

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_M\DATA\BM012318\
 Data File : BM013756.D
 Acq On : 24 Jan 2018 07:00
 Operator : SJ/JU
 Sample : J1220-13
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 F6A15

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-BM011018.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	3.157	25	29	35	rVB2	45085	56315	1.33%	0.089%
2	6.292	554	562	570	rBV	259033	399925	9.42%	0.633%
3	6.787	640	646	660	rBV2	848296	1542593	36.35%	2.442%
4	6.951	667	674	682	rBV	757526	1146437	27.01%	1.815%
5	7.139	699	706	719	rVB	1041303	1634641	38.52%	2.588%
6	7.604	779	785	796	rVB	887155	1367242	32.22%	2.165%
7	8.316	900	906	924	rVB	827830	1346154	31.72%	2.131%
8	8.704	965	972	976	rBV	142148	257744	6.07%	0.408%
9	8.763	976	982	989	rVB	978749	1651742	38.92%	2.615%
10	8.886	999	1003	1011	rVB2	49896	79410	1.87%	0.126%
11	9.163	1045	1050	1054	rBV3	30739	48843	1.15%	0.077%
12	9.475	1096	1103	1113	rBV	865144	1456922	34.33%	2.307%
13	9.657	1129	1134	1142	rBV4	31498	80866	1.91%	0.128%
14	10.010	1187	1194	1205	rBV	1152366	2023760	47.69%	3.204%
15	10.380	1250	1257	1264	rBV	1141401	1831996	43.17%	2.900%
16	10.527	1275	1282	1295	rBV	681019	1258487	29.65%	1.992%
17	11.339	1414	1420	1427	rBV5	19977	50532	1.19%	0.080%
18	13.080	1711	1716	1719	rBV5	30372	43591	1.03%	0.069%
19	13.121	1719	1723	1732	rVB7	54083	81853	1.93%	0.130%
20	13.386	1762	1768	1771	rBV6	25771	53795	1.27%	0.085%
21	13.563	1795	1798	1803	rBV4	47775	64243	1.51%	0.102%
22	13.674	1810	1817	1821	rBV	2031807	2763752	65.12%	4.375%
23	13.721	1821	1825	1833	rVV	535564	771282	18.17%	1.221%
24	13.786	1833	1836	1840	rVB5	37200	51164	1.21%	0.081%
25	13.945	1853	1863	1871	rBV	2344028	3625230	85.42%	5.739%
26	14.098	1882	1889	1893	rBV6	27454	44147	1.04%	0.070%
27	14.163	1893	1900	1903	rVV2	101572	137066	3.23%	0.217%
28	14.210	1903	1908	1912	rVV	782156	1134715	26.74%	1.796%
29	14.257	1912	1916	1924	rVB3	1410339	2159650	50.89%	3.419%
30	14.486	1949	1955	1969	rBV2	463679	999548	23.55%	1.582%
31	14.786	2001	2006	2014	rVB9	23288	45873	1.08%	0.073%
32	15.068	2049	2054	2056	rBV2	42418	57539	1.36%	0.091%
33	15.121	2059	2063	2069	rVB	128561	167395	3.94%	0.265%
34	15.257	2080	2086	2094	rBV	2869290	3920968	92.39%	6.208%

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Integration Parameters: LSCINT.P

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35	15.392	2103	2109	2119	rBV	616921	939468	22.14%	1.487%
36	15.504	2124	2128	2136	rVB5	45052	107472	2.53%	0.170%
37	15.627	2145	2149	2154	rVB3	73108	102101	2.41%	0.162%
38	15.739	2162	2168	2172	rVB7	34788	58656	1.38%	0.093%
39	15.898	2191	2195	2199	rBV	70105	93014	2.19%	0.147%
40	15.986	2205	2210	2216	rBV	121516	180073	4.24%	0.285%
41	16.774	2341	2344	2348	rBV2	74362	89717	2.11%	0.142%
42	16.827	2350	2353	2357	rVB3	36108	43627	1.03%	0.069%
43	16.915	2364	2368	2375	rBV7	42294	67394	1.59%	0.107%
44	17.009	2376	2384	2390	rVV	1751981	2412065	56.84%	3.819%
45	17.109	2395	2401	2412	rVB	2850959	3963626	93.39%	6.275%
46	17.403	2449	2451	2462	rVB	20594	51596	1.22%	0.082%
47	17.515	2467	2470	2475	rVB3	43849	50639	1.19%	0.080%
48	17.798	2514	2518	2523	rBV6	56855	80812	1.90%	0.128%
49	17.856	2524	2528	2534	rVV7	49972	78850	1.86%	0.125%
50	17.915	2534	2538	2553	rVV	573131	859943	20.26%	1.361%
51	18.209	2584	2588	2592	rBV6	47636	74572	1.76%	0.118%
52	19.050	2728	2731	2734	rBV3	65249	96651	2.28%	0.153%
53	19.086	2734	2737	2740	rBV5	37614	60656	1.43%	0.096%
54	19.203	2753	2757	2763	rBV2	280049	387208	9.12%	0.613%
55	19.409	2786	2792	2798	rBV	2992049	4218551	99.40%	6.679%
56	19.468	2800	2802	2807	rVB6	48333	56934	1.34%	0.090%
57	19.815	2855	2861	2866	rBV2	87002	134958	3.18%	0.214%
58	19.962	2883	2886	2889	rVB3	67715	89252	2.10%	0.141%
59	20.003	2890	2893	2899	rVB7	111623	153620	3.62%	0.243%
60	20.339	2946	2950	2951	rBV4	104049	119430	2.81%	0.189%
61	20.609	2994	2996	2999	rBV3	63306	67103	1.58%	0.106%
62	20.768	3019	3023	3029	rVB	2707372	3037658	71.58%	4.809%
63	20.933	3047	3051	3058	rVB	421688	559935	13.19%	0.886%
64	21.209	3093	3098	3105	rBV	1991619	2512773	59.21%	3.978%
65	21.815	3198	3201	3210	rVB	552172	813887	19.18%	1.289%
66	22.744	3356	3359	3362	rVB2	170190	186621	4.40%	0.295%
67	22.791	3364	3367	3373	rVB4	220795	339530	8.00%	0.538%
68	23.274	3442	3449	3458	rVB	2363784	4243956	100.00%	6.719%
69	23.409	3466	3472	3479	rVB	1346833	2453679	57.82%	3.885%
70	23.921	3555	3559	3565	rVB3	227738	374453	8.8	

LSC Area Percent Report

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Integrator: RTE
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Sampling : 1 Min Area: 1 % of largest Peak
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Stop Thrs : 0 Peak Location: TOP

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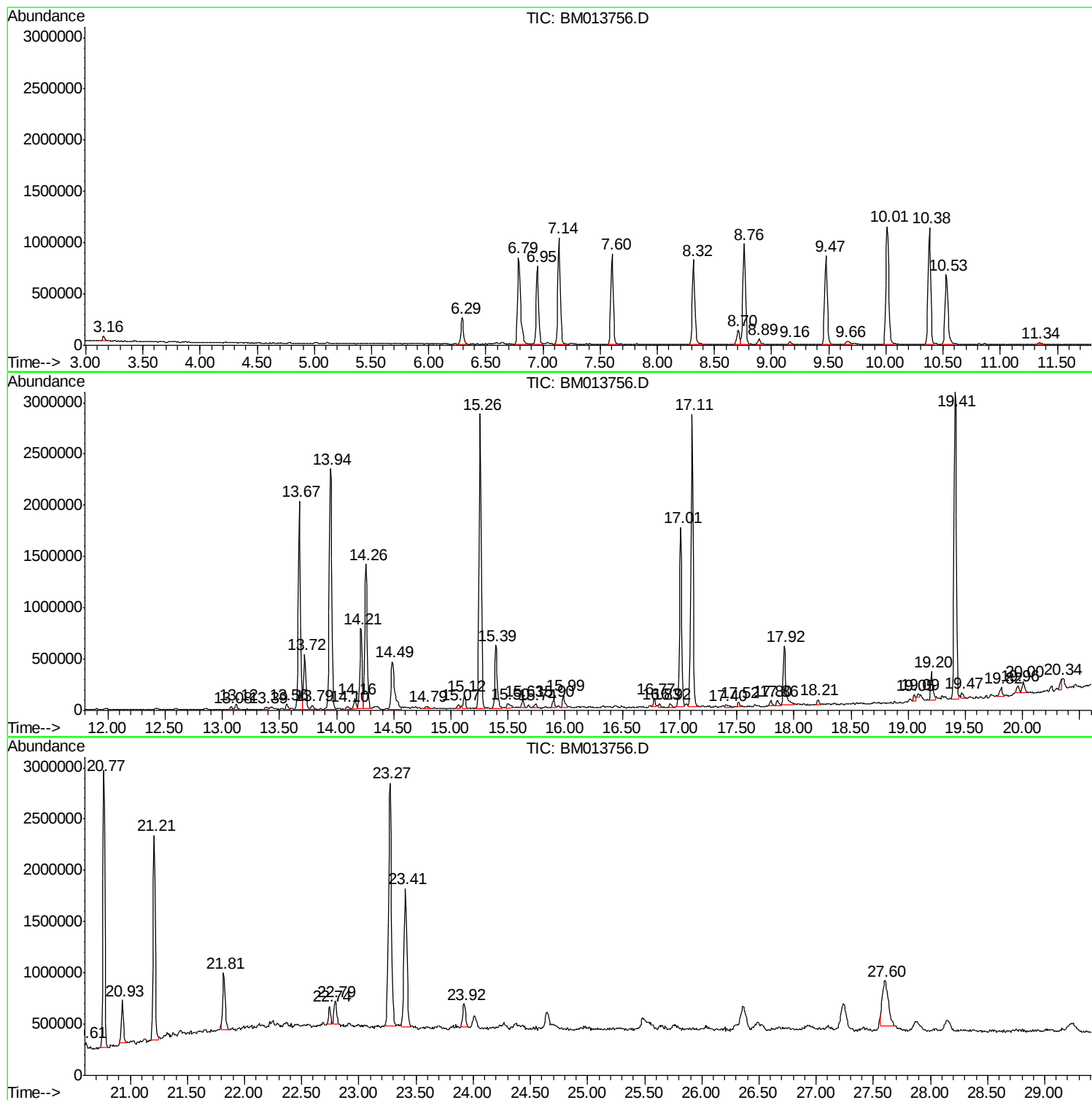
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM011018.M
Quant Title : SVOA CALIBRATION

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



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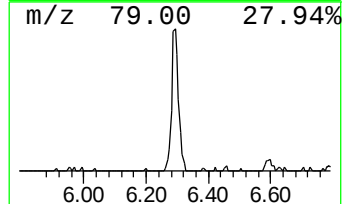
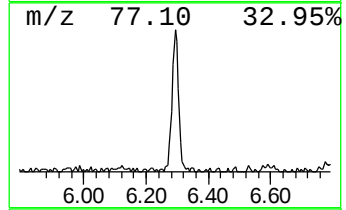
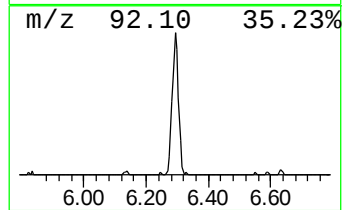
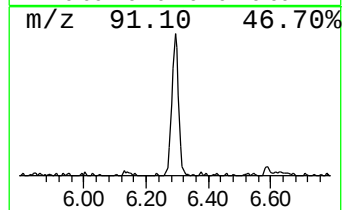
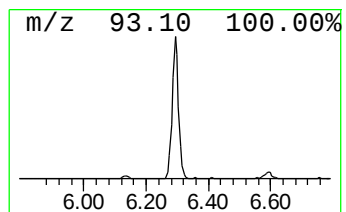
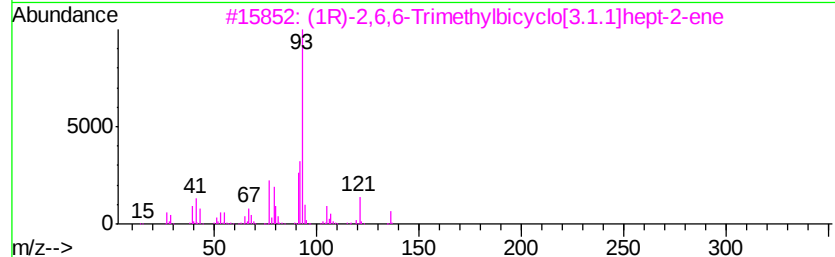
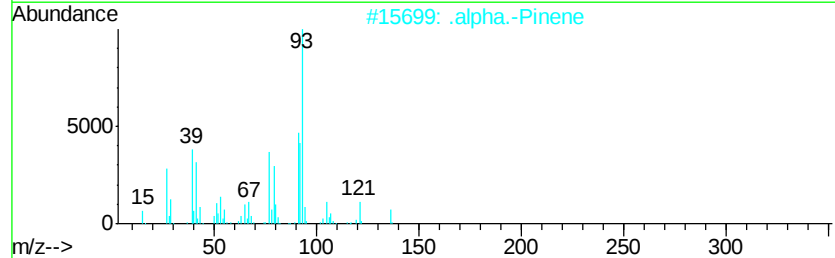
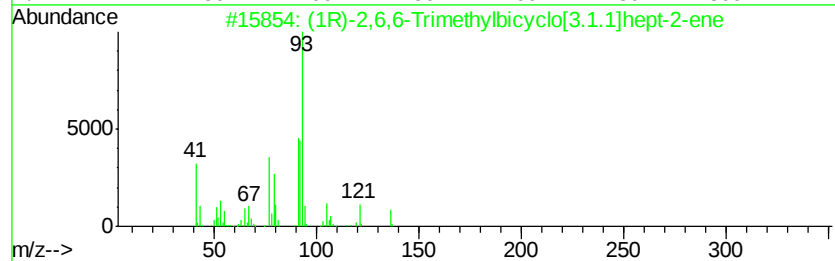
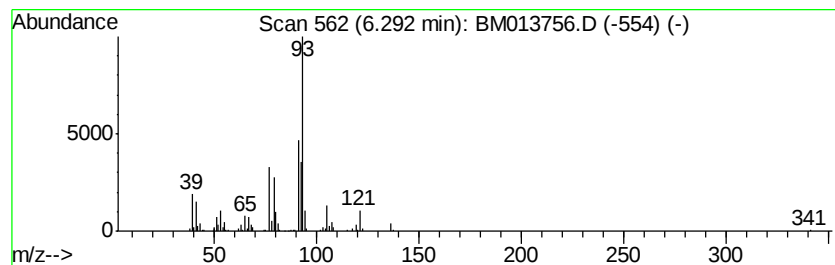
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM011018.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 (1R)-2,6,6-Trimethylbicyclo... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.29	5.85 ng/ul	399925	1,4-Dichlorobenzene-d4	7.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(1R)-2,6,6-Trimethylbicyclo[3.1.1]hept-2-ene	136	C10H16	007785-70-8	95
2		.alpha.-Pinene	136	C10H16	000080-56-8	95
3		(1R)-2,6,6-Trimethylbicyclo[3.1.1]hept-2-ene	136	C10H16	007785-70-8	91
4		Bicyclo[3.1.1]hept-2-ene, 3,6,6-trimethyl-	136	C10H16	004889-83-2	91
5		(+)-3-Carene	136	C10H16	000498-15-7	91



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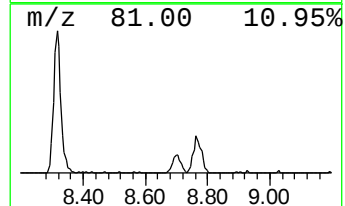
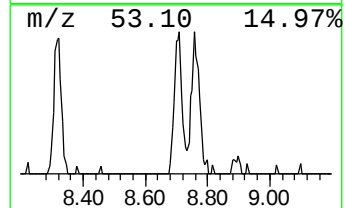
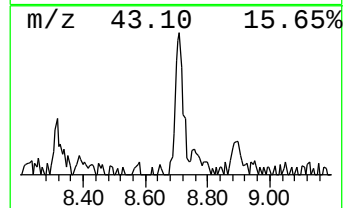
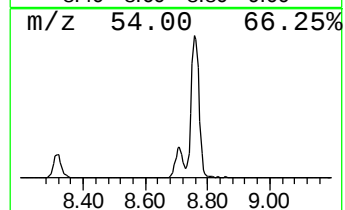
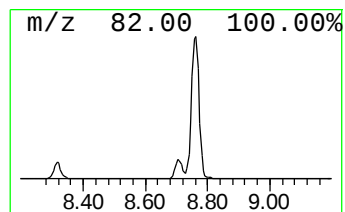
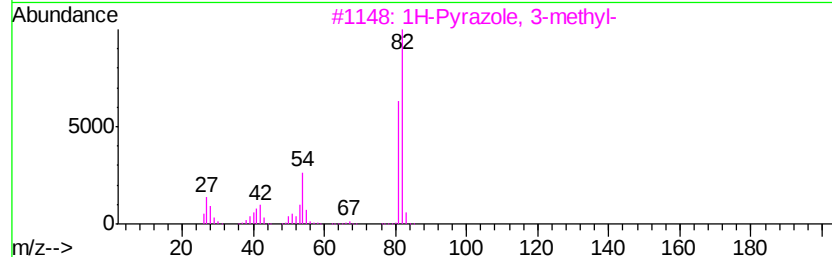
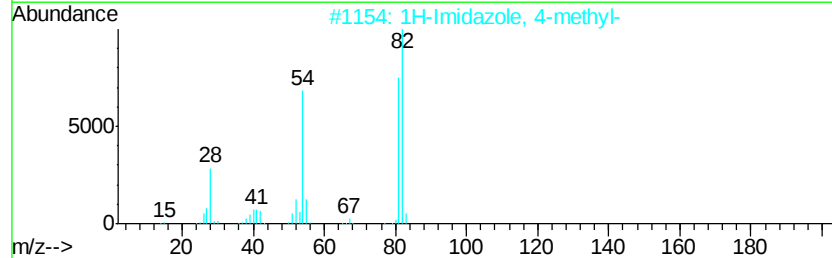
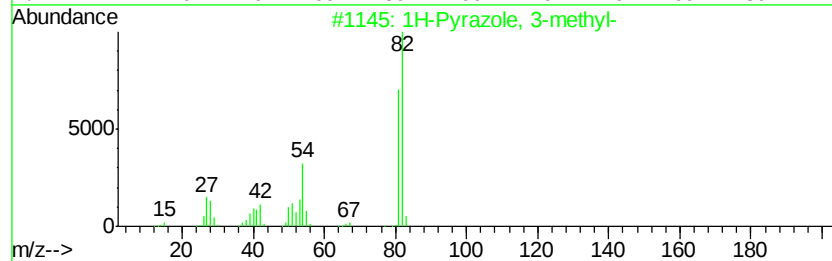
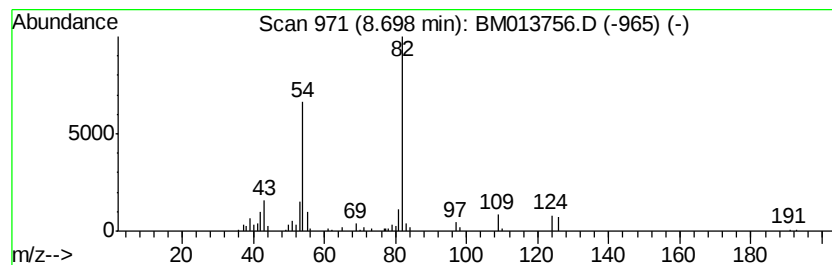
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM011018.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 1H-Pyrazole, 3-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.70	3.77 ng/ul	257744	1,4-Dichlorobenzene-d4	7.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Pyrazole, 3-methyl-	82	C4H6N2	001453-58-3	59
2		1H-Imidazole, 4-methyl-	82	C4H6N2	000822-36-6	56
3		1H-Pyrazole, 3-methyl-	82	C4H6N2	001453-58-3	53
4		2(5H)-Furanone, 5-(2-hydroxyethy...	126	C6H6O3	089291-42-9	50
5		1H-Pyrazole, 1-methyl-	82	C4H6N2	000930-36-9	50



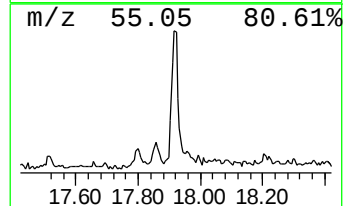
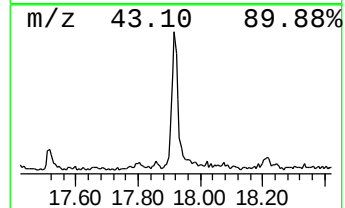
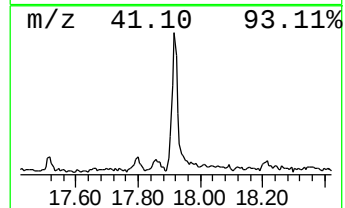
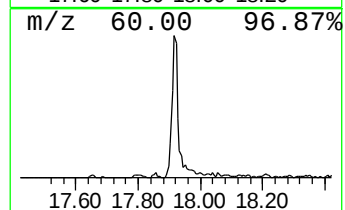
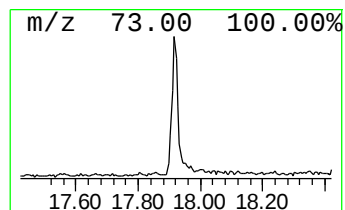
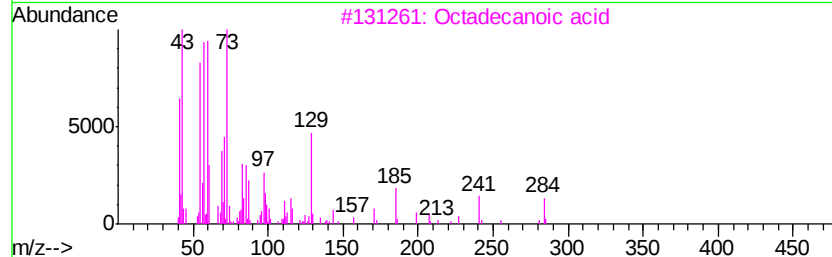
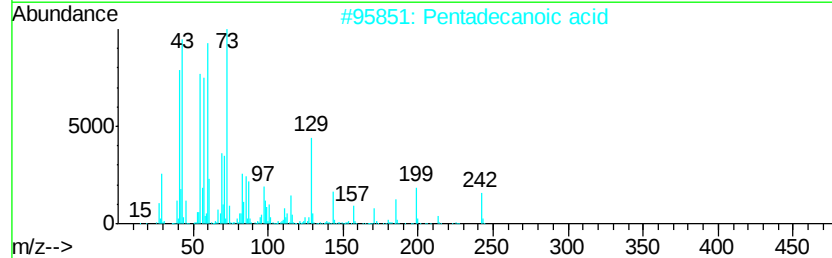
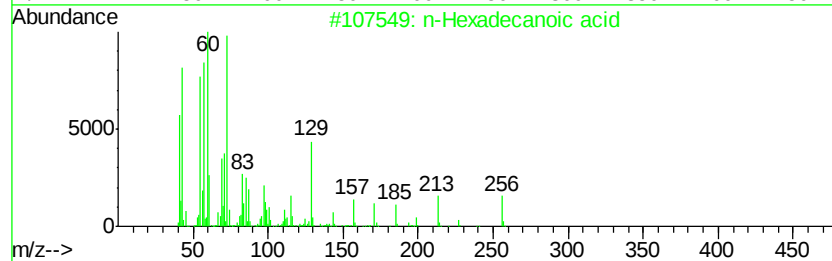
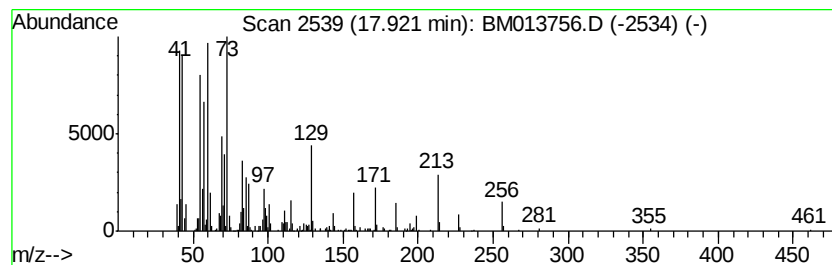
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 n-Hexadecanoic acid Concentration Rank 3



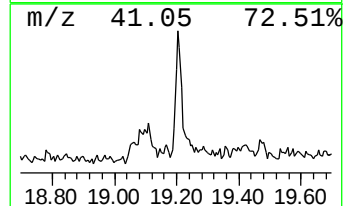
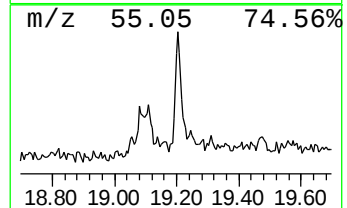
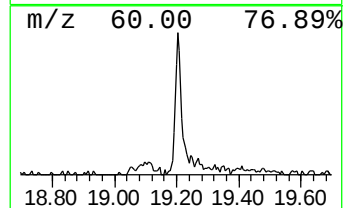
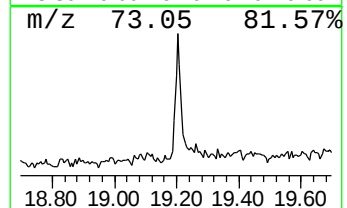
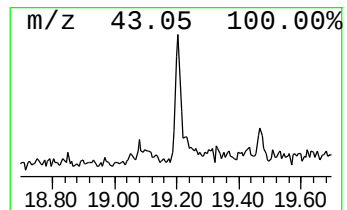
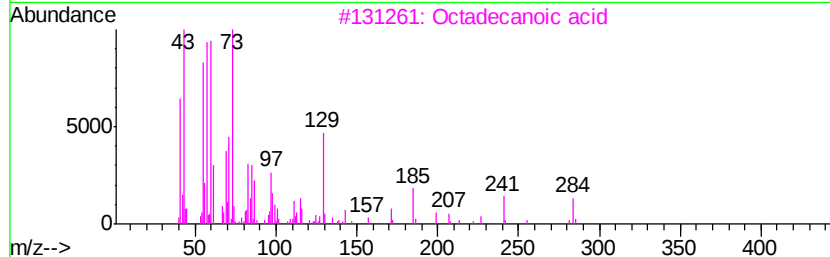
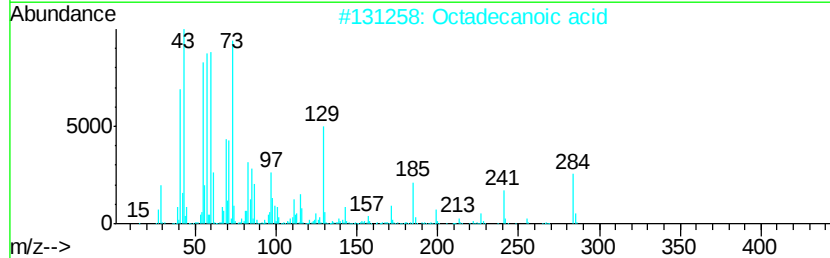
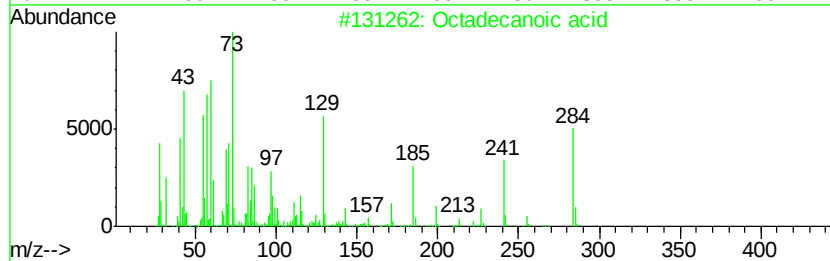
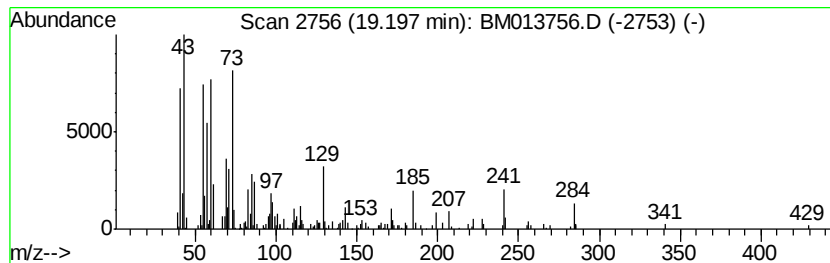
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Peak Number 4 Octadecanoic acid Concentration Rank 8



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Operator : SJ/JU
Sample : J1220-13
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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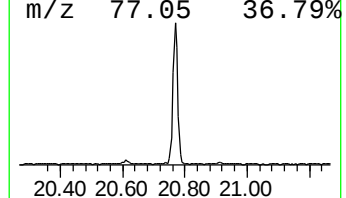
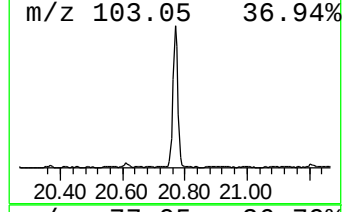
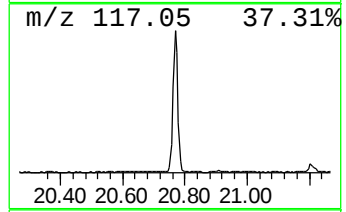
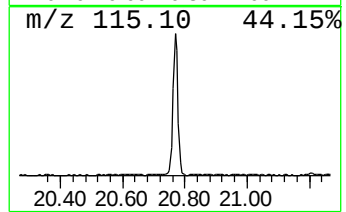
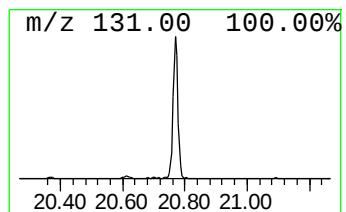
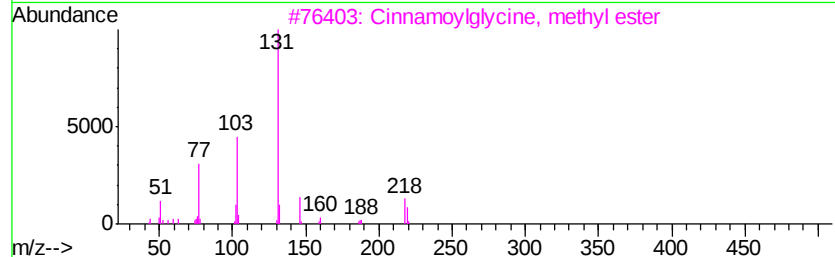
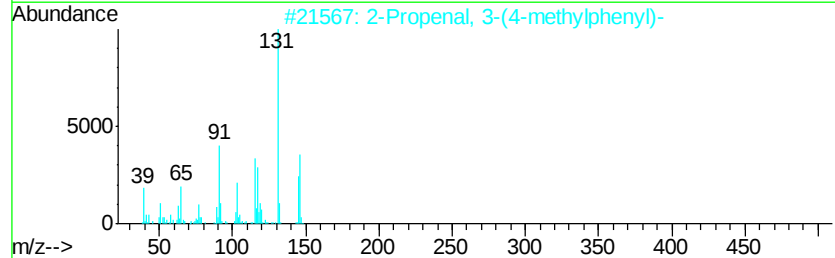
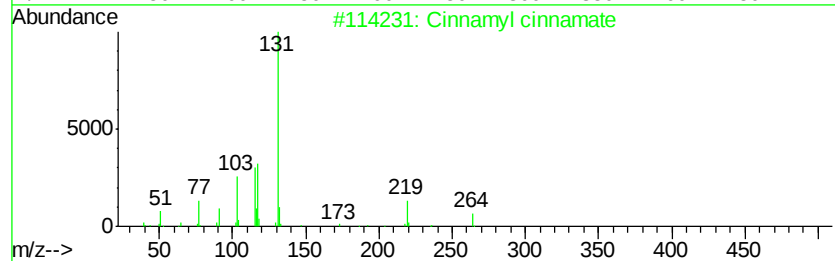
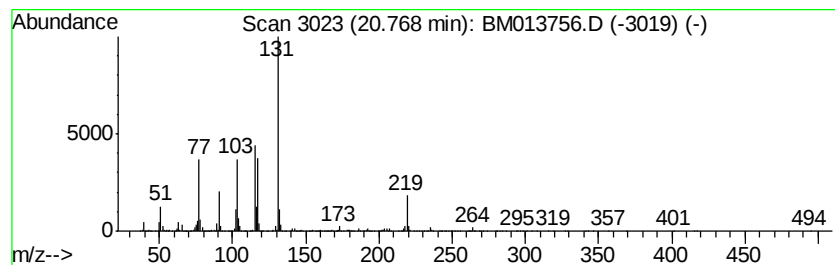
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Cinnamyl cinnamate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.77	24.18 ng/ul	3037660	Chrysene-d12	21.21

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cinnamyl cinnamate	264	C18H16O2	000122-69-0	64
2		2-Propenal, 3-(4-methylphenyl)-	146	C10H10O	001504-75-2	56
3		Cinnamoylglycine, methyl ester	219	C12H13NO3	040778-04-9	50
4		p-Vinylbenzohydrazide	162	C9H10N2O	1000244-74-9	47
5		N-(p-Vinylbenzoyl)-l-alanine	219	C12H13NO3	139184-60-4	47



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Data File : BM013756.D
Acq On : 24 Jan 2018 07:00
Operator : SJ/JU
Sample : J1220-13
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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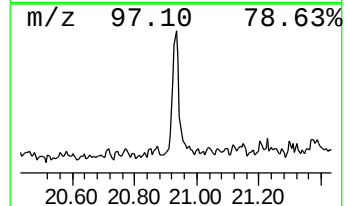
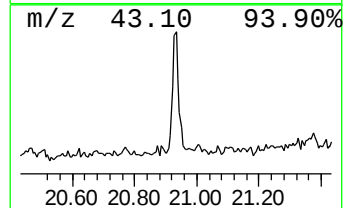
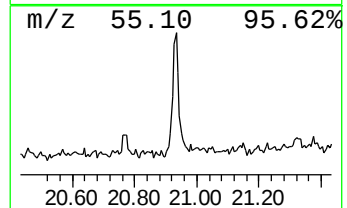
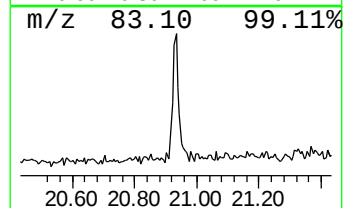
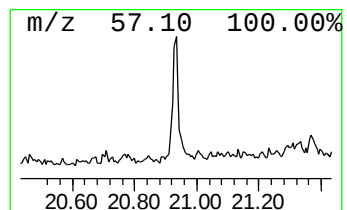
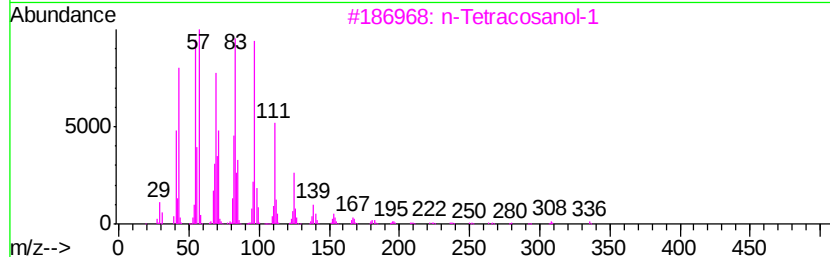
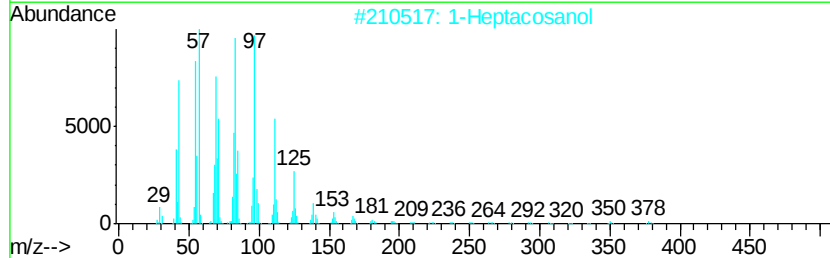
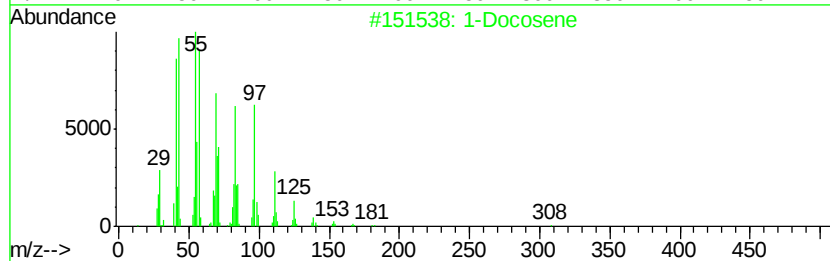
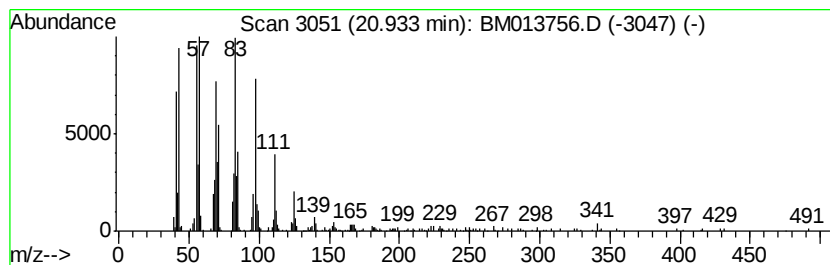
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Cyclic20.93 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.93	4.46 ng/ul	559935	Chrysene-d12	21.21

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Docosene	308	C22H44	001599-67-3	91
2		1-Heptacosanol	396	C27H56O	002004-39-9	91
3		n-Tetracosanol-1	354	C24H50O	000506-51-4	91
4		Nonadecyl trifluoroacetate	380	C21H39F3O2	1000351-76-3	90
5		Cyclopentadecane	210	C15H30	000295-48-7	87



Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA_M\DATA\BM012318\
Data File : BM013756.D
Acq On : 24 Jan 2018 07:00
Operator : SJ/JU
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Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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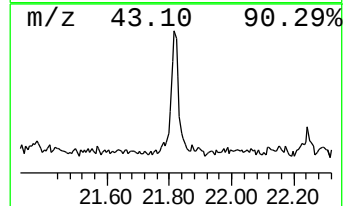
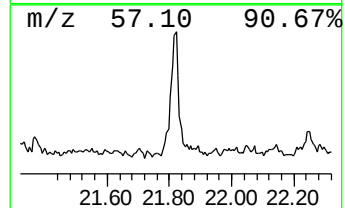
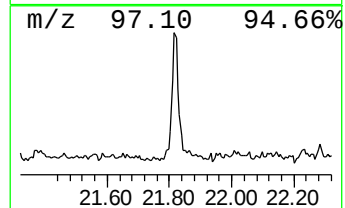
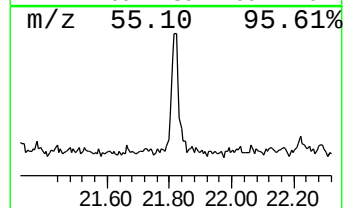
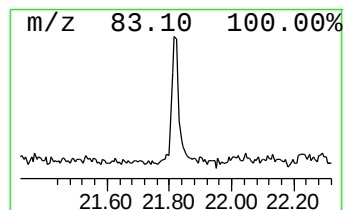
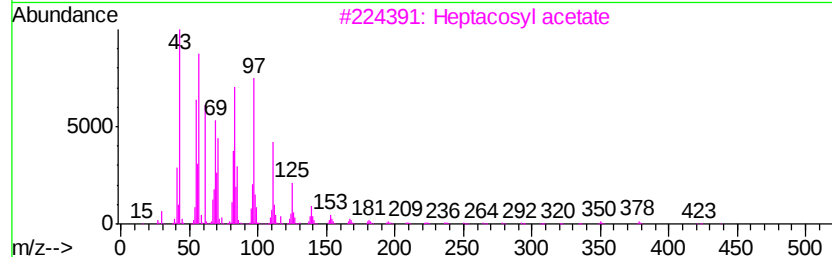
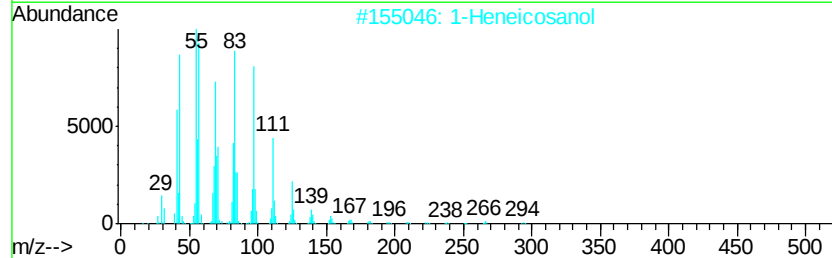
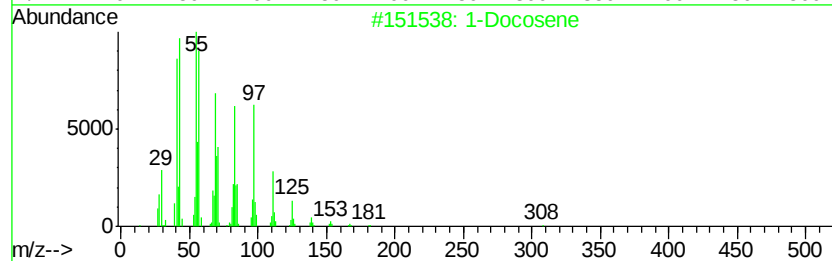
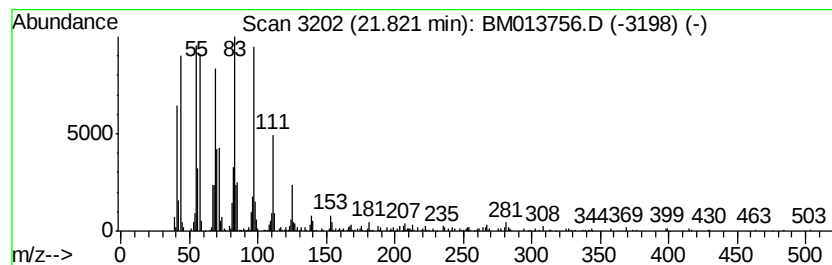
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Cyclic21.82 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.82	6.48 ng/ul	813887	Chrysene-d12	21.21

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Docosene	308	C22H44	001599-67-3	94
2			1-Heneicosanol	312	C21H44O	015594-90-8	94
3			Heptacosyl acetate	438	C29H58O2	1000351-78-2	93
4			n-Tetracosanol-1	354	C24H50O	000506-51-4	93
5			Cyclotetracosane	336	C24H48	000297-03-0	93



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Data File : BM013756.D
Acq On : 24 Jan 2018 07:00
Operator : SJ/JU
Sample : J1220-13
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
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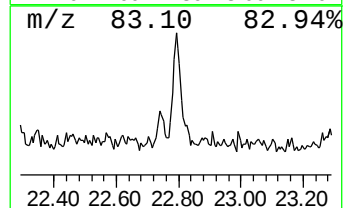
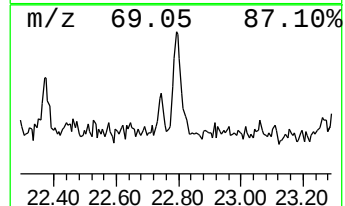
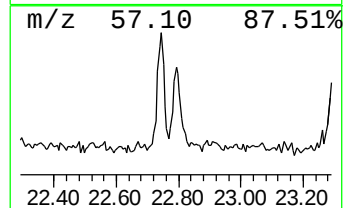
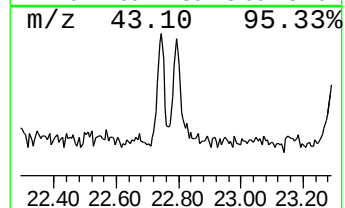
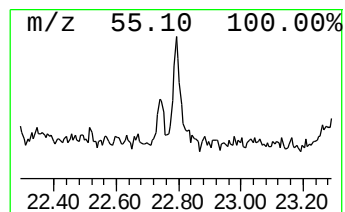
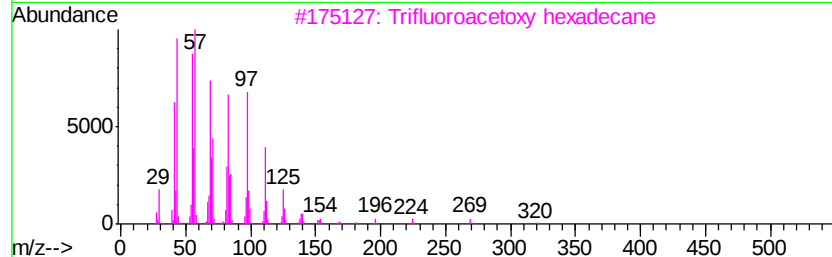
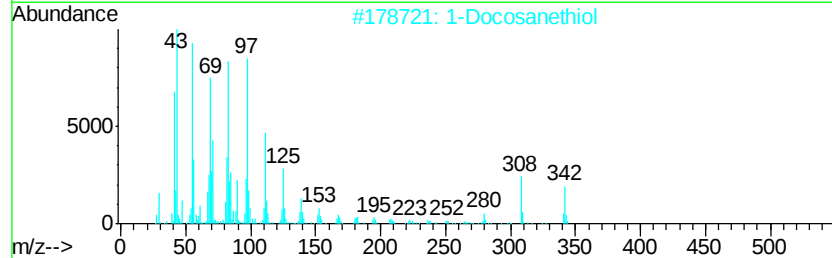
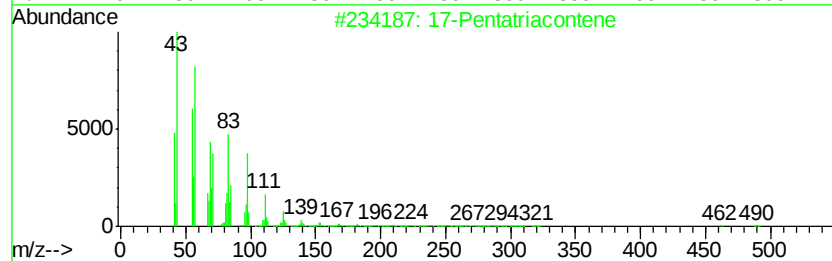
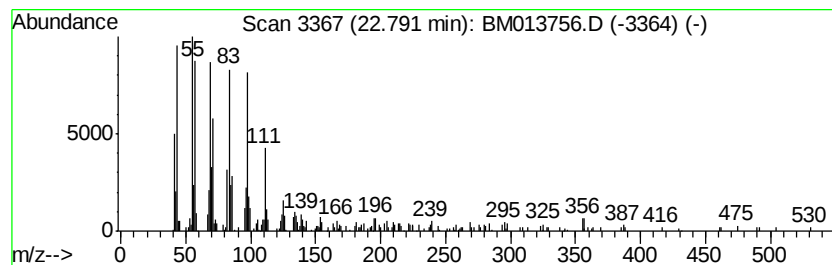
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Cyclic22.79 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.79	2.77 ng/ul	339530	Perylene-d12	23.41

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	17-Pentatriacontene	491	C35H70	006971-40-0	94
2		1-Docosanethiol	342	C22H46S	007773-83-3	91
3		Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	91
4		Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	90
5		Carbonic acid, octadecyl 2,2,2-t...	444	C21H39Cl3O3	1000314-56-3	89



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_M\DATA\BM012318\
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Acq On : 24 Jan 2018 07:00
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Misc :
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Instrument :
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ClientSampleId :
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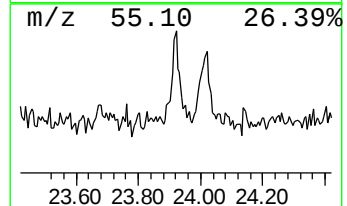
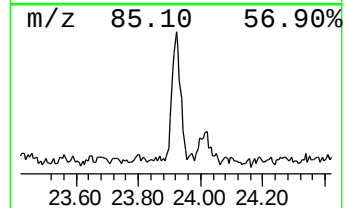
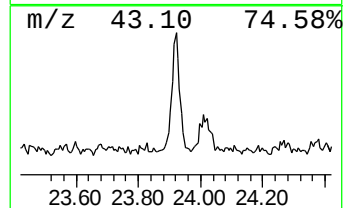
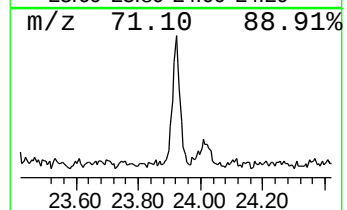
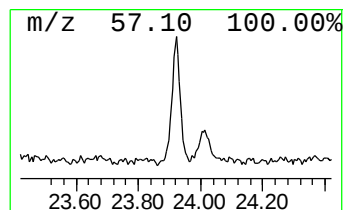
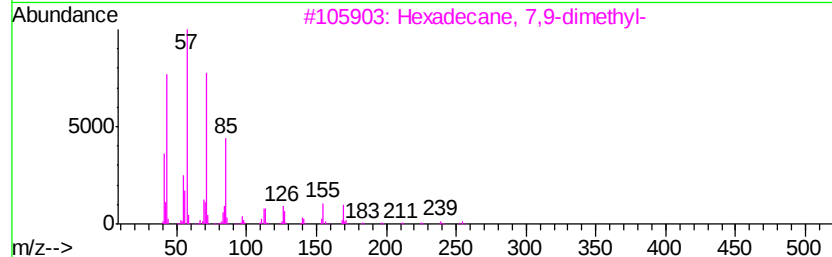
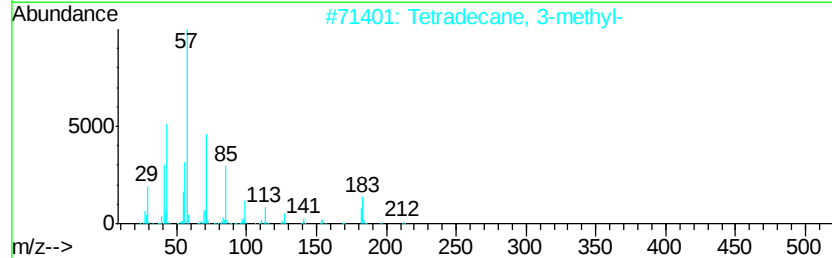
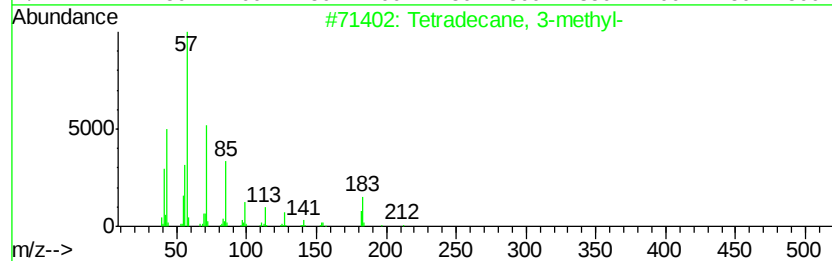
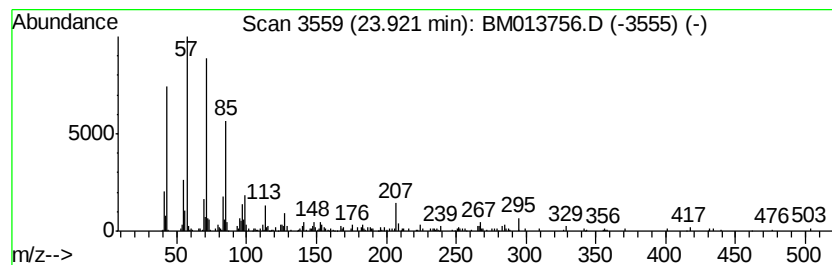
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.92	3.05 ng/ul	374453	Perylene-d12	23.41

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetradecane, 3-methyl-	212	C15H32	018435-22-8	76
2		Tetradecane, 3-methyl-	212	C15H32	018435-22-8	68
3		Hexadecane, 7,9-dimethyl-	254	C18H38	021164-95-4	68
4		Pentadecane, 4-methyl-	226	C16H34	002801-87-8	68
5		Hexadecane	226	C16H34	000544-76-3	64



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_M\DATA\BM012318\
Data File : BM013756.D
Acq On : 24 Jan 2018 07:00
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Misc :
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Instrument :
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ClientSampled :
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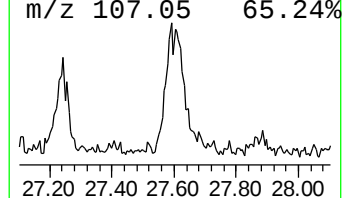
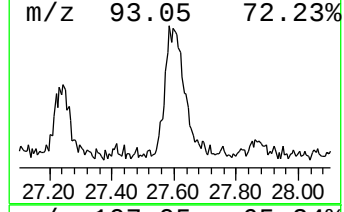
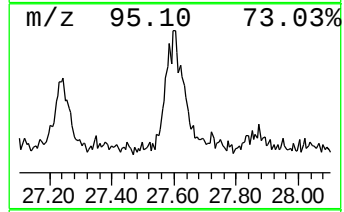
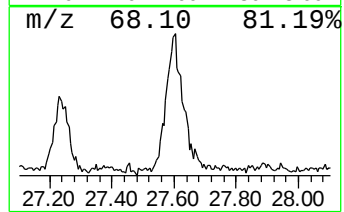
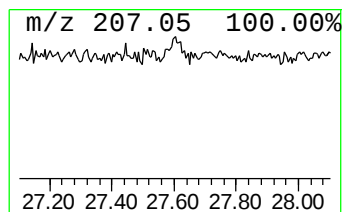
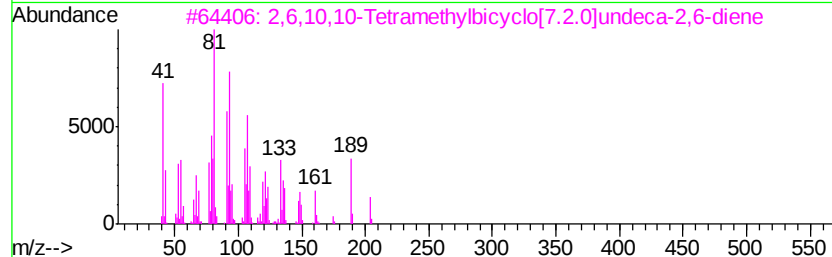
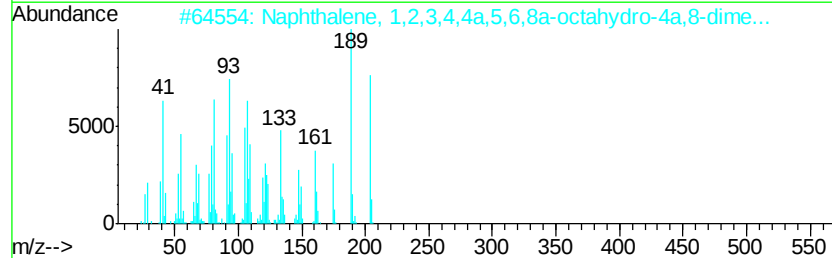
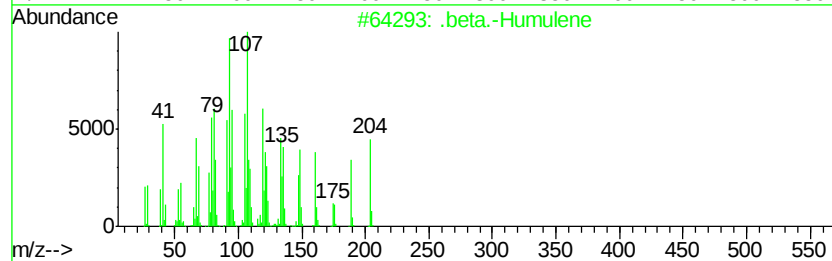
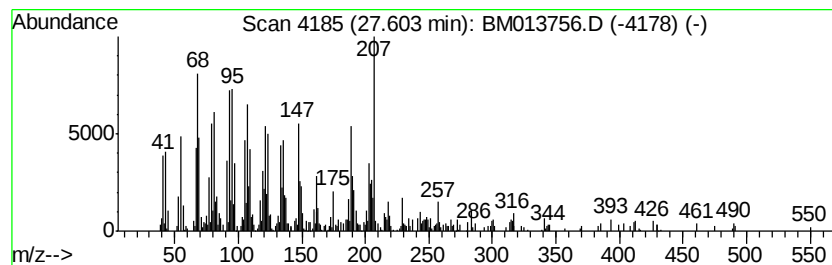
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 .beta.-Humulene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.60	13.44 ng/ul	1648960	Perylene-d12	23.41

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	.beta.-Humulene	204	C15H24	000116-04-1	90
2		Naphthalene, 1,2,3,4,4a,5,6,8a-o...	204	C15H24	000473-13-2	83
3		2,6,10,10-Tetramethylbicyclo[7.2...	204	C15H24	136296-37-2	56
4		Cyclohexane, 1-ethenyl-1-methyl-...	204	C15H24	000515-13-9	43
5		.alpha.-Bulnesene	204	C15H24	1000374-19-9	42



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_M\DATA\BM012318\
 Data File : BM013756.D
 Acq On : 24 Jan 2018 07:00
 Operator : SJ/JU
 Sample : J1220-13
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM011018.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
(1R)-2,6,6-Trimet...	6.29	5.8	ng/ul	399925	1	7.60	1367240	20.0
1H-Pyrazole, 3-me...	8.70	3.8	ng/ul	257744	1	7.60	1367240	20.0
n-Hexadecanoic acid	17.92	7.1	ng/ul	859943	4	17.01	2412070	20.0
Octadecanoic acid	19.20	3.1	ng/ul	387208	5	21.21	2512770	20.0
Cinnamyl cinnamate	20.77	24.2	ng/ul	3037660	5	21.21	2512770	20.0
(DEL) Alkane: Cyc...	20.93	4.5	ng/ul	559935	5	21.21	2512770	20.0
(DEL) Alkane: Cyc...	21.82	6.5	ng/ul	813887	5	21.21	2512770	20.0
(DEL) Alkane: Cyc...	22.79	2.8	ng/ul	339530	6	23.41	2453680	20.0
(DEL) Alkane: Str...	23.92	3.0	ng/ul	374453	6	23.41	2453680	20.0
.beta.-Humulene	27.60	13.4	ng/ul	1648960	6	23.41	2453680	20.0