Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\061505\G8368.D
Acq On : 15 Jun 05 18:17
Sample : 0506092-06a
Misc : samp asp_8260s
MS Integration Params: rteint.p
Quant Time: Jun 16 9:22 19105

Vial: 13
Operator: xl
Inst : inst g
Multiplr: 1.00

HA-6

Quant Results File: 8260BASP.RE

Quant Method : C:\HPCHEM\1\DATA\061505\8260BASP.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Fri May 20 09:55:33 2005
Response via : Initial Calibration
DataAcq Meth : 8260BASP

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) pentafluorobenzene	7.28	168	128700	50.00	ug/Kg	0.03
34) 1,4-difluorobenzene	8.41	114	299680	50.00	ug/Kg	0.05
52) chlorobenzene-d5	13.17	117	255469	50.00	ug/Kg	0.03
66) 1,4-dichlorobenzene-d4	17.31	152	96392	50.00	ug/Kg	0.04
System Monitoring Compounds						
31) Dibromofluoromethane	7.13	113	92345	55.00	ug/Kg	0.04
Spiked Amount 50.000	Range 80 - 120		Recovery = 110.00%			
48) toluene-d8	10.72	98	370254	50.30	ug/Kg	0.04
Spiked Amount 50.000	Range 81 - 120		Recovery = 100.60%			
65) 4-Bromofluorobenzene	15.27	95	120009	45.27	ug/Kg	0.04
Spiked Amount 50.000	Range 74 - 121		Recovery = 90.54%			
Target Compounds						Qvalue
54) tetrachloroethene	11.70	166	8279	4.35	ug/Kg	96

(#) = qualifier out of range (m) = manual integration

000087

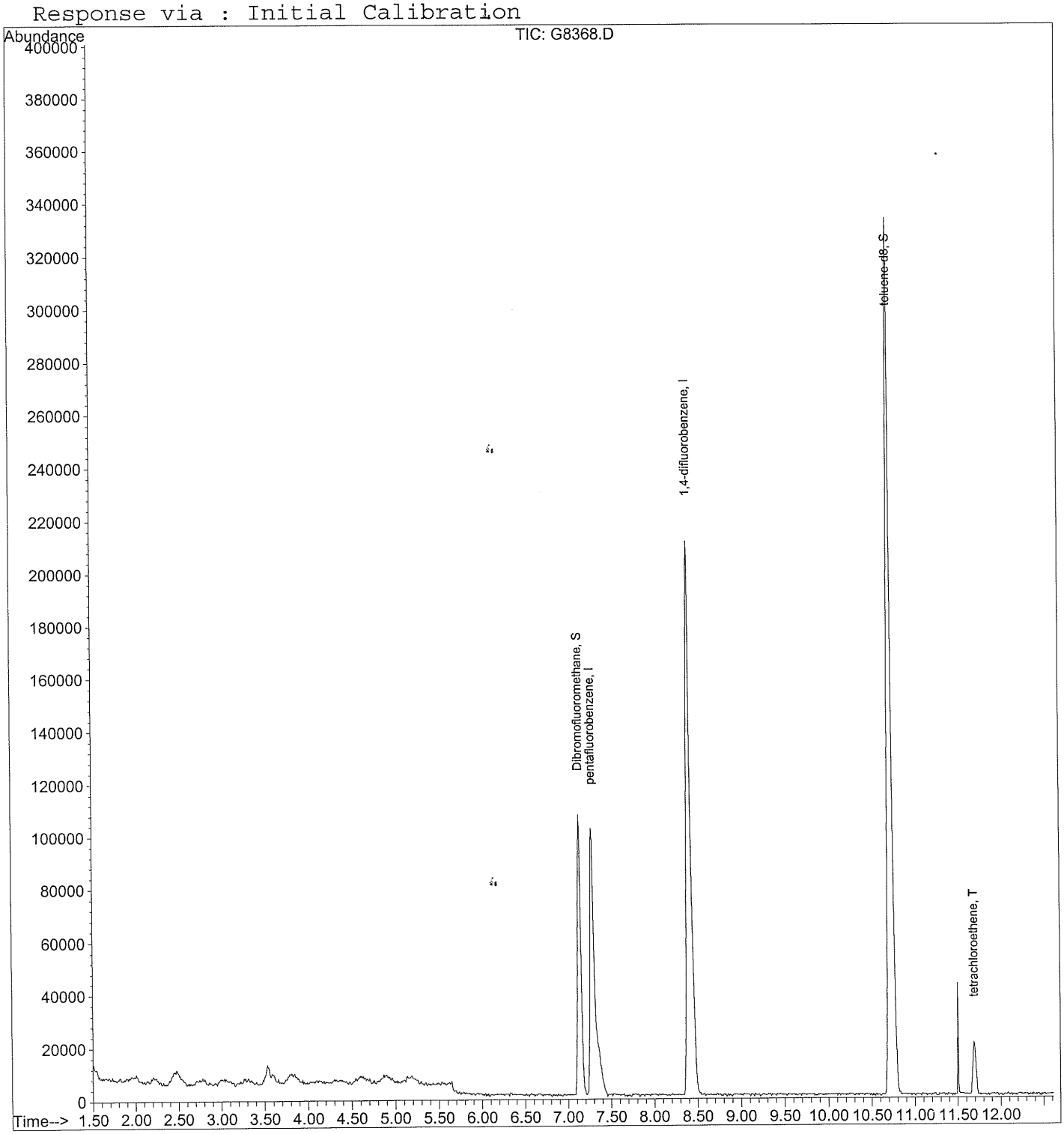
Quantitation Report

Data File : C:\HPCHEM\1\DATA\061505\G8368.D
Acq On : 15 Jun 05 18:17
Sample : 0506092-06a
Misc : samp asp_8260s
MS Integration Params: rteint.p
Quant Time: Jun 16 9:22 19105

Vial: 13
Operator: xl
Inst : inst g
Multiplr: 1.00

Quant Results File: 8260BASP.RE

Method : C:\HPCHEM\1\DATA\081605\8260BW.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Tue Aug 16 18:32:56 2005
Response via : Initial Calibration



000088

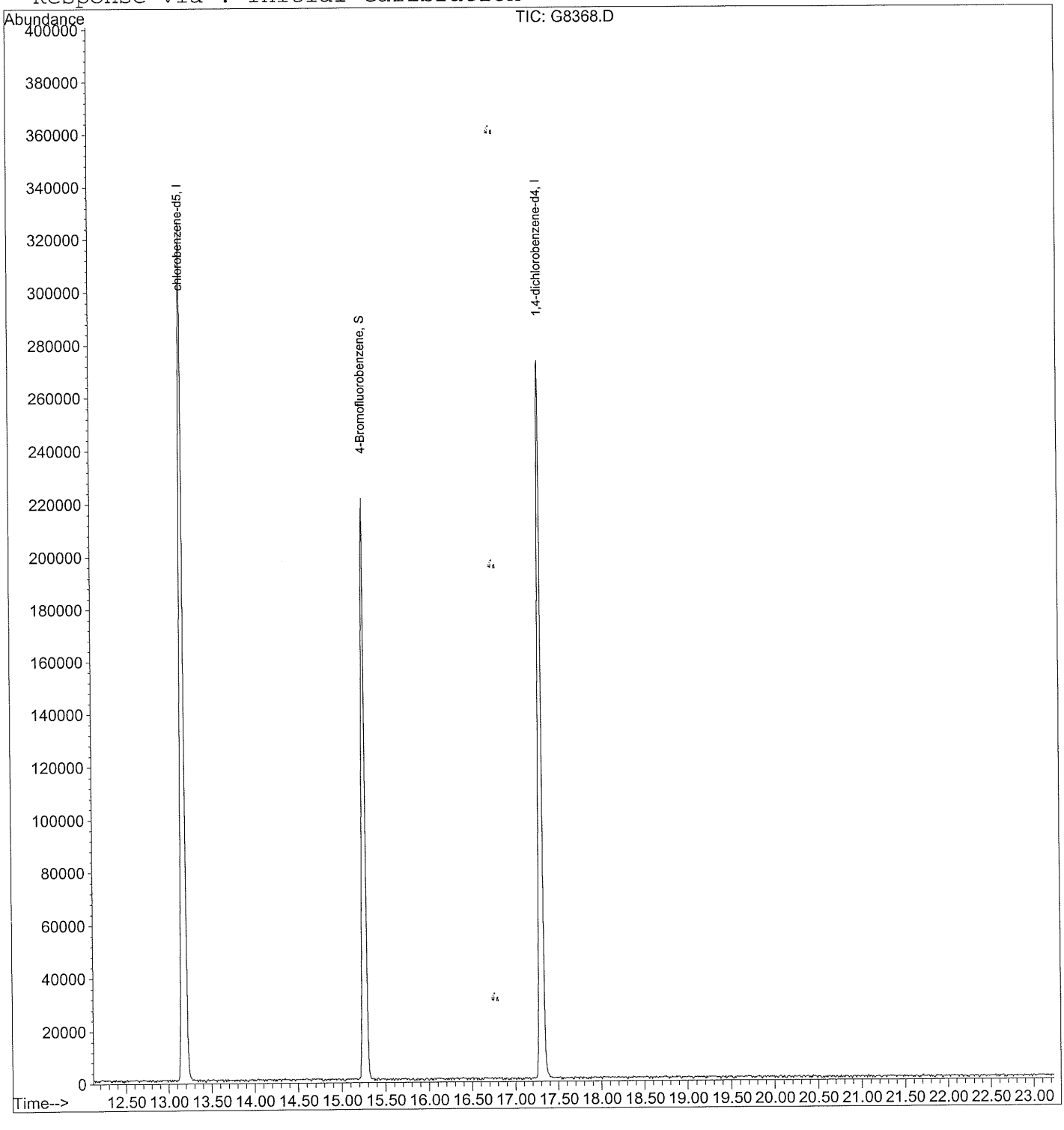
Quantitation Report

Data File : C:\HPCHEM\1\DATA\061505\G8368.D
Acq On : 15 Jun 05 18:17
Sample : 0506092-06a
Misc : samp asp_8260s
MS Integration Params: rteint.p
Quant Time: Jun 16 9:22 19105

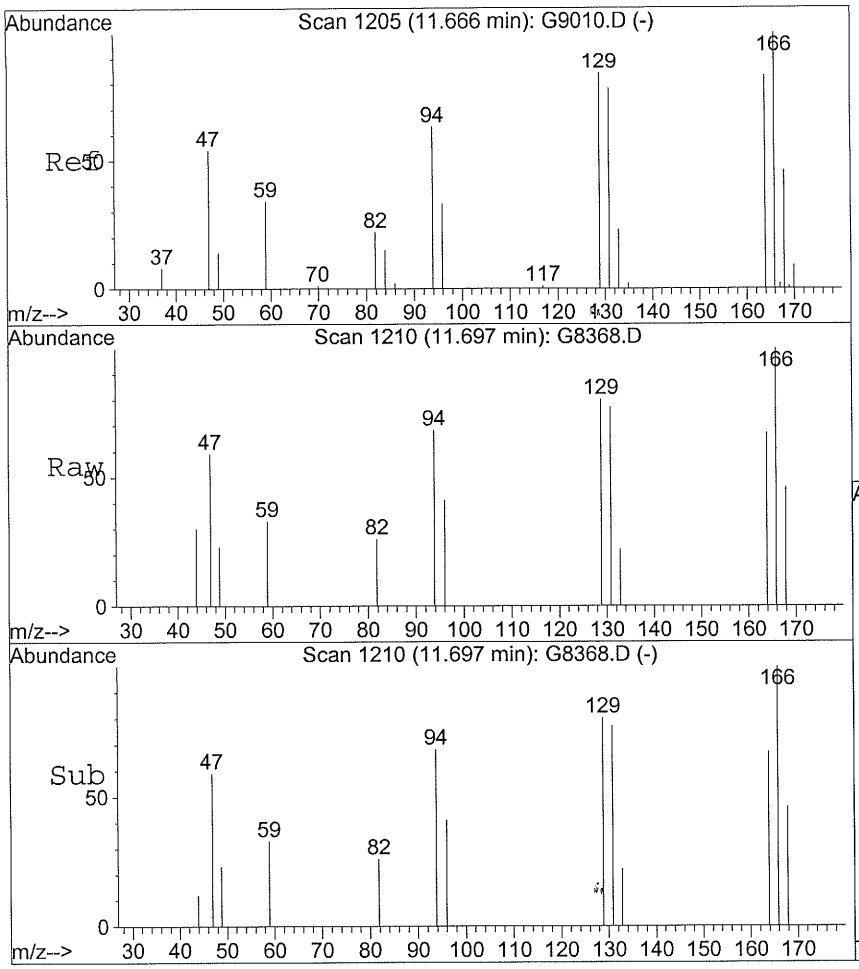
Vial: 13
Operator: xl
Inst : inst g
Multiplr: 1.00

Quant Results File: 8260BASP.RE

Method : C:\HPCHEM\1\DATA\081605\8260BW.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Tue Aug 16 18:32:56 2005
Response via : Initial Calibration



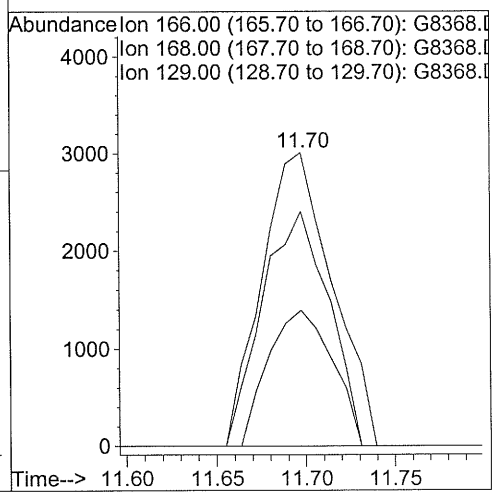
000089



#54
tetrachloroethene
Concen: 4.35 ug/Kg
RT: 11.70 min Scan# 1210
Delta R.T. 0.04 min
Lab File: G8368.D
Acq: 15 Jun 05 18:17

Tgt Ion:166 Resp: 8279

Ion	Ratio	Lower	Upper
166	100		
168	46.4	25.9	65.9
129	79.9	54.8	94.8



000090

STANDARD DATA
QUANT REPORT
CHROMATOGRAMS

Response Factor Report inst g

Method : C:\HPCHEM\1\DATA\051205\8260BASP.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Tue May 17 11:42:05 2005
Response via : Initial Calibration

Calibration Files
10 =G7889.D 20 =G7890.D 50 =G7891.D
100 =G7894.D 200 =G7893.D =

		Compound	10	20	50	100	200	Avg	%RSD
1)	I	pentafluorobenzene	----------ISTD----------						
2)	T	dichlorodifluoromet	1.007	1.070	1.067	1.117	0.953	1.043	6.10
3)	P	chloromethane	1.694	1.805	1.763	1.700	1.788	1.750	2.90
4)	C	vinyl chloride	1.779	1.877	1.874	1.960	1.882	1.874	3.41
5)	T	bromomethane	0.952	0.882	1.011	0.633	0.544	0.804	25.46
6)	T	chloroethane	1.005	1.024	1.032	1.074	1.001	1.027	2.85
7)	T	trichlorofluorometh	1.768	1.867	1.828	1.928	1.849	1.848	3.15
8)	T	Diethyl Ether	0.547	0.501	0.448	0.438	0.414	0.469	11.44
9)	T	Acrolein	0.104	0.091	0.101	0.094	0.089	0.096	6.84
10)	T	Acetone	0.366	0.376	0.265	0.252	0.239	0.300	21.97
11)	CM	1,1-dichloroethene	1.071	1.129	1.118	1.170	1.119	1.121	3.14
12)	T	Iodomethane	1.573	1.593	1.594	1.759	1.682	1.640	4.80
13)	T	methylene chloride	1.259	1.309	1.382	1.304	1.228	1.296	4.50
14)	T	Carbon Disulfide	4.495	4.726	4.805	5.155	4.981	4.832	5.19
15)		Isopropyl Alcohol		4.726	4.805	5.155	4.981	0.000	-1.00
16)		Acetonitrile		4.726	4.805	5.155	4.981	0.000	-1.00
17)	T	Acrylonitrile	0.313	0.326	0.308	0.272	0.263	0.296	9.26
18)	T	tert-Butyl alcohol	0.085	0.098	0.081	0.071	0.063	0.080	17.35
19)	T	methyl tert-butyl e	2.801	2.949	2.826	2.690	2.641	2.782	4.35
20)	T	trans-1,2-dichloroe	1.194	1.241	1.255	1.320	1.272	1.257	3.64
21)		n-Hexane	2.078	2.101	2.154	2.178	2.084	2.119	2.10
22)	P	1,1-dichloroethane	2.211	2.292	2.274	2.353	2.205	2.267	2.71
23)	T	di-isopropyl ether	4.292	4.317	4.300	4.275	4.151	4.267	1.55
24)	T	Vinyl Acetate	2.255	2.324	2.337	2.282	1.971	2.234	6.74
25)	T	ethyl tert-butyl et	2.694	2.785	2.745	2.766	2.826	2.763	1.77
26)	T	2-Butanone	0.415	0.452	0.401	0.364	0.363	0.399	9.44
27)	T	2,2-dichloropropane	1.349	1.379	1.464	1.547	1.549	1.458	6.36
28)	T	cis-1,2-dichloroeth	0.808	0.840	0.867	0.918	0.925	0.872	5.75
29)	T	bromochloromethane	0.310	0.326	0.322	0.332	0.333	0.325	2.92
30)	C	chloroform	1.648	1.661	1.704	1.785	1.763	1.712	3.54
31)	S	Dibromofluoromethan	0.678	0.666	0.652	0.641	0.624	0.652	3.19
32)	T	Tetrahydrofuran	0.209	0.238	0.215	0.193	0.187	0.208	9.69
33)	T	1,1,1-trichloroetha	1.181	1.185	1.237	1.297	1.314	1.243	4.97
34)	I	1,4-difluorobenzene	----------ISTD----------						
35)	T	carbon tetrachlorid	0.496	0.468	0.432	0.462	0.469	0.465	4.95
36)	T	1,1-dichloropropene	0.969	0.817	0.755	0.772	0.777	0.818	10.67
37)	M	benzene	1.771	1.834	1.922	2.053	2.168	1.950	8.28
38)	T	1,2-dichloroethane	0.605	0.624	0.630	0.648	0.649	0.631	2.90

#) = Out of Range

000092

Form VI

Response Factor Report inst g

Method : C:\HPCHEM\1\DATA\051205\8260BASP.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Tue May 17 11:42:05 2005
Response via : Initial Calibration

Calibration Files
10 =G7889.D 20 =G7890.D 50 =G7891.D
100 =G7894.D 200 =G7893.D =

		Compound	10	20	50	100	200	Avg	%RSD
39)	M	tert amyl methyl et	1.041	1.069	1.105	1.092	1.152	1.092	3.78
40)	M	trichloroethene	0.364	0.390	0.398	0.429	0.449	0.406	8.20
41)	C	1,2-dichloropropane	0.415	0.424	0.443	0.470	0.482	0.447	6.49
42)	T	dibromomethane	0.225	0.239	0.231	0.239	0.245	0.236	3.27
43)		1,4-Dioxane	0.000		0.231	0.239	0.245	0.000	0.00
44)	T	bromodichloromethan	0.510	0.541	0.559	0.604	0.619	0.567	7.89
45)	T	2-Chloroethyl vinyl	0.132	0.135	0.171	0.169	0.181	0.158	14.24
46)	T	4-Methyl-2-Pentanon	0.459	0.451	0.426	0.387	0.400	0.425	7.37
47)	T	cis-1,3-dichloropro	0.642	0.675	0.715	0.760	0.800	0.719	8.82
48)	S	toluene-d8	1.236	1.223	1.218	1.241	1.223	1.228	0.80
49)	CM	toluene	0.947	0.989	1.023	1.137	1.177	1.055	9.33
50)	T	trans-1,3-dichlorop	0.558	0.586	0.620	0.648	0.681	0.619	7.90
51)	T	1,1,2-trichloroetha	0.251	0.272	0.276	0.281	0.294	0.275	5.65
52)	I	chlorobenzene-d5	----------		ISTD	----------			
53)	T	2-Hexanone	0.287	0.316	0.322	0.297	0.315	0.307	4.83
54)	T	tetrachloroethene	0.333	0.348	0.364	0.395	0.423	0.373	9.77
55)	T	1,3-dichloropropane	0.682	0.714	0.729	0.731	0.757	0.723	3.80
56)	T	dibromochloromethan	0.296	0.307	0.336	0.346	0.367	0.330	8.79
57)	T	1,2-dibromoethane	0.288	0.314	0.324	0.319	0.334	0.316	5.45
58)	PM	chlorobenzene	1.032	1.065	1.109	1.195	1.252	1.131	8.08
59)	T	1,1,1,2-tetrachloro	0.310	0.319	0.345	0.382	0.426	0.356	13.47
60)	C	ethylbenzene	2.107	2.248	2.414	2.711	2.962	2.488	13.95
61)	T	m,p-xylene	0.698	0.751	0.818	0.926	1.029	0.844	15.82
62)	T	o-xylene	0.637	0.672	0.712	0.797	0.848	0.733	11.96
63)	T	styrene	1.127	1.221	1.321	1.488	1.600	1.352	14.28
64)	P	bromoform	0.144	0.164	0.172	0.179	0.188	0.169	9.89
65)	S	4-Bromofluorobenzen	0.511	0.517	0.524	0.522	0.519	0.519	0.96
66)	I	1,4-dichlorobenzene-d	----------		ISTD	----------			
67)	T	isopropylbenzene	4.138	4.335	4.538	5.142	5.423	4.715	11.57
68)	T	bromobenzene	0.794	0.825	0.851	0.921	0.948	0.868	7.49
69)	P	1,1,2,2-tetrachloro	0.983	1.056	1.025	0.944	1.035	1.008	4.45
70)	T	1,2,3-trichloroprop	0.783	0.856	0.814	0.788	0.835	0.815	3.79
71)	T	n-propylbenzene	5.990	6.213	6.480	7.327	7.862	6.774	11.68
72)	T	2-chlorotoluene	3.205	3.219	3.328	3.688	3.808	3.450	8.11
73)	T	4-chlorotoluene	3.703	3.903	4.043	4.557	4.860	4.213	11.39
74)	T	1,3,5-trimethylbenz	3.223	3.381	3.590	4.089	4.415	3.740	13.35
75)	T	tert-butylbenzene	2.053	2.131	2.271	2.544	2.616	2.323	10.70
76)	T	1,2,4-trimethylbenz	3.151	3.351	3.535	3.979	4.199	3.643	11.97

#) = Out of Range

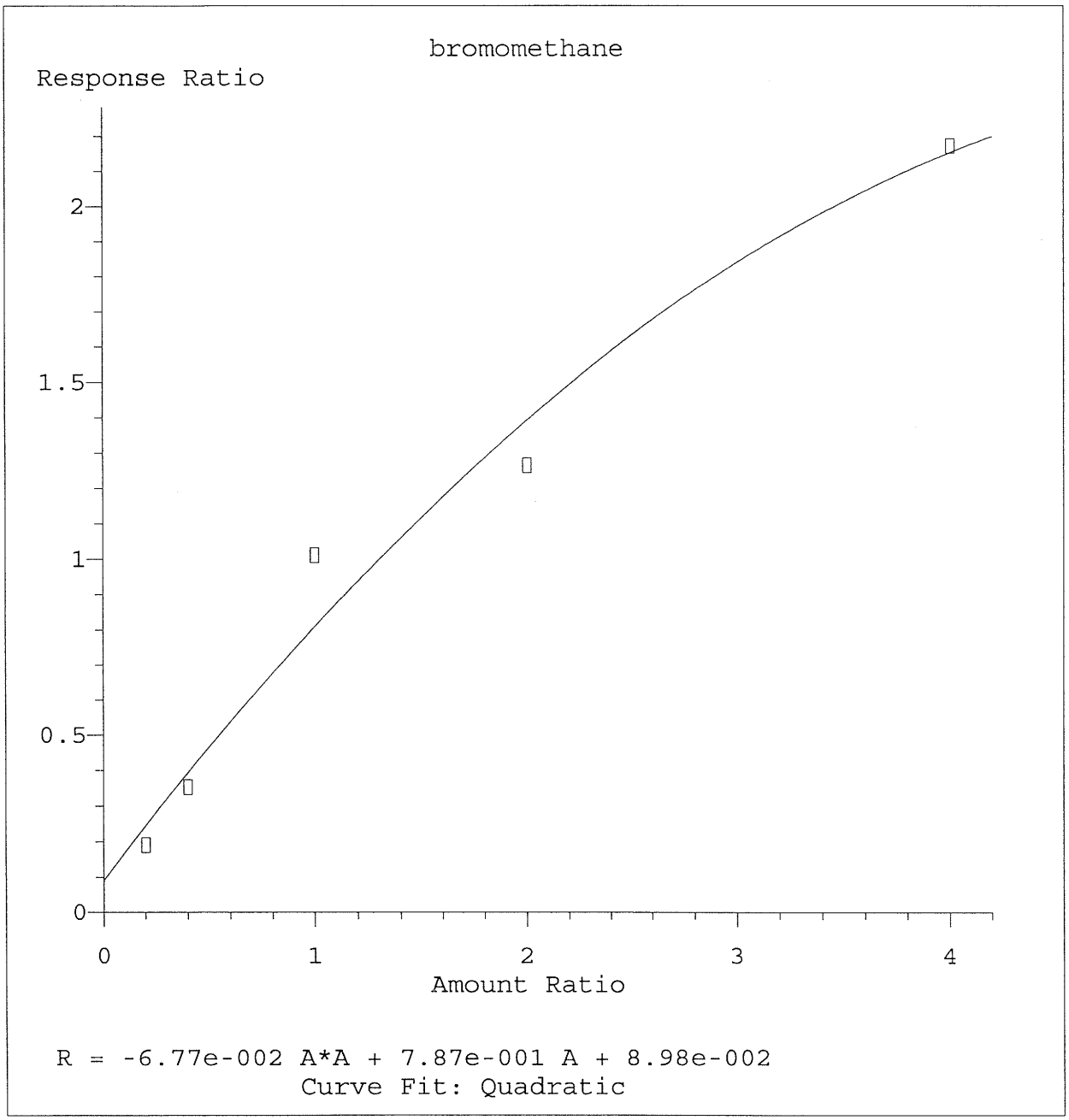
Response Factor Report inst g

Method : C:\HPCHEM\1\DATA\051205\8260BASP.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Tue May 17 11:42:05 2005
Response via : Initial Calibration

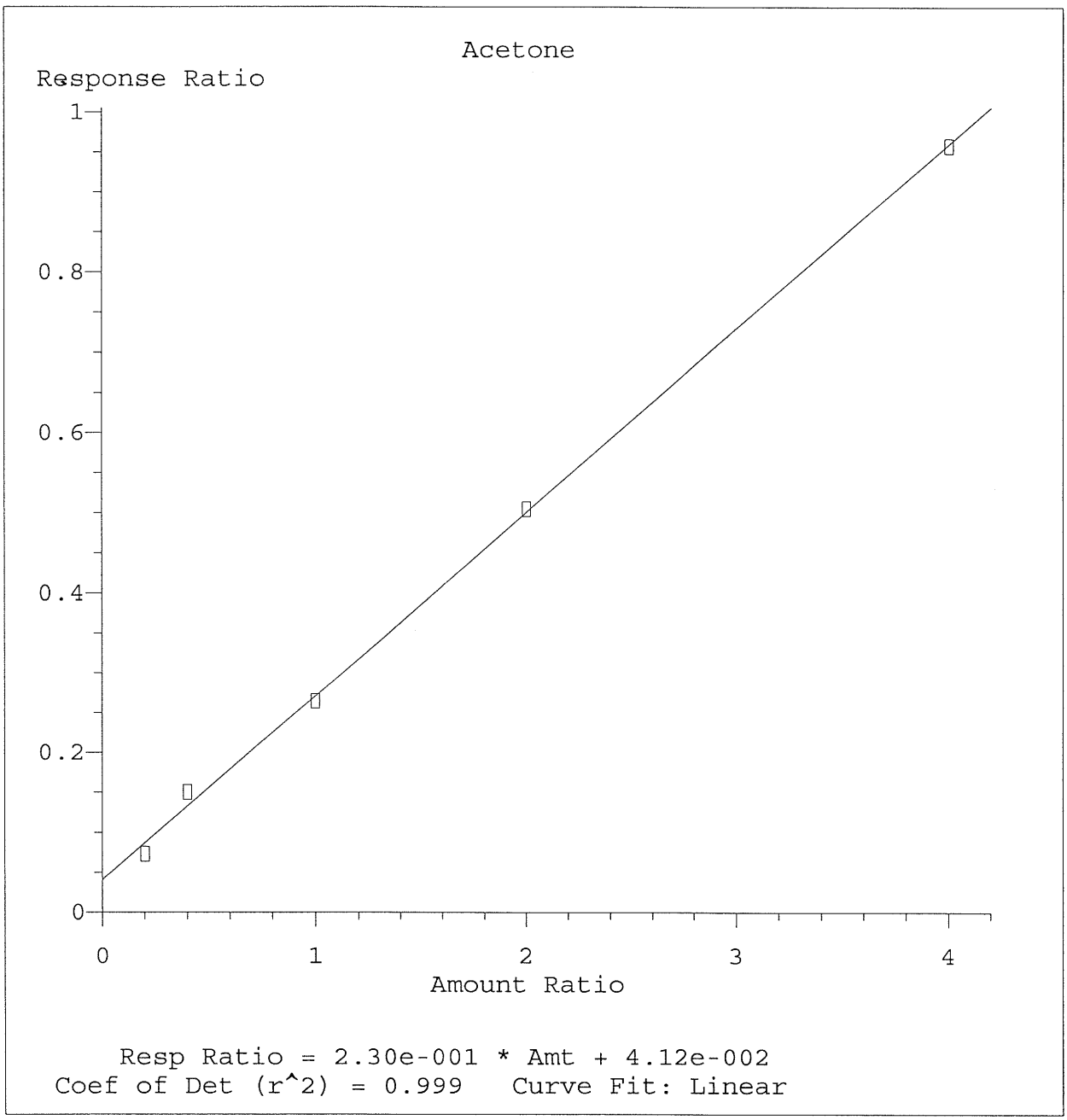
Calibration Files
10 =G7889.D 20 =G7890.D 50 =G7891.D
100 =G7894.D 200 =G7893.D =

		Compound	10	20	50	100	200	Avg	%RSD
77)	T	sec-butylbenzene	4.471	4.637	5.028	5.672	6.029	5.167	12.92
78)	T	1,3-dichlorobenzene	1.647	1.688	1.775	1.954	2.073	1.827	9.92
79)	T	4-isopropyltoluene	3.440	3.500	3.864	4.389	4.701	3.979	13.89
80)	T	1,4-dichlorobenzene	1.711	1.764	1.800	1.962	2.078	1.863	8.17
81)	T	1,2-dichlorobenzene	1.462	1.554	1.620	1.770	1.917	1.664	10.85
82)	T	n-butylbenzene	4.158	4.198	4.647	5.368	5.698	4.814	14.42
83)	T	1,2-dibromo-3-chlor	0.123	0.146	0.142	0.124	0.140	0.135	7.99
84)	T	1,2,4-trichlorobenz	0.796	0.828	0.882	0.960	1.049	0.903	11.36
85)	T	hexachlorobutadiene	0.363	0.348	0.391	0.456	0.488	0.409	14.77
86)	T	naphthalene	1.844	2.051	2.156	1.946	2.250	2.049	7.88
87)	T	1,2,3-trichlorobenz	0.702	0.730	0.777	0.782	0.877	0.774	8.63

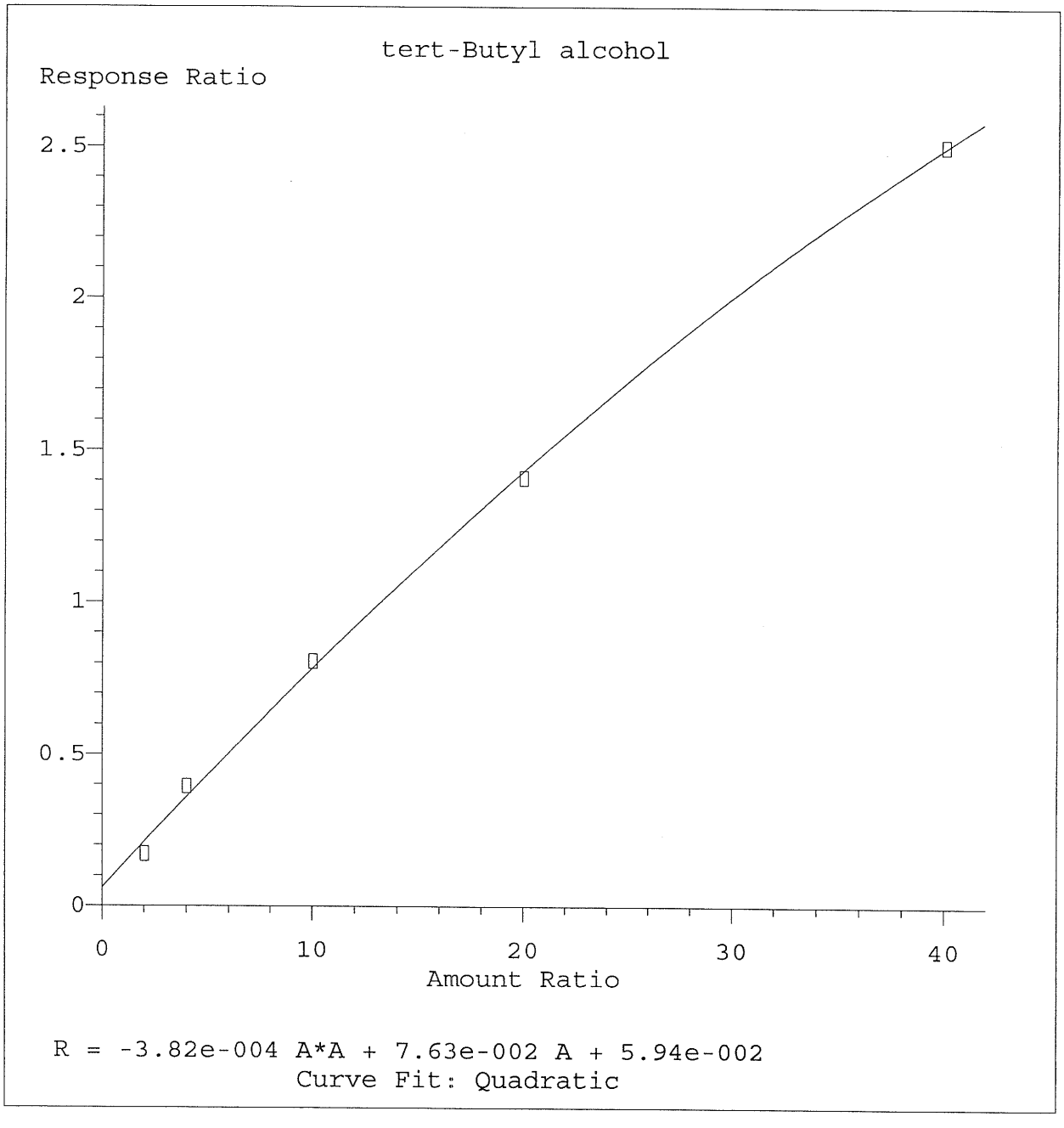
(#) = Out of Range



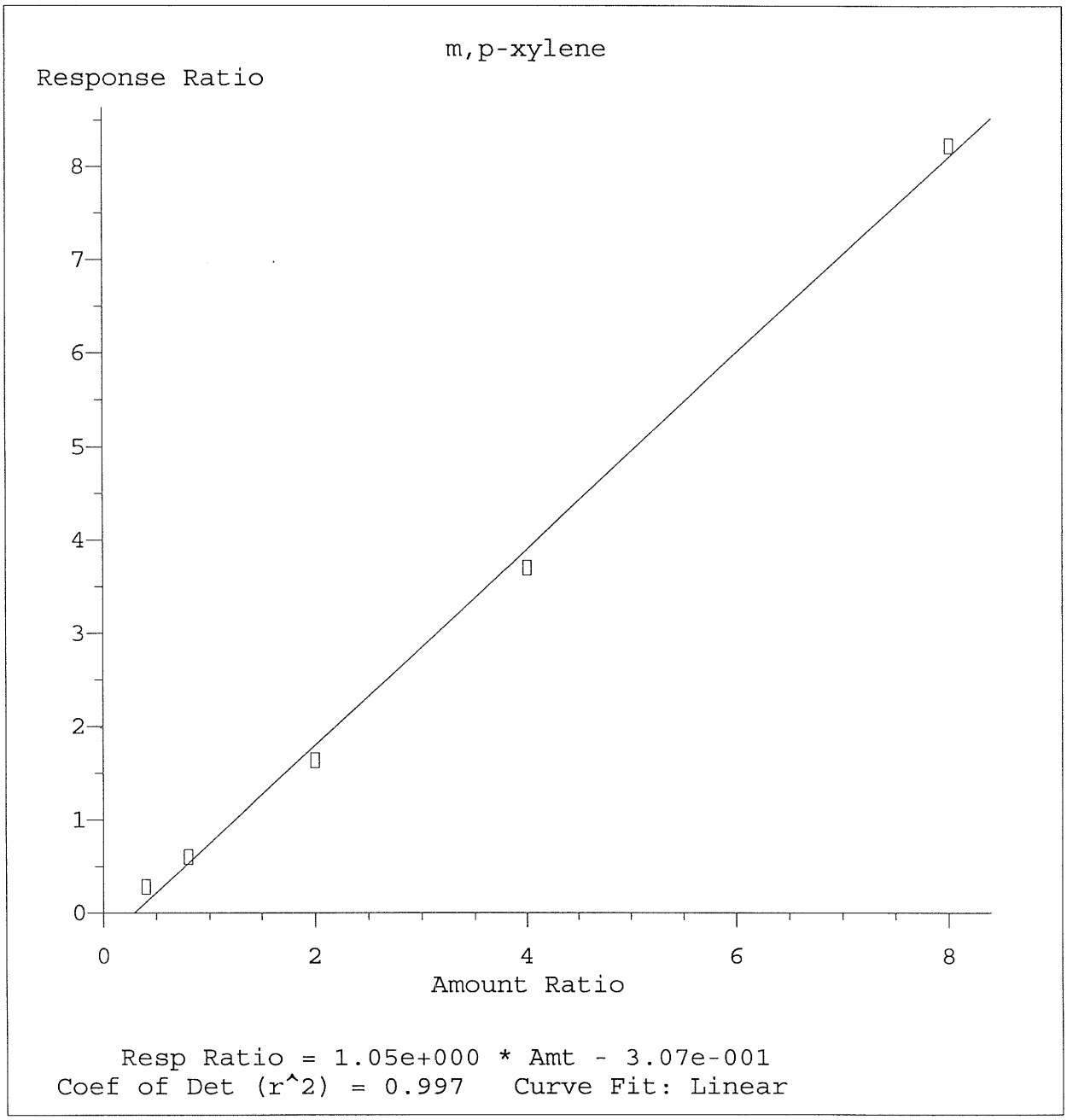
Method Name: C:\HPCHEM\1\DATA\051205\8260BASP.M
Calibration Table Last Updated: Fri May 20 09:55:33 2005



Method Name: C:\HPCHEM\1\DATA\051205\8260BASP.M
Calibration Table Last Updated: Fri May 20 09:52:41 2005



Method Name: C:\HPCHEM\1\DATA\051205\8260BASP.M
Calibration Table Last Updated: Fri May 20 09:53:26 2005



Method Name: C:\HPCHEM\1\DATA\051205\8260BASP.M
Calibration Table Last Updated: Fri May 20 09:54:22 2005

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\051205\G7888.D Vial: 2
Acq On : 12 May 05 9:10 Operator: XL
Sample : 5PPB 8260 STD Inst : inst g
Misc : Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 12 9:59 19105 Quant Results File: 8260BS.RES

Quant Method : C:\HPCHEM\1\DATA\051205\8260BS.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Tue May 03 11:15:26 2005
Response via : Initial Calibration
DataAcq Meth : 8260BS

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) pentafluorobenzene	7.25	168	332903	50.00	ug/L	-0.03
34) 1,4-difluorobenzene	8.36	114	702070	50.00	ug/L	-0.03
52) chlorobenzene-d5	13.14	117	589700	50.00	ug/L	-0.03
66) 1,4-dichlorobenzene-d4	17.27	152	250782	50.00	ug/L	-0.03
System Monitoring Compounds						
31) Dibromofluoromethane	7.10	113	222006	53.77	ug/L	-0.03
Spiked Amount 50.000	Range 86 - 118		Recovery = 107.54%			
48) toluene-d8	10.68	98	843224	48.64	ug/L	-0.03
Spiked Amount 50.000	Range 88 - 110		Recovery = 97.28%			
65) 4-Bromofluorobenzene	15.23	95	296832	49.15	ug/L	-0.03
Spiked Amount 50.000	Range 86 - 115		Recovery = 98.30%			
Target Compounds						Qvalue
2) dichlorodifluoromethane	1.54	85	30133	4.34	ug/L	100
3) chloromethane	1.74	50	48866	4.70	ug/L	99
4) vinyl chloride	1.85	62	52993	4.41	ug/L	97
5) bromomethane	2.22	96	29405	4.14	ug/L	93
6) chloroethane	2.34	64	30359	4.56	ug/L	96
7) trichlorofluoromethane	2.62	101	52837	4.45	ug/L	98
8) Diethyl Ether	3.01	45	16289	5.03	ug/L	95
9) Acrolein	3.27	56	15572	26.58	ug/L	97
10) Acetone	3.51	43	15947	1.35	ug/L	99
11) 1,1-dichloroethene	3.31	96	30928	4.22	ug/L	86
12) Iodomethane	3.53	142	49551	4.78	ug/L	91
13) methylene chloride	4.15	84	39954	4.96	ug/L	90
14) Carbon Disulfide	3.57	76	130398	4.17	ug/L	87
17) Acrylonitrile	4.68	53	9066	5.34	ug/L	83
18) tert-Butyl alcohol	4.48	59	22981	52.90	ug/L	100
19) methyl tert-butyl ether	4.53	73	78905	4.76	ug/L	100
20) trans-1,2-dichloroethene	4.53	96	35903	4.32	ug/L	90
22) 1,1-dichloroethane	5.30	63	64049	4.40	ug/L	97
23) di-isopropyl ether	5.37	45	129105	4.76	ug/L	98
24) Vinyl Acetate	5.42	43	102634	7.70	ug/L	96
25) ethyl tert-butyl ether	5.98	59	74947	4.61	ug/L	98
26) 2-Butanone	6.38	43	11813	5.33	ug/L	94
27) 2,2-dichloropropane	6.21	77	41062	4.55	ug/L	100

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\051205\G7889.D
Acq On : 12 May 05 9:42
Sample : vstd010
Misc : ical
MS Integration Params: rteint.p
Quant Time: May 12 14:38 19105

Vial: 3
Operator: XL
Inst : inst g
Multiplr: 1.00

Quant Results File: 8260BS.RES

Quant Method : C:\HPCHEM\1\DATA\051205\8260BS.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Tue May 03 11:15:26 2005
Response via : Initial Calibration
DataAcq Meth : 8260BS

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) pentafluorobenzene	7.24	168	317756	50.00	ug/L	-0.03
34) 1,4-difluorobenzene	8.36	114	662478	50.00	ug/L	-0.03
52) chlorobenzene-d5	13.14	117	574458	50.00	ug/L	-0.03
66) 1,4-dichlorobenzene-d4	17.27	152	251544	50.00	ug/L	-0.02
System Monitoring Compounds						
31) Dibromofluoromethane	7.10	113	215392	54.66	ug/L	-0.03
Spiked Amount 50.000	Range 86 - 118		Recovery = 109.32%			
48) toluene-d8	10.68	98	819071	50.07	ug/L	-0.02
Spiked Amount 50.000	Range 88 - 110		Recovery = 100.14%			
65) 4-Bromofluorobenzene	15.22	95	293786	49.94	ug/L	-0.03
Spiked Amount 50.000	Range 86 - 115		Recovery = 99.88%			
Target Compounds						Qvalue
2) dichlorodifluoromethane	1.54	85	64005	9.66	ug/L	99
3) chloromethane	1.74	50	107684	10.85	ug/L	97
4) vinyl chloride	1.85	62	113089	9.85	ug/L	98
5) bromomethane	2.21	96	60493	8.93	ug/L	97
6) chloroethane	2.34	64	63855	10.05	ug/L	92
7) trichlorofluoromethane	2.61	101	112342	9.91	ug/L	98
8) Diethyl Ether	3.02	45	34765	11.26	ug/L	95
9) Acrolein	3.26	56	33129	59.25	ug/L	88
10) Acetone	3.51	43	23272	8.72	ug/L	99
11) 1,1-dichloroethene	3.31	96	68058	9.72	ug/L	79
12) Iodomethane	3.53	142	99942	10.10	ug/L	89
13) methylene chloride	4.15	84	79992	10.40	ug/L	98
14) Carbon Disulfide	3.57	76	285673	9.57	ug/L	98
16) Acetonitrile	3.94	41	542	1.18	ug/L	99
17) Acrylonitrile	4.69	53	19909	12.29	ug/L	87
18) tert-Butyl alcohol	4.48	59	54287	130.92	ug/L	100
19) methyl tert-butyl ether	4.53	73	178032	11.26	ug/L	100
20) trans-1,2-dichloroethene	4.53	96	75911	9.56	ug/L	88
22) 1,1-dichloroethane	5.30	63	140510	10.12	ug/L	98
23) di-isopropyl ether	5.38	45	272730	10.53	ug/L	94
24) Vinyl Acetate	5.42	43	143296m	11.26	ug/L	96
25) ethyl tert-butyl ether	5.97	59	171177	11.04	ug/L	97
26) 2-Butanone	6.39	43	26400	12.49	ug/L	88

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\051205\G7889.D
Acq On : 12 May 05 9:42
Sample : vstd010
Misc : ical
MS Integration Params: rteint.p
Quant Time: May 12 14:38 19105

Vial: 3
Operator: XL
Inst : inst g
Multiplr: 1.00

Quant Results File: 8260BS.RES

Quant Method : C:\HPCHEM\1\DATA\051205\8260BS.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Tue May 03 11:15:26 2005
Response via : Initial Calibration
DataAcq Meth : 8260BS

	Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
27)	2,2-dichloropropane	6.21	77	85699	9.95	ug/L	94
28)	cis-1,2-dichloroethene	6.29	96	51377	9.98	ug/L	88
29)	bromochloromethane	6.68	128	19678m	10.57	ug/L	88
30)	chloroform	6.84	83	104715	10.53	ug/L	94
32)	Tetrahydrofuran	6.71	42	13272	12.26	ug/L	91
33)	1,1,1-trichloroethane	7.01	97	75081	10.19	ug/L	95
35)	carbon tetrachloride	7.23	117	65752m	10.60	ug/L	99
36)	1,1-dichloropropene	7.29	75	128350	12.56	ug/L	96
37)	benzene	7.60	78	234702	9.58	ug/L	96
38)	1,2-dichloroethane	7.77	62	80140	10.14	ug/L	99
39)	tert amyl methyl ether	7.82	73	137978	10.65	ug/L	97
40)	trichloroethene	8.69	95	48238m	9.14	ug/L	86
41)	1,2-dichloropropane	9.12	63	55037	9.98	ug/L	100
42)	dibromomethane	9.31	93	29839	10.55	ug/L	95
44)	bromodichloromethane	9.60	83	67620	9.51	ug/L	99
45)	2-Chloroethyl vinyl ether	10.15	63	17488	11.05	ug/L	93
46)	4-Methyl-2-Pentanone	10.63	43	60791	12.39	ug/L	84
47)	cis-1,3-dichloropropene	10.35	75	85091	9.80	ug/L	96
49)	toluene	10.79	92	125496	9.35	ug/L	86
50)	trans-1,3-dichloropropene	11.33	75	73891	10.16	ug/L	94
51)	1,1,2-trichloroethane	11.61	83	33256	10.20	ug/L	96
53)	2-Hexanone	12.03	43	32963	11.23	ug/L	99
54)	tetrachloroethene	11.65	166	38205	9.02	ug/L	99
55)	1,3-dichloropropane	11.89	76	78334	10.33	ug/L	90
56)	dibromochloromethane	12.20	129	33977	9.69	ug/L	95
57)	1,2-dibromoethane	12.37	107	33100	10.01	ug/L	97
58)	chlorobenzene	13.19	112	118536	9.52	ug/L	95
59)	1,1,1,2-tetrachloroethane	13.37	131	35583	9.40	ug/L	96
60)	ethylbenzene	13.36	91	242113	8.99	ug/L	93
61)	m,p-xylene	13.58	106	160354	17.45	ug/L	96
62)	o-xylene	14.25	106	73134	9.25	ug/L	96
63)	styrene	14.31	104	129511	8.96	ug/L	99
64)	bromoform	14.63	173	16543	11.03	ug/L	88
67)	isopropylbenzene	14.89	105	208197	8.93	ug/L	99
68)	bromobenzene	15.44	156	39928m	9.39	ug/L	89
69)	1,1,2,2-tetrachloroethane	15.59	83	49442m	10.86	ug/L	96
70)	1,2,3-trichloropropane	15.64	75	39395	10.22	ug/L	97

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\051205\G7889.D
Acq On : 12 May 05 9:42
Sample : vstd010
Misc : ical
MS Integration Params: rteint.p
Quant Time: May 12 14:38 19105

Vial: 3
Operator: XL
Inst : inst g
Multiplr: 1.00

Quant Results File: 8260BS.RES

Quant Method : C:\HPCHEM\1\DATA\051205\8260BS.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Tue May 03 11:15:26 2005
Response via : Initial Calibration
DataAcq Meth : 8260BS

	Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
71)	n-propylbenzene	15.63	91	301328	9.12	ug/L	100
72)	2-chlorotoluene	15.79	91	161230	9.55	ug/L	96
73)	4-chlorotoluene	16.00	91	186306	9.04	ug/L	95
74)	1,3,5-trimethylbenzene	15.97	105	162142	8.87	ug/L	100
75)	tert-butylbenzene	16.51	91	103262	8.92	ug/L	92
76)	1,2,4-trimethylbenzene	16.63	105	158528	8.92	ug/L	100
77)	sec-butylbenzene	16.91	105	224921	8.88	ug/L	97
78)	1,3-dichlorobenzene	17.13	146	82842	9.39	ug/L	97
79)	4-isopropyltoluene	17.19	119	173066	8.95	ug/L	98
80)	1,4-dichlorobenzene	17.31	146	86078	9.59	ug/L	95
81)	1,2-dichlorobenzene	17.96	146	73535	9.20	ug/L	96
82)	n-butylbenzene	17.93	91	209180	9.00	ug/L	94
83)	1,2-dibromo-3-chloropropan	19.42	75	6186m	10.64	ug/L	78
84)	1,2,4-trichlorobenzene	20.87	180	40068	9.18	ug/L	87
85)	hexachlorobutadiene	21.14	225	18287	8.86	ug/L	97
86)	naphthalene	21.31	128	92749	9.59	ug/L	100
87)	1,2,3-trichlorobenzene	21.76	180	35315	9.38	ug/L	96

(#) = qualifier out of range (m) = manual integration

000102

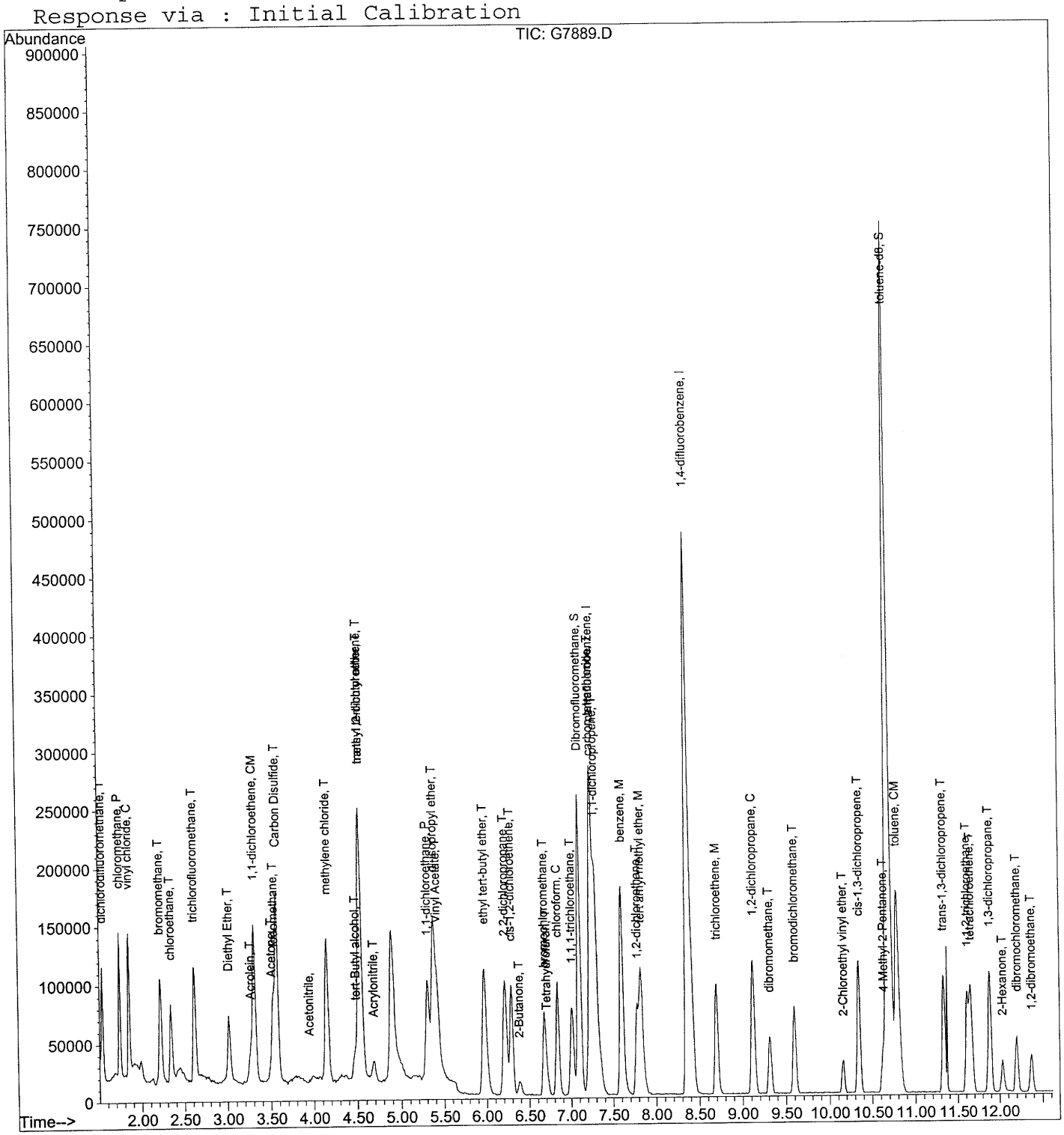
Quantitation Report

Data File : C:\HPCHEM\1\DATA\051205\G7889.D
Acq On : 12 May 05 9:42
Sample : vstd010
Misc : ical
MS Integration Params: rteint.p
Quant Time: May 12 14:38 19105

Vial: 3
Operator: XL
Inst : inst g
Multiplr: 1.00

Quant Results File: 8260BS.RES

Method : C:\HPCHEM\1\DATA\051705\8260BASP.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Fri May 20 09:55:33 2005
Response via : Initial Calibration



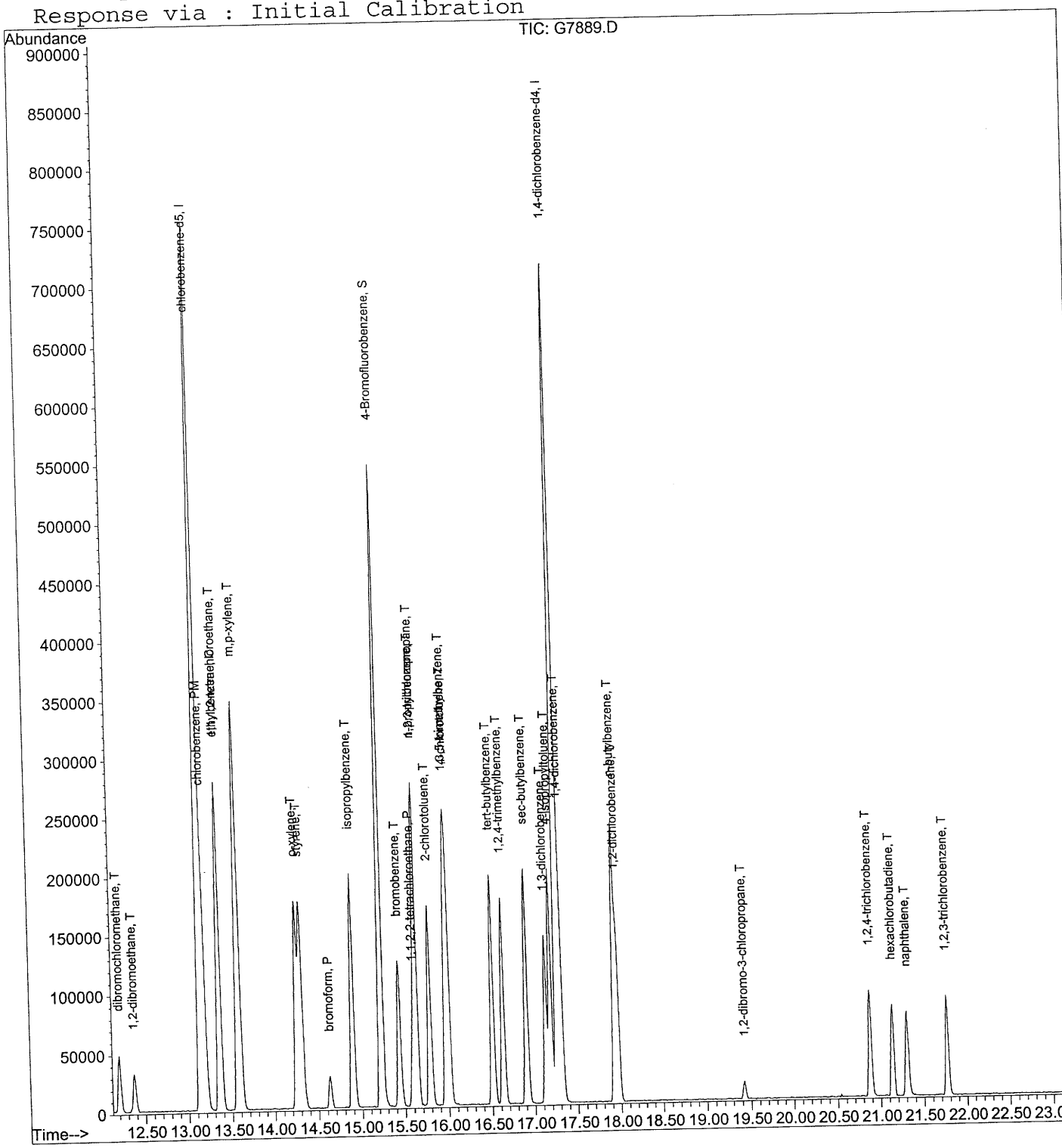
Quantitation Report

Data File : C:\HPCHEM\1\DATA\051205\G7889.D
Acq On : 12 May 05 9:42
Sample : vstd010
Misc : ical
MS Integration Params: rteint.p
Quant Time: May 12 14:38 19105

Vial: 3
Operator: XL
Inst : inst g
Multiplr: 1.00

Quant Results File: 8260BS.RES

Method : C:\HPCHEM\1\DATA\051705\8260BASP.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Fri May 20 09:55:33 2005
Response via : Initial Calibration



000104

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\051205\G7890.D
Acq On : 12 May 05 10:14
Sample : vstd020
Misc : ical
MS Integration Params: rteint.p
Quant Time: May 12 11:03 19105

Vial: 4
Operator: XL
Inst : inst g
Multiplr: 1.00

Quant Results File: 8260BS.RES

Quant Method : C:\HPCHEM\1\DATA\051205\8260BS.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Tue May 03 11:15:26 2005
Response via : Initial Calibration
DataAcq Meth : 8260BS

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) pentafluorobenzene	7.24	168	300171	50.00	ug/L	-0.03
34) 1,4-difluorobenzene	8.36	114	629579	50.00	ug/L	-0.03
52) chlorobenzene-d5	13.15	117	541713	50.00	ug/L	-0.02
66) 1,4-dichlorobenzene-d4	17.27	152	241714	50.00	ug/L	-0.02
System Monitoring Compounds						
31) Dibromofluoromethane	7.10	113	199776	53.66	ug/L	-0.03
Spiked Amount 50.000	Range 86 - 118		Recovery = 107.32%			
48) toluene-d8	10.68	98	769812	49.51	ug/L	-0.02
Spiked Amount 50.000	Range 88 - 110		Recovery = 99.02%			
65) 4-Bromofluorobenzene	15.23	95	279965	50.47	ug/L	-0.02
Spiked Amount 50.000	Range 86 - 115		Recovery = 100.94%			
Target Compounds						Qvalue
2) dichlorodifluoromethane	1.54	85	128431	20.53	ug/L	96
3) chloromethane	1.74	50	216754	23.11	ug/L	100
4) vinyl chloride	1.85	62	225311	20.77	ug/L	99
5) bromomethane	2.21	96	105912	16.55	ug/L	99
6) chloroethane	2.34	64	122989	20.48	ug/L	93
7) trichlorofluoromethane	2.62	101	224216	20.95	ug/L	99
8) Diethyl Ether	3.01	45	60102	20.60	ug/L	97
9) Acrolein	3.27	56	54390	102.97	ug/L	98
10) Acetone	3.50	43	45150	30.40	ug/L	99
11) 1,1-dichloroethene	3.31	96	135512	20.48	ug/L	89
12) Iodomethane	3.53	142	191302	20.46	ug/L	91
13) methylene chloride	4.15	84	157178	21.62	ug/L	98
14) Carbon Disulfide	3.57	76	567404	20.11	ug/L	93
17) Acrylonitrile	4.68	53	39179	25.61	ug/L	95
18) tert-Butyl alcohol	4.47	59	118189	301.73	ug/L	100
19) methyl tert-butyl ether	4.53	73	354110	23.71	ug/L	100
20) trans-1,2-dichloroethene	4.53	96	148999	19.86	ug/L	88
22) 1,1-dichloroethane	5.30	63	275168	20.97	ug/L	97
23) di-isopropyl ether	5.38	45	518286	21.18	ug/L	93
24) Vinyl Acetate	5.42	43	279006m	23.21	ug/L	100
25) ethyl tert-butyl ether	5.98	59	334355	22.83	ug/L	98
26) 2-Butanone	6.39	43	54330	27.21	ug/L	97
27) 2,2-dichloropropane	6.21	77	165612	20.36	ug/L	99

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\051205\G7890.D
Acq On : 12 May 05 10:14
Sample : vstd020
Misc : ical
MS Integration Params: rteint.p
Quant Time: May 12 11:03 19105

Vial: 4
Operator: XL
Inst : inst g
Multiplr: 1.00

Quant Results File: 8260BS.RES

Quant Method : C:\HPCHEM\1\DATA\051205\8260BS.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Tue May 03 11:15:26 2005
Response via : Initial Calibration
DataAcq Meth : 8260BS

Compound	R.T.	QIon	Response	Conc Unit	Qvalue
28) cis-1,2-dichloroethene	6.30	96	100858	20.74 ug/L	93
29) bromochloromethane	6.68	128	39119	22.24 ug/L	92
30) chloroform	6.84	83	199378	21.22 ug/L	98
32) Tetrahydrofuran	6.71	42	28597	27.97 ug/L	94
33) 1,1,1-trichloroethane	7.01	97	142230	20.43 ug/L	92
35) carbon tetrachloride	7.22	117	117889	20.84 ug/L	98
36) 1,1-dichloropropene	7.29	75	205799	21.20 ug/L	95
37) benzene	7.60	78	461911	19.84 ug/L	99
38) 1,2-dichloroethane	7.77	62	157262	20.94 ug/L	99
39) tert amyl methyl ether	7.82	73	269306	21.88 ug/L	98
40) trichloroethene	8.70	95	98250	19.60 ug/L	95
41) 1,2-dichloropropane	9.12	63	106693	20.35 ug/L	100
42) dibromomethane	9.31	93	60114	22.37 ug/L	96
44) bromodichloromethane	9.59	83	136168	20.15 ug/L	98
45) 2-Chloroethyl vinyl ether	10.15	63	34122	23.20 ug/L	92
46) 4-Methyl-2-Pentanone	10.62	43	113655	24.38 ug/L	82
47) cis-1,3-dichloropropene	10.35	75	170085	20.62 ug/L	100
49) toluene	10.78	92	249009	19.53 ug/L	98
50) trans-1,3-dichloropropene	11.32	75	147638	21.37 ug/L	94
51) 1,1,2-trichloroethane	11.61	83	68502	22.10 ug/L	97
53) 2-Hexanone	12.03	43	68444	24.74 ug/L	91
54) tetrachloroethene	11.65	166	75465	18.88 ug/L	98
55) 1,3-dichloropropane	11.89	76	154819	21.65 ug/L	95
56) dibromochloromethane	12.20	129	66443	20.10 ug/L	98
57) 1,2-dibromoethane	12.37	107	68115	21.84 ug/L	95
58) chlorobenzene	13.19	112	230840	19.65 ug/L	98
59) 1,1,1,2-tetrachloroethane	13.37	131	69112	19.36 ug/L	97
60) ethylbenzene	13.36	91	487009	19.18 ug/L	92
61) m,p-xylene	13.58	106	325599	37.57 ug/L	99
62) o-xylene	14.26	106	145594	19.53 ug/L	93
63) styrene	14.31	104	264574	19.40 ug/L	95
64) bromoform	14.63	173	35526	24.27 ug/L	95
67) isopropylbenzene	14.89	105	419138	18.70 ug/L	98
68) bromobenzene	15.43	156	79757	19.52 ug/L	92
69) 1,1,2,2-tetrachloroethane	15.58	83	102076m	23.34 ug/L	99
70) 1,2,3-trichloropropane	15.63	75	82726	22.34 ug/L	90
71) n-propylbenzene	15.63	91	600734	18.92 ug/L	99

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\051205\G7890.D
Acq On : 12 May 05 10:14
Sample : vstd020
Misc : ical
MS Integration Params: rteint.p
Quant Time: May 12 11:03 19105

Vial: 4
Operator: XL
Inst : inst g
Multiplr: 1.00

Quant Results File: 8260BS.RES

Quant Method : C:\HPCHEM\1\DATA\051205\8260BS.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Tue May 03 11:15:26 2005
Response via : Initial Calibration
DataAcq Meth : 8260BS

Compound	R.T.	QIon	Response	Conc Unit	Qvalue
72) 2-chlorotoluene	15.79	91	311244	19.19 ug/L	98
73) 4-chlorotoluene	16.00	91	377407	19.06 ug/L	94
74) 1,3,5-trimethylbenzene	15.97	105	326855	18.62 ug/L	99
75) tert-butylbenzene	16.50	91	206077	18.53 ug/L	93
76) 1,2,4-trimethylbenzene	16.63	105	324019	18.98 ug/L	96
77) sec-butylbenzene	16.91	105	448340	18.43 ug/L	96
78) 1,3-dichlorobenzene	17.13	146	163213	19.26 ug/L	95
79) 4-isopropyltoluene	17.19	119	338378	18.22 ug/L	97
80) 1,4-dichlorobenzene	17.31	146	170533	19.77 ug/L	94
81) 1,2-dichlorobenzene	17.96	146	150237	19.57 ug/L	97
82) n-butylbenzene	17.93	91	405898	18.17 ug/L	95
83) 1,2-dibromo-3-chloropropan	19.42	75	14158	25.34 ug/L	87
84) 1,2,4-trichlorobenzene	20.88	180	80043	19.09 ug/L	91
85) hexachlorobutadiene	21.14	225	33629m	16.95 ug/L	92
86) naphthalene	21.30	128	198301	21.33 ug/L	100
87) 1,2,3-trichlorobenzene	21.77	180	70592	19.52 ug/L	94

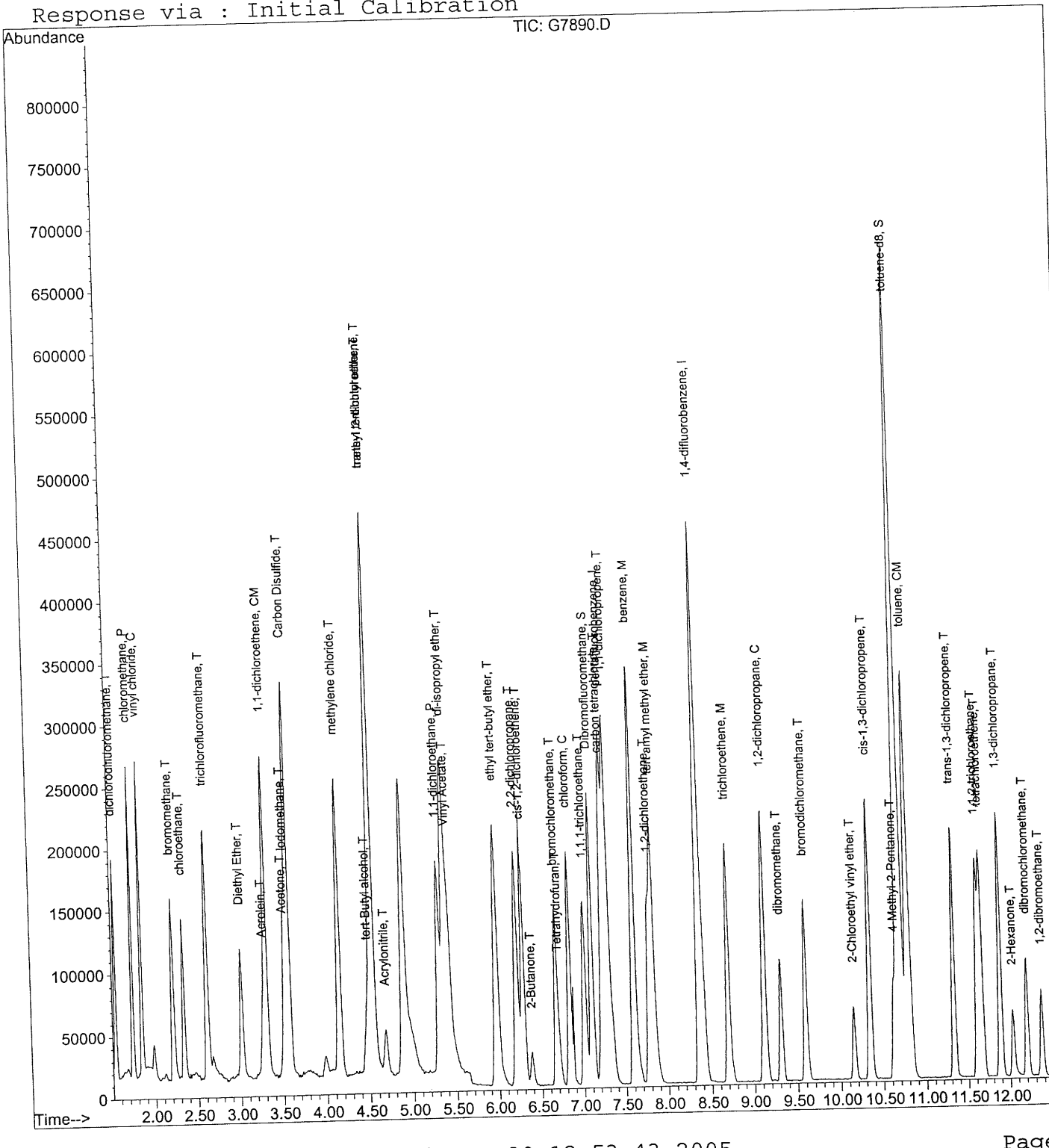
Quantitation Report

Data File : C:\HPCHEM\1\DATA\051205\G7890.D
Acq On : 12 May 05 10:14
Sample : vstd020
Misc : ical
MS Integration Params: rteint.p
Quant Time: May 12 11:03 19105

Vial: 4
Operator: XL
Inst : inst g
Multiplr: 1.00

Quant Results File: 8260BS.RES

Method : C:\HPCHEM\1\DATA\051705\8260BASP.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Fri May 20 09:55:33 2005
Response via : Initial Calibration



000108

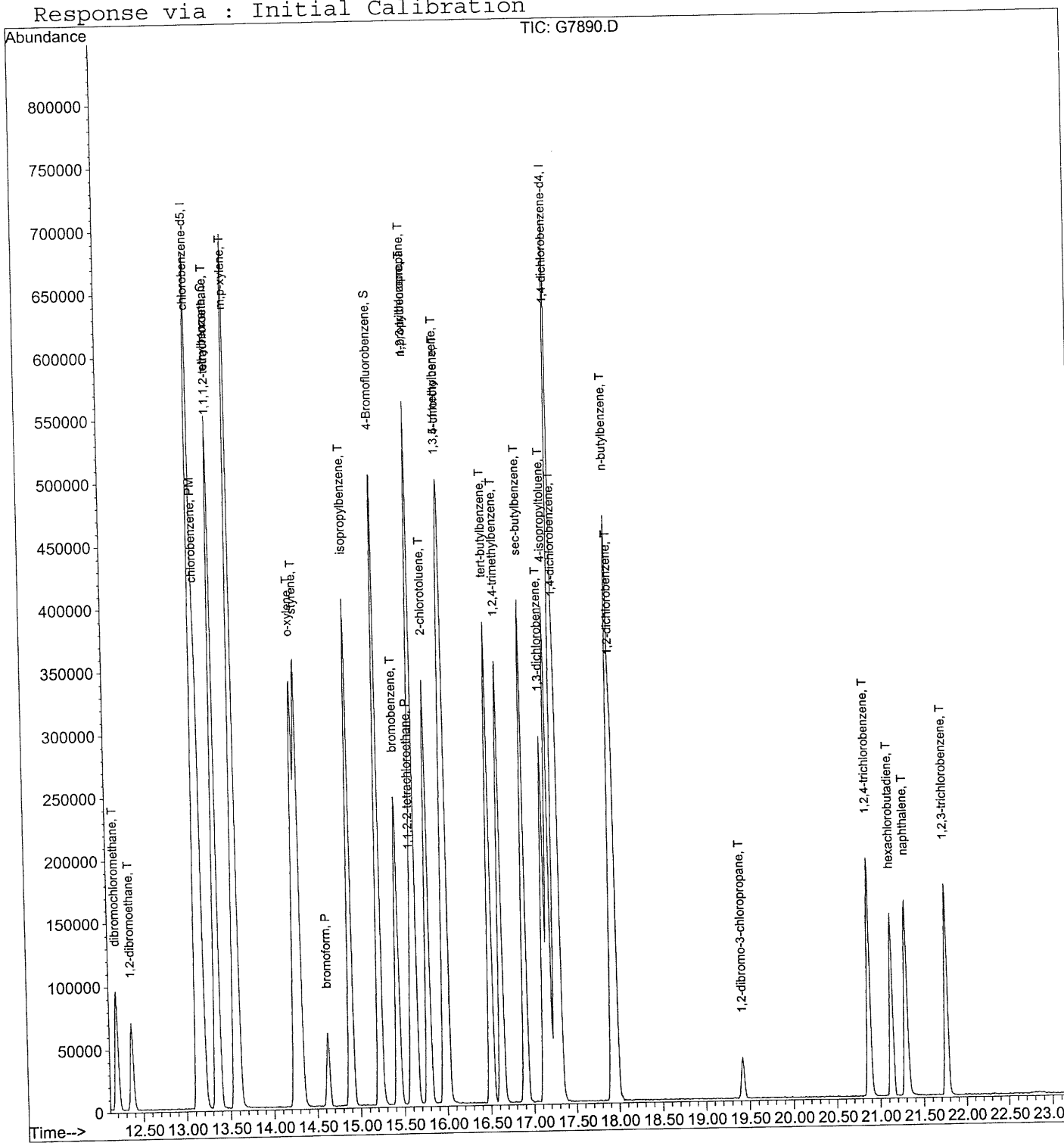
Quantitation Report

Data File : C:\HPCHEM\1\DATA\051205\G7890.D
Acq On : 12 May 05 10:14
Sample : vstd020
Misc : ical
MS Integration Params: rteint.p
Quant Time: May 12 11:03 19105

Vial: 4
Operator: XL
Inst : inst g
Multiplr: 1.00

Quant Results File: 8260BS.RES

Method : C:\HPCHEM\1\DATA\051705\8260BASP.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Fri May 20 09:55:33 2005
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\051205\G7891.D
Acq On : 12 May 05 10:46
Sample : vstd050
Misc : ical
MS Integration Params: rteint.p
Quant Time: May 12 11:48 19105

Vial: 2
Operator: XL
Inst : inst g
Multiplr: 1.00

Quant Results File: 8260BS.RES

Quant Method : C:\HPCHEM\1\DATA\051205\8260BS.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Tue May 03 11:15:26 2005
Response via : Initial Calibration
DataAcq Meth : 8260BS

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) pentafluorobenzene	7.24	168	312366	50.00	ug/L	-0.03
34) 1,4-difluorobenzene	8.37	114	653985	50.00	ug/L	-0.03
52) chlorobenzene-d5	13.14	117	551428	50.00	ug/L	-0.03
66) 1,4-dichlorobenzene-d4	17.27	152	254839	50.00	ug/L	-0.03
System Monitoring Compounds						
31) Dibromofluoromethane	7.10	113	203743	52.59	ug/L	-0.03
Spiked Amount 50.000	Range 86 - 118		Recovery = 105.18%			
48) toluene-d8	10.68	98	796690	49.33	ug/L	-0.03
Spiked Amount 50.000	Range 88 - 110		Recovery = 98.66%			
65) 4-Bromofluorobenzene	15.22	95	289045	51.18	ug/L	-0.03
Spiked Amount 50.000	Range 86 - 115		Recovery = 102.36%			
Target Compounds						Qvalue
2) dichlorodifluoromethane	1.53	85	333378	51.20	ug/L	97
3) chloromethane	1.74	50	550677	56.42	ug/L	99
4) vinyl chloride	1.85	62	585299	51.85	ug/L	99
5) bromomethane	2.22	96	315877	47.42	ug/L	97
6) chloroethane	2.33	64	322258	51.58	ug/L	92
7) trichlorofluoromethane	2.61	101	571018	51.26	ug/L	99
8) Diethyl Ether	3.01	45	139933	46.10	ug/L	92
9) Acrolein	3.26	56	156986	285.60	ug/L	95
10) Acetone	3.50	43	82658	60.66	ug/L	88
11) 1,1-dichloroethene	3.30	96	349285	50.74	ug/L	86
12) Iodomethane	3.52	142	497895	51.18	ug/L	90
13) methylene chloride	4.15	84	431794	57.08	ug/L	95
14) Carbon Disulfide	3.57	76	1500934	51.13	ug/L	92
17) Acrylonitrile	4.69	53	96071	60.34	ug/L	92
18) tert-Butyl alcohol	4.48	59	251837	617.82	ug/L	100
19) methyl tert-butyl ether	4.53	73	882652	56.79	ug/L	100
20) trans-1,2-dichloroethene	4.53	96	392053m	50.23	ug/L	86
22) 1,1-dichloroethane	5.30	63	710168	52.01	ug/L	97
23) di-isopropyl ether	5.37	45	1343039	52.74	ug/L	93
24) Vinyl Acetate	5.42	43	730092m	58.37	ug/L	99
25) ethyl tert-butyl ether	5.97	59	857518	56.27	ug/L	99
26) 2-Butanone	6.38	43	125144	60.23	ug/L	94
27) 2,2-dichloropropane	6.22	77	457439	54.04	ug/L	98

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\051205\G7891.D
Acq On : 12 May 05 10:46
Sample : vstd050
Misc : ical
MS Integration Params: rteint.p
Quant Time: May 12 11:48 19105

Vial: 2
Operator: XL
Inst : inst g
Multiplr: 1.00

Quant Results File: 8260BS.RES

Quant Method : C:\HPCHEM\1\DATA\051205\8260BS.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Tue May 03 11:15:26 2005
Response via : Initial Calibration
DataAcq Meth : 8260BS

	Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
28)	cis-1,2-dichloroethene	6.29	96	270730	53.51	ug/L	91
29)	bromochloromethane	6.68	128	100666	54.99	ug/L	98
30)	chloroform	6.83	83	532324	54.43	ug/L	98
32)	Tetrahydrofuran	6.70	42	67163	63.12	ug/L	97
33)	1,1,1-trichloroethane	7.01	97	386329	53.33	ug/L	96
35)	carbon tetrachloride	7.21	117	282267	49.28	ug/L	98
36)	1,1-dichloropropene	7.29	75	493652	48.95	ug/L	95
37)	benzene	7.60	78	1256900	51.97	ug/L	98
38)	1,2-dichloroethane	7.78	62	412049	52.83	ug/L	99
39)	tert amyl methyl ether	7.82	73	722538	56.50	ug/L	94
40)	trichloroethene	8.70	95	260324	49.99	ug/L	93
41)	1,2-dichloropropane	9.12	63	289737	53.20	ug/L	100
42)	dibromomethane	9.31	93	151169	54.15	ug/L	99
44)	bromodichloromethane	9.60	83	365830	52.12	ug/L	97
45)	2-Chloroethyl vinyl ether	10.15	63	112066	66.78	ug/L	98
46)	4-Methyl-2-Pentanone	10.62	43	278618	57.52	ug/L	81
47)	cis-1,3-dichloropropene	10.34	75	467902	54.61	ug/L	99
49)	toluene	10.79	92	668892	50.50	ug/L	93
50)	trans-1,3-dichloropropene	11.33	75	405547	56.51	ug/L	96
51)	1,1,2-trichloroethane	11.61	83	180428	56.05	ug/L	94
53)	2-Hexanone	12.04	43	177681	63.09	ug/L	97
54)	tetrachloroethene	11.66	166	200807	49.36	ug/L	98
55)	1,3-dichloropropane	11.88	76	402235	55.27	ug/L	93
56)	dibromochloromethane	12.20	129	185328	55.07	ug/L	95
57)	1,2-dibromoethane	12.36	107	178695	56.28	ug/L	96
58)	chlorobenzene	13.18	112	611660	51.15	ug/L	97
59)	1,1,1,2-tetrachloroethane	13.37	131	190320	52.36	ug/L	91
60)	ethylbenzene	13.35	91	1331390	51.51	ug/L	92
61)	m,p-xylene	13.57	106	902161	102.26	ug/L	100
62)	o-xylene	14.25	106	392784	51.77	ug/L	98
63)	styrene	14.31	104	728277	52.47	ug/L	95
64)	bromoform	14.63	173	94908	59.11	ug/L	92
67)	isopropylbenzene	14.90	105	1156550	48.94	ug/L	99
68)	bromobenzene	15.44	156	216780	50.32	ug/L	96
69)	1,1,2,2-tetrachloroethane	15.59	83	261282m	56.67	ug/L	99
70)	1,2,3-trichloropropane	15.64	75	207477	53.15	ug/L	97
71)	n-propylbenzene	15.63	91	1651457	49.32	ug/L	100

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\051205\G7891.D
Acq On : 12 May 05 10:46
Sample : vstd050
Misc : ical
MS Integration Params: rteint.p
Quant Time: May 12 11:48 19105

Vial: 2
Operator: XL
Inst : inst g
Multiplr: 1.00

Quant Results File: 8260BS.RES

Quant Method : C:\HPCHEM\1\DATA\051205\8260BS.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Tue May 03 11:15:26 2005
Response via : Initial Calibration
DataAcq Meth : 8260BS

Compound	R.T.	QIon	Response	Conc Unit	Qvalue
72) 2-chlorotoluene	15.78	91	847987	49.59 ug/L	98
73) 4-chlorotoluene	16.00	91	1030375	49.36 ug/L	94
74) 1,3,5-trimethylbenzene	15.98	105	914997	49.43 ug/L	100
75) tert-butylbenzene	16.51	91	578796	49.36 ug/L	96
76) 1,2,4-trimethylbenzene	16.63	105	900845	50.06 ug/L	99
77) sec-butylbenzene	16.90	105	1281279	49.94 ug/L	96
78) 1,3-dichlorobenzene	17.13	146	452249	50.62 ug/L	95
79) 4-isopropyltoluene	17.19	119	984672	50.28 ug/L	99
80) 1,4-dichlorobenzene	17.31	146	458725	50.45 ug/L	97
81) 1,2-dichlorobenzene	17.97	146	412732	50.99 ug/L	95
82) n-butylbenzene	17.92	91	1184168	50.29 ug/L	94
83) 1,2-dibromo-3-chloropropan	19.42	75	36241	61.52 ug/L	84
84) 1,2,4-trichlorobenzene	20.88	180	224679	50.84 ug/L	94
85) hexachlorobutadiene	21.14	225	99709m	47.68 ug/L	46
86) naphthalene	21.30	128	549327	56.05 ug/L	100
87) 1,2,3-trichlorobenzene	21.76	180	198103	51.96 ug/L	97

(#) = qualifier out of range (m) = manual integration

000112

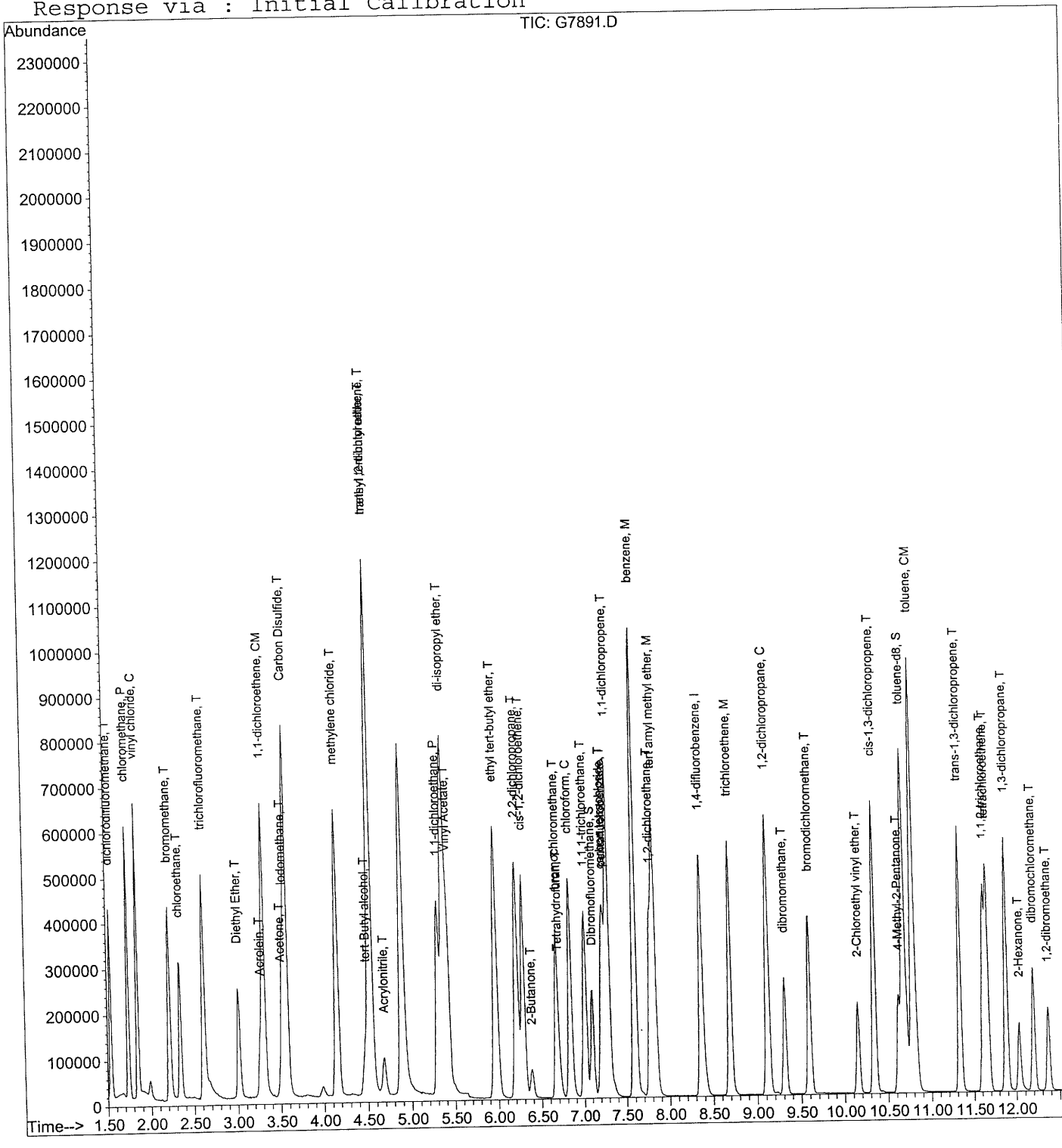
Quantitation Report

Data File : C:\HPCHEM\1\DATA\051205\G7891.D
Acq On : 12 May 05 10:46
Sample : vstd050
Misc : ical
MS Integration Params: rteint.p
Quant Time: May 12 11:48 19105

Vial: 2
Operator: XL
Inst : inst g
Multiplr: 1.00

Quant Results File: 8260BS.RES

Method : C:\HPCHEM\1\DATA\051705\8260BASP.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Fri May 20 09:55:33 2005
Response via : Initial Calibration



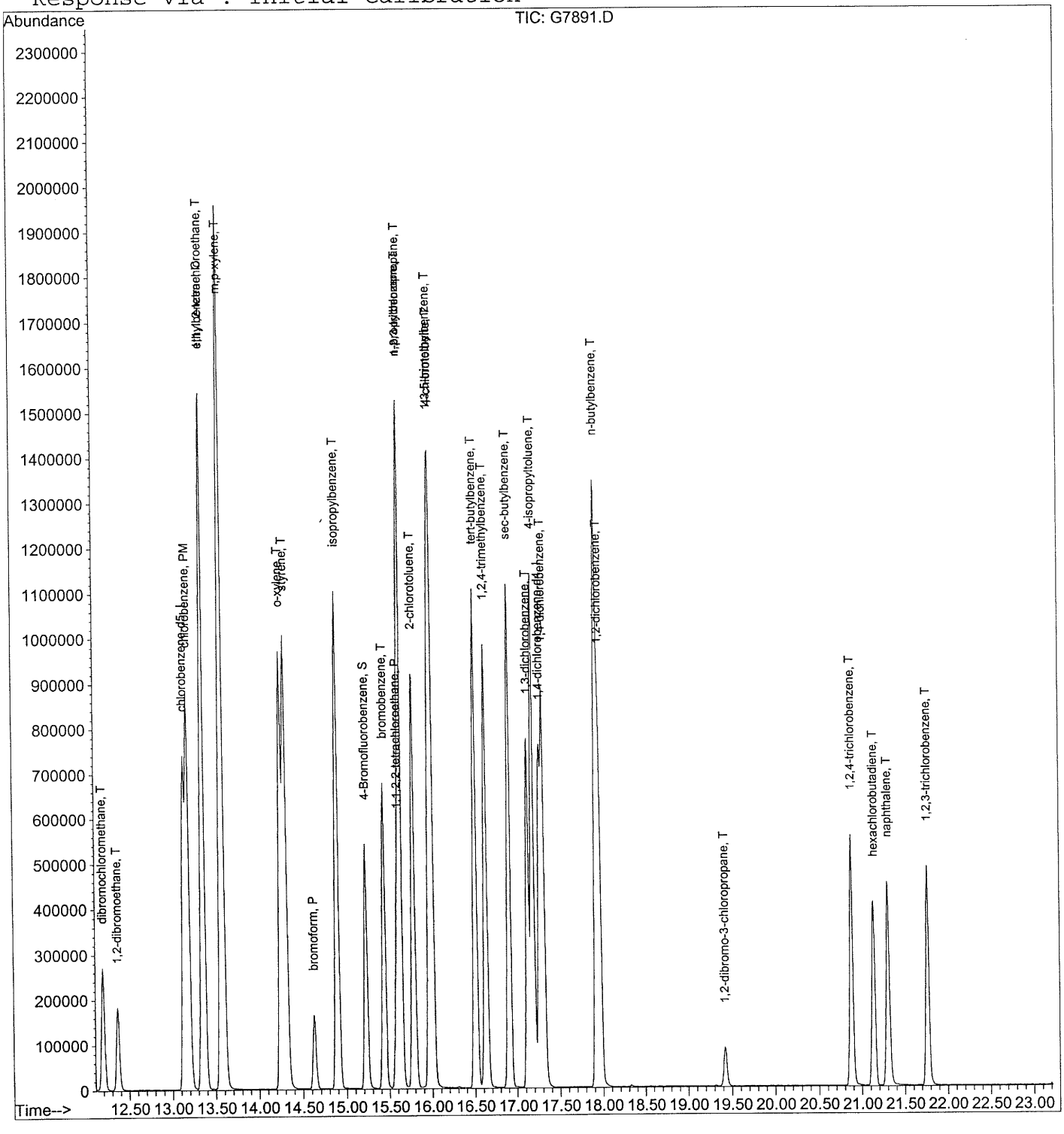
Quantitation Report

Data File : C:\HPCHEM\1\DATA\051205\G7891.D
Acq On : 12 May 05 10:46
Sample : vstd050
Misc : ical
MS Integration Params: rteint.p
Quant Time: May 12 11:48 19105

Vial: 2
Operator: XL
Inst : inst g
Multiplr: 1.00

Quant Results File: 8260BS.RES

Method : C:\HPCHEM\1\DATA\051705\8260BASP.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Fri May 20 09:55:33 2005
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\051205\G7894.D
Acq On : 12 May 05 13:15
Sample : vstd100
Misc : ical
MS Integration Params: rteint.p
Quant Time: May 12 16:02 19105

Vial: 1
Operator: XL
Inst : inst g
Multiplr: 1.00

Quant Results File: 8260BS.RES

Quant Method : C:\HPCHEM\1\DATA\051205\8260BS.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Tue May 03 11:15:26 2005
Response via : Initial Calibration
DataAcq Meth : 8260BS

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) pentafluorobenzene	7.24	168	305178	50.00	ug/L	-0.03
34) 1,4-difluorobenzene	8.36	114	619312	50.00	ug/L	-0.04
52) chlorobenzene-d5	13.14	117	532750	50.00	ug/L	-0.03
66) 1,4-dichlorobenzene-d4	17.27	152	244581	50.00	ug/L	-0.03
System Monitoring Compounds						
31) Dibromofluoromethane	7.09	113	195669	51.70	ug/L	-0.04
Spiked Amount 50.000	Range 86 - 118		Recovery = 103.40%			
48) toluene-d8	10.68	98	768441	50.25	ug/L	-0.03
Spiked Amount 50.000	Range 88 - 110		Recovery = 100.50%			
65) 4-Bromofluorobenzene	15.22	95	278234	51.00	ug/L	-0.03
Spiked Amount 50.000	Range 86 - 115		Recovery = 102.00%			
Target Compounds						Qvalue
2) dichlorodifluoromethane	1.53	85	681787	107.18	ug/L	99
3) chloromethane	1.74	50	1037496	108.81	ug/L	100
4) vinyl chloride	1.85	62	1196003	108.45	ug/L	100
5) bromomethane	2.20	96	386559	59.40	ug/L	99
6) chloroethane	2.33	64	655597	107.40	ug/L	92
7) trichlorofluoromethane	2.61	101	1176671	108.12	ug/L	99
8) Diethyl Ether	3.01	45	267097	90.06	ug/L	97
9) Acrolein	3.26	56	287289	534.97	ug/L	98
10) Acetone	3.50	43	154047	118.67	ug/L	99
11) 1,1-dichloroethene	3.30	96	713970	106.15	ug/L	86
12) Iodomethane	3.52	142	1073714	112.97	ug/L	91
13) methylene chloride	4.15	84	795833	107.68	ug/L	98
14) Carbon Disulfide	3.57	76	3146179	109.69	ug/L	91
17) Acrylonitrile	4.68	53	165945	106.68	ug/L	90
18) tert-Butyl alcohol	4.47	59	430420	1080.80	ug/L	100
19) methyl tert-butyl ether	4.53	73	1642038	108.13	ug/L	100
20) trans-1,2-dichloroethene	4.53	96	805694	105.65	ug/L	90
22) 1,1-dichloroethane	5.30	63	1436109	107.66	ug/L	98
23) di-isopropyl ether	5.37	45	2609397	104.89	ug/L	93
24) Vinyl Acetate	5.41	43	1392951m	113.99	ug/L	100
25) ethyl tert-butyl ether	5.97	59	1687945	113.38	ug/L	99
26) 2-Butanone	6.38	43	222107	109.41	ug/L	94
27) 2,2-dichloropropane	6.21	77	943991	114.16	ug/L	96

(#) = qualifier out of range (m) = manual integration

000115

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\051205\G7894.D
Acq On : 12 May 05 13:15
Sample : vstd100
Misc : ical
MS Integration Params: rteint.p
Quant Time: May 12 16:02 19105

Vial: 1
Operator: XL
Inst : inst g
Multiplr: 1.00

Quant Results File: 8260BS.RES

Quant Method : C:\HPCHEM\1\DATA\051205\8260BS.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Tue May 03 11:15:26 2005
Response via : Initial Calibration
DataAcq Meth : 8260BS

Compound	R.T.	QIon	Response	Conc Unit	Qvalue
28) cis-1,2-dichloroethene	6.28	96	560600	113.40 ug/L	89
29) bromochloromethane	6.67	128	202524m	113.24 ug/L	90
30) chloroform	6.83	83	1089336	114.01 ug/L	99
32) Tetrahydrofuran	6.70	42	117658	113.18 ug/L	95
33) 1,1,1-trichloroethane	7.01	97	791930	111.90 ug/L	94
35) carbon tetrachloride	7.21	117	572072	106.54 ug/L	97
36) 1,1-dichloropropene	7.29	75	956221	100.13 ug/L	97
37) benzene	7.60	78	2543232	111.04 ug/L	98
38) 1,2-dichloroethane	7.77	62	802537	108.65 ug/L	100
39) tert amyl methyl ether	7.82	73	1352843	111.72 ug/L	96
40) trichloroethene	8.70	95	531070	107.69 ug/L	92
41) 1,2-dichloropropane	9.12	63	582550	112.96 ug/L	100
42) dibromomethane	9.30	93	296601	112.20 ug/L	95
44) bromodichloromethane	9.59	83	747961	112.53 ug/L	97
45) 2-Chloroethyl vinyl ether	10.15	63	208888	116.23 ug/L	97
46) 4-Methyl-2-Pentanone	10.62	43	479431	104.53 ug/L	80
47) cis-1,3-dichloropropene	10.34	75	941008	115.98 ug/L	97
49) toluene	10.79	92	1408187	112.27 ug/L	92
50) trans-1,3-dichloropropene	11.33	75	803157	118.17 ug/L	97
51) 1,1,2-trichloroethane	11.61	83	348064	114.17 ug/L	93
53) 2-Hexanone	12.03	43	316367	116.27 ug/L	98
54) tetrachloroethene	11.66	166	420543	107.01 ug/L	99
55) 1,3-dichloropropane	11.88	76	778791	110.76 ug/L	95
56) dibromochloromethane	12.20	129	368685	113.40 ug/L	96
57) 1,2-dibromoethane	12.36	107	339532	110.68 ug/L	96
58) chlorobenzene	13.18	112	1272840	110.17 ug/L	98
59) 1,1,1,2-tetrachloroethane	13.36	131	406651	115.80 ug/L	93
60) ethylbenzene	13.35	91	2888303	115.67 ug/L	93
61) m,p-xylene	13.57	106	1973337	231.51 ug/L	99
62) o-xylene	14.25	106	849481	115.90 ug/L	97
63) styrene	14.31	104	1585963	118.26 ug/L	98
64) bromoform	14.63	173	190339	111.05 ug/L	94
67) isopropylbenzene	14.90	105	2515424	110.92 ug/L	98
68) bromobenzene	15.44	156	450638	108.99 ug/L	96
69) 1,1,2,2-tetrachloroethane	15.59	83	461666	104.33 ug/L	99
70) 1,2,3-trichloropropane	15.63	75	385295	102.85 ug/L	86
71) n-propylbenzene	15.62	91	3584172	111.53 ug/L	98

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\051205\G7894.D Vial: 1
Acq On : 12 May 05 13:15 Operator: XL
Sample : vstd100 Inst : inst g
Misc : ical Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 12 16:02 19105 Quant Results File: 8260BS.RES

Quant Method : C:\HPCHEM\1\DATA\051205\8260BS.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Tue May 03 11:15:26 2005
Response via : Initial Calibration
DataAcq Meth : 8260BS

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
72) 2-chlorotoluene	15.78	91	1804269	109.93	ug/L	98
73) 4-chlorotoluene	16.00	91	2229348	111.28	ug/L	97
74) 1,3,5-trimethylbenzene	15.98	105	2000042	112.59	ug/L	98
75) tert-butylbenzene	16.51	91	1244461	110.58	ug/L	98
76) 1,2,4-trimethylbenzene	16.63	105	1946408	112.69	ug/L	98
77) sec-butylbenzene	16.90	105	2774328	112.68	ug/L	96
78) 1,3-dichlorobenzene	17.12	146	956036	111.50	ug/L	95
79) 4-isopropyltoluene	17.19	119	2146959	114.23	ug/L	98
80) 1,4-dichlorobenzene	17.31	146	959515	109.94	ug/L	96
81) 1,2-dichlorobenzene	17.97	146	865707	111.44	ug/L	94
82) n-butylbenzene	17.92	91	2625885	116.19	ug/L	94
83) 1,2-dibromo-3-chloropropan	19.42	75	60814	107.57	ug/L	86
84) 1,2,4-trichlorobenzene	20.88	180	469481	110.68	ug/L	93
85) hexachlorobutadiene	21.14	225	223044m	111.12	ug/L	46
86) naphthalene	21.30	128	952036	101.22	ug/L	100
87) 1,2,3-trichlorobenzene	21.76	180	382330	104.48	ug/L	96

(#) = qualifier out of range (m) = manual integration

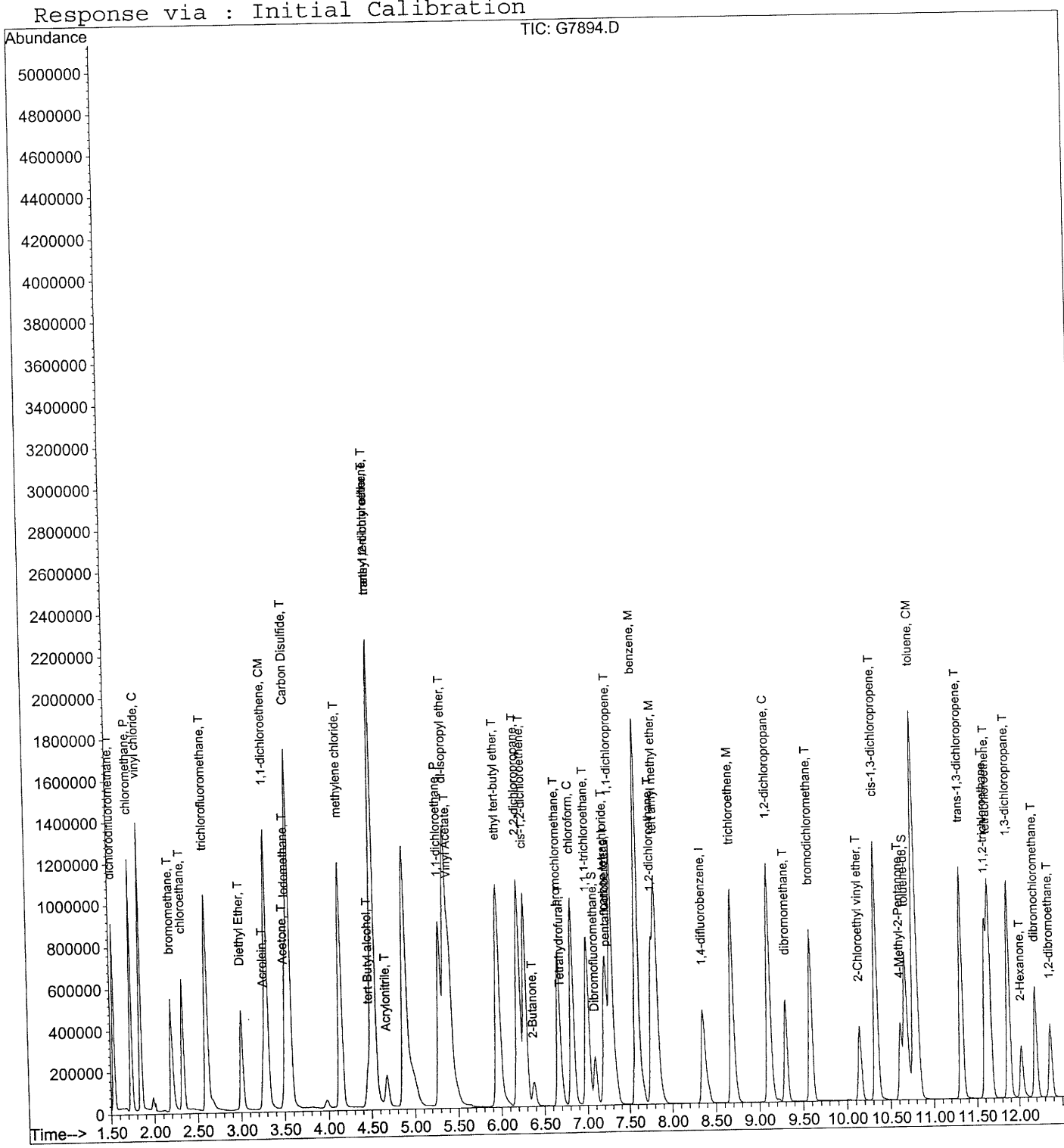
Quantitation Report

Data File : C:\HPCHEM\1\DATA\051205\G7894.D
Acq On : 12 May 05 13:15
Sample : vstd100
Misc : ical
MS Integration Params: rteint.p
Quant Time: May 12 16:02 19105

Vial: 1
Operator: XL
Inst : inst g
Multiplr: 1.00

Quant Results File: 8260BS.RES

Method : C:\HPCHEM\1\DATA\051705\8260BASP.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Fri May 20 09:55:33 2005
Response via : Initial Calibration



000118

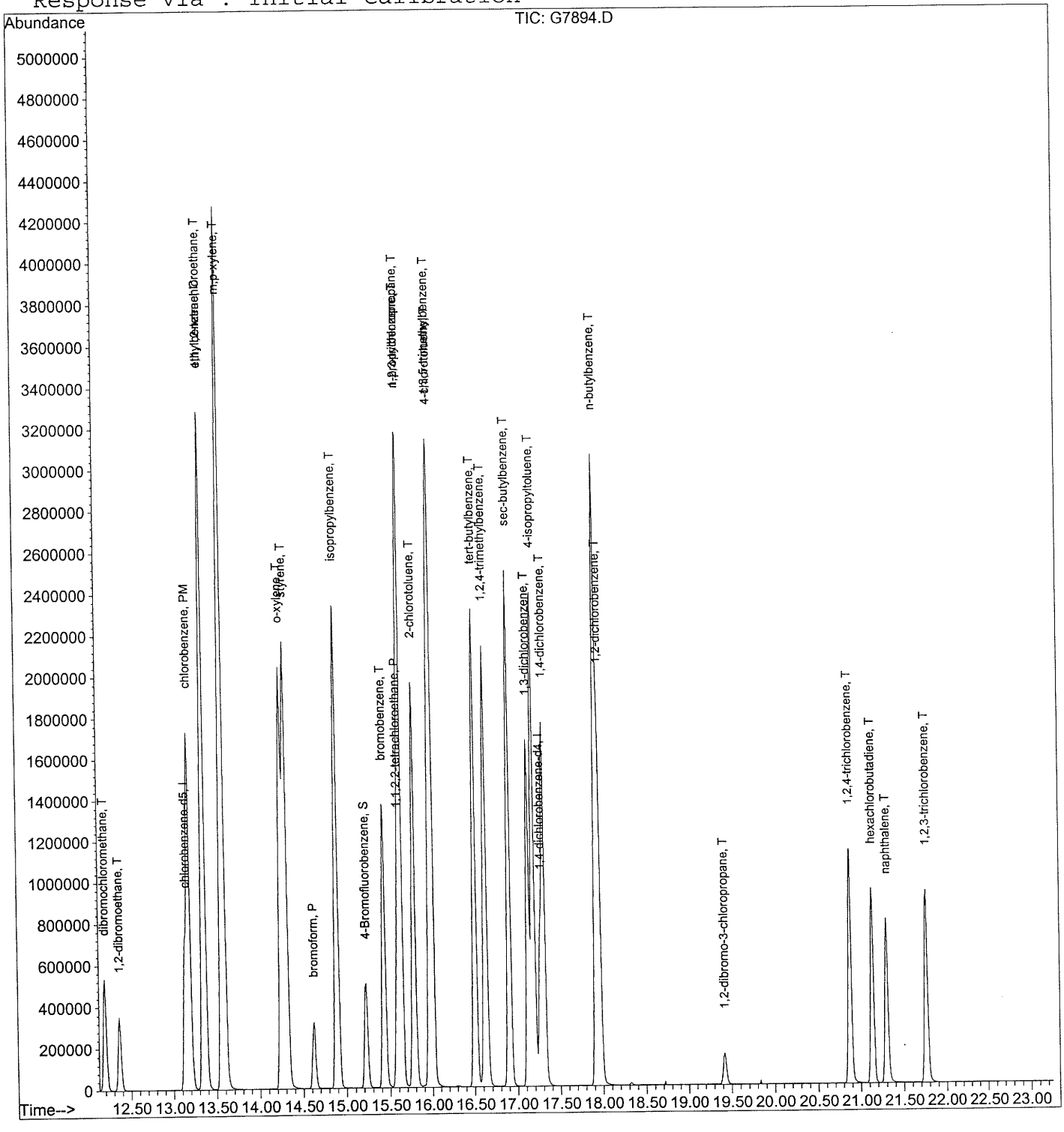
Quantitation Report

Data File : C:\HPCHEM\1\DATA\051205\G7894.D
Acq On : 12 May 05 13:15
Sample : vstd100
Misc : ical
MS Integration Params: rteint.p
Quant Time: May 12 16:02 19105

Vial: 1
Operator: XL
Inst : inst g
Multiplr: 1.00

Quant Results File: 8260BS.RES

Method : C:\HPCHEM\1\DATA\051705\8260BASP.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Fri May 20 09:55:33 2005
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\051205\G7893.D
Acq On : 12 May 05 11:54
Sample : vstd200
Misc : ical
MS Integration Params: rteint.p
Quant Time: May 12 12:24 19105

Vial: 4
Operator: XL
Inst : inst g
Multiplr: 1.00

Quant Results File: 8260BS.RES

Quant Method : C:\HPCHEM\1\DATA\051205\8260BS.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Tue May 03 11:15:26 2005
Response via : Initial Calibration
DataAcq Meth : 8260BS

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) pentafluorobenzene	7.24	168	339737	50.00	ug/L	-0.03
34) 1,4-difluorobenzene	8.36	114	678879	50.00	ug/L	-0.03
52) chlorobenzene-d5	13.14	117	582874	50.00	ug/L	-0.03
66) 1,4-dichlorobenzene-d4	17.27	152	271815	50.00	ug/L	-0.02
System Monitoring Compounds						
31) Dibromofluoromethane	7.10	113	212129	50.35	ug/L	-0.03
Spiked Amount 50.000	Range 86 - 118		Recovery = 100.70%			
48) toluene-d8	10.67	98	829942	49.51	ug/L	-0.03
Spiked Amount 50.000	Range 88 - 110		Recovery = 99.02%			
65) 4-Bromofluorobenzene	15.22	95	302729	50.72	ug/L	-0.03
Spiked Amount 50.000	Range 86 - 115		Recovery = 101.44%			
Target Compounds						Qvalue
2) dichlorodifluoromethane	1.53	85	1295073	182.89	ug/L	98
3) chloromethane	1.73	50	2429909	228.92	ug/L	100
4) vinyl chloride	1.84	62	2558149	208.38	ug/L	99
5) bromomethane	2.20	96	738723	101.97	ug/L	100
6) chloroethane	2.33	64	1360073	200.14	ug/L	91
7) trichlorofluoromethane	2.61	101	2512923	207.42	ug/L	100
8) Diethyl Ether	3.01	45	562515	170.37	ug/L	95
9) Acrolein	3.26	56	603972	1010.27	ug/L	98
10) Acetone	3.49	43	325437	213.58	ug/L	96
11) 1,1-dichloroethene	3.30	96	1520903	203.13	ug/L	88
12) Iodomethane	3.52	142	2285567	216.00	ug/L	93
13) methylene chloride	4.14	84	1669463	202.92	ug/L	97
14) Carbon Disulfide	3.56	76	6769087	212.00	ug/L	89
17) Acrylonitrile	4.68	53	357552	206.47	ug/L	87
18) tert-Butyl alcohol	4.46	59	850647	1918.73	ug/L	100
19) methyl tert-butyl ether	4.52	73	3589291	212.32	ug/L	100
20) trans-1,2-dichloroethene	4.52	96	1729024	203.67	ug/L	90
22) 1,1-dichloroethane	5.30	63	2995881	201.75	ug/L	96
23) di-isopropyl ether	5.37	45	5641606	203.70	ug/L	95
24) Vinyl Acetate	5.41	43	2678550	196.89	ug/L	100
25) ethyl tert-butyl ether	5.97	59	3839738	231.68	ug/L	98
26) 2-Butanone	6.38	43	493058	218.17	ug/L	95
27) 2,2-dichloropropane	6.21	77	2105090	228.67	ug/L	97

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\051205\G7893.D
Acq On : 12 May 05 11:54
Sample : vstd200
Misc : ical
MS Integration Params: rteint.p
Quant Time: May 12 12:24 19105

Vial: 4
Operator: XL
Inst : inst g
Multiplr: 1.00

Quant Results File: 8260BS.RES

Quant Method : C:\HPCHEM\1\DATA\051205\8260BS.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Tue May 03 11:15:26 2005
Response via : Initial Calibration
DataAcq Meth : 8260BS

Compound	R.T.	QIon	Response	Conc Unit	Qvalue
28) cis-1,2-dichloroethene	6.29	96	1256673	228.36 ug/L	91
29) bromochloromethane	6.67	128	453013	227.53 ug/L	99
30) chloroform	6.83	83	2395765	225.24 ug/L	99
32) Tetrahydrofuran	6.70	42	254216	219.67 ug/L	93
33) 1,1,1-trichloroethane	7.00	97	1785607	226.65 ug/L	96
35) carbon tetrachloride	7.21	117	1272713	217.19 ug/L	97
36) 1,1-dichloropropene	7.28	75	2110818	201.64 ug/L	96
37) benzene	7.59	78	5887668	234.50 ug/L	97
38) 1,2-dichloroethane	7.77	62	1762935	217.74 ug/L	99
39) tert amyl methyl ether	7.81	73	3127358	235.60 ug/L	96
40) trichloroethene	8.69	95	1219058	225.51 ug/L	93
41) 1,2-dichloropropane	9.11	63	1310172	231.75 ug/L	100
42) dibromomethane	9.31	93	664911	229.45 ug/L	97
44) bromodichloromethane	9.59	83	1680373	230.62 ug/L	98
45) 2-Chloroethyl vinyl ether	10.15	63	492371	205.39 ug/L	100
46) 4-Methyl-2-Pentanone	10.61	43	1085165m	215.83 ug/L	80
47) cis-1,3-dichloropropene	10.34	75	2173014	244.32 ug/L	97
49) toluene	10.78	92	3196721	232.51 ug/L	91
50) trans-1,3-dichloropropene	11.32	75	1849456	248.24 ug/L	95
51) 1,1,2-trichloroethane	11.61	83	797079	238.52 ug/L	95
53) 2-Hexanone	12.03	43	734673	246.78 ug/L	97
54) tetrachloroethene	11.65	166	986265	229.37 ug/L	99
55) 1,3-dichloropropane	11.88	76	1764451	229.37 ug/L	93
56) dibromochloromethane	12.19	129	855330	240.45 ug/L	97
57) 1,2-dibromoethane	12.37	107	779130	232.13 ug/L	96
58) chlorobenzene	13.19	112	2919511	230.97 ug/L	95
59) 1,1,1,2-tetrachloroethane	13.36	131	993580	258.61 ug/L	92
60) ethylbenzene	13.36	91	6904759	252.73 ug/L	94
61) m,p-xylene	13.58	106	4796089	514.29 ug/L	99
62) o-xylene	14.25	106	1976268	246.44 ug/L	98
63) styrene	14.31	104	3731395	254.31 ug/L	99
64) bromoform	14.63	173	438840	201.11 ug/L	94
67) isopropylbenzene	14.89	105	5895920	233.93 ug/L	99
68) bromobenzene	15.43	156	1030709	224.31 ug/L	93
69) 1,1,2,2-tetrachloroethane	15.58	83	1125209m	228.81 ug/L	98
70) 1,2,3-trichloropropane	15.63	75	907666	218.01 ug/L	89
71) n-propylbenzene	15.63	91	8547560	239.33 ug/L	98

(#) = qualifier out of range (m) = manual integration

000121

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\051205\G7893.D
Acq On : 12 May 05 11:54
Sample : vstd200
Misc : ical
MS Integration Params: rteint.p
Quant Time: May 12 12:24 19105

Vial: 4
Operator: XL
Inst : inst g
Multiplr: 1.00

Quant Results File: 8260BS.RES

Quant Method : C:\HPCHEM\1\DATA\051205\8260BS.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Tue May 03 11:15:26 2005
Response via : Initial Calibration
DataAcq Meth : 8260BS

Compound	R.T.	QIon	Response	Conc Unit	Qvalue
72) 2-chlorotoluene	15.79	91	4140169	226.98 ug/L	98
73) 4-chlorotoluene	16.00	91	5283593	237.31 ug/L	95
74) 1,3,5-trimethylbenzene	15.97	105	4800649	243.16 ug/L	100
75) tert-butylbenzene	16.50	91	2843796	227.37 ug/L	99
76) 1,2,4-trimethylbenzene	16.63	105	4565359	237.83 ug/L	100
77) sec-butylbenzene	16.91	105	6555331	239.57 ug/L	96
78) 1,3-dichlorobenzene	17.13	146	2254197	236.57 ug/L	95
79) 4-isopropyltoluene	17.19	119	5110808	244.68 ug/L	98
80) 1,4-dichlorobenzene	17.31	146	2258972	232.90 ug/L	97
81) 1,2-dichlorobenzene	17.96	146	2084652	241.46 ug/L	93
82) n-butylbenzene	17.93	91	6195660	246.68 ug/L	95
83) 1,2-dibromo-3-chloropropan	19.42	75	152292	242.38 ug/L	85
84) 1,2,4-trichlorobenzene	20.87	180	1140867	242.02 ug/L	92
85) hexachlorobutadiene	21.13	225	530981m	238.04 ug/L	46
86) naphthalene	21.30	128	2446038	234.01 ug/L	100
87) 1,2,3-trichlorobenzene	21.77	180	953621	234.49 ug/L	95

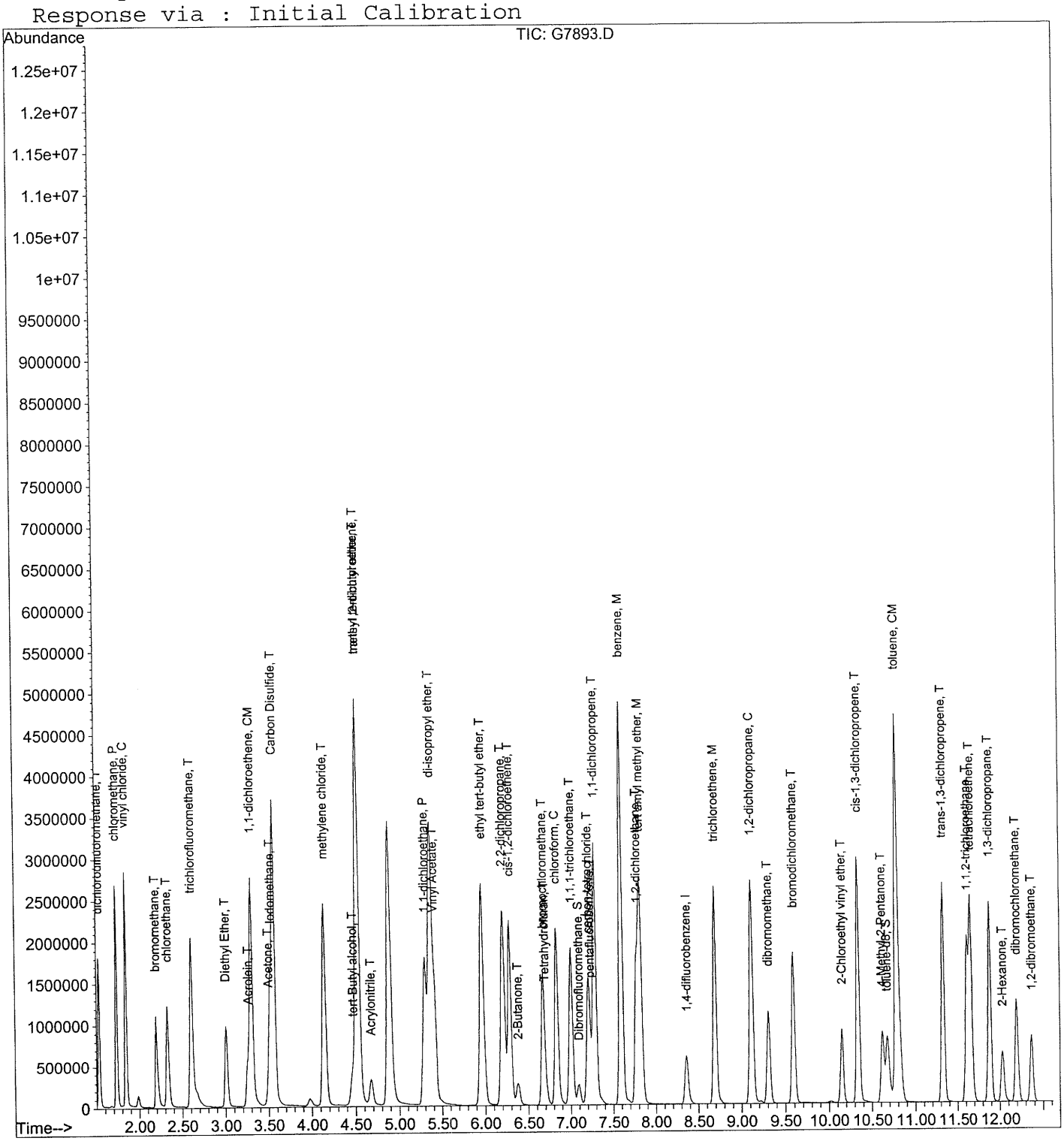
Quantitation Report

Data File : C:\HPCHEM\1\DATA\051205\G7893.D
Acq On : 12 May 05 11:54
Sample : vstd200
Misc : ical
MS Integration Params: rteint.p
Quant Time: May 12 12:24 19105

Vial: 4
Operator: XL
Inst : inst g
Multiplr: 1.00

Quant Results File: 8260BS.RES

Method : C:\HPCHEM\1\DATA\051705\8260BASP.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Fri May 20 09:55:33 2005
Response via : Initial Calibration



000123

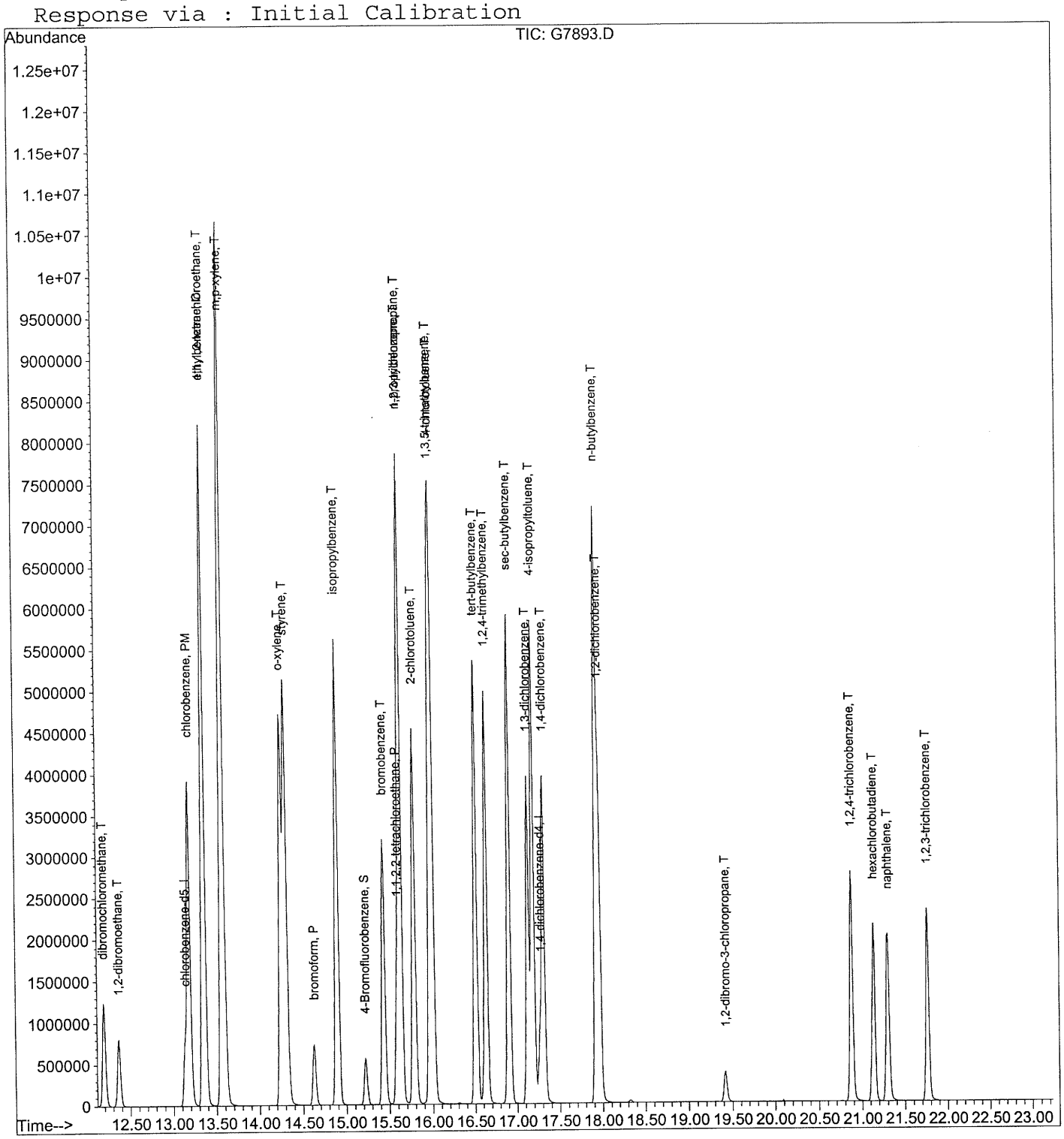
Quantitation Report

Data File : C:\HPCHEM\1\DATA\051205\G7893.D
Acq On : 12 May 05 11:54
Sample : vstd200
Misc : ical
MS Integration Params: rteint.p
Quant Time: May 12 12:24 19105

Vial: 4
Operator: XL
Inst : inst g
Multiplr: 1.00

Quant Results File: 8260BS.RES

Method : C:\HPCHEM\1\DATA\051705\8260BASP.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Fri May 20 09:55:33 2005
Response via : Initial Calibration



Method : C:\HPCHEM\1\DATA\061505\8260BASP.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Fri May 20 09:55:33 2005
Response via : Initial Calibration

Continuing Calibration File: G8358.D

Min. RRF : 0.000 Min. Rel. Area : 50%
Max. RRF Dev : 20% Max. Rel. Area : 200%

		Compound	AvgRF	CCRF	%Dev	Area%
39	M	tert amyl methyl ether	1.092	1.007	7.8	57
40	M	trichloroethene	0.406	0.442	-8.9	70
41	C	1,2-dichloropropane	0.447	0.490	-9.7	70
42	T	dibromomethane	0.236	0.234	0.7	64
43		1,4-Dioxane	0.000	-0.000#	100.5#	0#
44	T	bromodichloromethane	0.567	0.592	-4.6	67
45	T	2-Chloroethyl vinyl ether	0.158	0.188	-19.4	69
46	T	4-Methyl-2-Pentanone	0.425	0.428	-0.9	63
47	T	cis-1,3-dichloropropene	0.719	0.756	-5.2	67
48	S	toluene-d8	1.228	1.252	-2.0	65
49	CM	toluene	1.055	1.140	-8.1	70
50	T	trans-1,3-dichloropropene	0.619	0.603	2.6	61
51	T	1,1,2-trichloroethane	0.275	0.285	-3.6	65
52	I	chlorobenzene-d5	1.000	1.000	0.0	63
53	T	2-Hexanone	0.307	0.301	2.0	59
54	T	tetrachloroethene	0.373	0.387	-3.9	67
55	T	1,3-dichloropropane	0.723	0.745	-3.1	65
56	T	dibromochloromethane	0.330	0.342	-3.5	64
57	T	1,2-dibromoethane	0.316	0.317	-0.4	62
58	PM	chlorobenzene	1.131	1.207	-6.7	69
59	T	1,1,1,2-tetrachloroethane	0.356	0.368	-3.4	68
60	C	ethylbenzene	2.488	2.567	-3.2	67
61	T	m,p-xylene	0.844	0.883	-4.6	68
62	T	o-xylene	0.733	0.765	-4.3	68
63	T	styrene	1.352	1.418	-4.9	68
64	P	bromoform	0.169	0.161	5.2	59
65	S	4-Bromofluorobenzene	0.519	0.507	2.2	61
66	I	1,4-dichlorobenzene-d4	1.000	1.000	0.0	62
67	T	isopropylbenzene	4.715	4.973	-5.5	68
68	T	bromobenzene	0.868	0.890	-2.6	65
69	P	1,1,2,2-tetrachloroethane	1.008	1.035	-2.6	63
70	T	1,2,3-trichloropropane	0.815	0.825	-1.2	63
71	T	n-propylbenzene	6.774	7.340	-8.4	70
72	T	2-chlorotoluene	3.450	3.446	0.1	64
73	T	4-chlorotoluene	4.213	4.523	-7.3	69
74	T	1,3,5-trimethylbenzene	3.740	3.943	-5.4	68
75	T	tert-butylbenzene	2.323	2.482	-6.8	68

(#) = Out of Range

Method : C:\HPCHEM\1\DATA\061505\8260BASP.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Fri May 20 09:55:33 2005
Response via : Initial Calibration

Continuing Calibration File: G8358.D

Min. RRF : 0.000 Min. Rel. Area : 50%
Max. RRF Dev : 20% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%
76 T	1,2,4-trimethylbenzene	3.643	3.888	-6.7	68
77 T	sec-butylbenzene	5.167	5.573	-7.8	69
78 T	1,3-dichlorobenzene	1.827	1.923	-5.2	67
79 T	4-isopropyltoluene	3.979	4.045	-1.7	65
80 T	1,4-dichlorobenzene	1.863	1.950	-4.7	67
81 T	1,2-dichlorobenzene	1.664	1.729	-3.9	66
82 T	n-butylbenzene	4.814	5.377	-11.7	72
83 T	1,2-dibromo-3-chloropropane	0.135	0.115	15.2	50#
84 T	1,2,4-trichlorobenzene	0.903	0.905	-0.2	64
85 T	hexachlorobutadiene	0.409	0.445	-8.7	71
86 T	naphthalene	2.049	2.068	-0.9	59
87 T	1,2,3-trichlorobenzene	0.774	0.806	-4.2	64

(#) = Out of Range SPCC's out = 0 CCC's out = 0

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\061505\G8358.D
Acq On : 15 Jun 05 12:12
Sample : vstd050
Misc : ccv
MS Integration Params: rteint.p
Quant Time: Jun 16 9:30 19105

Vial: 2
Operator: xl
Inst : inst g
Multiplr: 1.00

Quant Results File: 8260BASP.RE

Quant Method : C:\HPCHEM\1\DATA\061505\8260BASP.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Fri May 20 09:55:33 2005
Response via : Initial Calibration
DataAcq Meth : 8260BS

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) pentafluorobenzene	7.28	168	188757	50.00	ug/Kg	0.03
34) 1,4-difluorobenzene	8.40	114	412244	50.00	ug/Kg	0.04
52) chlorobenzene-d5	13.17	117	349176	50.00	ug/Kg	0.03
66) 1,4-dichlorobenzene-d4	17.30	152	157999	50.00	ug/Kg	0.03
System Monitoring Compounds						
31) Dibromofluoromethane	7.13	113	128576	52.22	ug/Kg	0.04
Spiked Amount 50.000	Range 80 - 120		Recovery = 104.44%			
48) toluene-d8	10.72	98	516262	50.98	ug/Kg	0.04
Spiked Amount 50.000	Range 81 - 120		Recovery = 101.96%			
65) 4-Bromofluorobenzene	15.26	95	177085	48.88	ug/Kg	0.03
Spiked Amount 50.000	Range 74 - 121		Recovery = 97.76%			
Target Compounds						Qvalue
2) dichlorodifluoromethane	1.55	85	185352	47.08	ug/Kg	100
3) chloromethane	1.75	50	366027	55.40	ug/Kg	99
4) vinyl chloride	1.86	62	390989	55.26	ug/Kg	100
5) bromomethane	2.23	96	210927	74.94	ug/Kg	99
6) chloroethane	2.36	64	223928	57.75	ug/Kg	100
7) trichlorofluoromethane	2.63	101	386899	55.46	ug/Kg	100
8) Diethyl Ether	3.03	45	102195	57.67	ug/Kg	97
9) Acrolein	3.29	56	73835	204.41	ug/Kg	97
10) Acetone	3.52	43	57558	57.43	ug/Kg	99
11) 1,1-dichloroethene	3.33	96	229347	54.18	ug/Kg	90
12) Iodomethane	3.56	142	330665	53.40	ug/Kg	96
13) methylene chloride	4.18	84	259185	52.95	ug/Kg	96
14) Carbon Disulfide	3.60	76	998704	54.75	ug/Kg	98
17) Acrylonitrile	4.71	53	54026	48.28	ug/Kg	99
18) tert-Butyl alcohol	4.52	59	98387	312.53	ug/Kg	100
19) methyl tert-butyl ether	4.56	73	518475	49.37	ug/Kg	100
20) trans-1,2-dichloroethene	4.56	96	256799	54.13	ug/Kg	89
21) n-Hexane	4.92	57	446514m	55.82	ug/Kg	17
22) 1,1-dichloroethane	5.34	63	481908	56.32	ug/Kg	99
23) di-isopropyl ether	5.40	45	841211	52.22	ug/Kg	99
24) Vinyl Acetate	5.45	43	434736m	51.55	ug/Kg	99
25) ethyl tert-butyl ether	6.01	59	512282	49.11	ug/Kg	100
26) 2-Butanone	6.42	43	79732	52.93	ug/Kg	100

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\061505\G8358.D Vial: 2
Acq On : 15 Jun 05 12:12 Operator: xl
Sample : vstd050 Inst : inst g
Misc : ccv Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: Jun 16 9:30 19105 Quant Results File: 8260BASP.RE

Quant Method : C:\HPCHEM\1\DATA\061505\8260BASP.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Fri May 20 09:55:33 2005
Response via : Initial Calibration
DataAcq Meth : 8260BS

	Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
27)	2,2-dichloropropane	6.25	77	296656	53.91	ug/Kg	99
28)	cis-1,2-dichloroethene	6.32	96	192215	58.41	ug/Kg	99
29)	bromochloromethane	6.71	128	65808	53.71	ug/Kg	98
30)	chloroform	6.87	83	366082	56.64	ug/Kg	100
32)	Tetrahydrofuran	6.74	42	38887	49.43	ug/Kg	95
33)	1,1,1-trichloroethane	7.04	97	252495	53.81	ug/Kg	98
35)	carbon tetrachloride	7.24	117	197165	51.39	ug/Kg	99
36)	1,1-dichloropropene	7.32	75	342645	50.80	ug/Kg	98
37)	benzene	7.63	78	898089	55.87	ug/Kg	98
38)	1,2-dichloroethane	7.81	62	250041	48.04	ug/Kg	100
39)	tert amyl methyl ether	7.86	73	415105	46.11	ug/Kg	100
40)	trichloroethene	8.73	95	182244	54.44	ug/Kg	95
41)	1,2-dichloropropane	9.15	63	202143	54.85	ug/Kg	99
42)	dibromomethane	9.34	93	96516	49.63	ug/Kg	99
44)	bromodichloromethane	9.63	83	244219	52.28	ug/Kg	98
45)	2-Chloroethyl vinyl ether	10.19	63	77627	59.68	ug/Kg	99
46)	4-Methyl-2-Pentanone	10.66	43	176609	50.45	ug/Kg	99
47)	cis-1,3-dichloropropene	10.37	75	311792	52.62	ug/Kg	100
49)	toluene	10.82	92	470059	54.06	ug/Kg	98
50)	trans-1,3-dichloropropene	11.36	75	248393	48.69	ug/Kg	99
51)	1,1,2-trichloroethane	11.65	83	117350	51.82	ug/Kg	100
53)	2-Hexanone	12.07	43	105232	49.02	ug/Kg	96
54)	tetrachloroethene	11.69	166	135184	51.96	ug/Kg	98
55)	1,3-dichloropropane	11.92	76	260184	51.55	ug/Kg	99
56)	dibromochloromethane	12.23	129	119313	51.73	ug/Kg	100
57)	1,2-dibromoethane	12.40	107	110718	50.19	ug/Kg	98
58)	chlorobenzene	13.22	112	421331	53.36	ug/Kg	99
59)	1,1,1,2-tetrachloroethane	13.40	131	128648	51.70	ug/Kg	99
60)	ethylbenzene	13.38	91	896283	51.58	ug/Kg	97
61)	m,p-xylene	13.60	106	616584	98.56	ug/Kg	96
62)	o-xylene	14.29	106	267080	52.17	ug/Kg	96
63)	styrene	14.34	104	495247	52.47	ug/Kg	95
64)	bromoform	14.66	173	56088	47.42	ug/Kg	94
67)	isopropylbenzene	14.93	105	785776	52.73	ug/Kg	98
68)	bromobenzene	15.47	156	140678	51.31	ug/Kg	92
69)	1,1,2,2-tetrachloroethane	15.62	83	163459m	51.29	ug/Kg	100
70)	1,2,3-trichloropropane	15.67	75	130372m	50.62	ug/Kg	93

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\061505\G8358.D Vial: 2
Acq On : 15 Jun 05 12:12 Operator: xl
Sample : vstd050 Inst : inst g
Misc : ccv Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: Jun 16 9:30 19105 Quant Results File: 8260BASP.RE

Quant Method : C:\HPCHEM\1\DATA\061505\8260BASP.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Fri May 20 09:55:33 2005
Response via : Initial Calibration
DataAcq Meth : 8260BS

	Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
71)	n-propylbenzene	15.65	91	1159746	54.18	ug/Kg	100
72)	2-chlorotoluene	15.81	91	544415	49.94	ug/Kg	99
73)	4-chlorotoluene	16.03	91	714637	53.67	ug/Kg	99
74)	1,3,5-trimethylbenzene	16.01	105	622955	52.72	ug/Kg	99
75)	tert-butylbenzene	16.54	91	392076	53.41	ug/Kg	99
76)	1,2,4-trimethylbenzene	16.66	105	614260	53.36	ug/Kg	98
77)	sec-butylbenzene	16.93	105	880462	53.92	ug/Kg	100
78)	1,3-dichlorobenzene	17.16	146	303840	52.62	ug/Kg	97
79)	4-isopropyltoluene	17.22	119	639142	50.84	ug/Kg	99
80)	1,4-dichlorobenzene	17.34	146	308114	52.34	ug/Kg	98
81)	1,2-dichlorobenzene	18.00	146	273237	51.95	ug/Kg	97
82)	n-butylbenzene	17.96	91	849494	55.84	ug/Kg	98
83)	1,2-dibromo-3-chloropropan	19.45	75	18115	42.40	ug/Kg	96
84)	1,2,4-trichlorobenzene	20.91	180	143003	50.12	ug/Kg	99
85)	hexachlorobutadiene	21.17	225	70319m	54.36	ug/Kg	100
86)	naphthalene	21.33	128	326665	50.45	ug/Kg	100
87)	1,2,3-trichlorobenzene	21.79	180	127338	52.09	ug/Kg	98

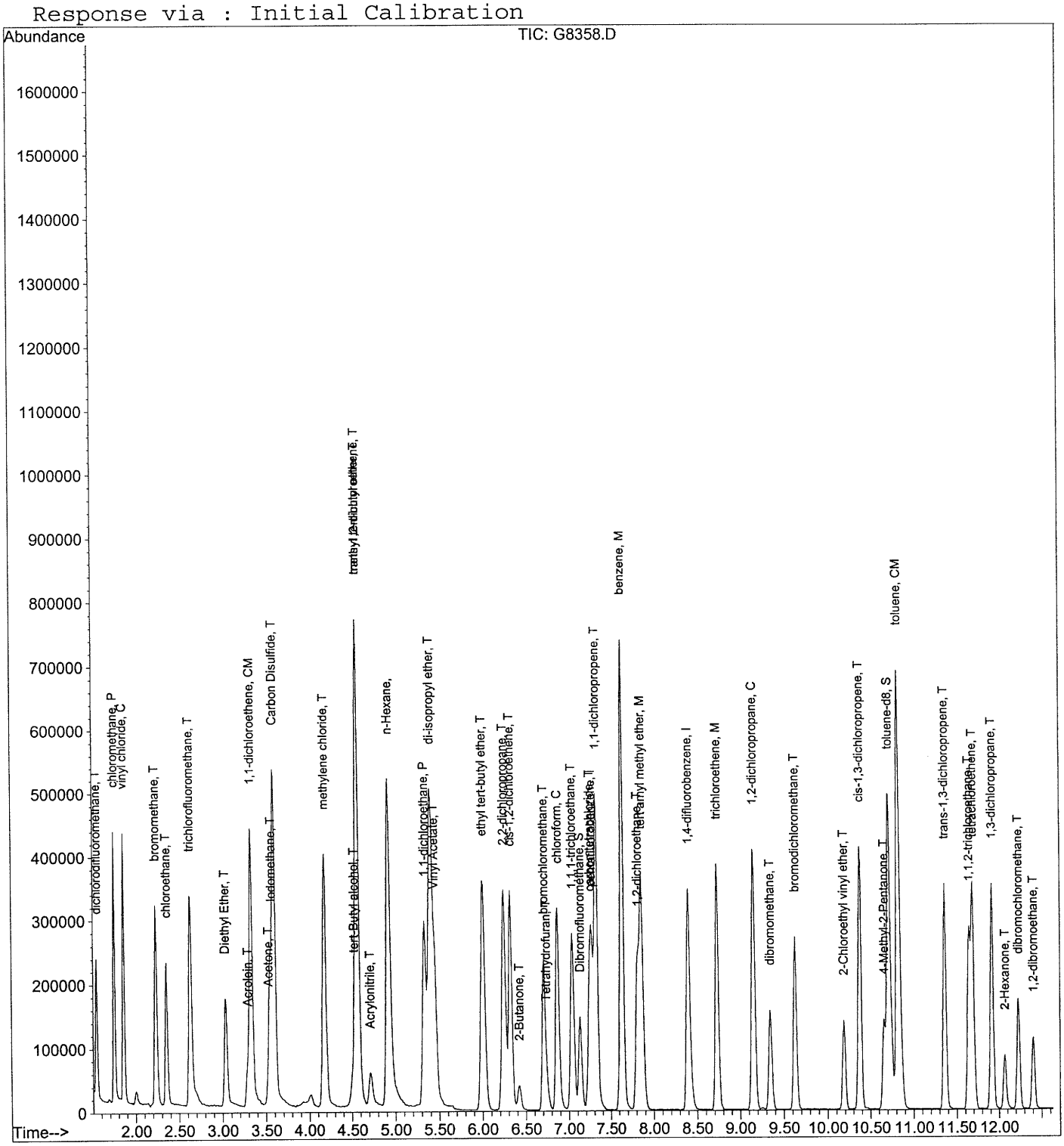
(#) = qualifier out of range (m) = manual integration

000129

Quantitation Report

Data File : C:\HPCHEM\1\DATA\061505\G8358.D Vial: 2
Acq On : 15 Jun 05 12:12 Operator: xl
Sample : vstd050 Inst : inst g
Misc : ccv Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: Jun 16 9:30 19105 Quant Results File: 8260BASP.RE

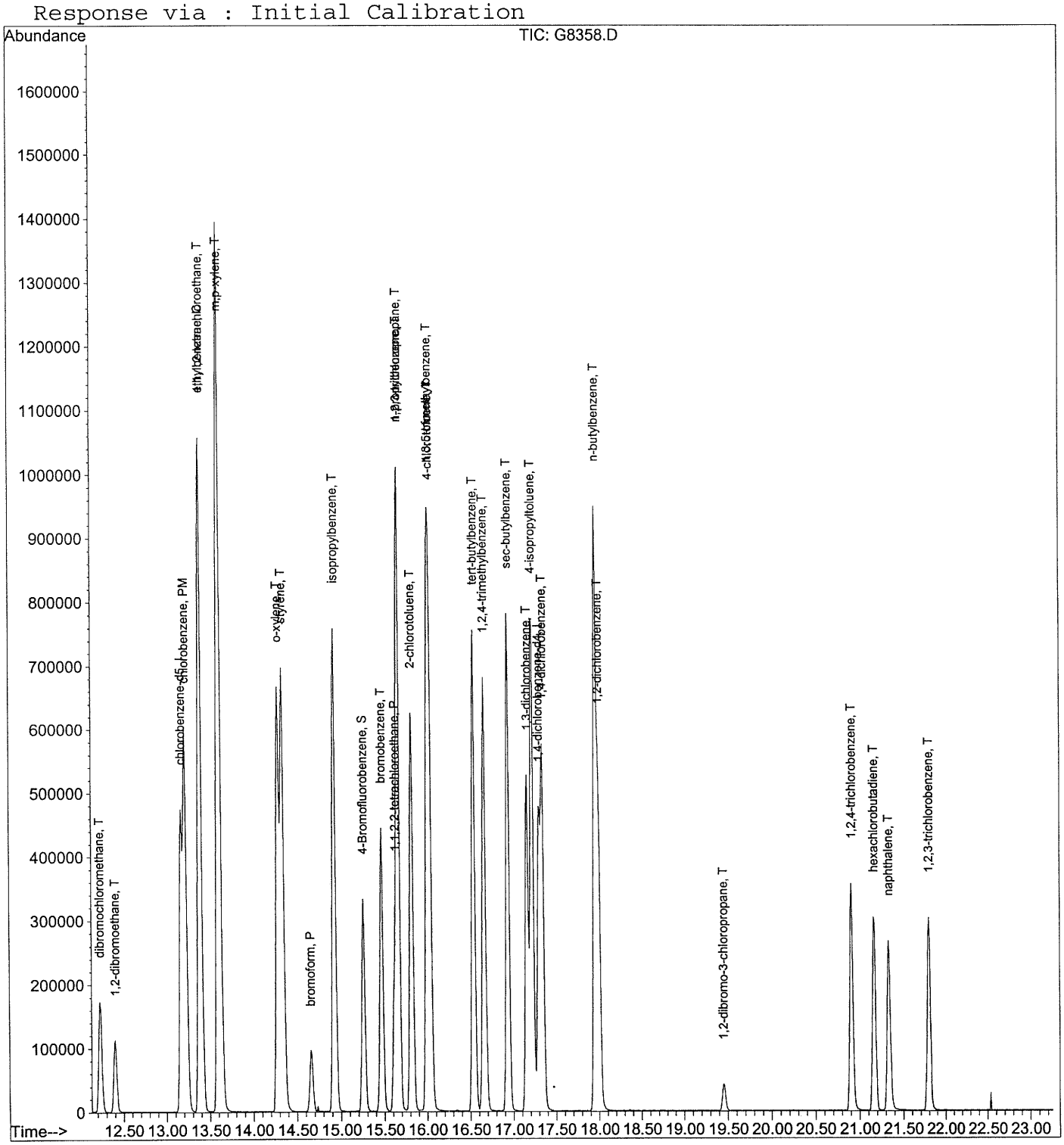
Method : C:\HPCHEM\1\DATA\061505\8260BASP.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Fri May 20 09:55:33 2005
Response via : Initial Calibration



Quantitation Report

Data File : C:\HPCHEM\1\DATA\061505\G8358.D Vial: 2
Acq On : 15 Jun 05 12:12 Operator: xl
Sample : vstd050 Inst : inst g
Misc : ccv Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: Jun 16 9:30 19105 Quant Results File: 8260BASP.RE

Method : C:\HPCHEM\1\DATA\061505\8260BASP.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Fri May 20 09:55:33 2005
Response via : Initial Calibration



000131

Continuing Calibration Report inst g

Method : C:\HPCHEM\1\DATA\061505\8260BASP.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Fri May 20 09:55:33 2005
Response via : Initial Calibration

Continuing Calibration File: G8358.D

Min. RRF : 0.000 Min. Rel. Area : 50%
Max. RRF Dev : 20% Max. Rel. Area : 200%

		Compound	AvgRF	CCRF	%Dev	Area%
1	I	pentafluorobenzene	1.000	1.000	0.0	60
2	T	dichlorodifluoromethane	1.043	0.982	5.8	56
3	P	chloromethane	1.750	1.939	-10.8	66
4	C	vinyl chloride	1.874	2.071	-10.5	67
5	T	bromomethane	0.804	1.117	-38.9#	67
6	T	chloroethane	1.027	1.186	-15.5	69
7	T	trichlorofluoromethane	1.848	2.050	-10.9	68
8	T	Diethyl Ether	0.469	0.541	-15.3	73
9	T	Acrolein	0.096	0.078	18.2	47#
10	T	Acetone	0.300	0.305	-1.7	70
11	CM	1,1-dichloroethene	1.121	1.215	-8.4	66
12	T	Iodomethane	1.640	1.752	-6.8	66
13	T	methylene chloride	1.296	1.373	-5.9	60
14	T	Carbon Disulfide	4.832	5.291	-9.5	67
15		Isopropyl Alcohol	0.000	0.054	0.0	0#
16		Acetonitrile	0.000	0.062	0.0	0#
17	T	Acrylonitrile	0.296	0.286	3.4	56
18	T	tert-Butyl alcohol	0.080	0.052	34.5#	39#
19	T	methyl tert-butyl ether	2.782	2.747	1.3	59
20	T	trans-1,2-dichloroethene	1.257	1.360	-8.3	66
21		n-Hexane	2.119	2.366	-11.6	66
22	P	1,1-dichloroethane	2.267	2.553	-12.6	68
23	T	di-isopropyl ether	4.267	4.457	-4.4	63
24	T	Vinyl Acetate	2.234	2.303	-3.1	60
25	T	ethyl tert-butyl ether	2.763	2.714	1.8	60
26	T	2-Butanone	0.399	0.422	-5.9	64
27	T	2,2-dichloropropane	1.458	1.572	-7.8	65
28	T	cis-1,2-dichloroethene	0.872	1.018	-16.8	71
29	T	bromochloromethane	0.325	0.349	-7.4	65
30	C	chloroform	1.712	1.939	-13.3	69
31	S	Dibromofluoromethane	0.652	0.681	-4.4	63
32	T	Tetrahydrofuran	0.208	0.206	1.1	58
33	T	1,1,1-trichloroethane	1.243	1.338	-7.6	65
34	I	1,4-difluorobenzene	1.000	1.000	0.0	63
35	T	carbon tetrachloride	0.465	0.478	-2.8	70
36	T	1,1-dichloropropene	0.818	0.831	-1.6	69
37	M	benzene	1.950	2.179	-11.7	71
38	T	1,2-dichloroethane	0.631	0.607	3.9	61

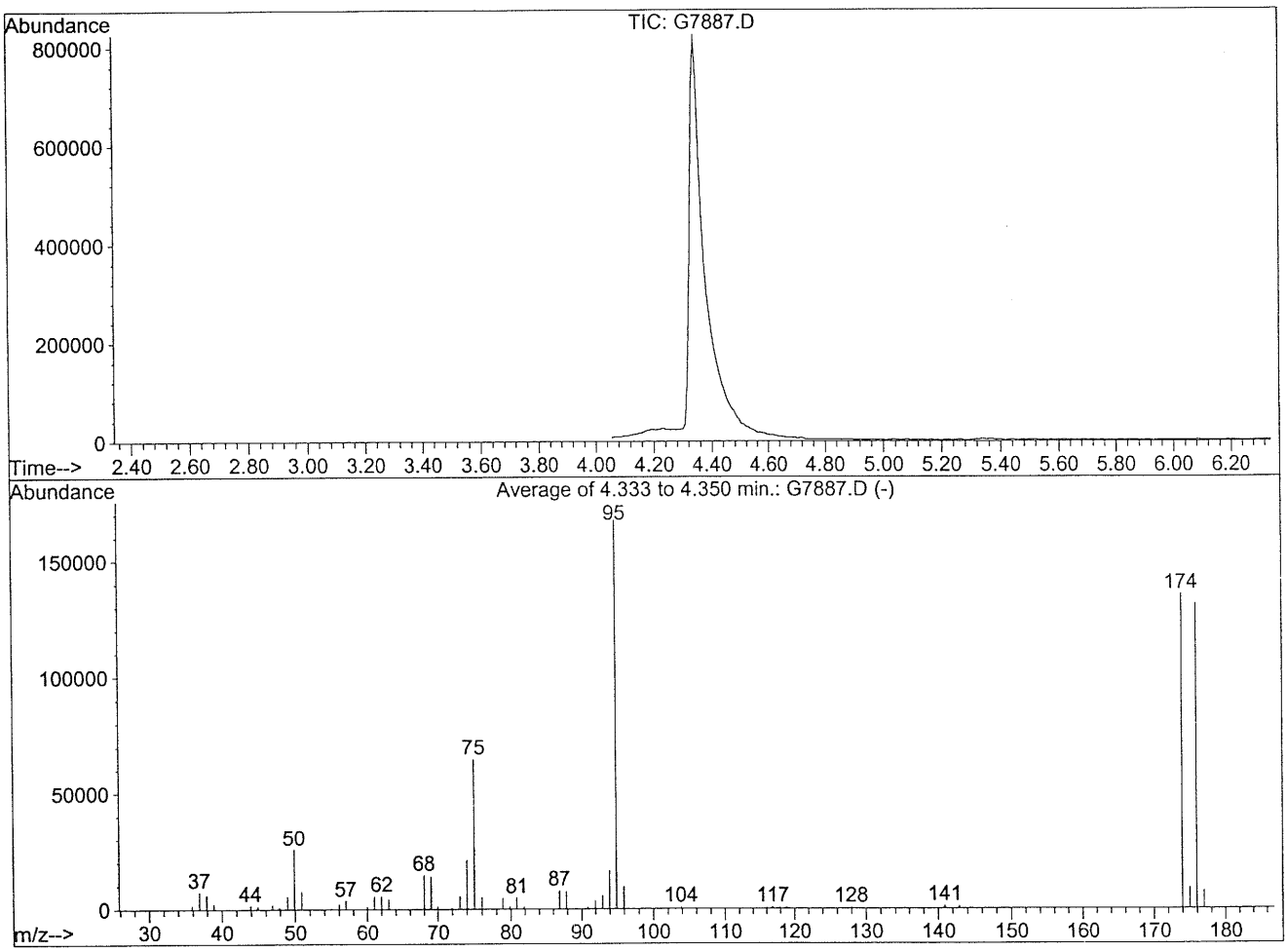
(#) = Out of Range

RAW QC DATA

SW-846 Method 8260

Data File : C:\HPCHEM\1\DATA\051205\G7887.D
Acq On : 12 May 05 8:47
Sample : 50ng bfb std
Misc : samp 8260_wst
MS Integration Params: rteint.p
Method : C:\HPCHEM\1\DATA\051205\8260BS.M (RTE Integrator)
Title : SW-846 Method 8260B

Vial: 1
Operator: XL
Inst : inst g
Multiplr: 1.00



AutoFind: Scans 34, 35, 36; Background Corrected with Scan 27

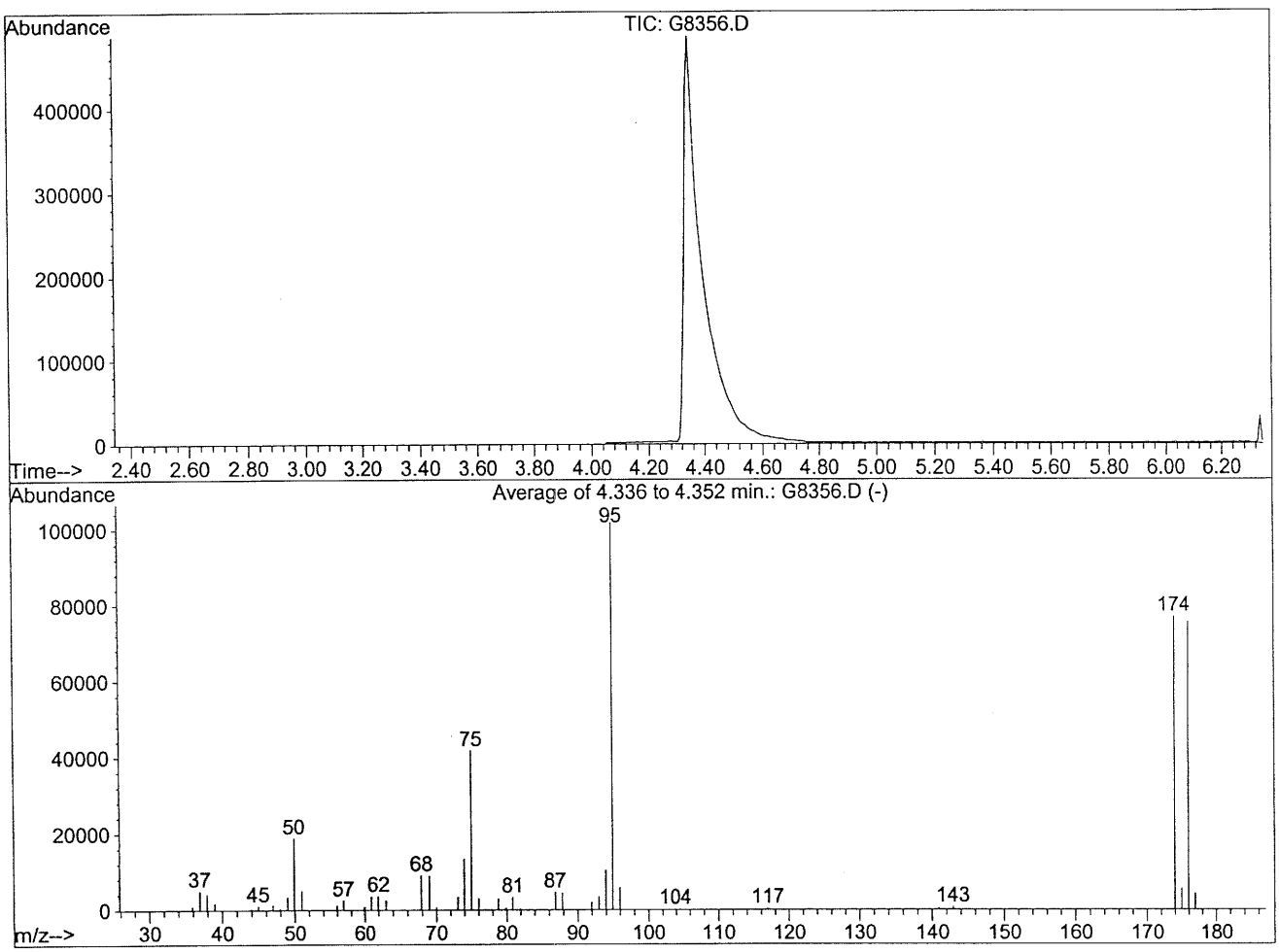
Target Mass	Rel. to Mass	Lower Limit%	Upper Limit%	Rel. Abn%	Raw Abn	Result Pass/Fail
50	95	15	40	15.4	25762	PASS
75	95	30	60	38.7	64781	PASS
95	95	100	100	100.0	167317	PASS
96	95	5	9	5.6	9336	PASS
173	174	0.00	2	0.0	0	PASS
174	95	50	100	80.9	135344	PASS
175	174	5	9	6.6	8895	PASS
176	174	95	101	96.8	131027	PASS
177	176	5	9	5.8	7549	PASS

G7887.D 8260BS.M Tue May 17 11:12:11 2005

BFB

Data File : C:\HPCHEM\1\DATA\061505\G8356.D
Acq On : 15 Jun 05 11:19
Sample : 50ng bfb std
Misc :
MS Integration Params: rteint.p
Method : C:\HPCHEM\1\DATA\061505\BFB.M (RTE Integrator)
Title :

Vial: 1
Operator: xl
Inst : inst g
Multiplr: 1.00

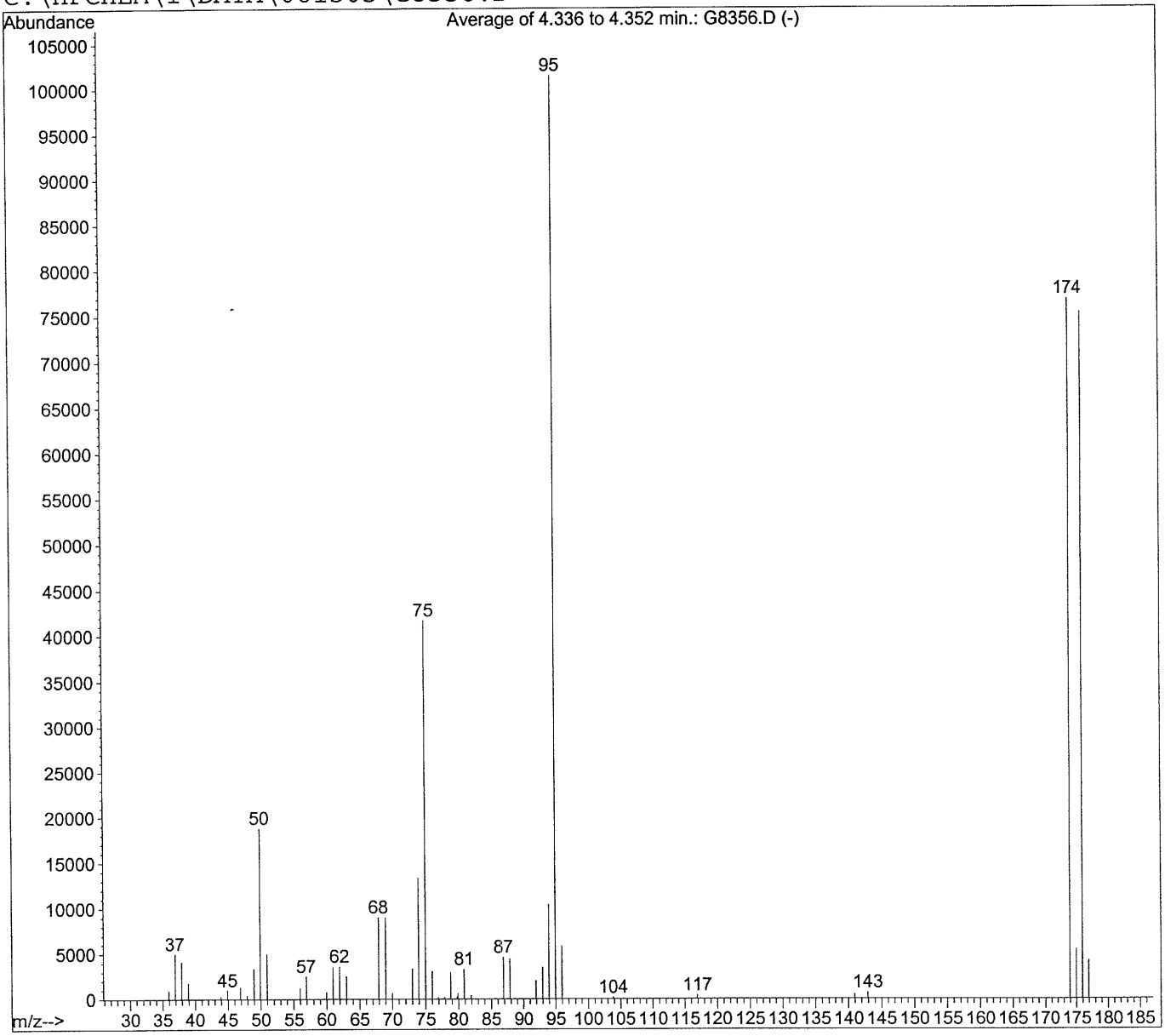


AutoFind: Scans 35, 36, 37; Background Corrected with Scan 29

Target Mass	Rel. to Mass	Lower Limit%	Upper Limit%	Rel. Abn%	Raw Abn	Result Pass/Fail
50	95	15	40	18.5	18803	PASS
75	95	30	60	41.1	41701	PASS
95	95	100	100	100.0	101501	PASS
96	95	5	9	5.8	5856	PASS
173	174	0.00	2	0.0	0	PASS
174	95	50	100	75.7	76827	PASS
175	174	5	9	7.1	5451	PASS
176	174	95	101	98.1	75403	PASS
177	176	5	9	5.6	4240	PASS

BFB 624 Results

C:\HPCHEM\1\DATA\061505\G8356.D Wed Jun 15 11:25:52 2005



Peak Apex is scan: 36
Average of 3 scans: 35,36,37 minus background scan 29

Target Mass	Comparison Mass	Lower Limit,%	Upper Limit,%	Relative Abundance,%	Result Pass/Fail
50	95	15	40	18.5	PASS
75	95	30	60	41.1	PASS
95	95	100	100	100.0	PASS
96	95	5	9	5.8	PASS
173	174	0	2	0.0	PASS
174	95	50	100	75.7	PASS
175	174	5	9	7.1	PASS
176	174	95	101	98.1	PASS
177	176	5	9	5.6	PASS

Form 1

CLIENT SAMPLE NO.

mblk061505

Lab Name: Toxikon Corporation Contract: _____

Lab Code: 10778 Case No.: LBG WHIT SAS No.: _____ SDG No.: 0506092

Matrix: (soil/water) Soil Lab Sample ID: mblk061505

Sample wt/vol: (g/mL) G Lab File ID: G8359.D

Level: (low/med) LOW Date Received:

% Moisture: not dec. Date Analyzed: 6/15/2005

GC Column: ID: (mm) Dilution Factor: 1.00

Extract Volume: _____ (µl)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) µg/Kg	Q
630-20-6	1,1,1,2-Tetrachloroethane	10	U
71-55-6	1,1,1-Trichloroethane	10	U
79-34-5	1,1,2,2-Tetrachloroethane	10	U
79-00-5	1,1,2-Trichloroethane	10	U
75-34-3	1,1-Dichloroethane	10	U
75-35-4	1,1-Dichloroethene	10	U
563-58-6	1,1-Dichloropropene	10	U
87-61-6	1,2,3-Trichlorobenzene	10	U
96-18-4	1,2,3-Trichloropropane	10	U
120-82-1	1,2,4-Trichlorobenzene	10	U
95-63-6	1,2,4-Trimethylbenzene	10	U
96-12-8	1,2-Dibromo-3-chloropropane	10	U
106-93-4	1,2-Dibromoethane	10	U
95-50-1	1,2-Dichlorobenzene	10	U
107-06-2	1,2-Dichloroethane	10	U
78-87-5	1,2-Dichloropropane	10	U
108-67-8	1,3,5-Trimethylbenzene	10	U
541-73-1	1,3-Dichlorobenzene	10	U
142-28-9	1,3-Dichloropropane	10	U
106-46-7	1,4-Dichlorobenzene	10	U
590-20-7	2,2-Dichloropropane	10	U
78-93-3	2-Butanone	10	U
110-75-8	2-Chloroethyl vinyl ether	10	U
95-49-8	2-Chlorotoluene	10	U
591-78-6	2-Hexanone	10	U
106-43-4	4-Chlorotoluene	10	U
99-87-6	4-Isopropyltoluene	10	U
108-10-1	4-Methyl-2-pentanone	10	U
67-64-1	Acetone	10	U
107-13-1	Acrylonitrile	100	U
71-43-2	Benzene	10	U
108-86-1	Bromobenzene	10	U
74-97-5	Bromochloromethane	10	U

Form 1

CLIENT SAMPLE NO.

mblk061505

Lab Name: Toxikon Corporation Contract: _____

Lab Code: 10778 Case No.: LBG WHIT SAS No.: _____ SDG No.: 0506092

Matrix: (soil/water) Soil Lab Sample ID: mblk061505

Sample wt/vol: (g/mL) G Lab File ID: G8359.D

Level: (low/med) LOW Date Received:

% Moisture: not dec. Date Analyzed: 6/15/2005

GC Column: ID: (mm) Dilution Factor: 1.00

Extract Volume: _____ (µl)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) µg/Kg	Q
75-27-4	Bromodichloromethane	10	U
75-25-2	Bromoform	10	U
74-83-9	Bromomethane	10	U
75-15-0	Carbon disulfide	10	U
56-23-5	Carbon tetrachloride	10	U
108-90-7	Chlorobenzene	10	U
75-00-3	Chloroethane	10	U
67-66-3	Chloroform	10	U
74-87-3	Chloromethane	10	U
156-59-2	cis-1,2-Dichloroethene	10	U
10061-01-5	cis-1,3-Dichloropropene	10	U
124-48-1	Dibromochloromethane	10	U
74-95-3	Dibromomethane	10	U
75-71-8	Dichlorodifluoromethane	10	U
100-41-4	Ethylbenzene	10	U
87-68-3	Hexachlorobutadiene	10	U
74-88-4	Iodomethane	10	U
98-82-8	Isopropylbenzene	10	U
1634-04-4	Methyl tert-butyl-ether	10	U
75-09-2	Methylene chloride	10	U
104-51-8	n-Butylbenzene	10	U
103-65-1	n-Propylbenzene	10	U
91-20-3	Naphthalene	10	U
135-98-8	sec-Butylbenzene	10	U
100-42-5	Styrene	10	U
98-06-6	tert-Butylbenzene	10	U
127-18-4	Tetrachloroethene	10	U
109-99-9	Tetrahydrofuran	10	U
108-88-3	Toluene	10	U
156-60-5	trans-1,2-Dichloroethene	10	U
10061-02-6	trans-1,3-Dichloropropene	10	U
79-01-6	Trichloroethene	10	U
75-69-4	Trichlorofluoromethane	10	U

CLIENT SAMPLE NO.

mblk061505

Lab Name: Toxikon Corporation Contract: _____

Lab Code: 10778 Case No.: LBG_WHIT SAS No.: _____ SDG No.: 0506092

Matrix: (soil/water) Soil Lab Sample ID: mblk061505

Sample wt/vol: (g/mL) G Lab File ID: G8359.D

Level: (low/med) LOW Date Received:

% Moisture: not dec. Date Analyzed: 6/15/2005

GC Column: ID: (mm) Dilution Factor: 1.00

Extract Volume: _____ (µl)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) µg/Kg	Q
108-05-4	Vinyl acetate	10	U
75-01-4	Vinyl chloride	10	U
1330-20-7	Xylenes, Total	10	U

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\061505\G8359.D
Acq On : 15 Jun 05 12:52
Sample : mblk061505
Misc : mblk asp_8260s
MS Integration Params: rteint.p
Quant Time: Jun 16 9:29 19105

Vial: 3
Operator: xl
Inst : inst g
Multiplr: 1.00

Quant Results File: 8260BASP.RE

Quant Method : C:\HPCHEM\1\DATA\061505\8260BASP.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Fri May 20 09:55:33 2005
Response via : Initial Calibration
DataAcq Meth : 8260BS

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) pentafluorobenzene	7.28	168	180195	50.00	ug/Kg	0.03
34) 1,4-difluorobenzene	8.40	114	398842	50.00	ug/Kg	0.04
52) chlorobenzene-d5	13.18	117	341565	50.00	ug/Kg	0.03
66) 1,4-dichlorobenzene-d4	17.30	152	137350	50.00	ug/Kg	0.03
System Monitoring Compounds						
31) Dibromofluoromethane	7.14	113	128843	54.81	ug/Kg	0.04
Spiked Amount 50.000	Range 80 - 120		Recovery = 109.62%			
48) toluene-d8	10.71	98	496763	50.71	ug/Kg	0.03
Spiked Amount 50.000	Range 81 - 120		Recovery = 101.42%			
65) 4-Bromofluorobenzene	15.26	95	169343	47.78	ug/Kg	0.03
Spiked Amount 50.000	Range 74 - 121		Recovery = 95.56%			

Target Compounds Qvalue

(#) = qualifier out of range (m) = manual integration

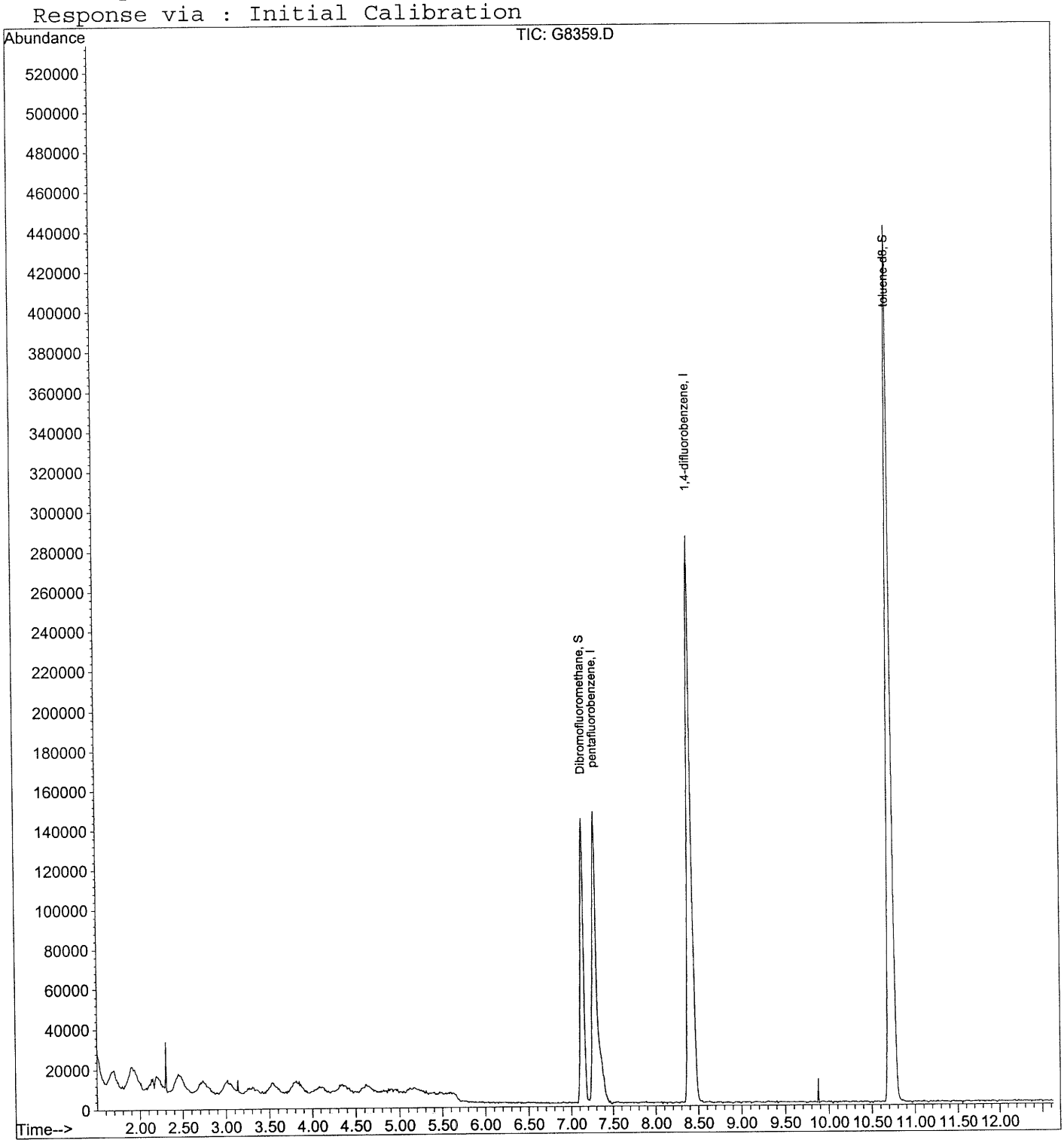
Quantitation Report

Data File : C:\HPCHEM\1\DATA\061505\G8359.D
Acq On : 15 Jun 05 12:52
Sample : mblk061505
Misc : mblk asp_8260s
MS Integration Params: rteint.p
Quant Time: Jun 16 9:29 19105

Vial: 3
Operator: xl
Inst : inst g
Multiplr: 1.00

Quant Results File: 8260BASP.RE

Method : C:\HPCHEM\1\DATA\061505\8260BASP.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Fri May 20 09:55:33 2005
Response via : Initial Calibration



000141

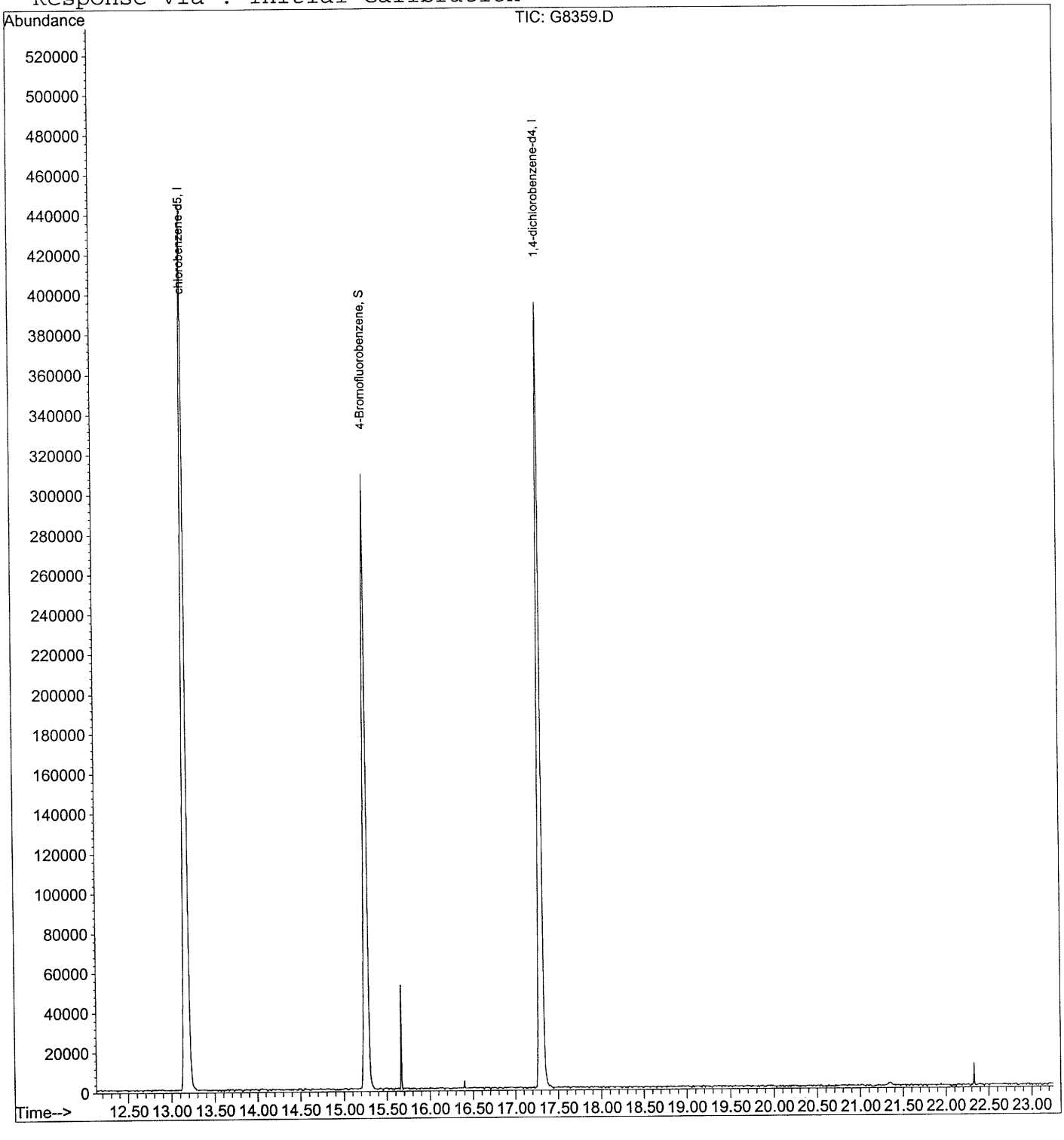
Quantitation Report

Data File : C:\HPCHEM\1\DATA\061505\G8359.D
Acq On : 15 Jun 05 12:52
Sample : mblk061505
Misc : mblk asp_8260s
MS Integration Params: rteint.p
Quant Time: Jun 16 9:29 19105

Vial: 3
Operator: xl
Inst : inst g
Multiplr: 1.00

Quant Results File: 8260BASP.RE

Method : C:\HPCHEM\1\DATA\061505\8260BASP.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Fri May 20 09:55:33 2005
Response via : Initial Calibration



000142

Form 1

CLIENT SAMPLE NO.

0506092-01ams

Lab Name: Toxikon Corporation Contract: _____

Lab Code: 10778 Case No.: LBG_WHIT SAS No.: _____ SDG No.: 0506092

Matrix: (soil/water) Soil Lab Sample ID: 0506092-01ams

Sample wt/vol: (g/mL) G Lab File ID: G8377.D

Level: (low/med) LOW Date Received: 6/10/2005

% Moisture: not dec. 14.6 Date Analyzed: 6/15/2005

GC Column: ID: (mm) Dilution Factor: 1.00

Extract Volume: _____ (µl)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) µg/Kg	Q
630-20-6	1,1,1,2-Tetrachloroethane	11.7	U
71-55-6	1,1,1-Trichloroethane	11.7	U
79-34-5	1,1,2,2-Tetrachloroethane	11.7	U
79-00-5	1,1,2-Trichloroethane	11.7	U
75-34-3	1,1-Dichloroethane	11.7	U
75-35-4	1,1-Dichloroethene	64.7	D
563-58-6	1,1-Dichloropropene	11.7	U
87-61-6	1,2,3-Trichlorobenzene	11.7	U
96-18-4	1,2,3-Trichloropropane	11.7	U
120-82-1	1,2,4-Trichlorobenzene	11.7	U
95-63-6	1,2,4-Trimethylbenzene	11.7	U
96-12-8	1,2-Dibromo-3-chloropropane	11.7	U
106-93-4	1,2-Dibromoethane	11.7	U
95-50-1	1,2-Dichlorobenzene	11.7	U
107-06-2	1,2-Dichloroethane	11.7	U
78-87-5	1,2-Dichloropropane	11.7	U
108-67-8	1,3,5-Trimethylbenzene	11.7	U
541-73-1	1,3-Dichlorobenzene	11.7	U
142-28-9	1,3-Dichloropropane	11.7	U
106-46-7	1,4-Dichlorobenzene	11.7	U
590-20-7	2,2-Dichloropropane	11.7	U
78-93-3	2-Butanone	11.7	U
110-75-8	2-Chloroethyl vinyl ether	11.7	U
95-49-8	2-Chlorotoluene	11.7	U
591-78-6	2-Hexanone	11.7	U
106-43-4	4-Chlorotoluene	11.7	U
99-87-6	4-Isopropyltoluene	11.7	U
108-10-1	4-Methyl-2-pentanone	11.7	U
67-64-1	Acetone	11.7	U
107-13-1	Acrylonitrile	117	U
71-43-2	Benzene	68.1	D
108-86-1	Bromobenzene	11.7	U
74-97-5	Bromochloromethane	11.7	U

Form 1

CLIENT SAMPLE NO.

0506092-01ams

Lab Name: Toxikon Corporation Contract: _____

Lab Code: 10778 Case No.: LBG WHIT SAS No.: _____ SDG No.: 0506092

Matrix: (soil/water) Soil Lab Sample ID: 0506092-01ams

Sample wt/vol: (g/mL) G Lab File ID: G8377.D

Level: (low/med) LOW Date Received: 6/10/2005

% Moisture: not dec. 14.6 Date Analyzed: 6/15/2005

GC Column: ID: (mm) Dilution Factor: 1.00

Extract Volume: _____ (µl)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) µg/Kg	Q
75-27-4	Bromodichloromethane	11.7	U
75-25-2	Bromoform	11.7	U
74-83-9	Bromomethane	11.7	U
75-15-0	Carbon disulfide	11.7	U
56-23-5	Carbon tetrachloride	11.7	U
108-90-7	Chlorobenzene	66.7	D
75-00-3	Chloroethane	11.7	U
67-66-3	Chloroform	11.7	U
74-87-3	Chloromethane	11.7	U
156-59-2	cis-1,2-Dichloroethene	11.7	U
10061-01-5	cis-1,3-Dichloropropene	11.7	U
124-48-1	Dibromochloromethane	11.7	U
74-95-3	Dibromomethane	11.7	U
75-71-8	Dichlorodifluoromethane	11.7	U
100-41-4	Ethylbenzene	11.7	U
87-68-3	Hexachlorobutadiene	11.7	U
74-88-4	Iodomethane	11.7	U
98-82-8	Isopropylbenzene	11.7	U
1634-04-4	Methyl tert-butyl-ether	11.7	U
75-09-2	Methylene chloride	11.7	U
104-51-8	n-Butylbenzene	11.7	U
103-65-1	n-Propylbenzene	11.7	U
91-20-3	Naphthalene	11.7	U
135-98-8	sec-Butylbenzene	11.7	U
100-42-5	Styrene	11.7	U
98-06-6	tert-Butylbenzene	11.7	U
127-18-4	Tetrachloroethene	11.7	U
109-99-9	Tetrahydrofuran	11.7	U
108-88-3	Toluene	67.1	D
156-60-5	trans-1,2-Dichloroethene	11.7	U
10061-02-6	trans-1,3-Dichloropropene	11.7	U
79-01-6	Trichloroethene	65.9	D
75-69-4	Trichlorofluoromethane	11.7	U

Form 1

CLIENT SAMPLE NO.

0506092-01ams

Lab Name: Toxikon Corporation Contract: _____

Lab Code: 10778 Case No.: LBG_WHIT SAS No.: _____ SDG No.: 0506092

Matrix: (soil/water) Soil Lab Sample ID: 0506092-01ams

Sample wt/vol: (g/mL) G Lab File ID: G8377.D

Level: (low/med) LOW Date Received: 6/10/2005

% Moisture: not dec. 14.6 Date Analyzed: 6/15/2005

GC Column: ID: (mm) Dilution Factor: 1.00

Extract Volume: _____ (µl)

CONCENTRATION UNITS:

CAS NO.	COMPOUND	(ug/L or ug/Kg) µg/Kg	Q
108-05-4	Vinyl acetate	11.7	U
75-01-4	Vinyl chloride	11.7	U
1330-20-7	Xylenes, Total	11.7	U

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\061505\G8377.D Vial: 15
Acq On : 15 Jun 05 23:00 Operator: xl
Sample : 0506092-01ams Inst : inst g
Misc : ms asp_8260s Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: Jun 16 9:26 19105 Quant Results File: 8260BASP.RE

Quant Method : C:\HPCHEM\1\DATA\061505\8260BASP.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Fri May 20 09:55:33 2005
Response via : Initial Calibration
DataAcq Meth : 8260BASP

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) pentafluorobenzene	7.28	168	141776	50.00	ug/Kg	0.04
34) 1,4-difluorobenzene	8.41	114	325131	50.00	ug/Kg	0.05
52) chlorobenzene-d5	13.18	117	279993	50.00	ug/Kg	0.04
66) 1,4-dichlorobenzene-d4	17.31	152	109569	50.00	ug/Kg	0.04
System Monitoring Compounds						
31) Dibromofluoromethane	7.14	113	101205	54.72	ug/Kg	0.05
Spiked Amount 50.000	Range 80 - 120		Recovery = 109.44%			
48) toluene-d8	10.72	98	406677	50.92	ug/Kg	0.04
Spiked Amount 50.000	Range 81 - 120		Recovery = 101.84%			
65) 4-Bromofluorobenzene	15.26	95	130426	44.89	ug/Kg	0.04
Spiked Amount 50.000	Range 74 - 121		Recovery = 89.78%			
Target Compounds						Qvalue
11) 1,1-dichloroethene	3.34	96	175592	55.23	ug/Kg	86
37) benzene	7.64	78	736832	58.12	ug/Kg	97
40) trichloroethene	8.74	95	148579	56.28	ug/Kg	97
49) toluene	10.83	92	392768	57.28	ug/Kg	99
58) chlorobenzene	13.22	112	360878	57.00	ug/Kg	97

(#) = qualifier out of range (m) = manual integration

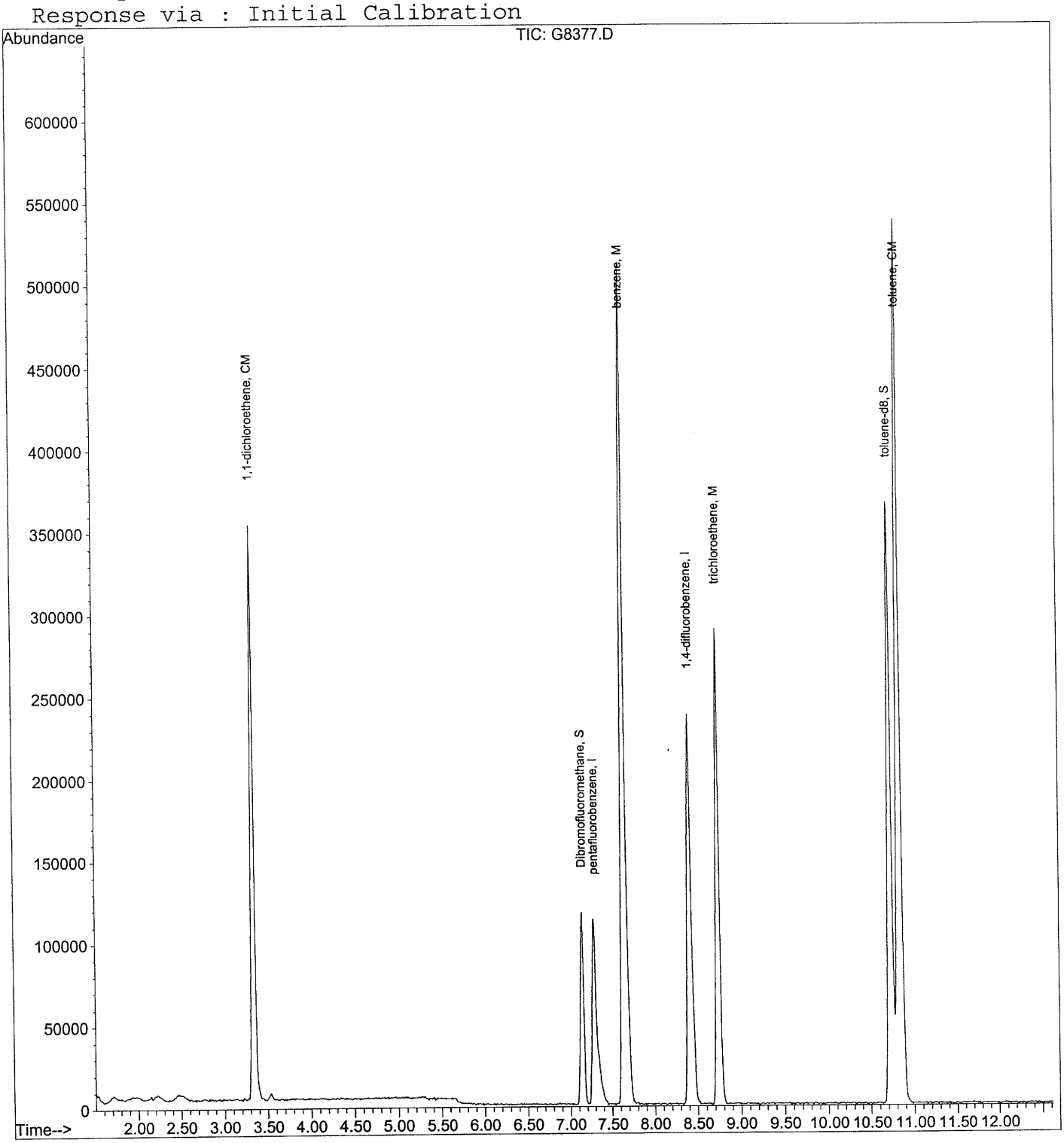
Quantitation Report

Data File : C:\HPCHEM\1\DATA\061505\G8377.D
Acq On : 15 Jun 05 23:00
Sample : 0506092-01ams
Misc : ms asp_8260s
MS Integration Params: rteint.p
Quant Time: Jun 16 9:26 19105

Vial: 15
Operator: xl
Inst : inst g
Multiplr: 1.00

Quant Results File: 8260BASP.RE

Method : C:\HPCHEM\1\DATA\061505\8260BASP.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Fri May 20 09:55:33 2005
Response via : Initial Calibration



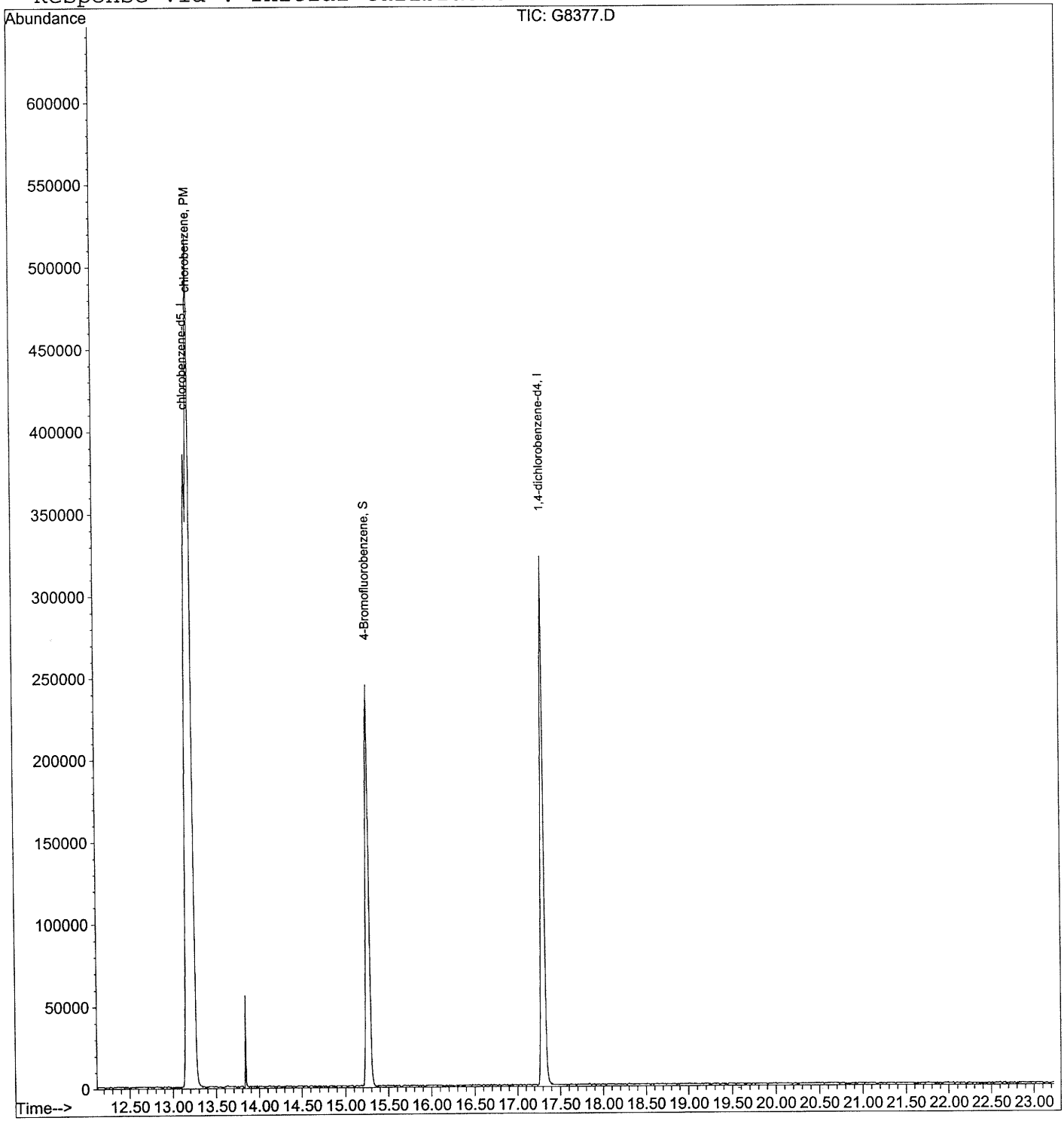
Quantitation Report

Data File : C:\HPCHEM\1\DATA\061505\G8377.D
Acq On : 15 Jun 05 23:00
Sample : 0506092-01ams
Misc : ms asp_8260s
MS Integration Params: rteint.p
Quant Time: Jun 16 9:26 19105

Vial: 15
Operator: xl
Inst : inst g
Multiplr: 1.00

Quant Results File: 8260BASP.RE

Method : C:\HPCHEM\1\DATA\061505\8260BASP.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Fri May 20 09:55:33 2005
Response via : Initial Calibration



Form 1

CLIENT SAMPLE NO.

0506092-01amsd

Lab Name: Toxikon Corporation Contract: _____

Lab Code: 10778 Case No.: LBG WHIT SAS No.: _____ SDG No.: 0506092

Matrix: (soil/water) Soil Lab Sample ID: 0506092-01amsd

Sample wt/vol: (g/mL) G Lab File ID: G8374.D

Level: (low/med) LOW Date Received: 6/10/2005

% Moisture: not dec. 14.6 Date Analyzed: 6/15/2005

GC Column: ID: (mm) Dilution Factor: 1.00

Extract Volume: _____ (µl)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) µg/Kg	Q
630-20-6	1,1,1,2-Tetrachloroethane	11.7	U
71-55-6	1,1,1-Trichloroethane	11.7	U
79-34-5	1,1,2,2-Tetrachloroethane	11.7	U
79-00-5	1,1,2-Trichloroethane	11.7	U
75-34-3	1,1-Dichloroethane	11.7	U
75-35-4	1,1-Dichloroethene	67.3	D
563-58-6	1,1-Dichloropropene	11.7	U
87-61-6	1,2,3-Trichlorobenzene	11.7	U
96-18-4	1,2,3-Trichloropropane	11.7	U
120-82-1	1,2,4-Trichlorobenzene	11.7	U
95-63-6	1,2,4-Trimethylbenzene	11.7	U
96-12-8	1,2-Dibromo-3-chloropropane	11.7	U
106-93-4	1,2-Dibromoethane	11.7	U
95-50-1	1,2-Dichlorobenzene	11.7	U
107-06-2	1,2-Dichloroethane	11.7	U
78-87-5	1,2-Dichloropropane	11.7	U
108-67-8	1,3,5-Trimethylbenzene	11.7	U
541-73-1	1,3-Dichlorobenzene	11.7	U
142-28-9	1,3-Dichloropropane	11.7	U
106-46-7	1,4-Dichlorobenzene	11.7	U
590-20-7	2,2-Dichloropropane	11.7	U
78-93-3	2-Butanone	11.7	U
110-75-8	2-Chloroethyl vinyl ether	11.7	U
95-49-8	2-Chlorotoluene	11.7	U
591-78-6	2-Hexanone	11.7	U
106-43-4	4-Chlorotoluene	11.7	U
99-87-6	4-Isopropyltoluene	11.7	U
108-10-1	4-Methyl-2-pentanone	11.7	U
67-64-1	Acetone	10	DJ
107-13-1	Acrylonitrile	117	U
71-43-2	Benzene	69.8	D
108-86-1	Bromobenzene	11.7	U
74-97-5	Bromochloromethane	11.7	U

Form 1

CLIENT SAMPLE NO.

0506092-01amsd

Lab Name: Toxikon Corporation Contract: _____

Lab Code: 10778 Case No.: LBG WHIT SAS No.: _____ SDG No.: 0506092

Matrix: (soil/water) Soil Lab Sample ID: 0506092-01amsd

Sample wt/vol: (g/mL) G Lab File ID: G8374.D

Level: (low/med) LOW Date Received: 6/10/2005

% Moisture: not dec. 14.6 Date Analyzed: 6/15/2005

GC Column: ID: (mm) Dilution Factor: 1.00

Extract Volume: _____ (µl)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) µg/Kg	Q
75-27-4	Bromodichloromethane	11.7	U
75-25-2	Bromoform	11.7	U
74-83-9	Bromomethane	11.7	U
75-15-0	Carbon disulfide	11.7	U
56-23-5	Carbon tetrachloride	11.7	U
108-90-7	Chlorobenzene	67.8	D
75-00-3	Chloroethane	11.7	U
67-66-3	Chloroform	11.7	U
74-87-3	Chloromethane	11.7	U
156-59-2	cis-1,2-Dichloroethene	11.7	U
10061-01-5	cis-1,3-Dichloropropene	11.7	U
124-48-1	Dibromochloromethane	11.7	U
74-95-3	Dibromomethane	11.7	U
75-71-8	Dichlorodifluoromethane	11.7	U
100-41-4	Ethylbenzene	11.7	U
87-68-3	Hexachlorobutadiene	11.7	U
74-88-4	Iodomethane	11.7	U
98-82-8	Isopropylbenzene	11.7	U
1634-04-4	Methyl tert-butyl-ether	11.7	U
75-09-2	Methylene chloride	11.7	U
104-51-8	n-Butylbenzene	11.7	U
103-65-1	n-Propylbenzene	11.7	U
91-20-3	Naphthalene	11.7	U
135-98-8	sec-Butylbenzene	11.7	U
100-42-5	Styrene	11.7	U
98-06-6	tert-Butylbenzene	11.7	U
127-18-4	Tetrachloroethene	3.6	DJ
109-99-9	Tetrahydrofuran	11.7	U
108-88-3	Toluene	68.4	D
156-60-5	trans-1,2-Dichloroethene	11.7	U
10061-02-6	trans-1,3-Dichloropropene	11.7	U
79-01-6	Trichloroethene	66.6	D
75-69-4	Trichlorofluoromethane	11.7	U

CLIENT SAMPLE NO.

0506092-01amsd

Lab Name: Toxikon Corporation Contract: _____

Lab Code: 10778 Case No.: LBG WHIT SAS No.: _____ SDG No.: 0506092

Matrix: (soil/water) Soil Lab Sample ID: 0506092-01amsd

Sample wt/vol: (g/mL) G Lab File ID: G8374.D

Level: (low/med) LOW Date Received: 6/10/2005

% Moisture: not dec. 14.6 Date Analyzed: 6/15/2005

GC Column: ID: (mm) Dilution Factor: 1.00

Extract Volume: _____ (µl)

CONCENTRATION UNITS:

CAS NO.	COMPOUND	(ug/L or ug/Kg) µg/Kg	Q
108-05-4	Vinyl acetate	11.7	U
75-01-4	Vinyl chloride	11.7	U
1330-20-7	Xylenes, Total	11.7	U

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\061505\G8374.D Vial: 15
Acq On : 15 Jun 05 21:27 Operator: xl
Sample : 0506092-01amsd Inst : inst g
Misc : msd asp_8260s Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: Jun 16 9:24 19105 Quant Results File: 8260BASP.RE

Quant Method : C:\HPCHEM\1\DATA\061505\8260BASP.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Fri May 20 09:55:33 2005
Response via : Initial Calibration
DataAcq Meth : 8260BASP

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) pentafluorobenzene	7.28	168	126496	50.00	ug/Kg	0.03
34) 1,4-difluorobenzene	8.40	114	291055	50.00	ug/Kg	0.04
52) chlorobenzene-d5	13.18	117	253535	50.00	ug/Kg	0.03
66) 1,4-dichlorobenzene-d4	17.30	152	96314	50.00	ug/Kg	0.03
System Monitoring Compounds						
31) Dibromofluoromethane	7.14	113	92587	56.11	ug/Kg	0.04
Spiked Amount 50.000	Range 80 - 120		Recovery = 112.22%			
48) toluene-d8	10.72	98	363507	50.85	ug/Kg	0.04
Spiked Amount 50.000	Range 81 - 120		Recovery = 101.70%			
65) 4-Bromofluorobenzene	15.26	95	121661	46.25	ug/Kg	0.03
Spiked Amount 50.000	Range 74 - 121		Recovery = 92.50%			
Target Compounds						Qvalue
10) Acetone	3.53	43	10240	8.66	ug/Kg	97
11) 1,1-dichloroethene	3.34	96	163012	57.46	ug/Kg	92
37) benzene	7.64	78	676712	59.62	ug/Kg	99
40) trichloroethene	8.74	95	134377	56.86	ug/Kg	97
49) toluene	10.82	92	358397	58.38	ug/Kg	99
54) tetrachloroethene	11.71	166	5867	3.11	ug/Kg	90
58) chlorobenzene	13.23	112	331861	57.89	ug/Kg	99

(#) = qualifier out of range (m) = manual integration

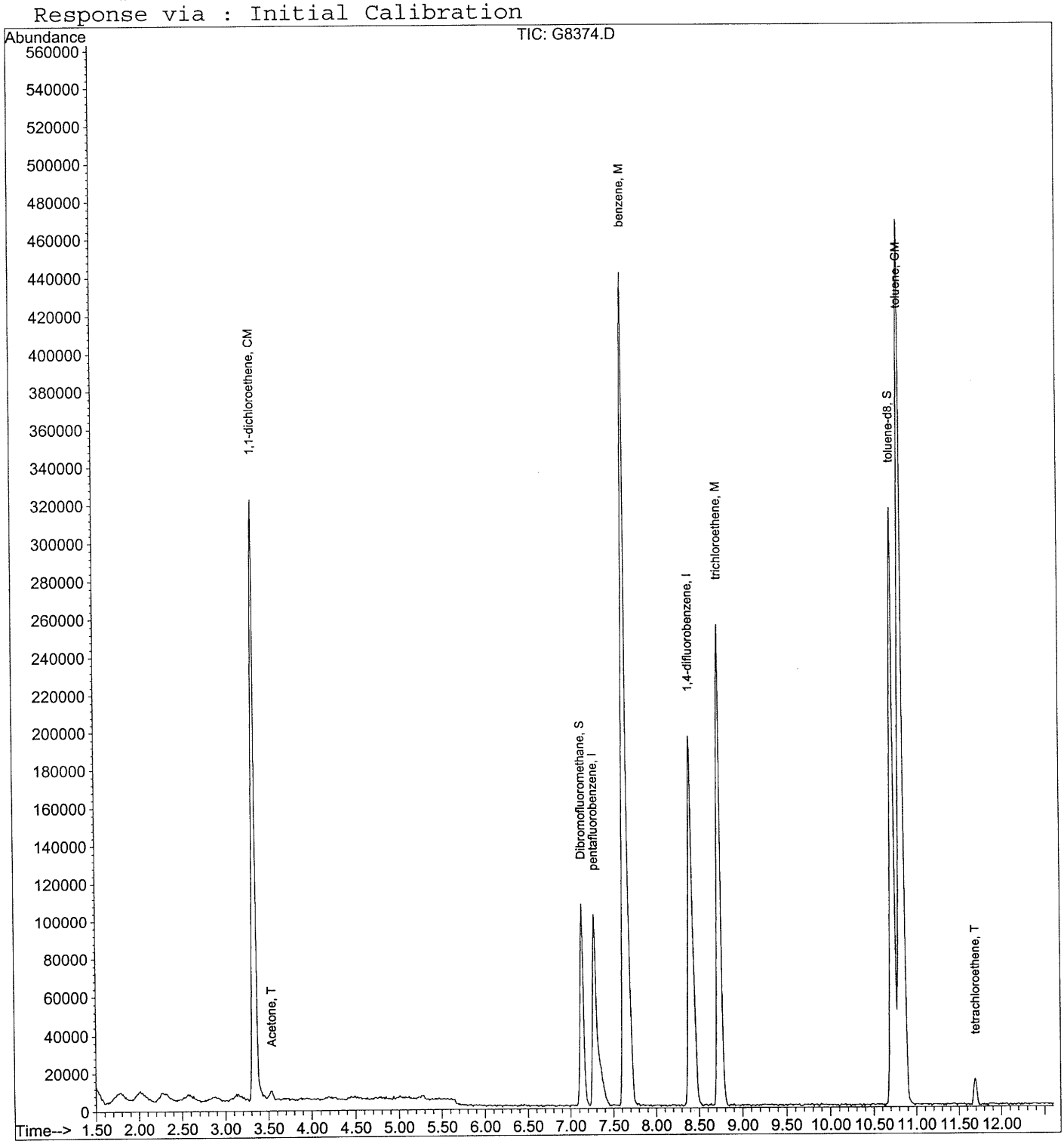
Quantitation Report

Data File : C:\HPCHEM\1\DATA\061505\G8374.D
Acq On : 15 Jun 05 21:27
Sample : 0506092-01amsd
Misc : msd asp_8260s
MS Integration Params: rteint.p
Quant Time: Jun 16 9:24 19105

Vial: 15
Operator: xl
Inst : inst g
Multiplr: 1.00

Quant Results File: 8260BASP.RE

Method : C:\HPCHEM\1\DATA\061505\8260BASP.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Fri May 20 09:55:33 2005
Response via : Initial Calibration

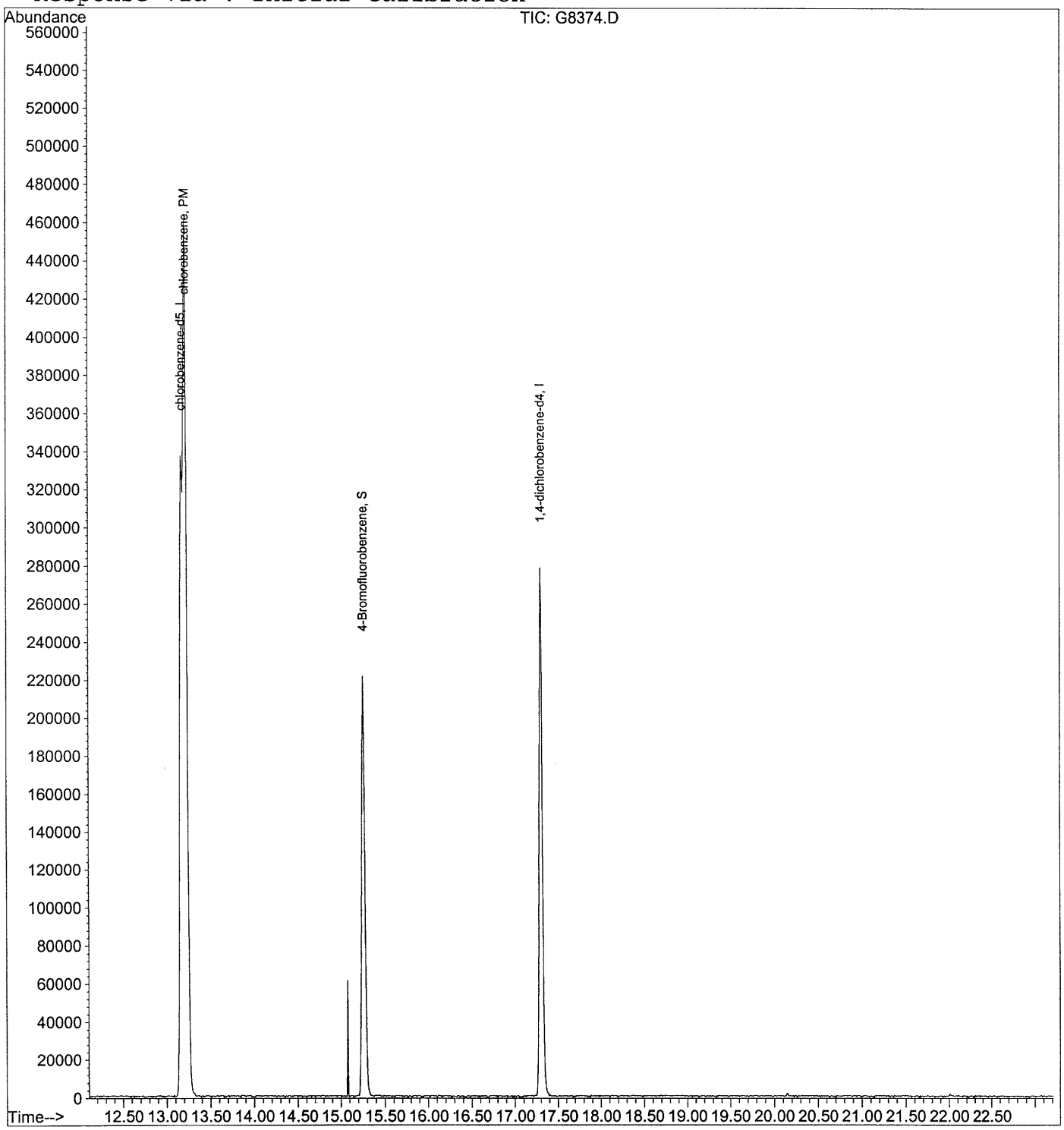


000153

Quantitation Report

Data File : C:\HPCHEM\1\DATA\061505\G8374.D Vial: 15
Acq On : 15 Jun 05 21:27 Operator: xl
Sample : 0506092-01amsd Inst : inst g
Misc : msd asp_8260s Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: Jun 16 9:24 19105 Quant Results File: 8260BASP.RE

Method : C:\HPCHEM\1\DATA\061505\8260BASP.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Fri May 20 09:55:33 2005
Response via : Initial Calibration



000154

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\061505\G8361.D
Acq On : 15 Jun 05 14:32
Sample : lcs061505
Misc : lcs asp_8260s
MS Integration Params: rteint.p
Quant Time: Jun 16 9:30 19105

Vial: 3
Operator: xl
Inst : inst g
Multiplr: 1.00

Quant Results File: 8260BS.RES

Quant Method : C:\HPCHEM\1\DATA\061505\8260BS.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Tue May 17 11:42:05 2005
Response via : Initial Calibration
DataAcq Meth : 8260BS

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) pentafluorobenzene	7.28	168	190618	50.00	ug/Kg	0.03
34) 1,4-difluorobenzene	8.40	114	418448	50.00	ug/Kg	0.04
52) chlorobenzene-d5	13.17	117	352958	50.00	ug/Kg	0.03
66) 1,4-dichlorobenzene-d4	17.30	152	161585	50.00	ug/Kg	0.03
System Monitoring Compounds						
31) Dibromofluoromethane	7.13	113	128680	51.56	ug/Kg	0.04
Spiked Amount 50.000	Range 80 - 120		Recovery = 103.12%			
48) toluene-d8	10.71	98	517154	50.50	ug/Kg	0.03
Spiked Amount 50.000	Range 81 - 120		Recovery = 101.00%			
65) 4-Bromofluorobenzene	15.26	95	181739	49.87	ug/Kg	0.03
Spiked Amount 50.000	Range 74 - 121		Recovery = 99.74%			
Target Compounds						Qvalue
2) dichlorodifluoromethane	1.54	85	170120	43.75	ug/Kg	98
3) chloromethane	1.75	50	343934	52.97	ug/Kg	99
4) vinyl chloride	1.86	62	367574	52.77	ug/Kg	99
5) bromomethane	2.23	96	208916	56.00	ug/Kg	100
6) chloroethane	2.36	64	212206	55.22	ug/Kg	97
7) trichlorofluoromethane	2.63	101	360787	52.44	ug/Kg	100
8) Diethyl Ether	3.03	45	101740	56.45	ug/Kg	97
9) Acrolein	3.29	56	73272	201.62	ug/Kg	94
10) Acetone	3.52	43	61632	62.04	ug/Kg	100
11) 1,1-dichloroethene	3.33	96	214281	51.60	ug/Kg	91
12) Iodomethane	3.56	142	310004	50.35	ug/Kg	94
13) methylene chloride	4.18	84	249261	51.06	ug/Kg	92
14) Carbon Disulfide	3.60	76	934602	52.39	ug/Kg	100
17) Acrylonitrile	4.71	53	58661	52.62	ug/Kg	98
18) tert-Butyl alcohol	4.51	59	121576	400.49	ug/Kg	100
19) methyl tert-butyl ether	4.57	73	506144	48.94	ug/Kg	100
20) trans-1,2-dichloroethene	4.56	96	244917	52.36	ug/Kg	91
21) n-Hexane	4.92	57	435179m	54.75	ug/Kg	16
22) 1,1-dichloroethane	5.34	63	462070	54.85	ug/Kg	99
23) di-isopropyl ether	5.41	45	835494	52.15	ug/Kg	98
24) Vinyl Acetate	5.46	43	403127	47.60	ug/Kg	100
25) ethyl tert-butyl ether	6.00	59	495395	48.53	ug/Kg	99
26) 2-Butanone	6.42	43	78906	52.84	ug/Kg	99

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\061505\G8361.D Vial: 3
Acq On : 15 Jun 05 14:32 Operator: xl
Sample : lcs061505 Inst : inst g
Misc : lcs asp_8260s Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: Jun 16 9:30 19105 Quant Results File: 8260BS.RES

Quant Method : C:\HPCHEM\1\DATA\061505\8260BS.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Tue May 17 11:42:05 2005
Response via : Initial Calibration
DataAcq Meth : 8260BS

	Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
27)	2,2-dichloropropane	6.25	77	278791	51.49	ug/Kg	97
28)	cis-1,2-dichloroethene	6.32	96	181108	56.57	ug/Kg	99
29)	bromochloromethane	6.71	128	63200	52.95	ug/Kg	99
30)	chloroform	6.87	83	348143	54.95	ug/Kg	100
32)	Tetrahydrofuran	6.74	42	40250	51.82	ug/Kg	95
33)	1,1,1-trichloroethane	7.04	97	238776	52.02	ug/Kg	98
35)	carbon tetrachloride	7.25	117	186902	45.66	ug/Kg	98
36)	1,1-dichloropropene	7.33	75	326816	48.30	ug/Kg	98
37)	benzene	7.63	78	865339	54.96	ug/Kg	99
38)	1,2-dichloroethane	7.81	62	242055	47.44	ug/Kg	98
39)	tert amyl methyl ether	7.86	73	399094	45.11	ug/Kg	99
40)	trichloroethene	8.73	95	172174	52.60	ug/Kg	98
41)	1,2-dichloropropane	9.16	63	193115	53.43	ug/Kg	100
42)	dibromomethane	9.34	93	94575	49.72	ug/Kg	96
44)	bromodichloromethane	9.63	83	230527	50.59	ug/Kg	100
45)	2-Chloroethyl vinyl ether	10.19	63	76119	55.40	ug/Kg	99
46)	4-Methyl-2-Pentanone	10.66	43	178291	50.88	ug/Kg	99
47)	cis-1,3-dichloropropene	10.37	75	291726	50.80	ug/Kg	99
49)	toluene	10.82	92	450491	53.20	ug/Kg	100
50)	trans-1,3-dichloropropene	11.36	75	238262	48.23	ug/Kg	96
51)	1,1,2-trichloroethane	11.65	83	112887	50.99	ug/Kg	96
53)	2-Hexanone	12.07	43	108999	52.60	ug/Kg	96
54)	tetrachloroethene	11.69	166	126997	49.98	ug/Kg	96
55)	1,3-dichloropropane	11.92	76	251532	51.08	ug/Kg	99
56)	dibromochloromethane	12.23	129	113392	50.78	ug/Kg	97
57)	1,2-dibromoethane	12.40	107	107469	49.80	ug/Kg	98
58)	chlorobenzene	13.22	112	403572	52.47	ug/Kg	97
59)	1,1,1,2-tetrachloroethane	13.39	131	124433	46.48	ug/Kg	99
60)	ethylbenzene	13.39	91	867551	46.25	ug/Kg	97
61)	m,p-xylene	13.61	106	592751	104.39	ug/Kg	96
62)	o-xylene	14.29	106	258945	47.23	ug/Kg	99
63)	styrene	14.34	104	475072	46.54	ug/Kg	96
64)	bromoform	14.66	173	54294	44.97	ug/Kg	98
67)	isopropylbenzene	14.93	105	750791	51.23	ug/Kg	99
68)	bromobenzene	15.47	156	135566	50.10	ug/Kg	95
69)	1,1,2,2-tetrachloroethane	15.62	83	164559m	52.18	ug/Kg	98
70)	1,2,3-trichloropropane	15.66	75	131019m	51.42	ug/Kg	99

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\061505\G8361.D
Acq On : 15 Jun 05 14:32
Sample : lcs061505
Misc : lcs asp_8260s
MS Integration Params: rteint.p
Quant Time: Jun 16 9:30 19105

Vial: 3
Operator: xl
Inst : inst g
Multiplr: 1.00

Quant Results File: 8260BS.RES

Quant Method : C:\HPCHEM\1\DATA\061505\8260BS.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Tue May 17 11:42:05 2005
Response via : Initial Calibration
DataAcq Meth : 8260BS

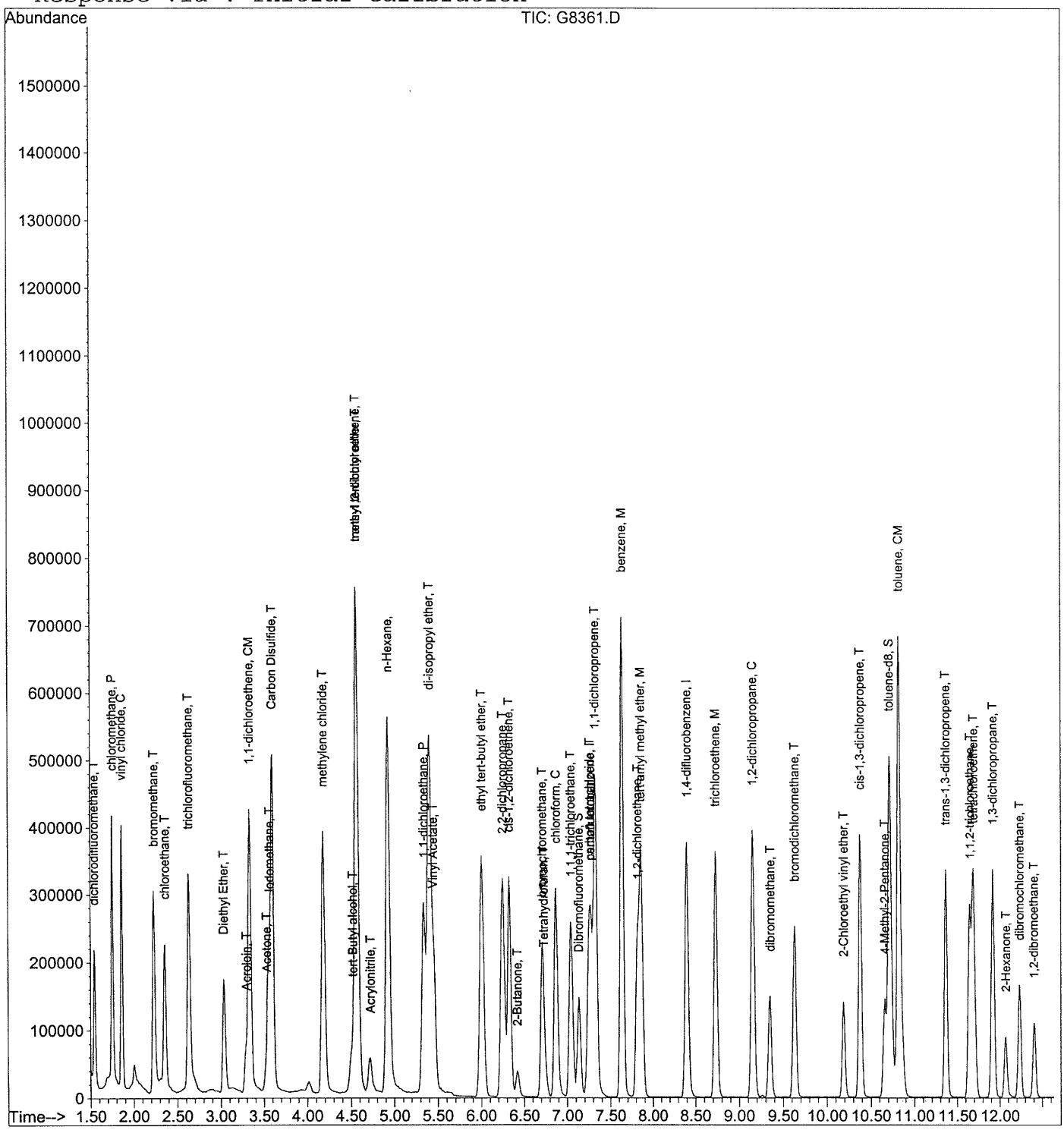
Compound	R.T.	QIon	Response	Conc Unit	Qvalue
71) n-propylbenzene	15.66	91	1120428	53.33 ug/Kg	99
72) 2-chlorotoluene	15.82	91	519428	48.10 ug/Kg	98
73) 4-chlorotoluene	16.03	91	686579	52.57 ug/Kg	98
74) 1,3,5-trimethylbenzene	16.00	105	596321	46.33 ug/Kg	98
75) tert-butylbenzene	16.54	91	376057	51.81 ug/Kg	99
76) 1,2,4-trimethylbenzene	16.66	105	591263	47.35 ug/Kg	98
77) sec-butylbenzene	16.94	105	842999	52.64 ug/Kg	99
78) 1,3-dichlorobenzene	17.17	146	289985	51.09 ug/Kg	99
79) 4-isopropyltoluene	17.22	119	615903	44.99 ug/Kg	99
80) 1,4-dichlorobenzene	17.34	146	298151	51.51 ug/Kg	99
81) 1,2-dichlorobenzene	18.00	146	263034	50.86 ug/Kg	98
82) n-butylbenzene	17.96	91	814568	54.61 ug/Kg	99
83) 1,2-dibromo-3-chloropropan	19.45	75	18632	43.06 ug/Kg	93
84) 1,2,4-trichlorobenzene	20.91	180	140733	50.21 ug/Kg	99
85) hexachlorobutadiene	21.17	225	66966m	47.11 ug/Kg	100
86) naphthalene	21.33	128	330783	52.23 ug/Kg	100
87) 1,2,3-trichlorobenzene	21.80	180	124336	51.87 ug/Kg	96

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : C:\HPCHEM\1\DATA\061505\G8361.D Vial: 3
Acq On : 15 Jun 05 14:32 Operator: xl
Sample : lcs061505 Inst : inst g
Misc : lcs asp_8260s Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: Jun 16 9:30 19105 Quant Results File: 8260BS.RES

Method : C:\HPCHEM\1\DATA\061505\8260BASP.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Fri May 20 09:55:33 2005
Response via : Initial Calibration



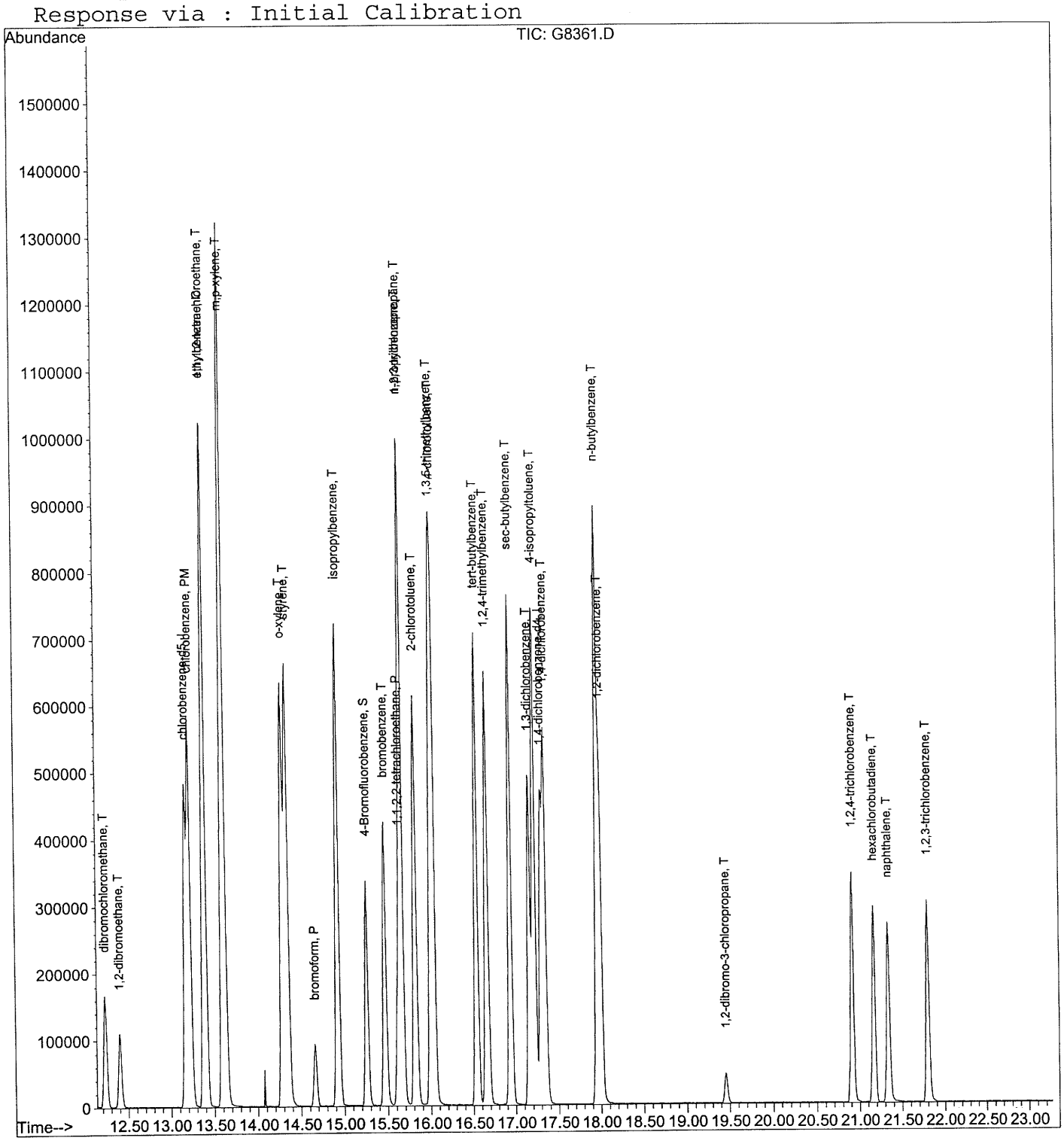
Quantitation Report

Data File : C:\HPCHEM\1\DATA\061505\G8361.D
Acq On : 15 Jun 05 14:32
Sample : lcs061505
Misc : lcs asp_8260s
MS Integration Params: rteint.p
Quant Time: Jun 16 9:30 19105

Vial: 3
Operator: xl
Inst : inst g
Multiplr: 1.00

Quant Results File: 8260BS.RES

Method : C:\HPCHEM\1\DATA\061505\8260BASP.M (RTE Integrator)
Title : SW-846 Method 8260B
Last Update : Fri May 20 09:55:33 2005
Response via : Initial Calibration



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LOGBOOK PAGES

TOXIKON GC/MS VOLATILE ANALYSIS LOG INSTRUMENT G

MAINTENANCE PERFORMED:

WORKING STANDARD ID (enter when a new standard is used):

	DATE	DATAFILE	CLIENT	LAB ID	SEQ	DILUTION	WT	ALS	COMMENTS	OPERATOR
1	6/15/05	G8356		50 ng BFB				DE		XL
2	6/15/05	G8357		50 ng 8260 std				2	8260-S, BSP-8260S	XL
3	6/15/05	G8358	✓	50 ng 8260 std				2		XL
4	6/15/05	G8359		MBLK 061505				3		XL
5	6/15/05	G8360		LCS 061505				2		XL
6	6/15/05	G8361		LCS 061505				2		XL
7	6/15/05	G8362		0506092-01A		NONE		3	8260-S	XL
8	6/15/05	G8363		0506092-01A		NONE		4	BSP-8260S	XL
9	6/15/05	G8364		0506092-02A		NONE		7	BSP-8260S	XL
10	6/15/05	G8365		0506092-03A		NONE		8	BSP-8260S	XL
11	6/15/05	G8366		0506092-04A		NONE		10	BSP-8260S	XL
12	6/15/05	G8367		0506092-05A		NONE		11	BSP-8260S	XL
13	6/15/05	G8368		0506092-06A		NONE		13	BSP-8260S	XL
14	6/15/05	G8369		0506092-01AMS		NONE		14	BSP-8260S	XL
15	6/15/05	G8370		0506092-01AMSD		NONE		15	BSP-8260S	XL
16	6/15/05	G8371		0506092-02A	RE	1:5		16	BSP-8260S	XL
17	6/15/05	G8372		0506092-03A	RE	NONE		3	BSP-8260S	XL
18	6/15/05	G8373		0506092-01AMS				14	BSP-8260S	XL
19	6/15/05	G8374		0506092-01AMSD				15	BSP-8260S	XL
20	6/15/05	G8375		0506092-06A	RE	NONE		16	BSP-8260S	XL
21	6/15/05	G8376		0506092-03A	RE	1:2		3	BSP-8260S	XL
22	6/15/05	G8377		0506092-01AMS		NONE		15	BSP-8260S	XL
23	6/15/05	G8378	out of time times	0506092-04A		NONE		4	BSP-8260S	XL
24	6/15/05	G8379		0506092-05A		NONE		7	BSP-8260S	XL
25										
26						XL				
27										

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Figure 1

DRY WEIGHT AND % MOISTURE DETERMINATION

Analyzed By/On: Am 6/9/05
Balance Used: 01/A 33

Sample ID #	Gross Wt. Initial (g)	Tare (g)	1 st Gross Wt. Final (g)	2nd Gross Wt. Final (g)	% Dry Weight	% Moisture
0506065 3	7.23	1.19	2.03	2.03	14.0	86.0
DUP 065 3	7.25	1.19	2.02	2.02	13.8	86.2
0506065.4	9.26	1.18	2.60	2.60	17.5	82.5
0506076.1	7.65	1.18	6.80	6.80	86.9	NT ~~13.0~~ 13.1
.2	6.11	1.18	5.90	5.90	95.7	4.3
.3	8.88	1.18	8.73	8.73	98.1	1.9
.4	6.80	1.18	6.64	6.64	97.2	2.8
0506086.1	9.21	1.19	8.16	8.16	86.9	13.1
.2	10.55	1.19	9.91	9.91	93.2	6.8
.3	7.85	1.20	6.91	6.91	85.9	14.1
.4	9.01	1.19	7.92	7.92	86.1	13.9
0506085.1	8.76	1.18	6.99	6.99	76.6	23.4
0506101.1	6.95	1.21	6.12	6.12	85.5	14.5
2	5.72	1.18	3.31	3.31	46.9	53.1
0506112.1	9.68	1.21	9.31	9.31	95.6	4.4
06077.1	7.12	1.18	5.51	5.51	72.9	27.1
06092.1	7.71	1.19	6.76	6.76	85.4	14.6
.2	9.31	1.19	7.99	7.99	83.7	16.3
.3	8.69	1.18	7.46	7.46	83.6	16.4
.4	9.78	1.19	8.44	8.44	84.4	15.6
.5	8.93	1.20	7.61	7.61	82.9	17.1
.6	6.93	1.18	6.17	6.17	86.8	13.2

Am 6/13/05 9 a.m

nm 6/15/05 9am

NT 6/15/05 @ 4:50 pm

NT 6/16/05 @ 3:45 pm

Calc % Dry Weight: [(Gross Wt. Final-Tare)/(Gross Wt. Initial-Tare)] X 100 = %
% Moisture is 100 - % Dry Weight