

Data Path : Z:\HPCHEM1\BNA G\DATA\BG032516\
 Data File : BG021508.D
 Acq On : 25 Mar 2016 10:21
 Operator : UM/SJ
 Sample : H1909-04
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 A4T40

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-BG031616.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	3.196	57	61	66	rVB2	3116	4632	1.51%	0.125%
2	3.243	66	69	73	rVB2	4229	5655	1.85%	0.152%
3	3.461	104	106	113	rVB5	1711	2603	0.85%	0.070%
4	3.860	172	174	179	rBV3	587	955	0.31%	0.026%
5	3.995	195	197	202	rVB2	1239	1394	0.46%	0.038%
6	4.548	286	291	299	rVV	15577	23381	7.64%	0.630%
7	5.153	388	394	403	rVB2	9205	15818	5.17%	0.426%
8	7.344	761	767	771	rBV	27659	50259	16.42%	1.354%
9	7.391	771	775	781	rVB	23001	37196	12.15%	1.002%
10	7.638	812	817	825	rBB	29675	49108	16.04%	1.323%
11	8.061	883	889	897	rBB	34859	58308	19.05%	1.571%
12	8.849	1017	1023	1030	rBV	34857	58557	19.13%	1.578%
13	9.230	1082	1088	1096	rBB	32214	54708	17.87%	1.474%
14	9.953	1205	1211	1218	rBB	28674	47570	15.54%	1.282%
15	10.570	1310	1316	1324	rBB	40297	67680	22.11%	1.823%
16	10.870	1360	1367	1373	rBV	58876	102033	33.33%	2.749%
17	10.923	1373	1376	1380	rVB	1468	1972	0.64%	0.053%
18	11.017	1385	1392	1400	rBV	29355	52611	17.18%	1.417%
19	12.374	1619	1623	1627	rBB2	1674	1995	0.65%	0.054%
20	14.078	1908	1913	1918	rBV	87197	122805	40.11%	3.308%
21	14.125	1918	1921	1926	rVB	19045	24251	7.92%	0.653%
22	14.371	1957	1963	1973	rBB	92810	144978	47.35%	3.906%
23	14.548	1989	1993	1996	rBB2	1926	1900	0.62%	0.051%
24	14.677	2009	2015	2022	rBB2	94450	152161	49.70%	4.099%
25	14.736	2022	2025	2029	rBB	985	1379	0.45%	0.037%
26	15.094	2080	2086	2095	rBV3	20822	38440	12.56%	1.036%
27	15.488	2150	2153	2157	rBB	3116	3985	1.30%	0.107%
28	15.664	2177	2183	2189	rBV	126194	182899	59.74%	4.927%
29	15.723	2189	2193	2197	rVB	1200	1869	0.61%	0.050%
30	15.787	2200	2204	2211	rBB2	31853	46461	15.18%	1.252%
31	15.899	2219	2223	2226	rBB2	1308	1519	0.50%	0.041%
32	16.011	2239	2242	2245	rBB2	891	1089	0.36%	0.029%
33	16.158	2264	2267	2272	rBB2	1051	1744	0.57%	0.047%
34	16.346	2297	2299	2307	rBV2	2909	5345	1.75%	0.144%

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35	17.074	2418	2423	2428	rBB	2564	3569	1.17%	0.096%
36	17.139	2431	2434	2440	rBB2	2420	3393	1.11%	0.091%
37	17.239	2447	2451	2455	rVB	787	1160	0.38%	0.031%
38	17.415	2475	2481	2485	rBV	129496	189741	61.98%	5.112%
39	17.456	2485	2488	2492	rVV	35801	50292	16.43%	1.355%
40	17.515	2492	2498	2509	rVV2	126560	200806	65.59%	5.410%
41	17.832	2548	2552	2556	rVV2	1688	2779	0.91%	0.075%
42	17.885	2556	2561	2563	rVV	747	1076	0.35%	0.029%
43	17.926	2565	2568	2571	rBB2	1067	1223	0.40%	0.033%
44	17.997	2577	2580	2584	rVB3	1154	1500	0.49%	0.040%
45	18.144	2602	2605	2609	rBV2	668	1031	0.34%	0.028%
46	18.285	2625	2629	2633	rVV	11263	14187	4.63%	0.382%
47	18.337	2633	2638	2646	rVV3	7166	12055	3.94%	0.325%
48	18.414	2646	2651	2656	rVV2	1890	2685	0.88%	0.072%
49	18.478	2656	2662	2666	rVV3	7902	14920	4.87%	0.402%
50	18.514	2666	2668	2673	rVB2	4306	5704	1.86%	0.154%
51	18.778	2707	2713	2717	rBV2	3689	5929	1.94%	0.160%
52	18.831	2717	2722	2726	rVB2	9213	13035	4.26%	0.351%
53	19.019	2751	2754	2757	rBV2	1646	2262	0.74%	0.061%
54	19.078	2760	2764	2766	rVB2	1489	2012	0.66%	0.054%
55	19.119	2769	2771	2773	rVB2	2079	1755	0.57%	0.047%
56	19.195	2775	2784	2789	rBV3	5556	10658	3.48%	0.287%
57	19.336	2802	2808	2815	rVB2	3877	8480	2.77%	0.228%
58	19.460	2820	2829	2835	rBV	79109	116613	38.09%	3.142%
59	19.518	2835	2839	2840	rBV2	1400	1564	0.51%	0.042%
60	19.548	2840	2844	2849	rVB7	2109	2953	0.96%	0.080%
61	19.607	2849	2854	2859	rVB3	2261	3365	1.10%	0.091%
62	19.654	2859	2862	2864	rBV3	2946	3610	1.18%	0.097%
63	19.689	2864	2868	2870	rVV3	919	1058	0.35%	0.029%
64	19.789	2877	2885	2889	rBV	170842	278903	91.10%	7.514%
65	19.818	2889	2890	2897	rVB	71272	65476	21.39%	1.764%
66	19.889	2897	2902	2906	rBV2	5410	7447	2.43%	0.201%
67	19.994	2918	2920	2925	rVB3	4614	5466	1.79%	0.147%
68	20.135	2938	2944	2951	rBV6	6644	16008	5.23%	0.431%
69	20.200	2951	2955	2956	rBV2	1810	2061	0.67%	0.056%
70	20.300	2963	2972	2980	rBV3	13583	34607	11.30%	0.932%
71	20.359	2980	2982	2984	rBV3	1698	1261	0.41%	0.034%

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72	20.406	2986	2990	2994	rBV	9313	14261	4.66%	0.384%
73	20.464	2994	3000	3004	rVV5	8317	14463	4.72%	0.390%
74	20.541	3010	3013	3017	rBV4	3061	3881	1.27%	0.105%
75	20.605	3020	3024	3027	rVV4	5121	6926	2.26%	0.187%
76	20.647	3027	3031	3035	rBV4	6468	9531	3.11%	0.257%
77	20.735	3044	3046	3050	rVB4	2677	3184	1.04%	0.086%
78	20.782	3050	3054	3059	rBV3	10169	11916	3.89%	0.321%
79	21.064	3098	3102	3107	rBV2	10497	17762	5.80%	0.479%
80	21.252	3129	3134	3136	rBV3	10324	15165	4.95%	0.409%
81	21.287	3136	3140	3145	rVB2	32942	46426	15.16%	1.251%
82	21.399	3155	3159	3168	rVB3	9412	20781	6.79%	0.560%
83	21.463	3168	3170	3171	rBV2	2579	1630	0.53%	0.044%
84	21.534	3180	3182	3187	rVB3	7997	9331	3.05%	0.251%
85	21.669	3197	3205	3210	rBV2	158962	306154	100.00%	8.248%
86	21.716	3210	3213	3218	rVB	39709	58317	19.05%	1.571%
87	21.827	3227	3232	3237	rBV3	23187	34340	11.22%	0.925%
88	22.068	3270	3273	3274	rBV2	4846	4167	1.36%	0.112%
89	22.392	3326	3328	3330	rBV2	4557	5248	1.71%	0.141%
90	22.427	3330	3334	3339	rVB2	19972	28305	9.25%	0.763%
91	23.108	3446	3450	3457	rVB4	18796	33935	11.08%	0.914%
92	23.296	3480	3482	3485	rBV4	3281	3394	1.11%	0.091%
93	23.866	3571	3579	3582	rBV2	29726	77519	25.32%	2.088%
94	24.589	3695	3702	3706	rBV	14953	35047	11.45%	0.944%
95	24.659	3707	3714	3721	rVV2	83486	217706	71.11%	5.865%
96	24.724	3722	3725	3733	rVB2	21351	50072	16.36%	1.349%
97	24.883	3744	3752	3761	rVB3	75608	201698	65.88%	5.434%
98	25.940	3930	3932	3939	rVB7	10978	20626	6.74%	0.556%
99	30.447	4697	4699	4700	rBV2	2491	1164	0.38%	0.031%
100	31.910	4947	4948	4949	rBV	2738	1151	0.38%	0.031%

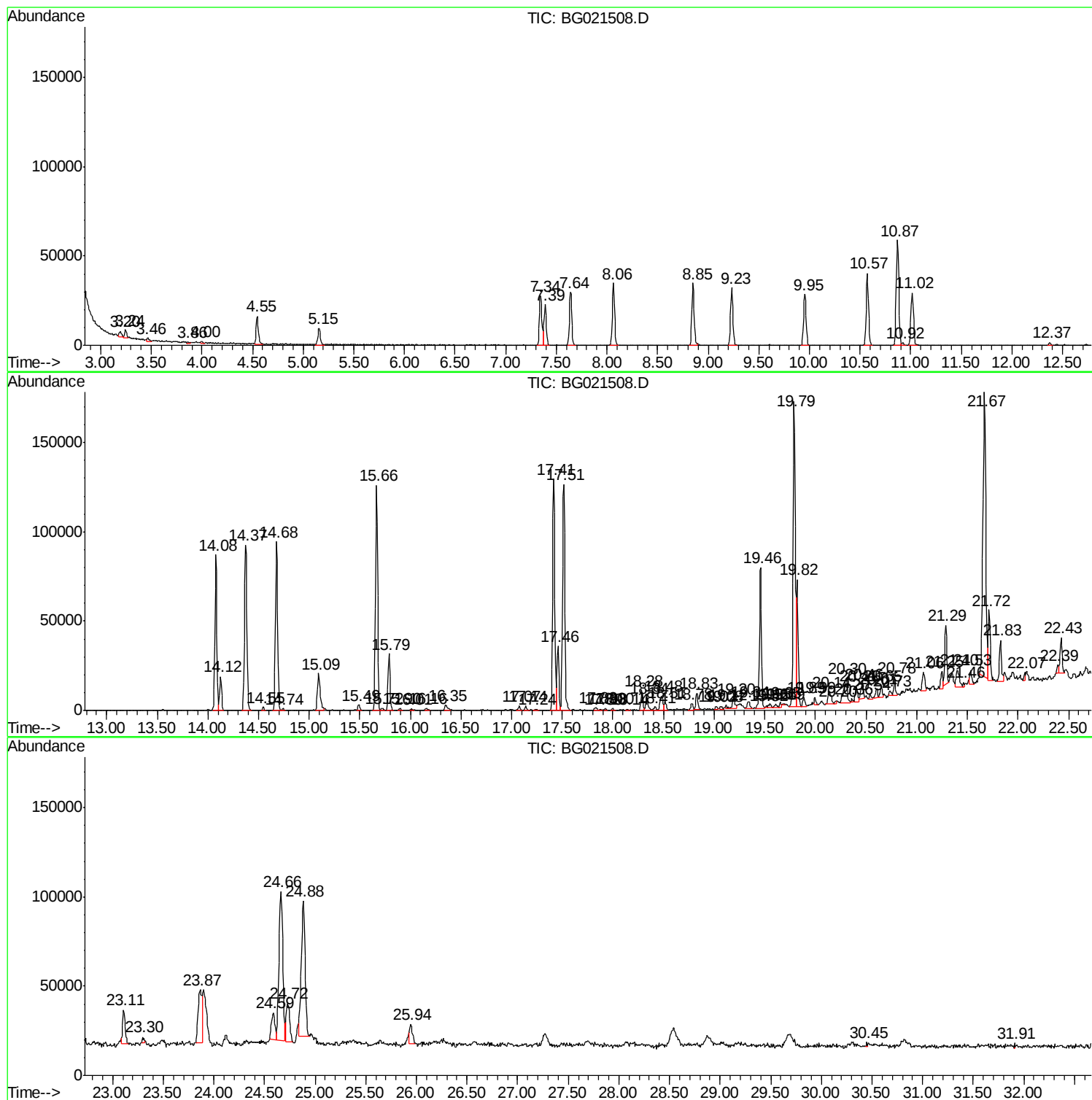
Sum of corrected areas: 3712004

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG031616.M
Quant Title : SVOA CALIBRATION

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



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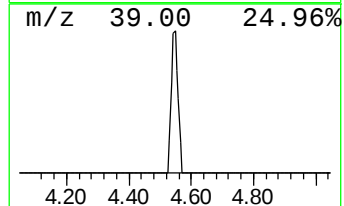
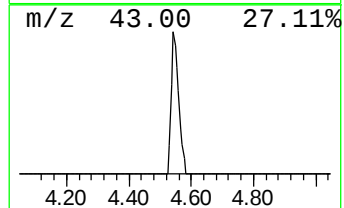
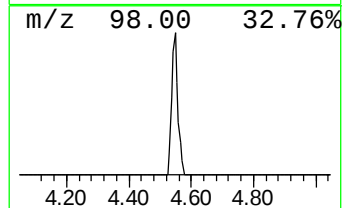
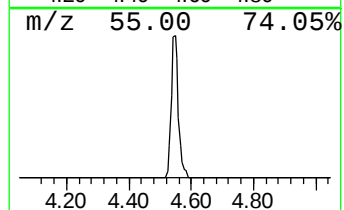
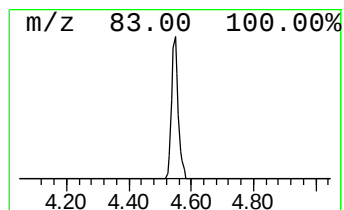
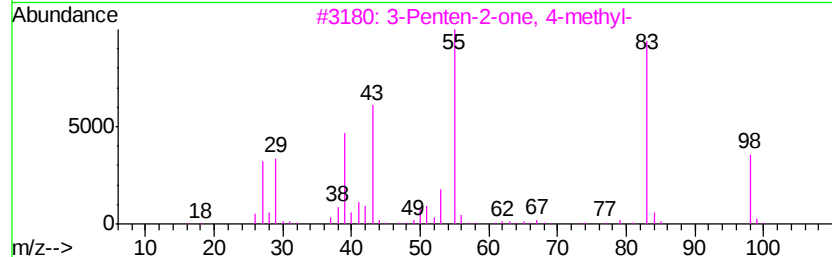
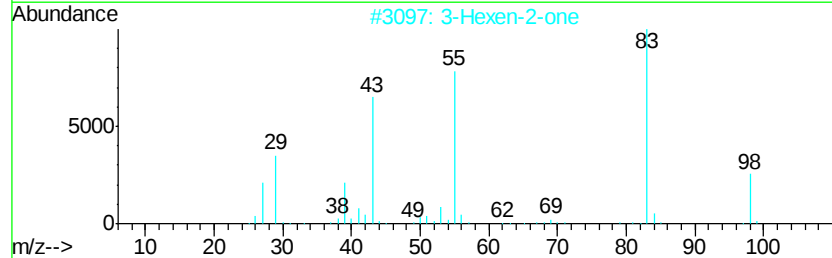
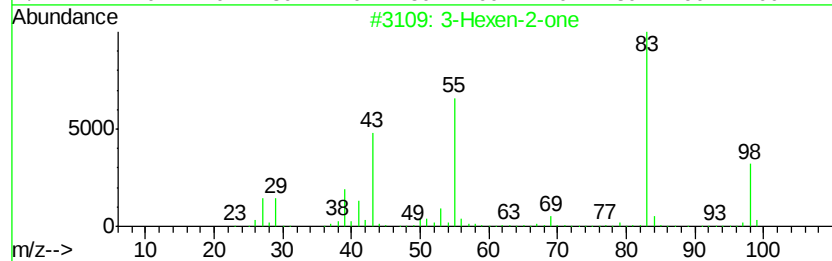
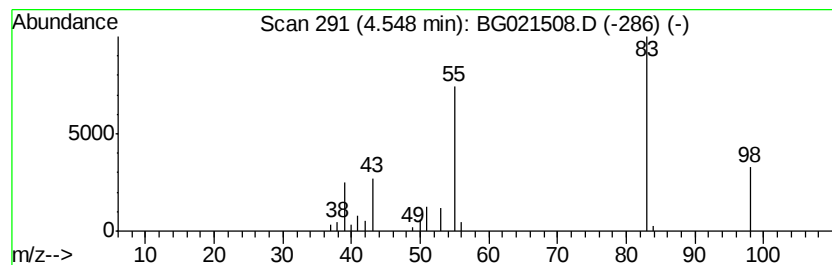
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG031616.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 3-Hexen-2-one Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.55	8.02 ng/ul	23381	1,4-Dichlorobenzene-d4	8.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Hexen-2-one	98	C6H10O	000763-93-9	86
2		3-Hexen-2-one	98	C6H10O	000763-93-9	78
3		3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	74
4		3-Hexen-2-one	98	C6H10O	000763-93-9	72
5		2-Pentene, 2,4-dimethyl-	98	C7H14	000625-65-0	72



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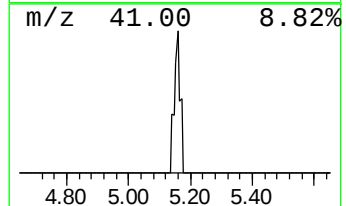
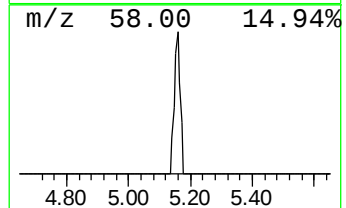
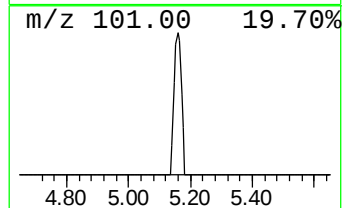
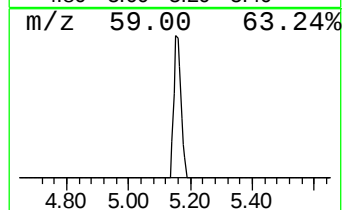
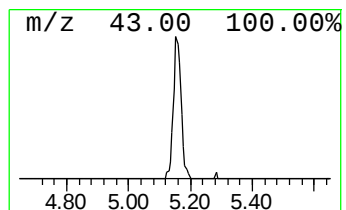
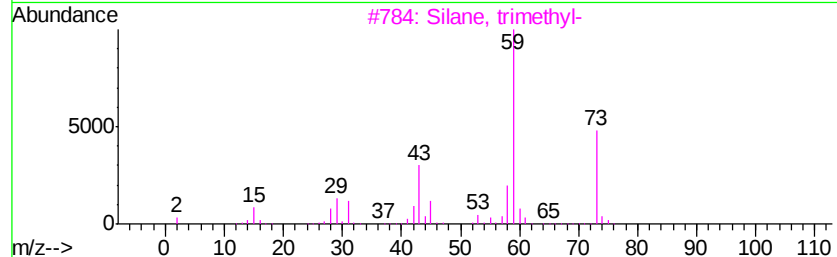
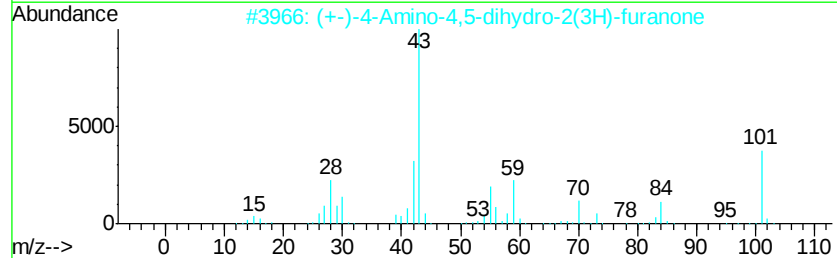
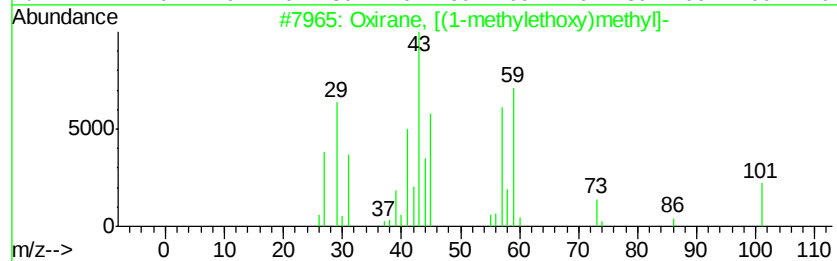
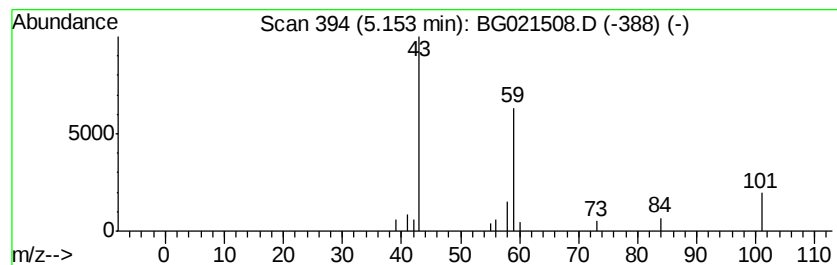
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown-01 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.15	5.43 ng/ul	15818	1,4-Dichlorobenzene-d4	8.07

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxirane, [(1-methylethoxy)methyl]-	116	C6H12O2	004016-14-2	43
2			(+)-4-Amino-4,5-dihydro-2(3H)-furanone	101	C4H7NO2	016504-58-8	9
3			Silane, trimethyl-	74	C3H10Si	000993-07-7	9
4			3-Octanol	130	C8H18O	000589-98-0	9
5			2-Nonanone	142	C9H18O	000821-55-6	9



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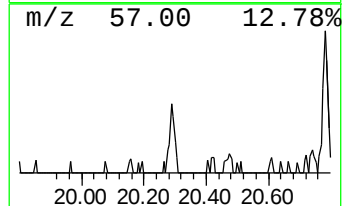
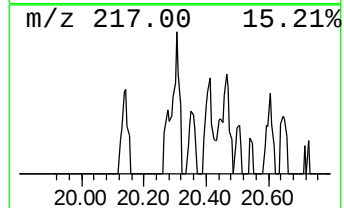
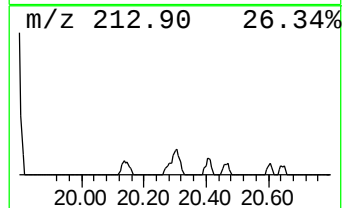
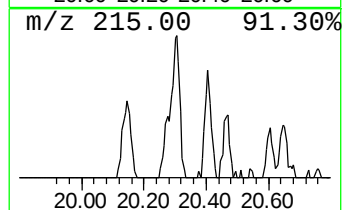
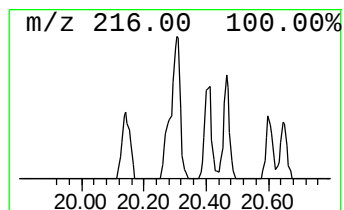
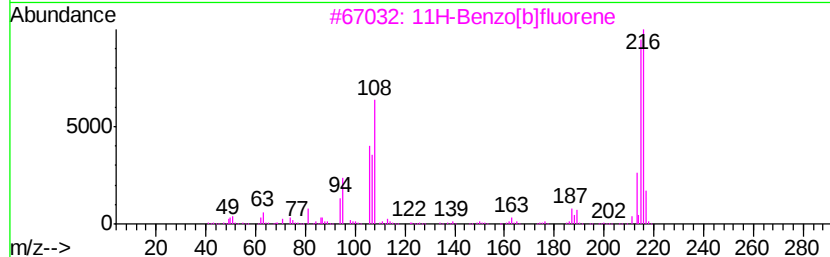
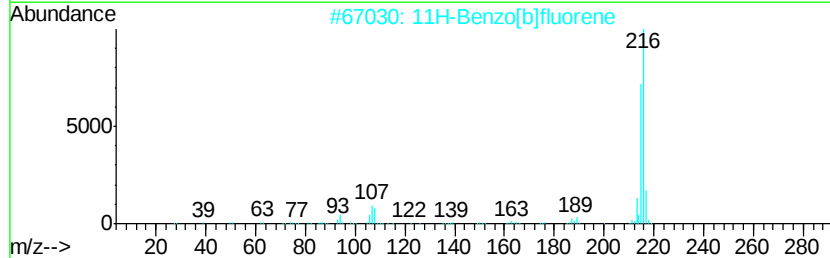
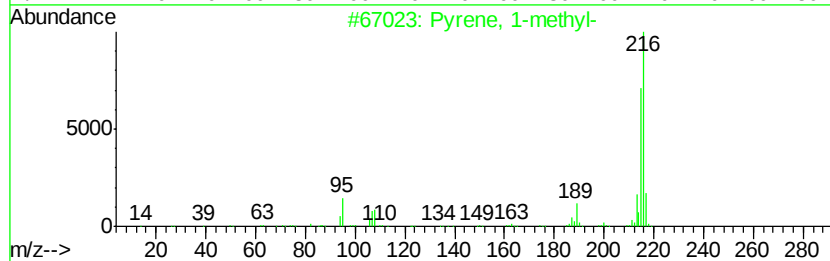
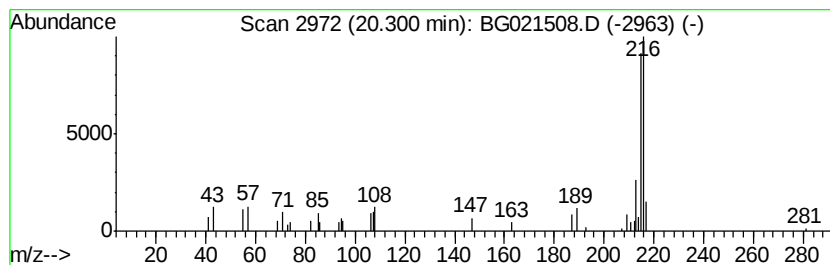
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Peak Number 3 Pyrene, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.30	2.26 ng/ul	34607	Chrysene-d12	21.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pyrene, 1-methyl-	216	C17H12	002381-21-7	90
2		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	87
3		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	87
4		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	87
5		7H-Benzo[c]fluorene	216	C17H12	000205-12-9	74



Data Path : Z:\HPCHEM1\BNA G\DATA\BG032516\
Data File : BG021508.D
Acq On : 25 Mar 2016 10:21
Operator : UM/SJ
Sample : H1909-04
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4T40

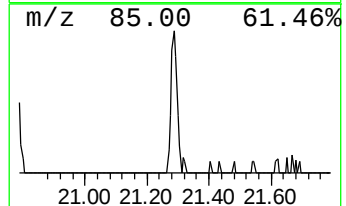
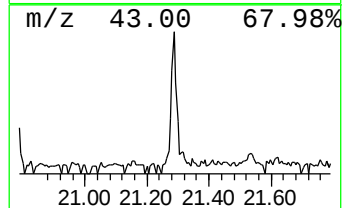
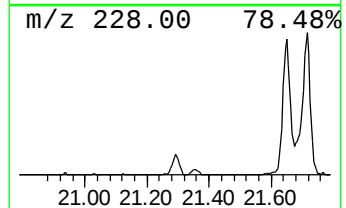
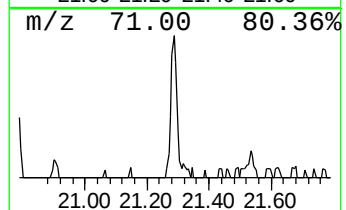
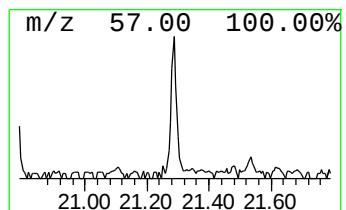
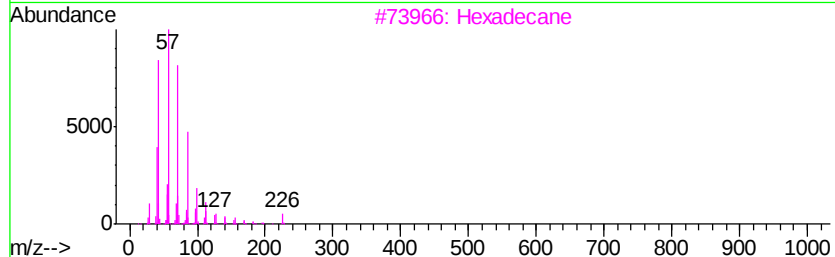
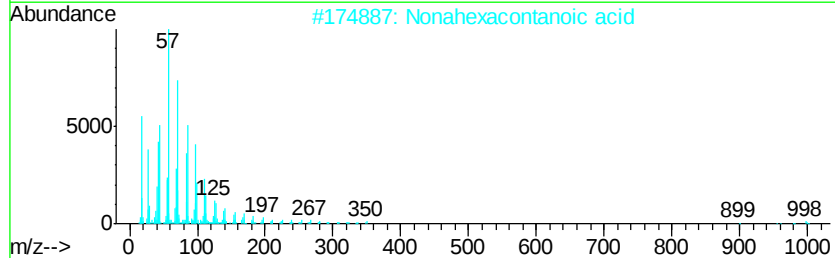
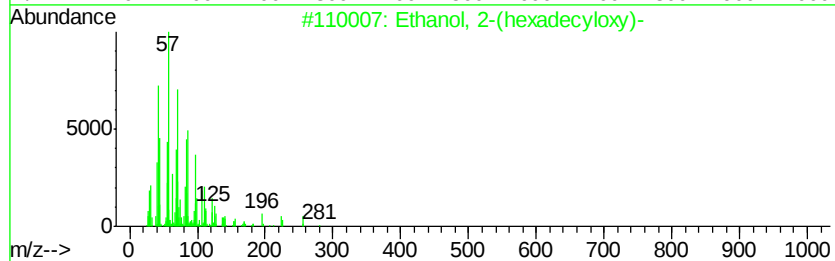
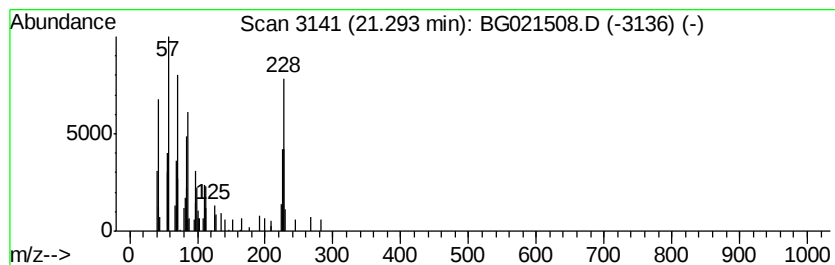
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Ethanol, 2-(hexadecyloxy)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.29	3.03 ng/ul	46426	Chrysene-d12	21.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanol, 2-(hexadecyloxy)-	286	C18H38O2	002136-71-2	52
2		Nonahexacontanoic acid	999	C69H138O2	040710-32-5	49
3		Hexadecane	226	C16H34	000544-76-3	42
4		Hexadecane	226	C16H34	000544-76-3	42
5		Trichloroacetic acid, tetradecyl...	358	C16H29Cl3O2	074339-52-9	42



Data Path : Z:\HPCHEM1\BNA G\DATA\BG032516\
Data File : BG021508.D
Acq On : 25 Mar 2016 10:21
Operator : UM/SJ
Sample : H1909-04
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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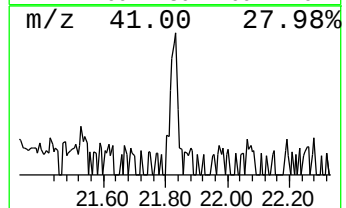
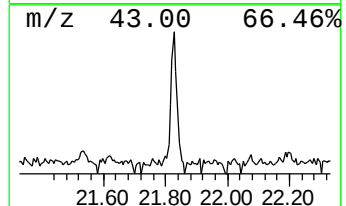
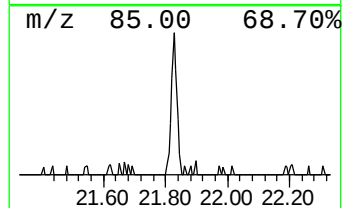
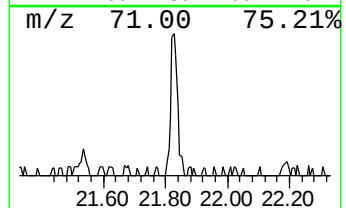
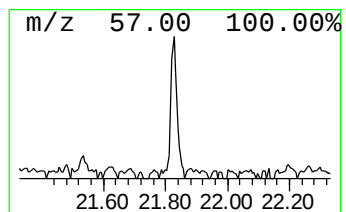
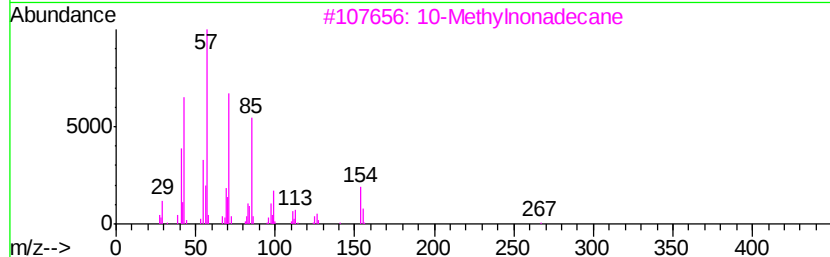
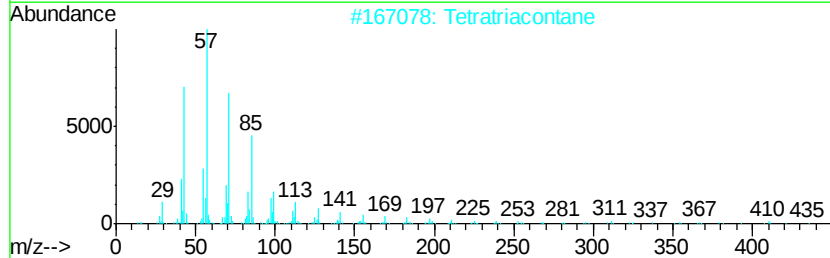
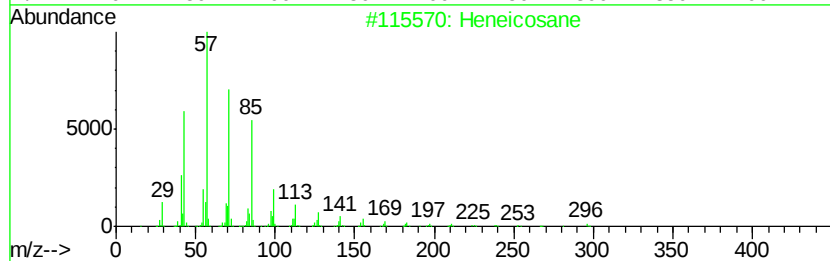
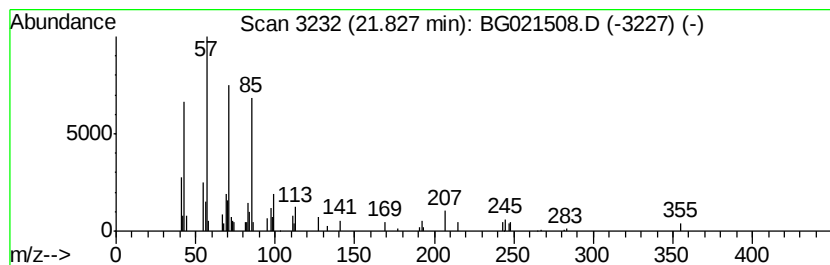
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.83	2.24 ng/ul	34340	Chrysene-d12	21.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	74
2		Tetratriacontane	479	C34H70	014167-59-0	74
3		10-Methylnonadecane	282	C20H42	056862-62-5	74
4		Octacosane	394	C28H58	000630-02-4	74
5		Heptacosane	380	C27H56	000593-49-7	74



Data Path : Z:\HPCHEM1\BNA G\DATA\BG032516\
Data File : BG021508.D
Acq On : 25 Mar 2016 10:21
Operator : UM/SJ
Sample : H1909-04
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4T40

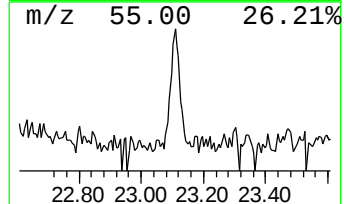
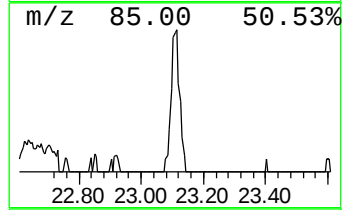
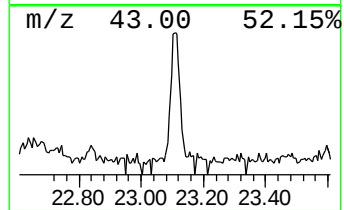
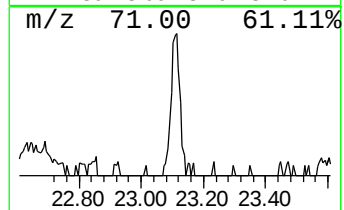
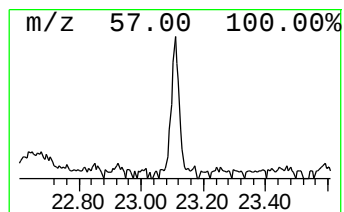
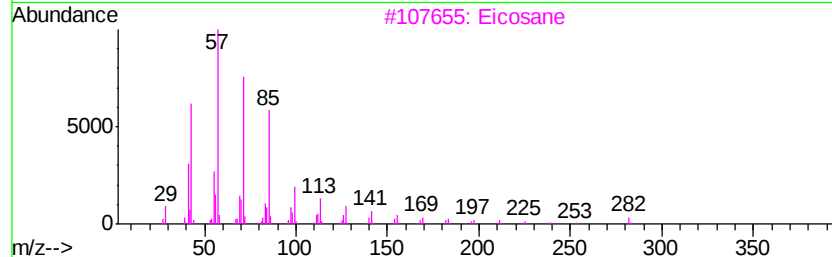
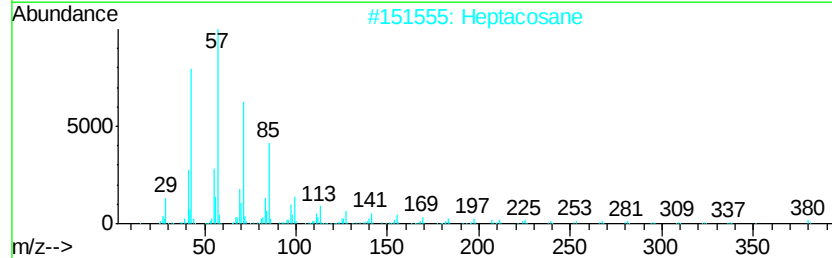
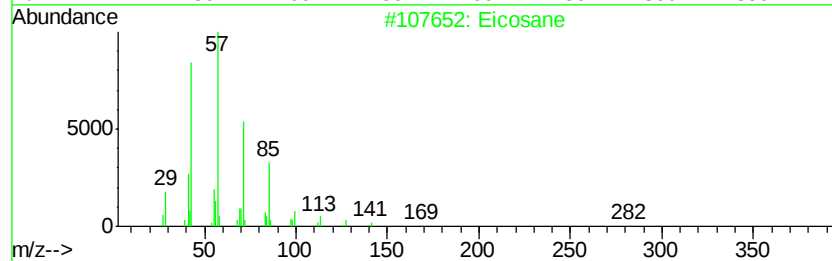
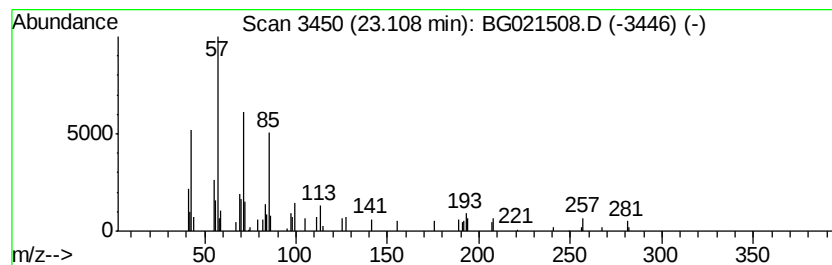
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.11	2.22 ng/ul	33935	Chrysene-d12	21.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	95
2		Heptacosane	380	C27H56	000593-49-7	81
3		Eicosane	282	C20H42	000112-95-8	76
4		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	74
5		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	74



Data Path : Z:\HPCHEM1\BNA G\DATA\BG032516\
Data File : BG021508.D
Acq On : 25 Mar 2016 10:21
Operator : UM/SJ
Sample : H1909-04
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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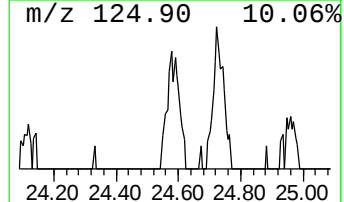
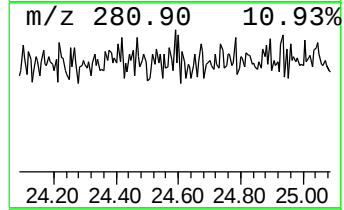
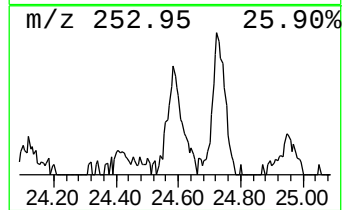
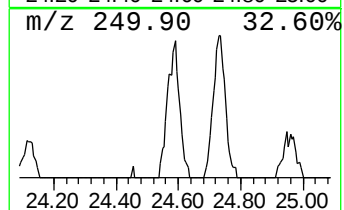
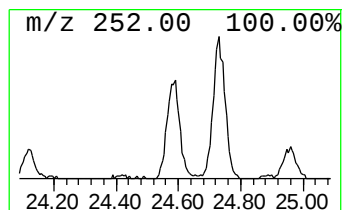
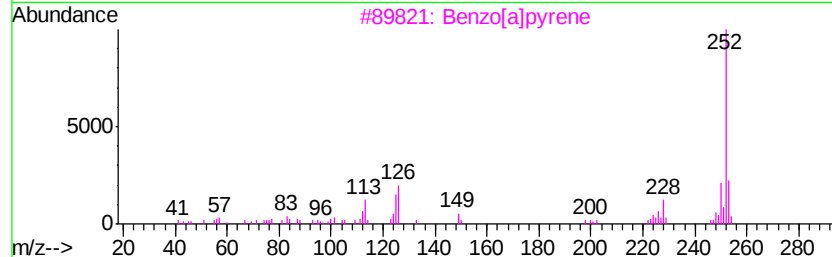
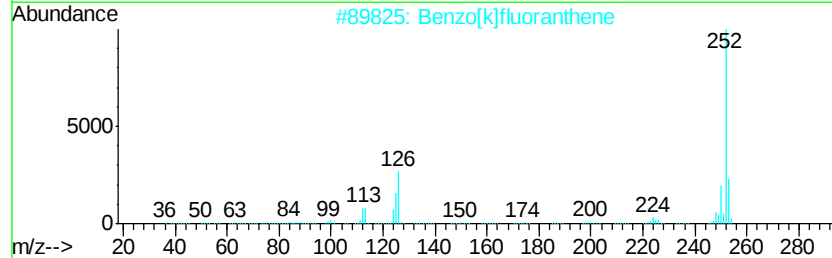
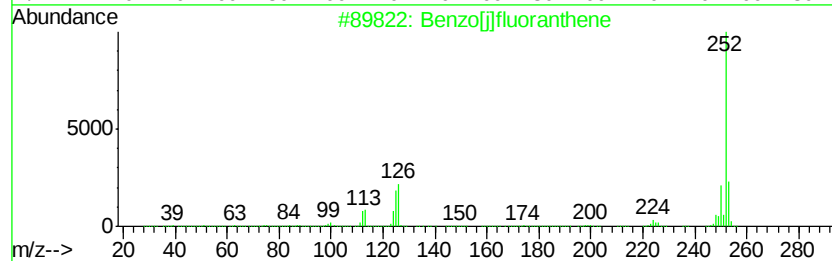
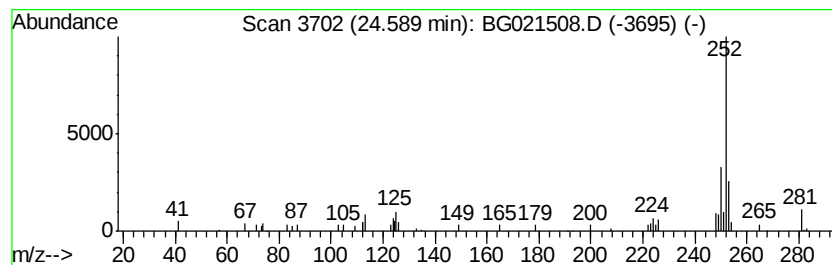
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Benzo[j]fluoranthene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.59	3.48 ng/ul	35047	Perylene-d12	24.88

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzo[i]fluoranthene	252	C20H12	000205-82-3	70
2		Benzo[k]fluoranthene	252	C20H12	000207-08-9	70
3		Benzo[a]pyrene	252	C20H12	000050-32-8	70
4		Perylene	252	C20H12	000198-55-0	70
5		Benz[e]acephenanthrylene	252	C20H12	000205-99-2	70



Data Path : Z:\HPCHEM1\BNA G\DATA\BG032516\
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Operator : UM/SJ
Sample : H1909-04
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4T40

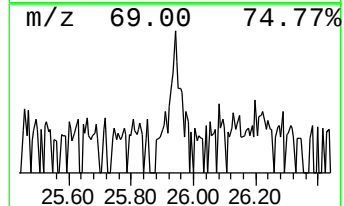
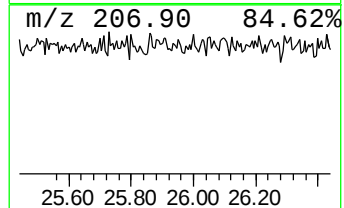
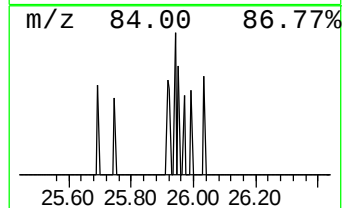
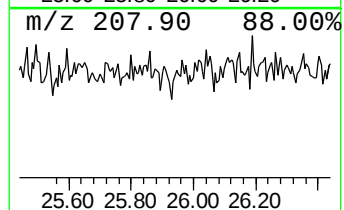
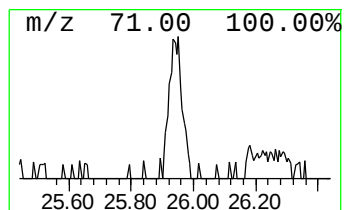
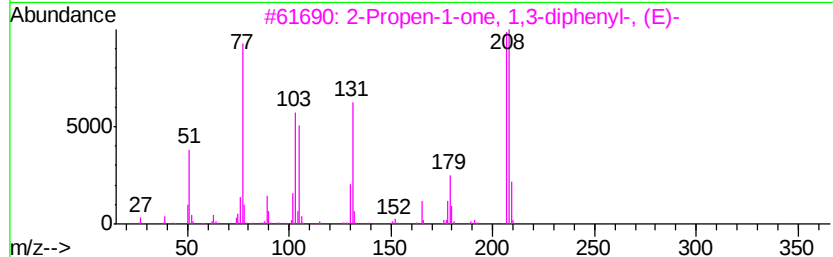
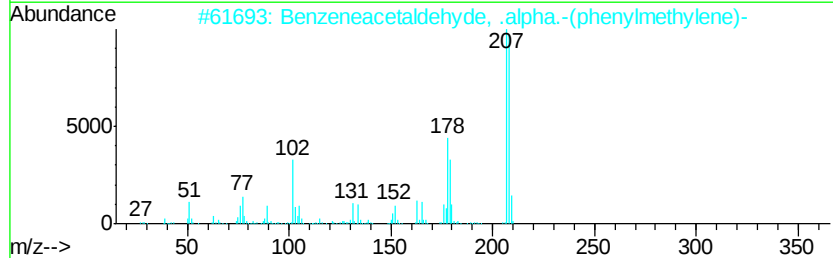
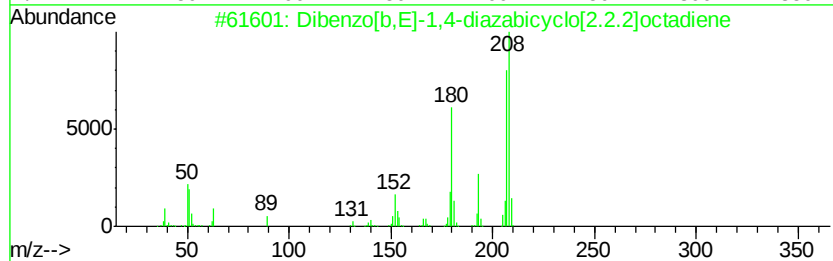
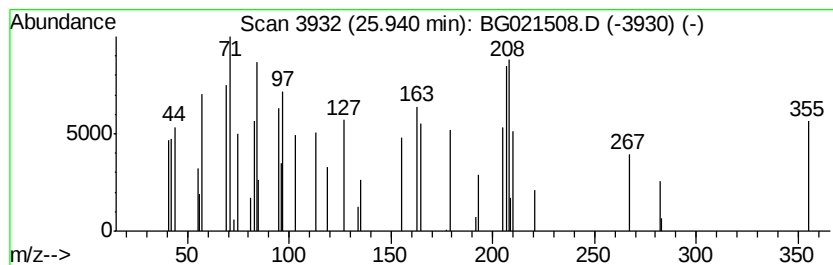
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-02 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.94	2.05 ng/ul	20626	Perylene-d12	24.88

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Dibenzo[b,E]-1,4-diazabicyclo[2....	208	C14H12N2	094509-68-9	25
2		Benzeneacetaldehyde, .alpha.-(ph...	208	C15H12O	013702-35-7	22
3		2-Propen-1-one, 1,3-diphenyl-, (E)-	208	C15H12O	000614-47-1	22
4		3-Phenyl-1-indanone	208	C15H12O	016618-72-7	18
5		Indole, 6-methyl-2-(2-pyridyl)-	208	C14H12N2	107919-89-1	14



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG032516\
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ALS Vial : 15 Sample Multiplier: 1

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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG031616.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
3-Hexen-2-one	4.55	8.0	ng/ul	23381	1	8.07	58308	20.0
unknown-01	5.15	5.4	ng/ul	15818	1	8.07	58308	20.0
Pyrene, 1-methyl-	20.30	2.3	ng/ul	34607	5	21.67	306154	20.0
Ethanol, 2-(hexad...	21.29	3.0	ng/ul	46426	5	21.67	306154	20.0
(DEL) Alkane: Str...	21.83	2.2	ng/ul	34340	5	21.67	306154	20.0
(DEL) Alkane: Str...	23.11	2.2	ng/ul	33935	5	21.67	306154	20.0
Benzo[j]fluoranthene	24.59	3.5	ng/ul	35047	6	24.88	201698	20.0
unknown-02	25.94	2.0	ng/ul	20626	6	24.88	201698	20.0