

Data Path : Z:\HPCHEM1\BNA M\DATA\BM050318\
 Data File : BM014950.D
 Acq On : 03 May 2018 19:28
 Operator : SJ/JU
 Sample : J2554-16
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AG4

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-BM041918.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	3.334	4	8	15	rVB	2415445	3078007	57.26%	4.033%
2	3.640	56	60	68	rVB	80391	109490	2.04%	0.143%
3	4.228	155	160	166	rVB	66433	95368	1.77%	0.125%
4	5.475	367	372	386	rBV	364287	560667	10.43%	0.735%
5	7.134	648	654	662	rVB	282134	458543	8.53%	0.601%
6	7.616	729	736	750	rBV2	1275259	2396476	44.59%	3.140%
7	7.792	757	766	780	rBV	1007561	1674021	31.14%	2.193%
8	7.916	780	787	796	rVB	70229	133080	2.48%	0.174%
9	8.010	796	803	811	rBV	1335928	2160089	40.19%	2.830%
10	8.492	878	885	900	rVB	1209510	1972043	36.69%	2.584%
11	8.704	915	921	930	rBV	371763	597247	11.11%	0.783%
12	9.186	996	1003	1015	rBV	1110158	1885330	35.08%	2.470%
13	9.675	1078	1086	1094	rBV	1172521	2003013	37.26%	2.625%
14	10.033	1141	1147	1153	rVB	46186	69593	1.29%	0.091%
15	10.404	1202	1210	1221	rBV	935954	1565655	29.13%	2.051%
16	10.939	1294	1301	1311	rBV	1405308	2386810	44.41%	3.127%
17	11.333	1360	1368	1382	rBV	1512561	2641374	49.14%	3.461%
18	11.457	1382	1389	1403	rBV	1188368	2137283	39.76%	2.801%
19	12.651	1587	1592	1598	rBV2	81966	122019	2.27%	0.160%
20	13.892	1796	1803	1809	rVB	55240	79326	1.48%	0.104%
21	14.416	1887	1892	1896	rBV2	45447	69975	1.30%	0.092%
22	14.486	1896	1904	1908	rVV	2333845	3170654	58.99%	4.155%
23	14.533	1908	1912	1917	rVB	579931	806329	15.00%	1.057%
24	14.798	1950	1957	1969	rBV	2599481	3938790	73.28%	5.161%
25	14.945	1973	1982	1987	rBV	200458	277465	5.16%	0.364%
26	15.039	1991	1998	2002	rBV4	40786	75863	1.41%	0.099%
27	15.098	2002	2008	2015	rVB2	1926313	2956221	55.00%	3.874%
28	15.257	2029	2035	2044	rBV	863038	1310153	24.37%	1.717%
29	15.445	2062	2067	2071	rVV5	37529	66009	1.23%	0.086%
30	15.557	2081	2086	2094	rBV3	40564	72233	1.34%	0.095%
31	15.886	2138	2142	2147	rVB	272677	331219	6.16%	0.434%
32	16.074	2167	2174	2181	rBV	3355605	4683194	87.13%	6.136%
33	16.174	2185	2191	2202	rVV	807127	1155038	21.49%	1.513%
34	16.286	2203	2210	2219	rVB2	67471	194616	3.62%	0.255%

Data Path : Z:\HPCHEM1\BNA M\DATA\BM050318\
Data File : BM014950.D
Acq On : 03 May 2018 19:28
Operator : SJ/JU
Sample : J2554-16
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0AG4

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-BM041918.M
Title : SVOA CALIBRATION

35	16.427	2229	2234	2238	rBV4	40071	67937	1.26%	0.089%
36	16.733	2281	2286	2295	rBV	217649	352193	6.55%	0.461%
37	17.333	2385	2388	2395	rVB3	52431	73618	1.37%	0.096%
38	17.521	2415	2420	2425	rBV	119612	149059	2.77%	0.195%
39	17.809	2463	2469	2475	rBV2	2224744	3205359	59.63%	4.200%
40	17.851	2475	2476	2480	rVV	62347	54963	1.02%	0.072%
41	17.909	2480	2486	2497	rVV	3120736	4625442	86.05%	6.061%
42	18.233	2537	2541	2548	rBV2	91357	158103	2.94%	0.207%
43	18.521	2585	2590	2596	rBV5	32654	67627	1.26%	0.089%
44	18.580	2596	2600	2605	rBV	108926	143113	2.66%	0.188%
45	18.639	2605	2610	2621	rBV	959808	1360044	25.30%	1.782%
46	19.633	2776	2779	2784	rBV2	55446	73399	1.37%	0.096%
47	19.739	2794	2797	2799	rBV2	59702	70919	1.32%	0.093%
48	19.815	2807	2810	2816	rVB	129926	175766	3.27%	0.230%
49	19.880	2816	2821	2831	rBV	555124	788251	14.66%	1.033%
50	20.145	2861	2866	2878	rVB	3650424	5265665	97.96%	6.900%
51	21.268	3055	3057	3065	rBV6	85801	126270	2.35%	0.165%
52	21.556	3102	3106	3114	rBV	641705	906695	16.87%	1.188%
53	21.897	3159	3164	3170	rBV	2555059	3264902	60.74%	4.278%
54	23.538	3440	3443	3449	rVB	365598	515511	9.59%	0.675%
55	24.215	3552	3558	3576	rVB	2548879	5375057	100.00%	7.043%
56	24.374	3579	3585	3597	rVB	1665912	3325690	61.87%	4.358%
57	24.909	3671	3676	3685	rVB	464749	938893	17.47%	1.230%

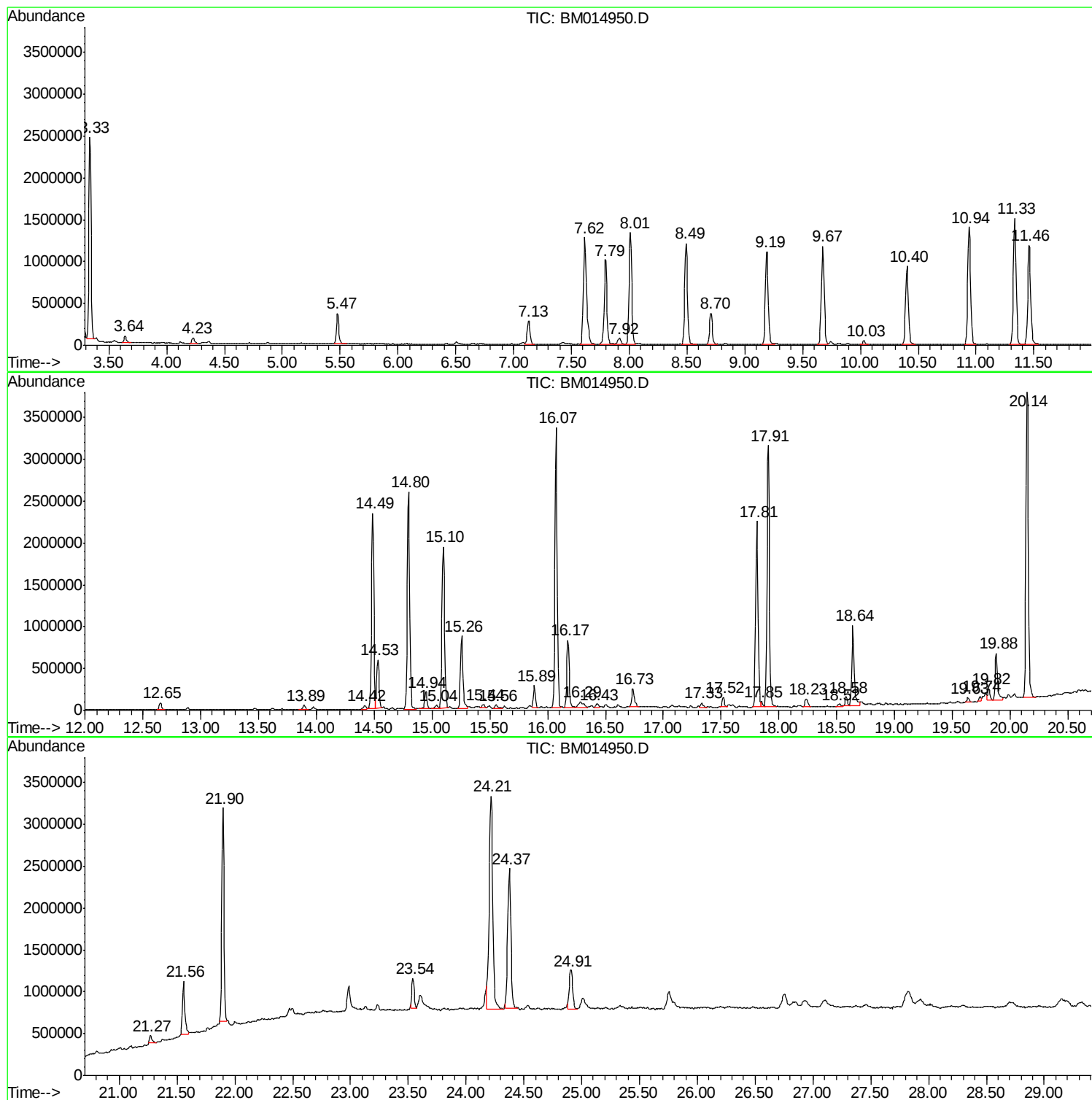
Sum of corrected areas: 76317669

Data Path : Z:\HPCHEM1\BNA M\DATA\BM050318\
Data File : BM014950.D
Acq On : 03 May 2018 19:28
Operator : SJ/JU
Sample : J2554-16
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0AG4

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM041918.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\HPCHEM1\BNA M\DATA\BM050318\
Data File : BM014950.D
Acq On : 03 May 2018 19:28
Operator : SJ/JU
Sample : J2554-16
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0AG4

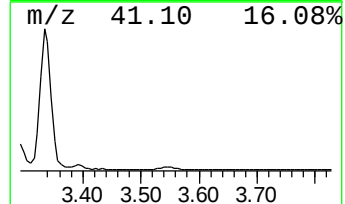
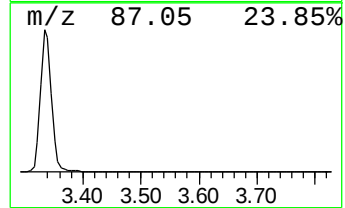
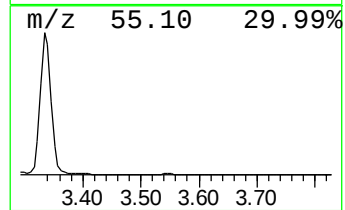
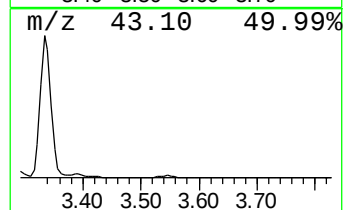
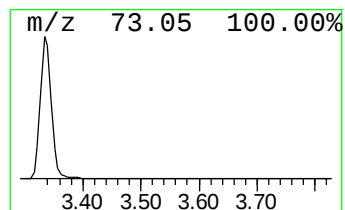
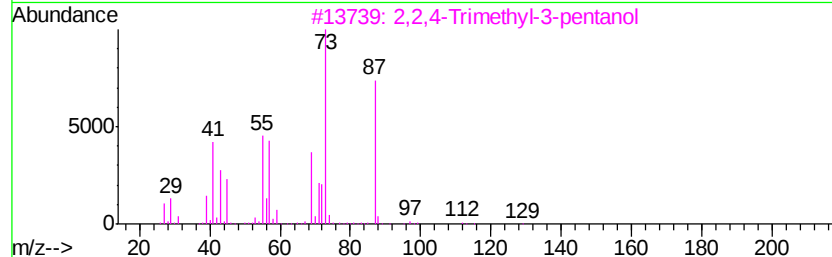
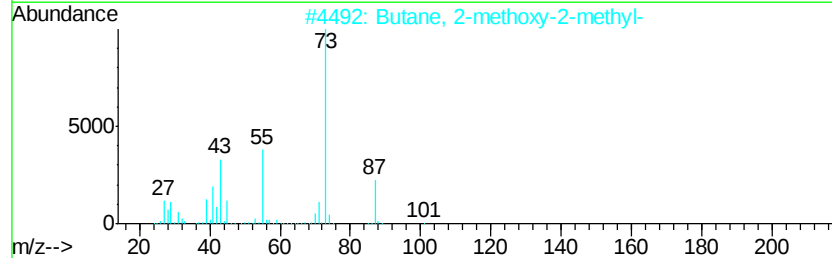
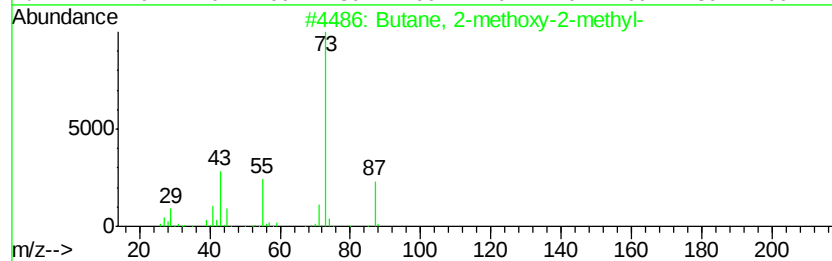
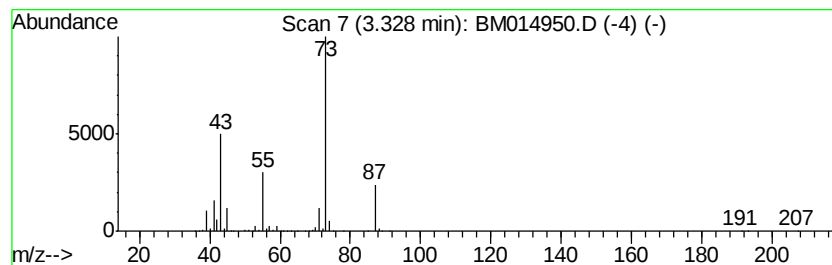
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM041918.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.33	31.22 ng/ul	3078010	1,4-Dichlorobenzene-d4	8.49

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	78
3		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
4		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
5		Silane, tetramethyl-	88	C4H12Si	000075-76-3	38



Data Path : Z:\HPCHEM1\BNA M\DATA\BM050318\
Data File : BM014950.D
Acq On : 03 May 2018 19:28
Operator : SJ/JU
Sample : J2554-16
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0AG4

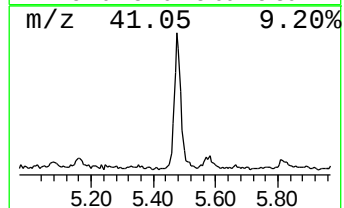
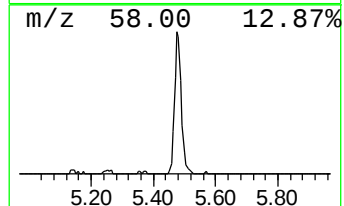
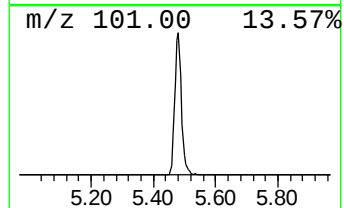
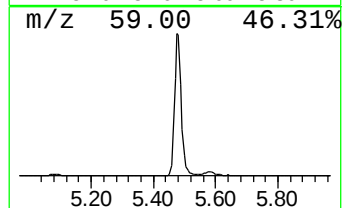
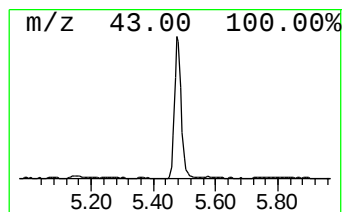
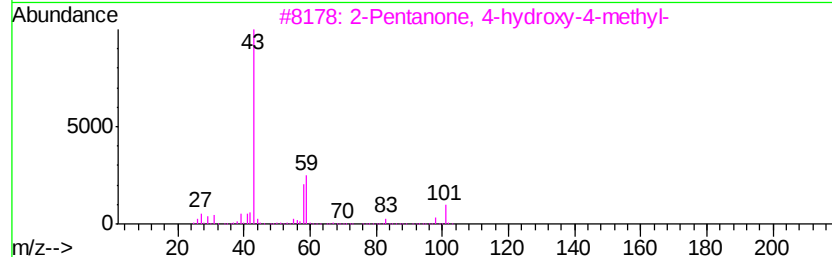
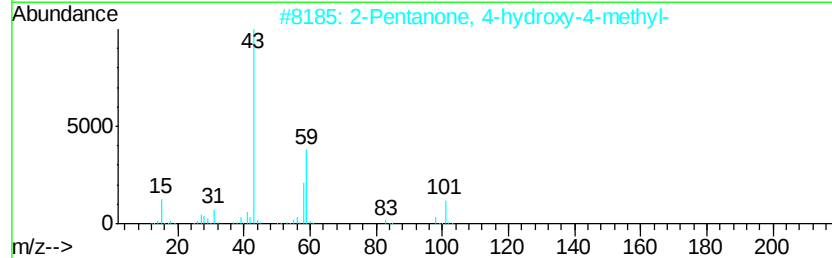
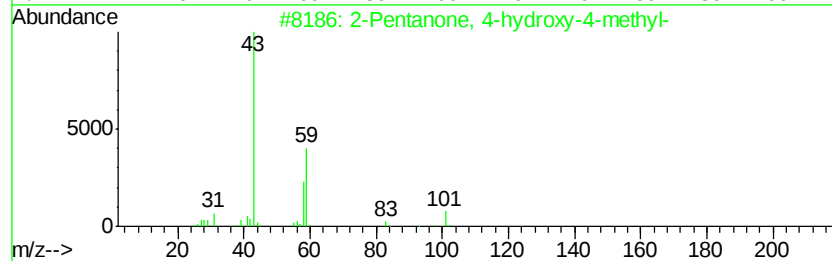
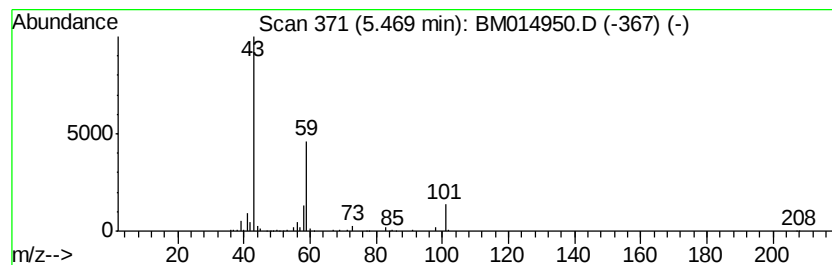
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM041918.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.47	5.69 ng/ul	560667	1,4-Dichlorobenzene-d4	8.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
3		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	47
4		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	25
5		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10



Data Path : Z:\HPCHEM1\BNA M\DATA\BM050318\
Data File : BM014950.D
Acq On : 03 May 2018 19:28
Operator : SJ/JU
Sample : J2554-16
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0AG4

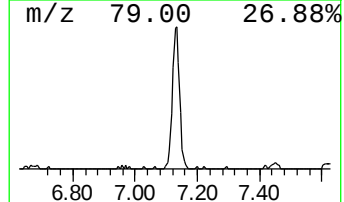
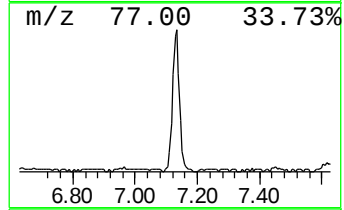
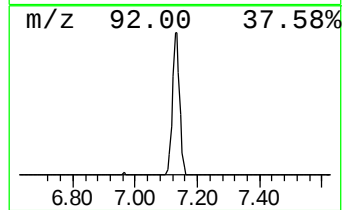
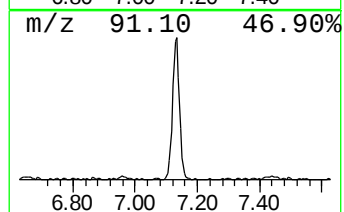
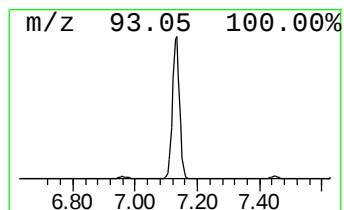
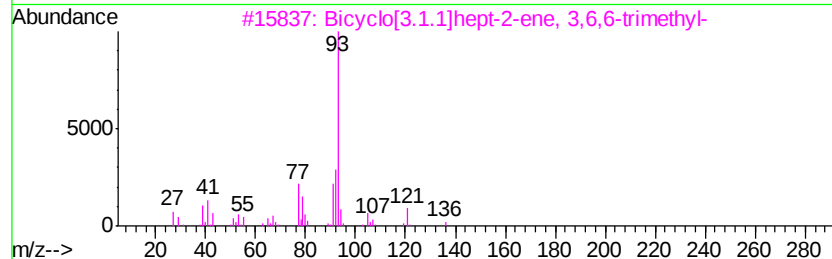
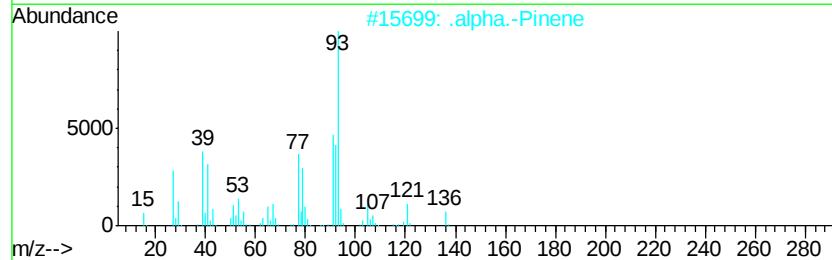
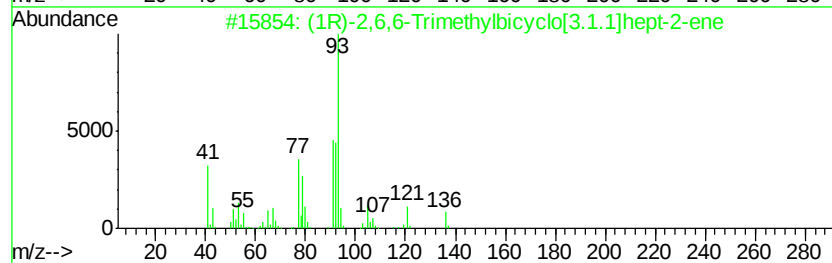
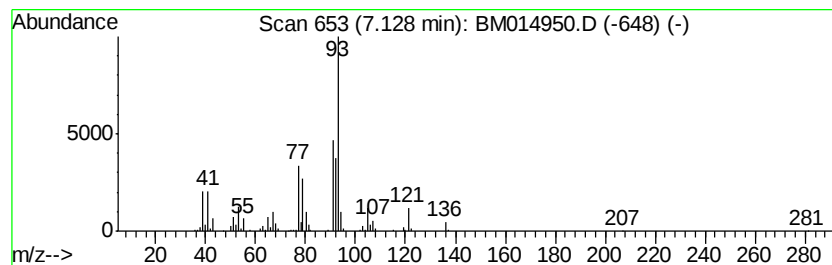
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM041918.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 (1R)-2,6,6-Trimethylbicyclo... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.13	4.65 ng/ul	458543	1,4-Dichlorobenzene-d4	8.49

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			(1R)-2,6,6-Trimethylbicyclo[3.1.1]hept-2-ene	136	C10H16	007785-70-8	96
2			.alpha.-Pinene	136	C10H16	000080-56-8	95
3			Bicyclo[3.1.1]hept-2-ene, 3,6,6-trimethyl-	136	C10H16	004889-83-2	94
4			.alpha.-Pinene	136	C10H16	000080-56-8	94
5			.gamma.-Terpinene	136	C10H16	000099-85-4	91



Data Path : Z:\HPCHEM1\BNA M\DATA\BM050318\
Data File : BM014950.D
Acq On : 03 May 2018 19:28
Operator : SJ/JU
Sample : J2554-16
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0AG4

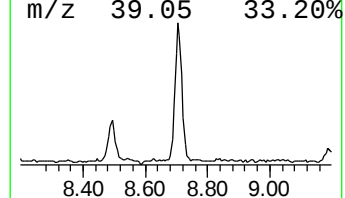
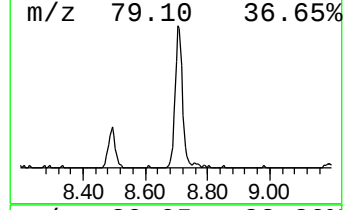
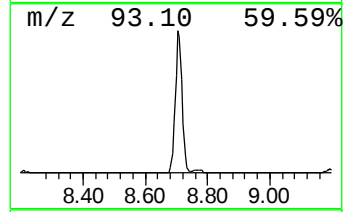
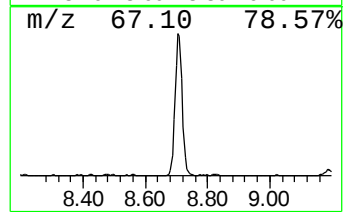
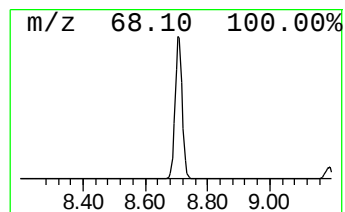
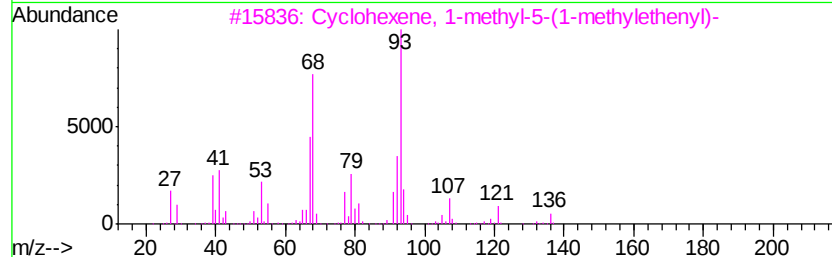
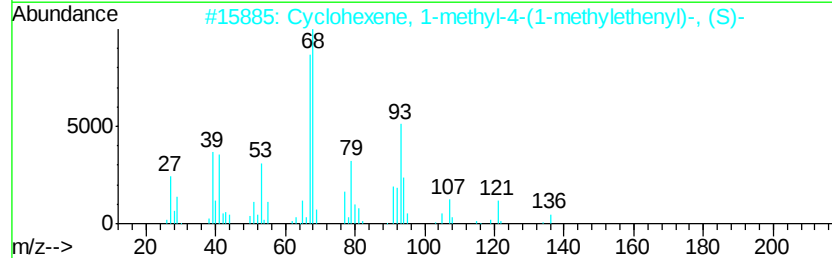
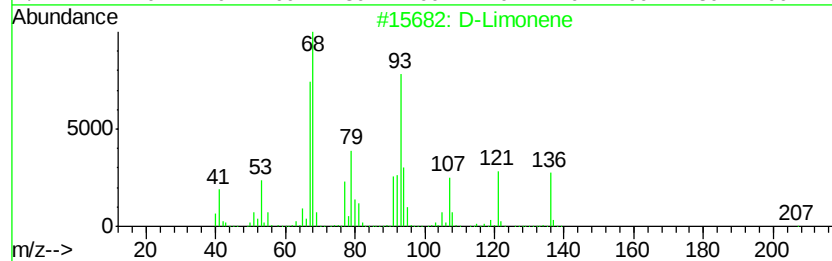
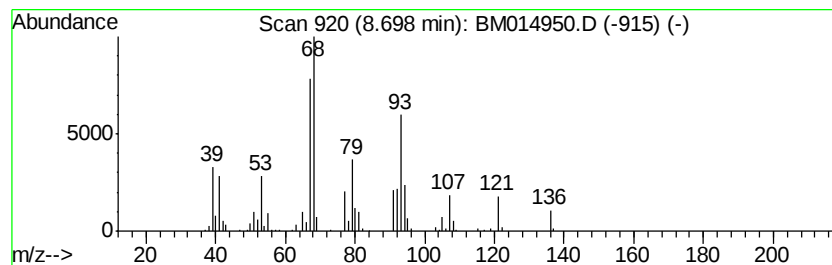
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM041918.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 D-Limonene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.70	6.06 ng/ul	597247	1,4-Dichlorobenzene-d4	8.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	D-Limonene	136	C10H16	005989-27-5	98
2		Cyclohexene, 1-methyl-4-(1-methyl-5-(1-methylethenyl)-2-yl)-	136	C10H16	005989-54-8	95
3		Cyclohexene, 1-methyl-5-(1-methyl-4-(1-methylethenyl)-2-yl)-	136	C10H16	013898-73-2	90
4		D-Limonene	136	C10H16	005989-27-5	87
5		Limonene	136	C10H16	000138-86-3	87



Data Path : Z:\HPCHEM1\BNA M\DATA\BM050318\
Data File : BM014950.D
Acq On : 03 May 2018 19:28
Operator : SJ/JU
Sample : J2554-16
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0AG4

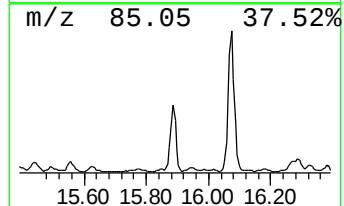
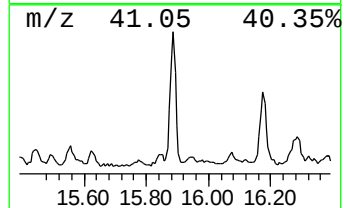
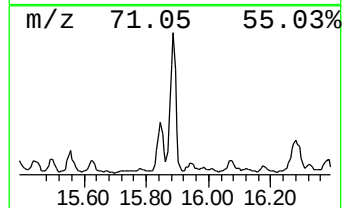
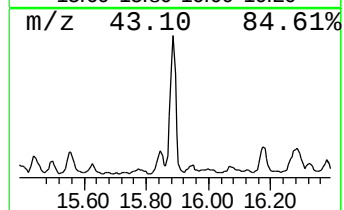
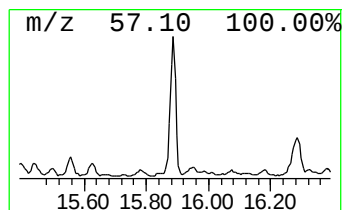
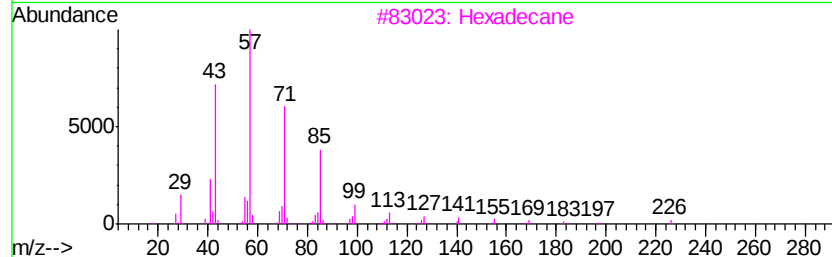
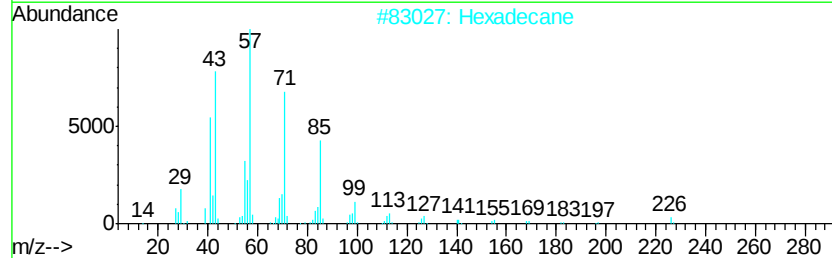
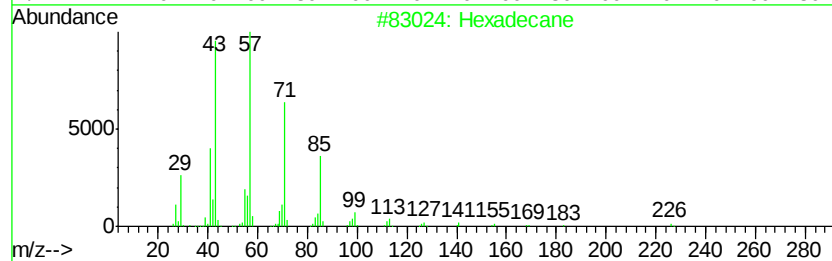
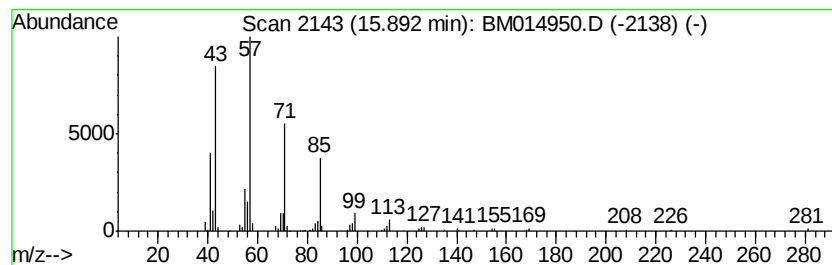
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM041918.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.89	2.24 ng/ul	331219	Acenaphthene-d10	15.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane	226	C16H34	000544-76-3	97
2		Hexadecane	226	C16H34	000544-76-3	94
3		Hexadecane	226	C16H34	000544-76-3	93
4		Hexadecane	226	C16H34	000544-76-3	91
5		Heneicosane	296	C21H44	000629-94-7	90



Data Path : Z:\HPCHEM1\BNA M\DATA\BM050318\
Data File : BM014950.D
Acq On : 03 May 2018 19:28
Operator : SJ/JU
Sample : J2554-16
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0AG4

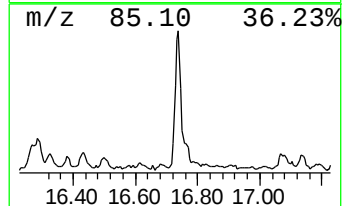
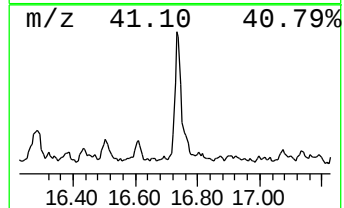
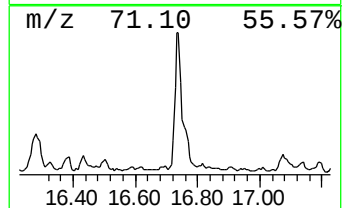
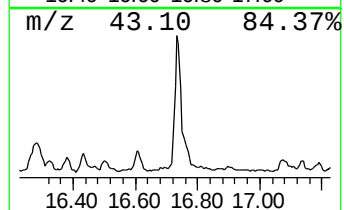
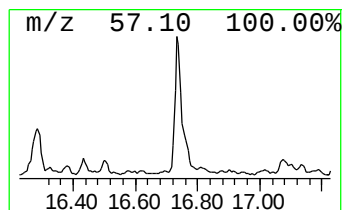
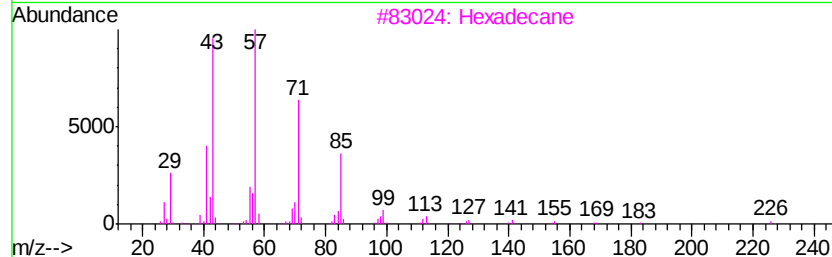
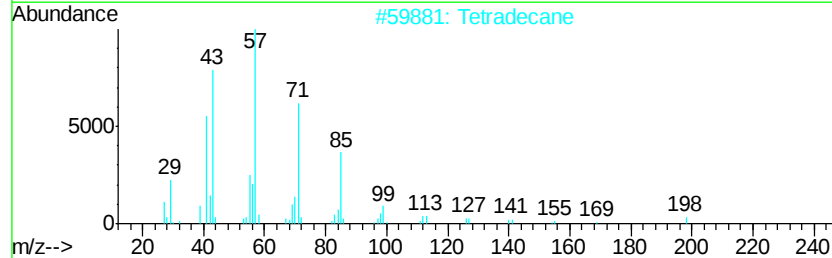
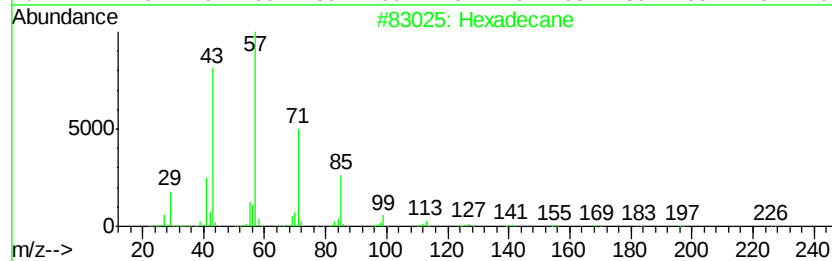
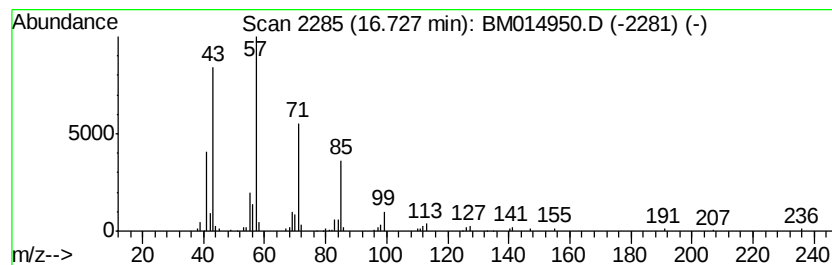
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM041918.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.73	2.20 ng/ul	352193	Phenanthrene-d10	17.81

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	91
2			Tetradecane	198	C14H30	000629-59-4	90
3			Hexadecane	226	C16H34	000544-76-3	87
4			Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	86
5			Tridecane	184	C13H28	000629-50-5	86



Data Path : Z:\HPCHEM1\BNA M\DATA\BM050318\
Data File : BM014950.D
Acq On : 03 May 2018 19:28
Operator : SJ/JU
Sample : J2554-16
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0AG4

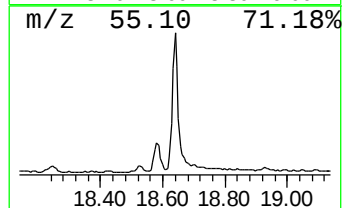
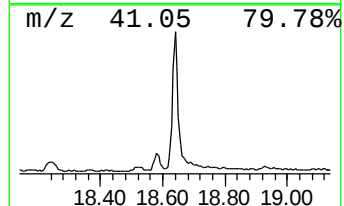
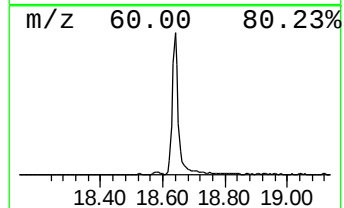
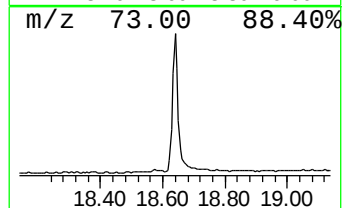
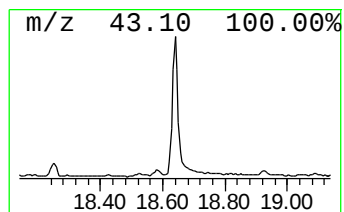
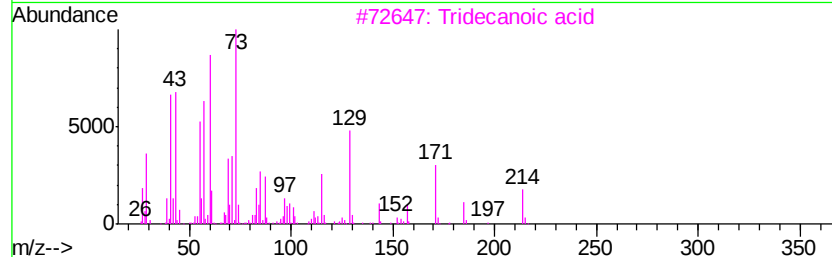
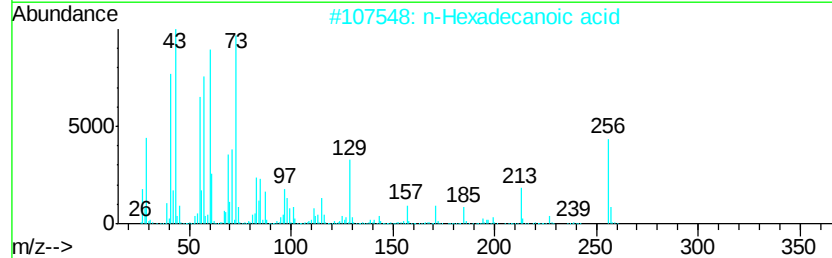
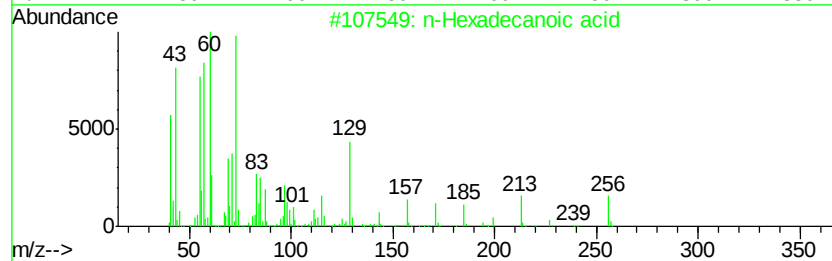
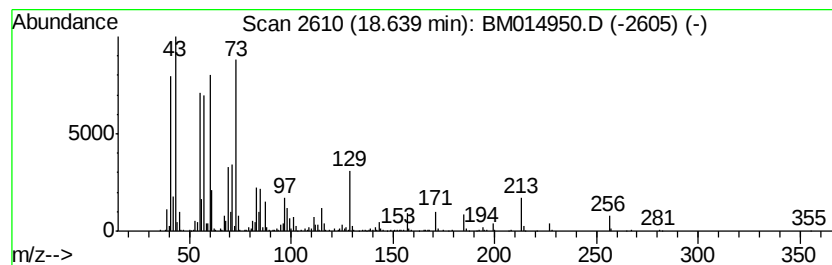
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM041918.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.64	8.49 ng/ul	1360040	Phenanthrene-d10	17.81

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	97
3		Tridecanoic acid	214	C13H26O2	000638-53-9	93
4		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91
5		Pentadecanoic acid	242	C15H30O2	001002-84-2	90



Data Path : Z:\HPCHEM1\BNA M\DATA\BM050318\
Data File : BM014950.D
Acq On : 03 May 2018 19:28
Operator : SJ/JU
Sample : J2554-16
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0AG4

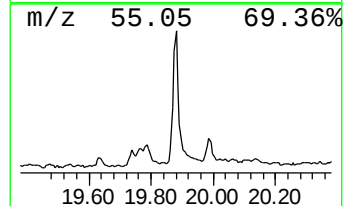
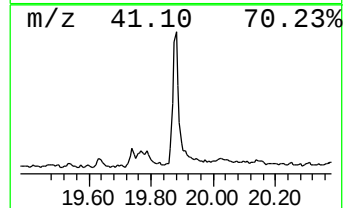
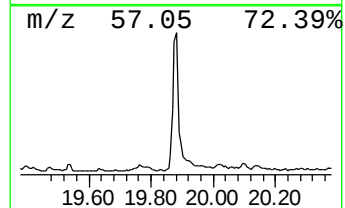
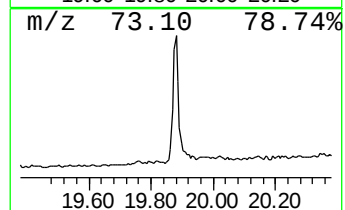
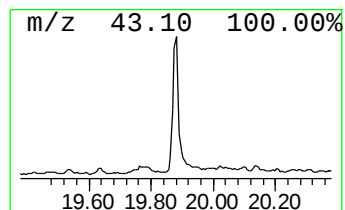
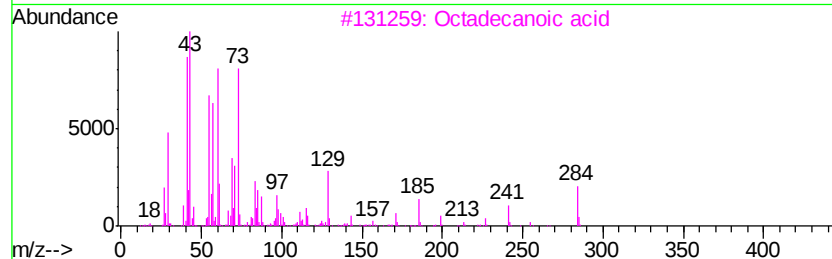
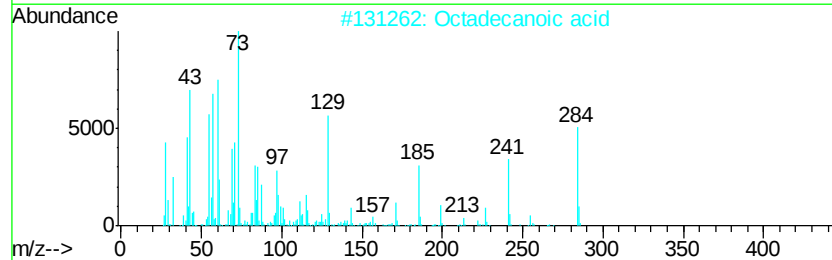
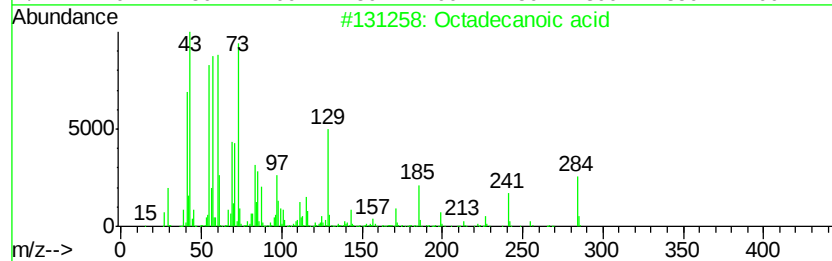
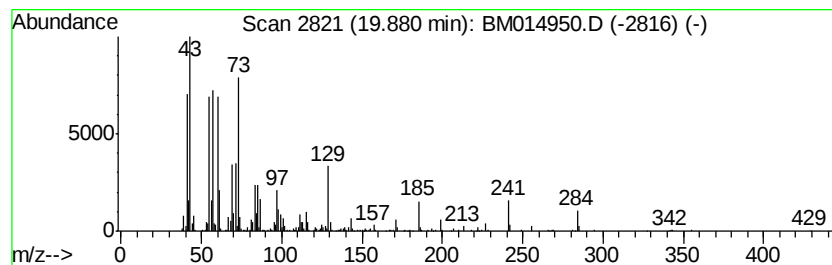
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM041918.M
Quant Title : SVOA CALIBRATION

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

Peak Number 8 Octadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.88	4.83 ng/ul	788251	Chrysene-d12	21.90

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			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	97
4			Octadecanoic acid	284	C18H36O2	000057-11-4	97
5			Octadecanoic acid	284	C18H36O2	000057-11-4	91



Data Path : Z:\HPCHEM1\BNA M\DATA\BM050318\
Data File : BM014950.D
Acq On : 03 May 2018 19:28
Operator : SJ/JU
Sample : J2554-16
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0AG4

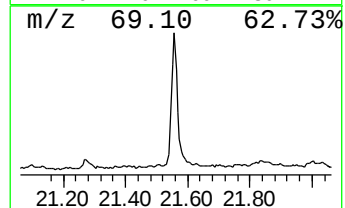
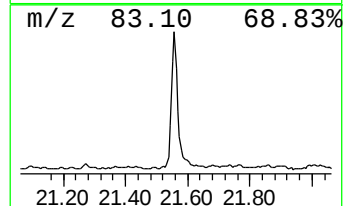
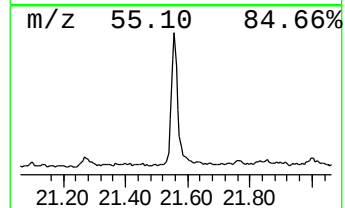
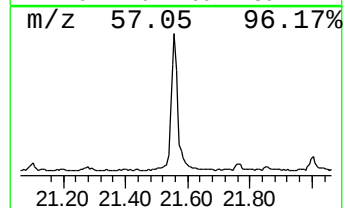
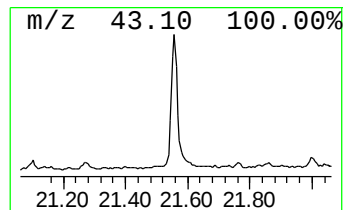
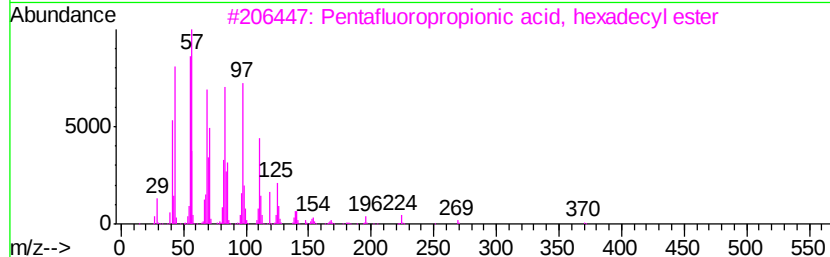
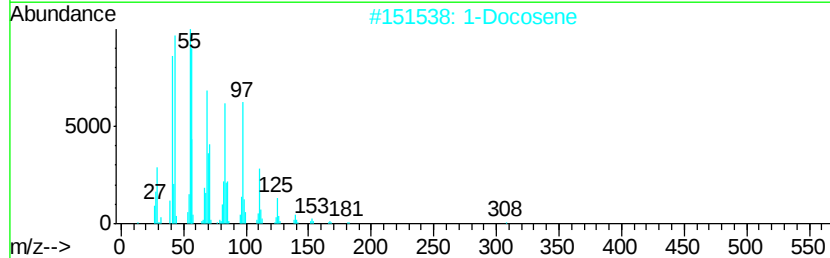
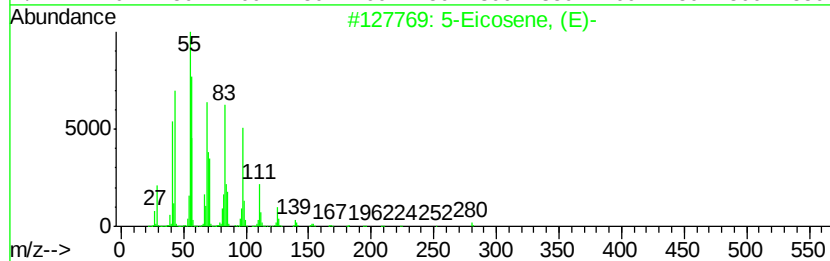
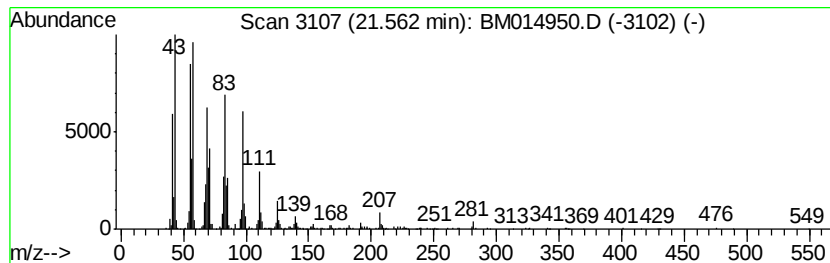
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM041918.M
Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Cyclic21.56 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.56	5.55 ng/ul	906695	Chrysene-d12	21.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Eicosene, (E)-	280	C20H40	074685-30-6	96
2		1-Docosene	308	C22H44	001599-67-3	94
3		Pentafluoropropionic acid, hexadecyl ester	388	C19H33F5O2	006222-07-7	91
4		Heptafluorobutyric acid, pentadecyl ester	424	C19H31F7O2	959261-23-5	91
5		Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	91



Data Path : Z:\HPCHEM1\BNA M\DATA\BM050318\
Data File : BM014950.D
Acq On : 03 May 2018 19:28
Operator : SJ/JU
Sample : J2554-16
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0AG4

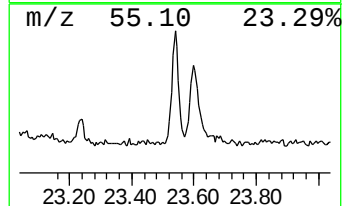
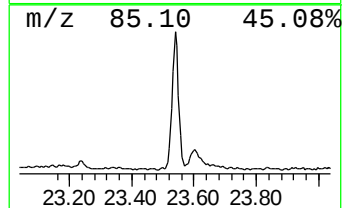
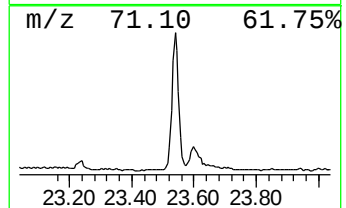
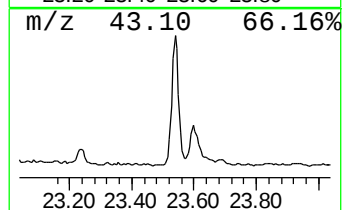
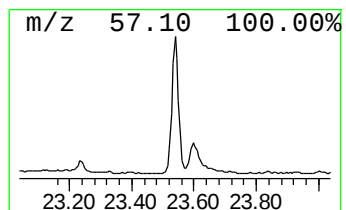
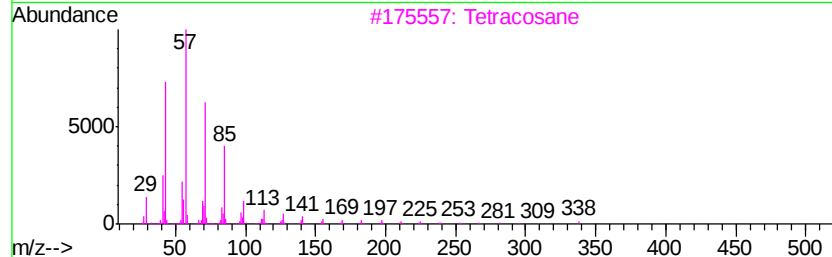
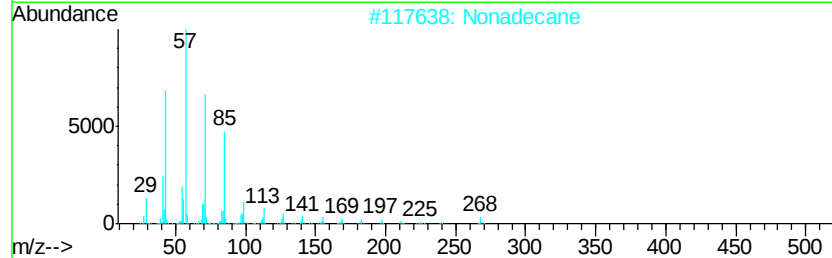
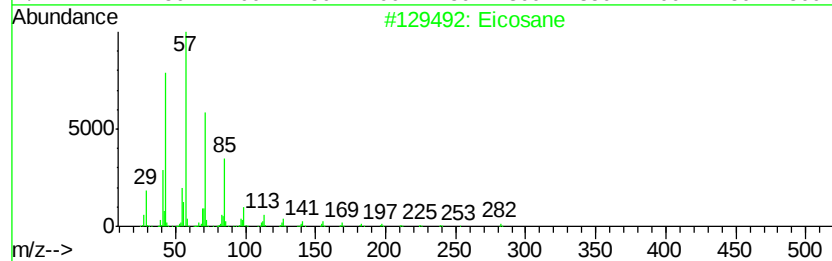
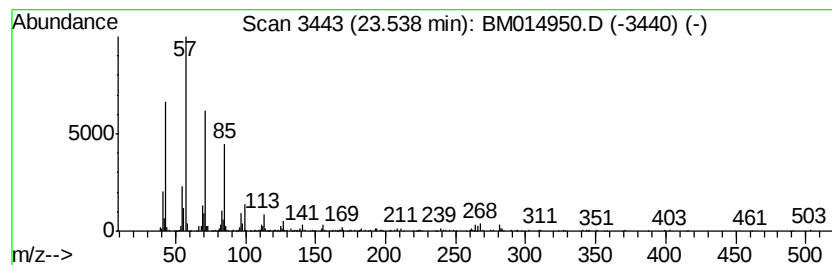
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM041918.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.54	3.10 ng/ul	515511	Perylene-d12	24.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	97
2		Nonadecane	268	C19H40	000629-92-5	97
3		Tetracosane	338	C24H50	000646-31-1	87
4		Heptacosane	380	C27H56	000593-49-7	87
5		Pentacosane	352	C25H52	000629-99-2	83



Data Path : Z:\HPCHEM1\BNA M\DATA\BM050318\
Data File : BM014950.D
Acq On : 03 May 2018 19:28
Operator : SJ/JU
Sample : J2554-16
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0AG4

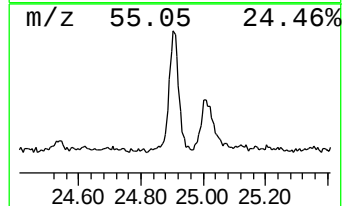
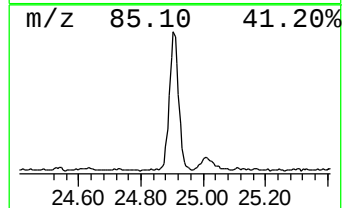
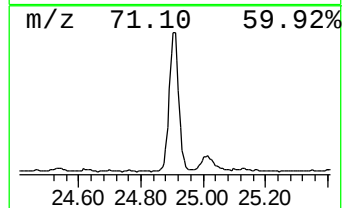
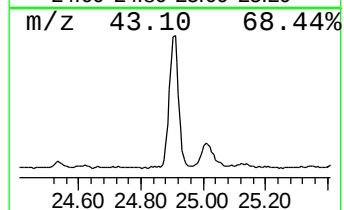
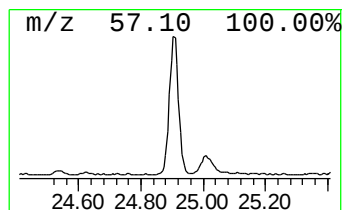
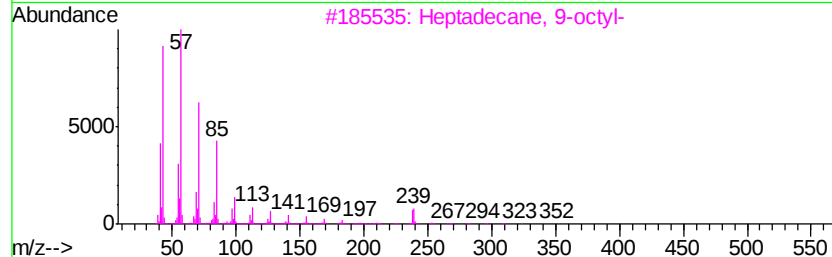
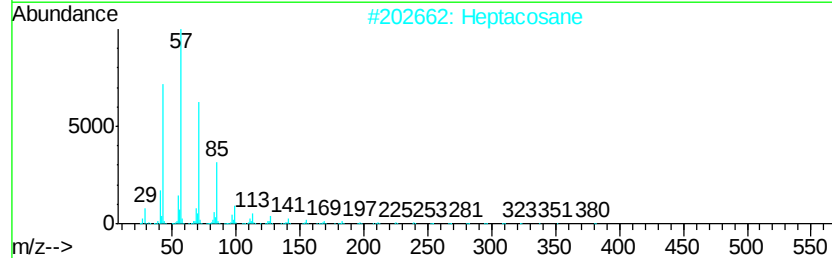
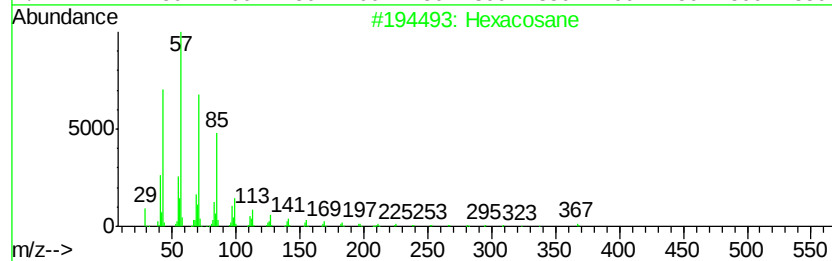
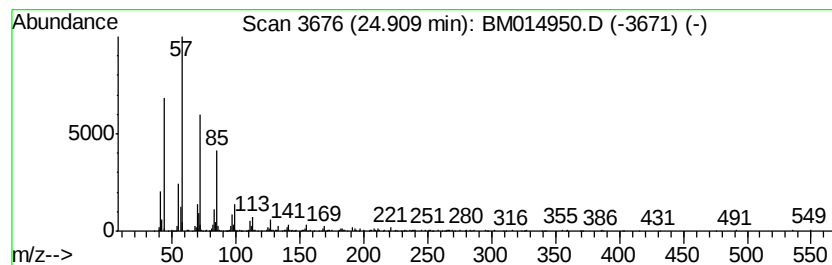
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM041918.M
Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.91	5.65 ng/ul	938893	Perylene-d12	24.37

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexacosane	366	C26H54	000630-01-3	90
2			Heptacosane	380	C27H56	000593-49-7	90
3			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	90
4			Tetracosane	338	C24H50	000646-31-1	90
5			Octacosane	394	C28H58	000630-02-4	90



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM050318\
Data File : BM014950.D
Acq On : 03 May 2018 19:28
Operator : SJ/JU
Sample : J2554-16
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0AG4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM041918.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
Butane, 2-methoxy...	3.33	31.2	ng/ul	3078010	1	8.49	1972040	20.0
2-Pentanone, 4-hy...	5.47	5.7	ng/ul	560667	1	8.49	1972040	20.0
(1R)-2,6,6-Trimet...	7.13	4.7	ng/ul	458543	1	8.49	1972040	20.0
D-Limonene	8.70	6.1	ng/ul	597247	1	8.49	1972040	20.0
(DEL) Alkane: Str...	15.89	2.2	ng/ul	331219	3	15.10	2956220	20.0
(DEL) Alkane: Str...	16.73	2.2	ng/ul	352193	4	17.81	3205360	20.0
n-Hexadecanoic acid	18.64	8.5	ng/ul	1360040	4	17.81	3205360	20.0
Octadecanoic acid	19.88	4.8	ng/ul	788251	5	21.90	3264900	20.0
(DEL) Alkane: Cyc...	21.56	5.5	ng/ul	906695	5	21.90	3264900	20.0
(DEL) Alkane: Str...	23.54	3.1	ng/ul	515511	6	24.37	3325690	20.0
(DEL) Alkane: Str...	24.91	5.7	ng/ul	938893	6	24.37	3325690	20.0