

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM032619\
 Data File : BM019477.D
 Acq On : 26 Mar 2019 19:15
 Operator : JU/SJ
 Sample : K2069-11
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C09D0

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 0.1 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.952	326	334	361	rBV	29111616	45239150	51.70%	16.830%
2	6.781	641	645	658	rBV	130944	236527	0.27%	0.088%
3	6.951	669	674	689	rBV	473764	793293	0.91%	0.295%
4	7.134	699	705	718	rBV	563052	893364	1.02%	0.332%
5	7.487	757	765	778	rBV	3873650	5877266	6.72%	2.187%
6	7.640	778	791	818	rVB	14980619	24623656	28.14%	9.161%
7	7.975	835	848	853	rBV2	38780332	87500939	100.00%	32.553%
8	8.316	900	906	922	rBV	343217	639198	0.73%	0.238%
9	8.769	976	983	998	rBV	546421	960216	1.10%	0.357%
10	9.087	1031	1037	1047	rBV2	237602	410013	0.47%	0.153%
11	9.310	1070	1075	1079	rBV2	67092	106413	0.12%	0.040%
12	9.357	1079	1083	1087	rVV	71950	148762	0.17%	0.055%
13	9.398	1087	1090	1097	rVV	107308	166881	0.19%	0.062%
14	9.481	1098	1104	1119	rVB	445943	815609	0.93%	0.303%
15	10.010	1188	1194	1211	rBV	660003	1198320	1.37%	0.446%
16	10.263	1226	1237	1251	rBV	29880381	55883568	63.87%	20.790%
17	10.386	1252	1258	1265	rVV	800631	1262773	1.44%	0.470%
18	10.763	1314	1322	1341	rBV	6558552	11083914	12.67%	4.124%
19	12.392	1591	1599	1600	rBV	160032	212334	0.24%	0.079%
20	12.427	1600	1605	1615	rVB	639819	1093330	1.25%	0.407%
21	13.045	1702	1710	1728	rBV	2427508	3831679	4.38%	1.426%
22	13.310	1749	1755	1766	rBV3	50052	121633	0.14%	0.045%
23	13.680	1812	1818	1823	rBV	1167447	1717141	1.96%	0.639%
24	13.951	1857	1864	1873	rBV	1549395	2189394	2.50%	0.815%
25	14.257	1910	1916	1926	rBV2	1048889	1590859	1.82%	0.592%
26	14.574	1965	1970	1979	rVB	237304	346264	0.40%	0.129%
27	15.257	2080	2086	2105	rBV	1926454	2737447	3.13%	1.018%
28	15.404	2105	2111	2124	rBV	405264	655420	0.75%	0.244%
29	17.015	2378	2385	2391	rBV2	1275606	1728246	1.98%	0.643%
30	17.115	2396	2402	2413	rBV	1867730	2755239	3.15%	1.025%
31	17.904	2530	2536	2545	rBV3	60704	114309	0.13%	0.043%
32	19.421	2788	2794	2805	rBV2	2333849	3114390	3.56%	1.159%
33	21.221	3095	3100	3109	rBV	1393848	1967489	2.25%	0.732%
34	21.786	3193	3196	3199	rBV2	80107	89383	0.10%	0.033%

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Title : SVOA CALIBRATION

35	22.239	3270	3273	3279	rVB	145425	175466	0.20%	0.065%
36	22.733	3353	3357	3362	rVB	181044	249965	0.29%	0.093%
37	23.291	3446	3452	3464	rBV2	1993149	3621229	4.14%	1.347%
38	23.439	3470	3477	3493	rVB	1057260	2106058	2.41%	0.784%
39	23.915	3554	3558	3565	rVB	169278	296807	0.34%	0.110%
40	24.644	3679	3682	3691	rVB	128647	240733	0.28%	0.090%

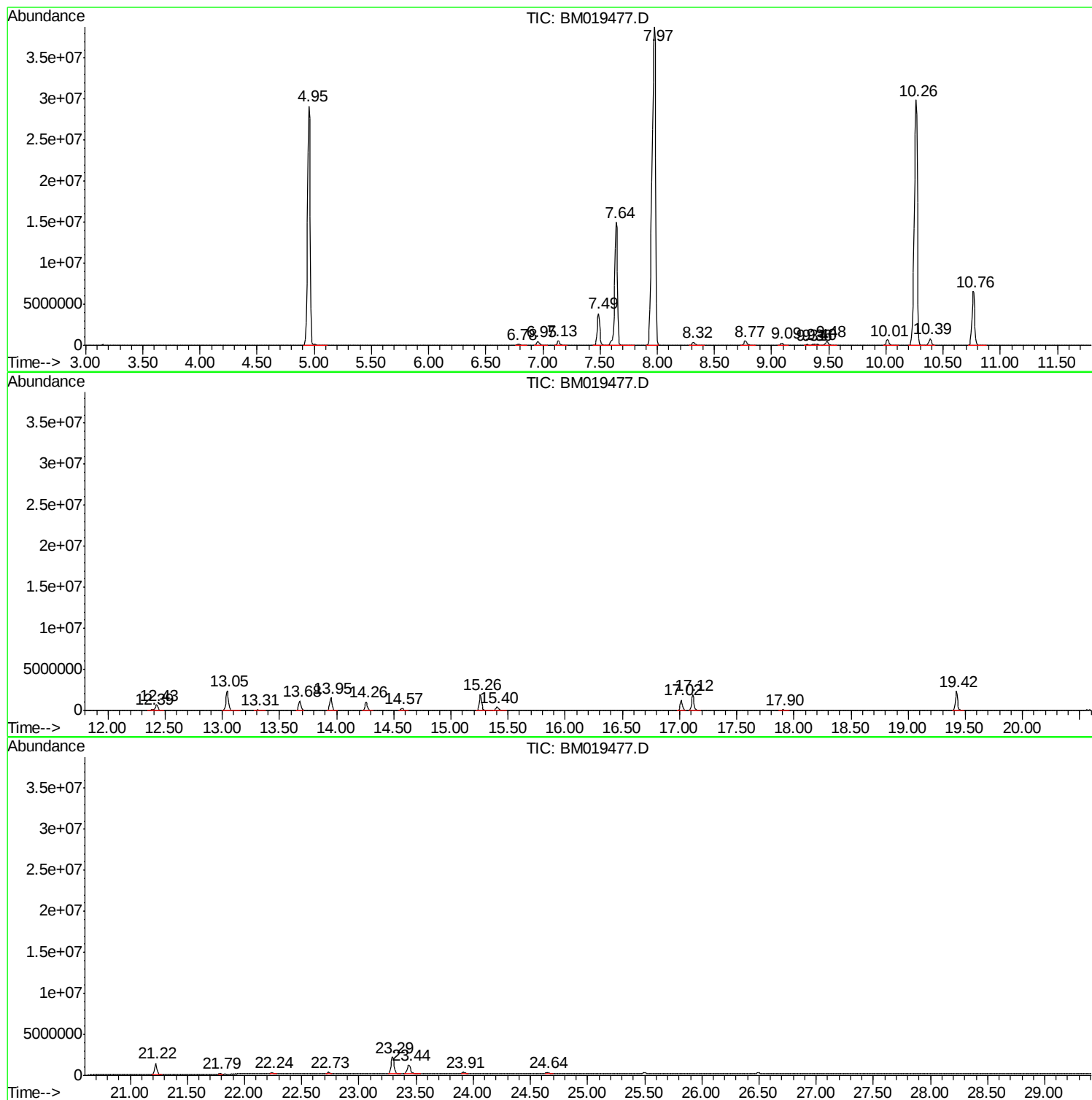
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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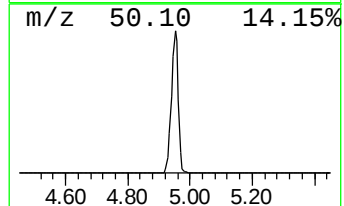
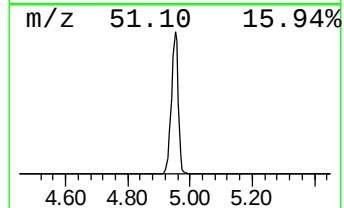
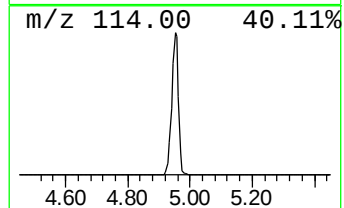
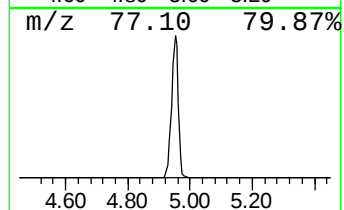
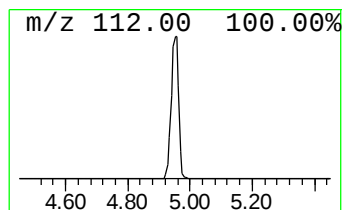
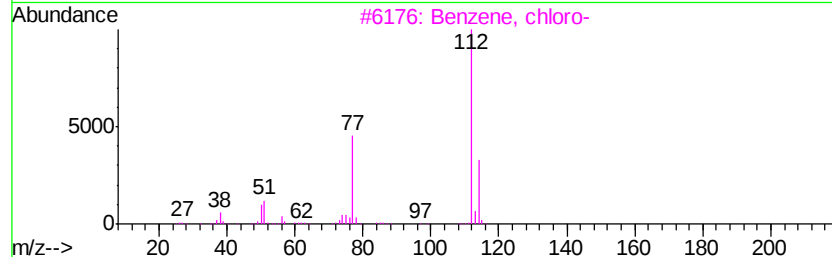
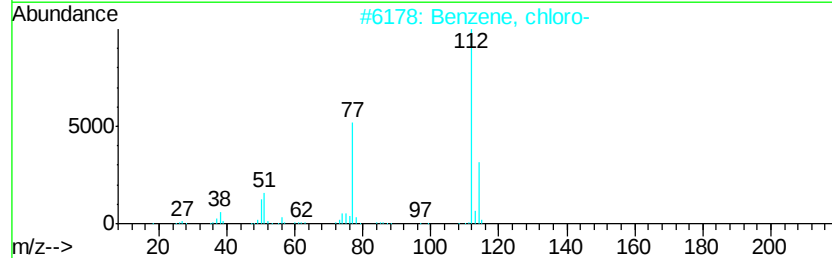
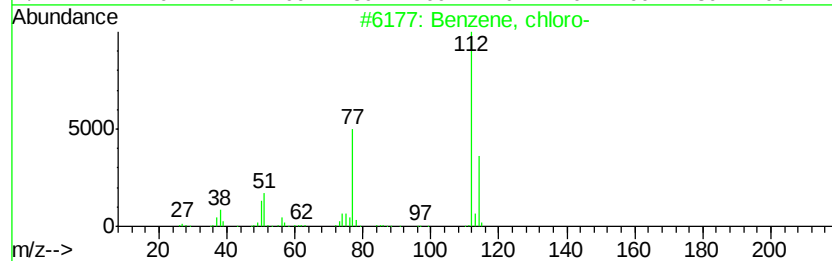
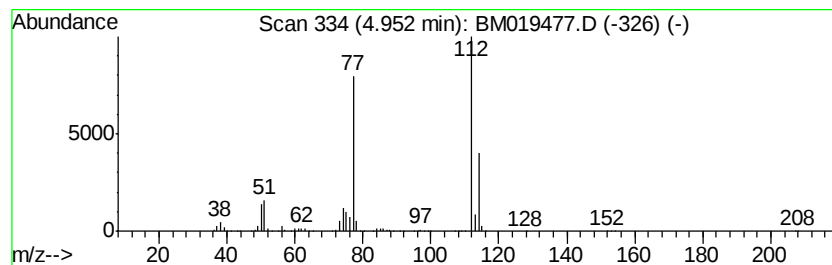
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Benzene, chloro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.95	36.74 ng/ul	45239200	1,4-Dichlorobenzene-d4	7.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, chloro-	112	C6H5Cl	000108-90-7	95
2		Benzene, chloro-	112	C6H5Cl	000108-90-7	95
3		Benzene, chloro-	112	C6H5Cl	000108-90-7	94
4		Benzene, chloro-	112	C6H5Cl	000108-90-7	94
5		Phenol, 3-fluoro-	112	C6H5FO	000372-20-3	22



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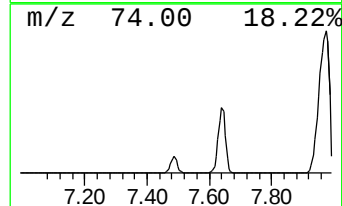
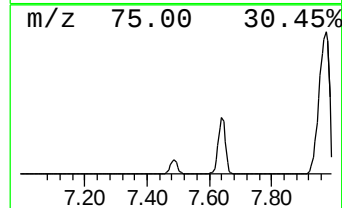
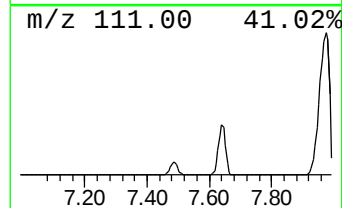
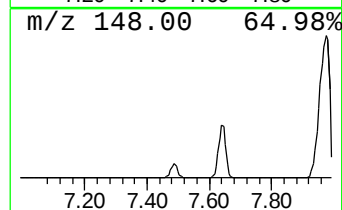
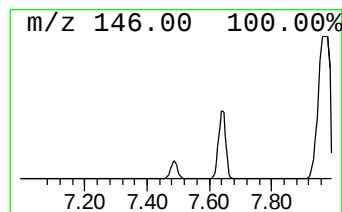
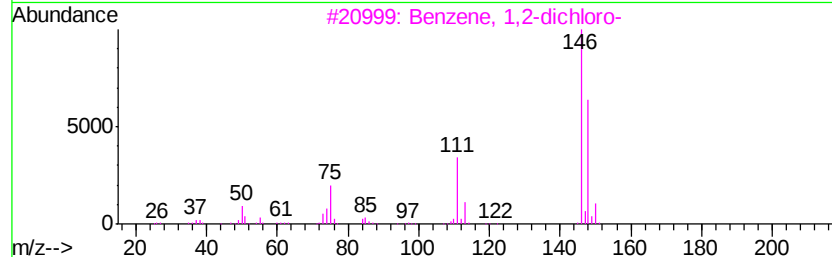
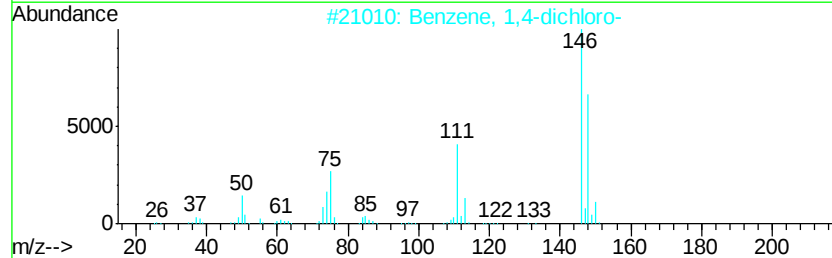
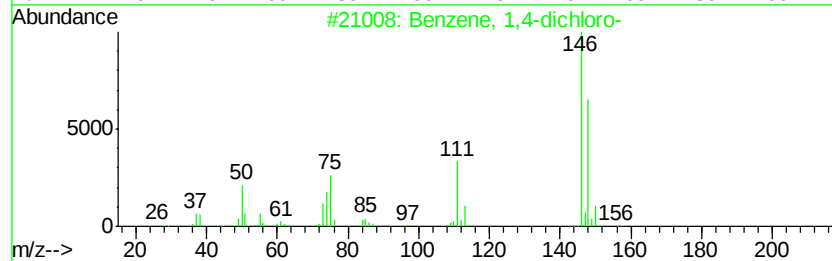
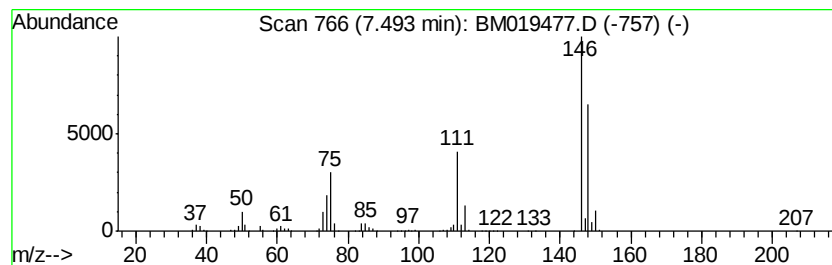
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Benzene, 1,4-dichloro- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.49	4.77 ng/ul	5877270	1,4-Dichlorobenzene-d4	7.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96
2		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96
3		Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	95
4		Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	95
5		Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	95



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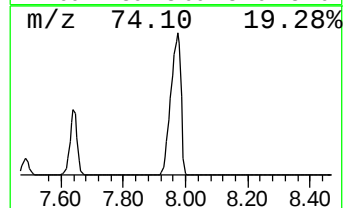
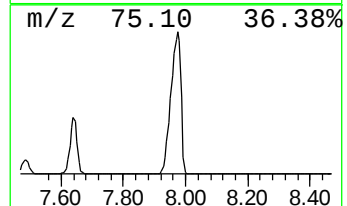
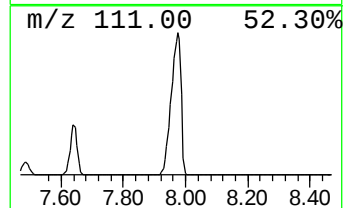
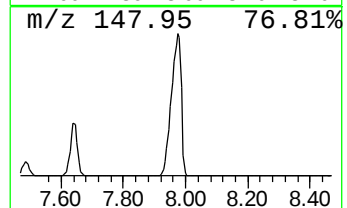
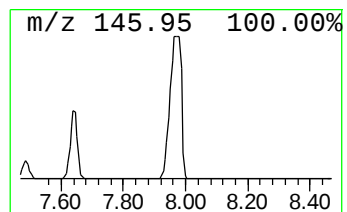
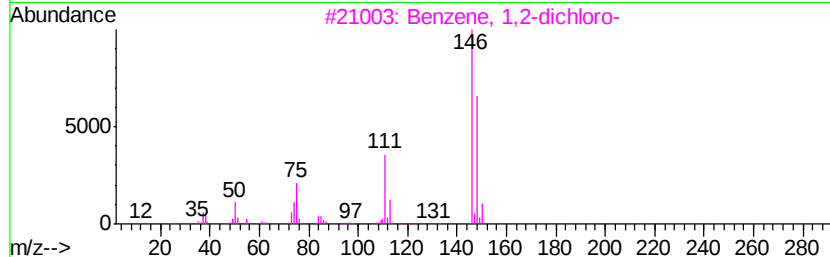
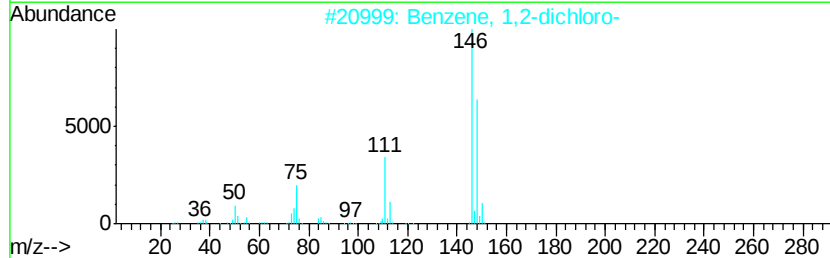
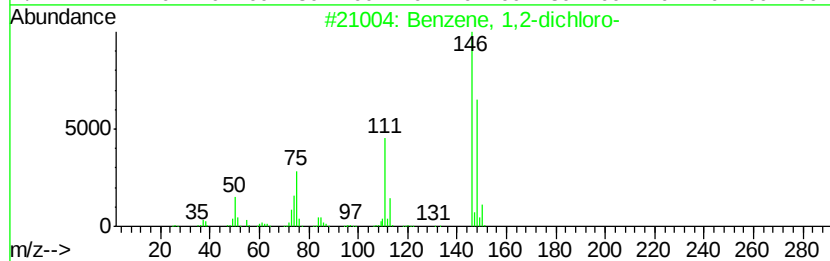
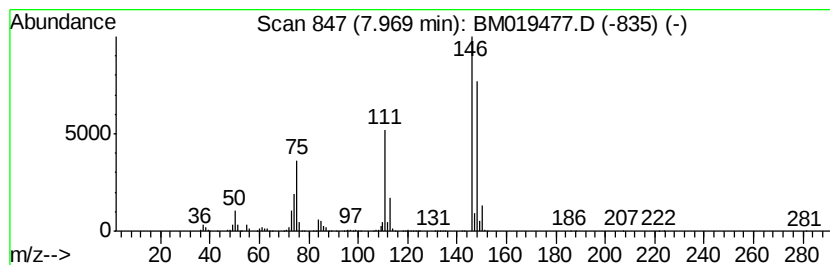
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
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Peak Number 3 Benzene, 1,2-dichloro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.97	71.07 ng/ul	87500900	1,4-Dichlorobenzene-d4	7.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	95
2		Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	94
3		Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	94
4		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	94
5		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	94



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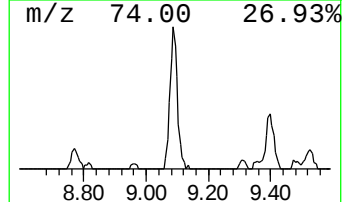
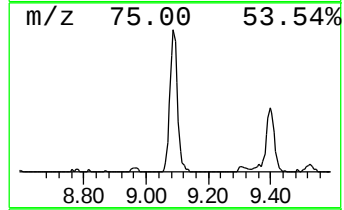
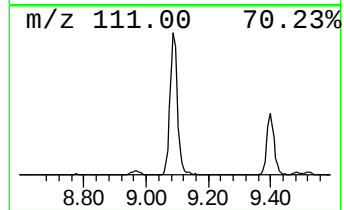
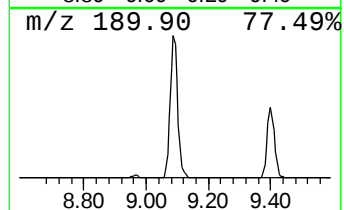
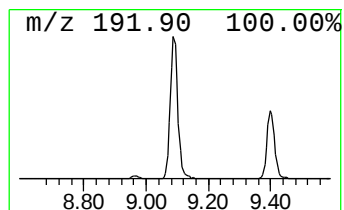
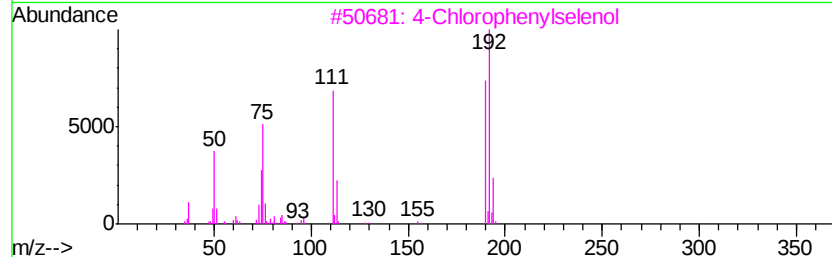
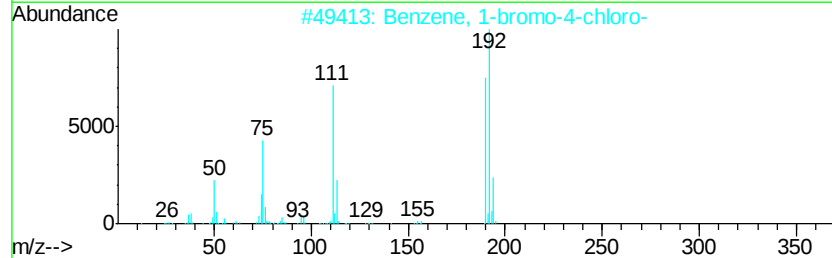
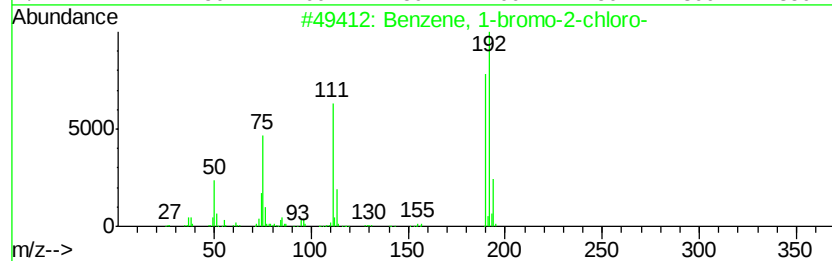
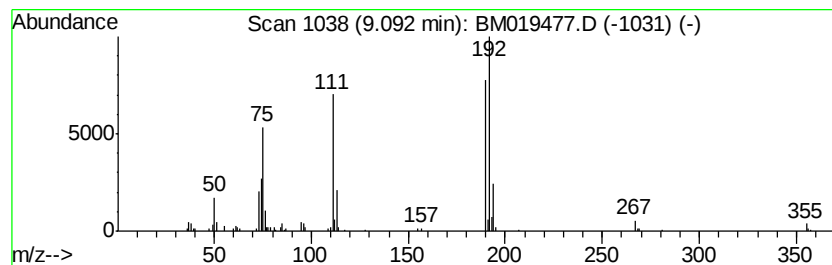
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Peak Number 4 Benzene, 1-bromo-2-chloro- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.09	6.49 ng/ul	410013	Napthalene-d8	10.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	98
2		Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	98
3		4-Chlorophenylselenol	192	C6H5ClSe	016645-10-6	97
4		Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	96
5		Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	95



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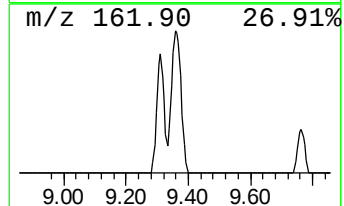
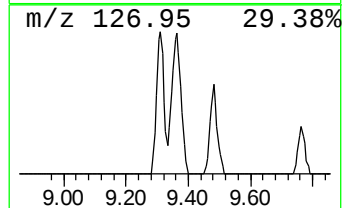
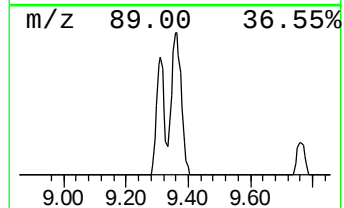
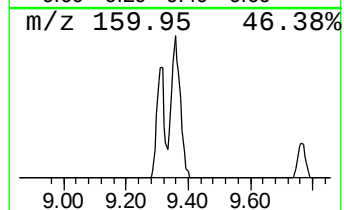
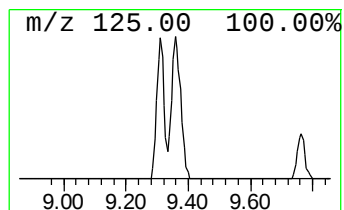
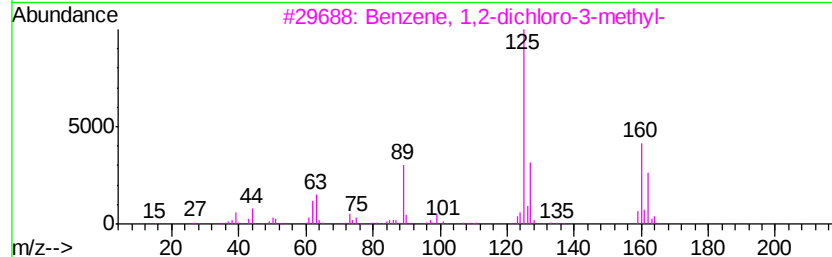
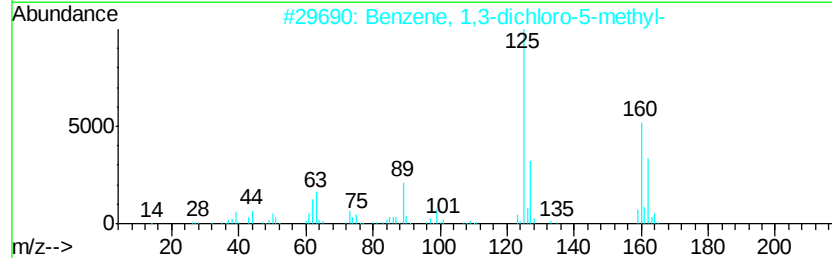
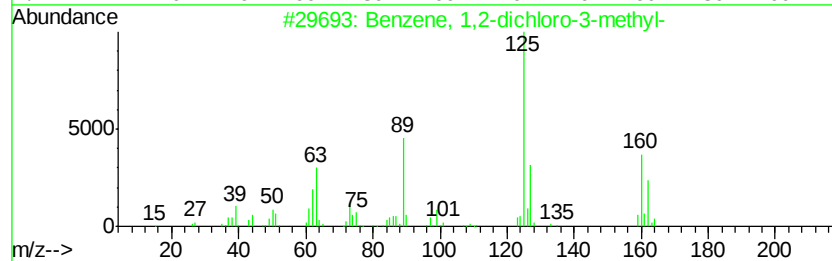
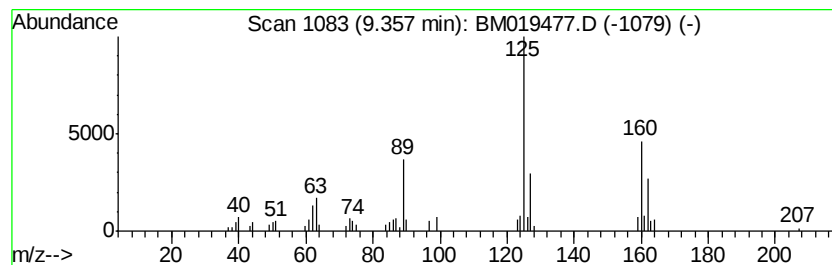
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, 1,2-dichloro-3-met... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.36	2.36 ng/ul	148762	Naphthalene-d8	10.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2-dichloro-3-methyl-	160	C7H6Cl2	032768-54-0	97
2		Benzene, 1,3-dichloro-5-methyl-	160	C7H6Cl2	025186-47-4	96
3		Benzene, 1,2-dichloro-3-methyl-	160	C7H6Cl2	032768-54-0	96
4		Benzene, 1,4-dichloro-2-methyl-	160	C7H6Cl2	019398-61-9	96
5		Benzene, 1,3-dichloro-2-methyl-	160	C7H6Cl2	000118-69-4	95



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM032619\
Data File : BM019477.D
Acq On : 26 Mar 2019 19:15
Operator : JU/SJ
Sample : K2069-11
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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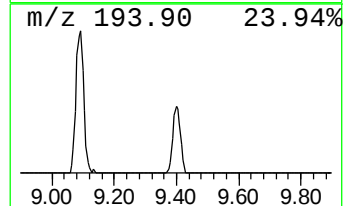
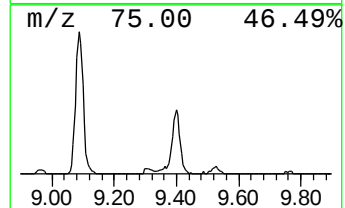
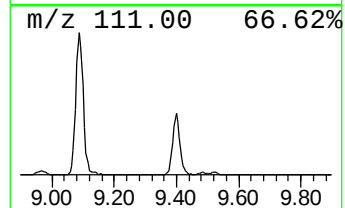
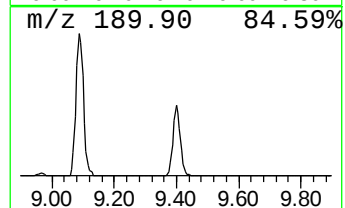
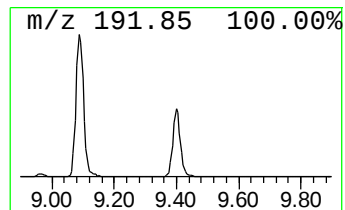
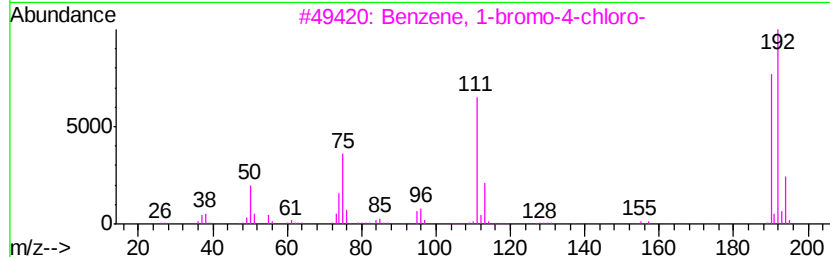
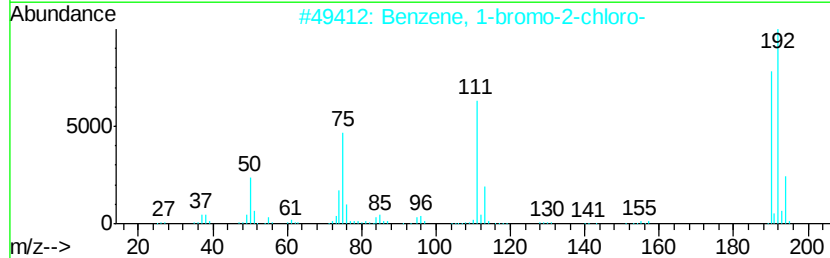
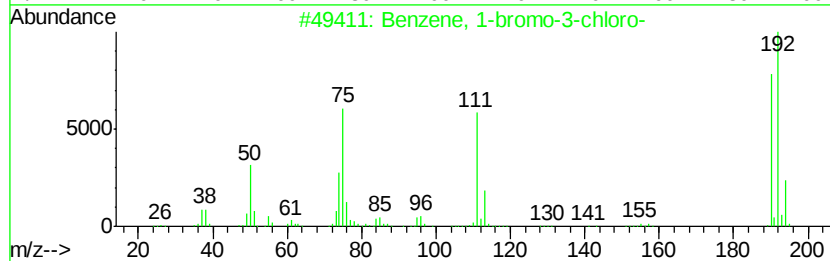
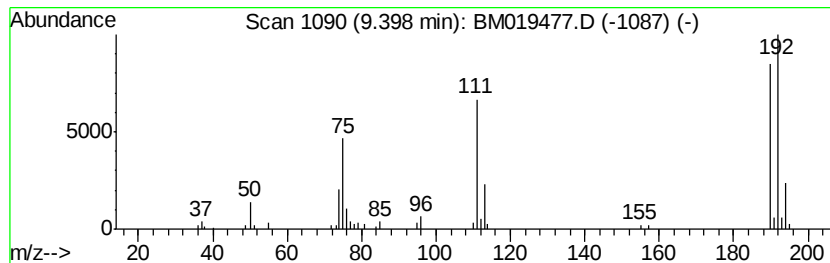
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzene, 1-bromo-3-chloro- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.40	2.64 ng/ul	166881	Napthalene-d8	10.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	97
2		Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	95
3		Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	94
4		Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	94
5		Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM032619\
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Operator : JU/SJ
Sample : K2069-11
Misc :
ALS Vial : 9 Sample Multiplier: 1

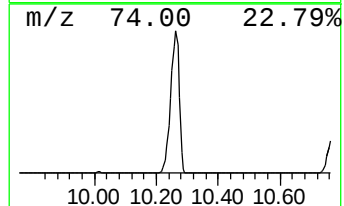
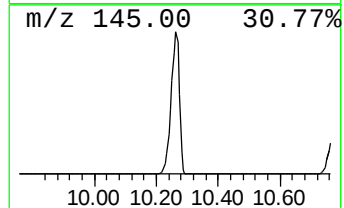
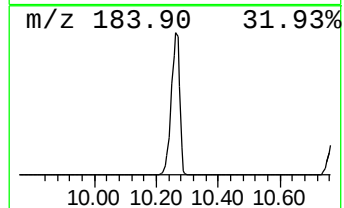
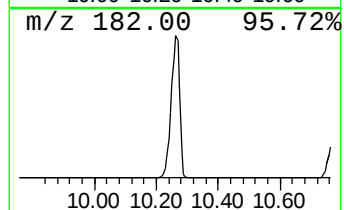
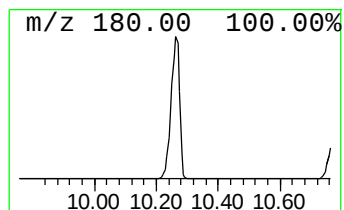
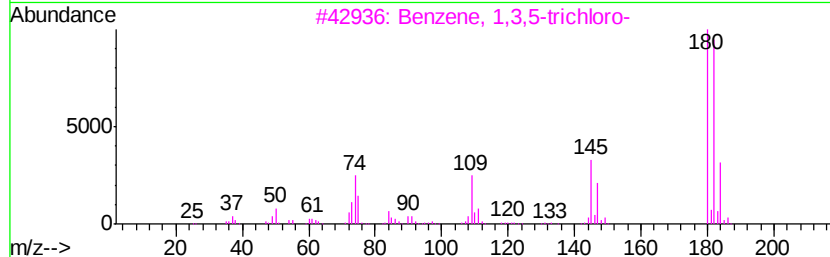
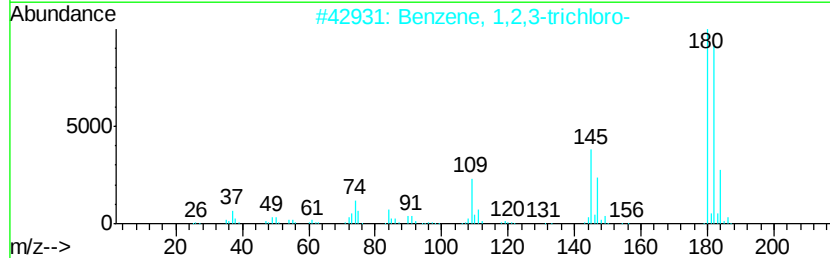
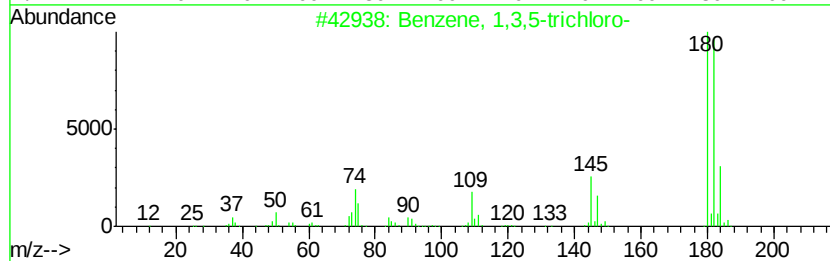
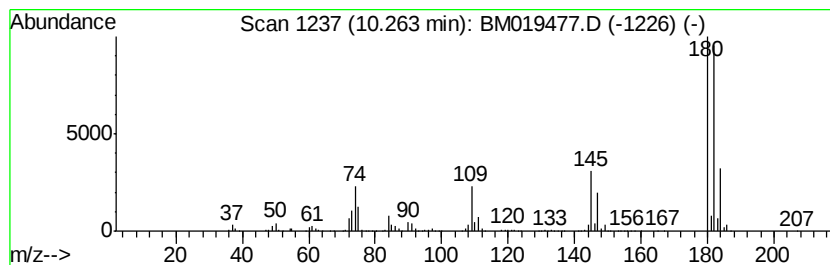
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Benzene, 1,3,5-trichloro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.26	885.09 ng/ul	55883600	Napthalene-d8	10.39
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1,3,5-trichloro-	180 C6H3Cl3	000108-70-3 98
2		Benzene, 1,2,3-trichloro-	180 C6H3Cl3	000087-61-6 98
3		Benzene, 1,3,5-trichloro-	180 C6H3Cl3	000108-70-3 96
4		Benzene, 1,2,4-trichloro-	180 C6H3Cl3	000120-82-1 95
5		Benzene, 1,3,5-trichloro-	180 C6H3Cl3	000108-70-3 95



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM032619\
Data File : BM019477.D
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Sample : K2069-11
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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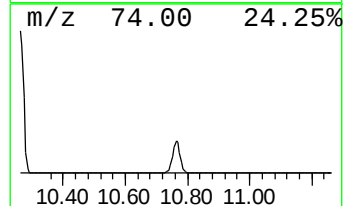
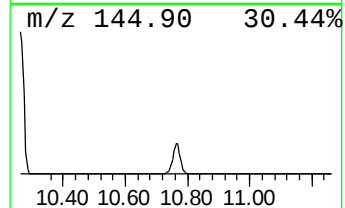
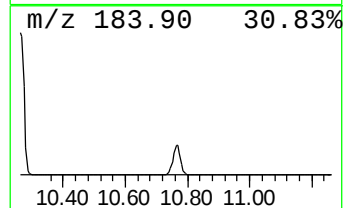
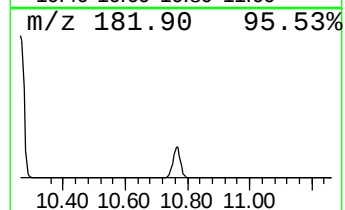
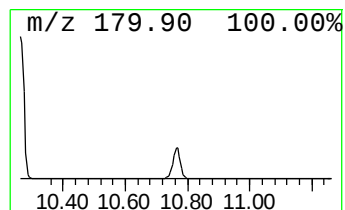
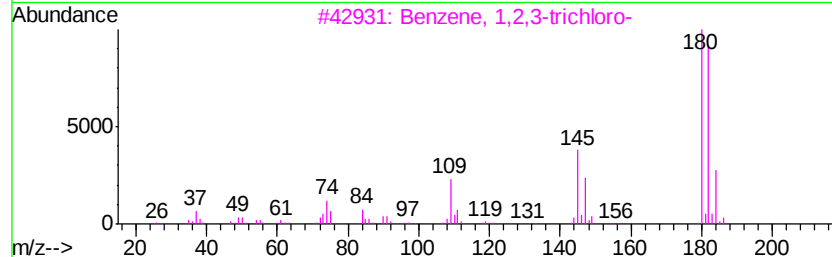
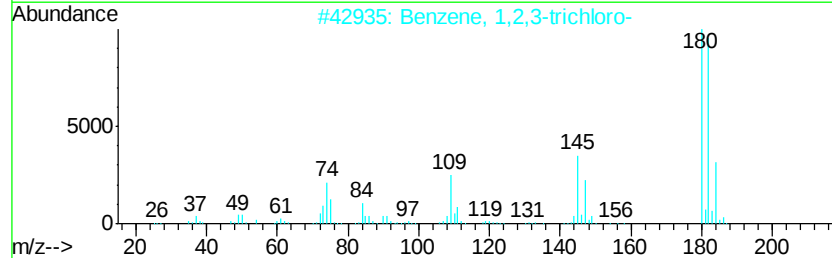
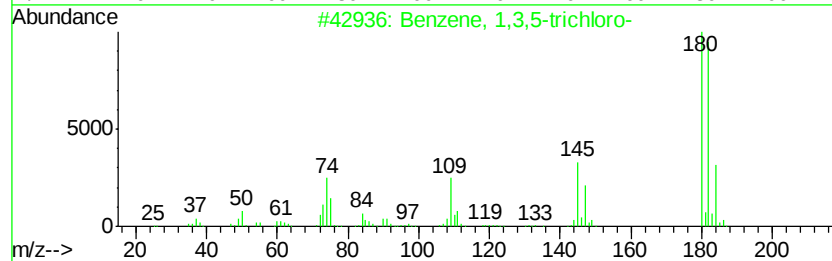
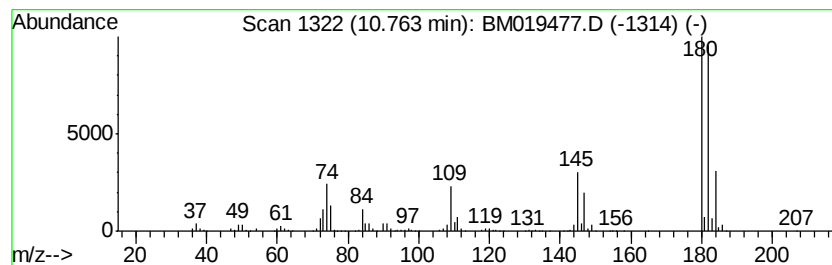
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Benzene, 1,2,3-trichloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.76	175.55 ng/ul	11083900	Napthalene-d8	10.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	98
2		Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	98
3		Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	98
4		Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	98
5		Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	97



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM032619\
Data File : BM019477.D
Acq On : 26 Mar 2019 19:15
Operator : JU/SJ
Sample : K2069-11
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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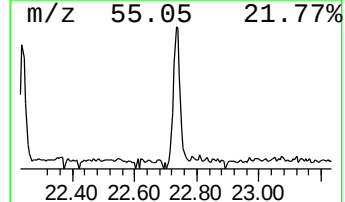
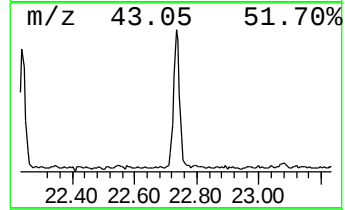
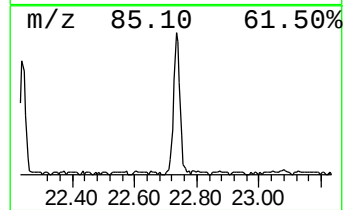
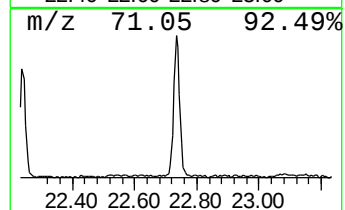
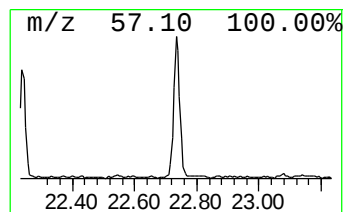
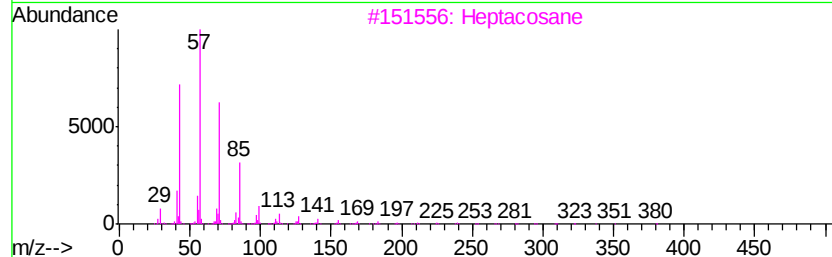
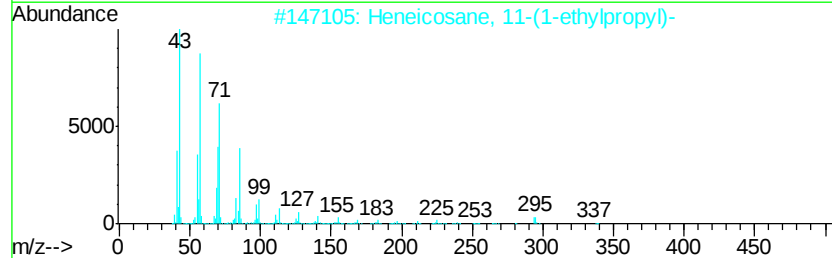
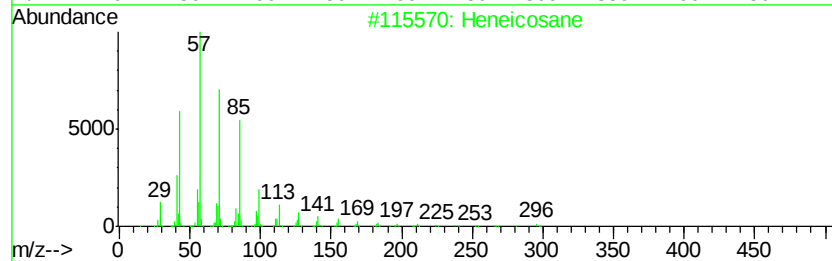
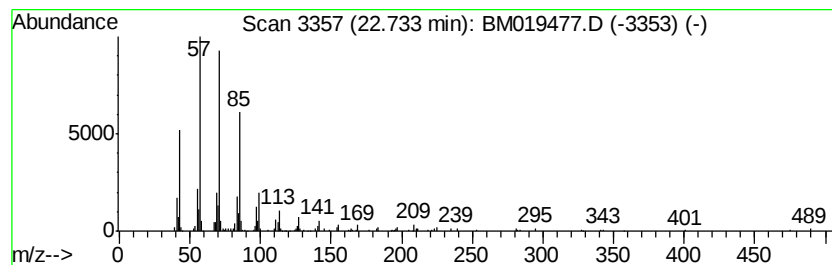
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.73	2.37 ng/ul	249965	Perylene-d12	23.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91
3		Heptacosane	380	C27H56	000593-49-7	91
4		Octacosane	394	C28H58	000630-02-4	87
5		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM032619\
Data File : BM019477.D
Acq On : 26 Mar 2019 19:15
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Sample : K2069-11
Misc :
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Instrument :
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ClientSampleId :
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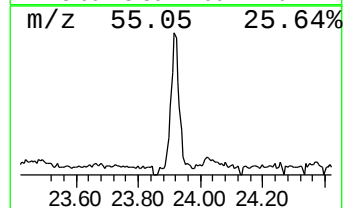
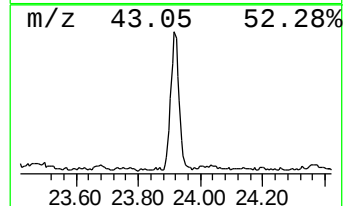
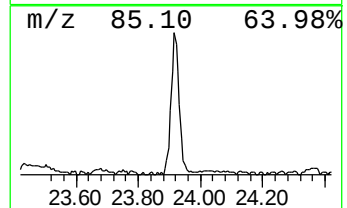
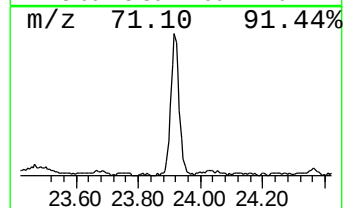
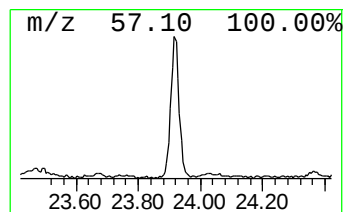
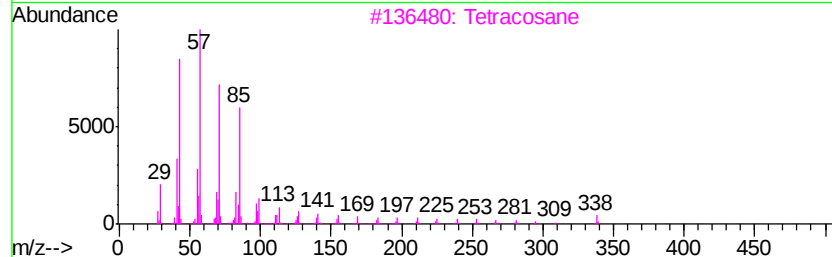
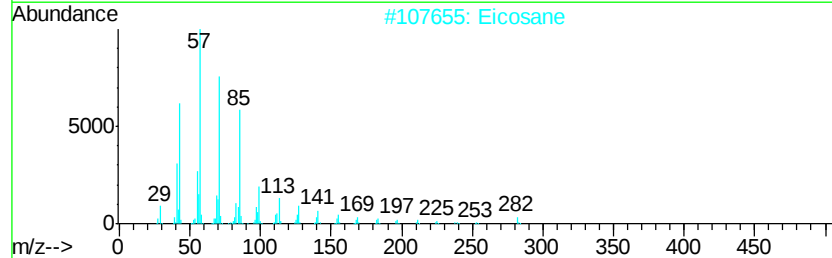
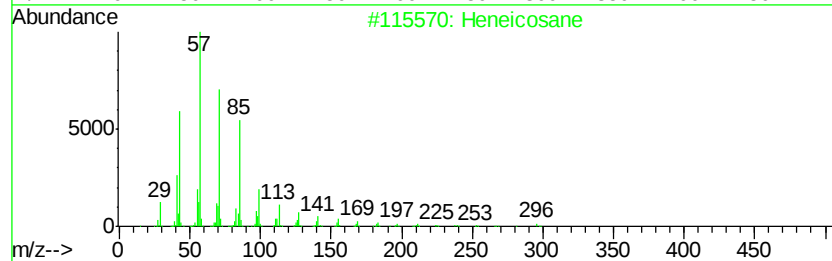
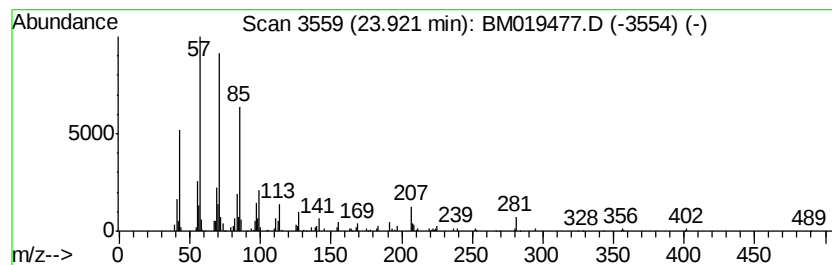
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.92	2.82 ng/ul	296807	Perylene-d12	23.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	96
2		Eicosane	282	C20H42	000112-95-8	91
3		Tetracosane	338	C24H50	000646-31-1	87
4		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	87
5		Octadecane	254	C18H38	000593-45-3	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM032619\
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Sample : K2069-11
Misc :
ALS Vial : 9 Sample Multiplier: 1

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ClientSampleId :
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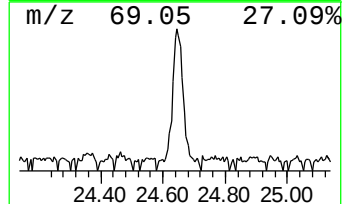
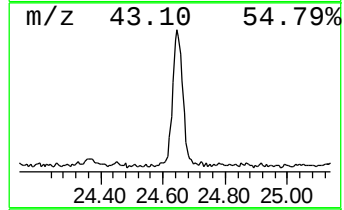
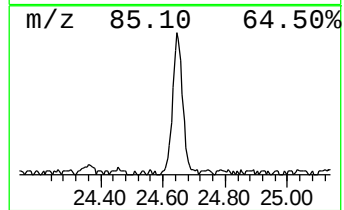
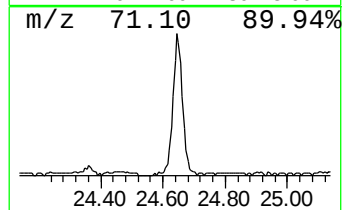
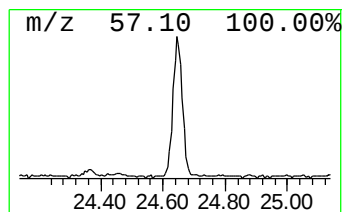
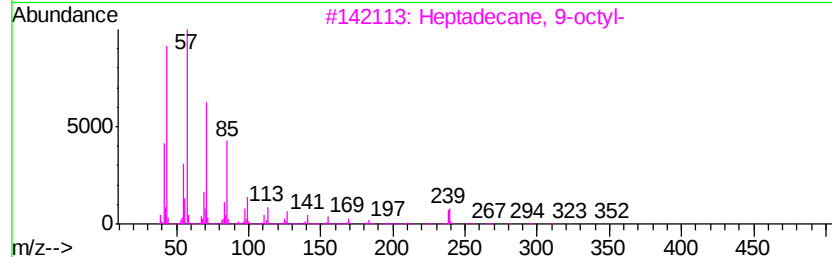
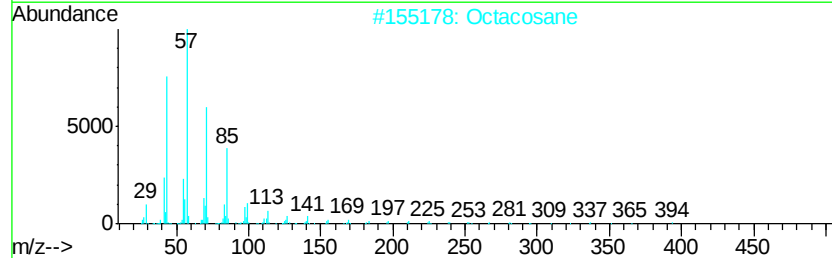
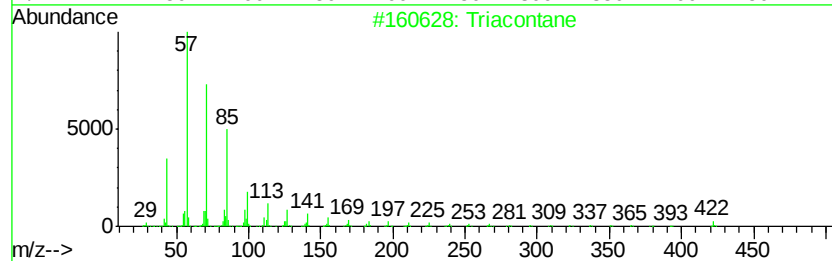
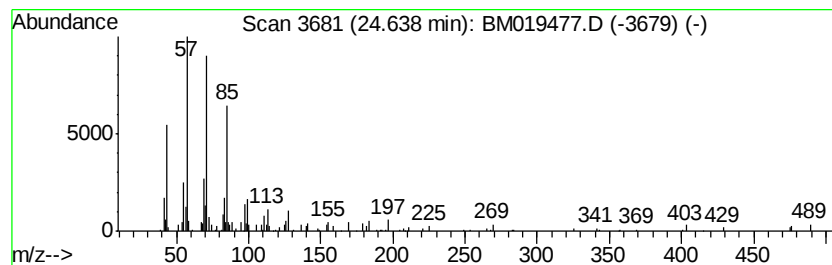
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.64	2.29 ng/ul	240733	Perylene-d12	23.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Triacontane	422	C30H62	000638-68-6	87
2		Octacosane	394	C28H58	000630-02-4	86
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	86
4		Hentriacontane	437	C31H64	000630-04-6	86
5		Hexacosane	366	C26H54	000630-01-3	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM032619\
 Data File : BM019477.D
 Acq On : 26 Mar 2019 19:15
 Operator : JU/SJ
 Sample : K2069-11
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
 Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Benzene, 1,4-dich...	7.49	4.8	ng/ul	5877270	1	7.60	24623700	20.0
Benzene, 1,2-dich...	7.97	71.1	ng/ul	87500900	1	7.60	24623700	20.0
Benzene, 1-bromo-...	9.09	6.5	ng/ul	410013	2	10.39	1262770	20.0
Benzene, 1,2-dich...	9.36	2.4	ng/ul	148762	2	10.39	1262770	20.0
Benzene, 1-bromo-...	9.40	2.6	ng/ul	166881	2	10.39	1262770	20.0
Benzene, 1,3,5-tr...	10.26	885.1	ng/ul	55883600	2	10.39	1262770	20.0
Benzene, 1,2,3-tr...	10.76	175.6	ng/ul	11083900	2	10.39	1262770	20.0
(DEL) Alkane: Str...	22.73	2.4	ng/ul	249965	6	23.44	2106060	20.0
(DEL) Alkane: Str...	23.92	2.8	ng/ul	296807	6	23.44	2106060	20.0
(DEL) Alkane: Str...	24.64	2.3	ng/ul	240733	6	23.44	2106060	20.0