

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM121517\
 Data File : BM013455.D
 Acq On : 16 Dec 2017 01:07
 Operator : SJ/JU
 Sample : I6779-04
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 F9A87

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-BM113017.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	3.128	4	7	15	rVB2	83450	106612	1.50%	0.126%
2	4.246	192	197	207	rVB	1756949	2267896	31.88%	2.677%
3	4.834	292	297	306	rVB	432793	587316	8.26%	0.693%
4	6.857	634	641	651	rVB	1138583	2008998	28.24%	2.372%
5	7.016	662	668	678	rVB	814652	1202654	16.91%	1.420%
6	7.210	694	701	709	rBV	1196342	1834362	25.79%	2.166%
7	7.675	772	780	788	rBV	808881	1282259	18.03%	1.514%
8	8.381	893	900	909	rBV	1480180	2371164	33.33%	2.799%
9	8.828	968	976	985	rBV	1090677	1772130	24.91%	2.092%
10	9.545	1091	1098	1106	rVB	924638	1531007	21.52%	1.807%
11	10.081	1181	1189	1198	rBV	1477150	2502454	35.18%	2.954%
12	10.451	1245	1252	1257	rBV	1144888	1846573	25.96%	2.180%
13	10.498	1257	1260	1269	rVV	150111	246127	3.46%	0.291%
14	10.598	1269	1277	1293	rVV	523665	1010251	14.20%	1.193%
15	11.986	1505	1513	1521	rBV	670851	1122392	15.78%	1.325%
16	12.345	1564	1574	1581	rVB	83881	146341	2.06%	0.173%
17	13.145	1703	1710	1717	rVB2	68262	109693	1.54%	0.130%
18	13.498	1758	1770	1777	rBV2	193783	362749	5.10%	0.428%
19	13.633	1788	1793	1798	rBV2	65052	106013	1.49%	0.125%
20	13.733	1804	1810	1814	rVV	2629168	3582061	50.35%	4.229%
21	13.780	1814	1818	1826	rVB	674334	927685	13.04%	1.095%
22	14.010	1850	1857	1871	rVB	2873067	4258095	59.86%	5.027%
23	14.221	1889	1893	1899	rVV2	76847	101647	1.43%	0.120%
24	14.316	1899	1909	1915	rVV2	1582003	2477649	34.83%	2.925%
25	14.380	1915	1920	1928	rVB	278377	445165	6.26%	0.526%
26	14.533	1939	1946	1958	rBV	1150699	1848239	25.98%	2.182%
27	14.716	1966	1977	1985	rBV	254915	427860	6.01%	0.505%
28	15.180	2052	2056	2063	rVB	115648	147902	2.08%	0.175%
29	15.315	2071	2079	2084	rBV	3567210	5011196	70.44%	5.916%
30	15.368	2084	2088	2093	rVB	365348	500649	7.04%	0.591%
31	15.445	2093	2101	2107	rBV	883757	1272086	17.88%	1.502%
32	15.533	2112	2116	2123	rBV4	40607	113454	1.59%	0.134%
33	15.663	2134	2138	2146	rVB4	50541	91692	1.29%	0.108%
34	15.810	2157	2163	2172	rVB4	55805	143867	2.02%	0.170%

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35	16.039	2197	2202	2210	rBV4	133164	223892	3.15%	0.264%
36	16.292	2240	2245	2250	rBV	113644	156913	2.21%	0.185%
37	16.833	2332	2337	2341	rBV3	64866	91672	1.29%	0.108%
38	16.886	2341	2346	2351	rVB	118029	190443	2.68%	0.225%
39	17.068	2370	2377	2380	rBV2	2030904	2951283	41.49%	3.484%
40	17.109	2380	2384	2388	rVV	1201321	1654017	23.25%	1.953%
41	17.168	2388	2394	2405	rVB	3856050	5678530	79.82%	6.704%
42	17.474	2436	2446	2451	rBV	202768	304974	4.29%	0.360%
43	17.574	2458	2463	2468	rVB4	59502	93355	1.31%	0.110%
44	17.939	2515	2525	2527	rBV3	88532	148077	2.08%	0.175%
45	17.974	2527	2531	2544	rVV3	767828	1281120	18.01%	1.512%
46	18.068	2544	2547	2552	rVV3	79495	137363	1.93%	0.162%
47	18.133	2553	2558	2563	rVV2	143330	266800	3.75%	0.315%
48	18.174	2563	2565	2570	rVB3	81996	102623	1.44%	0.121%
49	18.356	2589	2596	2605	rBV7	100973	210482	2.96%	0.248%
50	18.445	2607	2611	2617	rVV	60174	85919	1.21%	0.101%
51	18.539	2621	2627	2631	rBV3	154856	201939	2.84%	0.238%
52	19.127	2718	2727	2733	rBV	671244	952672	13.39%	1.125%
53	19.186	2733	2737	2742	rVB	269317	325889	4.58%	0.385%
54	19.256	2742	2749	2759	rBV	425136	663787	9.33%	0.784%
55	19.350	2763	2765	2772	rVB4	42126	79392	1.12%	0.094%
56	19.468	2778	2785	2798	rBV	4678098	7113728	100.00%	8.398%
57	19.909	2855	2860	2865	rBV	2144544	2712951	38.14%	3.203%
58	19.992	2870	2874	2882	rVB	129377	190864	2.68%	0.225%
59	20.092	2886	2891	2894	rBV	96431	127511	1.79%	0.151%
60	20.974	3037	3041	3052	rVB	1015753	1312608	18.45%	1.550%
61	21.256	3082	3089	3094	rBV	2807081	3705896	52.09%	4.375%
62	21.292	3094	3095	3100	rVB	131012	126470	1.78%	0.149%
63	21.415	3113	3116	3123	rVV6	62117	91772	1.29%	0.108%
64	22.309	3265	3268	3272	rVB2	87242	98234	1.38%	0.116%
65	22.809	3349	3353	3359	rBV2	150668	252135	3.54%	0.298%
66	23.338	3436	3443	3456	rVB2	2972189	5721388	80.43%	6.755%
67	23.480	3460	3467	3481	rVB	1495036	2854075	40.12%	3.369%
68	24.009	3553	3557	3564	rVB5	107272	201854	2.84%	0.238%
69	24.091	3566	3571	3583	rVB5	184981	437401	6.15%	0.516%
70	24.744	3678	3682	3688	rVB7	94359	189159	2.66%	0.223%

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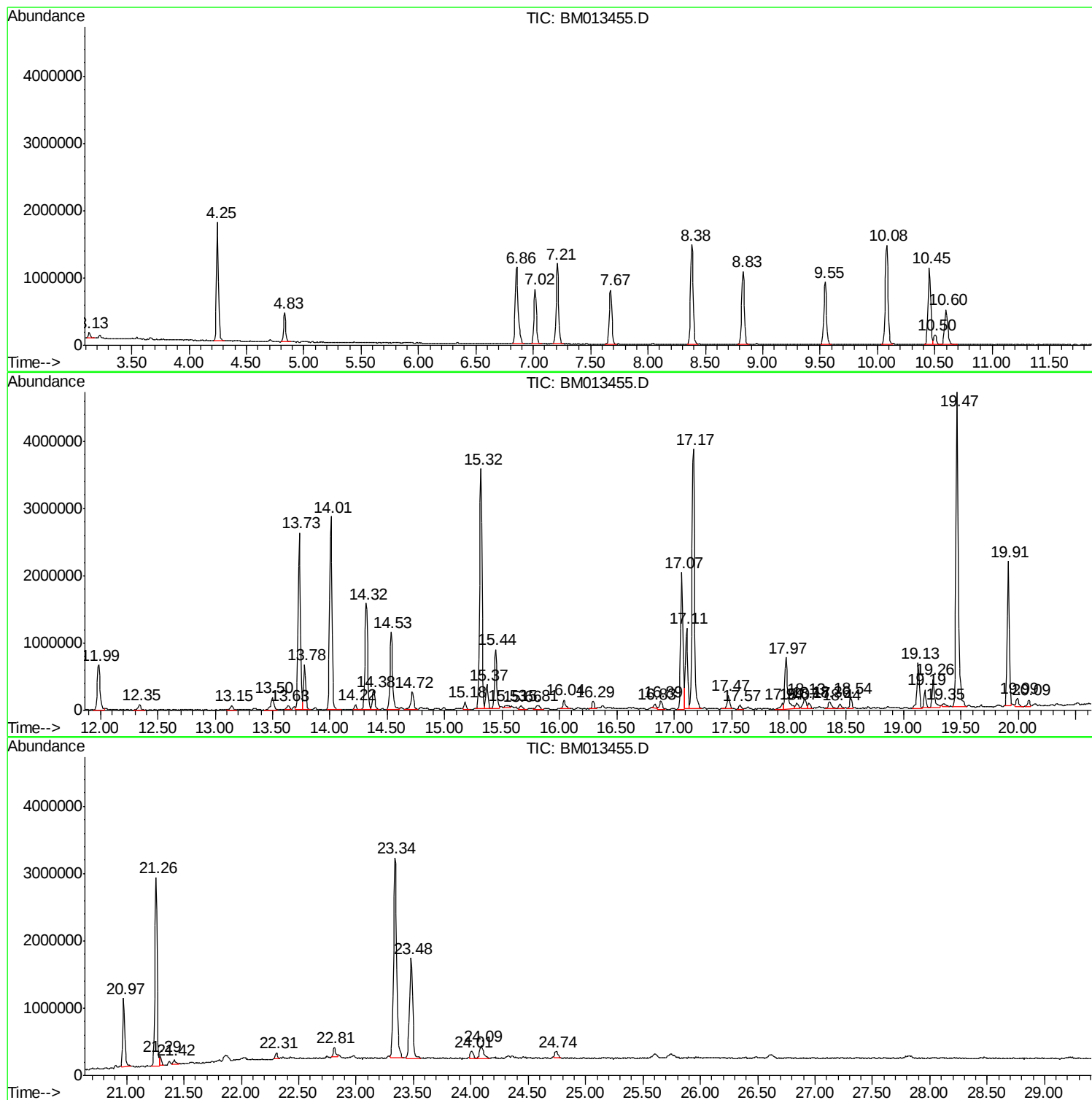
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM113017.M
Quant Title : SVOA CALIBRATION

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



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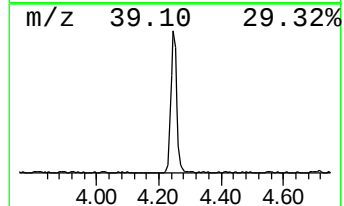
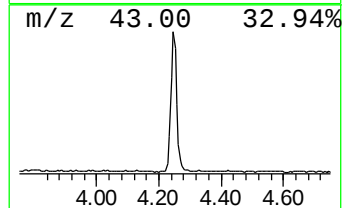
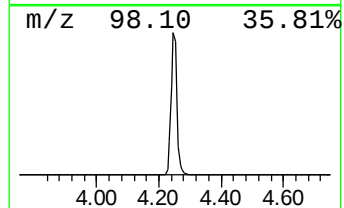
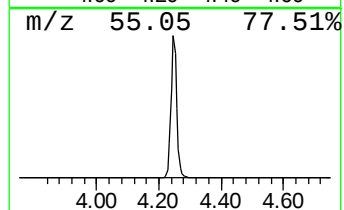
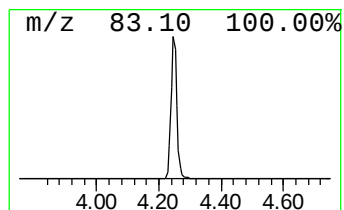
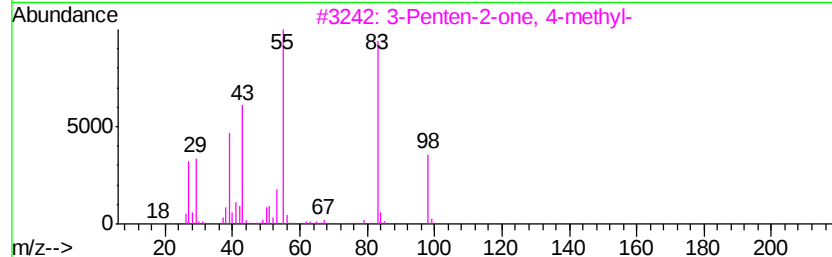
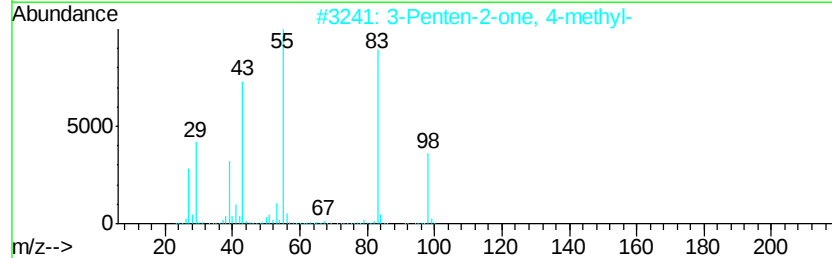
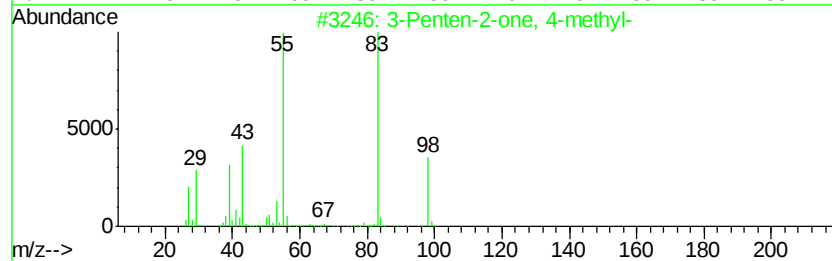
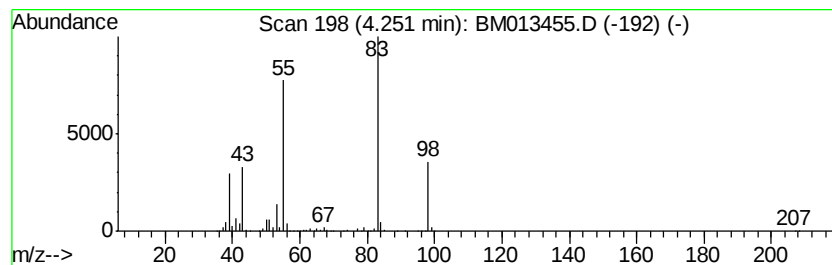
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Quant Title : SVOA CALIBRATION

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

Peak Number 1 3-Penten-2-one, 4-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.25	35.37 ng/ul	2267900	1,4-Dichlorobenzene-d4	7.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	94
2		3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	91
3		3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	91
4		3-Hexen-2-one	98	C6H10O	000763-93-9	90
5		2-Pentene, 3,4-dimethyl-, (E)-	98	C7H14	004914-92-5	86



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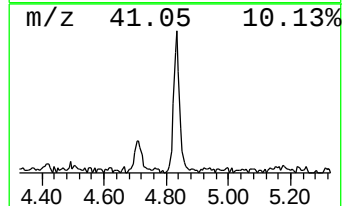
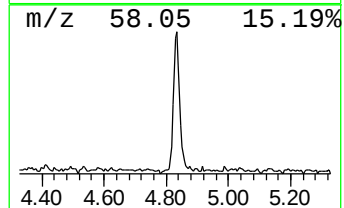
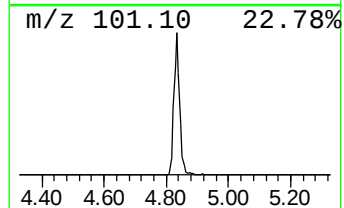
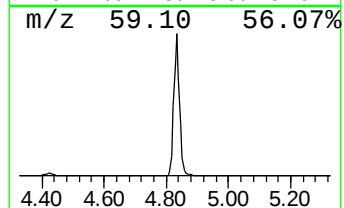
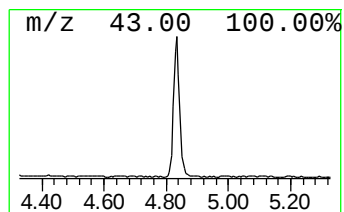
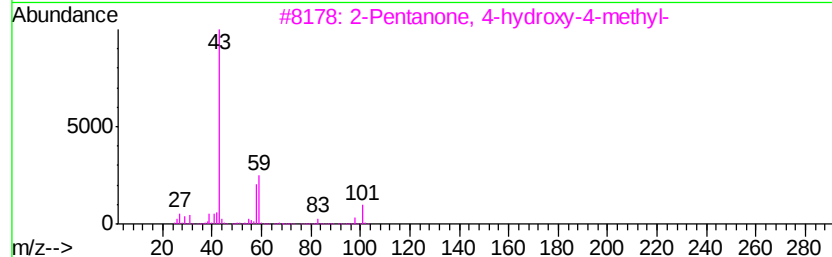
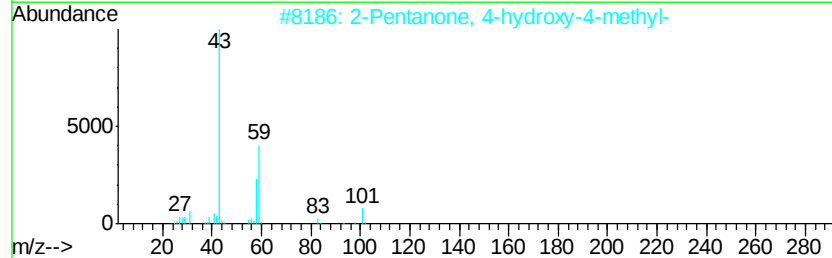
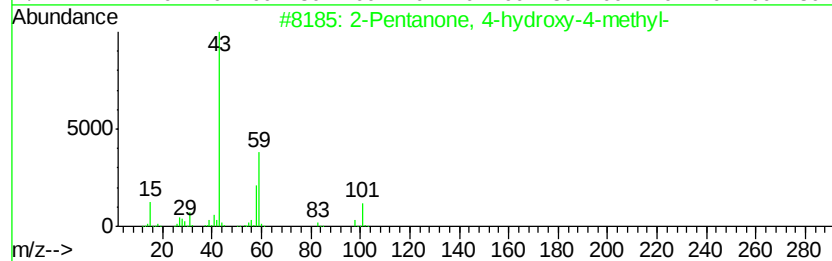
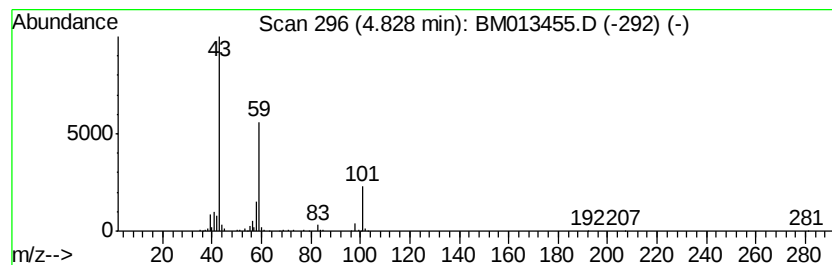
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM113017.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.83	9.16 ng/ul	587316	1,4-Dichlorobenzene-d4	7.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
3		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
4		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	38
5		3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	25



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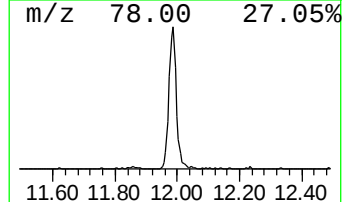
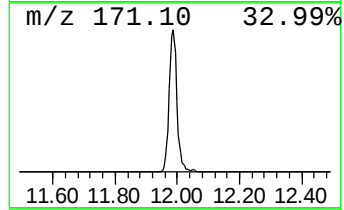
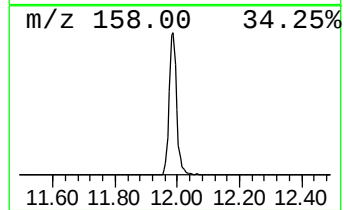
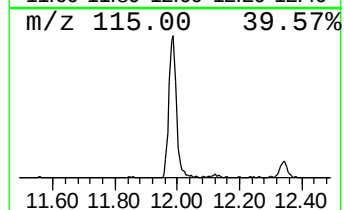
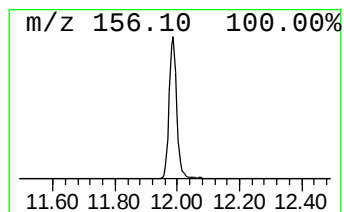
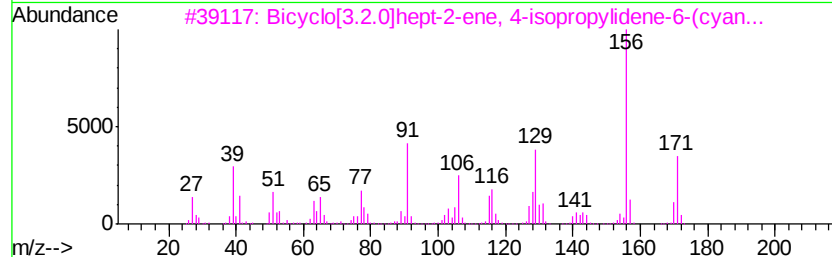
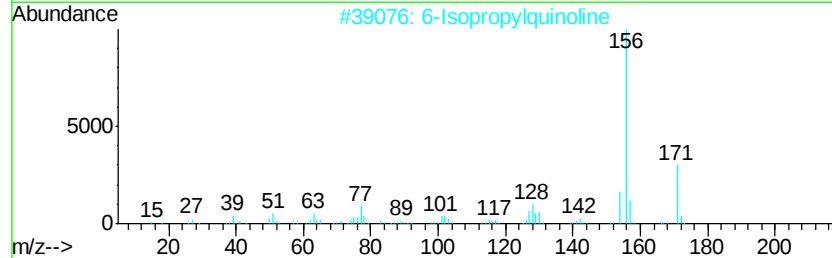
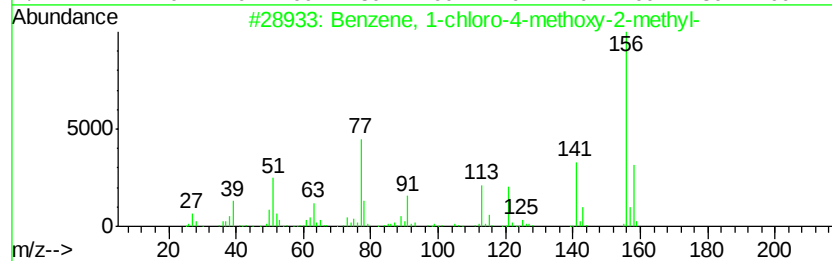
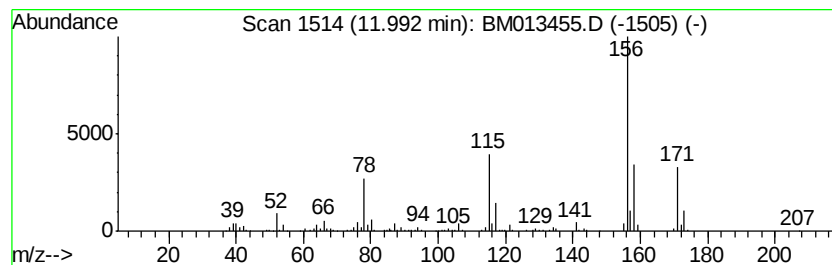
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown-01 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.99	12.16 ng/ul	1122390	Naphthalene-d8	10.45

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-chloro-4-methoxy-2-me...	156	C8H9ClO	013334-71-9	38
2		6-Isopropylquinoline	171	C12H13N	000135-79-5	32
3		Bicyclo[3.2.0]hept-2-ene, 4-isop...	171	C12H13N	1000217-15-6	32
4		Benzene, 2,5-cyclohexadien-1-yl-	156	C12H12	004794-05-2	28
5		2-(2-Chloro-pyridin-3-yl)-propan...	171	C8H10ClNO	267003-35-0	25



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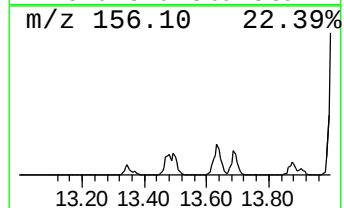
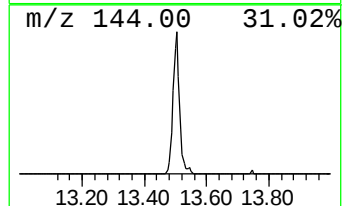
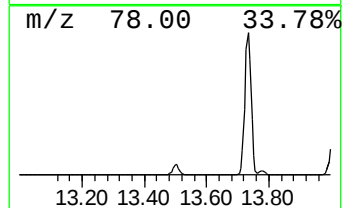
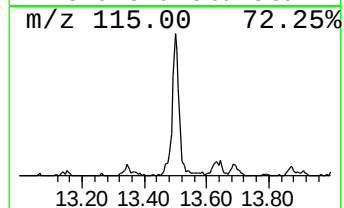
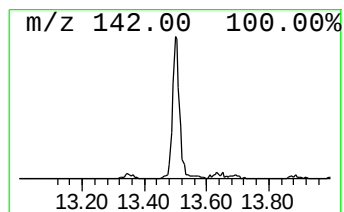
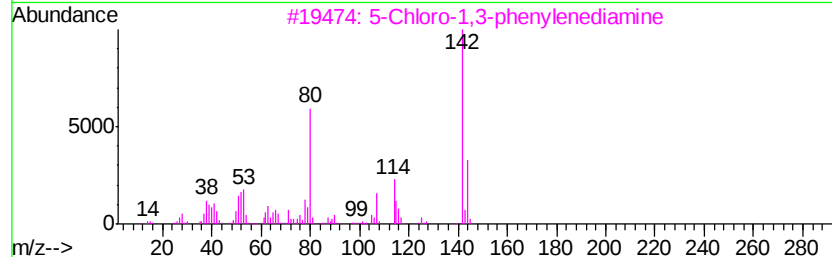
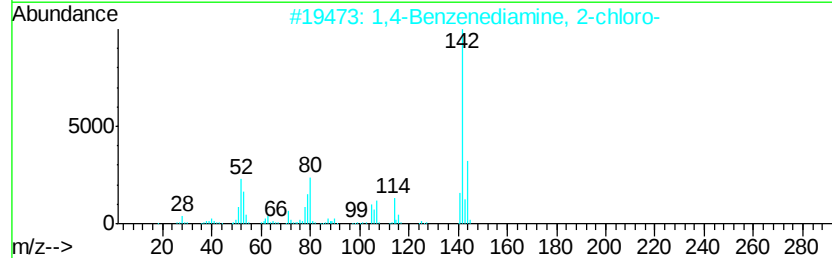
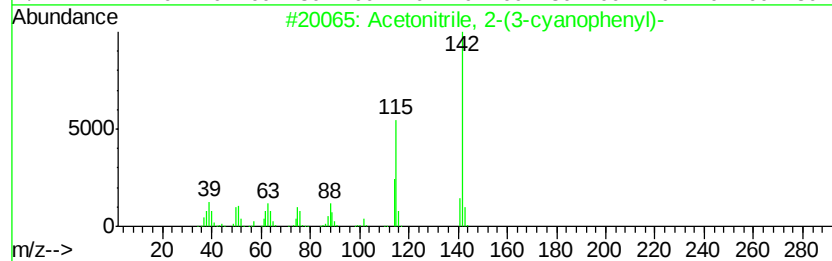
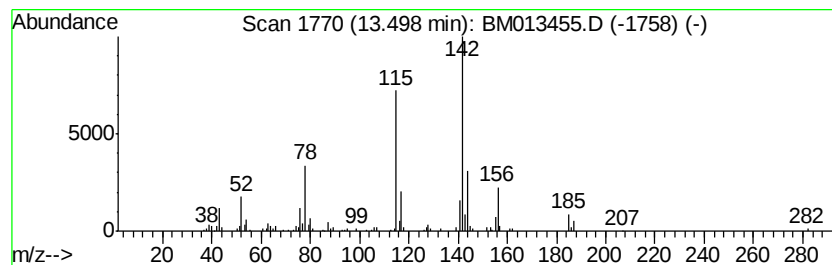
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM113017.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 unknown-02 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.50	2.93 ng/ul	362749	Acenaphthene-d10	14.32
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Acetonitrile, 2-(3-cyanophenyl)-	142 C9H6N2	016532-78-8 47
2		1,4-Benzenediamine, 2-chloro-	142 C6H7ClN2	000615-66-7 38
3		5-Chloro-1,3-phenylenediamine	142 C6H7ClN2	033786-89-9 38
4		Benzeneacetonitrile, 2-cyano-	142 C9H6N2	003759-28-2 38
5		2,6-Dicyanotoluene	142 C9H6N2	002317-22-8 38



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM121517\
Data File : BM013455.D
Acq On : 16 Dec 2017 01:07
Operator : SJ/JU
Sample : I6779-04
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
F9A87

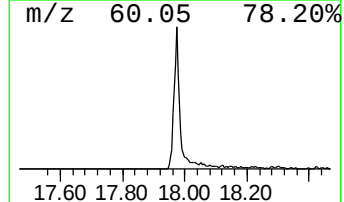
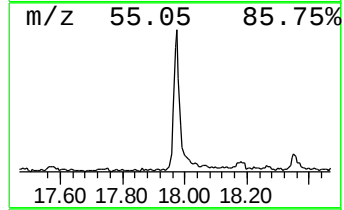
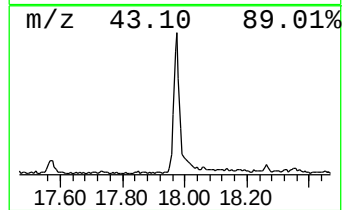
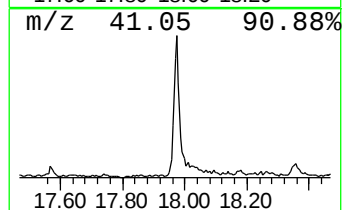
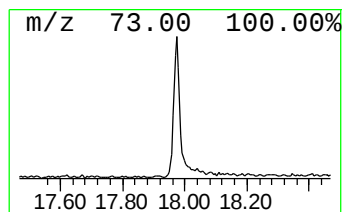
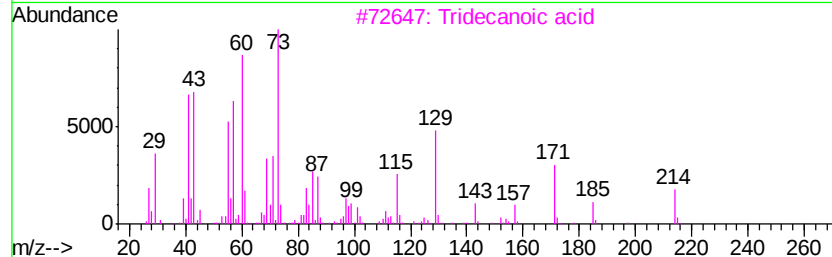
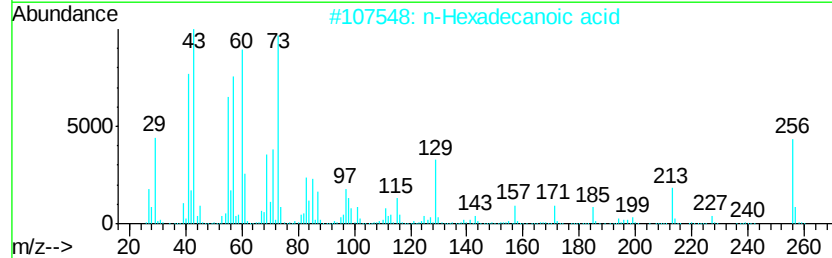
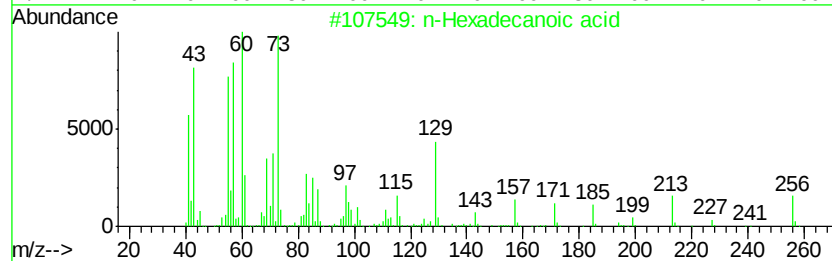
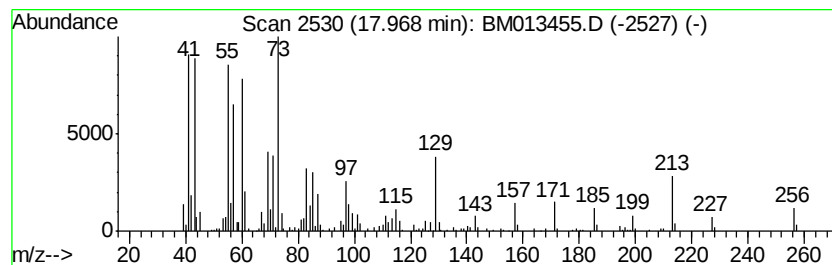
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.97	8.68 ng/ul	1281120	Phenanthrene-d10	17.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94
3		Tridecanoic acid	214	C13H26O2	000638-53-9	91
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	91
5		Pentadecanoic acid	242	C15H30O2	001002-84-2	87



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM121517\
Data File : BM013455.D
Acq On : 16 Dec 2017 01:07
Operator : SJ/JU
Sample : I6779-04
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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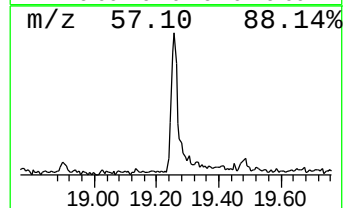
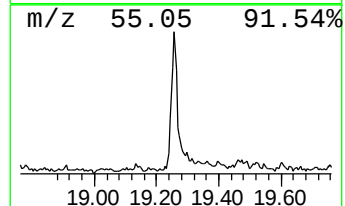
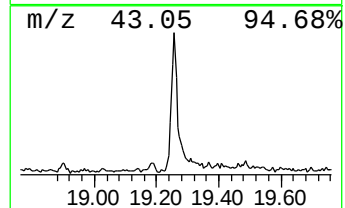
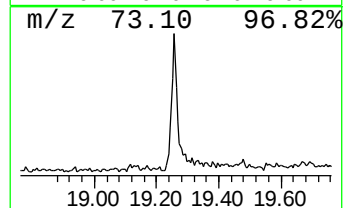
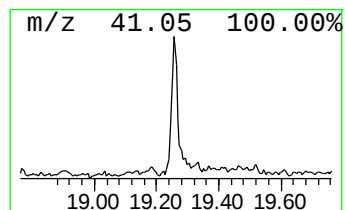
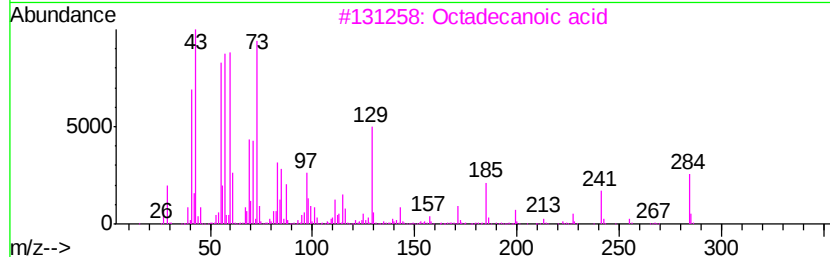
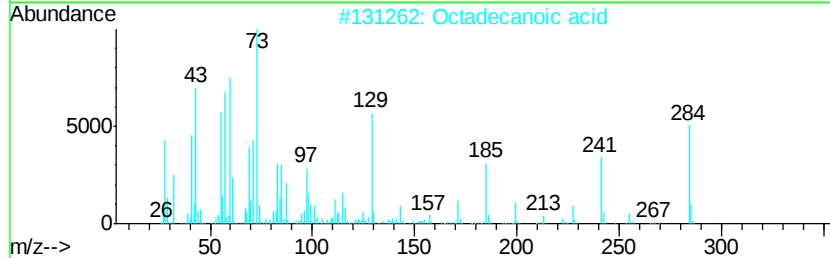
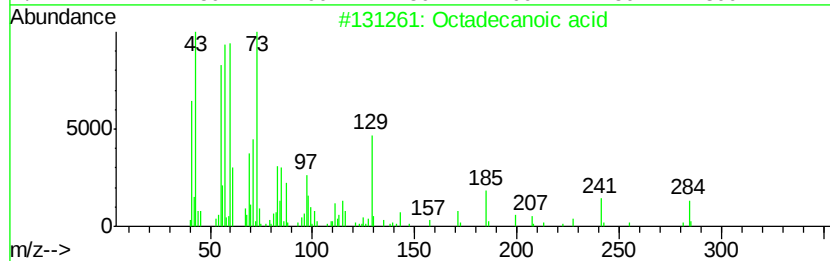
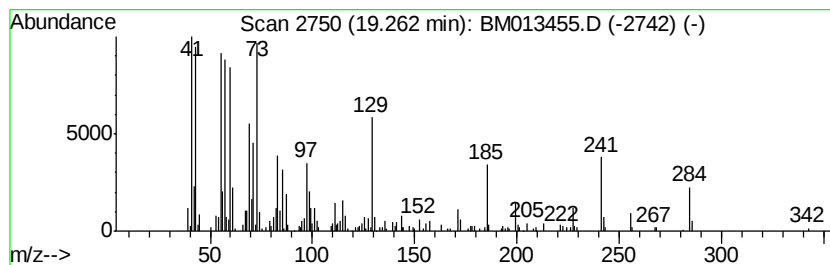
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Octadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.26	3.58 ng/ul	663787	Chrysene-d12	21.26

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Octadecanoic acid	284	C18H36O2	000057-11-4	99
3			Octadecanoic acid	284	C18H36O2	000057-11-4	99
4			Octadecanoic acid	284	C18H36O2	000057-11-4	95
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	90



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM121517\
Data File : BM013455.D
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Operator : SJ/JU
Sample : I6779-04
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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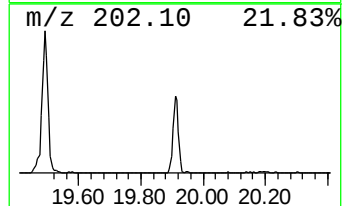
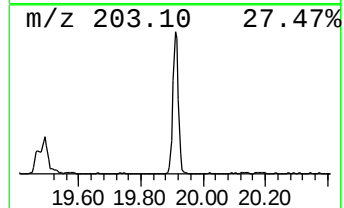
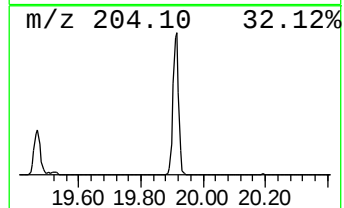
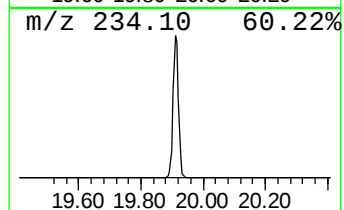
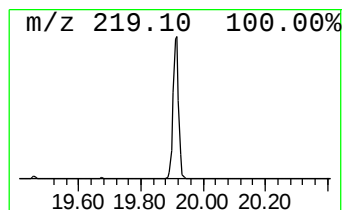
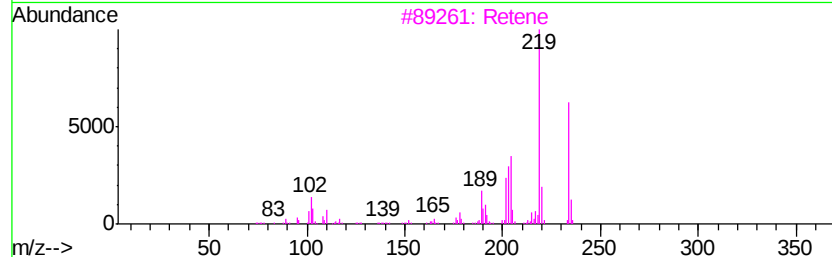
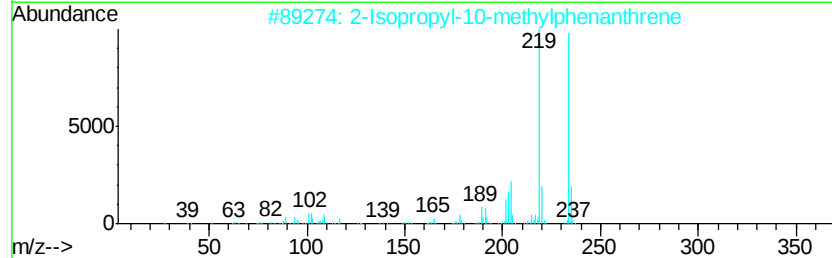
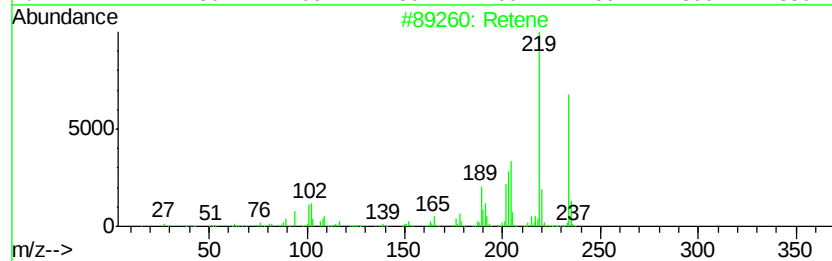
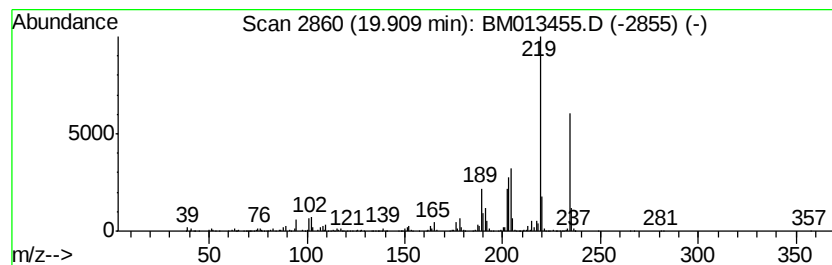
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Retene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.91	14.64 ng/ul	2712950	Chrysene-d12	21.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Retene	234	C18H18	000483-65-8	98
2		2-Isopropyl-10-methylphenanthrene	234	C18H18	066552-97-4	95
3		Retene	234	C18H18	000483-65-8	94
4		Retene	234	C18H18	000483-65-8	93
5		Phenanthrene, 2,4,5,7-tetramethyl-	234	C18H18	007396-38-5	76



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM121517\
Data File : BM013455.D
Acq On : 16 Dec 2017 01:07
Operator : SJ/JU
Sample : I6779-04
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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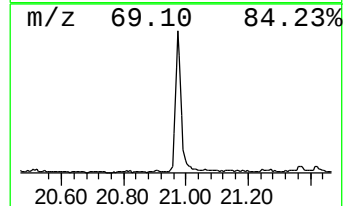
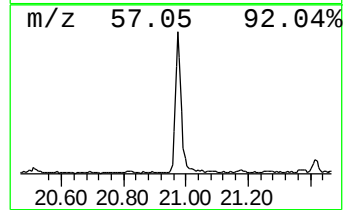
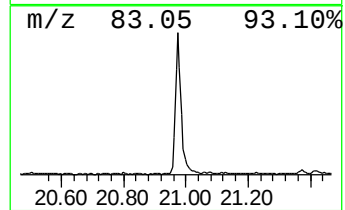
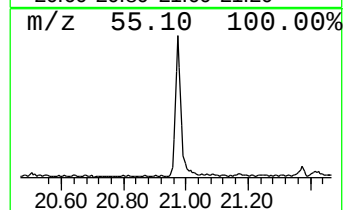
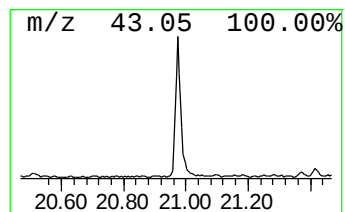
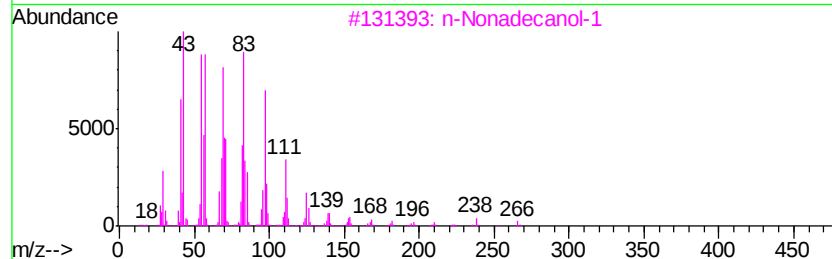
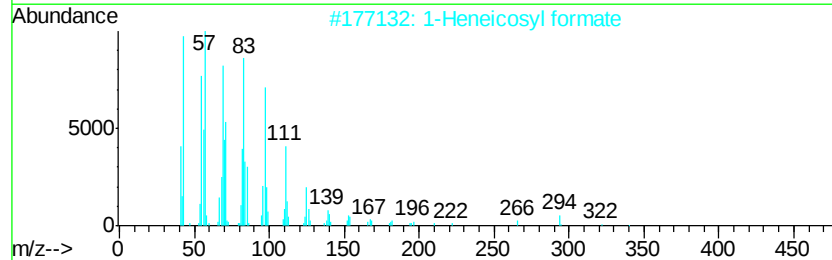
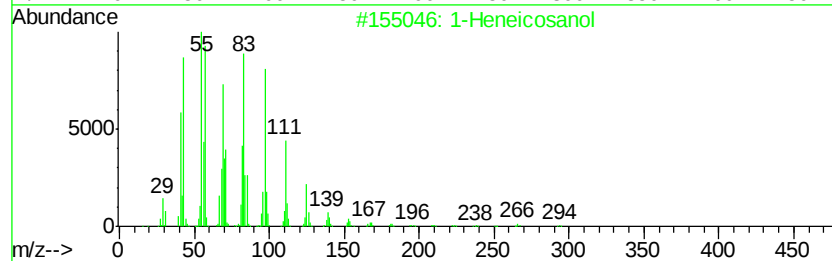
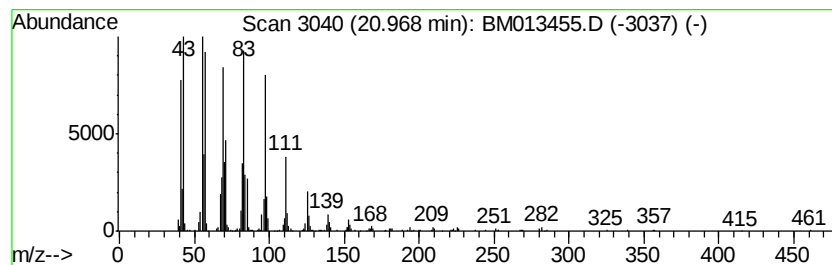
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 1-Heneicosanol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.97	7.08 ng/ul	1312610	Chrysene-d12	21.26

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1	1-Heneicosanol	312	C21H44O	015594-90-8	95
2	1	1-Heneicosyl formate	340	C22H44O2	077899-03-7	94
3	1	n-Nonadecanol-1	284	C19H40O	001454-84-8	94
4	1	Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	94
5	1	Trifluoroacetic acid, pentadecyl...	324	C17H31F3O2	959010-23-2	94



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Misc :
ALS Vial : 12 Sample Multiplier: 1

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ClientSampleId :
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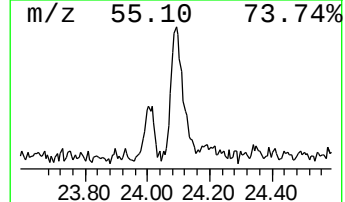
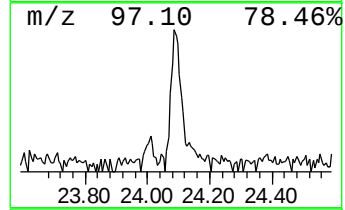
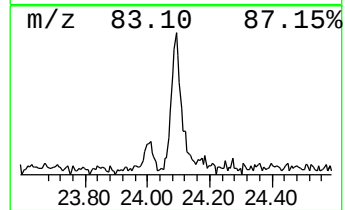
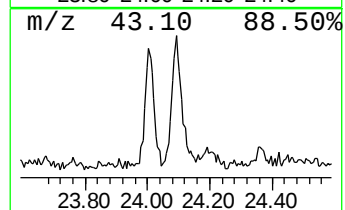
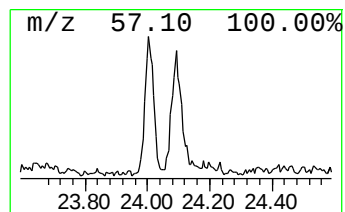
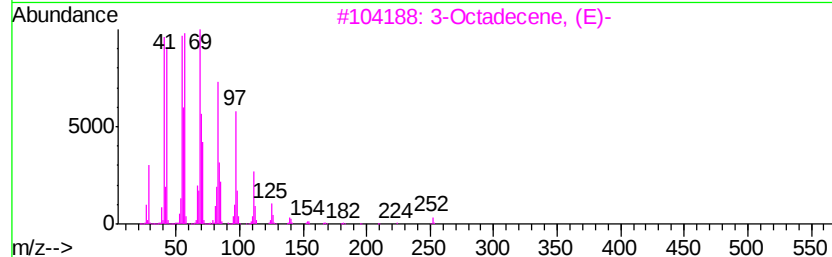
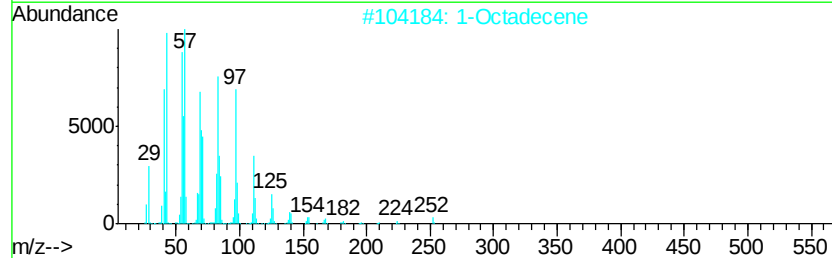
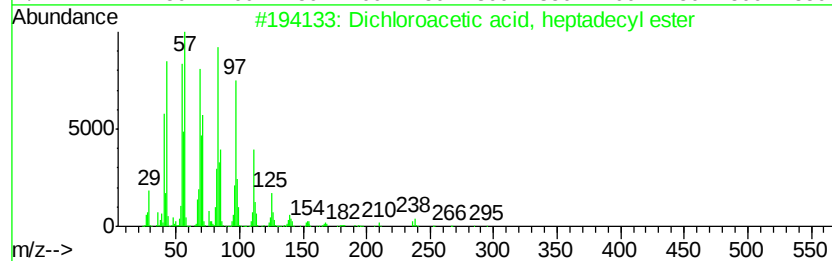
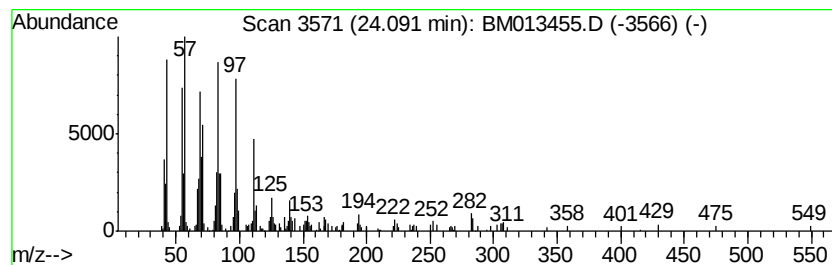
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Dichloroacetic acid, heptad... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.09	3.07 ng/ul	437401	Perylene-d12	23.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	90
2		1-Octadecene	252	C18H36	000112-88-9	90
3		3-Octadecene, (E)-	252	C18H36	007206-19-1	90
4		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	87
5		Cyclohexadecane	224	C16H32	000295-65-8	87



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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
3-Penten-2-one, 4...	4.25	35.4	ng/ul	2267900	1	7.67	1282260	20.0
2-Pentanone, 4-hy...	4.83	9.2	ng/ul	587316	1	7.67	1282260	20.0
unknown-01	11.99	12.2	ng/ul	1122390	2	10.45	1846570	20.0
unknown-02	13.50	2.9	ng/ul	362749	3	14.32	2477650	20.0
n-Hexadecanoic acid	17.97	8.7	ng/ul	1281120	4	17.07	2951280	20.0
Octadecanoic acid	19.26	3.6	ng/ul	663787	5	21.26	3705900	20.0
Retene	19.91	14.6	ng/ul	2712950	5	21.26	3705900	20.0
1-Heneicosanol	20.97	7.1	ng/ul	1312610	5	21.26	3705900	20.0
Dichloroacetic ac...	24.09	3.1	ng/ul	437401	6	23.48	2854080	20.0