

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA M\DATA\BM101417\
 Data File : BM012186.D
 Acq On : 15 Oct 2017 00:41
 Operator : SJ/JU
 Sample : I5522-16
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AC1

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-BM101217.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	6.981	656	662	676	rBV	165188	338527	6.05%	0.520%
2	7.139	683	689	700	rBV	126290	214448	3.84%	0.330%
3	7.339	716	723	739	rBV	198835	368995	6.60%	0.567%
4	7.810	795	803	816	rBV	468650	805943	14.41%	1.239%
5	8.528	919	925	936	rBV	202921	398765	7.13%	0.613%
6	8.981	995	1002	1014	rBV	182344	344943	6.17%	0.530%
7	9.704	1118	1125	1134	rBV	165638	319606	5.72%	0.491%
8	10.245	1210	1217	1234	rBV	252932	546721	9.78%	0.840%
9	10.610	1272	1279	1294	rBV	716944	1221226	21.84%	1.877%
10	10.769	1299	1306	1319	rBV2	91593	231527	4.14%	0.356%
11	13.868	1825	1833	1837	rBV	671657	959812	17.17%	1.475%
12	13.916	1837	1841	1850	rVB	267645	399120	7.14%	0.613%
13	14.157	1876	1882	1897	rVB	736629	1167107	20.87%	1.794%
14	14.463	1927	1934	1941	rBV2	1244535	1862510	33.31%	2.863%
15	14.698	1966	1974	1986	rBV	79395	230826	4.13%	0.355%
16	15.457	2097	2103	2109	rBV	995564	1495729	26.75%	2.299%
17	15.598	2120	2127	2136	rBV	217015	368659	6.59%	0.567%
18	16.151	2217	2221	2224	rBV2	37519	58409	1.04%	0.090%
19	16.945	2349	2356	2362	rBV7	39044	86375	1.54%	0.133%
20	17.092	2376	2381	2387	rBV	143962	214890	3.84%	0.330%
21	17.156	2387	2392	2395	rBV2	83821	109688	1.96%	0.169%
22	17.209	2395	2401	2405	rBV2	1776064	2556378	45.72%	3.929%
23	17.251	2405	2408	2412	rVB	515269	638806	11.43%	0.982%
24	17.309	2412	2418	2422	rBV2	1129488	1620102	28.98%	2.490%
25	17.968	2524	2530	2535	rBV	139970	216077	3.86%	0.332%
26	18.027	2536	2540	2545	rVV2	92770	134104	2.40%	0.206%
27	18.086	2545	2550	2555	rVV2	971454	1321917	23.64%	2.032%
28	18.127	2555	2557	2567	rVB2	170520	277594	4.96%	0.427%
29	18.215	2568	2572	2577	rVB	50341	68110	1.22%	0.105%
30	18.280	2577	2583	2587	rBV2	165507	312286	5.59%	0.480%
31	18.315	2587	2589	2595	rVB	79639	97043	1.74%	0.149%
32	18.386	2595	2601	2606	rBV2	224772	371887	6.65%	0.572%
33	18.580	2630	2634	2639	rVV	100160	153148	2.74%	0.235%
34	18.633	2639	2643	2649	rVV2	84179	141480	2.53%	0.217%

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Integration Parameters: LSCINT.P

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35	18.692	2650	2653	2659	rVB7	45952	81192	1.45%	0.125%
36	19.003	2701	2706	2710	rBV2	121240	203907	3.65%	0.313%
37	19.150	2725	2731	2737	rBV3	85504	188841	3.38%	0.290%
38	19.268	2739	2751	2758	rBV	2093822	3226247	57.70%	4.959%
39	19.362	2763	2767	2780	rVV	741789	1063972	19.03%	1.635%
40	19.603	2802	2808	2810	rBV2	1922323	2958805	52.92%	4.548%
41	19.627	2810	2812	2820	rVB	1750741	2206270	39.46%	3.391%
42	19.698	2821	2824	2831	rVB2	92933	138267	2.47%	0.213%
43	19.809	2840	2843	2847	rVB	99077	121971	2.18%	0.187%
44	19.950	2862	2867	2869	rBV2	137771	239656	4.29%	0.368%
45	20.092	2888	2891	2893	rBV4	129823	182820	3.27%	0.281%
46	20.121	2893	2896	2902	rVB	318882	434695	7.77%	0.668%
47	20.221	2910	2913	2917	rBV	155210	203383	3.64%	0.313%
48	20.421	2945	2947	2950	rBV	80507	85416	1.53%	0.131%
49	20.874	3021	3024	3029	rVB	174226	211280	3.78%	0.325%
50	21.033	3048	3051	3054	rBV2	169837	236099	4.22%	0.363%
51	21.074	3054	3058	3062	rVV	877063	966137	17.28%	1.485%
52	21.280	3090	3093	3098	rVB2	350878	379068	6.78%	0.583%
53	21.386	3105	3111	3115	rBV2	3261081	5591276	100.00%	8.594%
54	21.427	3115	3118	3127	rVB	1191947	1536605	27.48%	2.362%
55	21.950	3204	3207	3209	rBV3	616463	763462	13.65%	1.173%
56	21.974	3209	3211	3221	rVB2	906516	1250983	22.37%	1.923%
57	22.662	3325	3328	3334	rVB2	397276	539141	9.64%	0.829%
58	22.944	3371	3376	3380	rBV	1767353	2484763	44.44%	3.819%
59	22.997	3380	3385	3397	rVB	1815669	4204705	75.20%	6.463%
60	23.497	3466	3470	3472	rBV	458710	749748	13.41%	1.152%
61	23.597	3483	3487	3494	rVB	887908	1576804	28.20%	2.424%
62	23.697	3498	3504	3510	rBV	1542929	2907725	52.00%	4.469%
63	23.856	3526	3531	3540	rVB2	884841	1624592	29.06%	2.497%
64	24.197	3583	3589	3597	rVB	2516120	5032289	90.00%	7.735%
65	25.438	3795	3800	3808	rVB2	769600	1689504	30.22%	2.597%
66	25.874	3868	3874	3883	rVB	893355	2327276	41.62%	3.577%

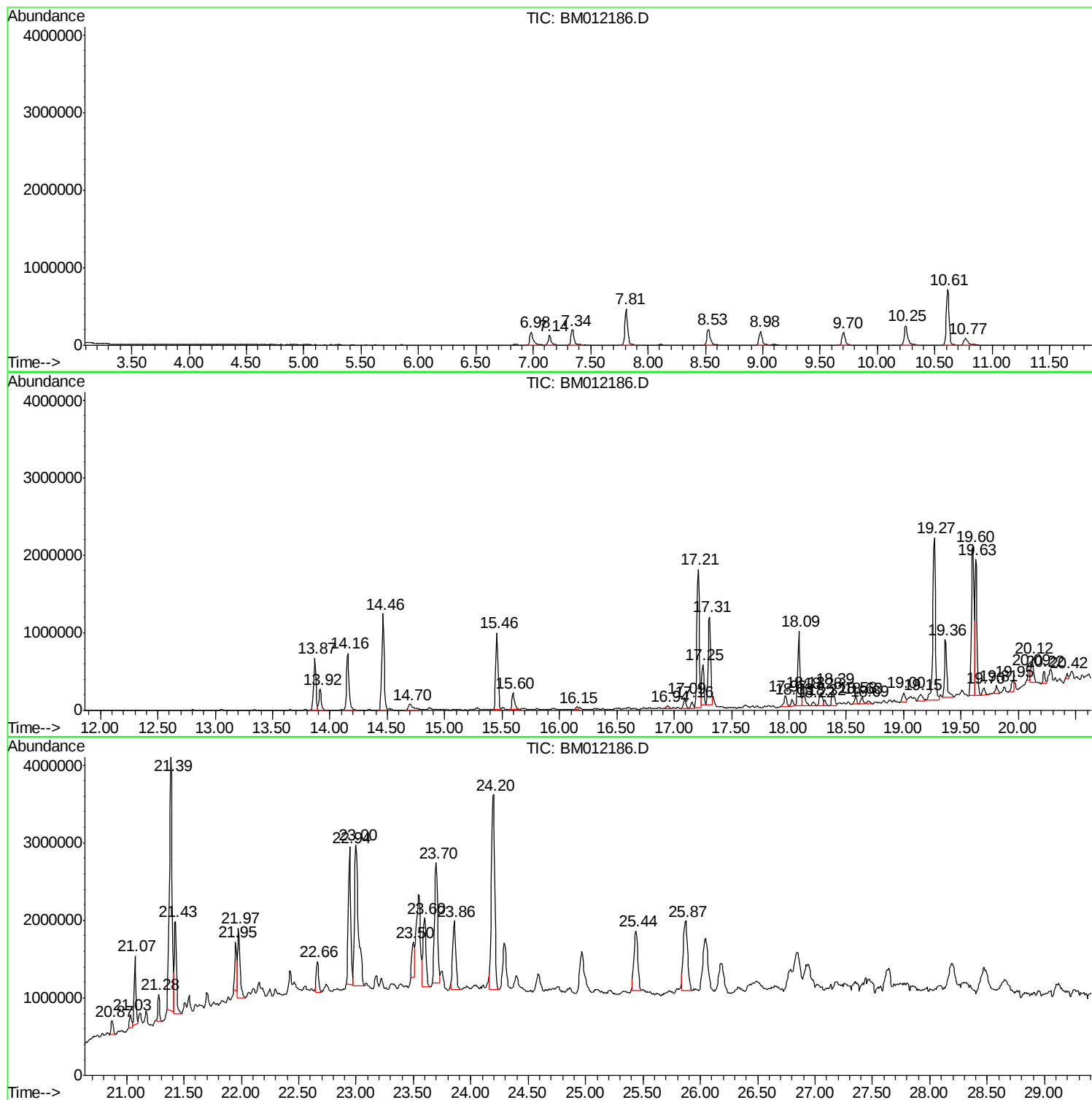
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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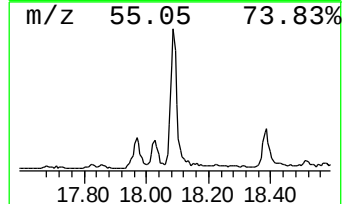
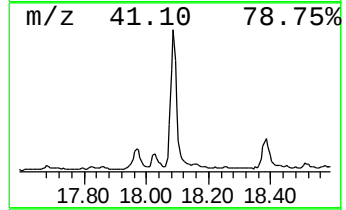
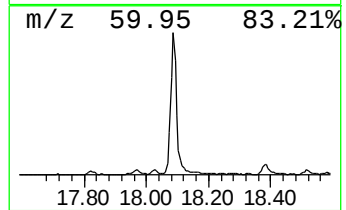
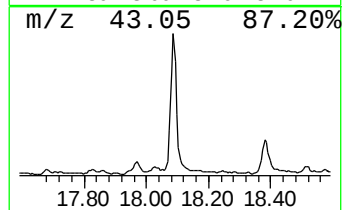
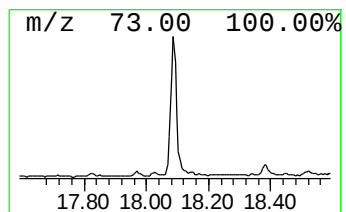
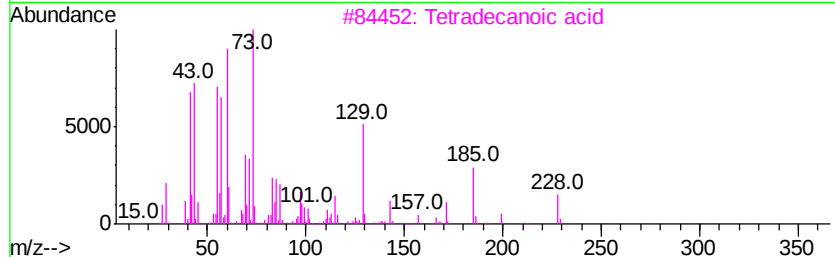
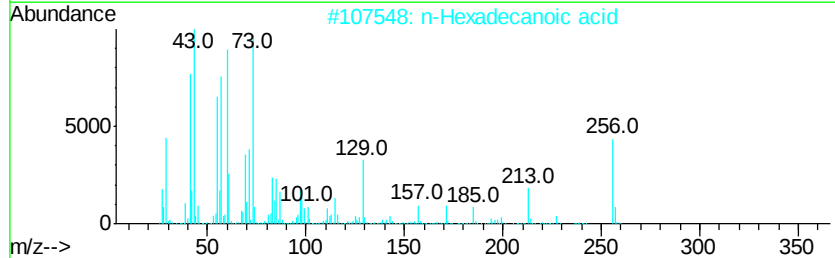
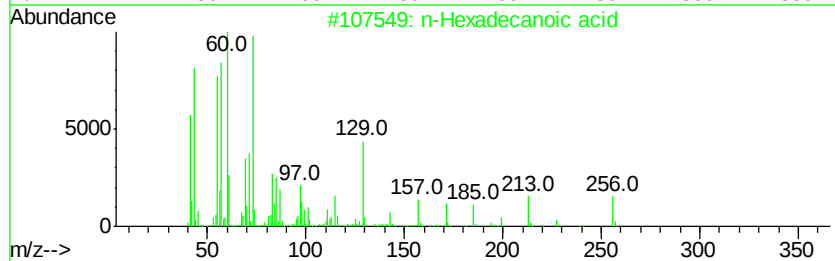
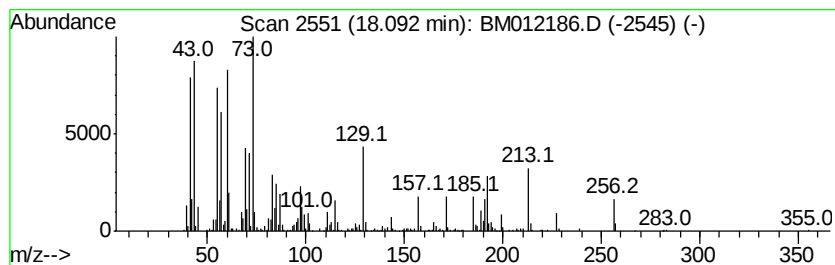
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM101217.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.09	10.34 ng/ul	1321920	Phenanthrene-d10	17.21

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		Tetradecanoic acid	228	C14H28O2	000544-63-8	93
4		Tridecanoic acid	214	C13H26O2	000638-53-9	86
5		Tridecanoic acid	214	C13H26O2	000638-53-9	86



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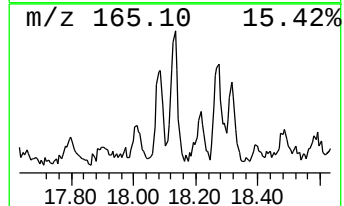
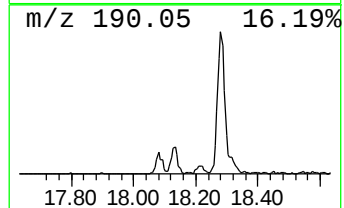
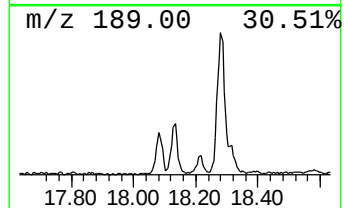
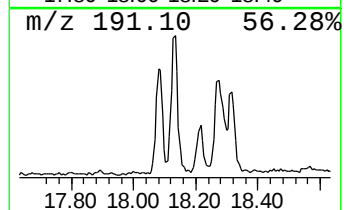
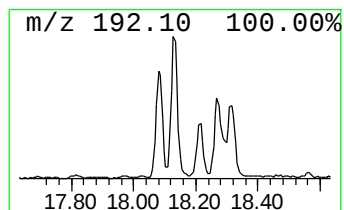
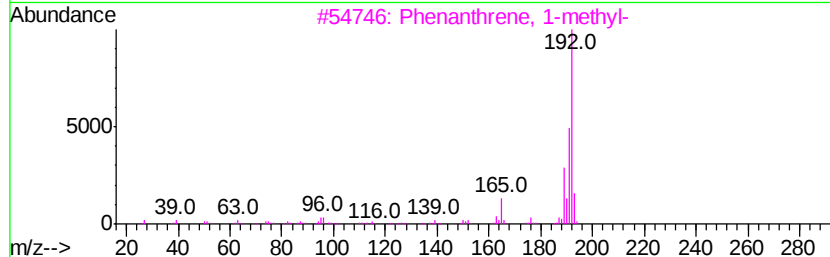
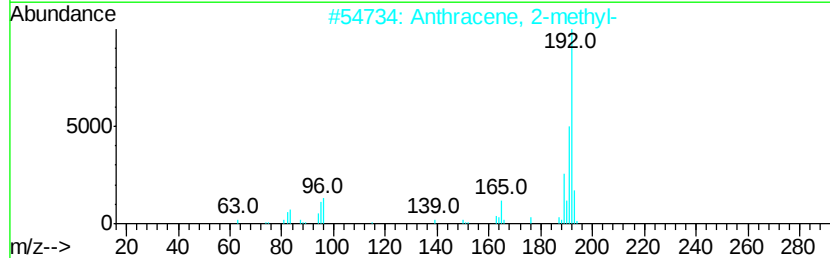
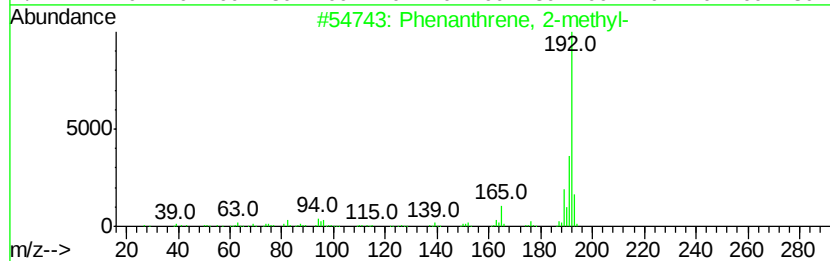
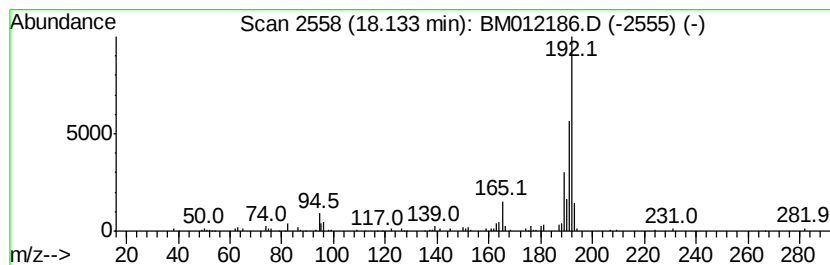
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM101217.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Phenanthrene, 2-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.13	2.17 ng/ul	277594	Phenanthrene-d10	17.21

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
2		Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
3		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	94
4		Anthracene, 1-methyl-	192	C15H12	000610-48-0	94
5		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93



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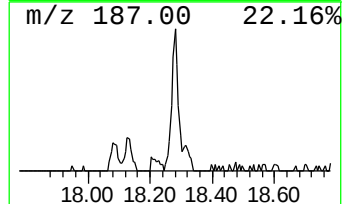
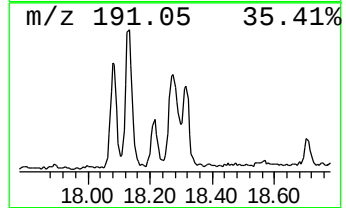
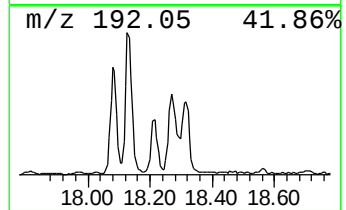
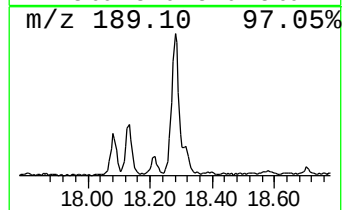
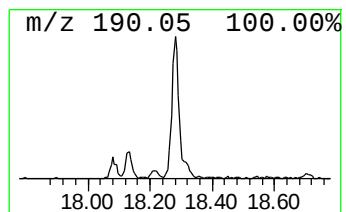
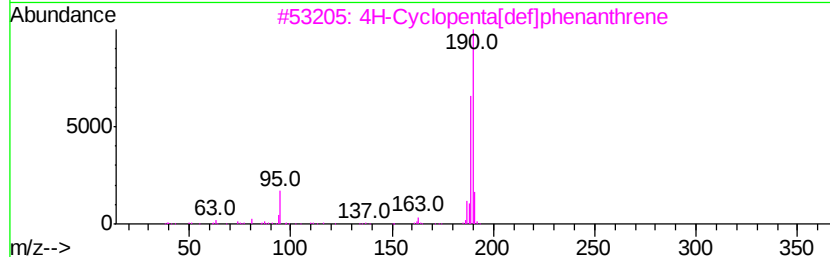
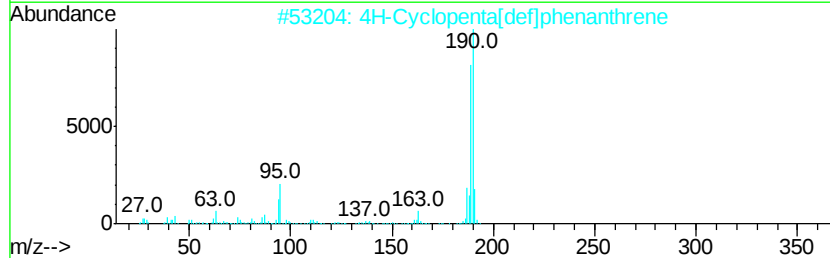
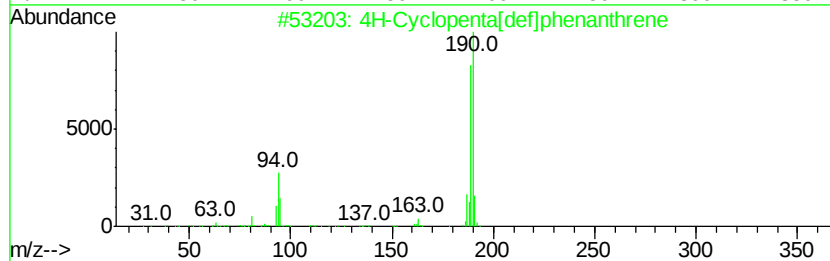
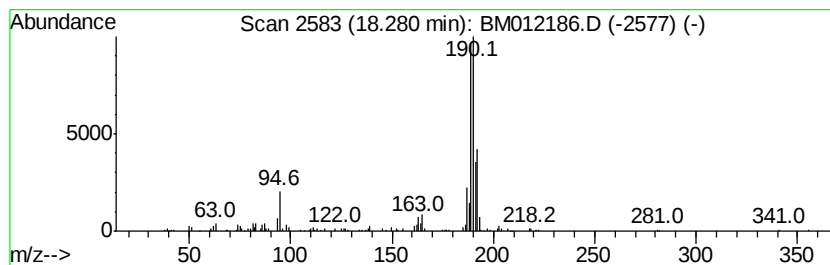
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 4H-Cyclopenta[def]phenanthrene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.28	2.44 ng/ul	312286	Phenanthrene-d10	17.21

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	74
2		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	70
3		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	58
4		6H-Cyclobuta[flk]phenanthrene	190	C15H10	083469-43-6	42
5		Carbonic acid, ethyl 2-formyl-4,...	262	C10H8Cl2O4	1000331-34-4	40



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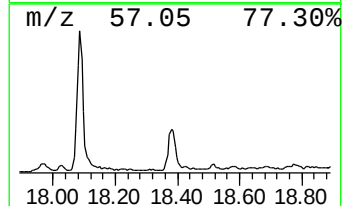
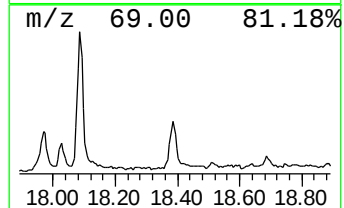
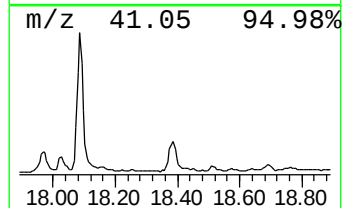
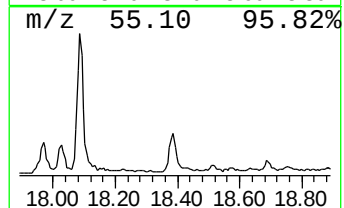
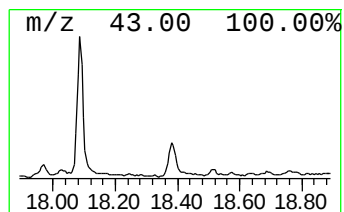
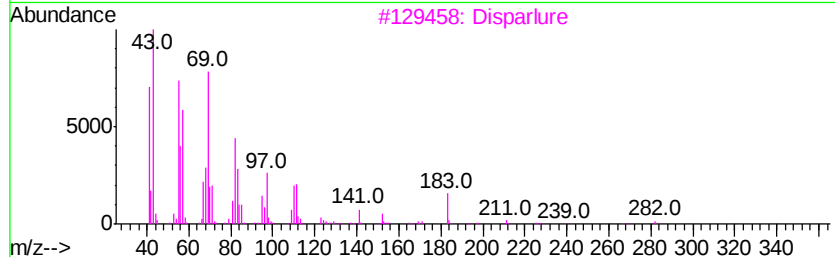
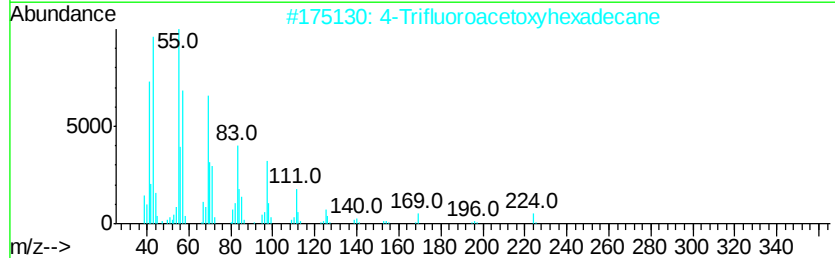
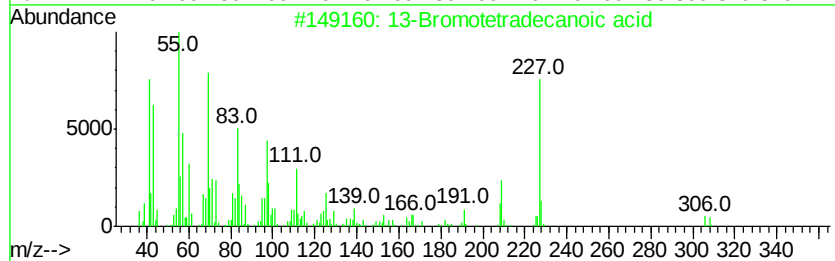
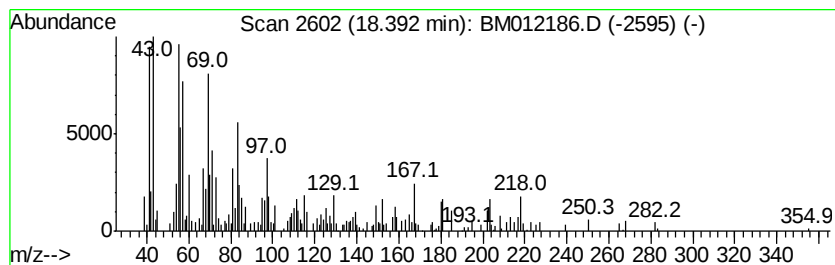
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Peak Number 4 13-Bromotetradecanoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.39	2.91 ng/ul	371887	Phenanthrene-d10	17.21

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		13-Bromotetradecanoic acid	306	C14H27BrO2	1000099-12-8	64
2		4-Trifluoroacetoxyhexadecane	338	C18H33F3O2	1000215-97-5	62
3		Disparlure	282	C19H38O	029804-22-6	60
4		Cyclododecane	168	C12H24	000294-62-2	58
5		Cyclopentane, 1-butyl-2-propyl-	168	C12H24	062199-50-2	55



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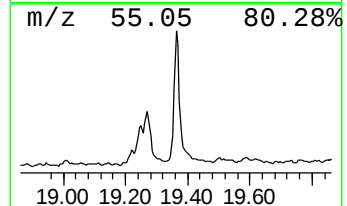
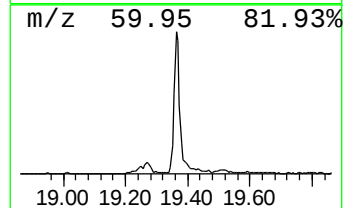
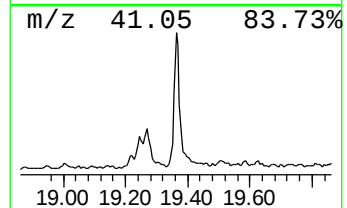
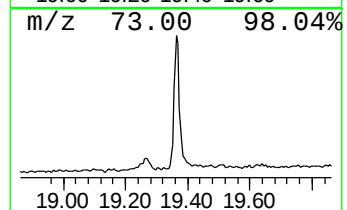
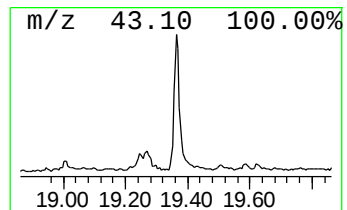
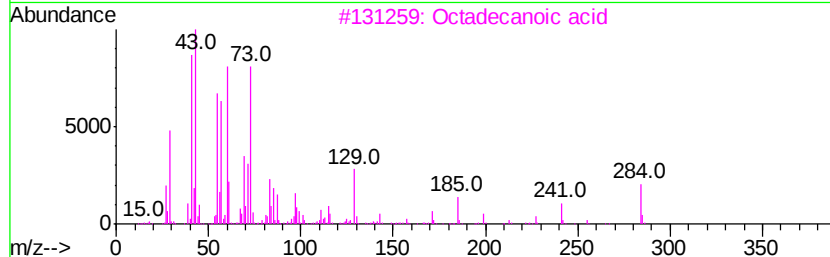
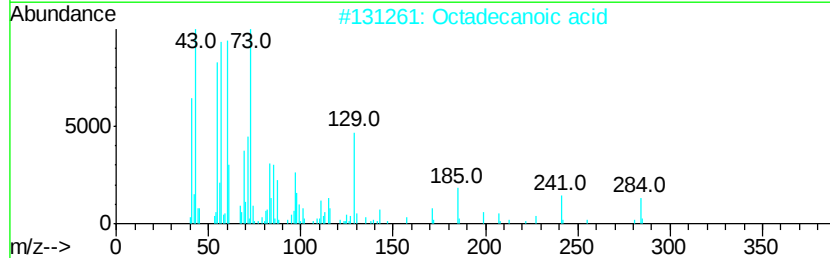
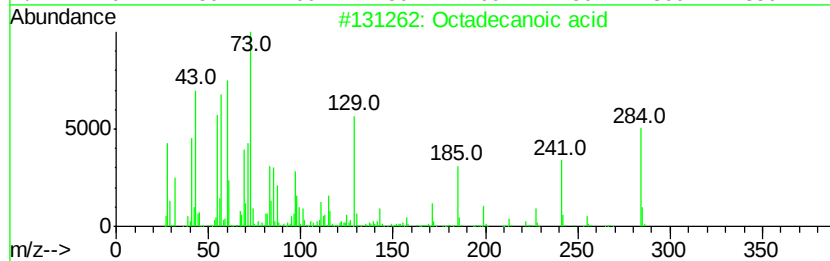
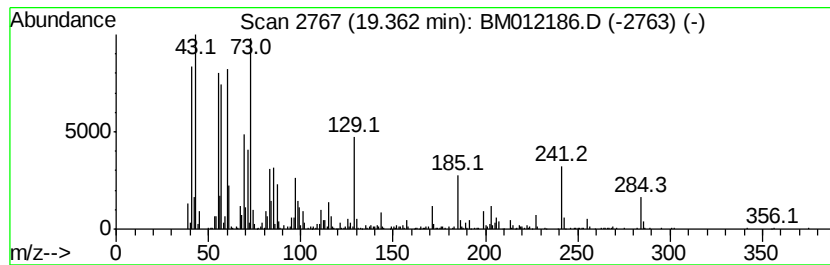
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Octadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.36	3.81 ng/ul	1063970	Chrysene-d12	21.39

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	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	98
4	Octadecanoic acid		284	C18H36O2	000057-11-4	98
5	Octadecanoic acid		284	C18H36O2	000057-11-4	95



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Data File : BM012186.D
Acq On : 15 Oct 2017 00:41
Operator : SJ/JU
Sample : I5522-16
Misc :
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ClientSampleId :
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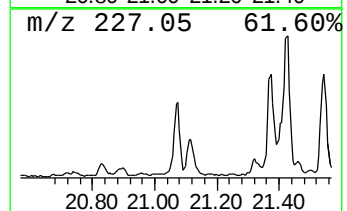
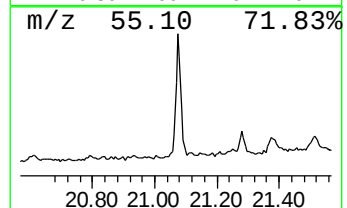
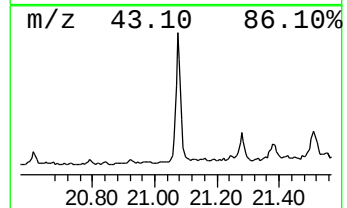
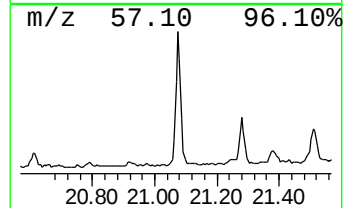
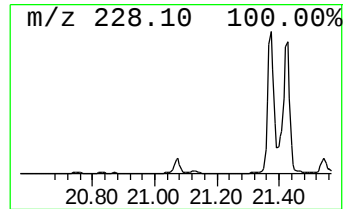
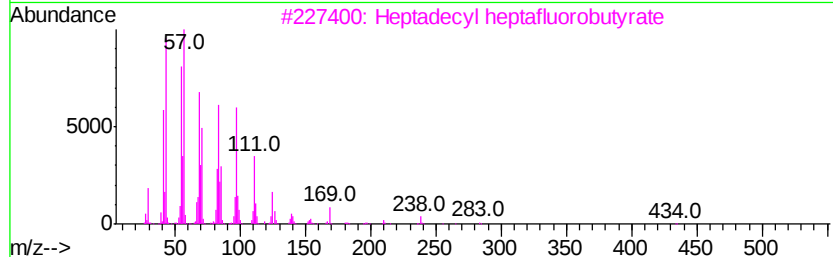
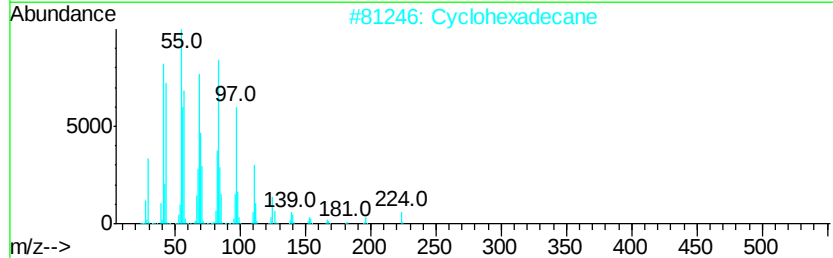
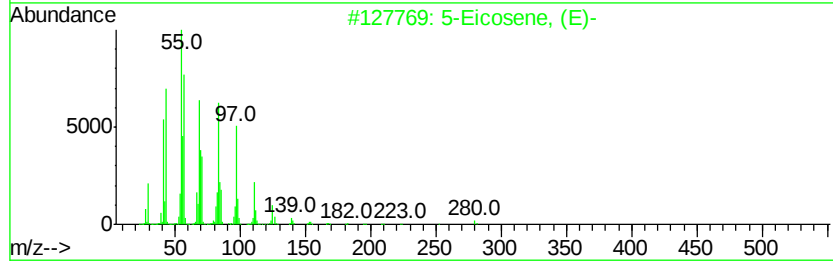
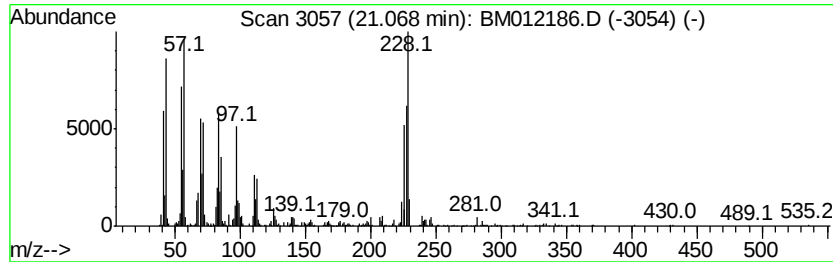
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Cyclic21.07 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.07	3.46 ng/ul	966137	Chrysene-d12	21.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Eicosene, (E)-	280	C20H40	074685-30-6	89
2		Cyclohexadecane	224	C16H32	000295-65-8	83
3		Heptadecyl heptafluorobutvrate	452	C21H35F7O2	959085-66-6	64
4		Cyclooctacosane	392	C28H56	000297-24-5	55
5		Eicosyl trifluoroacetate	394	C22H41F3O2	1000351-75-2	55



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Instrument :
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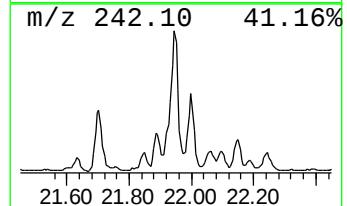
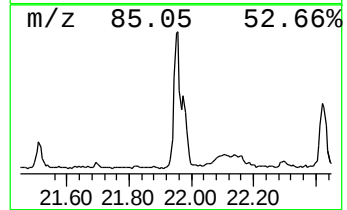
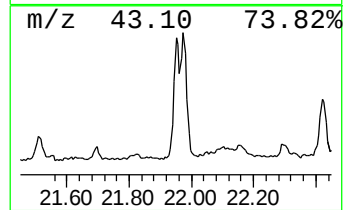
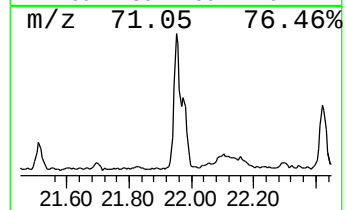
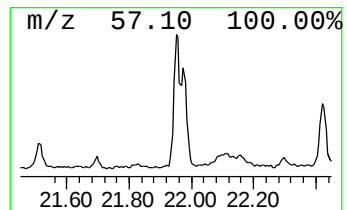
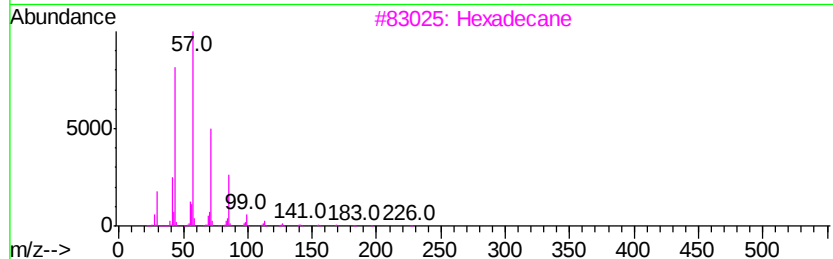
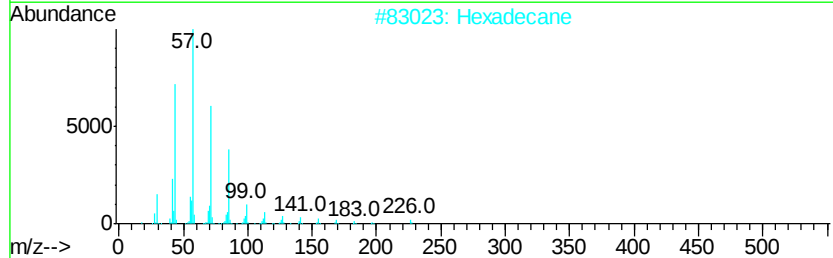
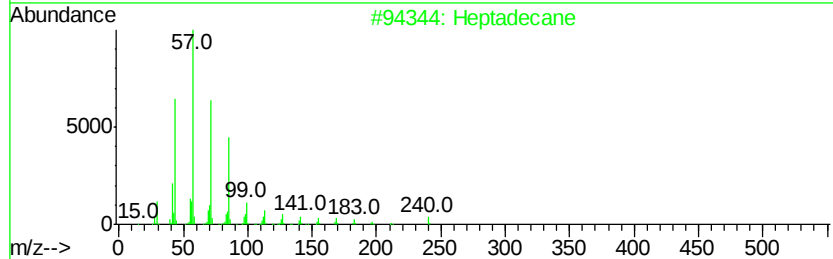
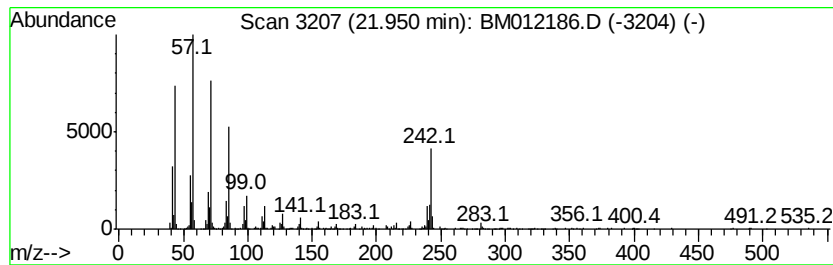
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.95	2.73 ng/ul	763462	Chrysene-d12	21.39

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane	240	C17H36	000629-78-7	97
2			Hexadecane	226	C16H34	000544-76-3	96
3			Hexadecane	226	C16H34	000544-76-3	93
4			Hexadecane	226	C16H34	000544-76-3	91
5			Hexadecane	226	C16H34	000544-76-3	83



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA M\DATA\BM101417\
Data File : BM012186.D
Acq On : 15 Oct 2017 00:41
Operator : SJ/JU
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Misc :
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Instrument :
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ClientSampleId :
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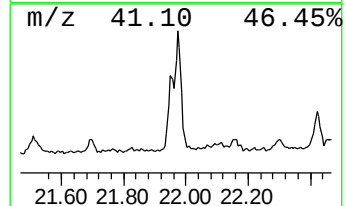
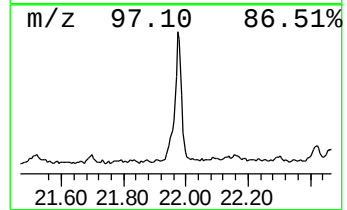
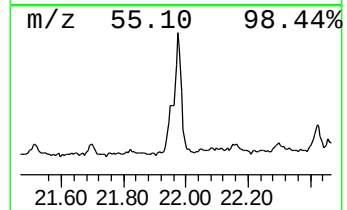
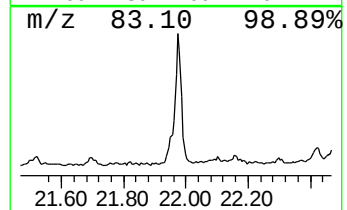
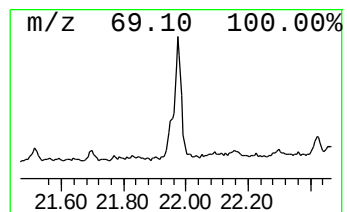
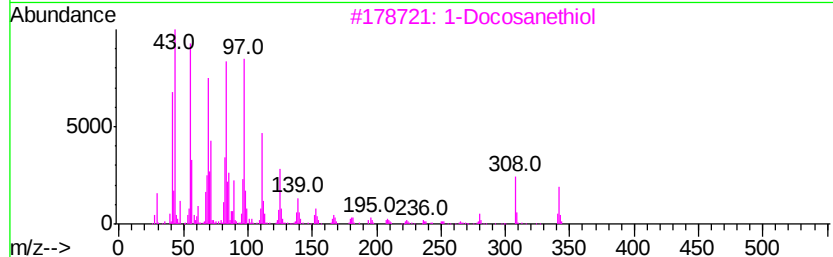
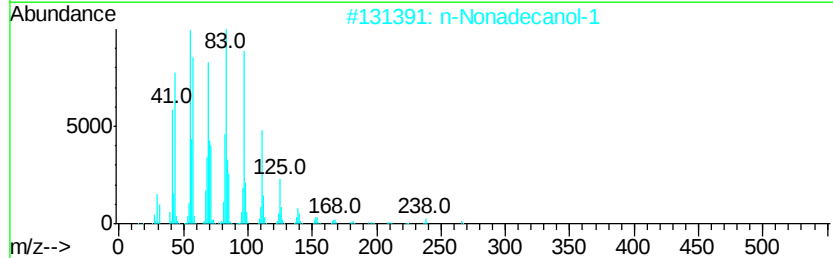
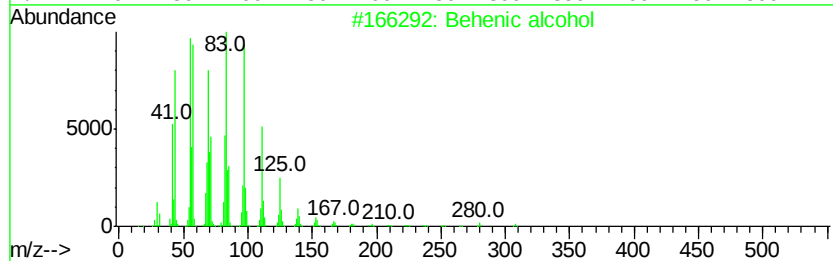
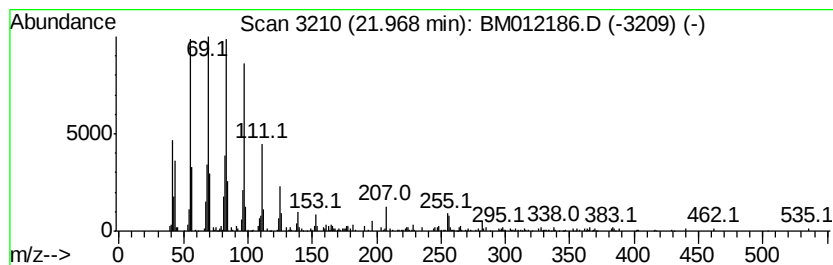
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Behenic alcohol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.97	4.47 ng/ul	1250980	Chrysene-d12	21.39

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Behenic alcohol		326	C22H46O	000661-19-8	94
2	n-Nonadecanol-1		284	C19H40O	001454-84-8	90
3	1-Docosanethiol		342	C22H46S	007773-83-3	80
4	Fumaric acid, octadecyl 2,2,3,3,...		582	C27H42F8O4	1000348-79-4	68
5	1-Octadecanol		270	C18H38O	000112-92-5	64



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA M\DATA\BM101417\
Data File : BM012186.D
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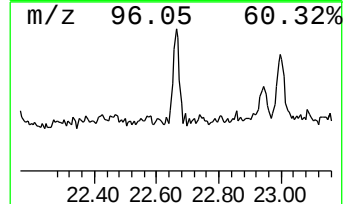
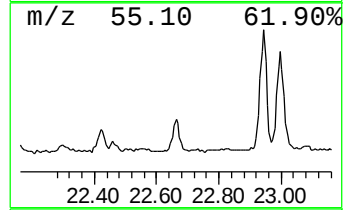
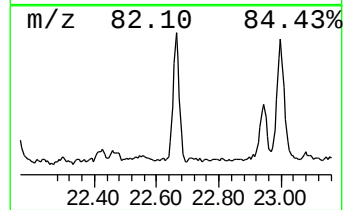
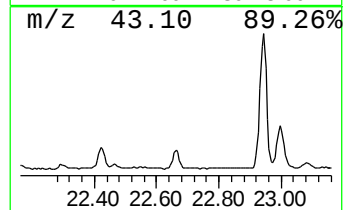
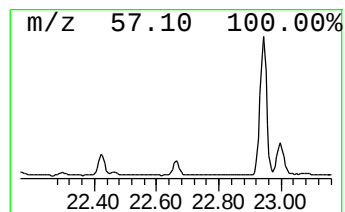
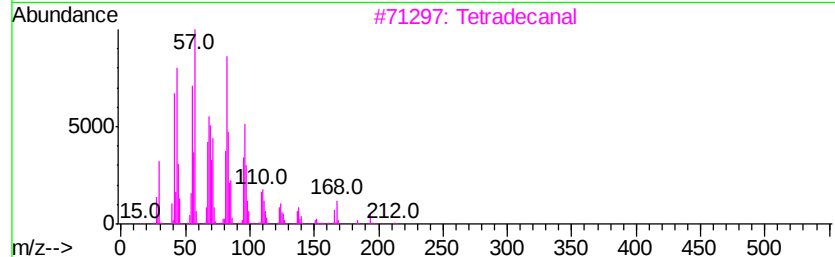
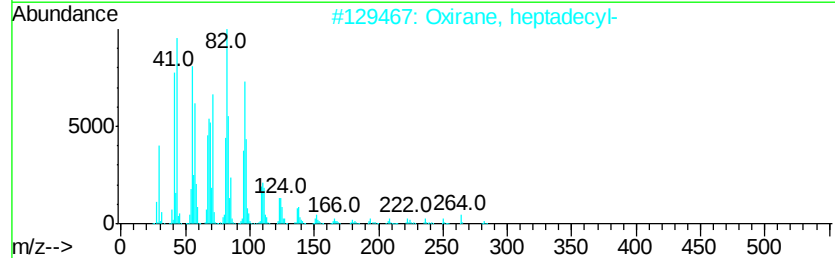
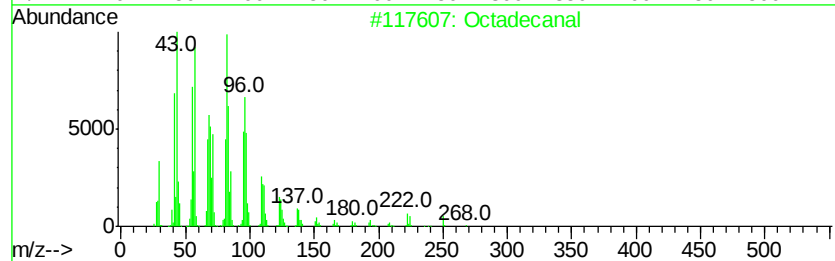
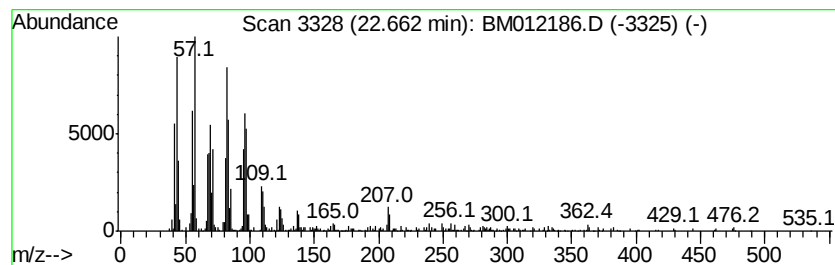
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Octadecanal Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.66	3.71 ng/ul	539141	Perylene-d12	23.70

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanal		268	C18H36O	000638-66-4	93
2	Oxirane, heptadecyl-		282	C19H38O	067860-04-2	93
3	Tetradecanal		212	C14H28O	000124-25-4	93
4	Oxirane, hexadecyl-		268	C18H36O	007390-81-0	87
5	Pentadecanal		226	C15H30O	002765-11-9	87



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA M\DATA\BM101417\
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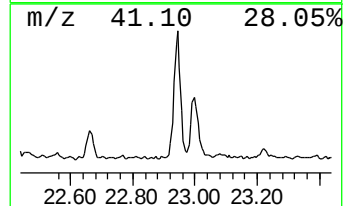
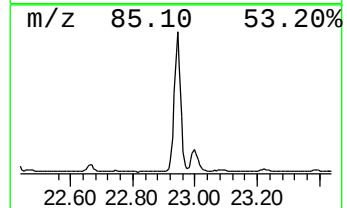
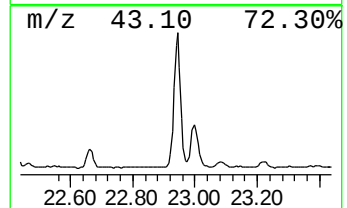
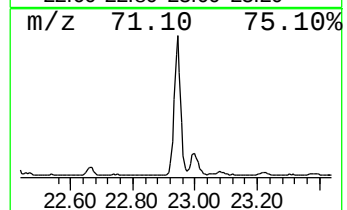
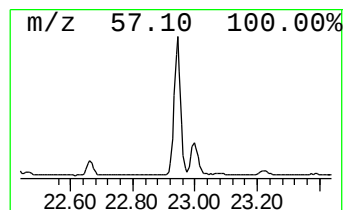
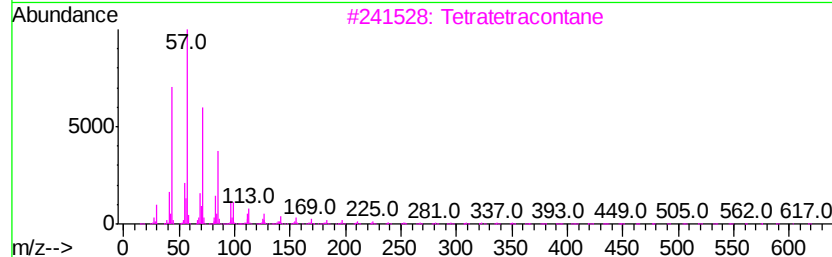
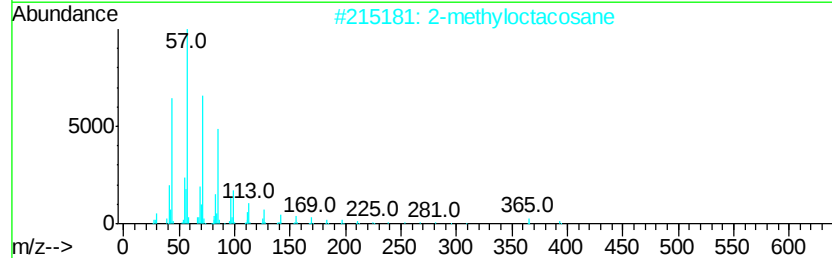
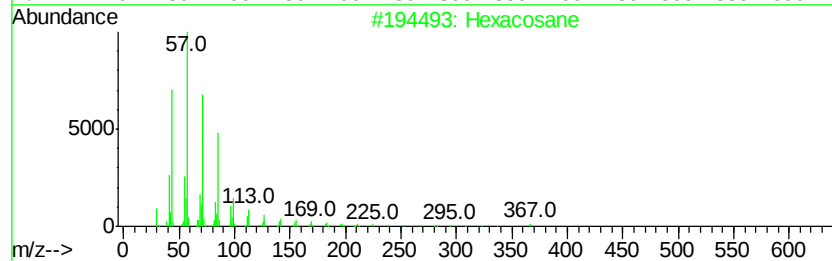
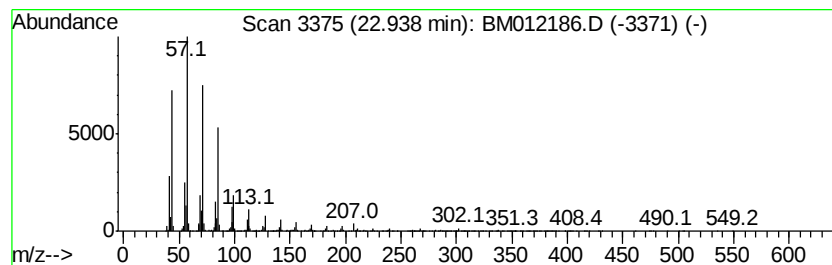
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.94	17.09 ng/ul	2484760	Perylene-d12	23.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosane	366	C26H54	000630-01-3	93
2		2-methyloctacosane	408	C29H60	1000376-72-8	91
3		Tetratetracontane	619	C44H90	007098-22-8	91
4		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
5		Heneicosane	296	C21H44	000629-94-7	90



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA M\DATA\BM101417\
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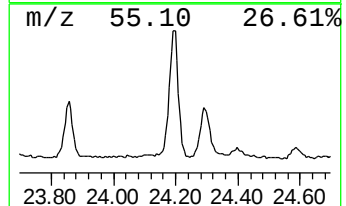
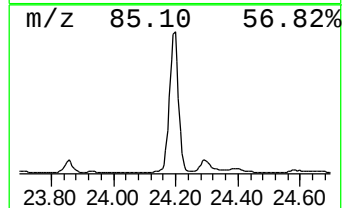
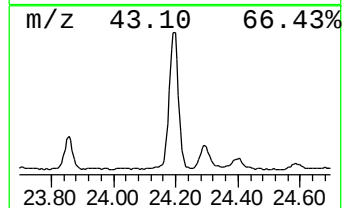
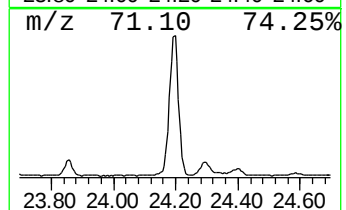
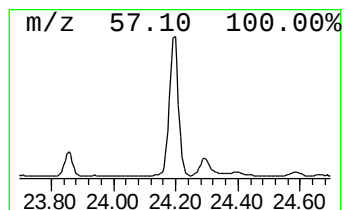
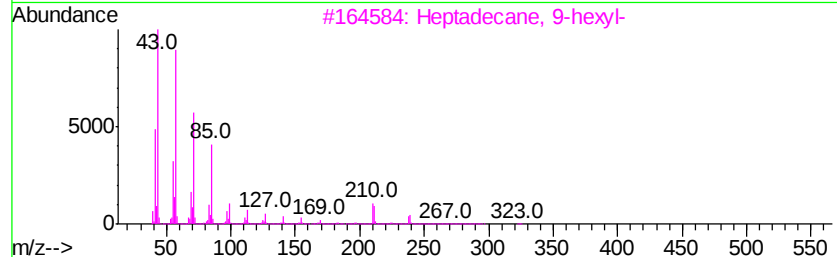
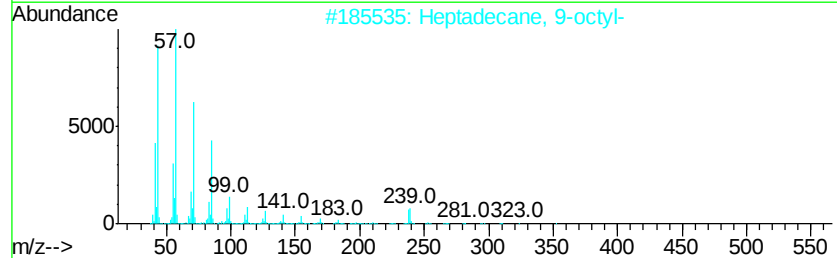
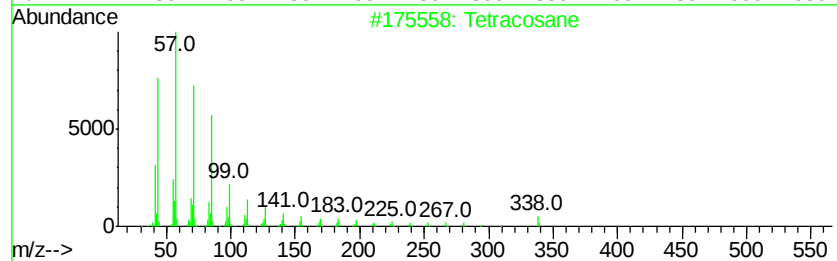
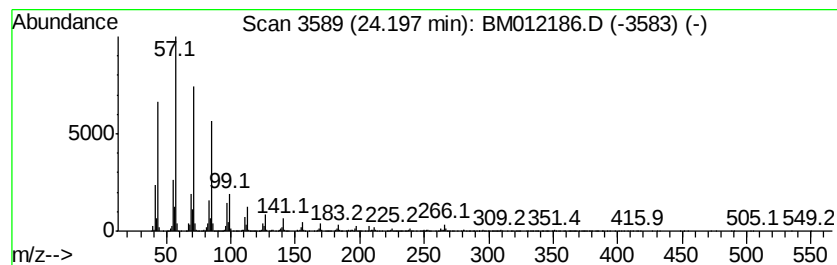
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.20	34.61 ng/ul	5032290	Perylene-d12	23.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetracosane	338	C24H50	000646-31-1	98
2		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
3		Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	93
4		Heptacosane	380	C27H56	000593-49-7	93
5		Docosane, 11-decyl-	451	C32H66	055401-55-3	91



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Acq On : 15 Oct 2017 00:41
Operator : SJ/JU
Sample : I5522-16
Misc :
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Instrument :
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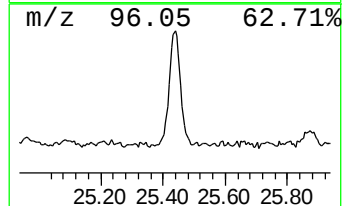
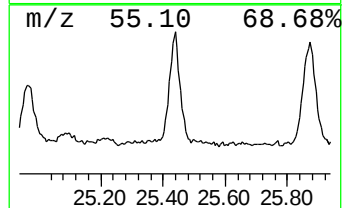
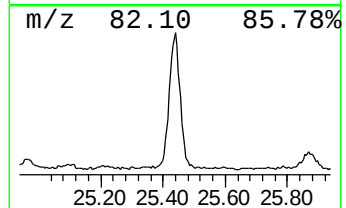
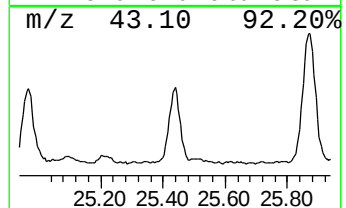
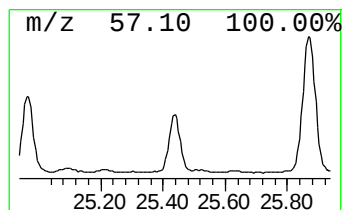
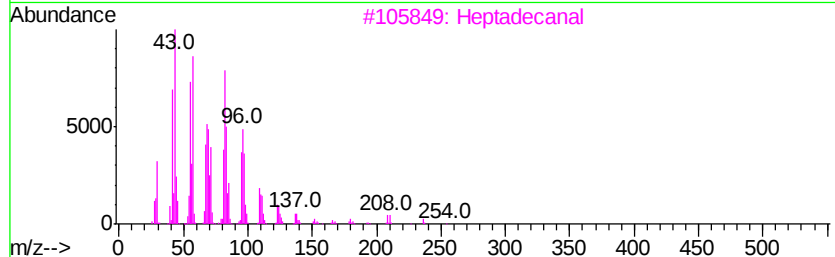
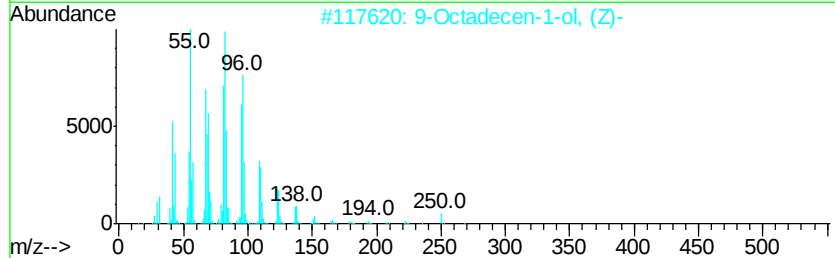
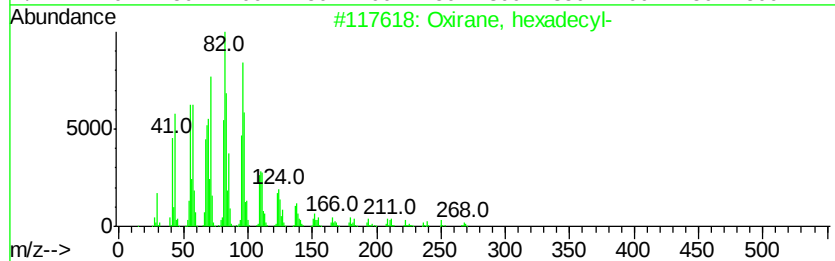
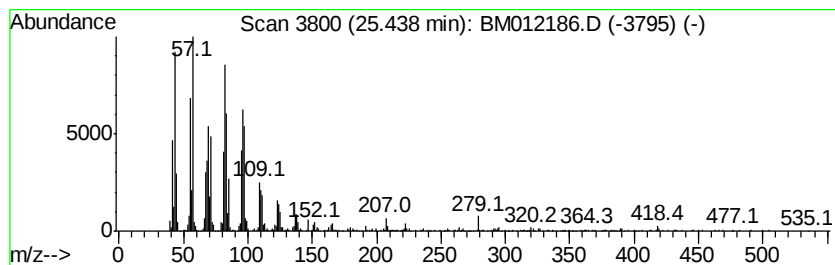
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 Oxirane, hexadecyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.44	11.62 ng/ul	1689500	Perylene-d12	23.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Oxirane, hexadecyl-	268	C18H36O	007390-81-0	94
2		9-Octadecen-1-ol, (Z)-	268	C18H36O	000143-28-2	92
3		Heptadecanal	254	C17H34O	1000376-70-0	91
4		Hexadecanal	240	C16H32O	000629-80-1	90
5		1,19-Eicosadiene	278	C20H38	014811-95-1	89



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA M\DATA\BM101417\
Data File : BM012186.D
Acq On : 15 Oct 2017 00:41
Operator : SJ/JU
Sample : I5522-16
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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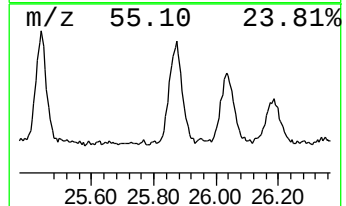
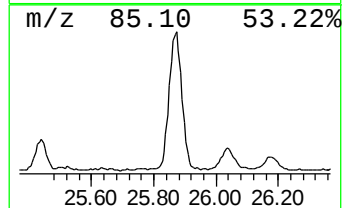
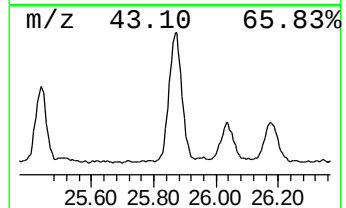
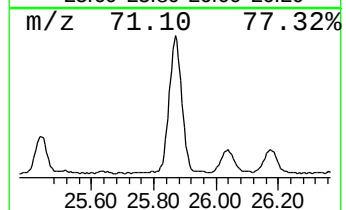
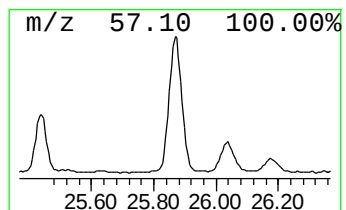
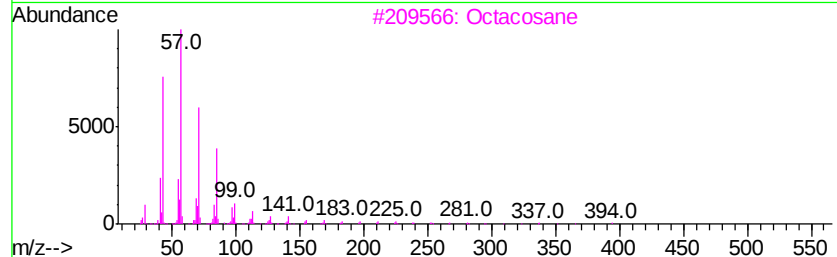
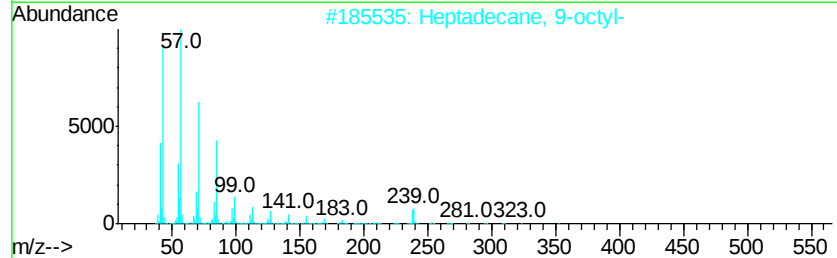
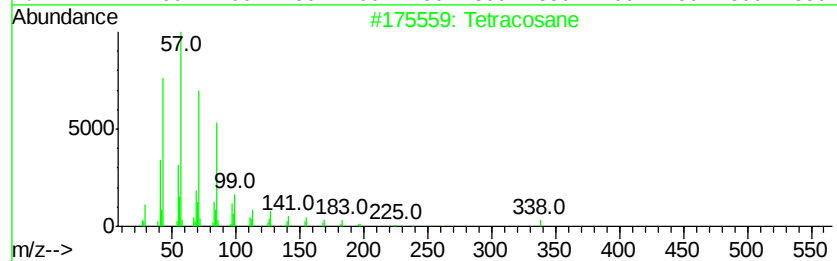
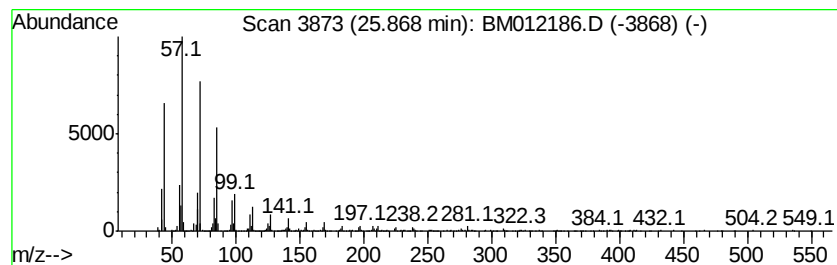
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.87	16.01 ng/ul	2327280	Perylene-d12	23.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetracosane	338	C24H50	000646-31-1	96
2		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	95
3		Octacosane	394	C28H58	000630-02-4	95
4		Docosane, 11-decyl-	451	C32H66	055401-55-3	91
5		2-methyloctacosane	408	C29H60	1000376-72-8	91



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_M\DATA\BM101417\
 Data File : BM012186.D
 Acq On : 15 Oct 2017 00:41
 Operator : SJ/JU
 Sample : I5522-16
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AC1

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM101217.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	18.09	10.3	ng/ul	1321920	4	17.21	2556380	20.0
Phenanthrene, 2-m...	18.13	2.2	ng/ul	277594	4	17.21	2556380	20.0
4H-Cyclopenta[def...	18.28	2.4	ng/ul	312286	4	17.21	2556380	20.0
13-Bromotetradeca...	18.39	2.9	ng/ul	371887	4	17.21	2556380	20.0
Octadecanoic acid	19.36	3.8	ng/ul	1063970	5	21.39	5591280	20.0
(DEL) Alkane: Cyc...	21.07	3.5	ng/ul	966137	5	21.39	5591280	20.0
(DEL) Alkane: Str...	21.95	2.7	ng/ul	763462	5	21.39	5591280	20.0
Behenic alcohol	21.97	4.5	ng/ul	1250980	5	21.39	5591280	20.0
Octadecanal	22.66	3.7	ng/ul	539141	6	23.70	2907730	20.0
(DEL) Alkane: Str...	22.94	17.1	ng/ul	2484760	6	23.70	2907730	20.0
(DEL) Alkane: Str...	24.20	34.6	ng/ul	5032290	6	23.70	2907730	20.0
Oxirane, hexadecyl-	25.44	11.6	ng/ul	1689500	6	23.70	2907730	20.0
(DEL) Alkane: Str...	25.87	16.0	ng/ul	2327280	6	23.70	2907730	20.0