

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM091917\
 Data File : BM011616.D
 Acq On : 19 Sep 2017 22:15
 Operator : SJ/JU
 Sample : I5271-16
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AJ1

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:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090817MA.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	5.287	331	340	344	rBV6	69074948	184863551	77.53%	16.435%
2	7.116	645	651	661	rBV	412413	674349	0.28%	0.060%
3	7.293	674	681	696	rBV	1589262	2592524	1.09%	0.230%
4	7.498	708	716	727	rBV2	1593203	3193356	1.34%	0.284%
5	7.869	768	779	790	rBV	47825653	89441459	37.51%	7.952%
6	8.040	790	808	812	rBV5	68340326	190156783	79.75%	16.906%
7	8.351	849	861	879	rVB3	67367100	169110487	70.92%	15.035%
8	8.663	907	914	929	rBV	1161107	1966576	0.82%	0.175%
9	8.963	959	965	978	rBV2	180009	372435	0.16%	0.033%
10	9.128	986	993	1012	rBV	1997868	3472487	1.46%	0.309%
11	9.469	1043	1051	1063	rBV	2487187	4214345	1.77%	0.375%
12	9.692	1082	1089	1093	rBV	673177	1173012	0.49%	0.104%
13	9.739	1093	1097	1099	rVV	644549	1071338	0.45%	0.095%
14	9.781	1099	1104	1111	rVV	1456258	2923104	1.23%	0.260%
15	9.857	1111	1117	1122	rVV	1559701	2758694	1.16%	0.245%
16	9.910	1122	1126	1138	rVB	373755	691590	0.29%	0.061%
17	10.151	1160	1167	1183	rVB2	232447	460511	0.19%	0.041%
18	10.392	1201	1208	1215	rBV	2171289	4118918	1.73%	0.366%
19	10.457	1215	1219	1229	rVB	296657	611052	0.26%	0.054%
20	10.681	1242	1257	1262	rBV5	71202395	238441145	100.00%	21.198%
21	10.786	1269	1275	1282	rBV	2190818	3507732	1.47%	0.312%
22	11.175	1330	1341	1360	rBV	48252531	93921375	39.39%	8.350%
23	11.975	1466	1477	1483	rBV2	244410	449844	0.19%	0.040%
24	12.063	1483	1492	1506	rVB5	149184	379970	0.16%	0.034%
25	12.333	1528	1538	1547	rBV8	116227	269634	0.11%	0.024%
26	12.510	1562	1568	1584	rBV3	132733	276487	0.12%	0.025%
27	12.757	1603	1610	1612	rBV	897448	1376893	0.58%	0.122%
28	12.792	1612	1616	1624	rVV	2106896	3313551	1.39%	0.295%
29	13.028	1646	1656	1664	rBV	990417	1640790	0.69%	0.146%
30	13.398	1712	1719	1733	rVB	12222268	18785604	7.88%	1.670%
31	13.598	1748	1753	1759	rBV	335610	538812	0.23%	0.048%
32	13.663	1759	1764	1772	rVB	782071	1190262	0.50%	0.106%
33	13.998	1812	1821	1826	rBV	4696082	6539531	2.74%	0.581%
34	14.057	1826	1831	1837	rVB	1462159	2154446	0.90%	0.192%

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Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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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

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35	14.292	1865	1871	1881	rVB	5201726	7458426	3.13%	0.663%
36	14.598	1914	1923	1940	rBV2	3251488	5596790	2.35%	0.498%
37	14.727	1940	1945	1950	rVV	276894	502162	0.21%	0.045%
38	14.780	1950	1954	1971	rVB	342867	672049	0.28%	0.060%
39	14.910	1971	1976	1982	rVB	617220	920185	0.39%	0.082%
40	15.074	1998	2004	2011	rBV	1069658	1603566	0.67%	0.143%
41	15.586	2085	2091	2098	rBV	6627999	9430888	3.96%	0.838%
42	15.698	2103	2110	2119	rBV	2423243	3497714	1.47%	0.311%
43	16.180	2187	2192	2196	rVV	244527	426668	0.18%	0.038%
44	16.221	2196	2199	2208	rVV2	193376	409887	0.17%	0.036%
45	16.521	2244	2250	2252	rBV3	154516	265601	0.11%	0.024%
46	16.551	2252	2255	2264	rVB2	250487	396248	0.17%	0.035%
47	17.339	2382	2389	2394	rBV	3865649	5693241	2.39%	0.506%
48	17.433	2399	2405	2424	rVB2	7350165	10707533	4.49%	0.952%
49	19.474	2748	2752	2760	rBV	155574	244182	0.10%	0.022%
50	19.721	2787	2794	2804	rBV	9300847	12473460	5.23%	1.109%
51	20.227	2875	2880	2887	rBV3	130535	280130	0.12%	0.025%
52	21.503	3091	3097	3108	rBV	4907603	6629157	2.78%	0.589%
53	23.103	3366	3369	3381	rVB	356824	555097	0.23%	0.049%
54	23.721	3466	3474	3489	rVB2	6050571	12331348	5.17%	1.096%
55	23.874	3493	3500	3510	rVB2	2850324	5810484	2.44%	0.517%
56	24.403	3585	3590	3600	rVB	494349	1024906	0.43%	0.091%
57	25.209	3722	3727	3742	rVB	510672	1233762	0.52%	0.110%

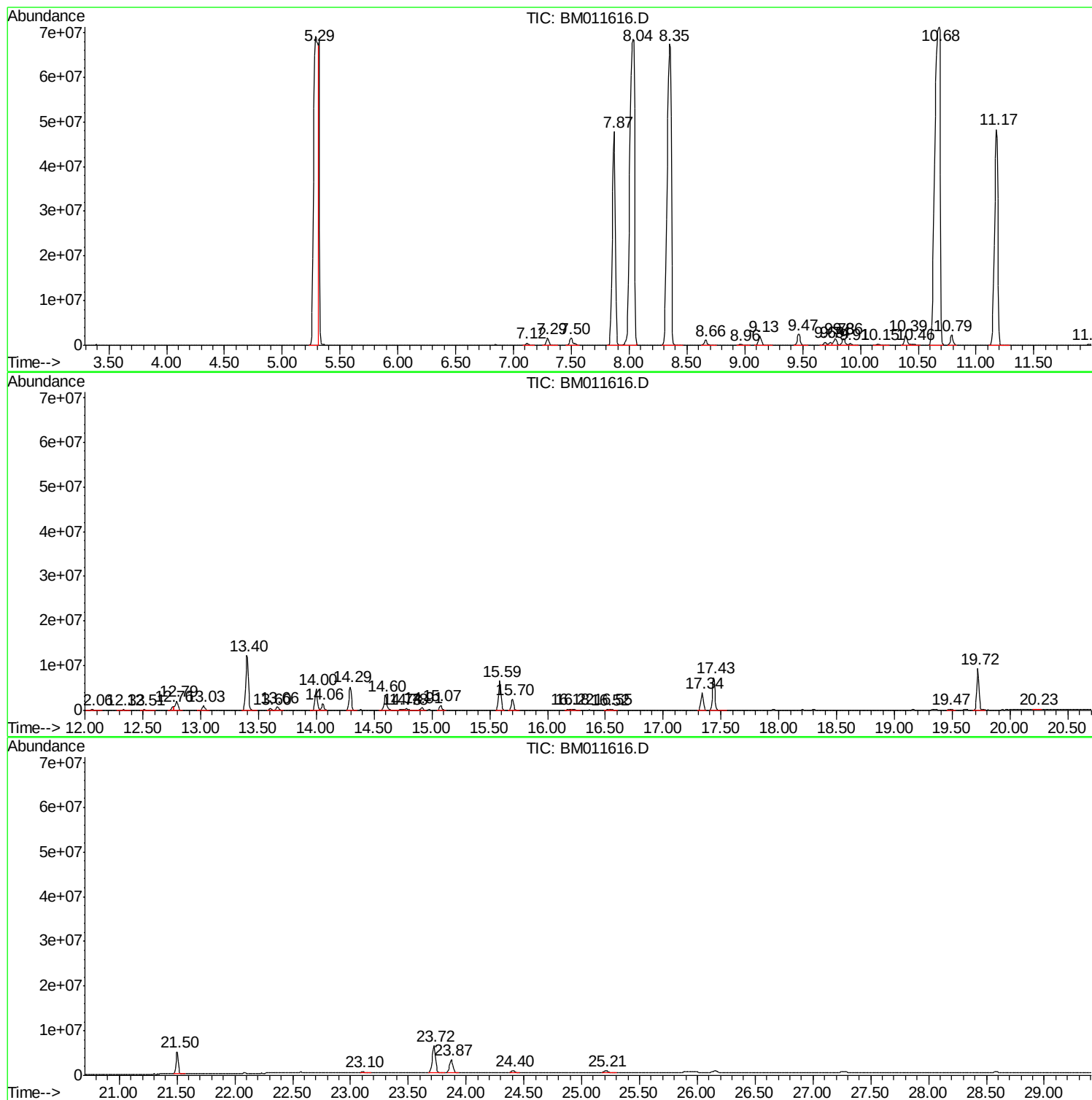
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090817MA.M
Quant Title : SVOA CALIBRATION

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



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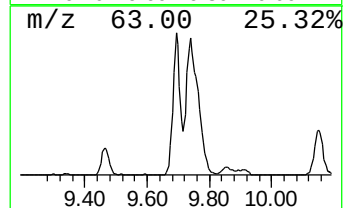
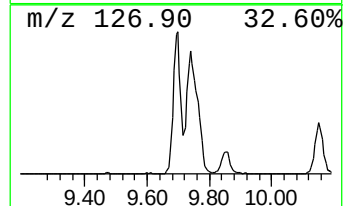
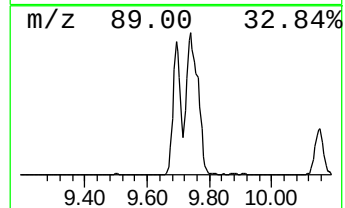
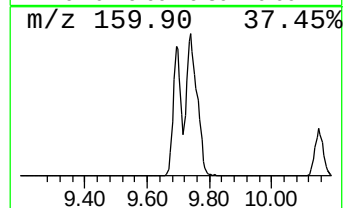
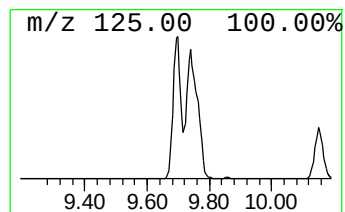
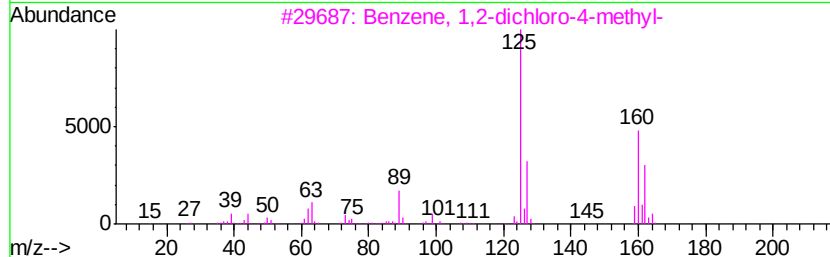
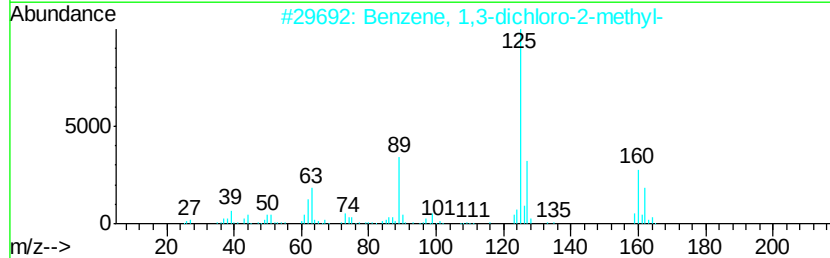
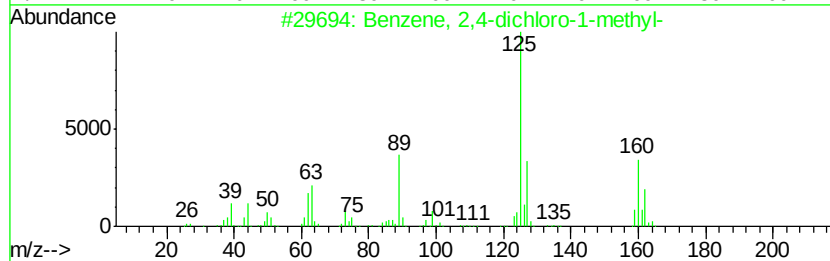
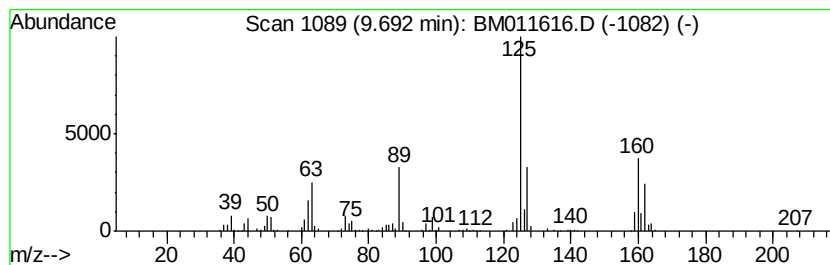
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090817MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzene, 2,4-dichloro-1-met... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.69	6.69 ng/ul	1173010	Napthalene-d8	10.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2,4-dichloro-1-methyl-	160	C7H6Cl2	000095-73-8	98
2		Benzene, 1,3-dichloro-2-methyl-	160	C7H6Cl2	000118-69-4	98
3		Benzene, 1,2-dichloro-4-methyl-	160	C7H6Cl2	000095-75-0	97
4		Benzene, 1,3-dichloro-2-methyl-	160	C7H6Cl2	000118-69-4	97
5		Benzene, 1,2-dichloro-3-methyl-	160	C7H6Cl2	032768-54-0	96



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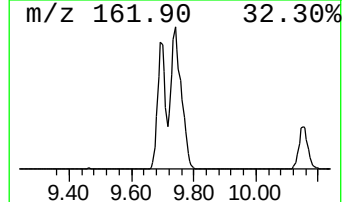
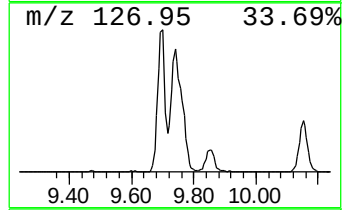
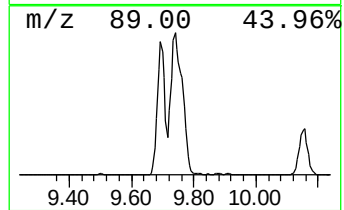
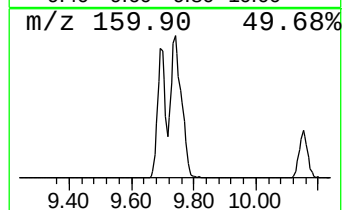
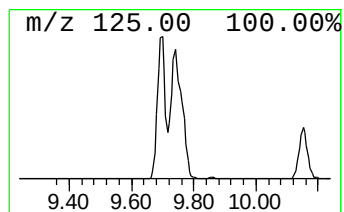
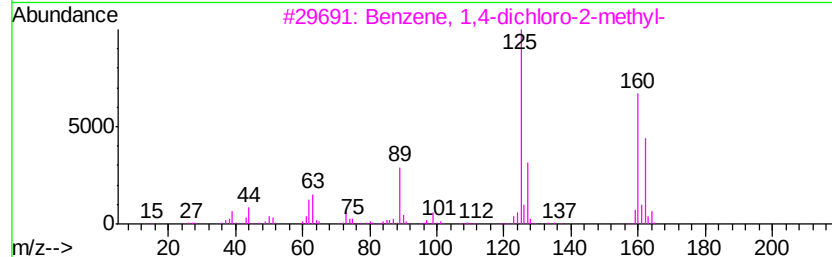
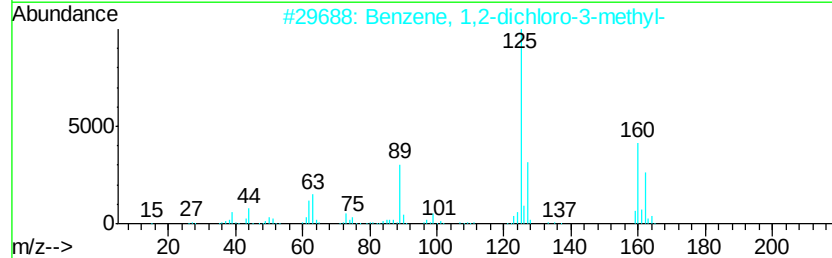
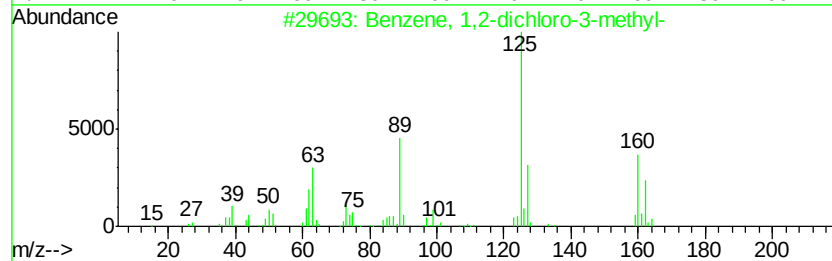
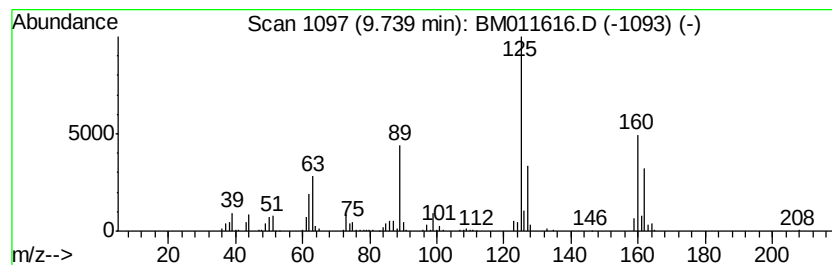
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090817MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 Benzene, 1,2-dichloro-3-met... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.74	6.11 ng/ul	1071340	Napthalene-d8	10.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2-dichloro-3-methyl-	160	C7H6Cl2	032768-54-0	97
2		Benzene, 1,2-dichloro-3-methyl-	160	C7H6Cl2	032768-54-0	96
3		Benzene, 1,4-dichloro-2-methyl-	160	C7H6Cl2	019398-61-9	95
4		Benzene, 1,2-dichloro-4-methyl-	160	C7H6Cl2	000095-75-0	95
5		Benzene, 1,2-dichloro-3-methyl-	160	C7H6Cl2	032768-54-0	95



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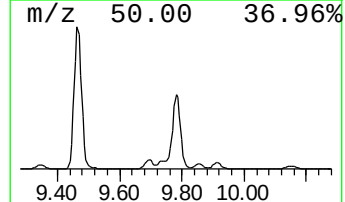
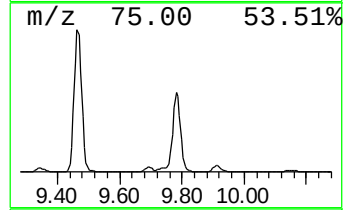
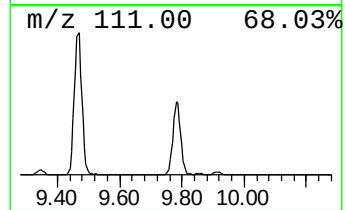
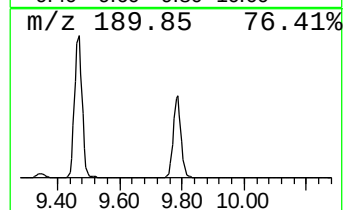
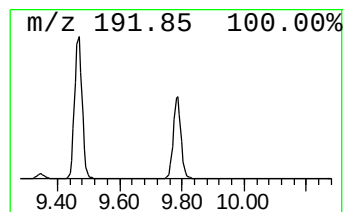
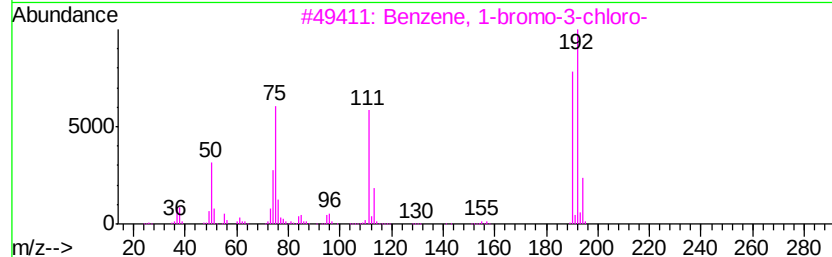
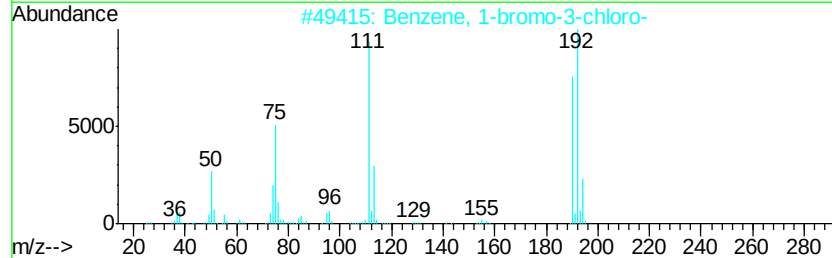
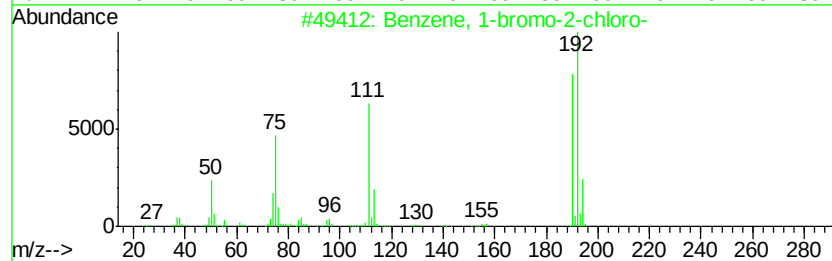
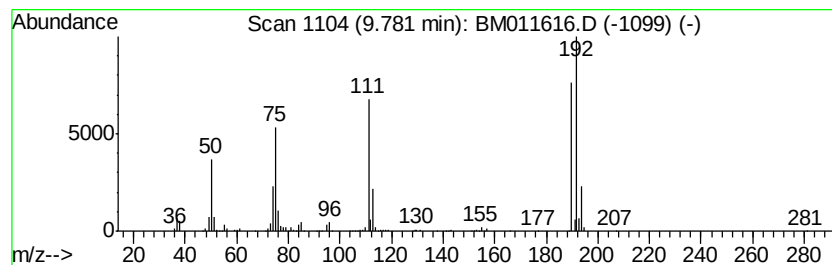
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090817MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 8 Benzene, 1-bromo-2-chloro- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.78	16.67 ng/ul	2923100	Naphthalene-d8	10.79

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	97
3		Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	97
4		Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	96
5		Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	96



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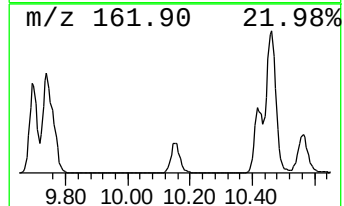
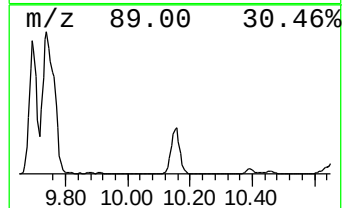
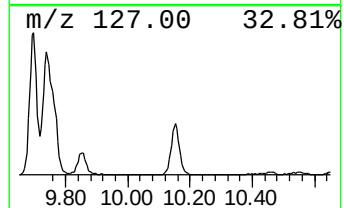
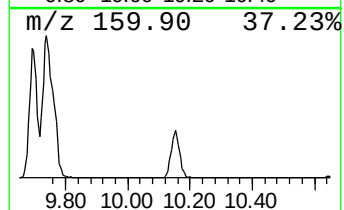
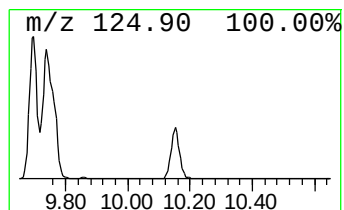
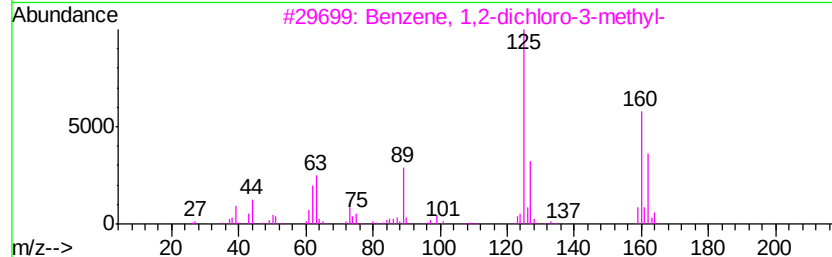
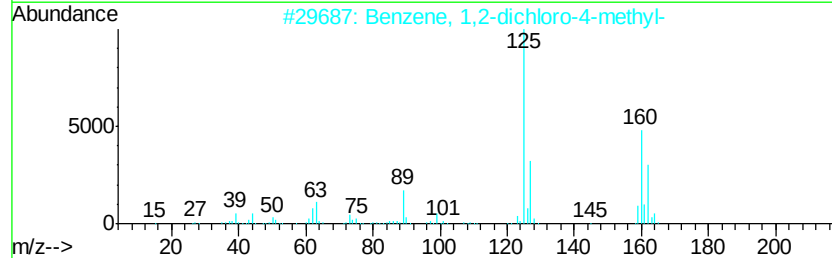
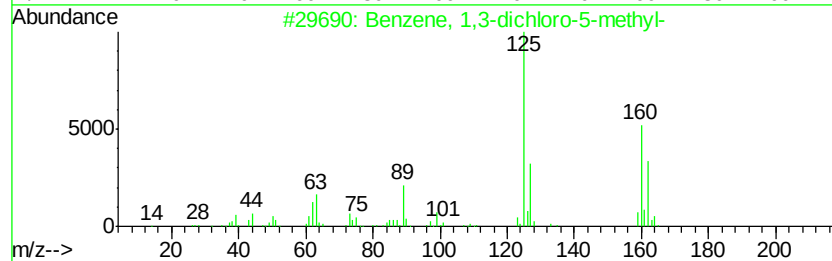
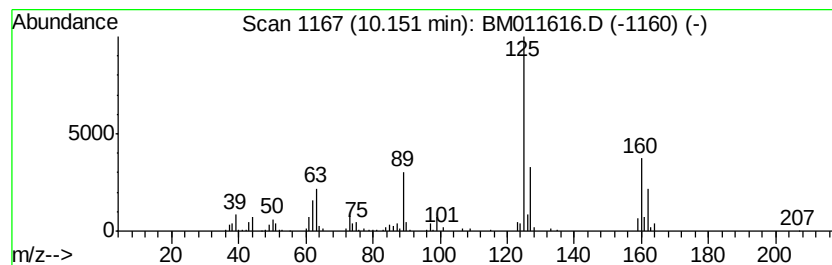
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090817MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 Benzene, 1,3-dichloro-5-met... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.15	2.63 ng/ul	460511	Napthalene-d8	10.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,3-dichloro-5-methyl-	160	C7H6Cl2	025186-47-4	97
2		Benzene, 1,2-dichloro-4-methyl-	160	C7H6Cl2	000095-75-0	97
3		Benzene, 1,2-dichloro-3-methyl-	160	C7H6Cl2	032768-54-0	96
4		Benzene, 1,2-dichloro-4-methyl-	160	C7H6Cl2	000095-75-0	96
5		Benzene, 1,2-dichloro-4-methyl-	160	C7H6Cl2	000095-75-0	95



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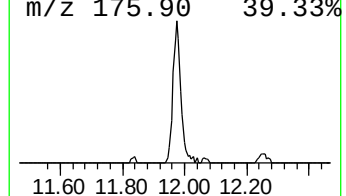
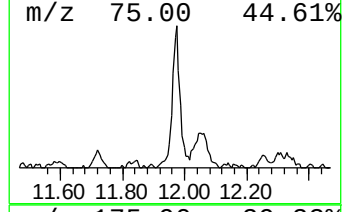
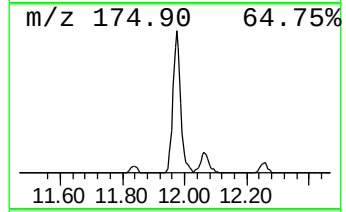
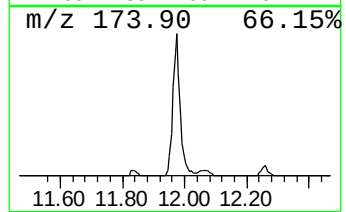
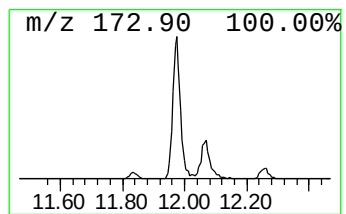
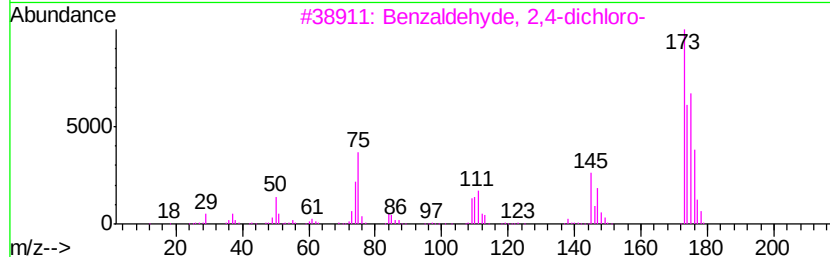
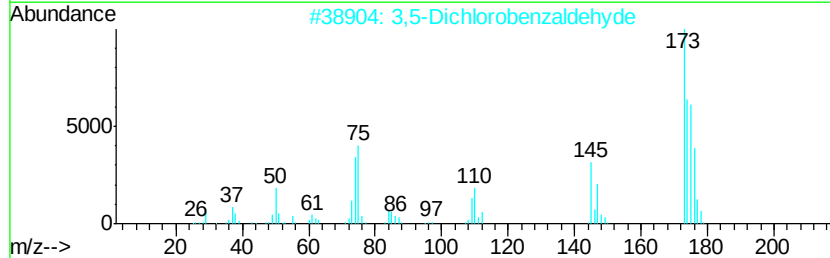
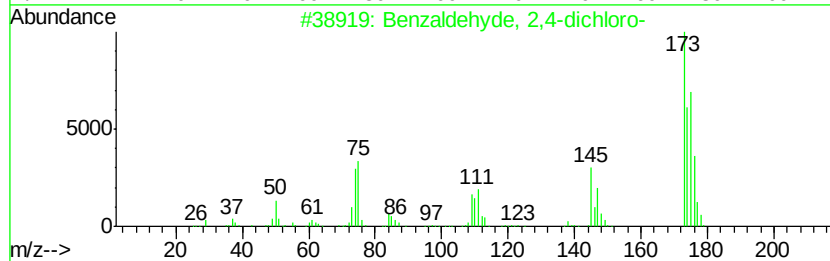
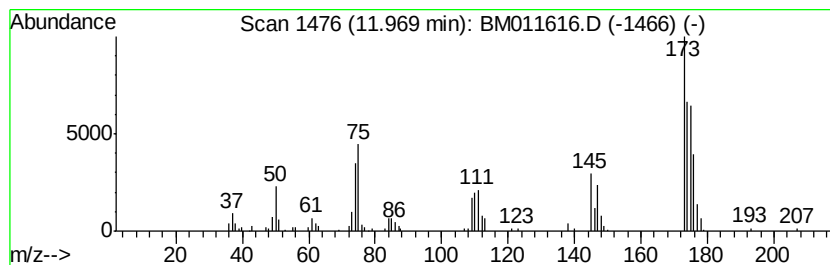
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 Benzaldehyde, 2,4-dichloro- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.97	2.56 ng/ul	449844	Napthalene-d8	10.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzaldehyde, 2,4-dichloro-	174	C7H4Cl2O	000874-42-0	99
2		3,5-Dichlorobenzaldehyde	174	C7H4Cl2O	010203-08-4	99
3		Benzaldehyde, 2,4-dichloro-	174	C7H4Cl2O	000874-42-0	98
4		3,5-Dichlorobenzaldehyde	174	C7H4Cl2O	010203-08-4	98
5		2,3-Dichlorobenzaldehyde	174	C7H4Cl2O	006334-18-5	98



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM091917\
Data File : BM011616.D
Acq On : 19 Sep 2017 22:15
Operator : SJ/JU
Sample : I5271-16
Misc :
ALS Vial : 20 Sample Multiplier: 1

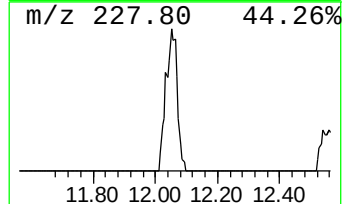
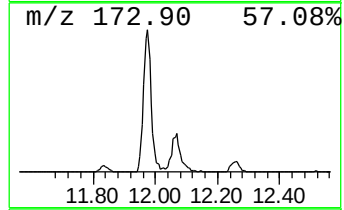
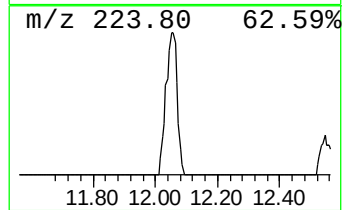
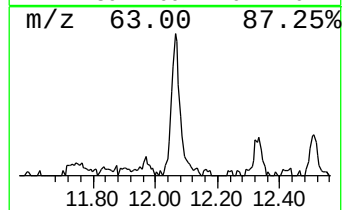
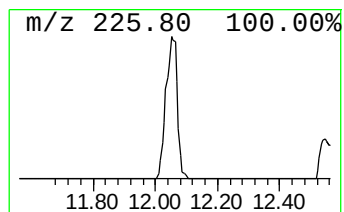
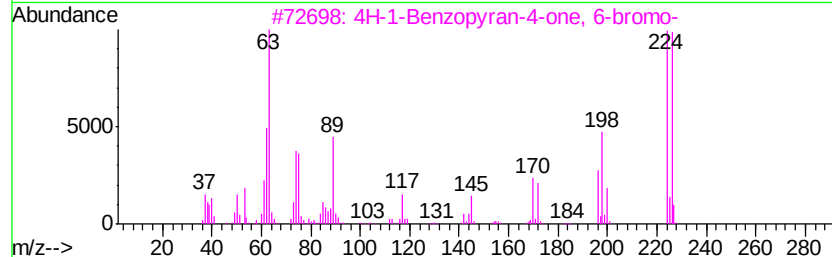
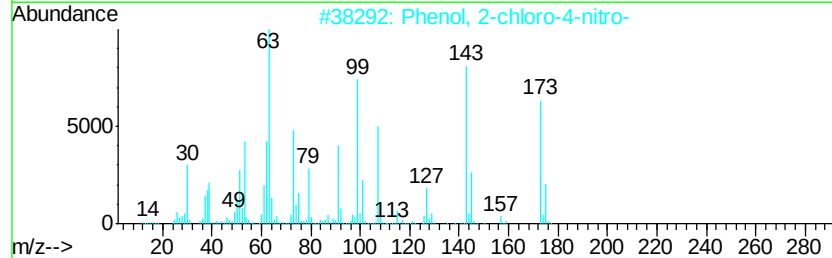
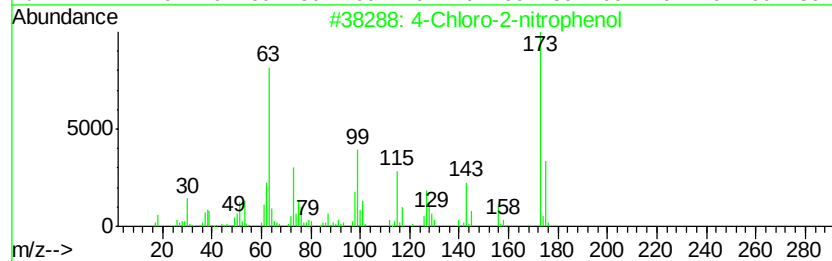
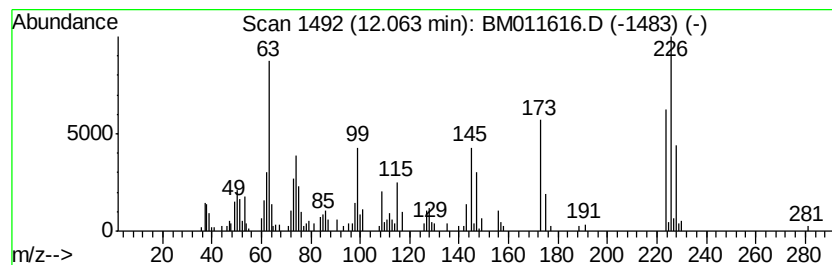
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 4-Chloro-2-nitrophenol Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.		
12.06	2.17 ng/ul	379970	Naphthalene-d8	10.79		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		4-Chloro-2-nitrophenol	173	C6H4ClNO3	000089-64-5	95
2		Phenol, 2-chloro-4-nitro-	173	C6H4ClNO3	000619-08-9	76
3		4H-1-Benzopyran-4-one, 6-bromo-	224	C9H5BrO2	051483-92-2	70
4		Phenol, 5-chloro-2-nitro-	173	C6H4ClNO3	000611-07-4	70
5		Benzene, 1-bromo-3,5-dichloro-	224	C6H3BrCl2	019752-55-7	64



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM091917\
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Acq On : 19 Sep 2017 22:15
Operator : SJ/JU
Sample : I5271-16
Misc :
ALS Vial : 20 Sample Multiplier: 1

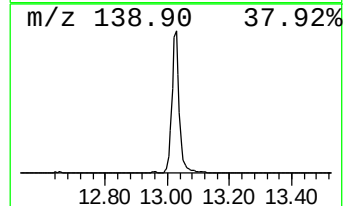
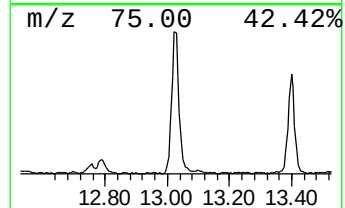
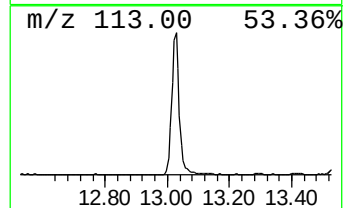
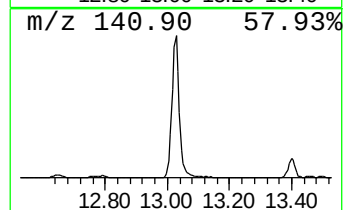
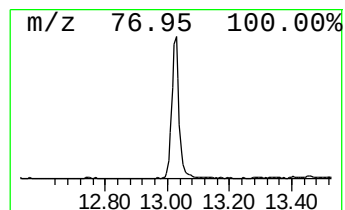
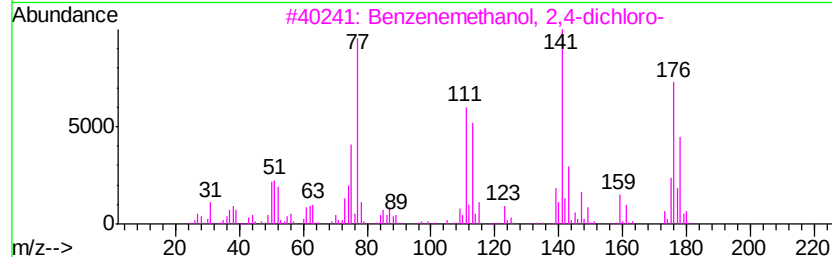
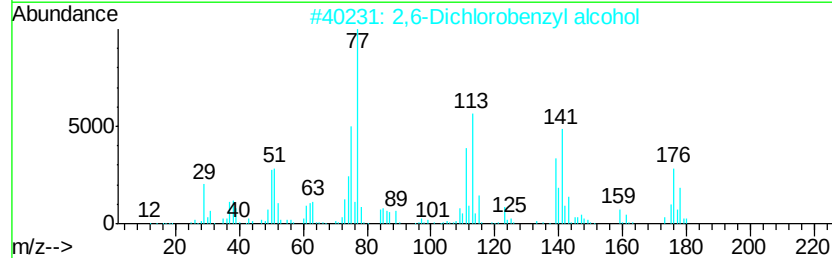
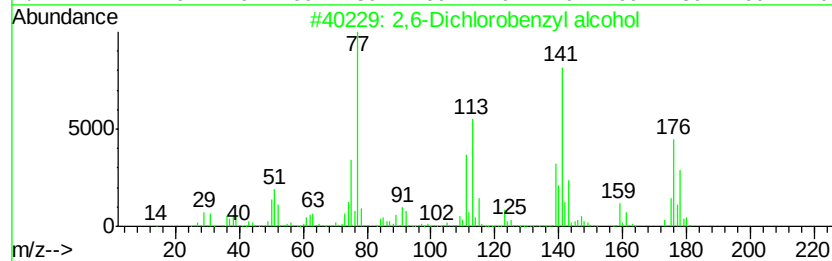
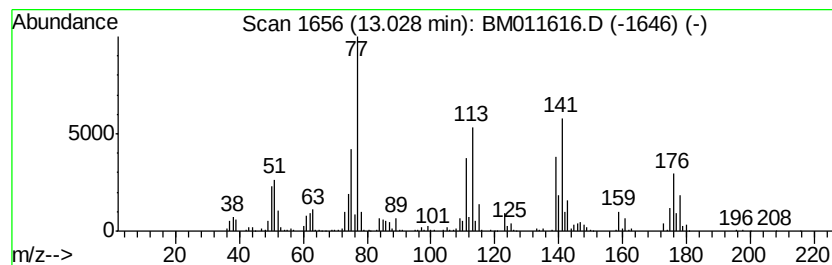
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090817MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 15 2,6-Dichlorobenzyl alcohol Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.03	5.86 ng/ul	1640790	Acenaphthene-d10	14.60
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	2,6-Dichlorobenzyl alcohol	176 C7H6Cl2O	015258-73-8	97
2	2,6-Dichlorobenzyl alcohol	176 C7H6Cl2O	015258-73-8	96
3	Benzenemethanol, 2,4-dichloro-	176 C7H6Cl2O	001777-82-8	91
4	3,5-Dichlorobenzyl alcohol	176 C7H6Cl2O	060211-57-6	86
5	3,5-Dichlorobenzyl alcohol	176 C7H6Cl2O	060211-57-6	83



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM091917\
Data File : BM011616.D
Acq On : 19 Sep 2017 22:15
Operator : SJ/JU
Sample : I5271-16
Misc :
ALS Vial : 20 Sample Multiplier: 1

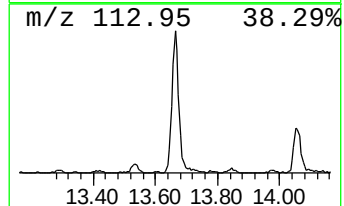
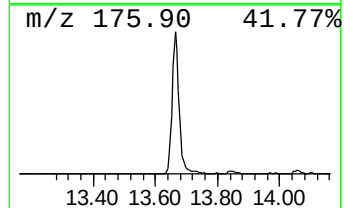
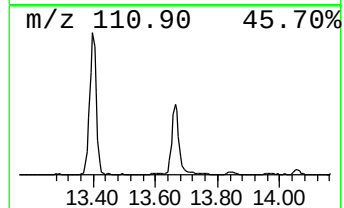
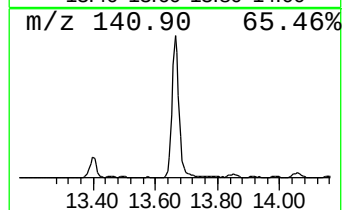
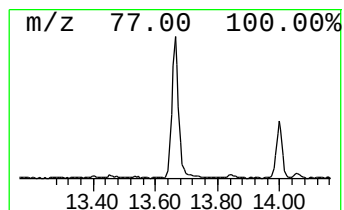
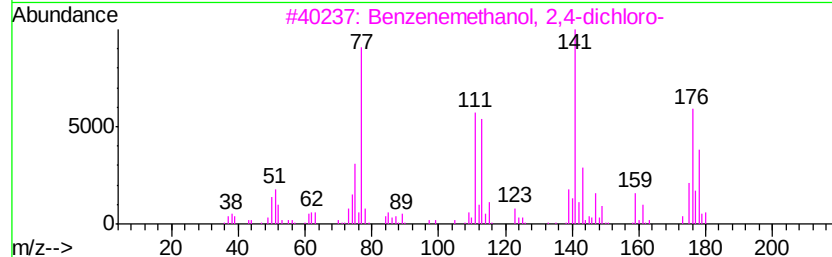
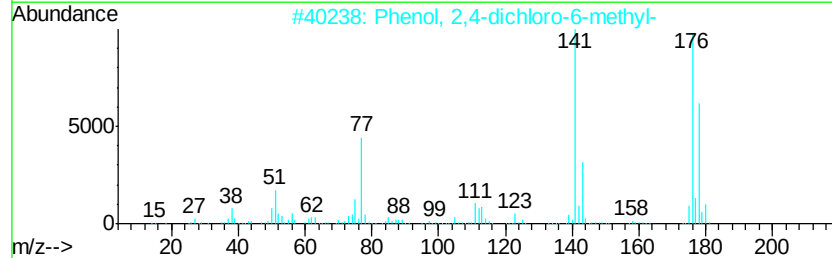
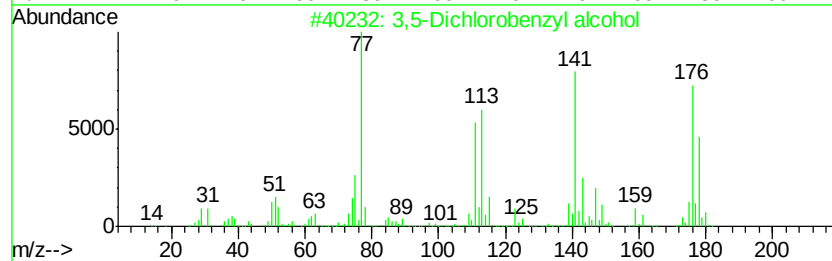
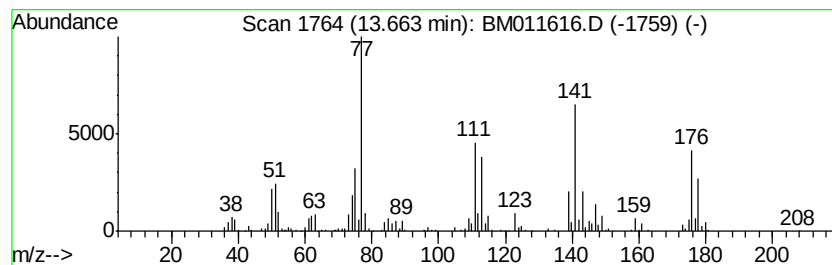
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 3,5-Dichlorobenzyl alcohol Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.66	4.25 ng/ul	1190260	Acenaphthene-d10	14.60
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		3,5-Dichlorobenzyl alcohol	176 C7H6Cl2O	060211-57-6 96
2		Phenol, 2,4-dichloro-6-methyl-	176 C7H6Cl2O	001570-65-6 93
3		Benzenemethanol, 2,4-dichloro-	176 C7H6Cl2O	001777-82-8 90
4		Phenol, 2,4-dichloro-6-methyl-	176 C7H6Cl2O	001570-65-6 86
5		Phenol, 2,4-dichloro-6-methyl-	176 C7H6Cl2O	001570-65-6 86



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM091917\
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Sample : I5271-16
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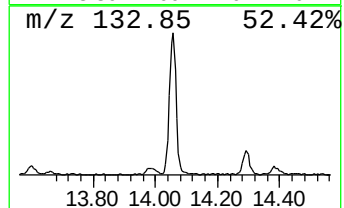
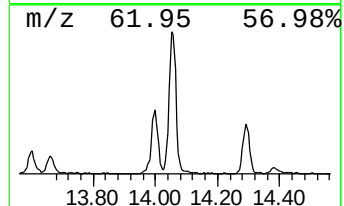
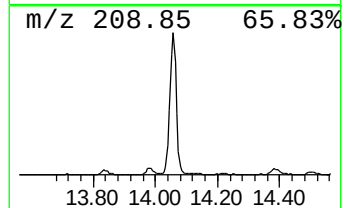
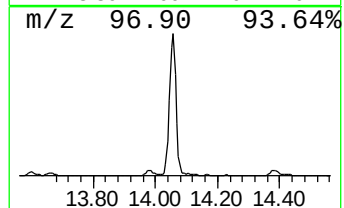
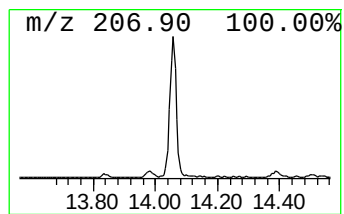
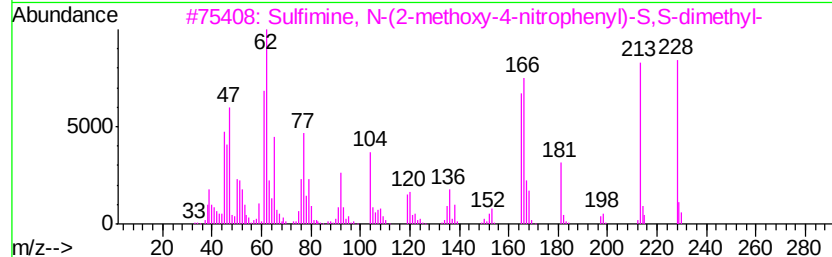
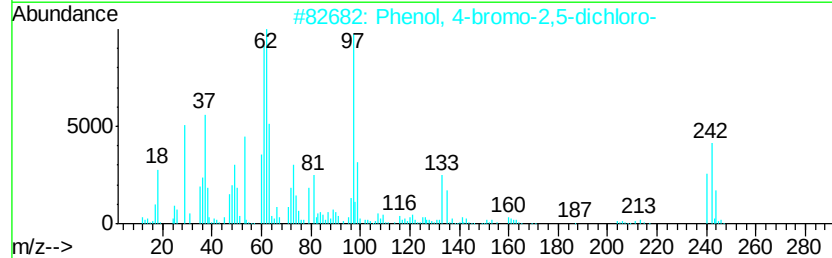
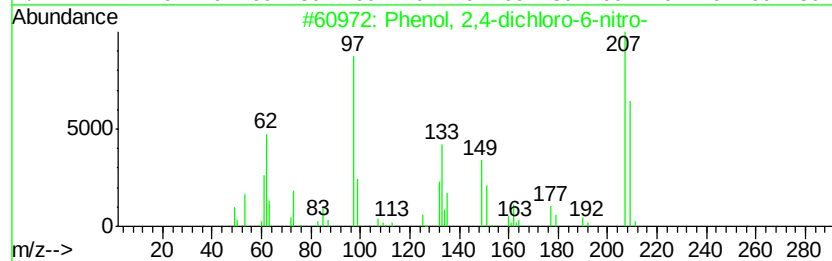
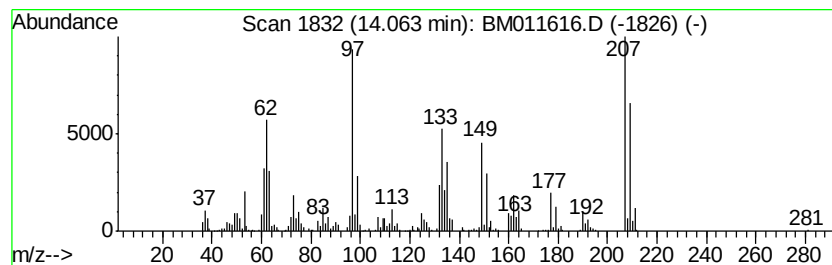
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090817MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 17 Phenol, 2,4-dichloro-6-nitro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.06	7.70 ng/ul	2154450	Acenaphthene-d10	14.60
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Phenol, 2,4-dichloro-6-nitro-	207	C6H3Cl2NO3	000609-89-2 99
2	Phenol, 4-bromo-2,5-dichloro-	240	C6H3BrCl2O	001940-42-7 32
3	Sulfimine, N-(2-methoxy-4-nitrop...	228	C9H12N2O3S	339246-70-7 14
4	2,3-Dichloropropionyl chloride	160	C3H3Cl3O	007623-13-4 13
5	Propanoic acid, 2,2-dichloro-	142	C3H4Cl2O2	000075-99-0 12



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM091917\
Data File : BM011616.D
Acq On : 19 Sep 2017 22:15
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Sample : I5271-16
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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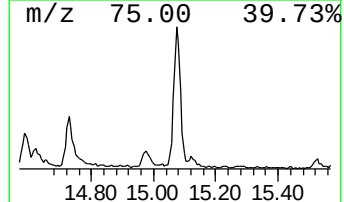
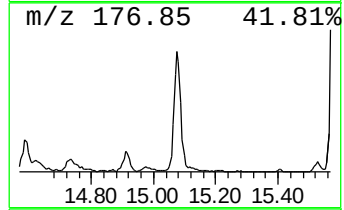
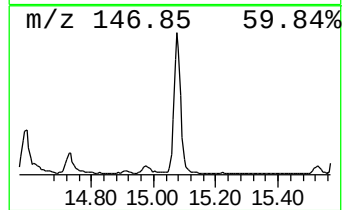
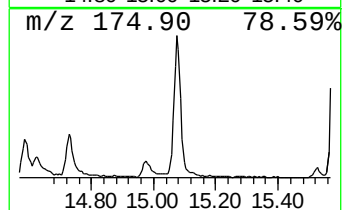
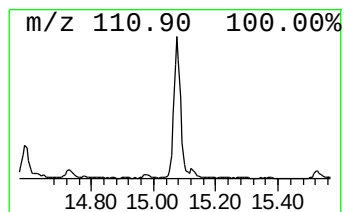
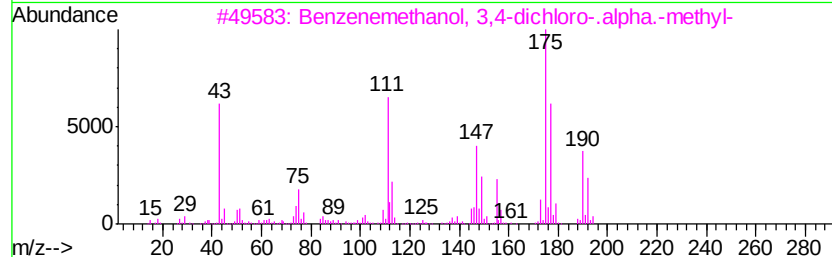
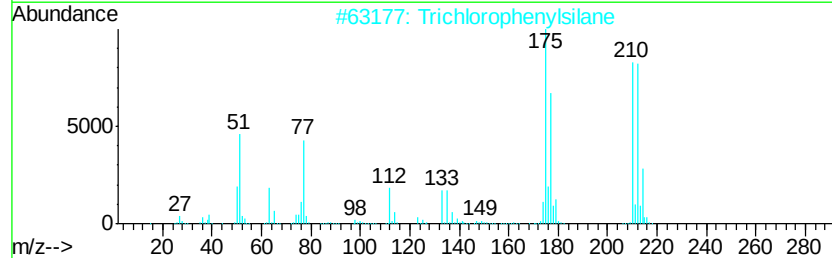
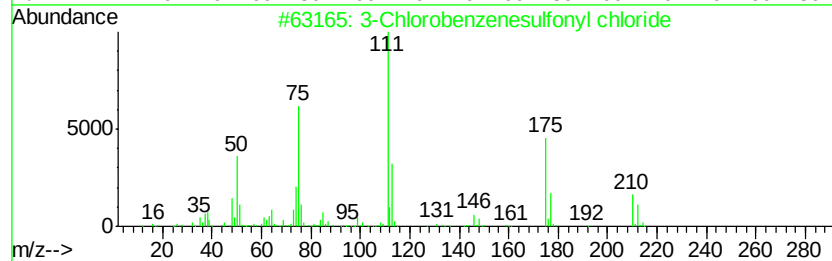
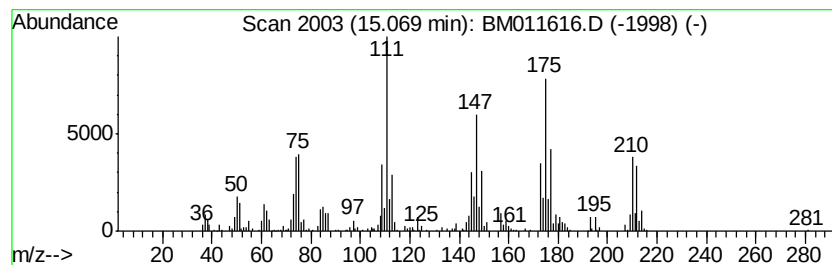
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 3-Chlorobenzenesulfonyl chl... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.07	5.73 ng/ul	1603570	Acenaphthene-d10	14.60

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chlorobenzenesulfonyl chloride	210	C6H4Cl2O2S	002888-06-4	64
2			Trichlorophenylsilane	210	C6H5Cl3Si	000098-13-5	50
3			Benzenemethanol, 3,4-dichloro-.a...	190	C8H8Cl2O	001475-11-2	47
4			Benzenemethanol, 2,3-dichloro-.a...	190	C8H8Cl2O	054798-91-3	47
5			Bicyclo[3.2.0]hept-2-en-6-one, 5...	254	C7H5BrCl2O	022374-10-3	37



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM091917\
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Acq On : 19 Sep 2017 22:15
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Misc :
ALS Vial : 20 Sample Multiplier: 1

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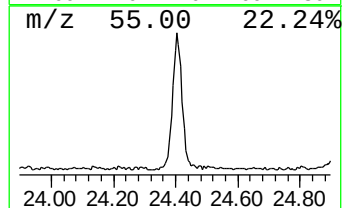
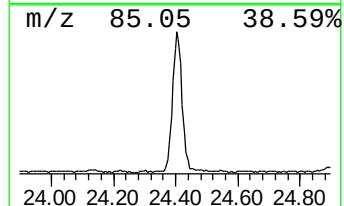
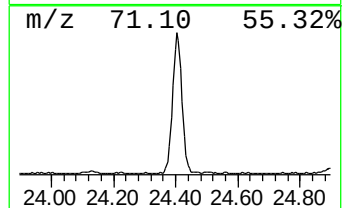
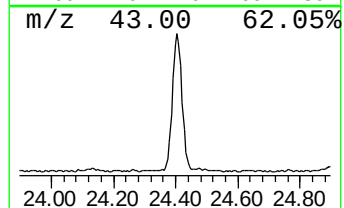
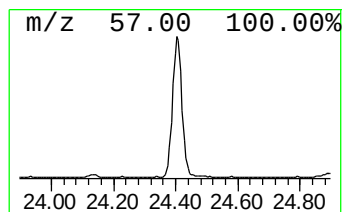
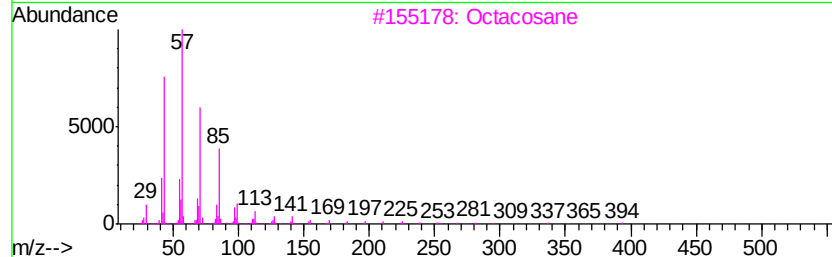
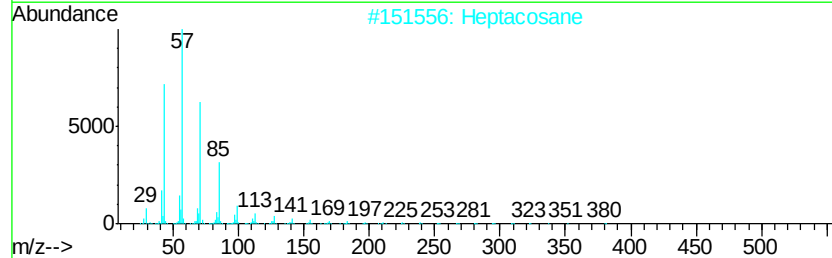
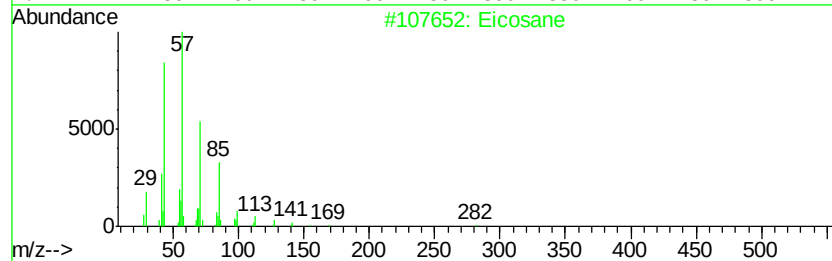
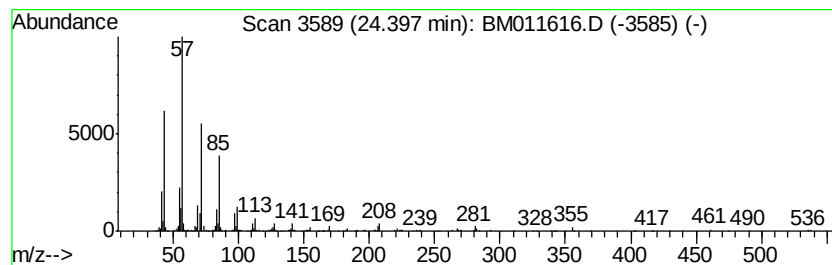
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.40	3.53 ng/ul	1024910	Perylene-d12	23.87

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	92
2			Heptacosane	380	C27H56	000593-49-7	87
3			Octacosane	394	C28H58	000630-02-4	87
4			Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	83
5			Hentriacontane	437	C31H64	000630-04-6	83



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Data File : BM011616.D
Acq On : 19 Sep 2017 22:15
Operator : SJ/JU
Sample : I5271-16
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0AJ1

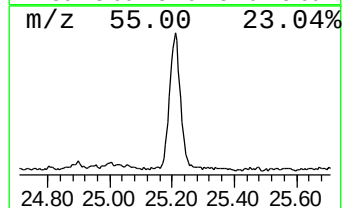
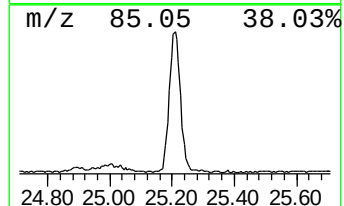
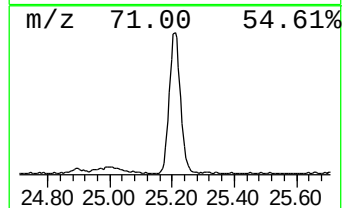
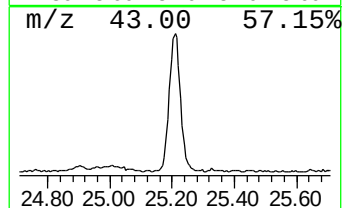
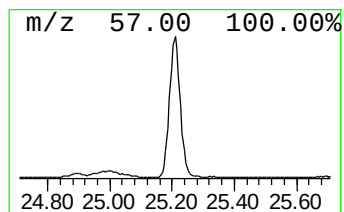
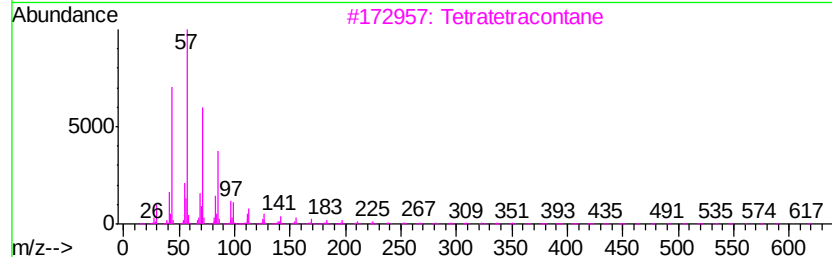
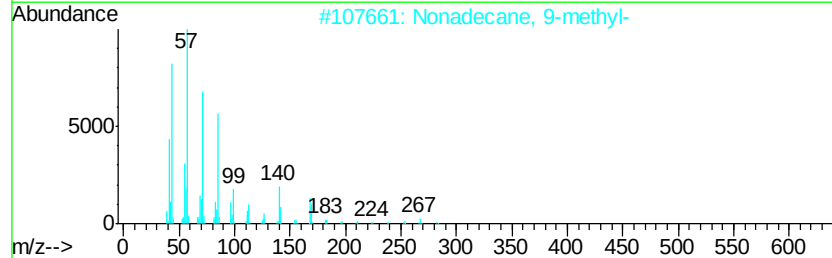
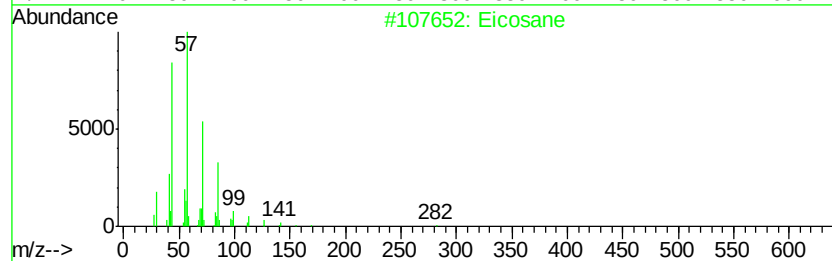
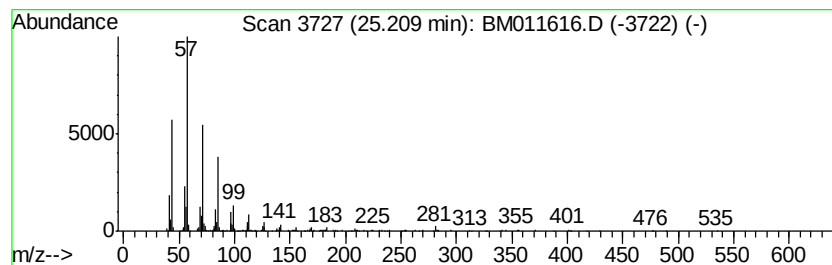
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.21	4.25 ng/ul	1233760	Perylene-d12	23.87

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	93
2			Nonadecane, 9-methyl-	282	C20H42	013287-24-6	93
3			Tetratetracontane	619	C44H90	007098-22-8	90
4			Octacosane	394	C28H58	000630-02-4	90
5			Tetracosane	338	C24H50	000646-31-1	90



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM091917\
 Data File : BM011616.D
 Acq On : 19 Sep 2017 22:15
 Operator : SJ/JU
 Sample : I5271-16
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AJ1

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090817MA.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, 2,4-dich...	9.69	6.7	ng/ul	1173010	2	10.79	3507730	20.0
Benzene, 1,2-dich...	9.74	6.1	ng/ul	1071340	2	10.79	3507730	20.0
Benzene, 1-bromo-...	9.78	16.7	ng/ul	2923100	2	10.79	3507730	20.0
Benzene, 1,3-dich...	10.15	2.6	ng/ul	460511	2	10.79	3507730	20.0
Benzaldehyde, 2,4...	11.97	2.6	ng/ul	449844	2	10.79	3507730	20.0
4-Chloro-2-nitrop...	12.06	2.2	ng/ul	379970	2	10.79	3507730	20.0
2,6-Dichlorobenz...	13.03	5.9	ng/ul	1640790	3	14.60	5596790	20.0
3,5-Dichlorobenz...	13.66	4.3	ng/ul	1190260	3	14.60	5596790	20.0
Phenol, 2,4-dichl...	14.06	7.7	ng/ul	2154450	3	14.60	5596790	20.0
3-Chlorobenzenesu...	15.07	5.7	ng/ul	1603570	3	14.60	5596790	20.0
(DEL) Alkane: Str...	24.40	3.5	ng/ul	1024910	6	23.87	5810480	20.0
(DEL) Alkane: Str...	25.21	4.3	ng/ul	1233760	6	23.87	5810480	20.0