

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM101217\
 Data File : BM012126.D
 Acq On : 13 Oct 2017 07:40
 Operator : SJ/JU
 Sample : I5568-02
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 B-51(1-2)-092817

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-BM101217.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	5.334	377	382	390	rBV2	249464	403268	7.13%	0.559%
2	5.469	398	405	415	rBV	667307	1327405	23.47%	1.839%
3	5.840	461	468	474	rBV2	348691	613591	10.85%	0.850%
4	7.022	664	669	681	rVB	625622	1269272	22.44%	1.759%
5	7.175	683	695	700	rBV	516425	838498	14.82%	1.162%
6	7.381	722	730	736	rBV	757356	1215672	21.49%	1.685%
7	7.445	736	741	744	rVV	361091	518426	9.17%	0.718%
8	7.487	744	748	753	rVB	219178	324046	5.73%	0.449%
9	7.845	804	809	815	rBV	1121313	1837816	32.49%	2.547%
10	7.898	815	818	824	rVB	192537	294308	5.20%	0.408%
11	8.387	893	901	909	rVB2	202012	464443	8.21%	0.644%
12	8.481	909	917	924	rBV3	364707	672998	11.90%	0.933%
13	8.569	925	932	940	rBV	807692	1438648	25.43%	1.994%
14	8.945	984	996	1001	rBV	225253	389492	6.89%	0.540%
15	9.022	1001	1009	1012	rVV	641059	1195547	21.14%	1.657%
16	9.051	1012	1014	1024	rVB	342399	475405	8.40%	0.659%
17	9.534	1090	1096	1103	rVB	135995	235723	4.17%	0.327%
18	9.739	1124	1131	1141	rVB2	516170	1008095	17.82%	1.397%
19	9.992	1165	1174	1178	rBV	128585	232263	4.11%	0.322%
20	10.286	1218	1224	1230	rBV	841956	1514254	26.77%	2.098%
21	10.345	1230	1234	1243	rVB2	121018	242864	4.29%	0.337%
22	10.657	1280	1287	1292	rBV	1491693	2549683	45.08%	3.533%
23	10.704	1292	1295	1303	rVB	649437	1037090	18.34%	1.437%
24	10.804	1303	1312	1318	rBV2	149035	333182	5.89%	0.462%
25	11.481	1418	1427	1435	rBV2	134485	354103	6.26%	0.491%
26	13.327	1732	1741	1752	rBV	1460859	2297028	40.61%	3.183%
27	13.545	1772	1778	1783	rBV	4013400	5656218	100.00%	7.838%
28	13.592	1783	1786	1791	rVB	277411	407312	7.20%	0.564%
29	13.904	1834	1839	1843	rBV	1499655	2081648	36.80%	2.885%
30	13.951	1843	1847	1852	rVB	700927	947805	16.76%	1.313%
31	14.198	1883	1889	1900	rBV	1882190	2831741	50.06%	3.924%
32	14.504	1935	1941	1948	rVB2	2212691	3290395	58.17%	4.560%
33	14.733	1971	1980	1987	rBV	340499	720755	12.74%	0.999%
34	14.804	1987	1992	2000	rVB	151102	324871	5.74%	0.450%

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Min Area: 1 % of largest Peak
Max Peaks: 100
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If leading or trailing edge < 100 prefer < Baseline drop else tangent >
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35	15.321	2076	2080	2086	rVB2	195765	300551	5.31%	0.416%
36	15.498	2104	2110	2116	rBV	2234949	3146601	55.63%	4.360%
37	16.186	2224	2227	2229	rBV2	176514	230204	4.07%	0.319%
38	16.980	2359	2362	2367	rBV2	152744	211195	3.73%	0.293%
39	17.251	2403	2408	2413	rBV2	2570397	3921451	69.33%	5.434%
40	17.298	2413	2416	2420	rVB	775000	966832	17.09%	1.340%
41	17.351	2420	2425	2430	rBV2	2455106	3407921	60.25%	4.722%
42	18.068	2543	2547	2553	rBV	987756	1449433	25.63%	2.009%
43	18.133	2553	2558	2562	rBV	2456824	3554634	62.84%	4.926%
44	19.309	2755	2758	2764	rVB	1947789	2712536	47.96%	3.759%
45	19.409	2771	2775	2781	rBV	2340219	3289491	58.16%	4.558%
46	19.651	2811	2816	2819	rBV	3400855	5220565	92.30%	7.234%
47	21.321	3097	3100	3106	rVB	3971113	4408807	77.95%	6.109%

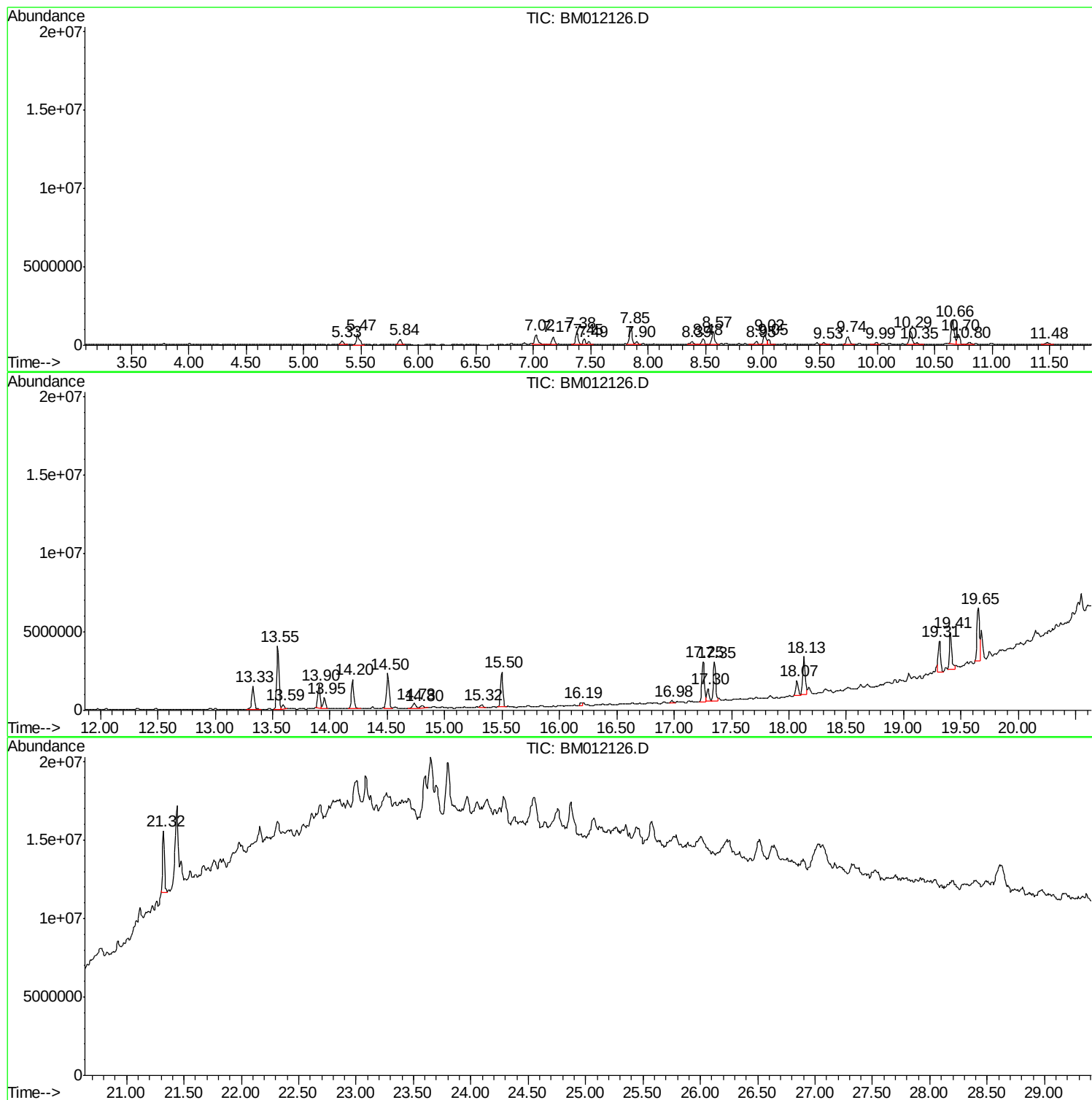
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM101217.M
Quant Title : SVOA CALIBRATION

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



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Sample : I5568-02
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
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ClientSampled :
B-51(1-2)-092817

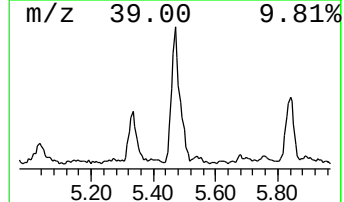
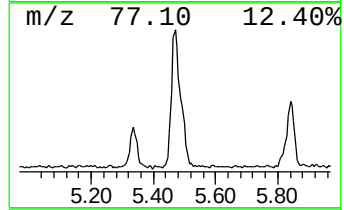
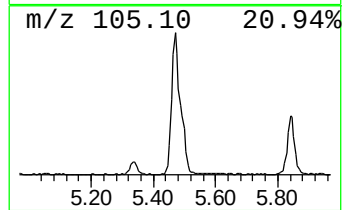
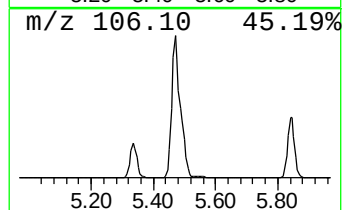
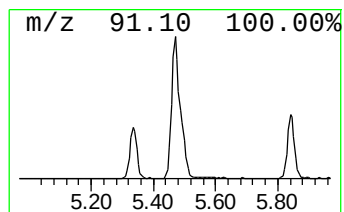
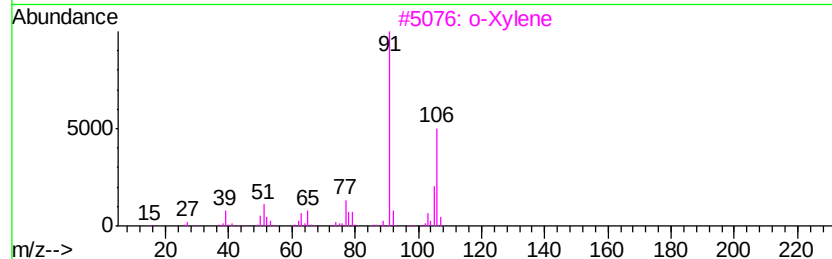
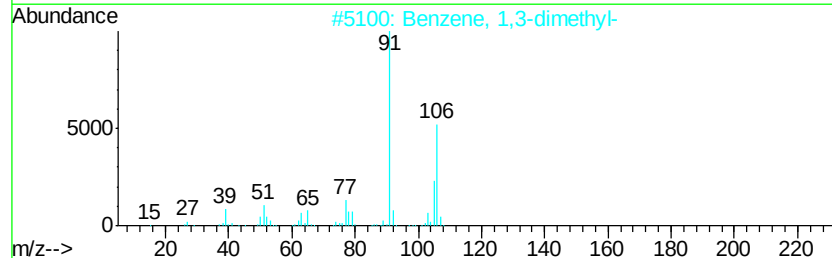
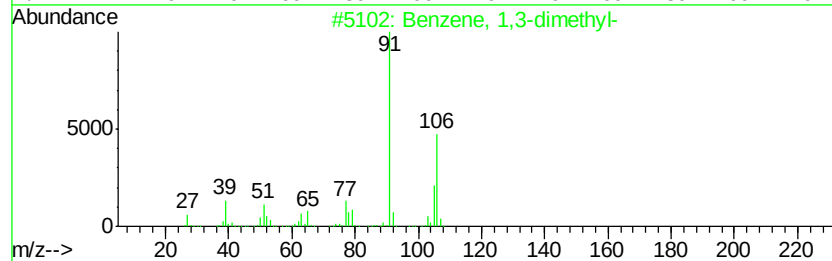
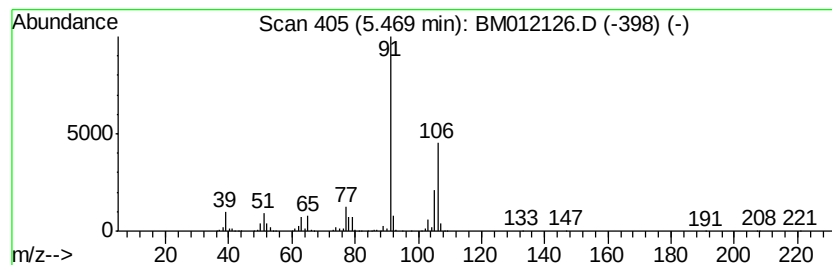
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM101217.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzene, 1,3-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.47	14.45 ng/ul	1327410	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		Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
3		o-Xylene	106	C8H10	000095-47-6	97
4		p-Xylene	106	C8H10	000106-42-3	97
5		p-Xylene	106	C8H10	000106-42-3	94



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Instrument :
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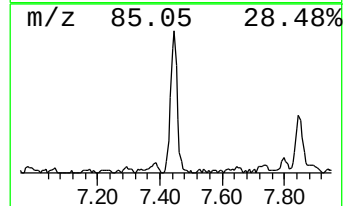
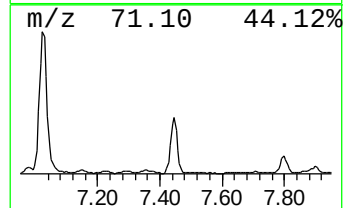
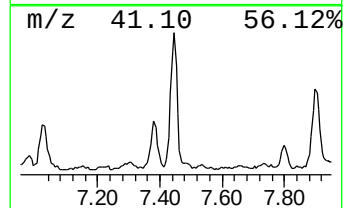
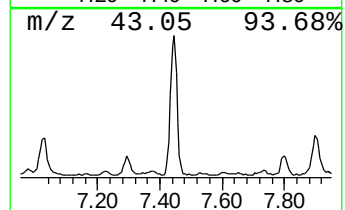
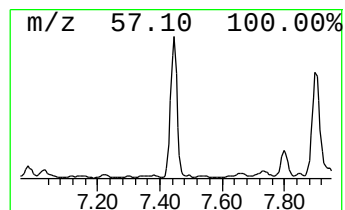
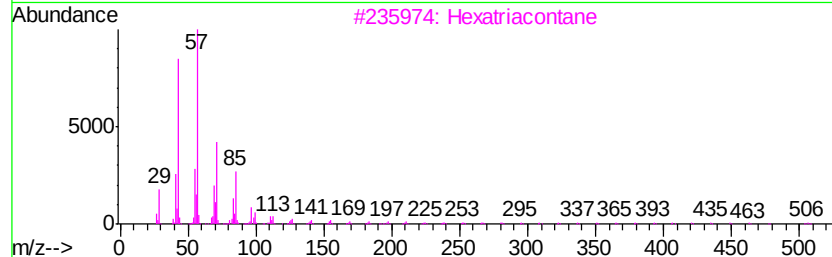
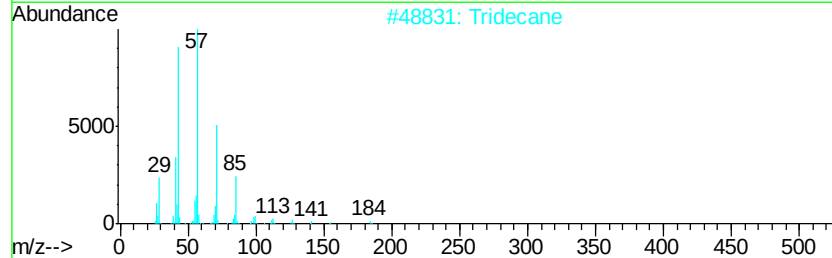
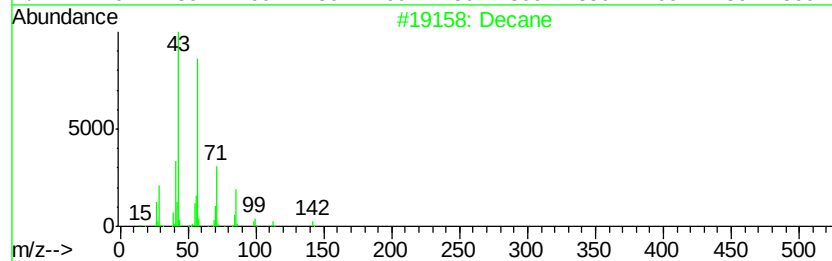
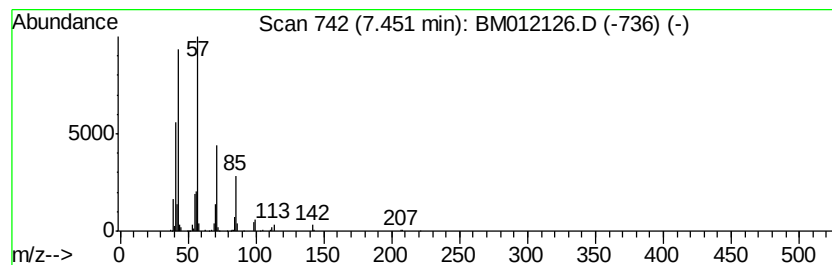
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM101217.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
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Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.45	5.64 ng/ul	518426	1,4-Dichlorobenzene-d4	7.85

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane	142	C10H22	000124-18-5	93
2			Tridecane	184	C13H28	000629-50-5	78
3			Hexatriacontane	507	C36H74	000630-06-8	78
4			Undecane, 2,3-dimethyl-	184	C13H28	017312-77-5	72
5			Decane	142	C10H22	000124-18-5	70



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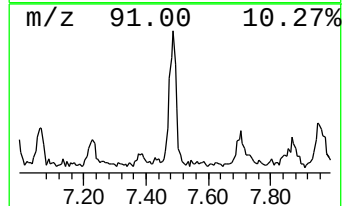
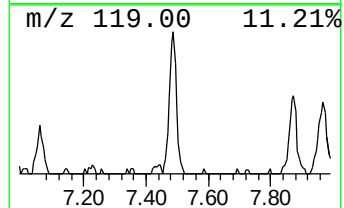
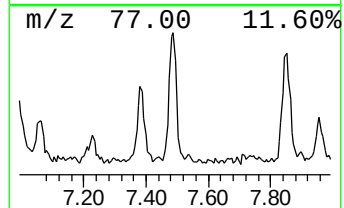
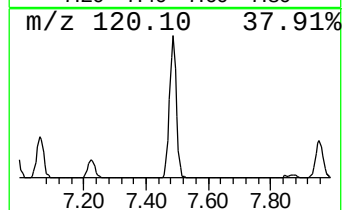
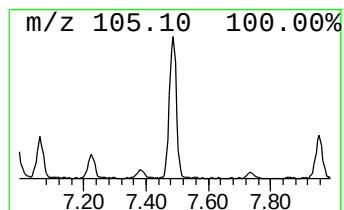
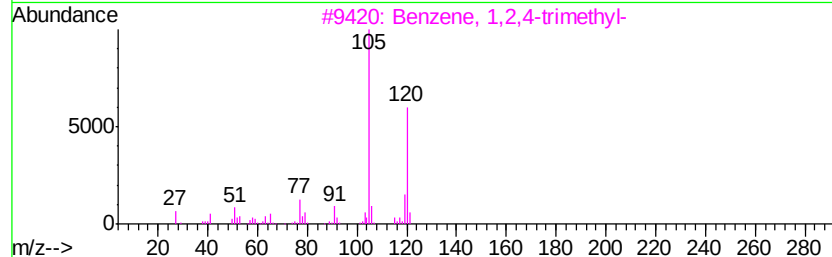
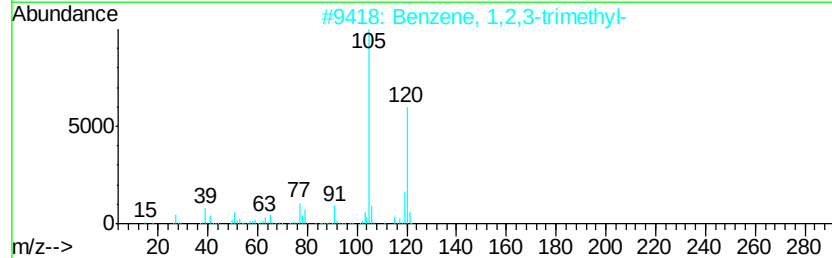
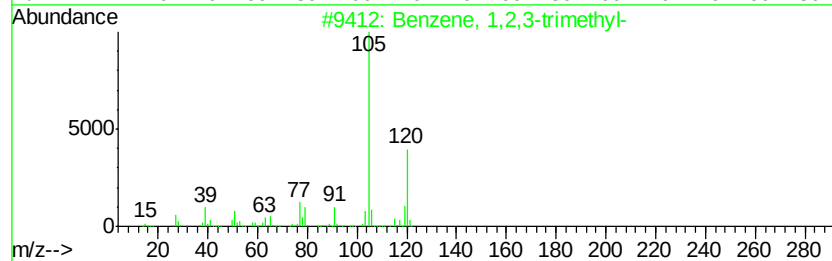
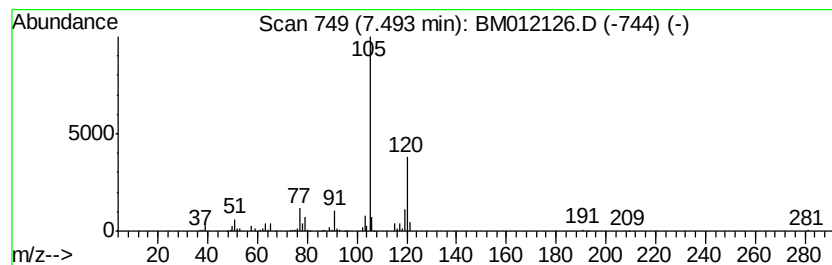
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM101217.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
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Peak Number 5 Benzene, 1,2,3-trimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.49	3.53 ng/ul	324046	1,4-Dichlorobenzene-d4	7.85

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5		Mesitylene	120	C9H12	000108-67-8	91



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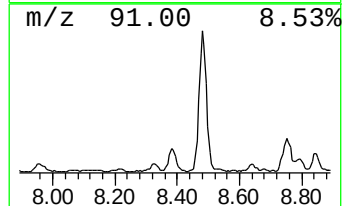
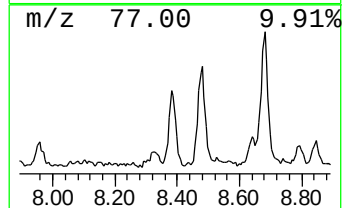
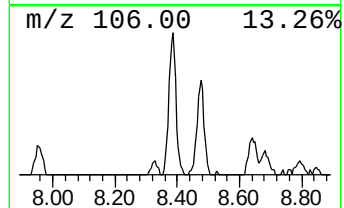
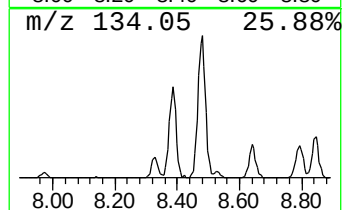
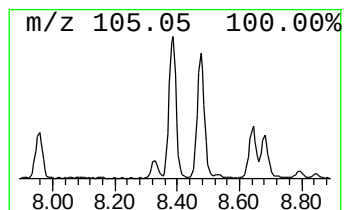
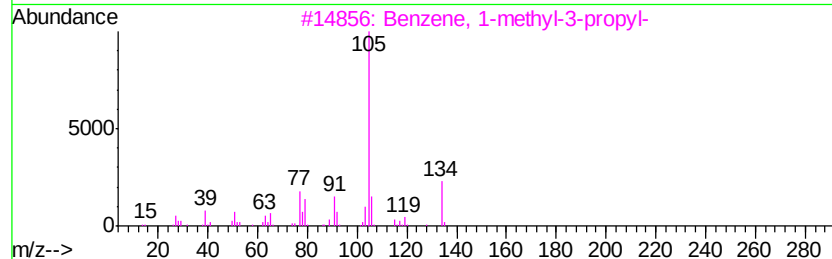
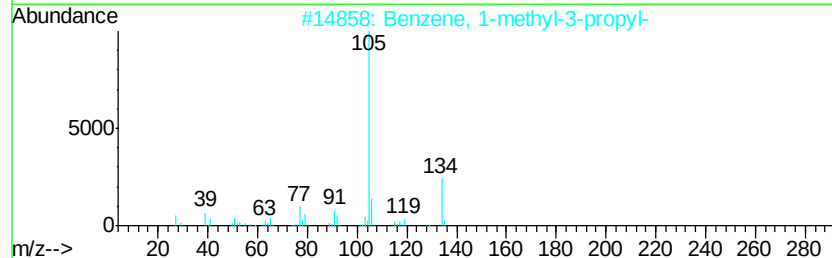
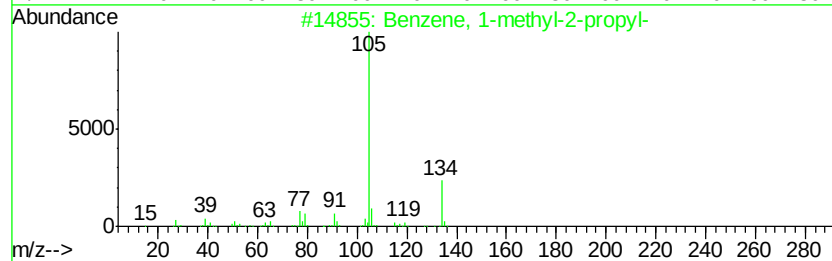
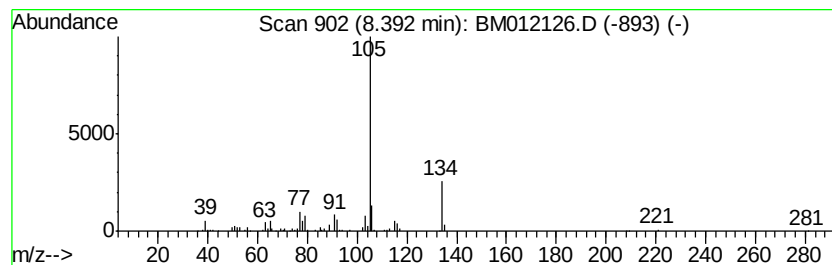
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzene, 1-methyl-2-propyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.39	5.05 ng/ul	464443	1,4-Dichlorobenzene-d4	7.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	91
2		Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	91
3		Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	91
4		Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	91
5		Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91



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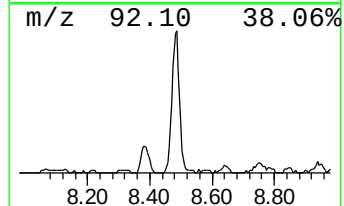
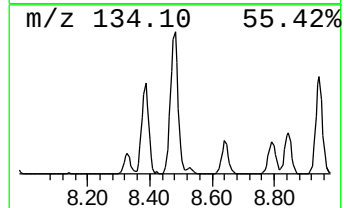
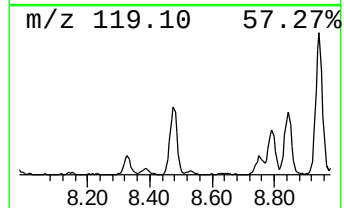
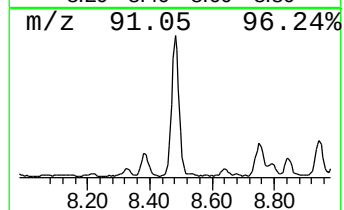
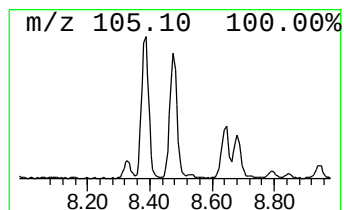
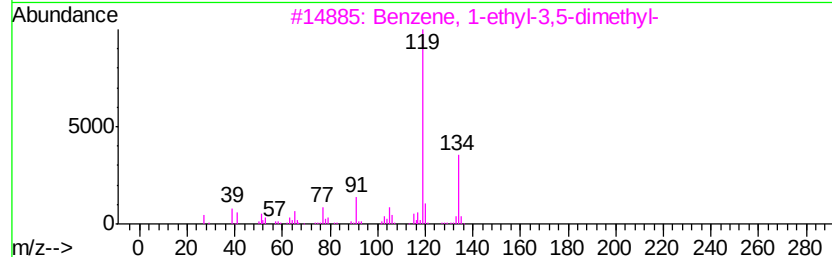
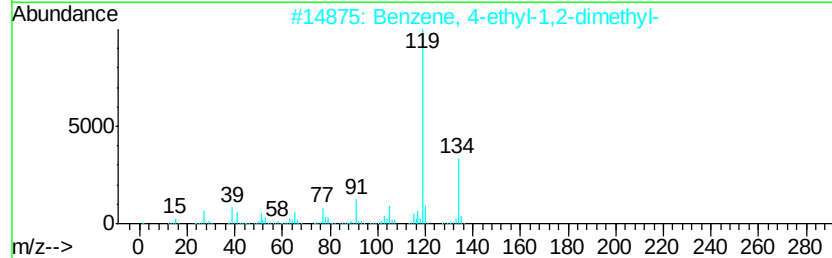
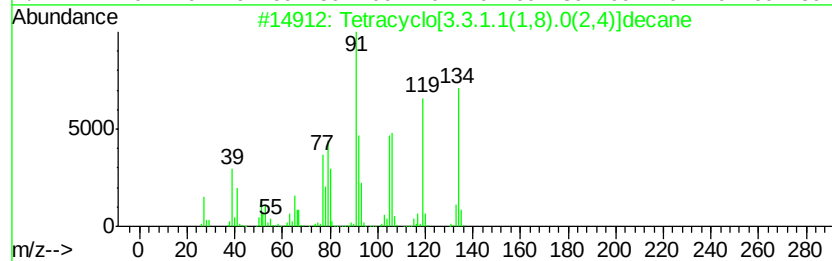
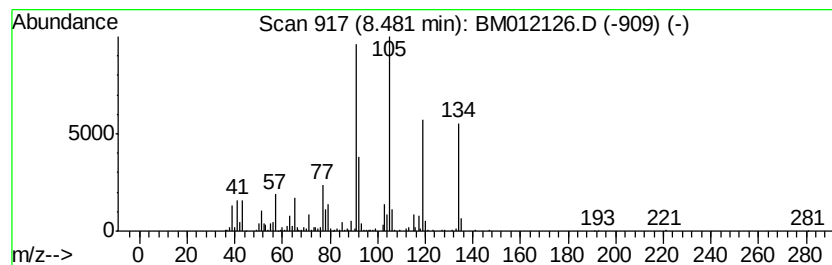
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Tetracyclo[3.3.1.1(1,8).0(2... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.48	7.32 ng/ul	672998	1,4-Dichlorobenzene-d4	7.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetracyclo[3.3.1.1(1,8).0(2,4)]d...	134	C10H14	1000185-58-7	74
2		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	70
3		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	70
4		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	70
5		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	70



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM101217\
Data File : BM012126.D
Acq On : 13 Oct 2017 07:40
Operator : SJ/JU
Sample : I5568-02
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
B-51(1-2)-092817

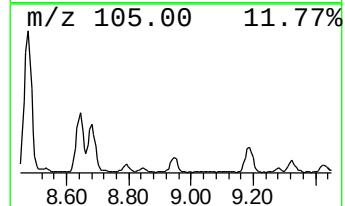
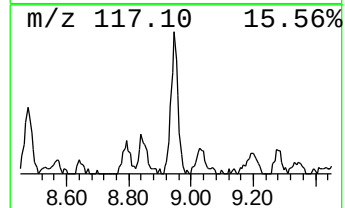
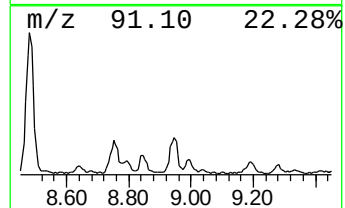
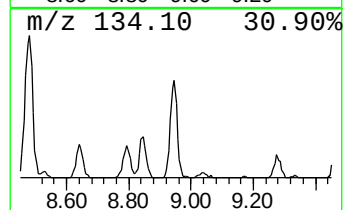
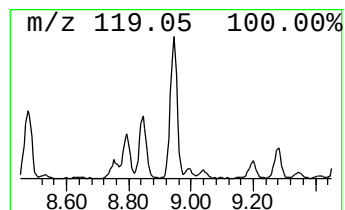
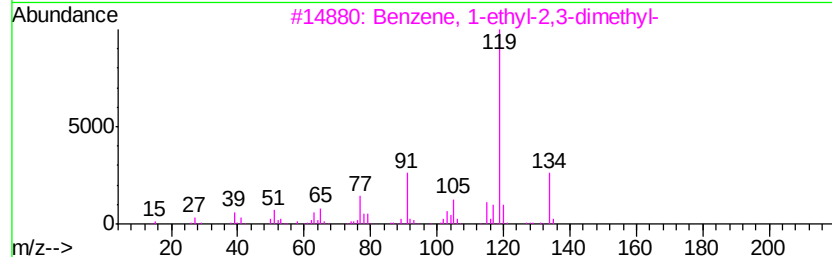
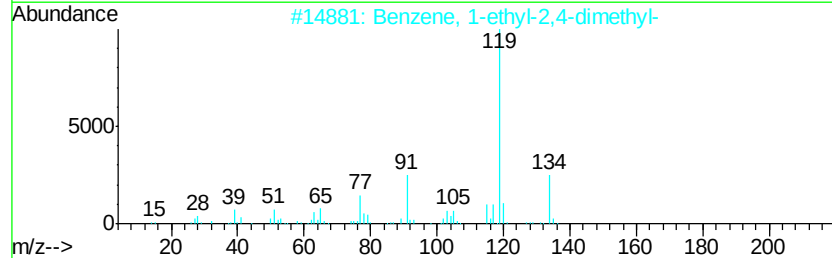
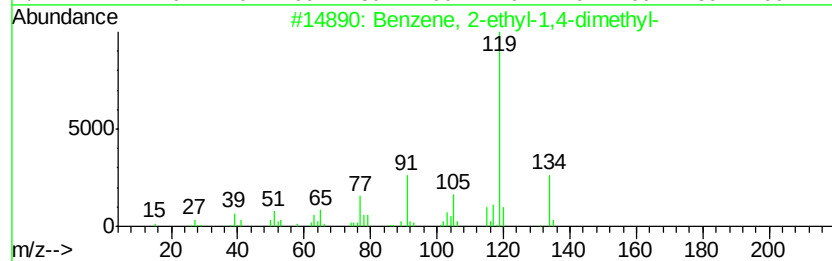
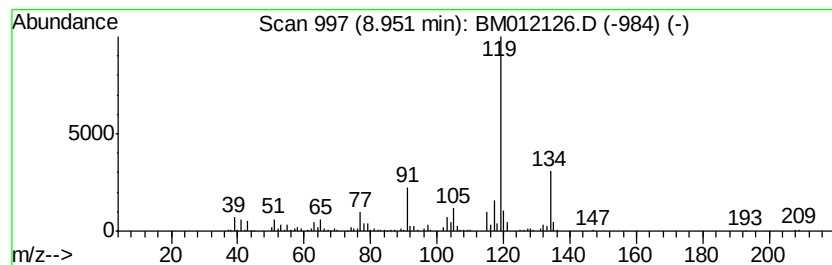
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.95	4.24 ng/ul	389492	1,4-Dichlorobenzene-d4	7.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	95
3		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
4		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94
5		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM101217\
Data File : BM012126.D
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Operator : SJ/JU
Sample : I5568-02
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ALS Vial : 26 Sample Multiplier: 1

Instrument :
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ClientSampleId :
B-51(1-2)-092817

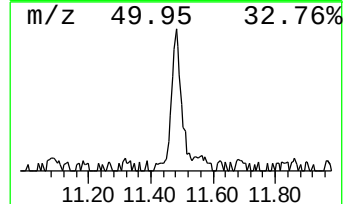
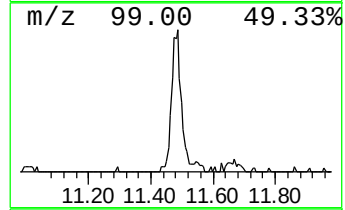
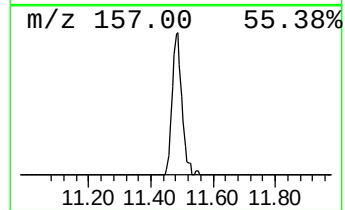
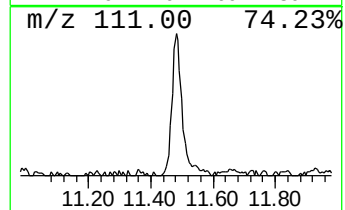
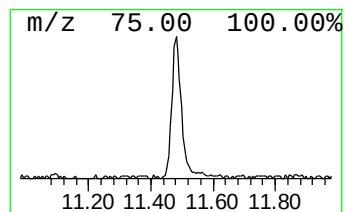
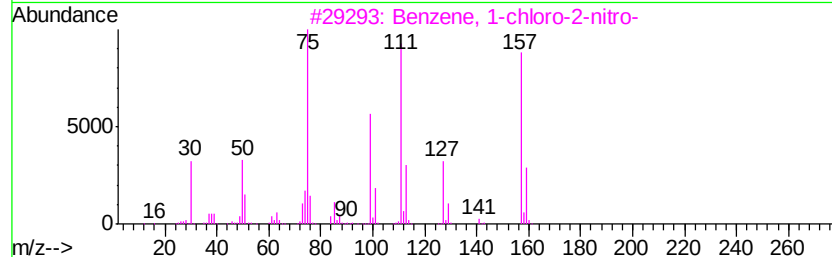
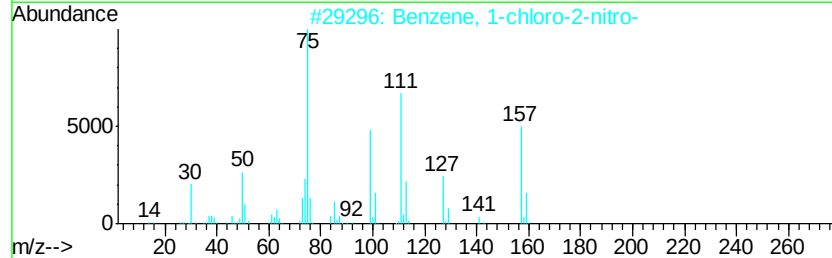
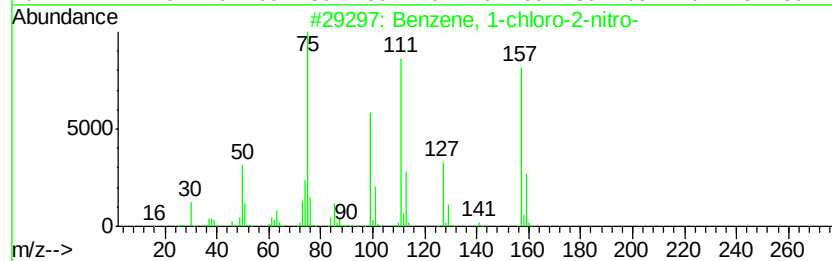
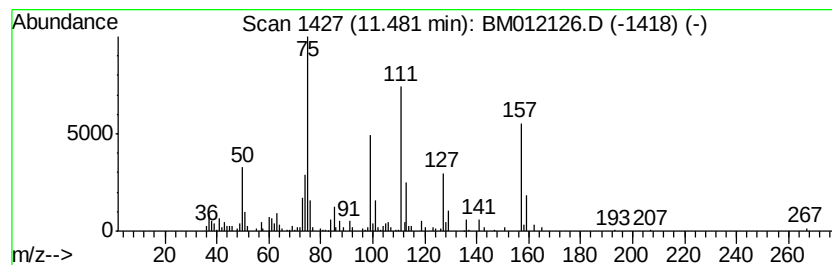
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Benzene, 1-chloro-2-nitro- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.48	2.78 ng/ul	354103	Naphthalene-d8	10.66

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



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM101217\
Data File : BM012126.D
Acq On : 13 Oct 2017 07:40
Operator : SJ/JU
Sample : I5568-02
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
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ClientSampleId :
B-51(1-2)-092817

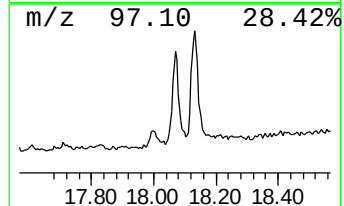
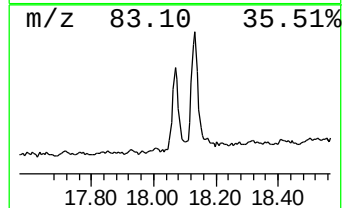
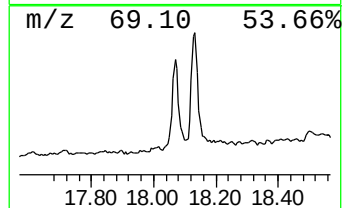
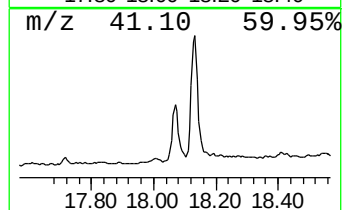
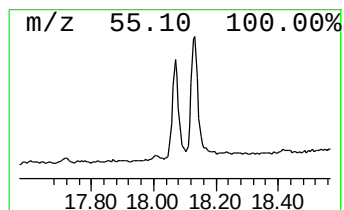
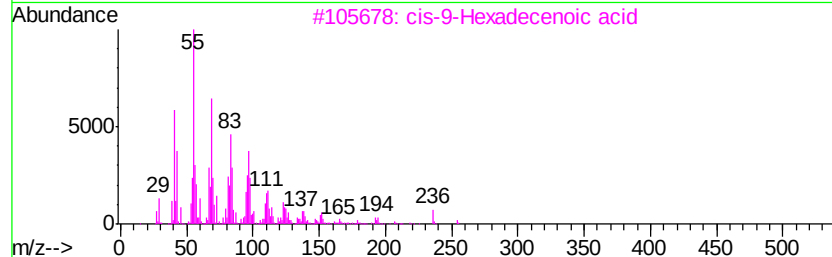
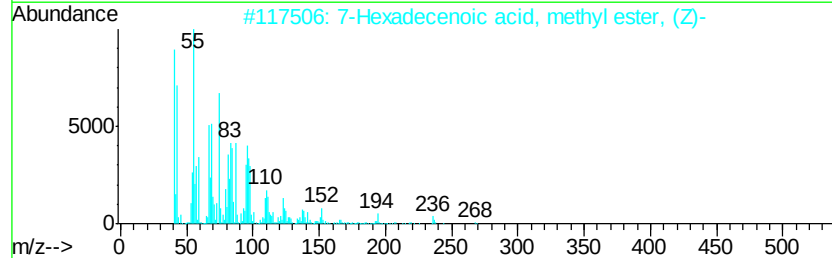
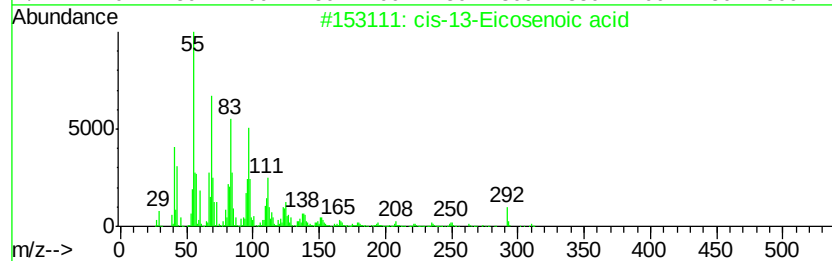
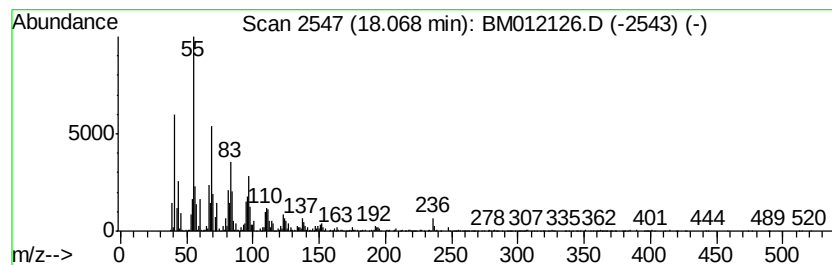
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 cis-13-Eicosenoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.07	7.39 ng/ul	1449430	Phenanthrene-d10	17.25

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	cis-13-Eicosenoic acid	310	C20H38O2	017735-94-3	97
2		7-Hexadecenoic acid, methyl este...	268	C17H32O2	056875-67-3	93
3		cis-9-Hexadecenoic acid	254	C16H30O2	1000333-19-5	90
4		cis-Vaccenic acid	282	C18H34O2	000506-17-2	87
5		cis-10-Nonadecenoic acid	296	C19H36O2	073033-09-7	83



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM101217\
Data File : BM012126.D
Acq On : 13 Oct 2017 07:40
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Sample : I5568-02
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
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ClientSampleId :
B-51(1-2)-092817

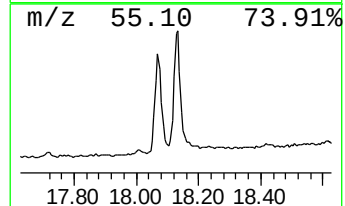
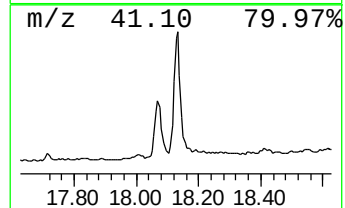
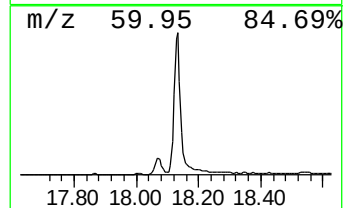
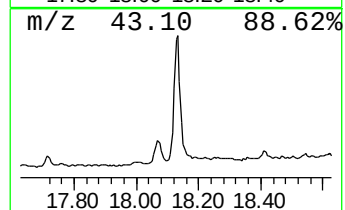
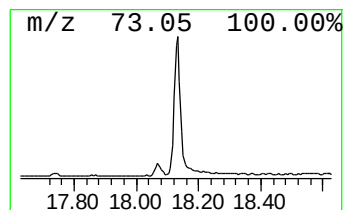
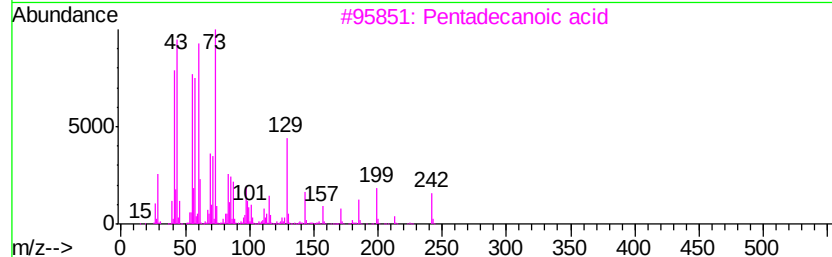
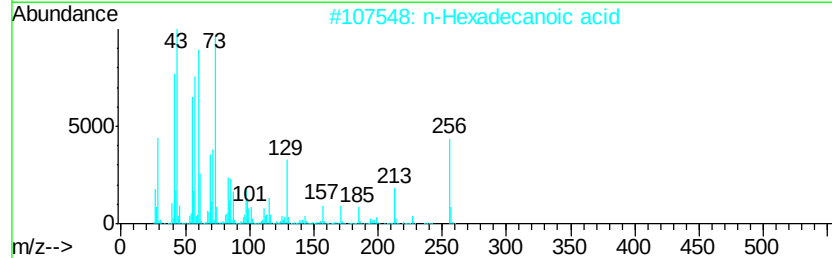
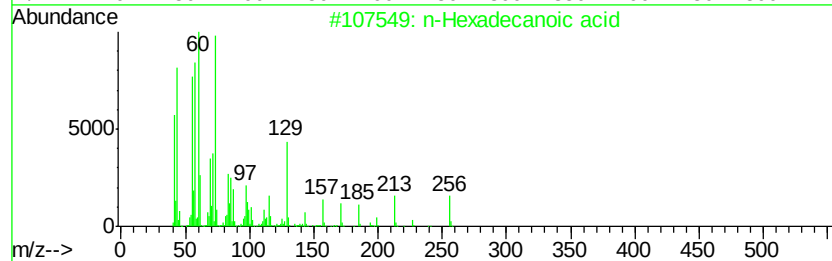
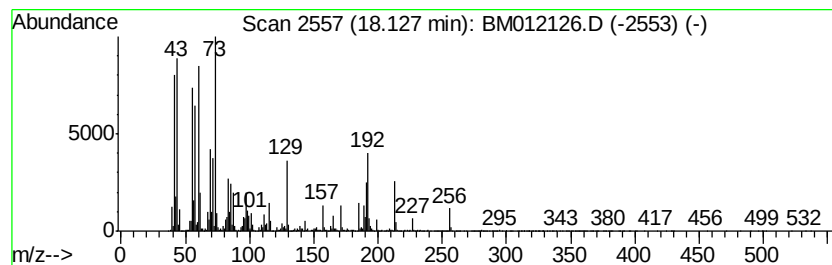
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.13	18.13 ng/ul	3554630	Phenanthrene-d10	17.25

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	90
4		Tridecanoic acid	214	C13H26O2	000638-53-9	86
5		Tetradecanoic acid	228	C14H28O2	000544-63-8	83



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM101217\
Data File : BM012126.D
Acq On : 13 Oct 2017 07:40
Operator : SJ/JU
Sample : I5568-02
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ClientSampleId :
B-51(1-2)-092817

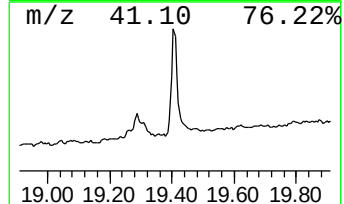
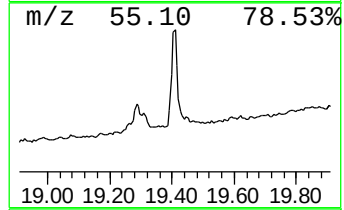
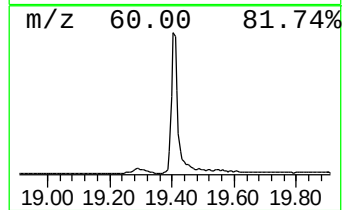
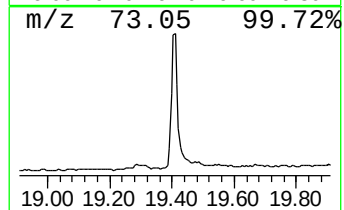
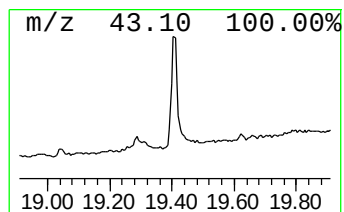
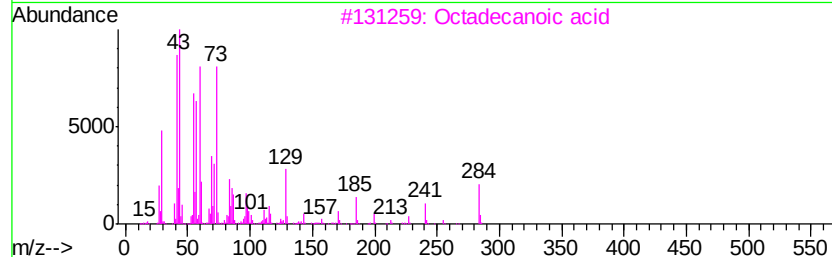
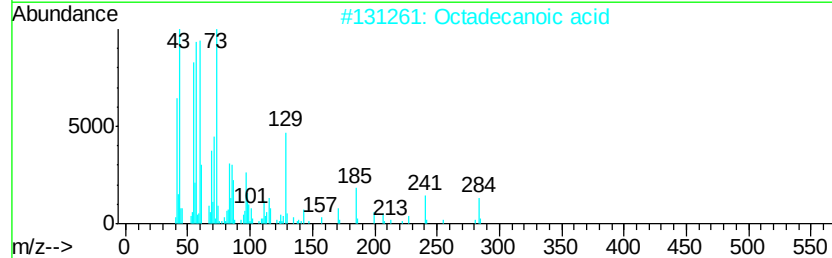
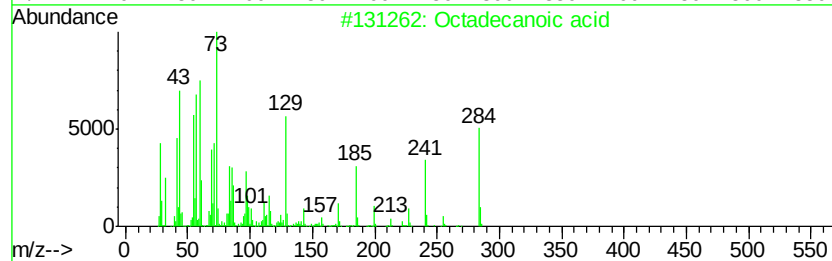
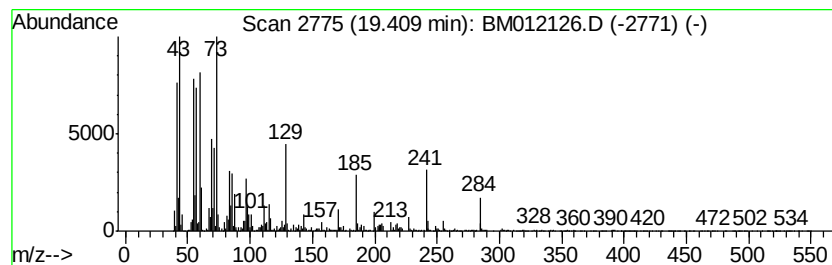
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Octadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.41	14.92 ng/ul	3289490	Chrysene-d12	21.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		Octadecanoic acid	284	C18H36O2	000057-11-4	99
3		Octadecanoic acid	284	C18H36O2	000057-11-4	95
4		Octadecanoic acid	284	C18H36O2	000057-11-4	93
5		Pentadecanoic acid	242	C15H30O2	001002-84-2	90



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM101217\
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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
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(DEL) Alkane: Str...	7.45	5.6	ng/ul	518426	1	7.85	1837820	20.0
Benzene, 1,2,3-tr...	7.49	3.5	ng/ul	324046	1	7.85	1837820	20.0
Benzene, 1-methyl...	8.39	5.0	ng/ul	464443	1	7.85	1837820	20.0
Tetracyclo[3.3.1....	8.48	7.3	ng/ul	672998	1	7.85	1837820	20.0
Benzene, 2-ethyl-...	8.95	4.2	ng/ul	389492	1	7.85	1837820	20.0
Benzene, 1-chloro...	11.48	2.8	ng/ul	354103	2	10.66	2549680	20.0
cis-13-Eicosenoic...	18.07	7.4	ng/ul	1449430	4	17.25	3921450	20.0
n-Hexadecanoic acid	18.13	18.1	ng/ul	3554630	4	17.25	3921450	20.0
Octadecanoic acid	19.41	14.9	ng/ul	3289490	5	21.44	4408810	20.0