

Data Path : Z:\HPCHEM1\BNA M\DATA\BM101117\
 Data File : BM012079.D
 Acq On : 11 Oct 2017 21:33
 Operator : SJ/JU
 Sample : I5649-09
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 B-57(2-4)-100317

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	3.258	25	29	36	rVB	68734	95399	1.81%	0.134%
2	3.787	114	119	124	rBV	99567	135262	2.56%	0.189%
3	5.334	376	382	391	rVB	454652	671584	12.72%	0.940%
4	7.004	656	666	681	rBV	980907	1765204	33.44%	2.472%
5	7.169	686	694	706	rVB	856529	1337803	25.34%	1.873%
6	7.369	721	728	741	rBV	1133700	1852413	35.09%	2.594%
7	7.840	799	808	816	rBV	1194406	1939468	36.74%	2.716%
8	7.969	824	830	839	rVB	98487	156667	2.97%	0.219%
9	8.051	839	844	850	rBV2	61063	98610	1.87%	0.138%
10	8.545	922	928	936	rBV	947580	1678824	31.80%	2.351%
11	9.004	998	1006	1021	rBV	1035851	1849423	35.03%	2.590%
12	9.728	1117	1129	1140	rBV	899850	1627938	30.84%	2.280%
13	10.269	1213	1221	1233	rBV	1267988	2331116	44.16%	3.265%
14	10.645	1278	1285	1291	rBV	1600772	2694466	51.04%	3.773%
15	10.698	1291	1294	1301	rVB	37997	59534	1.13%	0.083%
16	10.786	1301	1309	1327	rBV	916294	1803016	34.15%	2.525%
17	12.051	1517	1524	1531	rBV3	32024	60561	1.15%	0.085%
18	13.292	1730	1735	1743	rBV3	39929	64308	1.22%	0.090%
19	13.810	1818	1823	1828	rBV5	27997	53422	1.01%	0.075%
20	13.892	1828	1837	1842	rBV	2341656	3456657	65.48%	4.841%
21	13.939	1842	1845	1852	rVB	1037696	1424785	26.99%	1.995%
22	14.180	1880	1886	1899	rBV	2863726	4162651	78.85%	5.829%
23	14.369	1914	1918	1927	rVB	83647	118716	2.25%	0.166%
24	14.492	1927	1939	1950	rBV2	2192386	3548768	67.22%	4.970%
25	14.698	1968	1974	1989	rBV	365250	820907	15.55%	1.150%
26	14.886	2000	2006	2010	rBV6	30902	56314	1.07%	0.079%
27	15.321	2075	2080	2085	rVB	119265	152528	2.89%	0.214%
28	15.480	2099	2107	2114	rBV	3371194	4814789	91.21%	6.743%
29	15.610	2124	2129	2141	rBV	447584	705999	13.37%	0.989%
30	15.721	2142	2148	2153	rVB2	92716	162648	3.08%	0.228%
31	15.821	2161	2165	2171	rBV9	35389	56922	1.08%	0.080%
32	16.180	2221	2226	2229	rBV	139953	216980	4.11%	0.304%
33	16.521	2278	2284	2286	rBV7	35498	64024	1.21%	0.090%
34	16.768	2321	2326	2332	rBV4	52256	94614	1.79%	0.132%

Data Path : Z:\HPCHEM1\BNA M\DATA\BM101117\
Data File : BM012079.D
Acq On : 11 Oct 2017 21:33
Operator : SJ/JU
Sample : I5649-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
B-57(2-4)-100317

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

35	16.968	2357	2360	2365	rBV	74663	91428	1.73%	0.128%
36	17.021	2365	2369	2373	rVB3	36649	55764	1.06%	0.078%
37	17.233	2399	2405	2410	rBV2	2564881	3668813	69.50%	5.138%
38	17.333	2416	2422	2433	rBV	3307510	4541844	86.04%	6.360%
39	17.510	2448	2452	2461	rBV7	40617	101049	1.91%	0.142%
40	18.109	2549	2554	2559	rBV	994481	1384881	26.23%	1.939%
41	18.404	2601	2604	2608	rBV6	32053	59956	1.14%	0.084%
42	19.033	2707	2711	2714	rBV2	88998	129240	2.45%	0.181%
43	19.268	2749	2751	2752	rBV2	92369	74557	1.41%	0.104%
44	19.386	2767	2771	2785	rBV	799141	1264607	23.96%	1.771%
45	19.627	2806	2812	2826	rBV	3608502	5278980	100.00%	7.393%
46	20.145	2898	2900	2904	rVB2	189644	184941	3.50%	0.259%
47	20.533	2962	2966	2972	rVB	590236	699084	13.24%	0.979%
48	21.098	3059	3062	3067	rVB	786702	963868	18.26%	1.350%
49	21.409	3111	3115	3120	rBV	2867144	3697811	70.05%	5.178%
50	23.580	3478	3484	3498	rVB2	2577273	5261801	99.67%	7.369%
51	23.727	3504	3509	3517	rVB2	1965193	3816232	72.29%	5.344%

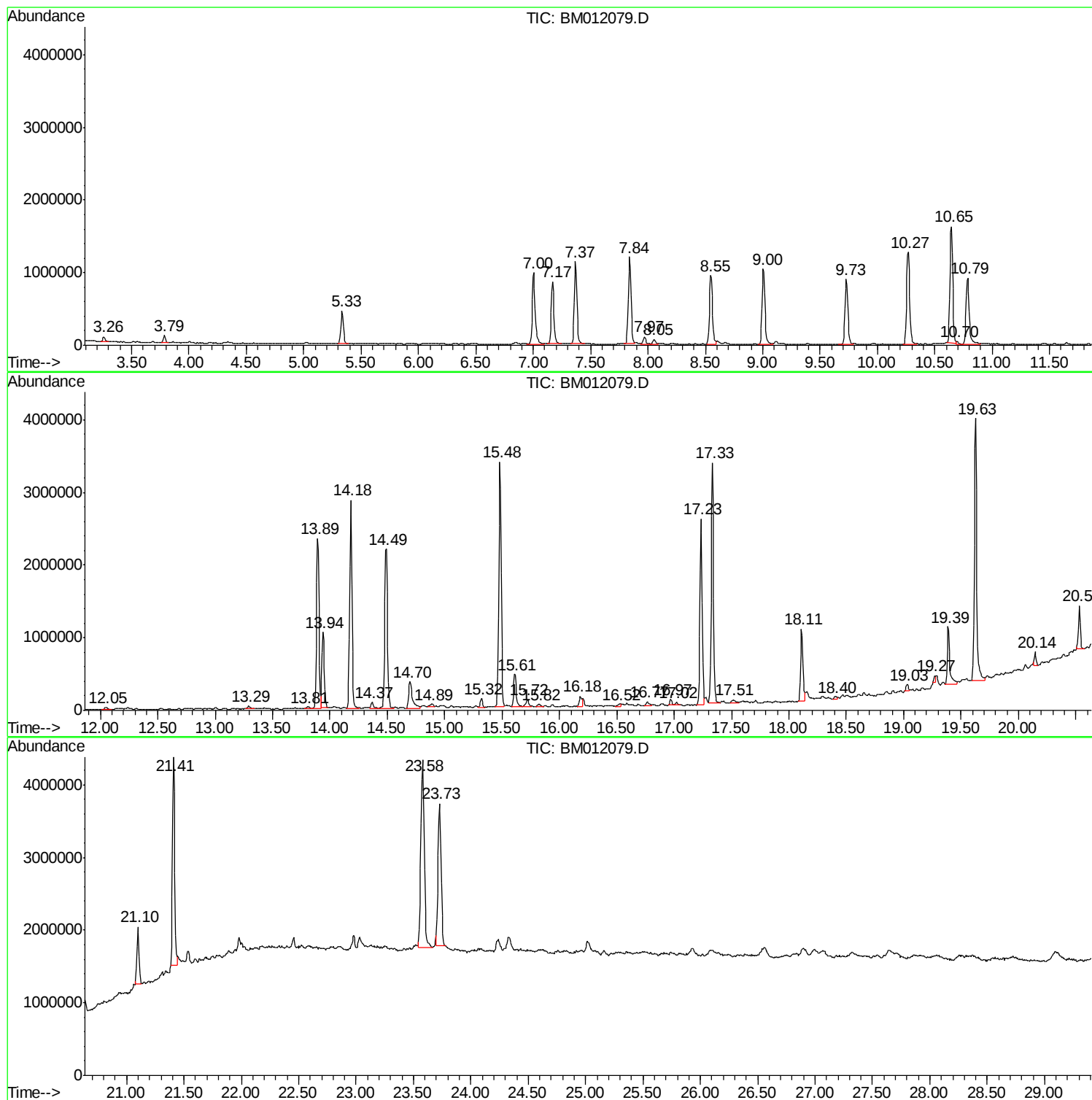
Sum of corrected areas: 71407176

Data Path : Z:\HPCHEM1\BNA M\DATA\BM101117\
Data File : BM012079.D
Acq On : 11 Oct 2017 21:33
Operator : SJ/JU
Sample : I5649-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampled :
B-57(2-4)-100317

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



Data Path : Z:\HPCHEM1\BNA M\DATA\BM101117\
Data File : BM012079.D
Acq On : 11 Oct 2017 21:33
Operator : SJ/JU
Sample : I5649-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
B-57(2-4)-100317

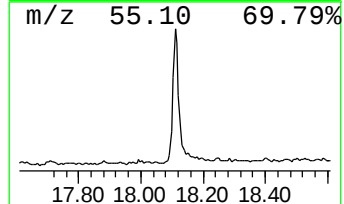
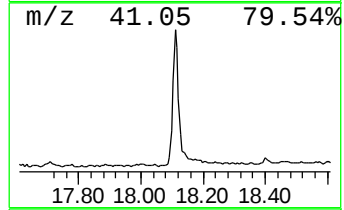
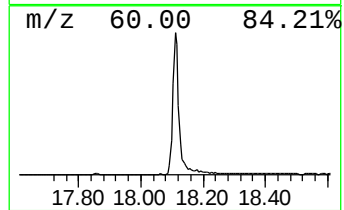
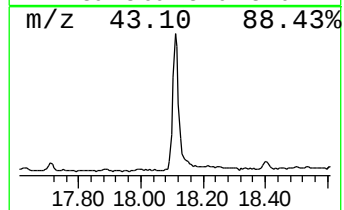
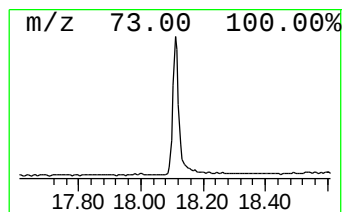
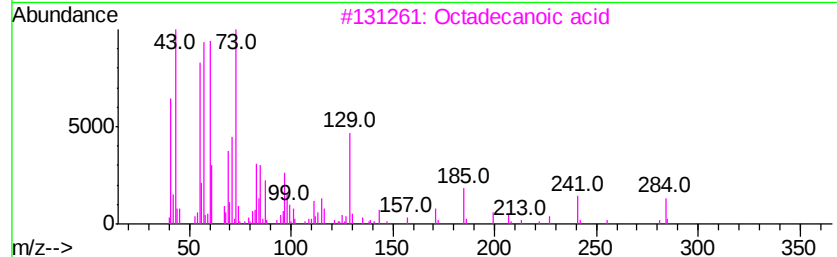
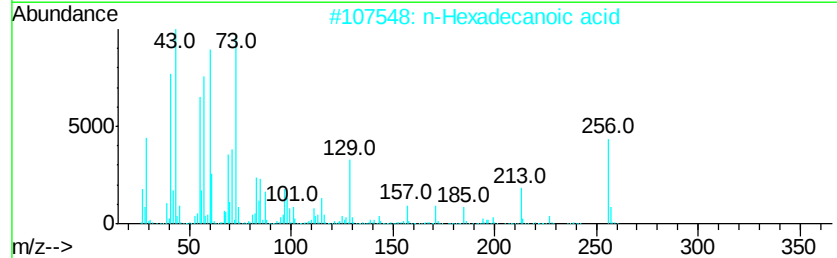
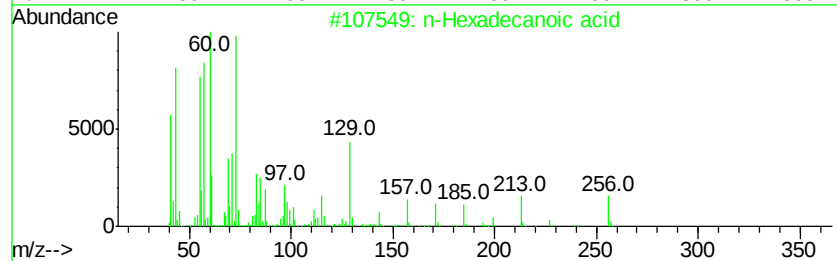
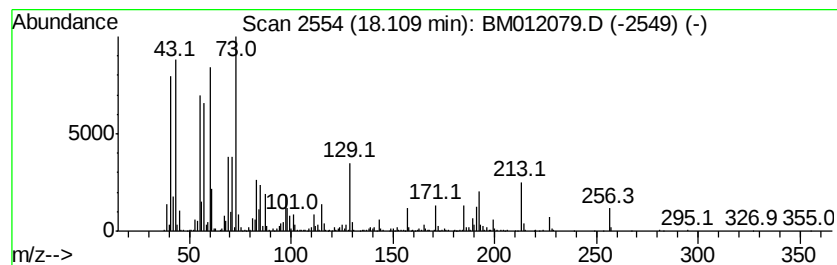
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM100317.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.11	7.55 ng/ul	1384880	Phenanthrene-d10	17.23

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		Octadecanoic acid	284	C18H36O2	000057-11-4	90
4		Tridecanoic acid	214	C13H26O2	000638-53-9	90
5		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	83



Data Path : Z:\HPCHEM1\BNA M\DATA\BM101117\
Data File : BM012079.D
Acq On : 11 Oct 2017 21:33
Operator : SJ/JU
Sample : I5649-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
B-57(2-4)-100317

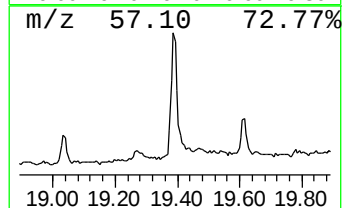
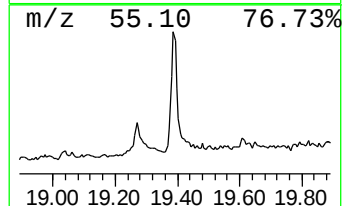
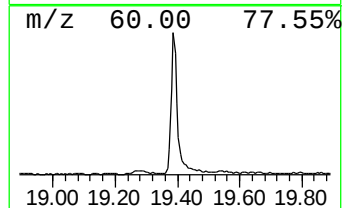
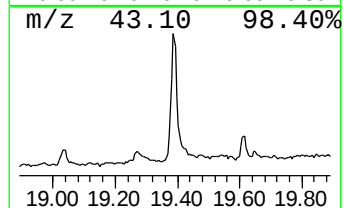
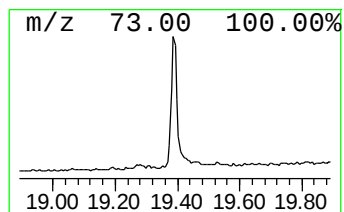
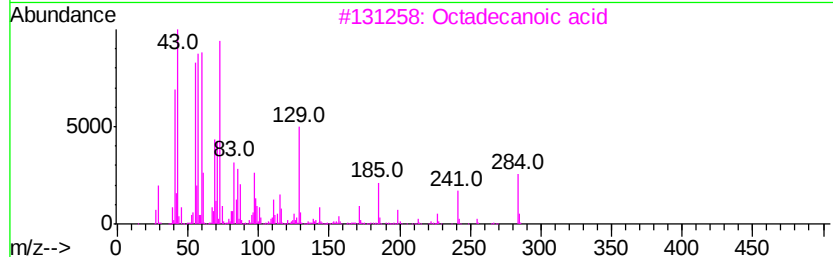
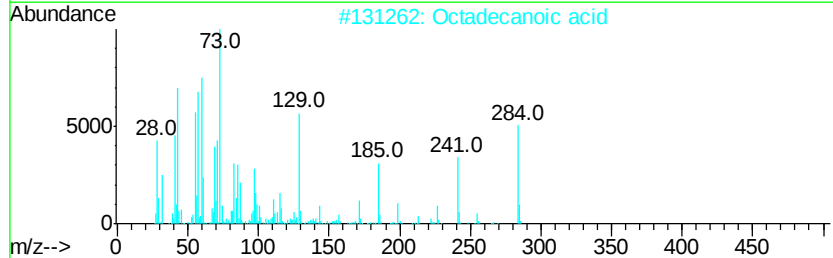
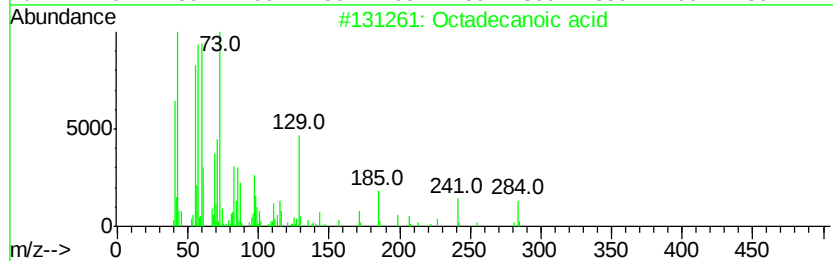
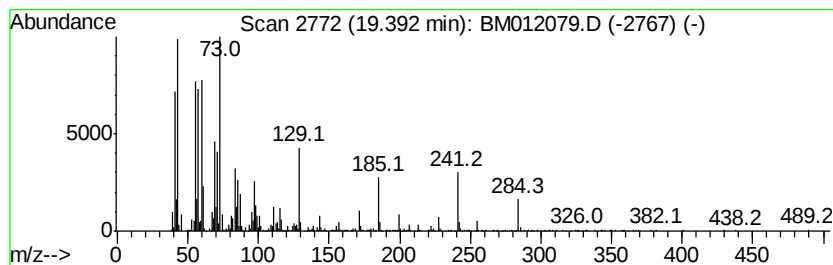
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM100317.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Octadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.39	6.84 ng/ul	1264610	Chrysene-d12	21.41

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	99
4		Octadecanoic acid	284	C18H36O2	000057-11-4	98
5		Octadecanoic acid	284	C18H36O2	000057-11-4	94



Data Path : Z:\HPCHEM1\BNA M\DATA\BM101117\
Data File : BM012079.D
Acq On : 11 Oct 2017 21:33
Operator : SJ/JU
Sample : I5649-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
B-57(2-4)-100317

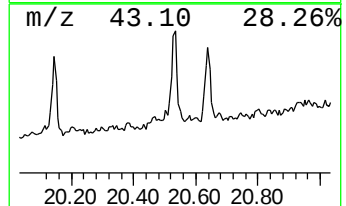
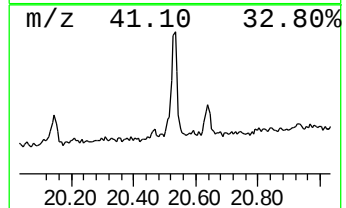
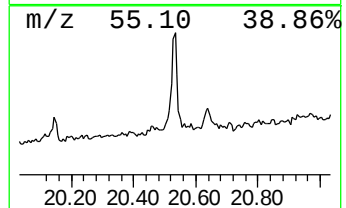
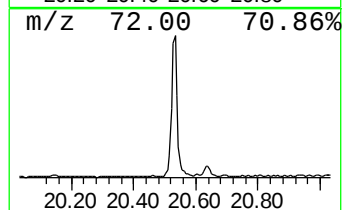
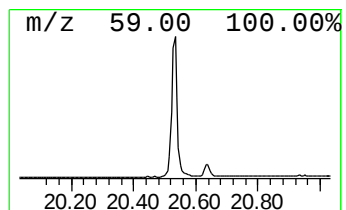
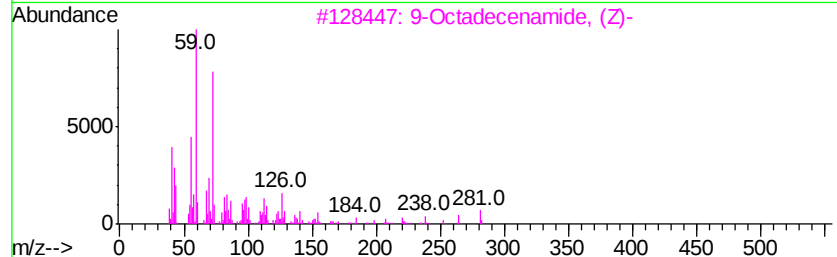
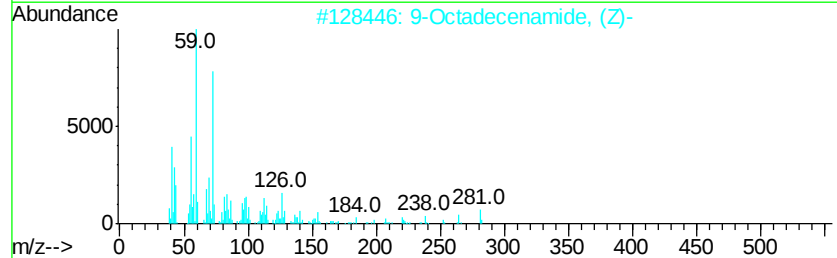
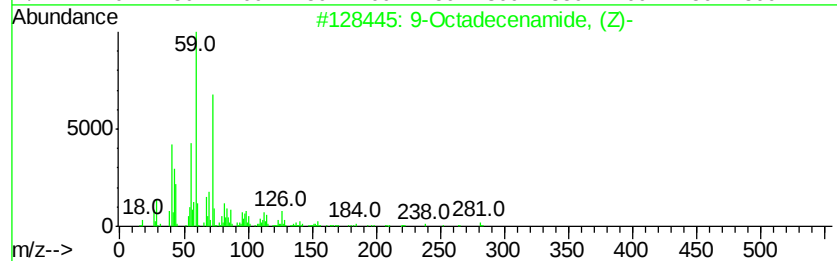
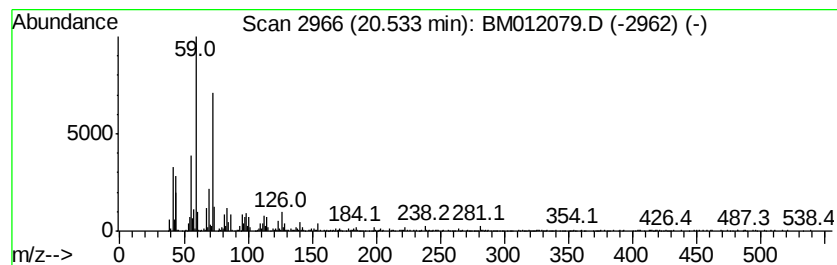
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM100317.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 9-Octadecenamide, (Z)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.53	3.78 ng/ul	699084	Chrysene-d12	21.41

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	99
2		9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	91
3		9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	91
4		7-Nonenamide	155	C9H17NO	090949-53-4	72
5		9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	72



Data Path : Z:\HPCHEM1\BNA M\DATA\BM101117\
Data File : BM012079.D
Acq On : 11 Oct 2017 21:33
Operator : SJ/JU
Sample : I5649-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
B-57(2-4)-100317

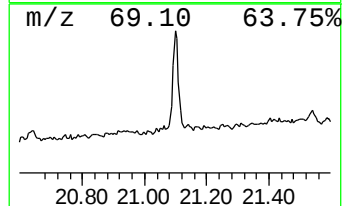
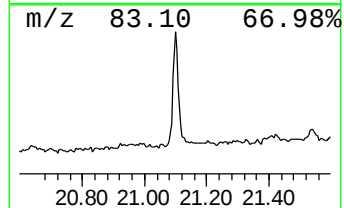
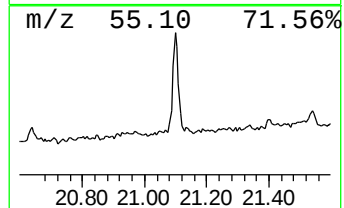
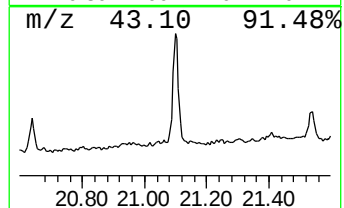
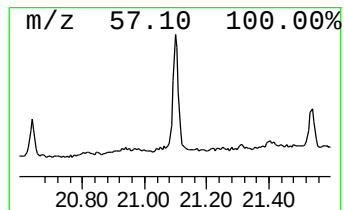
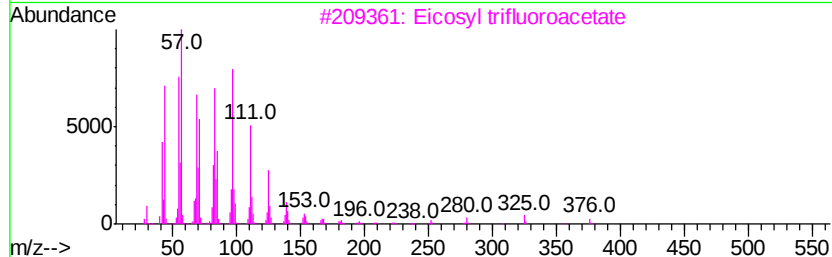
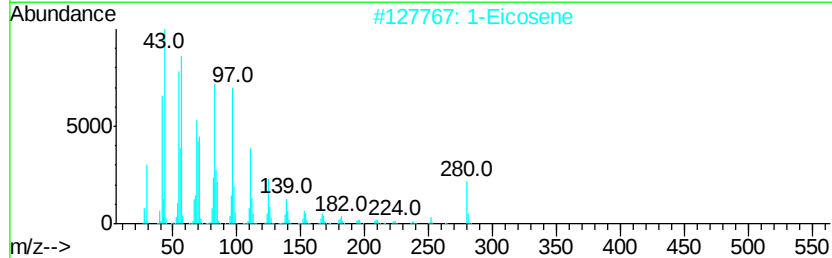
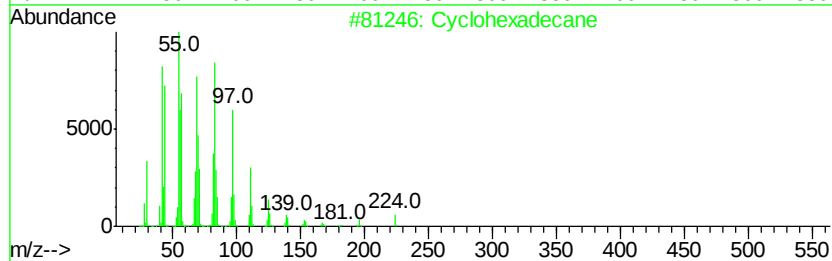
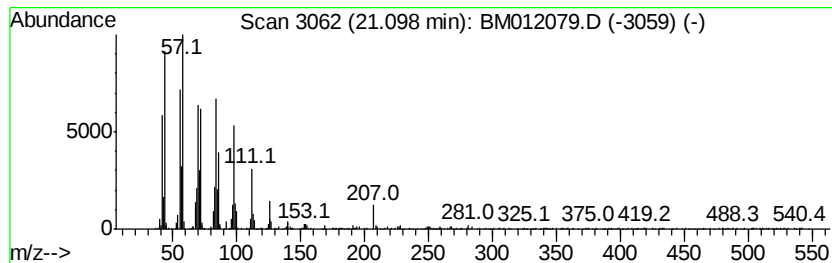
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM100317.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Cyclic21.10 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.10	5.21 ng/ul	963868	Chrysene-d12	21.41

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexadecane	224	C16H32	000295-65-8	96
2		1-Eicosene	280	C20H40	003452-07-1	93
3		Eicosyl trifluoroacetate	394	C22H41F3O2	1000351-75-2	91
4		Tricosyl trifluoroacetate	436	C25H47F3O2	1000351-75-1	91
5		Cyclotriacontane	420	C30H60	000297-35-8	91



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM101117\
Data File : BM012079.D
Acq On : 11 Oct 2017 21:33
Operator : SJ/JU
Sample : I5649-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
B-57(2-4)-100317

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
n-Hexadecanoic acid	18.11	7.5	ng/ul	1384880	4	17.23	3668810	20.0
Octadecanoic acid	19.39	6.8	ng/ul	1264610	5	21.41	3697810	20.0
9-Octadecenamide, ...	20.53	3.8	ng/ul	699084	5	21.41	3697810	20.0
(DEL) Alkane: Cyc...	21.10	5.2	ng/ul	963868	5	21.41	3697810	20.0