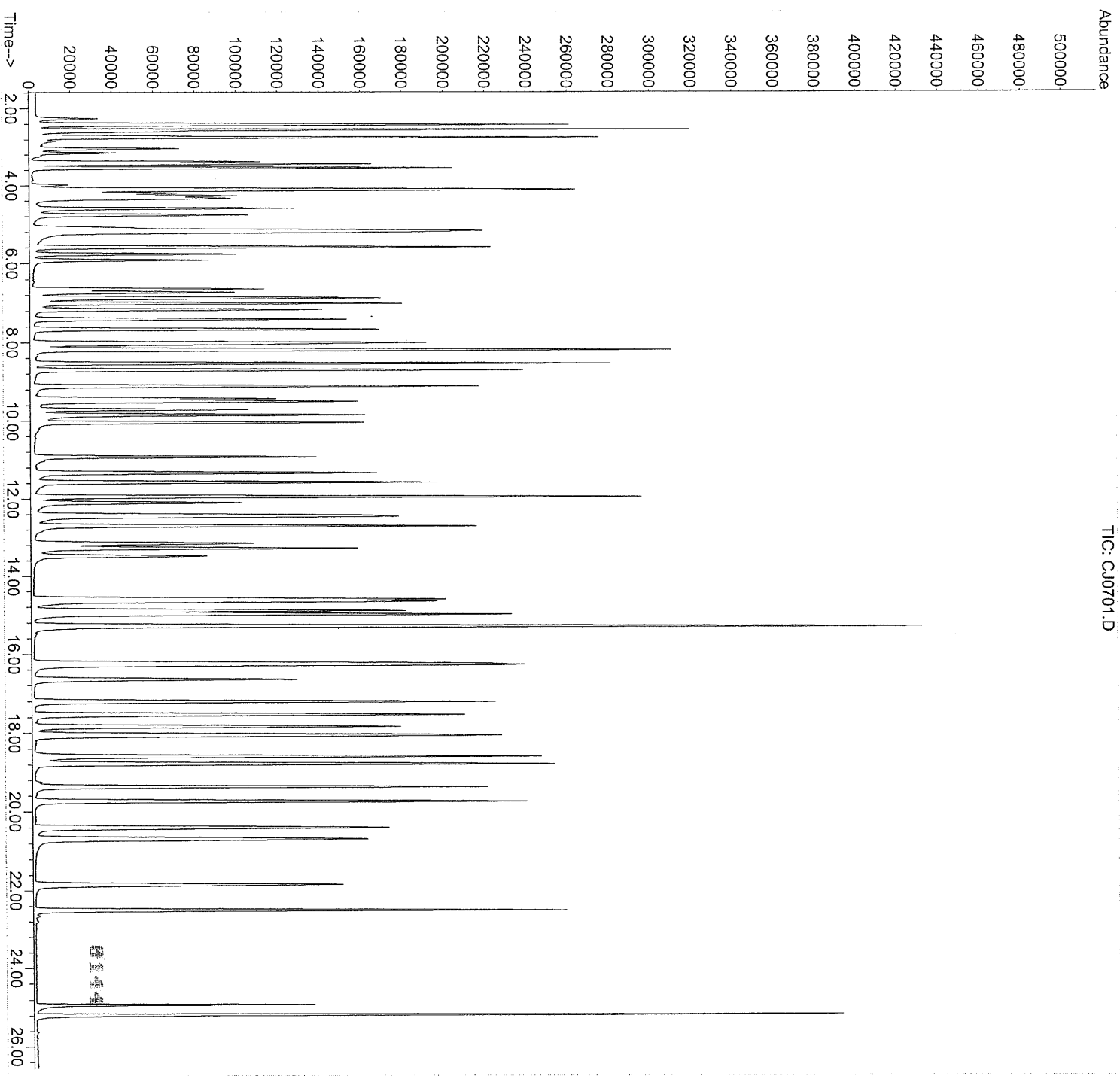
Instrument ID: HP09464
EPA SAMPLE NO.: VSTD010

Data File: C:\MSDCHEM\1\DATA\OCT28\CJ0701.D
DATE INJECTED: 10/28/05 TIME INJECTED: 08:05

COMPOUND NAME	ICAL MEAN RRF	CCAL RRF	%D
Dichlorodifluoromethane	2.928	2.424	-20
Freon 114	3.324	2.694	-20
Chloromethane	1.882	1.547	-20
Vinyl Chloride	1.320	1.050	-20
Bromomethane	1.001	0.820	-20
Chloroethane	0.655	0.557	-10
Trichlorofluoromethane	2.869	2.414	-20
1,1-Dichloroethene	2.058	1.724	-20
Freon 113	1.378	1.196	-10
3-Chloropropene	0.519	0.399	-20
Methylene Chloride	0.840	0.743	-10
1,1-Dichloroethane	2.326	2.044	-10
cis-1,2-Dichloroethene	1.637	1.441	-10
Chloroform	2.131	1.897	-10
1,1,1-Trichloroethane	2.186	1.954	-10
Carbon tetrachloride	2.376	2.121	-10
1,2-Dichloroethane	0.402	0.402	0
Benzene	0.817	0.812	-1
Trichloroethene	0.447	0.403	-10
1,2-Dichloropropane	0.366	0.355	-3
cis-1,3-dichloropropene	0.468	0.469	0
Toluene	1.074	1.050	-2
trans-1,3-Dichloropropene	0.417	0.450	8
1,1,2-Trichloroethane	0.394	0.377	-4
Tetrachloroethene	0.494	0.459	-7
1,2-Dibromoethane	0.543	0.551	1
Chlorobenzene	0.886	0.875	-1
Ethylbenzene	1.503	1.486	-1
m/p-Xylene	1.212	1.222	1
o-Xylene	1.176	1.175	0
Styrene	0.852	0.876	3
1,1,2,2-Tetrachloroethane	0.546	0.623	14
4-Ethyltoluene	1.588	1.642	3
1,3,5-Trimethylbenzene	1.362	1.390	2
1,2,4-Trimethylbenzene	1.347	1.415	5
1,3-Dichlorobenzene	0.771	0.799	4
1,4-Dichlorobenzene	0.738	0.759	3
1,2-Dichlorobenzene	0.736	0.698	-5
1,2,4-Trichlorobenzene	0.250	0.291	16

Data Path : C:\MSDCHEM\1\DATA\OCT28\
Data File : C10701.D
Acq On : 28 Oct 2005 8:05
Operator : JBS
Sample : VSTD010
Misc :
ALS Vial : 27 Sample Multiplier: 1

Quant Time: Oct 28 08:32:42 2005
Quant Method : C:\MSDCHEM\1\METHODS\ALLT015.M
Quant Title :
Quant Update : Thu Oct 27 13:03:08 2005
Response via : Initial Calibration



Data Path : C:\MSDCHEM\1\DATA\OCT28\
 Data File : CJ0701.D
 Acq On : 28 Oct 2005 8:05
 Operator : JBS
 Sample : VSTD010
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Quant Time : Oct 28 08:32:42 2005
 Quant Method : C:\MSDCHEM\1\METHODS\ALLT015.M
 Quant Title :
 Quant Update : Thu Oct 27 13:03:08 2005
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
1) Bromochloromethane	7.01	130	784315	10.000	0.00
37) 1,4-Difluorobenzene	8.70	114	2732019	10.000	0.00
51) Chlorobenzene d5	14.58	117	2282851	10.000	0.00

Target Compounds	R.T.	QIon	Response	Conc Units	Dev(Min)	Qvalue
2) Propene	2.39	41	869036	8.038	PPB	97
3) Dichlorodifluoromethane	2.42	85	1910824	8.322	PPB	99
4) Chlorodifluoromethane	2.43	51	1610847	8.290	PPB	96
5) Freon 114	2.54	85	2112858	8.103	PPB	#
6) Chloromethane	2.59	50	1213238	8.222	PPB	96
7) Vinyl Chloride	2.70	62	823696	7.955	PPB	97
8) 1,3-Butadiene	2.74	54	1553694	8.382	PPB	97
9) Bromomethane	3.04	94	643381	8.194	PPB	94
10) Chloroethane	3.15	64	436645	8.501	PPB	97
11) Dichlorofluoromethane	3.39	67	1501115	8.543	PPB	98
12) Trichlorofluoromethane	3.44	101	1893109	8.414	PPB	98
13) Pentane	3.55	43	1991243	8.677	PPB	98
14) Acrolein	3.98	56	252538	10.233	PPB	88
15) 1,1-Dichloroethene	4.09	61	1352364	8.377	PPB	86
16) Freon 113	4.12	103	938018	8.680	PPB	94
17) Acetone	4.20	43	1531277	9.677	PPB	100
18) Methyl Iodide	4.27	142	1642324	8.812	PPB	98
19) Carbon Disulfide	4.35	76	1919281	9.010	PPB	93
20) Acetonitrile	4.59	41	1648196	10.214	PPB	94
21) 3-Chloropropene	4.59	76	312836	7.680	PPB	94
22) Methylene Chloride	4.76	84	582376	8.844	PPB	74
23) tert-Butyl Alcohol	5.10	59	1706718	9.402	PPB	100
24) Acrylonitrile	5.15	53	1026938	9.810	PPB	97
25) trans-1,2-Dichloroethene	5.16	61	1158970	9.362	PPB	77
26) Methyl t-Butyl Ether	5.20	73	1966121	8.952	PPB	89
27) Hexane	5.58	57	1583166	8.996	PPB	88
28) 1,1-Dichloroethane	5.77	63	1523308	8.350	PPB	99
29) Vinyl Acetate	5.92	43	2126519	11.453	PPB	#
30) cis-1,2-Dichloroethene	6.65	61	1073404	8.359	PPB	78
31) 2-Butanone	6.73	72	340138	9.800	PPB	100
32) Ethyl Acetate	6.86	70	159028	10.210	PPB	#
33) Methyl Acrylate	6.89	55	1643237	9.910	PPB	94
34) Chloroform	7.16	83	1487954	8.904	PPB	96
35) 1,1,1-Trichloroethane	7.41	97	1532804	8.941	PPB	93
36) Carbon Tetrachloride	7.66	117	1580260	8.481	PPB	99
38) 1,2-Dichloroethane	8.00	62	1098425	9.995	PPB	99
39) Benzene	8.00	78	2217172	9.932	PPB	95
40) Isooctane	8.18	57	5072792	10.265	PPB	97
41) Heptane	8.53	43	2264175	10.264	PPB	92
42) Trichloroethene	9.11	130	1099758	8.999	PPB	88
43) Ethyl Acrylate	9.43	55	2268063	11.448	PPB	88
44) 1,2-Dichloropropane	9.50	63	968781	9.681	PPB	89
45) Methyl Methacrylate	9.85	69	746583	11.044	PPB	96
46) Dibromomethane	9.71	174	638108	10.487	PPB	77
47) 1,4-Dioxane	9.87	88	311123	10.423	PPB	76
48) Bromodichloromethane	10.04	83	1628626	10.592	PPB	38

0145

Data Path : C:\MSDCHEM\1\DATA\OCT28\
 Data File : CJ0701.D
 Acq On : 28 Oct 2005 8:05
 Operator : JBS
 Sample : VSTD010
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Quant Time : Oct 28 08:32:42 2005
 Quant Method : C:\MSDCHEM\1\METHODS\ALLT015.M
 Quant Title :
 Quant Update : Thu Oct 27 13:03:08 2005
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev	(Min)
49) cis-1,3-Dichloropropene	10.92	75	1153188	9.028	PPB	#	71
50) 4-Methyl-2-Pentanone	11.33	43	2755225	10.422	PPB	#	94
52) Toluene	11.57	91	2398074	9.784	PPB	#	98
53) Octane	11.94	43	3043789	10.414	PPB	#	88
54) trans-1,3-Dichloropropene	12.10	75	924589	9.703	PPB	#	87
55) Ethyl Methacrylate	12.41	69	1155578	9.940	PPB	#	76
56) 1,1,2-Trichloroethane	12.46	97	818589	9.107	PPB	#	87
57) Tetrachloroethene	12.70	166	1047121	9.294	PPB	#	97
58) 2-Hexanone	13.15	43	2256090	10.266	PPB	#	77
59) Dibromochloromethane	13.28	127	1272951	10.069	PPB	#	99
60) 1,2-Dibromoethane	13.48	107	1258205	10.141	PPB	#	98
61) Chlorobenzene	14.64	112	1996614	9.872	PPB	#	85
62) 1,1,1,2-Tetrachloroethane	14.88	131	1138353	9.267	PPB	#	98
63) Ethylbenzene	14.97	91	3391644	9.885	PPB	#	92
64) m/p-Xylene	15.27	91	5440256	19.656	PPB	#	90
65) o-Xylene	16.22	91	2682776	9.996	PPB	#	90
66) Styrene	16.27	104	1999016	10.283	PPB	#	100
67) Bromoform	16.65	173	1358466	10.413	PPB	#	97
68) Cumene	17.20	105	3321402	10.683	PPB	#	93
69) 1,1,2,2-Tetrachloroethane	18.03	83	1492168	11.977	PPB	#	96
70) 1,2,3-Trichloropropane	18.06	110	501028	10.386	PPB	#	64
71) Bromobenzene	17.83	156	1071707	10.567	PPB	#	99
72) 4-Ethyltoluene	18.59	105	3561208	9.825	PPB	#	94
73) 1,3,5-Trimethylbenzene	18.77	105	3173295	10.204	PPB	#	91
74) Alpha Methyl Styrene	19.36	118	1519857	10.933	PPB	#	95
75) 1,2,4-Trimethylbenzene	19.73	105	3068856	9.982	PPB	#	91
76) 1,3-Dichlorobenzene	20.40	146	1823098	10.357	PPB	#	97
77) 1,4-Dichlorobenzene	20.69	146	1732848	10.281	PPB	#	97
78) Benzyl Chloride	16.22	91	2682776	9.996	PPB	#	100
79) 1,2-Dichlorobenzene	21.84	146	1593743	9.485	PPB	#	96
80) Hexachloroethane	22.50	117	1225547	10.677	PPB	#	97
81) 1,2,4-Trichlorobenzene	24.92	180	664223	8.957	PPB	#	99
82) Hexachlorobutadiene	25.18	225	952028	9.365	PPB	#	97

(#) = qualifier out of range (m) = manual integration (+) = signals summed

0146

FORM7
LANCASTER LABORATORIES
VOLATILE ORGANICS IN AIR

CONTINUING CALIBRATION DATA SHEET

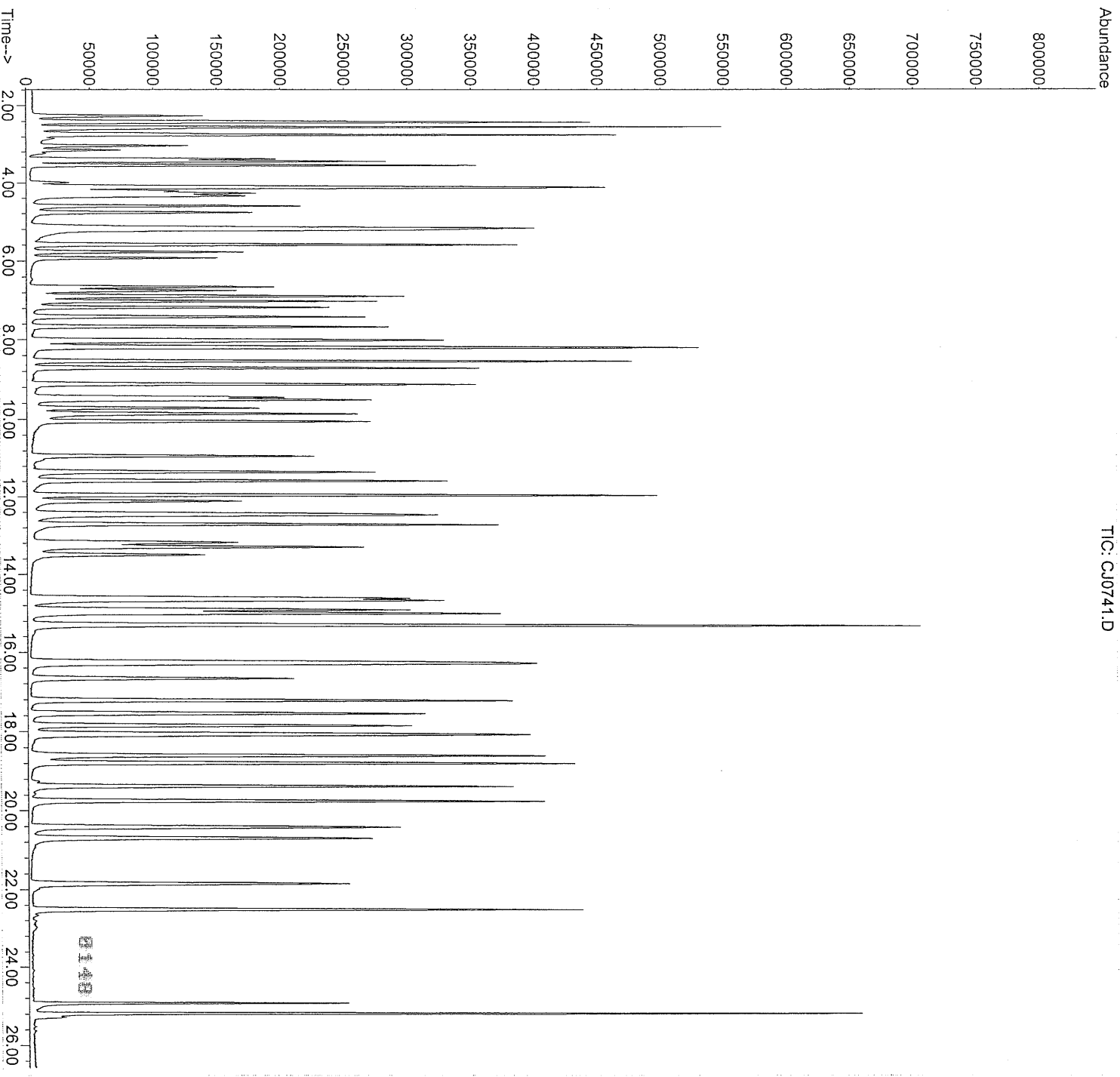
Instrument ID: HP09464
EPA SAMPLE NO.: VSTD010

Data File: C:\MSDCHEM\1\DATA\OCT29\CJ0741.D
DATE INJECTED: 10/29/05 TIME INJECTED: 12:47

COMPOUND NAME	ICAL MEAN RRF	CCAL RRF	%D
Dichlorodifluoromethane	2.928	2.610	-10
Freon 114	3.324	2.941	-10
Chloromethane	1.882	1.650	-10
Vinyl Chloride	1.320	1.177	-10
Bromomethane	1.001	0.899	-10
Chloroethane	0.655	0.603	-8
Trichlorofluoromethane	2.869	2.580	-10
1,1-Dichloroethene	2.058	1.893	-8
Freon 113	1.378	1.331	-3
3-Chloropropene	0.519	0.465	-10
Methylene Chloride	0.840	0.813	-3
1,1-Dichloroethane	2.326	2.252	-3
cis-1,2-Dichloroethene	1.637	1.643	0
Chloroform	2.131	2.076	-3
1,1,1-Trichloroethane	2.186	2.121	-3
Carbon Tetrachloride	2.376	2.308	-3
1,2-Dichloroethane	0.402	0.435	8
Benzene	0.817	0.909	11
Trichloroethene	0.447	0.431	-4
1,2-Dichloropropane	0.366	0.401	10
cis-1,3-Dichloropropene	0.468	0.538	15
Toluene	1.074	1.210	13
trans-1,3-Dichloropropene	0.417	0.494	18
1,1,2-Trichloroethane	0.394	0.420	7
Tetrachloroethane	0.494	0.541	10
1,2-Dibromoethane	0.543	0.606	11
Chlorobenzene	0.886	0.964	9
Ethylbenzene	1.503	1.656	10
m/p-Xylene	1.212	1.338	10
o-Xylene	1.176	1.298	10
Styrene	0.852	0.961	13
1,1,2,2-Tetrachloroethane	0.546	0.750	38
4-Ethyltoluene	1.588	1.813	14
1,3,5-Trimethylbenzene	1.362	1.514	11
1,2,4-Trimethylbenzene	1.347	1.528	13
1,3-Dichlorobenzene	0.771	0.876	14
1,4-Dichlorobenzene	0.738	0.837	13
1,2-Dichlorobenzene	0.736	0.765	4
1,2,4-Trichlorobenzene	0.250	0.328	31

Data Path : C:\MSDCHEM\1\DATA\OCT29\
Data File : CJ0741.D
Acq On : 29 Oct 2005 12:47
Operator : JBS
Sample : VSTD010
Misc :
ALS Vial : 27 Sample Multiplier: 1

Quant Time: Oct 29 13:13:52 2005
Quant Method : C:\MSDCHEM\1\METHODS\ALLTO15.M
Quant Title :
Quant Update : Fri Oct 28 08:49:53 2005
Response via : Initial Calibration



Data Path : C:\MSDCHEM\1\DATA\OCT29\
 Data File : CJ0741.D
 Acq On : 29 Oct 2005 12:47
 Operator : JBS
 Sample : VSTD010
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Quant Time: Oct 29 13:13:52 2005
 Quant Method : C:\MSDCHEM\1\METHODS\ALLT015.M
 Quant Title :
 Qlast Update : Fri Oct 28 08:49:53 2005
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	7.01	130	1240511	10.000	PPB	0.00
37) 1,4-Difluorobenzene	8.70	114	4279276	10.000	PPB	0.00
51) Chlorobenzene d5	14.59	117	3559350	10.000	PPB	0.00

Target Compounds

2) Propene	2.39	41	1499754	8.770	PPB	96
3) Dichlorodifluoromethane	2.42	85	3253540	8.959	PPB	98
4) Chlorodifluoromethane	2.43	51	2750655	8.950	PPB	96
5) Freon 114	2.54	85	3648650	8.847	PPB	80
6) Chloromethane	2.59	50	2046442	8.768	PPB	96
7) Vinyl Chloride	2.70	62	1459849	8.914	PPB	98
8) 1,3-Butadiene	2.74	54	2648630	9.034	PPB	96
9) Bromomethane	3.04	94	1115109	8.979	PPB	96
10) Chloroethane	3.15	64	747669	9.203	PPB	99
11) Dichlorofluoromethane	3.39	67	2639890	9.498	PPB	98
12) Trichlorofluoromethane	3.44	101	3200752	8.995	PPB	99
13) Pentane	3.55	43	3420180	9.423	PPB	97
14) Acrolein	3.99	56	468953	12.014	PPB	90
15) 1,1-Dichloroethene	4.09	61	2348051	9.196	PPB	87
16) Freon 113	4.12	103	1651499	9.662	PPB	94
17) Acetone	4.21	43	2640401	10.549	PPB	98
18) Methyl Iodide	4.27	142	2901627	9.843	PPB	99
19) Carbon Disulfide	4.35	76	3376956	10.023	PPB	92
20) Acetonitrile	4.60	41	2794458	10.948	PPB	94
21) 3-Chloropropene	4.60	76	576897	8.955	PPB	84
22) Methylene Chloride	4.77	84	1008022	9.679	PPB	77
23) tert-Butyl Alcohol	5.14	59	3120622	10.869	PPB	100
24) Acrylonitrile	5.15	53	1833409	11.073	PPB	98
25) trans-1,2-Dichloroethene	5.16	61	2045724	10.448	PPB	77
26) Methyl t-Butyl Ether	5.21	73	3563393	10.257	PPB	79
27) Hexane	5.59	57	2777091	9.978	PPB	86
28) 1,1-Dichloroethane	5.77	63	2654530	9.200	PPB	99
29) Vinyl Acetate	5.92	43	3807127	12.964	PPB	95
30) cis-1,2-Dichloroethene	6.66	61	1936276	9.533	PPB	77
31) 2-Butanone	6.74	72	620258	11.299	PPB	100
32) Ethyl Acetate	6.87	70	291660	11.839	PPB	100
33) Methyl Acrylate	6.90	55	2926302	11.158	PPB	94
34) Chloroform	7.17	83	2575529	9.744	PPB	96
35) 1,1,1-Trichloroethane	7.41	97	2631265	9.704	PPB	93
36) Carbon Tetrachloride	7.67	117	2720021	9.230	PPB	100
38) 1,2-Dichloroethane	8.05	62	1860162	10.806	PPB	100
39) Benzene	8.01	78	3888784	11.121	PPB	96
40) Isooctane	8.18	57	8916500	11.519	PPB	96
41) Heptane	8.53	43	3891872	11.264	PPB	93
42) Trichloroethene	9.12	130	1842776	9.627	PPB	88
43) Ethyl Acrylate	9.44	55	4034975	13.003	PPB	93
44) 1,2-Dichloropropane	9.51	63	1717764	10.959	PPB	97
45) Methyl Methacrylate	9.86	69	1314562	12.415	PPB	78
46) Dibromomethane	9.72	174	1137611	11.936	PPB	78
47) 1,4-Dioxane	9.89	88	438104	9.370	PPB	38
48) Bromodichloromethane	10.05	83	2805792	11.650	PPB	93

8149

Data Path : C:\MSDCHEM\1\DATA\OCT29\
 Data File : C070741.D
 Acq On : 29 Oct 2005 12:47
 Operator : JBS
 Sample : VSTD010
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Quant Time: Oct 29 13:13:52 2005
 Quant Method : C:\MSDCHEM\1\METHODS\ALLTO15.M
 Quant Title :
 QIast Update : Fri Oct 28 08:49:53 2005
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev	(Min)
49) cis-1,3-Dichloropropene	10.93	75	2070838	10.350	PPB	#	71
50) 4-Methyl-2-Pentanone	11.34	43	4641866	11.210	PPB	#	95
52) Toluene	11.58	91	4305104	11.265	PPB	#	99
53) Octane	11.95	43	5254735	11.531	PPB	#	89
54) trans-1,3-Dichloropropene	12.11	75	1583226	10.656	PPB	#	90
55) Ethyl Methacrylate	12.43	69	2116476	11.677	PPB	#	79
56) 1,1,2-Trichloroethane	12.47	97	1420768	10.137	PPB	#	86
57) Tetrachloroethene	12.70	166	1924760	10.958	PPB	#	98
58) 2-Hexanone	13.17	43	3829686	11.177	PPB	#	78
59) Dibromochloromethane	13.29	127	2086089	10.583	PPB	#	98
60) 1,2-Dibromoethane	13.49	107	2156796	11.149	PPB	#	97
61) Chlorobenzene	14.65	112	3432162	10.884	PPB	#	86
62) 1,1,1,2-Tetrachloroethane	14.89	131	1937484	10.116	PPB	#	98
63) Ethylbenzene	14.99	91	5893567	11.017	PPB	#	92
64) m/p-Xylene	15.29	91	9288802	21.525	PPB	#	91
65) o-Xylene	16.23	91	4621778	11.045	PPB	#	89
66) Styrene	16.28	104	3420745	11.286	PPB	#	100
67) Bromoform	16.66	173	2281105	11.214	PPB	#	97
68) Cumene	17.21	105	5762941	11.888	PPB	#	93
69) 1,1,2,2-Tetrachloroethane	18.04	83	2804623	14.438	PPB	#	96
70) 1,2,3-Trichloropropane	18.07	110	858096	11.408	PPB	#	64
71) Bromobenzene	17.84	156	1827886	11.560	PPB	#	98
72) 4-Ethyltoluene	18.59	105	6129986	10.847	PPB	#	93
73) 1,3,5-Trimethylbenzene	18.78	105	5389513	11.115	PPB	#	92
74) Alpha Methyl Styrene	19.37	118	2609376	12.039	PPB	#	94
75) 1,2,4-Trimethylbenzene	19.74	105	5166606	10.778	PPB	#	90
76) 1,3-Dichlorobenzene	20.40	146	3119547	11.366	PPB	#	96
77) 1,4-Dichlorobenzene	20.69	146	2979498	11.338	PPB	#	97
78) Benzyl Chloride	16.23	91	4621778	11.045	PPB	#	100
79) 1,2-Dichlorobenzene	21.85	146	2722938	10.394	PPB	#	96
80) Hexachloroethane	22.50	117	2090535	11.682	PPB	#	97
81) 1,2,4-Trichlorobenzene	24.92	180	1168915	10.110	PPB	#	98
82) Hexachlorobutadiene	25.18	225	1547505	9.763	PPB	#	96

(#) = qualifier out of range (m) = manual integration (+) = signals summed

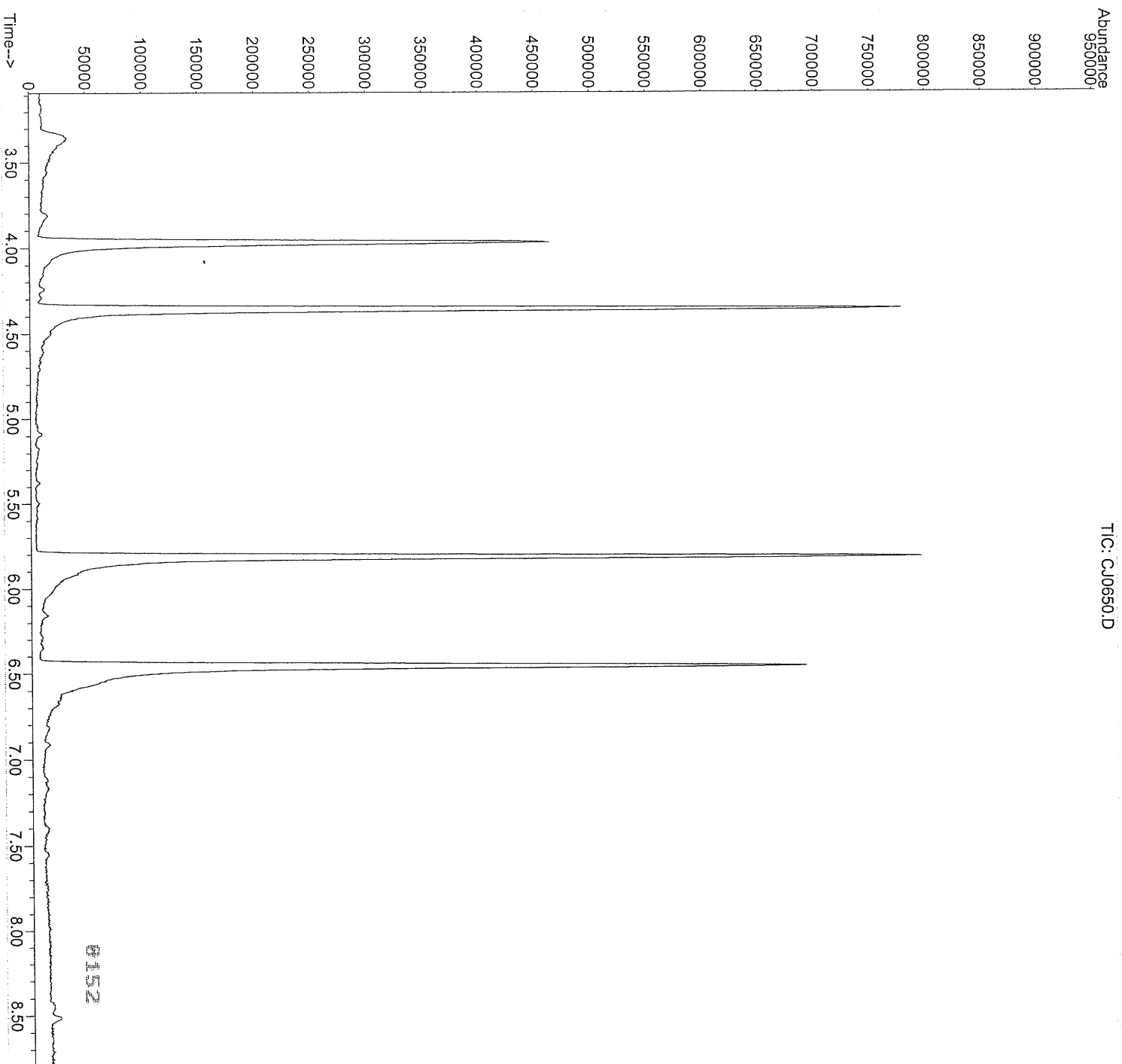
2158

Raw QC Data

0151

Data Path : C:\MSDCHEM\1\DATA\OCT27\
Data File : C:\MSDCHEM\1\DATA\OCT27\
Acq On : 27 Oct 2005 7:28
Operator : JBS
Sample : BFB
Misc :
ALS Vial : 21 Sample Multiplier: 1

Quant Time: Oct 27 07:37:29 2005
Quant Method : C:\MSDCHEM\1\METHODS\BFB5973.M
Quant Title :
Quant Update : Thu Feb 17 14:59:05 2005
Response via : Initial Calibration



Data Path : C:\MSDCHEM\1\DATA\OCT27\
 Data File : CJ0650.D
 Acq On : 27 Oct 2005 7:28
 Operator : JBS
 Sample : BFB
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Quant Time: Oct 27 07:37:29 2005
 Quant Method : C:\MSDCHEM\1\METHODS\BFB5973.M
 Quant Title :
 Qlast Update : Thu Feb 17 14:59:05 2005
 Response via : Initial Calibration

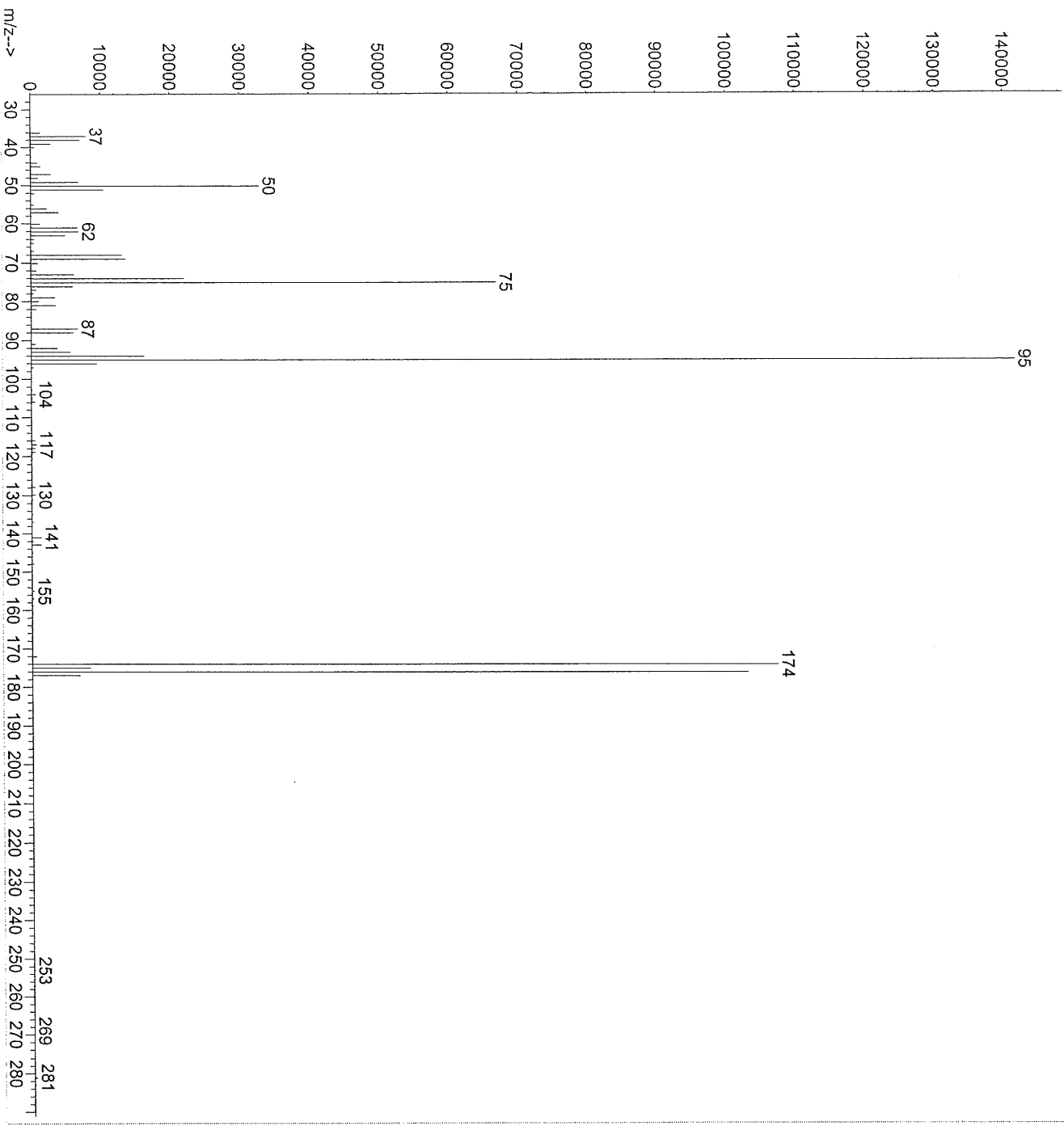
Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)

Target Compounds						Qvalue
1) BFB	6.46	95	3645735	No	Calib	

(#) = qualifier out of range (m) = manual integration (+) = signals summed

0153

Average of 6.451 to 6.458 min.: CJ0650.D (-)



Peak Apex is scan: 1051

Average of 3 scans: 1050,1051,1052 minus background scan 1036

Target Mass	Comparison Mass	Lower Limit, %	Upper Limit, %	Relative Abundance, %	Result
50	95	15	40	23.2	PASS
75	95	30	60	47.2	PASS
95	95	100	100	100.0	PASS
96	95	5	9	6.6	PASS
173	174	0	2	0.0	PASS
174	95	50	100	75.8	PASS
175	174	5	9	7.8	PASS
176	174	95	101	96.0	PASS
177	176	5	9	6.7	PASS

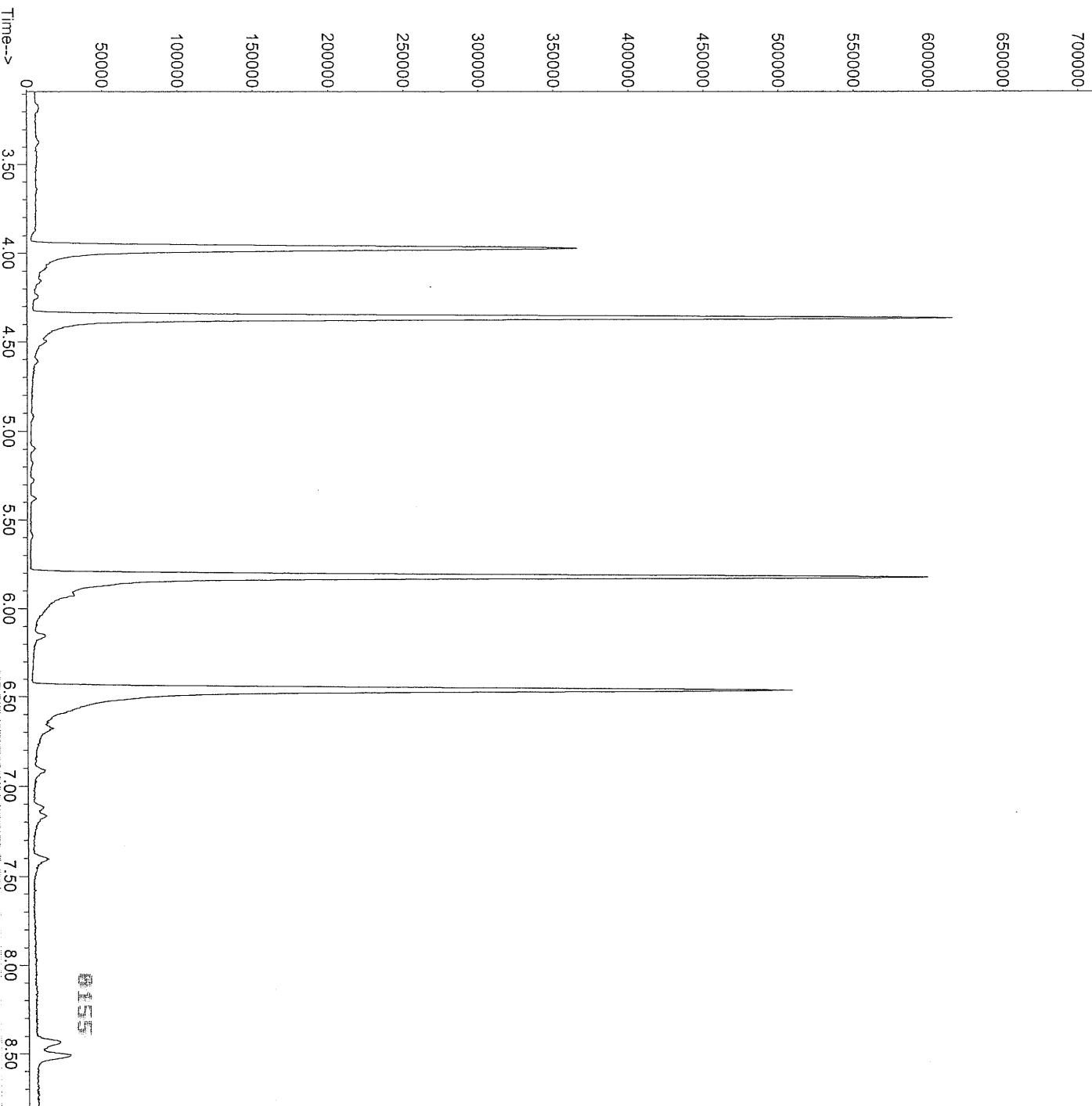
0154

Data Path : C:\MSDCHEM\1\DATA\OCT28\
Data File : CJ0700.D
Acq On : 28 Oct 2005 7:39
Operator : JBS
Sample : BFB
Misc :
ALS Vial : 21 Sample Multiplier: 1

Quant Time: Oct 28 07:48:44 2005
Quant Method : C:\MSDCHEM\1\METHODS\BFB5973.M
Quant Title :
Quant Update : Thu Feb 17 14:59:05 2005
Response via : Initial Calibration

Abundance

TIC: CJ0700.D



Quantitation Report (Not Reviewed)

Data Path : C:\MSDCHEM\1\DATA\OCT28\
Data File : CU0700.D
Acq On : 28 Oct 2005 7:39
Operator : JBS
Sample : BFB
Misc :
ALS Vial : 21 Sample Multiplier: 1

Quant Time: Oct 28 07:48:44 2005
Quant Method : C:\MSDCHEM\1\METHODS\BFB5973.M
Quant Title :
Qlast Update : Thu Feb 17 14:59:05 2005
Response via : Initial Calibration

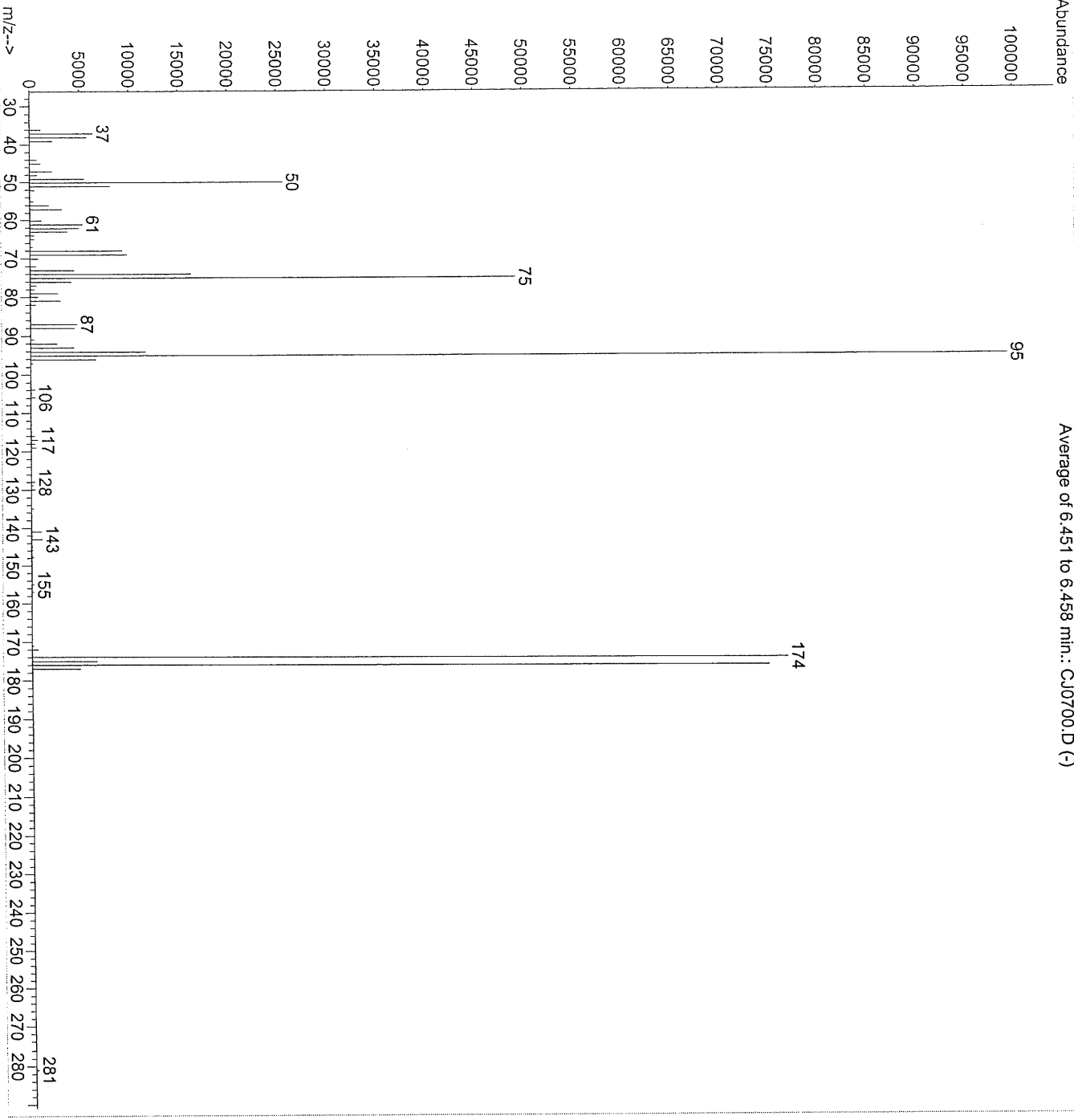
Internal Standards	R.T.	QIon	Response	Conc	Units	Dev	(Min)

Target Compounds							Qvalue
1) BFB	6.46	95	2574147	No	Calib		

(#) = qualifier out of range (m) = manual integration (+) = signals summed

0156

Average of 6.451 to 6.458 min.: CJ0700.D (-)



Peak Apex is scan: 1051

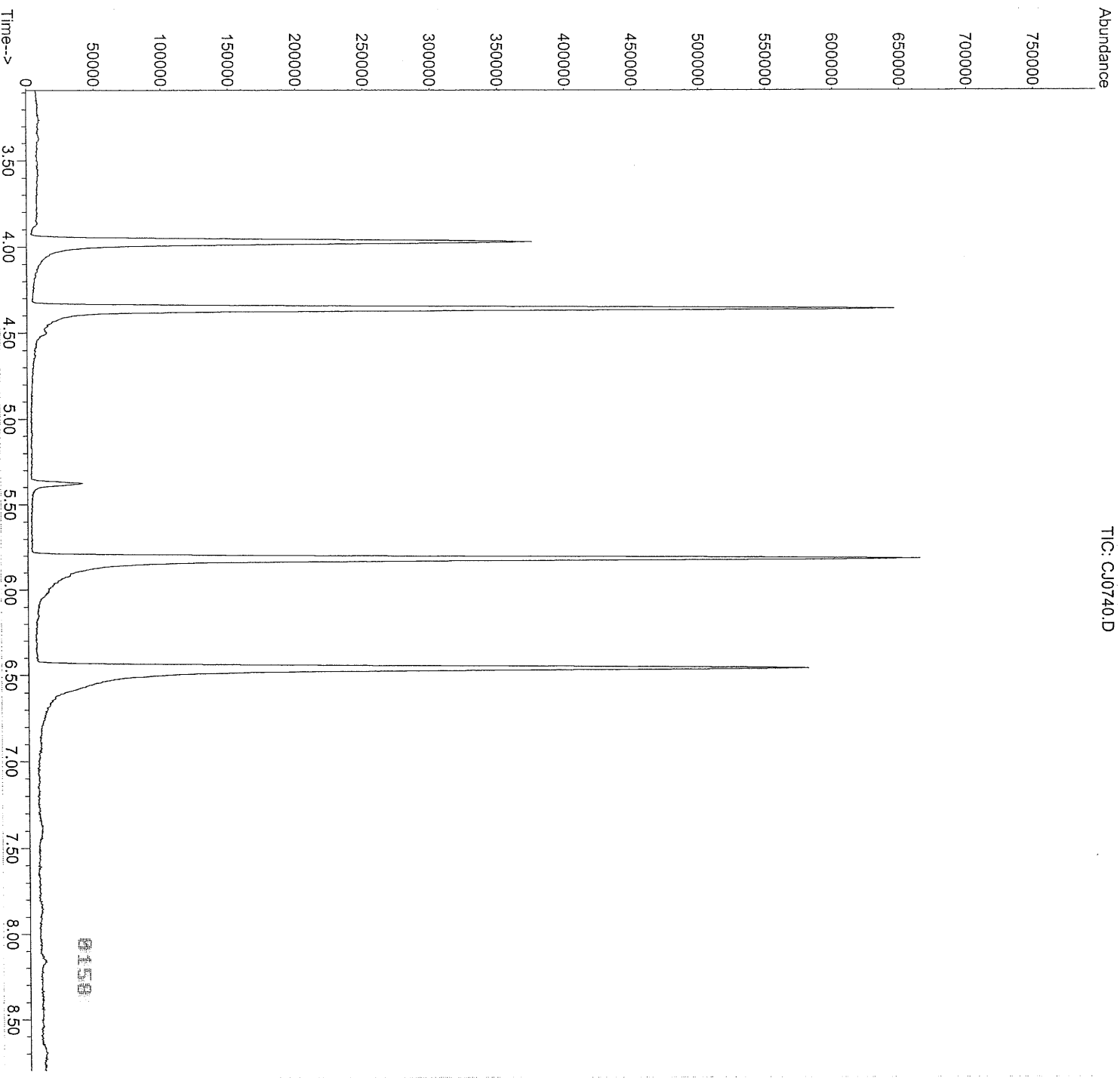
Average of 3 scans: 1050,1051,1052 minus background scan 1038

Target Mass	Comparison Mass	Lower Limit, %	Upper Limit, %	Relative Abundance, %	Result
50	95	15	40	25.9	PASS
75	95	30	60	49.6	PASS
95	95	100	100	100.0	PASS
96	95	5	9	6.6	PASS
173	174	0	2	0.0	PASS
174	95	50	100	77.4	PASS
175	174	5	9	8.5	PASS
176	174	95	101	97.5	PASS
177	176	5	9	6.5	PASS

0157

Data Path : C:\MSDCHEM\1\DATA\OCT29\
Data File : CJ0740.D
Acq On : 29 Oct 2005 12:20
Operator : JBS
Sample : BFB
Misc :
ALS Vial : 21 Sample Multiplier: 1

Quant Time: Oct 29 12:29:50 2005
Quant Method : C:\MSDCHEM\1\METHODS\BFB5973.M
Quant Title :
Quant Update : Thu Feb 17 14:59:05 2005
Response via : Initial Calibration



Quantitation Report (Not Reviewed)

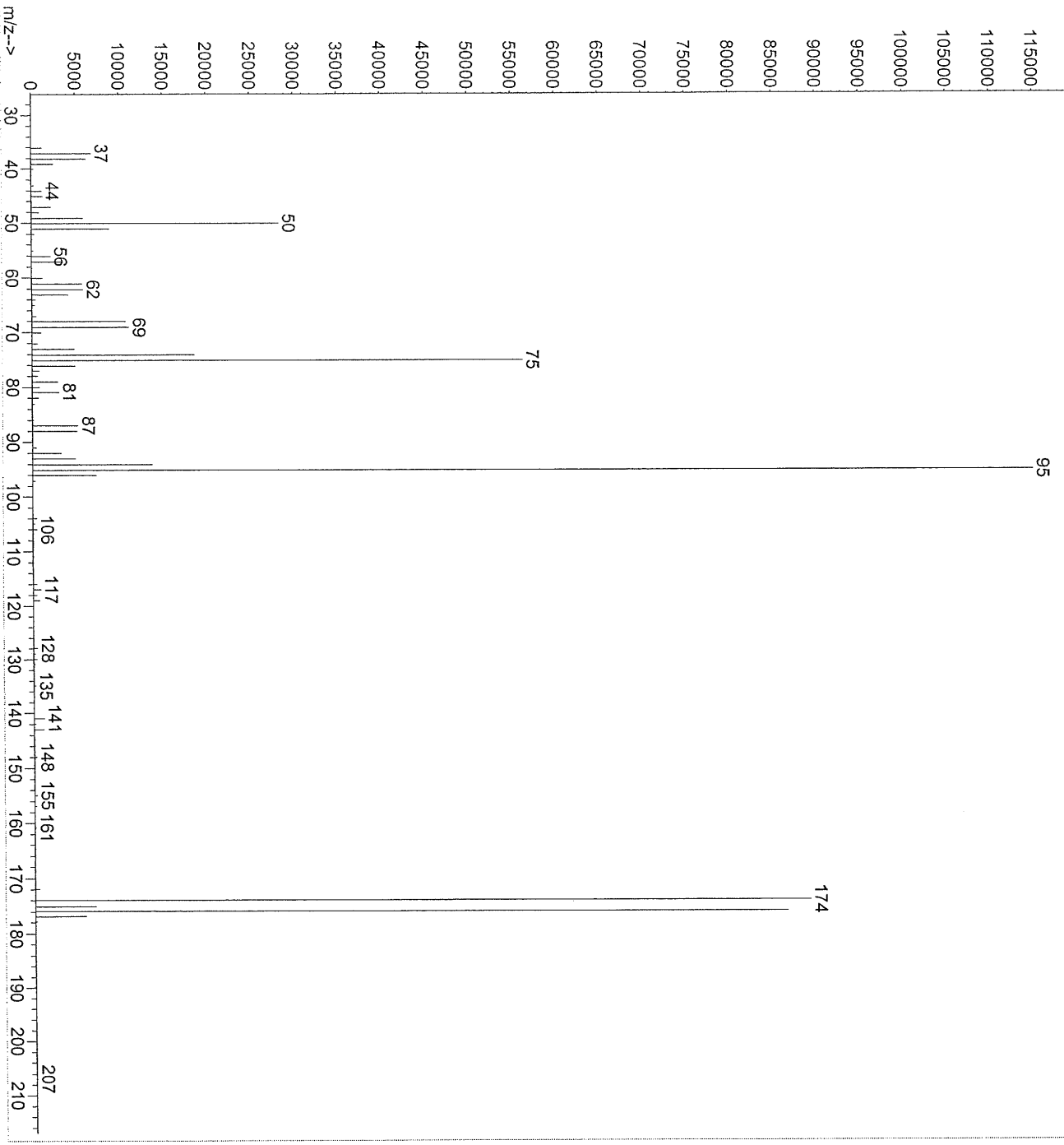
Data Path : C:\MSDCHEM\1\DATA\OCT29\
 Data File : CU0740.D
 Acq On : 29 Oct 2005 12:20
 Operator : JBS
 Sample : BFB
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Quant Time: Oct 29 12:29:50 2005
 Quant Method : C:\MSDCHEM\1\METHODS\BFB5973.M
 Quant Title :
 Qlast Update : Thu Feb 17 14:59:05 2005
 Response via : Initial Calibration

Target Compounds	R.T.	QIon	Response	Conc	Units	Dev (Min)	Qvalue
1) BFB	6.46	95	2942573	No	Calib		

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Average of 6.454 to 6.461 min.: CJ0740.D (-)



Peak Apex is scan: 1052

Average of 3 scans: 1051, 1052, 1053 minus background scan 1037

Target Mass	Comparison Mass	Lower Limit, %	Upper Limit, %	Relative Abundance, %	Result
50	95	15	40	24.7	PASS
75	95	30	60	49.0	PASS
95	95	100	100	100.0	PASS
96	95	5	9	6.4	PASS
173	174	0	2	0.0	PASS
174	95	50	100	77.5	PASS
175	174	5	9	7.8	PASS
176	174	95	101	97.0	PASS
177	176	5	9	6.7	PASS

0160

LANCASTER LABORATORIES
 VOLATILE ORGANICS IN AIR
 SUMMA CANISTER SAMPLE
 ANALYSIS DATA SHEET

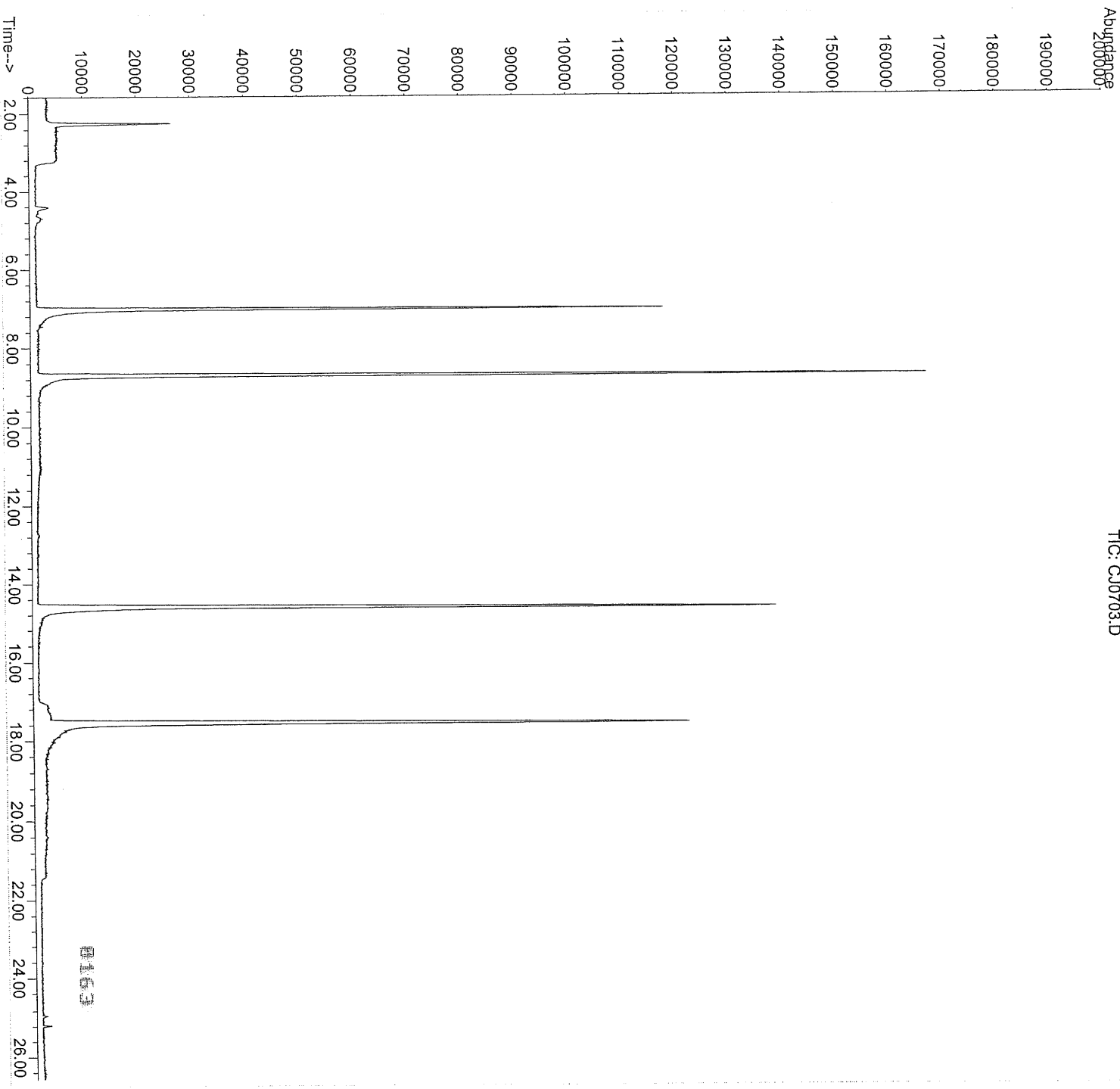
Sample No.: VBLKC23 Date Collected: Date Received: Time Analyzed: 09:42
 Lab Sample ID: VBLKC23 Date Analyzed: 10/28/05
 Canister ID: SUMMA Pressure Rec'd: 14.7 psia Final Pressure: 14.7 psia
 Injection Volume: 250 cc Nominal Volume: 250 cc Dilution Factor: 1.0
 Instrument ID: HP09464 ^{Sum} Lab File ID: C:\MSDCHEM\1\DATA\OCT28\CJ0703.D
 11/6/05

CAS RN	COMPOUND NAME	CONCENTRATION UNITS:	MDL	ppb (v)	Q
75-71-8	Dichlorodifluoromethane			0.2	U
76-14-2	Freon 114			0.2	U
74-87-3	Chloromethane			0.2	U
75-01-4	Vinyl Chloride			0.2	U
74-83-9	Bromomethane			0.2	U
75-00-3	Chloroethane			0.2	U
75-69-4	Trichlorofluoromethane			0.2	U
75-35-4	1,1-Dichloroethane			0.2	U
76-13-1	Freon 113			0.5	U
107-05-1	3-Chloropropene			0.5	U
75-09-2	Methylene Chloride			0.5	U
75-34-3	1,1-Dichloroethane			0.2	U
156-59-2	cis-1,2-Dichloroethene			0.2	U
67-66-3	Chloroform			0.2	U
71-55-6	1,1,1-Trichloroethane			0.2	U
56-23-5	Carbon Tetrachloride			0.2	U
107-06-2	1,2-Dichloroethane			0.2	U
71-43-2	Benzene			0.2	U
79-01-6	Trichloroethene			0.2	U
78-87-5	1,2-Dichloropropane			0.2	U
10061-01-5	cis-1,3-Dichloropropene			0.2	U
108-88-3	Toluene			0.2	U
10061-02-6	trans-1,3-Dichloropropene			0.2	U
79-00-5	1,1,2-Trichloroethane			0.2	U
127-18-4	Tetrachloroethene			0.2	U
106-93-4	1,2-Dibromoethane			0.2	U
108-90-7	Chlorobenzene			0.2	U
100-41-4	Ethylbenzene			0.2	U
1330-20-7	m/p-Xylene			0.2	U
95-47-6	o-Xylene			0.2	U
100-42-5	Styrene			0.2	U
79-34-5	1,1,2,2-Tetrachloroethane			0.2	U
622-96-8	4-Ethyltoluene			0.2	U
108-67-8	1,3,5-Trimethylbenzene			0.2	U

U = Compound was undetected at the specified limit of quantitation.
 B = Compound was found in method blank. D = analysis of diluted sample.
 J = Compound was detected, but below the limit of quantitation.

Data Path : C:\MSDCHEM\1\DATA\OCT28\
Data File : CJ0703.D
Acq On : 28 Oct 2005 9:42
Operator : JBS
Sample : VBKIC23
Misc :
ALS Vial : 34 Sample Multiplier: 1
Quant Time: Oct 28 10:08:48 2005
Quant Method : C:\MSDCHEM\1\METHODS\ALLT015.M
Quant Title :
Quant Update : Fri Oct 28 08:49:53 2005
Response via : Initial Calibration

TIC: CJ0703.D



Data Path : C:\MSDCHEM\1\DATA\OCT28\
 Data File : CJ0703.D
 Acq On : 28 Oct 2005 9:42
 Operator : JBS
 Sample : VBLKC23
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Quant Time: Oct 28 10:08:48 2005
 Quant Method : C:\MSDCHEM\1\METHODS\ALLTO15.M
 Quant Title :
 QLast Update : Fri Oct 28 08:49:53 2005
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	7.01	130	566273	10.000		0.00
37) 1,4-Difluorobenzene	8.69	114	2061811	10.000		0.00
51) Chlorobenzene d5	14.57	117	1826290	10.000		0.00

Target Compounds

Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

LANCASTER LABORATORIES
VOLATILE ORGANICS IN AIR
SUMMA CANISTER SAMPLE
ANALYSIS DATA SHEET

Sample No.: VBLKC25 Date Collected: Date Received: 14:57
 Lab Sample ID: VBLKC25 Date Analyzed: 10/29/05 Time Analyzed: 14:57
 Canister ID: SUMMA Pressure Rec'd: 14.7 psia Final Pressure: 14.7 psia
 Injection Volume: 250 cc Nominal Volume: 250 cc Dilution Factor: 1.0
 Instrument ID: HP09464 ^{11/6/05} Lab File ID: C:\MSDCHEM\1\DATA\OCT29\CJ0744.D

CAS RN	COMPOUND NAME	CONCENTRATION	UNITS: MDL	ppb (v)	Q
75-71-8	Dichlorodifluoromethane	0.2			U
76-14-2	Freon 114	0.2			U
74-87-3	Chloromethane	0.2			U
75-01-4	Vinyl Chloride	0.2			U
74-83-9	Bromomethane	0.2			U
75-00-3	Chloroethane	0.2			U
75-69-4	Trichlorofluoromethane	0.2			U
75-35-4	1,1-Dichloroethene	0.2			U
76-13-1	Freon 113	0.5			U
107-05-1	3-Chloropropene	0.5			U
75-09-2	Methylene Chloride	0.5			U
75-34-3	1,1-Dichloroethane	0.2			U
156-59-2	cis-1,2-Dichloroethene	0.2			U
67-66-3	Chloroform	0.2			U
71-55-6	1,1,1-Trichloroethane	0.2			U
56-23-5	Carbon Tetrachloride	0.2			U
107-06-2	1,2-Dichloroethane	0.2			U
71-43-2	Benzene	0.2			U
79-01-6	Trichloroethene	0.2			U
78-87-5	1,2-Dichloropropane	0.2			U
10061-01-5	cis-1,3-Dichloropropene	0.2			U
108-88-3	Toluene	0.2			U
10061-02-6	trans-1,3-Dichloropropene	0.2			U
79-00-5	1,1,2-Trichloroethane	0.2			U
127-18-4	Tetrachloroethene	0.2			U
106-93-4	1,2-Dibromoethane	0.2			U
108-90-7	Chlorobenzene	0.2			U
100-41-4	Ethylbenzene	0.2			U
1330-20-7	m/p-Xylene	0.2			U
95-47-6	o-Xylene	0.2			U
100-42-5	Styrene	0.2			U
79-34-5	1,1,2,2-Tetrachloroethane	0.2			U
622-96-8	4-Ethyltoluene	0.2			U
108-67-8	1,3,5-Trimethylbenzene	0.2			U

U = Compound was undetected at the specified limit of quantitation.
 B = Compound was found in method blank. D = analysis of diluted sample.
 J = Compound was detected, but below the limit of quantitation.

LANCASTER LABORATORIES
 VOLATILE ORGANICS IN AIR
 SUMMA CANISTER SAMPLE
 ANALYSIS DATA SHEET

Sample No.: VBLKC25 Date Collected: Date Received: Time Analyzed: 14:57
 Lab Sample ID: VBLKC25 Date Analyzed: 10/29/05
 Canister ID: SUMMA Pressure Rec'd: 14.7 psia Final Pressure: 14.7 psia
 Injection Volume: 250 cc Nominal Volume: 250 cc Dilution Factor: 1.0
 Instrument ID: HP09464 ^{4/14/05} Lab File ID: C:\MSDCHEM\1\DATA\OCT29\CJ0744.D

CAS RN	COMPOUND NAME	CONCENTRATION	UNITS	MDL	ppb (v)	Q
95-63-6	1,2,4-Trimethylbenzene				0.2	U
541-73-1	1,3-Dichlorobenzene				0.5	U
106-46-7	1,4-Dichlorobenzene				0.5	U
95-50-1	1,2-Dichlorobenzene				0.5	U
120-82-1	1,2,4-Trichlorobenzene				1	U

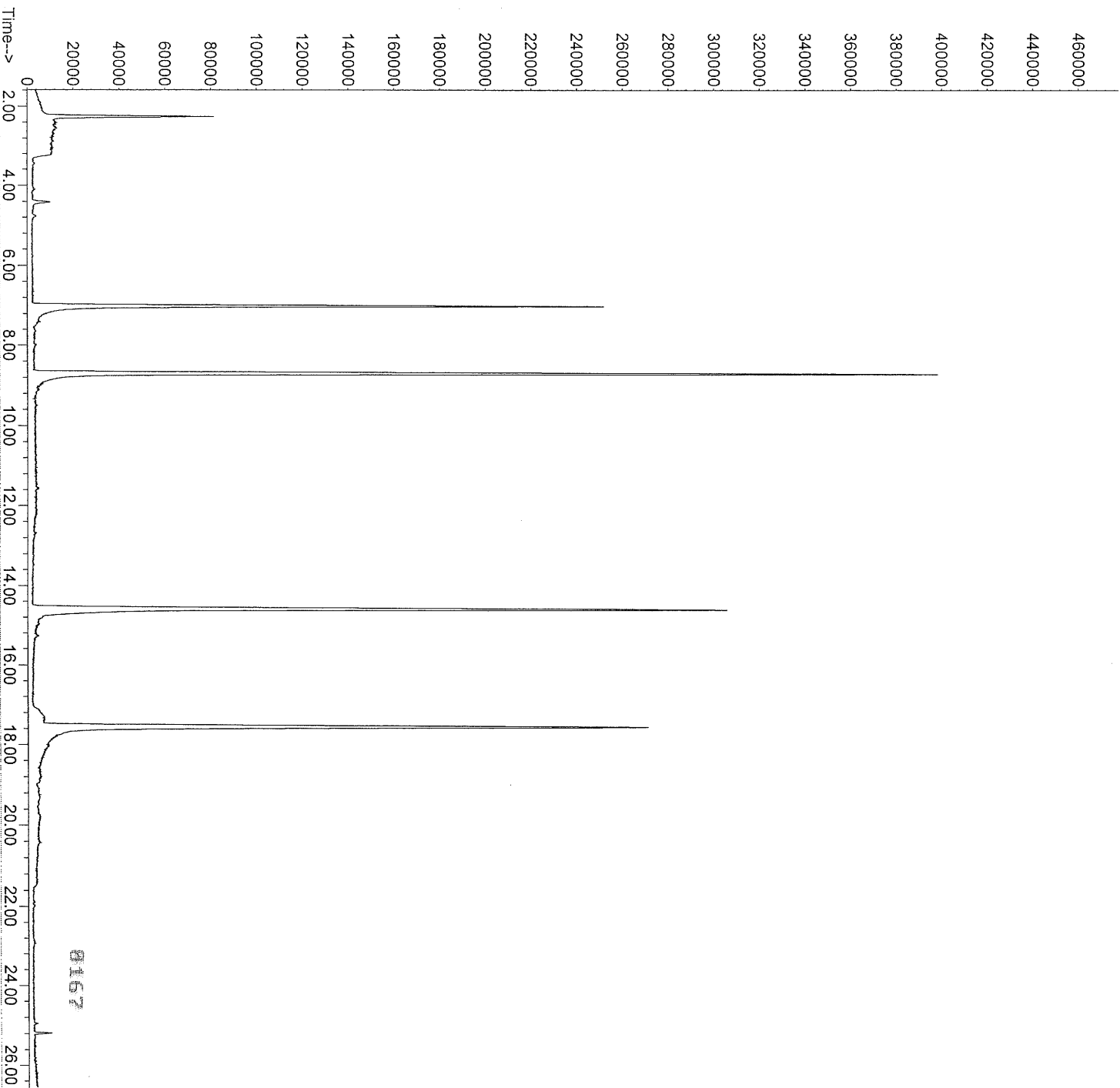
U = Compound was undetected at the specified limit of quantitation.
 B = Compound was found in method blank. D = analysis of diluted sample.
 J = Compound was detected, but below the limit of quantitation.

Data Path : C:\MSDCHEM\1\DATA\OCT29\
Data File : CJ0744.D
Acq On : 29 Oct 2005 14:57
Operator : JBS
Sample : VBLKC25
Misc :
ALS Vial : 35 Sample Multiplier: 1

Quant Time: Oct 29 15:24:11 2005
Quant Method : C:\MSDCHEM\1\METHODS\ALLT015.M
Quant Title :
Quant Update : Sat Oct 29 13:19:44 2005
Response via : Initial Calibration

Abundance

TIC: CJ0744.D



Data Path : C:\MSDCHEM\1\DATA\OCT29\
 Data File : CJ0744.D
 Acq On : 29 Oct 2005 14:57
 Operator : JBS
 Sample : VBLKC25
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Quant Time: Oct 29 15:24:11 2005
 Quant Method : C:\MSDCHEM\1\METHODS\ALLTO15.M
 Quant Title :
 QLast Update : Sat Oct 29 13:19:44 2005
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	7.01	130	1282140	10.000		0.00
37) 1,4-Difluorobenzene	8.69	114	5110543	10.000		-0.01
51) Chlorobenzene d5	14.57	117	4104204	10.000		-0.01
Target Compounds						Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

0158

LANCASTER LABORATORIES
VOLATILE ORGANICS IN AIR
SUMMA CANISTER SAMPLE
ANALYSIS DATA SHEET

Sample No.: LCSC23 Date Collected: Date Received: 10/28/05 Time Analyzed: 10:18
 Lab Sample ID: LCSC23 Date Analyzed: 10/28/05 Pressure Rec'd: 14.7 psia Final Pressure: 14.7 psia
 Canister ID: SUMMA Injection Volume: 250 cc Nominal Volume: 250 cc Dilution Factor: 1.0
 Instrument ID: HP09464 ¹⁹² _{11/4/05} Lab File ID: C:\MSDCHEM\1\DATA\OCT28\CJ0704.D

CAS RN	COMPOUND NAME	CONCENTRATION UNITS: MDL	ug/m3	Q
75-71-8	Dichlorodifluoromethane		37	
76-14-2	Freon 114		52	
74-87-3	Chloromethane		15	
75-01-4	Vinyl Chloride		18	
74-83-9	Bromomethane		29	
75-00-3	Chloroethane		20	
75-69-4	Trichlorofluoromethane		44	
75-35-4	1,1-Dichloroethene		31	
76-13-1	Freon 113		65	
107-05-1	3-Chloropropene		2	
75-09-2	Methylene Chloride		28	
75-34-3	1,1-Dichloroethane		33	
156-59-2	cis-1,2-Dichloroethene		32	
67-66-3	Chloroform		43	
71-55-6	1,1,1-Trichloroethane		48	
56-23-5	Carbon Tetrachloride		53	
107-06-2	1,2-Dichloroethane		38	
71-43-2	Benzene		27	
79-01-6	Trichloroethene		42	
78-87-5	1,2-Dichloropropane		39	
10061-01-5	cis-1,3-Dichloropropene		34	
108-88-3	Toluene		33	
10061-02-6	trans-1,3-Dichloropropene		40	
79-00-5	1,1,2-Trichloroethane		46	
127-18-4	Tetrachloroethene		55	
106-93-4	1,2-Dibromoethane		66	
108-90-7	Chlorobenzene		38	
100-41-4	Ethylbenzene		36	
1330-20-7	m/p-Xylene		73	
95-47-6	o-Xylene		39	
100-42-5	Styrene		40	
79-34-5	1,1,2,2-Tetrachloroethane		77	
622-96-8	4-Ethyltoluene		41	
108-67-8	1,3,5-Trimethylbenzene		48	

U = Compound was undetected at the specified limit of quantitation.
 B = Compound was found in method blank. D = analysis of diluted sample.
 J = Compound was detected, but below the limit of quantitation.

LANCASTER LABORATORIES
 VOLATILE ORGANICS IN AIR
 SUMMA CANISTER SAMPLE
 ANALYSIS DATA SHEET

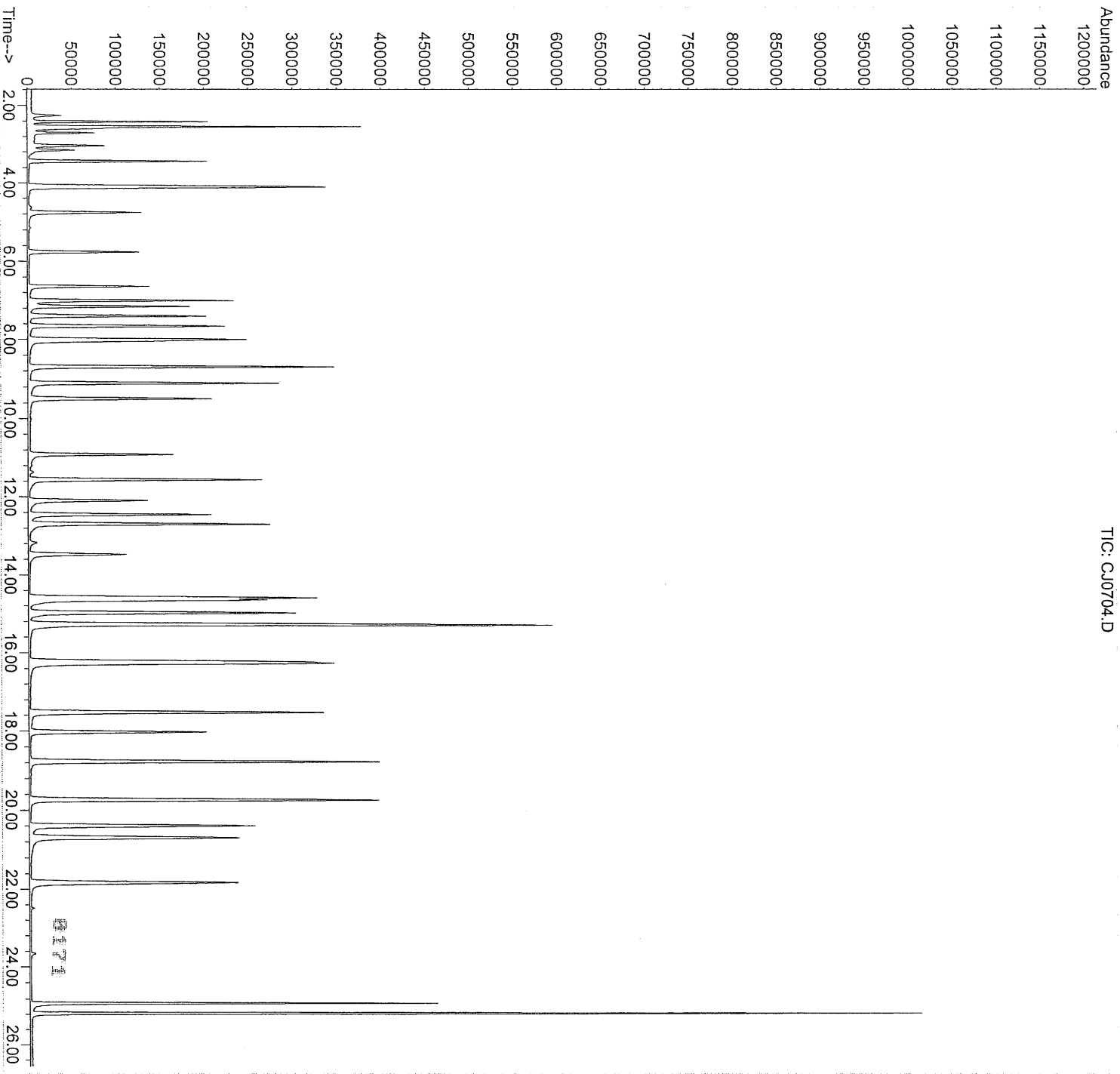
Sample No.: LCSC23 Date Collected: Date Received: Time Analyzed: 10:18
 Lab Sample ID: LCSC23 Date Analyzed: 10/28/05 Final Pressure: 14.7 psia
 Canister ID: SUMMA Pressure Rec'd: 14.7 psia Dilution Factor: 1.0
 Injection Volume: 250 cc Nominal Volume: 250 cc
 Instrument ID: HP09464 ^{11/4/05} Lab File ID: C:\MSDCHEM\1\DATA\OCT28\CJ0704.D

CAS RN	COMPOUND NAME	CONCENTRATION	UNITS	MDL	ug/m3	Q
95-63-6	1,2,4-Trimethylbenzene				50	
541-73-1	1,3-Dichlorobenzene				57	
106-46-7	1,4-Dichlorobenzene				59	
95-50-1	1,2-Dichlorobenzene				56	
120-82-1	1,2,4-Trichlorobenzene				120	

U = Compound was undetected at the specified limit of quantitation.
 B = Compound was found in method blank. D = analysis of diluted sample.
 J = Compound was detected, but below the limit of quantitation.

Data Path : C:\MSDCHEM\1\DATA\OCT28\
Data File : CJ0704.D
Acq On : 28 Oct 2005 10:18
Operator : JBS
Sample : LCSC23
Misc :
ALS Vial : 35 Sample Multiplier: 1
Quant Time: Oct 28 10:45:25 2005
Quant Method : C:\MSDCHEM\1\METHODS\ALLT015.M
Quant Title :
Quant Update : Fri Oct 28 08:49:53 2005
Response via : Initial Calibration

TIC: CJ0704.D



Data Path : C:\MSDCHEM\1\DATA\OCT28\
 Data File : CJ0704.D
 Acq On : 28 Oct 2005 10:18
 Operator : JBS
 Sample : LCSC23
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Quant Time: Oct 28 10:45:25 2005
 Quant Method : C:\MSDCHEM\1\METHODS\ALLT015.M
 Quant Title :
 Quant Update : Fri Oct 28 08:49:53 2005
 Response via : Initial Calibration

Internal Standards	R.T.	Qion	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	7.01	130	1023334	10.000	PPB	0.00
37) 1,4-Difluorobenzene	8.69	114	3957263	10.000	PPB	0.00
51) Chlorobenzene d5	14.58	117	3645200	10.000	PPB	0.00

Target Compounds	R.T.	Qion	Response	Conc	Units	Dev(Min)	Qvalue
3) Dichlorodifluoromethane	2.42	85	2243514	7.488	PPB		100
5) Freon 114	2.54	85	2552697	7.504	PPB	#	80
6) Chloromethane	2.59	50	1432922	7.442	PPB		96
7) Vinyl Chloride	2.70	62	964855	7.142	PPB		99
9) Bromomethane	3.04	94	758755	7.406	PPB		96
10) Chloroethane	3.15	64	519673	7.754	PPB		99
12) Trichlorofluoromethane	3.44	101	2290542	7.803	PPB		98
15) 1,1-Dichloroethene	4.08	61	1639890	7.786	PPB	#	86
16) Freon 113	4.11	103	1195514	8.478	PPB		93
22) Methylene Chloride	4.76	84	703143	8.184	PPB	#	70
28) 1,1-Dichloroethane	5.76	63	1929975	8.108	PPB		100
30) cis-1,2-Dichloroethene	6.65	61	1350221	8.059	PPB	#	75
34) Chloroform	7.16	83	1912082	8.769	PPB		94
35) 1,1,1-Trichloroethane	7.40	97	1971257	8.813	PPB	#	92
36) Carbon Tetrachloride	7.66	117	2046138	8.417	PPB		99
38) 1,2-Dichloroethane	8.04	62	1482453	9.313	PPB		99
39) Benzene	8.00	78	2726239	8.431	PPB	#	93
42) Trichloroethene	9.10	130	1387041	7.836	PPB	#	86
44) 1,2-Dichloropropane	9.50	63	1236740	8.532	PPB		97
49) cis-1,3-Dichloropropene	10.91	75	1403670	7.586	PPB	#	72
50) 4-Methyl-2-Pentanone	11.37	43	77938	0.204	PPB	#	49
52) Toluene	11.56	91	3376257	8.626	PPB		100
54) trans-1,3-Dichloropropene	12.09	75	1333806	8.766	PPB		97
56) 1,1,2-Trichloroethane	12.45	97	1219411	8.496	PPB		88
57) Tetrachloroethene	12.69	166	1455357	8.090	PPB		99
58) 2-Hexanone	13.18	43	168999	0.482	PPB	#	87
60) 1,2-Dibromoethane	13.47	107	1689947	8.530	PPB		98
61) Chlorobenzene	14.64	112	2641298	8.179	PPB	#	86
63) Ethylbenzene	14.97	91	4560282	8.324	PPB		92
64) m/p-Xylene	15.27	91	7478820	16.923	PPB		91
65) o-Xylene	16.22	91	3811061	8.893	PPB		89
66) Styrene	16.26	104	2949851	9.503	PPB	#	100
69) 1,1,2,2-Tetrachloroethane	18.02	83	2235612	11.238	PPB		96
72) 4-Ethyltoluene	18.76	105	4814286	8.318	PPB	#	74
73) 1,3,5-Trimethylbenzene	18.76	105	4809353	9.685	PPB		92
74) Alpha Methyl Styrene	19.73	118	63780	0.287	PPB	#	1
75) 1,2,4-Trimethylbenzene	19.73	105	4968238	10.120	PPB		90
76) 1,3-Dichlorobenzene	20.39	146	2672400	9.507	PPB		97
77) 1,4-Dichlorobenzene	20.68	146	2634489	9.789	PPB		96
78) Benzyl Chloride	16.22	91	3811061	8.893	PPB	#	100
79) 1,2-Dichlorobenzene	21.84	146	2509671	9.354	PPB		96
81) 1,2,4-Trichlorobenzene	24.91	180	1989343	16.800	PPB		99
82) Hexachlorobutadiene	25.18	225	2286802	14.087	PPB		97

0172

(#) = qualifier out of range (m) = manual integration (+) = signals summed