

Data Path : Z:\HPCHEM1\BNA M\DATA\BM100817\
 Data File : BM011985.D
 Acq On : 08 Oct 2017 20:10
 Operator : SJ/JU
 Sample : I5695-02
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 TWP-B-44

Integration Parameters: LSCINT.P

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

Filtering: 5
 Min Area: 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-BM100317.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	3.263	26	30	43	rVB	69841	118105	1.12%	0.142%
2	5.475	399	406	422	rVB	390110	796962	7.55%	0.957%
3	5.851	463	470	477	rBV	189168	305514	2.89%	0.367%
4	6.934	649	654	659	rBV	70865	116844	1.11%	0.140%
5	7.010	659	667	673	rVV	349603	669163	6.34%	0.803%
6	7.063	673	676	683	rVB	113631	177128	1.68%	0.213%
7	7.181	688	696	702	rBV	903948	1494980	14.16%	1.794%
8	7.381	723	730	744	rBV	1115456	1911346	18.10%	2.294%
9	7.492	744	749	756	rVB	255224	395901	3.75%	0.475%
10	7.851	802	810	823	rVV	702579	1193025	11.30%	1.432%
11	7.963	823	829	837	rVB	129404	227825	2.16%	0.273%
12	8.487	912	918	923	rBV	84291	143397	1.36%	0.172%
13	8.557	923	930	942	rBV	1008653	1799945	17.04%	2.160%
14	8.951	992	997	1001	rVB2	111013	156280	1.48%	0.188%
15	9.016	1001	1008	1019	rBV	1260907	2204678	20.88%	2.646%
16	9.281	1047	1053	1062	rBV5	107910	225284	2.13%	0.270%
17	9.475	1081	1086	1091	rVB	81415	126110	1.19%	0.151%
18	9.534	1091	1096	1101	rBV	87764	136323	1.29%	0.164%
19	9.739	1123	1131	1140	rBV	1138278	1979758	18.75%	2.376%
20	9.828	1140	1146	1155	rVB2	125055	262077	2.48%	0.315%
21	10.051	1175	1184	1190	rBV2	113572	210146	1.99%	0.252%
22	10.275	1215	1222	1233	rBV	1627023	2978497	28.21%	3.575%
23	10.651	1278	1286	1293	rBV	1034619	1851332	17.53%	2.222%
24	13.904	1832	1839	1850	rBV	3450427	4841999	45.85%	5.812%
25	14.192	1881	1888	1900	rBV	3827384	5727435	54.24%	6.874%
26	14.498	1931	1940	1947	rBV2	1667465	2626092	24.87%	3.152%
27	14.704	1969	1975	1985	rBV2	210154	543650	5.15%	0.653%
28	15.492	2102	2109	2115	rBV	4912918	7053923	66.80%	8.466%
29	15.616	2124	2130	2142	rBV	1704251	2547490	24.12%	3.058%
30	15.727	2145	2149	2156	rVB2	64838	120740	1.14%	0.145%
31	17.245	2400	2407	2413	rBV2	2182095	3262446	30.89%	3.916%
32	17.339	2417	2423	2436	rVV	5557360	7693271	72.85%	9.234%
33	18.115	2551	2555	2561	rBV2	104523	146667	1.39%	0.176%
34	19.268	2747	2751	2761	rVB2	74058	150008	1.42%	0.180%

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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: 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	19.392	2768	2772	2781	rBV	227919	383909	3.64%	0.461%
36	19.633	2807	2813	2825	rBV	7119081	9756585	92.39%	11.710%
37	21.415	3111	3116	3124	rBV	3149447	4151406	39.31%	4.983%
38	23.591	3478	3486	3501	rBV2	5132055	10559959	100.00%	12.675%
39	23.733	3504	3510	3521	rVB	2128084	4270120	40.44%	5.125%

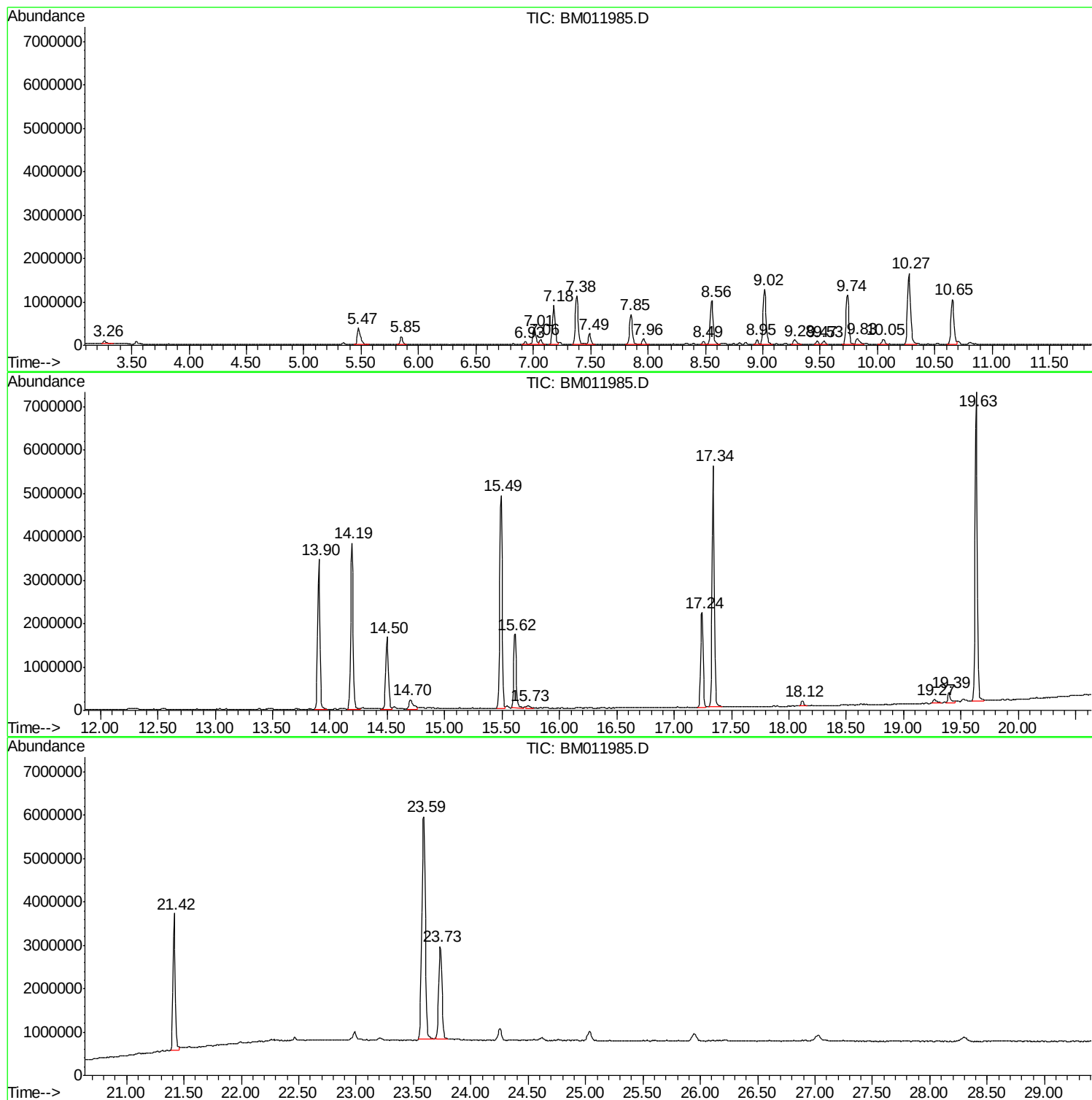
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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM100317.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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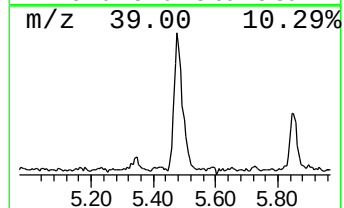
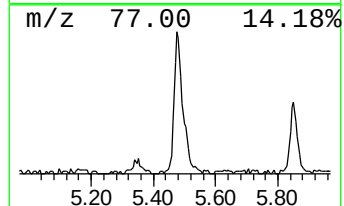
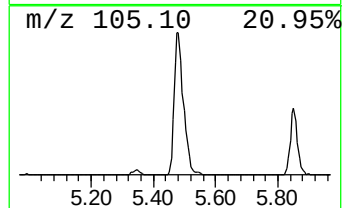
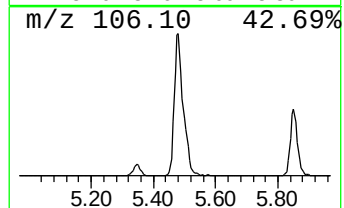
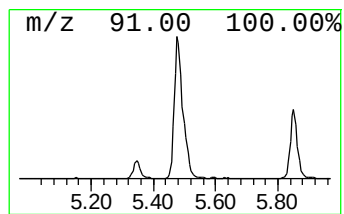
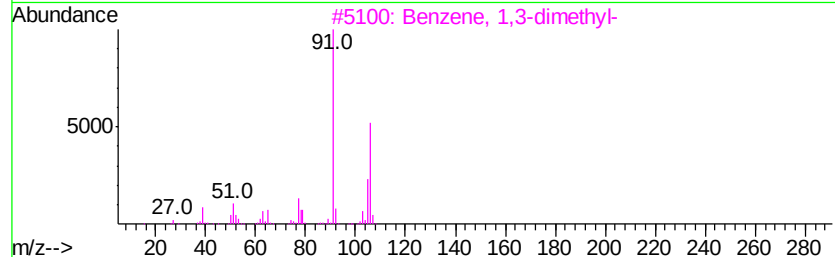
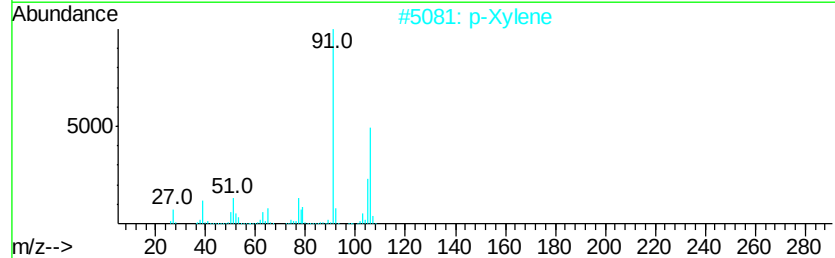
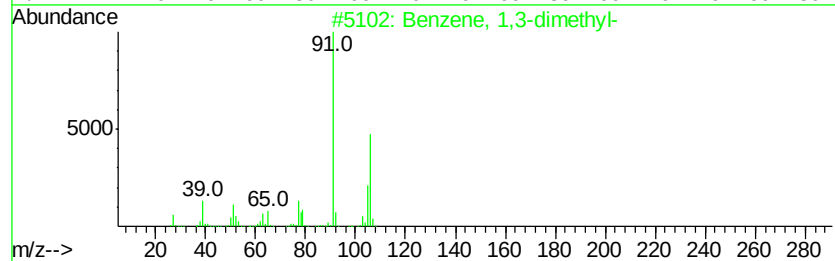
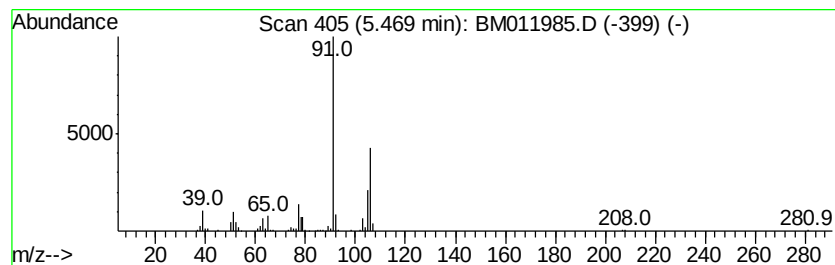
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.47	13.36 ng/ul	796962	1,4-Dichlorobenzene-d4	7.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
2		p-Xylene	106	C8H10	000106-42-3	97
3		Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	95
4		Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	95
5		p-Xylene	106	C8H10	000106-42-3	94



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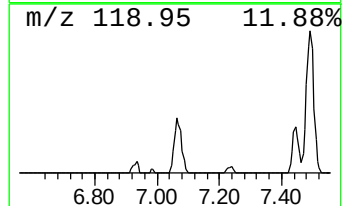
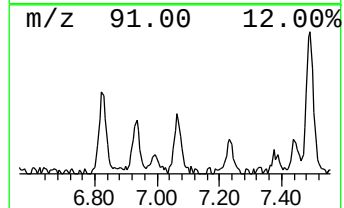
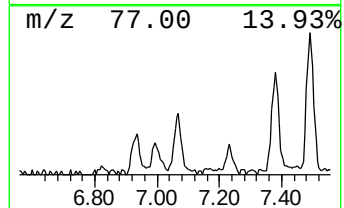
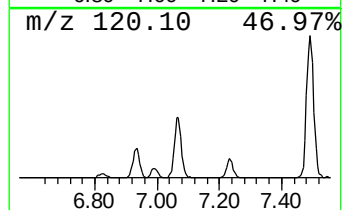
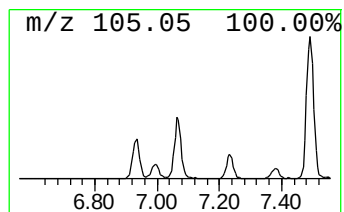
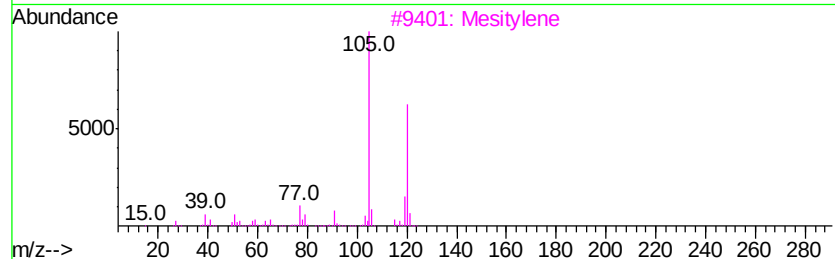
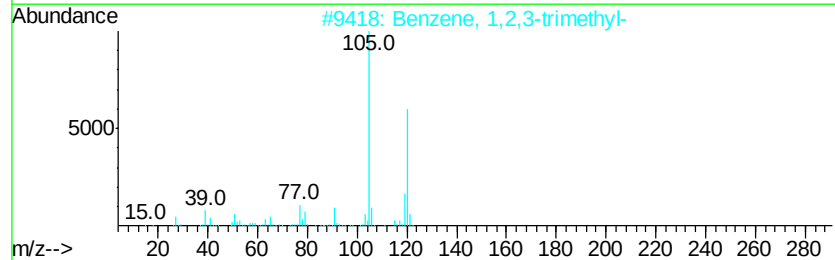
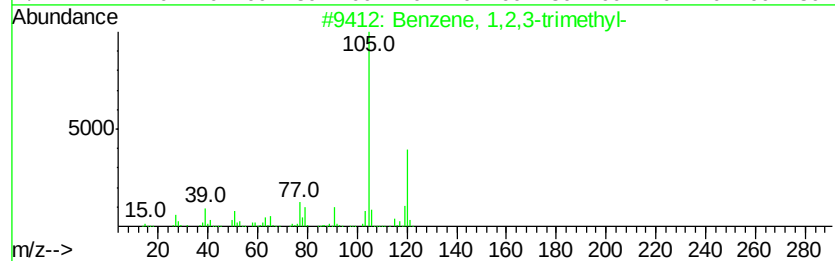
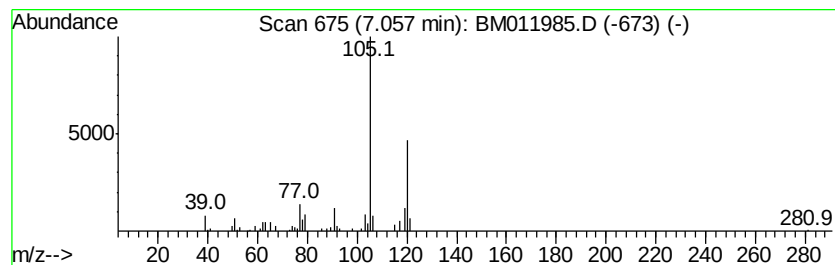
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM100317.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.06	2.97 ng/ul	177128	1,4-Dichlorobenzene-d4	7.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
3		Mesitylene	120	C9H12	000108-67-8	94
4		Mesitylene	120	C9H12	000108-67-8	91
5		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



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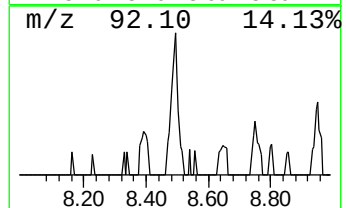
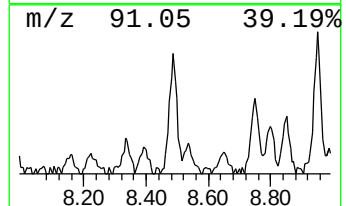
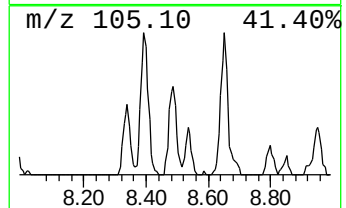
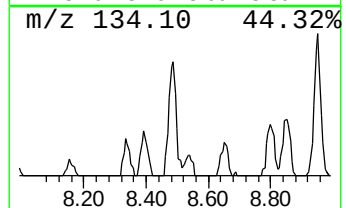
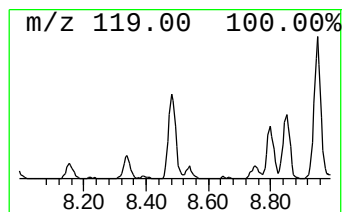
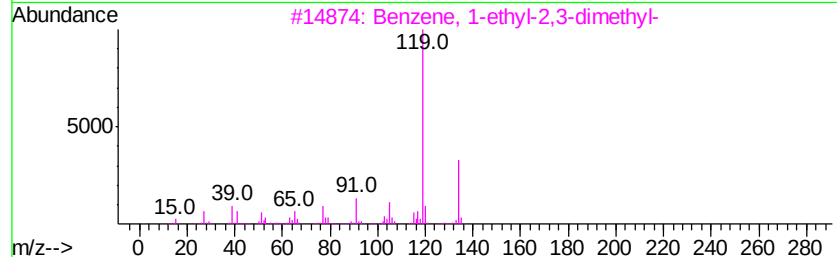
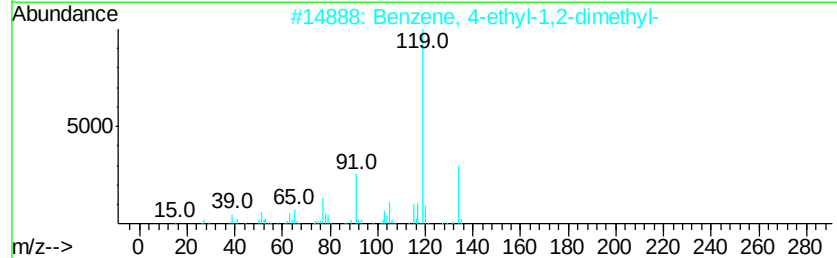
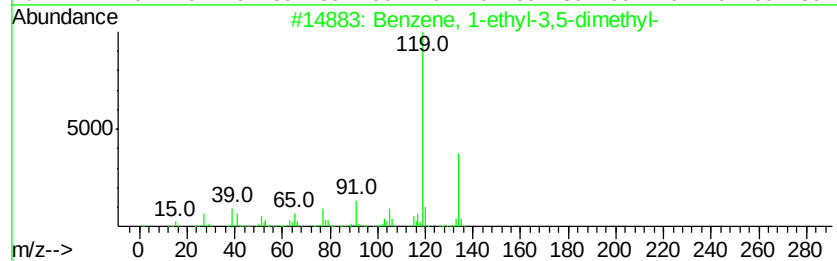
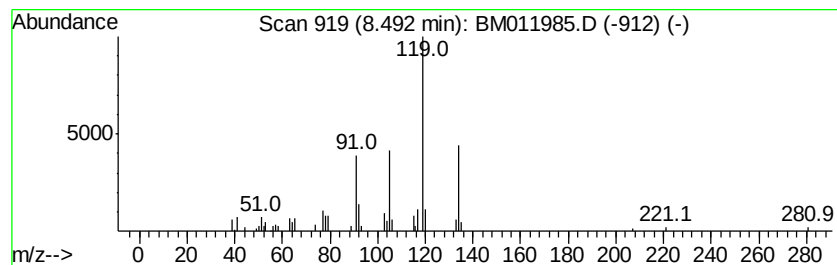
Instrument :
BNA_M
ClientSampled :
TWP-B-44

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM100317.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzene, 1-ethyl-3,5-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.		
8.49	2.40 ng/ul	143397	1,4-Dichlorobenzene-d4	7.85		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	94	
2	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	93	
3	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	93	
4	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	91	
5	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	90	



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Instrument :
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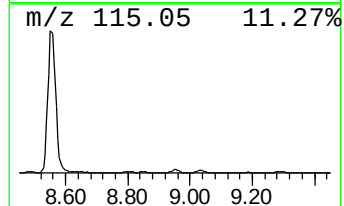
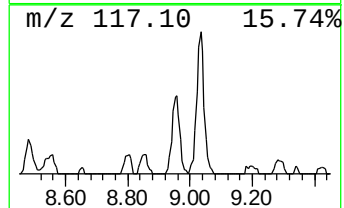
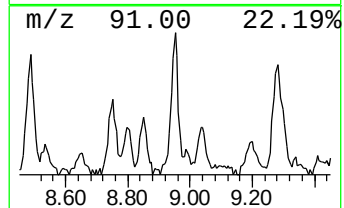
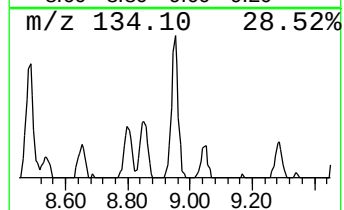
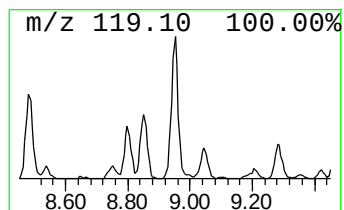
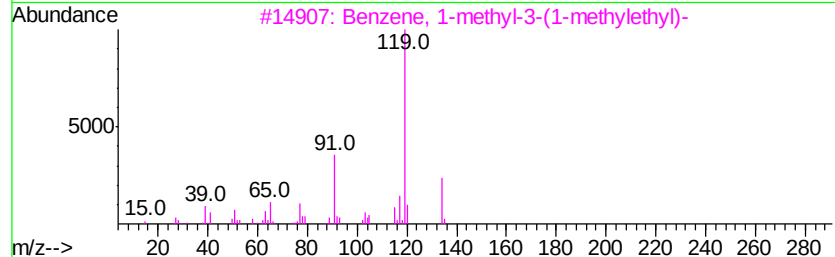
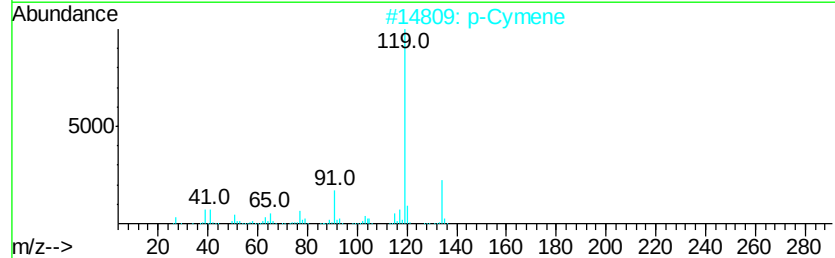
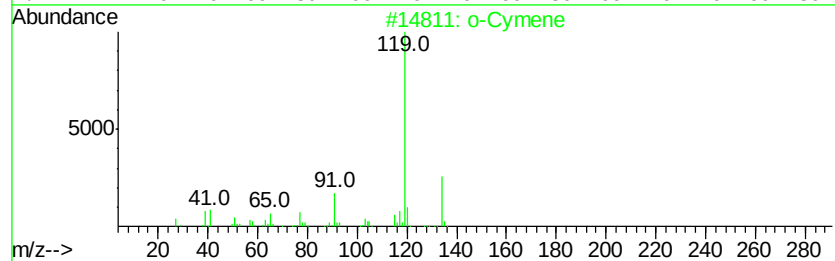
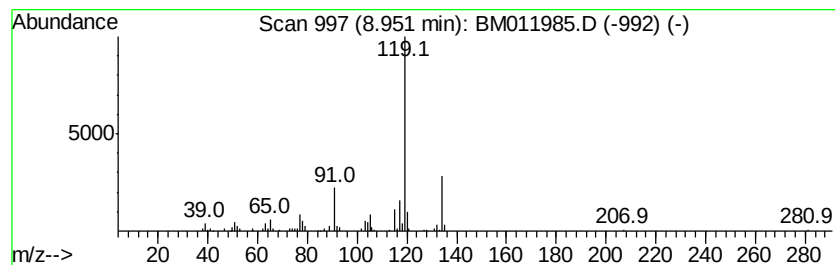
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM100317.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 o-Cymene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.95	2.62 ng/ul	156280	1,4-Dichlorobenzene-d4	7.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	o-Cymene	134	C10H14	000527-84-4	95
2		p-Cymene	134	C10H14	000099-87-6	94
3		Benzene, 1-methyl-3-(1-methylethyl-...	134	C10H14	000535-77-3	91
4		o-Cymene	134	C10H14	000527-84-4	91
5		o-Cymene	134	C10H14	000527-84-4	90



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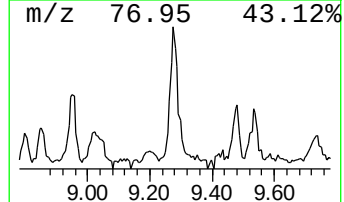
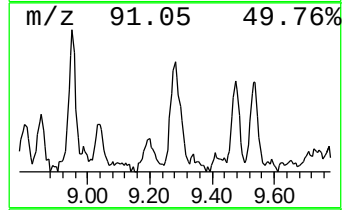
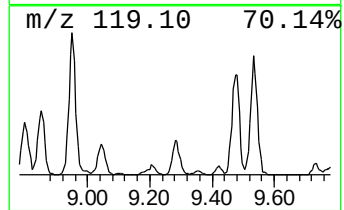
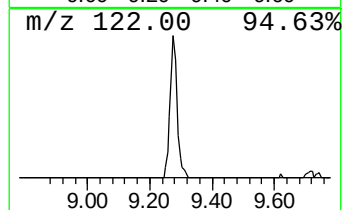
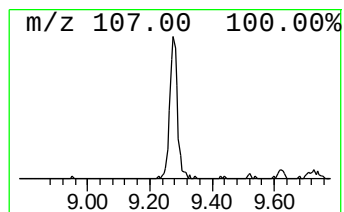
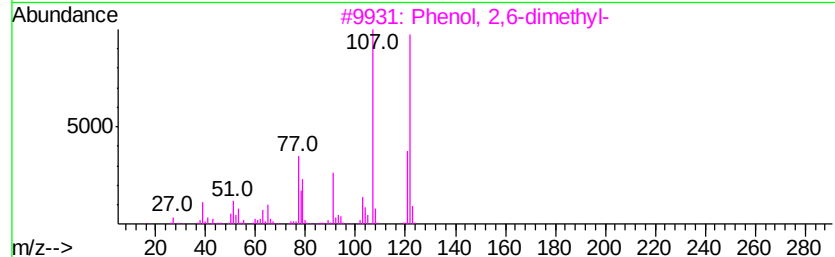
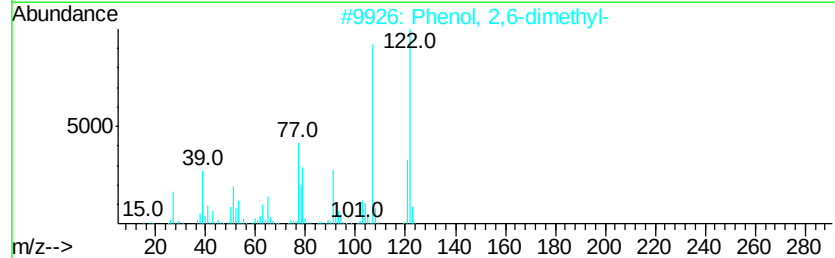
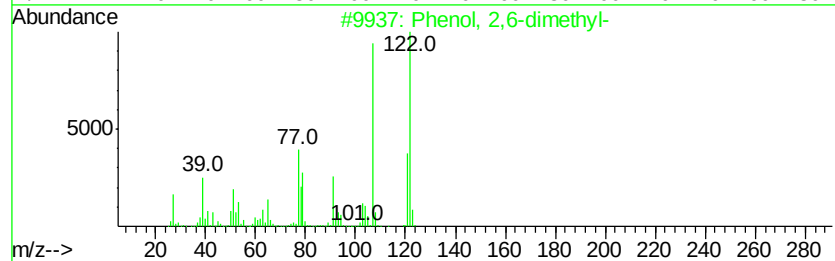
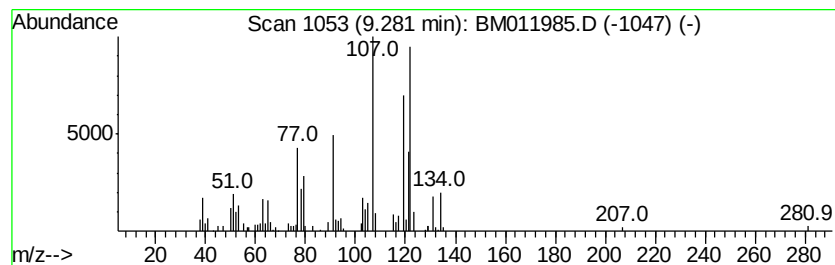
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TIC Integration Parameters: LSCINT.P

Peak Number 8 Phenol, 2,6-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.28	2.43 ng/ul	225284	Napthalene-d8	10.65

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,6-dimethyl-		122	C8H10O	000576-26-1	98
2	Phenol, 2,6-dimethyl-		122	C8H10O	000576-26-1	96
3	Phenol, 2,6-dimethyl-		122	C8H10O	000576-26-1	94
4	Phenol, 2,3-dimethyl-		122	C8H10O	000526-75-0	94
5	Phenol, 2,6-dimethyl-		122	C8H10O	000576-26-1	94



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ClientSampleId :
TWP-B-44

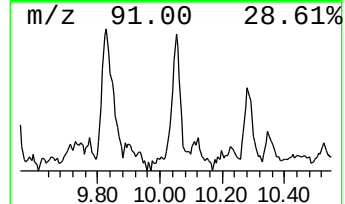
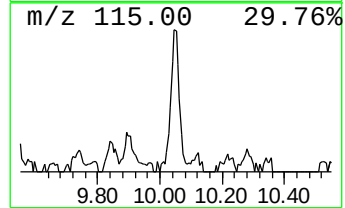
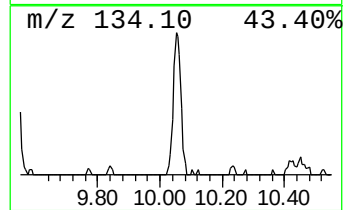
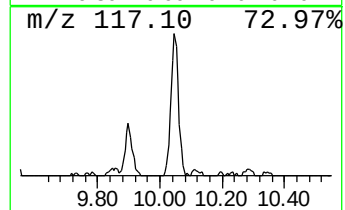
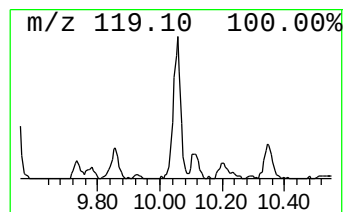
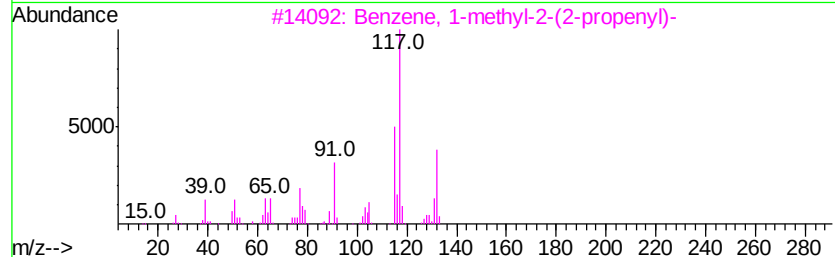
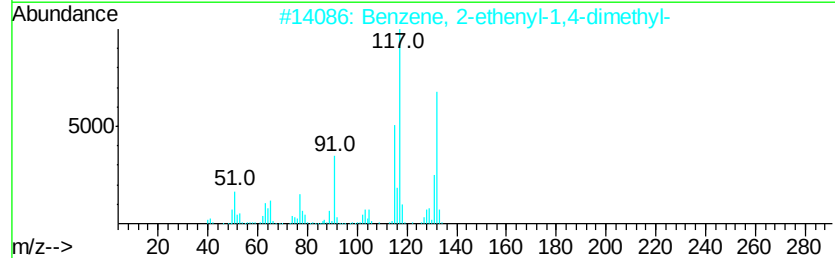
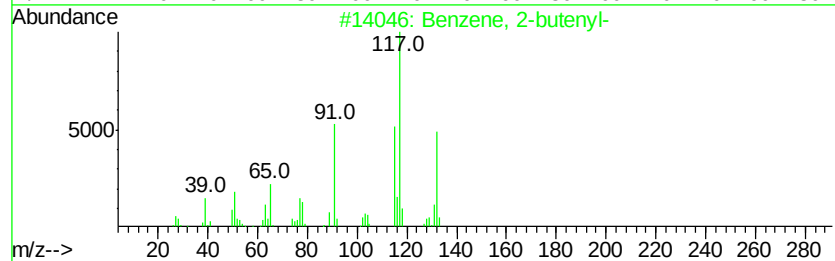
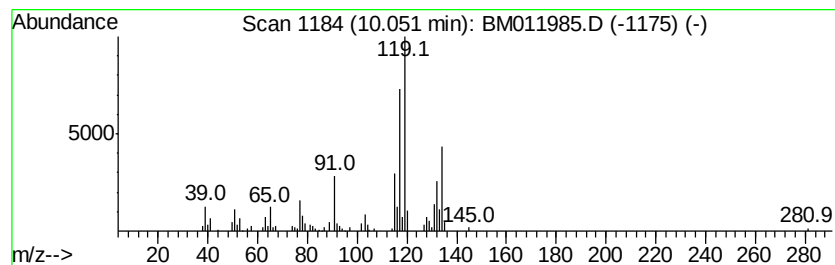
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Benzene, 2-butenyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.05	2.27 ng/ul	210146	Naphthalene-d8	10.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-butenyl-	132	C10H12	001560-06-1	83
2		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	83
3		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	80
4		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	56
5		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	55



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM100817\
Data File : BM011985.D
Acq On : 08 Oct 2017 20:10
Operator : SJ/JU
Sample : I5695-02
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
TWP-B-44

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100317.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, 1,3-dime...	5.47	13.4	ng/ul	796962	1	7.85	1193030	20.0
Benzene, 1,2,3-tr...	7.06	3.0	ng/ul	177128	1	7.85	1193030	20.0
Benzene, 1-ethyl-...	8.49	2.4	ng/ul	143397	1	7.85	1193030	20.0
o-Cymene	8.95	2.6	ng/ul	156280	1	7.85	1193030	20.0
Phenol, 2,6-dimet...	9.28	2.4	ng/ul	225284	2	10.65	1851330	20.0
Benzene, 2-butenyl-	10.05	2.3	ng/ul	210146	2	10.65	1851330	20.0