

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM060117\
 Data File : BM010362.D
 Acq On : 01 Jun 2017 19:12
 Operator : SJ/MA
 Sample : I3163-05
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 JHSF5

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-BM052717.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	7.104	673	683	693	rBV	970945	1609450	23.42%	1.634%
2	7.257	701	709	721	rBV	716255	1193821	17.37%	1.212%
3	7.457	735	743	755	rVB	1166069	1854430	26.99%	1.883%
4	7.916	813	821	830	rBV	1560369	2458577	35.78%	2.496%
5	8.640	936	944	957	rBV	935954	1562996	22.75%	1.587%
6	9.104	1014	1023	1031	rBV	1030903	1736284	25.27%	1.763%
7	9.816	1137	1144	1154	rBV	904630	1532604	22.30%	1.556%
8	10.345	1225	1234	1248	rBV	1356936	2386122	34.73%	2.423%
9	10.722	1291	1298	1308	rBV	1854520	3111017	45.27%	3.159%
10	10.886	1319	1326	1338	rBV	431400	843193	12.27%	0.856%
11	13.969	1843	1850	1854	rBV	2502999	3396854	49.43%	3.449%
12	14.010	1854	1857	1864	rVB	451740	626166	9.11%	0.636%
13	14.251	1891	1898	1906	rBV	2675480	3885529	56.55%	3.945%
14	14.551	1943	1949	1959	rBV2	2650641	4149739	60.39%	4.213%
15	14.763	1980	1985	1986	rBV3	93966	121160	1.76%	0.123%
16	14.792	1986	1990	2004	rVB2	413114	865134	12.59%	0.878%
17	15.545	2111	2118	2126	rVB	3408300	4892428	71.20%	4.967%
18	15.669	2133	2139	2145	rBV	843223	1151061	16.75%	1.169%
19	15.780	2154	2158	2164	rVB5	48396	78869	1.15%	0.080%
20	16.198	2223	2229	2235	rBV2	103000	197586	2.88%	0.201%
21	16.657	2304	2307	2311	rBV4	75252	96311	1.40%	0.098%
22	16.986	2359	2363	2367	rBV6	54775	74929	1.09%	0.076%
23	17.151	2387	2391	2392	rBV3	94289	97569	1.42%	0.099%
24	17.174	2392	2395	2399	rVV4	135144	178878	2.60%	0.182%
25	17.215	2399	2402	2408	rVB7	66762	88525	1.29%	0.090%
26	17.298	2408	2416	2427	rBV	3400746	4988180	72.59%	5.064%
27	17.398	2427	2433	2440	rBV	3738358	5217329	75.93%	5.297%
28	18.033	2535	2541	2546	rBV6	124375	218117	3.17%	0.221%
29	18.092	2547	2551	2555	rBV7	59126	78270	1.14%	0.079%
30	18.145	2556	2560	2567	rBV	775987	1147324	16.70%	1.165%
31	18.327	2587	2591	2596	rVB8	110208	137449	2.00%	0.140%
32	18.386	2597	2601	2604	rBV5	101462	122314	1.78%	0.124%
33	18.457	2606	2613	2622	rVB2	660146	1011550	14.72%	1.027%
34	18.557	2625	2630	2636	rVB10	89589	157165	2.29%	0.160%

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Integration Parameters: LSCINT.P

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Max Peaks: 100
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35	18.610	2636	2639	2641	rBV4	63541	69461	1.01%	0.071%
36	18.727	2655	2659	2663	rBV4	187758	257148	3.74%	0.261%
37	18.804	2668	2672	2673	rBV4	48500	75241	1.09%	0.076%
38	18.880	2681	2685	2691	rBV	678156	855027	12.44%	0.868%
39	19.045	2710	2713	2715	rBV4	72728	101614	1.48%	0.103%
40	19.080	2715	2719	2723	rVV3	543167	686952	10.00%	0.697%
41	19.121	2723	2726	2735	rVB7	138646	246229	3.58%	0.250%
42	19.415	2773	2776	2778	rBV2	325497	393296	5.72%	0.399%
43	19.445	2778	2781	2788	rVB2	1059712	1364194	19.85%	1.385%
44	19.686	2816	2822	2835	rBV	5028495	6507230	94.70%	6.607%
45	20.151	2897	2901	2906	rBV4	257423	330932	4.82%	0.336%
46	20.209	2907	2911	2915	rVB6	127682	194953	2.84%	0.198%
47	20.392	2937	2942	2950	rBV	749199	1207180	17.57%	1.226%
48	20.492	2956	2959	2972	rBV7	529603	1126062	16.39%	1.143%
49	20.962	3036	3039	3043	rBV5	225102	327702	4.77%	0.333%
50	21.009	3045	3047	3053	rVB6	220796	278681	4.06%	0.283%
51	21.103	3060	3063	3064	rBV2	466766	469804	6.84%	0.477%
52	21.133	3064	3068	3083	rVB2	2442053	4413631	64.23%	4.481%
53	21.427	3115	3118	3120	rBV3	474520	620082	9.02%	0.630%
54	21.462	3120	3124	3132	rVB	5269047	6638282	96.61%	6.740%
55	21.980	3209	3212	3216	rBV	454773	525326	7.65%	0.533%
56	22.027	3217	3220	3228	rVV6	485844	963731	14.03%	0.978%
57	22.703	3333	3335	3341	rVB4	487889	642496	9.35%	0.652%
58	22.974	3378	3381	3386	rVB	643352	827023	12.04%	0.840%
59	23.656	3491	3497	3504	rVB	3788944	6871468	100.00%	6.977%
60	23.809	3517	3523	3531	rVB	3478642	6826536	99.35%	6.931%
61	24.327	3606	3611	3627	rBV	1504848	4406756	64.13%	4.474%

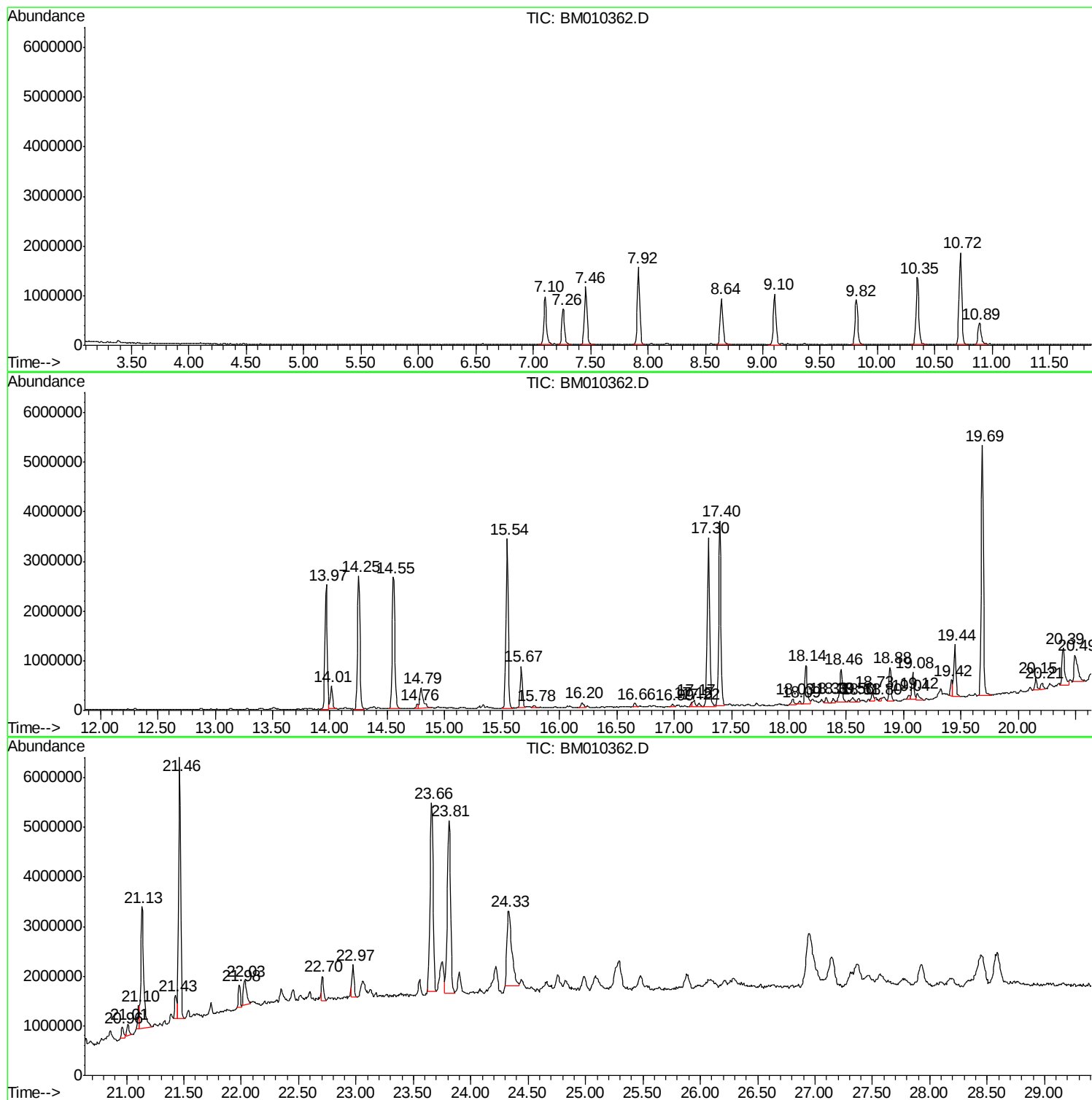
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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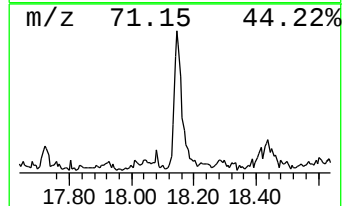
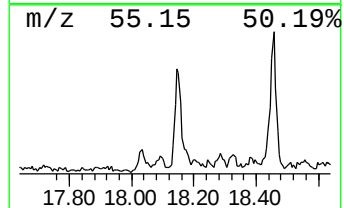
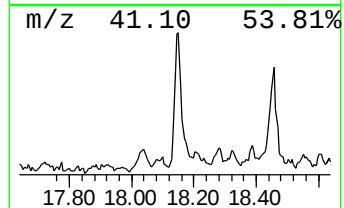
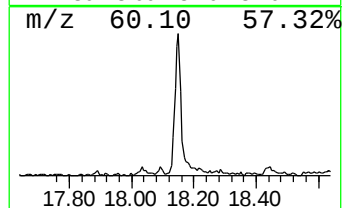
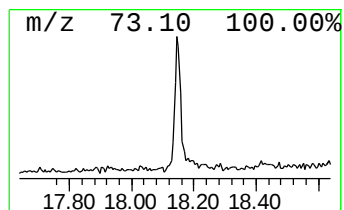
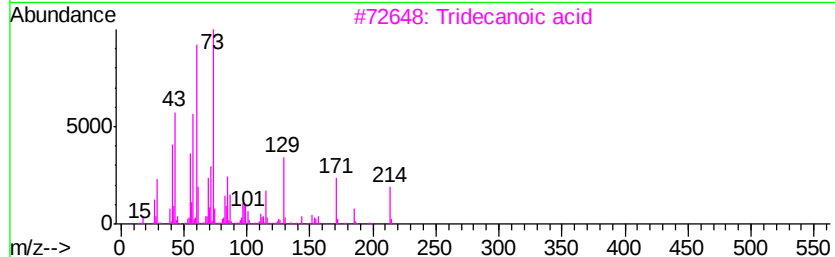
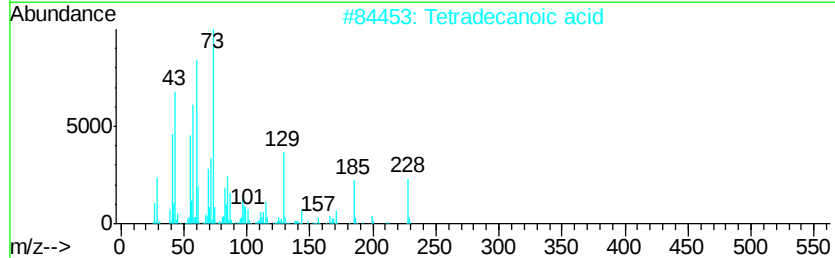
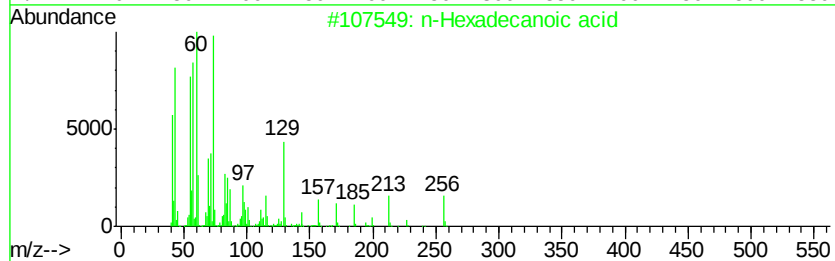
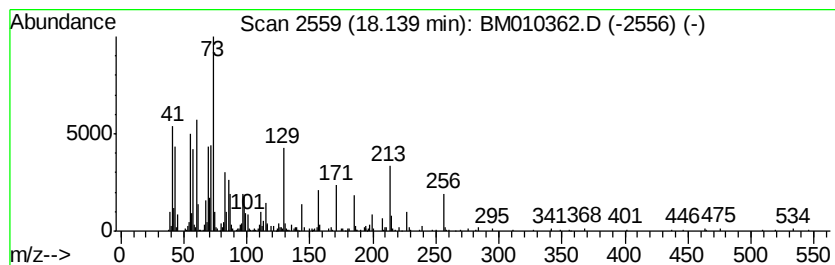
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM052717.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.14	4.60 ng/ul	1147320	Phenanthrene-d10	17.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	83
3		Tridecanoic acid	214	C13H26O2	000638-53-9	64
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	60
5		Tridecanoic acid	214	C13H26O2	000638-53-9	52



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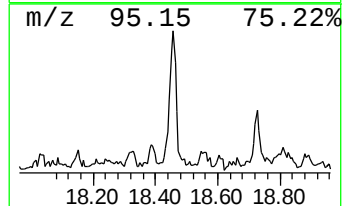
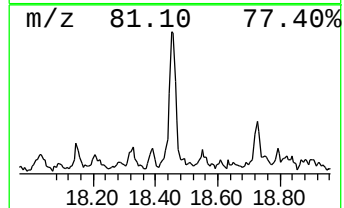
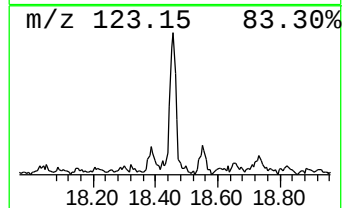
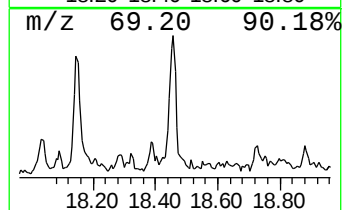
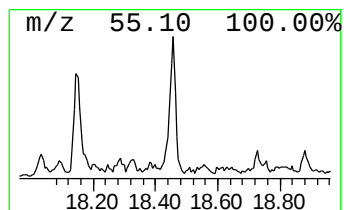
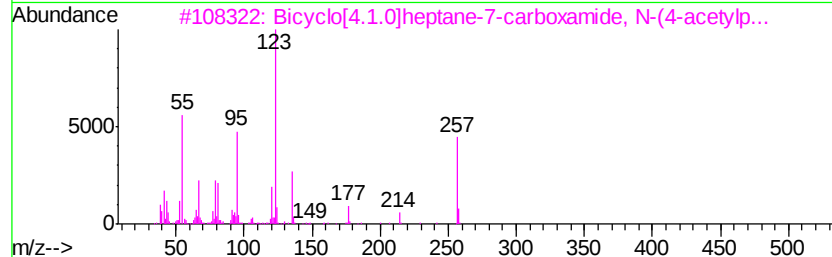
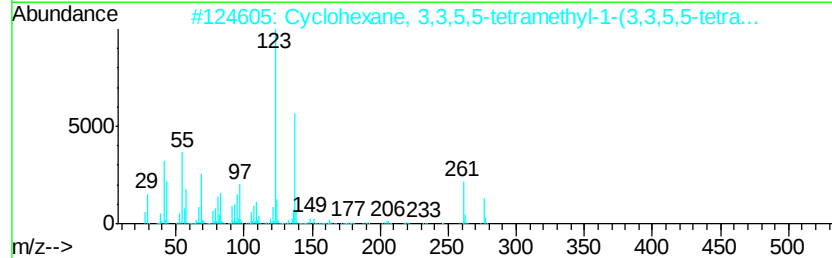
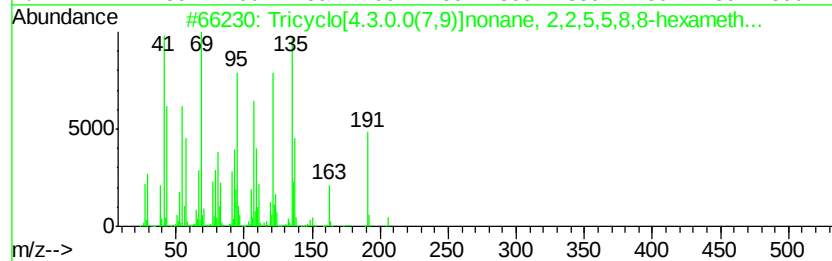
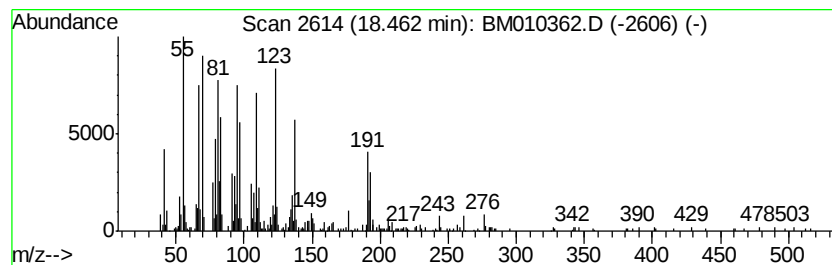
Instrument :
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM052717.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 unknown-01 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.46	4.06 ng/ul	1011550	Phenanthrene-d10	17.30
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Tricyclo[4.3.0.0(7,9)]nonane, 2,...	206 C15H26	054832-82-5 25
2		Cyclohexane, 3,3,5,5-tetramethyl...	276 C20H36	1000157-88-6 25
3		Bicyclo[4.1.0]heptane-7-carboxam...	257 C16H19N02	1000278-06-2 15
4		Divinylbis(cyclopropyl)silane	164 C10H16Si	1000109-95-2 14
5		Silane, [(11-fluoroundecyl)oxy]t...	262 C14H31FOSi	026305-81-7 14



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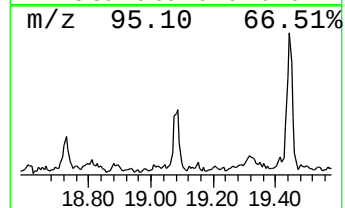
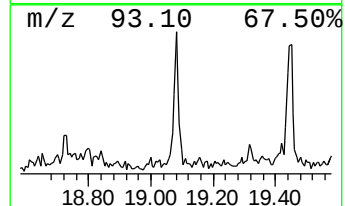
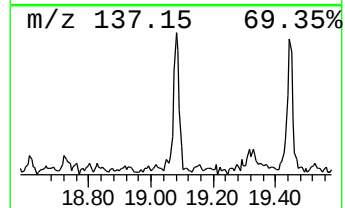
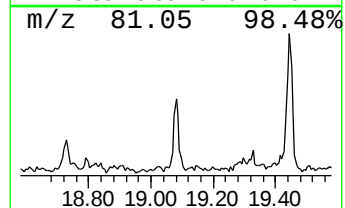
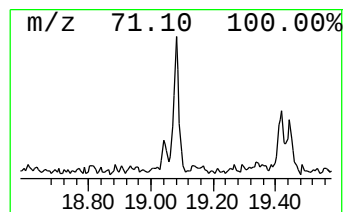
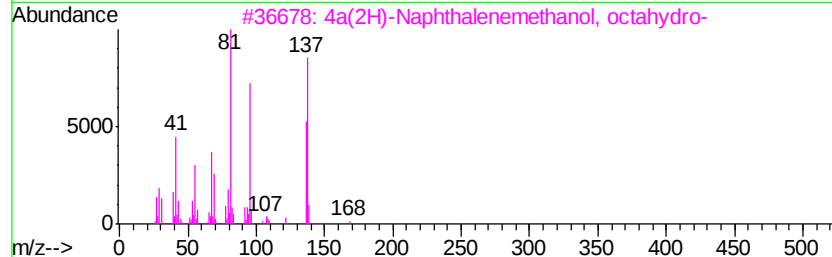
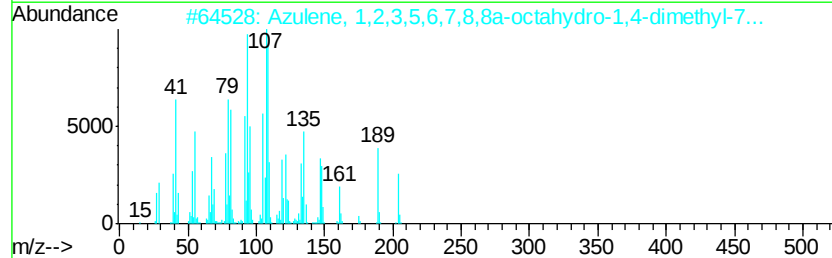
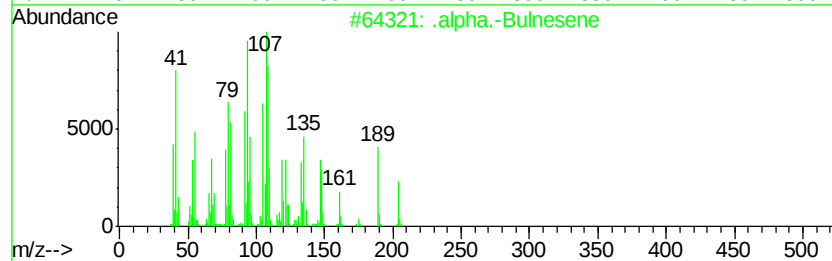
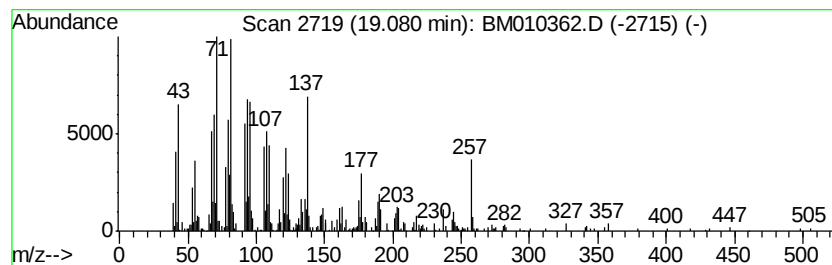
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown-02 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.08	2.75 ng/ul	686952	Phenanthrene-d10	17.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	.alpha.-Bulnesene	204	C15H24	1000374-19-9	45
2		Azulene, 1,2,3,5,6,7,8,8a-octahy...	204	C15H24	003691-11-0	44
3		4a(2H)-Naphthalenemethanol, octa...	168	C11H20O	099992-19-5	38
4		2-Bromomethyl-1-isopropenyl-3-me...	216	C10H17Br	1000187-07-8	38
5		1,3,3-Trimethyl-2-hydroxymethyl-...	222	C15H26O	1000144-10-7	38



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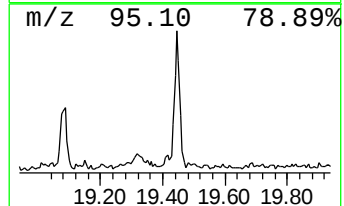
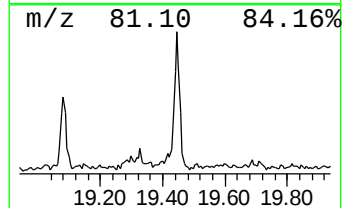
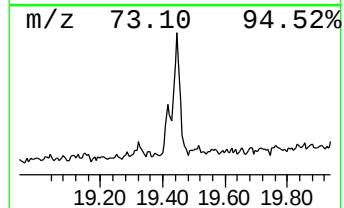
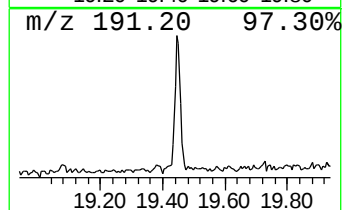
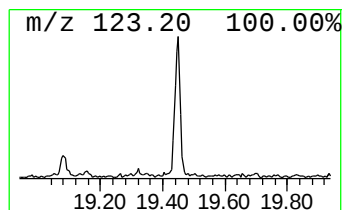
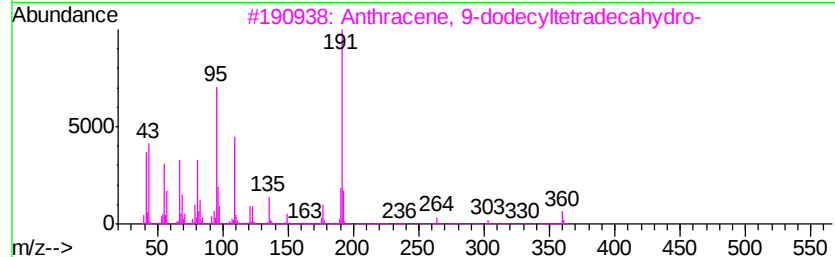
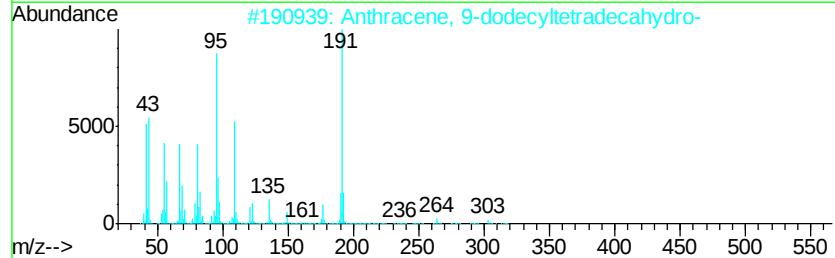
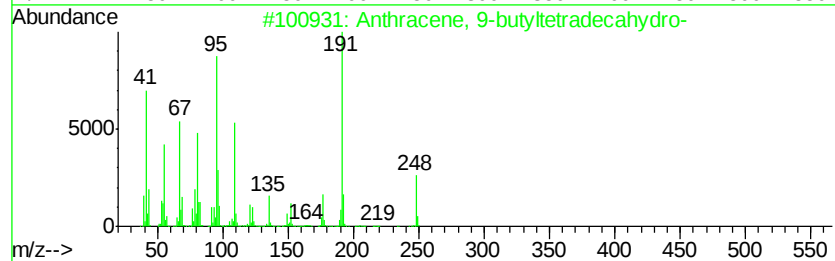
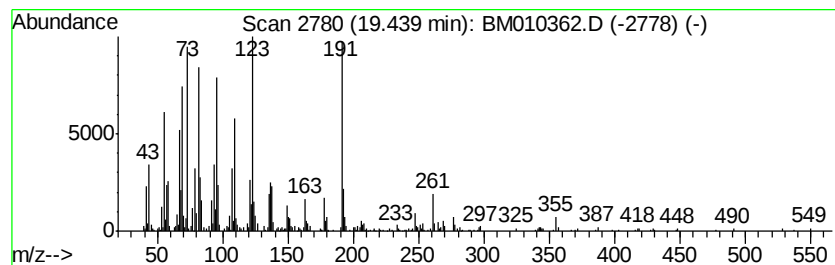
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 5 Anthracene, 9-butyltetradec... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.44	4.11 ng/ul	1364190	Chrysene-d12	21.46

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Anthracene, 9-butyltetradecahydro-	248	C18H32	055133-89-6	64
2		Anthracene, 9-dodecyltetradecahy...	360	C26H48	055401-75-7	43
3		Anthracene, 9-dodecyltetradecahy...	360	C26H48	055401-75-7	43
4		1H-Indene, 1-(1,5-dimethyl-2-hex...	248	C18H32	054411-95-9	30
5		Cyclodecene, 3-bromo-	216	C10H17Br	056325-56-5	22



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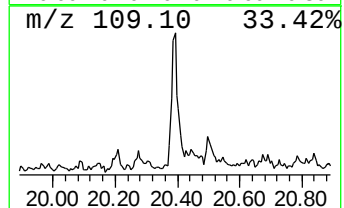
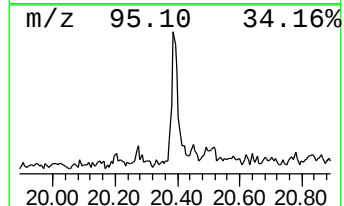
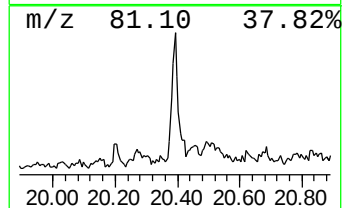
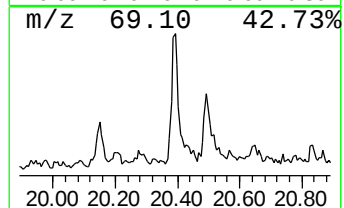
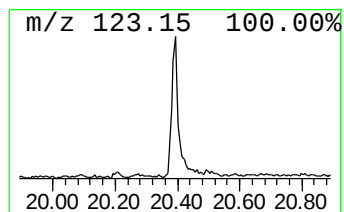
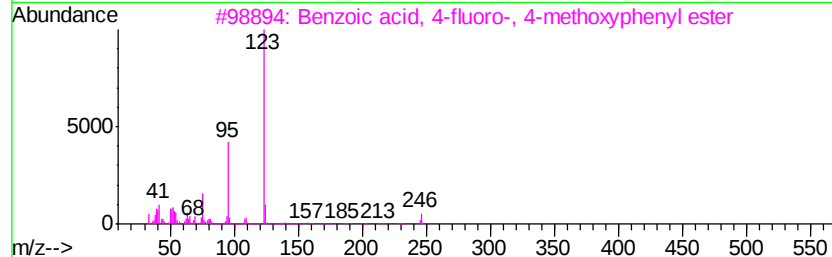
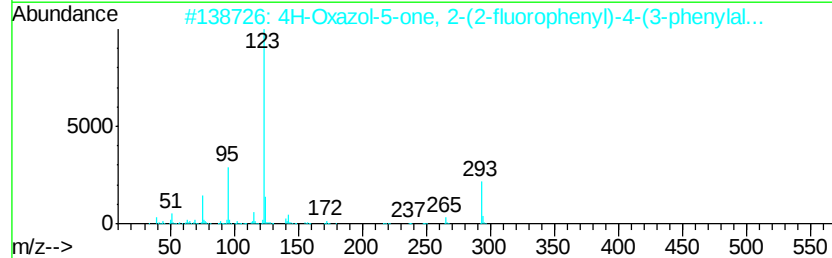
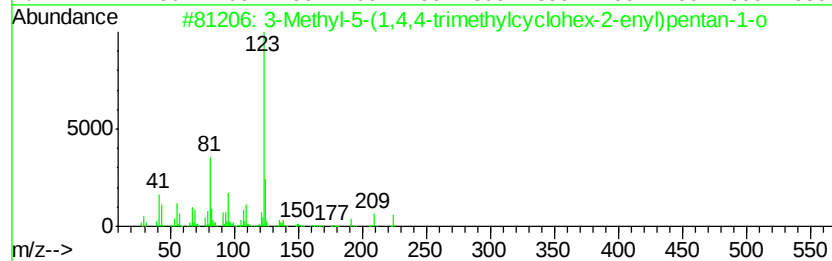
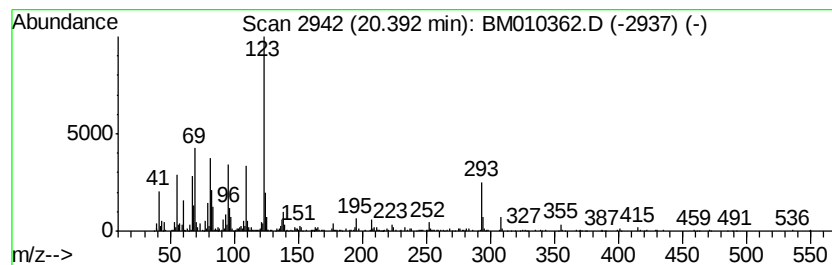
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 3-Methyl-5-(1,4,4-trimethyl... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.39	3.64 ng/ul	1207180	Chrysene-d12	21.46

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Methyl-5-(1,4,4-trimethylcyclo...	224	C15H28O	1000187-09-1	64
2		4H-Oxazol-5-one, 2-(2-fluorophen...	293	C18H12FN02	1000310-91-1	64
3		Benzoic acid, 4-fluoro-, 4-metho...	246	C14H11F03	080079-00-1	50
4		3,3,5,5-Tetramethylcyclohexyl me...	236	C11H22F02P	1000298-38-5	50
5		Ether, [1-(p-fluorophenyl)ethyl]...	212	C12H17F02	1000149-23-5	50



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM060117\
Data File : BM010362.D
Acq On : 01 Jun 2017 19:12
Operator : SJ/MA
Sample : I3163-05
Misc :
ALS Vial : 16 Sample Multiplier: 1

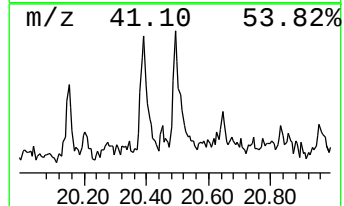
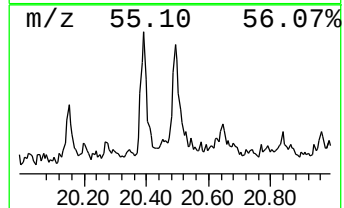
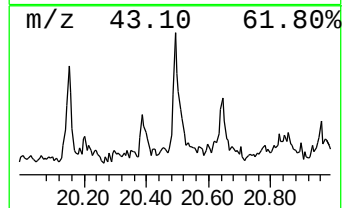
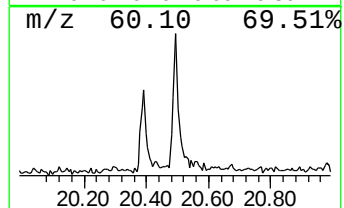
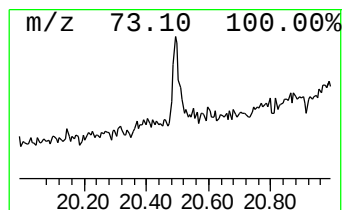
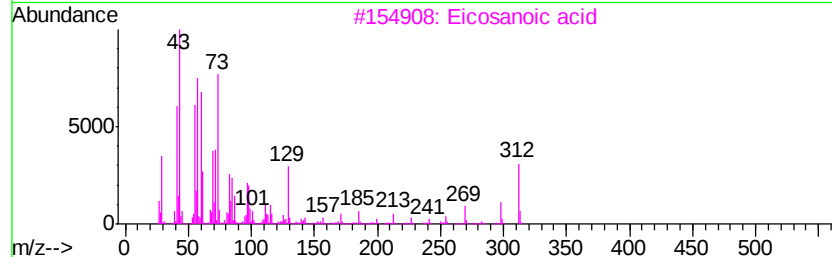
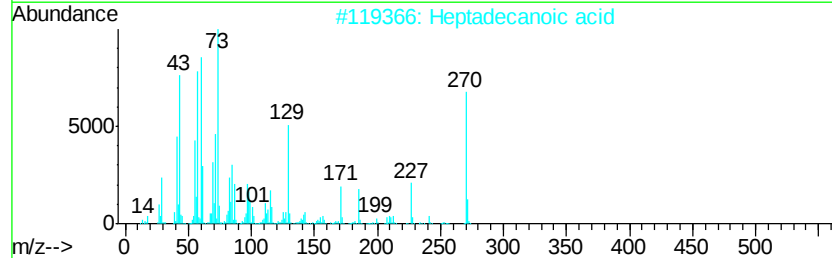
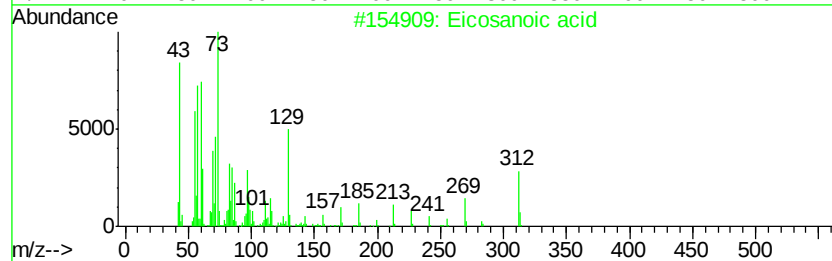
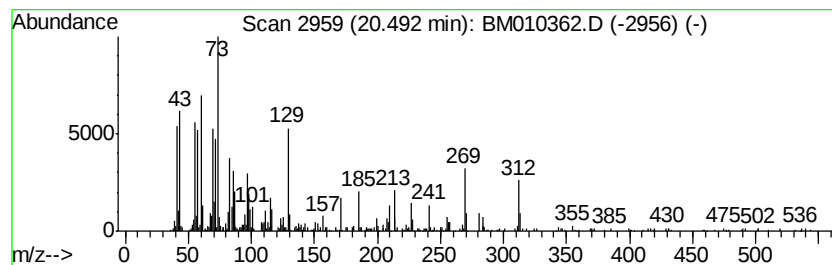
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM052717.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 Eicosanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.49	3.39 ng/ul	1126060	Chrysene-d12	21.46
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Eicosanoic acid	312 C20H40O2	000506-30-9	94
2	Heptadecanoic acid	270 C17H34O2	000506-12-7	89
3	Eicosanoic acid	312 C20H40O2	000506-30-9	89
4	Eicosanoic acid	312 C20H40O2	000506-30-9	89
5	Heptadecanoic acid	270 C17H34O2	000506-12-7	70



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM060117\
Data File : BM010362.D
Acq On : 01 Jun 2017 19:12
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Sample : I3163-05
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Instrument :
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ClientSampleId :
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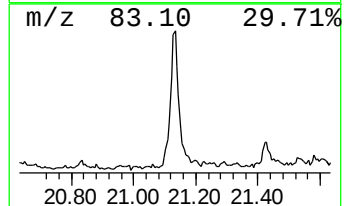
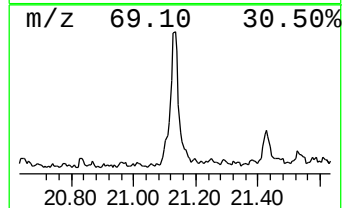
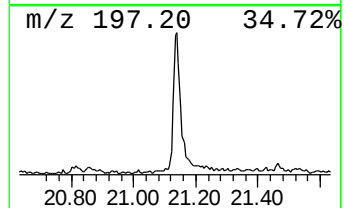
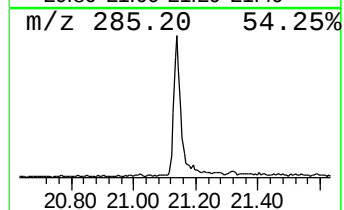
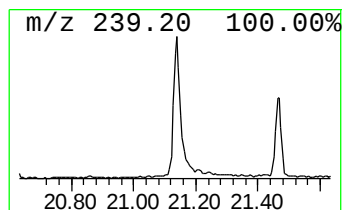
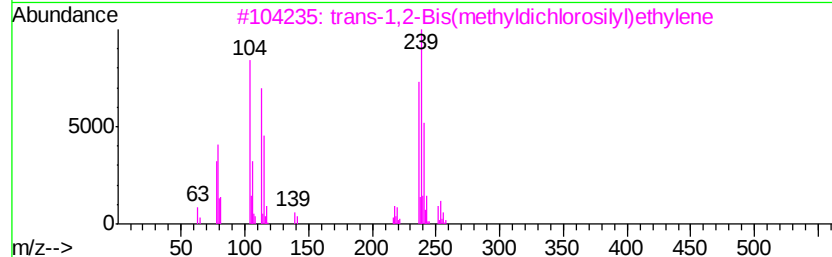
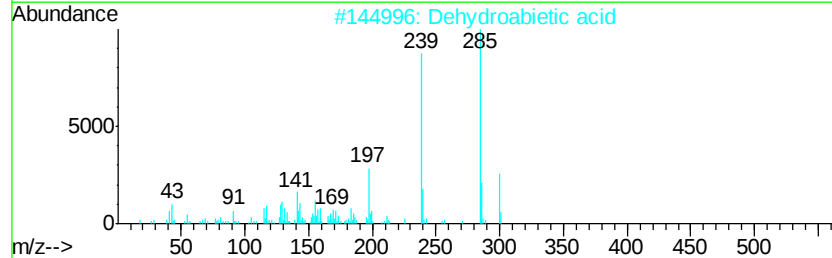
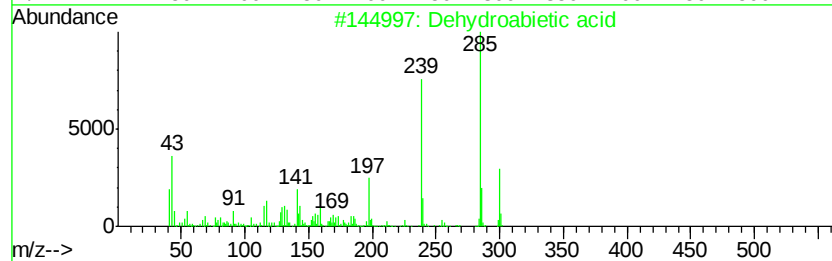
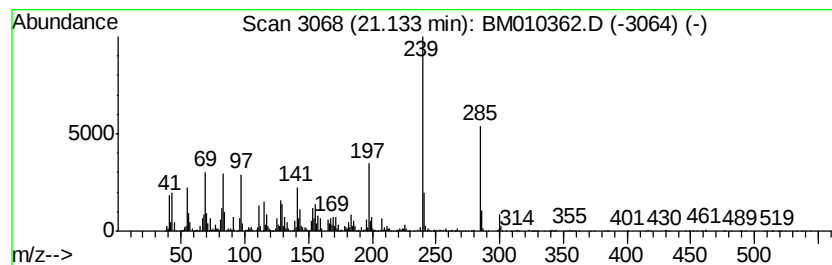
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Dehydroabietic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.13	13.30 ng/ul	4413630	Chrysene-d12	21.46

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dehydroabietic acid	300	C20H28O2	001740-19-8	93	
2	Dehydroabietic acid	300	C20H28O2	001740-19-8	89	
3	trans-1,2-Bis(methyldichlorosilyl)...	252	C4H8Cl4Si2	065899-10-7	74	
4	1-Phenanthrenecarboxylic acid, 1...	300	C20H28O2	005155-70-4	74	
5	Butanoic acid, 2-(cyano)(2,4,6-t...	300	C17H20N2O3	1000267-71-7	68	



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM060117\
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Acq On : 01 Jun 2017 19:12
Operator : SJ/MA
Sample : I3163-05
Misc :
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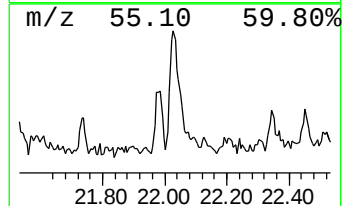
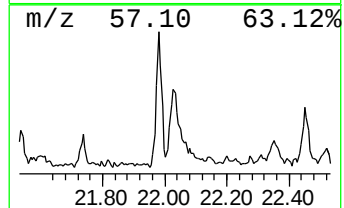
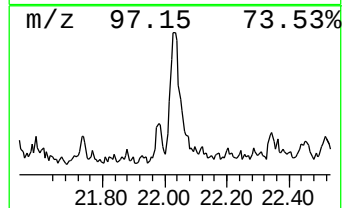
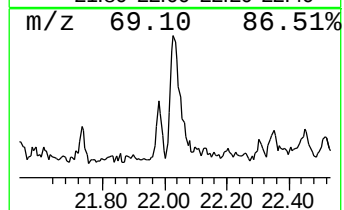
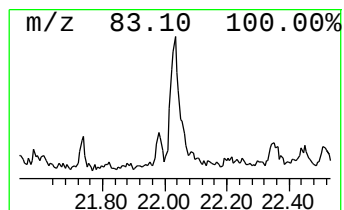
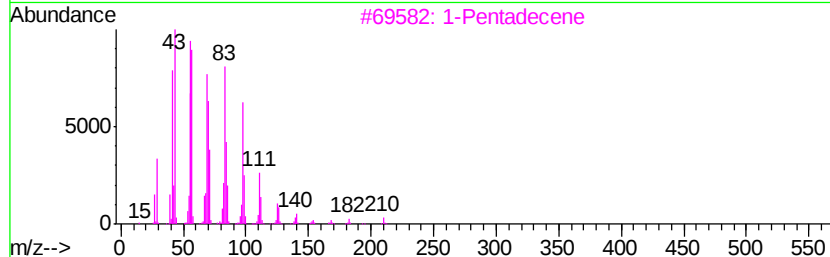
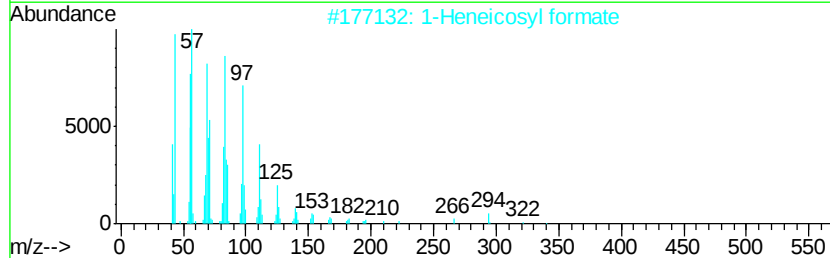
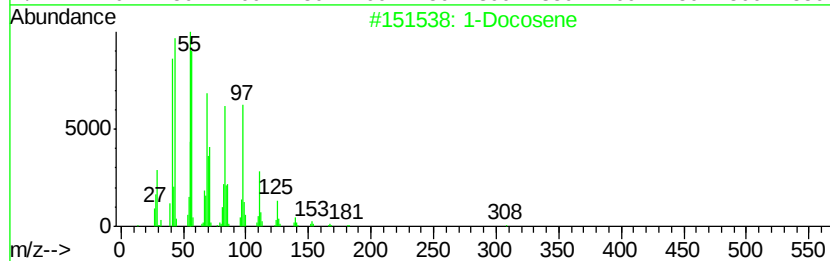
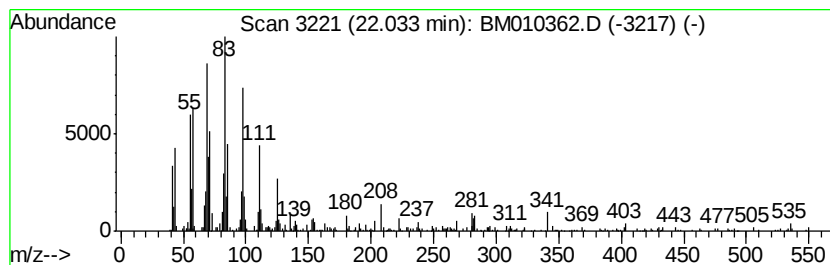
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Cyclic22.03 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.03	2.90 ng/ul	963731	Chrysene-d12	21.46
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1-Docosene		308 C22H44	001599-67-3 91
2	1-Heneicosyl formate		340 C22H44O2	077899-03-7 90
3	1-Pentadecene		210 C15H30	013360-61-7 90
4	3-Eicosene, (E)-		280 C20H40	074685-33-9 87
5	Heptadecyl heptafluorobutyrate		452 C21H35F7O2	959085-66-6 87



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM060117\
Data File : BM010362.D
Acq On : 01 Jun 2017 19:12
Operator : SJ/MA
Sample : I3163-05
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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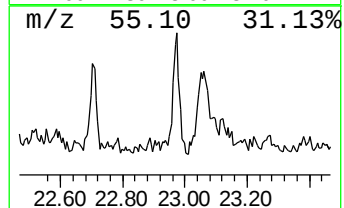
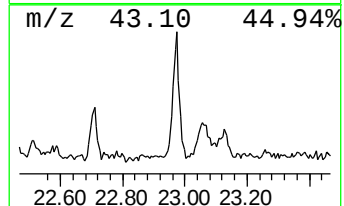
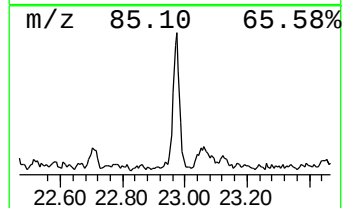
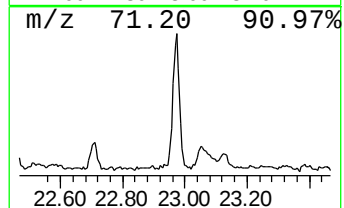
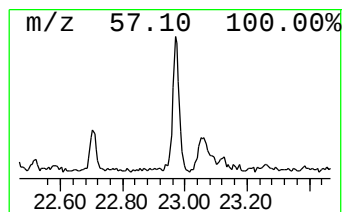
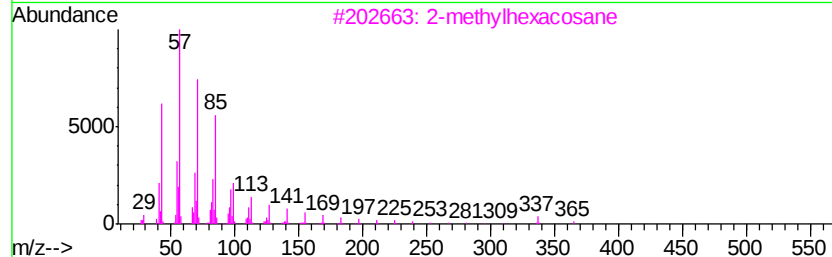
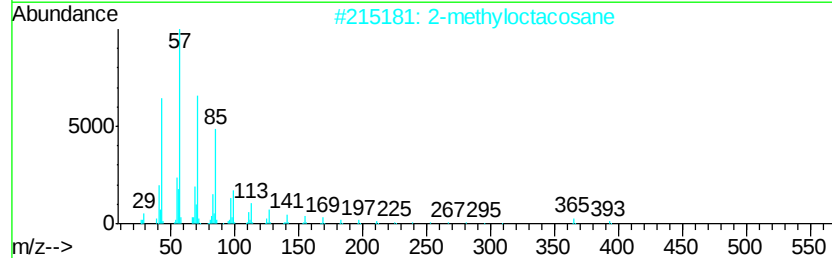
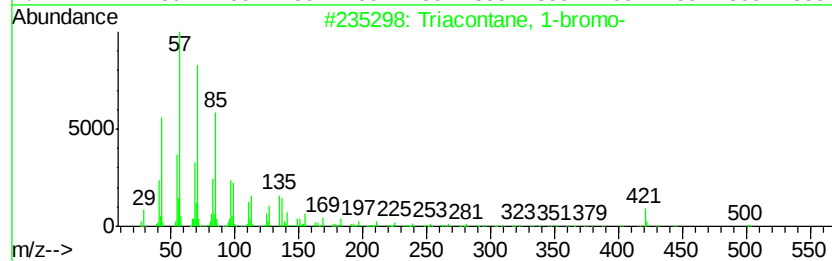
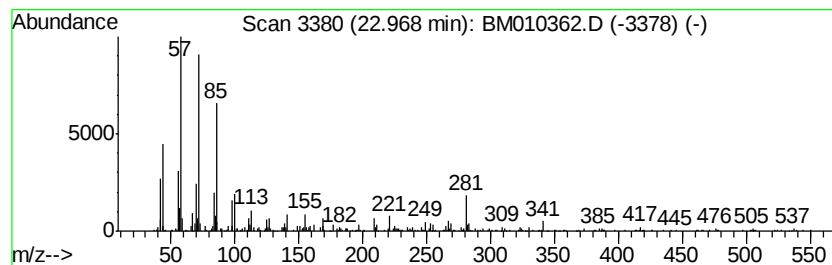
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Triacontane, 1-bromo- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.97	2.42 ng/ul	827023	Perylene-d12	23.81

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Triacontane, 1-bromo-	500	C30H61Br	004209-22-7	87
2			2-methyloctacosane	408	C29H60	1000376-72-8	87
3			2-methylhexacosane	380	C27H56	1000376-72-7	87
4			Docosane, 11-decyl-	451	C32H66	055401-55-3	86
5			Hexacosane	366	C26H54	000630-01-3	86



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM060117\
Data File : BM010362.D
Acq On : 01 Jun 2017 19:12
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Sample : I3163-05
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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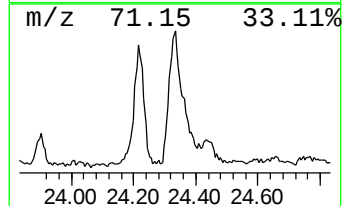
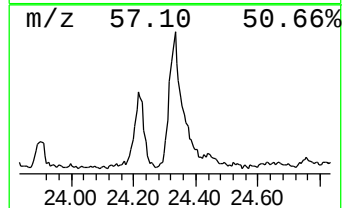
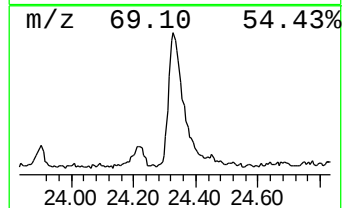
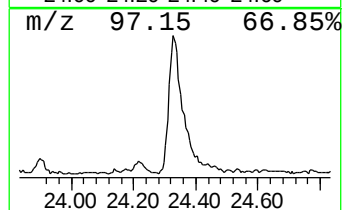
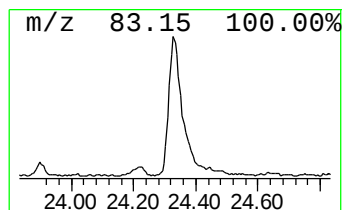
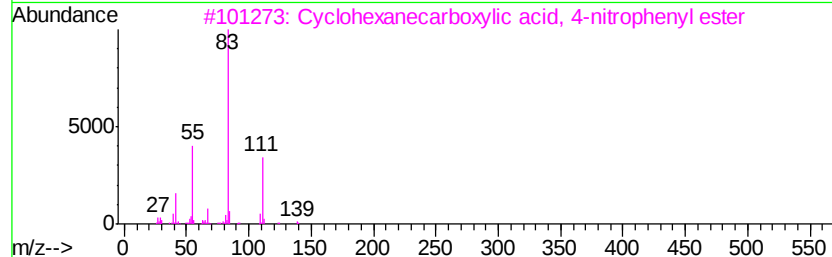
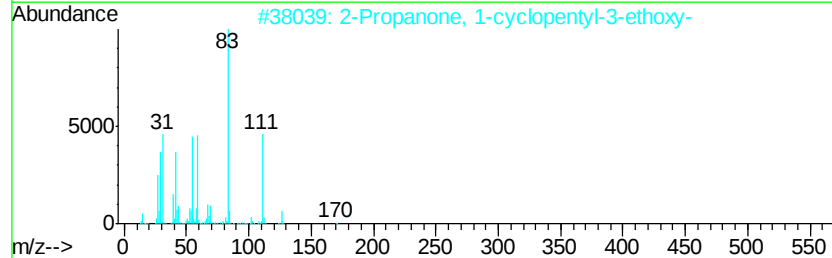
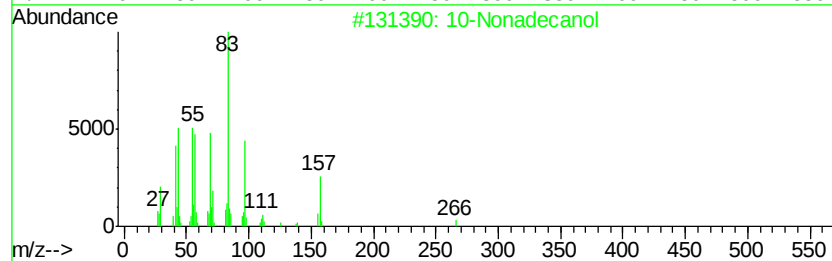
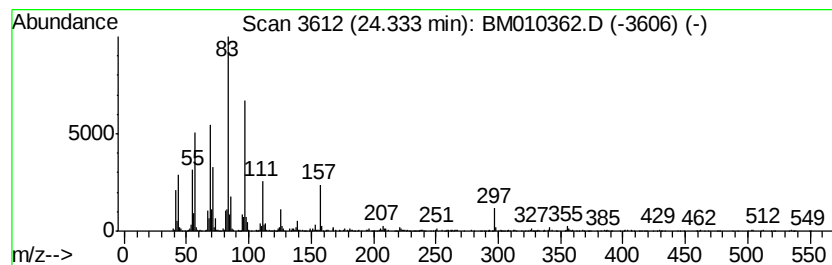
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 10-Nonadecanol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.33	12.91 ng/ul	4406760	Perylene-d12	23.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	10-Nonadecanol	284	C19H40O	016840-84-9	52
2		2-Propanone, 1-cyclopentyl-3-eth...	170	C10H18O2	051149-71-4	38
3		Cyclohexanecarboxylic acid, 4-ni...	249	C13H15NO4	1000307-70-8	38
4		Isocitronellol	156	C10H20O	018479-52-2	35
5		2-Hexene, 4,4,5-trimethyl-	126	C9H18	055702-61-9	30



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM060117\
Data File : BM010362.D
Acq On : 01 Jun 2017 19:12
Operator : SJ/MA
Sample : I3163-05
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM052717.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	18.14	4.6	ng/ul	1147320	4	17.30	4988180	20.0
unknown-01	18.46	4.1	ng/ul	1011550	4	17.30	4988180	20.0
Naphthalene, 6-et...	18.88	3.4	ng/ul	855027	4	17.30	4988180	20.0
unknown-02	19.08	2.8	ng/ul	686952	4	17.30	4988180	20.0
Anthracene, 9-but...	19.44	4.1	ng/ul	1364190	5	21.46	6638280	20.0
3-Methyl-5-(1,4,4...	20.39	3.6	ng/ul	1207180	5	21.46	6638280	20.0
Eicosanoic acid	20.49	3.4	ng/ul	1126060	5	21.46	6638280	20.0
Dehydroabietic acid	21.13	13.3	ng/ul	4413630	5	21.46	6638280	20.0
(DEL) Alkane: Cyc...	22.03	2.9	ng/ul	963731	5	21.46	6638280	20.0
Triacontane, 1-br...	22.97	2.4	ng/ul	827023	6	23.81	6826540	20.0
10-Nonadecanol	24.33	12.9	ng/ul	4406760	6	23.81	6826540	20.0