

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM041415\
 Data File : BM000987.D
 Acq On : 14 Apr 2015 17:14
 Operator : TP/IZ
 Sample : G1858-11
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 P001-SD008-0006-02

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 0.1 % 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\SOM01.2-EPA-BM041415.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	2.734	3	8	18	rBV	768203	1486481	23.88%	1.467%
2	2.816	18	22	28	rVB	600074	748827	12.03%	0.739%
3	3.357	111	114	122	rBV	45522	68881	1.11%	0.068%
4	3.628	157	160	165	rVB5	31469	40339	0.65%	0.040%
5	4.028	226	228	235	rVB7	15731	25650	0.41%	0.025%
6	4.710	340	344	353	rBV	46393	82523	1.33%	0.081%
7	5.628	495	500	503	rBV7	11631	21551	0.35%	0.021%
8	6.781	689	696	711	rVB	1243599	2151643	34.57%	2.123%
9	6.904	712	717	718	rVV3	25107	34920	0.56%	0.034%
10	6.940	718	723	734	rVV	954768	1466686	23.56%	1.447%
11	7.134	749	756	767	rBV	1233860	1890613	30.37%	1.866%
12	7.228	767	772	777	rVV4	30011	48874	0.79%	0.048%
13	7.528	819	823	828	rVV6	25991	54744	0.88%	0.054%
14	7.598	828	835	843	rVV	1406949	2222813	35.71%	2.193%
15	7.810	866	871	872	rBV4	18041	21118	0.34%	0.021%
16	8.051	908	912	917	rBV8	18812	35047	0.56%	0.035%
17	8.310	948	956	963	rBV	1550166	2513101	40.37%	2.480%
18	8.757	1024	1032	1041	rBV	1095166	1749040	28.10%	1.726%
19	9.181	1099	1104	1111	rVB5	22620	34739	0.56%	0.034%
20	9.475	1145	1154	1166	rBV	827435	1363520	21.91%	1.346%
21	10.010	1237	1245	1254	rBV	1508751	2487689	39.97%	2.455%
22	10.181	1268	1274	1281	rVB	12029	24464	0.39%	0.024%
23	10.381	1301	1308	1315	rBV	1848774	3118532	50.10%	3.077%
24	10.528	1324	1333	1345	rBV	1271011	2258650	36.29%	2.229%
25	11.469	1488	1493	1496	rVV7	17785	28296	0.45%	0.028%
26	11.710	1525	1534	1540	rVV	348333	569307	9.15%	0.562%
27	11.980	1574	1580	1586	rVV4	32387	50020	0.80%	0.049%
28	12.057	1588	1593	1602	rVB6	19143	29582	0.48%	0.029%
29	12.280	1627	1631	1635	rVB6	18888	27845	0.45%	0.027%
30	12.598	1679	1685	1690	rBV2	54521	92961	1.49%	0.092%
31	12.857	1722	1729	1737	rVB	75240	119687	1.92%	0.118%
32	13.080	1761	1767	1772	rVB4	48345	77465	1.24%	0.076%
33	13.157	1772	1780	1787	rBV	199627	342846	5.51%	0.338%
34	13.669	1862	1867	1872	rVV	2982247	4085933	65.64%	4.032%

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Integration Parameters: LSCINT.P

Integrator: RTE
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35	13.716	1872	1875	1885	rVB	942778	1304047	20.95%	1.287%
36	13.945	1905	1914	1927	rBV	2897970	4338131	69.69%	4.281%
37	14.163	1946	1951	1956	rVB2	127971	173785	2.79%	0.171%
38	14.257	1961	1967	1976	rVV2	2512829	3859897	62.01%	3.809%
39	14.469	1995	2003	2011	rBV	1006485	1659901	26.67%	1.638%
40	14.533	2011	2014	2020	rVB3	115628	189521	3.04%	0.187%
41	14.663	2033	2036	2039	rBV5	11868	19904	0.32%	0.020%
42	14.698	2039	2042	2045	rVB4	21447	28742	0.46%	0.028%
43	14.792	2054	2058	2063	rVB7	26600	40098	0.64%	0.040%
44	14.927	2076	2081	2084	rVB7	14083	22646	0.36%	0.022%
45	15.033	2094	2099	2106	rBV	86095	141767	2.28%	0.140%
46	15.116	2109	2113	2118	rVB	151166	190886	3.07%	0.188%
47	15.251	2130	2136	2143	rBV	3700995	5462306	87.75%	5.390%
48	15.304	2143	2145	2150	rVB6	25458	28930	0.46%	0.029%
49	15.380	2152	2158	2171	rBV	851180	1202945	19.33%	1.187%
50	15.492	2174	2177	2179	rBV3	46425	62324	1.00%	0.062%
51	15.521	2180	2182	2187	rVB2	61465	84296	1.35%	0.083%
52	15.586	2190	2193	2194	rBV	47372	46220	0.74%	0.046%
53	15.621	2194	2199	2206	rVB	2415258	3225506	51.82%	3.183%
54	15.839	2228	2236	2244	rBV	35513	88654	1.42%	0.087%
55	15.980	2256	2260	2262	rBV	174299	228516	3.67%	0.225%
56	16.004	2262	2264	2269	rVB	155477	181028	2.91%	0.179%
57	16.257	2302	2307	2311	rVV	110418	152909	2.46%	0.151%
58	16.345	2313	2322	2327	rVV	567283	963990	15.49%	0.951%
59	16.398	2327	2331	2333	rVV2	113288	199842	3.21%	0.197%
60	16.421	2333	2335	2339	rVV	136582	185149	2.97%	0.183%
61	16.457	2339	2341	2343	rVV3	38763	46643	0.75%	0.046%
62	16.480	2343	2345	2353	rVB5	43231	62829	1.01%	0.062%
63	16.557	2353	2358	2361	rBV7	19503	28931	0.46%	0.029%
64	16.768	2390	2394	2399	rVV	155066	212012	3.41%	0.209%
65	16.821	2399	2403	2406	rVV	77372	113263	1.82%	0.112%
66	16.862	2406	2410	2416	rVB	243310	294348	4.73%	0.290%
67	16.968	2424	2428	2430	rBV	254166	334378	5.37%	0.330%
68	17.004	2430	2434	2445	rVB	3199014	4381765	70.40%	4.324%
69	17.104	2445	2451	2457	rBV	4415254	6024989	96.79%	5.945%
70	17.204	2463	2468	2472	rVB8	43858	70085	1.13%	0.069%
71	17.274	2473	2480	2490	rVB	521817	939220	15.09%	0.927%

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72	17.392	2497	2500	2504	rBV4	48913	50152	0.81%	0.049%
73	17.509	2516	2520	2523	rVB	143594	170778	2.74%	0.169%
74	17.680	2547	2549	2554	rVB5	21908	25712	0.41%	0.025%
75	17.792	2564	2568	2573	rVB8	21707	34096	0.55%	0.034%
76	17.909	2583	2588	2597	rBV	888480	1384624	22.24%	1.366%
77	18.198	2634	2637	2641	rBV	126487	147026	2.36%	0.145%
78	18.668	2715	2717	2721	rVB5	24327	21583	0.35%	0.021%
79	18.768	2731	2734	2737	rBV4	36895	53237	0.86%	0.053%
80	18.839	2743	2746	2753	rVB	130369	166466	2.67%	0.164%
81	19.045	2776	2781	2783	rBV	63316	95044	1.53%	0.094%
82	19.098	2787	2790	2793	rVB4	81779	99542	1.60%	0.098%
83	19.198	2800	2807	2818	rBV2	293119	607678	9.76%	0.600%
84	19.292	2821	2823	2827	rVB4	43650	53088	0.85%	0.052%
85	19.403	2836	2842	2850	rBV2	4448490	6224511	100.00%	6.142%
86	19.462	2850	2852	2856	rVB5	37082	40406	0.65%	0.040%
87	19.933	2928	2932	2934	rBV4	56942	91237	1.47%	0.090%
88	19.962	2934	2937	2941	rVB2	102624	123695	1.99%	0.122%
89	20.339	2997	3001	3005	rVB3	77624	99723	1.60%	0.098%
90	20.456	3016	3021	3025	rBV4	83791	128119	2.06%	0.126%
91	20.915	3094	3099	3110	rBV	4602428	5773429	92.75%	5.697%
92	21.203	3142	3148	3158	rBV	3967015	5101381	81.96%	5.034%
93	21.321	3164	3168	3171	rBV2	261222	316905	5.09%	0.313%
94	21.780	3243	3246	3247	rBV2	203602	206189	3.31%	0.203%
95	22.715	3400	3405	3410	rBV	758574	1068203	17.16%	1.054%
96	23.244	3488	3495	3510	rBV2	2914202	5413603	86.97%	5.342%
97	23.380	3511	3518	3529	rVB2	2706250	4932658	79.25%	4.868%
98	23.880	3597	3603	3609	rBV	482927	882278	14.17%	0.871%
99	23.962	3611	3617	3628	rBV2	778113	1685821	27.08%	1.664%
100	25.427	3862	3866	3874	rVB5	180370	376318	6.05%	0.371%

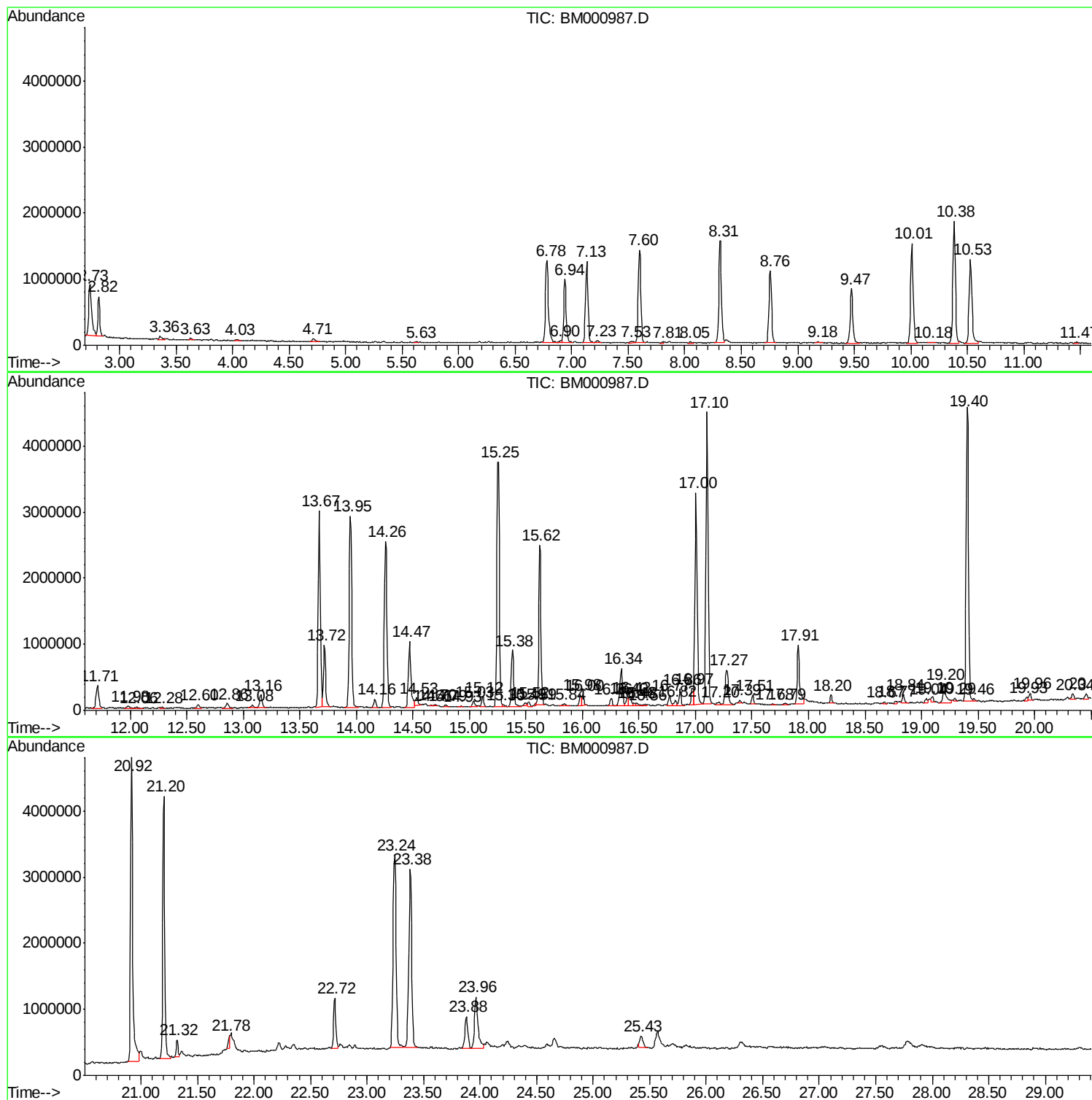
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM01.2-EPA-BM041415.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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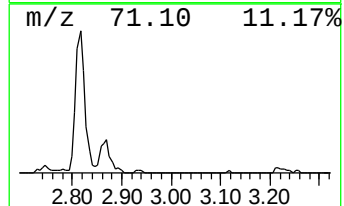
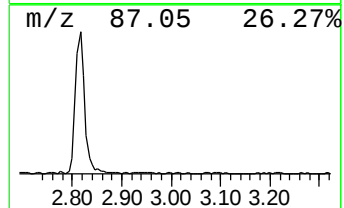
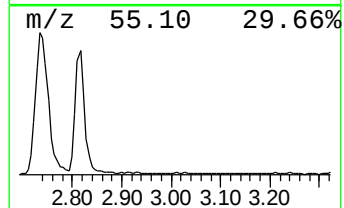
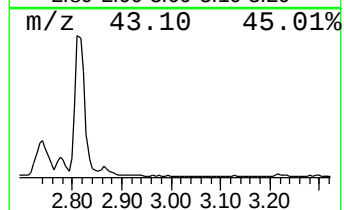
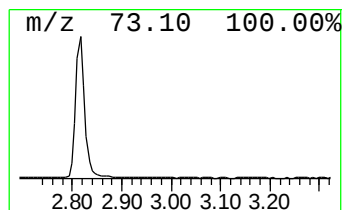
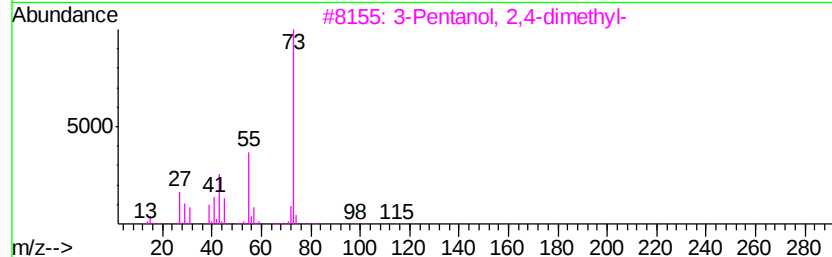
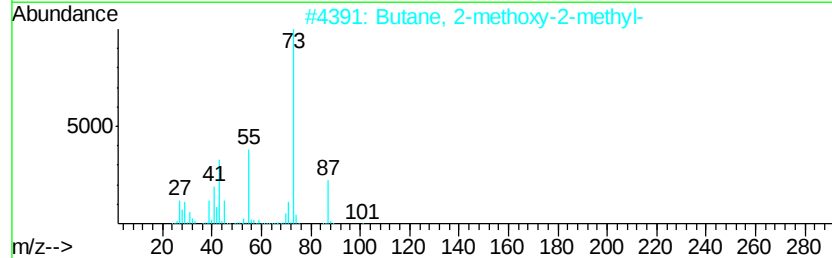
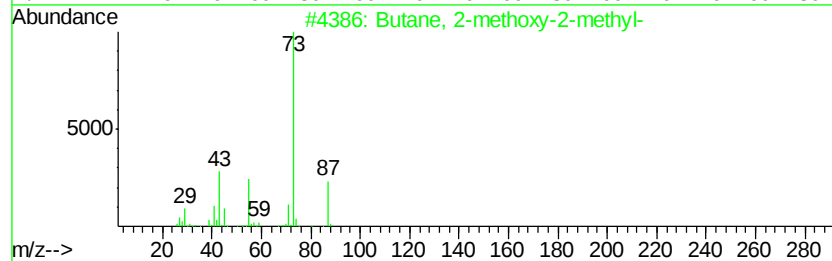
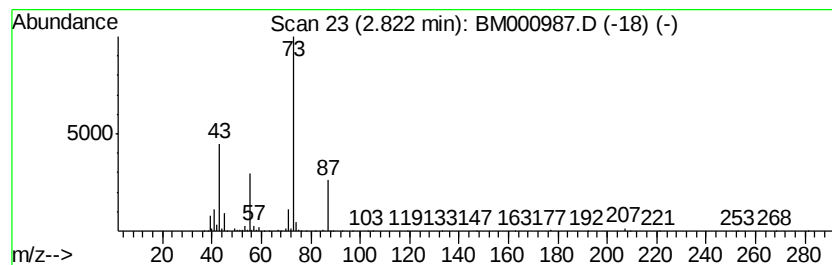
Instrument :
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM01.2-EPA-BM041415.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
2.82	6.74 ng/ul	748827	1,4-Dichlorobenzene-d4	7.60	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64
3	3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	37
4	Silane, tetramethyl-	88	C4H12Si	000075-76-3	35
5	3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	23



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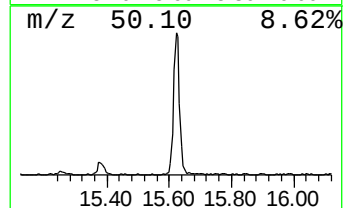
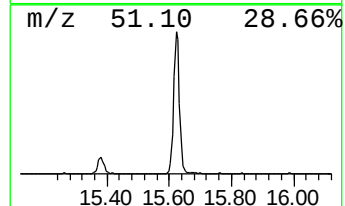
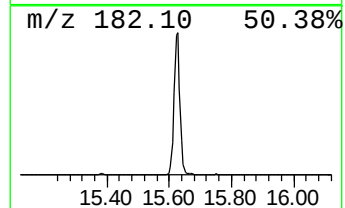
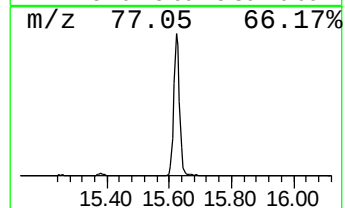
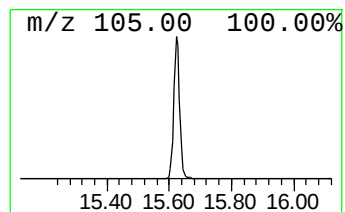
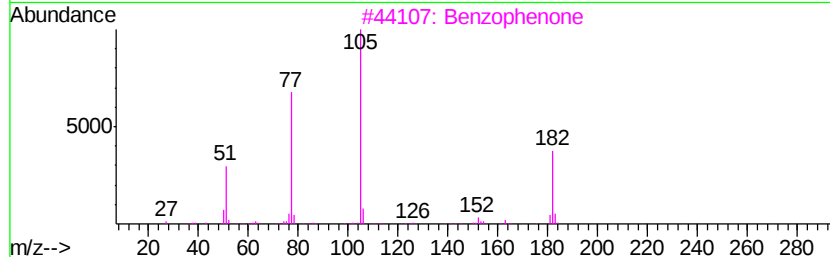
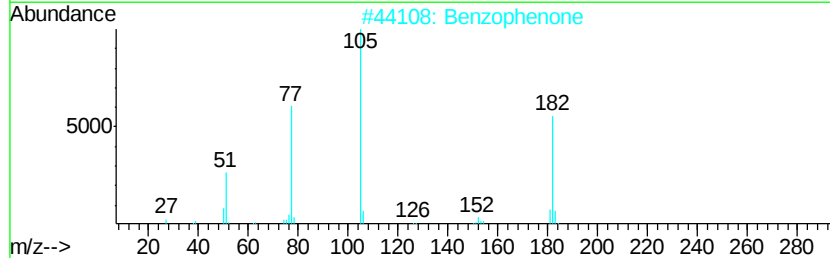
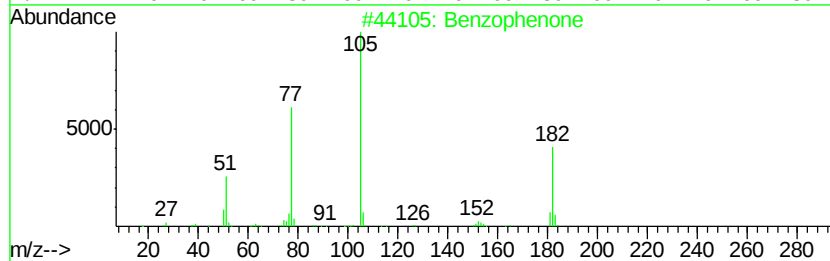
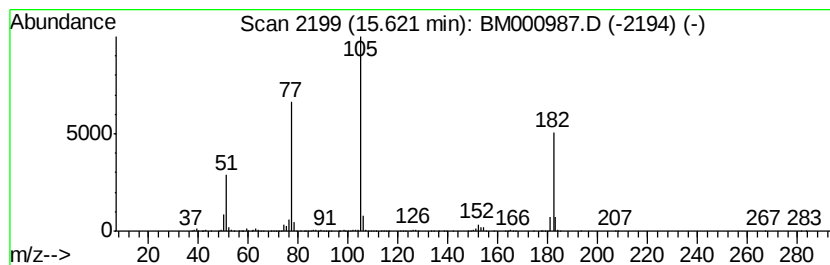
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM01.2-EPA-BM041415.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzophenone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.62	16.71 ng/ul	3225510	Acenaphthene-d10	14.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzophenone	182	C13H10O	000119-61-9	96
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	90



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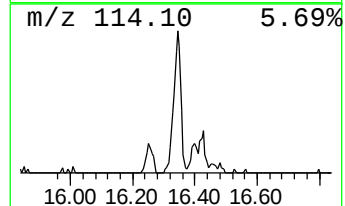
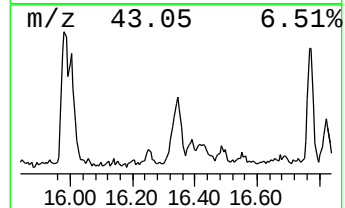
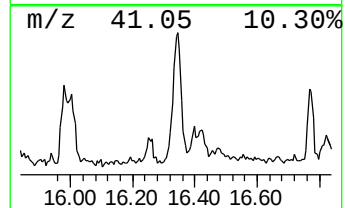
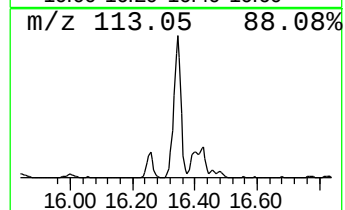
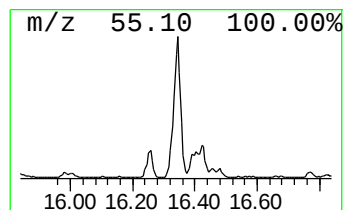
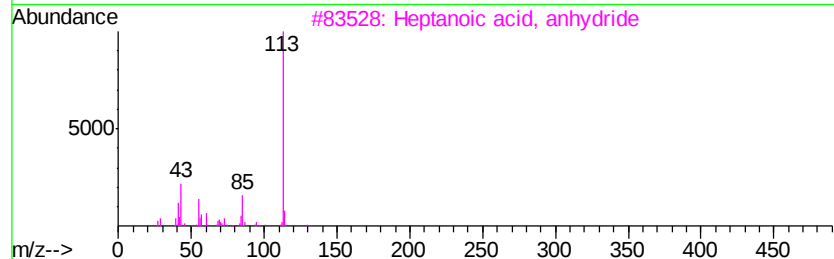
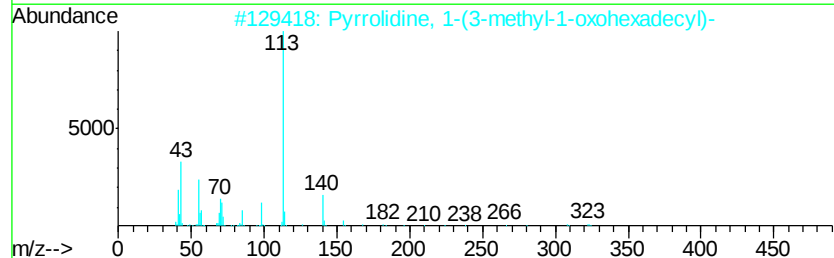
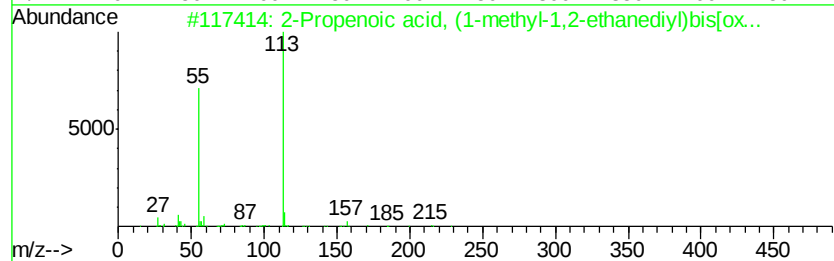
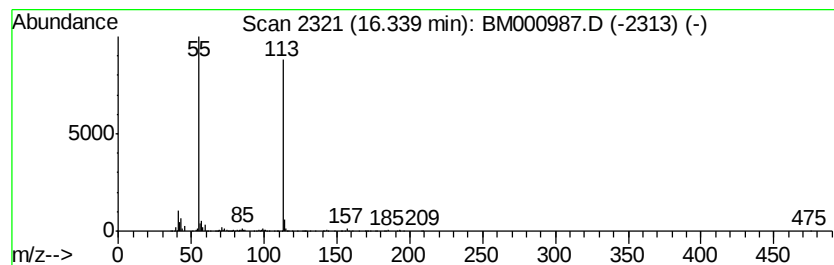
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 2-Propenoic acid, (1-methyl... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.34	4.40 ng/ul	963990	Phenanthrene-d10	17.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Propenoic acid, (1-methyl-1,2-...	300	C15H24O6	042978-66-5	64
2		Pyrrolidine, 1-(3-methyl-1-oxohe...	323	C21H41NO	056630-57-0	38
3		Heptanoic acid, anhydride	242	C14H26O3	000626-27-7	38
4		2-Propenoic acid, (1-methyl-1,2-...	300	C15H24O6	042978-66-5	37
5		1H-Imidazole, 4-nitro-	113	C3H3N3O2	003034-38-6	9



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM041415\
Data File : BM000987.D
Acq On : 14 Apr 2015 17:14
Operator : TP/IZ
Sample : G1858-11
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
P001-SD008-0006-02

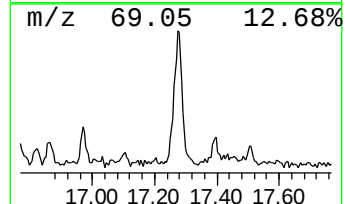
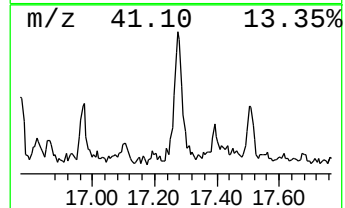
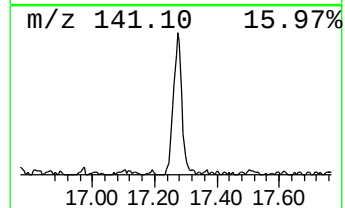
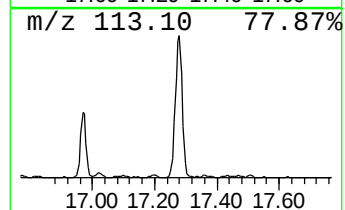
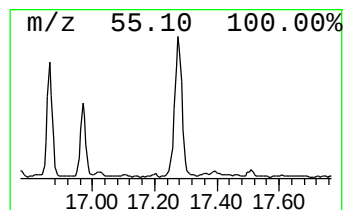
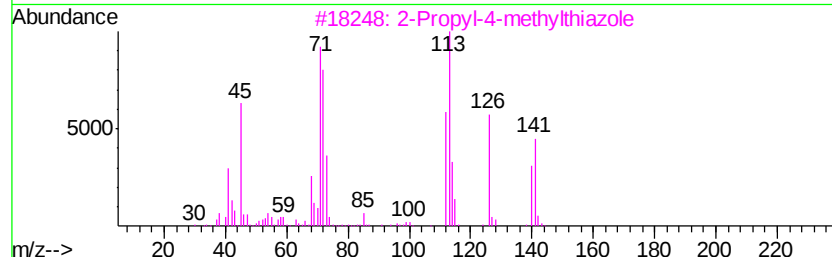
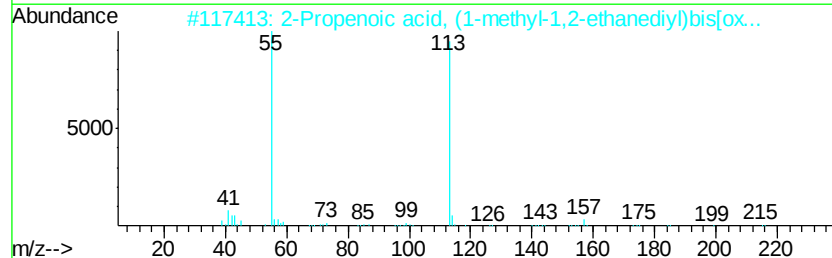
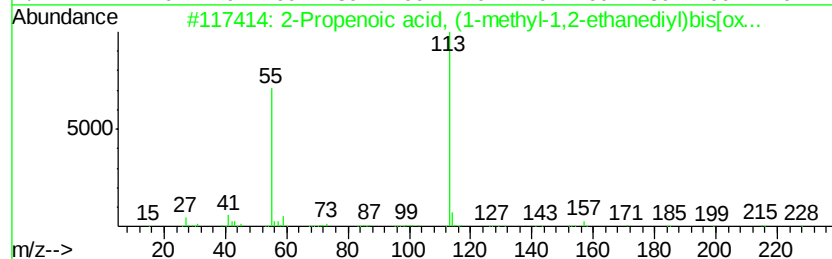
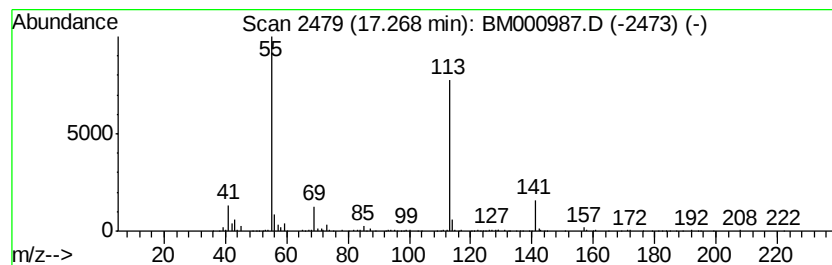
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 unknown-01 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.27	4.29 ng/ul	939220	Phenanthrene-d10	17.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Propenoic acid, (1-methyl-1,2-...	300	C15H24O6	042978-66-5	45
2		2-Propenoic acid, (1-methyl-1,2-...	300	C15H24O6	042978-66-5	42
3		2-Propyl-4-methylthiazole	141	C7H11NS	052414-87-6	40
4		2,5-Pyrrolidinedione, 3-methyl-4...	155	C8H13NO2	016493-22-4	36
5		3-Methylpentyl ethylphosphonoflu...	196	C8H18FO2P	1000298-39-1	36



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM041415\
Data File : BM000987.D
Acq On : 14 Apr 2015 17:14
Operator : TP/IZ
Sample : G1858-11
Misc :
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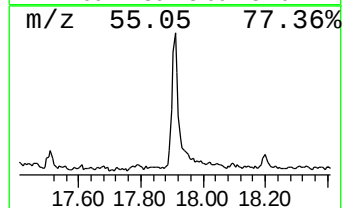
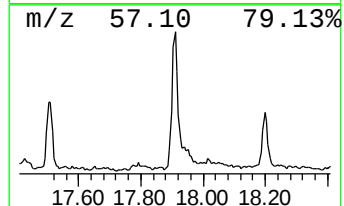
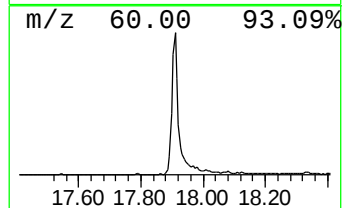
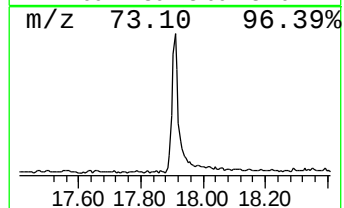
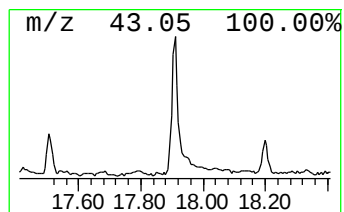
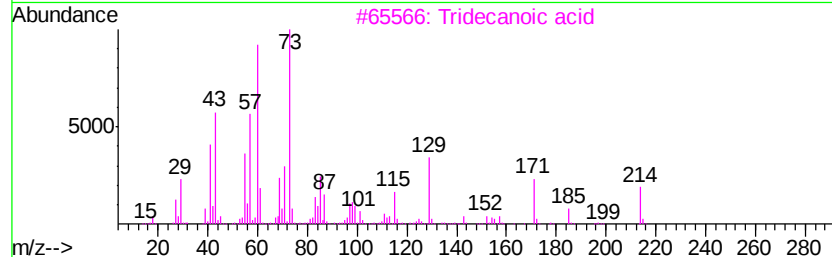
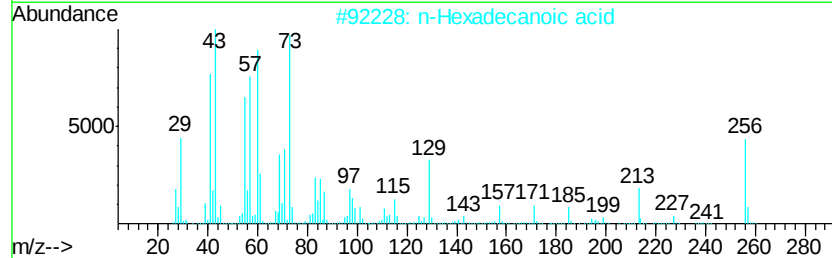
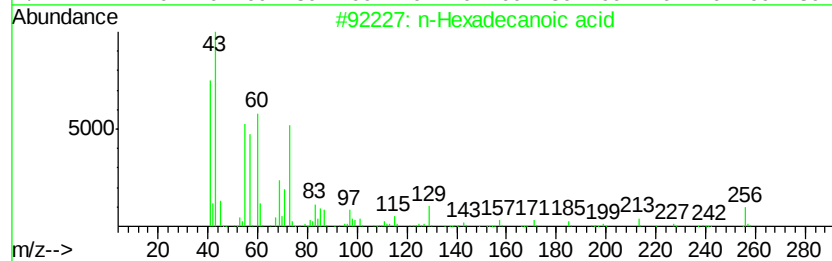
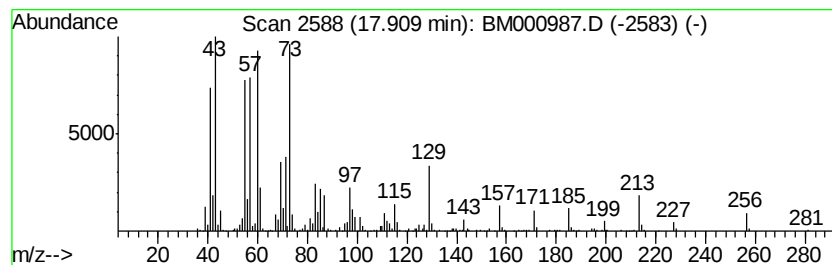
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM01.2-EPA-BM041415.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.91	6.32 ng/ul	1384620	Phenanthrene-d10	17.00	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
3	Tridecanoic acid	214	C13H26O2	000638-53-9	87
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	86
5	Tetradecanoic acid	228	C14H28O2	000544-63-8	83



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM041415\
Data File : BM000987.D
Acq On : 14 Apr 2015 17:14
Operator : TP/IZ
Sample : G1858-11
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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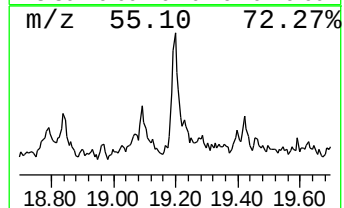
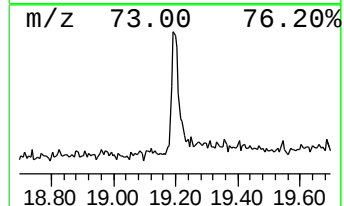
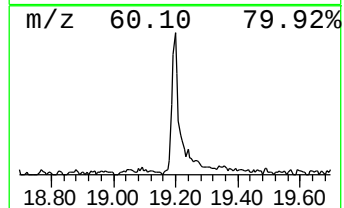
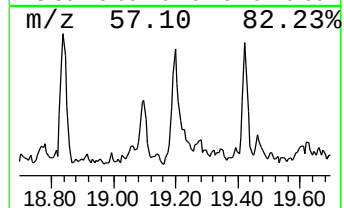
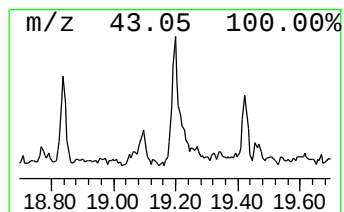
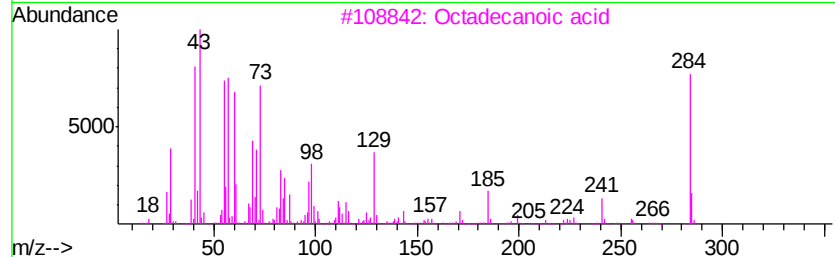
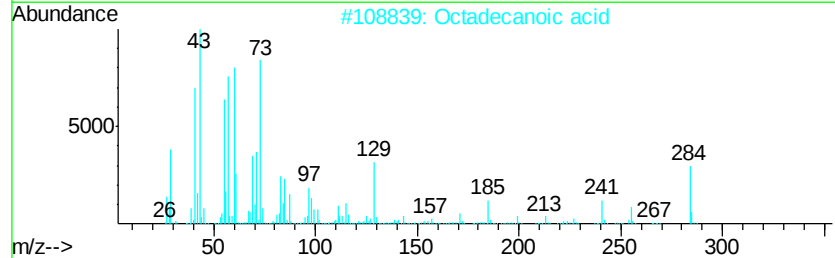
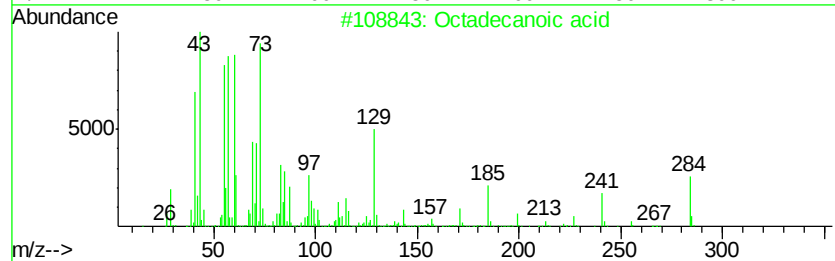
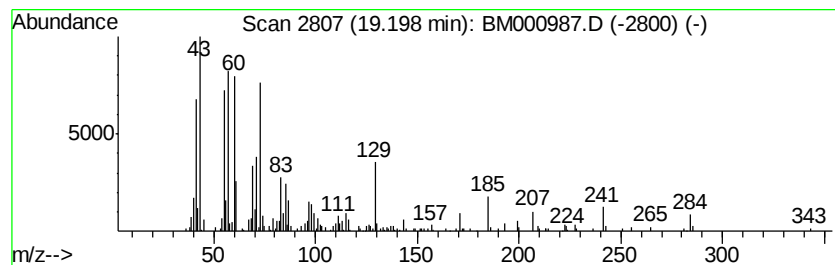
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Octadecanoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.20	2.38 ng/ul	607678	Chrysene-d12	21.20

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	94
3		Octadecanoic acid	284	C18H36O2	000057-11-4	83
4		Octadecanoic acid	284	C18H36O2	000057-11-4	76
5		Tridecanoic acid	214	C13H26O2	000638-53-9	59



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM041415\
Data File : BM000987.D
Acq On : 14 Apr 2015 17:14
Operator : TP/IZ
Sample : G1858-11
Misc :
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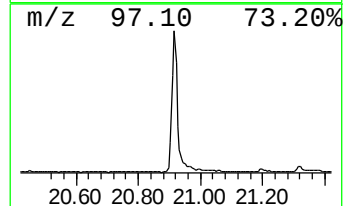
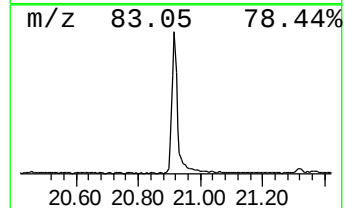
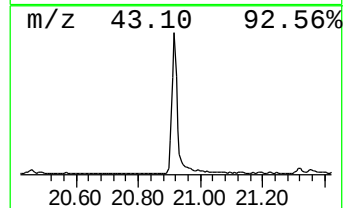
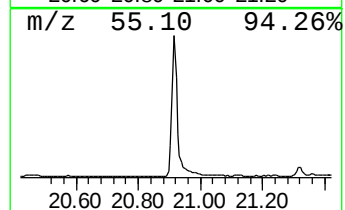
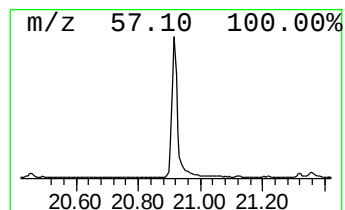
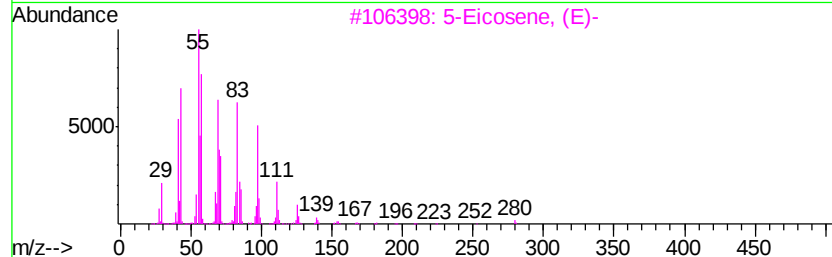
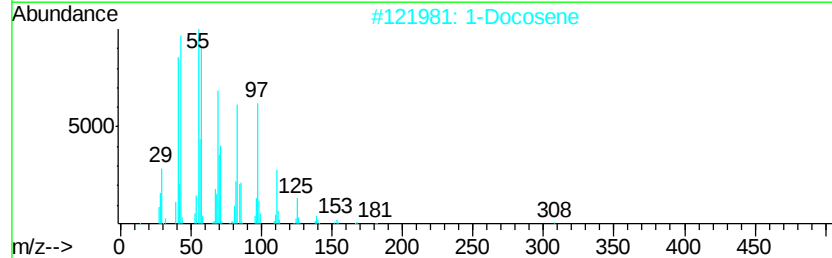
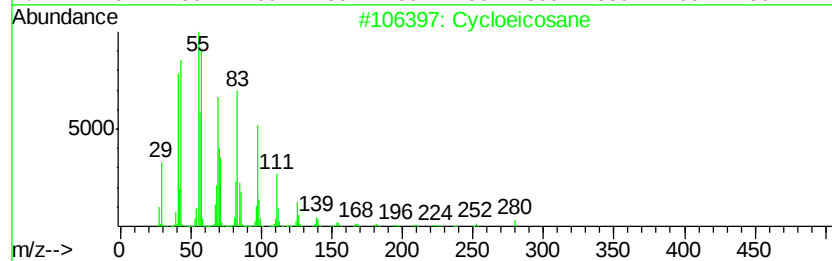
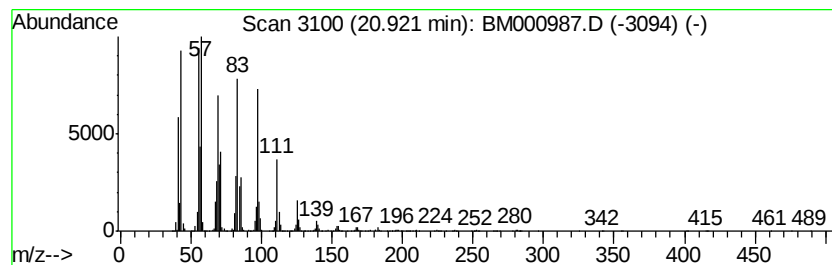
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 (DEL) Alkane: Straight-Chain Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.92	22.63 ng/ul	5773430	Chrysene-d12	21.20
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Cycloeicosane	280 C20H40	000296-56-0 94
2		1-Docosene	308 C22H44	001599-67-3 94
3		5-Eicosene, (E)-	280 C20H40	074685-30-6 91
4		2- Chloropropionic acid, hexadec...	332 C19H37ClO2	086711-81-1 91
5		2- Chloropropionic acid, pentade...	318 C18H35ClO2	1000292-44-1 91



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM041415\
Data File : BM000987.D
Acq On : 14 Apr 2015 17:14
Operator : TP/IZ
Sample : G1858-11
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
P001-SD008-0006-02

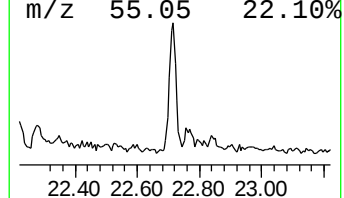
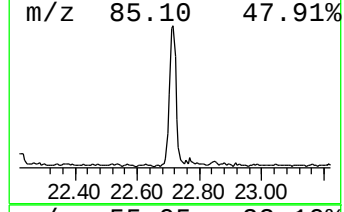
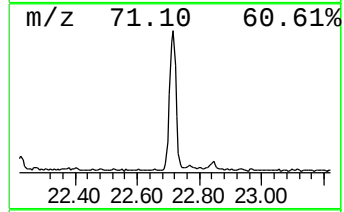
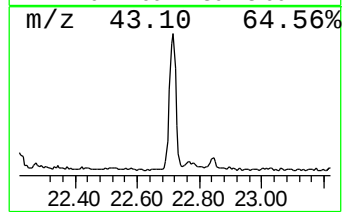
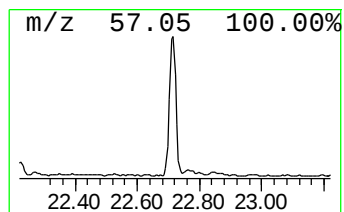
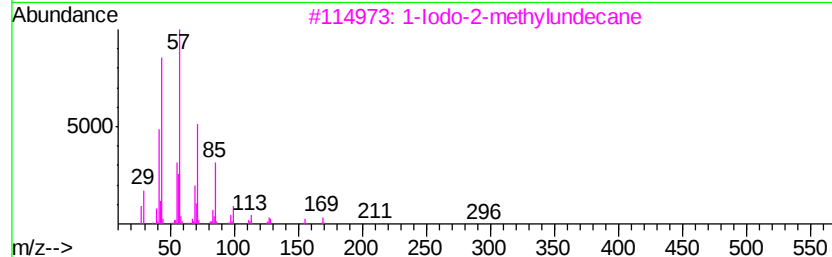
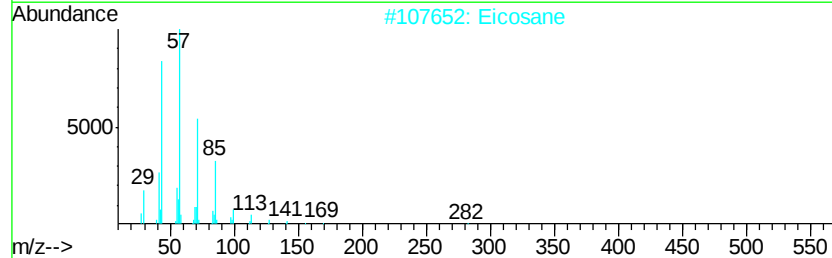
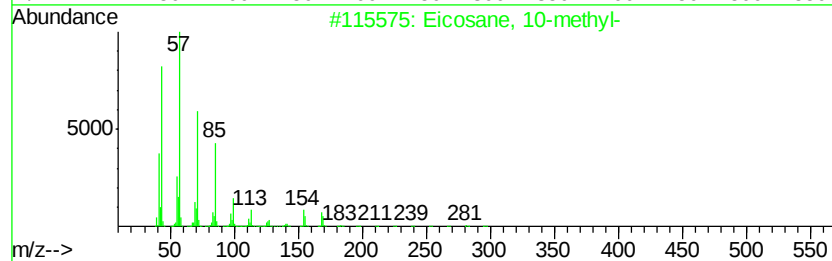
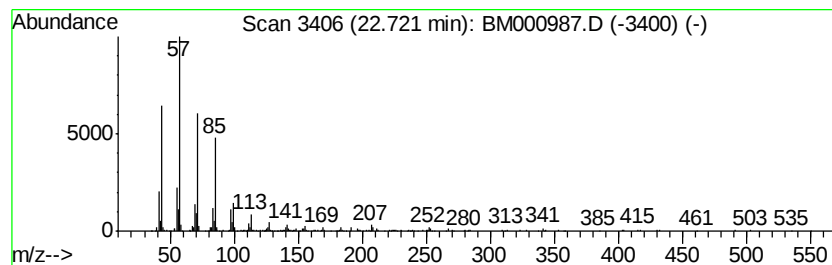
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 (DEL) Alkane: Straight-Chain Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.72	4.33 ng/ul	1068200	Perylene-d12	23.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
2		Eicosane	282	C20H42	000112-95-8	91
3		1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	90
4		Heptacosane	380	C27H56	000593-49-7	90
5		Tetracosane	338	C24H50	000646-31-1	87



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM041415\
Data File : BM000987.D
Acq On : 14 Apr 2015 17:14
Operator : TP/IZ
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Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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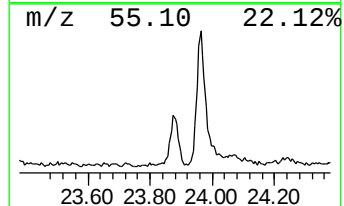
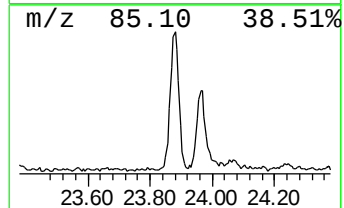
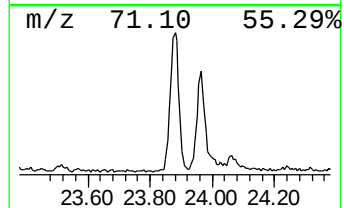
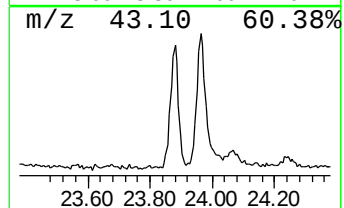
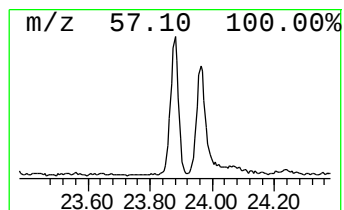
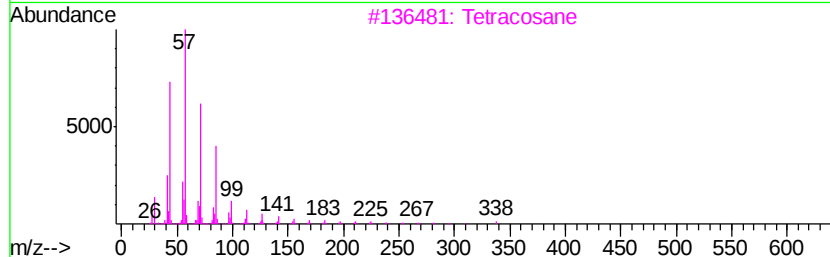
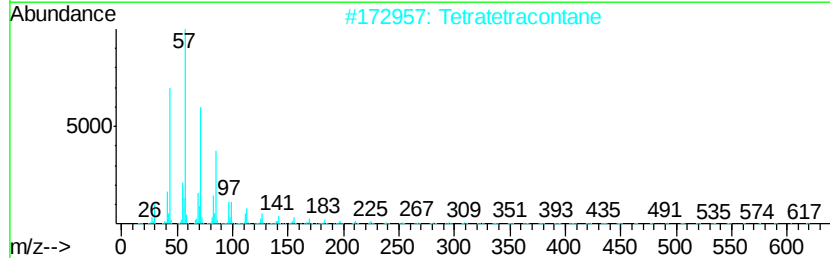
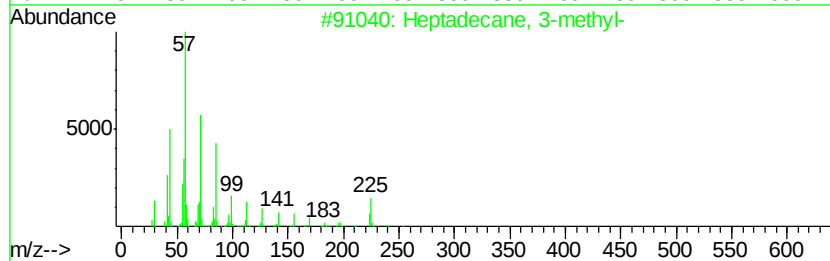
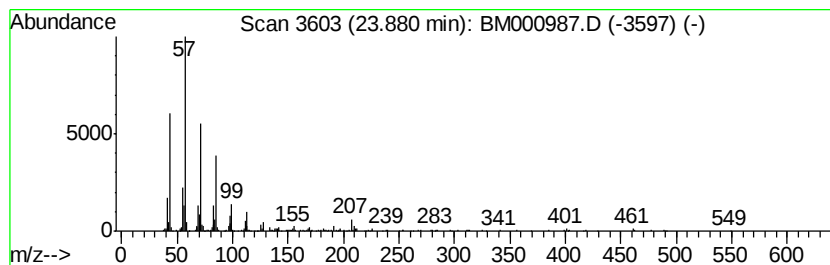
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 (DEL) Alkane: Straight-Chain Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.88	3.58 ng/ul	882278	Perylene-d12	23.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 3-methyl-	254	C18H38	006418-44-6	93
2		Tetratetracontane	619	C44H90	007098-22-8	90
3		Tetracosane	338	C24H50	000646-31-1	90
4		Octacosane	394	C28H58	000630-02-4	90
5		Heneicosane	296	C21H44	000629-94-7	87



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM041415\
Data File : BM000987.D
Acq On : 14 Apr 2015 17:14
Operator : TP/IZ
Sample : G1858-11
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
P001-SD008-0006-02

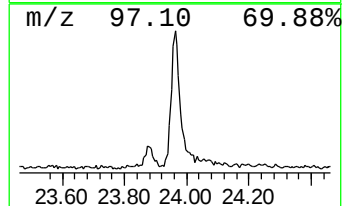
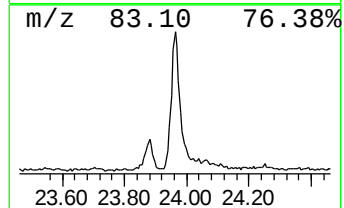
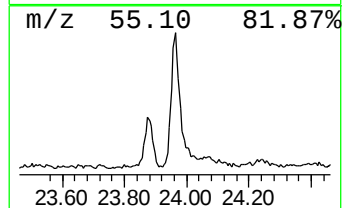
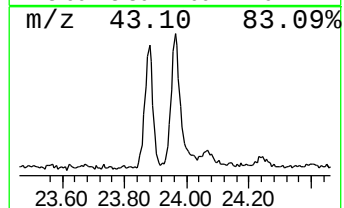
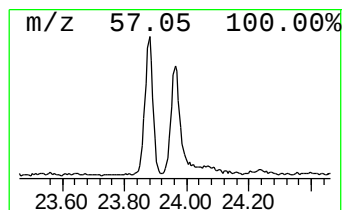
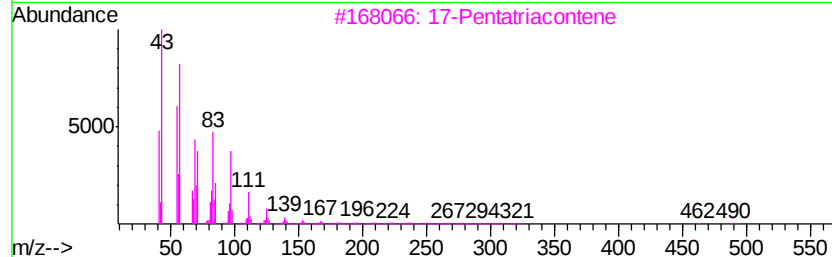
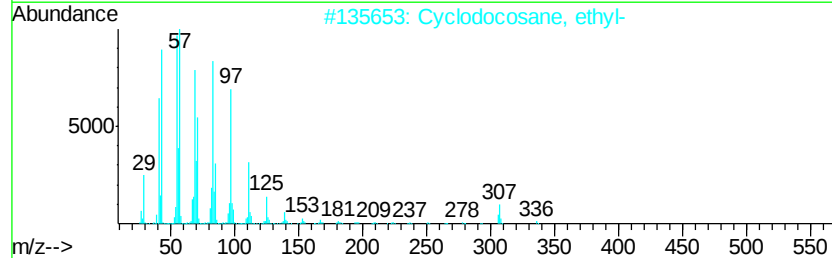
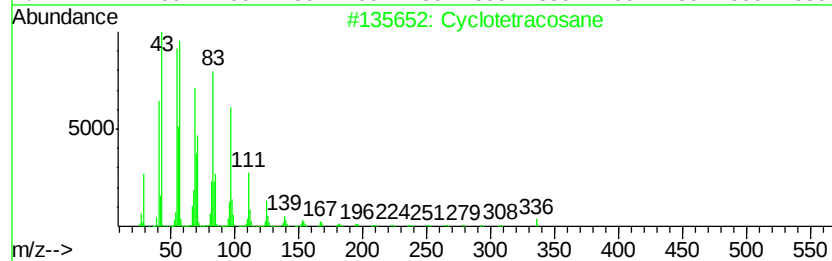
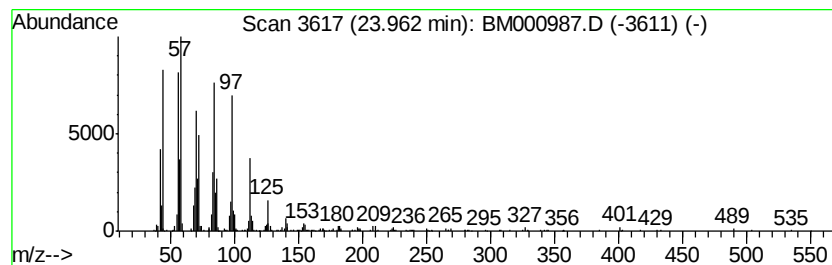
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM01.2-EPA-BM041415.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.96	6.84 ng/ul	1685820	Perylene-d12	23.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclotetracosane	336	C24H48	000297-03-0	91
2		Cyclodocosane, ethyl-	336	C24H48	1000151-22-6	91
3		17-Pentatriacontene	491	C35H70	006971-40-0	91
4		Z-5-Nonadecene	266	C19H38	1000131-11-8	83
5		E-14-Hexadecenal	238	C16H30O	330207-53-9	80



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM041415\
Data File : BM000987.D
Acq On : 14 Apr 2015 17:14
Operator : TP/IZ
Sample : G1858-11
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
P001-SD008-0006-02

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM01.2-EPA-BM041415.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	2.82	6.7	ng/ul	748827	1	7.60	2222810	20.0
Benzophenone	15.62	16.7	ng/ul	3225510	3	14.26	3859900	20.0
2-Propenoic acid, ...	16.34	4.4	ng/ul	963990	4	17.00	4381770	20.0
unknown-01	17.27	4.3	ng/ul	939220	4	17.00	4381770	20.0
n-Hexadecanoic acid	17.91	6.3	ng/ul	1384620	4	17.00	4381770	20.0
Octadecanoic acid	19.20	2.4	ng/ul	607678	5	21.20	5101380	20.0
(DEL) Alkane: Str...	20.92	22.6	ng/ul	5773430	5	21.20	5101380	20.0
(DEL) Alkane: Str...	22.72	4.3	ng/ul	1068200	6	23.38	4932660	20.0
(DEL) Alkane: Str...	23.88	3.6	ng/ul	882278	6	23.38	4932660	20.0
(DEL) Alkane: Str...	23.96	6.8	ng/ul	1685820	6	23.38	4932660	20.0