

Data Path : Z:\HPCHEM1\BNA M\DATA\BM101717\
 Data File : BM012277.D
 Acq On : 18 Oct 2017 08:20
 Operator : SJ/JU
 Sample : I5747-02
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 B-23(1-3)-100517

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	3.169	9	14	18	rBV	155260	223961	1.59%	0.249%
2	3.551	74	79	89	rVB	133219	188840	1.34%	0.210%
3	3.946	140	146	168	rVB	9878806	14068895	100.00%	15.665%
4	4.475	230	236	252	rBV	4613879	6865634	48.80%	7.645%
5	5.269	365	371	390	rBV	612018	1007480	7.16%	1.122%
6	6.881	628	645	649	rBV	320389	1052548	7.48%	1.172%
7	6.957	649	658	676	rVB2	895471	2127607	15.12%	2.369%
8	7.116	678	685	696	rVV	678170	1076490	7.65%	1.199%
9	7.316	711	719	727	rBV	1052406	1922832	13.67%	2.141%
10	7.781	792	798	813	rVB	742014	1261564	8.97%	1.405%
11	8.498	913	920	934	rBV	766044	1416883	10.07%	1.578%
12	8.951	990	997	1013	rBV	672401	1222900	8.69%	1.362%
13	9.675	1113	1120	1130	rBV	601146	1098793	7.81%	1.223%
14	9.957	1156	1168	1182	rBV2	326752	822928	5.85%	0.916%
15	10.216	1205	1212	1233	rBV	932218	1768615	12.57%	1.969%
16	10.380	1233	1240	1248	rVB3	111686	206551	1.47%	0.230%
17	10.586	1268	1275	1290	rVB	911355	1593523	11.33%	1.774%
18	10.733	1290	1300	1314	rBV	710533	1422239	10.11%	1.584%
19	11.904	1493	1499	1510	rBV	169612	313288	2.23%	0.349%
20	13.057	1683	1695	1706	rBV	549863	822186	5.84%	0.915%
21	13.845	1819	1829	1833	rBV	2043248	2885962	20.51%	3.213%
22	13.892	1833	1837	1851	rVB	1217397	1740317	12.37%	1.938%
23	14.133	1871	1878	1891	rVB	2322323	3477453	24.72%	3.872%
24	14.439	1921	1930	1938	rBV2	1532276	2362798	16.79%	2.631%
25	14.668	1961	1969	1990	rBV	329216	852588	6.06%	0.949%
26	15.098	2038	2042	2048	rVB3	106339	156203	1.11%	0.174%
27	15.433	2092	2099	2105	rBV	3118808	4350069	30.92%	4.844%
28	15.580	2118	2124	2133	rBV	547256	913540	6.49%	1.017%
29	15.798	2154	2161	2171	rVB	1271911	1754365	12.47%	1.953%
30	17.186	2391	2397	2401	rBV2	2046905	2941098	20.90%	3.275%
31	17.227	2401	2404	2408	rVV	351552	496110	3.53%	0.552%
32	17.286	2408	2414	2430	rVB	3278791	4668515	33.18%	5.198%
33	18.062	2541	2546	2551	rBV	641112	885351	6.29%	0.986%
34	18.115	2551	2555	2564	rVB2	164663	259749	1.85%	0.289%

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35	18.250	2574	2578	2583	rBV2	96183	171545	1.22%	0.191%
36	18.980	2697	2702	2705	rBV2	108741	164387	1.17%	0.183%
37	19.245	2741	2747	2752	rVB	894515	1284783	9.13%	1.431%
38	19.345	2760	2764	2769	rBV2	353010	554135	3.94%	0.617%
39	19.580	2798	2804	2808	rBV	4733305	7013682	49.85%	7.810%
40	20.486	2955	2958	2965	rVB	584208	675750	4.80%	0.752%
41	21.056	3051	3055	3058	rBV2	437356	508882	3.62%	0.567%
42	21.368	3102	3108	3112	rBV2	2925476	4285899	30.46%	4.772%
43	21.403	3112	3114	3121	rVB	592704	754374	5.36%	0.840%
44	21.927	3200	3203	3208	rBV3	301253	539455	3.83%	0.601%
45	22.721	3337	3338	3342	rBV3	224575	342813	2.44%	0.382%
46	23.521	3468	3474	3479	rBV2	1801694	3197711	22.73%	3.561%
47	23.662	3493	3498	3505	rBV	1139629	2088060	14.84%	2.325%

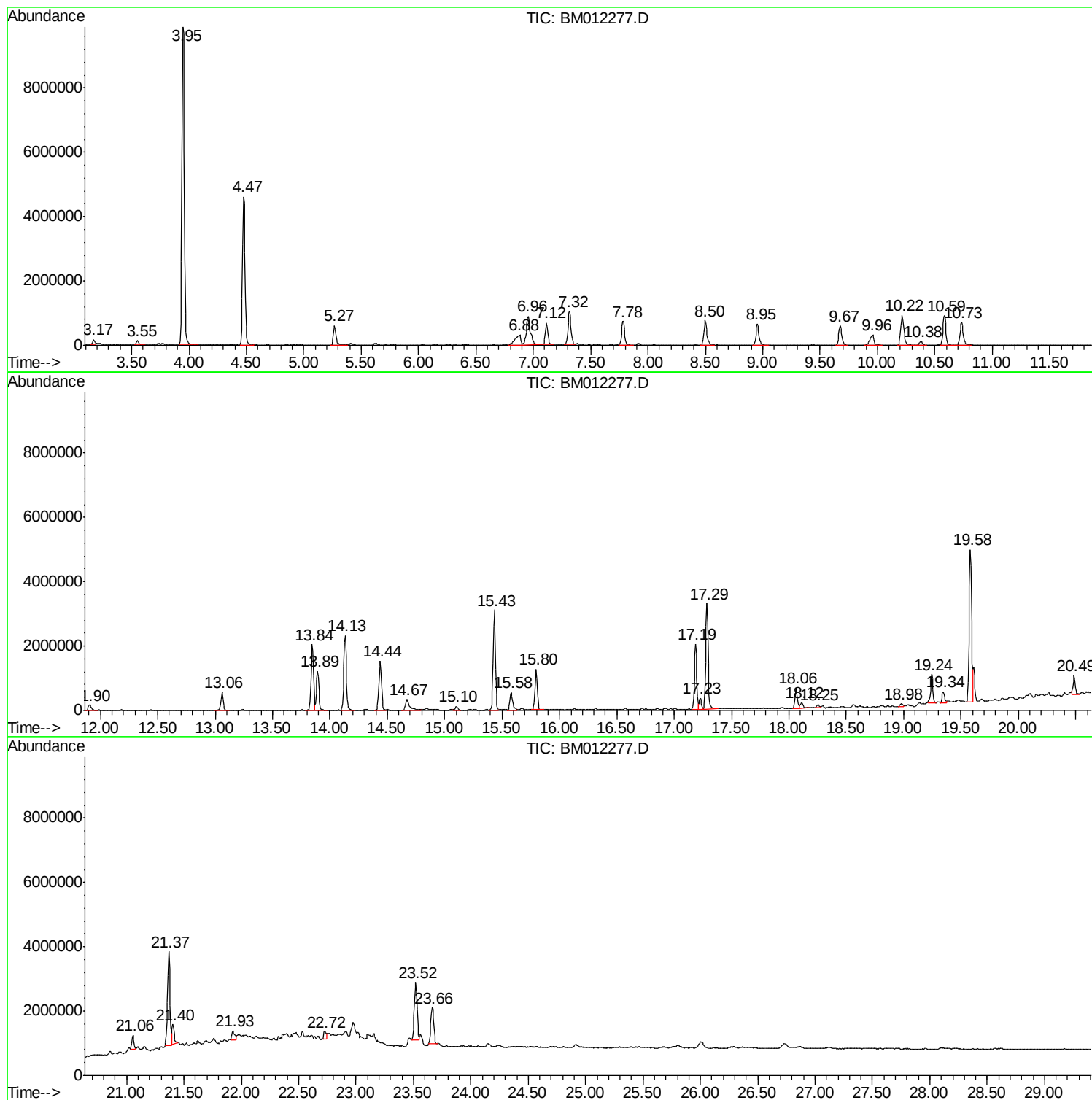
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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM101217.M
Quant Title : SVOA CALIBRATION

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



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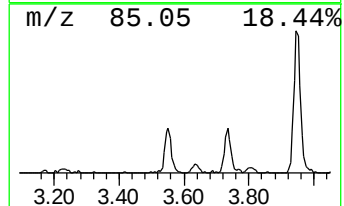
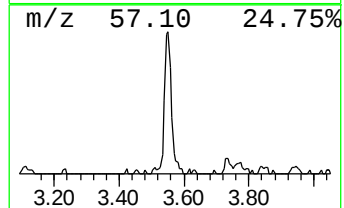
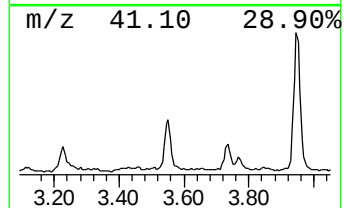
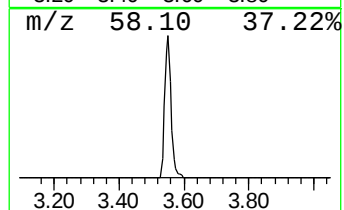
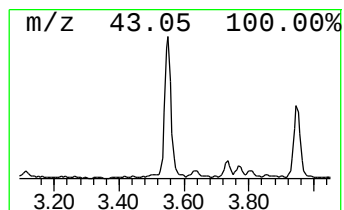
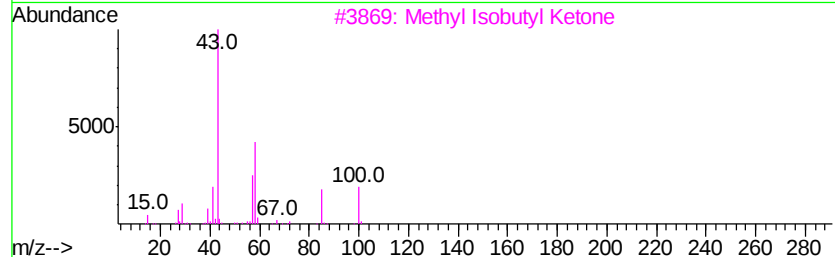
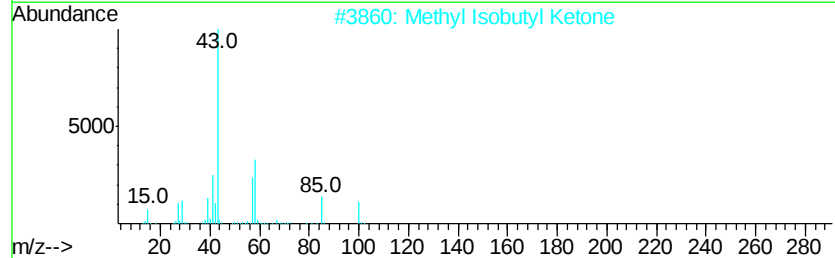
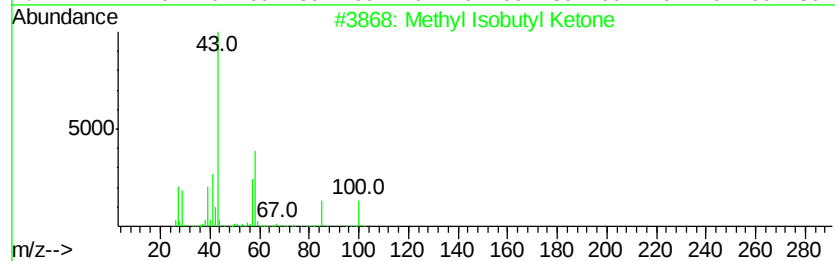
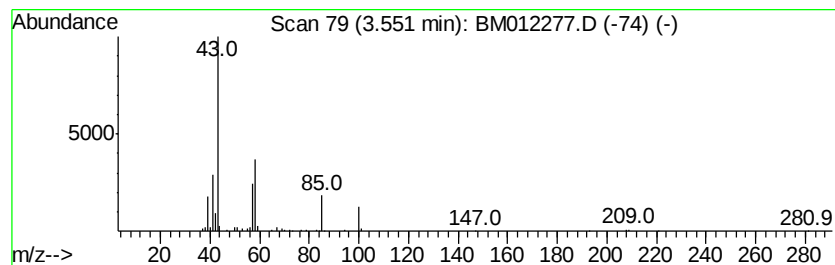
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Quant Title : SVOA CALIBRATION

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

Peak Number 1 Methyl Isobutyl Ketone Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.55	2.99 ng/ul	188840	1,4-Dichlorobenzene-d4	7.78

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl Isobutyl Ketone	100	C6H12O	000108-10-1	87
2			Methyl Isobutyl Ketone	100	C6H12O	000108-10-1	83
3			Methyl Isobutyl Ketone	100	C6H12O	000108-10-1	72
4			2-Hexanone	100	C6H12O	000591-78-6	72
5			2-Hexanone	100	C6H12O	000591-78-6	59



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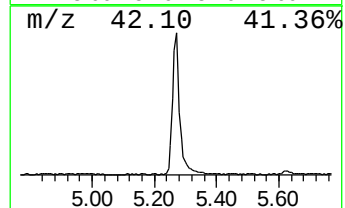
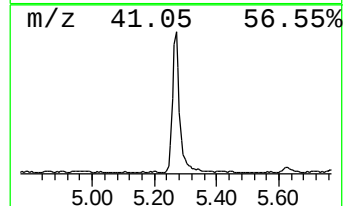
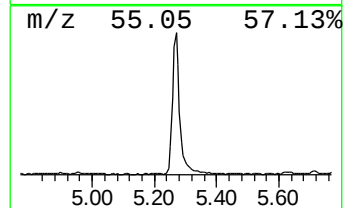
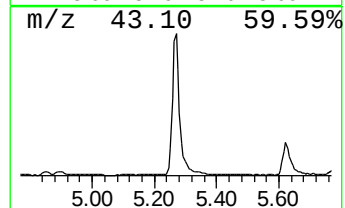
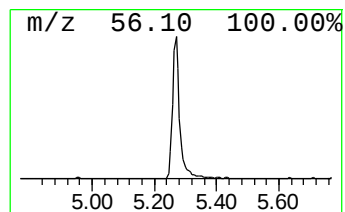
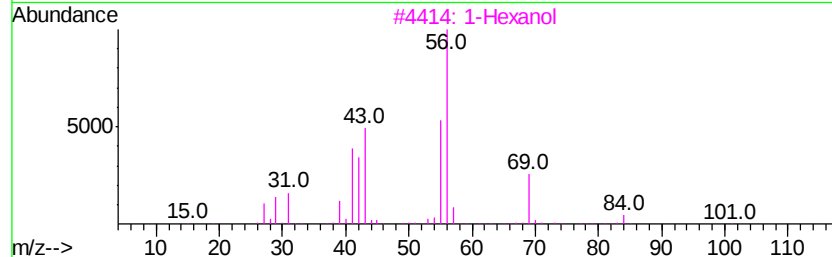
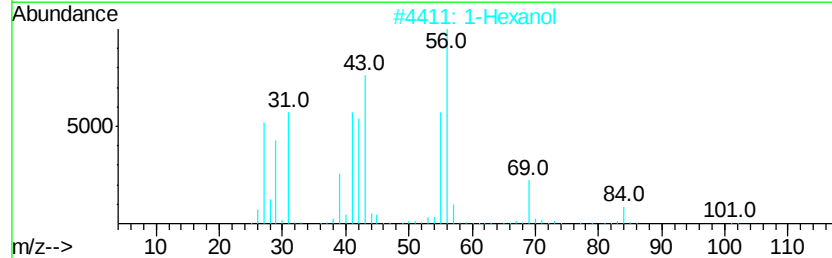
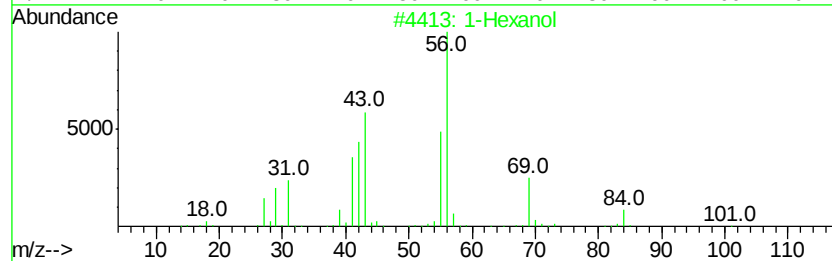
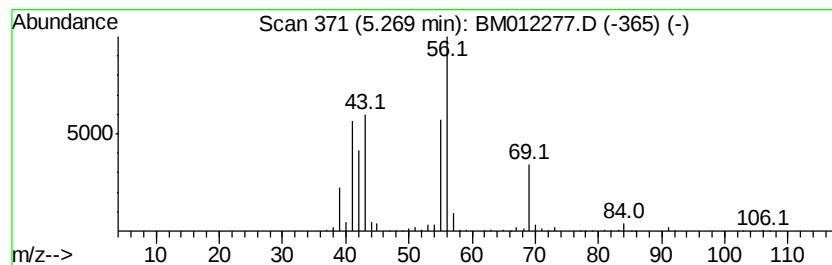
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 4 1-Hexanol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.27	15.97 ng/ul	1007480	1,4-Dichlorobenzene-d4	7.78

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Hexanol	102	C6H14O	000111-27-3	83
2		1-Hexanol	102	C6H14O	000111-27-3	83
3		1-Hexanol	102	C6H14O	000111-27-3	83
4		2H-Pyran-2-one, tetrahydro-5,6-d...	128	C7H12O2	024405-16-1	78
5		Formic acid, hexyl ester	130	C7H14O2	000629-33-4	78



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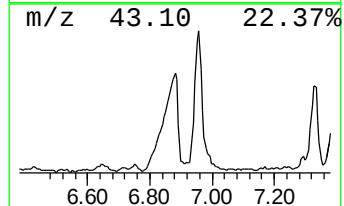
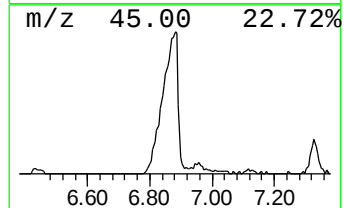
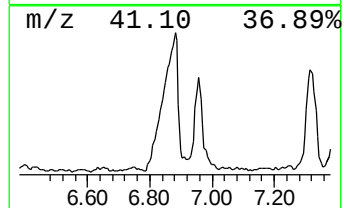
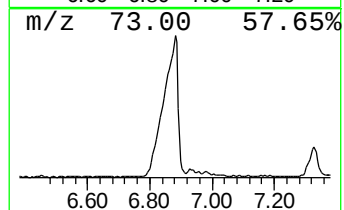
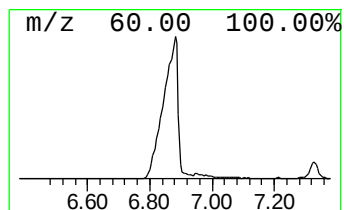
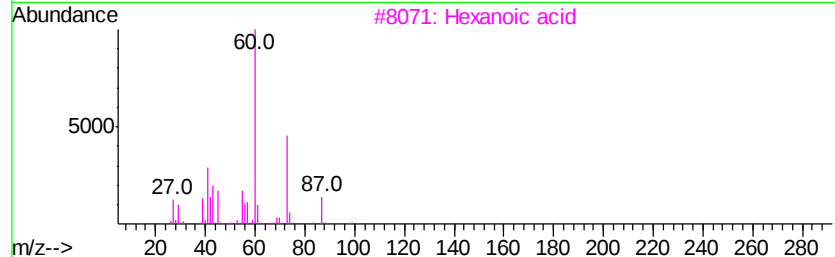
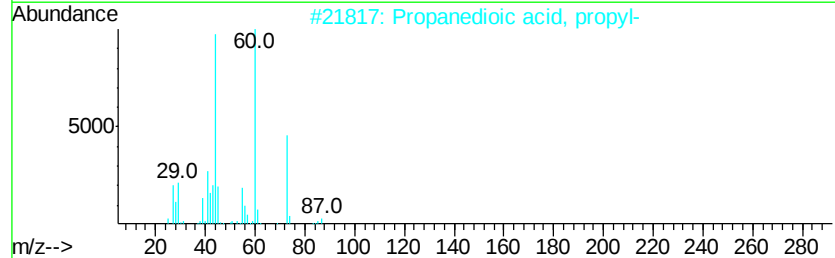
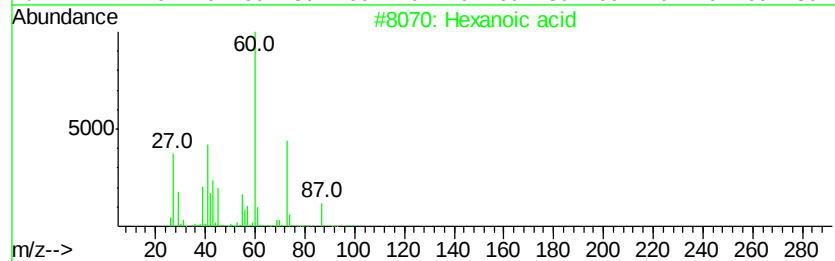
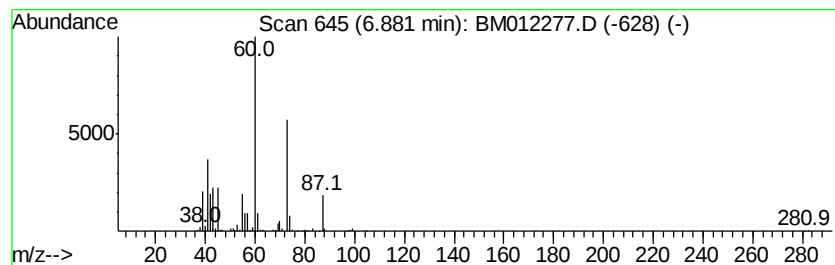
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 5 Hexanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.88	16.69 ng/ul	1052550	1,4-Dichlorobenzene-d4	7.78

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanoic acid		116	C6H12O2	000142-62-1	90
2	Propanedioic acid, propyl-		146	C6H10O4	000616-62-6	78
3	Hexanoic acid		116	C6H12O2	000142-62-1	72
4	Hexanoic acid		116	C6H12O2	000142-62-1	64
5	Hexanoic acid		116	C6H12O2	000142-62-1	64



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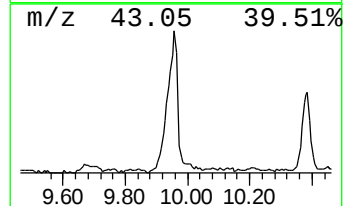
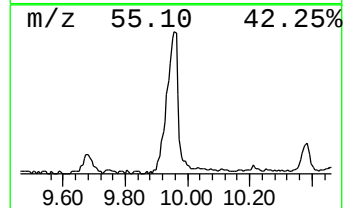
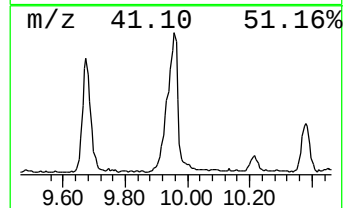
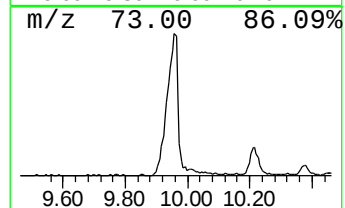
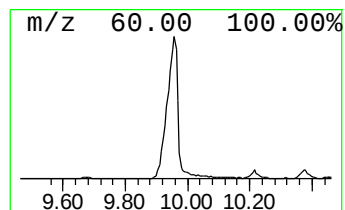
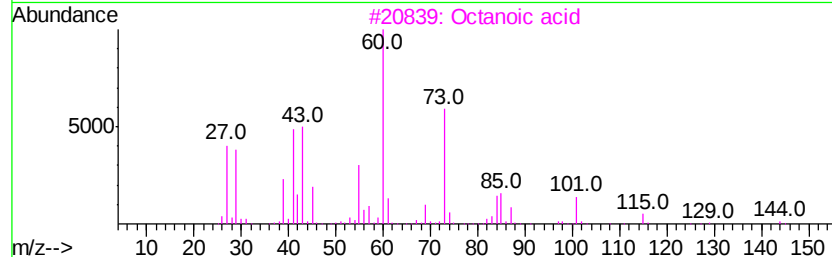
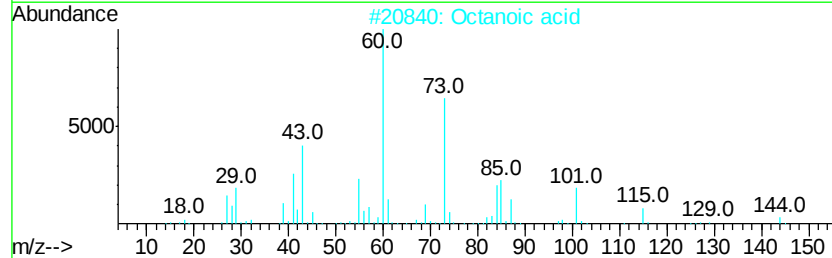
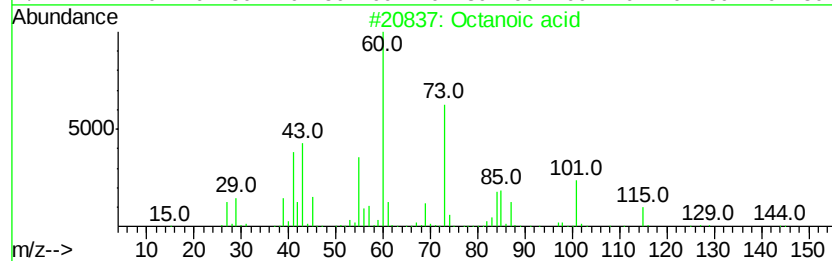
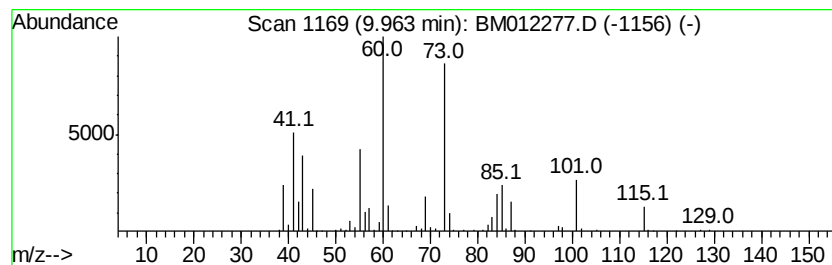
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Octanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.96	10.33 ng/ul	822928	Naphthalene-d8	10.59

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octanoic acid	144	C8H16O2	000124-07-2	91
2		Octanoic acid	144	C8H16O2	000124-07-2	86
3		Octanoic acid	144	C8H16O2	000124-07-2	83
4		Nonanoic acid	158	C9H18O2	000112-05-0	53
5		Hexanoic acid	116	C6H12O2	000142-62-1	50



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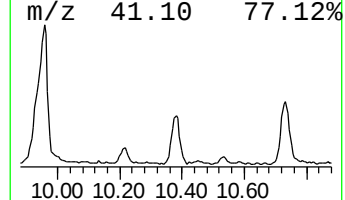
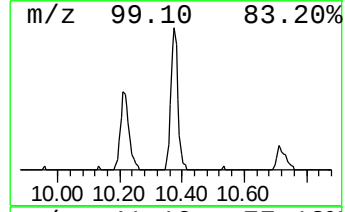
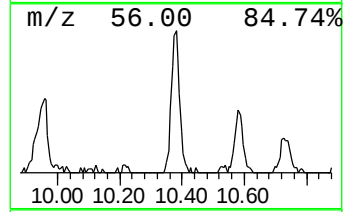
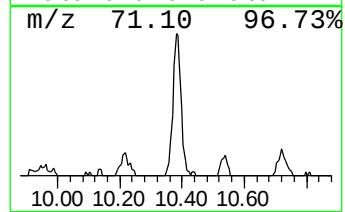
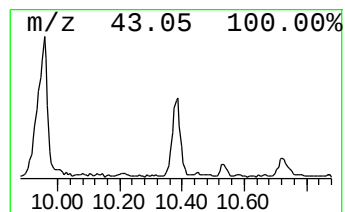
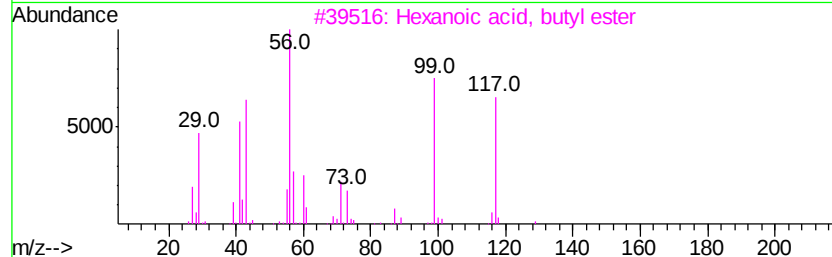
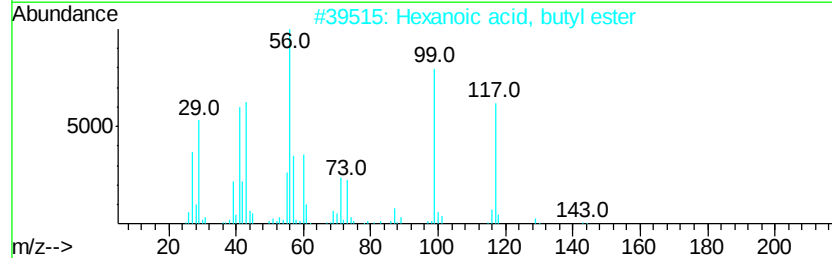
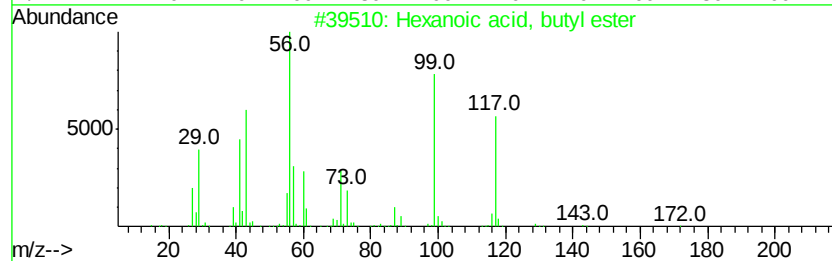
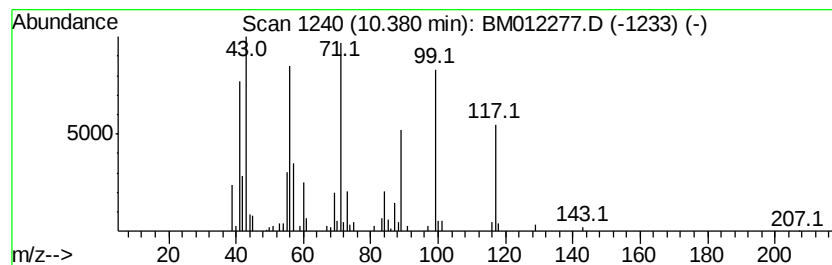
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Hexanoic acid, butyl ester Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.		
10.38	2.59 ng/ul	206551	Naphthalene-d8	10.59		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanoic acid, butyl ester	172	C10H20O2	000626-82-4	50	
2	Hexanoic acid, butyl ester	172	C10H20O2	000626-82-4	50	
3	Hexanoic acid, butyl ester	172	C10H20O2	000626-82-4	47	
4	Hexanoic acid, hexyl ester	200	C12H24O2	006378-65-0	47	
5	4H-Imidazol-4-one, 2-amino-1,5-d...	99	C3H5N3O	000503-86-6	43	



Data Path : Z:\HPCHEM1\BNA M\DATA\BM101717\
Data File : BM012277.D
Acq On : 18 Oct 2017 08:20
Operator : SJ/JU
Sample : I5747-02
Misc :
ALS Vial : 28 Sample Multiplier: 1

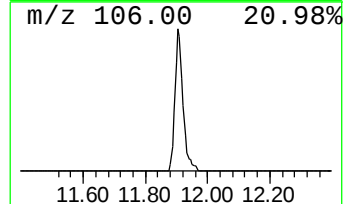
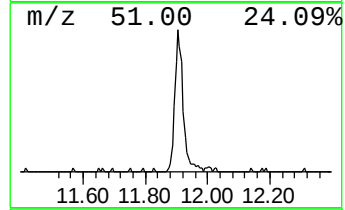
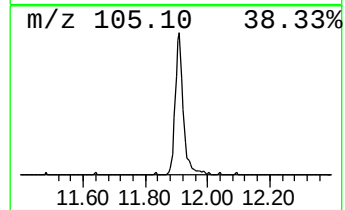
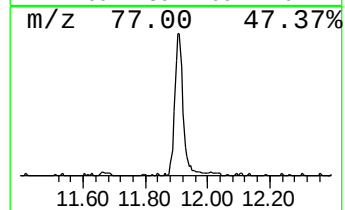
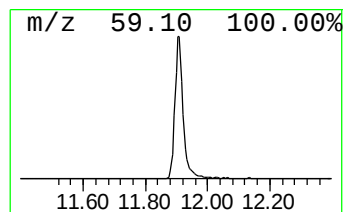
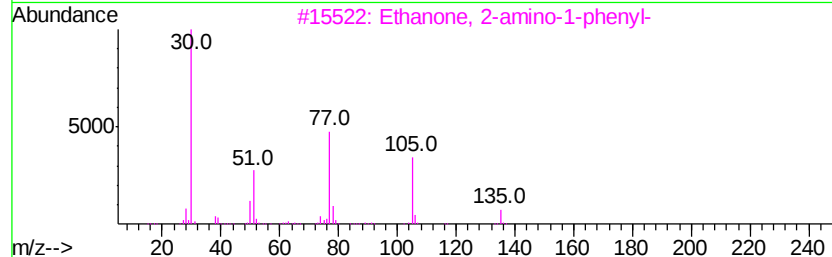
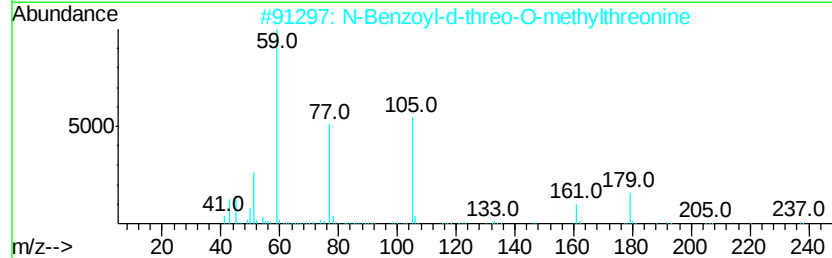
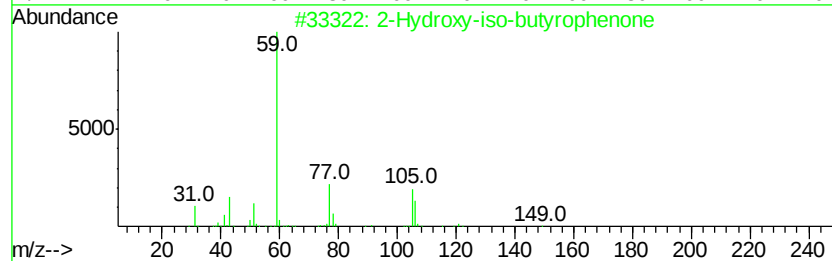
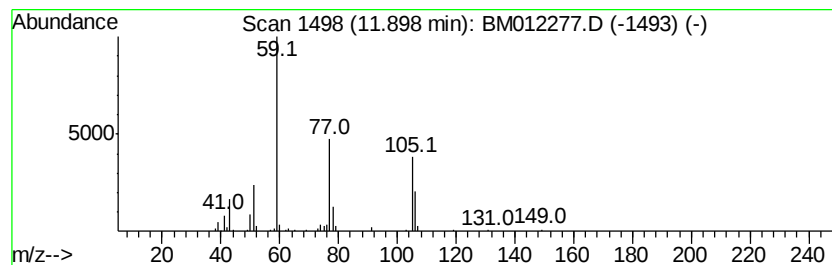
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM101217.M
Quant Title : SVOA CALIBRATION

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

Peak Number 8 2-Hydroxy-iso-butyrophenone Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.		
11.90	3.93 ng/ul	313288	Napthalene-d8	10.59		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hydroxy-iso-butvropfenone	164	C10H12O2	007473-98-5	64	
2	N-Benzoyl-d-threo-O-methylthreonine	237	C12H15NO4	131644-54-7	39	
3	Ethanone, 2-amino-1-phenyl-	135	C8H9NO	000613-89-8	35	
4	Propanedioic acid, (benzoylhvdr...	282	C12H14N2O6	1000142-99-9	25	
5	Guanidine	59	CH5N3	000113-00-8	9	



Data Path : Z:\HPCHEM1\BNA M\DATA\BM101717\
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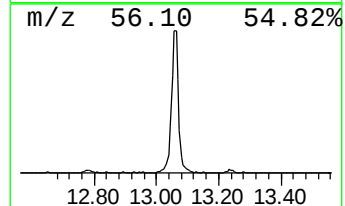
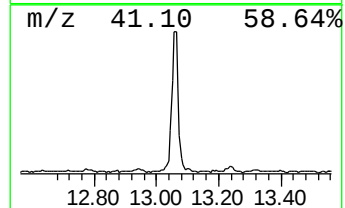
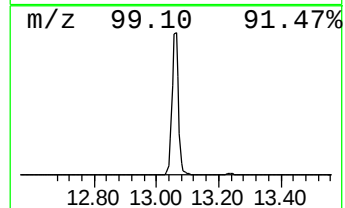
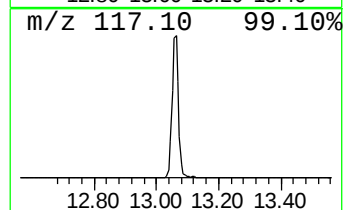
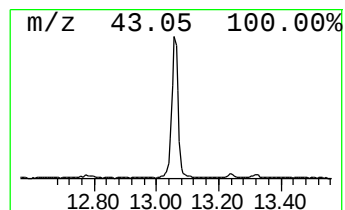
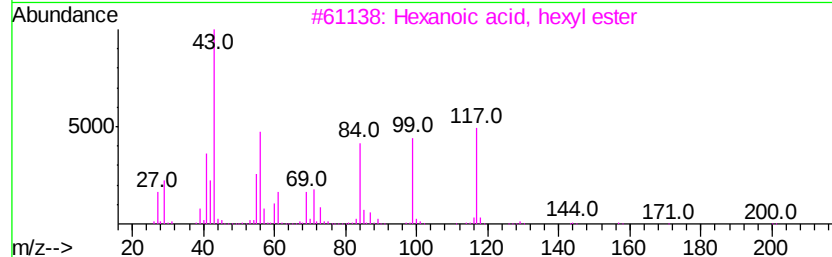
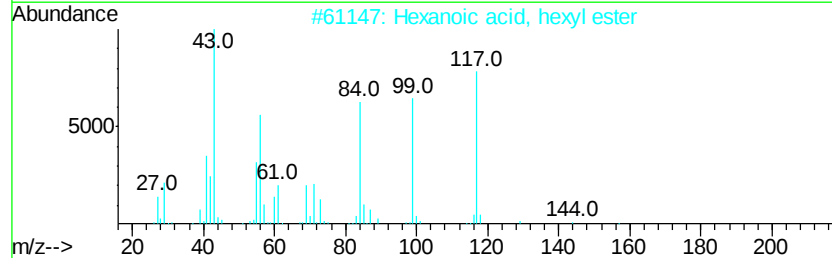
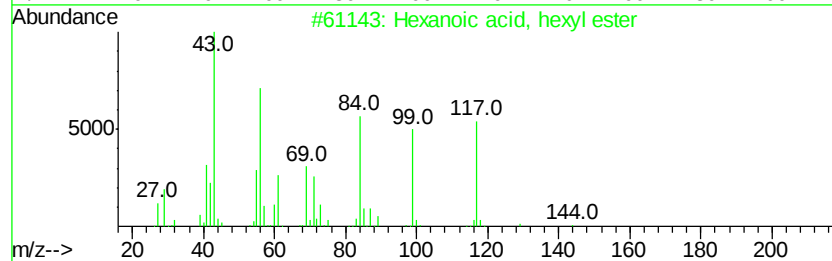
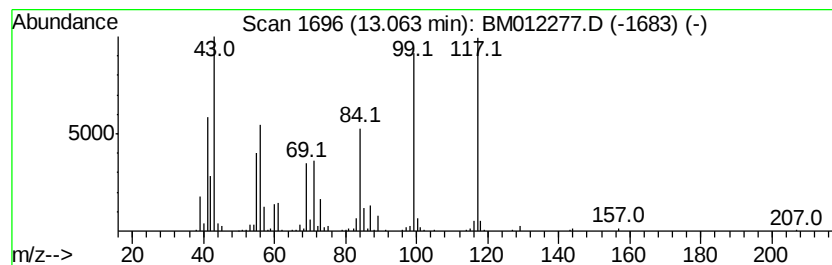
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Hexanoic acid, hexyl ester Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.06	6.96 ng/ul	822186	Acenaphthene-d10	14.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexanoic acid, hexyl ester	200	C12H24O2	006378-65-0	72
2		Hexanoic acid, hexyl ester	200	C12H24O2	006378-65-0	58
3		Hexanoic acid, hexyl ester	200	C12H24O2	006378-65-0	58
4		Hexanoic acid, pentyl ester	186	C11H22O2	000540-07-8	50
5		Hexanoic acid, butyl ester	172	C10H20O2	000626-82-4	50



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Misc :
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ClientSampleId :
B-23(1-3)-100517

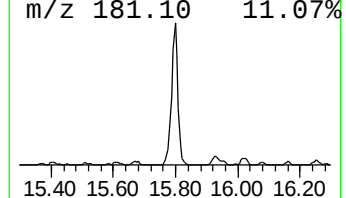
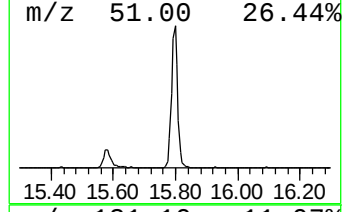
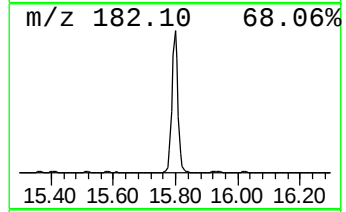
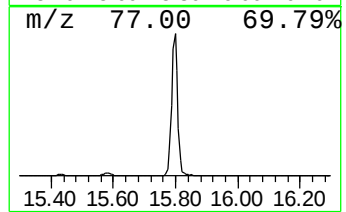
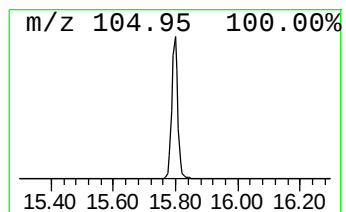
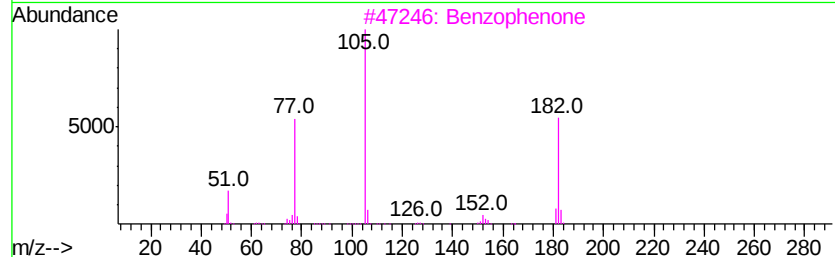
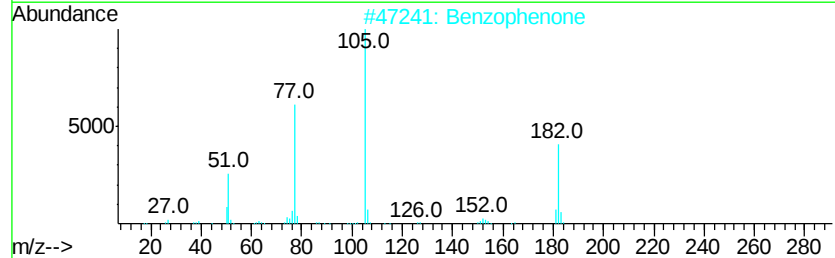
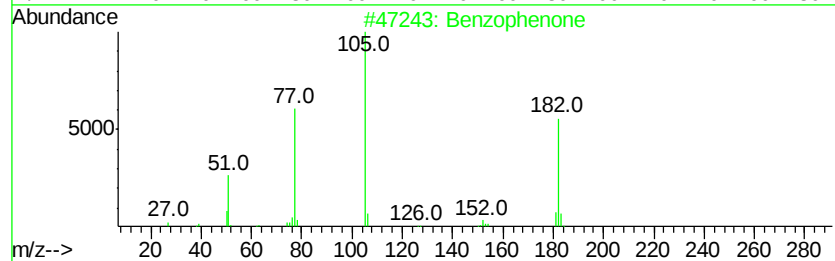
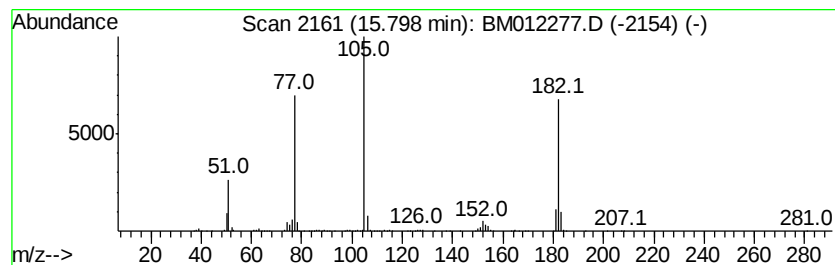
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Benzophenone Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.80	14.85 ng/ul	1754370	Acenaphthene-d10	14.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzophenone	182	C13H10O	000119-61-9	96
2		Benzophenone	182	C13H10O	000119-61-9	95
3		Benzophenone	182	C13H10O	000119-61-9	95
4		Benzophenone	182	C13H10O	000119-61-9	95
5		Benzophenone	182	C13H10O	000119-61-9	91



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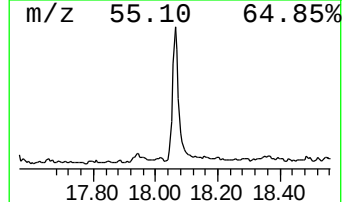
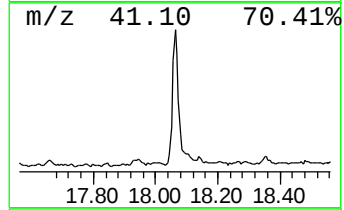
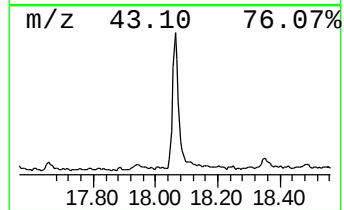
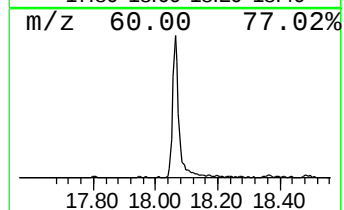
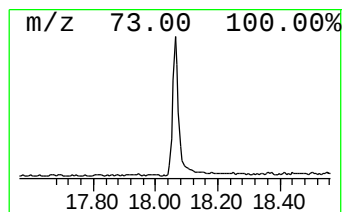
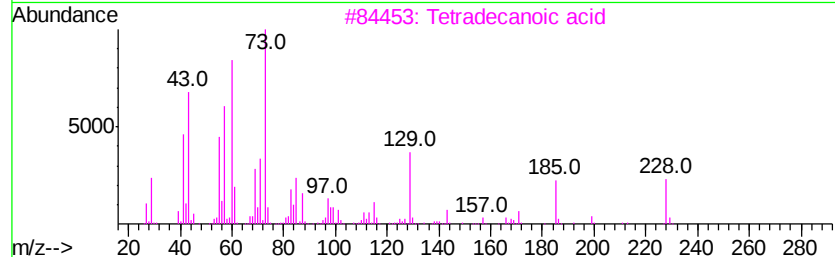
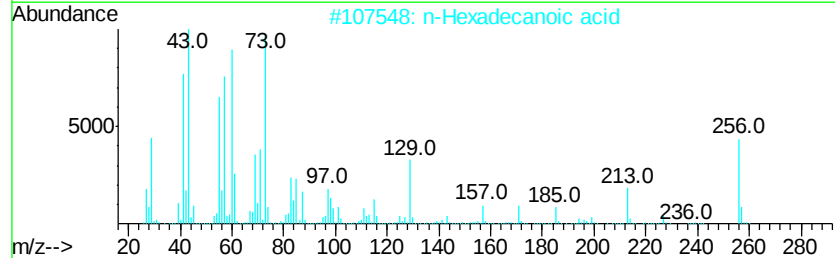
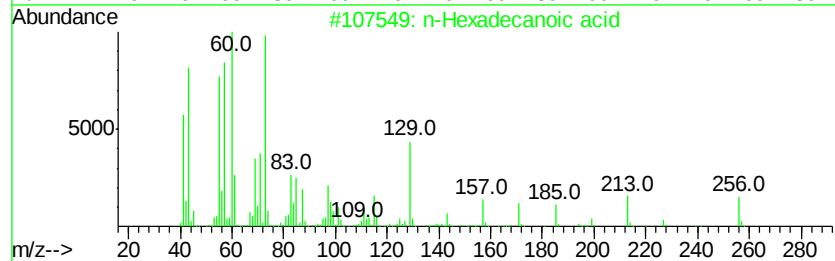
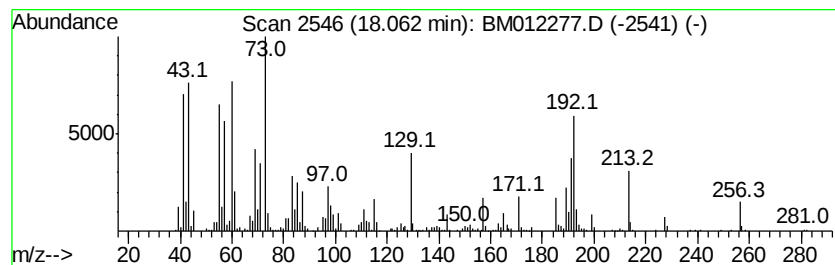
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 n-Hexadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD		R.T.	
18.06	6.02 ng/ul	885351	Phenanthrene-d10		17.19	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99	
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97	
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	91	
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	90	
5	Tridecanoic acid	214	C13H26O2	000638-53-9	78	



Data Path : Z:\HPCHEM1\BNA M\DATA\BM101717\
Data File : BM012277.D
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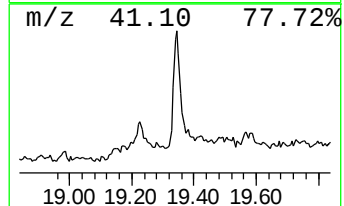
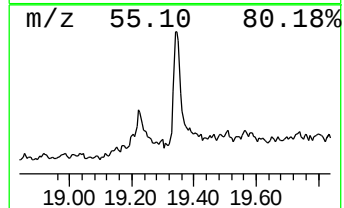
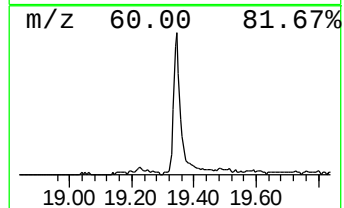
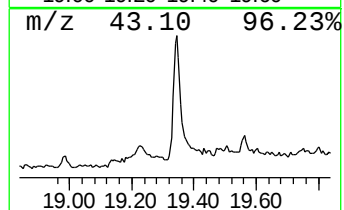
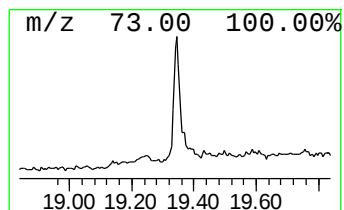
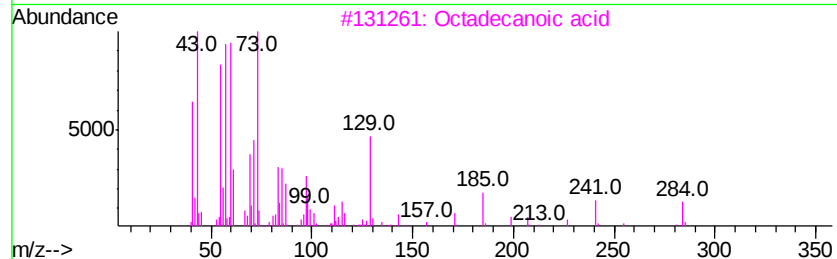
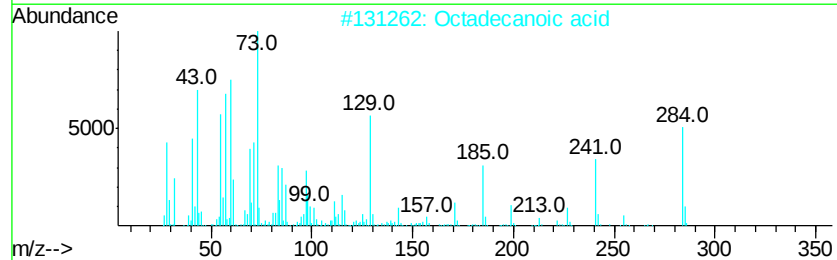
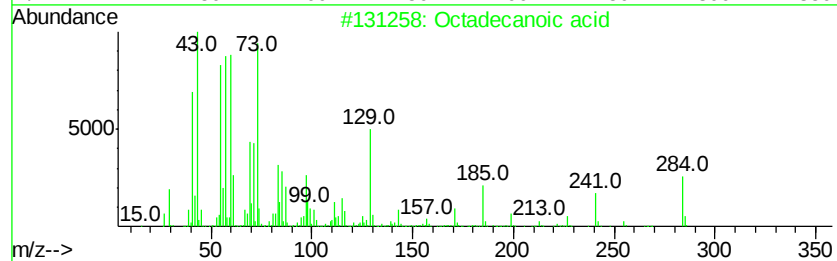
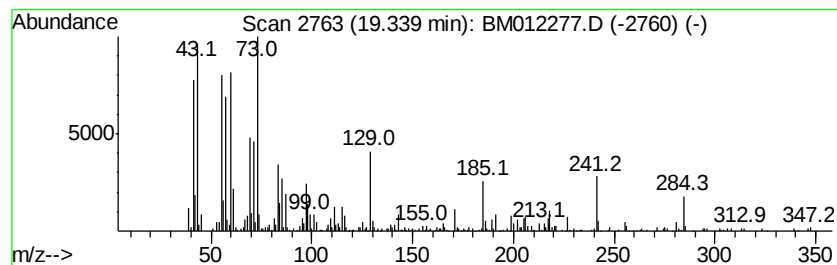
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Octadecanoic acid Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.34	2.59 ng/ul	554135	Chrysene-d12	21.37

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	96
3			Octadecanoic acid	284	C18H36O2	000057-11-4	96
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94
5			Octadecanoic acid	284	C18H36O2	000057-11-4	91



Data Path : Z:\HPCHEM1\BNA M\DATA\BM101717\
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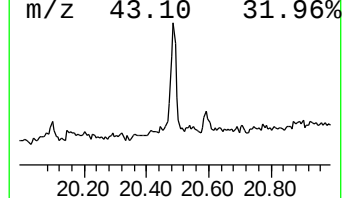
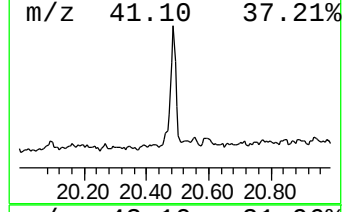
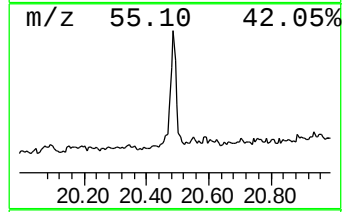
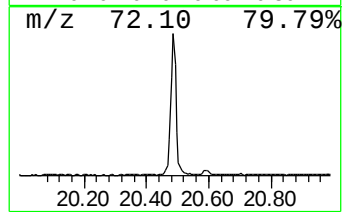
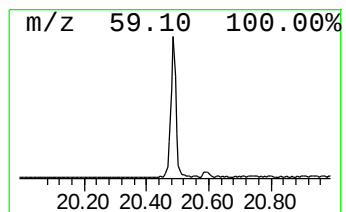
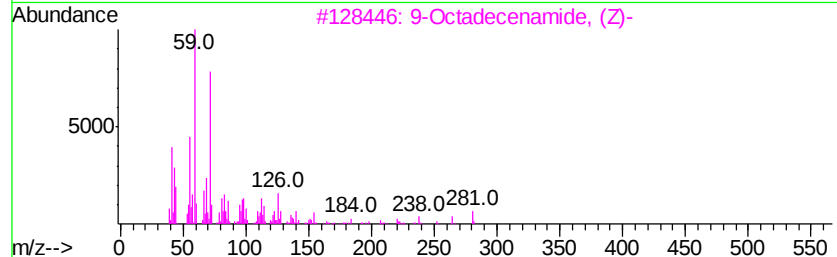
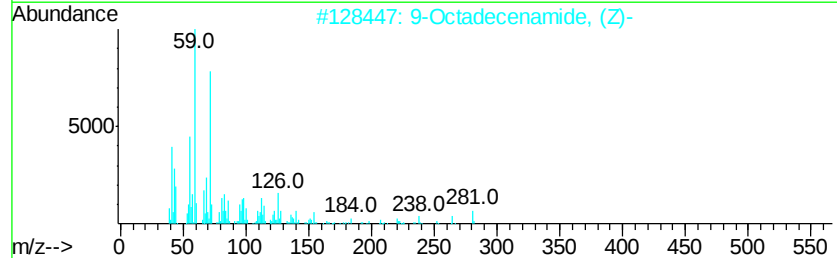
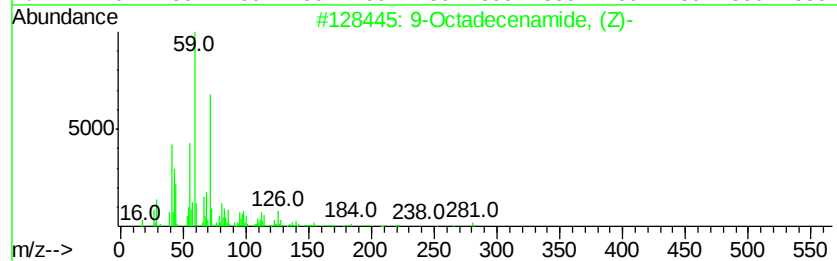
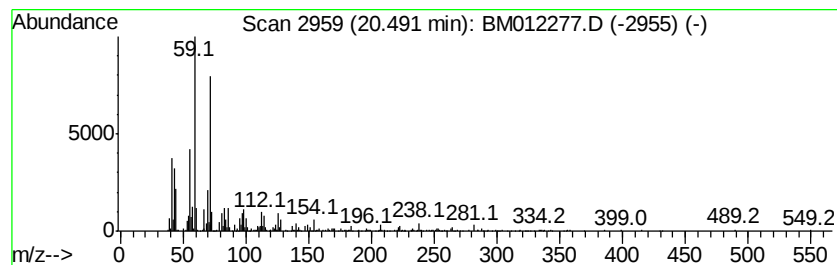
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 9-Octadecenamide, (Z)- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.49	3.15 ng/ul	675750	Chrysene-d12	21.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	99
2		9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	97
3		9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	97
4		Tetradecanamide	227	C14H29NO	000638-58-4	72
5		7-Nonenamide	155	C9H17NO	090949-53-4	59



Data Path : Z:\HPCHEM1\BNA M\DATA\BM101717\
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ClientSampleId :
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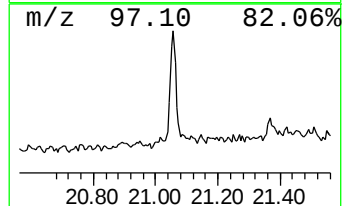
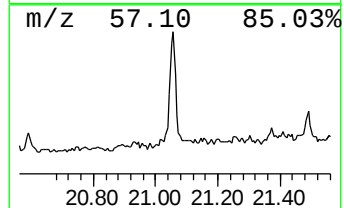
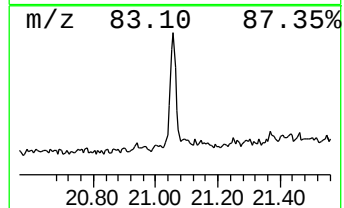
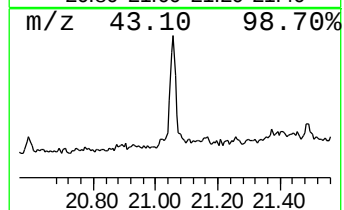
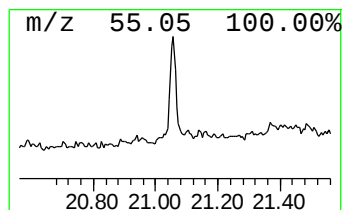
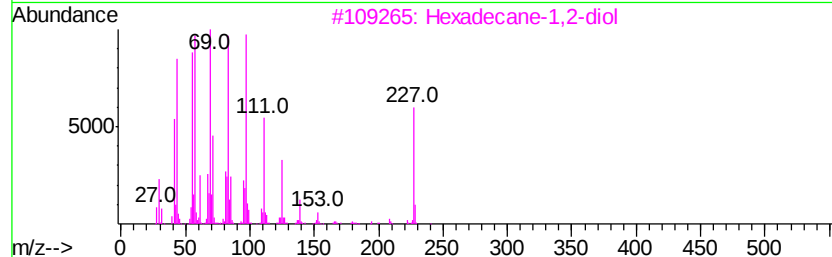
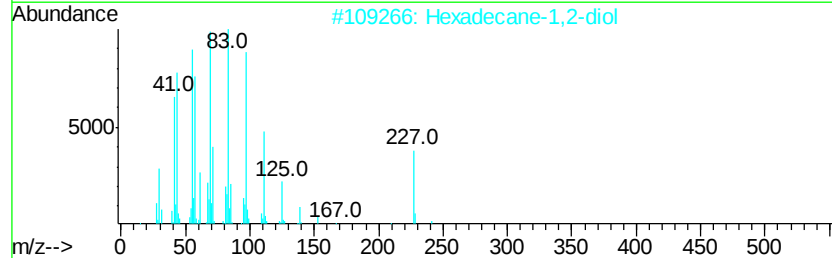
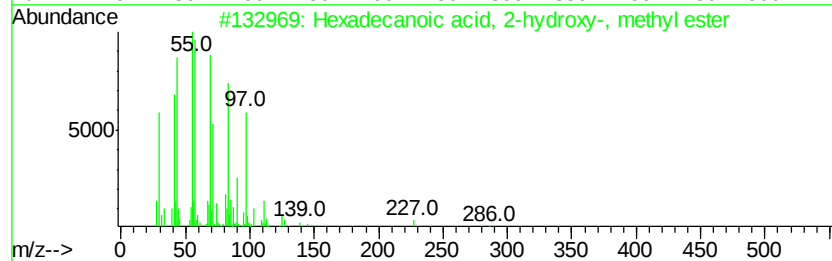
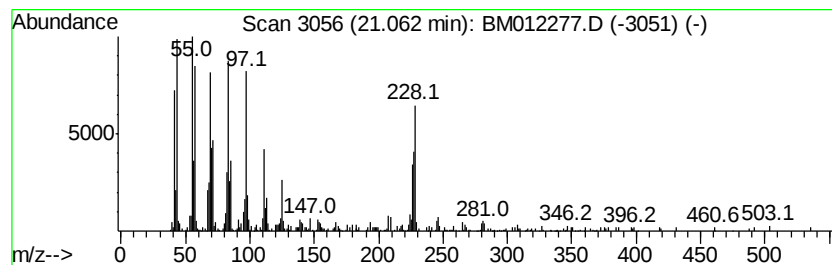
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 Hexadecanoic acid, 2-hydrox... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.06	2.37 ng/ul	508882	Chrysene-d12	21.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecanoic acid, 2-hydroxy-, m...	286	C17H34O3	016742-51-1	50
2		Hexadecane-1,2-diol	258	C16H34O2	006920-24-7	49
3		Hexadecane-1,2-diol	258	C16H34O2	006920-24-7	38
4		Trifluoroacetic acid,n-tridecyl ...	296	C15H27F3O2	053800-02-5	30
5		erythro-9,10-Dibromopentacosane	508	C25H50Br2	059907-02-7	25



Data Path : Z:\HPCHEM1\BNA M\DATA\BM101717\
Data File : BM012277.D
Acq On : 18 Oct 2017 08:20
Operator : SJ/JU
Sample : I5747-02
Misc :
ALS Vial : 28 Sample Multiplier: 1

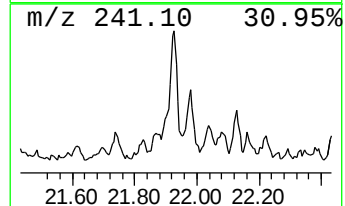
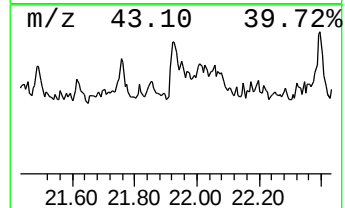
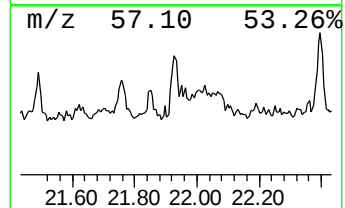
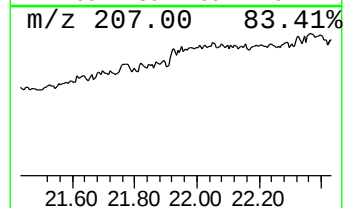
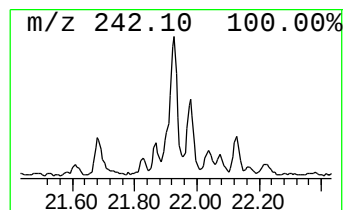
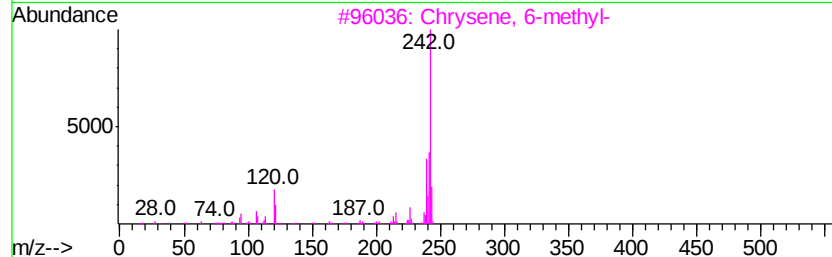
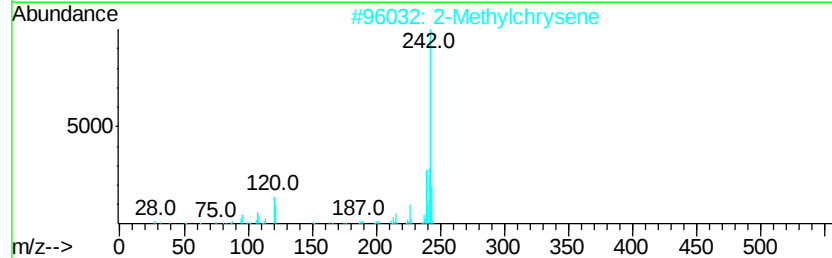
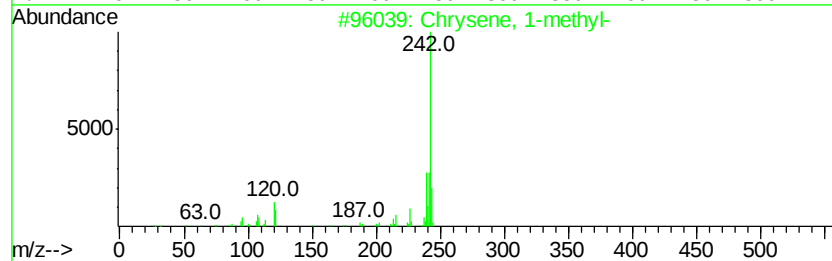
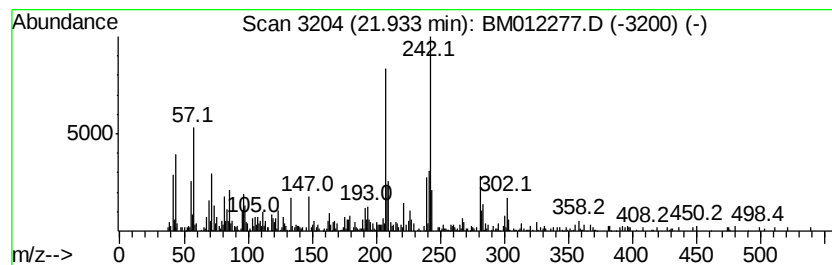
Instrument :
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ClientSampleId :
B-23(1-3)-100517

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 15 Chrysene, 1-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD		R.T.		
21.93	2.52 ng/ul	539455	Chrysene-d12		21.37		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Chrysene, 1-methyl-	242	C19H14	003351-28-8	84
2			2-Methylchrysene	242	C19H14	003351-32-4	83
3			Chrysene, 6-methyl-	242	C19H14	001705-85-7	59
4			Benz[<i>a</i>]anthracene, 7-methyl-	242	C19H14	002541-69-7	46
5			Benz[<i>a</i>]anthracene, 7-methyl-	242	C19H14	002541-69-7	46



Data Path : Z:\HPCHEM1\BNA M\DATA\BM101717\
Data File : BM012277.D
Acq On : 18 Oct 2017 08:20
Operator : SJ/JU
Sample : I5747-02
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
B-23(1-3)-100517

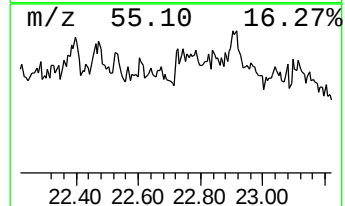
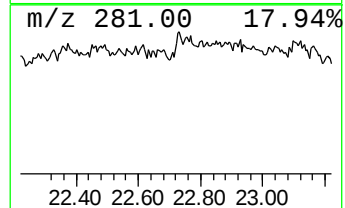
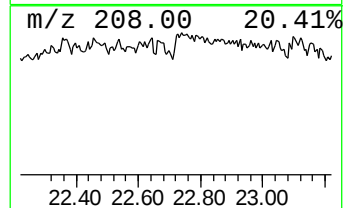
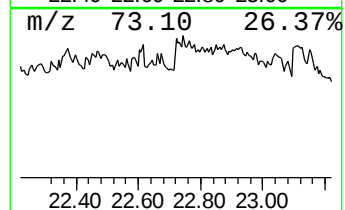
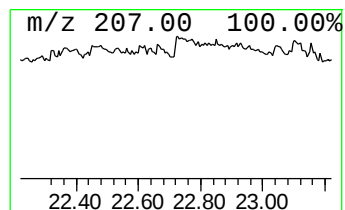
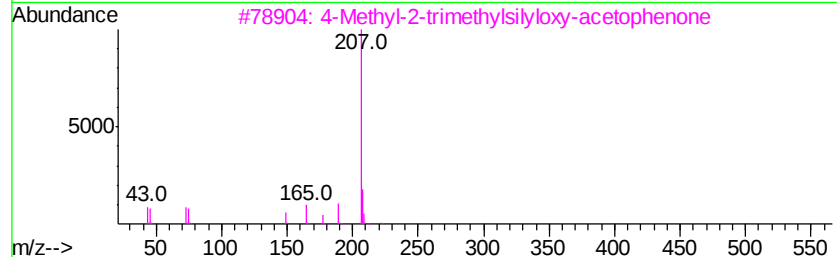
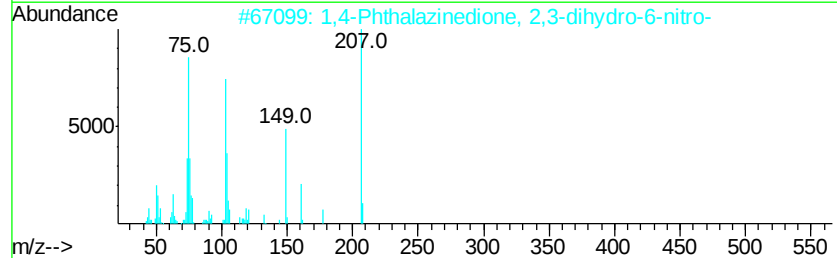
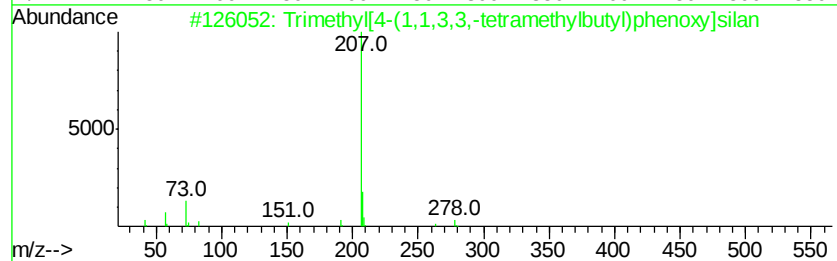
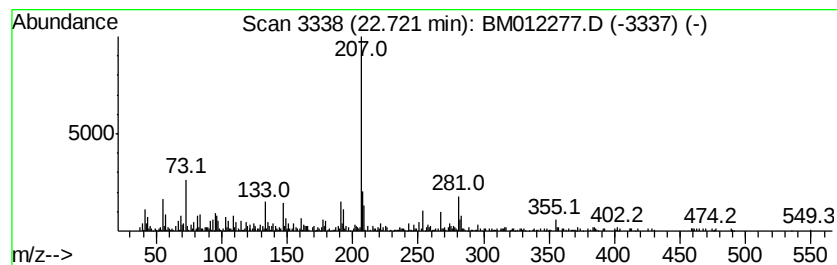
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM101217.M
Quant Title : SVOA CALIBRATION

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

Peak Number 16 Trimethyl[4-(1,1,3,3,-tetra... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.72	3.28 ng/ul	342813	Perylene-d12	23.66

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Trimethyl[4-(1,1,3,3,-tetramethy...	278	C17H30OSi	078721-87-6	72
2			1,4-Phthalazinedione, 2,3-dihydr...	207	C8H5N3O4	003682-19-7	60
3			4-Methyl-2-trimethylsilyloxy-ace...	222	C12H18O2Si	097389-70-3	53
4			1,2-Benzisothiazol-3-amine tbdms	264	C13H20N2SSi	1000332-57-2	53
5			1H-Benzo[4,5]furo[3,2-f]indole	207	C14H9NO	000242-97-7	52



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM101717\
 Data File : BM012277.D
 Acq On : 18 Oct 2017 08:20
 Operator : SJ/JU
 Sample : I5747-02
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 B-23(1-3)-100517

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
Methyl Isobutyl K...	3.55	3.0	ng/ul	188840	1	7.78	1261560	20.0
1-Hexanol	5.27	16.0	ng/ul	1007480	1	7.78	1261560	20.0
Hexanoic acid	6.88	16.7	ng/ul	1052550	1	7.78	1261560	20.0
Octanoic acid	9.96	10.3	ng/ul	822928	2	10.59	1593520	20.0
Hexanoic acid, bu...	10.38	2.6	ng/ul	206551	2	10.59	1593520	20.0
2-Hydroxy-iso-but...	11.90	3.9	ng/ul	313288	2	10.59	1593520	20.0
Hexanoic acid, he...	13.06	7.0	ng/ul	822186	3	14.44	2362800	20.0
Benzophenone	15.80	14.9	ng/ul	1754370	3	14.44	2362800	20.0
n-Hexadecanoic acid	18.06	6.0	ng/ul	885351	4	17.19	2941100	20.0
Octadecanoic acid	19.34	2.6	ng/ul	554135	5	21.37	4285900	20.0
9-Octadecenamide,...	20.49	3.1	ng/ul	675750	5	21.37	4285900	20.0
Hexadecanoic acid...	21.06	2.4	ng/ul	508882	5	21.37	4285900	20.0
Chrysene, 1-methyl-	21.93	2.5	ng/ul	539455	5	21.37	4285900	20.0
Trimethyl[4-(1,1,...	22.72	3.3	ng/ul	342813	6	23.66	2088060	20.0