

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM012617\
 Data File : BM008900.D
 Acq On : 27 Jan 2017 06:57
 Operator : SJ/MA
 Sample : I1426-19
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BDMY7

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\SOM02.2-EPA-BM012317.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	2.716	20	22	27	rVB4	16043	20731	1.16%	0.082%
2	2.834	38	42	47	rVB	64035	78618	4.39%	0.312%
3	2.916	52	56	60	rVV2	75037	98723	5.51%	0.392%
4	2.963	60	64	71	rVB	249630	283074	15.81%	1.124%
5	3.046	73	78	82	rBV2	104013	132715	7.41%	0.527%
6	3.387	132	136	139	rBV3	22013	28008	1.56%	0.111%
7	3.869	213	218	228	rVB	116904	167482	9.36%	0.665%
8	5.392	472	477	483	rVB2	19985	28654	1.60%	0.114%
9	5.528	495	500	507	rVB2	22593	40981	2.29%	0.163%
10	6.175	606	610	615	rBV4	12617	19723	1.10%	0.078%
11	6.257	620	624	637	rVB7	18460	61594	3.44%	0.244%
12	6.416	645	651	659	rBV3	40039	67245	3.76%	0.267%
13	6.516	659	668	672	rBV3	44086	84138	4.70%	0.334%
14	6.657	687	692	696	rVB4	12208	22147	1.24%	0.088%
15	6.716	696	702	707	rBV4	26986	44413	2.48%	0.176%
16	6.898	728	733	738	rVB4	15393	23624	1.32%	0.094%
17	7.075	754	763	776	rBV2	472384	853230	47.66%	3.386%
18	7.251	787	793	802	rVB	272934	437367	24.43%	1.736%
19	7.451	820	827	835	rBV	449378	696056	38.88%	2.763%
20	7.928	897	908	914	rBV	463829	763852	42.67%	3.032%
21	8.616	1019	1025	1032	rBV	454218	709943	39.66%	2.818%
22	9.086	1098	1105	1114	rVB	453822	750536	41.92%	2.979%
23	9.463	1162	1169	1173	rVB4	13729	22514	1.26%	0.089%
24	9.810	1221	1228	1235	rBV	395975	644640	36.01%	2.559%
25	10.339	1311	1318	1325	rBV	526133	846575	47.29%	3.360%
26	10.727	1378	1384	1391	rVB	579444	945735	52.83%	3.754%
27	10.863	1401	1407	1417	rBV2	307129	519419	29.01%	2.062%
28	13.968	1927	1935	1939	rBV	841951	1166145	65.14%	4.628%
29	14.010	1939	1942	1947	rVB	169873	220756	12.33%	0.876%
30	14.257	1978	1984	1991	rVB	821889	1176081	65.69%	4.668%
31	14.439	2008	2015	2018	rBV3	15992	22375	1.25%	0.089%
32	14.562	2029	2036	2043	rVB	805157	1150653	64.27%	4.567%
33	14.745	2059	2067	2076	rBV	470290	685311	38.28%	2.720%
34	14.957	2094	2103	2107	rBV3	67029	92613	5.17%	0.368%

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35	15.392	2172	2177	2183	rVB	38670	53937	3.01%	0.214%
36	15.557	2198	2205	2211	rBV	1019772	1506230	84.14%	5.978%
37	15.662	2216	2223	2235	rBV	413429	561928	31.39%	2.230%
38	15.792	2240	2245	2249	rVB5	20084	35150	1.96%	0.140%
39	16.251	2317	2323	2331	rBV2	60586	100293	5.60%	0.398%
40	16.845	2416	2424	2426	rBV6	16202	27131	1.52%	0.108%
41	17.045	2454	2458	2462	rBV	39442	44324	2.48%	0.176%
42	17.092	2462	2466	2473	rVB8	14565	26114	1.46%	0.104%
43	17.309	2495	2503	2510	rBV2	944273	1385744	77.41%	5.500%
44	17.403	2513	2519	2526	rVV	1125012	1630857	91.10%	6.473%
45	18.174	2645	2650	2656	rBV2	130966	173622	9.70%	0.689%
46	19.327	2840	2846	2854	rVB8	18175	46626	2.60%	0.185%
47	19.450	2861	2867	2871	rBV4	89625	119058	6.65%	0.473%
48	19.692	2903	2908	2914	rBV	1298797	1790244	100.00%	7.105%
49	21.162	3155	3158	3164	rBV5	106674	142258	7.95%	0.565%
50	21.474	3206	3211	3216	rBV	1259981	1598067	89.27%	6.343%
51	23.680	3580	3586	3594	rVB2	840335	1655942	92.50%	6.572%
52	23.832	3606	3612	3620	rVB2	742545	1392376	77.78%	5.526%

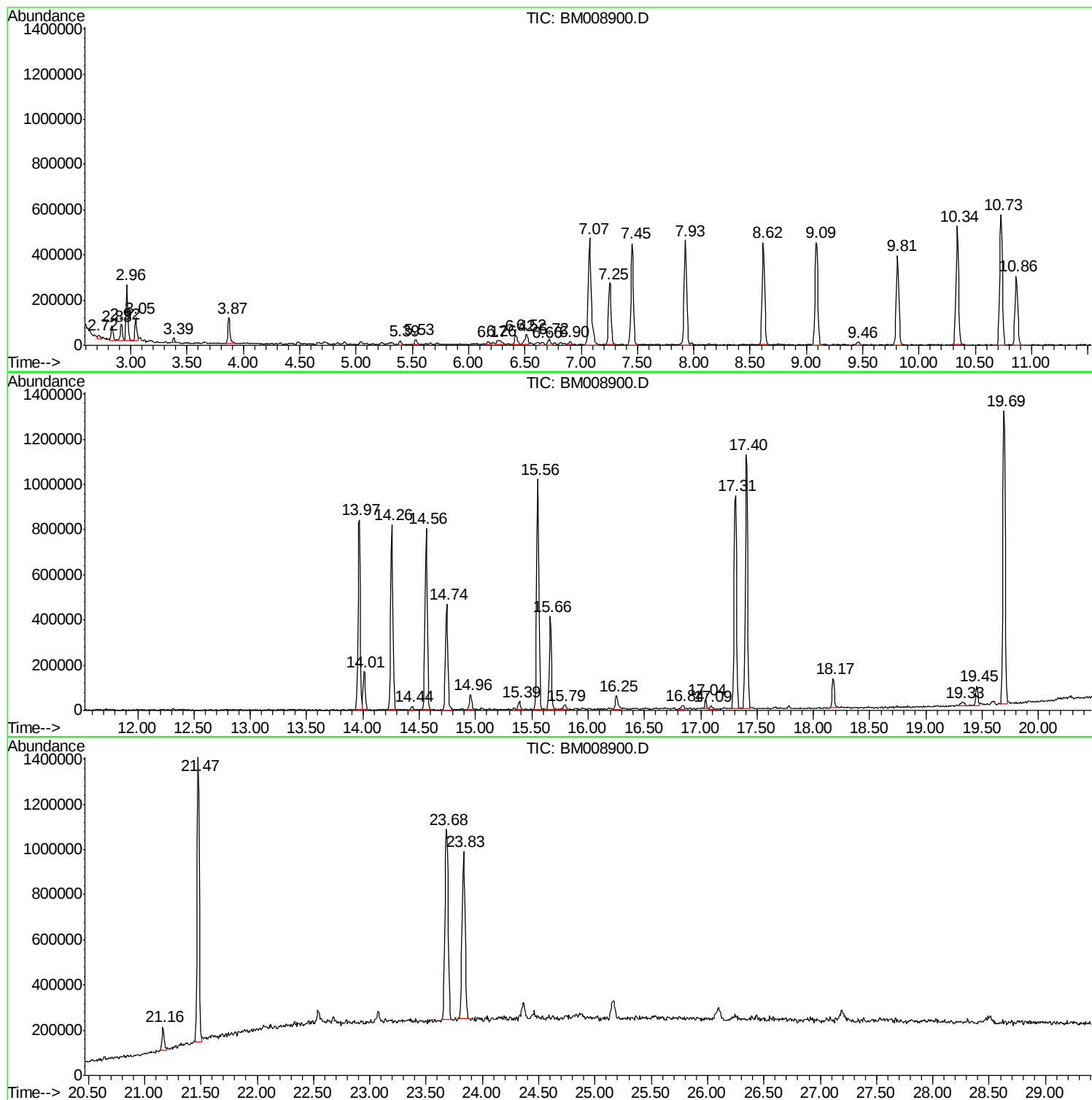
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM012317.M
Quant Title : SVOA CALIBRATION

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



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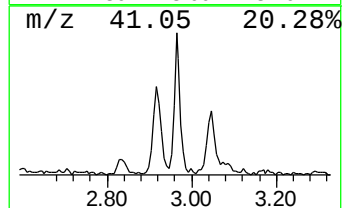
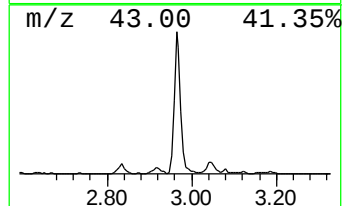
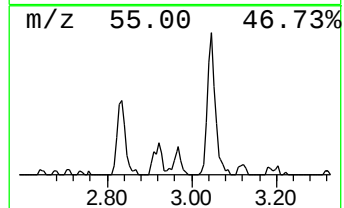
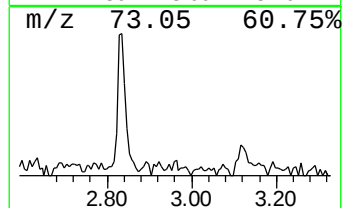
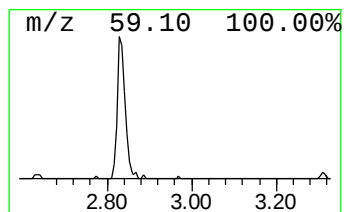
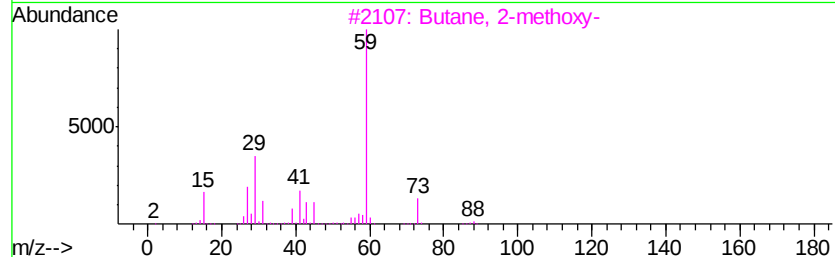
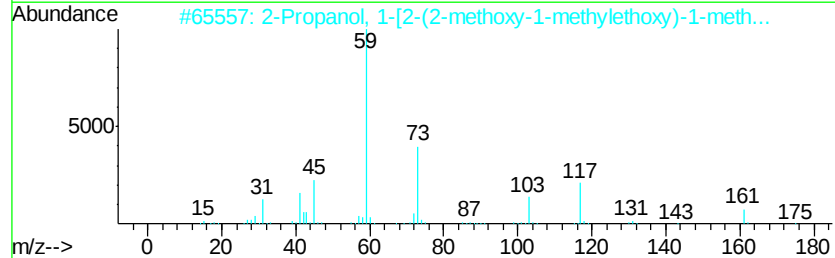
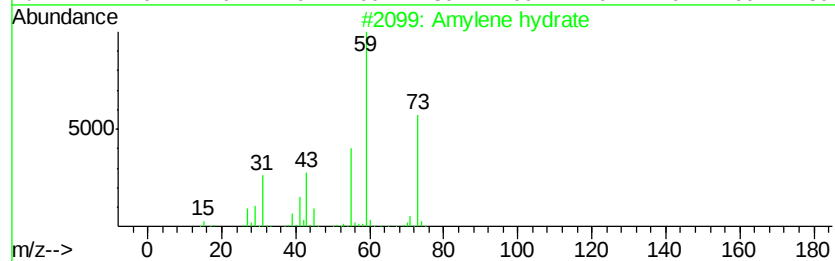
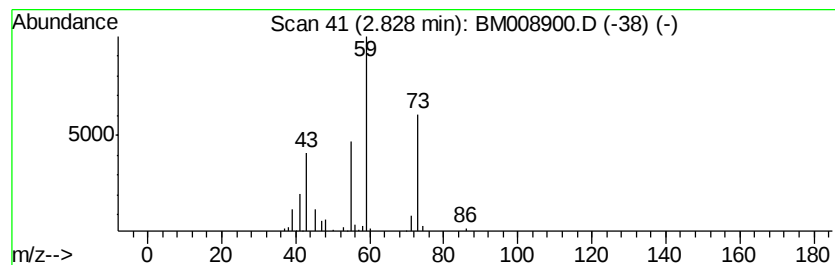
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM012317.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Amylene hydrate Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.83	2.06 ng/ul	78618	1,4-Dichlorobenzene-d4	7.92

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Amylene hydrate	88	C5H12O	000075-85-4	64
2			2-Propanol, 1-[2-(2-methoxy-1-me...	206	C10H22O4	020324-33-8	59
3			Butane, 2-methoxy-	88	C5H12O	006795-87-5	53
4			2,5,8,11-Tetraoxatetradecan-13-o...	264	C13H28O5	020324-34-9	45
5			Hydrazine, 1,2-diethyl-	88	C4H12N2	001615-80-1	40



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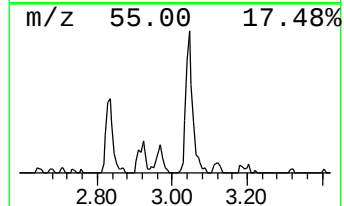
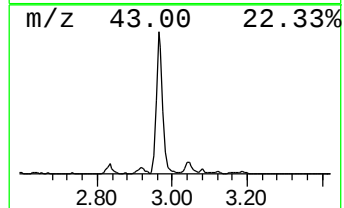
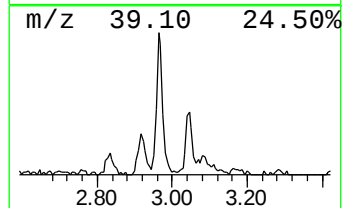
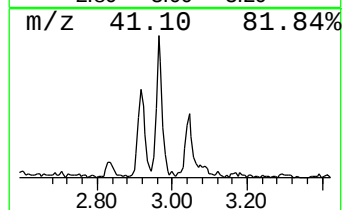
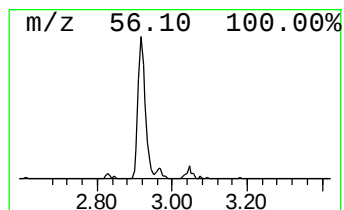
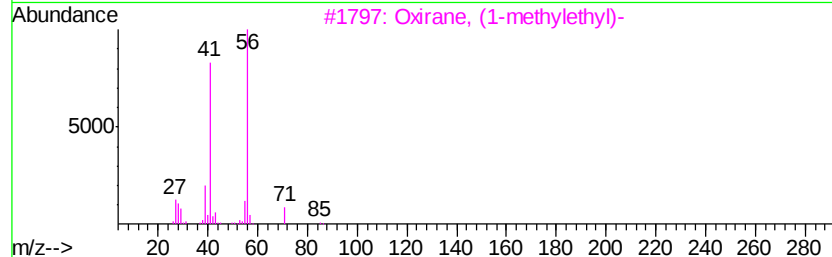
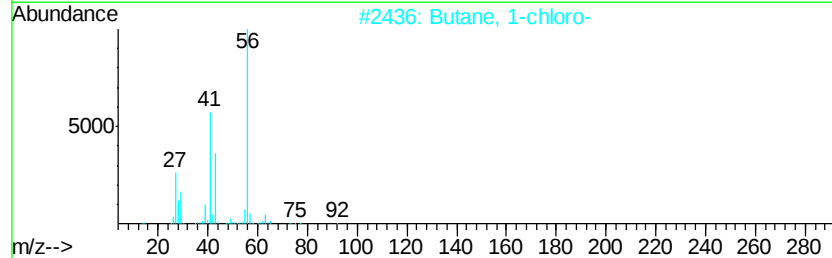
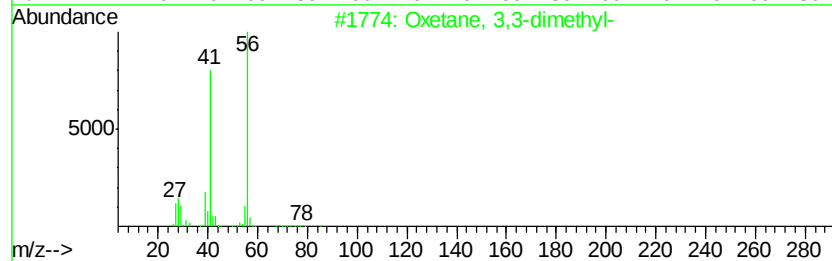
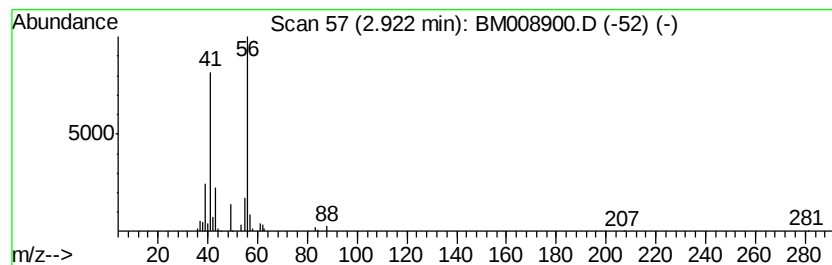
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Oxetane, 3,3-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.92	2.58 ng/ul	98723	1,4-Dichlorobenzene-d4	7.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Oxetane, 3,3-dimethyl-	86	C5H10O	006921-35-3	64
2		Butane, 1-chloro-	92	C4H9Cl	000109-69-3	64
3		Oxirane, (1-methylethyl)-	86	C5H10O	001438-14-8	64
4		Butane, 1-chloro-	92	C4H9Cl	000109-69-3	53
5		3,4-Dimethyldihydrofuran-2,5-dione	128	C6H8O3	007475-92-5	50



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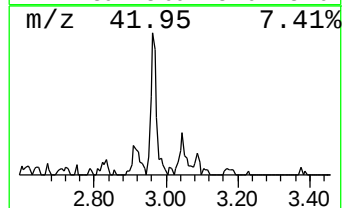
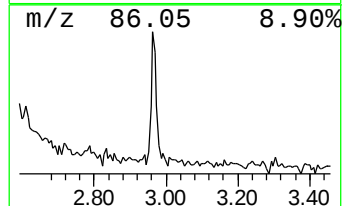
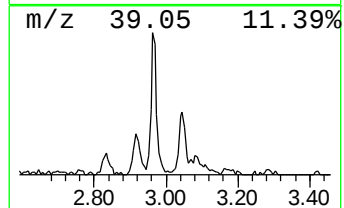
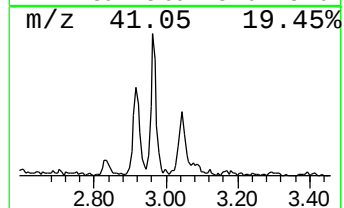
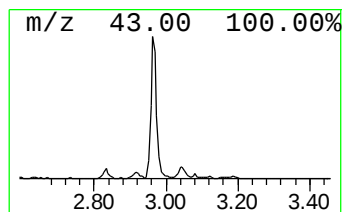
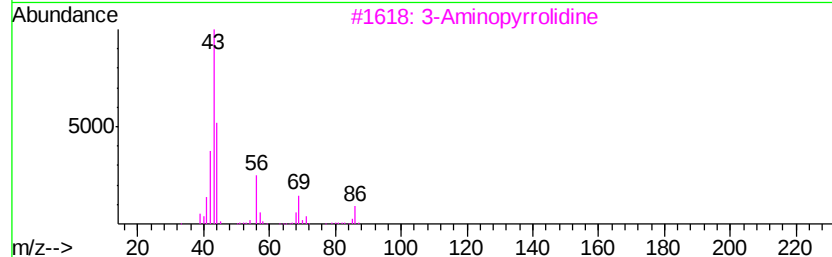
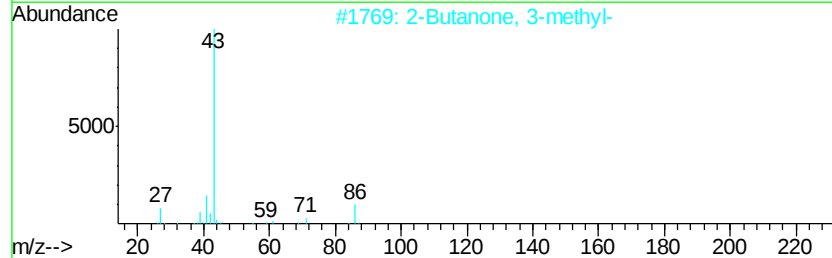
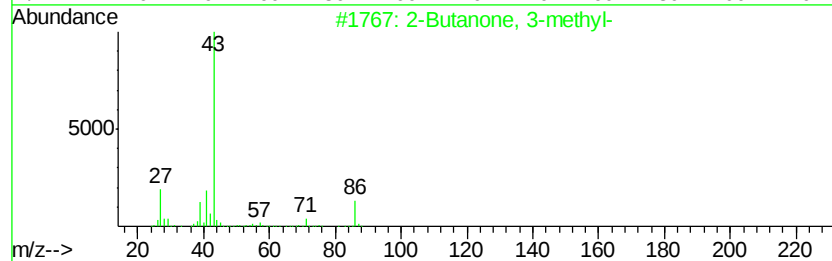
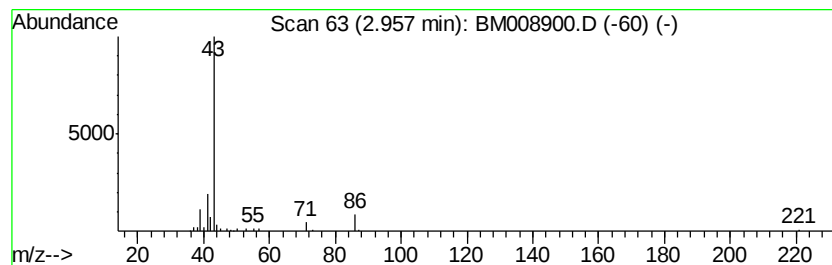
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.96	7.41 ng/ul	283074	1,4-Dichlorobenzene-d4	7.92

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butanone, 3-methyl-			86	C5H10O	000563-80-4	47
2	2-Butanone, 3-methyl-			86	C5H10O	000563-80-4	9
3	3-Aminopyrrolidine			86	C4H10N2	079286-79-6	9
4	Azetidine, 1-nitroso-			86	C3H6N2O	015216-10-1	9
5	2-Pentanone			86	C5H10O	000107-87-9	9



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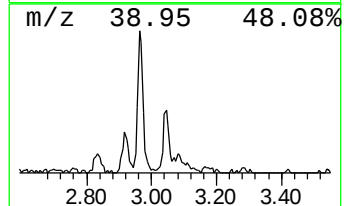
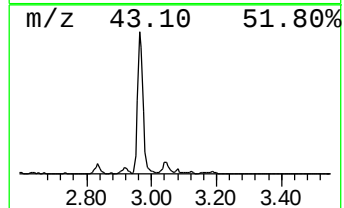
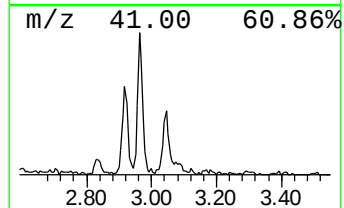
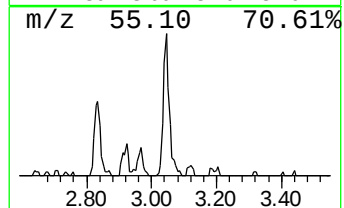
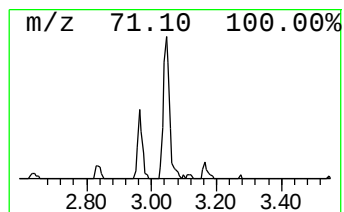
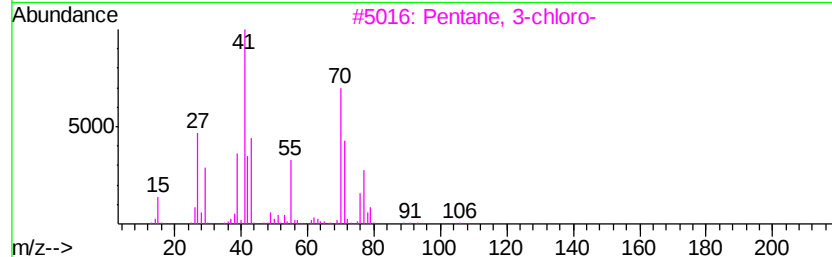
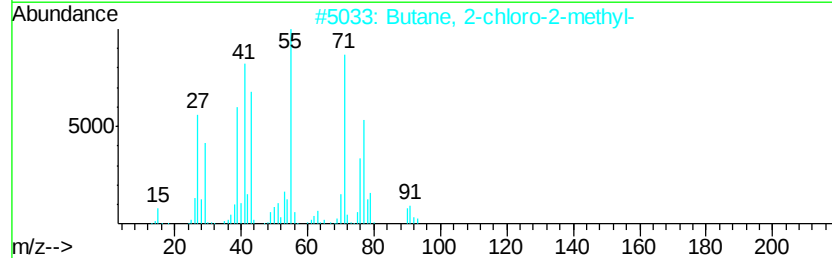
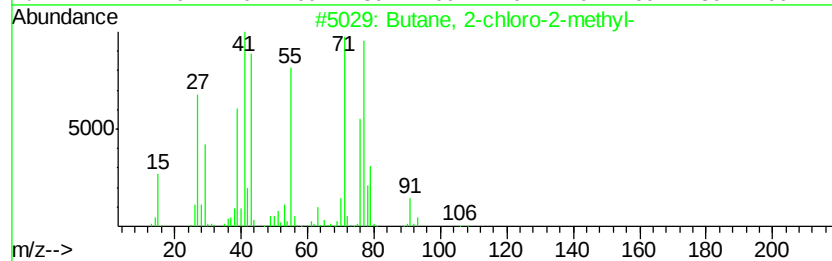
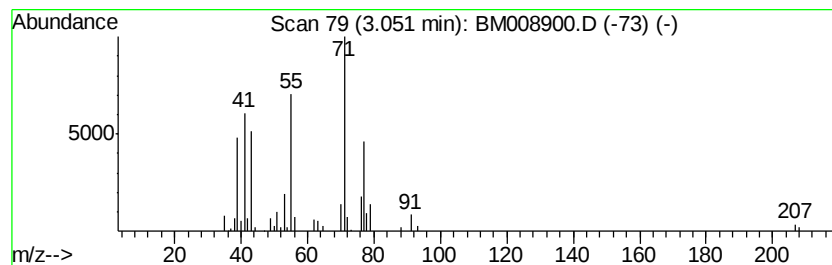
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 4 unknown-02 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.05	3.47 ng/ul	132715	1,4-Dichlorobenzene-d4	7.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-chloro-2-methyl-	106	C5H11Cl	000594-36-5	40
2		Butane, 2-chloro-2-methyl-	106	C5H11Cl	000594-36-5	38
3		Pentane, 3-chloro-	106	C5H11Cl	000616-20-6	38
4		Trans-3-methylpent-3-ene-5-ol	100	C6H12O	030801-95-7	23
5		1-Butene, 3-methyl-	70	C5H10	000563-45-1	22



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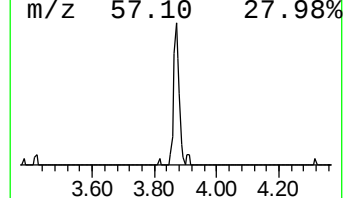
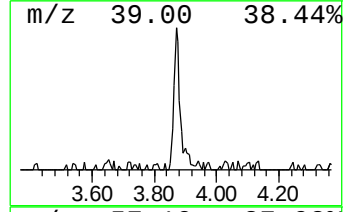
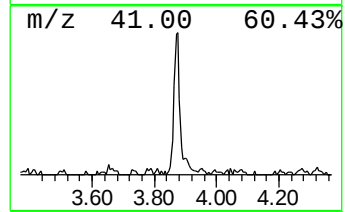
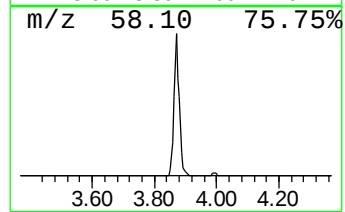
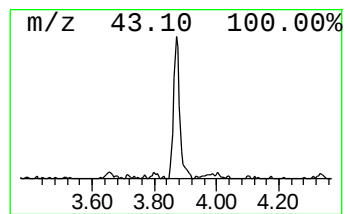
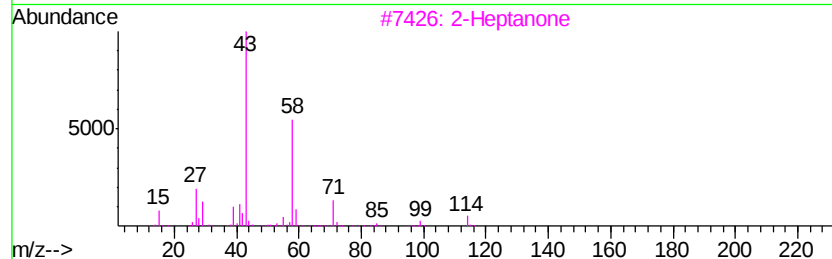
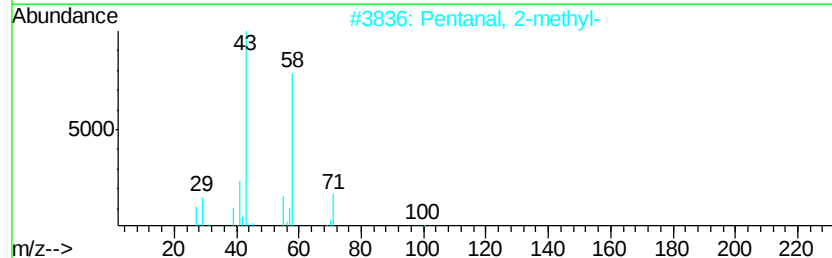
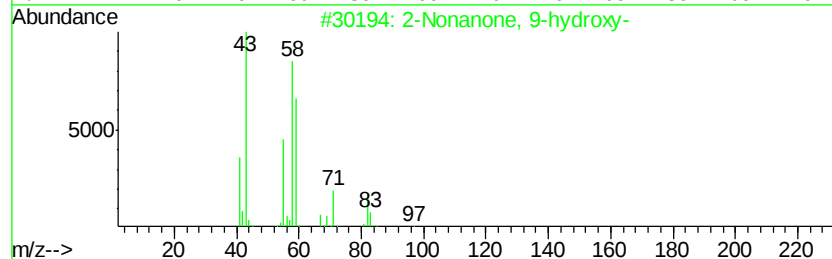
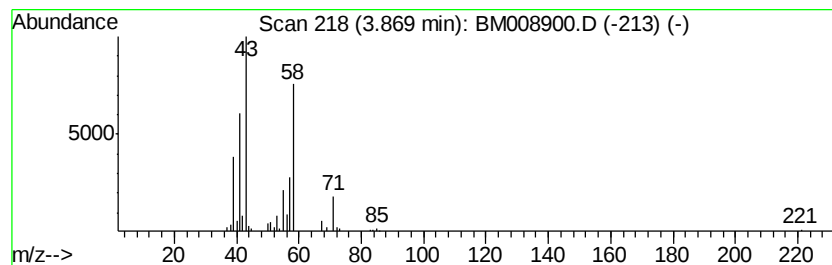
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 unknown-03 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.87	4.39 ng/ul	167482	1,4-Dichlorobenzene-d4	7.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Nonanone, 9-hydroxy-	158	C9H18O2	025368-56-3	43
2		Pentanal, 2-methyl-	100	C6H12O	000123-15-9	40
3		2-Heptanone	114	C7H14O	000110-43-0	40
4		S-[2-[N,N-Dimethylamino]ethyl]mo...	261	C10H19N3O3S	1000212-14-3	39
5		2-Hexanone, 5-methyl-	114	C7H14O	000110-12-3	36



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM012617\
Data File : BM008900.D
Acq On : 27 Jan 2017 06:57
Operator : SJ/MA
Sample : I1426-19
Misc :
ALS Vial : 23 Sample Multiplier: 1

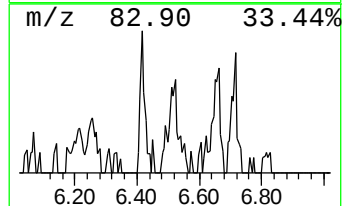
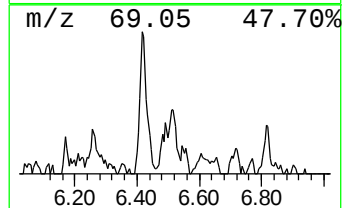
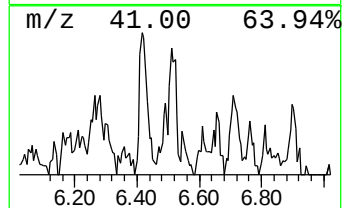
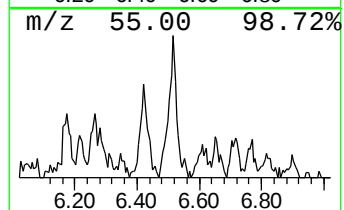
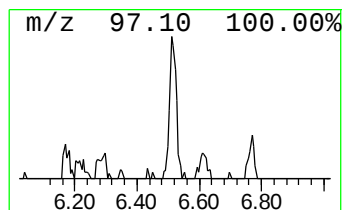
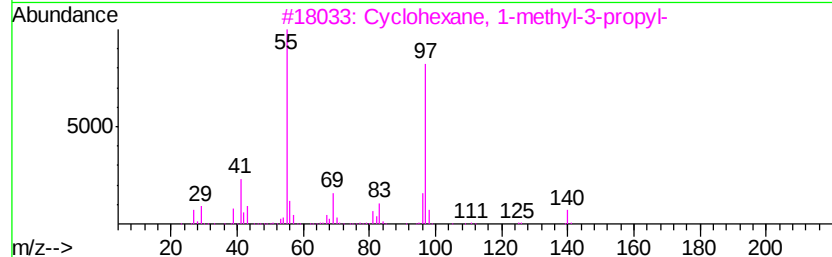
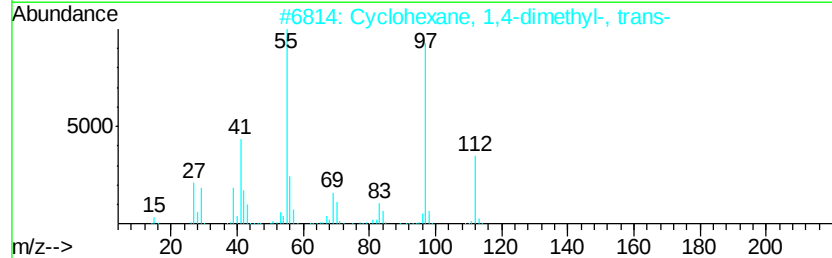
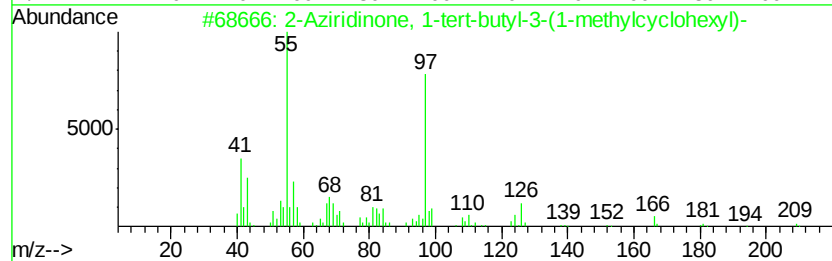
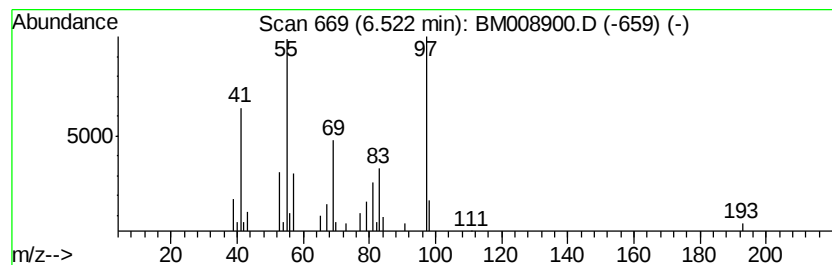
Instrument :
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ClientSampleId :
BDMY7

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM012317.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 2-Aziridinone, 1-tert-butyl... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.		
6.52	2.20 ng/ul	84138	1,4-Dichlorobenzene-d4	7.92		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Aziridinone, 1-tert-butyl-3-(1...	209	C13H23NO	024161-49-7	50	
2	Cyclohexane, 1,4-dimethyl-, trans-	112	C8H16	002207-04-7	42	
3	Cyclohexane, 1-methyl-3-propyl-	140	C10H20	004291-80-9	40	
4	Cyclopropane, 3-chloro-1,1,2,2-t...	132	C7H13Cl	014123-41-2	40	
5	Oxalic acid, butyl cyclohexylmet...	242	C13H22O4	1000309-68-2	40	



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM012617\
Data File : BM008900.D
Acq On : 27 Jan 2017 06:57
Operator : SJ/MA
Sample : I1426-19
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BDMY7

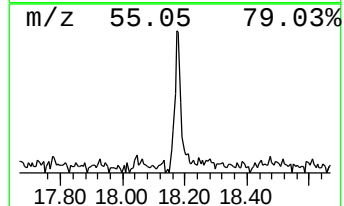
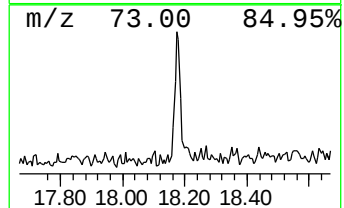
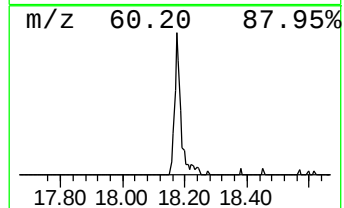
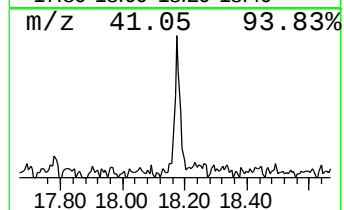
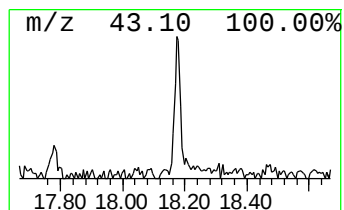
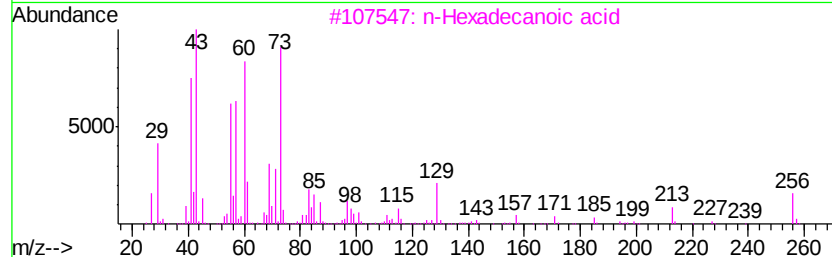
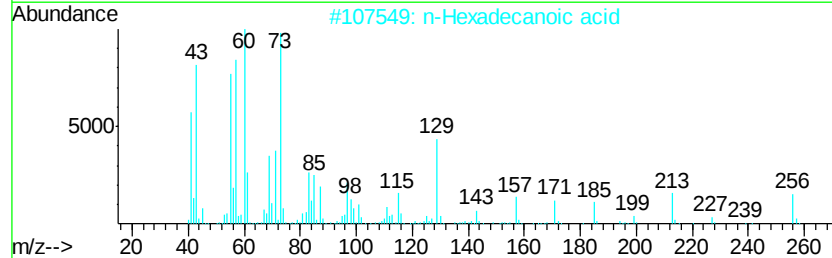
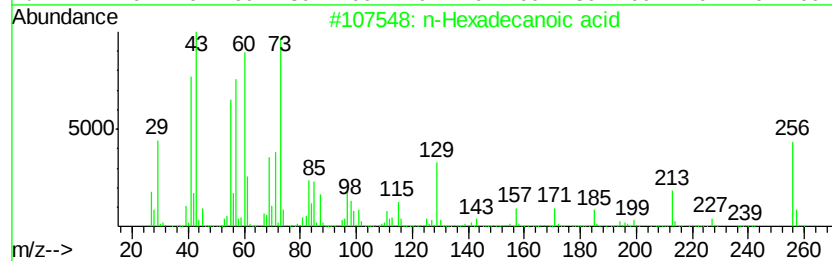
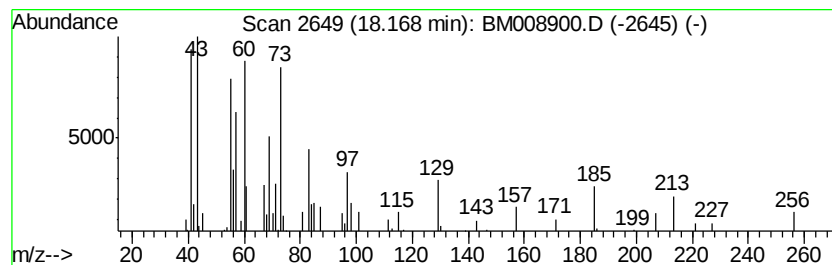
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.17	2.51 ng/ul	173622	Phenanthrene-d10	17.31

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	89
3		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	76
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	52
5		Pentadecanoic acid	242	C15H30O2	001002-84-2	50



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM012617\
Data File : BM008900.D
Acq On : 27 Jan 2017 06:57
Operator : SJ/MA
Sample : I1426-19
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BDMY7

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM012317.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
Amylene hydrate	2.83	2.1	ng/ul	78618	1	7.92	763852	20.0
Oxetane, 3,3-dime...	2.92	2.6	ng/ul	98723	1	7.92	763852	20.0
unknown-01	2.96	7.4	ng/ul	283074	1	7.92	763852	20.0
unknown-02	3.05	3.5	ng/ul	132715	1	7.92	763852	20.0
unknown-03	3.87	4.4	ng/ul	167482	1	7.92	763852	20.0
2-Aziridinone, 1-...	6.52	2.2	ng/ul	84138	1	7.92	763852	20.0
n-Hexadecanoic acid	18.17	2.5	ng/ul	173622	4	17.31	1385740	20.0