

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM032619\
 Data File : BM019478.D
 Acq On : 26 Mar 2019 19:52
 Operator : JU/SJ
 Sample : K2069-12
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C09D9

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	5.069	326	354	358	RBV5	53759446	414036182	100.00%	26.918%
2	6.299	557	563	574	RBV	353721	589321	0.14%	0.038%
3	6.987	669	680	688	RBV	350723	977378	0.24%	0.064%
4	7.146	700	707	720	RVB	476866	1064680	0.26%	0.069%
5	7.510	757	769	784	RBV	21565055	87040232	21.02%	5.659%
6	7.751	784	810	814	RVV6	51548954	346171803	83.61%	22.506%
7	8.104	840	870	873	RBV7	51981498	405996523	98.06%	26.395%
8	8.328	902	908	920	RVB	458926	818033	0.20%	0.053%
9	8.781	978	985	988	RBV	689451	1188760	0.29%	0.077%
10	8.834	988	994	1010	RVB	6403080	10965913	2.65%	0.713%
11	9.098	1032	1039	1049	RBV	846744	1342561	0.32%	0.087%
12	9.404	1087	1091	1100	RVB	400176	685959	0.17%	0.045%
13	9.492	1100	1106	1110	RBV	606995	1089294	0.26%	0.071%
14	10.022	1190	1196	1207	RBV	852426	1555653	0.38%	0.101%
15	10.304	1227	1244	1248	RBV2	48321657	151708445	36.64%	9.863%
16	10.404	1255	1261	1266	RBV	941691	1361347	0.33%	0.089%
17	10.786	1311	1326	1332	RBV	26019373	50037488	12.09%	3.253%
18	11.010	1354	1364	1379	RBV	6528927	10942062	2.64%	0.711%
19	11.222	1391	1400	1414	RBV	1070225	2114619	0.51%	0.137%
20	12.398	1593	1600	1601	RBV	421520	637696	0.15%	0.041%
21	12.428	1601	1605	1614	RVB	1256585	1996795	0.48%	0.130%
22	13.045	1698	1710	1722	RBV	4553754	6659579	1.61%	0.433%
23	13.680	1810	1818	1825	RBV	1689305	2648458	0.64%	0.172%
24	13.951	1858	1864	1872	RVB	1972868	2836663	0.69%	0.184%
25	14.257	1905	1916	1923	RBV2	1153938	1737142	0.42%	0.113%
26	15.257	2077	2086	2094	RBV	2398330	3402903	0.82%	0.221%
27	15.398	2105	2110	2121	RVB	702660	1034674	0.25%	0.067%
28	17.015	2379	2385	2395	RBV2	1324643	1880780	0.45%	0.122%
29	17.115	2395	2402	2414	RBV	2530121	3531876	0.85%	0.230%
30	18.398	2615	2620	2629	RVV	6832951	9110438	2.20%	0.592%
31	18.521	2636	2641	2652	RVB	278687	417526	0.10%	0.027%
32	18.733	2673	2677	2680	RVV	319281	463854	0.11%	0.030%
33	19.421	2788	2794	2804	RBV	2892097	3887916	0.94%	0.253%
34	21.221	3095	3100	3107	RBV	1561155	2058604	0.50%	0.134%

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35	23.297	3446	3453	3464	rVB2	2220972	4051291	0.98%	0.263%
36	23.433	3470	3476	3493	rVB	1062165	2114646	0.51%	0.137%

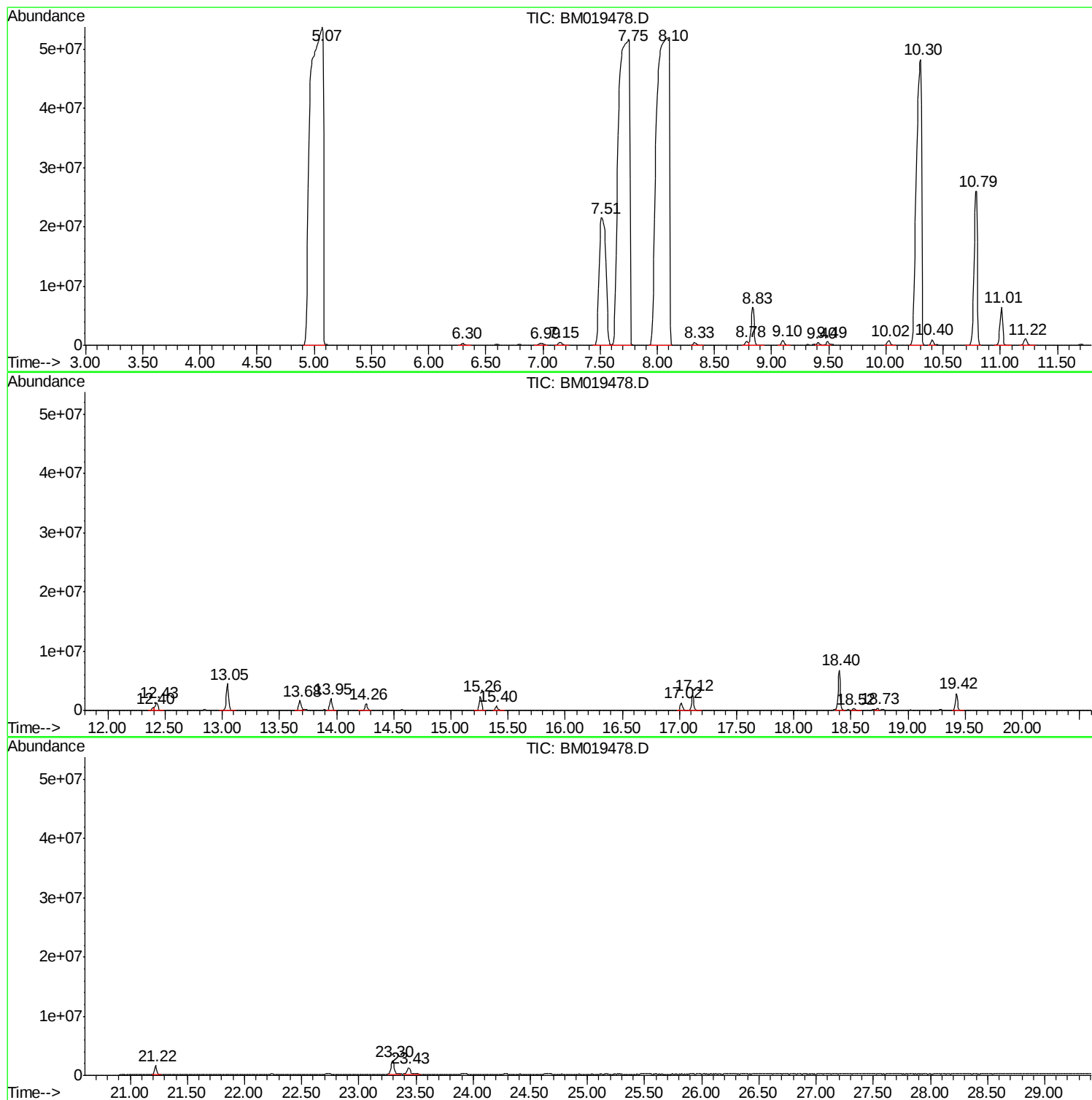
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Quant Title : SVOA CALIBRATION

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



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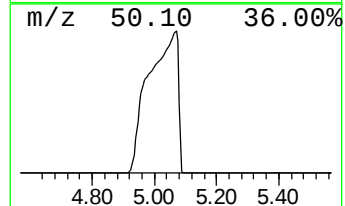
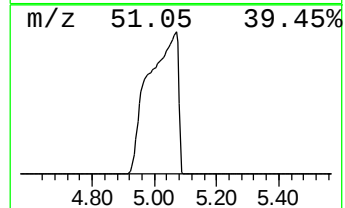
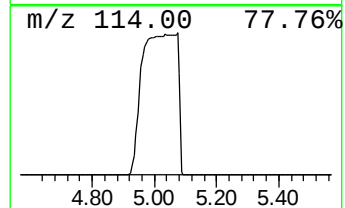
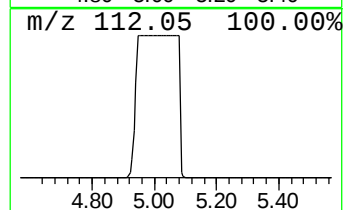
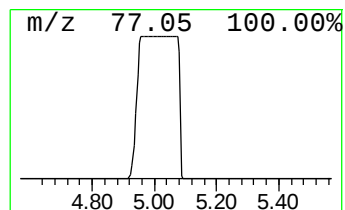
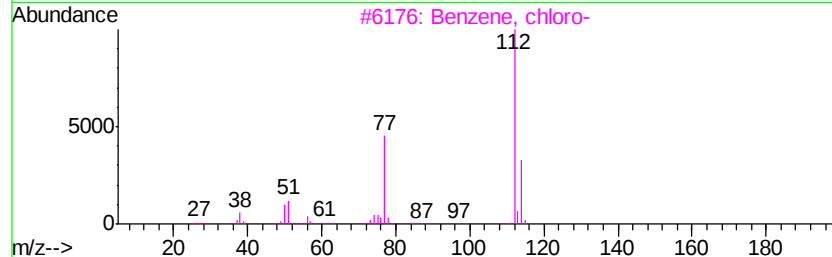
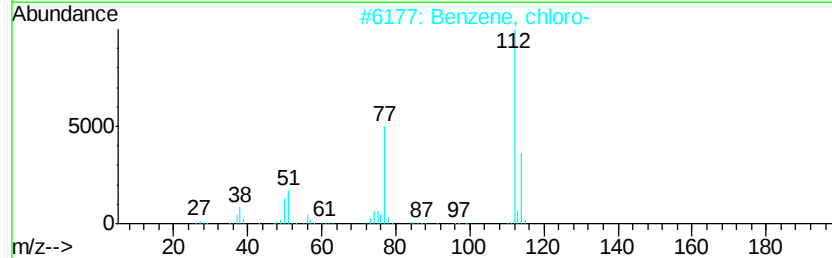
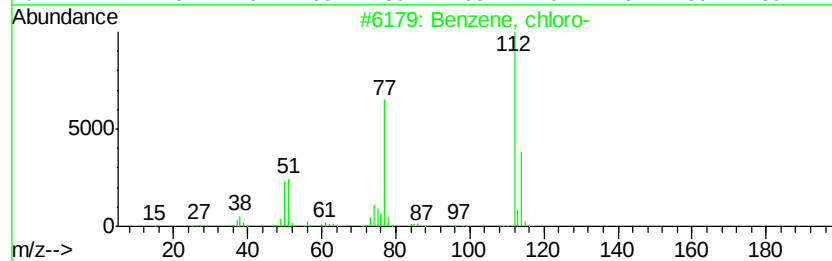
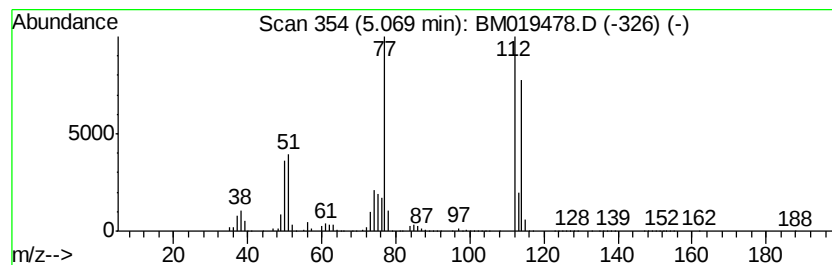
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Benzene, chloro- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.07	23.92 ng/ul	414036000	1,4-Dichlorobenzene-d4	7.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, chloro-	112	C6H5Cl	000108-90-7	76
2		Benzene, chloro-	112	C6H5Cl	000108-90-7	53
3		Benzene, chloro-	112	C6H5Cl	000108-90-7	42
4		Benzene, chloro-	112	C6H5Cl	000108-90-7	42
5		Diazobenzene, 4-nitro-	151	C6H5N3O2	022719-27-3	22



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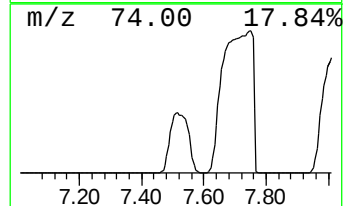
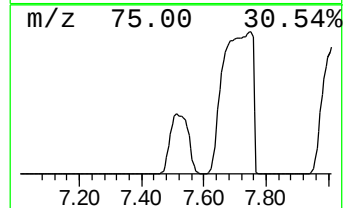
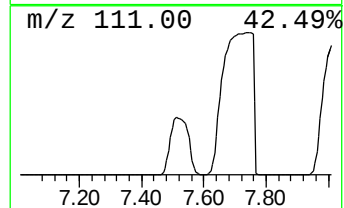
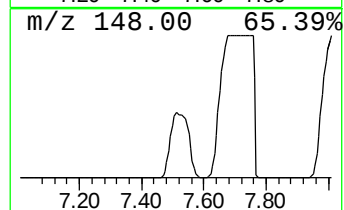
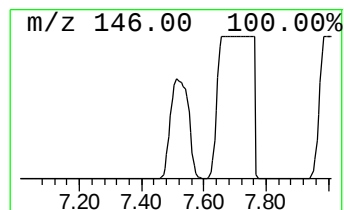
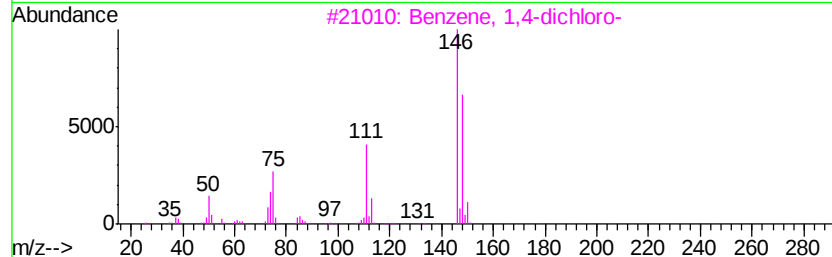
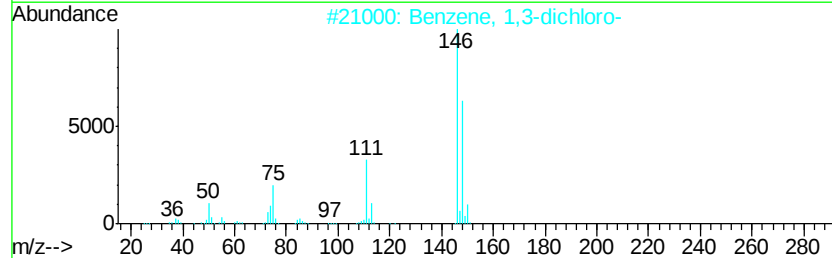
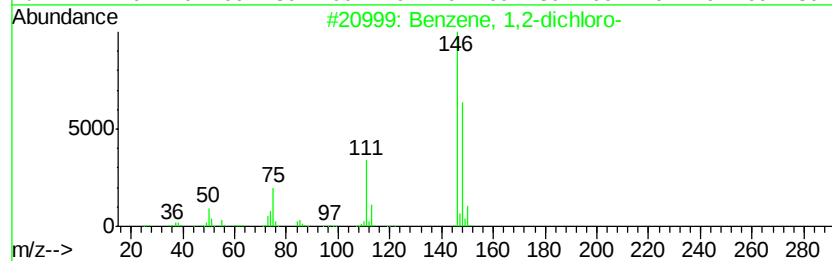
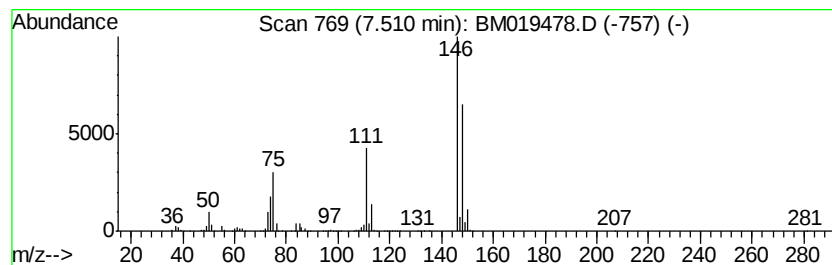
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzene, 1,2-dichloro- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.51	5.03 ng/ul	87040200	1,4-Dichlorobenzene-d4	7.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	97
2		Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	97
3		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96
4		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	95
5		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	95



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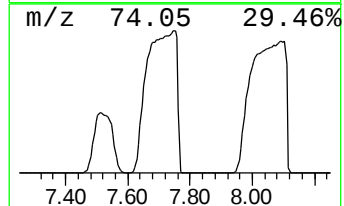
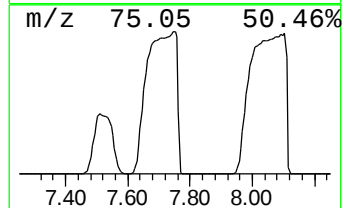
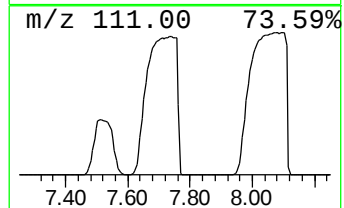
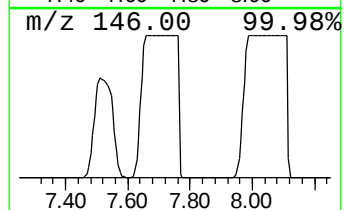
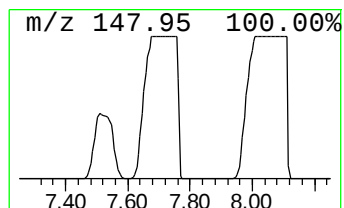
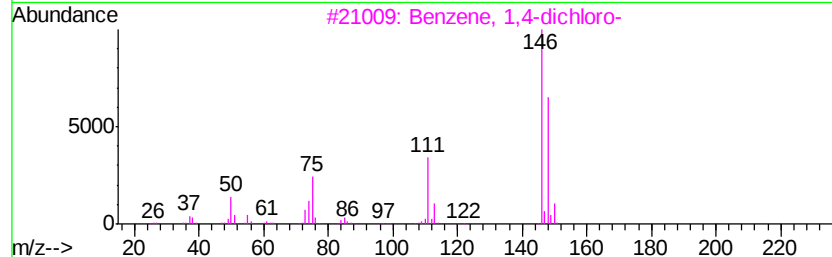
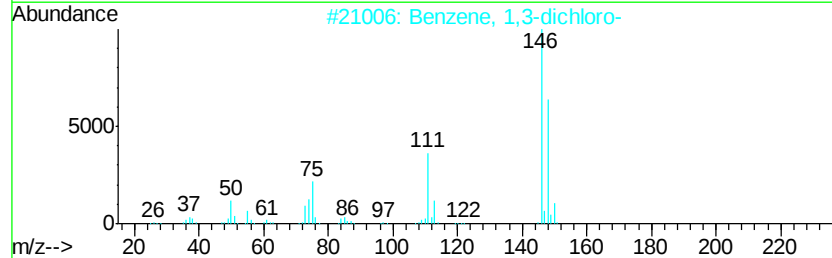
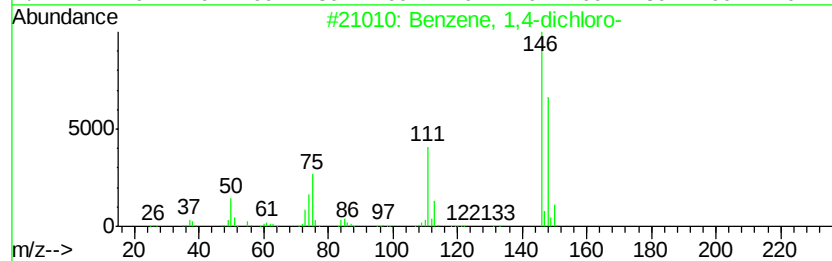
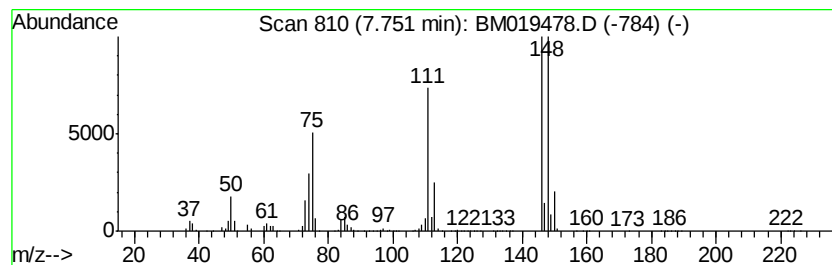
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 3 Benzene, 1,4-dichloro- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.75	20.00 ng/ul	346172000	1,4-Dichlorobenzene-d4	7.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	91
2		Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	87
3		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	87
4		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	81
5		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	81



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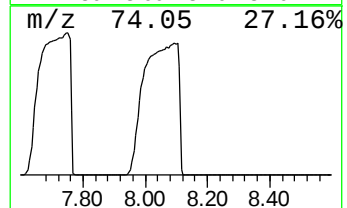
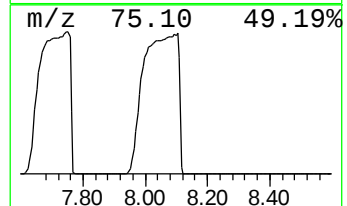
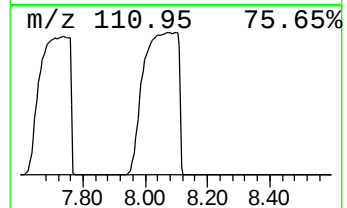
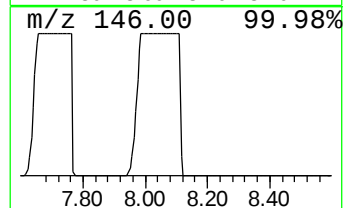
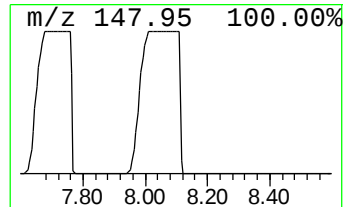
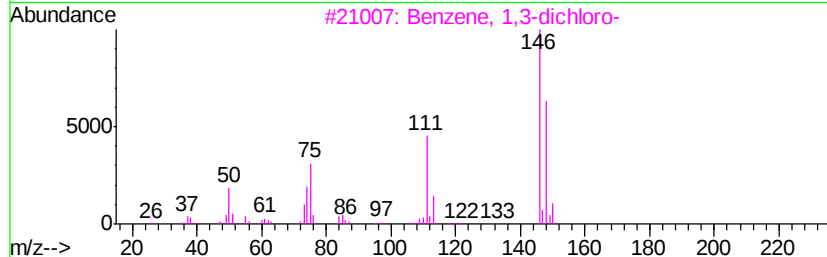
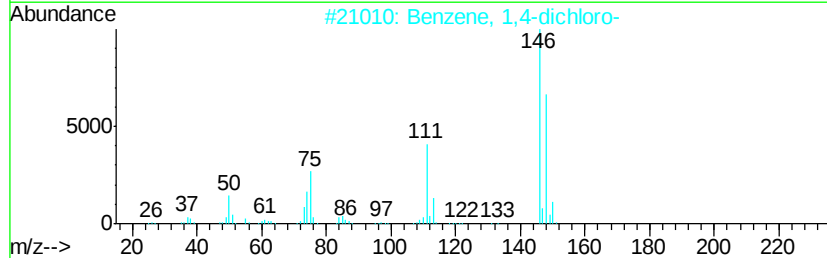
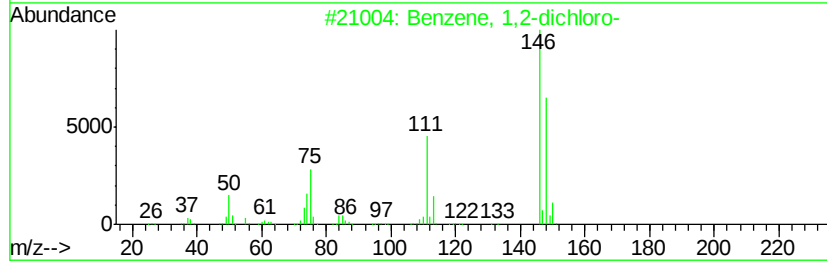
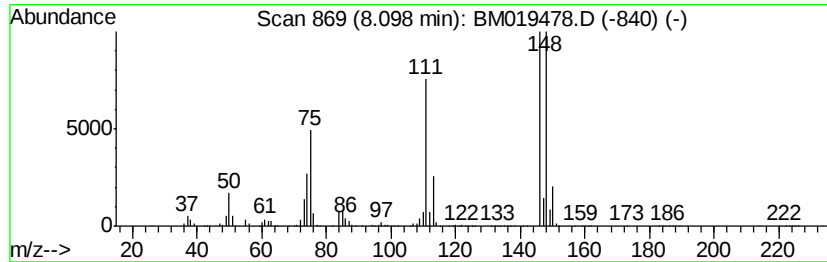
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Peak Number 4 Benzene, 1,3-dichloro- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.10	23.46 ng/ul	405997000	1,4-Dichlorobenzene-d4	7.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	93
2		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	91
3		Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	90
4		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	87
5		Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	81



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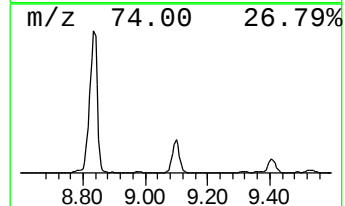
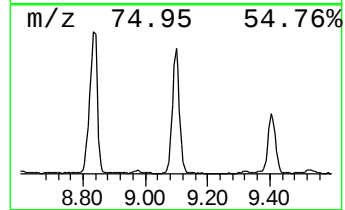
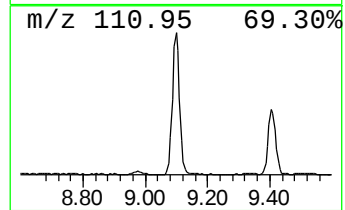
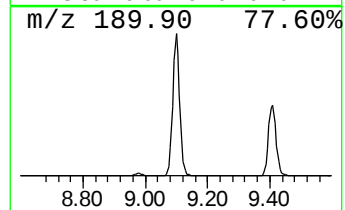
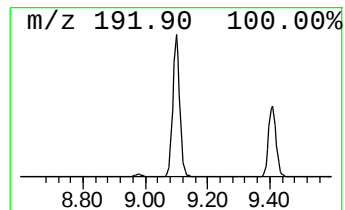
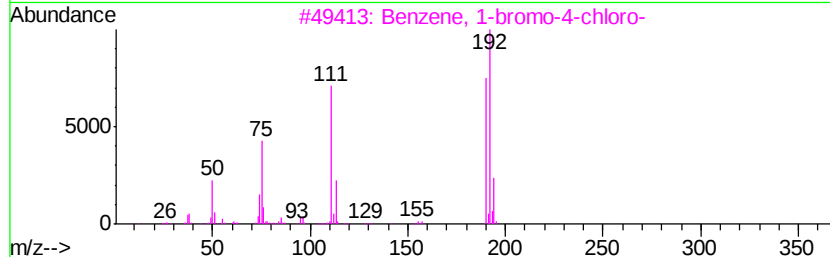
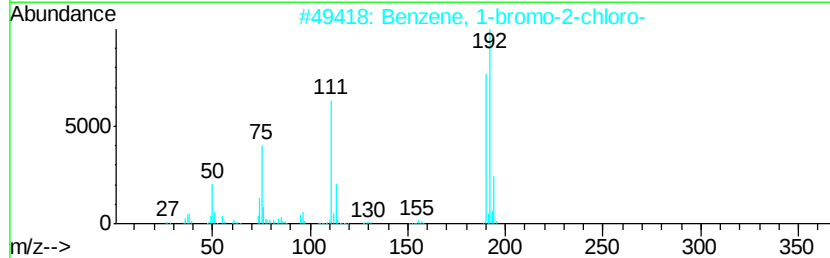
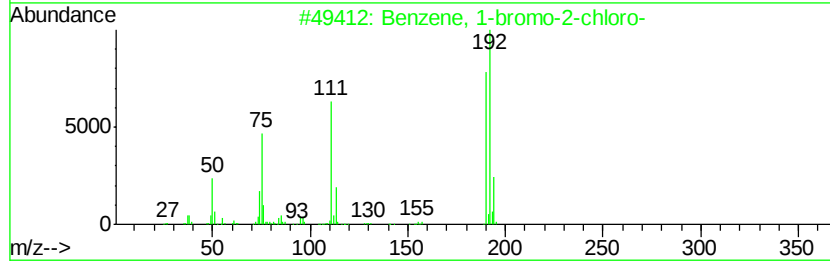
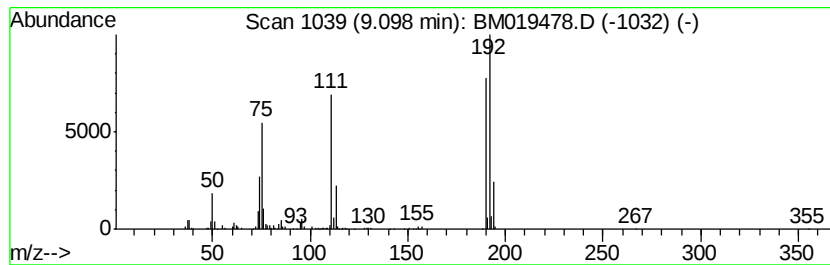
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Peak Number 5 Benzene, 1-bromo-2-chloro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.10	19.72 ng/ul	1342560	Napthalene-d8	10.40

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-2-chloro-	190	C6H4BrCl	000694-80-4	98
3		Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	97
4		Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	97
5		4-Chlorophenylselenol	192	C6H5ClSe	016645-10-6	95



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM032619\
Data File : BM019478.D
Acq On : 26 Mar 2019 19:52
Operator : JU/SJ
Sample : K2069-12
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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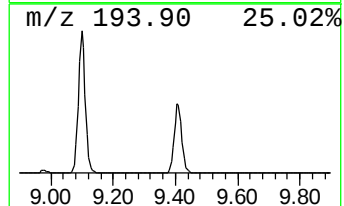
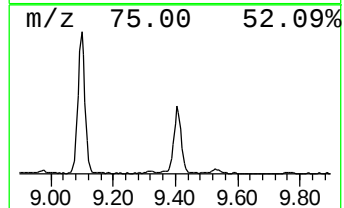
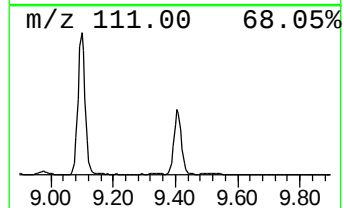
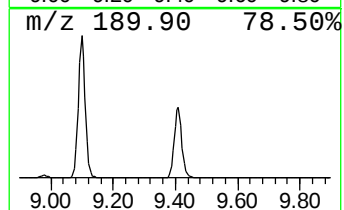
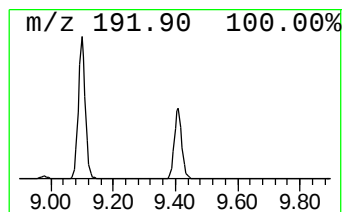
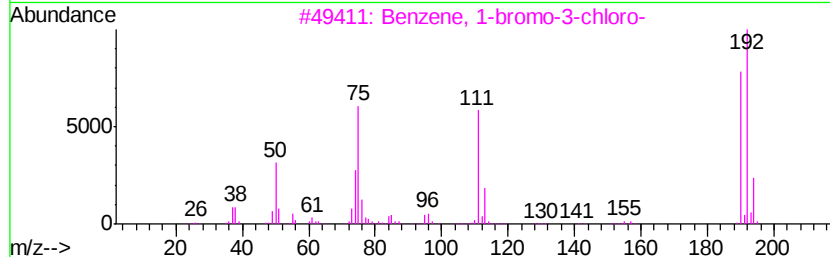
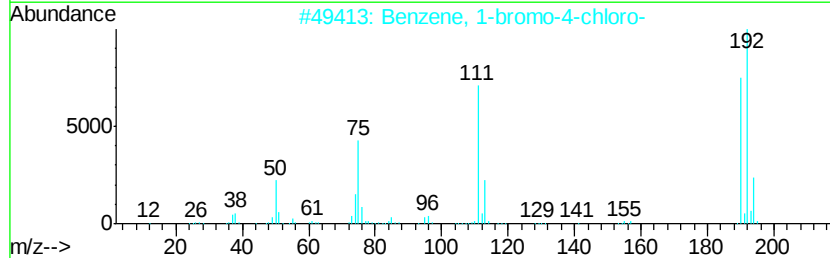
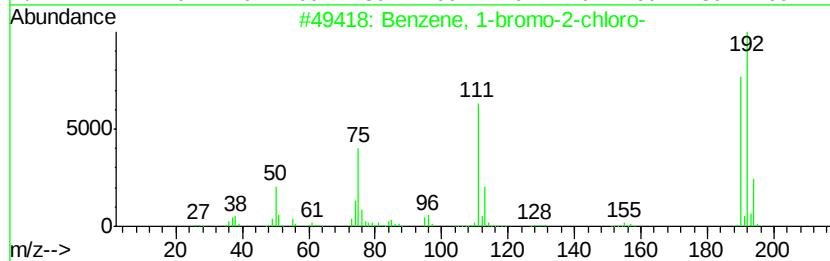
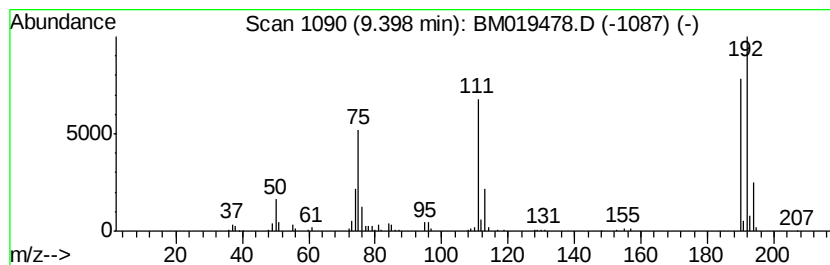
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzene, 1-bromo-4-chloro- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.40	10.08 ng/ul	685959	Napthalene-d8	10.40

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	97
3		Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	97
4		Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	95
5		Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	95



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM032619\
Data File : BM019478.D
Acq On : 26 Mar 2019 19:52
Operator : JU/SJ
Sample : K2069-12
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C09D9

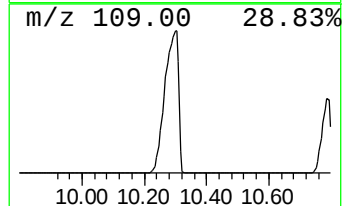
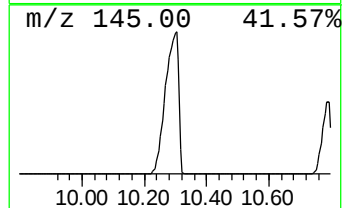
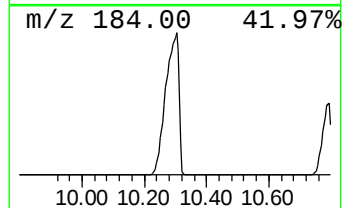
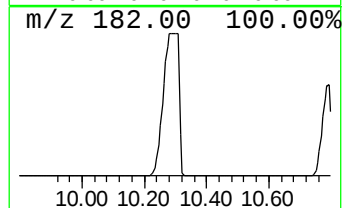
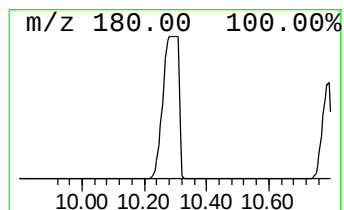
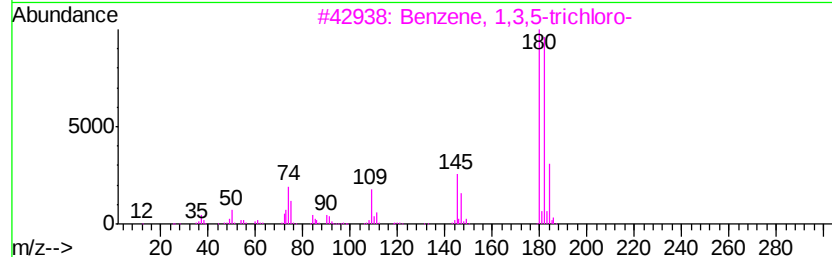
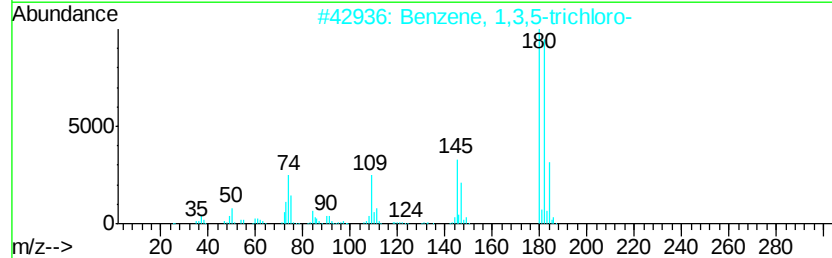
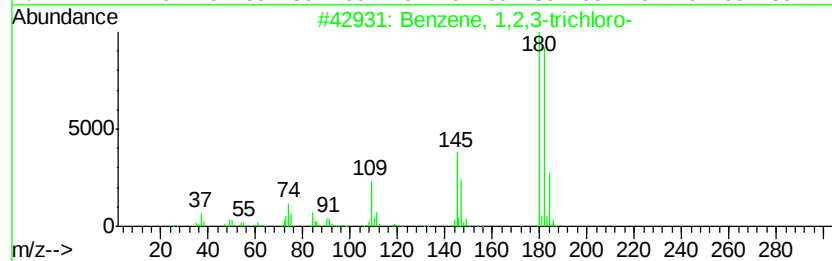
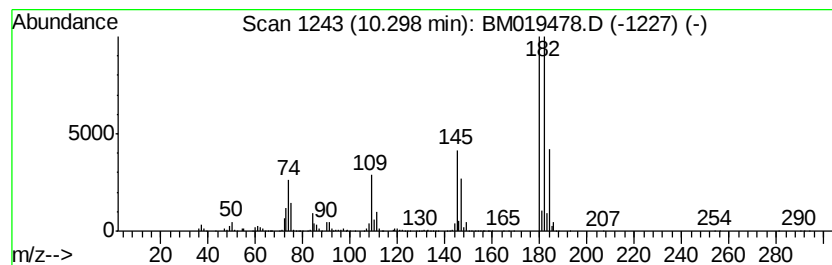
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Benzene, 1,2,3-trichloro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.30	2228.80 ng/ul	151708000	Napthalene-d8	10.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	96
2		Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	96
3		Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	95
4		Benzene, 1,2,4-trichloro-	180	C6H3Cl3	000120-82-1	95
5		Benzene, 1,2,4-trichloro-	180	C6H3Cl3	000120-82-1	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM032619\
Data File : BM019478.D
Acq On : 26 Mar 2019 19:52
Operator : JU/SJ
Sample : K2069-12
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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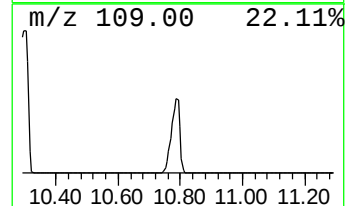
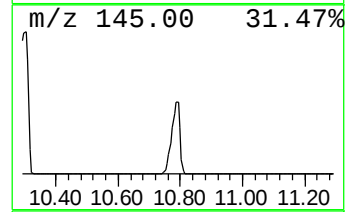
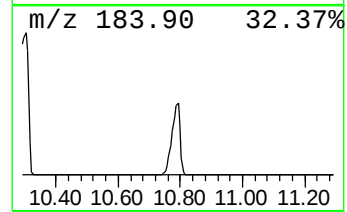
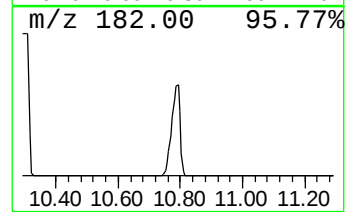
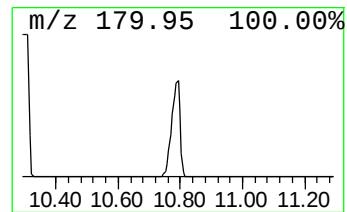
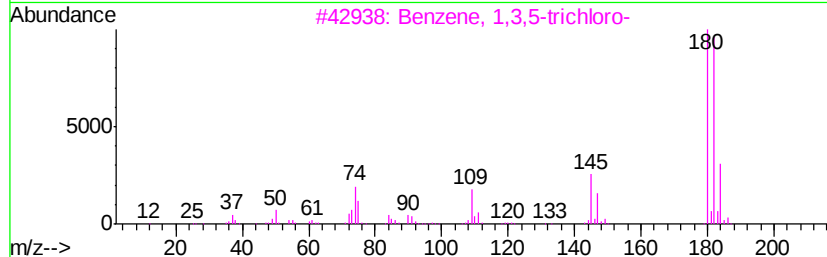
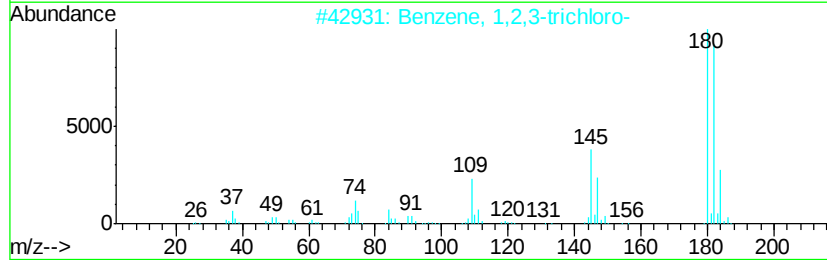
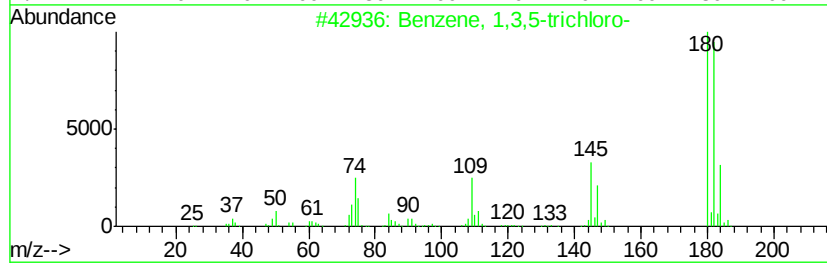
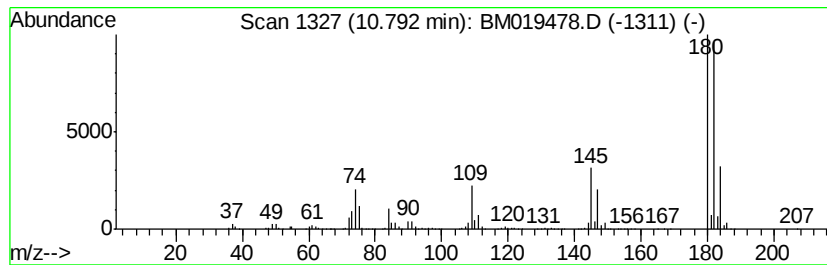
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Benzene, 1,3,5-trichloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.79	735.12 ng/ul	50037500	Naphthalene-d8	10.40

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, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	98
4		Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	97
5		Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	96



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM032619\
Data File : BM019478.D
Acq On : 26 Mar 2019 19:52
Operator : JU/SJ
Sample : K2069-12
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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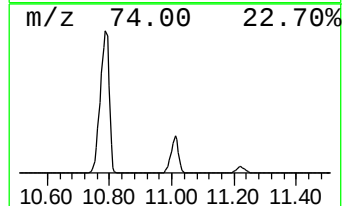
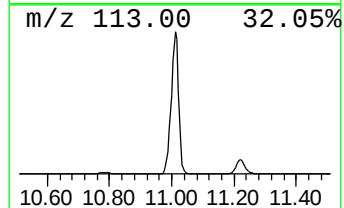
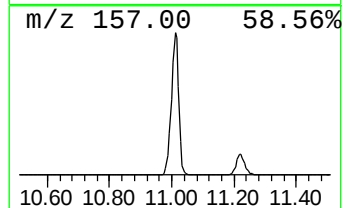
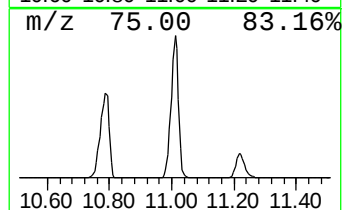
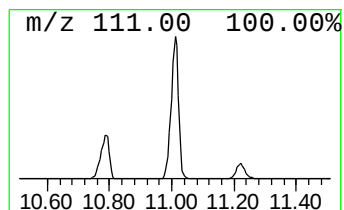
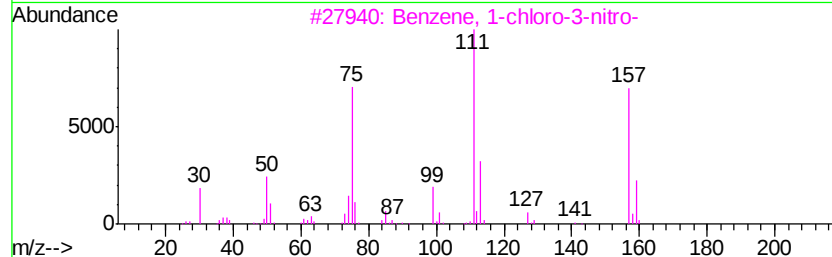
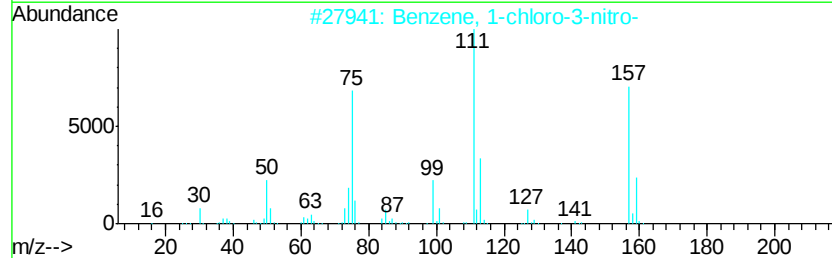
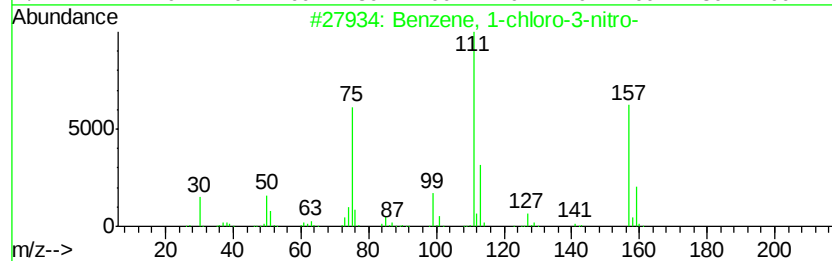
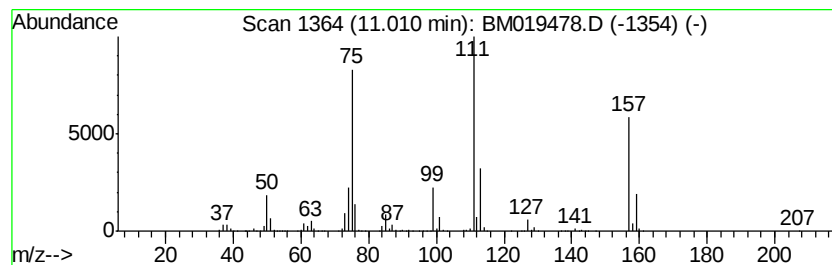
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Benzene, 1-chloro-3-nitro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.01	160.75 ng/ul	10942100	Naphthalene-d8	10.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	96
2		Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	96
3		Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	95
4		Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	94
5		Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM032619\
Data File : BM019478.D
Acq On : 26 Mar 2019 19:52
Operator : JU/SJ
Sample : K2069-12
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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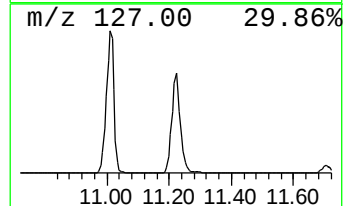
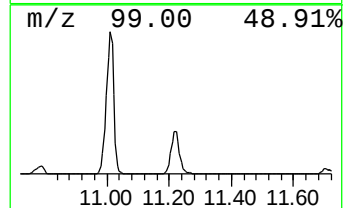
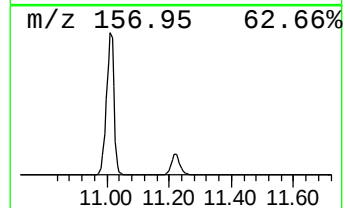
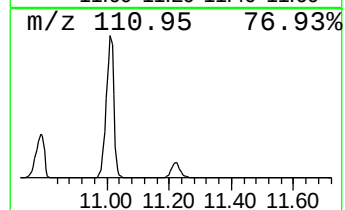
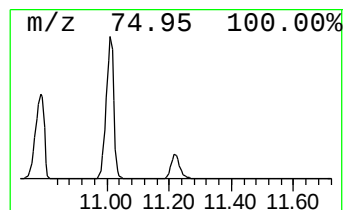
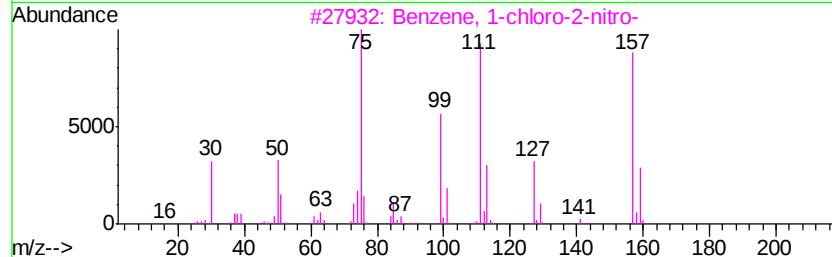
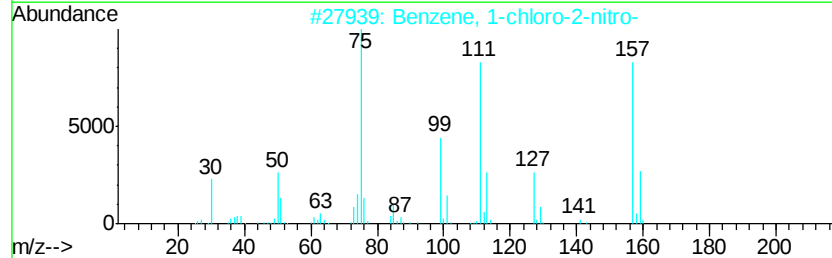
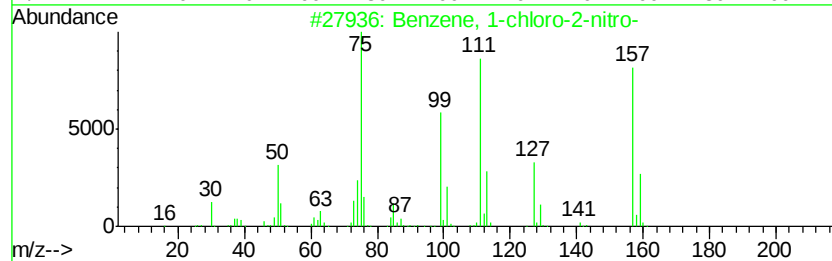
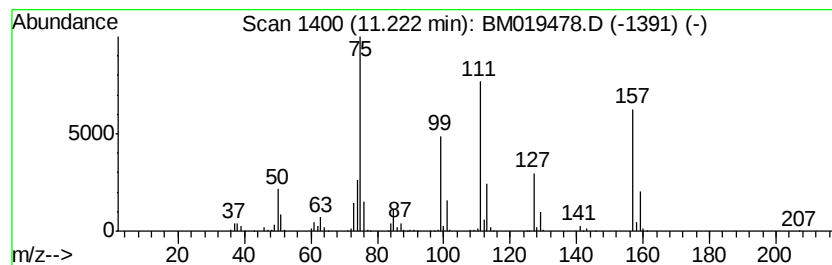
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Benzene, 1-chloro-2-nitro- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.22	31.07 ng/ul	2114620	Napthalene-d8	10.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	98
2		Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	97
3		Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	96
4		Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	95
5		Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM032619\
Data File : BM019478.D
Acq On : 26 Mar 2019 19:52
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Misc :
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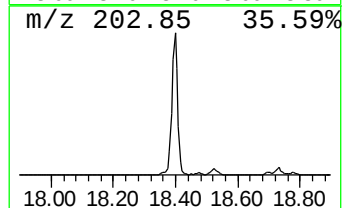
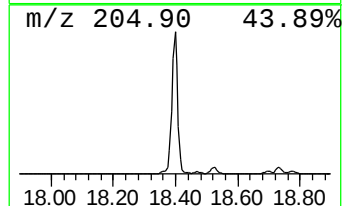
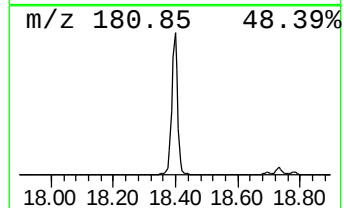
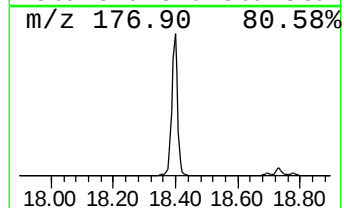
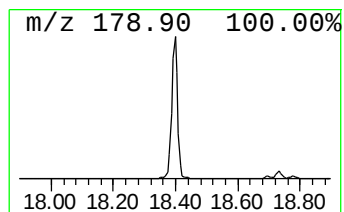
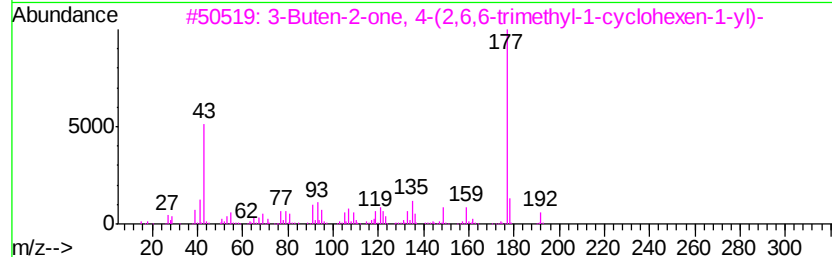
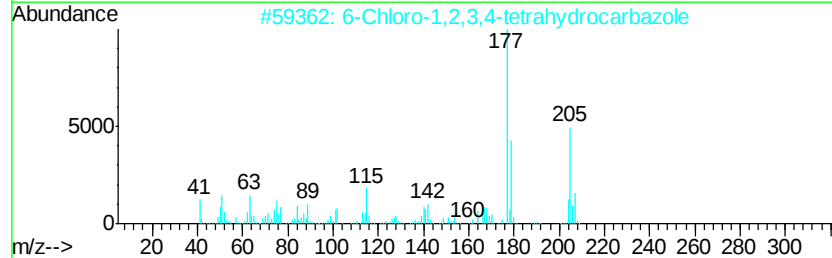
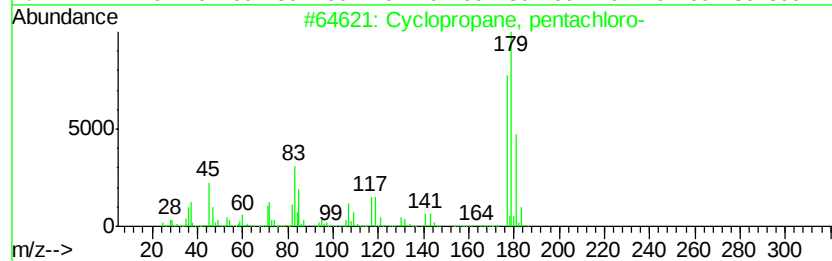
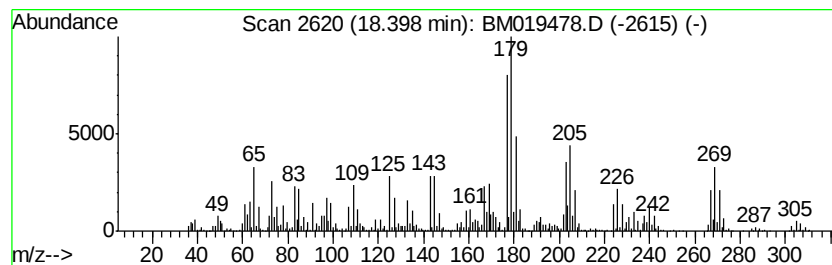
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 unknown-01 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.40	96.88 ng/ul	9110440	Phenanthrene-d10	17.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopropane, pentachloro-	212	C3HCl5	006262-51-7	37
2		6-Chloro-1,2,3,4-tetrahydrocarba...	205	C12H12ClN	036684-65-8	30
3		3-Buten-2-one, 4-(2,6,6-trimethy...	192	C13H20O	014901-07-6	15
4		4-Butylbenzoic acid, nonyl ester	304	C20H32O2	1000292-20-6	15
5		3-Buten-2-one, 4-(2,6,6-trimethy...	192	C13H20O	014901-07-6	15



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM032619\
Data File : BM019478.D
Acq On : 26 Mar 2019 19:52
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Sample : K2069-12
Misc :
ALS Vial : 10 Sample Multiplier: 1

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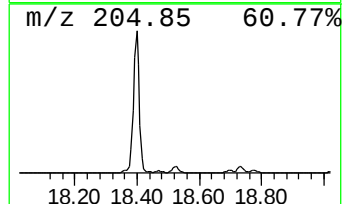
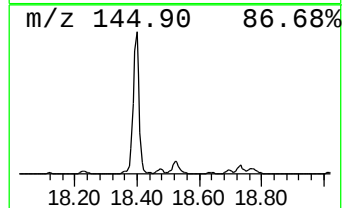
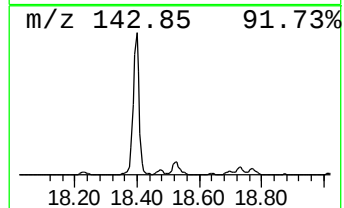
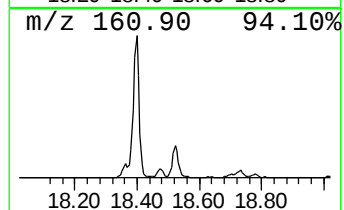
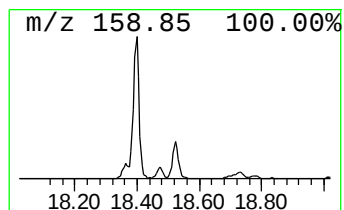
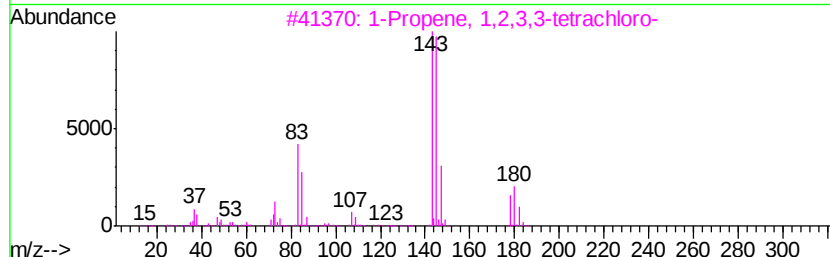
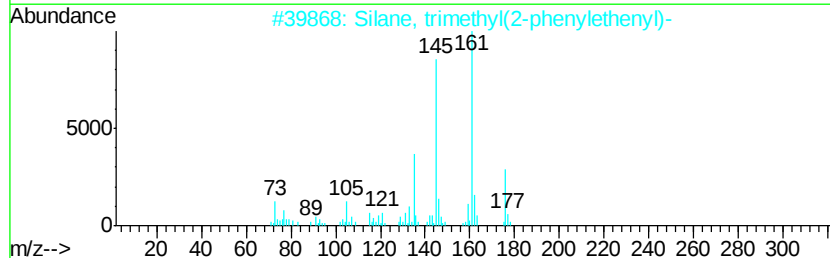
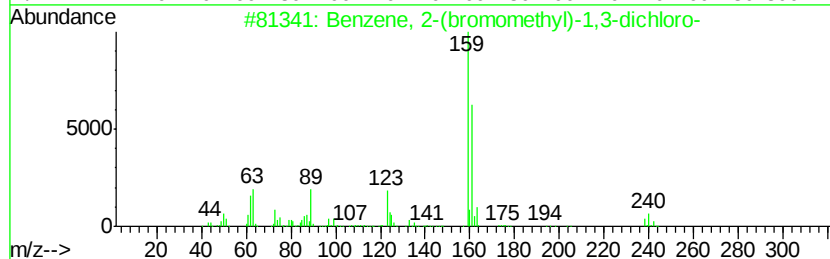
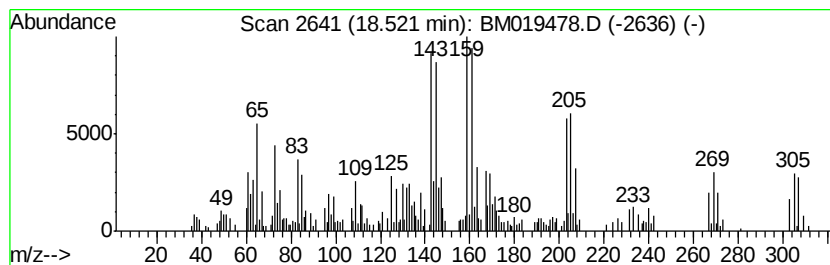
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 unknown-02 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.52	4.44 ng/ul	417526	Phenanthrene-d10	17.02

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-(bromomethyl)-1,3-dic...	238	C7H5BrCl2	020443-98-5	20
2			Silane, trimethyl(2-phenylethenyl)-	176	C11H16Si	018001-47-3	16
3			1-Propene, 1,2,3,3-tetrachloro-	178	C3H2Cl4	020589-85-9	15
4			Quinoline, 5-methoxy-	159	C10H9NO	006931-19-7	15
5			Benzene, 1,2,3-trichloro-4-methyl-	194	C7H5Cl3	007359-72-0	12



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM032619\
Data File : BM019478.D
Acq On : 26 Mar 2019 19:52
Operator : JU/SJ
Sample : K2069-12
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C09D9

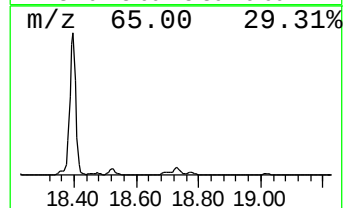
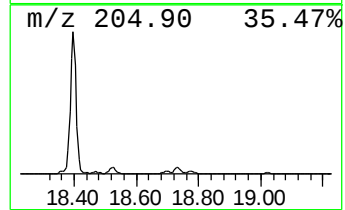
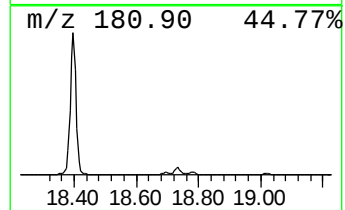
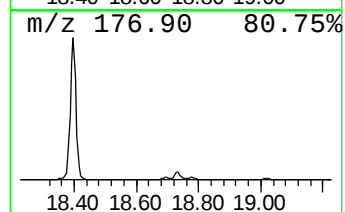
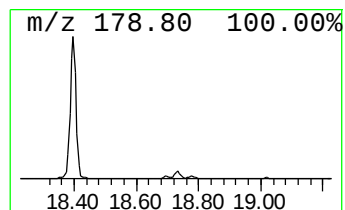
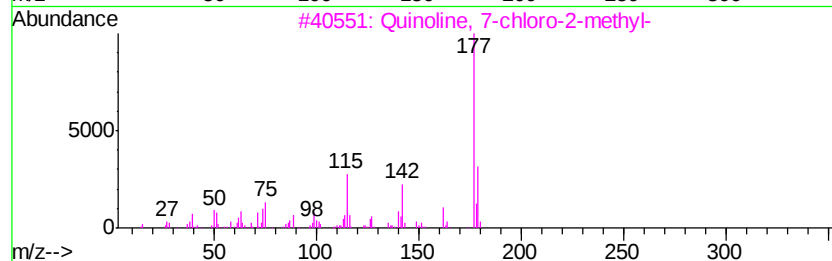
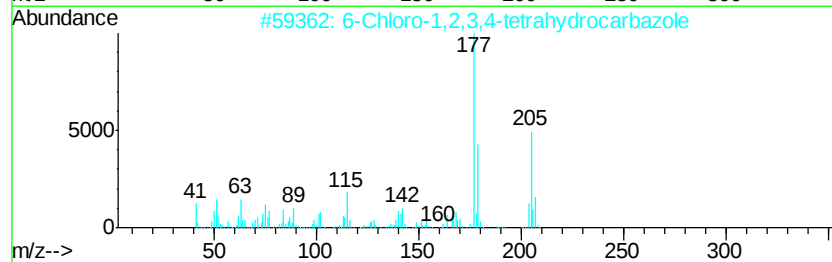
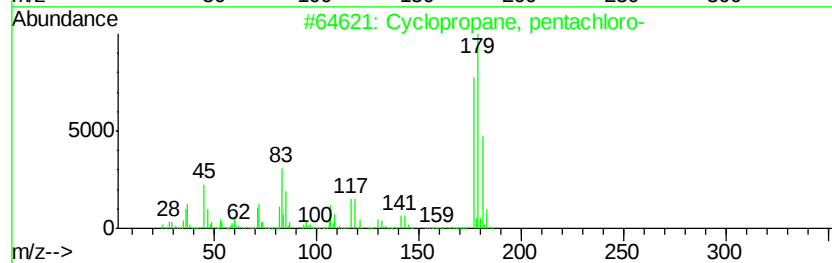
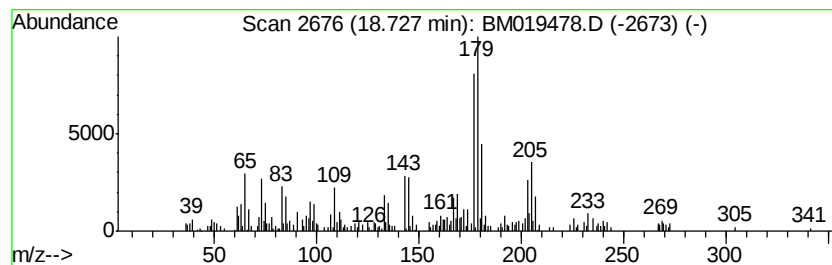
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 unknown-03 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.73	4.93 ng/ul	463854	Phenanthrene-d10	17.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopropane, pentachloro-	212	C3HCl5	006262-51-7	38
2		6-Chloro-1,2,3,4-tetrahydrocarba...	205	C12H12ClN	036684-65-8	27
3		Quinoline, 7-chloro-2-methyl-	177	C10H8ClN	004965-33-7	14
4		Methyl 2,5-dichlorothiophene-3-c...	210	C6H4Cl2O2S	145129-54-0	12
5		3-Buten-2-one, 4-(2,6,6-trimethy...	192	C13H20O	000079-77-6	11



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM032619\
 Data File : BM019478.D
 Acq On : 26 Mar 2019 19:52
 Operator : JU/SJ
 Sample : K2069-12
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C09D9

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
Benzene, chloro-	5.07	23.9	ng/ul	414036000	1	7.66	346172000	20.0
Benzene, 1,2-dich...	7.51	5.0	ng/ul	87040200	1	7.66	346172000	20.0
Benzene, 1,4-dich...	7.75	20.0	ng/ul	346172000	1	7.66	346172000	20.0
Benzene, 1,3-dich...	8.10	23.5	ng/ul	405997000	1	7.66	346172000	20.0
Benzene, 1-bromo-...	9.10	19.7	ng/ul	1342560	2	10.40	1361350	20.0
Benzene, 1-bromo-...	9.40	10.1	ng/ul	685959	2	10.40	1361350	20.0
Benzene, 1,2,3-tr...	10.30	2228.8	ng/ul	151708000	2	10.40	1361350	20.0
Benzene, 1,3,5-tr...	10.79	735.1	ng/ul	50037500	2	10.40	1361350	20.0
Benzene, 1-chloro...	11.01	160.8	ng/ul	10942100	2	10.40	1361350	20.0
Benzene, 1-chloro...	11.22	31.1	ng/ul	2114620	2	10.40	1361350	20.0
unknown-01	18.40	96.9	ng/ul	9110440	4	17.02	1880780	20.0
unknown-02	18.52	4.4	ng/ul	417526	4	17.02	1880780	20.0
unknown-03	18.73	4.9	ng/ul	463854	4	17.02	1880780	20.0