

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM071918\
 Data File : BM015992.D
 Acq On : 19 Jul 2018 16:22
 Operator : SJ/JU
 Sample : J4024-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DB1P0

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:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM071318MA.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	5.687	403	408	416	rBV	74688	121354	4.64%	0.534%
2	5.822	425	431	448	rBV	164125	369700	14.13%	1.626%
3	6.210	492	497	507	rBV	34574	55469	2.12%	0.244%
4	7.422	698	703	717	rBV2	48191	105312	4.03%	0.463%
5	7.575	724	729	741	rBV	194537	309870	11.84%	1.363%
6	7.781	759	764	780	rBV	220173	392596	15.01%	1.727%
7	8.257	839	845	860	rBV	212303	367395	14.04%	1.616%
8	8.975	962	967	980	rBV	165657	293049	11.20%	1.289%
9	9.439	1041	1046	1064	rBV	284471	507341	19.39%	2.231%
10	10.163	1164	1169	1186	rBV	257878	462821	17.69%	2.035%
11	10.704	1256	1261	1284	rBV	356283	662246	25.31%	2.913%
12	11.081	1318	1325	1341	rVB	344873	586377	22.41%	2.579%
13	13.722	1770	1774	1783	rBB	28655	42677	1.63%	0.188%
14	13.927	1804	1809	1819	rBB	638555	916690	35.04%	4.032%
15	14.280	1863	1869	1874	rBV	876061	1212955	46.36%	5.335%
16	14.327	1874	1877	1886	rVB	192782	257712	9.85%	1.133%
17	14.580	1913	1920	1928	rBV	884065	1261865	48.23%	5.550%
18	14.880	1964	1971	1979	rBV2	561753	842283	32.20%	3.704%
19	15.110	2005	2010	2027	rBV3	35702	99329	3.80%	0.437%
20	15.863	2132	2138	2146	rBV	1182070	1661160	63.50%	7.306%
21	15.998	2156	2161	2175	rBV	458555	648329	24.78%	2.851%
22	17.610	2429	2435	2445	rBV	770329	1070716	40.93%	4.709%
23	17.710	2445	2452	2463	rBV	1418459	2026721	77.47%	8.913%
24	18.451	2569	2578	2584	rBV	143783	198543	7.59%	0.873%
25	19.698	2786	2790	2799	rBV	68476	96043	3.67%	0.422%
26	19.968	2830	2836	2844	rBV	1843431	2432661	92.99%	10.699%
27	21.398	3074	3079	3091	rBV2	159507	307525	11.76%	1.352%
28	21.721	3129	3134	3141	rVB	1067189	1375366	52.57%	6.049%
29	23.891	3500	3503	3507	rVV3	58773	86873	3.32%	0.382%
30	23.962	3507	3515	3524	rVB	1350070	2616120	100.00%	11.506%
31	24.109	3533	3540	3551	rVB	708847	1350645	51.63%	5.940%

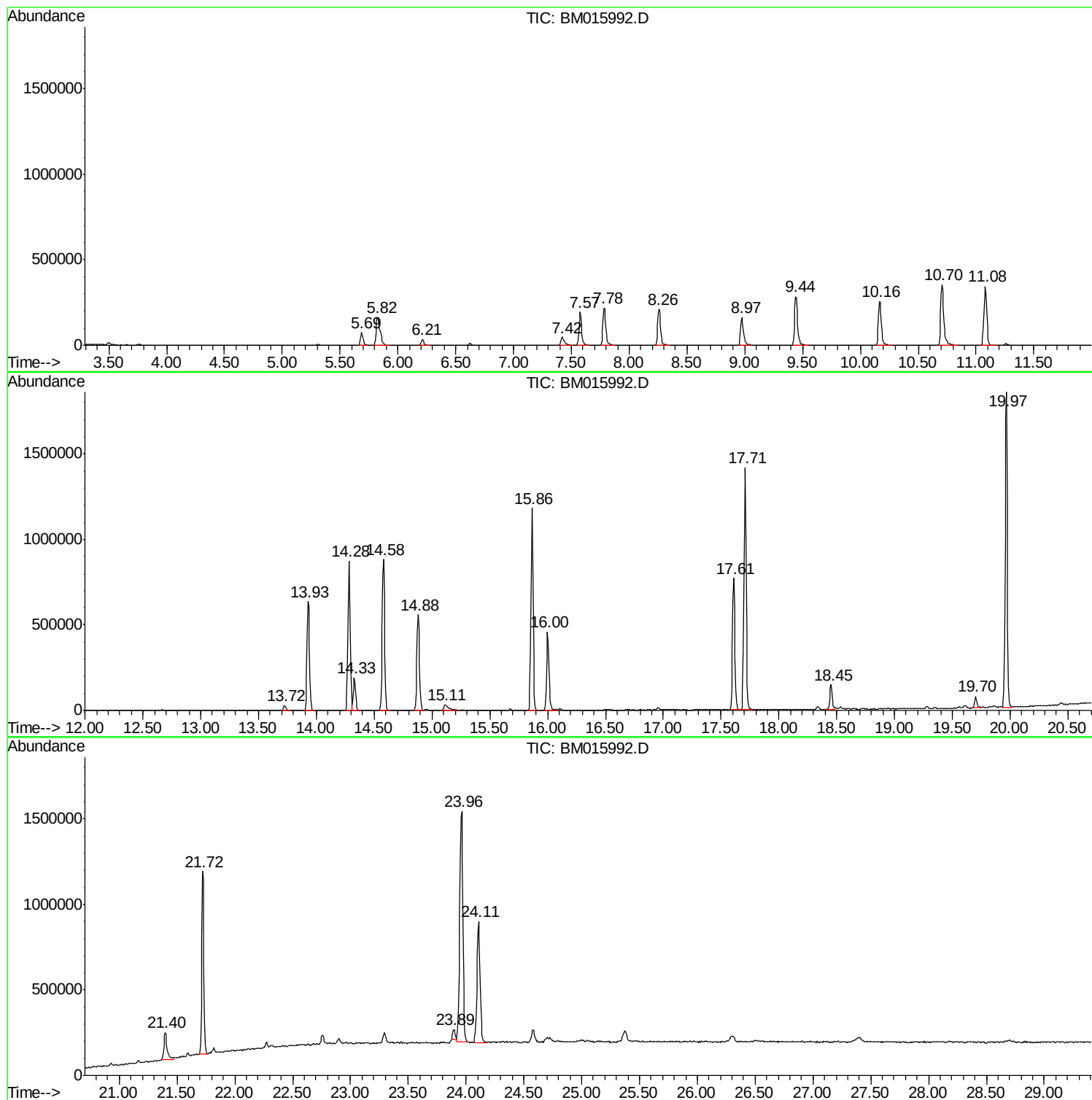
Sum of corrected areas: 22737743

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM071918\
Data File : BM015992.D
Acq On : 19 Jul 2018 16:22
Operator : SJ/JU
Sample : J4024-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DB1P0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM071318MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM071918\
Data File : BM015992.D
Acq On : 19 Jul 2018 16:22
Operator : SJ/JU
Sample : J4024-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DB1P0

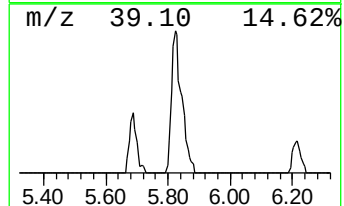
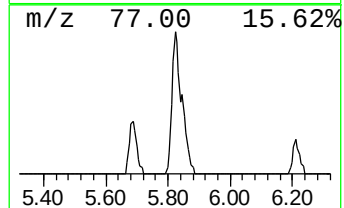
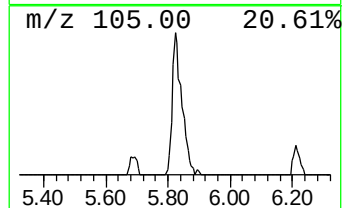
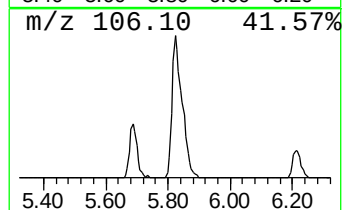
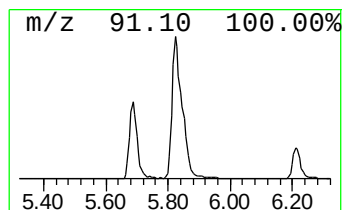
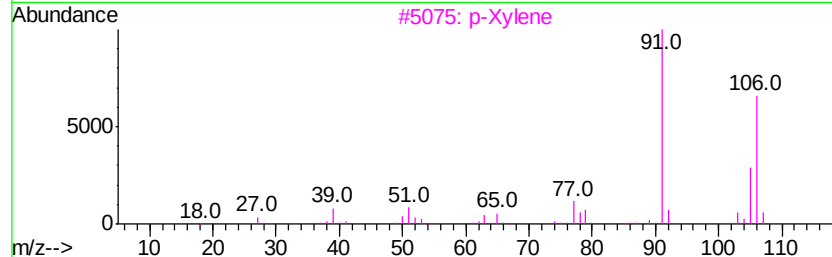
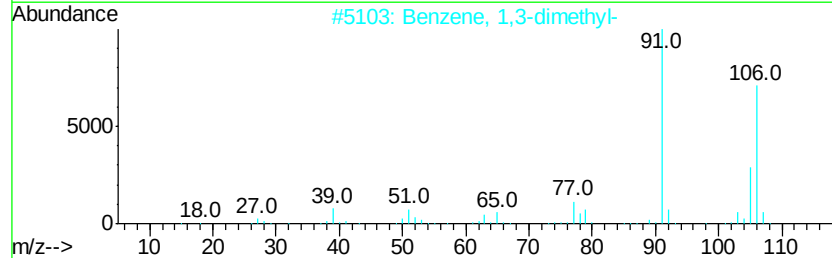
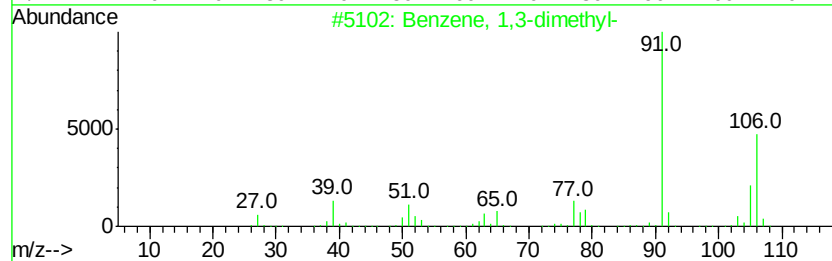
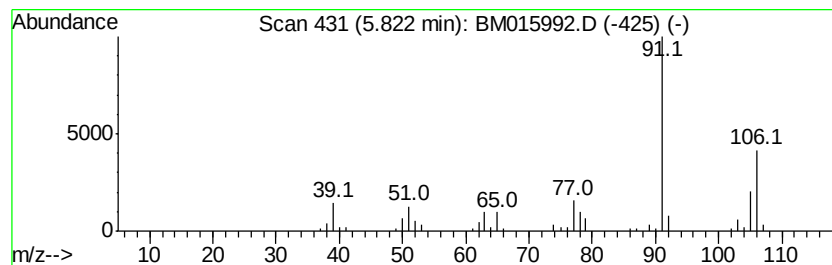
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM071318MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.82	20.13 ng/ul	369700	1,4-Dichlorobenzene-d4	8.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	94
2		Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	94
3		p-Xylene	106	C8H10	000106-42-3	94
4		p-Xylene	106	C8H10	000106-42-3	93
5		o-Xylene	106	C8H10	000095-47-6	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM071918\
Data File : BM015992.D
Acq On : 19 Jul 2018 16:22
Operator : SJ/JU
Sample : J4024-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DB1P0

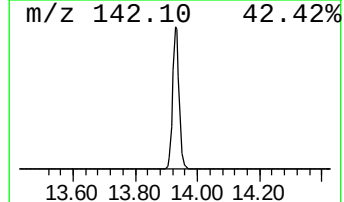
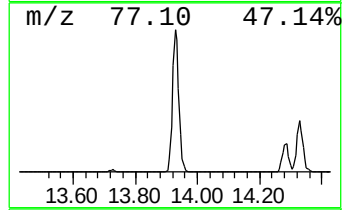
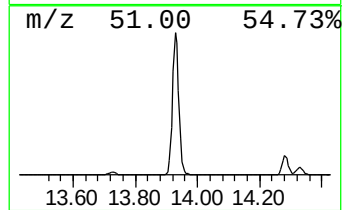
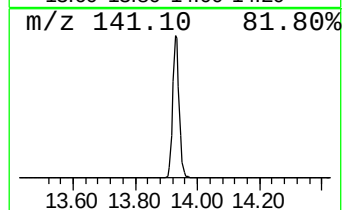
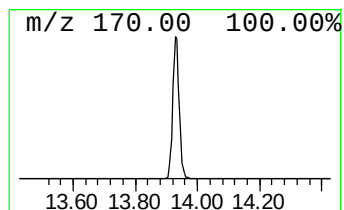
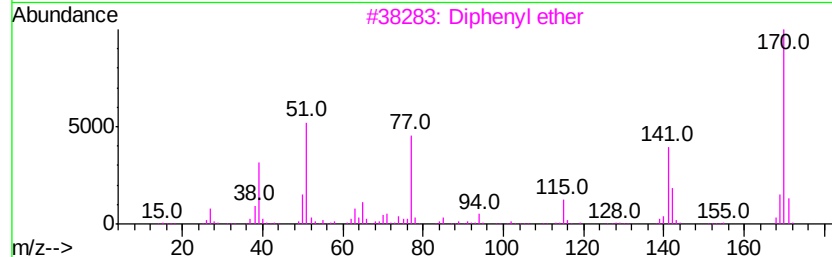
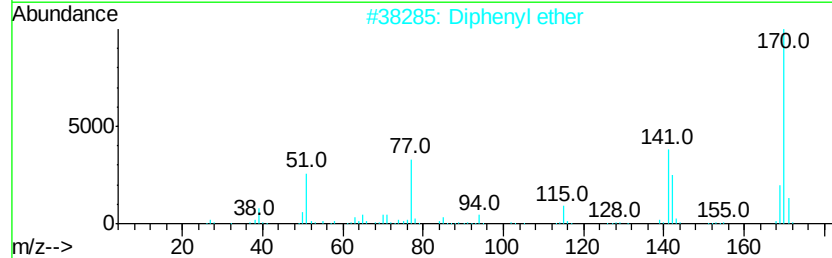
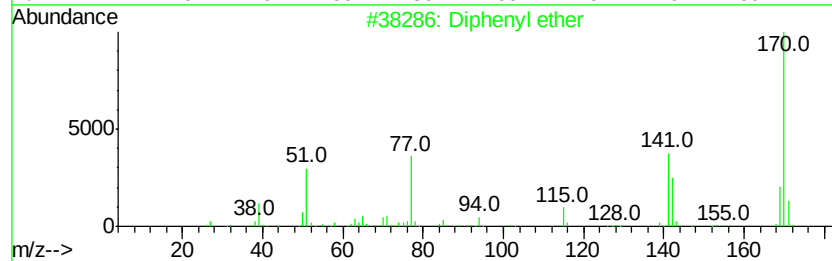
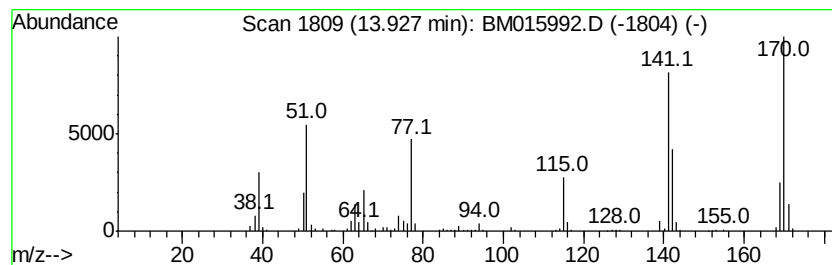
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM071318MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 Diphenyl ether Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.93	21.77 ng/ul	916690	Acenaphthene-d10	14.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Diphenyl ether	170	C12H10O	000101-84-8	76
2		Diphenyl ether	170	C12H10O	000101-84-8	76
3		Diphenyl ether	170	C12H10O	000101-84-8	52
4		Spiro[cyclopropane-1,1'(4'H)-nap...	170	C12H10O	033498-24-7	47
5		Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM071918\
Data File : BM015992.D
Acq On : 19 Jul 2018 16:22
Operator : SJ/JU
Sample : J4024-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DB1P0

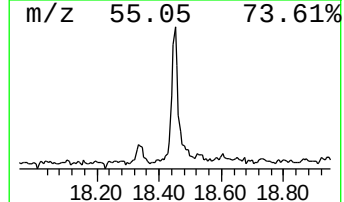
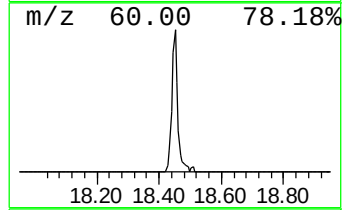
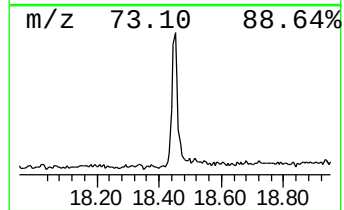
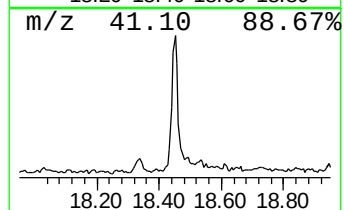
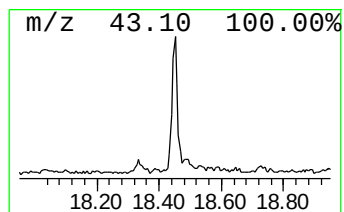
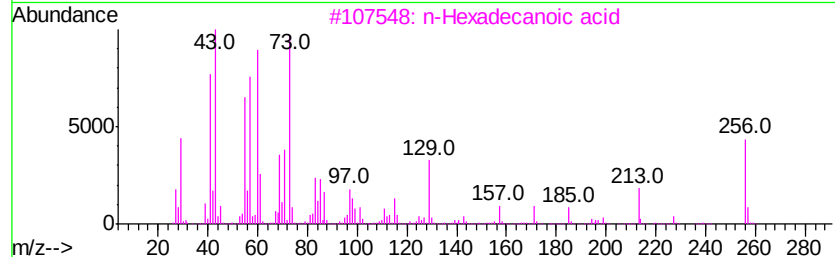
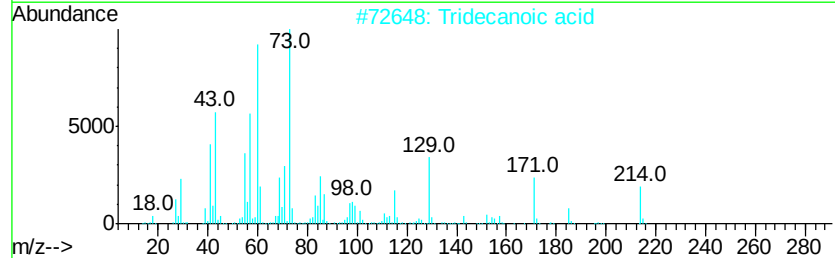
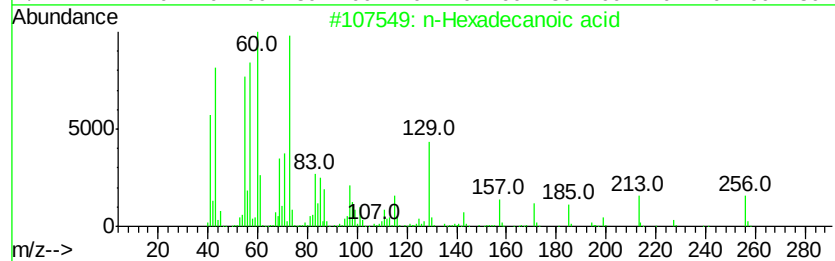
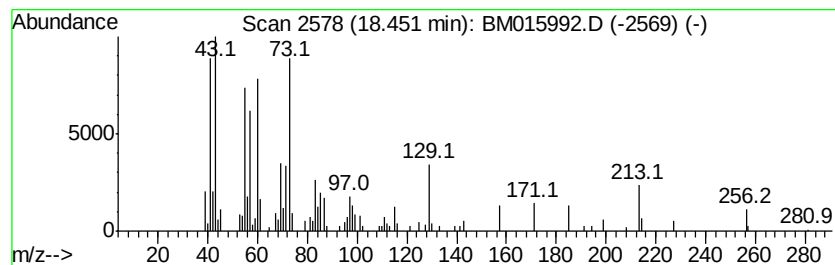
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM071318MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.45	3.71 ng/ul	198543	Phenanthrene-d10	17.61

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Tridecanoic acid	214	C13H26O2	000638-53-9	90
3		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	81
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	76
5		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	70



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM071918\
Data File : BM015992.D
Acq On : 19 Jul 2018 16:22
Operator : SJ/JU
Sample : J4024-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DB1P0

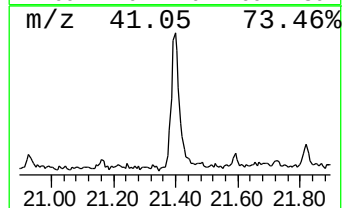
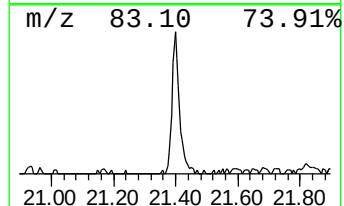
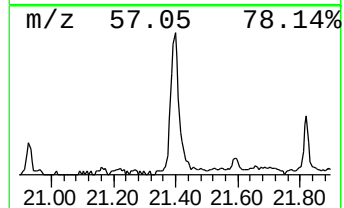
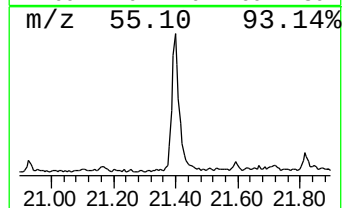
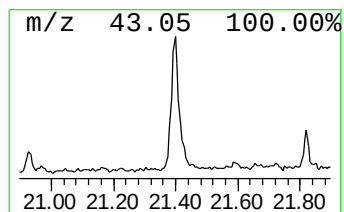
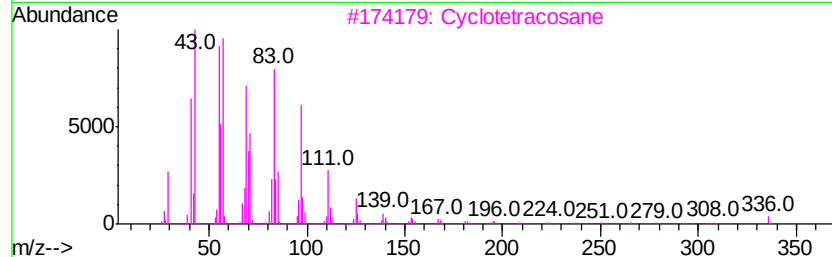
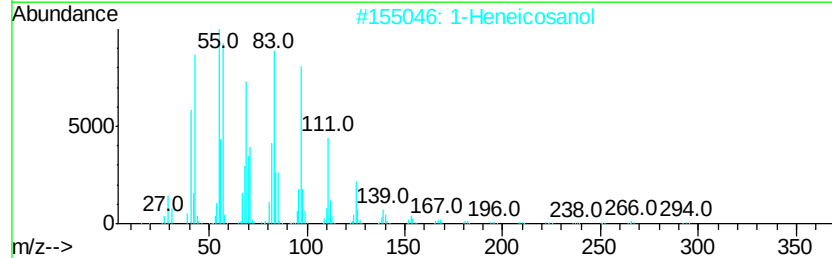
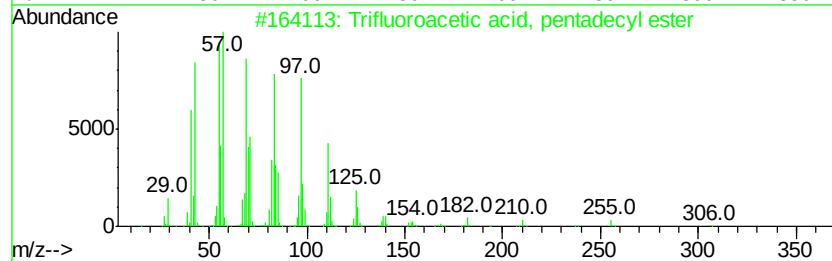
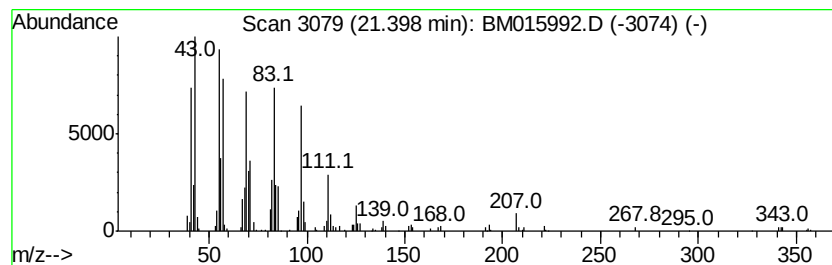
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM071318MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 Trifluoroacetic acid, penta... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.40	4.47 ng/ul	307525	Chrysene-d12	21.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Trifluoroacetic acid, pentadecyl...	324	C17H31F3O2	959010-23-2	94
2		1-Heneicosanol	312	C21H44O	015594-90-8	94
3		Cyclotetracosane	336	C24H48	000297-03-0	93
4		Bromoacetic acid, hexadecyl ester	362	C18H35BrO2	005454-48-8	93
5		1-Tricosene	322	C23H46	018835-32-0	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM071918\
Data File : BM015992.D
Acq On : 19 Jul 2018 16:22
Operator : SJ/JU
Sample : J4024-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DB1P0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM071318MA.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, 1,3-dime...	5.82	20.1	ng/ul	369700	1	8.26	367395	20.0
Diphenyl ether	13.93	21.8	ng/ul	916690	3	14.88	842283	20.0
n-Hexadecanoic acid	18.45	3.7	ng/ul	198543	4	17.61	1070720	20.0
Trifluoroacetic a...	21.40	4.5	ng/ul	307525	5	21.72	1375370	20.0