

Data Path : Z:\HPCHEM1\BNA M\DATA\BM030518\
 Data File : BM014470.D
 Acq On : 05 Mar 2018 13:23
 Operator : SJ/JU
 Sample : J1678-22
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BE5Z5

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-BM021418.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	4.663	282	285	288	rBV	38638	42221	1.48%	0.110%
2	6.728	630	636	653	rVB	347657	718779	25.20%	1.866%
3	6.863	653	659	670	rBV	343196	538898	18.89%	1.399%
4	7.057	686	692	700	rBV	475980	772771	27.09%	2.006%
5	7.510	763	769	777	rBV	340945	556933	19.53%	1.446%
6	7.798	813	818	824	rVB6	14206	29854	1.05%	0.078%
7	8.251	889	895	908	rBV	373972	750643	26.32%	1.949%
8	8.675	960	967	977	rBV	501665	879803	30.84%	2.284%
9	9.392	1082	1089	1106	rBV	446338	846885	29.69%	2.199%
10	9.945	1176	1183	1203	rBV	402632	1062191	37.24%	2.758%
11	10.286	1233	1241	1247	rBV	453780	804304	28.20%	2.088%
12	10.451	1263	1269	1283	rBV	240352	616445	21.61%	1.600%
13	12.980	1693	1699	1704	rBV5	26630	42451	1.49%	0.110%
14	13.145	1721	1727	1733	rBV9	12839	29955	1.05%	0.078%
15	13.469	1777	1782	1788	rBV4	78326	126768	4.44%	0.329%
16	13.598	1798	1804	1809	rBV	1158251	1653983	57.99%	4.294%
17	13.645	1809	1812	1824	rVB	216292	339476	11.90%	0.881%
18	13.863	1842	1849	1860	rBV	1230497	1983380	69.53%	5.149%
19	14.069	1880	1884	1889	rBV2	62090	86358	3.03%	0.224%
20	14.174	1895	1902	1909	rBV2	655770	1053502	36.93%	2.735%
21	14.239	1909	1913	1919	rVB5	22253	42519	1.49%	0.110%
22	14.492	1949	1956	1969	rBV2	174543	530696	18.61%	1.378%
23	14.621	1975	1978	1984	rVV8	20890	36748	1.29%	0.095%
24	15.033	2041	2048	2053	rVB2	75717	111754	3.92%	0.290%
25	15.180	2066	2073	2080	rBV	1557103	2277707	79.85%	5.913%
26	15.239	2081	2083	2088	rVB4	22772	37862	1.33%	0.098%
27	15.357	2097	2103	2111	rBV	389582	667672	23.41%	1.733%
28	15.898	2190	2195	2204	rVB	68094	118645	4.16%	0.308%
29	16.363	2270	2274	2276	rBV5	18651	32543	1.14%	0.084%
30	16.692	2325	2330	2335	rBV7	34672	62002	2.17%	0.161%
31	16.851	2349	2357	2364	rBV6	53862	98376	3.45%	0.255%
32	16.939	2364	2372	2376	rVV2	879042	1335146	46.81%	3.466%
33	16.980	2376	2379	2384	rVV	430974	628960	22.05%	1.633%
34	17.039	2384	2389	2400	rVB2	1636735	2470054	86.60%	6.413%

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35	17.333	2434	2439	2441	rBV6	39778	60767	2.13%	0.158%
36	17.368	2442	2445	2453	rVB3	46045	75563	2.65%	0.196%
37	17.662	2490	2495	2499	rBV8	15578	30214	1.06%	0.078%
38	17.727	2499	2506	2511	rBV9	57661	107715	3.78%	0.280%
39	17.821	2518	2522	2523	rBV	53031	68343	2.40%	0.177%
40	17.851	2523	2527	2536	rVB2	366130	637959	22.37%	1.656%
41	17.951	2541	2544	2547	rVB5	28305	39409	1.38%	0.102%
42	18.009	2549	2554	2558	rBV2	86751	150276	5.27%	0.390%
43	18.051	2559	2561	2568	rVB3	59701	76832	2.69%	0.199%
44	18.151	2571	2578	2579	rBV7	35929	65593	2.30%	0.170%
45	18.268	2594	2598	2600	rBV4	38034	52333	1.83%	0.136%
46	18.327	2604	2608	2614	rVV3	52574	101748	3.57%	0.264%
47	18.386	2615	2618	2625	rVB8	52718	103279	3.62%	0.268%
48	18.580	2644	2651	2653	rVB7	23722	39915	1.40%	0.104%
49	18.633	2655	2660	2662	rBV5	29158	47038	1.65%	0.122%
50	18.768	2677	2683	2686	rVV3	126034	187618	6.58%	0.487%
51	18.904	2704	2706	2710	rVB4	36472	35848	1.26%	0.093%
52	19.015	2715	2725	2731	rBV	829735	1204238	42.22%	3.126%
53	19.139	2742	2746	2751	rBV3	182882	256149	8.98%	0.665%
54	19.251	2762	2765	2768	rBV5	38033	58412	2.05%	0.152%
55	19.351	2777	2782	2786	rBV	1898504	2852367	100.00%	7.405%
56	19.380	2786	2787	2796	rVV	693880	623910	21.87%	1.620%
57	19.451	2798	2799	2806	rVB7	46091	70678	2.48%	0.183%
58	19.586	2816	2822	2827	rBV2	110320	176381	6.18%	0.458%
59	19.715	2839	2844	2845	rBV5	55027	90007	3.16%	0.234%
60	19.780	2849	2855	2863	rVV	763105	1077673	37.78%	2.798%
61	19.886	2866	2873	2878	rVB4	117849	334528	11.73%	0.868%
62	19.980	2887	2889	2893	rVB2	45562	53987	1.89%	0.140%
63	20.186	2921	2924	2928	rBV5	47962	60411	2.12%	0.157%
64	20.762	3019	3022	3026	rBV4	115898	136979	4.80%	0.356%
65	20.798	3026	3028	3032	rBV3	67363	77852	2.73%	0.202%
66	20.862	3035	3039	3046	rVB2	665118	827125	29.00%	2.147%
67	21.062	3071	3073	3079	rVB2	107010	121205	4.25%	0.315%
68	21.156	3083	3089	3093	rVV2	1025532	1722301	60.38%	4.471%
69	21.192	3093	3095	3100	rVB	328795	378797	13.28%	0.983%
70	22.650	3340	3343	3347	rBV	453726	573426	20.10%	1.489%
71	22.697	3347	3351	3363	rVB	344465	866759	30.39%	2.250%

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72	23.203	3432	3437	3442	rBV2	901973	1567971	54.97%	4.071%
73	23.339	3455	3460	3466	rBV2	494394	838237	29.39%	2.176%
74	23.803	3534	3539	3547	rVB	729349	1282128	44.95%	3.329%

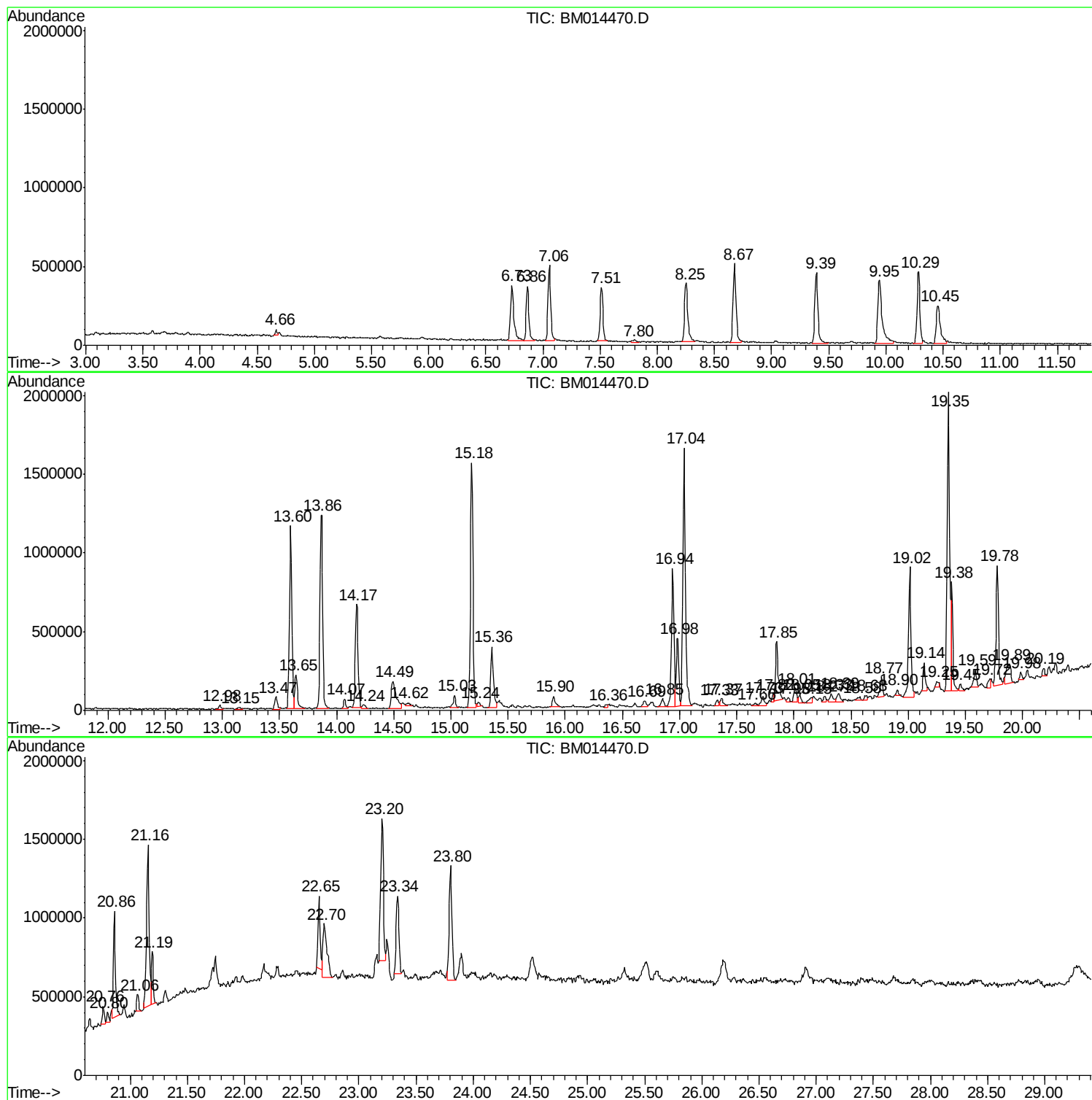
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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM021418.M
Quant Title : SVOA CALIBRATION

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



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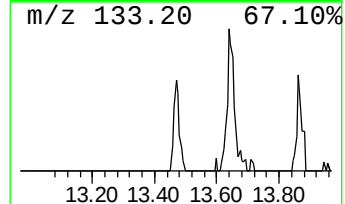
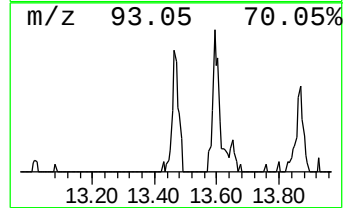
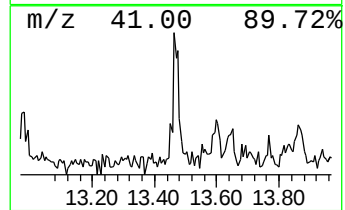
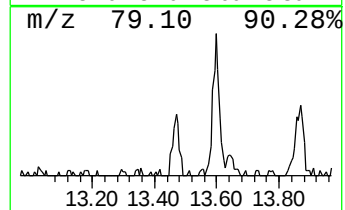
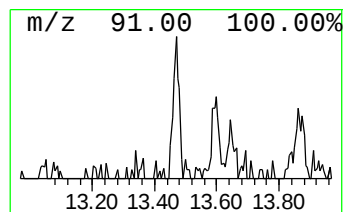
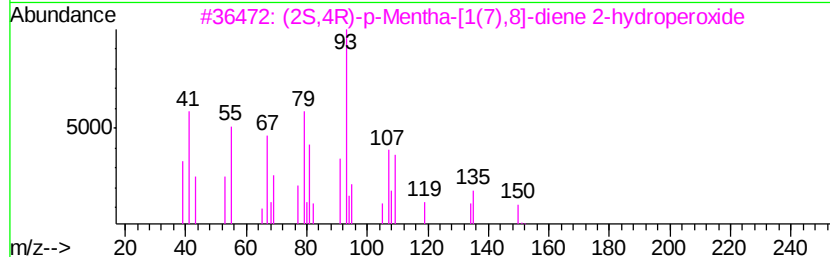
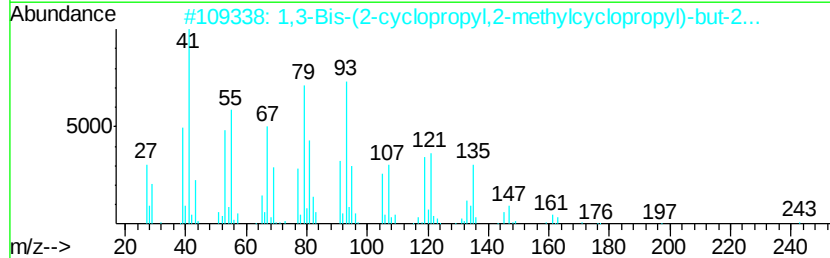
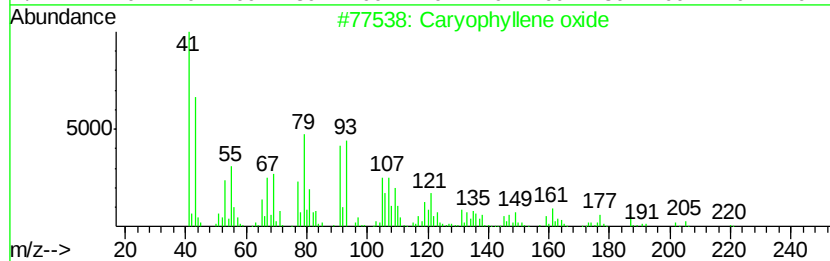
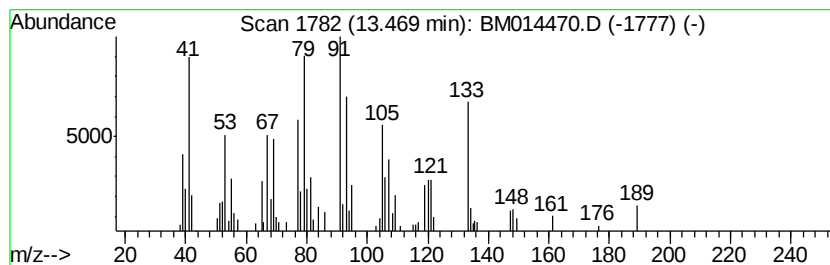
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Caryophyllene oxide Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.47	2.41 ng/ul	126768	Acenaphthene-d10	14.17
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Caryophyllene oxide	220 C15H24O	001139-30-6 59
2		1,3-Bis-(2-cyclopropyl,2-methylcyclopropyl)-but-2-ene	258 C18H26O	1000222-08-6 52
3		(2S,4R)-p-Mentha-[1(7),8]-diene	168 C10H16O2	1000292-74-4 50
4		1,3,7-Octatriene, 3,7-dimethyl-	136 C10H16	000502-99-8 38
5		4-Tridecen-6-yne, (E)-	178 C13H22	074744-43-7 27



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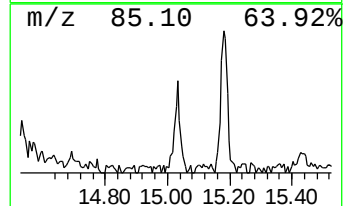
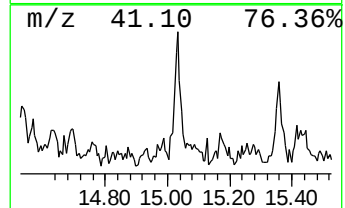
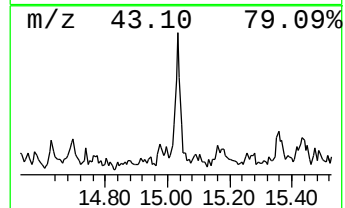
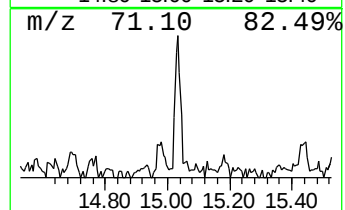
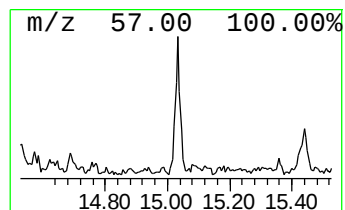
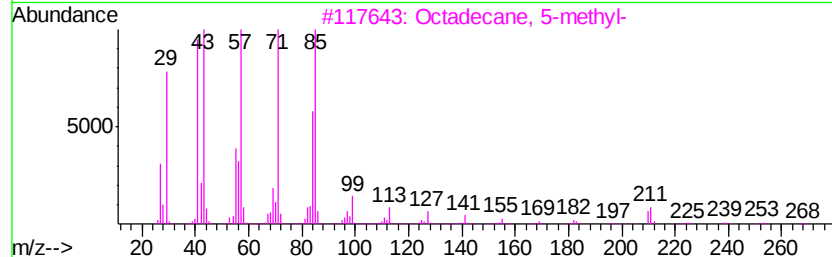
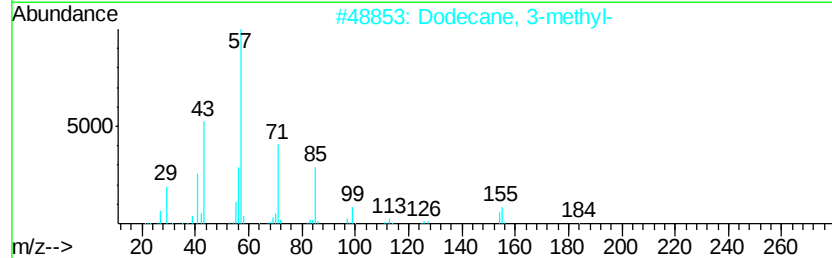
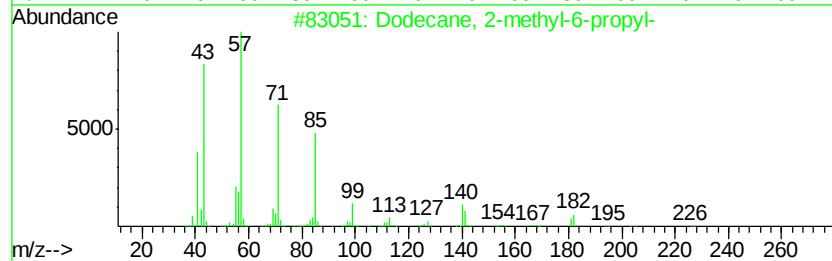
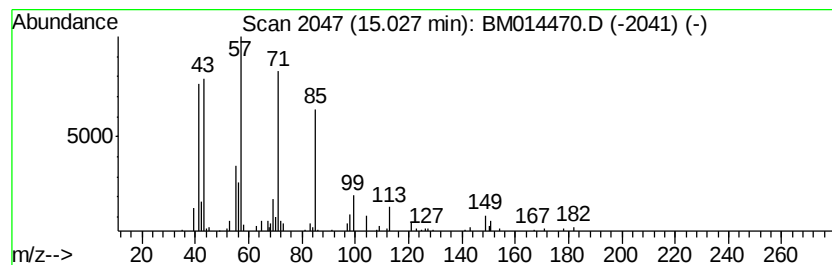
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.03	2.12 ng/ul	111754	Acenaphthene-d10	14.17

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	78
2			Dodecane, 3-methyl-	184	C13H28	017312-57-1	72
3			Octadecane, 5-methyl-	268	C19H40	025117-35-5	72
4			Dodecane, 5-methyl-	184	C13H28	017453-93-9	64
5			1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	59



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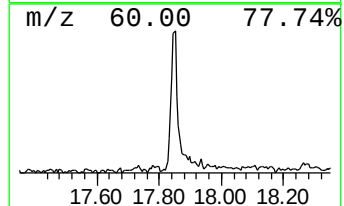
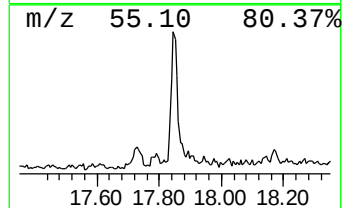
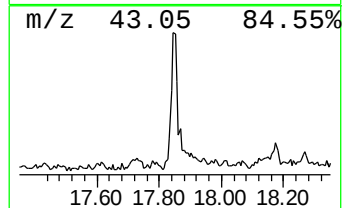
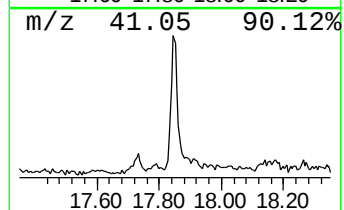
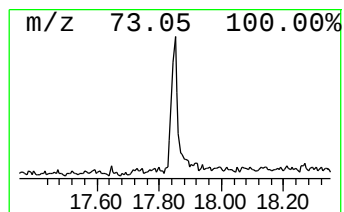
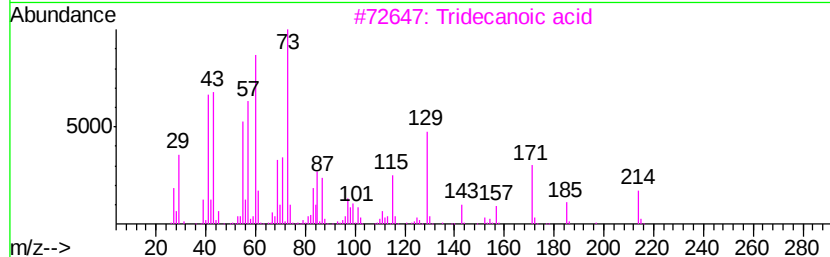
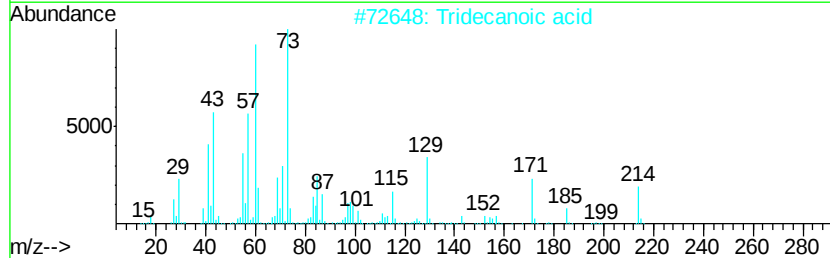
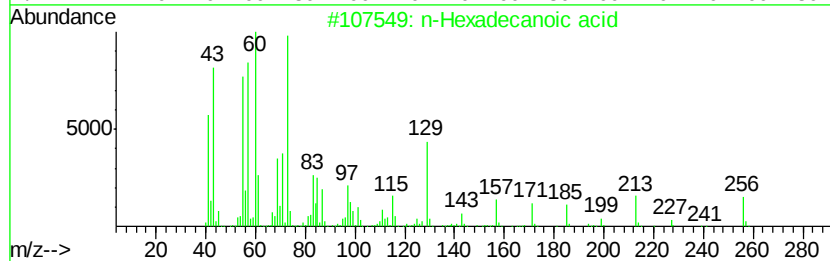
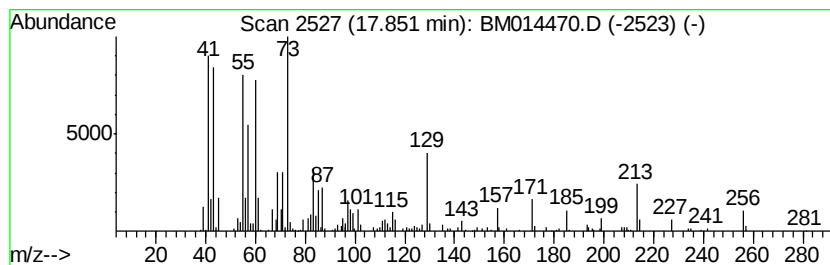
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Peak Number 3 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD		R.T.		
17.85	9.56 ng/ul	637959	Phenanthrene-d10		16.94		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2			Tridecanoic acid	214	C13H26O2	000638-53-9	93
3			Tridecanoic acid	214	C13H26O2	000638-53-9	93
4			Tridecanoic acid	214	C13H26O2	000638-53-9	90
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	74



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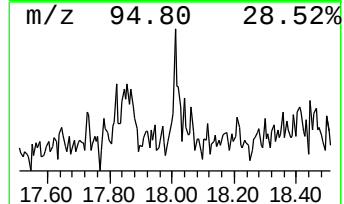
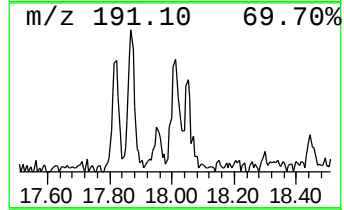
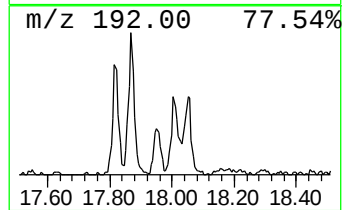
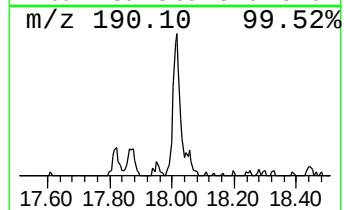
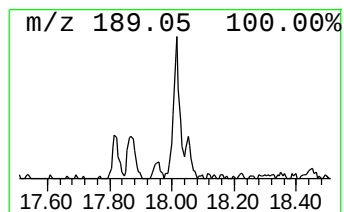
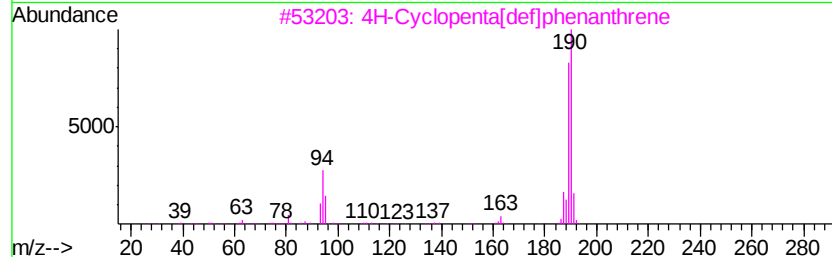
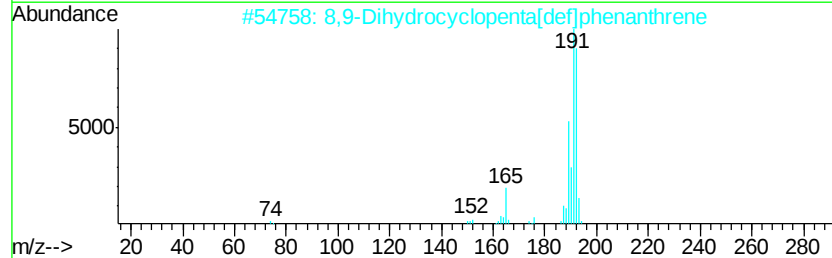
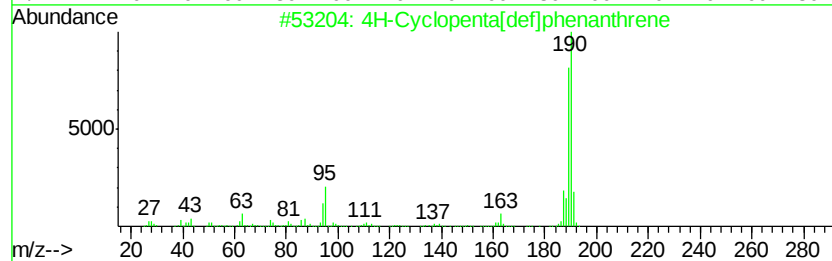
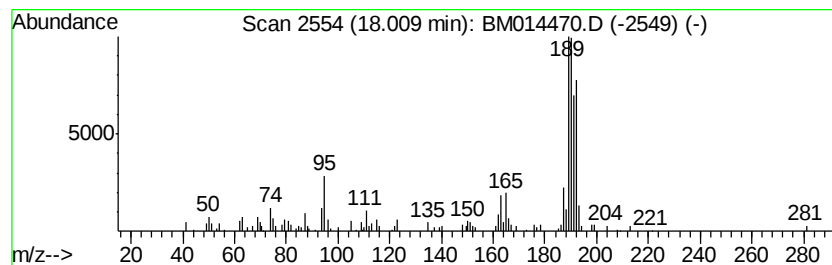
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 unknown-01 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.01	2.25 ng/ul	150276	Phenanthrene-d10	16.94

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	49
2			8,9-Dihydrocyclopenta[def]phenan...	192	C15H12	027410-55-5	46
3			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	46
4			Benzene, 1,1'-(2-cyclopropen-1-v...	192	C15H12	022825-21-4	41
5			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	38



Data Path : Z:\HPCHEM1\BNA M\DATA\BM030518\
Data File : BM014470.D
Acq On : 05 Mar 2018 13:23
Operator : SJ/JU
Sample : J1678-22
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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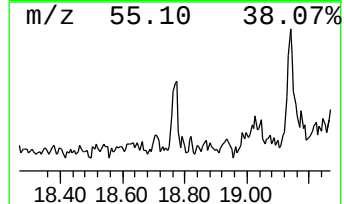
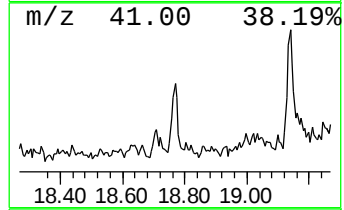
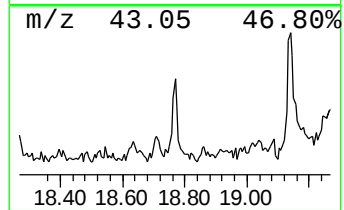
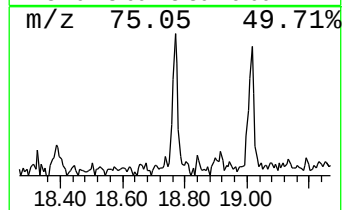
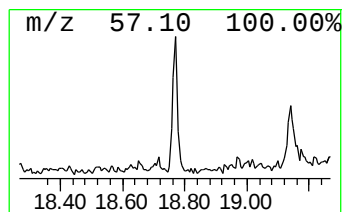
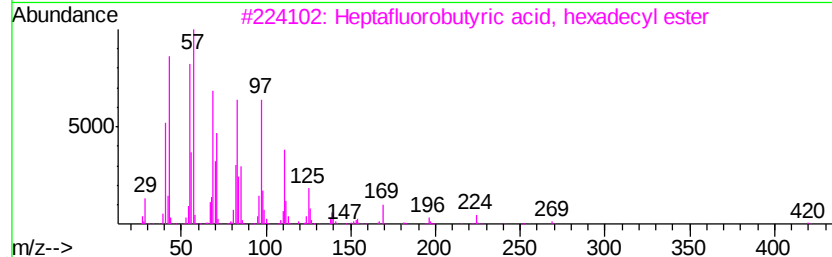
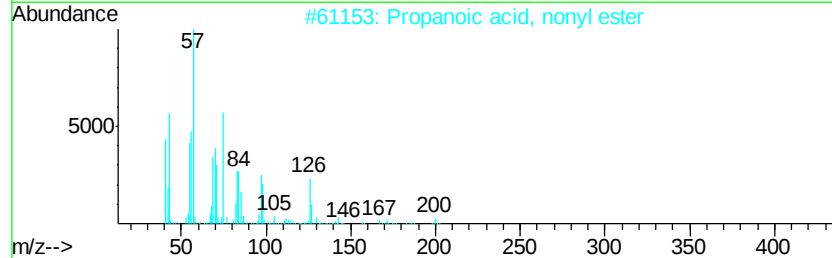
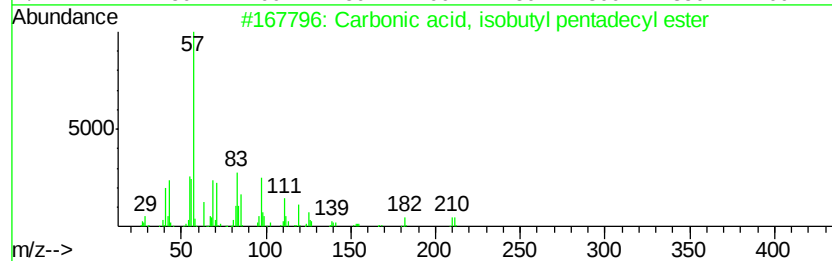
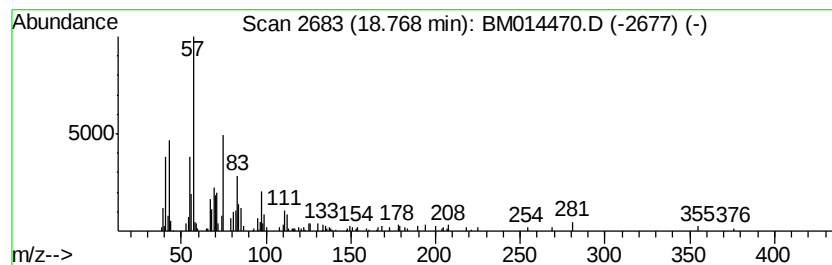
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Carbonic acid, isobutyl pen... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.77	2.81 ng/ul	187618	Phenanthrene-d10	16.94

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Carbonic acid, isobutyl pentadec...	328	C20H40O3	1000314-61-2	53
2		Propanoic acid, nonyl ester	200	C12H24O2	053184-67-1	53
3		Heptafluorobutyric acid, hexadec...	438	C20H33F7O2	006385-15-5	47
4		Heptafluorobutyric acid, n-octad...	466	C22H37F7O2	000400-57-7	47
5		Pentafluoropropionic acid, hepta...	402	C20H35F5O2	959218-78-1	47



Data Path : Z:\HPCHEM1\BNA M\DATA\BM030518\
Data File : BM014470.D
Acq On : 05 Mar 2018 13:23
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Sample : J1678-22
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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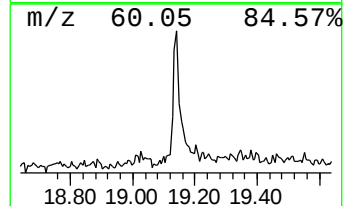
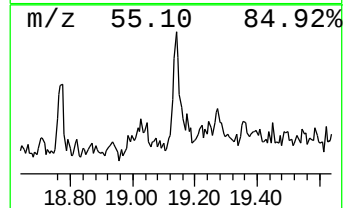
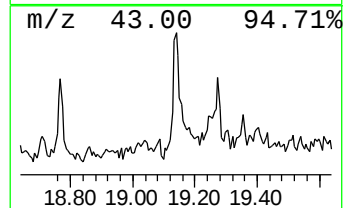
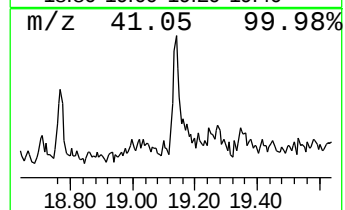
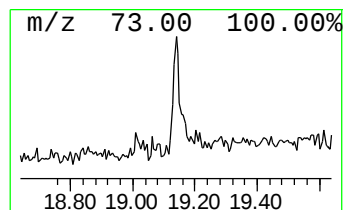
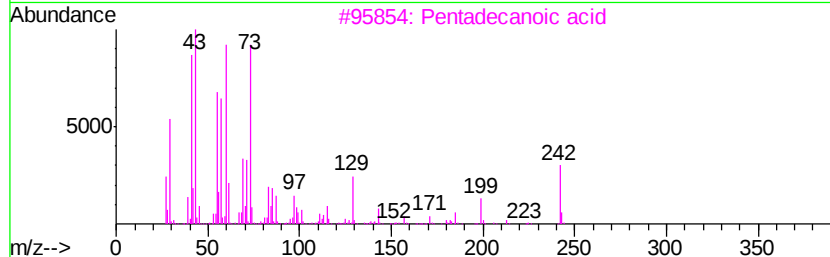
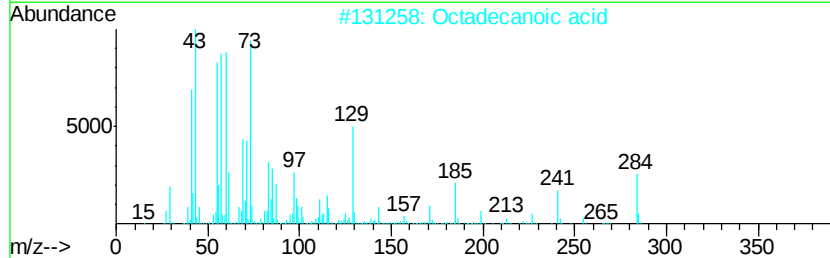
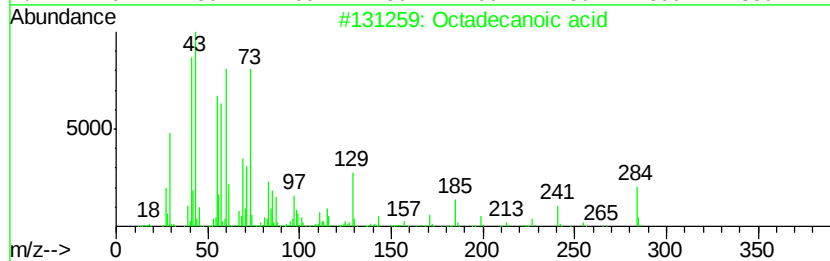
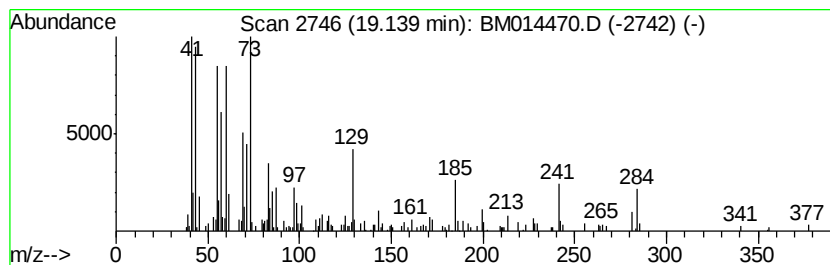
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Octadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.14	2.97 ng/ul	256149	Chrysene-d12	21.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	96
2		Octadecanoic acid	284	C18H36O2	000057-11-4	93
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	92
4		Octadecanoic acid	284	C18H36O2	000057-11-4	91
5		Octadecanoic acid	284	C18H36O2	000057-11-4	86



Data Path : Z:\HPCHEM1\BNA M\DATA\BM030518\
Data File : BM014470.D
Acq On : 05 Mar 2018 13:23
Operator : SJ/JU
Sample : J1678-22
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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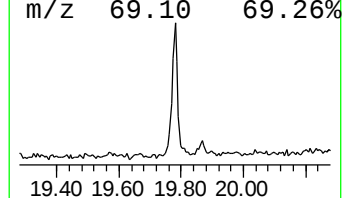
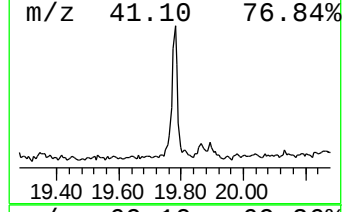
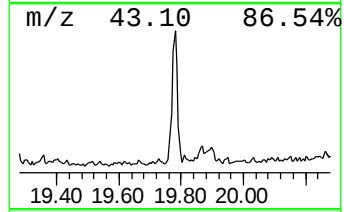
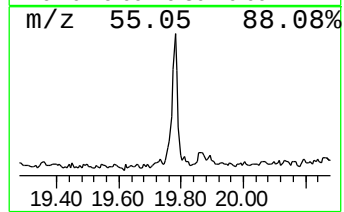
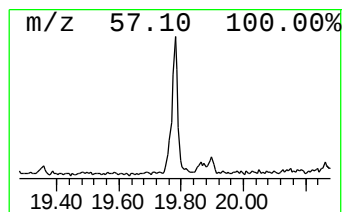
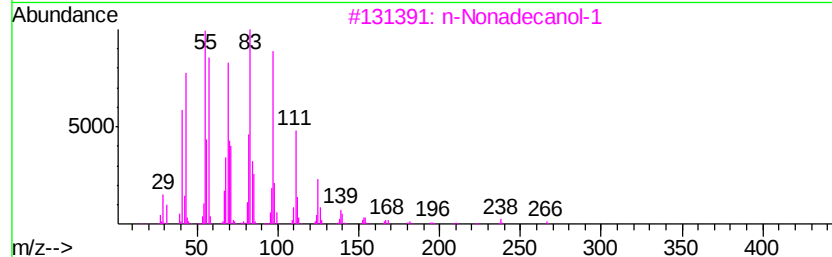
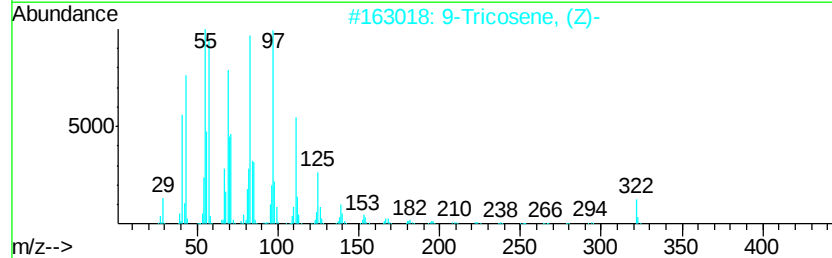
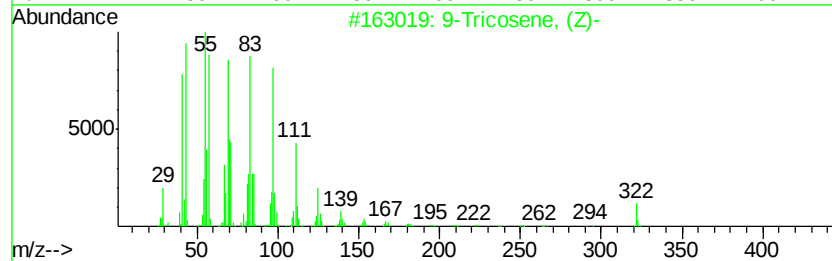
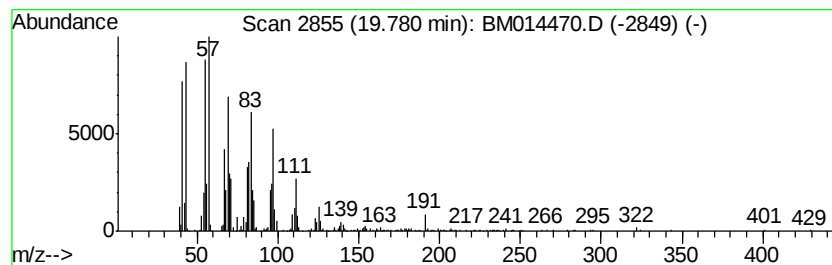
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Cyclic19.78 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.78	12.51 ng/ul	1077670	Chrysene-d12	21.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Tricosene, (Z)-	322	C23H46	027519-02-4	92
2		9-Tricosene, (Z)-	322	C23H46	027519-02-4	68
3		n-Nonadecanol-1	284	C19H40O	001454-84-8	64
4		Bromoacetic acid, octadecyl ester	390	C20H39BrO2	018992-03-5	64
5		1-Eicosanol	298	C20H42O	000629-96-9	64



Data Path : Z:\HPCHEM1\BNA M\DATA\BM030518\
Data File : BM014470.D
Acq On : 05 Mar 2018 13:23
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Sample : J1678-22
Misc :
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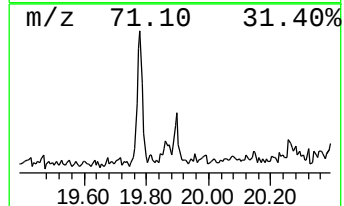
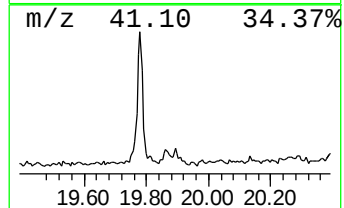
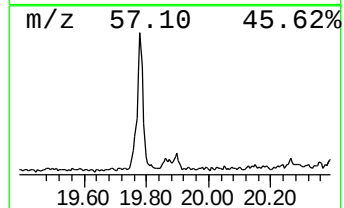
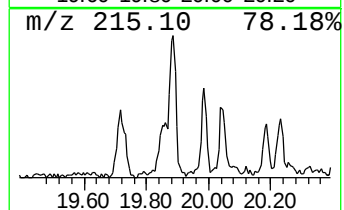
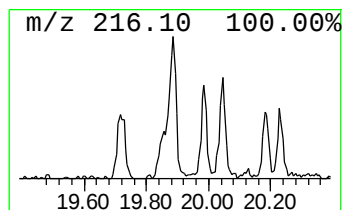
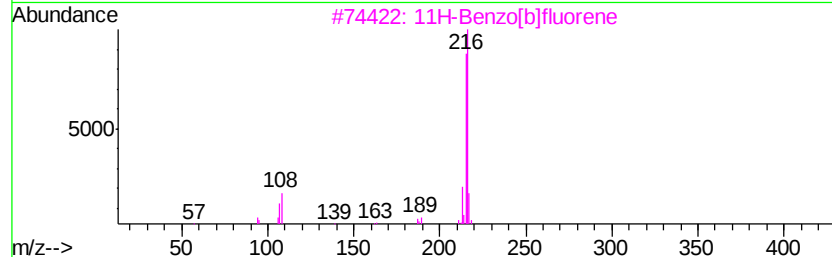
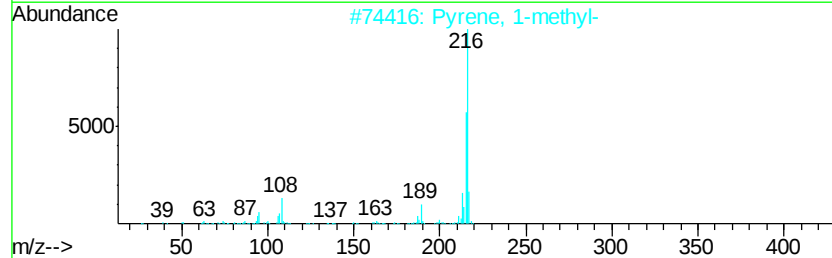
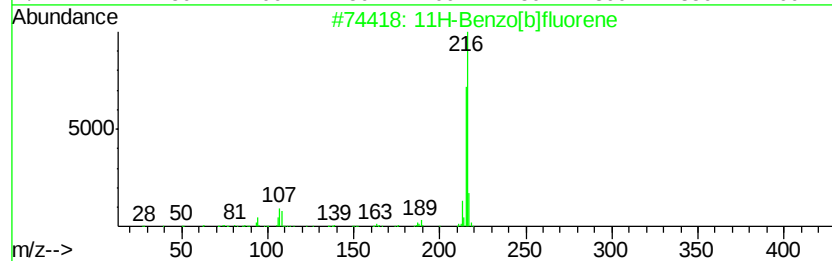
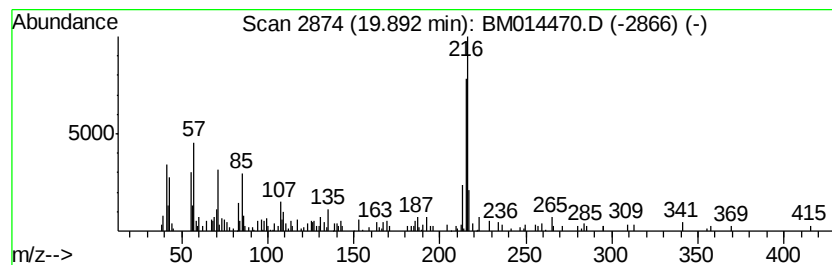
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 11H-Benzo[b]fluorene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.89	3.88 ng/ul	334528	Chrysene-d12	21.16

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	87
2		Pyrene, 1-methyl-	216	C17H12	002381-21-7	83
3		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	76
4		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	76
5		Pyrene, 1-methyl-	216	C17H12	002381-21-7	74



Data Path : Z:\HPCHEM1\BNA M\DATA\BM030518\
Data File : BM014470.D
Acq On : 05 Mar 2018 13:23
Operator : SJ/JU
Sample : J1678-22
Misc :
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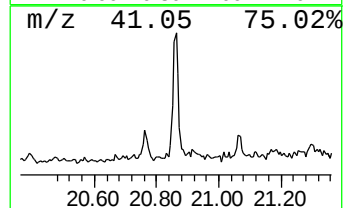
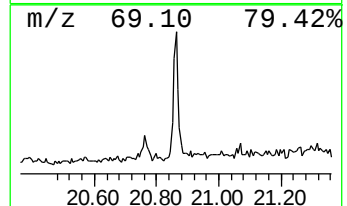
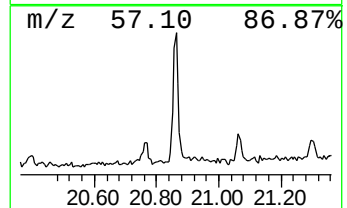
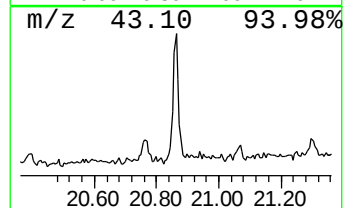
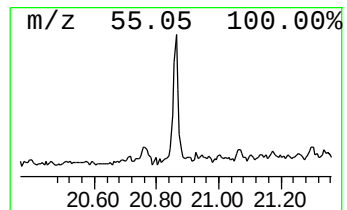
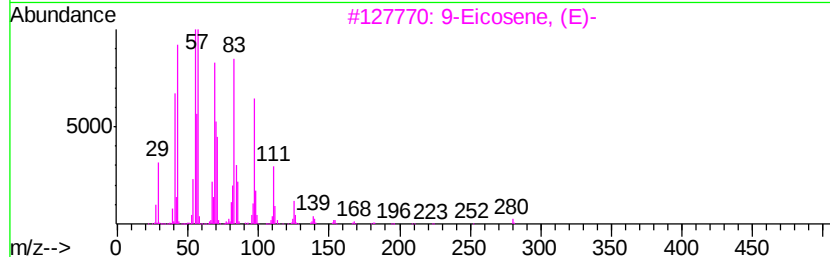
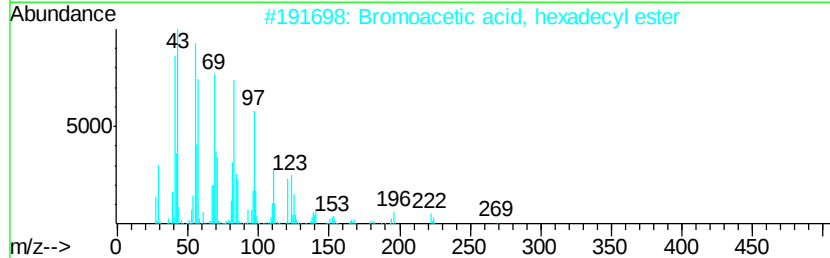
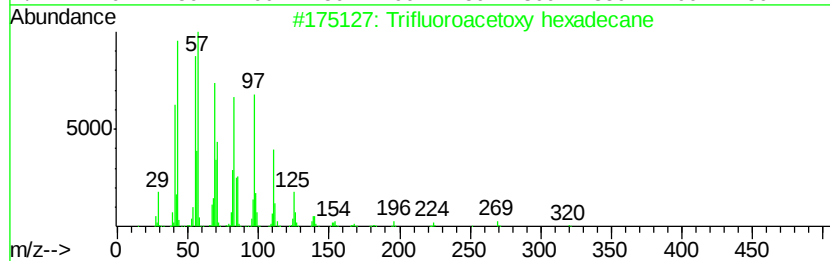
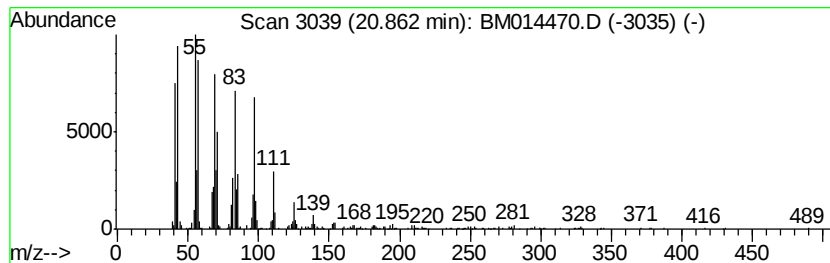
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Trifluoroacetoxy hexadecane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.86	9.60 ng/ul	827125	Chrysene-d12	21.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	91
2		Bromoacetic acid, hexadecyl ester	362	C18H35BrO2	005454-48-8	90
3		9-Eicosene, (E)-	280	C20H40	074685-29-3	87
4		n-Tetracosanol-1	354	C24H50O	000506-51-4	83
5		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	70



Data Path : Z:\HPCHEM1\BNA M\DATA\BM030518\
Data File : BM014470.D
Acq On : 05 Mar 2018 13:23
Operator : SJ/JU
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Misc :
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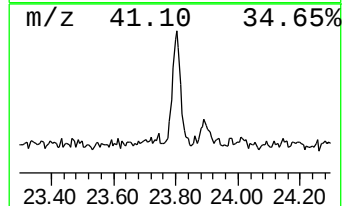
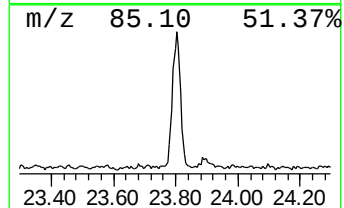
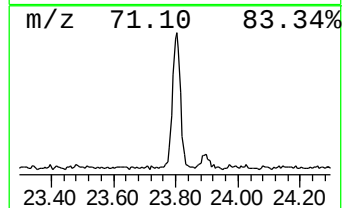
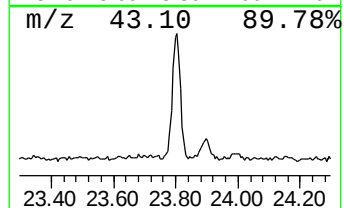
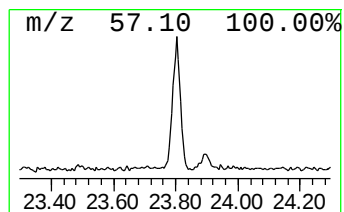
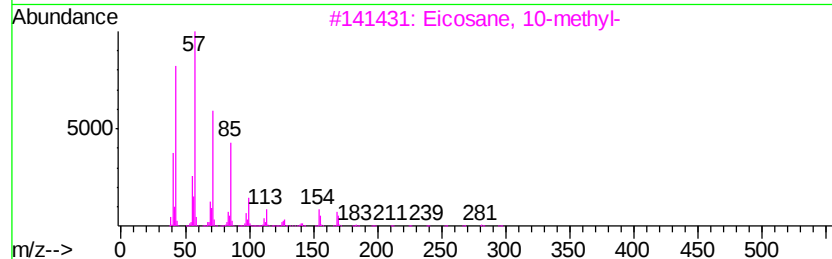
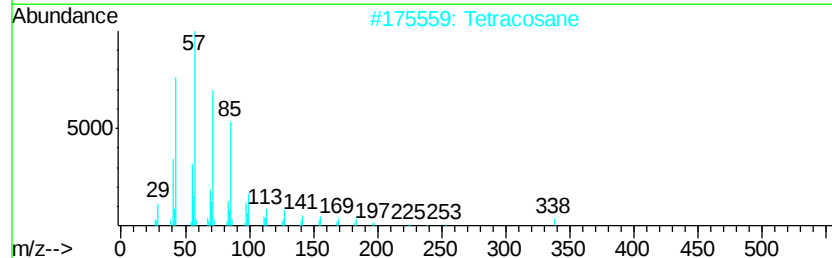
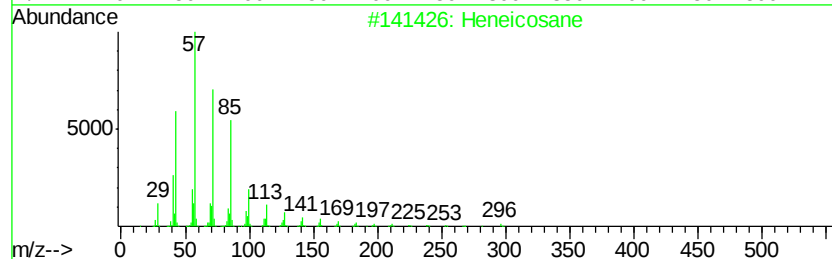
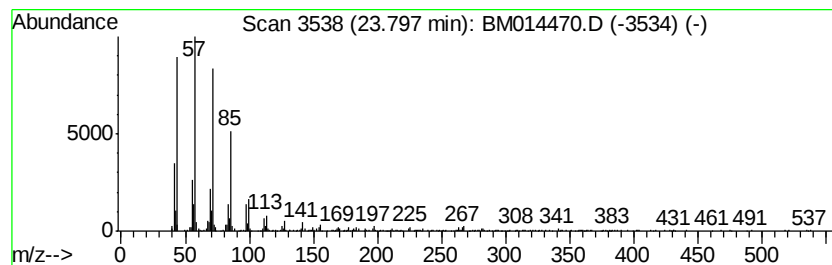
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.80	30.59 ng/ul	1282130	Perylene-d12	23.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	97
2		Tetracosane	338	C24H50	000646-31-1	94
3		Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
4		Heneicosane	296	C21H44	000629-94-7	91
5		Tetracosane, 11-decyl-	479	C34H70	055429-84-0	90



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM030518\
Data File : BM014470.D
Acq On : 05 Mar 2018 13:23
Operator : SJ/JU
Sample : J1678-22
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM021418.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
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(DEL) Alkane: Str...	15.03	2.1	ng/ul	111754	3	14.17	1053500	20.0
n-Hexadecanoic acid	17.85	9.6	ng/ul	637959	4	16.94	1335150	20.0
unknown-01	18.01	2.3	ng/ul	150276	4	16.94	1335150	20.0
Carbonic acid, is...	18.77	2.8	ng/ul	187618	4	16.94	1335150	20.0
Octadecanoic acid	19.14	3.0	ng/ul	256149	5	21.16	1722300	20.0
(DEL) Alkane: Cyc...	19.78	12.5	ng/ul	1077670	5	21.16	1722300	20.0
11H-Benzo[b]fluorene	19.89	3.9	ng/ul	334528	5	21.16	1722300	20.0
Trifluoroacetoxy ...	20.86	9.6	ng/ul	827125	5	21.16	1722300	20.0
(DEL) Alkane: Str...	23.80	30.6	ng/ul	1282130	6	23.34	838237	20.0