

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM032719\
 Data File : BM019517.D
 Acq On : 27 Mar 2019 22:35
 Operator : JU/SJ
 Sample : K2069-05
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C09D7

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	5.005	323	343	347	rBV3	49107590	266249384	100.00%	29.520%
2	6.281	554	560	568	rBV	175149	296134	0.11%	0.033%
3	6.957	668	675	684	rBV	230601	470734	0.18%	0.052%
4	7.128	698	704	718	rVB	283475	560331	0.21%	0.062%
5	7.493	755	766	778	rBV	16781905	42338368	15.90%	4.694%
6	7.669	778	796	800	rBV4	49147494	200058621	75.14%	22.181%
7	8.010	833	854	855	rBV4	49585904	198697693	74.63%	22.031%
8	8.310	899	905	915	rVB	225161	398820	0.15%	0.044%
9	8.763	975	982	984	rBV	380104	596149	0.22%	0.066%
10	8.810	984	990	1005	rVB	4625481	7506428	2.82%	0.832%
11	9.081	1030	1036	1046	rVB	314712	502408	0.19%	0.056%
12	9.475	1097	1103	1107	rBV	317765	584691	0.22%	0.065%
13	10.010	1187	1194	1204	rBV	415766	772619	0.29%	0.086%
14	10.269	1224	1238	1242	rBV	44745079	101468510	38.11%	11.250%
15	10.381	1251	1257	1263	rVB	536107	779427	0.29%	0.086%
16	10.763	1309	1322	1335	rBV	18644020	31130995	11.69%	3.452%
17	10.992	1353	1361	1375	rBV	4411048	7105551	2.67%	0.788%
18	11.204	1390	1397	1412	rBV	634653	1352605	0.51%	0.150%
19	12.410	1592	1602	1613	rBV	840448	1629683	0.61%	0.181%
20	13.033	1701	1708	1720	rVB	2467126	3762644	1.41%	0.417%
21	13.669	1808	1816	1825	rBV	899347	1513008	0.57%	0.168%
22	13.939	1857	1862	1872	rVB	1011623	1451438	0.55%	0.161%
23	14.245	1907	1914	1922	rBV2	680668	1027324	0.39%	0.114%
24	15.245	2077	2084	2093	rBV	1184656	1771469	0.67%	0.196%
25	15.392	2104	2109	2121	rVB	394631	624385	0.23%	0.069%
26	17.004	2378	2383	2392	rVB2	872984	1228105	0.46%	0.136%
27	17.104	2395	2400	2410	rVV2	1427928	2005060	0.75%	0.222%
28	18.392	2613	2619	2627	rVV	9636068	12695405	4.77%	1.408%
29	18.515	2635	2640	2648	rVB	339358	468354	0.18%	0.052%
30	18.686	2664	2669	2672	rBV2	168556	272995	0.10%	0.030%
31	18.721	2672	2675	2679	rVV	419409	555951	0.21%	0.062%
32	19.415	2786	2793	2807	rBV	1915179	2516951	0.95%	0.279%
33	21.215	3094	3099	3106	rBV	1258440	1769026	0.66%	0.196%
34	22.721	3352	3355	3360	rVB	199071	277863	0.10%	0.031%

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35	23.280	3443	3450	3462	rBV	2244252	4095355	1.54%	0.454%
36	23.421	3468	3474	3488	rVB	1205179	2280928	0.86%	0.253%
37	23.897	3551	3555	3566	rVB	268396	499808	0.19%	0.055%
38	24.627	3674	3679	3688	rVB	276157	604620	0.23%	0.067%

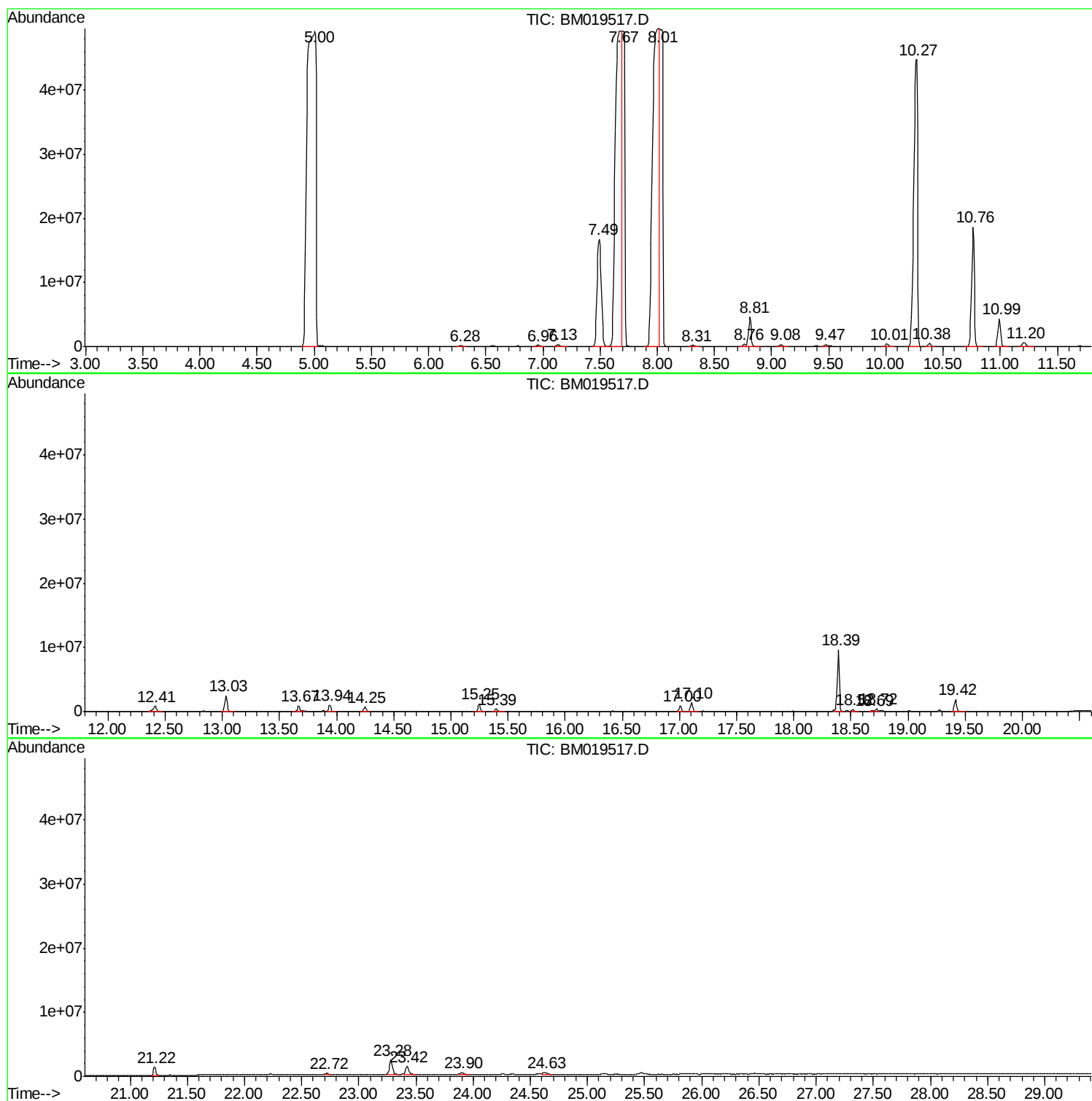
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
Quant Title : SVOA CALIBRATION

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



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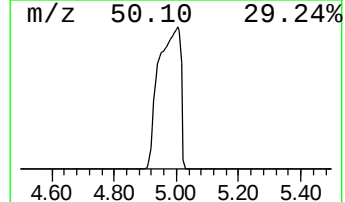
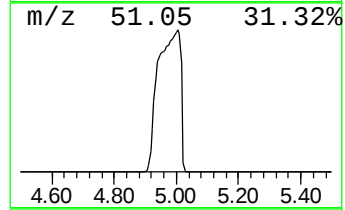
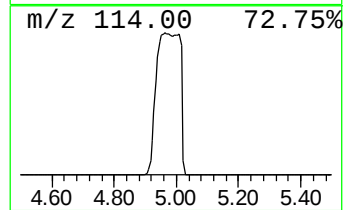
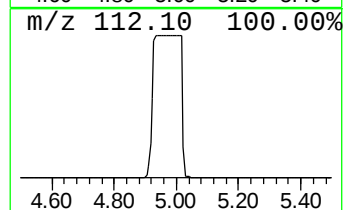
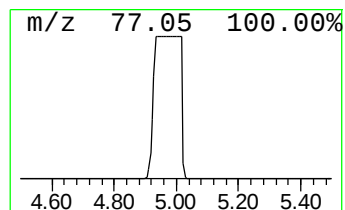
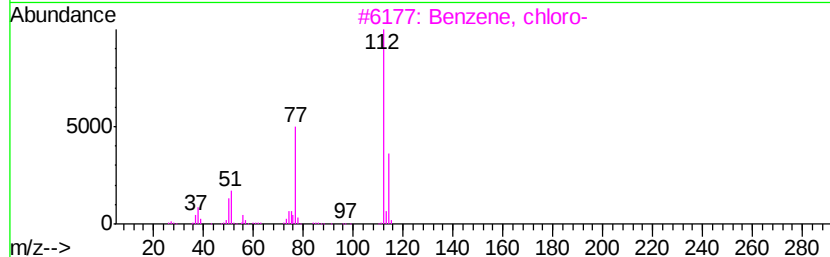
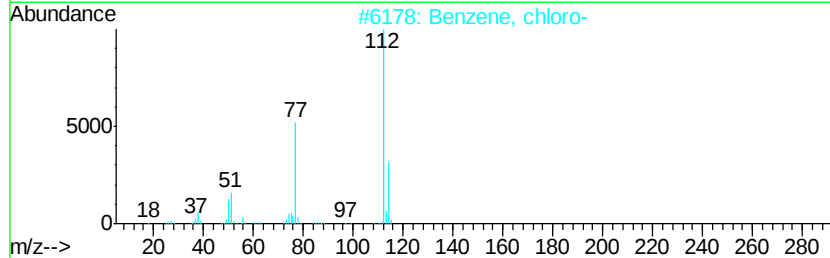
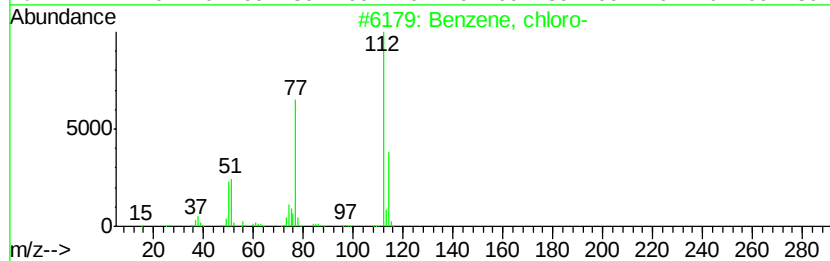
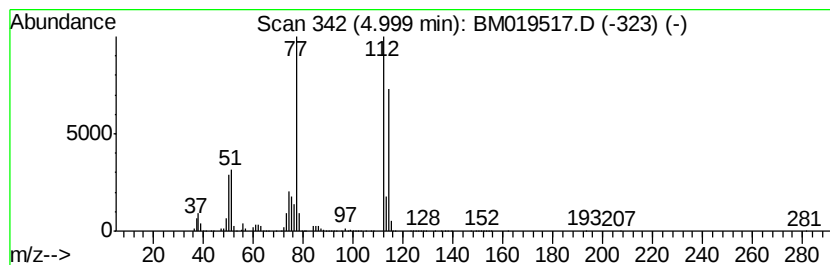
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Quant Title : SVOA CALIBRATION

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.
5.00	26.62 ng/ul	266249000	1,4-Dichlorobenzene-d4	7.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, chloro-	112	C6H5Cl	000108-90-7	90
2		Benzene, chloro-	112	C6H5Cl	000108-90-7	68
3		Benzene, chloro-	112	C6H5Cl	000108-90-7	64
4		Benzene, chloro-	112	C6H5Cl	000108-90-7	53
5		Benzene, 1,4-difluoro-	114	C6H4F2	000540-36-3	9



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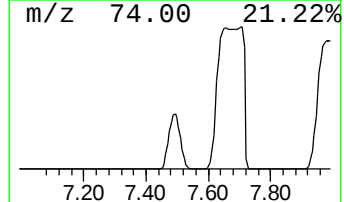
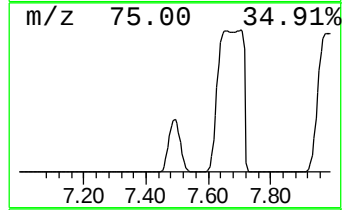
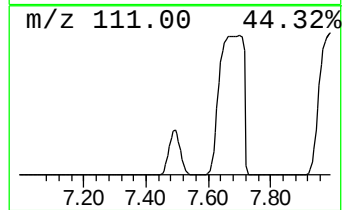
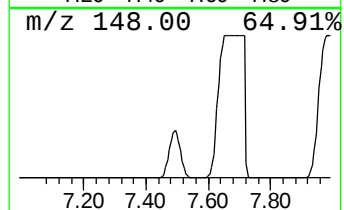
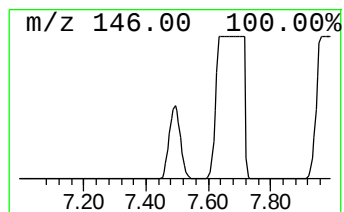
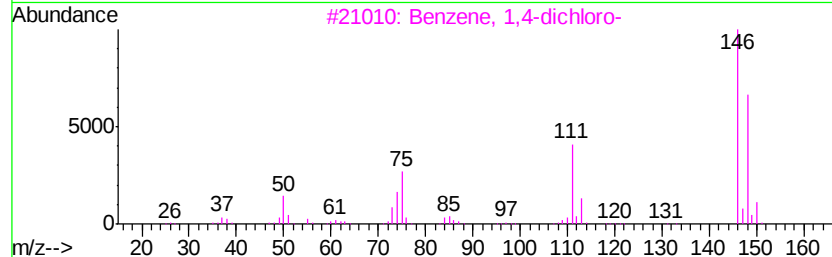
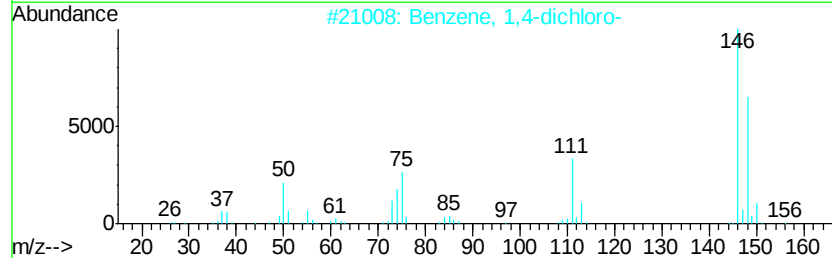
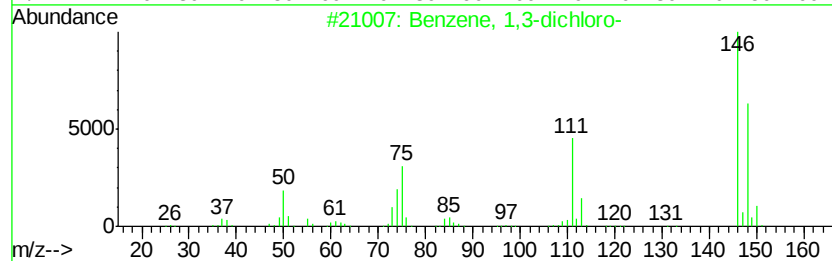
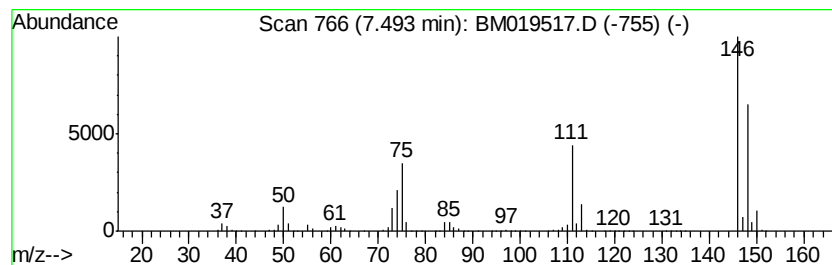
Instrument :
BNA_M
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 2 Benzene, 1,3-dichloro- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.49	4.23 ng/ul	42338400	1,4-Dichlorobenzene-d4	7.63
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Benzene, 1,3-dichloro-	146 C6H4Cl2	000541-73-1	96
2	Benzene, 1,4-dichloro-	146 C6H4Cl2	000106-46-7	96
3	Benzene, 1,4-dichloro-	146 C6H4Cl2	000106-46-7	96
4	Benzene, 1,3-dichloro-	146 C6H4Cl2	000541-73-1	95
5	Benzene, 1,3-dichloro-	146 C6H4Cl2	000541-73-1	95



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Instrument :
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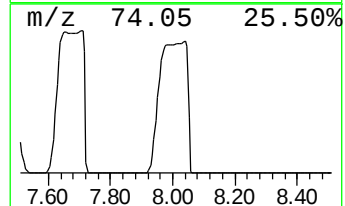
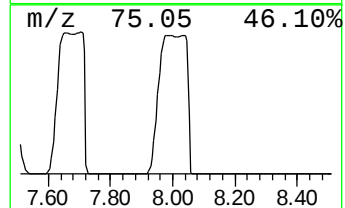
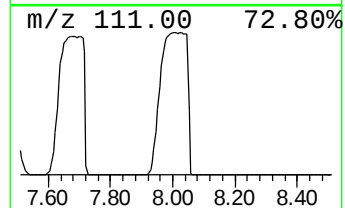
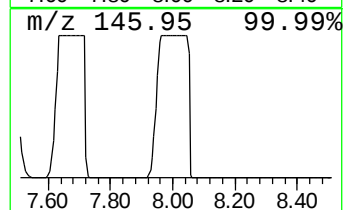
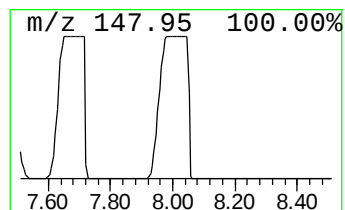
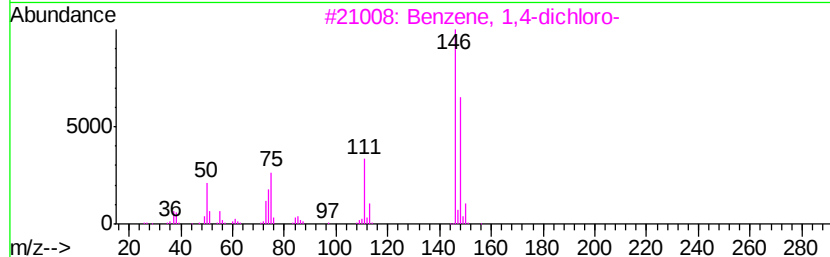
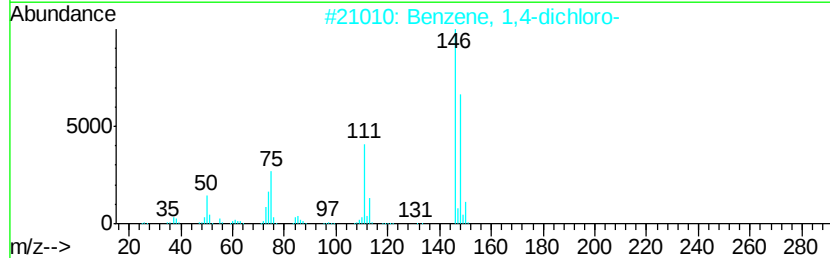
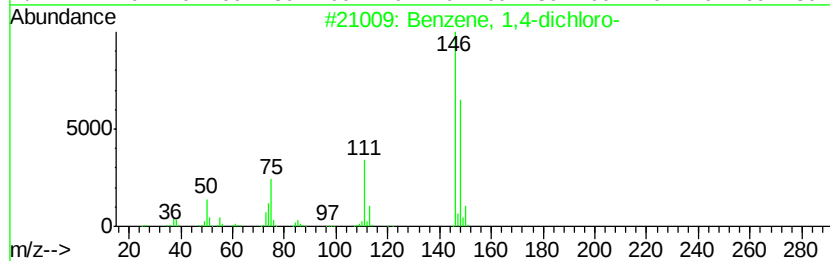
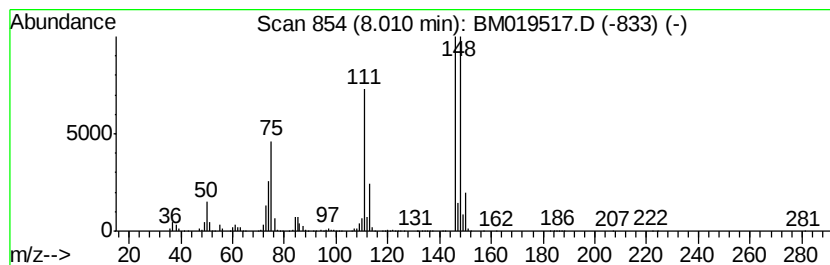
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
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Peak Number 3 Benzene, 1,4-dichloro- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.01	19.86 ng/ul	198698000	1,4-Dichlorobenzene-d4	7.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	94
2		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	91
3		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	87
4		Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	81
5		Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	81



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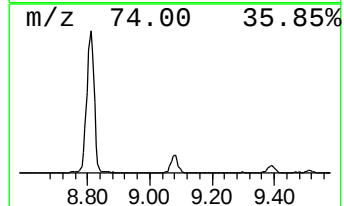
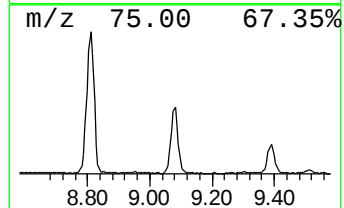
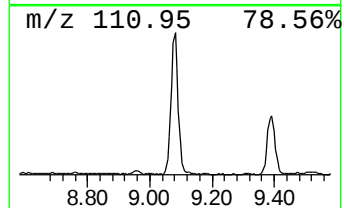
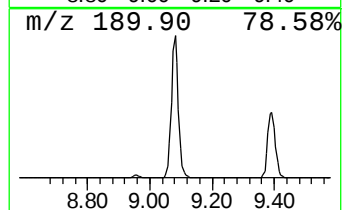
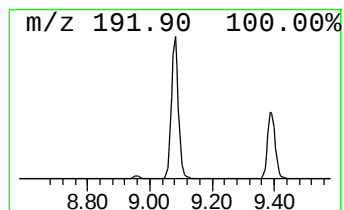
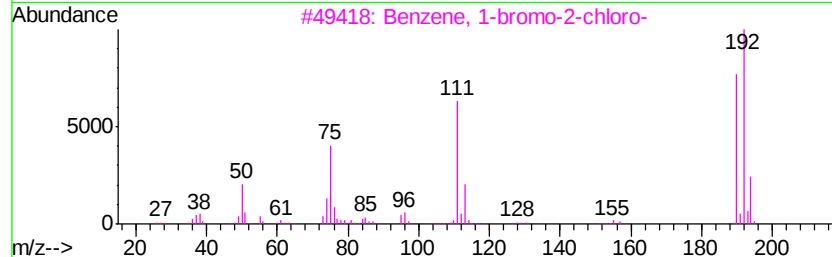
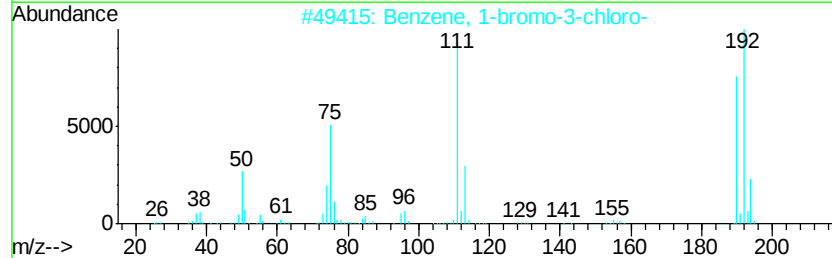
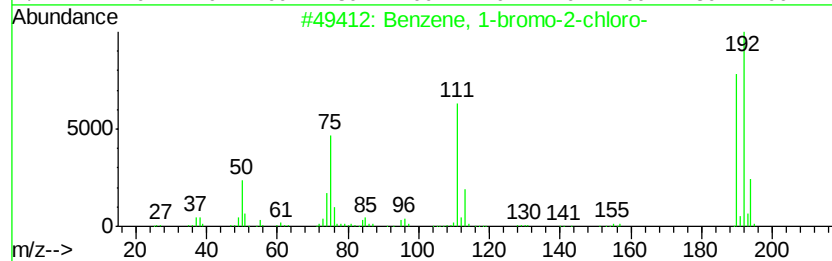
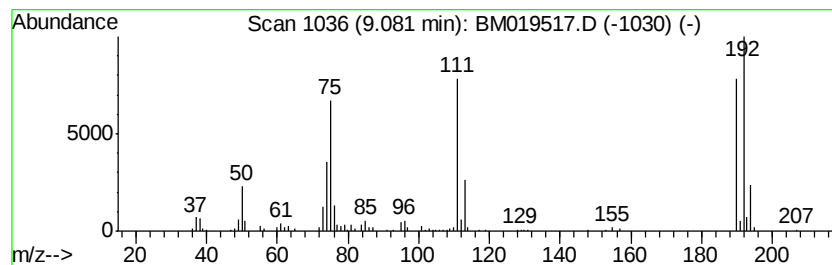
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Peak Number 4 Benzene, 1-bromo-2-chloro- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.08	12.89 ng/ul	502408	Napthalene-d8	10.38

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-3-chloro-	190	C6H4BrCl	000108-37-2	98
3		Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	98
4		Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	97
5		Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	97



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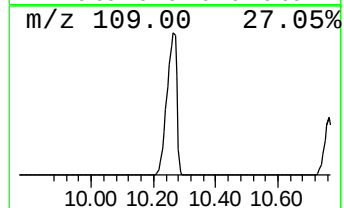
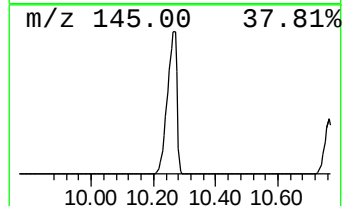
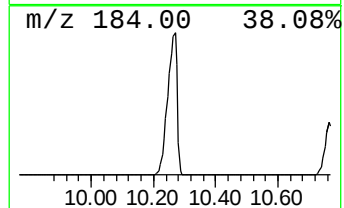
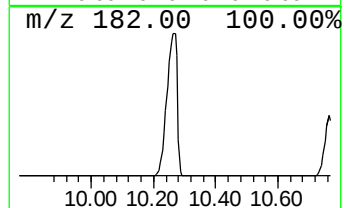
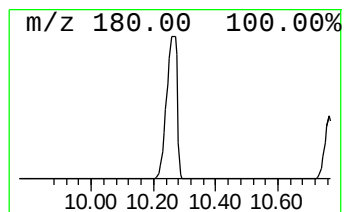
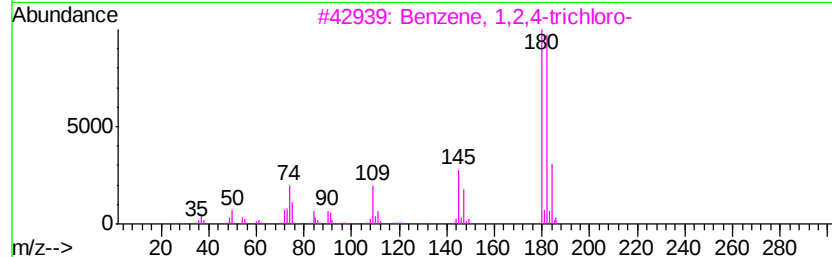
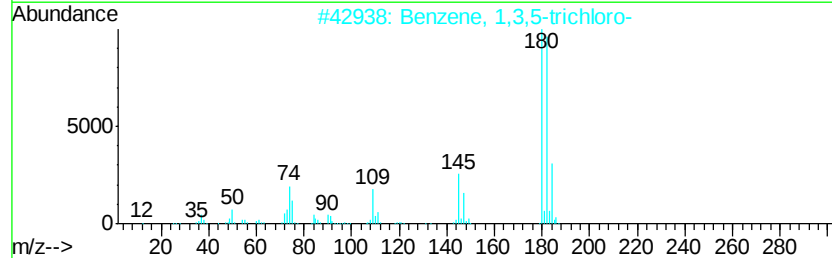
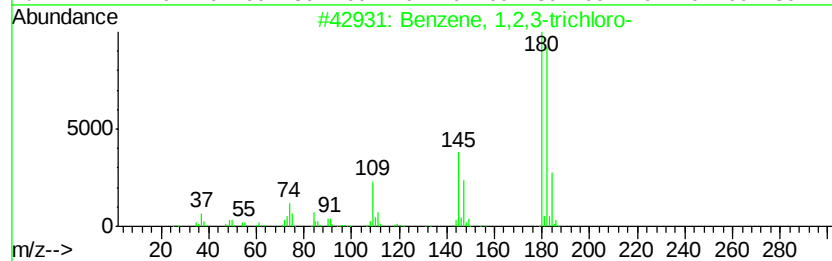
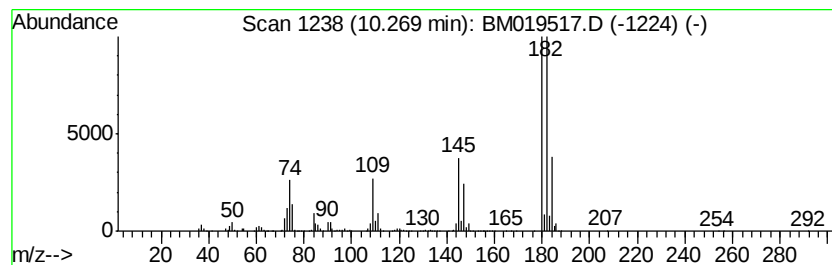
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.27	2603.67 ng/ul	101469000	Napthalene-d8	10.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	98
2		Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	96
3		Benzene, 1,2,4-trichloro-	180	C6H3Cl3	000120-82-1	95
4		Benzene, 1,2,4-trichloro-	180	C6H3Cl3	000120-82-1	95
5		Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	95



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Data File : BM019517.D
Acq On : 27 Mar 2019 22:35
Operator : JU/SJ
Sample : K2069-05
Misc :
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Instrument :
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ClientSampleId :
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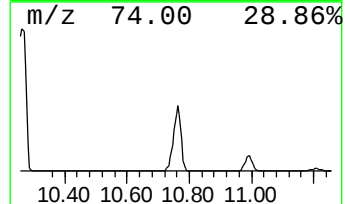
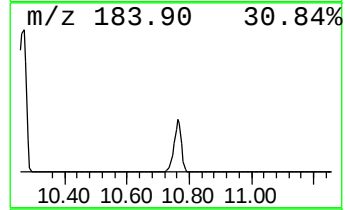
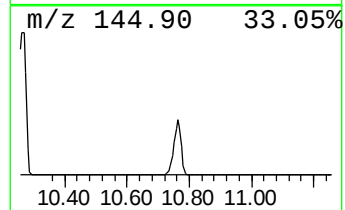
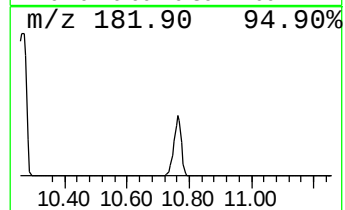
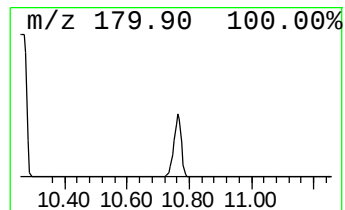
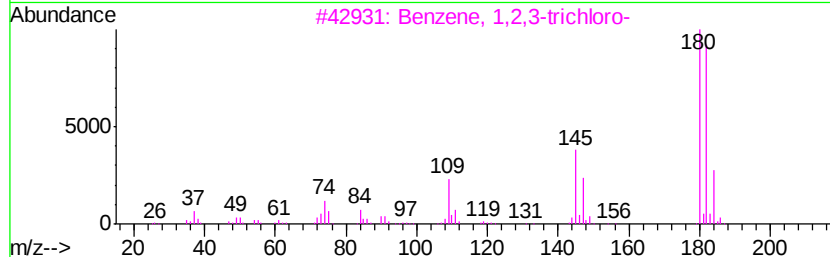
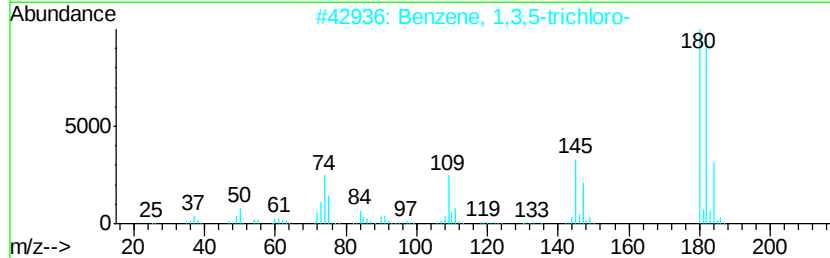
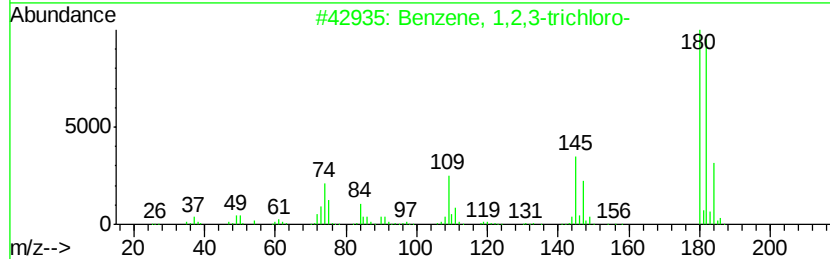
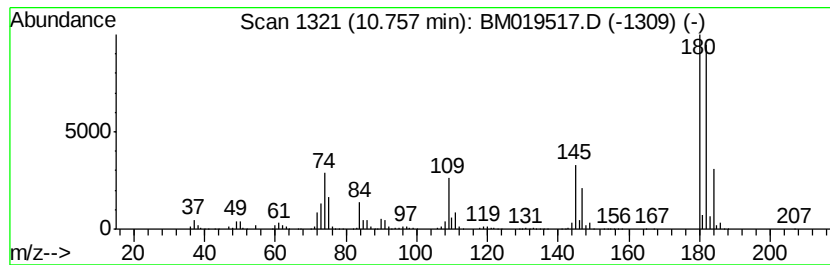
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzene, 1,3,5-trichloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.76	798.82 ng/ul	31131000	Napthalene-d8	10.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	98
2		Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	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\BM032719\
Data File : BM019517.D
Acq On : 27 Mar 2019 22:35
Operator : JU/SJ
Sample : K2069-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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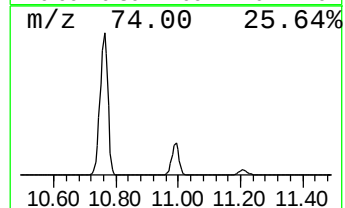
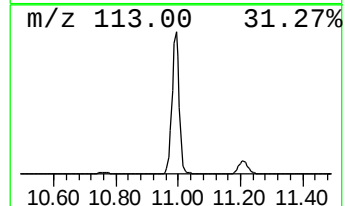
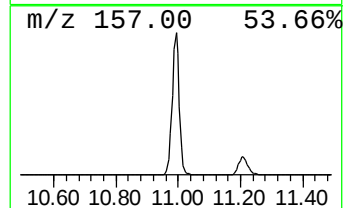
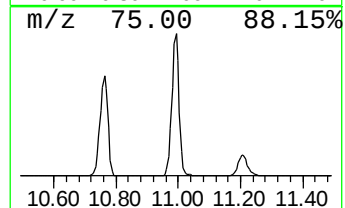
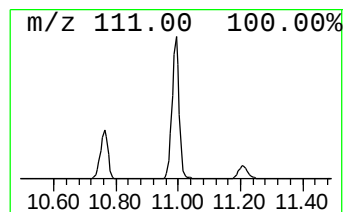
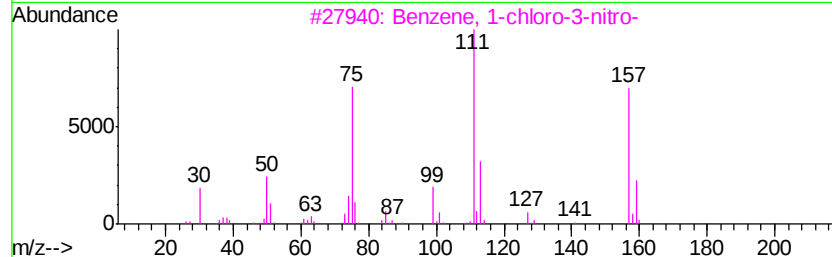
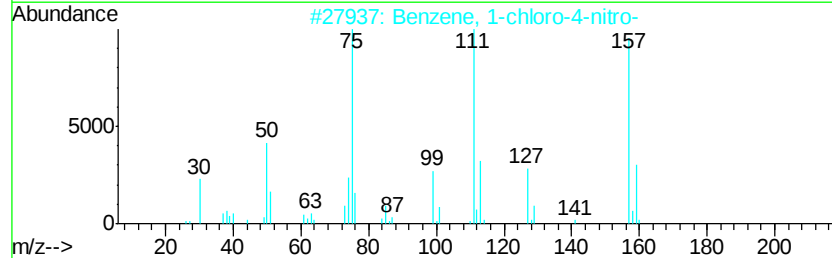
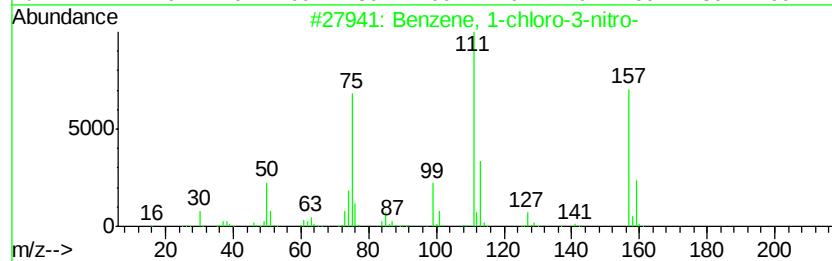
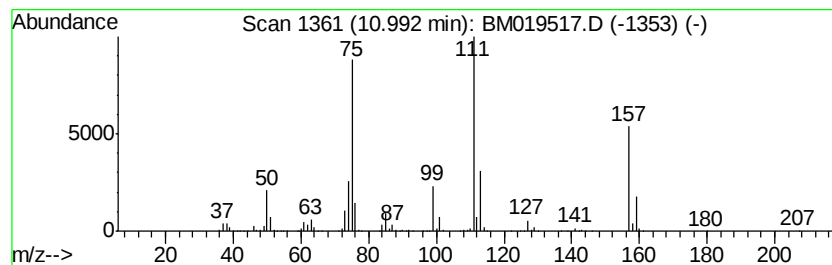
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Benzene, 1-chloro-3-nitro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.99	182.33 ng/ul	7105550	Naphthalene-d8	10.38

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-4-nitro-	157	C6H4ClNO2	000100-00-5	94
3		Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	93
4		Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	93
5		Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM032719\
Data File : BM019517.D
Acq On : 27 Mar 2019 22:35
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Sample : K2069-05
Misc :
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ClientSampleId :
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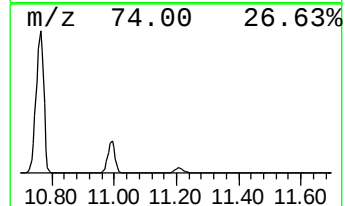
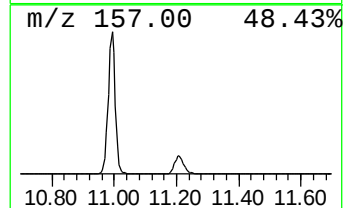
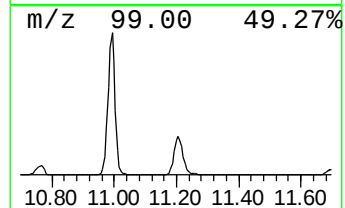
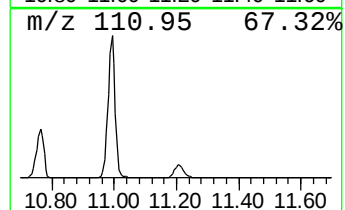
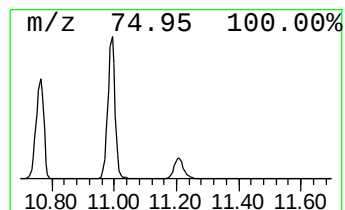
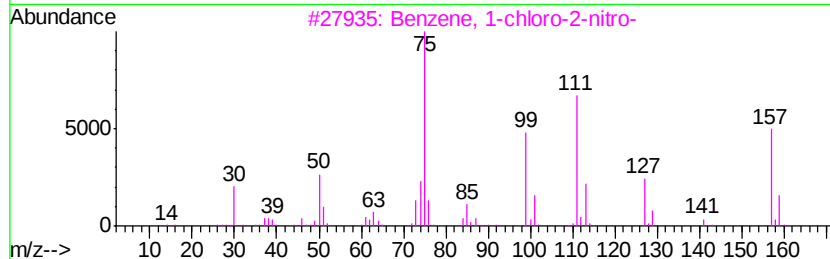
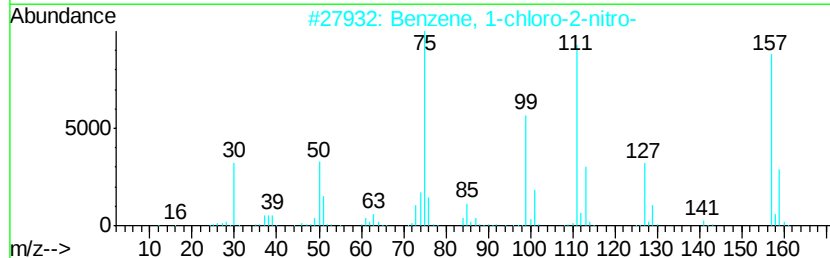
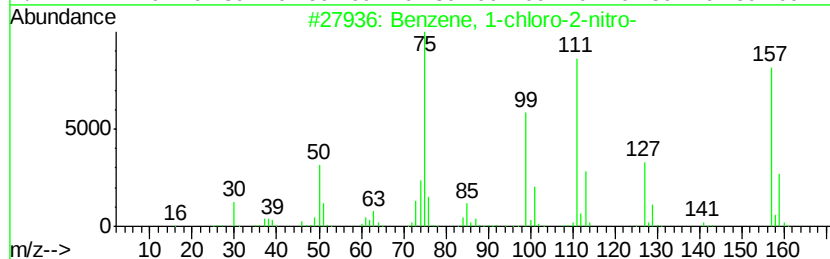
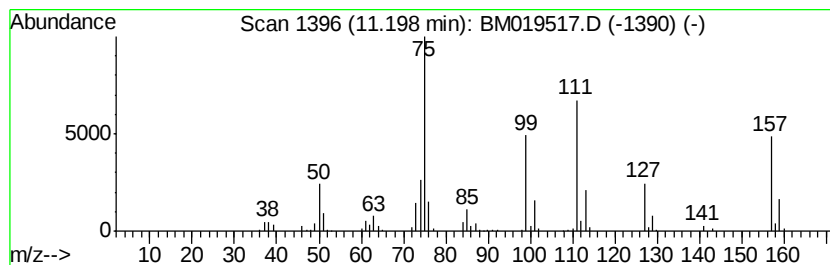
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Benzene, 1-chloro-2-nitro- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.20	34.71 ng/ul	1352610	Napthalene-d8	10.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	97
2		Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	96
3		Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	96
4		Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	94
5		Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM032719\
Data File : BM019517.D
Acq On : 27 Mar 2019 22:35
Operator : JU/SJ
Sample : K2069-05
Misc :
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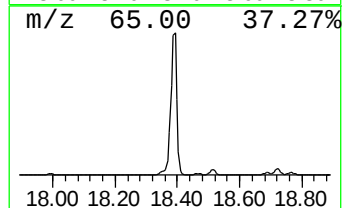
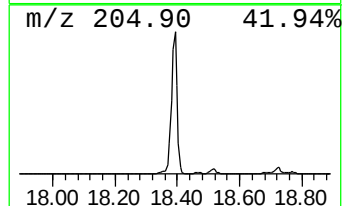
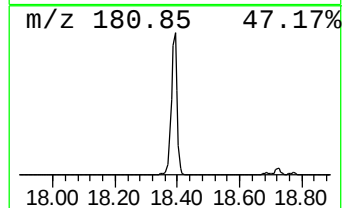
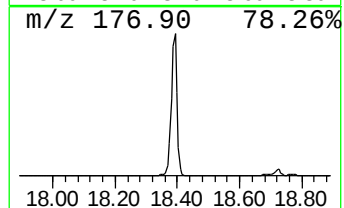
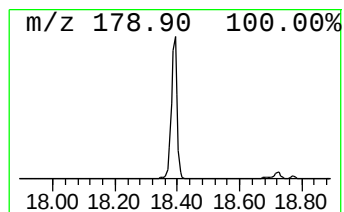
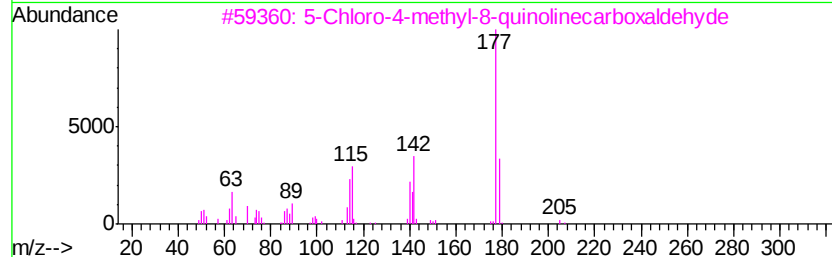
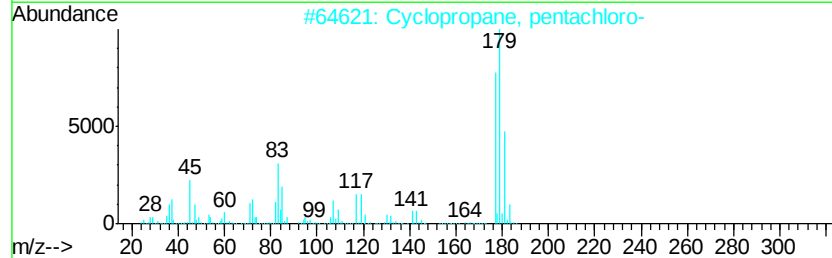
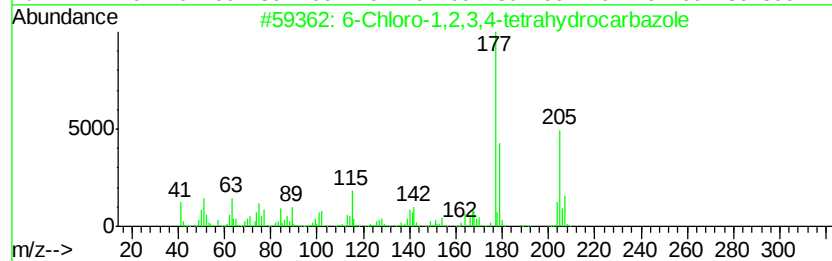
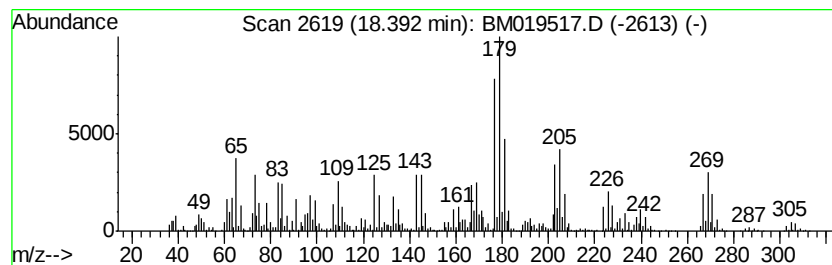
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 unknown-01 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.39	206.75 ng/ul	12695400	Phenanthrene-d10	17.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	6-Chloro-1,2,3,4-tetrahydrocarba...	205	C12H12ClN	036684-65-8	38
2		Cyclopropane, pentachloro-	212	C3HCl5	006262-51-7	37
3		5-Chloro-4-methyl-8-quinolinecar...	205	C11H8ClNO	028662-47-7	27
4		Octamethyl-1,4-cyclohexadiene	192	C14H24	172684-93-4	25
5		2H-1-Benzopyran, 3,4,4a,5,6,8a-h...	194	C13H22O	063335-66-0	11



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM032719\
Data File : BM019517.D
Acq On : 27 Mar 2019 22:35
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Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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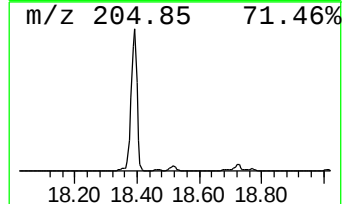
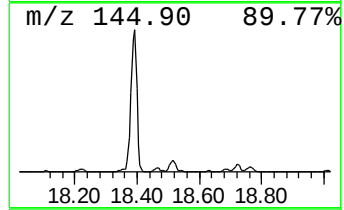
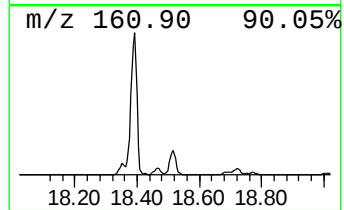
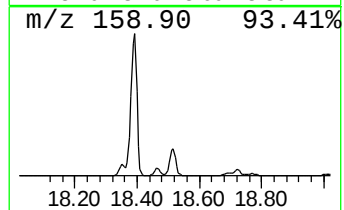
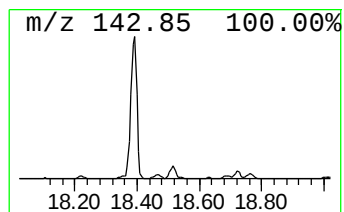
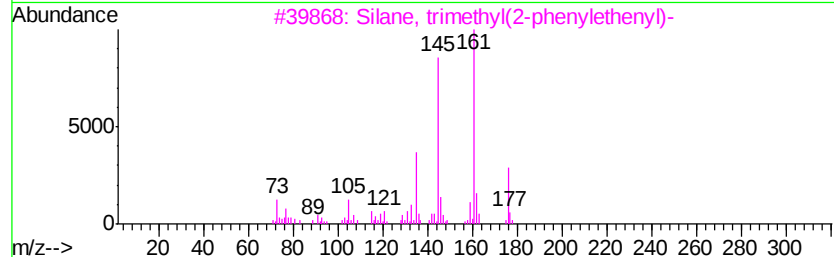
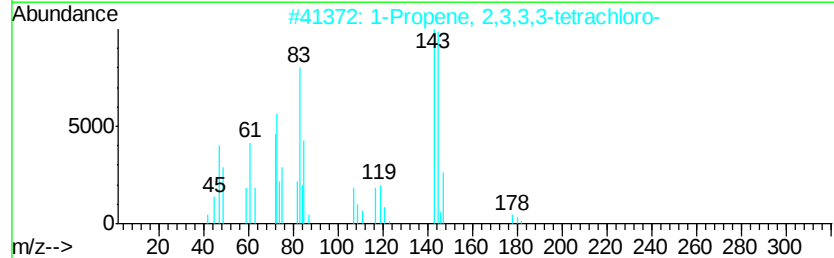
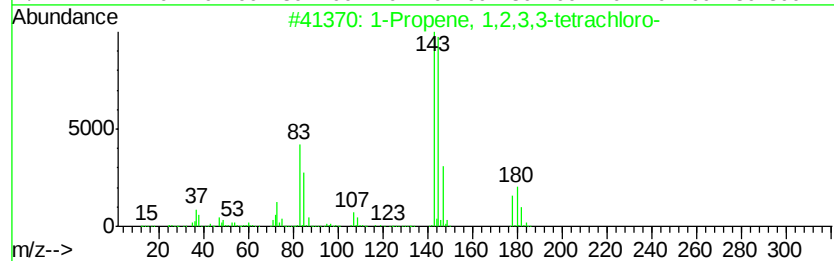
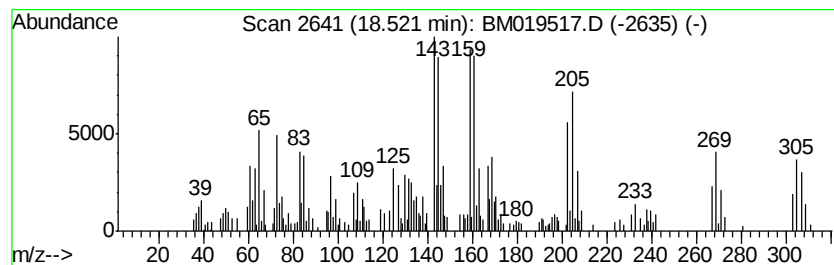
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 unknown-02 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.52	7.63 ng/ul	468354	Phenanthrene-d10	17.00

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Propene, 1,2,3,3-tetrachloro-	178	C3H2Cl4	020589-85-9	18
2			1-Propene, 2,3,3,3-tetrachloro-	178	C3H2Cl4	016500-91-7	14
3			Silane, trimethyl(2-phenylethenyl)-	176	C11H16Si	018001-47-3	12
4			2-(5-Aminopentyl)benzimidazole	203	C12H17N3	039650-63-0	11
5			Hydrazinecarboxamide, 2-(1-methy...	203	C11H13N3O	005468-31-5	10



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM032719\
Data File : BM019517.D
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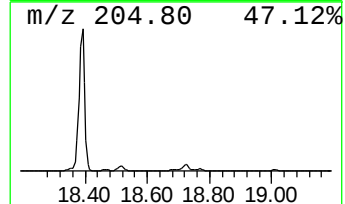
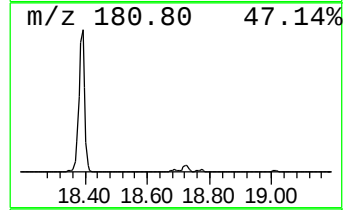
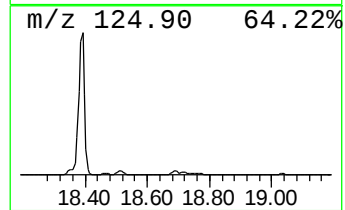
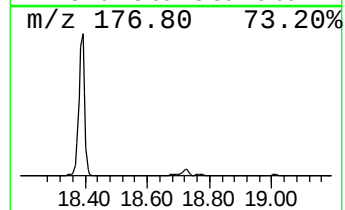
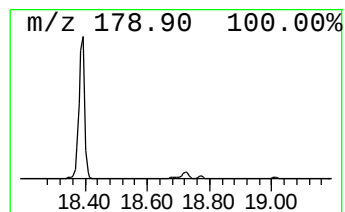
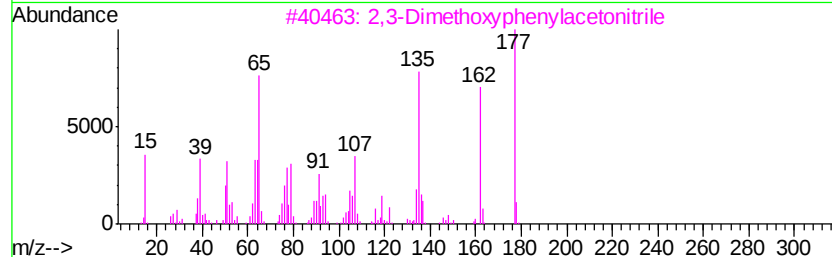
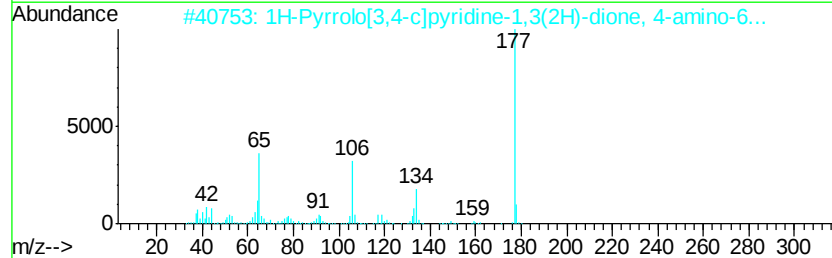
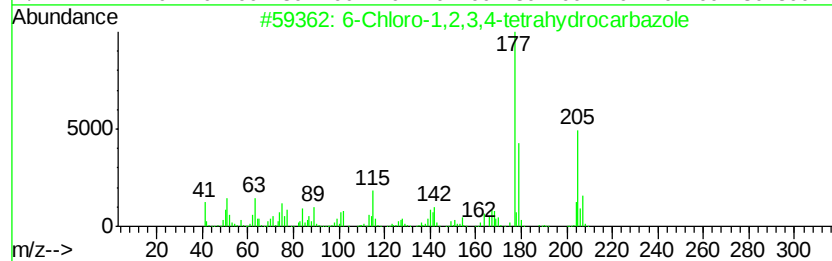
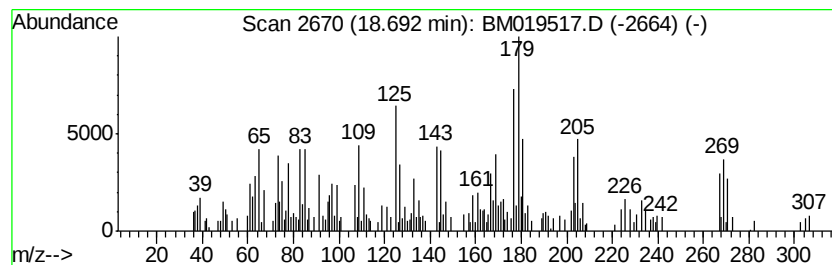
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 6-Chloro-1,2,3,4-tetrahydro... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.69	4.45 ng/ul	272995	Phenanthrene-d10	17.00

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			6-Chloro-1,2,3,4-tetrahydrocarba...	205	C12H12ClN	036684-65-8	53
2			1H-Pyrrolo[3,4-c]pyridine-1,3(2H)...	177	C8H7N3O2	090004-20-9	16
3			2,3-Dimethoxyphenylacetoneitrile	177	C10H11NO2	004468-57-9	16
4			Cyclopropane, pentachloro-	212	C3HCl5	006262-51-7	12
5			Germanium tetrachloride	214	Cl4Ge	010038-98-9	12



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM032719\
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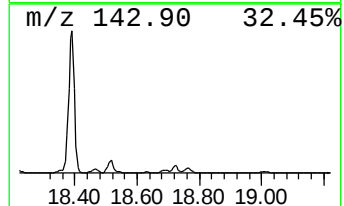
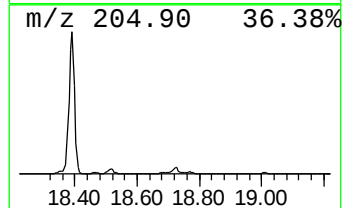
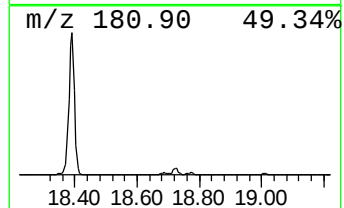
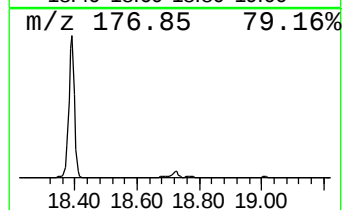
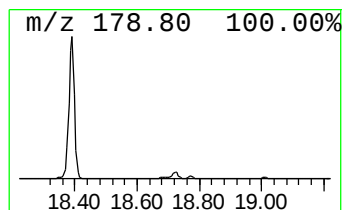
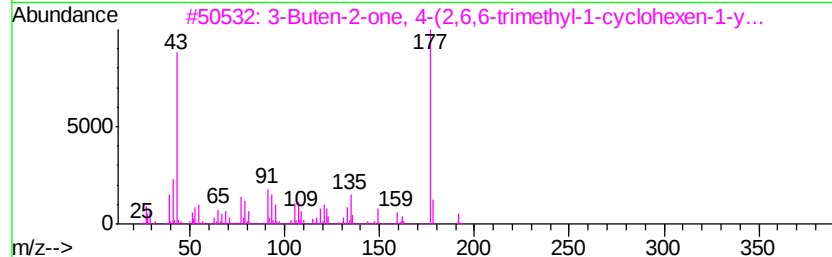
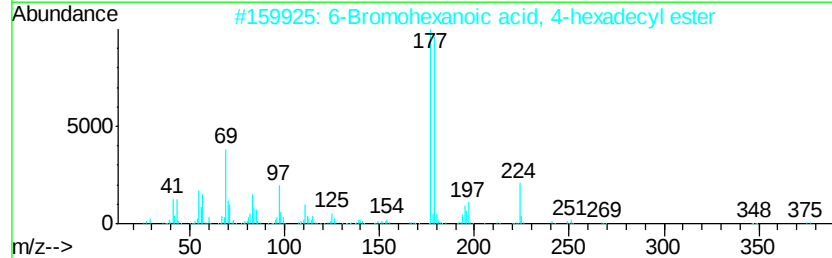
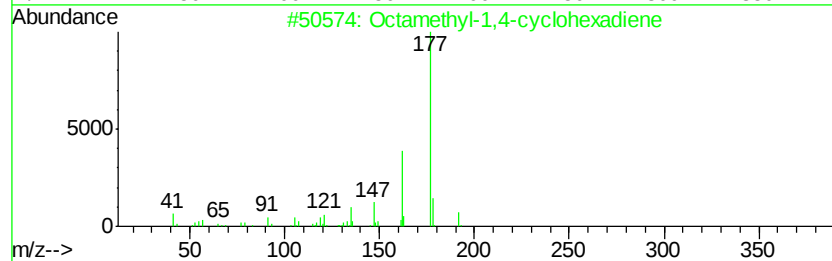
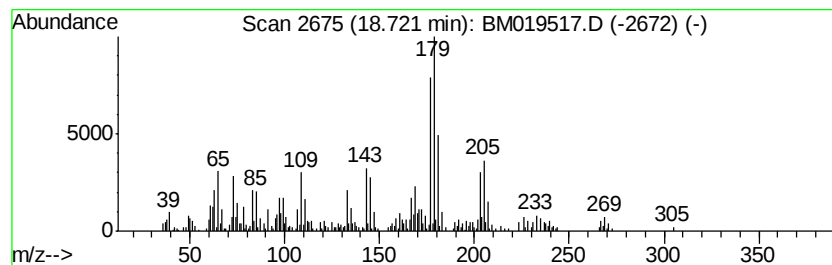
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 unknown-03 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.72	9.05 ng/ul	555951	Phenanthrene-d10	17.00

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octamethyl-1,4-cyclohexadiene	192	C14H24	172684-93-4	25
2			6-Bromohexanoic acid, 4-hexadecy...	418	C22H43BrO2	1000282-90-6	22
3			3-Buten-2-one, 4-(2,6,6-trimethy...	192	C13H20O	000079-77-6	15
4			Benzenamine, 4-fluoro-3-(trifluo...	179	C7H5F4N	002357-47-3	12
5			Silane, trimethyl(1-phenylethoxy)-	194	C11H18OSi	014856-75-8	11



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM032719\
Data File : BM019517.D
Acq On : 27 Mar 2019 22:35
Operator : JU/SJ
Sample : K2069-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C09D7

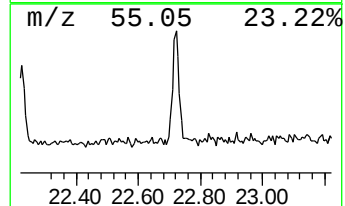
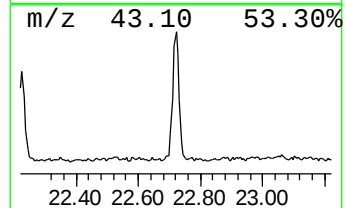
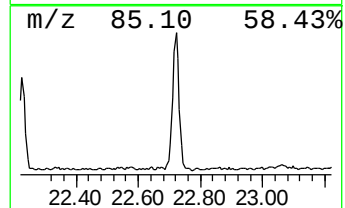
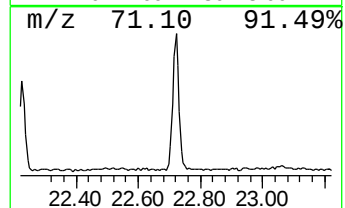
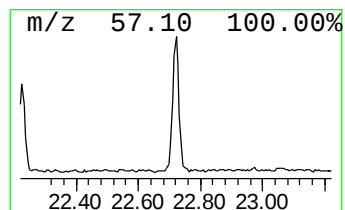
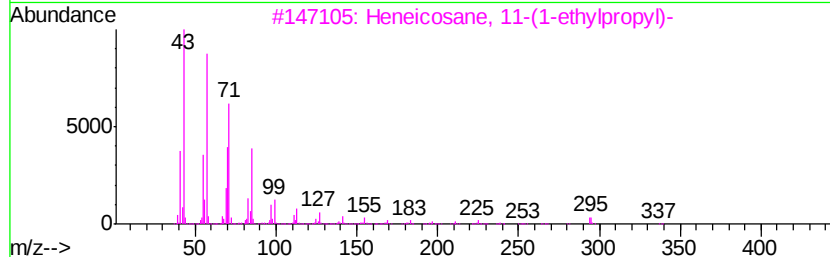
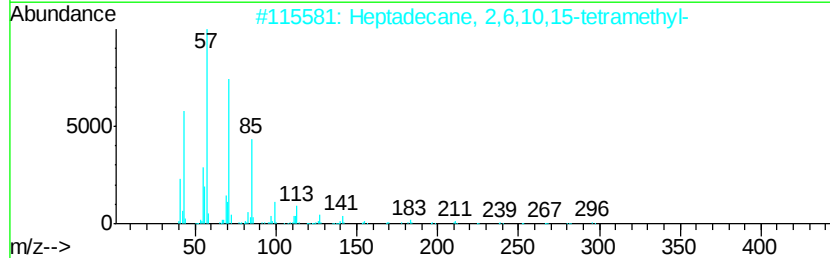
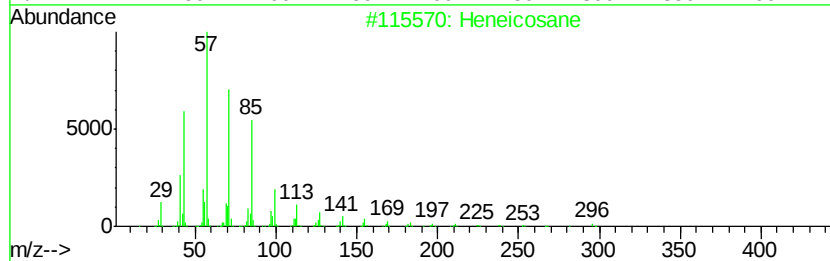
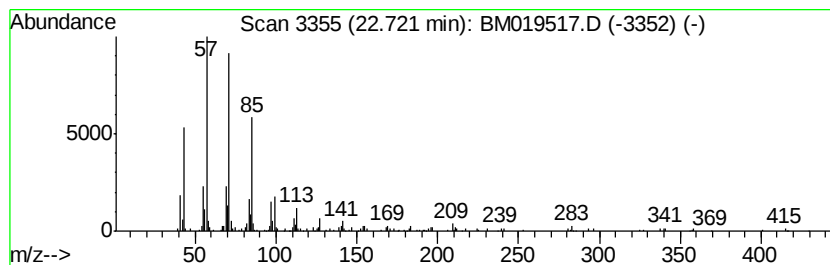
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.72	2.44 ng/ul	277863	Perylene-d12	23.42

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	96
2			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	93
3			Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91
4			Octadecane, 2-methyl-	268	C19H40	001560-88-9	90
5			Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM032719\
Data File : BM019517.D
Acq On : 27 Mar 2019 22:35
Operator : JU/SJ
Sample : K2069-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C09D7

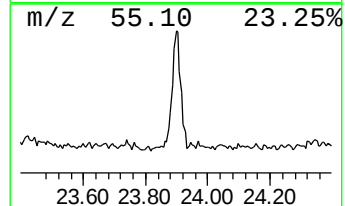
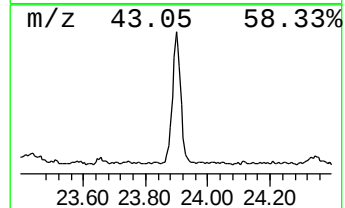
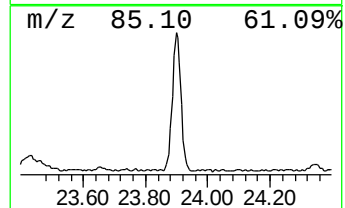
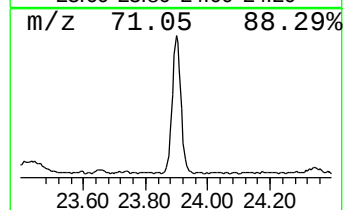
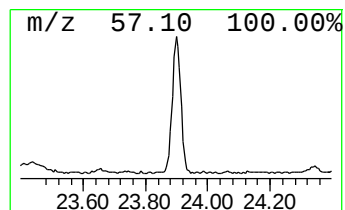
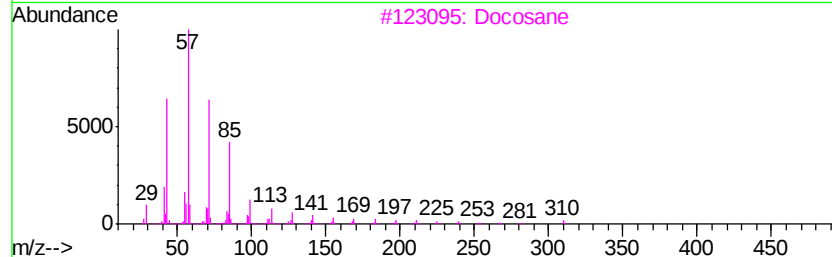
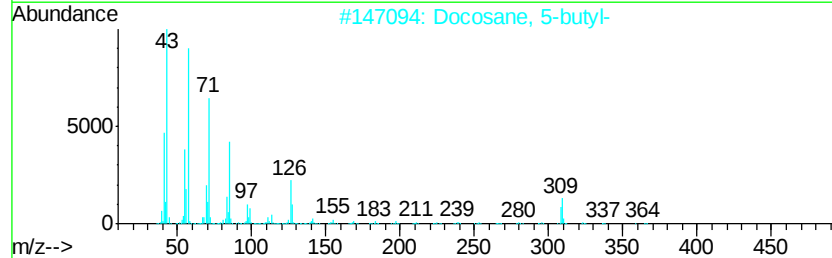
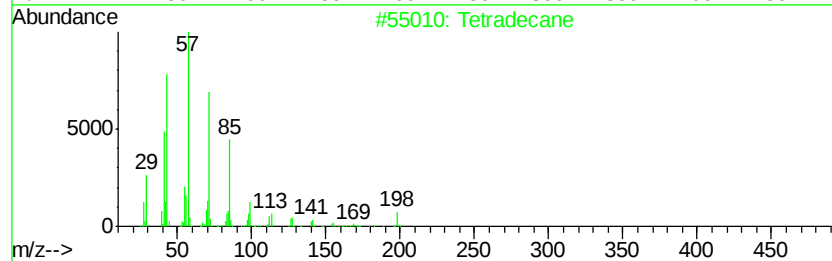
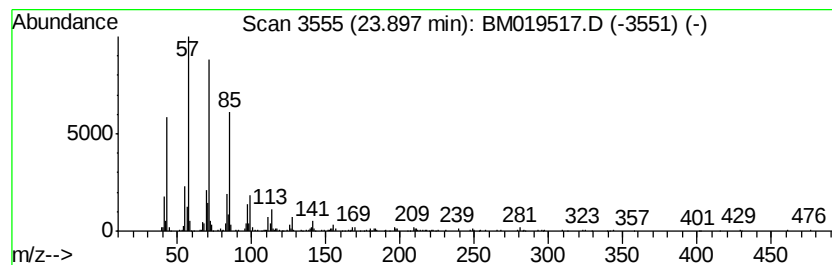
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 (DEL) Alkane: Straight-Chai... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.90	4.38 ng/ul	499808	Perylene-d12	23.42

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecane	198	C14H30	000629-59-4	93
2			Docosane, 5-butyl-	366	C26H54	055282-16-1	90
3			Docosane	310	C22H46	000629-97-0	86
4			Octadecane	254	C18H38	000593-45-3	83
5			Tetradecane, 4-methyl-	212	C15H32	025117-24-2	74



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM032719\
Data File : BM019517.D
Acq On : 27 Mar 2019 22:35
Operator : JU/SJ
Sample : K2069-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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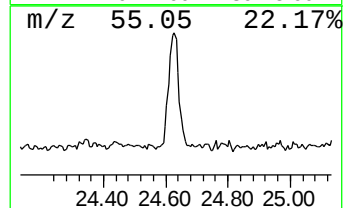
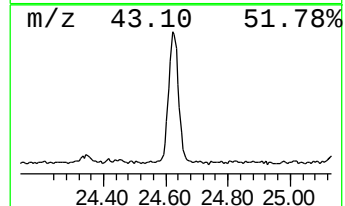
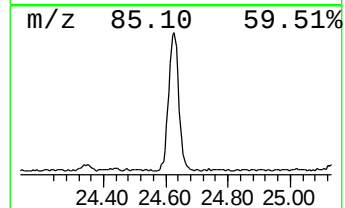
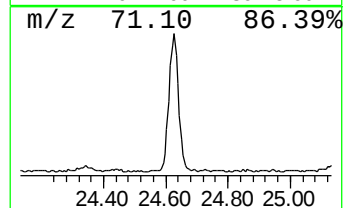
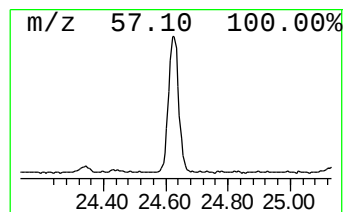
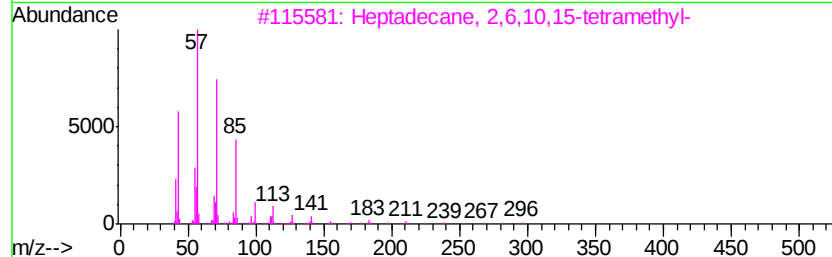
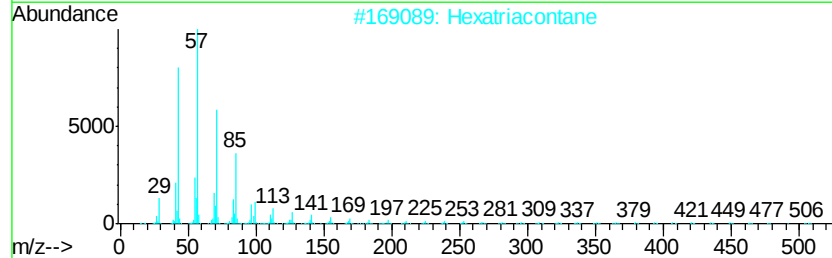
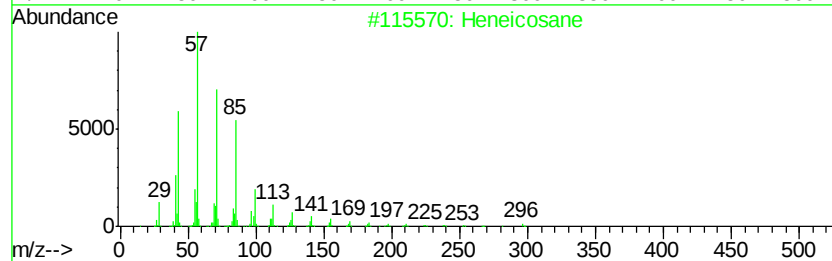
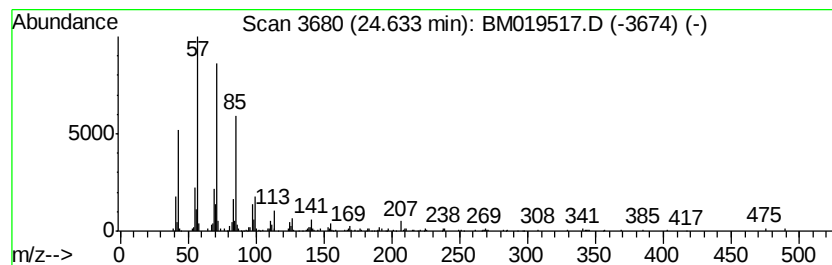
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.63	5.30 ng/ul	604620	Perylene-d12	23.42

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	83
2			Hexatriacontane	507	C36H74	000630-06-8	80
3			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	68
4			Heptacosane	380	C27H56	000593-49-7	64
5			Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	49



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM032719\
 Data File : BM019517.D
 Acq On : 27 Mar 2019 22:35
 Operator : JU/SJ
 Sample : K2069-05
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C09D7

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-	5.00	26.6	ng/ul	266249000	1	7.63	200059000	20.0
Benzene, 1,3-dich...	7.49	4.2	ng/ul	42338400	1	7.63	200059000	20.0
Benzene, 1,4-dich...	8.01	19.9	ng/ul	198698000	1	7.63	200059000	20.0
Benzene, 1-bromo-...	9.08	12.9	ng/ul	502408	2	10.38	779427	20.0
Benzene, 1,2,3-tr...	10.27	2603.7	ng/ul	101469000	2	10.38	779427	20.0
Benzene, 1,3,5-tr...	10.76	798.8	ng/ul	31131000	2	10.38	779427	20.0
Benzene, 1-chloro...	10.99	182.3	ng/ul	7105550	2	10.38	779427	20.0
Benzene, 1-chloro...	11.20	34.7	ng/ul	1352610	2	10.38	779427	20.0
unknown-01	18.39	206.8	ng/ul	12695400	4	17.00	1228110	20.0
unknown-02	18.52	7.6	ng/ul	468354	4	17.00	1228110	20.0
6-Chloro-1,2,3,4-...	18.69	4.5	ng/ul	272995	4	17.00	1228110	20.0
unknown-03	18.72	9.1	ng/ul	555951	4	17.00	1228110	20.0
(DEL) Alkane: Str...	22.72	2.4	ng/ul	277863	6	23.42	2280930	20.0
(DEL) Alkane: Str...	23.90	4.4	ng/ul	499808	6	23.42	2280930	20.0
(DEL) Alkane: Str...	24.63	5.3	ng/ul	604620	6	23.42	2280930	20.0