

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM072718\
 Data File : BM016097.D
 Acq On : 27 Jul 2018 11:43
 Operator : SJ/JU
 Sample : J4172-05
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0D03

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	3.463	27	30	40	rVB	13786	20414	1.74%	0.126%
2	4.146	138	146	154	rVB3	11516	19271	1.65%	0.119%
3	7.363	687	693	711	rBV	216050	389829	33.32%	2.408%
4	7.528	715	721	734	rVV	185262	286329	24.47%	1.769%
5	7.728	750	755	772	rBB	221650	381179	32.58%	2.355%
6	8.204	830	836	849	rBB	260982	431590	36.89%	2.666%
7	8.922	952	958	973	rBV	269253	458162	39.16%	2.831%
8	9.392	1032	1038	1059	rBB	236330	422048	36.07%	2.607%
9	10.116	1155	1161	1174	rBB	145445	262057	22.40%	1.619%
10	10.651	1247	1252	1266	rBV2	218490	396612	33.90%	2.450%
11	11.022	1309	1315	1324	rBV	347877	585097	50.01%	3.615%
12	11.169	1335	1340	1355	rBV	236402	445181	38.05%	2.750%
13	14.133	1839	1844	1849	rBV2	82955	112171	9.59%	0.693%
14	14.233	1856	1861	1866	rBV	548028	728546	62.27%	4.501%
15	14.280	1866	1869	1877	rVB	47638	68137	5.82%	0.421%
16	14.527	1905	1911	1919	rBV	573183	826656	70.66%	5.107%
17	14.833	1956	1963	1970	rBV2	445595	697910	59.65%	4.312%
18	14.898	1970	1974	1979	rVB	17364	24853	2.12%	0.154%
19	15.816	2124	2130	2137	rBV	705276	986133	84.29%	6.092%
20	15.874	2137	2140	2145	rVB	10633	12976	1.11%	0.080%
21	17.562	2421	2427	2431	rBV	510487	732026	62.57%	4.522%
22	17.604	2431	2434	2438	rVV	159895	204398	17.47%	1.263%
23	17.657	2438	2443	2455	rVB	684235	1039152	88.82%	6.420%
24	17.980	2492	2498	2507	rVB3	11537	21595	1.85%	0.133%
25	18.404	2566	2570	2579	rBV2	84955	139501	11.92%	0.862%
26	18.474	2579	2582	2588	rVB2	20411	32447	2.77%	0.200%
27	18.627	2601	2608	2612	rBV	34908	60677	5.19%	0.375%
28	18.915	2653	2657	2661	rBV2	12425	16654	1.42%	0.103%
29	19.515	2757	2759	2763	rBV4	8689	15518	1.33%	0.096%
30	19.592	2767	2772	2779	rVB	357753	485701	41.51%	3.001%
31	19.656	2779	2783	2789	rBV7	41063	58809	5.03%	0.363%
32	19.921	2822	2828	2831	rBV	849538	1169967	100.00%	7.228%
33	19.951	2831	2833	2841	rVB	293175	327925	28.03%	2.026%
34	20.021	2841	2845	2850	rBV2	13489	23588	2.02%	0.146%

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35	20.127	2860	2863	2868	rVB	18277	20389	1.74%	0.126%
36	20.398	2906	2909	2912	rBV4	30380	38523	3.29%	0.238%
37	20.433	2912	2915	2922	rVB	52864	72690	6.21%	0.449%
38	20.533	2928	2932	2936	rBV	42691	60091	5.14%	0.371%
39	20.886	2989	2992	2996	rBV2	38164	40304	3.44%	0.249%
40	21.345	3066	3070	3078	rBV4	101913	208637	17.83%	1.289%
41	21.680	3120	3127	3131	rBV3	602986	1008284	86.18%	6.229%
42	21.715	3131	3133	3139	rVB	148375	183650	15.70%	1.135%
43	21.780	3141	3144	3150	rVB	64091	67850	5.80%	0.419%
44	21.839	3152	3154	3159	rVB3	42178	40401	3.45%	0.250%
45	22.227	3217	3220	3223	rBV	61742	83785	7.16%	0.518%
46	22.721	3299	3304	3310	rBV2	272930	434166	37.11%	2.682%
47	23.239	3388	3392	3397	rVB3	88885	139311	11.91%	0.861%
48	23.327	3403	3407	3412	rBV	128964	272386	23.28%	1.683%
49	23.833	3489	3493	3497	rBV2	136501	216990	18.55%	1.341%
50	23.891	3497	3503	3508	rBV	448399	824162	70.44%	5.092%
51	24.039	3523	3528	3535	rVB	332670	591710	50.57%	3.656%

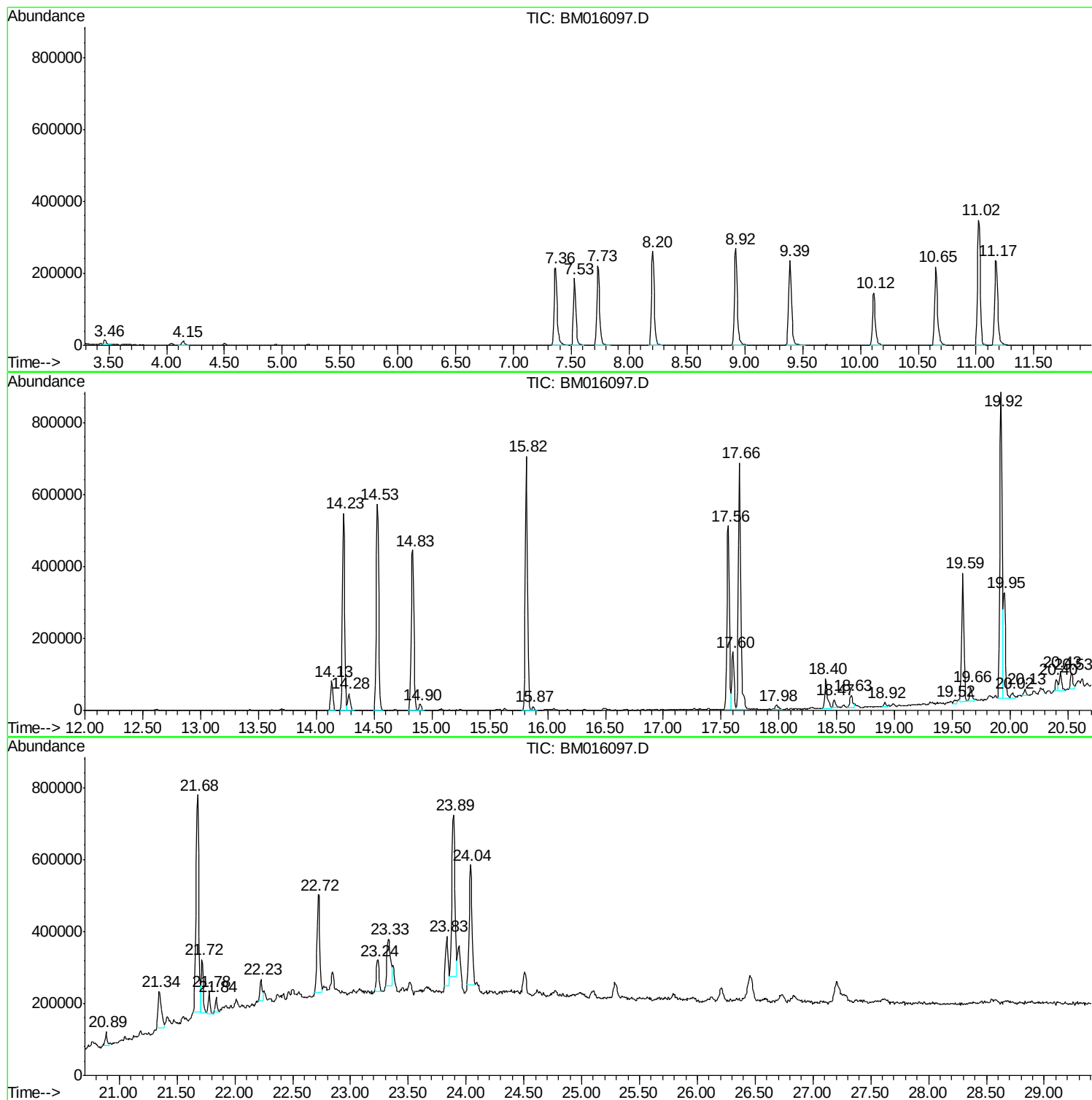
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Quant Title : SVOA CALIBRATION

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



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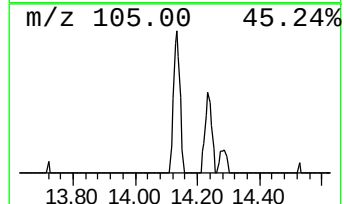
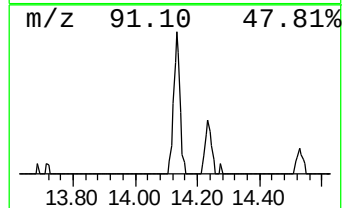
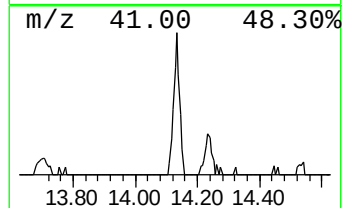
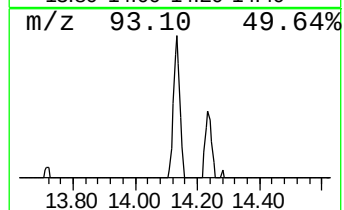
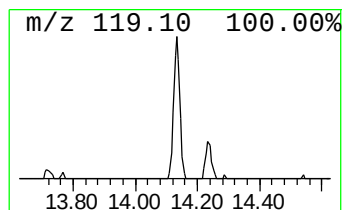
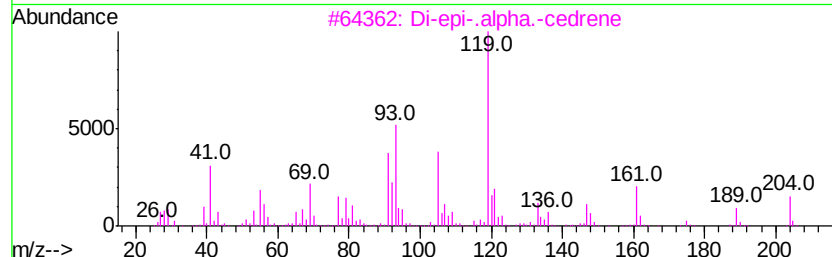
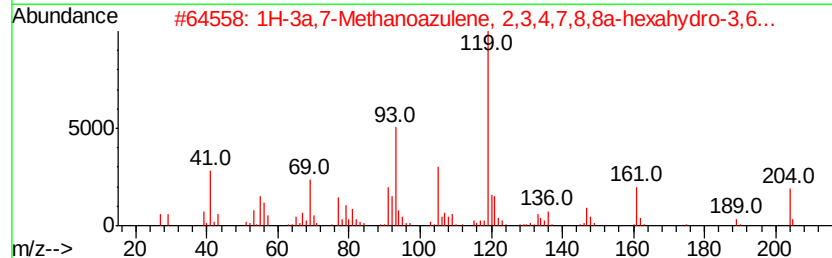
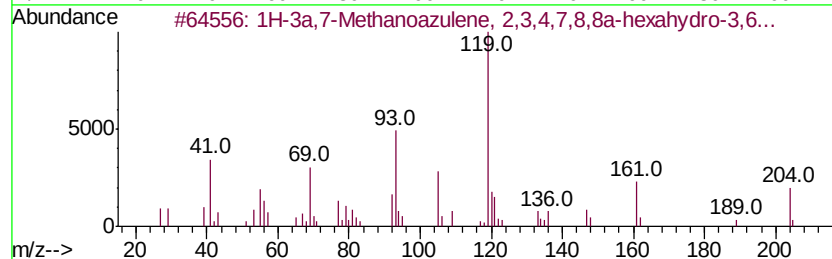
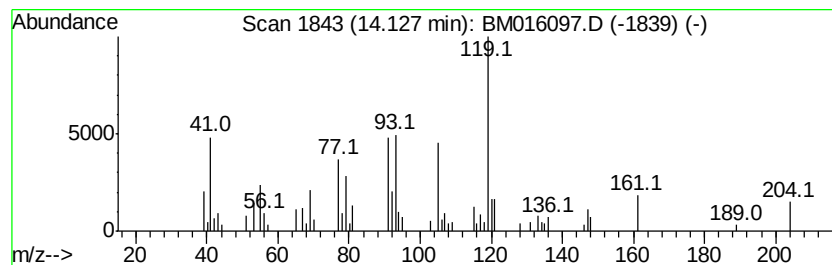
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM071318MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 1H-3a,7-Methanoazulene, 2,3... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.13	3.21 ng/ul	112171	Acenaphthene-d10	14.83
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1H-3a,7-Methanoazulene, 2,3,4,7,...	204 C15H24	000469-61-4	93
2	1H-3a,7-Methanoazulene, 2,3,4,7,...	204 C15H24	000469-61-4	93
3	Di-epi-.alpha.-cedrene	204 C15H24	050894-66-1	90
4	Copaene	204 C15H24	003856-25-5	49
5	s-Triazolo[4,3-a]pyridine, 3,5,7...	161 C9H11N3	004919-15-7	47



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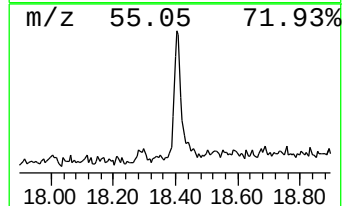
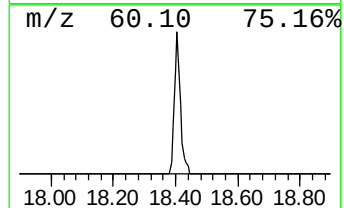
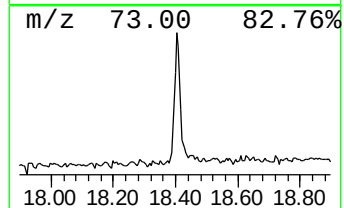
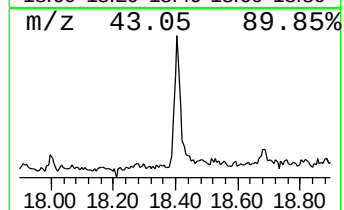
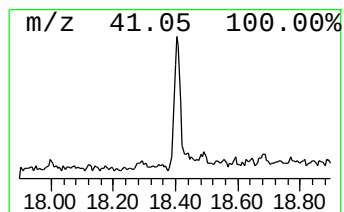
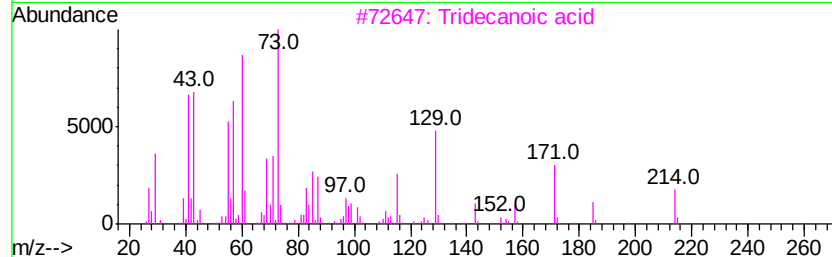
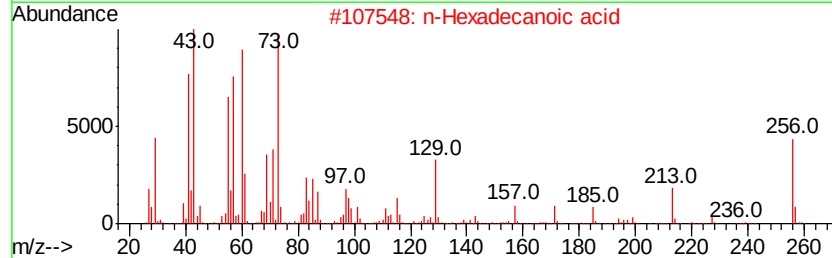
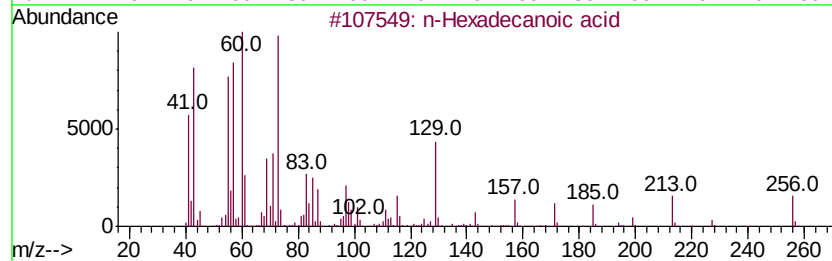
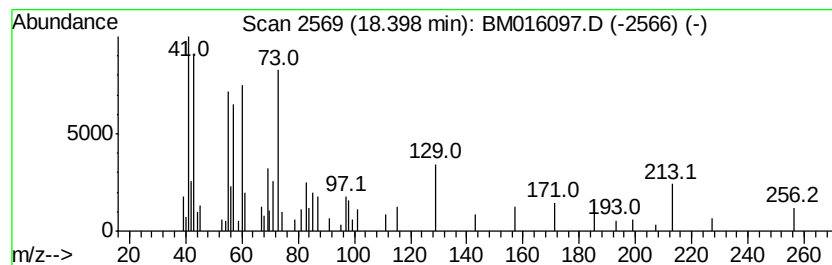
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.40	3.81 ng/ul	139501	Phenanthrene-d10	17.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94
3		Tridecanoic acid	214	C13H26O2	000638-53-9	76
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	74
5		Tetradecanoic acid	228	C14H28O2	000544-63-8	72



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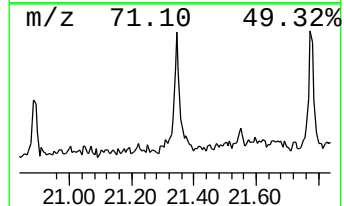
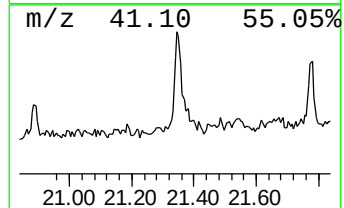
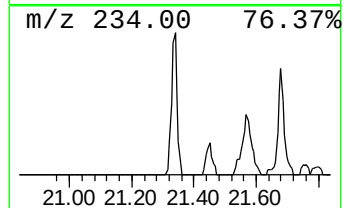
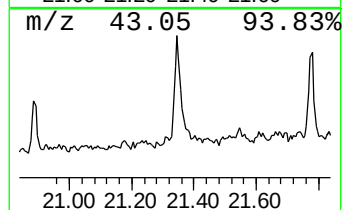
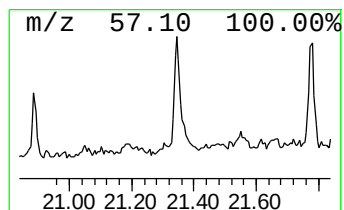
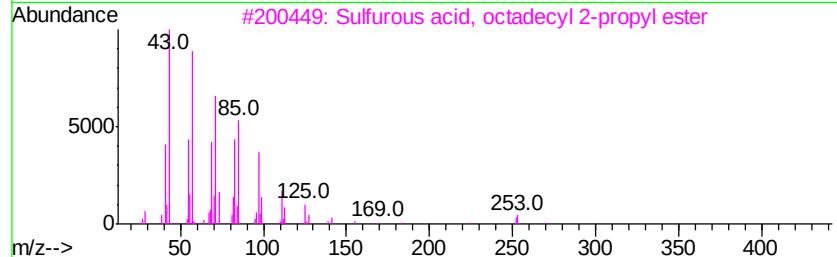
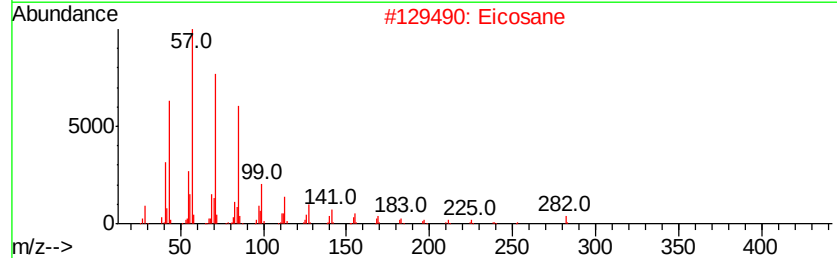
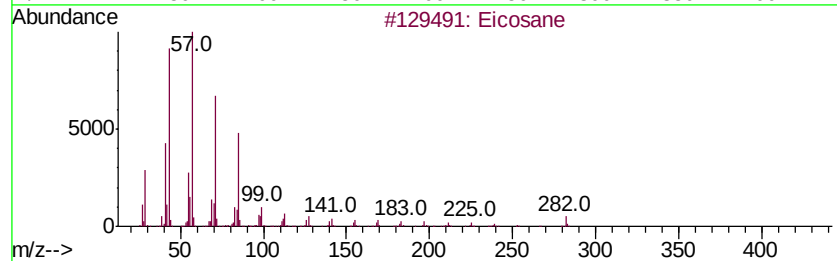
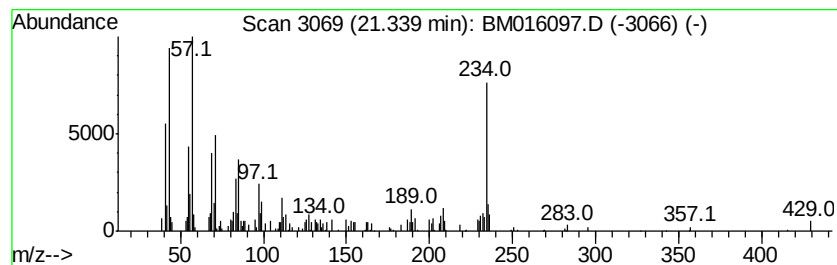
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TIC Library : C:\Database\NIST11.L
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Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.34	4.14 ng/ul	208637	Chrysene-d12	21.68
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Eicosane		282 C20H42	000112-95-8 78
2	Eicosane		282 C20H42	000112-95-8 60
3	Sulfurous acid, octadecyl 2-propyl...		376 C21H44O3S	1000309-12-7 52
4	Oxalic acid, cyclobutyl pentadecyl...		354 C21H38O4	1000309-70-5 52
5	Oxalic acid, isobutyl pentadecyl...		356 C21H40O4	1000309-38-0 52



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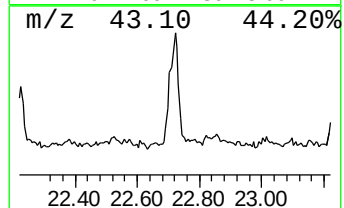
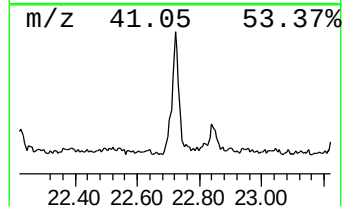
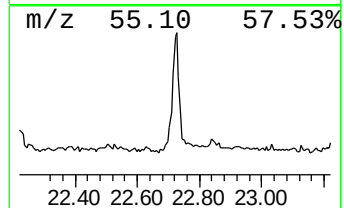
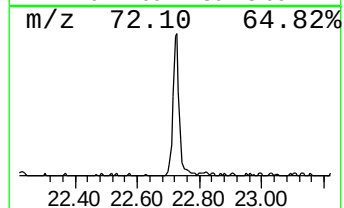
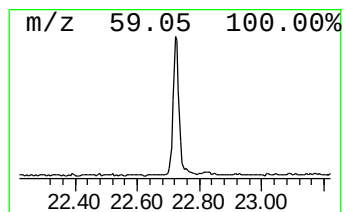
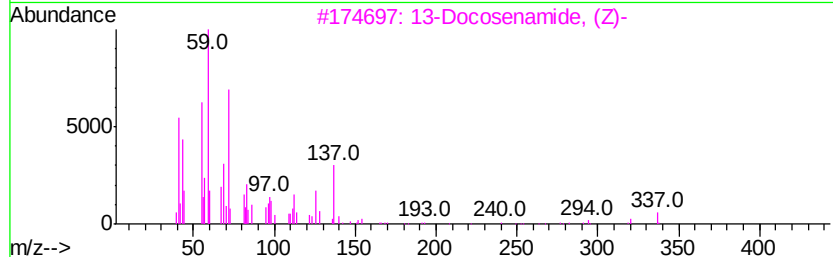
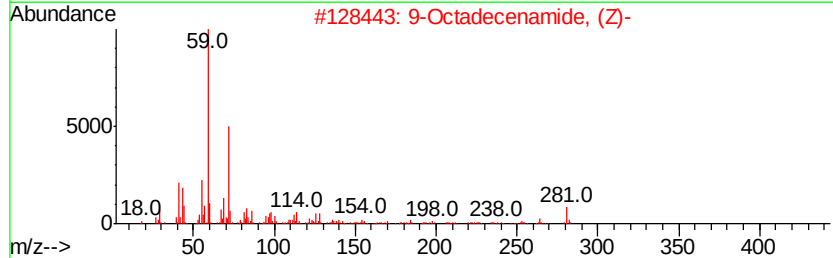
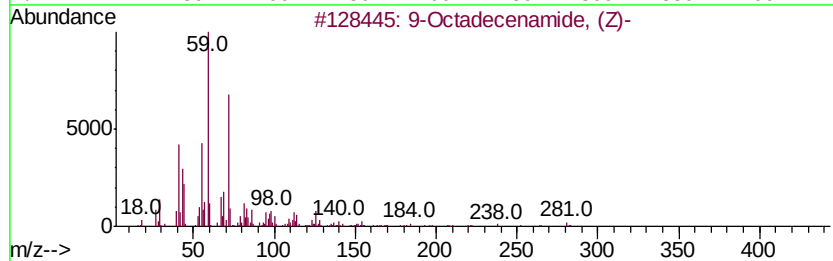
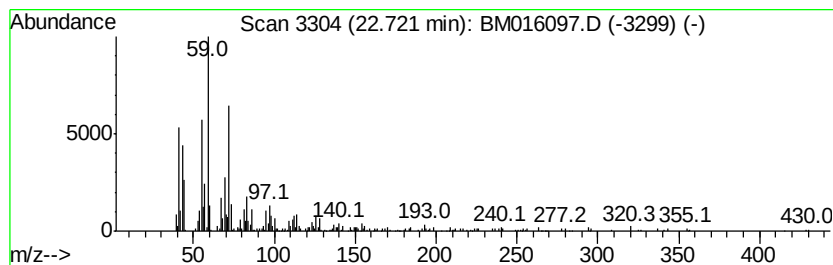
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Peak Number 4 9-Octadecenamide, (Z)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.72	8.61 ng/ul	434166	Chrysene-d12	21.68

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	91		
2	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	74		
3	13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	70		
4	Heptanamide, 4-ethyl-5-methyl-	171	C10H21NO	054789-40-1	64		
5	Dodecanamide	199	C12H25NO	001120-16-7	53		



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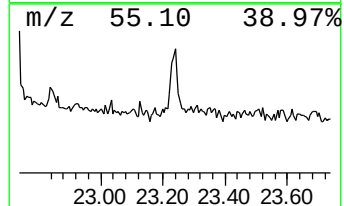
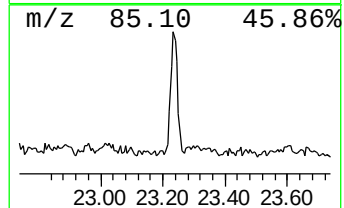
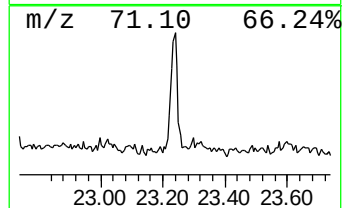
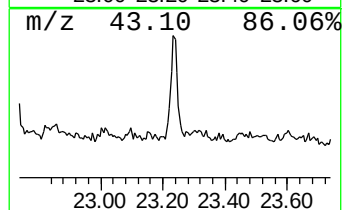
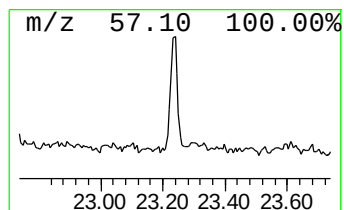
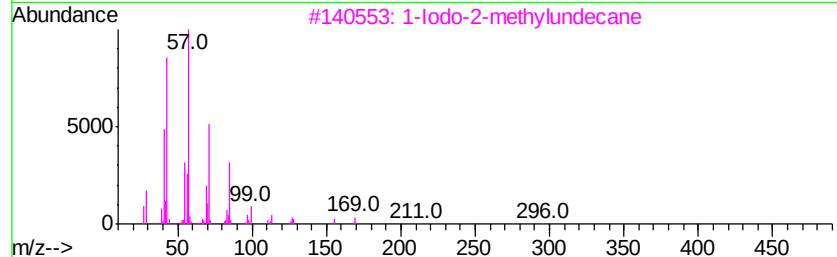
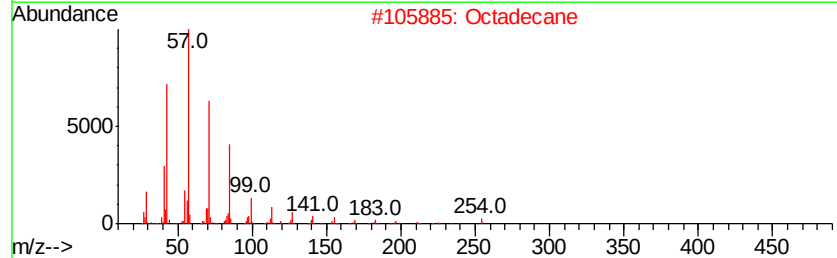
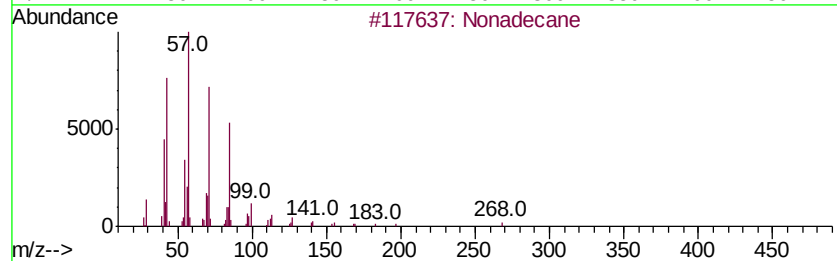
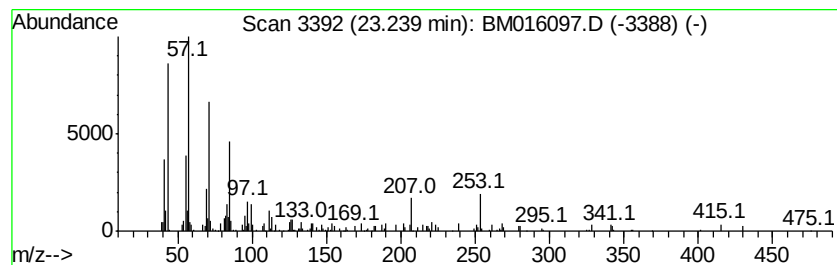
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.24	4.71 ng/ul	139311	Perylene-d12	24.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane	268	C19H40	000629-92-5	78
2		Octadecane	254	C18H38	000593-45-3	70
3		1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	60
4		Nonadecane	268	C19H40	000629-92-5	60
5		Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	53



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM072718\
Data File : BM016097.D
Acq On : 27 Jul 2018 11:43
Operator : SJ/JU
Sample : J4172-05
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0D03

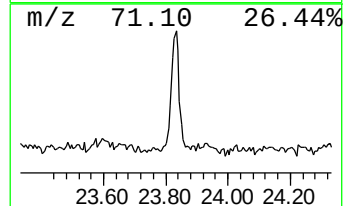
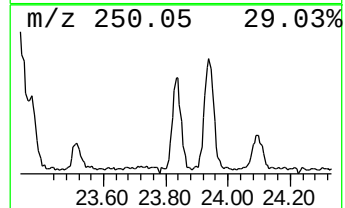
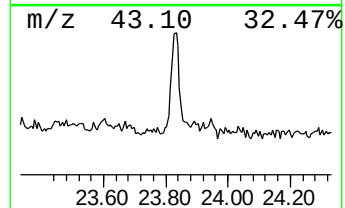
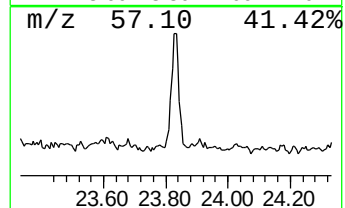
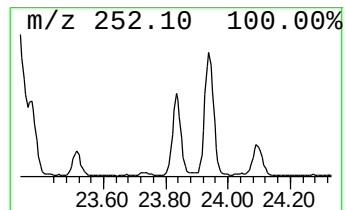
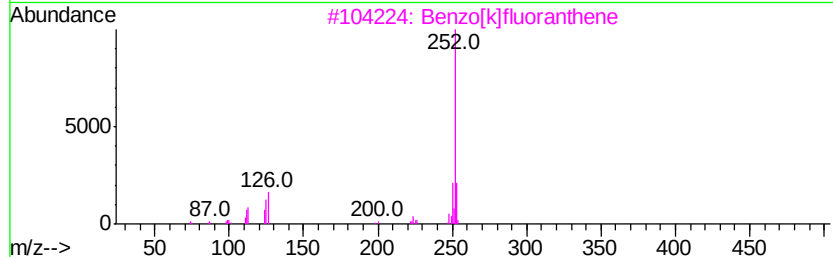
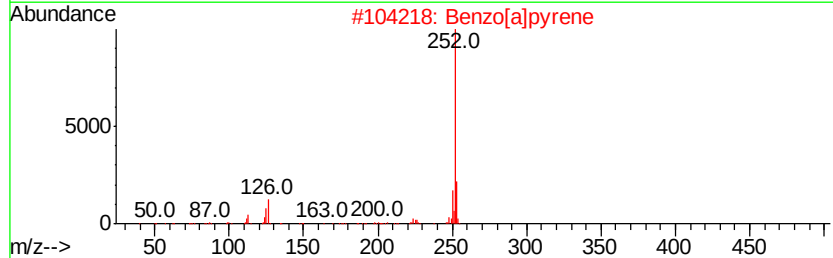
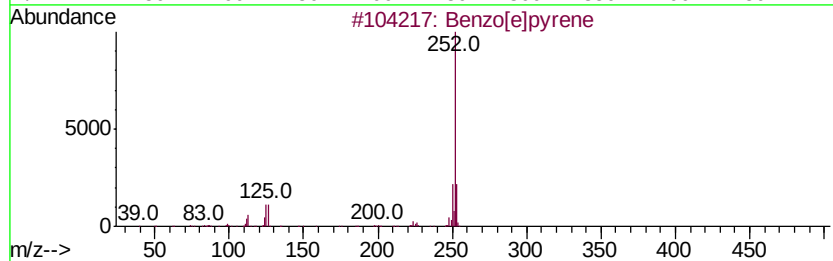
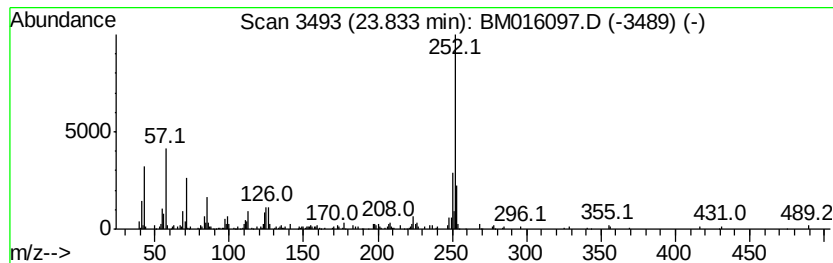
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzo[e]pyrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.83	7.33 ng/ul	216990	Perylene-d12	24.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[e]pyrene	252	C20H12	000192-97-2	96
2		Benzo[a]pyrene	252	C20H12	000050-32-8	93
3		Benzo[k]fluoranthene	252	C20H12	000207-08-9	93
4		Benzo[e]acephenanthrylene	252	C20H12	000205-99-2	90
5		Benzo[k]fluoranthene	252	C20H12	000207-08-9	89



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM072718\
Data File : BM016097.D
Acq On : 27 Jul 2018 11:43
Operator : SJ/JU
Sample : J4172-05
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0D03

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
1H-3a,7-Methanoaz...	14.13	3.2	ng/ul	112171	3	14.83	697910	20.0
n-Hexadecanoic acid	18.40	3.8	ng/ul	139501	4	17.56	732026	20.0
(DEL) Alkane: Str...	21.34	4.1	ng/ul	208637	5	21.68	1008280	20.0
9-Octadecenamide,...	22.72	8.6	ng/ul	434166	5	21.68	1008280	20.0
(DEL) Alkane: Str...	23.24	4.7	ng/ul	139311	6	24.04	591710	20.0
Benzo[e]pyrene	23.83	7.3	ng/ul	216990	6	24.04	591710	20.0