

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG030716\
 Data File : BG021340.D
 Acq On : 7 Mar 2016 15:23
 Operator : UM/SJ
 Sample : H1625-09
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 JHFG1

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_G\METHODS\SOM02.2-EPA-BG030316.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	3.190	55	60	68	rBV3	3391	6892	2.50%	0.183%
2	3.537	113	119	122	rVB2	1694	2264	0.82%	0.060%
3	7.274	749	755	765	rBB	45418	79601	28.91%	2.111%
4	7.444	779	784	792	rBB	34484	55441	20.13%	1.470%
5	7.655	812	820	826	rBB	45115	72684	26.40%	1.928%
6	8.125	894	900	906	rBB	43969	70920	25.76%	1.881%
7	8.819	1012	1018	1026	rBV	57513	97237	35.31%	2.579%
8	9.289	1091	1098	1107	rBV	48393	81989	29.78%	2.175%
9	9.377	1110	1113	1117	rBB2	986	1144	0.42%	0.030%
10	10.012	1215	1221	1228	rBB	39536	67318	24.45%	1.785%
11	10.546	1306	1312	1319	rBB	56772	94707	34.39%	2.512%
12	10.934	1371	1378	1384	rBV	66981	113556	41.24%	3.012%
13	11.063	1394	1400	1409	rBB2	39139	67331	24.45%	1.786%
14	13.084	1742	1744	1748	rBB	579	780	0.28%	0.021%
15	13.613	1830	1834	1841	rBB	1205	1757	0.64%	0.047%
16	14.136	1916	1923	1927	rBV	95774	140351	50.97%	3.722%
17	14.177	1927	1930	1936	rVB	44017	59491	21.61%	1.578%
18	14.430	1967	1973	1980	rVB	114713	176108	63.96%	4.671%
19	14.606	1999	2003	2008	rBB2	5702	6674	2.42%	0.177%
20	14.735	2019	2025	2032	rBB2	94993	151323	54.96%	4.013%
21	14.918	2050	2056	2066	rVB	42194	69594	25.27%	1.846%
22	15.041	2075	2077	2078	rBV	1293	848	0.31%	0.022%
23	15.059	2078	2080	2083	rVV	1256	1359	0.49%	0.036%
24	15.106	2085	2088	2093	rVB2	1259	1562	0.57%	0.041%
25	15.164	2095	2098	2101	rBB	973	890	0.32%	0.024%
26	15.223	2102	2108	2111	rBB	1813	2290	0.83%	0.061%
27	15.294	2117	2120	2122	rBB2	1234	1257	0.46%	0.033%
28	15.505	2152	2156	2160	rBV	1886	2752	1.00%	0.073%
29	15.552	2160	2164	2169	rVB	12460	16087	5.84%	0.427%
30	15.617	2170	2175	2179	rBV2	1580	2240	0.81%	0.059%
31	15.722	2187	2193	2199	rVB	149514	219375	79.67%	5.818%
32	15.799	2202	2206	2207	rBV	629	766	0.28%	0.020%
33	15.828	2207	2211	2217	rVB	35155	48213	17.51%	1.279%
34	15.957	2223	2233	2237	rBV2	4729	9657	3.51%	0.256%

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Integration Parameters: LSCINT.P

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 Title : SVOA CALIBRATION

35	15.999	2237	2240	2243	rBV2	1786	1968	0.71%	0.052%
36	16.057	2246	2250	2254	rVB	1700	2550	0.93%	0.068%
37	16.110	2254	2259	2263	rBB	2453	3214	1.17%	0.085%
38	16.145	2263	2265	2266	rBV	699	732	0.27%	0.019%
39	16.169	2266	2269	2274	rVB2	2153	2383	0.87%	0.063%
40	16.240	2279	2281	2284	rBV2	579	753	0.27%	0.020%
41	16.375	2297	2304	2306	rBV2	891	1752	0.64%	0.046%
42	16.410	2306	2310	2314	rBV	13856	20237	7.35%	0.537%
43	16.604	2341	2343	2345	rVB2	913	899	0.33%	0.024%
44	16.757	2365	2369	2372	rBV4	2267	3563	1.29%	0.094%
45	16.809	2376	2378	2383	rVB2	1594	1838	0.67%	0.049%
46	16.868	2385	2388	2393	rVB2	1299	1860	0.68%	0.049%
47	16.915	2393	2396	2402	rVB3	1259	2129	0.77%	0.056%
48	16.980	2404	2407	2410	rBV	1550	2137	0.78%	0.057%
49	17.080	2421	2424	2428	rVB2	740	1202	0.44%	0.032%
50	17.203	2441	2445	2449	rBV	8282	10157	3.69%	0.269%
51	17.256	2449	2454	2458	rVB3	2720	4628	1.68%	0.123%
52	17.362	2469	2472	2476	rVB4	841	1237	0.45%	0.033%
53	17.473	2485	2491	2497	rBV	121057	181143	65.79%	4.804%
54	17.567	2501	2507	2515	rVB2	159307	237868	86.39%	6.309%
55	17.673	2522	2525	2526	rBV	1196	1081	0.39%	0.029%
56	17.738	2532	2536	2539	rBV4	1323	1518	0.55%	0.040%
57	17.826	2545	2551	2555	rBV3	1359	2511	0.91%	0.067%
58	17.938	2565	2570	2576	rBV	4532	5713	2.07%	0.152%
59	18.214	2615	2617	2619	rBV2	1105	1267	0.46%	0.034%
60	18.267	2622	2626	2629	rBV4	872	1547	0.56%	0.041%
61	18.302	2629	2632	2633	rBV2	779	750	0.27%	0.020%
62	18.331	2633	2637	2642	rBV3	7860	12484	4.53%	0.331%
63	18.402	2645	2649	2650	rBV4	1844	2174	0.79%	0.058%
64	18.425	2650	2653	2654	rBV3	2153	2150	0.78%	0.057%
65	18.443	2654	2656	2659	rBV4	2251	2953	1.07%	0.078%
66	18.637	2675	2689	2690	rBV6	20159	55468	20.14%	1.471%
67	18.919	2735	2737	2740	rVB4	1729	1626	0.59%	0.043%
68	18.995	2748	2750	2751	rVB	1358	794	0.29%	0.021%
69	19.077	2760	2764	2765	rBV3	1085	1293	0.47%	0.034%
70	19.130	2772	2773	2777	rVV3	777	875	0.32%	0.023%
71	19.189	2779	2783	2786	rVV5	1609	2389	0.87%	0.063%

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72	19.254	2791	2794	2796	rVV3	1394	1469	0.53%	0.039%
73	19.477	2829	2832	2836	rBV5	2444	3952	1.44%	0.105%
74	19.547	2842	2844	2849	rVB4	1537	1847	0.67%	0.049%
75	19.589	2849	2851	2857	rBV4	2883	4904	1.78%	0.130%
76	19.735	2871	2876	2878	rVB6	1824	2526	0.92%	0.067%
77	19.759	2878	2880	2881	rBV2	1230	798	0.29%	0.021%
78	19.841	2889	2894	2901	rBV2	188706	275355	100.00%	7.303%
79	20.012	2921	2923	2925	rBV3	1429	1166	0.42%	0.031%
80	20.106	2938	2939	2940	rBV	1989	1176	0.43%	0.031%
81	20.276	2966	2968	2972	rVB5	3202	2646	0.96%	0.070%
82	20.323	2973	2976	2979	rVB5	2789	2926	1.06%	0.078%
83	20.411	2989	2991	2995	rBV4	2128	1936	0.70%	0.051%
84	20.493	3002	3005	3008	rBV4	2317	2605	0.95%	0.069%
85	20.840	3061	3064	3068	rVB6	3773	4568	1.66%	0.121%
86	21.351	3146	3151	3157	rVB2	43328	65981	23.96%	1.750%
87	21.727	3209	3215	3222	rBV	132488	214374	77.85%	5.686%
88	22.538	3350	3353	3363	rVB9	11875	26211	9.52%	0.695%
89	23.572	3515	3529	3548	rVB5	42392	242612	88.11%	6.435%
90	24.759	3722	3731	3740	rVB	92746	248897	90.39%	6.601%
91	24.982	3760	3769	3779	rBV2	75522	218707	79.43%	5.801%
92	25.399	3838	3840	3841	rBV2	2172	1251	0.45%	0.033%
93	26.104	3952	3960	3963	rBV7	9945	21882	7.95%	0.580%
94	27.086	4125	4127	4128	rBV2	2062	1142	0.41%	0.030%
95	27.468	4186	4192	4201	rVB10	6534	20202	7.34%	0.536%
96	30.123	4642	4644	4645	rBV2	2191	1826	0.66%	0.048%
97	30.670	4735	4737	4738	rBV2	2034	1105	0.40%	0.029%
98	31.533	4882	4884	4885	rBV2	1944	1668	0.61%	0.044%
99	31.692	4909	4911	4912	rBV2	2413	1662	0.60%	0.044%
100	32.009	4943	4965	4966	rBV7	19763	91677	33.29%	2.432%

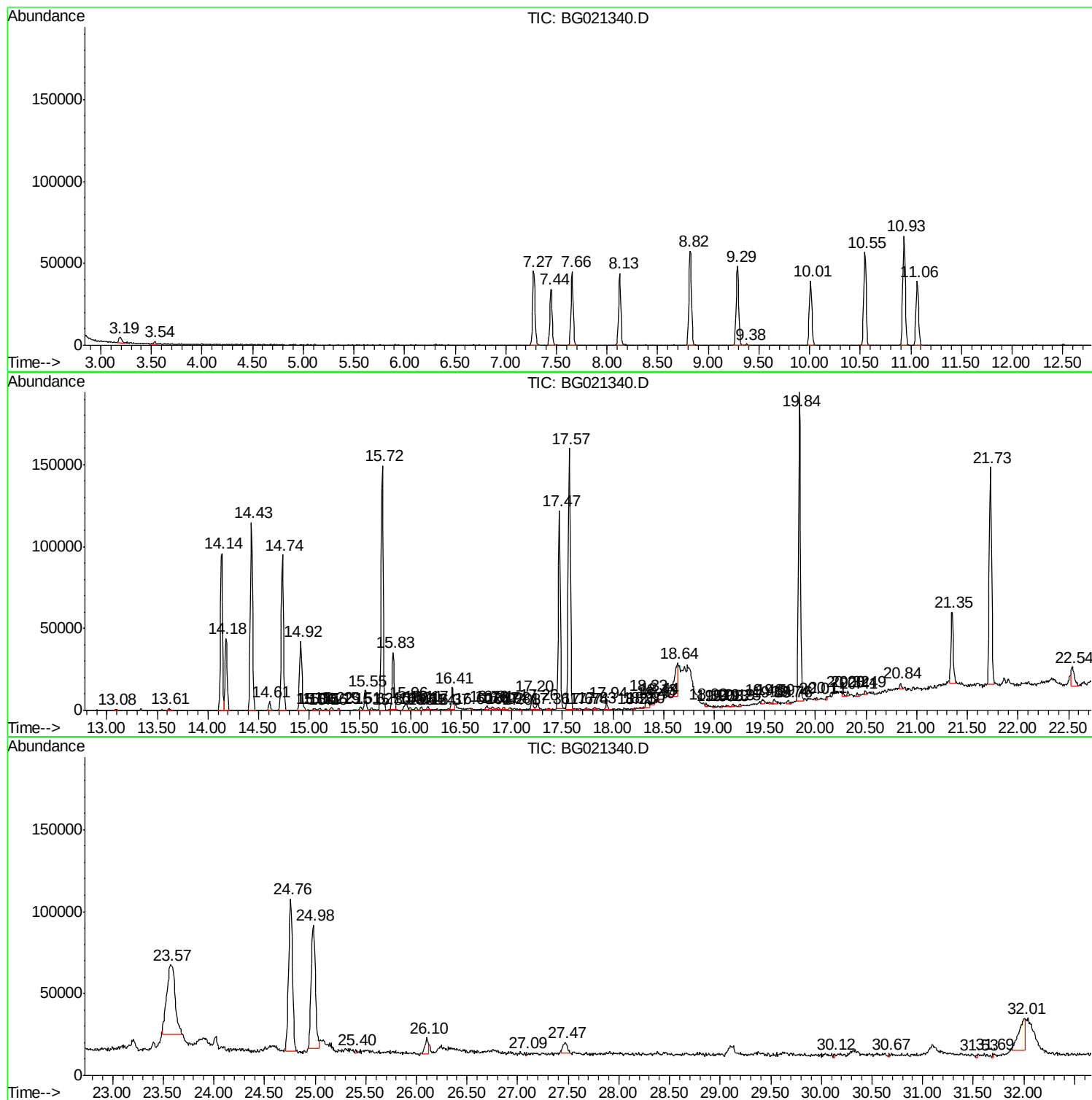
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Instrument :
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG030316.M
Quant Title : SVOA CALIBRATION

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



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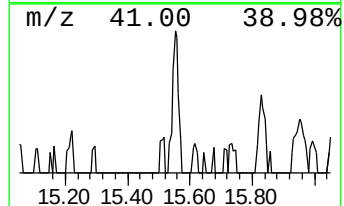
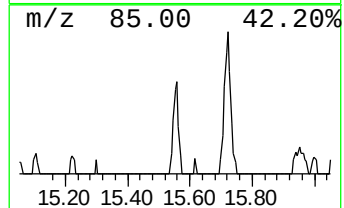
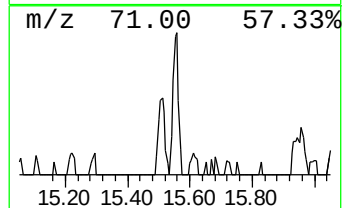
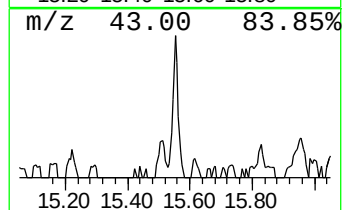
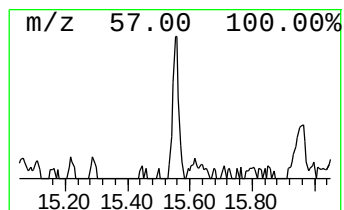
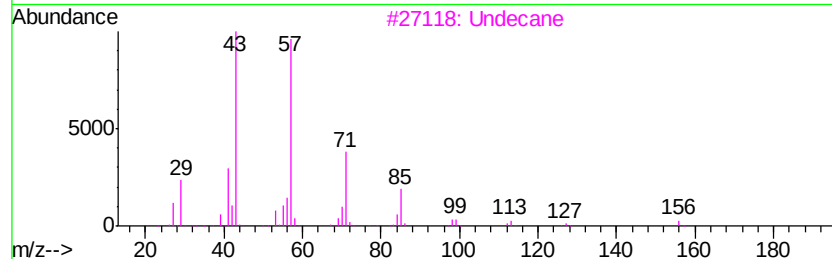
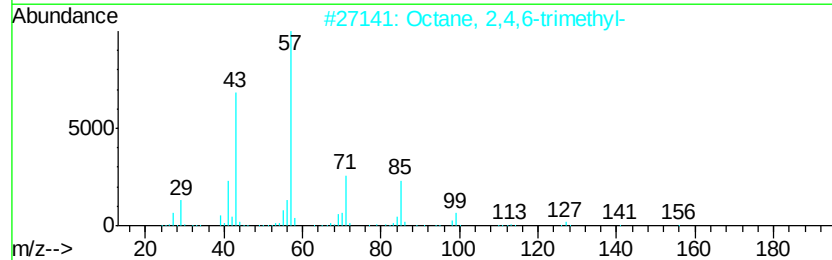
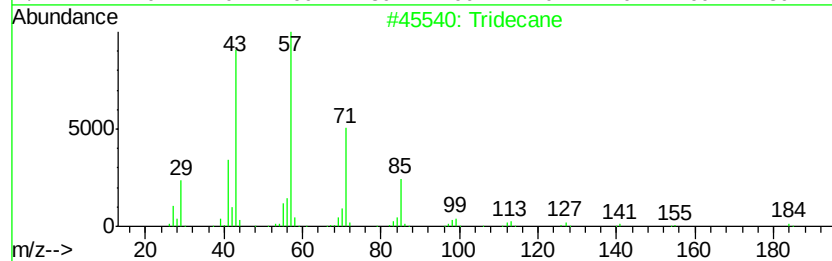
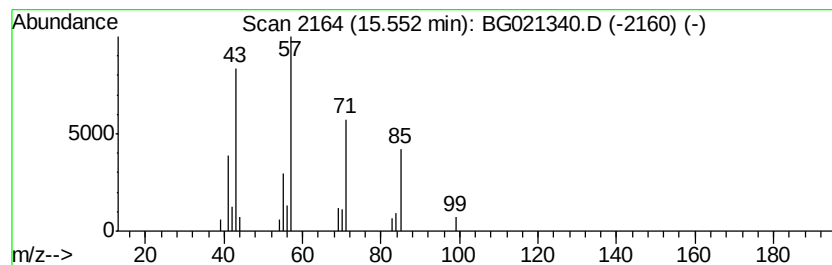
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG030316.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.55	2.13 ng/ul	16087	Acenaphthene-d10	14.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane	184	C13H28	000629-50-5	72
2		Octane, 2,4,6-trimethyl-	156	C11H24	062016-37-9	64
3		Undecane	156	C11H24	001120-21-4	64
4		Hexadecane	226	C16H34	000544-76-3	56
5		Decane	142	C10H22	000124-18-5	53



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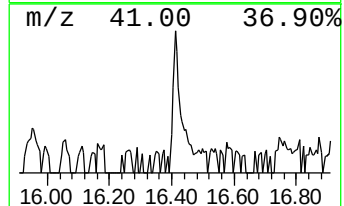
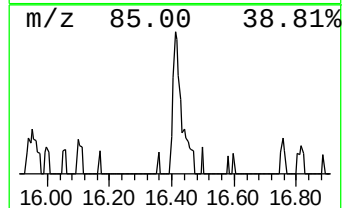
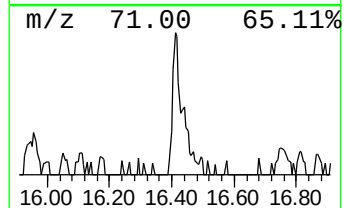
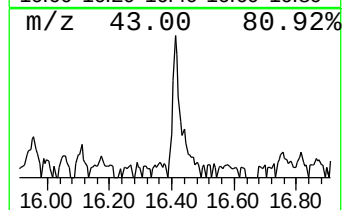
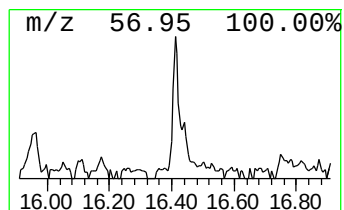
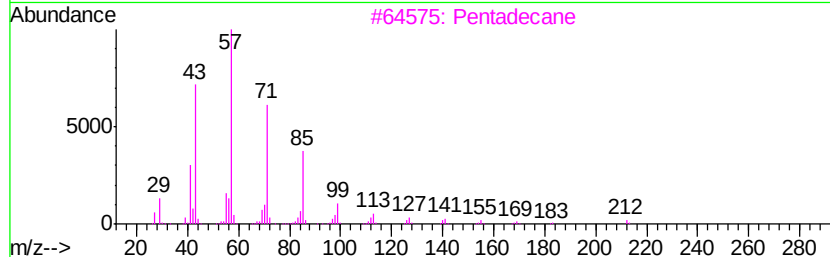
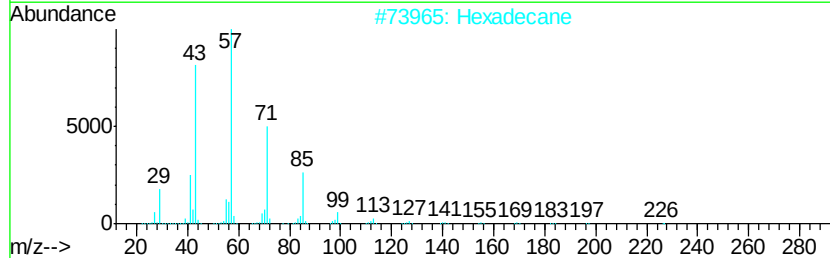
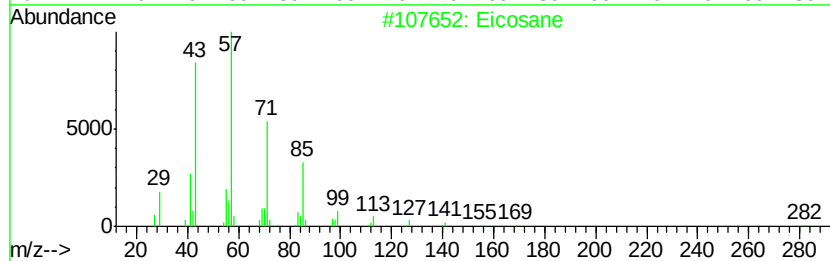
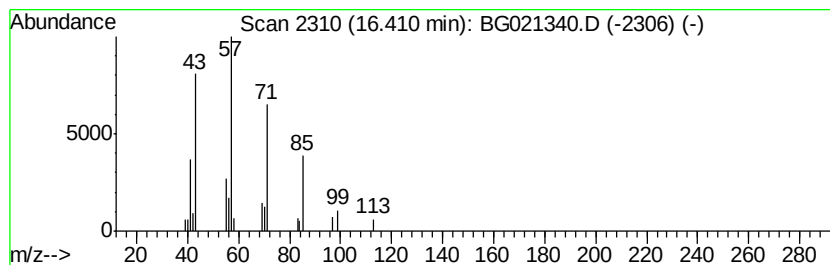
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG030316.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.41	2.23 ng/ul	20237	Phenanthrene-d10	17.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	90
2		Hexadecane	226	C16H34	000544-76-3	83
3		Pentadecane	212	C15H32	000629-62-9	78
4		Hexadecane	226	C16H34	000544-76-3	78
5		Tridecane	184	C13H28	000629-50-5	78



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ALS Vial : 11 Sample Multiplier: 1

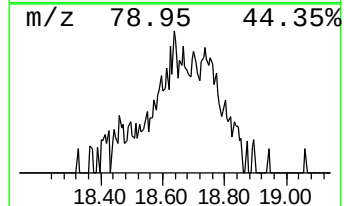
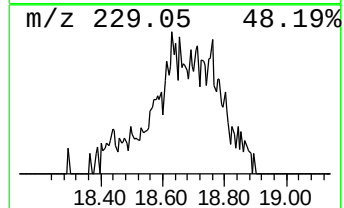
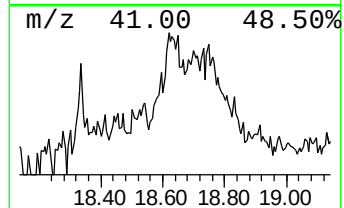
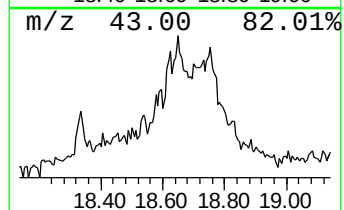
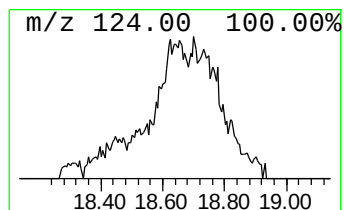
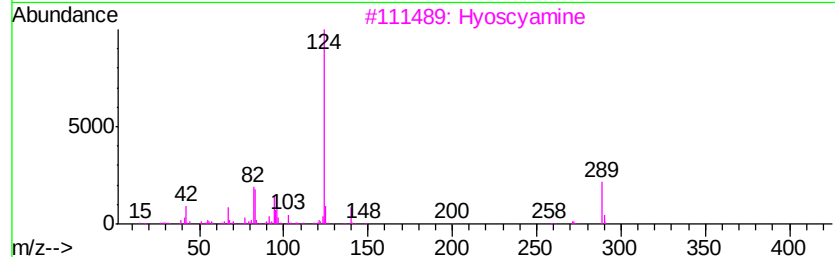
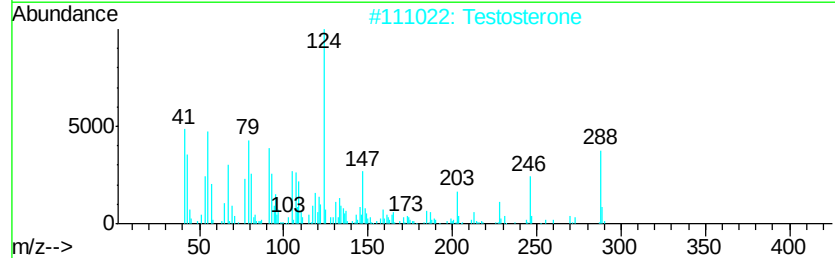
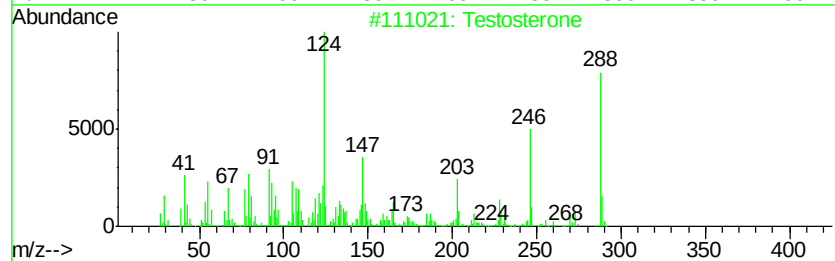
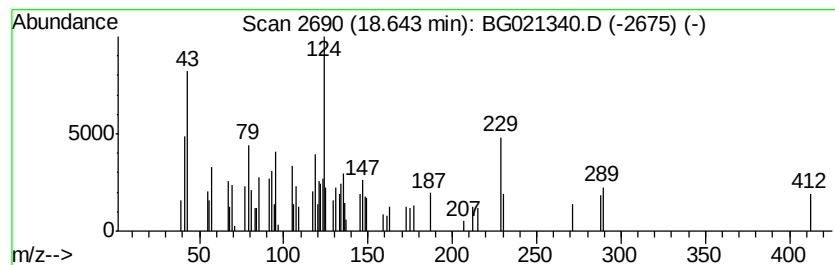
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 unknown-01 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD		R.T.	
18.64	6.12 ng/ul	55468	Phenanthrene-d10		17.47	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Testosterone	288	C19H28O2	000058-22-0	35
2		Testosterone	288	C19H28O2	000058-22-0	27
3		Hyoscyamine	289	C17H23NO3	000101-31-5	22
4		Hyoscyamine	289	C17H23NO3	000101-31-5	22
5		Hyoscyamine	289	C17H23NO3	000101-31-5	22



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Misc :
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ClientSampleId :
JHFG1

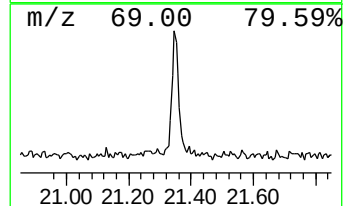
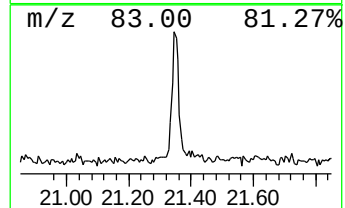
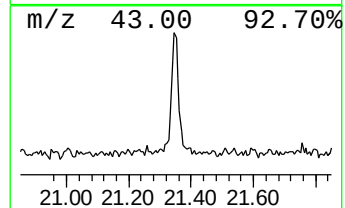
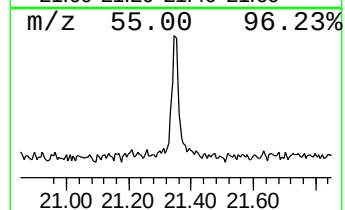
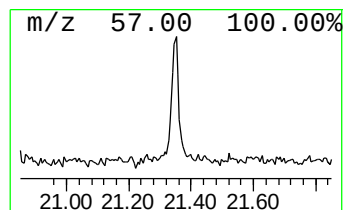
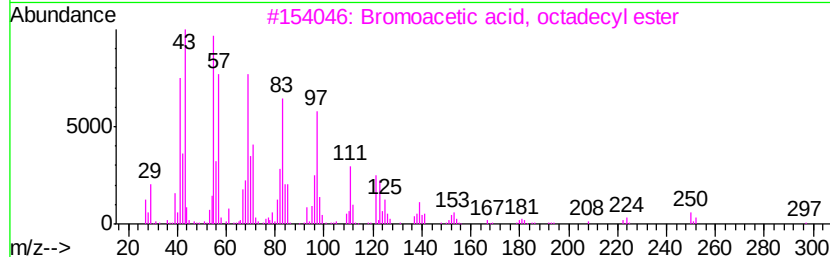
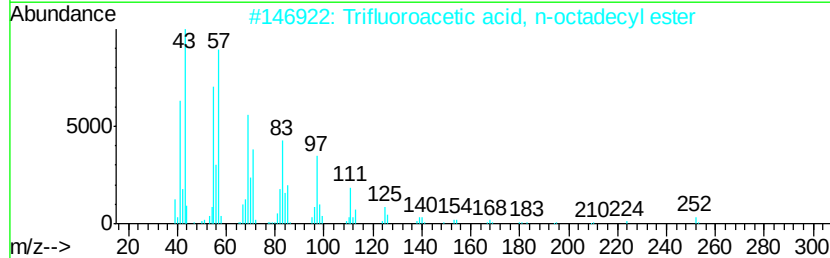
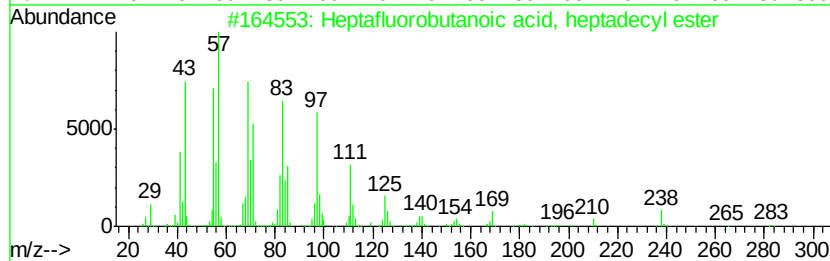
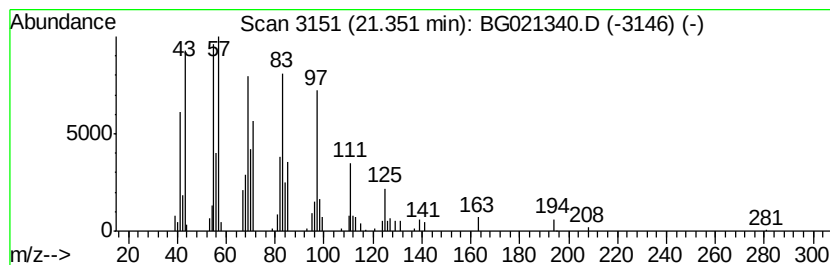
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Heptafluorobutanoic acid, h... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.35	6.16 ng/ul	65981	Chrysene-d12	21.73

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	91
2		Trifluoroacetic acid, n-octadecy...	366	C20H37F3O2	079392-43-1	91
3		Bromoacetic acid, octadecyl ester	390	C20H39BrO2	018992-03-5	91
4		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	90
5		1-Heptadecanol	256	C17H36O	001454-85-9	90



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG030716\
Data File : BG021340.D
Acq On : 7 Mar 2016 15:23
Operator : UM/SJ
Sample : H1625-09
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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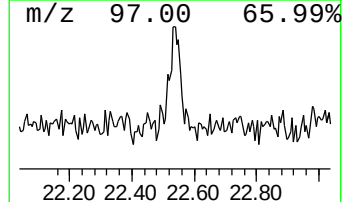
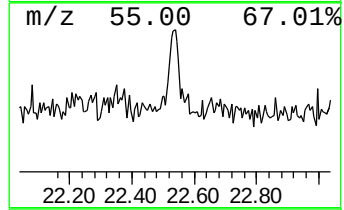
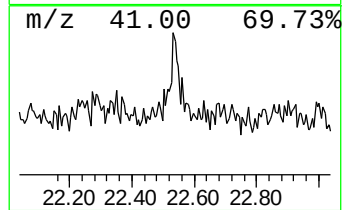
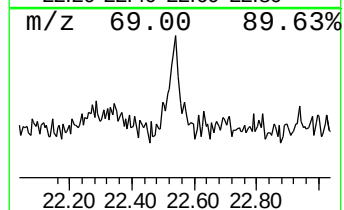
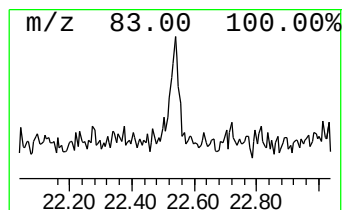
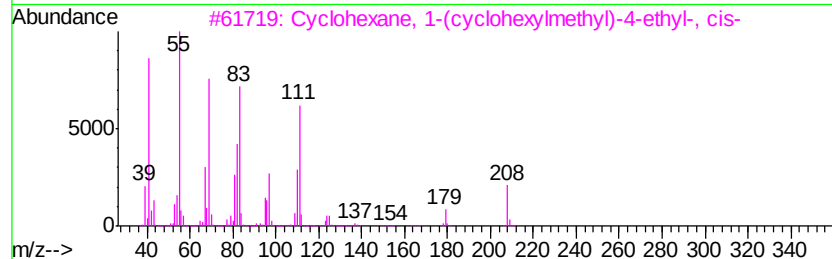
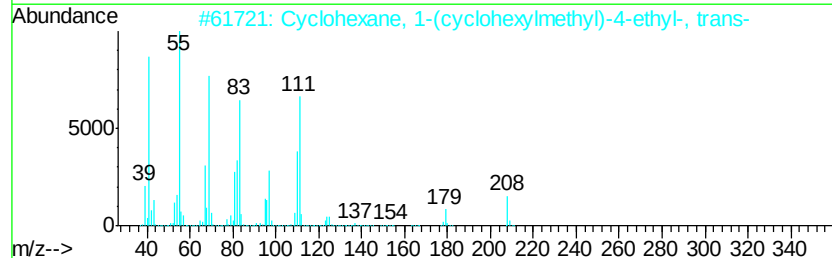
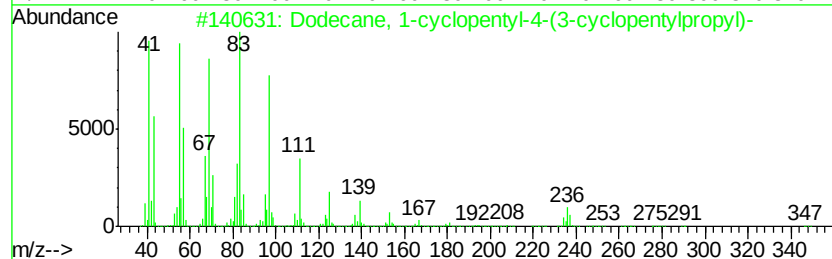
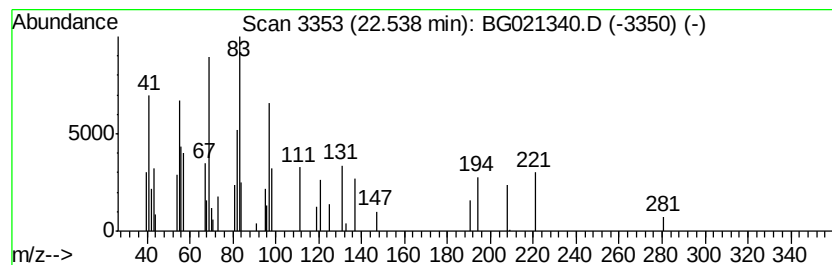
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Branched22.54 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.54	2.45 ng/ul	26211	Chrysene-d12	21.73

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dodecane, 1-cyclopentyl-4-(3-cyc...	348	C25H48	007225-68-5	50
2		Cyclohexane, 1-(cyclohexylmethyl...	208	C15H28	054934-94-0	46
3		Cyclohexane, 1-(cyclohexylmethyl...	208	C15H28	054934-95-1	43
4		1,1'-Bicyclohexyl, 2-propyl-, tr...	208	C15H28	054934-89-3	35
5		9-Nonadecene	266	C19H38	031035-07-1	35



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG030716\
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Acq On : 7 Mar 2016 15:23
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Sample : H1625-09
Misc :
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Instrument :
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ClientSampleId :
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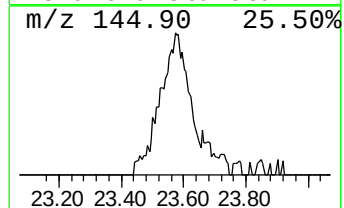
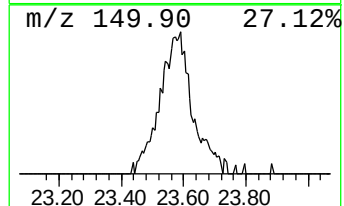
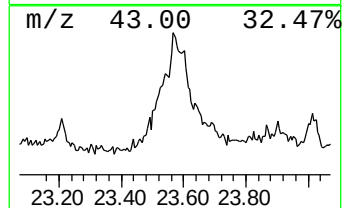
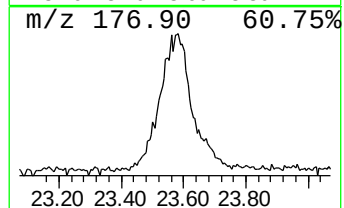
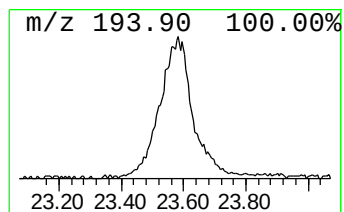
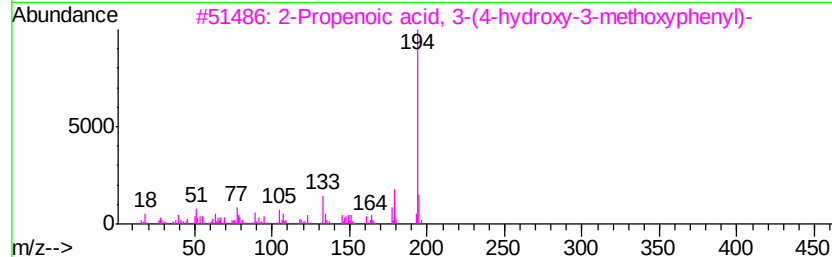
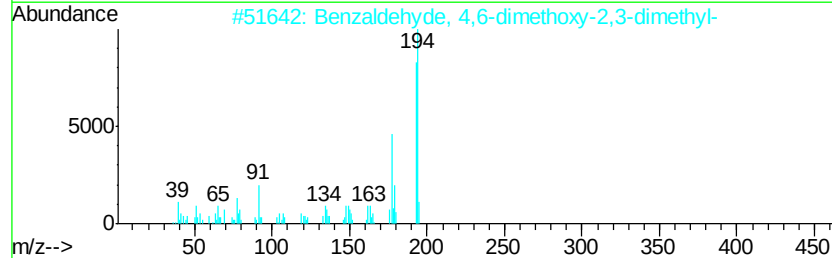
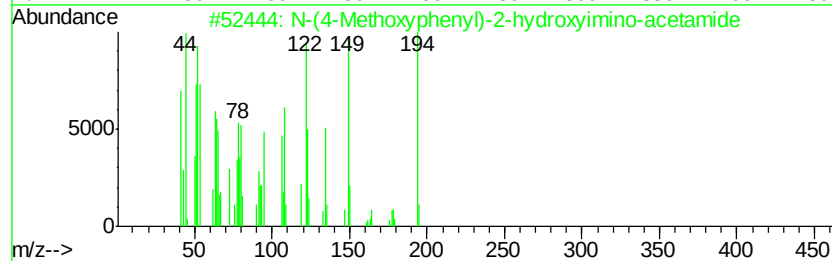
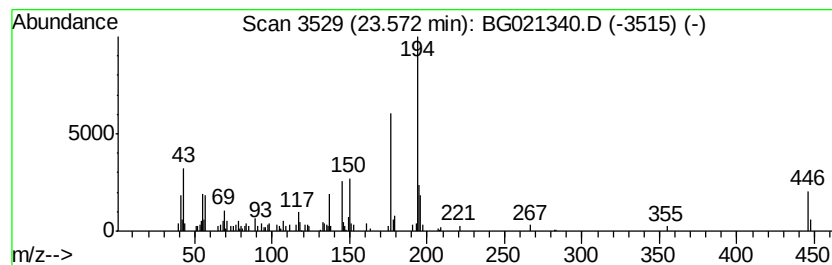
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 unknown-02 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.57	22.19 ng/ul	242612	Perylene-d12	24.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	N-(4-Methoxyphenyl)-2-hydroxyimi...	194	C9H10N2O3	1000143-61-3	35
2		Benzaldehyde, 4,6-dimethoxy-2,3-...	194	C11H14O3	034883-13-1	35
3		2-Propenoic acid, 3-(4-hydroxy-3...	194	C10H10O4	001135-24-6	27
4		3,3'-Dimethyl-5-methoxy-4,5'-bii...	194	C9H10N2O3	019668-80-5	22
5		N-[2-(2-Isopropyl-phenoxy)-ethyl...	329	C19H23NO2S	1000275-37-1	17



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG030716\
Data File : BG021340.D
Acq On : 7 Mar 2016 15:23
Operator : UM/SJ
Sample : H1625-09
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
JHFG1

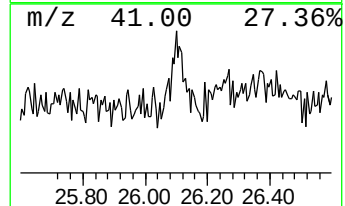
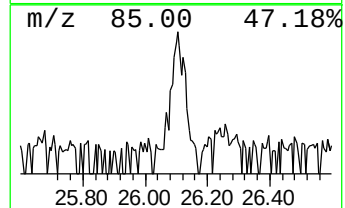
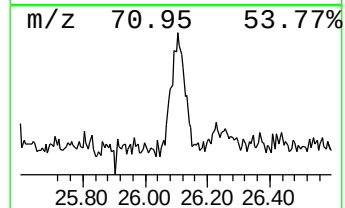
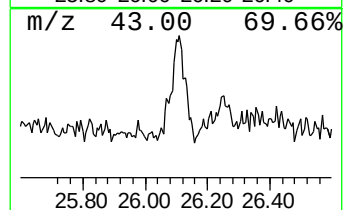
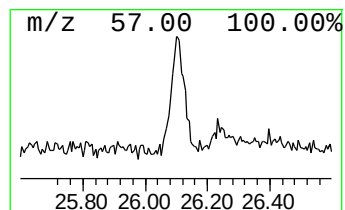
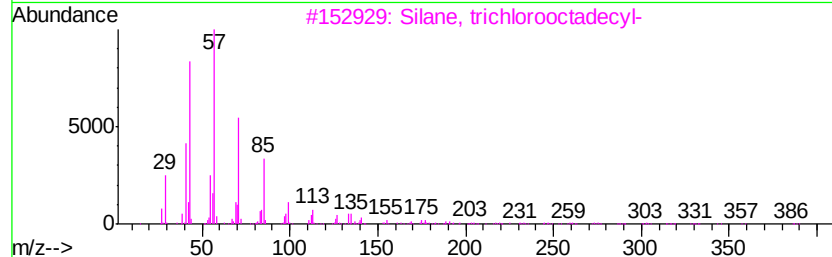
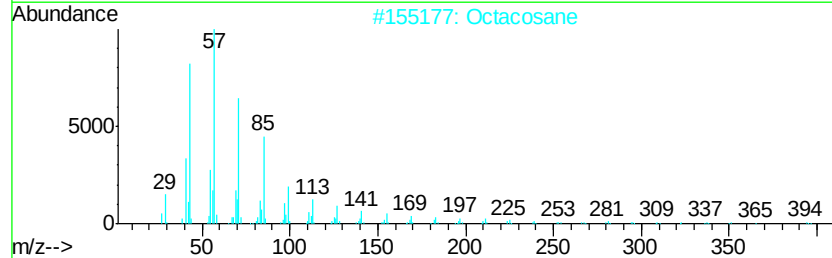
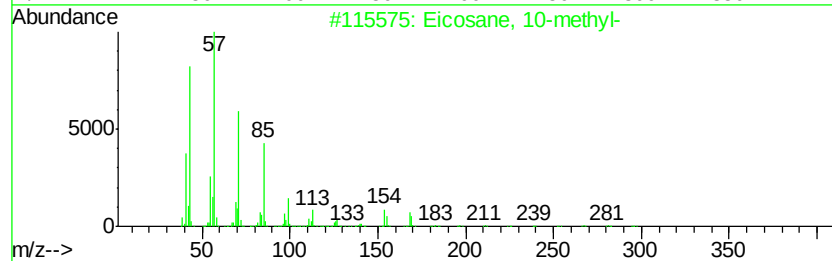
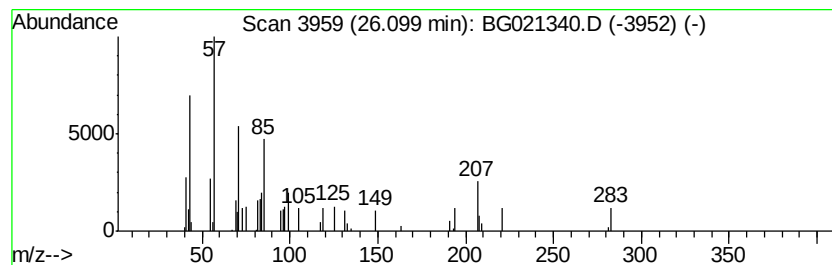
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.10	2.00 ng/ul	21882	Perylene-d12	24.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane, 10-methyl-	296	C21H44	054833-23-7	43
2		Octacosane	394	C28H58	000630-02-4	43
3		Silane, trichlorooctadecyl-	386	C18H37Cl3Si	000112-04-9	43
4		Pentadecane, 2-methyl-	226	C16H34	001560-93-6	43
5		Pentacosane	352	C25H52	000629-99-2	43



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG030716\
Data File : BG021340.D
Acq On : 7 Mar 2016 15:23
Operator : UM/SJ
Sample : H1625-09
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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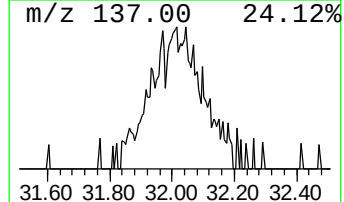
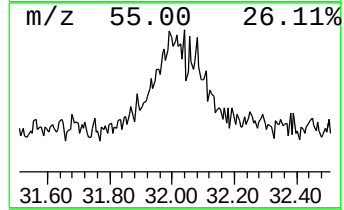
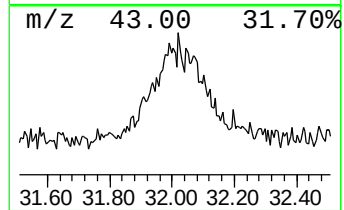
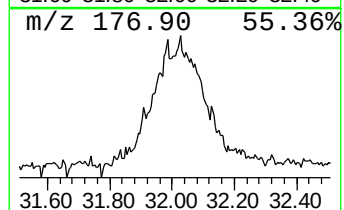
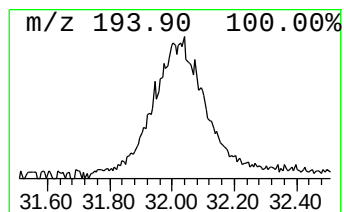
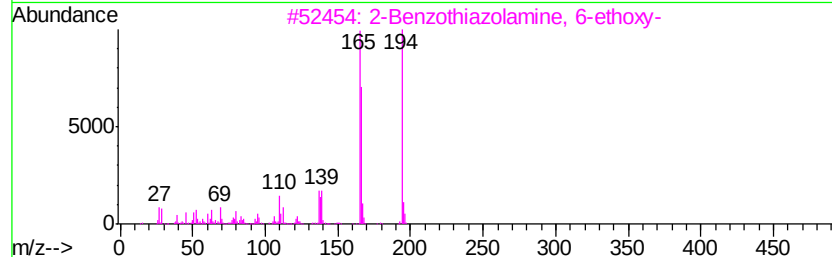
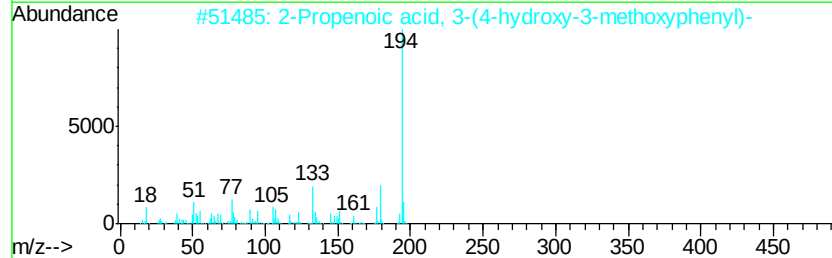
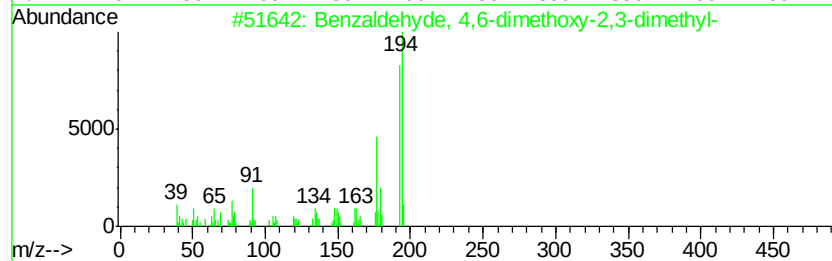
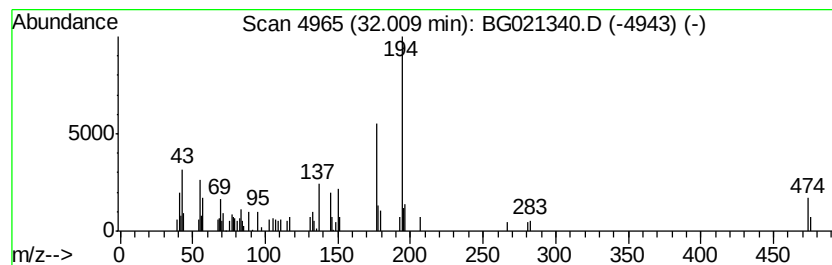
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-03 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
32.01	8.38 ng/ul	91677	Perylene-d12	24.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzaldehyde, 4,6-dimethoxy-2,3-...	194	C11H14O3	034883-13-1	40
2		2-Propenoic acid, 3-(4-hydroxy-3-...	194	C10H10O4	001135-24-6	38
3		2-Benzothiazolamine, 6-ethoxy-	194	C9H10N2OS	000094-45-1	38
4		3-Hydroxy-4-methoxycinnamic acid	194	C10H10O4	000537-73-5	35
5		1,1-Dimethyl-3-(p-tolyl)thiourea	194	C10H14N2S	002741-13-1	35



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG030716\
Data File : BG021340.D
Acq On : 7 Mar 2016 15:23
Operator : UM/SJ
Sample : H1625-09
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG030316.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
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(DEL) Alkane: Str...	16.41	2.2	ng/ul	20237	4	17.47	181143	20.0
unknown-01	18.64	6.1	ng/ul	55468	4	17.47	181143	20.0
Heptafluorobutano...	21.35	6.2	ng/ul	65981	5	21.73	214374	20.0
(DEL) Alkane: Bra...	22.54	2.5	ng/ul	26211	5	21.73	214374	20.0
unknown-02	23.57	22.2	ng/ul	242612	6	24.98	218707	20.0
(DEL) Alkane: Str...	26.10	2.0	ng/ul	21882	6	24.98	218707	20.0
unknown-03	32.01	8.4	ng/ul	91677	6	24.98	218707	20.0