

Data Path : Z:\HPCHEM1\BNA M\DATA\BM112217\
 Data File : BM013098.D
 Acq On : 22 Nov 2017 22:42
 Operator : SJ/JU
 Sample : I6514-09
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AD0

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-BM111617.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.246	23	27	35	rBV	98229	125546	1.48%	0.179%
2	6.887	640	646	659	rVB	1028419	1739277	20.53%	2.480%
3	7.051	665	674	683	rBV	832704	1247510	14.72%	1.779%
4	7.245	700	707	716	rVB	1106703	1722022	20.32%	2.455%
5	7.710	779	786	794	rBV	1028500	1599932	18.88%	2.281%
6	8.416	900	906	913	rBV	1324240	2091074	24.68%	2.981%
7	8.863	975	982	991	rBV	1085493	1769664	20.88%	2.523%
8	9.581	1097	1104	1111	rBV	922292	1489566	17.58%	2.124%
9	10.116	1185	1195	1206	rBV	1397029	2249291	26.54%	3.207%
10	10.492	1252	1259	1267	rVB	1385509	2226947	26.28%	3.175%
11	10.634	1273	1283	1295	rBV	1018083	1735583	20.48%	2.475%
12	13.769	1809	1816	1820	rBV	2335611	3324543	39.23%	4.740%
13	13.810	1820	1823	1828	rVB	402354	508365	6.00%	0.725%
14	14.039	1853	1862	1870	rBV	2479953	3826420	45.16%	5.456%
15	14.263	1894	1900	1906	rBV	107006	144039	1.70%	0.205%
16	14.351	1908	1915	1922	rBV2	1903074	2822818	33.31%	4.025%
17	14.557	1941	1950	1962	rBV	686468	1061679	12.53%	1.514%
18	15.180	2051	2056	2059	rBV	103532	148373	1.75%	0.212%
19	15.216	2059	2062	2067	rVB	123739	150291	1.77%	0.214%
20	15.345	2075	2084	2091	rBV	3144107	4428097	52.26%	6.313%
21	15.463	2097	2104	2112	rBV	802778	1079912	12.74%	1.540%
22	16.080	2204	2209	2216	rBV2	72638	117993	1.39%	0.168%
23	17.098	2376	2382	2388	rBV2	2165162	3030159	35.76%	4.320%
24	17.198	2393	2399	2407	rVB	3060576	4355498	51.40%	6.210%
25	18.004	2531	2536	2544	rBV2	368884	519871	6.14%	0.741%
26	19.286	2750	2754	2762	rBV	312163	443047	5.23%	0.632%
27	19.492	2783	2789	2797	rBV	3154898	4310723	50.87%	6.146%
28	21.009	3042	3047	3056	rBV2	615206	966924	11.41%	1.379%
29	21.139	3065	3069	3074	rBV4	135354	188037	2.22%	0.268%
30	21.198	3076	3079	3082	rBV3	93395	121917	1.44%	0.174%
31	21.280	3088	3093	3098	rVB	2247445	2818553	33.26%	4.019%
32	21.886	3192	3196	3205	rBV3	319516	575211	6.79%	0.820%
33	22.045	3220	3223	3228	rVB	265096	327528	3.87%	0.467%
34	22.192	3245	3248	3252	rBV6	123455	143081	1.69%	0.204%

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Integration Parameters: LSCINT.P

Integrator: RTE
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35	22.256	3257	3259	3264	rVB6	141077	187392	2.21%	0.267%
36	22.321	3264	3270	3277	rBV	6331992	8473808	100.00%	12.082%
37	22.850	3357	3360	3364	rBV	214216	281053	3.32%	0.401%
38	22.897	3364	3368	3374	rVB2	287124	427229	5.04%	0.609%
39	23.368	3441	3448	3454	rBV2	1962392	3637325	42.92%	5.186%
40	23.509	3466	3472	3479	rVB2	1321131	2496521	29.46%	3.559%
41	24.068	3562	3567	3575	rVB3	291219	559321	6.60%	0.797%
42	24.150	3577	3581	3589	rVB7	171734	342409	4.04%	0.488%
43	24.815	3690	3694	3701	rVB4	161581	323956	3.82%	0.462%

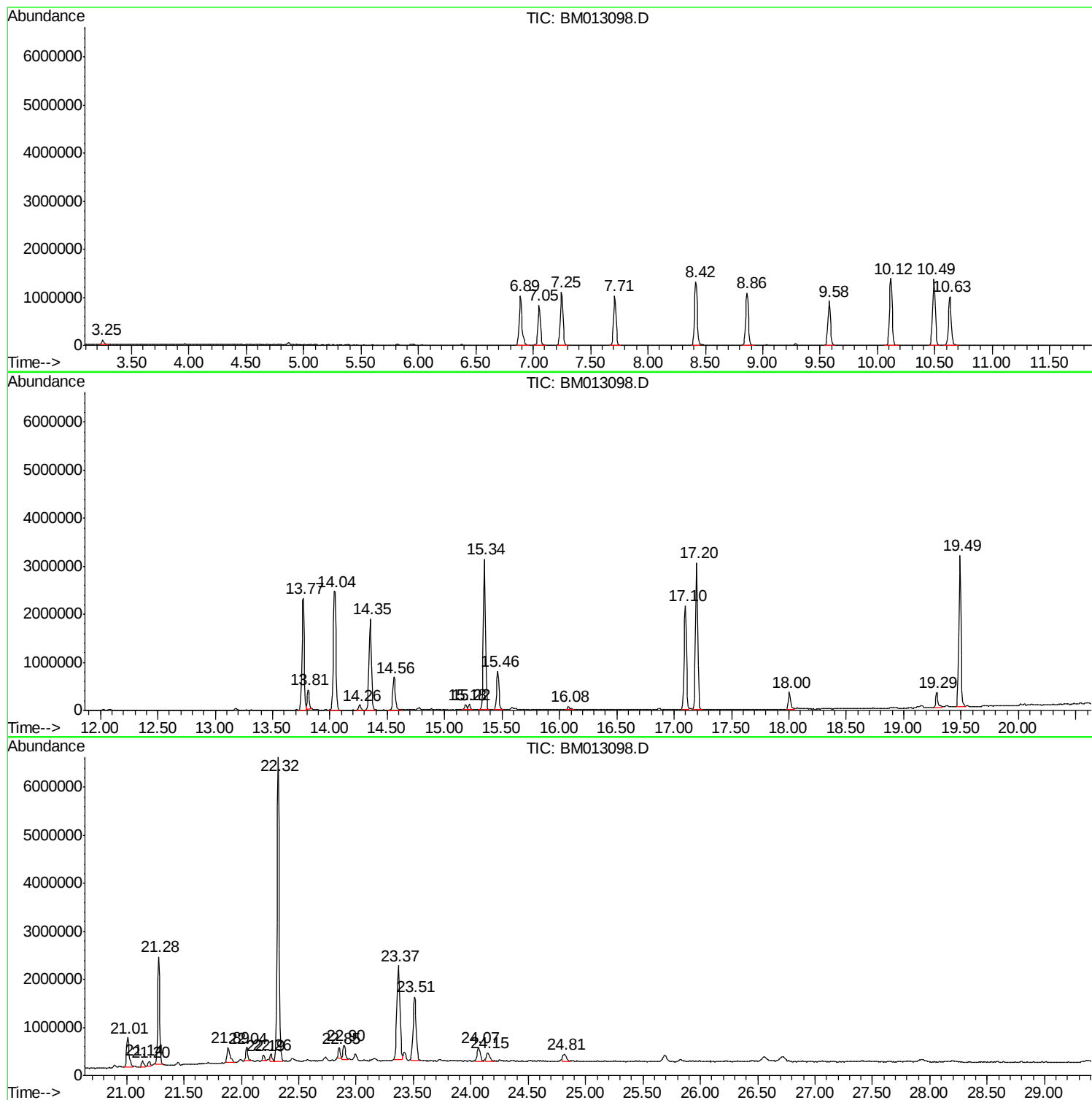
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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM111617.M
Quant Title : SVOA CALIBRATION

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



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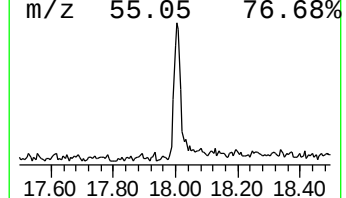
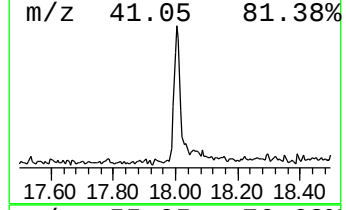
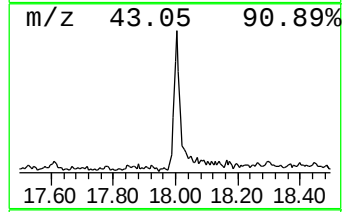
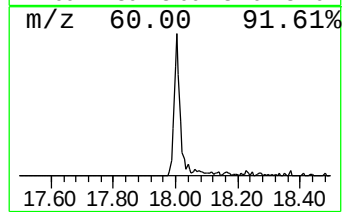
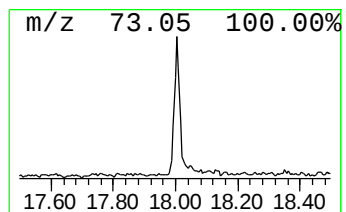
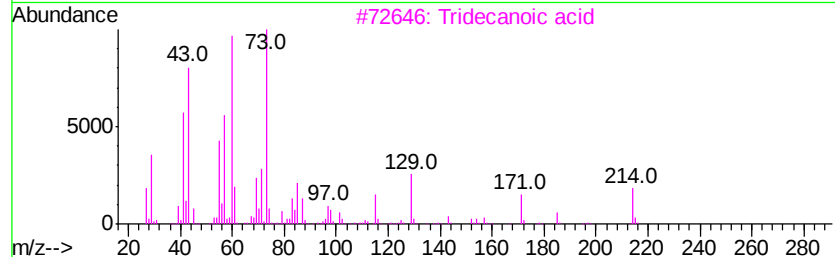
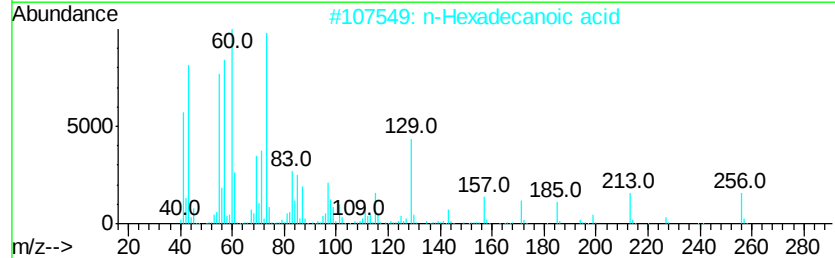
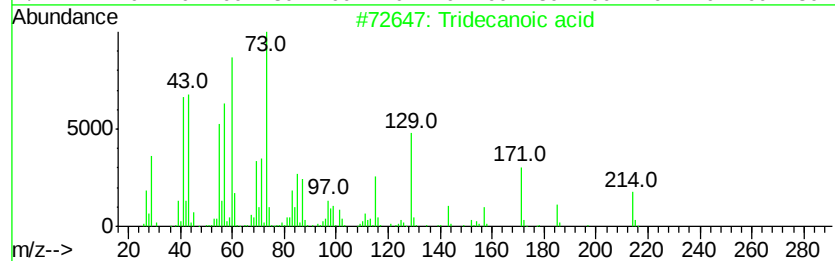
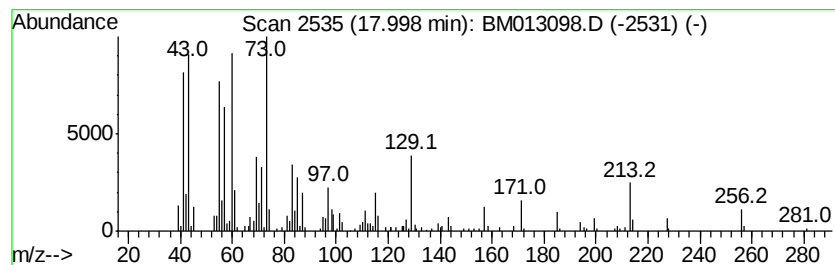
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM111617.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Tridecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.00	3.43 ng/ul	519871	Phenanthrene-d10	17.10

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tridecanoic acid	214	C13H26O2	000638-53-9	97
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
3		Tridecanoic acid	214	C13H26O2	000638-53-9	95
4		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
5		Tridecanoic acid	214	C13H26O2	000638-53-9	93



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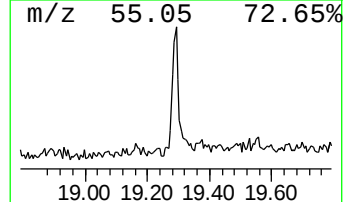
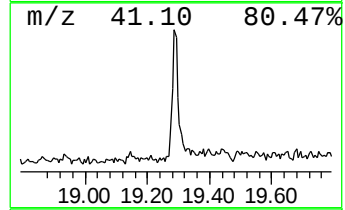
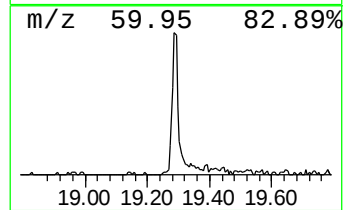
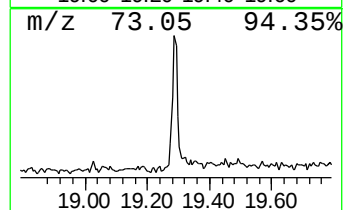
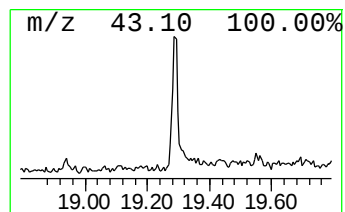
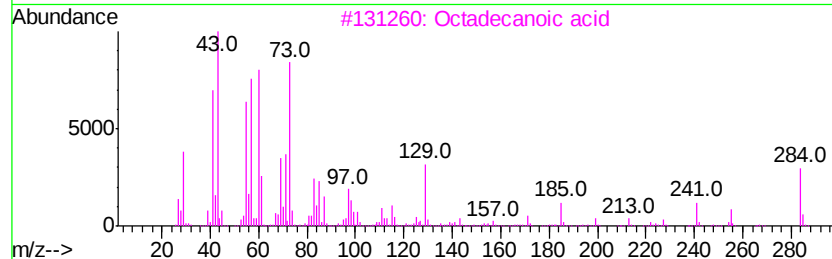
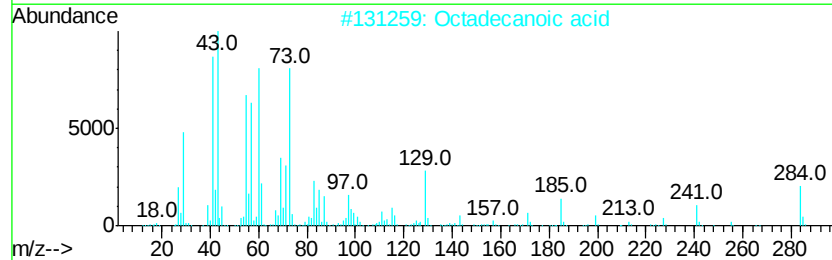
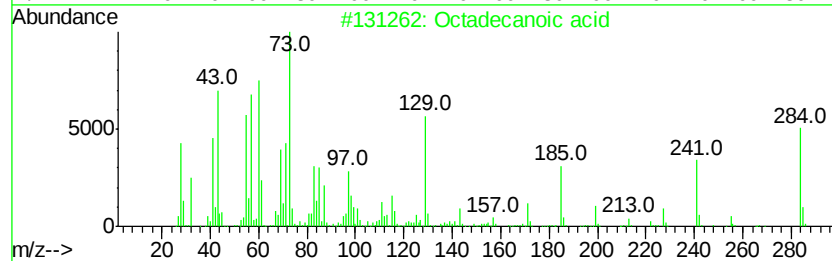
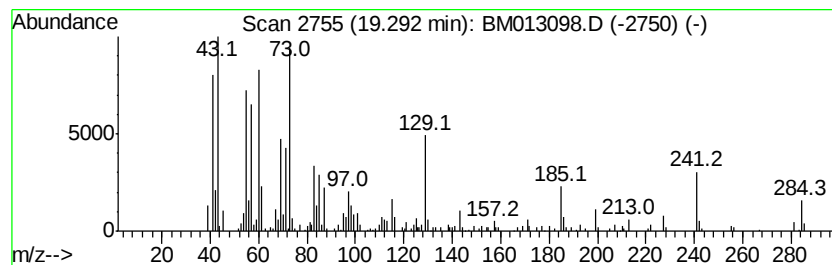
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Peak Number 2 Octadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.
19.29	3.14 ng/ul	443047	Chrysene-d12		21.28

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	98
2			Octadecanoic acid	284	C18H36O2	000057-11-4	98
3			Octadecanoic acid	284	C18H36O2	000057-11-4	97
4			Octadecanoic acid	284	C18H36O2	000057-11-4	97
5			Octadecanoic acid	284	C18H36O2	000057-11-4	95



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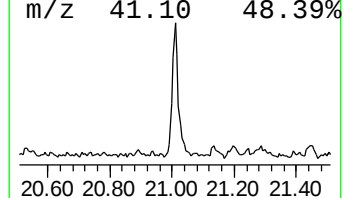
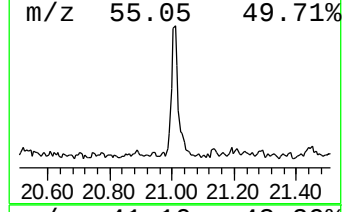
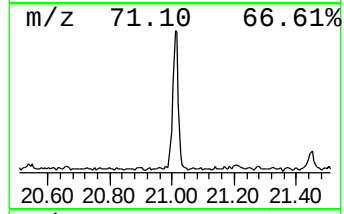
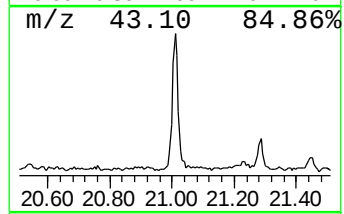
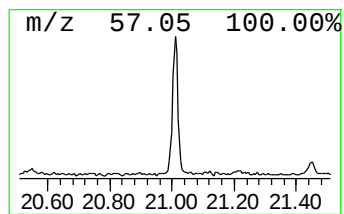
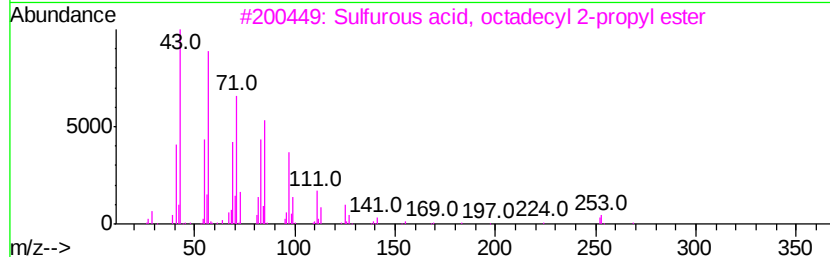
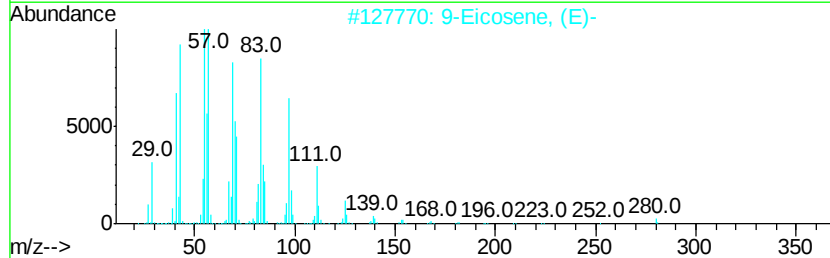
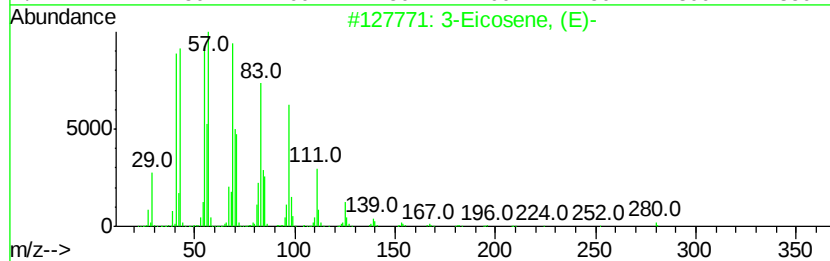
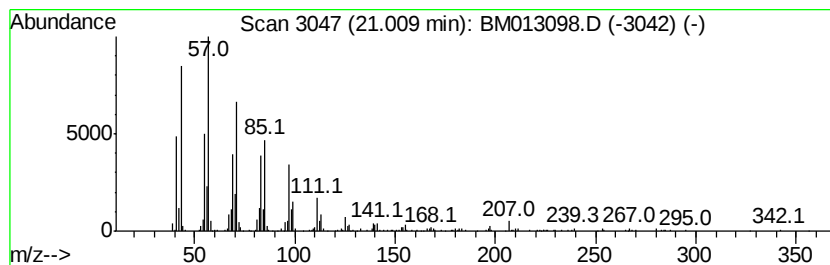
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 (DEL) Alkane: Cyclic21.01 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.01	6.86 ng/ul	966924	Chrysene-d12	21.28

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Eicosene, (E)-	280	C20H40	074685-33-9	96
2			9-Eicosene, (E)-	280	C20H40	074685-29-3	90
3			Sulfurous acid, octadecyl 2-propyl ...	376	C21H44O3S	1000309-12-7	87
4			Oxalic acid, isobutyl hexadecyl ...	370	C22H42O4	1000309-38-1	74
5			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	64



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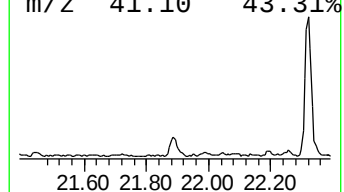
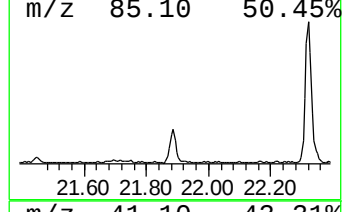
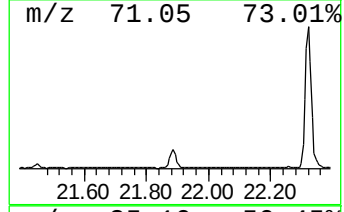
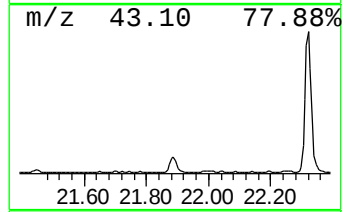
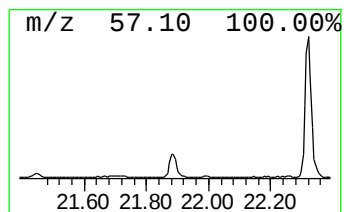
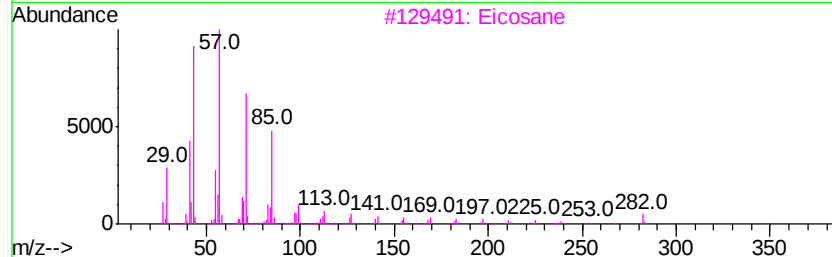
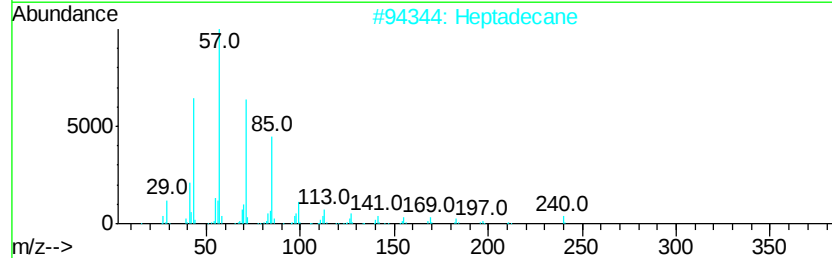
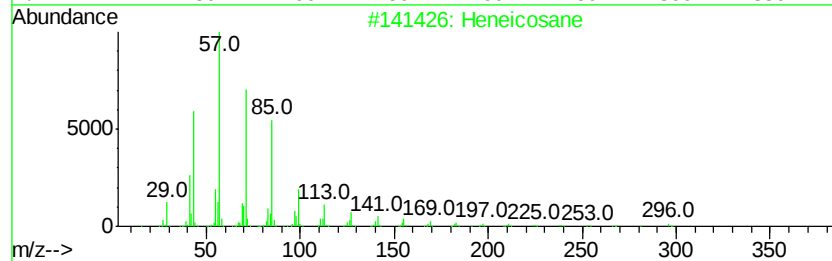
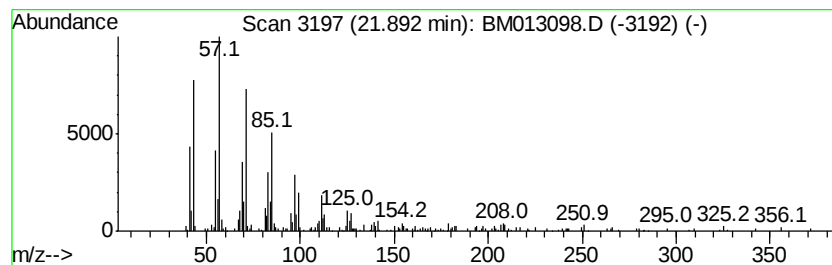
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Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.89	4.08 ng/ul	575211	Chrysene-d12	21.28

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	98
2			Heptadecane	240	C17H36	000629-78-7	94
3			Eicosane	282	C20H42	000112-95-8	94
4			Nonadecane	268	C19H40	000629-92-5	93
5			Heneicosane	296	C21H44	000629-94-7	93



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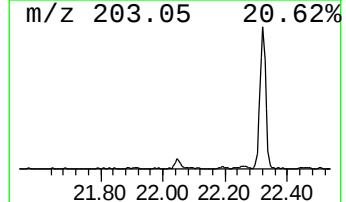
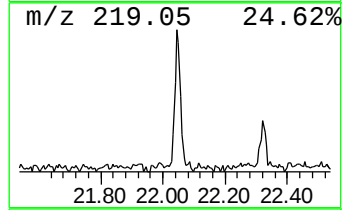
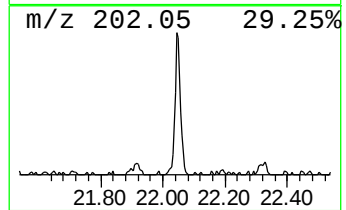
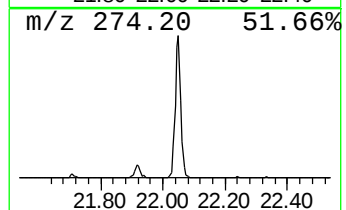
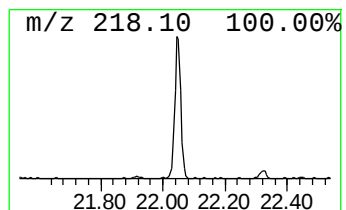
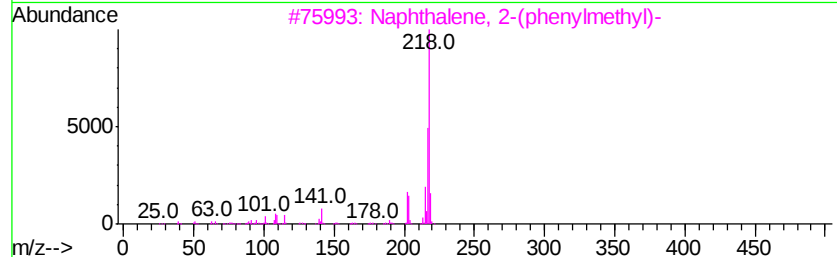
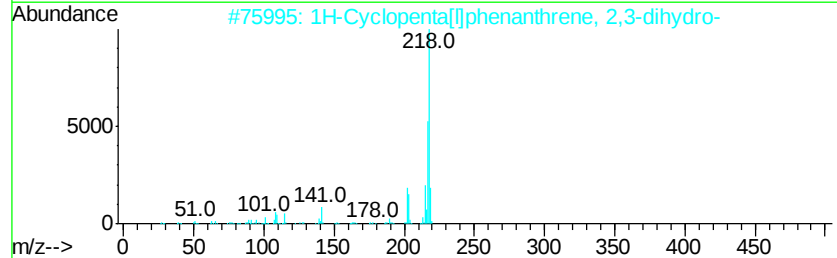
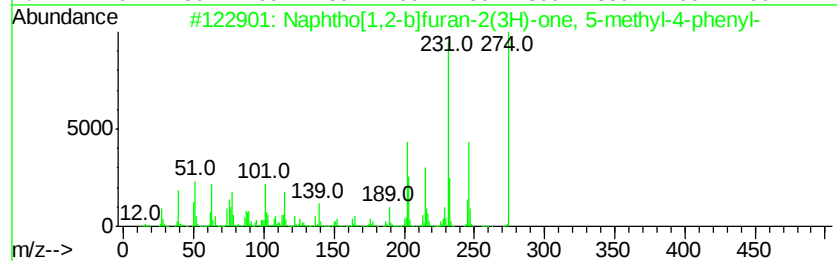
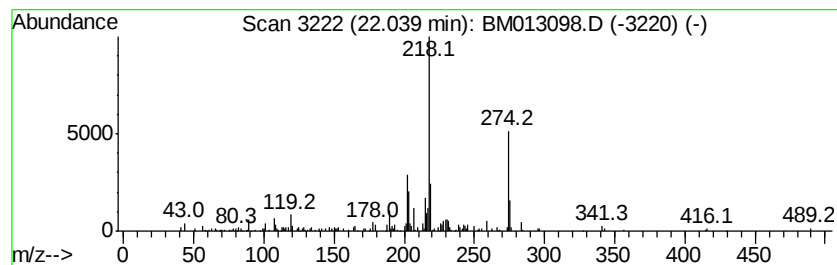
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TIC Integration Parameters: LSCINT.P

Peak Number 5 Naphtho[1,2-b]furan-2(3H)-o... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.04	2.32 ng/ul	327528	Chrysene-d12	21.28

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphtho[1,2-b]furan-2(3H)-one, 5...	274	C19H14O2	077573-41-2	60
2		1H-Cyclopenta[1]phenanthrene, 2,...	218	C17H14	000723-98-8	49
3		Naphthalene, 2-(phenylmethyl)-	218	C17H14	000613-59-2	47
4		2,4-Diethoxyquinazoline	218	C12H14N2O2	000611-65-4	46
5		7-Isopropenyl-1,4a-dimethyl-4,4a...	218	C15H22O	000473-08-5	41



Data Path : Z:\HPCHEM1\BNA M\DATA\BM112217\
Data File : BM013098.D
Acq On : 22 Nov 2017 22:42
Operator : SJ/JU
Sample : I6514-09
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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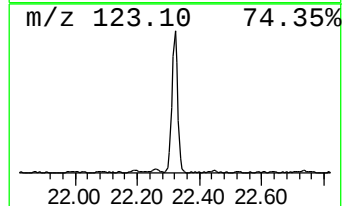
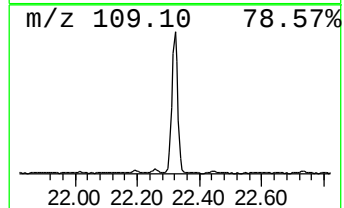
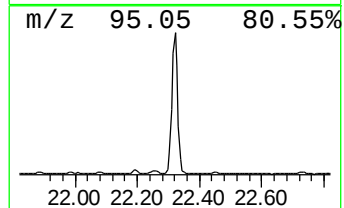
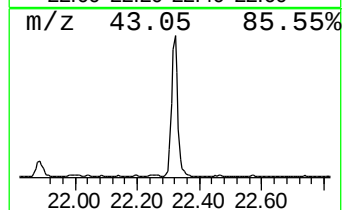
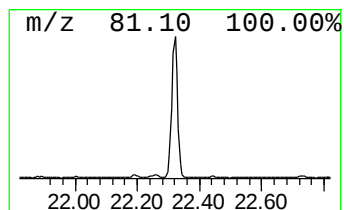
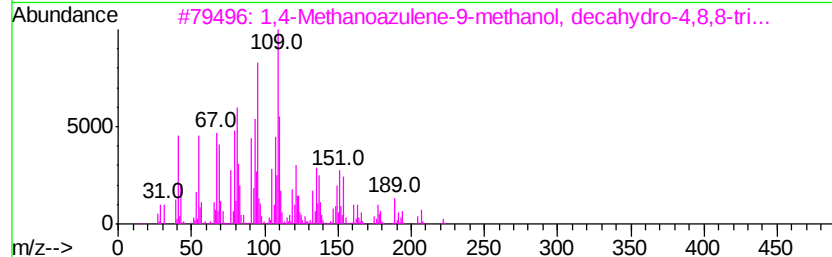
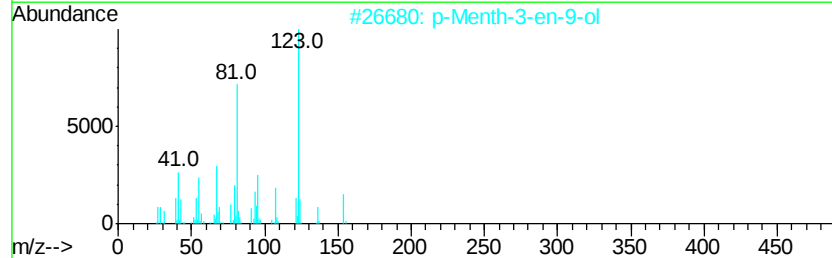
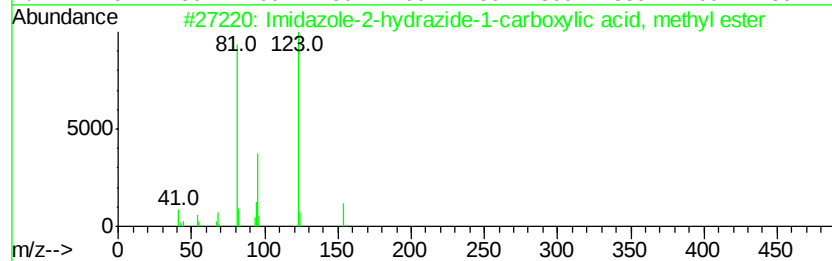
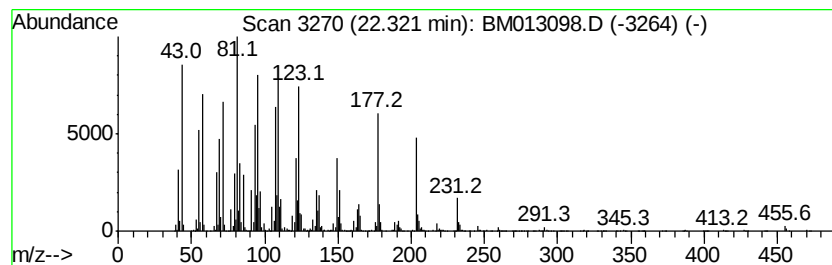
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.32	60.13 ng/ul	8473810	Chrysene-d12	21.28

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Imidazole-2-hydrazide-1-carboxyl...	154	C5H6N4O2	1000126-42-2	27
2		p-Menth-3-en-9-ol	154	C10H18O	015714-10-0	27
3		1,4-Methanoazulene-9-methanol, d...	222	C15H26O	001139-17-9	25
4		Vinyltris(cyclopropyl)silane	178	C11H18Si	1000109-97-3	22
5		Bicyclo[2.2.2]octane, 1-methyl-4...	202	C10H18O2S	069855-48-7	22



Data Path : Z:\HPCHEM1\BNA M\DATA\BM112217\
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Acq On : 22 Nov 2017 22:42
Operator : SJ/JU
Sample : I6514-09
Misc :
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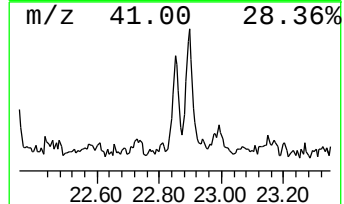
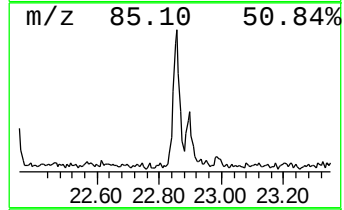
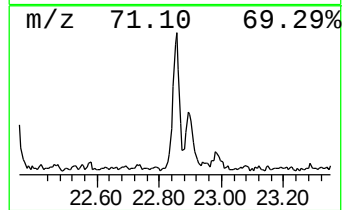
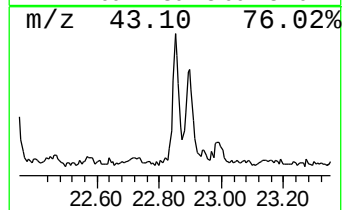
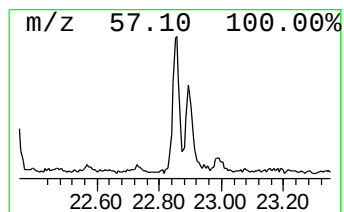
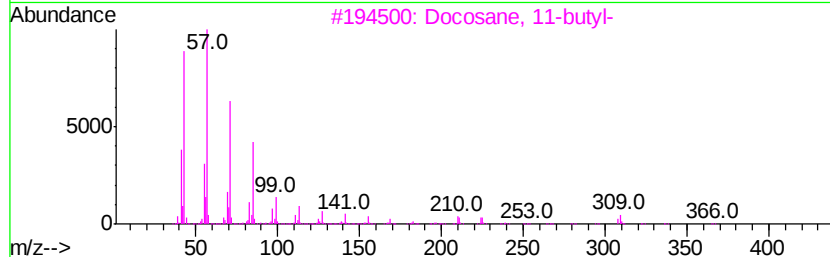
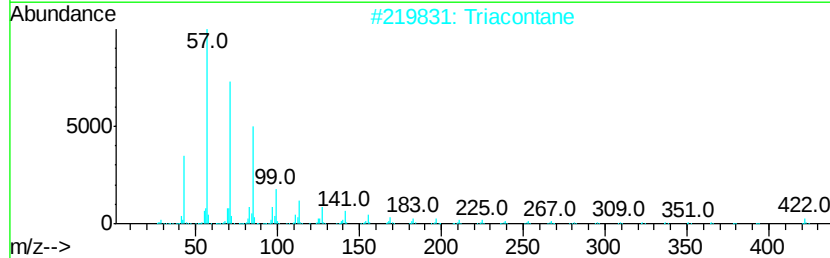
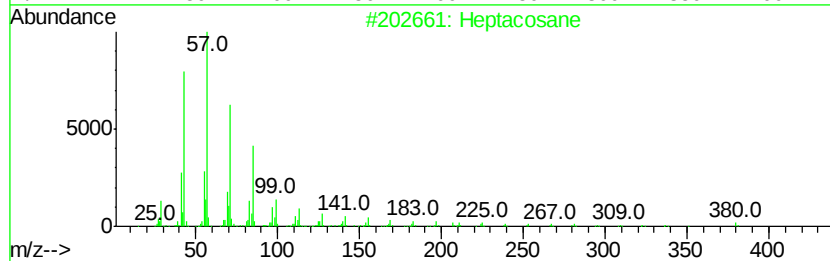
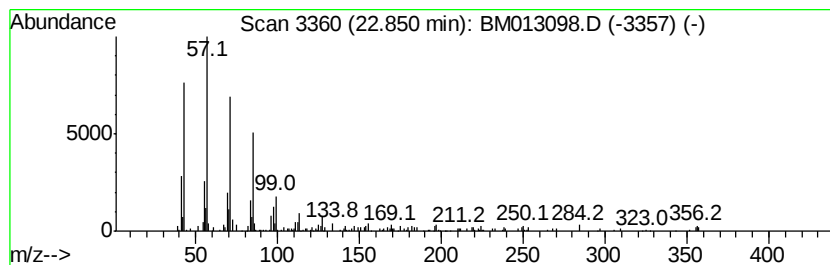
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.85	2.25 ng/ul	281053	Perylene-d12	23.51
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Heptacosane	380	C27H56	000593-49-7 90
2	Triacosane	422	C30H62	000638-68-6 87
3	Docosane, 11-butyl-	366	C26H54	013475-76-8 87
4	Heptadecane, 9-octyl-	352	C25H52	007225-64-1 87
5	Tetracosane	338	C24H50	000646-31-1 87



Data Path : Z:\HPCHEM1\BNA M\DATA\BM112217\
Data File : BM013098.D
Acq On : 22 Nov 2017 22:42
Operator : SJ/JU
Sample : I6514-09
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Instrument :
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ClientSampleId :
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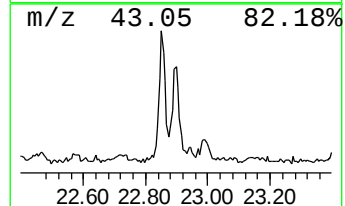
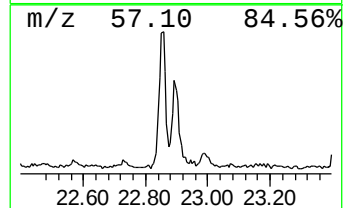
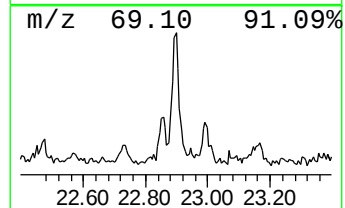
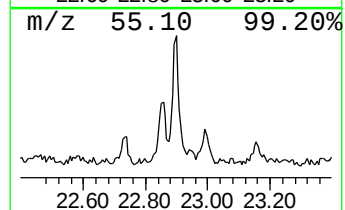
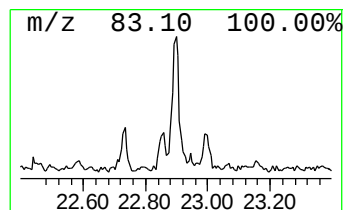
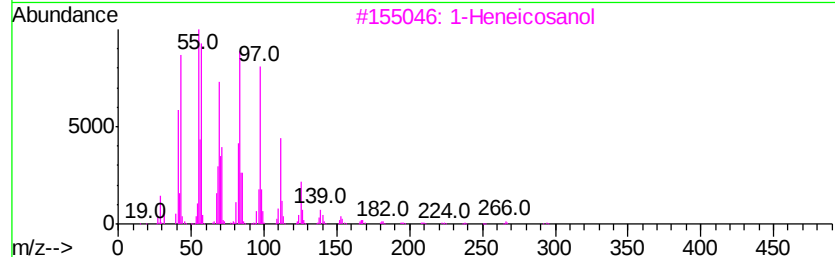
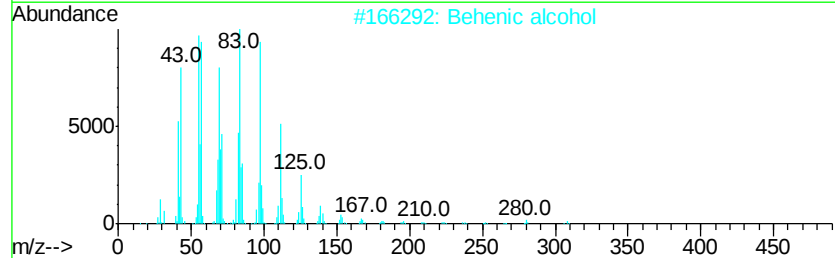
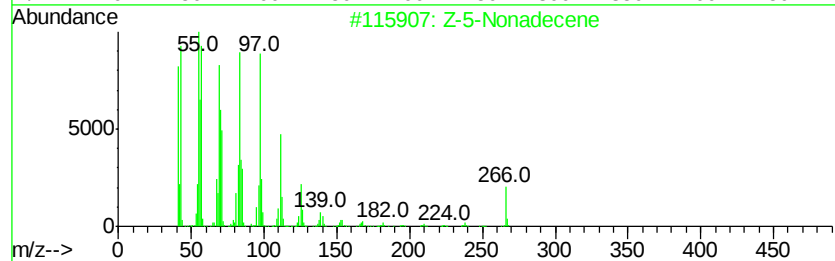
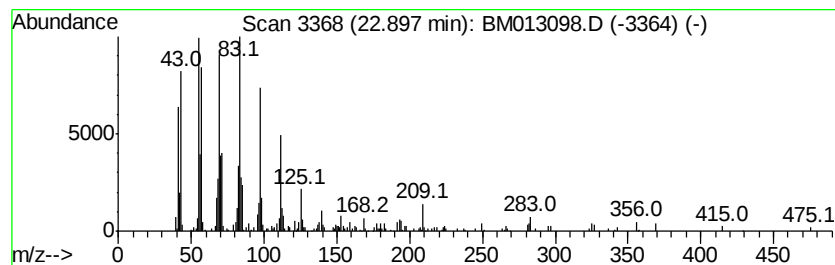
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Cyclic22.90 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.90	3.42 ng/ul	427229	Perylene-d12	23.51

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Z-5-Nonadecene	266	C19H38	1000131-11-8	95
2		Behenic alcohol	326	C22H46O	000661-19-8	93
3		1-Heneicosanol	312	C21H44O	015594-90-8	91
4		Fumaric acid, 2-chloropropyl dod...	360	C19H33ClO4	1000348-57-0	87
5		n-Pentadecanol	228	C15H32O	000629-76-5	87



Data Path : Z:\HPCHEM1\BNA M\DATA\BM112217\
Data File : BM013098.D
Acq On : 22 Nov 2017 22:42
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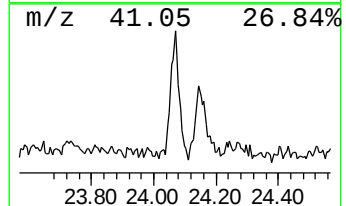
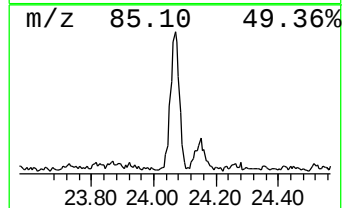
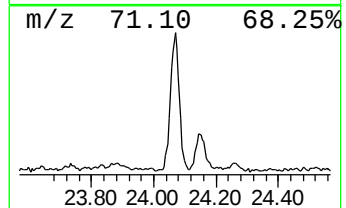
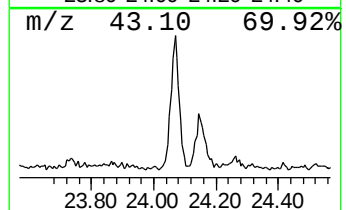
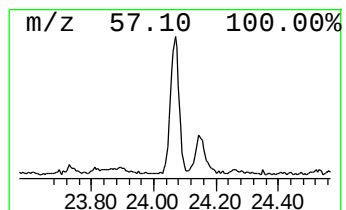
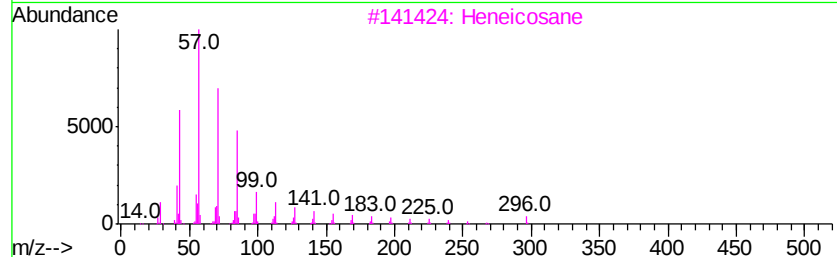
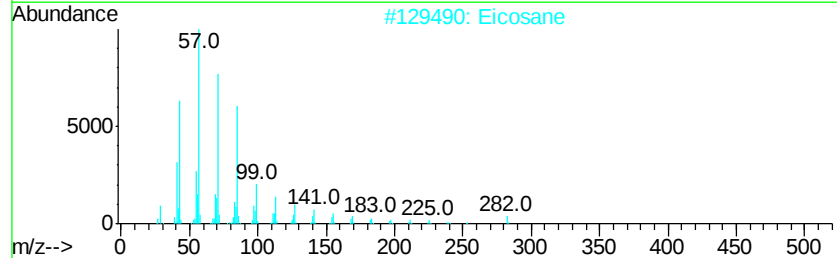
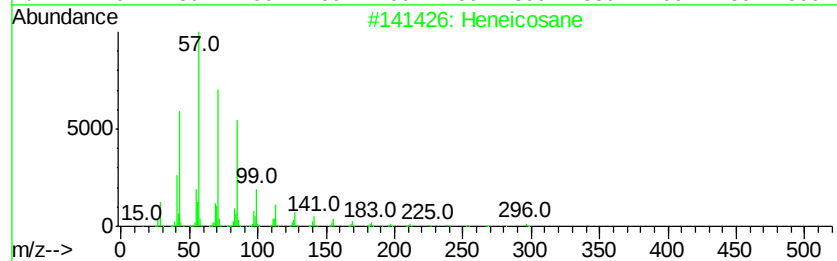
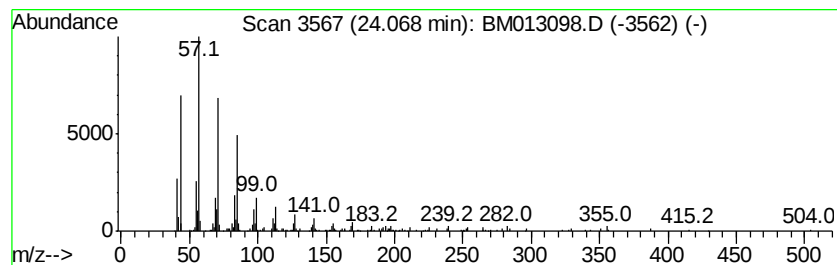
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.07	4.48 ng/ul	559321	Perylene-d12	23.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	98
2		Eicosane	282	C20H42	000112-95-8	97
3		Heneicosane	296	C21H44	000629-94-7	95
4		Eicosane	282	C20H42	000112-95-8	93
5		Eicosane, 2-methyl-	296	C21H44	001560-84-5	93



Data Path : Z:\HPCHEM1\BNA M\DATA\BM112217\
Data File : BM013098.D
Acq On : 22 Nov 2017 22:42
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Misc :
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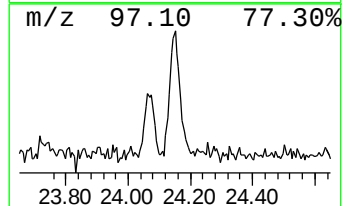
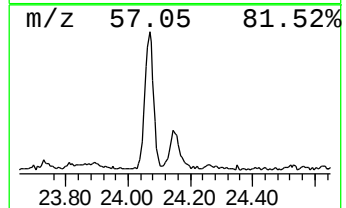
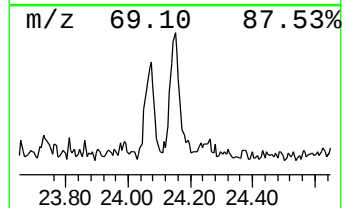
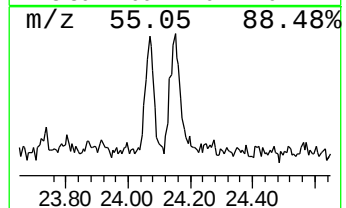
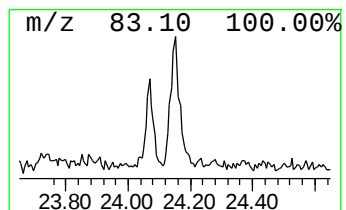
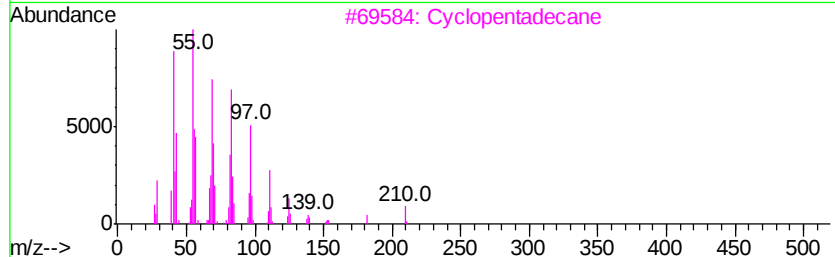
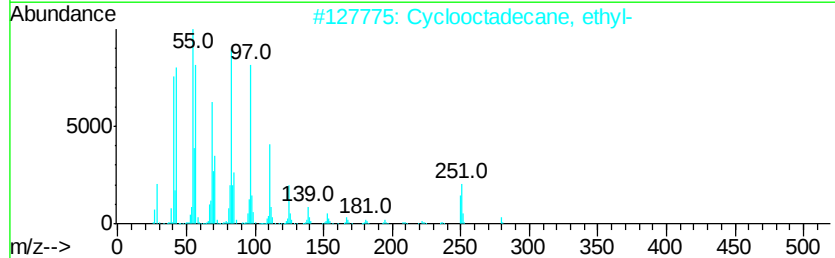
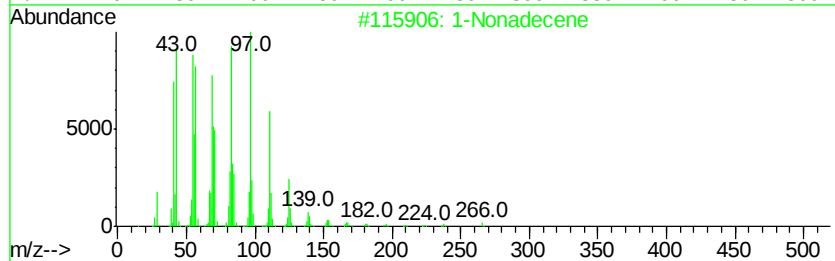
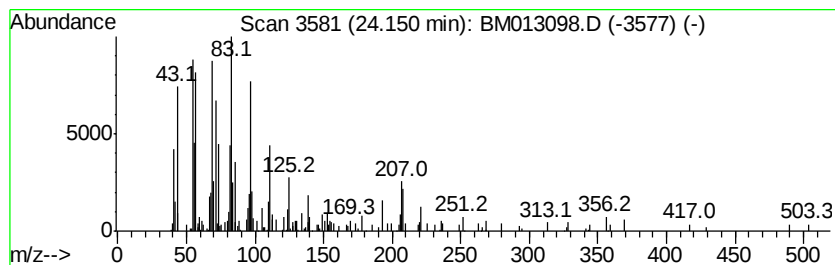
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Cyclic24.15 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.15	2.74 ng/ul	342409	Perylene-d12	23.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Nonadecene	266	C19H38	018435-45-5	83
2		Cyclooctadecane, ethyl-	280	C20H40	1000151-22-5	78
3		Cyclopentadecane	210	C15H30	000295-48-7	72
4		Bromoacetic acid, octadecyl ester	390	C20H39BrO2	018992-03-5	72
5		9-Eicosene, (E)-	280	C20H40	074685-29-3	64



Data Path : Z:\HPCHEM1\BNA M\DATA\BM112217\
Data File : BM013098.D
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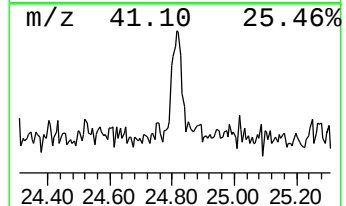
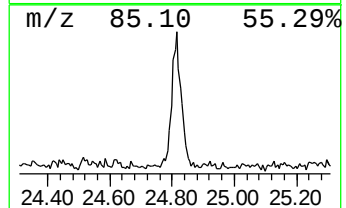
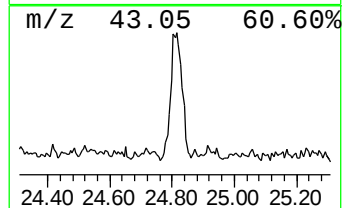
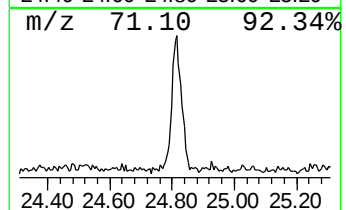
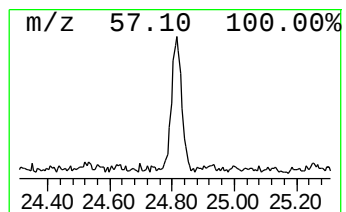
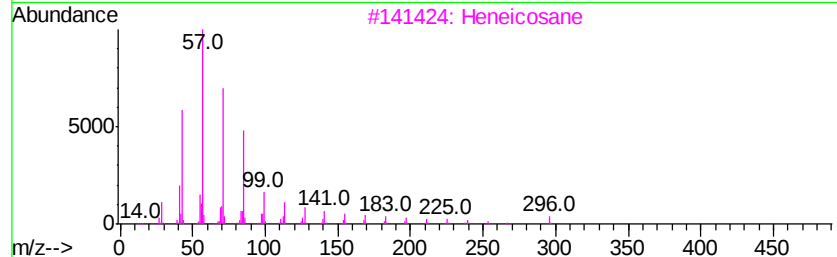
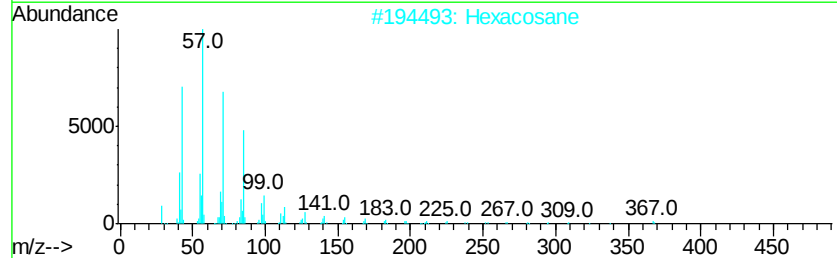
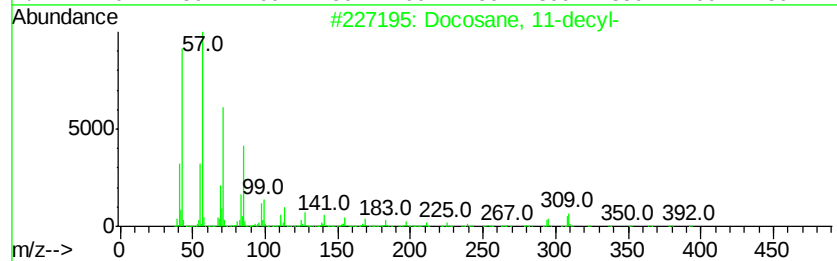
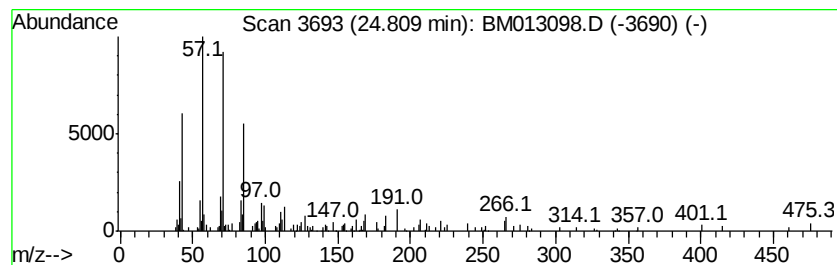
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.81	2.60 ng/ul	323956	Perylene-d12	23.51

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Docosane, 11-decyl-	451	C32H66	055401-55-3	72
2			Hexacosane	366	C26H54	000630-01-3	64
3			Heneicosane	296	C21H44	000629-94-7	64
4			Tricosane	324	C23H48	000638-67-5	64
5			Hexatriacontane	507	C36H74	000630-06-8	64



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM112217\
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Acq On : 22 Nov 2017 22:42
Operator : SJ/JU
Sample : I6514-09
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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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
Tridecanoic acid	18.00	3.4	ng/ul	519871	4	17.10	3030160	20.0
Octadecanoic acid	19.29	3.1	ng/ul	443047	5	21.28	2818550	20.0
(DEL) Alkane: Cyc...	21.01	6.9	ng/ul	966924	5	21.28	2818550	20.0
(DEL) Alkane: Str...	21.89	4.1	ng/ul	575211	5	21.28	2818550	20.0
Naphtho[1,2-b]fur...	22.04	2.3	ng/ul	327528	5	21.28	2818550	20.0
unknown-01	22.32	60.1	ng/ul	8473810	5	21.28	2818550	20.0
(DEL) Alkane: Str...	22.85	2.3	ng/ul	281053	6	23.51	2496520	20.0
(DEL) Alkane: Cyc...	22.90	3.4	ng/ul	427229	6	23.51	2496520	20.0
(DEL) Alkane: Str...	24.07	4.5	ng/ul	559321	6	23.51	2496520	20.0
(DEL) Alkane: Cyc...	24.15	2.7	ng/ul	342409	6	23.51	2496520	20.0
(DEL) Alkane: Str...	24.81	2.6	ng/ul	323956	6	23.51	2496520	20.0