

Data Path : Z:\HPCHEM1\BNA G\DATA\BG110915\
 Data File : BG019535.D
 Acq On : 9 Nov 2015 16:31
 Operator : UM/NP
 Sample : G4245-17DL2 50X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BCY52DL2

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-BG102815.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	59	61	64	rVB3	9910	11372	0.50%	0.117%
2	5.634	470	476	482	rBV	21746	33288	1.45%	0.341%
3	5.769	494	499	501	rBV2	13528	24080	1.05%	0.247%
4	6.151	558	564	570	rVB2	18548	31361	1.37%	0.322%
5	6.633	641	646	650	rBV4	4077	6321	0.28%	0.065%
6	7.244	745	750	753	rBV3	7590	11390	0.50%	0.117%
7	7.303	756	760	766	rVB4	6712	12175	0.53%	0.125%
8	7.379	768	773	781	rBV3	13267	20330	0.89%	0.209%
9	7.556	798	803	806	rVB3	3717	5233	0.23%	0.054%
10	7.808	840	846	852	rVB3	33425	57012	2.48%	0.585%
11	7.890	854	860	866	rBV	27220	45468	1.98%	0.466%
12	8.172	901	908	915	rBV	120817	201232	8.76%	2.064%
13	8.290	923	928	934	rVB3	8936	15820	0.69%	0.162%
14	8.549	964	972	979	rBV	124075	211104	9.19%	2.165%
15	8.713	995	1000	1009	rVB	110205	190981	8.32%	1.959%
16	9.019	1047	1052	1058	rVV4	4447	8424	0.37%	0.086%
17	9.277	1092	1096	1098	rBV2	5896	6603	0.29%	0.068%
18	9.365	1103	1111	1117	rVB5	8594	20312	0.88%	0.208%
19	9.541	1135	1141	1146	rBV2	10903	19100	0.83%	0.196%
20	9.659	1155	1161	1168	rBV4	22686	48200	2.10%	0.494%
21	9.729	1169	1173	1179	rVB4	7568	12158	0.53%	0.125%
22	9.865	1194	1196	1206	rVB6	3305	6722	0.29%	0.069%
23	10.241	1255	1260	1264	rBV3	12136	19593	0.85%	0.201%
24	10.388	1279	1285	1296	rBV7	21794	54532	2.37%	0.559%
25	10.487	1296	1302	1308	rBV5	9809	17994	0.78%	0.185%
26	10.623	1319	1325	1328	rBV6	4853	9486	0.41%	0.097%
27	11.004	1382	1390	1393	rBV	165601	297046	12.94%	3.047%
28	11.057	1393	1399	1407	rVV	1284940	2296350	100.00%	23.554%
29	11.175	1410	1419	1427	rVB	97711	183400	7.99%	1.881%
30	11.410	1455	1459	1467	rBV3	3729	7156	0.31%	0.073%
31	12.097	1573	1576	1585	rVB5	4416	8604	0.37%	0.088%
32	12.373	1619	1623	1626	rBV3	5288	6917	0.30%	0.071%
33	12.544	1648	1652	1656	rVB2	7376	9347	0.41%	0.096%
34	12.644	1661	1669	1676	rBV	53073	86452	3.76%	0.887%

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35	12.761	1686	1689	1696	rVV2	5828	8744	0.38%	0.090%
36	12.867	1696	1707	1714	rVB	138370	240717	10.48%	2.469%
37	13.643	1832	1839	1845	rBV2	45184	71550	3.12%	0.734%
38	13.842	1867	1873	1876	rBV4	9400	15748	0.69%	0.162%
39	13.966	1890	1894	1895	rBV	12967	12088	0.53%	0.124%
40	14.124	1914	1921	1926	rBV3	25815	46419	2.02%	0.476%
41	14.189	1926	1932	1937	rVB3	15296	34993	1.52%	0.359%
42	14.359	1956	1961	1964	rBV4	10187	15040	0.65%	0.154%
43	14.495	1979	1984	1987	rBV3	9941	16004	0.70%	0.164%
44	14.806	2023	2037	2043	rBV2	307517	511356	22.27%	5.245%
45	14.871	2043	2048	2054	rVB	153183	232560	10.13%	2.385%
46	14.929	2054	2058	2062	rVB3	12145	16927	0.74%	0.174%
47	14.982	2062	2067	2074	rBV	26951	40197	1.75%	0.412%
48	15.070	2078	2082	2086	rBV3	13866	22448	0.98%	0.230%
49	15.200	2098	2104	2113	rBV	113996	171977	7.49%	1.764%
50	15.270	2113	2116	2119	rBV4	6980	9178	0.40%	0.094%
51	15.570	2165	2167	2172	rBV	3354	4942	0.22%	0.051%
52	15.793	2200	2205	2209	rVV3	16934	24494	1.07%	0.251%
53	15.846	2209	2214	2220	rVV2	95081	141731	6.17%	1.454%
54	15.975	2227	2236	2245	rVV8	25939	67429	2.94%	0.692%
55	16.069	2247	2252	2260	rVV8	11657	29866	1.30%	0.306%
56	16.134	2260	2263	2268	rVB2	9037	10826	0.47%	0.111%
57	16.222	2274	2278	2281	rBV4	3933	5780	0.25%	0.059%
58	16.281	2281	2288	2296	rVB4	9439	23833	1.04%	0.244%
59	16.516	2322	2328	2333	rVB4	17783	29666	1.29%	0.304%
60	16.721	2355	2363	2371	rBV5	17931	34720	1.51%	0.356%
61	16.792	2371	2375	2380	rBV2	16296	27823	1.21%	0.285%
62	16.898	2388	2393	2397	rVB5	7713	15275	0.67%	0.157%
63	16.997	2407	2410	2411	rBV2	5301	4841	0.21%	0.050%
64	17.109	2425	2429	2434	rVB5	3928	6948	0.30%	0.071%
65	17.321	2460	2465	2470	rBV4	28429	47580	2.07%	0.488%
66	17.368	2470	2473	2476	rVB2	9420	11076	0.48%	0.114%
67	17.426	2478	2483	2485	rBV3	6103	8195	0.36%	0.084%
68	17.491	2490	2494	2495	rVV2	8038	9832	0.43%	0.101%
69	17.544	2497	2503	2508	rVV	481131	767459	33.42%	7.872%
70	17.591	2508	2511	2516	rVV	121924	167876	7.31%	1.722%
71	17.644	2516	2520	2523	rVV	22551	34280	1.49%	0.352%

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72	17.679	2523	2526	2530	rVV3	11783	15890	0.69%	0.163%
73	17.750	2533	2538	2544	rVB6	7964	14414	0.63%	0.148%
74	17.902	2560	2564	2566	rVV3	14681	19127	0.83%	0.196%
75	17.943	2566	2571	2578	rVV	139553	218445	9.51%	2.241%
76	18.032	2581	2586	2593	rVV3	26888	58992	2.57%	0.605%
77	18.096	2593	2597	2602	rVV2	30646	46686	2.03%	0.479%
78	18.190	2605	2613	2618	rBV9	12298	26388	1.15%	0.271%
79	18.237	2618	2621	2624	rVV4	6037	6758	0.29%	0.069%
80	18.290	2624	2630	2633	rVV6	9722	15262	0.66%	0.157%
81	18.337	2633	2638	2641	rVV6	4610	10254	0.45%	0.105%
82	18.378	2641	2645	2648	rVV4	16533	25033	1.09%	0.257%
83	18.460	2655	2659	2666	rVB5	37306	64911	2.83%	0.666%
84	18.531	2666	2671	2677	rBV2	23210	36700	1.60%	0.376%
85	18.607	2682	2684	2689	rBV2	6430	10835	0.47%	0.111%
86	18.684	2694	2697	2699	rBV	9364	9973	0.43%	0.102%
87	18.784	2712	2714	2718	rVV3	5866	6003	0.26%	0.062%
88	18.848	2721	2725	2730	rVV4	16592	26909	1.17%	0.276%
89	18.901	2731	2734	2739	rVB2	11424	16201	0.71%	0.166%
90	18.983	2746	2748	2753	rBV5	4630	6457	0.28%	0.066%
91	19.107	2766	2769	2772	rBV4	6492	7944	0.35%	0.081%
92	19.295	2799	2801	2806	rVV5	4503	5658	0.25%	0.058%
93	19.506	2835	2837	2841	rVB3	6882	6346	0.28%	0.065%
94	19.589	2847	2851	2854	rVB3	9641	13608	0.59%	0.140%
95	19.918	2903	2907	2910	rBV2	24260	30372	1.32%	0.312%
96	20.370	2979	2984	2990	rBV2	74699	115953	5.05%	1.189%
97	21.275	3136	3138	3142	rBV4	6568	9853	0.43%	0.101%
98	21.827	3225	3232	3240	rVB	612138	1025156	44.64%	10.515%
99	24.377	3664	3666	3668	rBV2	7180	5605	0.24%	0.057%
100	25.176	3792	3802	3814	rVB	328989	957402	41.69%	9.820%

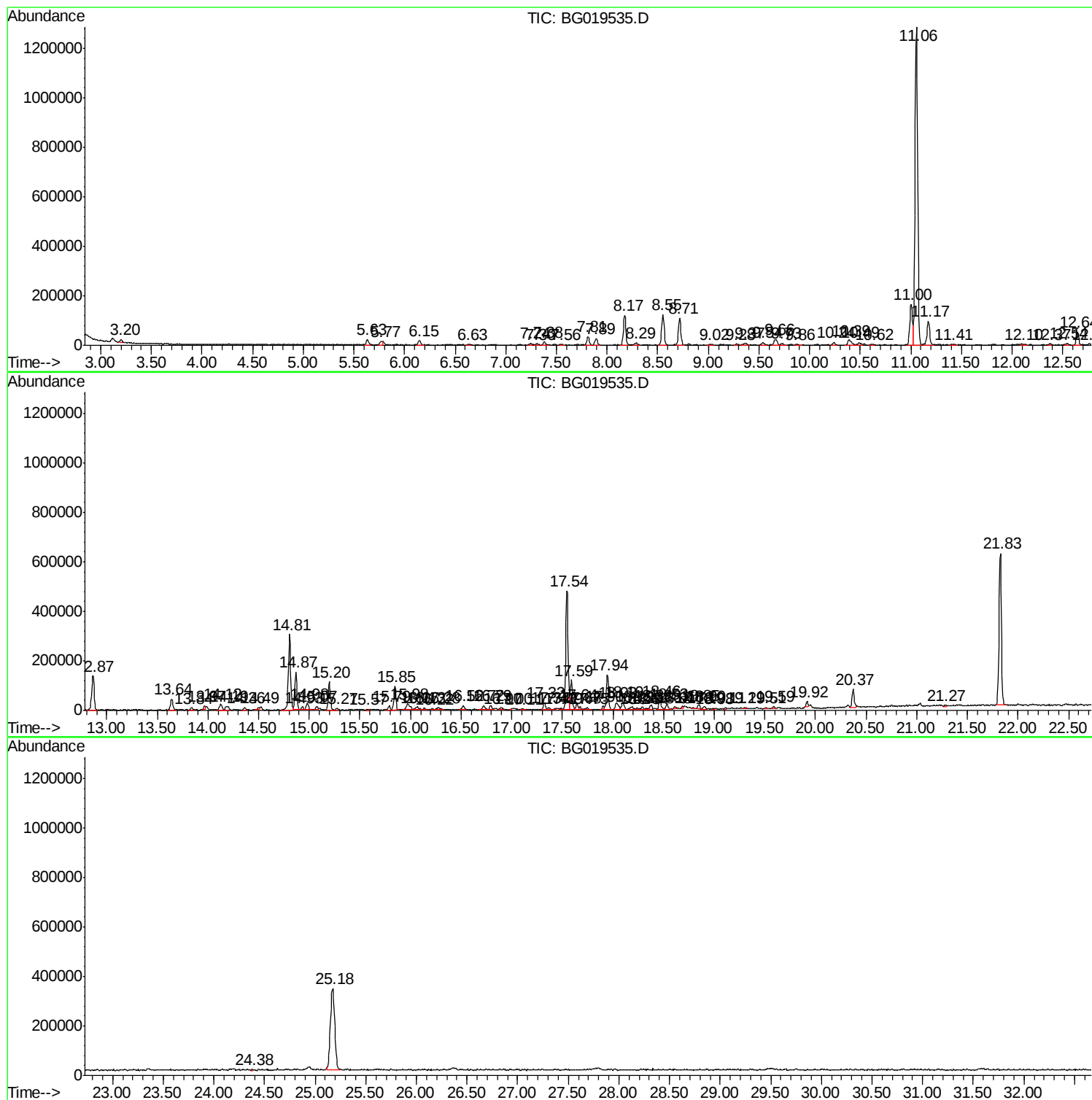
Sum of corrected areas: 9749116

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG102815.M
Quant Title : SVOA CALIBRATION

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



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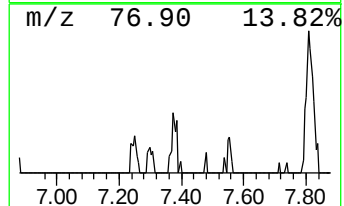
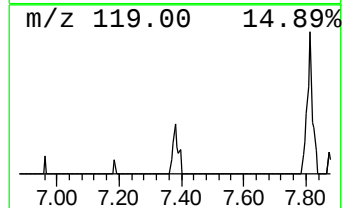
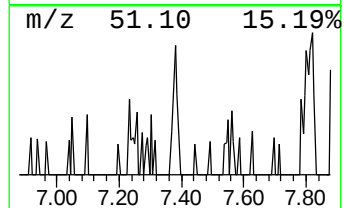
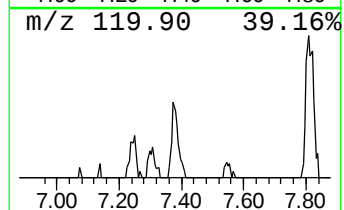
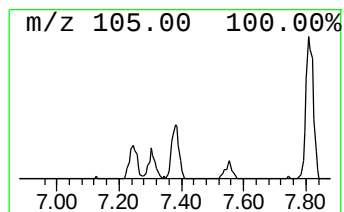
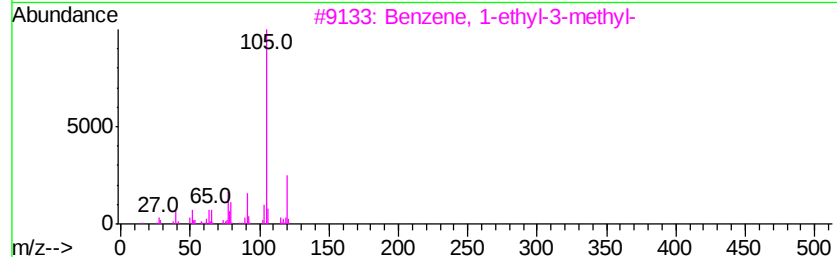
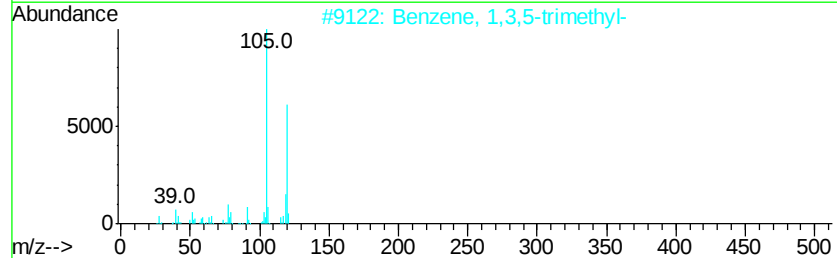
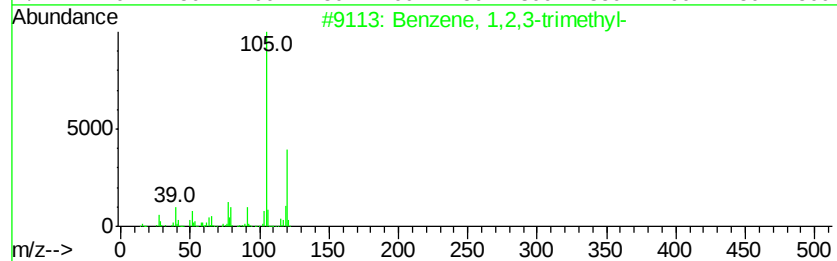
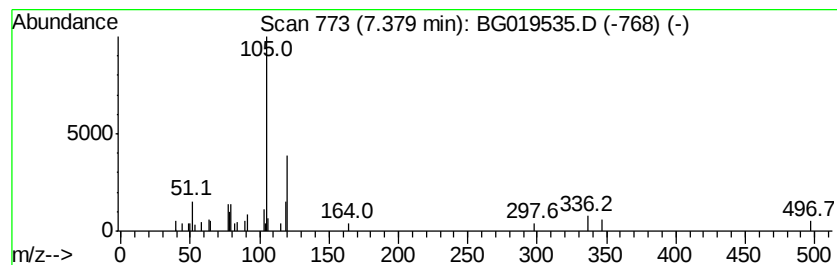
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG102815.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 Benzene, 1,2,3-trimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.38	2.02 ng/ul	20330	1,4-Dichlorobenzene-d4	8.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	58
2		Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	52
3		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	52
4		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	52
5		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	50



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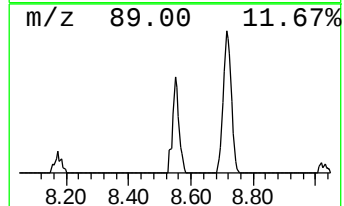
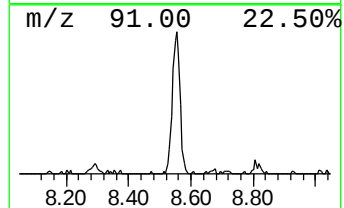
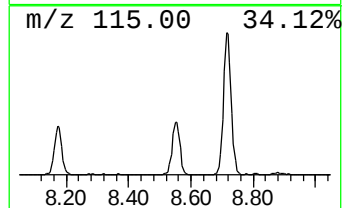
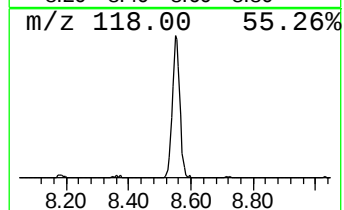
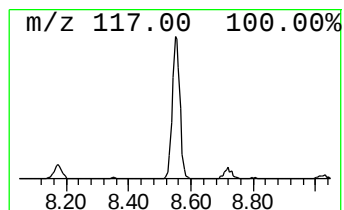
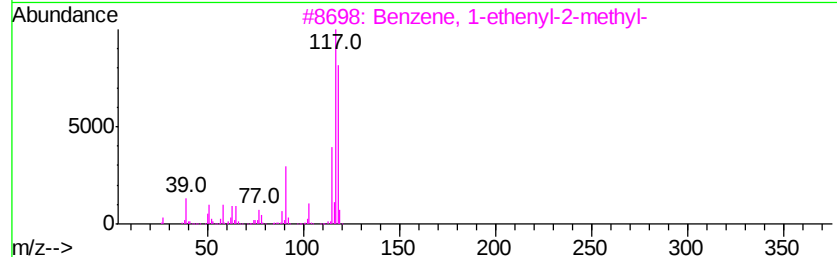
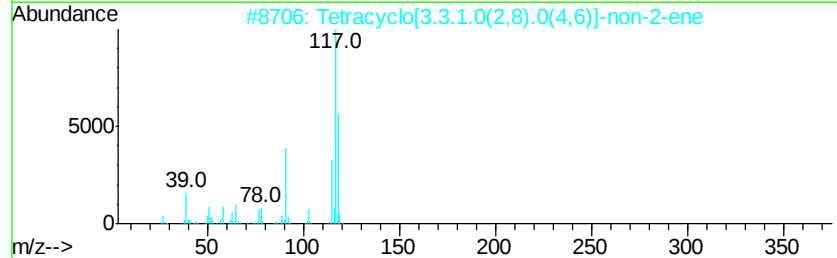
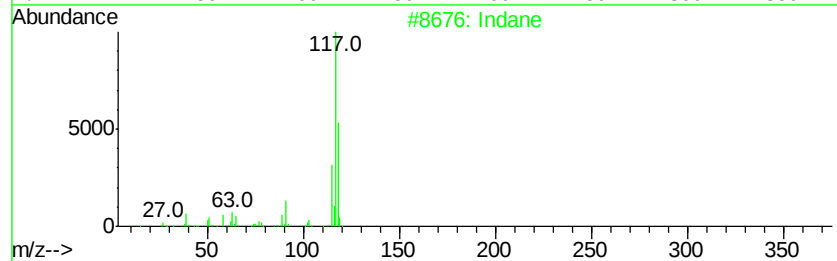
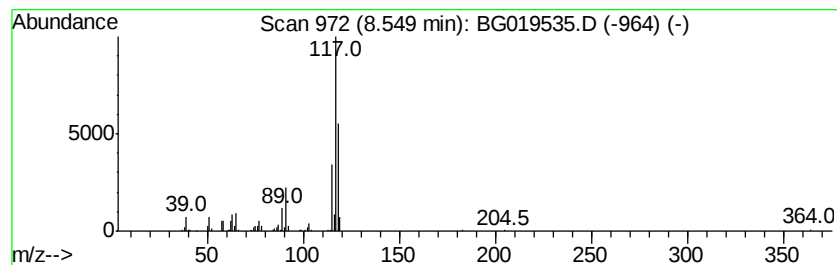
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG102815.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.55	20.98 ng/ul	211104	1,4-Dichlorobenzene-d4	8.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indane	118	C9H10	000496-11-7	87
2		Tetracyclo[3.3.1.0(2,8).0(4,6)]-	118	C9H10	1000191-13-7	81
3		Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	81
4		Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	81
5		Indane	118	C9H10	000496-11-7	81



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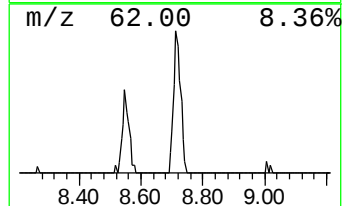
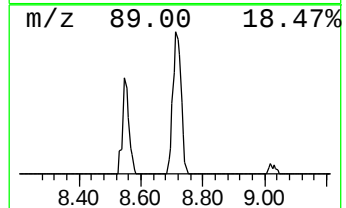
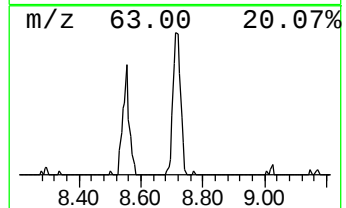
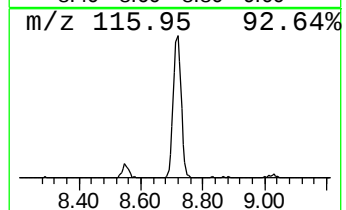
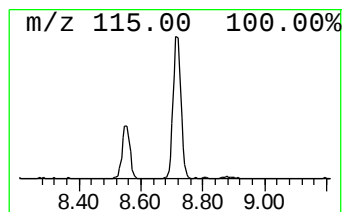
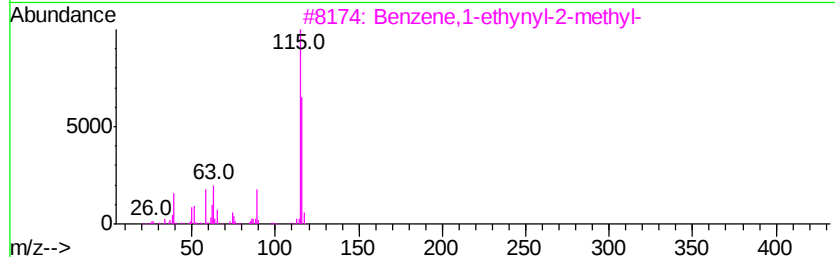
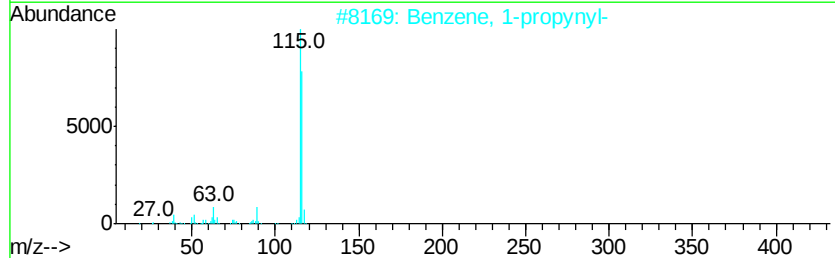
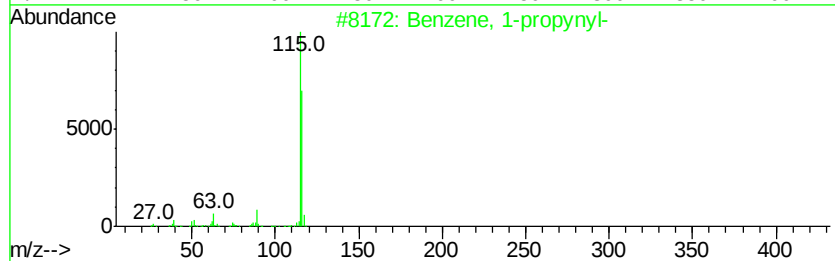
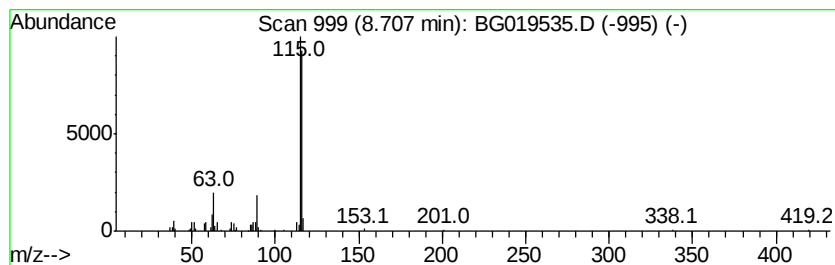
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Peak Number 8 Benzene, 1-propynyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.71	18.98 ng/ul	190981	1,4-Dichlorobenzene-d4	8.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-propynyl-	116	C9H8	000673-32-5	94
2		Benzene, 1-propynyl-	116	C9H8	000673-32-5	91
3		Benzene,1-ethynyl-2-methyl-	116	C9H8	000766-47-2	91
4		Indene	116	C9H8	000095-13-6	91
5		Indene	116	C9H8	000095-13-6	91



Data Path : Z:\HPCHEM1\BNA G\DATA\BG110915\
Data File : BG019535.D
Acq On : 9 Nov 2015 16:31
Operator : UM/NP
Sample : G4245-17DL2 50X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BCY52DL2

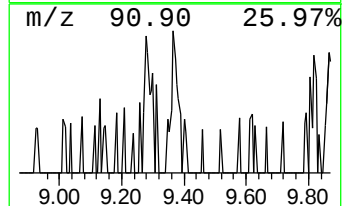
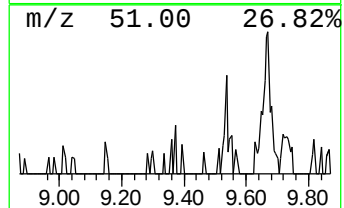
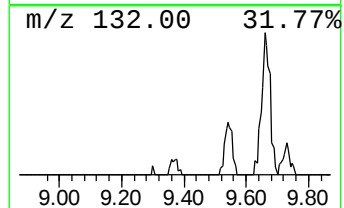
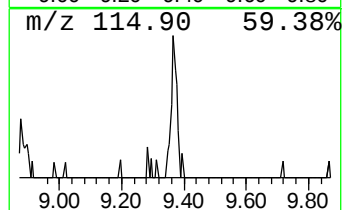
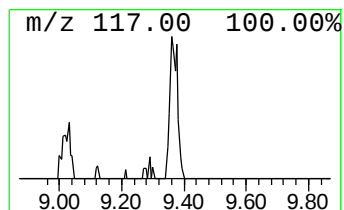
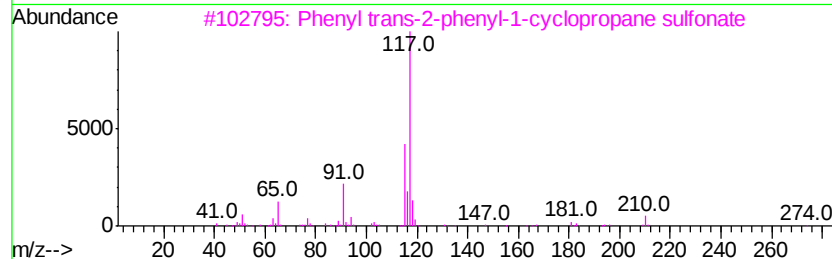
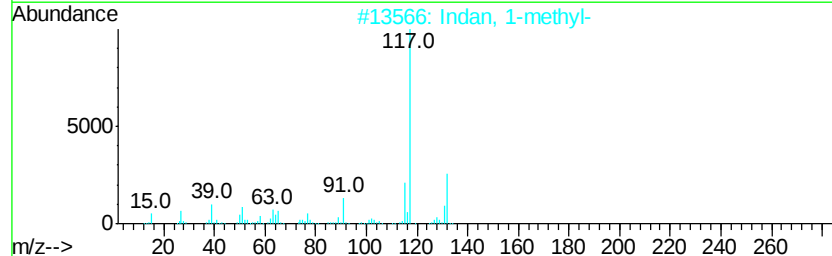
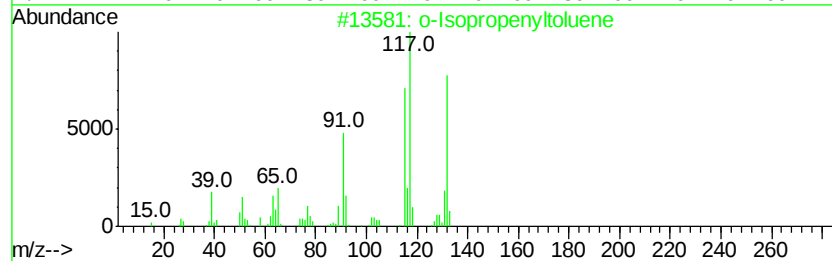
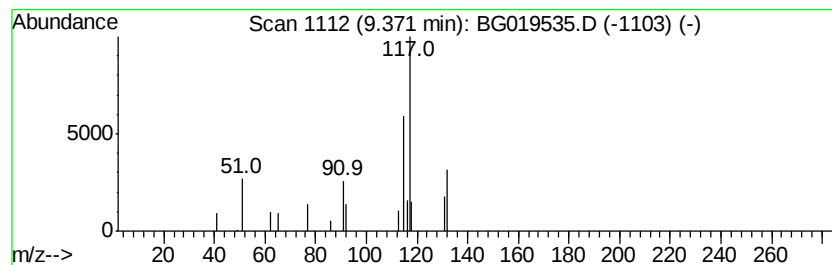
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 unknown-01 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.37	2.02 ng/ul	20312	1,4-Dichlorobenzene-d4	8.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	o-Isopropenyltoluene	132	C10H12	007399-49-7	47
2		Indan, 1-methyl-	132	C10H12	000767-58-8	38
3		Phenyl trans-2-phenyl-1-cyclopro...	274	C15H14O3S	017299-18-2	38
4		Benzene, (3-chloro-1-propenyl)-	152	C9H9Cl	002687-12-9	38
5		Benzene, 1-(chloromethyl)-2-ethe...	152	C9H9Cl	022570-84-9	23



Data Path : Z:\HPCHEM1\BNA G\DATA\BG110915\
Data File : BG019535.D
Acq On : 9 Nov 2015 16:31
Operator : UM/NP
Sample : G4245-17DL2 50X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BCY52DL2

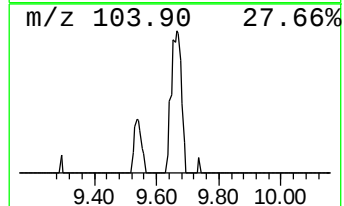
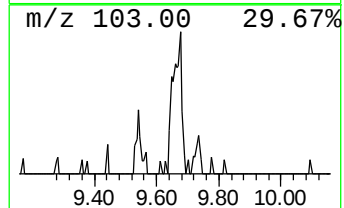
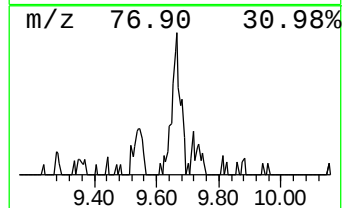
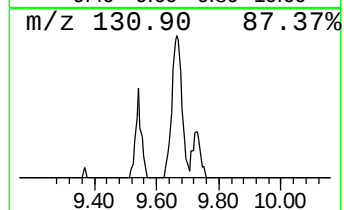
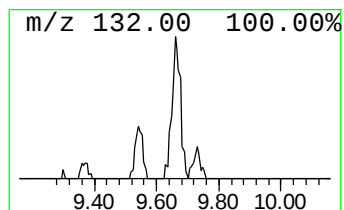
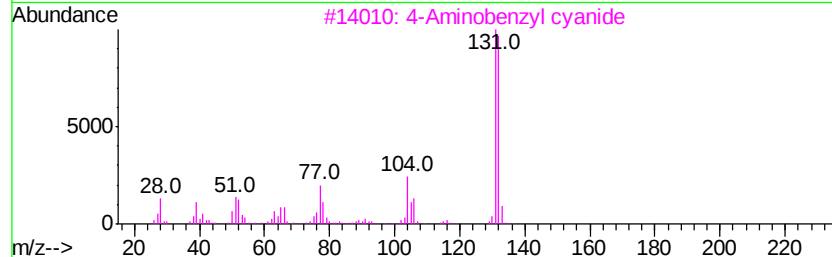
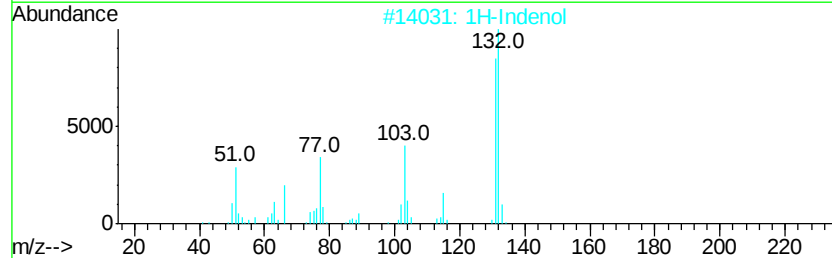
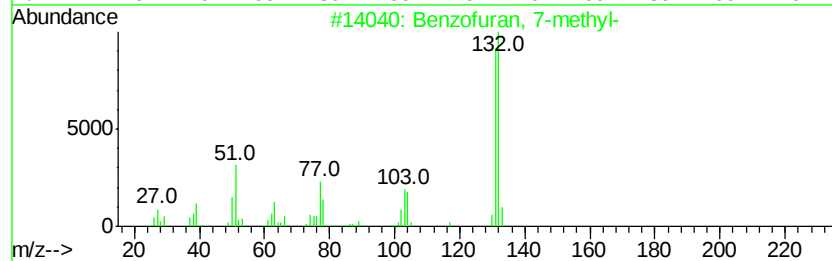
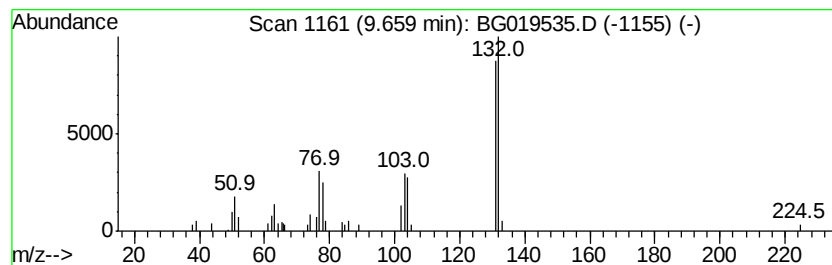
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Benzofuran, 7-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.66	3.25 ng/ul	48200	Naphthalene-d8	11.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzofuran, 7-methyl-	132	C9H8O	017059-52-8	94
2		1H-Indenol	132	C9H8O	056631-57-3	74
3		4-Aminobenzyl cyanide	132	C8H8N2	003544-25-0	59
4		3-Phenyl-2-propyn-1-ol	132	C9H8O	001504-58-1	59
5		Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	50



Data Path : Z:\HPCHEM1\BNA G\DATA\BG110915\
Data File : BG019535.D
Acq On : 9 Nov 2015 16:31
Operator : UM/NP
Sample : G4245-17DL2 50X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BCY52DL2

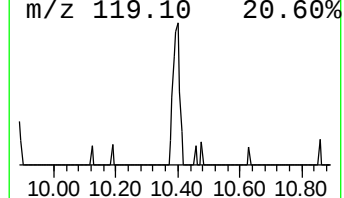
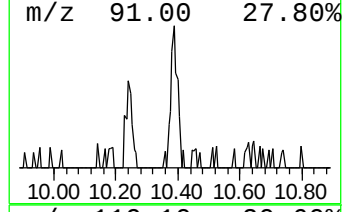
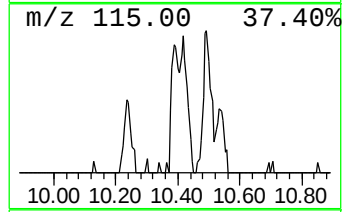
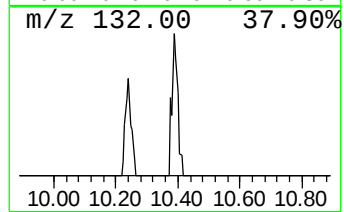
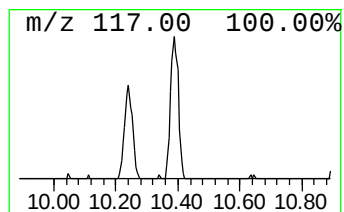
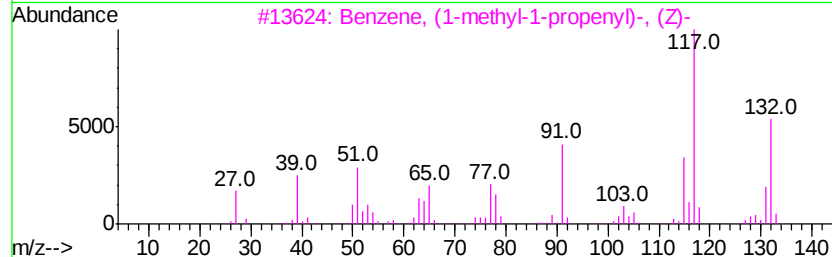
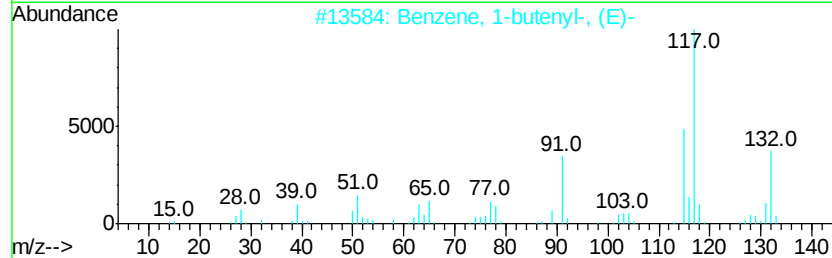
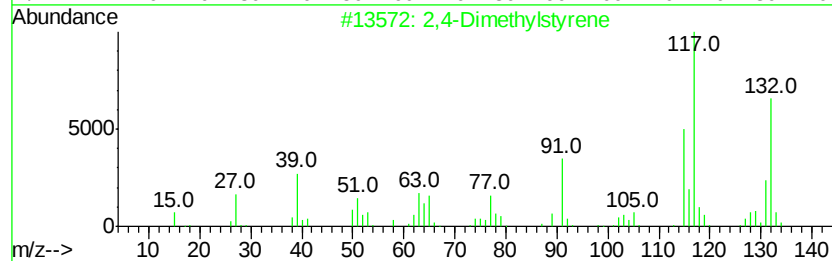
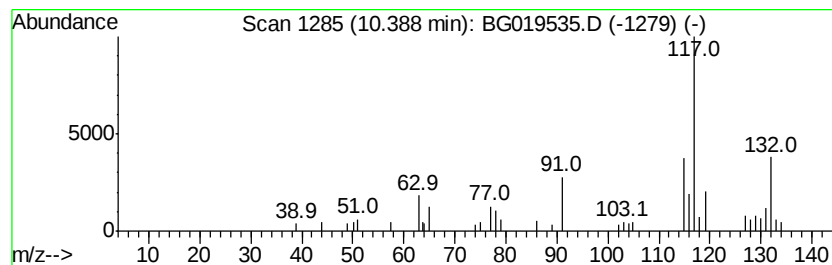
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 2,4-Dimethylstyrene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.39	3.67 ng/ul	54532	Naphthalene-d8	11.00

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,4-Dimethylstvrene	132	C10H12	002234-20-0	90
2			Benzene, 1-butenyl-, (E)-	132	C10H12	001005-64-7	81
3			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	76
4			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	76
5			Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	72



Data Path : Z:\HPCHEM1\BNA G\DATA\BG110915\
Data File : BG019535.D
Acq On : 9 Nov 2015 16:31
Operator : UM/NP
Sample : G4245-17DL2 50X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BCY52DL2

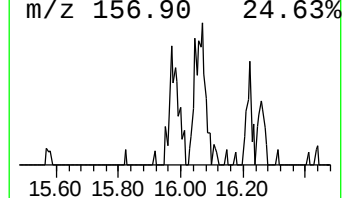
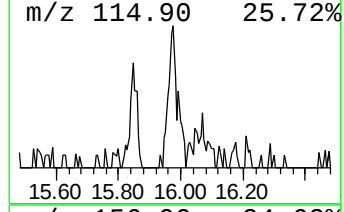
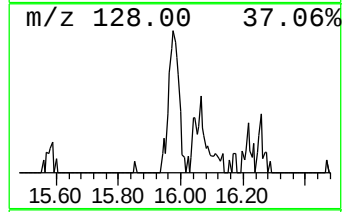
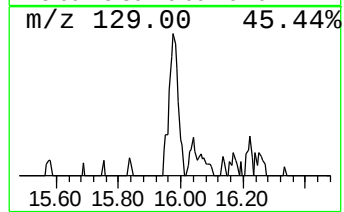
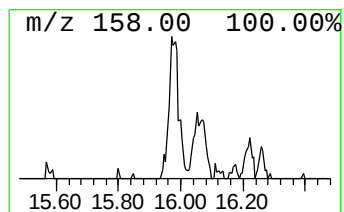
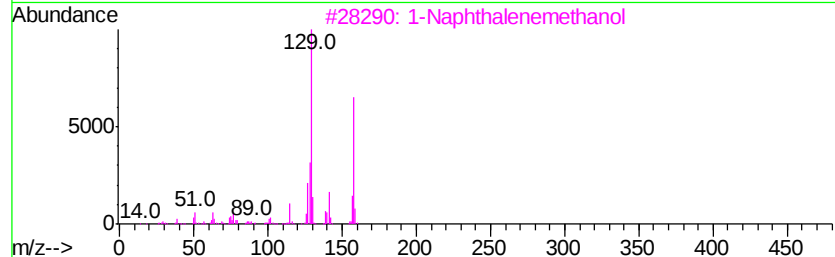
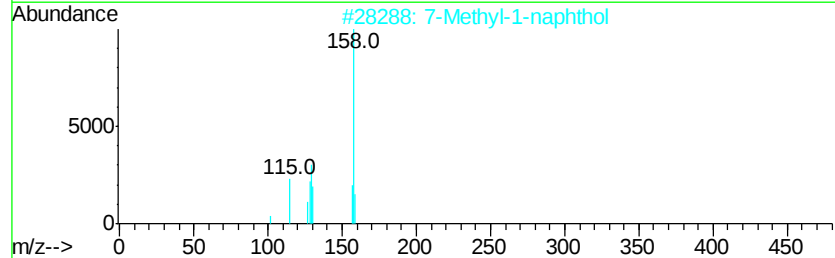
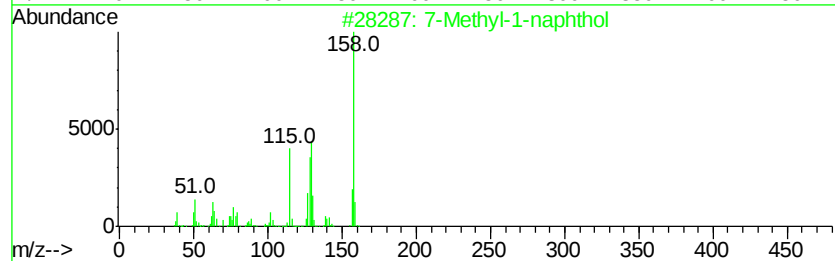
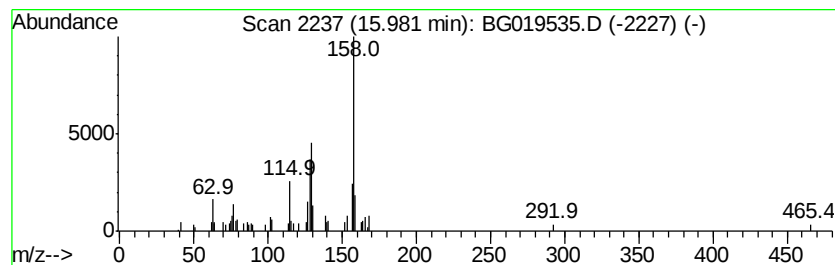
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 7-Methyl-1-naphthol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.98	2.64 ng/ul	67429	Acenaphthene-d10	14.81

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			7-Methyl-1-naphthol	158	C11H10O	006939-33-9	94
2			7-Methyl-1-naphthol	158	C11H10O	006939-33-9	90
3			1-Naphthalenemethanol	158	C11H10O	004780-79-4	72
4			1-Naphthalenol, 2-methyl-	158	C11H10O	007469-77-4	68
5			1-Naphthalenol, 4-methyl-	158	C11H10O	010240-08-1	62



Data Path : Z:\HPCHEM1\BNA G\DATA\BG110915\
Data File : BG019535.D
Acq On : 9 Nov 2015 16:31
Operator : UM/NP
Sample : G4245-17DL2 50X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampled :
BCY52DL2

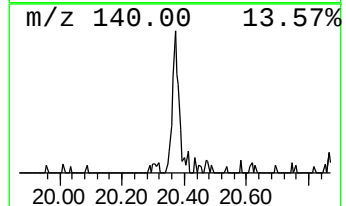
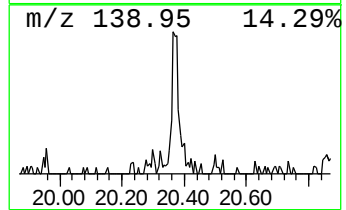
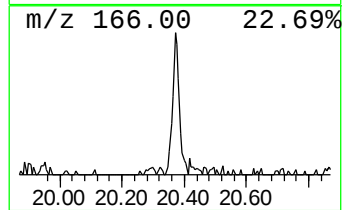
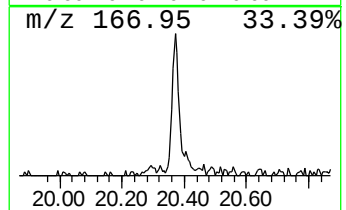
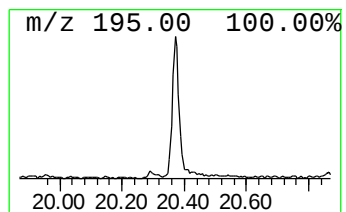
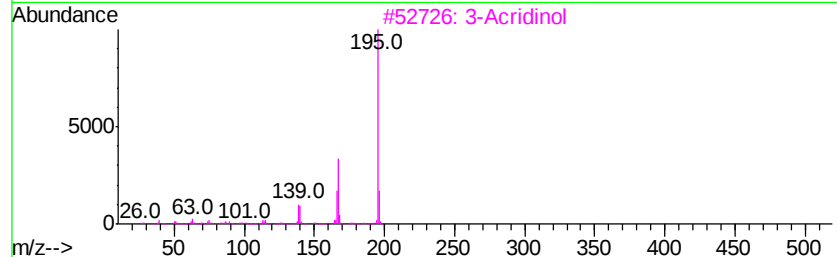
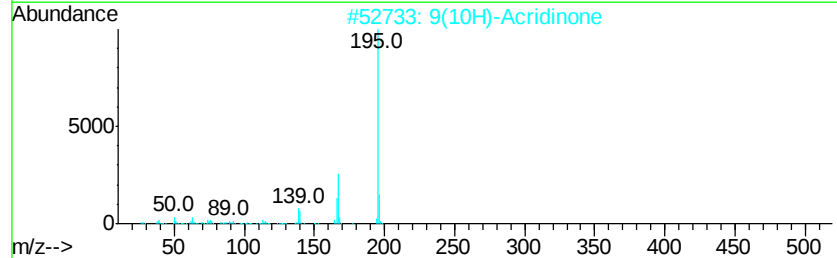
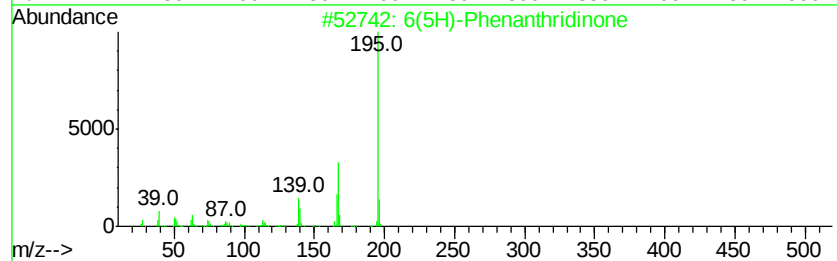
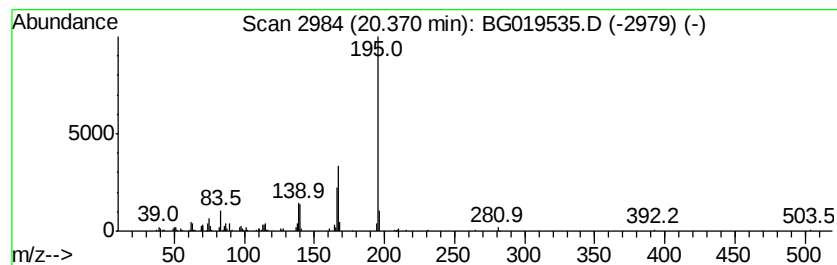
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 6(5H)-Phenanthridinone Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.37	2.26 ng/ul	115953	Chrysene-d12	21.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	6(5H)-Phenanthridinone	195	C13H9NO	001015-89-0	93
2		9(10H)-Acridinone	195	C13H9NO	000578-95-0	90
3		3-Acridinol	195	C13H9NO	007132-70-9	90
4		9(10H)-Acridinone	195	C13H9NO	000578-95-0	90
5		4-Cyano-4'-hydroxybiphenyl	195	C13H9NO	019812-93-2	83



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG110915\
Data File : BG019535.D
Acq On : 9 Nov 2015 16:31
Operator : UM/NP
Sample : G4245-17DL2 50X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BCY52DL2

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG102815.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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Indane	8.55	21.0	ng/ul	211104	1	8.17	201232	20.0
Benzene, 1-propynyl-	8.71	19.0	ng/ul	190981	1	8.17	201232	20.0
unknown-01	9.37	2.0	ng/ul	20312	1	8.17	201232	20.0
Benzofuran, 7-met...	9.66	3.3	ng/ul	48200	2	11.00	297046	20.0
2,4-Dimethylstyrene	10.39	3.7	ng/ul	54532	2	11.00	297046	20.0
7-Methyl-1-naphthol	15.98	2.6	ng/ul	67429	3	14.81	511356	20.0
6(5H)-Phenanthrid...	20.37	2.3	ng/ul	115953	5	21.83	1025160	20.0