

Data Path : Z:\HPCHEM1\BNA M\DATA\BM102717\
 Data File : BM012507.D
 Acq On : 28 Oct 2017 04:17
 Operator : SJ/JU
 Sample : I5786-08
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 B-7(0.5-1.5)_101017

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-BM102417.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.093	165	171	182	rVB	921688	1361379	1.68%	0.362%
2	5.528	409	415	425	rBV	2059463	4396108	5.42%	1.168%
3	5.893	472	477	487	rVB	1266189	1957392	2.41%	0.520%
4	6.869	632	643	649	rBV5	394440	1031947	1.27%	0.274%
5	7.034	664	671	679	rVV2	2953315	6277291	7.74%	1.668%
6	7.198	694	699	705	rVB	2006641	3080751	3.80%	0.819%
7	7.404	723	734	745	rVB	2636651	4609778	5.68%	1.225%
8	7.498	745	750	752	rBV	1262121	1882453	2.32%	0.500%
9	7.798	790	801	804	rBV7	315474	873530	1.08%	0.232%
10	7.869	804	813	818	rVB2	2747254	5266241	6.49%	1.399%
11	8.422	900	907	917	rVB6	616321	1630103	2.01%	0.433%
12	8.569	927	932	939	rVB	2695748	4362809	5.38%	1.159%
13	8.975	990	1001	1004	rBV3	613265	1397387	1.72%	0.371%
14	9.022	1004	1009	1015	rVV	2283034	3771582	4.65%	1.002%
15	9.098	1017	1022	1026	rVB	844754	1205864	1.49%	0.320%
16	9.745	1124	1132	1138	rBV	2081928	3597361	4.43%	0.956%
17	10.004	1169	1176	1182	rVB2	551171	1102476	1.36%	0.293%
18	10.280	1213	1223	1232	rVB	3168926	5426151	6.69%	1.442%
19	10.663	1278	1288	1293	rBV	3454019	6149147	7.58%	1.634%
20	10.792	1304	1310	1321	rBV	1104031	2086986	2.57%	0.555%
21	11.463	1414	1424	1437	rBV	462542	965847	1.19%	0.257%
22	13.910	1831	1840	1844	rBV	4986419	7117083	8.77%	1.891%
23	14.192	1882	1888	1900	rVB	5844752	8423774	10.38%	2.239%
24	14.498	1932	1940	1948	rBV2	4744296	7224777	8.90%	1.920%
25	14.692	1967	1973	1986	rBV	2047987	3252758	4.01%	0.864%
26	15.492	2099	2109	2115	rBV	7447746	10237580	12.62%	2.721%
27	15.610	2123	2129	2135	rVB	1935295	2784631	3.43%	0.740%
28	15.851	2166	2170	2176	rVB	1501127	2079825	2.56%	0.553%
29	17.239	2400	2406	2411	rBV2	5637381	7847407	9.67%	2.085%
30	17.339	2417	2423	2432	rVB	7296021	10328027	12.73%	2.745%
31	18.162	2554	2563	2567	rBV	25208776	43065026	53.07%	11.444%
32	18.209	2567	2571	2576	rVB2	27585781	37797215	46.58%	10.045%
33	18.921	2685	2692	2700	rVB	1477306	3002070	3.70%	0.798%
34	19.039	2708	2712	2714	rBV2	746043	1048770	1.29%	0.279%

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35	19.068	2714	2717	2724	rVB3	858595	1498876	1.85%	0.398%
36	19.327	2751	2761	2770	rBV2	30629415	81151525	100.00%	21.566%
37	19.439	2775	2780	2786	rVV	16336620	21893234	26.98%	5.818%
38	19.627	2807	2812	2823	rVB2	8686614	13161620	16.22%	3.498%
39	20.550	2963	2969	2974	rVB	6131733	7839994	9.66%	2.083%
40	21.121	3064	3066	3070	rVB	2480811	2589271	3.19%	0.688%
41	21.327	3097	3101	3108	rVB	12346751	14553854	17.93%	3.868%
42	21.409	3111	3115	3120	rBV	6251802	8536849	10.52%	2.269%
43	23.574	3479	3483	3494	rVB2	5466287	9957466	12.27%	2.646%
44	23.721	3504	3508	3514	rVB	4517086	8473035	10.44%	2.252%

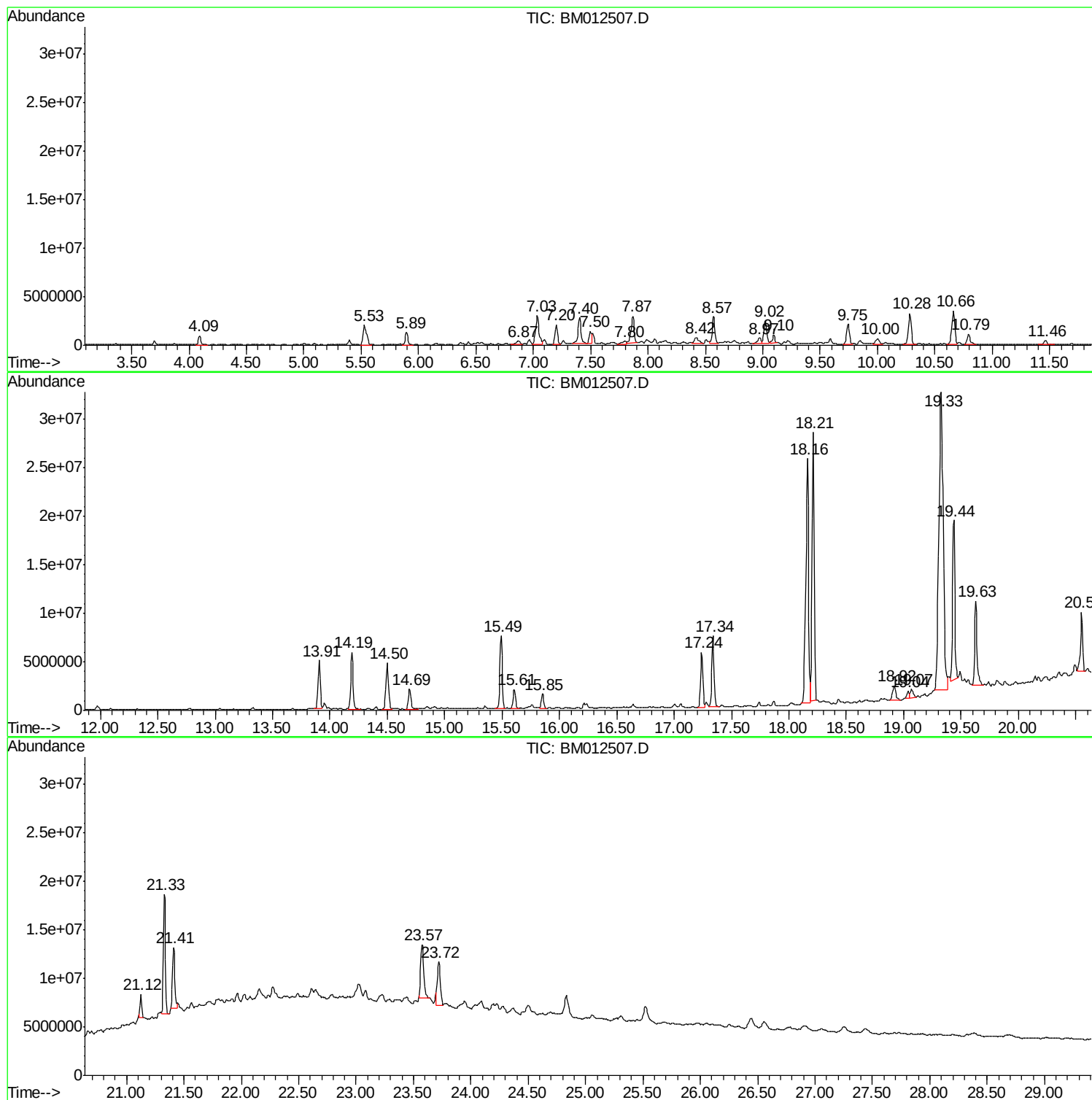
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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM102417.M
Quant Title : SVOA CALIBRATION

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



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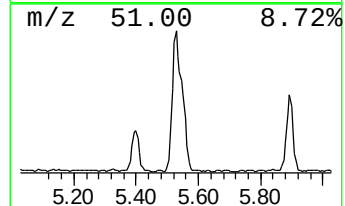
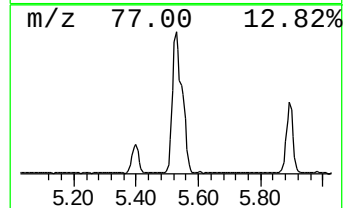
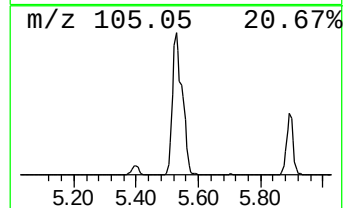
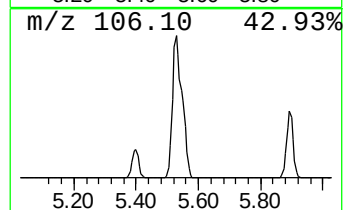
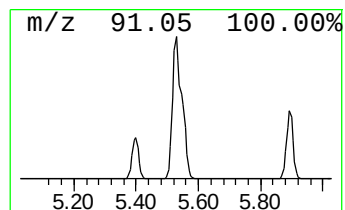
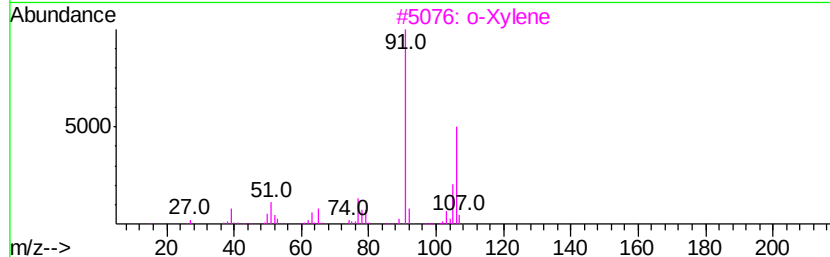
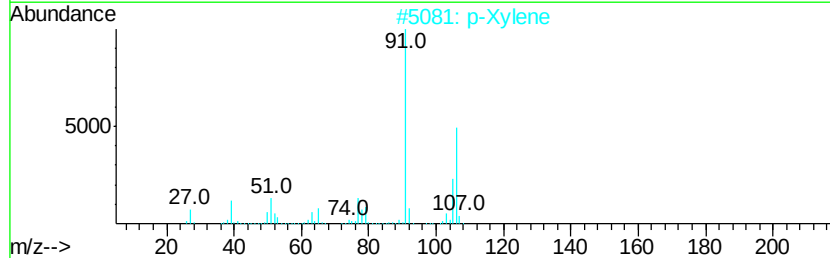
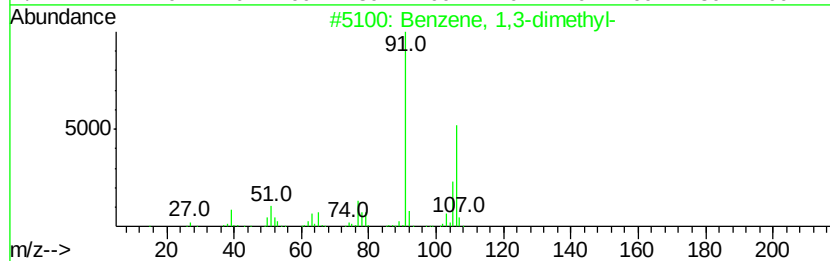
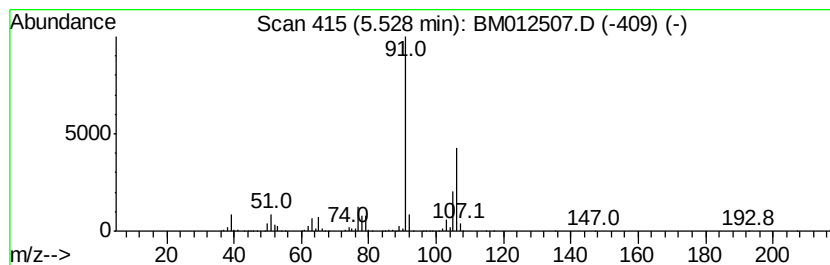
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzene, 1,3-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.53	16.70 ng/ul	4396110	1,4-Dichlorobenzene-d4	7.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
2		p-Xylene	106	C8H10	000106-42-3	97
3		o-Xylene	106	C8H10	000095-47-6	97
4		Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
5		p-Xylene	106	C8H10	000106-42-3	95



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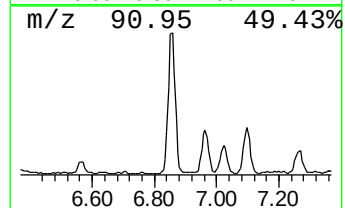
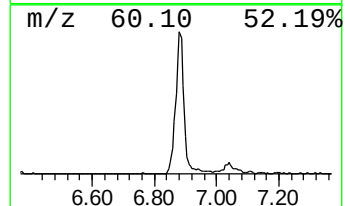
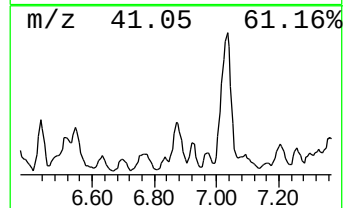
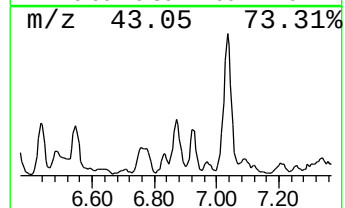
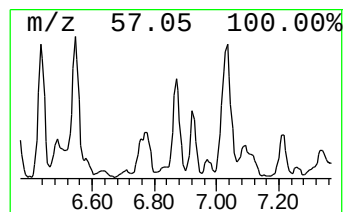
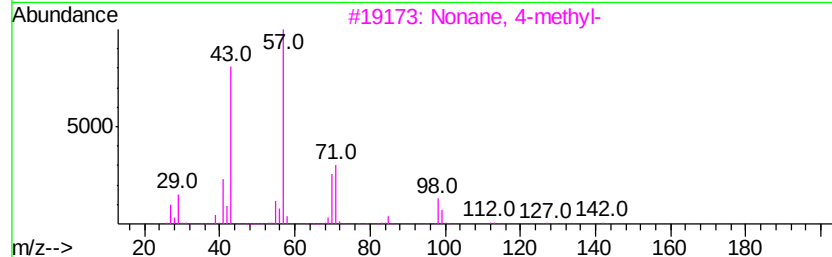
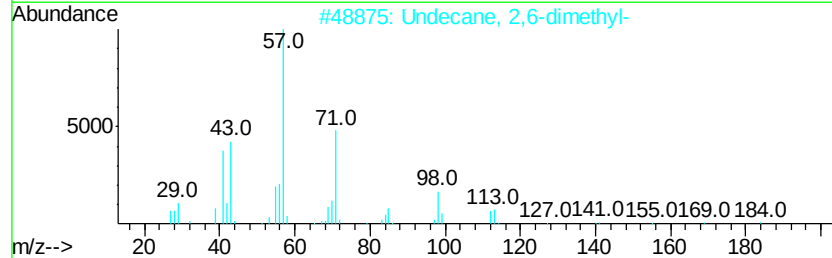
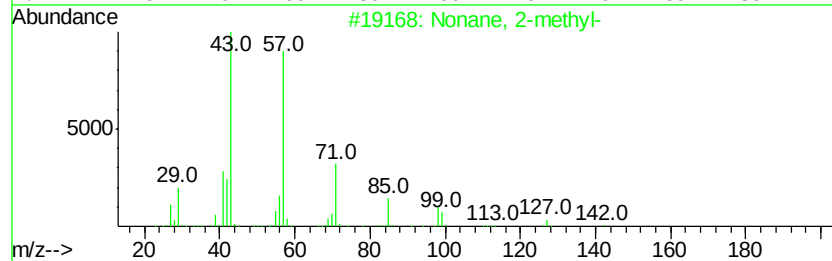
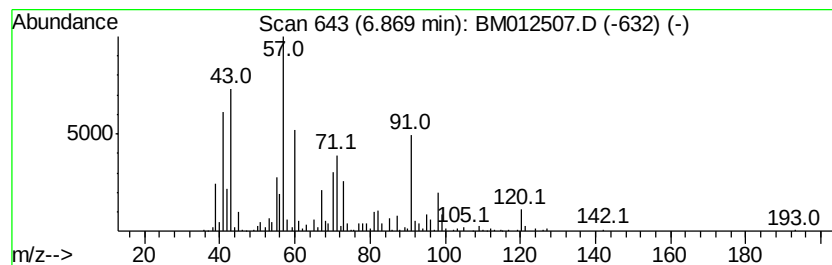
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Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.87	3.92 ng/ul	1031950	1,4-Dichlorobenzene-d4	7.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonane, 2-methyl-	142	C10H22	000871-83-0	38
2		Undecane, 2,6-dimethyl-	184	C13H28	017301-23-4	38
3		Nonane, 4-methyl-	142	C10H22	017301-94-9	30
4		Heptane, 2,3,4-trimethyl-	142	C10H22	052896-95-4	27
5		Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	27



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Instrument :
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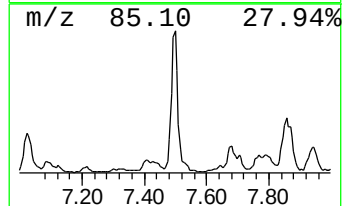
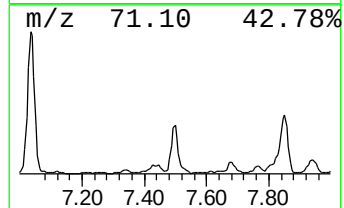
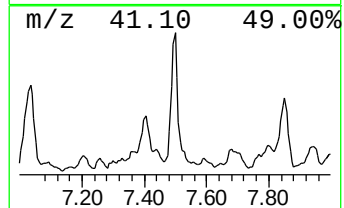
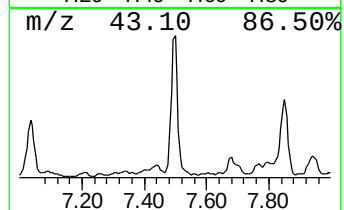
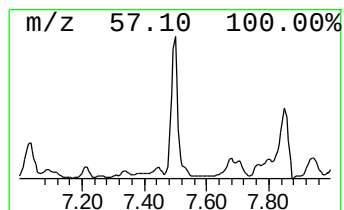
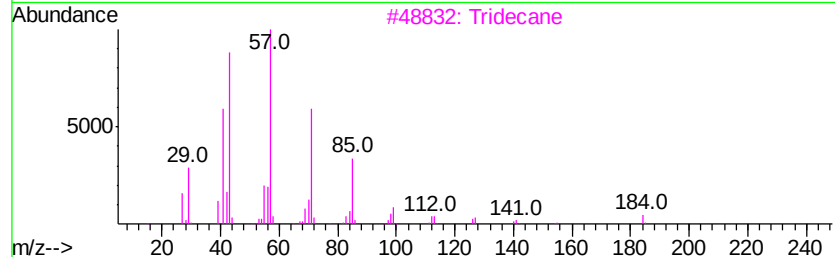
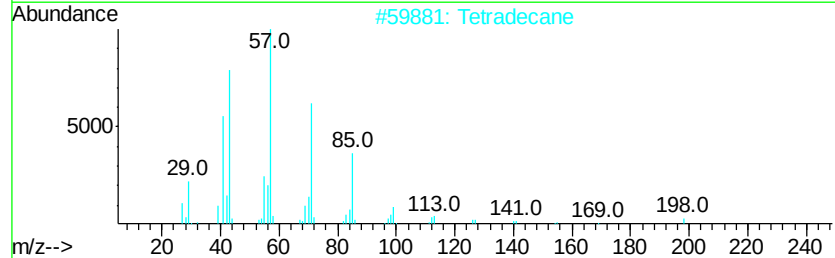
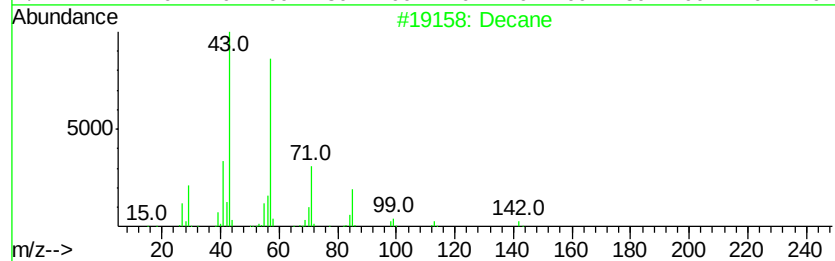
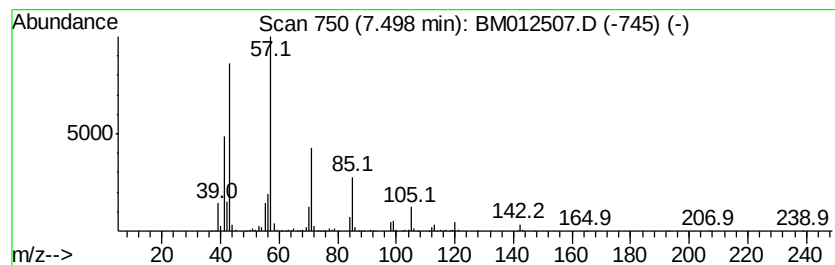
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM102417.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
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Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.50	7.15 ng/ul	1882450	1,4-Dichlorobenzene-d4	7.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Decane	142	C10H22	000124-18-5	91
2		Tetradecane	198	C14H30	000629-59-4	72
3		Tridecane	184	C13H28	000629-50-5	64
4		Hexatriacontane	507	C36H74	000630-06-8	64
5		Decane	142	C10H22	000124-18-5	64



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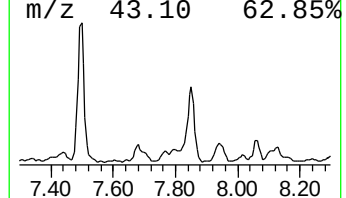
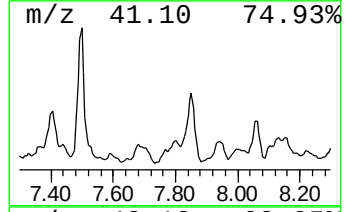
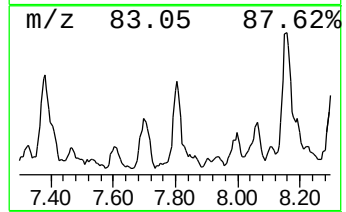
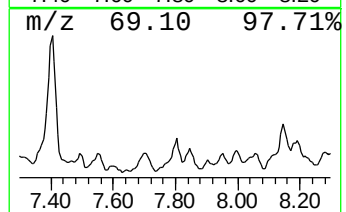
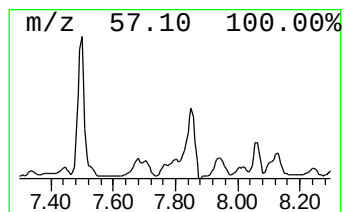
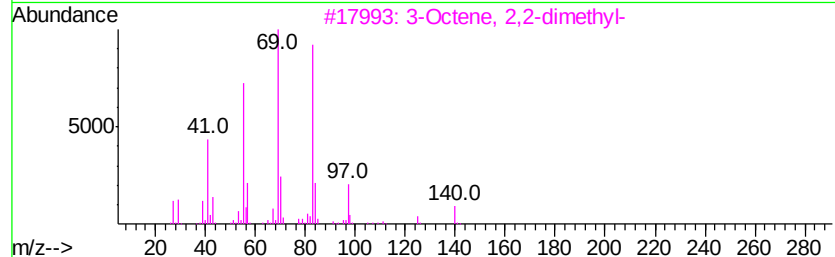
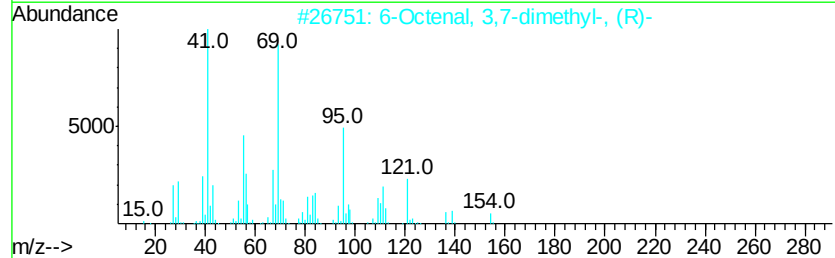
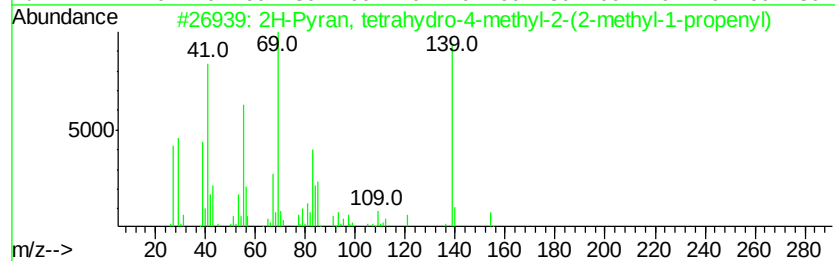
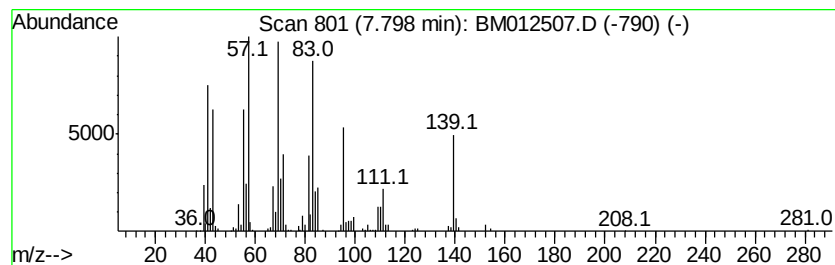
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Peak Number 6 2H-Pyran, tetrahydro-4-meth... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.		
7.80	3.32 ng/ul	873530	1,4-Dichlorobenzene-d4	7.87		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		2H-Pyran, tetrahydro-4-methyl-2-...	154	C10H18O	016409-43-1	50
2		6-Octenal, 3,7-dimethyl-, (R)-	154	C10H18O	002385-77-5	50
3		3-Octene, 2,2-dimethyl-	140	C10H20	086869-76-3	38
4		5-Undecene, 7-methyl-, (Z)-	168	C12H24	074630-62-9	22
5		Methanone, dicyclopropyl-	110	C7H10O	001121-37-5	18



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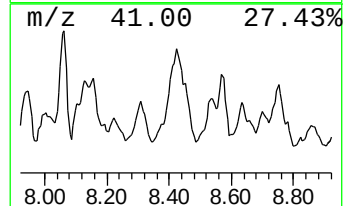
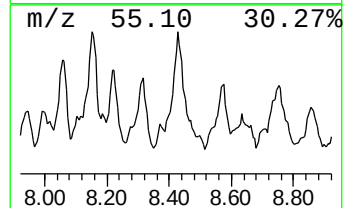
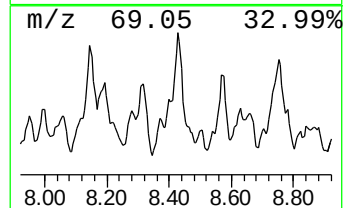
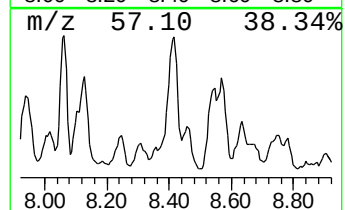
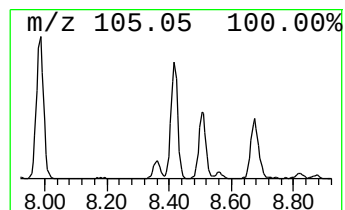
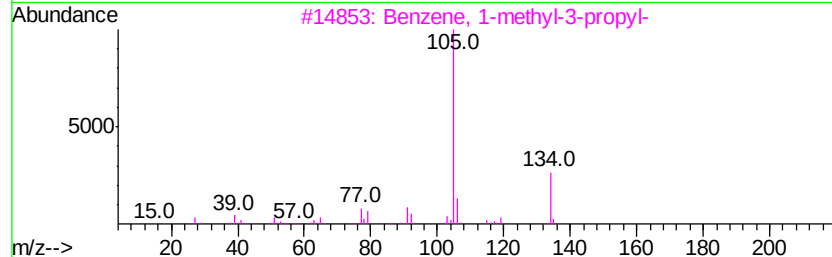
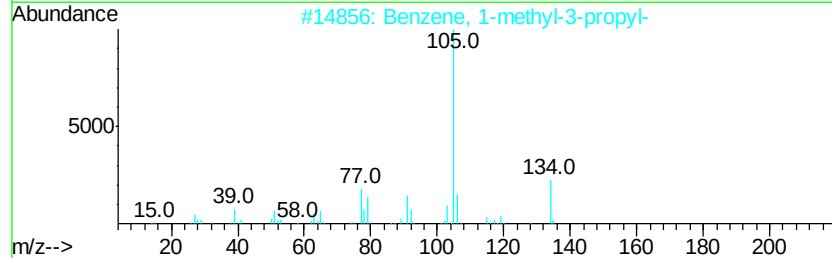
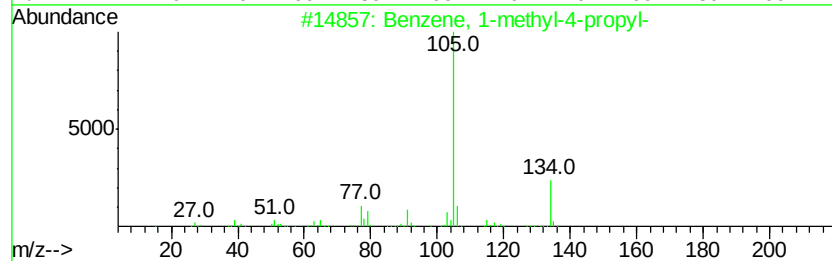
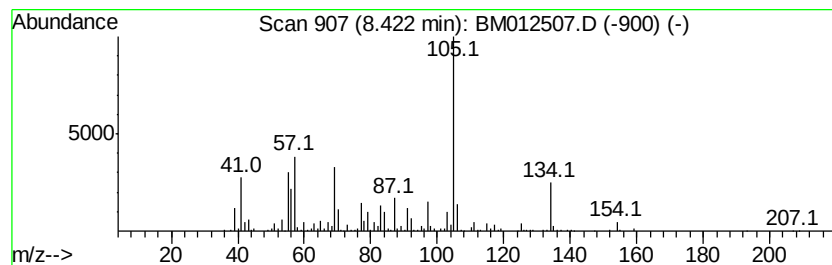
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TIC Integration Parameters: LSCINT.P

Peak Number 7 Benzene, 1-methyl-4-propyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.42	6.19 ng/ul	1630100	1,4-Dichlorobenzene-d4	7.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	86
2		Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	50
3		Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	46
4		Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	46
5		Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	46



Data Path : Z:\HPCHEM1\BNA M\DATA\BM102717\
Data File : BM012507.D
Acq On : 28 Oct 2017 04:17
Operator : SJ/JU
Sample : I5786-08
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
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ClientSampleId :
B-7(0.5-1.5)_101017

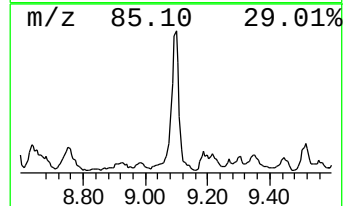
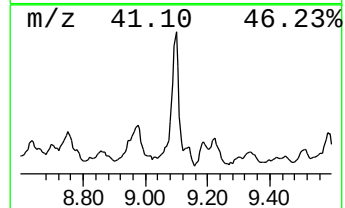
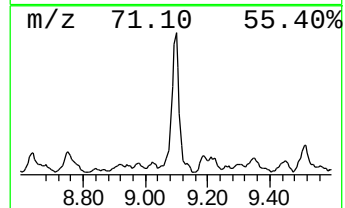
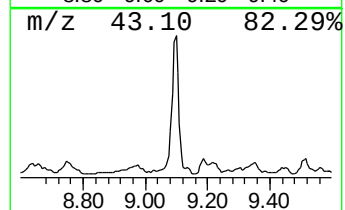
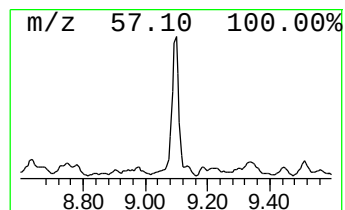
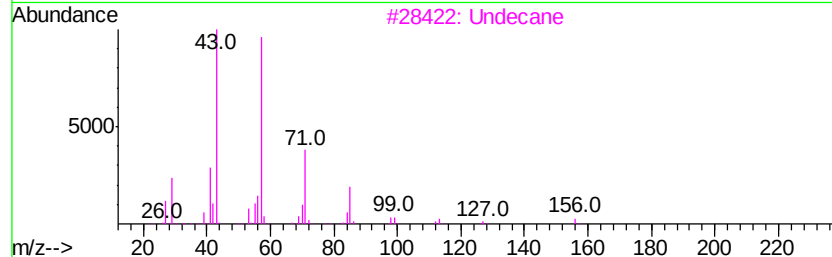
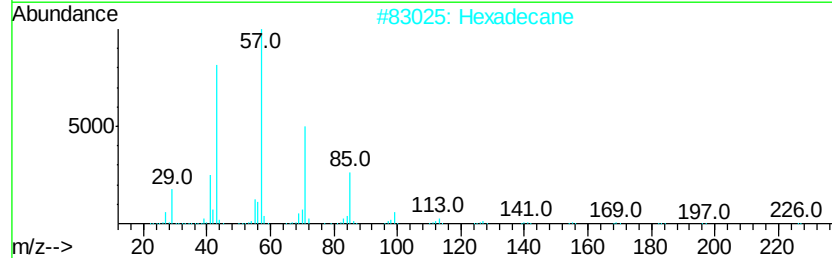
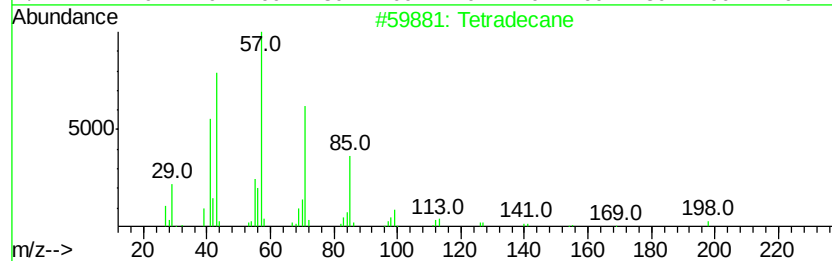
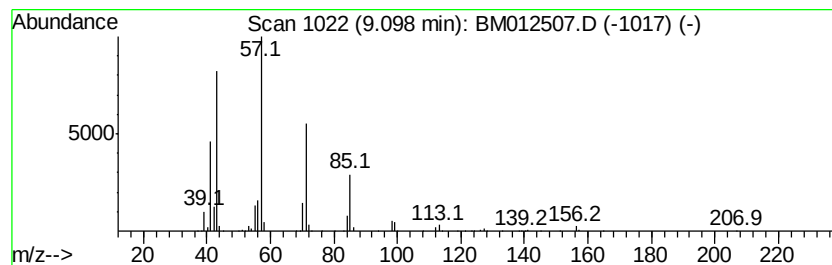
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.10	4.58 ng/ul	1205860	1,4-Dichlorobenzene-d4	7.87

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecane	198	C14H30	000629-59-4	90
2			Hexadecane	226	C16H34	000544-76-3	78
3			Undecane	156	C11H24	001120-21-4	76
4			1-Iodo-2-methylnonane	268	C10H21I	1000101-47-9	64
5			Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	64



Data Path : Z:\HPCHEM1\BNA M\DATA\BM102717\
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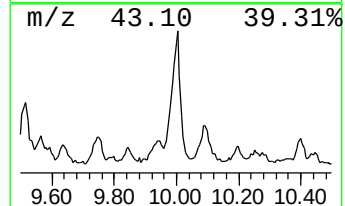
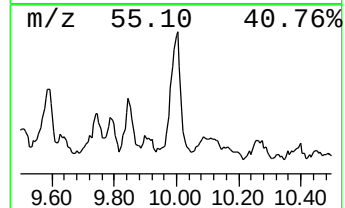
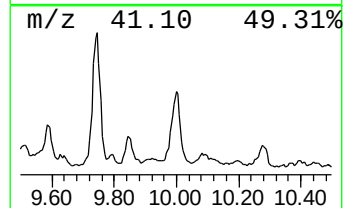
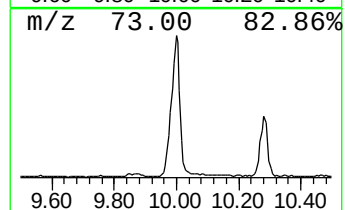
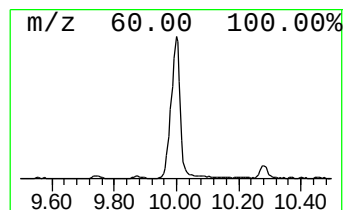
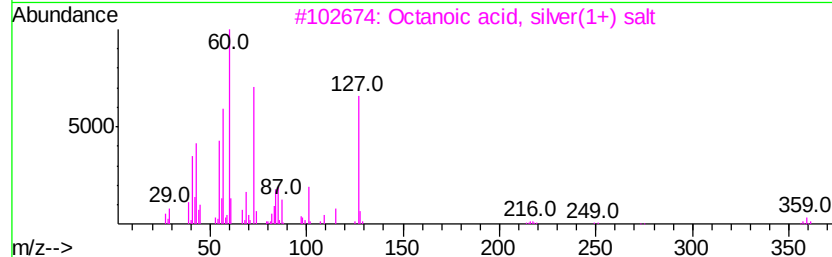
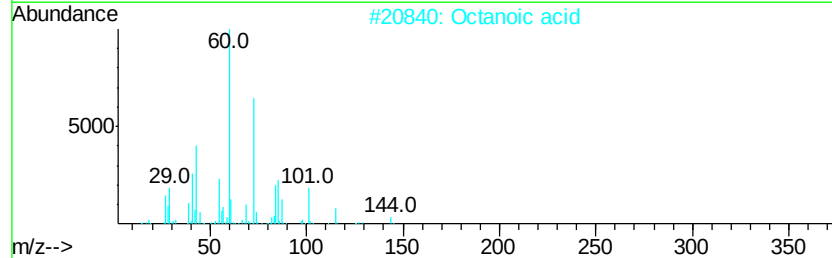
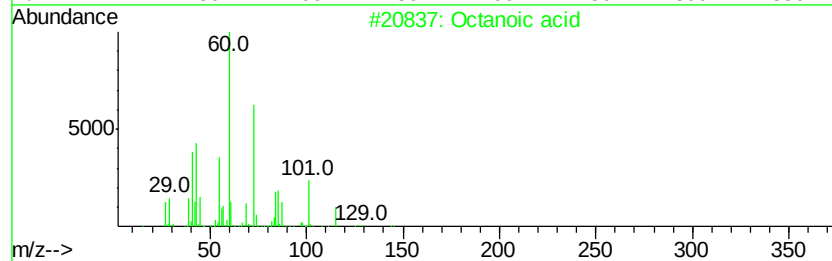
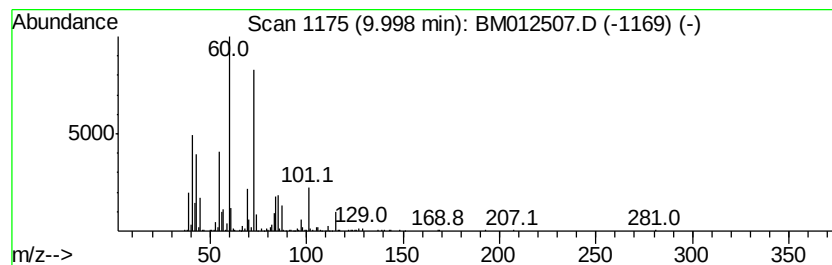
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Octanoic acid Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.00	3.59 ng/ul	1102480	Naphthalene-d8	10.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octanoic acid	144	C8H16O2	000124-07-2	96
2		Octanoic acid	144	C8H16O2	000124-07-2	90
3		Octanoic acid, silver(1+) salt	250	C8H15AgO2	024927-67-1	86
4		Octanoic acid	144	C8H16O2	000124-07-2	86
5		Hexanoic acid	116	C6H12O2	000142-62-1	46



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Sample : I5786-08
Misc :
ALS Vial : 31 Sample Multiplier: 1

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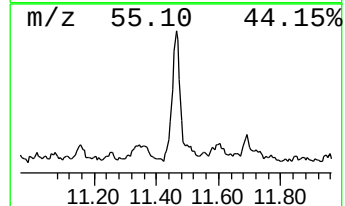
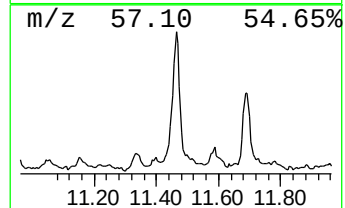
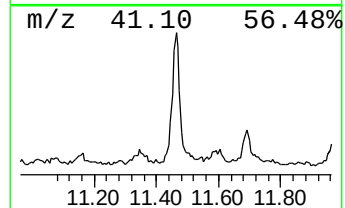
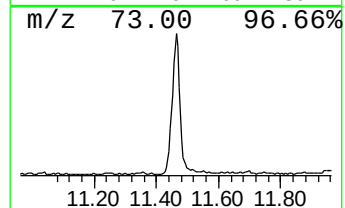
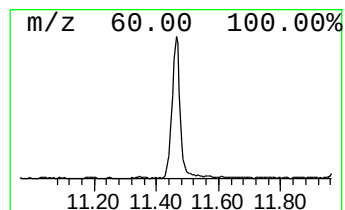
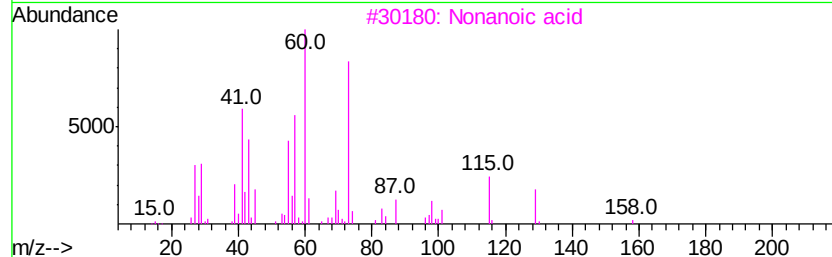
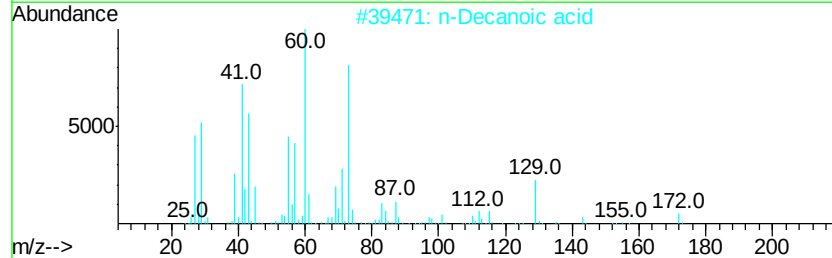
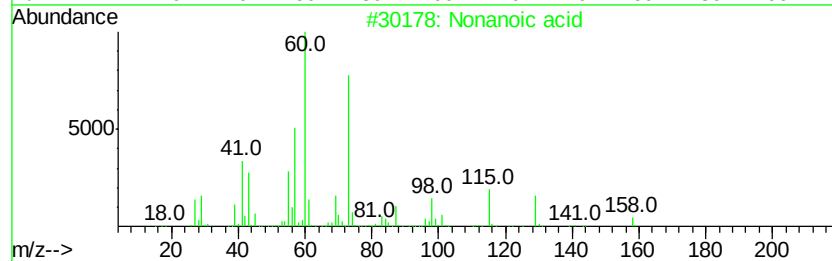
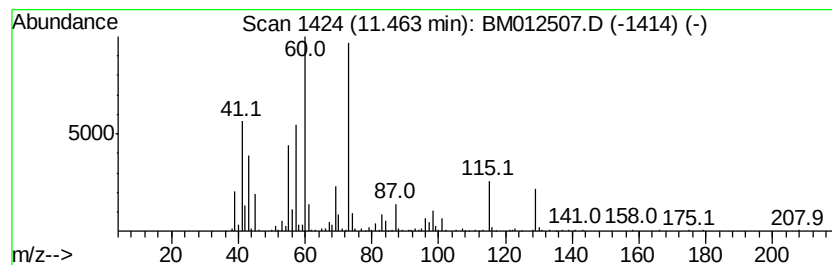
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Nonanoic acid Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.46	3.14 ng/ul	965847	Naphthalene-d8	10.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonanoic acid	158	C9H18O2	000112-05-0	72
2		n-Decanoic acid	172	C10H20O2	000334-48-5	72
3		Nonanoic acid	158	C9H18O2	000112-05-0	64
4		n-Decanoic acid	172	C10H20O2	000334-48-5	59
5		n-Decanoic acid	172	C10H20O2	000334-48-5	59



Data Path : Z:\HPCHEM1\BNA M\DATA\BM102717\
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Acq On : 28 Oct 2017 04:17
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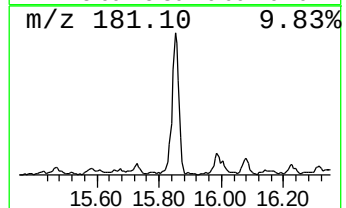
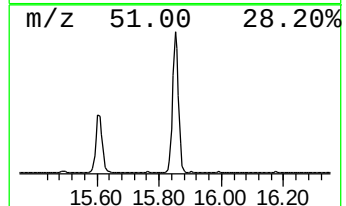
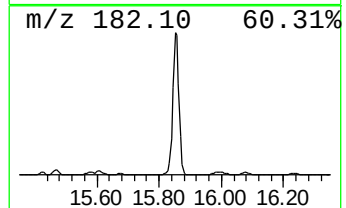
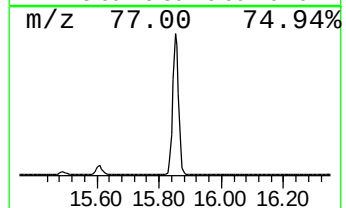
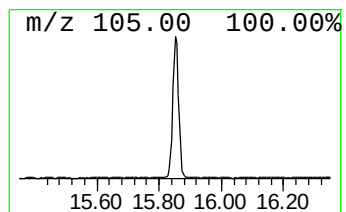
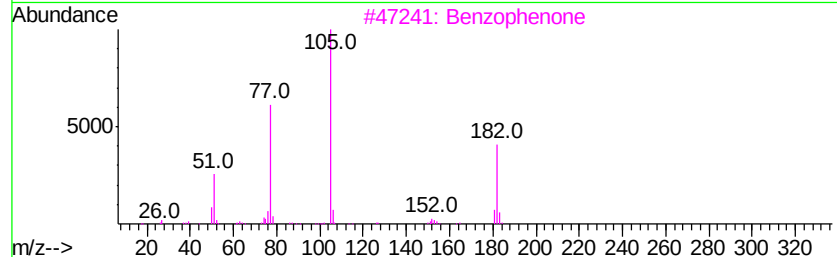
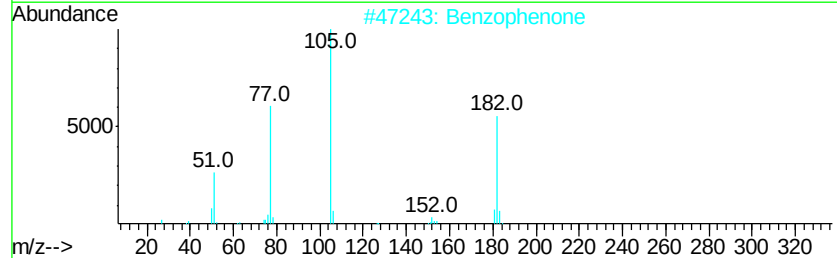
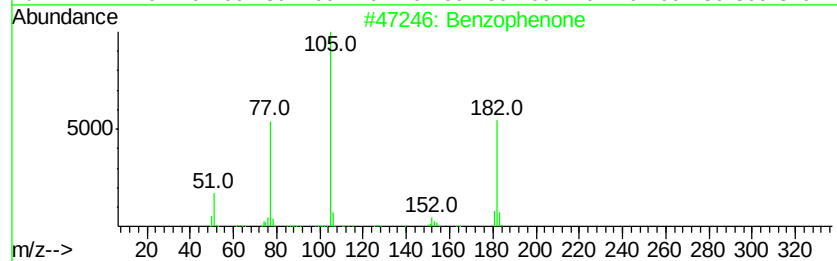
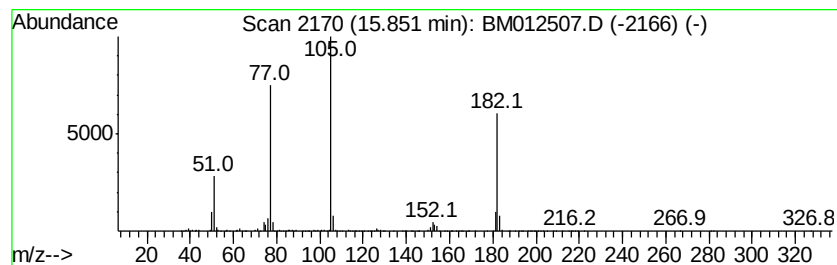
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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM102417.M
Quant Title : SVOA CALIBRATION

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

Peak Number 11 Benzophenone Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.85	5.76 ng/ul	2079830	Acenaphthene-d10	14.50
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Benzophenone	182 C13H10O	000119-61-9	97
2	Benzophenone	182 C13H10O	000119-61-9	96
3	Benzophenone	182 C13H10O	000119-61-9	96
4	Benzophenone	182 C13H10O	000119-61-9	95
5	Benzophenone	182 C13H10O	000119-61-9	91



Data Path : Z:\HPCHEM1\BNA M\DATA\BM102717\
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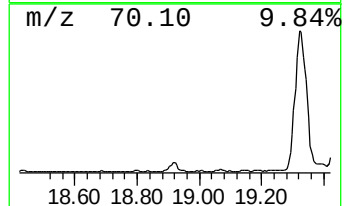
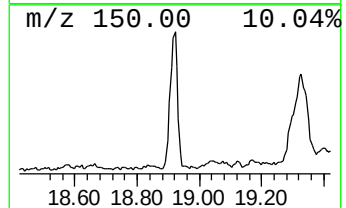
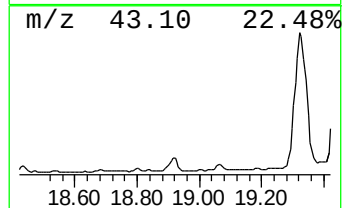
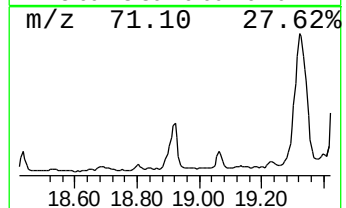
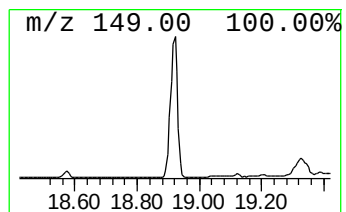
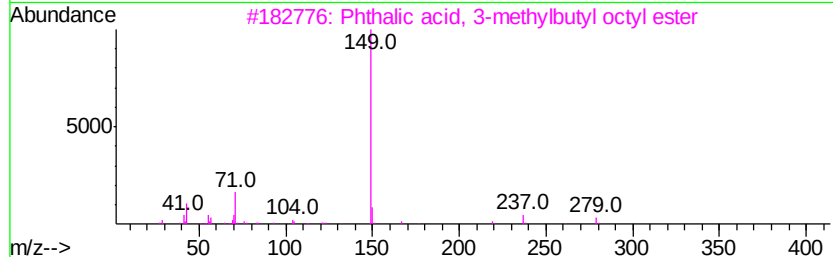
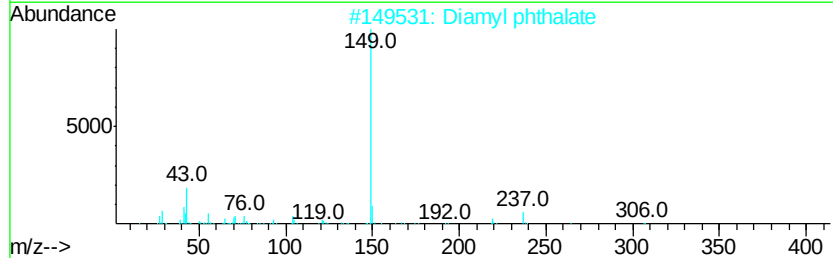
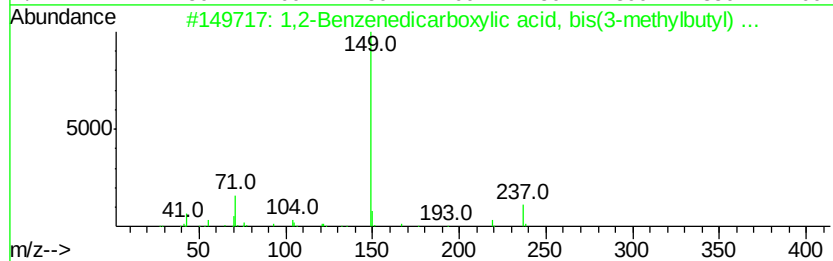
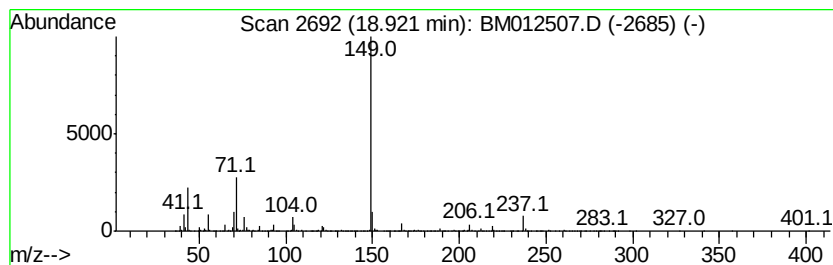
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 1,2-Benzenedicarboxylic aci... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.92	7.65 ng/ul	3002070	Phenanthrene-d10	17.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,2-Benzenedicarboxylic acid, bi...	306	C18H26O4	000605-50-5	86
2		Diamyl phthalate	306	C18H26O4	000131-18-0	72
3		Phthalic acid, 3-methylbutyl oct...	348	C21H32O4	1000356-80-9	64
4		Phthalic acid, isohexyl pentyl e...	320	C19H28O4	1000308-95-8	64
5		Diamyl phthalate	306	C18H26O4	000131-18-0	64



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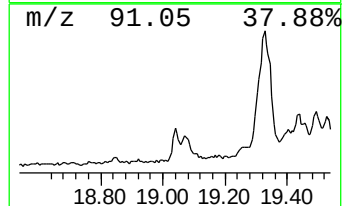
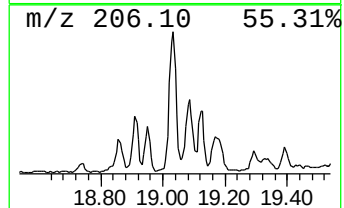
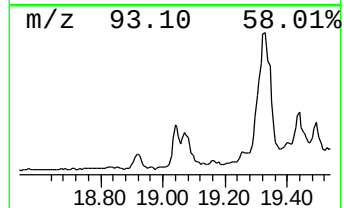
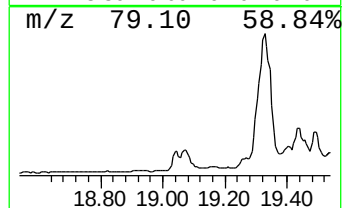
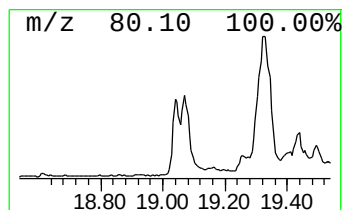
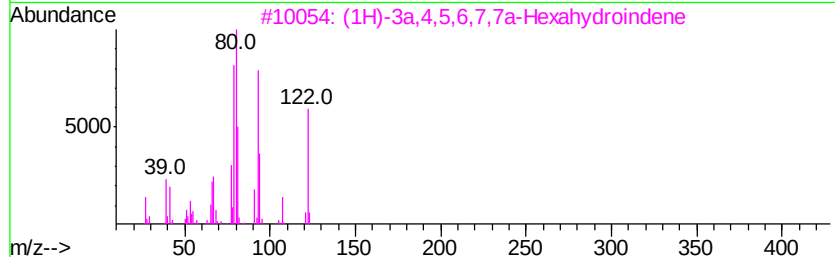
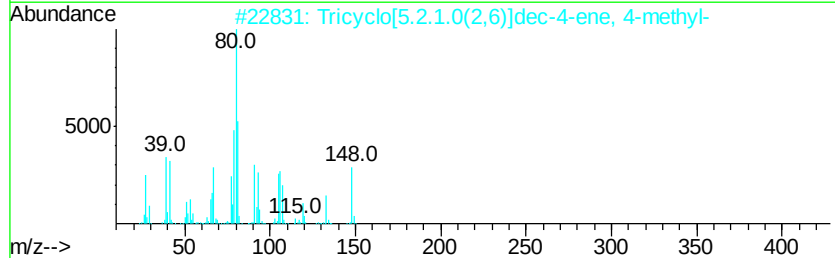
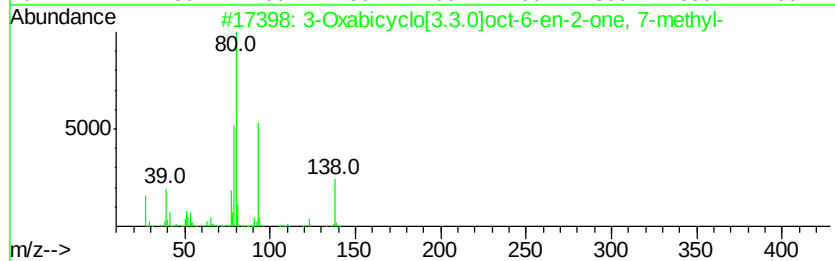
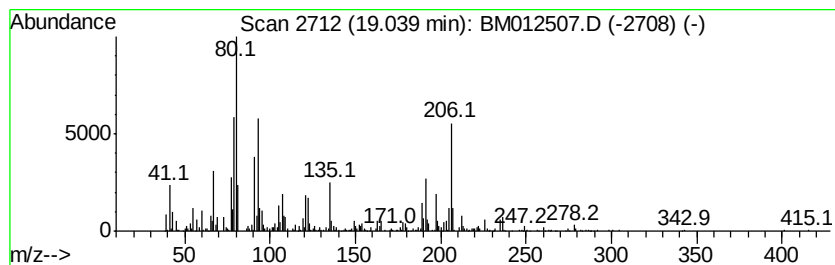
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 3-Oxabicyclo[3.3.0]oct-6-en... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.		
19.04	2.67 ng/ul	1048770	Phenanthrene-d10	17.24		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		3-Oxabicyclo[3.3.0]oct-6-en-2-on...	138	C8H10O2	1000155-62-3	59
2		Tricyclo[5.2.1.0(2,6)]dec-4-ene,...	148	C11H16	1000150-03-0	59
3		(1H)-3a,4,5,6,7,7a-Hexahydroindene	122	C9H14	1000196-59-5	59
4		Bicyclo[3.2.0]hept-2-ene, 2-methyl-	108	C8H12	094122-58-4	53
5		4,7-Methanoisobenzofuran-1,3-dio...	178	C10H10O3	000117-40-8	52



Data Path : Z:\HPCHEM1\BNA M\DATA\BM102717\
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ALS Vial : 31 Sample Multiplier: 1

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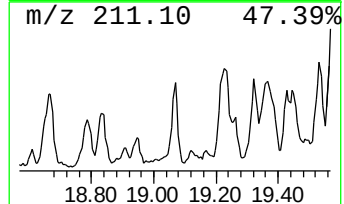
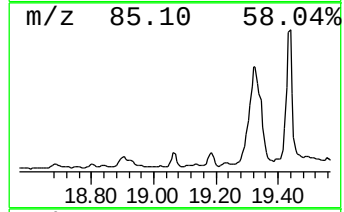
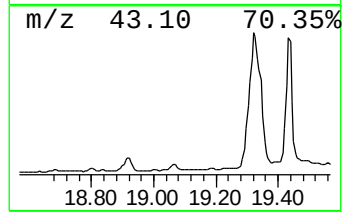
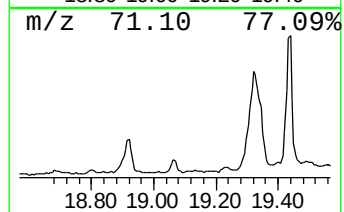
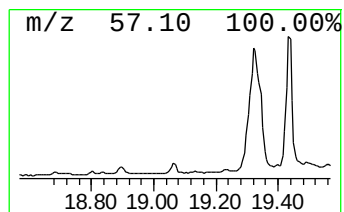
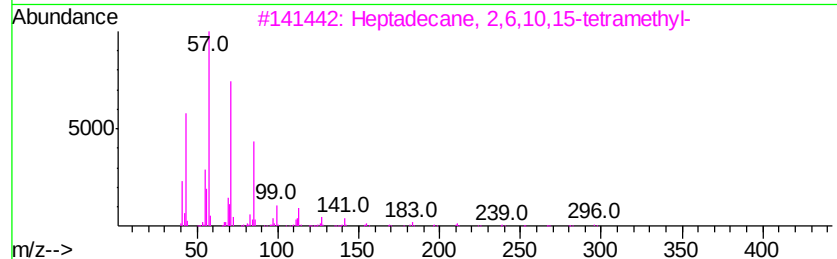
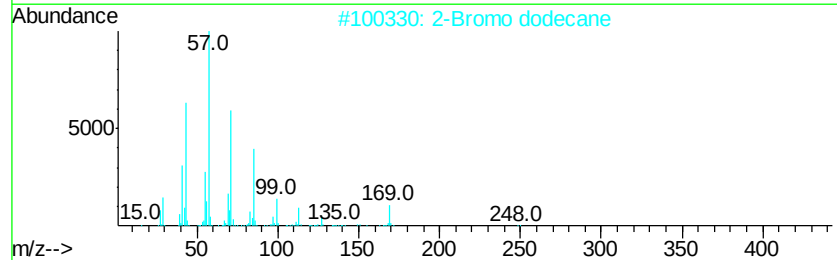
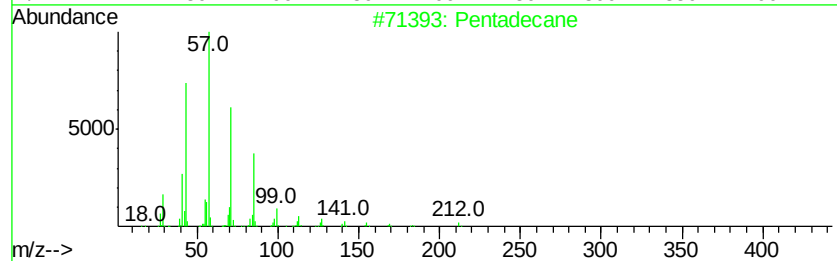
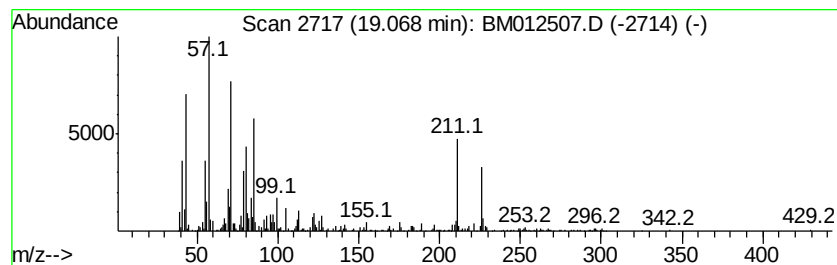
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 (DEL) Alkane: Straight-Chai... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.07	3.82 ng/ul	1498880	Phenanthrene-d10	17.24

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecane	212	C15H32	000629-62-9	93
2			2-Bromo dodecane	248	C12H25Br	013187-99-0	89
3			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	83
4			Heneicosane	296	C21H44	000629-94-7	64
5			Octadecane, 2-methyl-	268	C19H40	001560-88-9	64



Data Path : Z:\HPCHEM1\BNA M\DATA\BM102717\
Data File : BM012507.D
Acq On : 28 Oct 2017 04:17
Operator : SJ/JU
Sample : I5786-08
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
B-7(0.5-1.5)_101017

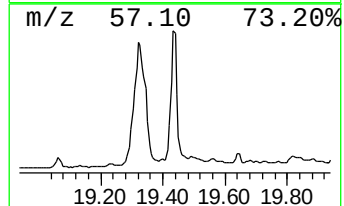
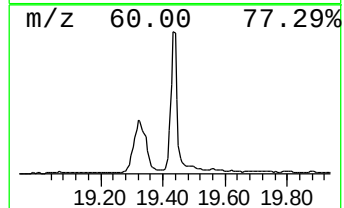
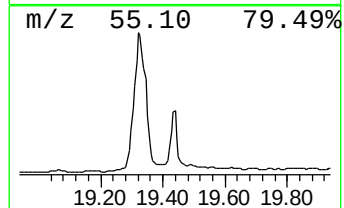
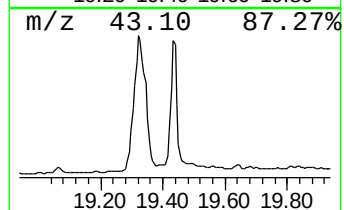
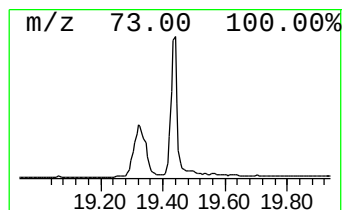
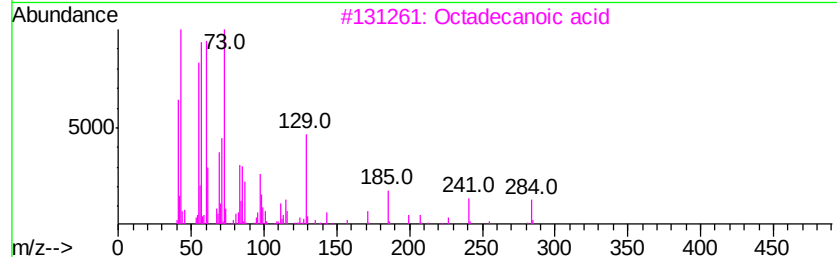
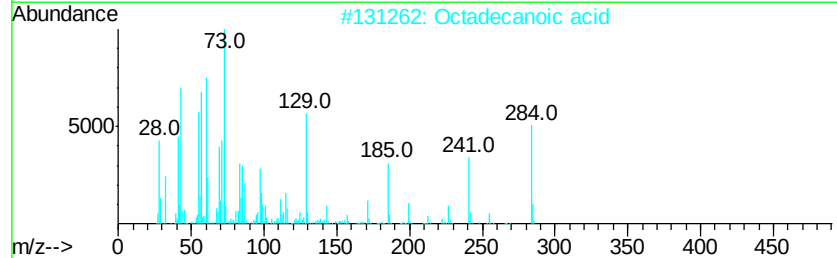
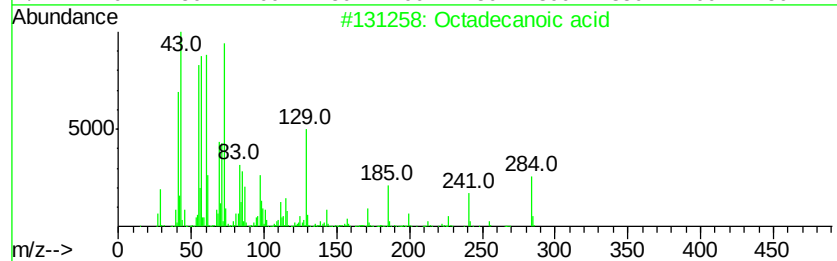
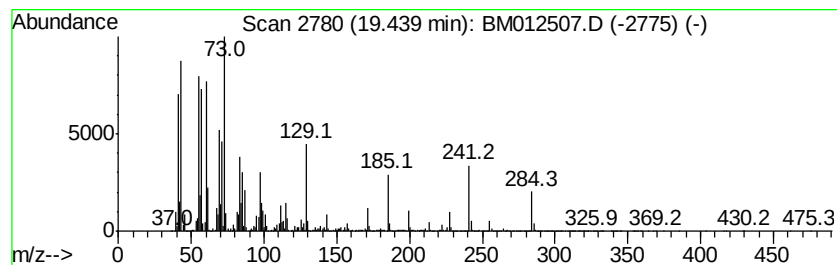
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 Octadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.44	51.29 ng/ul	21893200	Chrysene-d12	21.41

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	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	99
4		Octadecanoic acid	284	C18H36O2	000057-11-4	97
5		Octadecanoic acid	284	C18H36O2	000057-11-4	93



Data Path : Z:\HPCHEM1\BNA M\DATA\BM102717\
Data File : BM012507.D
Acq On : 28 Oct 2017 04:17
Operator : SJ/JU
Sample : I5786-08
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
B-7(0.5-1.5)_101017

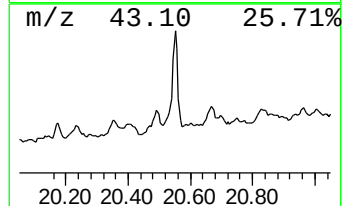
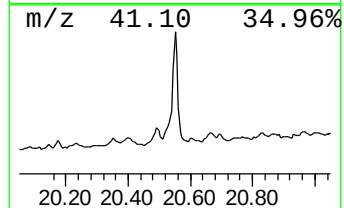
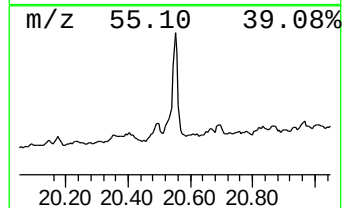
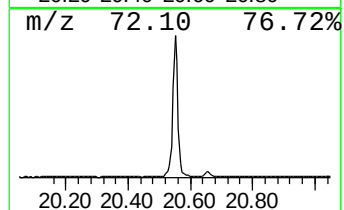
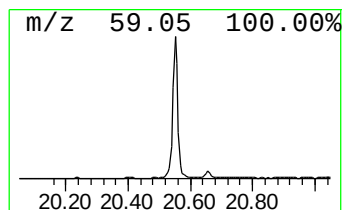
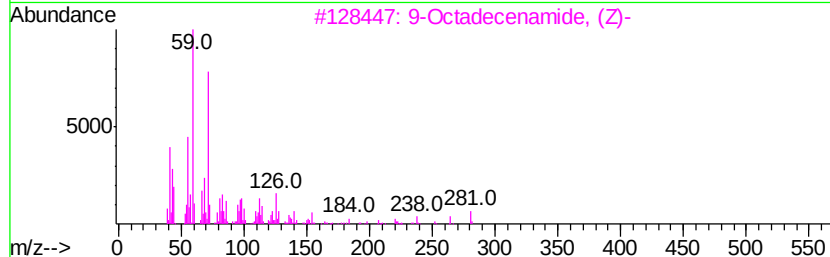
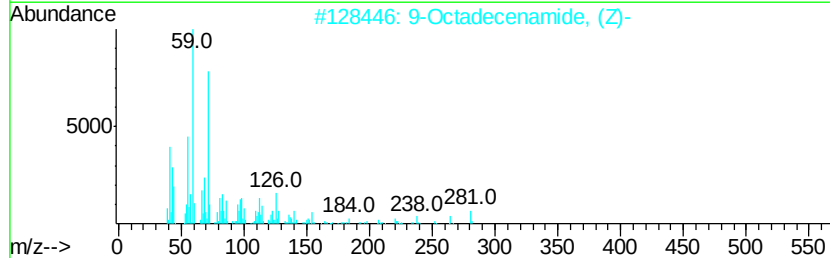
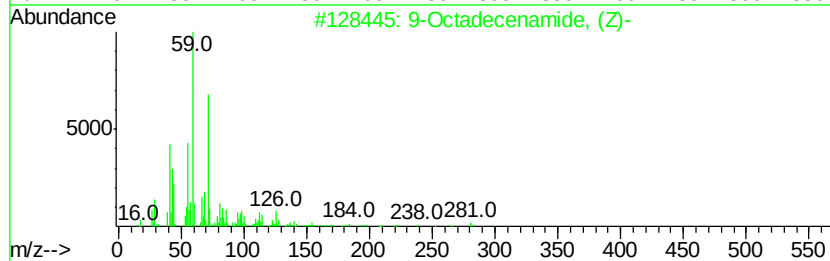
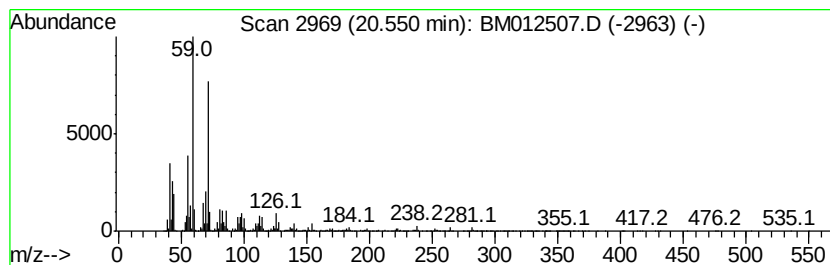
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 9-Octadecenamide, (Z)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.55	18.37 ng/ul	7839990	Chrysene-d12	21.41

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	94
3		9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	94
4		9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	53
5		9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	52



Data Path : Z:\HPCHEM1\BNA M\DATA\BM102717\
Data File : BM012507.D
Acq On : 28 Oct 2017 04:17
Operator : SJ/JU
Sample : I5786-08
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
B-7(0.5-1.5)_101017

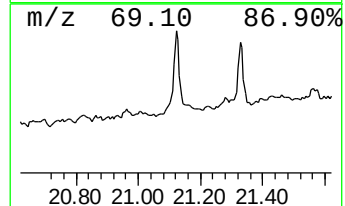
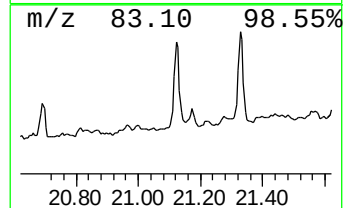
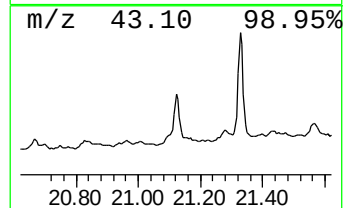
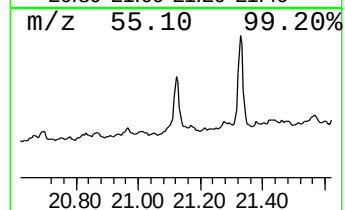
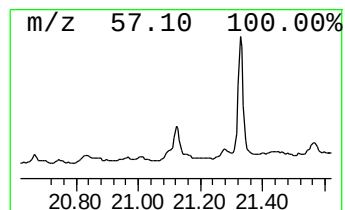
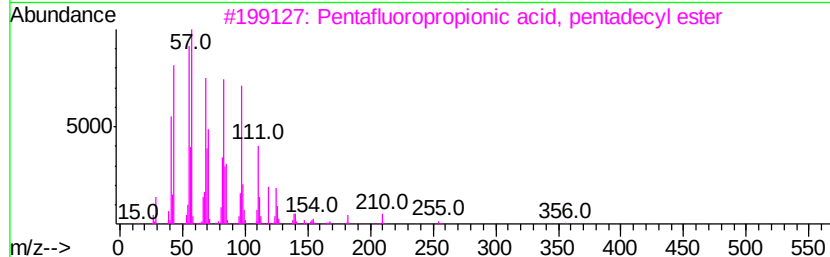
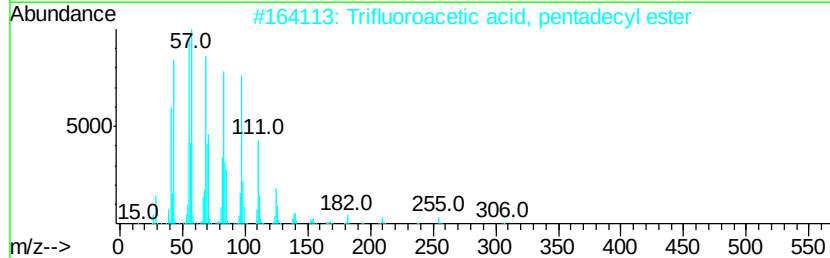
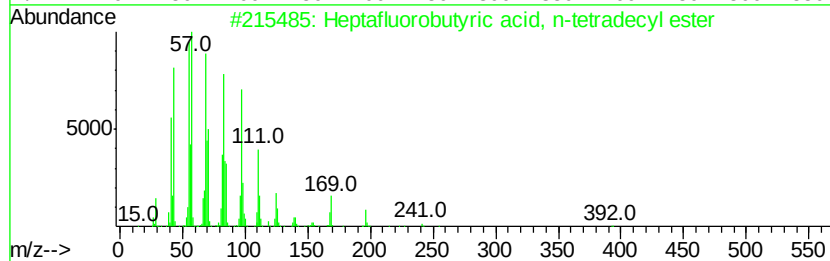
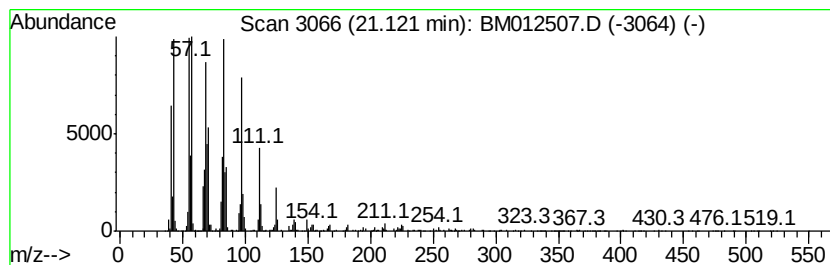
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 Heptafluorobutyric acid, n-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.12	6.07 ng/ul	2589270	Chrysene-d12	21.41

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptafluorobutyric acid, n-tetra...	410	C18H29F7O2	007365-36-8	94
2			Trifluoroacetic acid, pentadecyl...	324	C17H31F3O2	959010-23-2	94
3			Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	94
4			Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	93
5			Trifluoroacetoxyl hexadecane	338	C18H33F3O2	006222-03-3	93



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM102717\
 Data File : BM012507.D
 Acq On : 28 Oct 2017 04:17
 Operator : SJ/JU
 Sample : I5786-08
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102417.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
Benzene, 1,3-dime...	5.53	16.7	ng/ul	4396110	1	7.87	5266240	20.0
(DEL) Alkane: Str...	6.87	3.9	ng/ul	1031950	1	7.87	5266240	20.0
(DEL) Alkane: Str...	7.50	7.2	ng/ul	1882450	1	7.87	5266240	20.0
2H-Pyran, tetrahy...	7.80	3.3	ng/ul	873530	1	7.87	5266240	20.0
Benzene, 1-methyl...	8.42	6.2	ng/ul	1630100	1	7.87	5266240	20.0
(DEL) Alkane: Str...	9.10	4.6	ng/ul	1205860	1	7.87	5266240	20.0
Octanoic acid	10.00	3.6	ng/ul	1102480	2	10.66	6149150	20.0
Nonanoic acid	11.46	3.1	ng/ul	965847	2	10.66	6149150	20.0
Benzophenone	15.85	5.8	ng/ul	2079830	3	14.50	7224780	20.0
1,2-Benzenedicarb...	18.92	7.7	ng/ul	3002070	4	17.24	7847410	20.0
3-Oxabicyclo[3.3...	19.04	2.7	ng/ul	1048770	4	17.24	7847410	20.0
(DEL) Alkane: Str...	19.07	3.8	ng/ul	1498880	4	17.24	7847410	20.0
Octadecanoic acid	19.44	51.3	ng/ul	21893200	5	21.41	8536850	20.0
9-Octadecenamide, ...	20.55	18.4	ng/ul	7839990	5	21.41	8536850	20.0
Heptafluorobutyri...	21.12	6.1	ng/ul	2589270	5	21.41	8536850	20.0