

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM091718\
 Data File : BM016764.D
 Acq On : 17 Sep 2018 11:30
 Operator : SJ/JU
 Sample : J4874-13
 Misc : PE
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A3Z82

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-BM091018MA.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	3.022	3	6	16	rVB	2830582	3155665	13.33%	1.152%
2	3.269	44	48	54	rBV	81762	96716	0.41%	0.035%
3	3.810	137	140	146	rVB	37355	49191	0.21%	0.018%
4	3.899	150	155	160	rBV2	20976	32759	0.14%	0.012%
5	4.887	318	323	330	rBV	60875	88647	0.37%	0.032%
6	5.204	372	377	381	rBV2	18031	28287	0.12%	0.010%
7	5.881	488	492	497	rBV	17983	25877	0.11%	0.009%
8	6.928	658	670	680	rBV2	2273573	5230523	22.09%	1.909%
9	7.075	688	695	703	rBV	1025251	1489857	6.29%	0.544%
10	7.163	706	710	716	rVB	19640	30606	0.13%	0.011%
11	7.269	721	728	730	rBV	1250252	1995054	8.43%	0.728%
12	7.298	730	733	741	rVV	2342263	3398533	14.35%	1.240%
13	7.363	741	744	751	rVB5	19405	32903	0.14%	0.012%
14	7.622	779	788	797	rVB	2032305	3140805	13.26%	1.146%
15	7.734	797	807	809	rBV	924602	1446662	6.11%	0.528%
16	7.769	809	813	822	rVB	2722444	4132772	17.45%	1.508%
17	8.081	859	866	874	rVB	2709258	4195272	17.72%	1.531%
18	8.263	889	897	902	rBV	615122	1539953	6.50%	0.562%
19	8.316	902	906	912	rVB	380854	597406	2.52%	0.218%
20	8.434	919	926	936	rBV	1429496	2198830	9.29%	0.802%
21	8.804	982	989	996	rBV	2097481	3328834	14.06%	1.215%
22	8.881	996	1002	1010	rVV	1090540	1760429	7.43%	0.642%
23	9.598	1116	1124	1132	rBV	689812	1116790	4.72%	0.408%
24	9.704	1137	1142	1145	rVV2	19925	31806	0.13%	0.012%
25	9.745	1145	1149	1154	rVB	34279	52667	0.22%	0.019%
26	10.157	1208	1219	1231	rBV2	2522333	6289791	26.56%	2.295%
27	10.375	1249	1256	1265	rVB	2614088	4290231	18.12%	1.566%
28	10.516	1269	1280	1283	rBV	1203286	2059610	8.70%	0.752%
29	10.563	1283	1288	1296	rVV	3620540	6056410	25.58%	2.210%
30	10.651	1296	1303	1316	rVB	1416333	2382290	10.06%	0.869%
31	10.851	1329	1337	1343	rBV	2523973	4048875	17.10%	1.478%
32	11.228	1394	1401	1407	rBV2	25167	40295	0.17%	0.015%
33	11.551	1446	1456	1464	rBV	1696642	2825958	11.94%	1.031%
34	12.045	1535	1540	1545	rBV3	18080	30109	0.13%	0.011%

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35	12.104	1545	1550	1556	rVB	24277	41864	0.18%	0.015%
36	12.551	1618	1626	1633	rBV	4396049	6948930	29.35%	2.536%
37	12.775	1657	1664	1667	rBV	766212	1398854	5.91%	0.511%
38	12.804	1667	1669	1686	rVB	875020	1326521	5.60%	0.484%
39	13.657	1810	1814	1818	rBV2	17543	24828	0.10%	0.009%
40	13.786	1830	1836	1845	rBV	2271229	3128480	13.21%	1.142%
41	13.951	1857	1864	1872	rBV	2521128	3403294	14.37%	1.242%
42	14.063	1876	1883	1891	rVB	2812300	4131350	17.45%	1.508%
43	14.369	1929	1935	1941	rBV2	1570598	2491111	10.52%	0.909%
44	14.433	1941	1946	1955	rVB	5772268	8396924	35.46%	3.064%
45	14.569	1962	1969	1986	rBV	998308	1482265	6.26%	0.541%
46	14.739	1991	1998	2008	rVV	2182717	3060487	12.93%	1.117%
47	15.210	2071	2078	2087	rBV	4653734	6065796	25.62%	2.214%
48	15.368	2098	2105	2109	rBV	3553916	5016339	21.19%	1.831%
49	15.421	2109	2114	2120	rVV	5807690	8338267	35.22%	3.043%
50	15.480	2120	2124	2137	rVV	728448	998480	4.22%	0.364%
51	15.610	2141	2146	2152	rVB2	36327	50869	0.21%	0.019%
52	15.674	2152	2157	2163	rBV	114214	156044	0.66%	0.057%
53	16.280	2255	2260	2264	rVB2	38200	48077	0.20%	0.018%
54	16.327	2264	2268	2275	rBV	36317	51348	0.22%	0.019%
55	16.421	2277	2284	2291	rBV	4478229	6135238	25.91%	2.239%
56	16.763	2336	2342	2354	rBV	2765053	3976295	16.79%	1.451%
57	16.857	2354	2358	2363	rVB5	15054	34390	0.15%	0.013%
58	17.115	2396	2402	2410	rBV2	1918703	2817477	11.90%	1.028%
59	17.215	2413	2419	2422	rBV	3653987	5395490	22.79%	1.969%
60	17.251	2422	2425	2435	rVB	6009497	8063397	34.05%	2.943%
61	17.739	2503	2508	2512	rBV2	137734	171377	0.72%	0.063%
62	18.533	2639	2643	2649	rBV5	16658	26585	0.11%	0.010%
63	19.145	2743	2747	2755	rVB	38460	60865	0.26%	0.022%
64	19.421	2786	2794	2798	rBV2	87624	154170	0.65%	0.056%
65	19.515	2803	2810	2812	rBV	4400684	6549377	27.66%	2.390%
66	19.545	2812	2815	2823	rVB	8677943	10722000	45.28%	3.913%
67	20.221	2928	2930	2936	rVB5	28987	31355	0.13%	0.011%
68	20.462	2965	2971	2985	rBV	6092820	6852131	28.94%	2.501%
69	20.609	2991	2996	3002	rBV	7581772	8978111	37.92%	3.277%
70	21.245	3099	3104	3108	rBV	8299402	8630519	36.45%	3.150%
71	21.298	3108	3113	3118	rVV2	7501909	12772881	53.94%	4.661%

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72	21.345	3118	3121	3130	rVB	7652492	9275137	39.17%	3.385%
73	22.880	3375	3382	3386	rBV	5483038	9484465	40.06%	3.461%
74	22.927	3386	3390	3400	rVB	4615308	7431664	31.39%	2.712%
75	23.409	3465	3472	3476	rBV	3460858	6774498	28.61%	2.472%
76	23.456	3476	3480	3489	rVV	4699878	8634224	36.47%	3.151%
77	23.550	3490	3496	3508	rVB	1888365	3580868	15.12%	1.307%
78	25.821	3868	3882	3901	rBV2	7240779	23677768	100.00%	8.641%

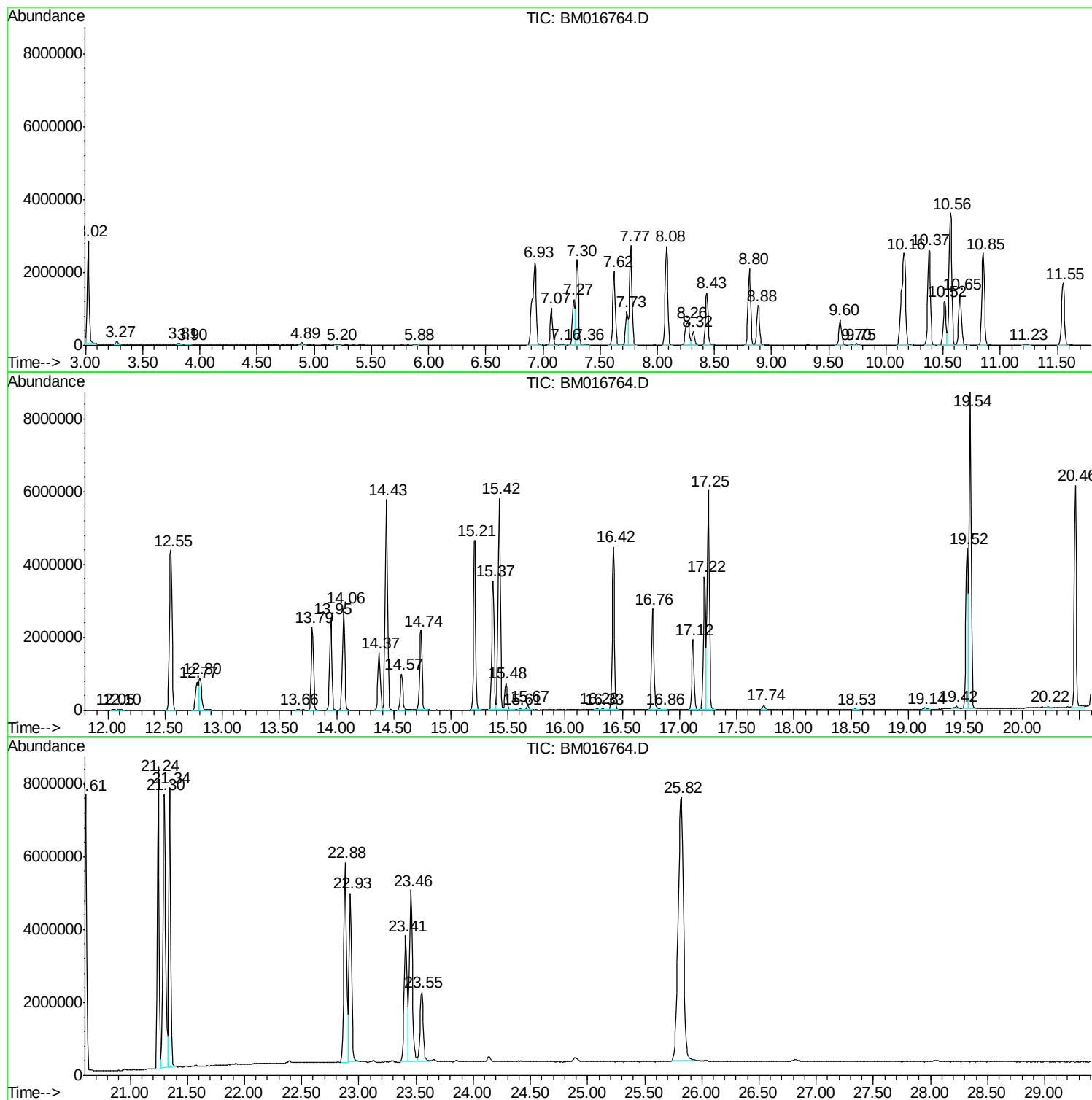
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Instrument :
BNA_M
ClientSampled :
A3Z82

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
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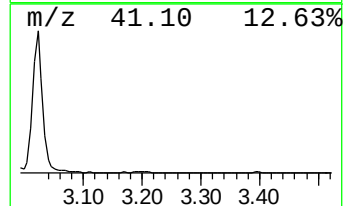
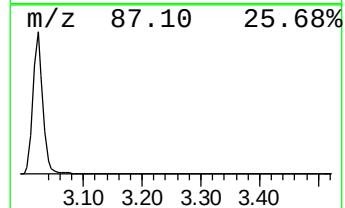
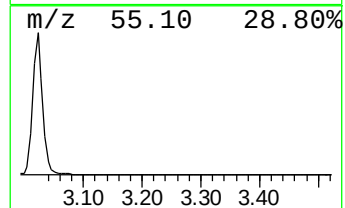
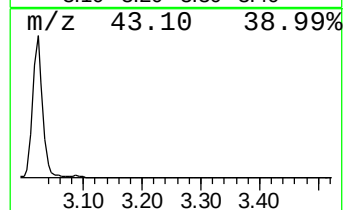
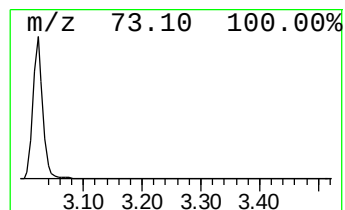
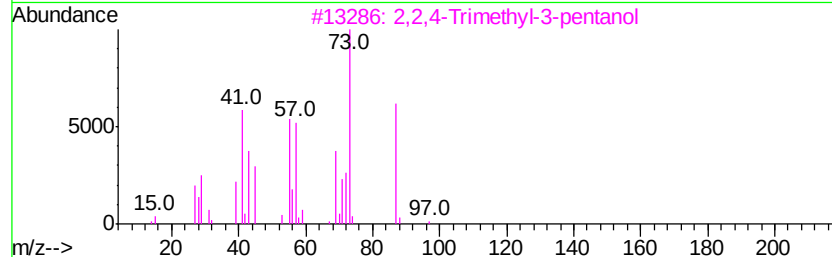
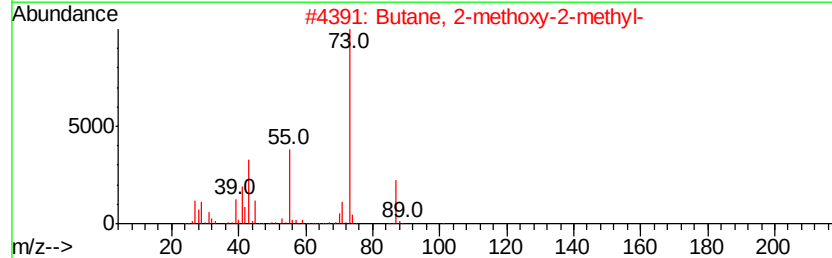
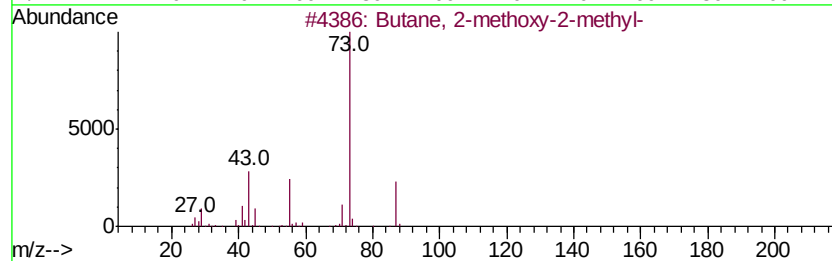
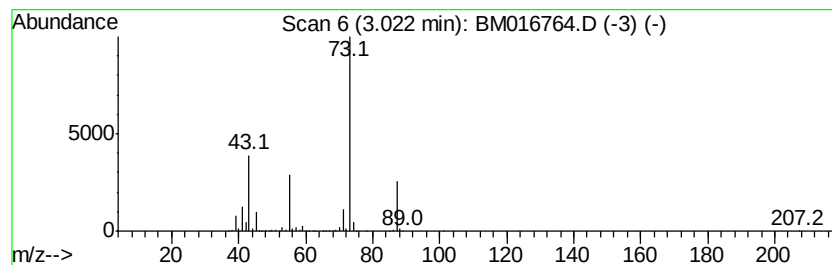
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.02	43.63 ng/ul	3155670	1,4-Dichlorobenzene-d4	7.73

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	56
3		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
4		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	27
5		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25



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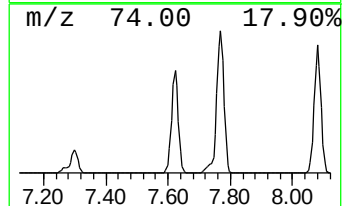
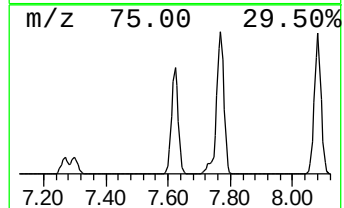
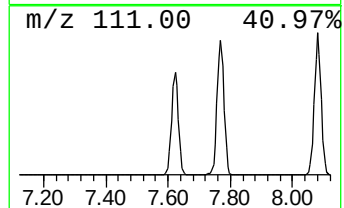
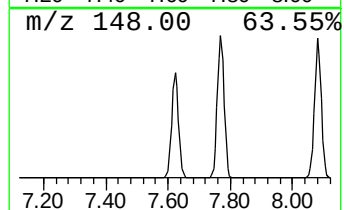
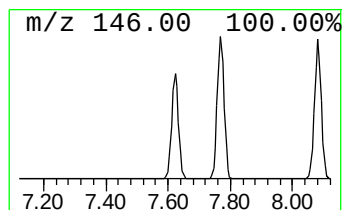
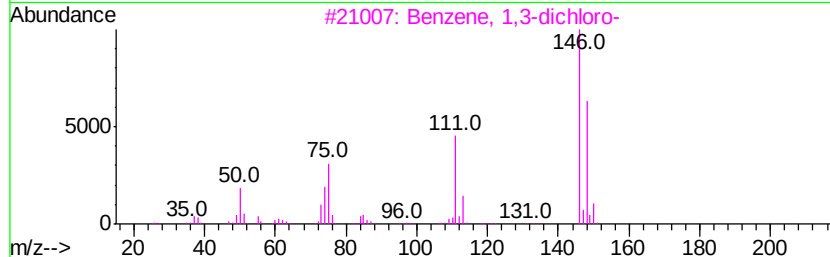
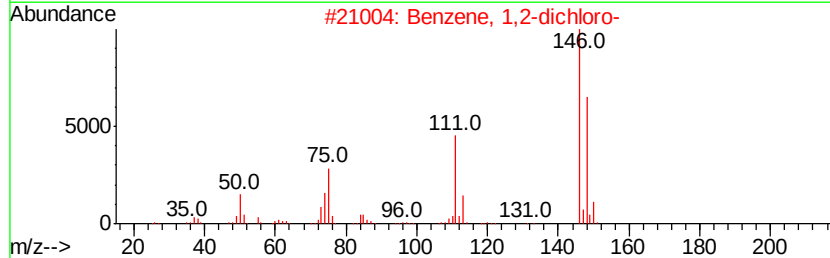
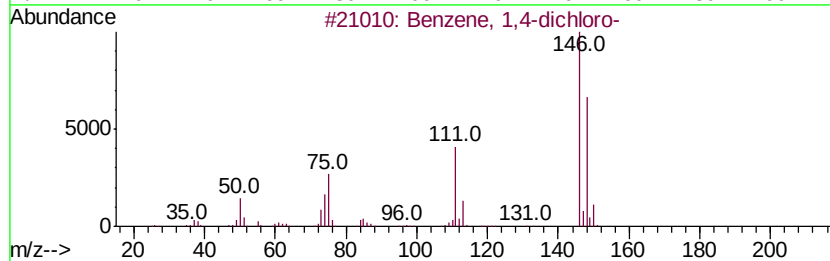
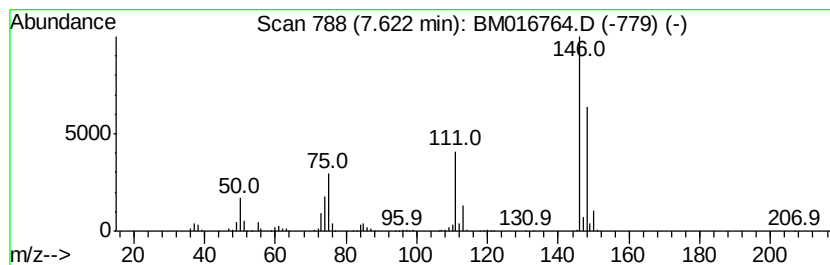
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzene, 1,4-dichloro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.62	43.42 ng/ul	3140810	1,4-Dichlorobenzene-d4	7.73

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	98
2		Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	97
3		Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	97
4		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96
5		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96



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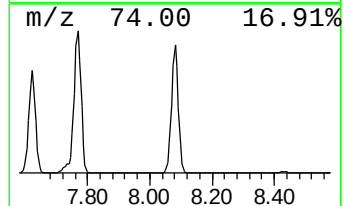
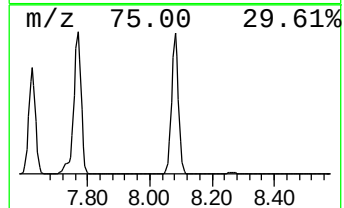
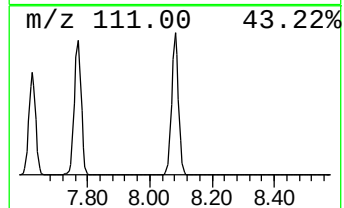
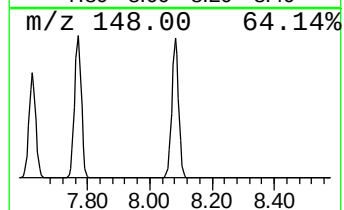
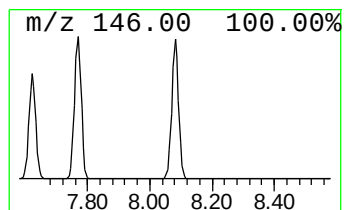
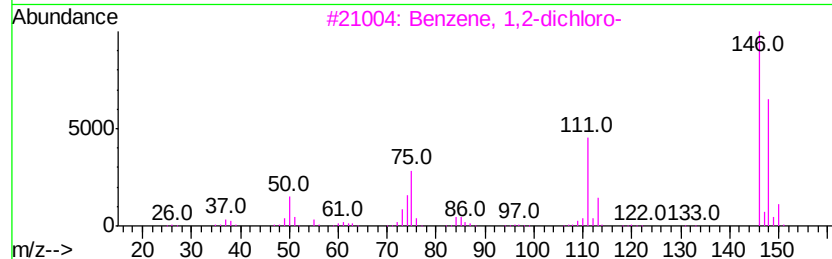
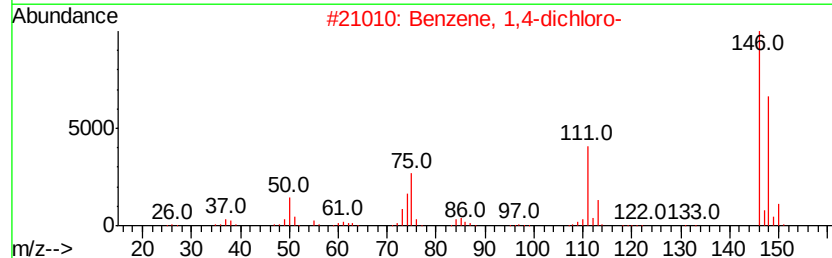
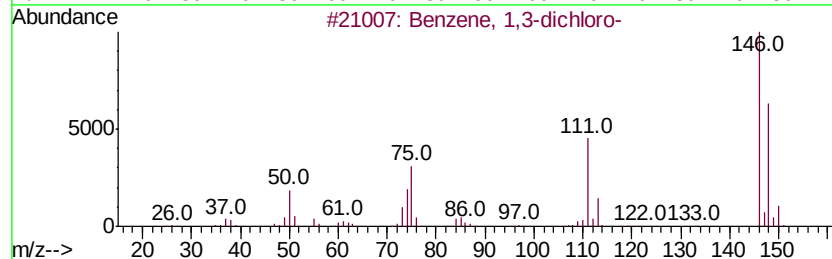
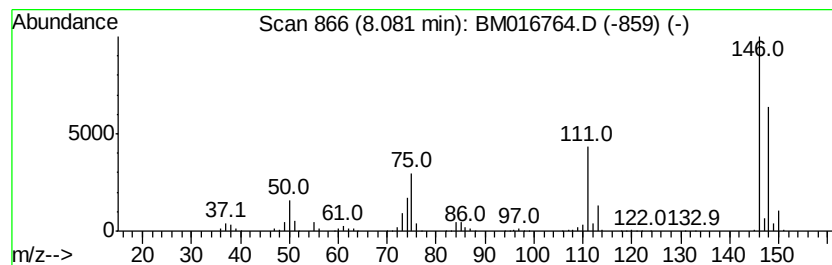
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.08	58.00 ng/ul	4195270	1,4-Dichlorobenzene-d4	7.73

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	98
2		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	98
3		Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	98
4		Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	97
5		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	97



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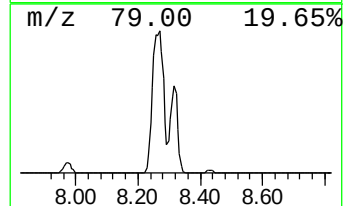
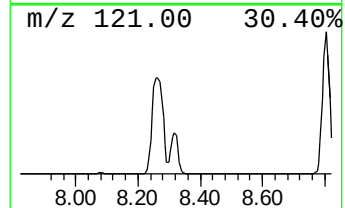
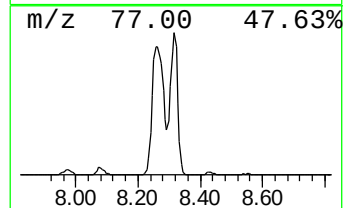
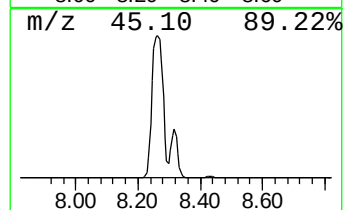
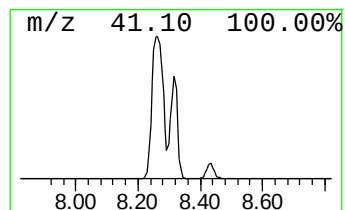
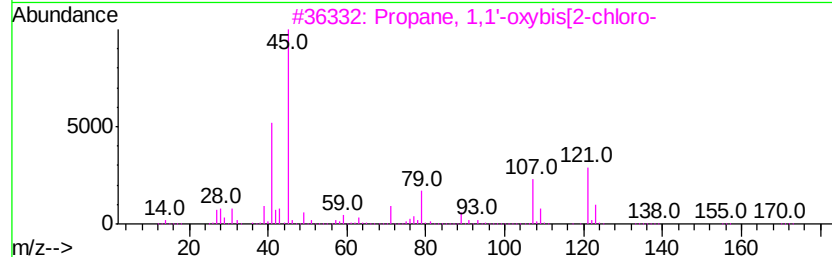
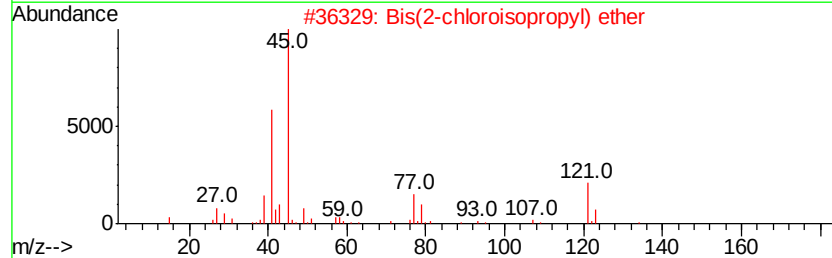
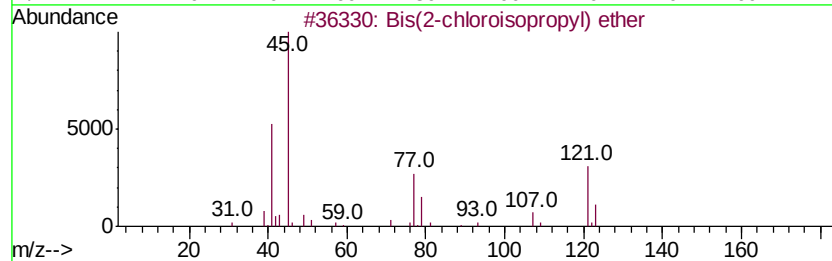
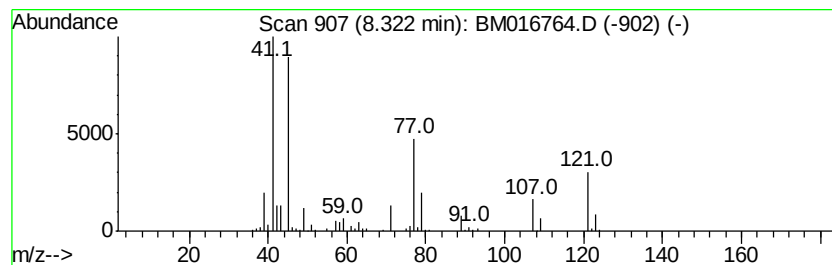
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Propane, 1,1'-oxybis[2-chloro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.32	8.26 ng/ul	597406	1,4-Dichlorobenzene-d4	7.73

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Bis(2-chloroisopropyl) ether	170	C6H12Cl2O	039638-32-9	59
2		Bis(2-chloroisopropyl) ether	170	C6H12Cl2O	039638-32-9	53
3		Propane, 1,1'-oxybis[2-chloro-	170	C6H12Cl2O	054460-96-7	53
4		Propane, 1,1'-oxybis[3-chloro-	170	C6H12Cl2O	000629-36-7	50
5		Propane, 2,2'-oxybis[1-chloro-	170	C6H12Cl2O	000108-60-1	40



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM091718\
Data File : BM016764.D
Acq On : 17 Sep 2018 11:30
Operator : SJ/JU
Sample : J4874-13
Misc : PE
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
A3Z82

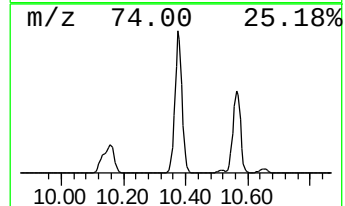
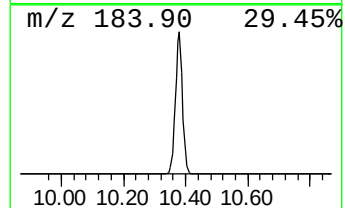
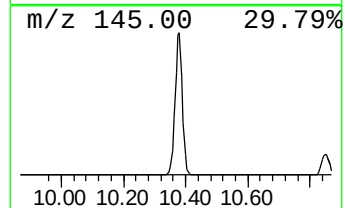
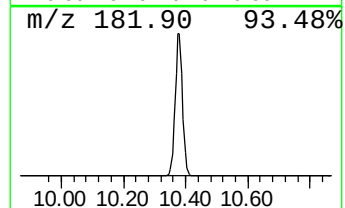
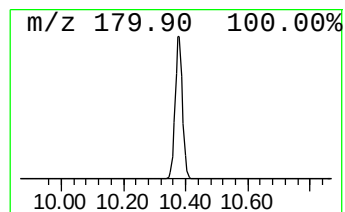
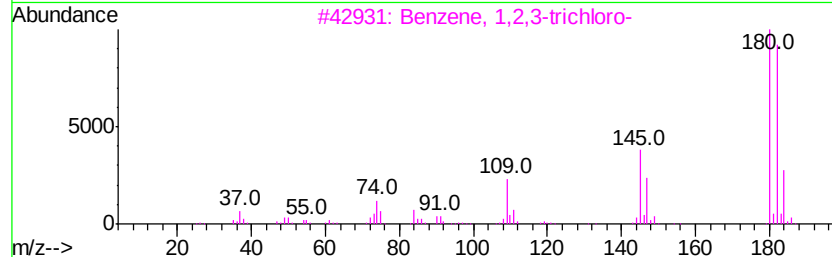
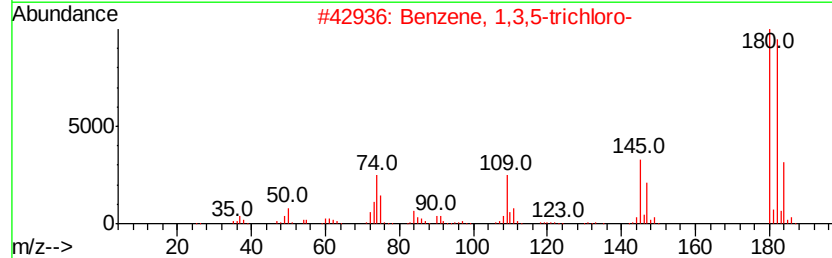
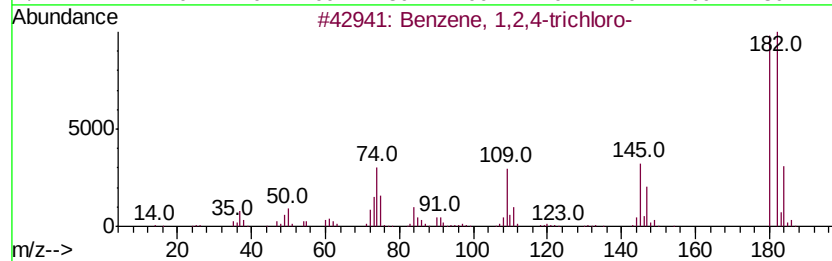
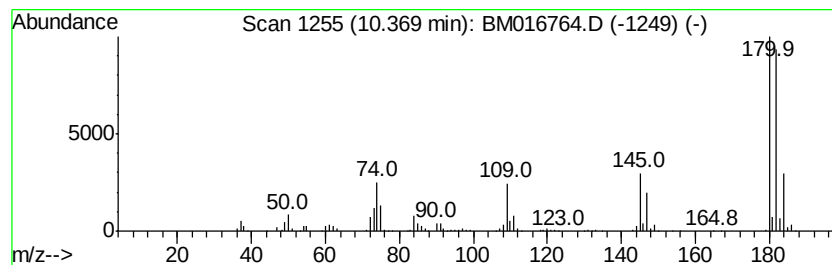
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benzene, 1,2,4-trichloro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.37	41.66 ng/ul	4290230	Naphthalene-d8	10.52

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM091718\
Data File : BM016764.D
Acq On : 17 Sep 2018 11:30
Operator : SJ/JU
Sample : J4874-13
Misc : PE
ALS Vial : 3 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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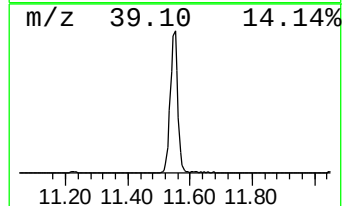
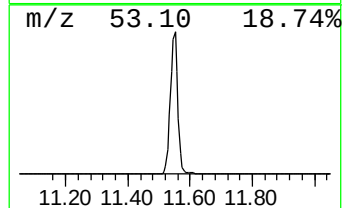
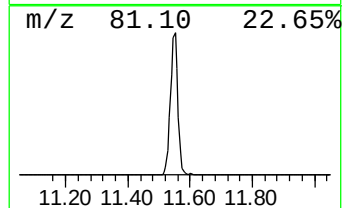
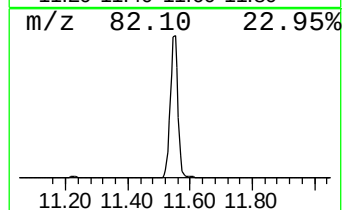
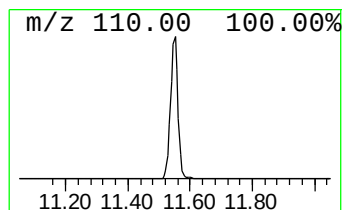
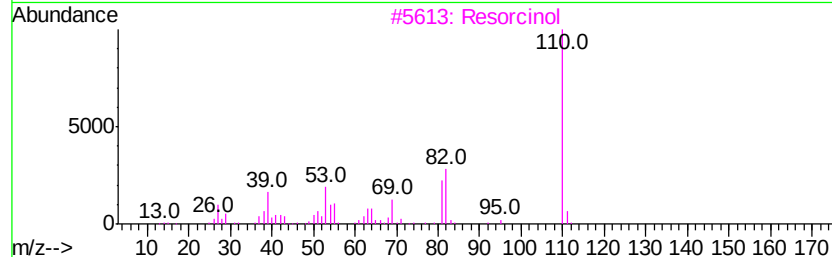
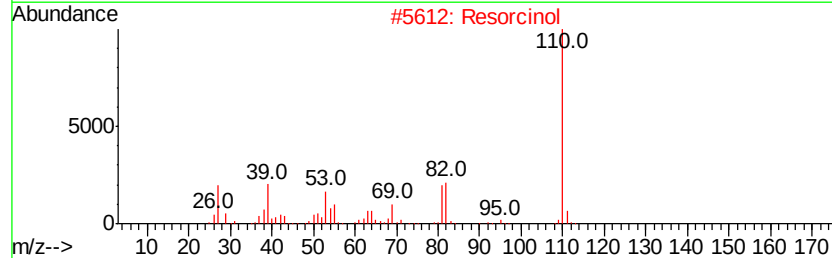
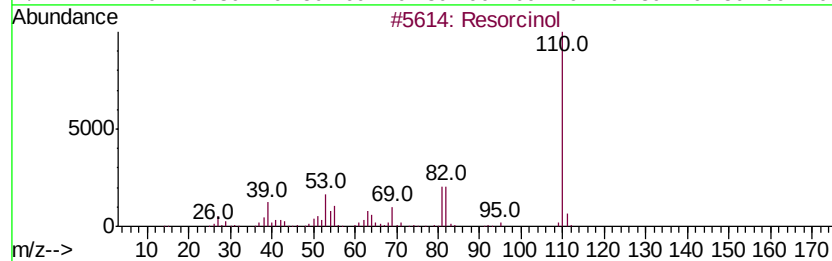
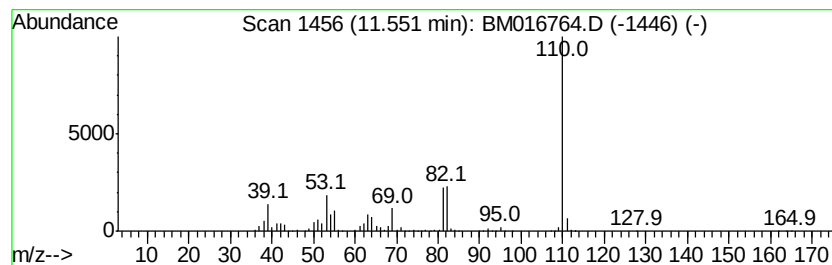
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Resorcinol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.55	27.44 ng/ul	2825960	Napthalene-d8	10.52

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Resorcinol	110	C6H6O2	000108-46-3	97
2		Resorcinol	110	C6H6O2	000108-46-3	95
3		Resorcinol	110	C6H6O2	000108-46-3	95
4		Resorcinol	110	C6H6O2	000108-46-3	91
5		Hydroquinone	110	C6H6O2	000123-31-9	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM091718\
Data File : BM016764.D
Acq On : 17 Sep 2018 11:30
Operator : SJ/JU
Sample : J4874-13
Misc : PE
ALS Vial : 3 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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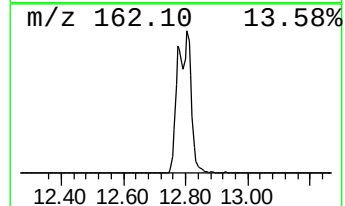
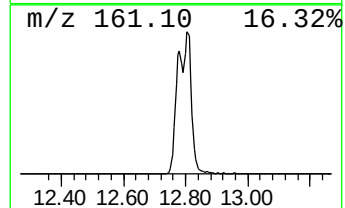
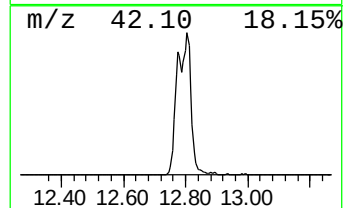
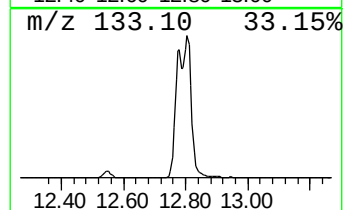
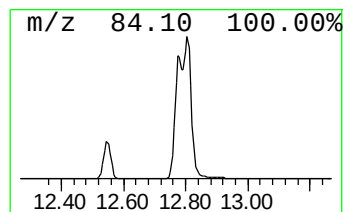
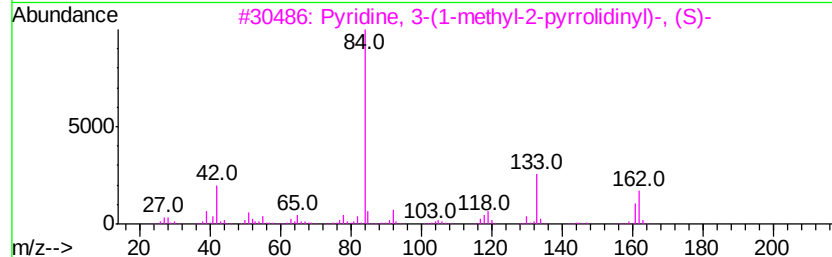
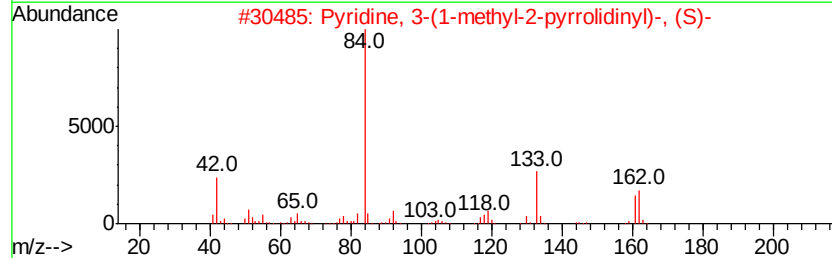
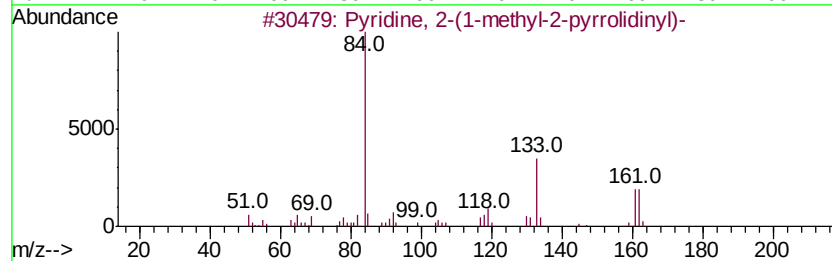
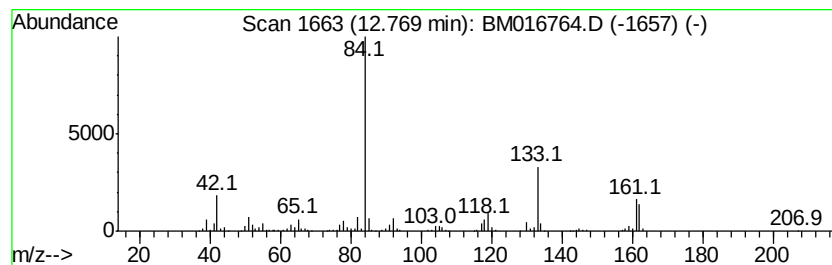
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Pyridine, 2-(1-methyl-2-pyr... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.77	11.23 ng/ul	1398850	Acenaphthene-d10	14.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pyridine, 2-(1-methyl-2-pyrrolid...	162	C10H14N2	023950-04-1	94
2		Pyridine, 3-(1-methyl-2-pyrrolid...	162	C10H14N2	000054-11-5	94
3		Pyridine, 3-(1-methyl-2-pyrrolid...	162	C10H14N2	000054-11-5	91
4		Pyridine, 3-(1-methyl-2-pyrrolid...	162	C10H14N2	000054-11-5	91
5		Pyridine, 3-(1-methyl-2-pyrrolid...	162	C10H14N2	000054-11-5	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM091718\
Data File : BM016764.D
Acq On : 17 Sep 2018 11:30
Operator : SJ/JU
Sample : J4874-13
Misc : PE
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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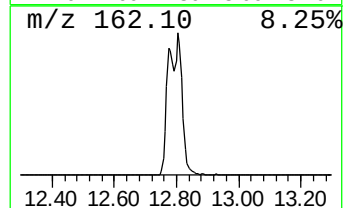
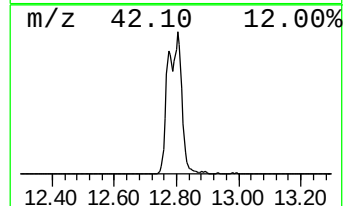
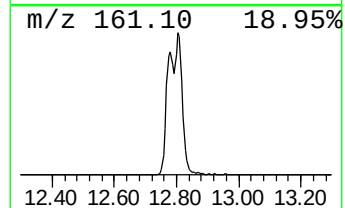
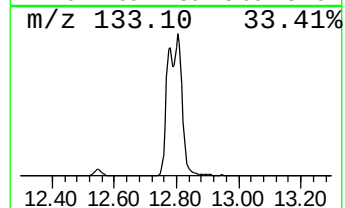
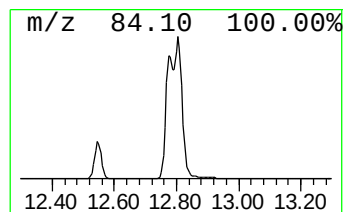
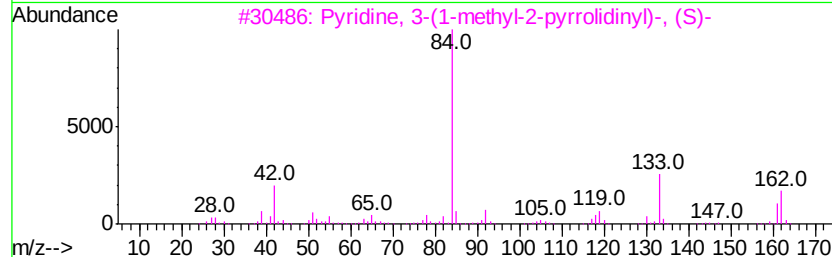
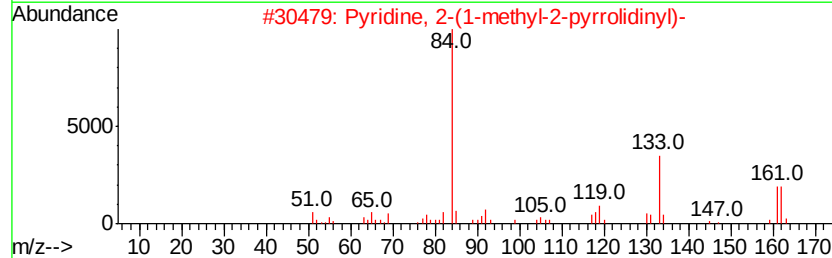
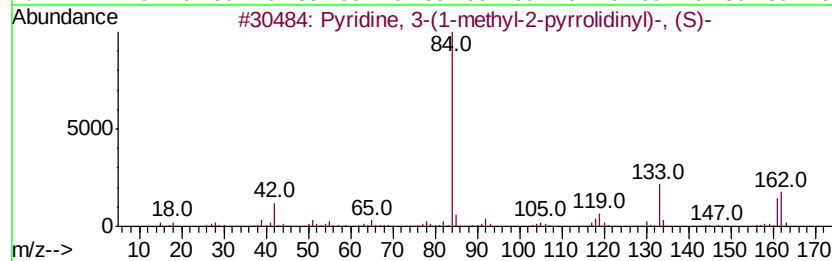
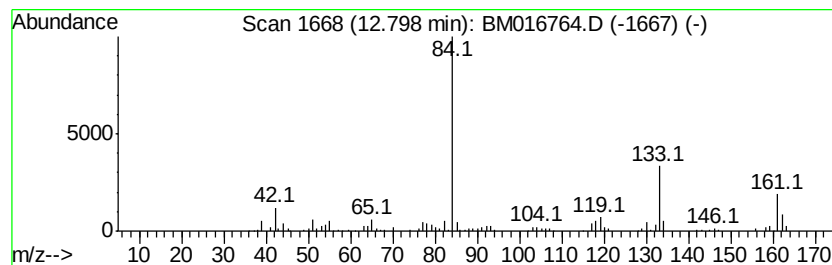
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Pyridine, 3-(1-methyl-2-pyr... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.80	10.65 ng/ul	1326520	Acenaphthene-d10	14.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pyridine, 3-(1-methyl-2-pyrrolid...	162	C10H14N2	000054-11-5	90
2		Pyridine, 2-(1-methyl-2-pyrrolid...	162	C10H14N2	023950-04-1	90
3		Pyridine, 3-(1-methyl-2-pyrrolid...	162	C10H14N2	000054-11-5	90
4		Pyridine, 3-(1-methyl-2-pyrrolid...	162	C10H14N2	000054-11-5	86
5		Pyridine, 3-(1-methyl-2-pyrrolid...	162	C10H14N2	000054-11-5	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM091718\
Data File : BM016764.D
Acq On : 17 Sep 2018 11:30
Operator : SJ/JU
Sample : J4874-13
Misc : PE
ALS Vial : 3 Sample Multiplier: 1

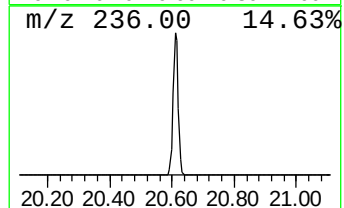
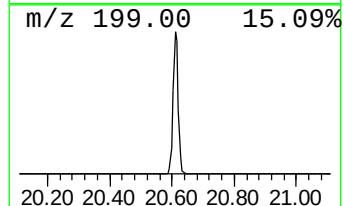
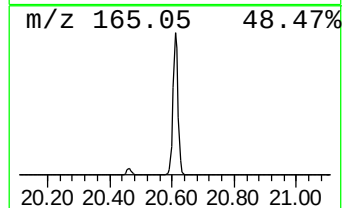
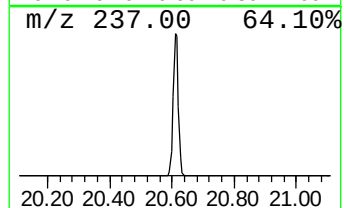
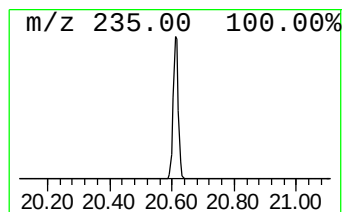
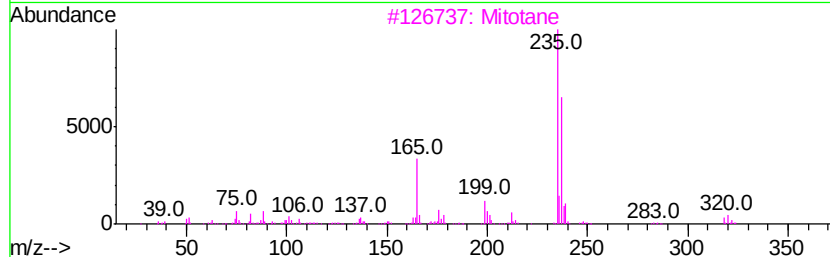
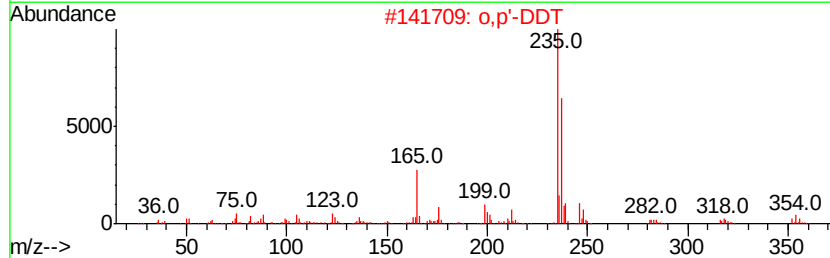
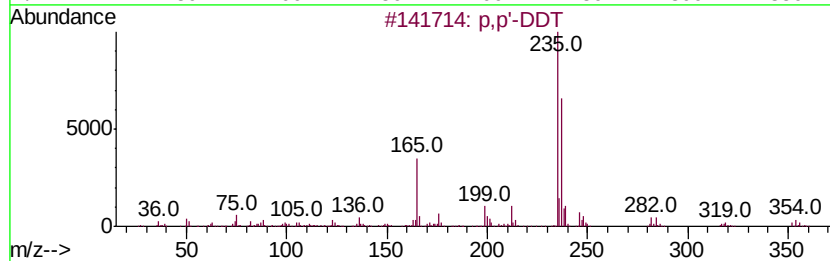
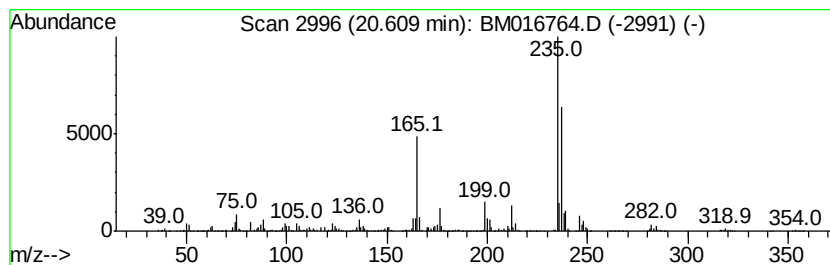
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 p,p'-DDT Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.61	14.06 ng/ul	8978110	Chrysene-d12	21.31
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	p,p'-DDT		352 C14H9Cl5	000050-29-3 99
2	o,p'-DDT		352 C14H9Cl5	000789-02-6 97
3	Mitotane		318 C14H10Cl4	000053-19-0 95
4	p,p'-DDT		352 C14H9Cl5	000050-29-3 93
5	3-Butanone, 1,1-bis(4-chlorophen...		320 C18H18Cl2O	1000158-20-4 93



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM091718\
Data File : BM016764.D
Acq On : 17 Sep 2018 11:30
Operator : SJ/JU
Sample : J4874-13
Misc : PE
ALS Vial : 3 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.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
Butane, 2-methoxy...	3.02	43.6	ng/ul	3155670	1	7.73	1446660	20.0
Benzene, 1,4-dich...	7.62	43.4	ng/ul	3140810	1	7.73	1446660	20.0
Benzene, 1,3-dich...	8.08	58.0	ng/ul	4195270	1	7.73	1446660	20.0
Propane, 1,1'-oxy...	8.32	8.3	ng/ul	597406	1	7.73	1446660	20.0
Benzene, 1,2,4-tr...	10.37	41.7	ng/ul	4290230	2	10.52	2059610	20.0
Resorcinol	11.55	27.4	ng/ul	2825960	2	10.52	2059610	20.0
Pyridine, 2-(1-me...	12.77	11.2	ng/ul	1398850	3	14.37	2491110	20.0
Pyridine, 3-(1-me...	12.80	10.7	ng/ul	1326520	3	14.37	2491110	20.0
p,p'-DDT	20.61	14.1	ng/ul	8978110	5	21.31	12772900	20.0