

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM052116\
 Data File : BM005570.D
 Acq On : 22 May 2016 03:01
 Operator : UM/SJ
 Sample : H2890-16
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 D9WY6

Integration Parameters: LSCINT.P

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

Filtering: 5
 Min Area: 1 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM052016.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	2.652	7	11	28	rVB3	48344	123517	6.52%	0.413%
2	2.987	64	68	70	rBV	28301	31117	1.64%	0.104%
3	4.040	242	247	256	rBV	15217	24049	1.27%	0.080%
4	5.846	546	554	561	rVB2	20446	29345	1.55%	0.098%
5	6.804	709	717	726	rBV	39517	67171	3.55%	0.225%
6	6.916	731	736	741	rVV	76867	111881	5.91%	0.374%
7	6.987	741	748	755	rVV	283679	477216	25.19%	1.597%
8	7.045	755	758	766	rVB	23166	34793	1.84%	0.116%
9	7.157	770	777	783	rBV	198478	309175	16.32%	1.035%
10	7.216	783	787	791	rVB	16259	24107	1.27%	0.081%
11	7.357	804	811	819	rBV	273718	435750	23.00%	1.458%
12	7.469	819	830	836	rVB2	65239	127151	6.71%	0.426%
13	7.822	880	890	897	rBV	464659	759893	40.11%	2.543%
14	7.881	897	900	905	rVV	26949	39457	2.08%	0.132%
15	7.934	905	909	915	rVB	21139	31696	1.67%	0.106%
16	8.522	1003	1009	1017	rVB	354995	573518	30.27%	1.919%
17	8.975	1080	1086	1094	rBV	246514	420056	22.17%	1.406%
18	9.045	1094	1098	1101	rVV	24831	36080	1.90%	0.121%
19	9.081	1101	1104	1109	rVB2	18842	26483	1.40%	0.089%
20	9.698	1201	1209	1220	rBV	199325	348063	18.37%	1.165%
21	10.233	1289	1300	1308	rBV	338798	581503	30.69%	1.946%
22	10.610	1355	1364	1370	rBV	631502	1128648	59.57%	3.777%
23	10.663	1370	1373	1380	rVB	39408	59490	3.14%	0.199%
24	10.745	1380	1387	1397	rBV	319480	558680	29.49%	1.870%
25	11.522	1512	1519	1525	rBV2	15917	25990	1.37%	0.087%
26	11.639	1530	1539	1544	rVV3	14810	28835	1.52%	0.097%
27	11.716	1548	1552	1557	rVB2	16134	25670	1.35%	0.086%
28	11.880	1572	1580	1584	rVB4	16353	30835	1.63%	0.103%
29	11.939	1584	1590	1595	rBV	75656	117433	6.20%	0.393%
30	11.992	1595	1599	1603	rBV	18780	27845	1.47%	0.093%
31	12.198	1625	1634	1640	rVV2	60106	101745	5.37%	0.341%
32	12.275	1640	1647	1654	rVB	415135	672904	35.52%	2.252%
33	12.492	1675	1684	1694	rVB	317115	505827	26.70%	1.693%
34	13.092	1782	1786	1793	rVB	61236	91039	4.81%	0.305%

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Integration Parameters: LSCINT.P

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35	13.286	1813	1819	1823	rBV2	88252	140948	7.44%	0.472%
36	13.380	1832	1835	1845	rVB	18874	31578	1.67%	0.106%
37	13.480	1845	1852	1864	rVB	63726	131634	6.95%	0.441%
38	13.616	1868	1875	1877	rBV	125169	228412	12.06%	0.764%
39	13.774	1895	1902	1907	rBV	234001	392104	20.70%	1.312%
40	13.827	1907	1911	1913	rVV	134199	195085	10.30%	0.653%
41	13.863	1913	1917	1921	rVV	660274	934745	49.34%	3.128%
42	13.910	1921	1925	1933	rVV	250275	365537	19.29%	1.223%
43	14.010	1937	1942	1946	rVV	114970	174405	9.21%	0.584%
44	14.045	1946	1948	1952	rVV3	32323	48240	2.55%	0.161%
45	14.098	1952	1957	1960	rVV2	57282	91677	4.84%	0.307%
46	14.145	1960	1965	1980	rVB	715951	1202499	63.47%	4.024%
47	14.351	1996	2000	2006	rBV	57553	79825	4.21%	0.267%
48	14.457	2006	2018	2024	rVV2	922114	1551272	81.88%	5.192%
49	14.510	2024	2027	2029	rVV2	24381	38460	2.03%	0.129%
50	14.539	2029	2032	2036	rVB2	33515	46363	2.45%	0.155%
51	14.657	2047	2052	2062	rBV3	54753	136569	7.21%	0.457%
52	14.745	2063	2067	2072	rVV3	24507	41154	2.17%	0.138%
53	14.804	2072	2077	2080	rVV4	27187	47223	2.49%	0.158%
54	14.851	2080	2085	2092	rVV	91757	159761	8.43%	0.535%
55	14.927	2092	2098	2102	rVV	67560	115833	6.11%	0.388%
56	14.968	2102	2105	2112	rVB6	21516	36566	1.93%	0.122%
57	15.068	2115	2122	2128	rBV3	46800	106113	5.60%	0.355%
58	15.121	2128	2131	2138	rVB2	54512	79145	4.18%	0.265%
59	15.245	2144	2152	2154	rBV3	50696	107335	5.67%	0.359%
60	15.304	2159	2162	2167	rVB	76079	91758	4.84%	0.307%
61	15.374	2167	2174	2176	rBV6	19947	35360	1.87%	0.118%
62	15.445	2177	2186	2191	rVV	932998	1389479	73.34%	4.650%
63	15.498	2191	2195	2198	rVB3	35040	47952	2.53%	0.160%
64	15.621	2208	2216	2220	rBV5	41682	77746	4.10%	0.260%
65	15.710	2228	2231	2236	rVB3	37745	49868	2.63%	0.167%
66	15.792	2236	2245	2252	rVB8	41069	112000	5.91%	0.375%
67	15.863	2253	2257	2262	rBV8	15087	23969	1.27%	0.080%
68	15.945	2263	2271	2275	rBV5	18895	48351	2.55%	0.162%
69	15.992	2275	2279	2293	rVB5	37522	84300	4.45%	0.282%
70	16.110	2294	2299	2301	rBV3	22190	34050	1.80%	0.114%
71	16.168	2304	2309	2311	rBV3	77469	116839	6.17%	0.391%

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Integration Parameters: LSCINT.P

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72	16.280	2323	2328	2330	rBV2	24114	36554	1.93%	0.122%
73	16.339	2334	2338	2343	rVB5	40563	69948	3.69%	0.234%
74	16.398	2344	2348	2352	rVB4	28082	40059	2.11%	0.134%
75	16.445	2352	2356	2359	rBV3	16387	25555	1.35%	0.086%
76	16.498	2362	2365	2371	rVV5	14401	26131	1.38%	0.087%
77	16.592	2377	2381	2389	rVB4	88692	153239	8.09%	0.513%
78	16.662	2389	2393	2399	rBV3	29500	52448	2.77%	0.176%
79	16.739	2402	2406	2410	rVB3	26867	35846	1.89%	0.120%
80	16.951	2439	2442	2447	rVB	75780	91835	4.85%	0.307%
81	17.009	2448	2452	2459	rVB3	35478	60551	3.20%	0.203%
82	17.198	2477	2484	2488	rBV2	1089929	1738398	91.76%	5.818%
83	17.239	2488	2491	2495	rVB	291148	362135	19.11%	1.212%
84	17.292	2495	2500	2505	rBV2	966688	1397437	73.76%	4.677%
85	17.574	2545	2548	2550	rBV3	30045	35225	1.86%	0.118%
86	17.692	2565	2568	2574	rVB3	48516	64017	3.38%	0.214%
87	18.086	2629	2635	2637	rBV2	82611	135973	7.18%	0.455%
88	18.162	2643	2648	2652	rBV	759107	967535	51.07%	3.238%
89	18.262	2660	2665	2669	rBV3	63115	129098	6.81%	0.432%
90	18.568	2714	2717	2720	rBV	37003	47488	2.51%	0.159%
91	18.698	2735	2739	2745	rBV	157524	236174	12.47%	0.790%
92	18.798	2752	2756	2761	rBV	86923	143995	7.60%	0.482%
93	18.945	2778	2781	2784	rVB	81209	85893	4.53%	0.287%
94	19.245	2828	2832	2837	rBV	585329	828338	43.72%	2.772%
95	19.356	2848	2851	2860	rBV4	68135	144651	7.64%	0.484%
96	19.580	2884	2889	2893	rBV	1136256	1884284	99.46%	6.306%
97	21.080	3140	3144	3151	rVB3	294200	401839	21.21%	1.345%
98	21.368	3187	3193	3197	rBV2	1256080	1894521	100.00%	6.341%
99	23.503	3552	3556	3561	rBV2	531144	906841	47.87%	3.035%
100	23.650	3576	3581	3587	rVB	692089	1302407	68.75%	4.359%

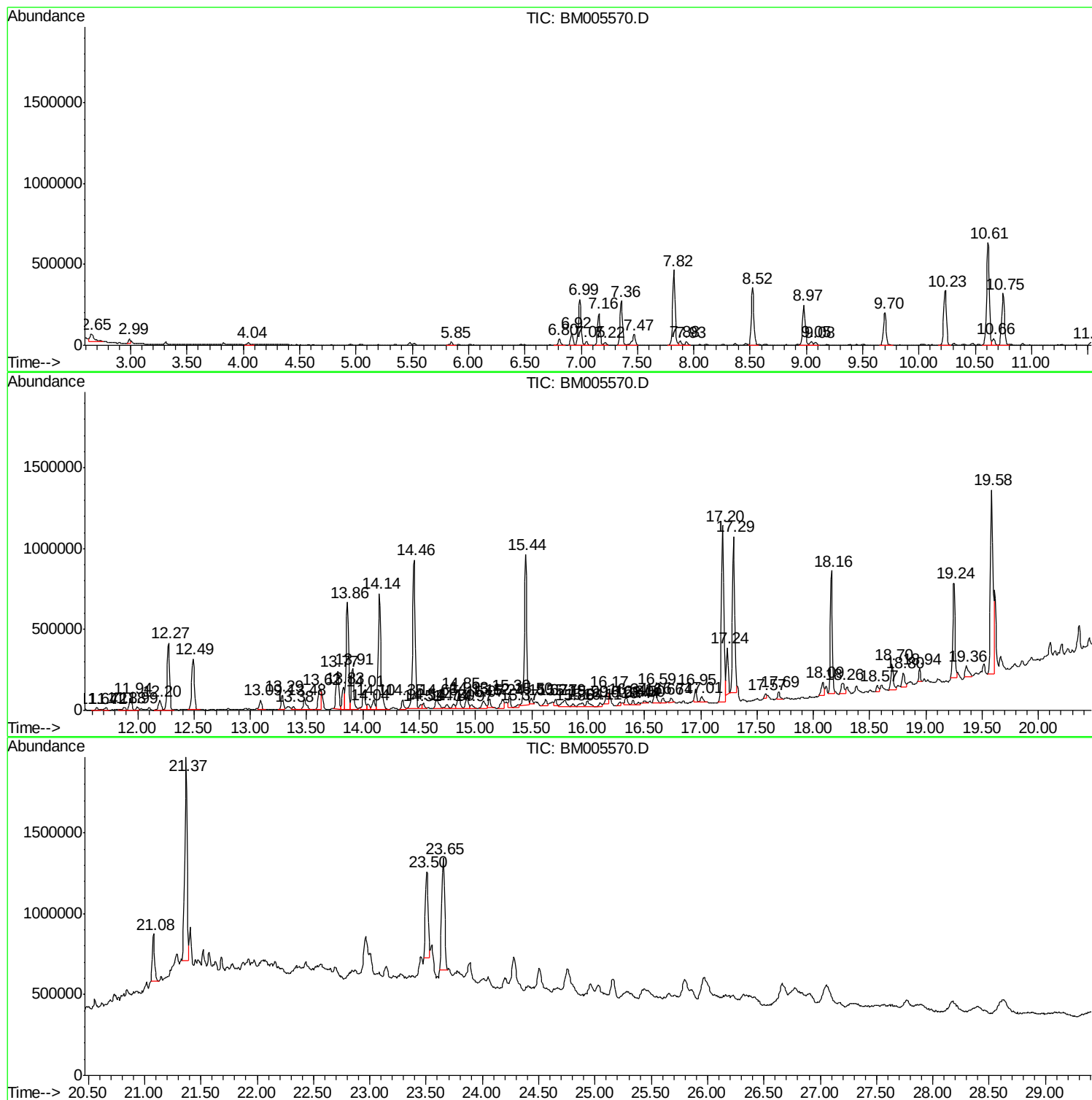
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM052016.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
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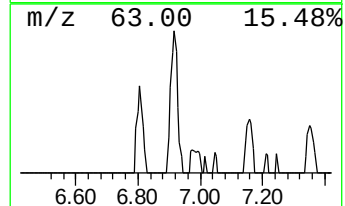
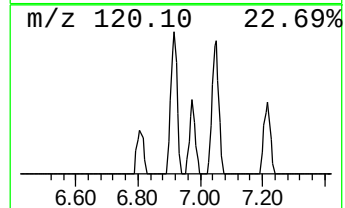
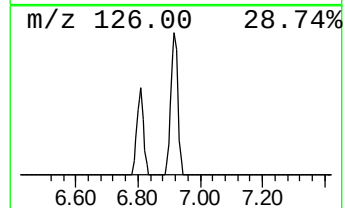
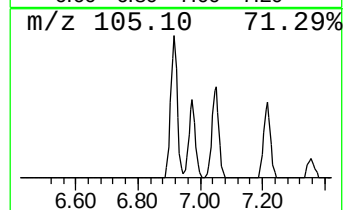
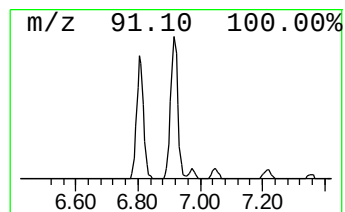
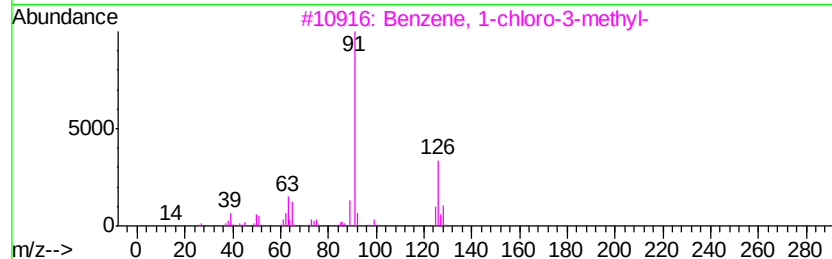
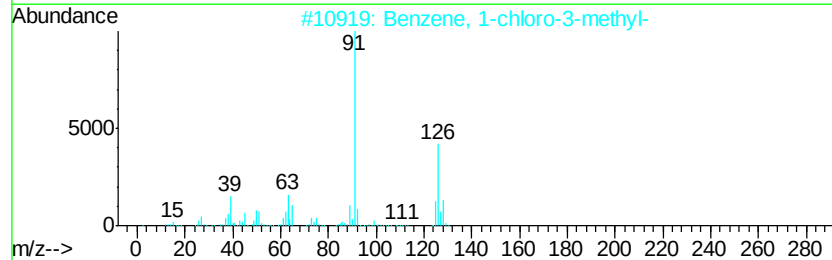
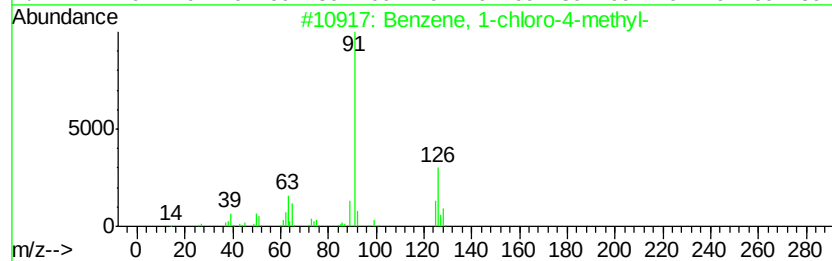
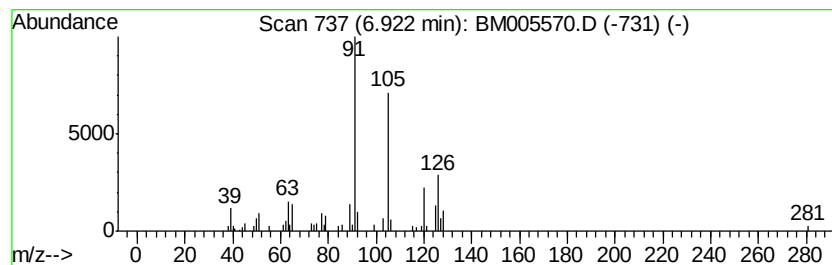
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM052016.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzene, 1-chloro-4-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.92	2.94 ng/ul	111881	1,4-Dichlorobenzene-d4	7.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-chloro-4-methyl-	126	C7H7Cl	000106-43-4	93
2		Benzene, 1-chloro-3-methyl-	126	C7H7Cl	000108-41-8	93
3		Benzene, 1-chloro-3-methyl-	126	C7H7Cl	000108-41-8	93
4		Benzene, 1-chloro-4-methyl-	126	C7H7Cl	000106-43-4	91
5		Benzene, 1-chloro-3-methyl-	126	C7H7Cl	000108-41-8	91



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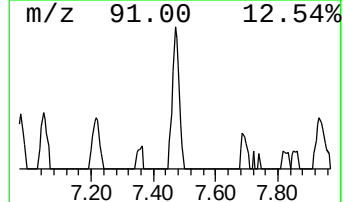
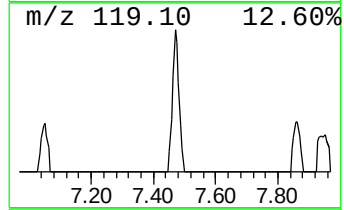
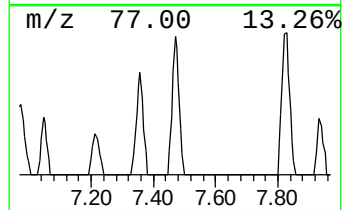
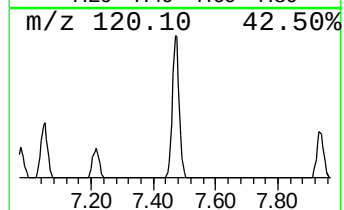
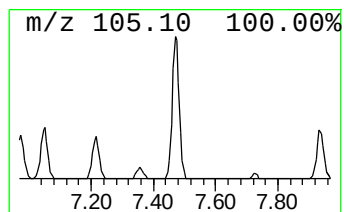
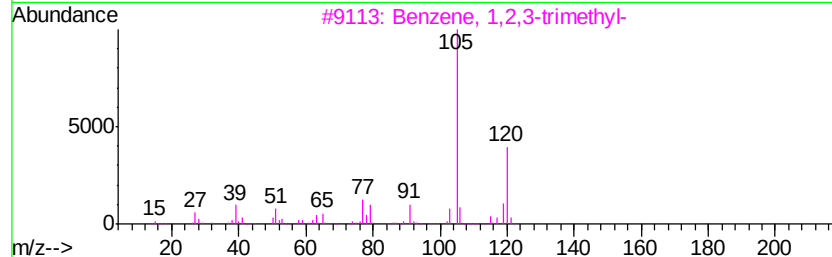
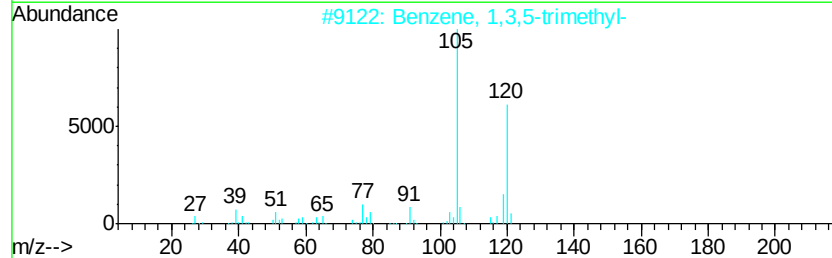
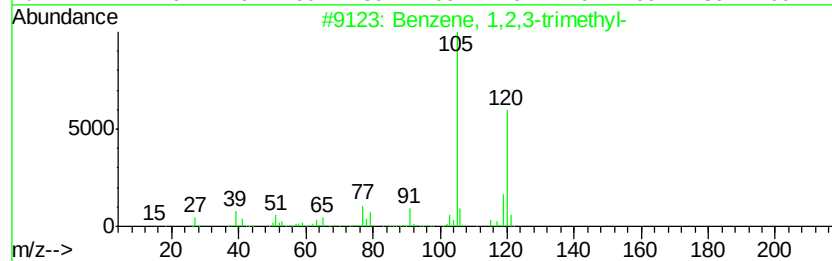
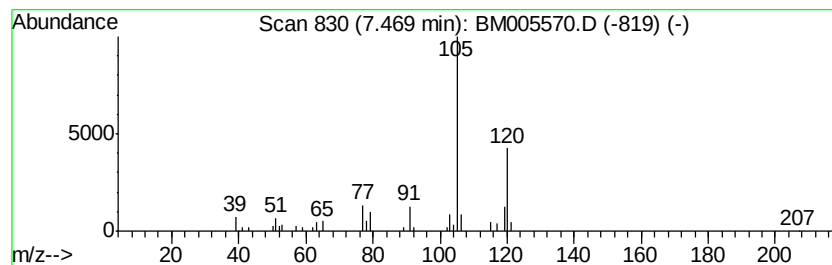
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM052016.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.47	3.35 ng/ul	127151	1,4-Dichlorobenzene-d4	7.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2		Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	95
3		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
4		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
5		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94



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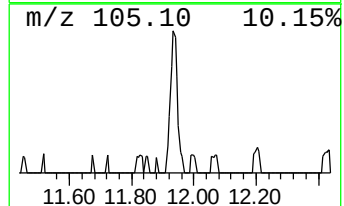
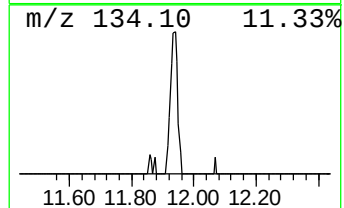
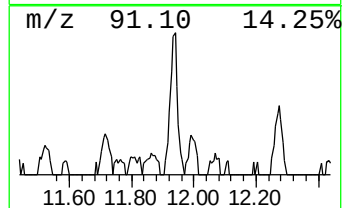
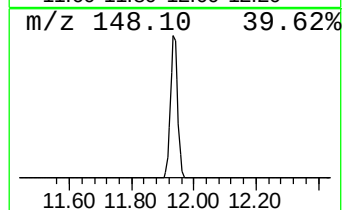
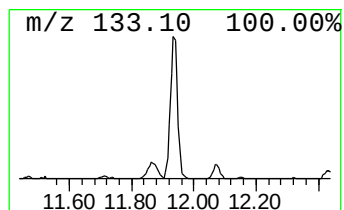
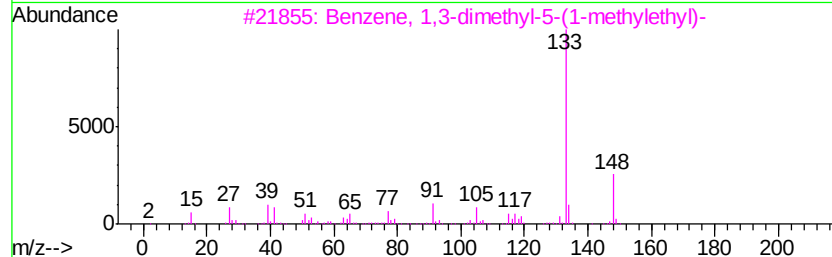
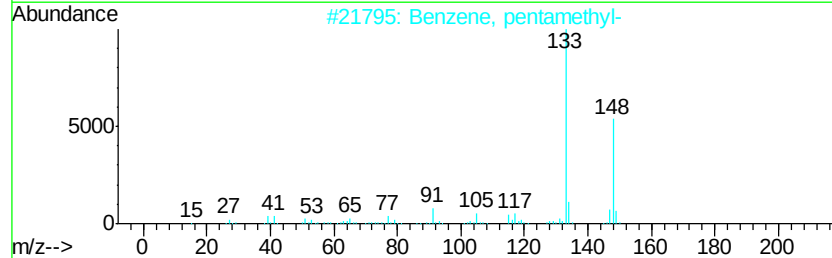
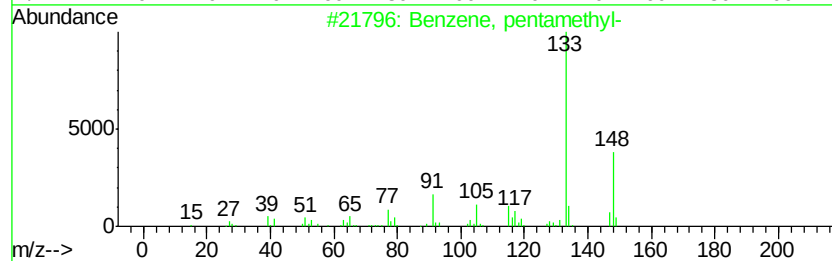
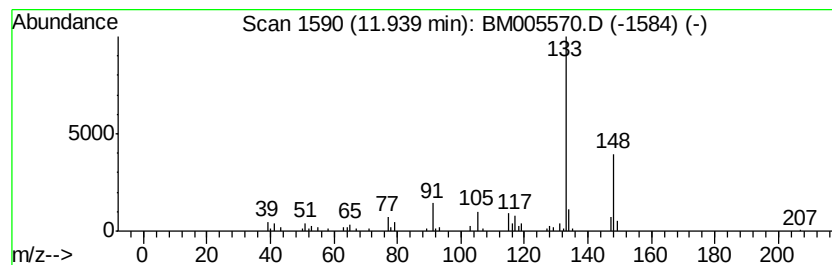
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzene, pentamethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.94	2.08 ng/ul	117433	Napthalene-d8	10.61

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, pentamethyl-	148	C11H16	000700-12-9	97
2		Benzene, pentamethyl-	148	C11H16	000700-12-9	94
3		Benzene, 1,3-dimethyl-5-(1-methy...	148	C11H16	004706-90-5	91
4		Benzene, pentamethyl-	148	C11H16	000700-12-9	91
5		Benzene, 1-ethyl-2,4,5-trimethyl-	148	C11H16	017851-27-3	90



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM052116\
Data File : BM005570.D
Acq On : 22 May 2016 03:01
Operator : UM/SJ
Sample : H2890-16
Misc :
ALS Vial : 20 Sample Multiplier: 1

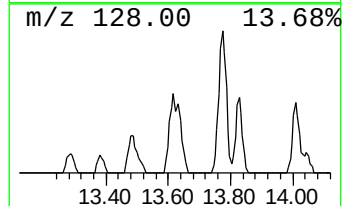
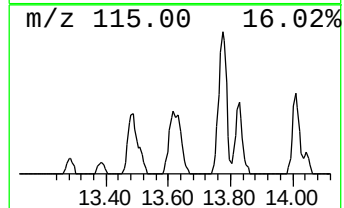
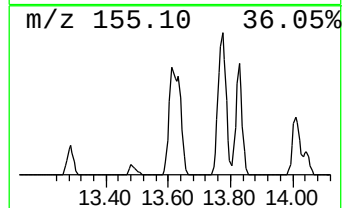
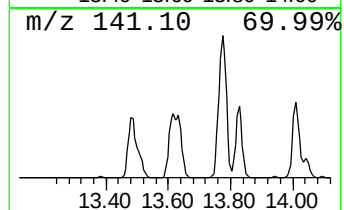
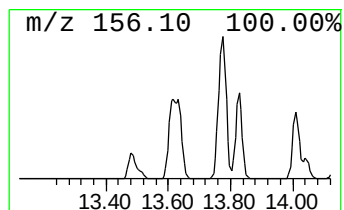
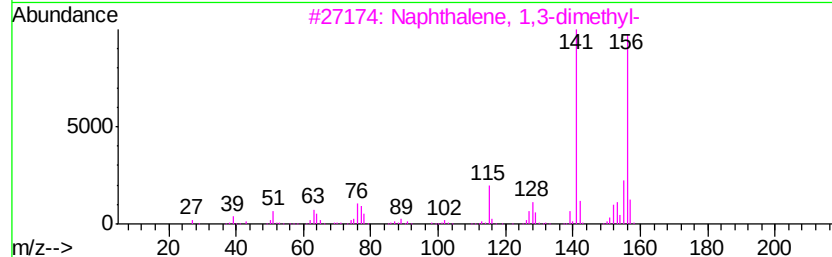
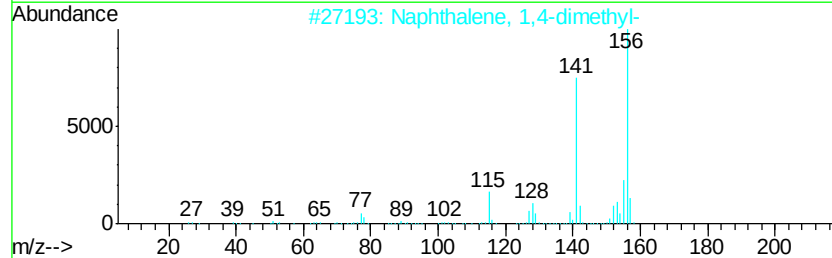
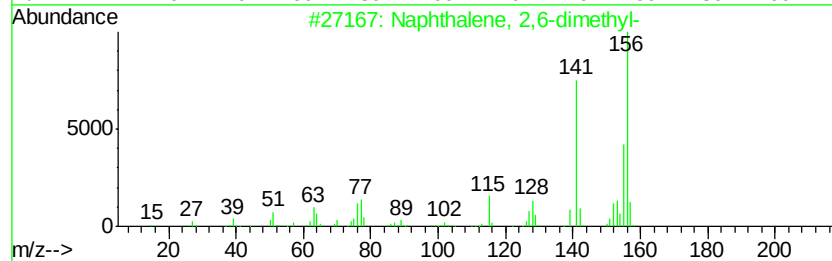
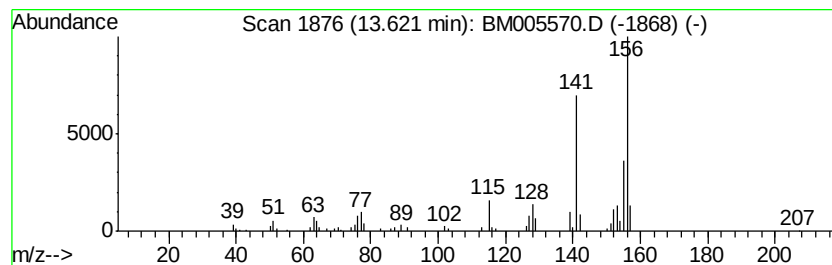
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM052016.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 Naphthalene, 2,6-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.62	2.94 ng/ul	228412	Acenaphthene-d10	14.46
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 2,6-dimethyl-	156 C12H12	000581-42-0 97
2		Naphthalene, 1,4-dimethyl-	156 C12H12	000571-58-4 97
3		Naphthalene, 1,3-dimethyl-	156 C12H12	000575-41-7 96
4		Naphthalene, 1,6-dimethyl-	156 C12H12	000575-43-9 96
5		Naphthalene, 2,7-dimethyl-	156 C12H12	000582-16-1 96



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM052116\
Data File : BM005570.D
Acq On : 22 May 2016 03:01
Operator : UM/SJ
Sample : H2890-16
Misc :
ALS Vial : 20 Sample Multiplier: 1

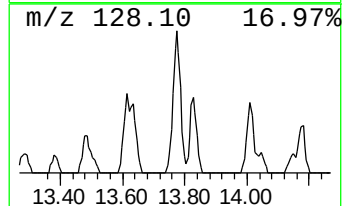
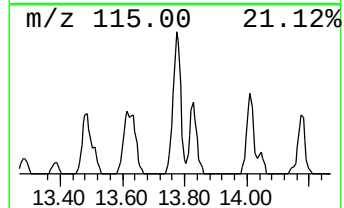
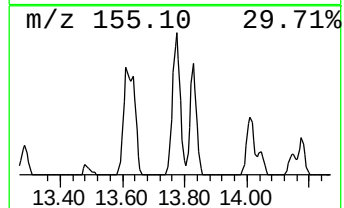
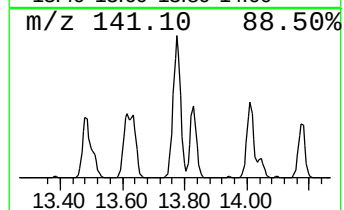
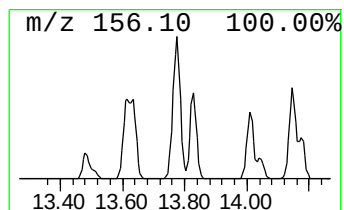
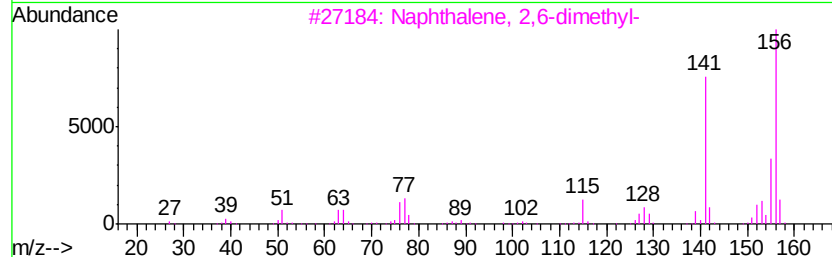
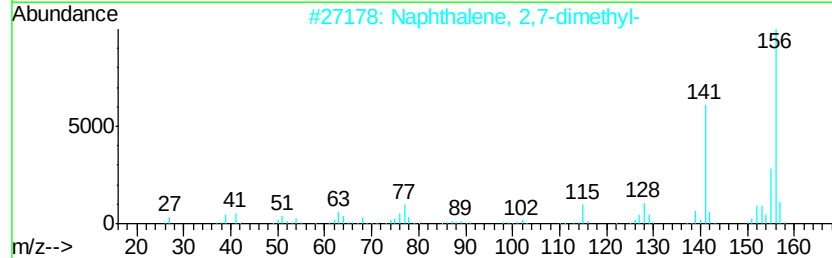
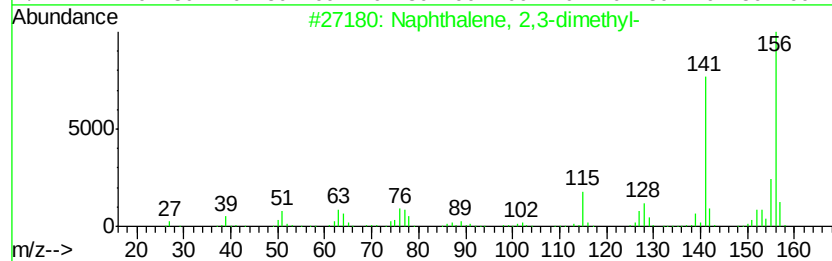
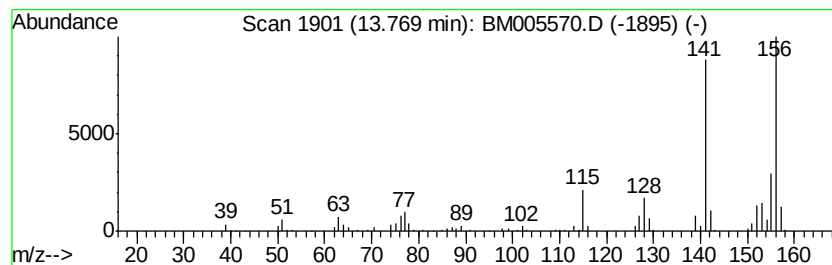
Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM052016.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 Naphthalene, 2,3-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.77	5.06 ng/ul	392104	Acenaphthene-d10	14.46
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 2,3-dimethyl-	156 C12H12	000581-40-8 97
2		Naphthalene, 2,7-dimethyl-	156 C12H12	000582-16-1 97
3		Naphthalene, 2,6-dimethyl-	156 C12H12	000581-42-0 97
4		Naphthalene, 1,6-dimethyl-	156 C12H12	000575-43-9 97
5		Naphthalene, 1,7-dimethyl-	156 C12H12	000575-37-1 97



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM052116\
Data File : BM005570.D
Acq On : 22 May 2016 03:01
Operator : UM/SJ
Sample : H2890-16
Misc :
ALS Vial : 20 Sample Multiplier: 1

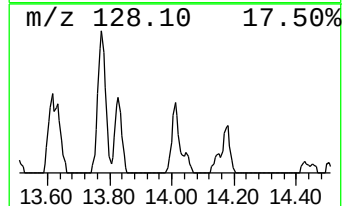
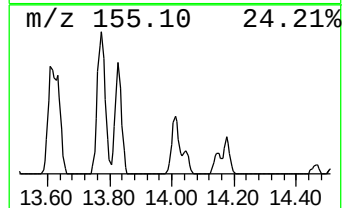
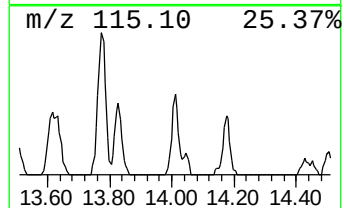
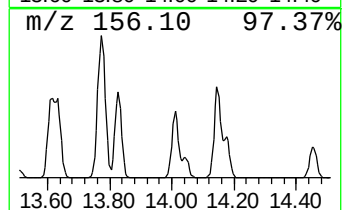
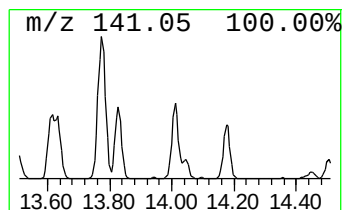
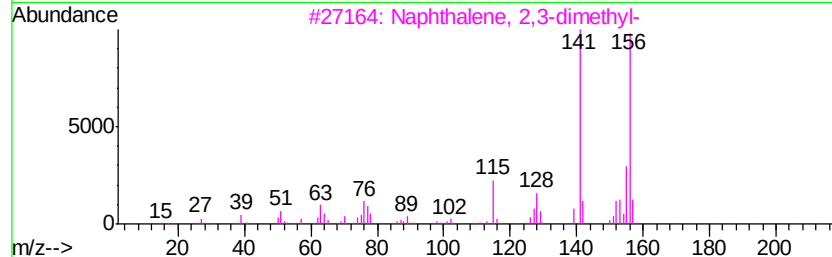
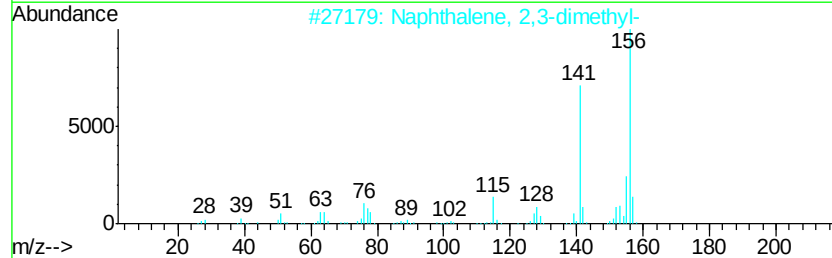
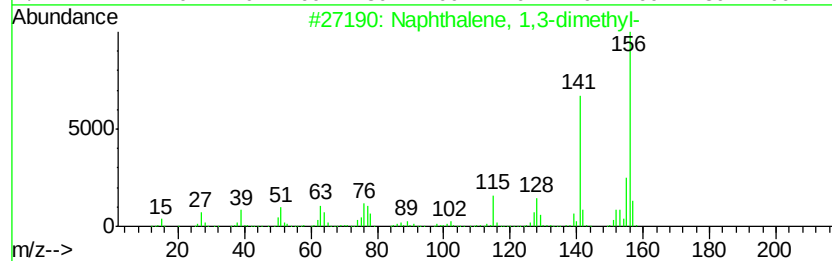
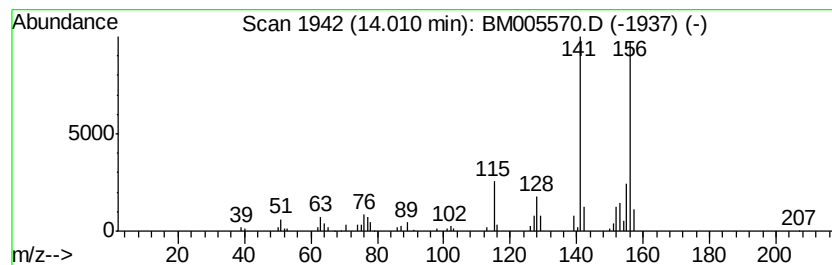
Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM052016.M
Quant Title : SVOA CALIBRATION

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

Peak Number 8 Naphthalene, 1,3-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.01	2.25 ng/ul	174405	Acenaphthene-d10	14.46
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,3-dimethyl-	156 C12H12	000575-41-7 97
2		Naphthalene, 2,3-dimethyl-	156 C12H12	000581-40-8 97
3		Naphthalene, 2,3-dimethyl-	156 C12H12	000581-40-8 97
4		Naphthalene, 2,3-dimethyl-	156 C12H12	000581-40-8 97
5		Naphthalene, 2,3-dimethyl-	156 C12H12	000581-40-8 97



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM052116\
Data File : BM005570.D
Acq On : 22 May 2016 03:01
Operator : UM/SJ
Sample : H2890-16
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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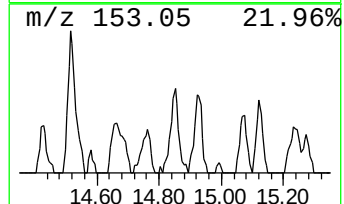
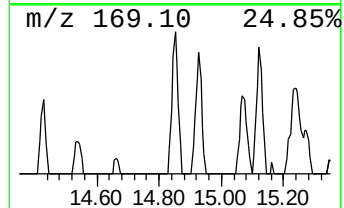
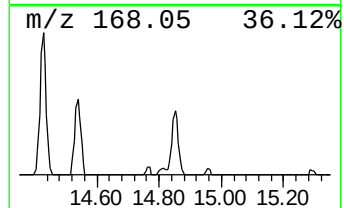
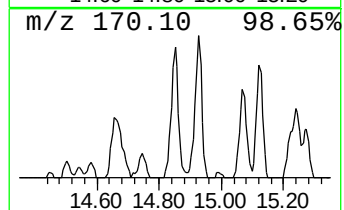
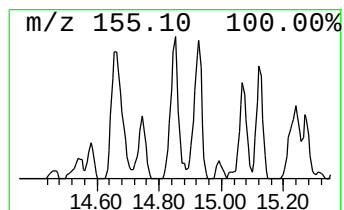
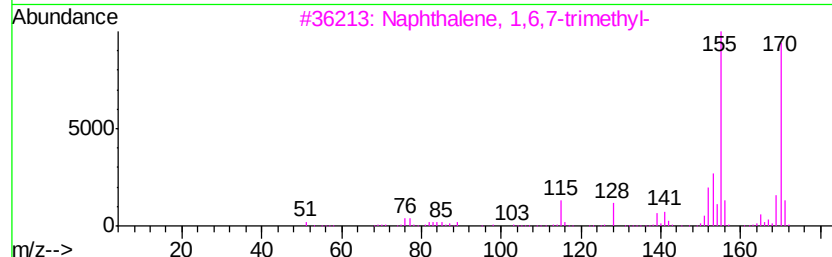
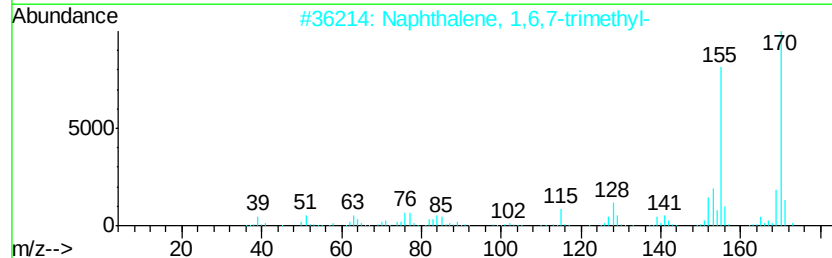
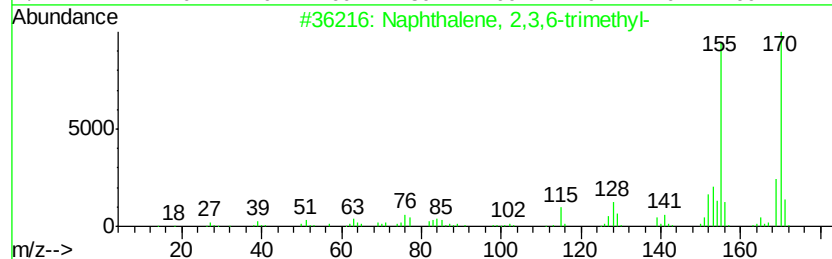
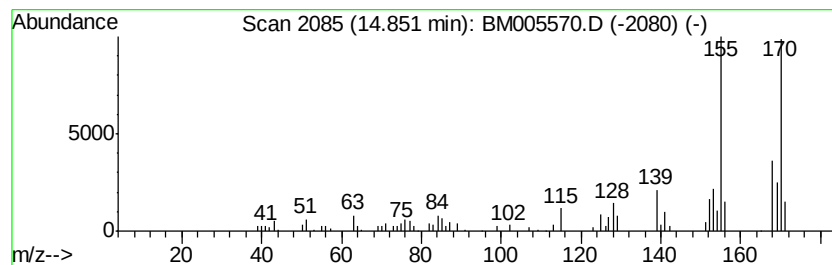
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Naphthalene, 2,3,6-trimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.85	2.06 ng/ul	159761	Acenaphthene-d10	14.46

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	96
2			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	96
3			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	95
4			Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	95
5			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	95



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM052116\
Data File : BM005570.D
Acq On : 22 May 2016 03:01
Operator : UM/SJ
Sample : H2890-16
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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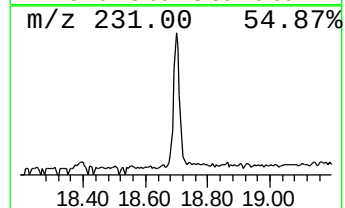
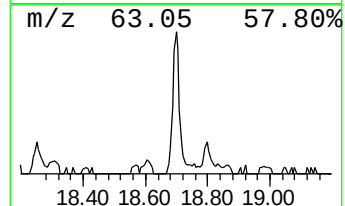
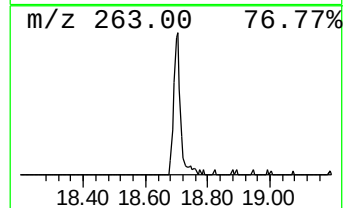
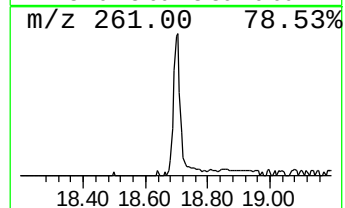
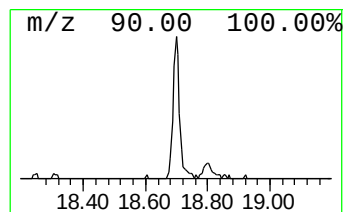
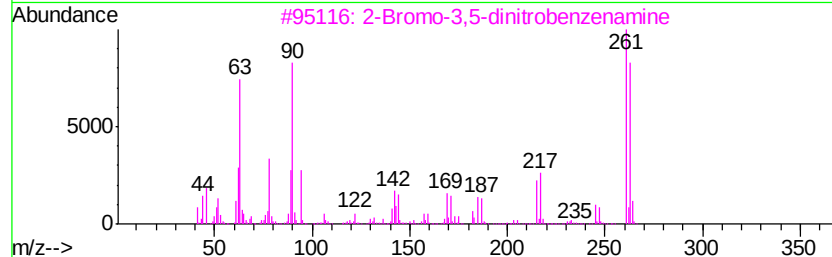
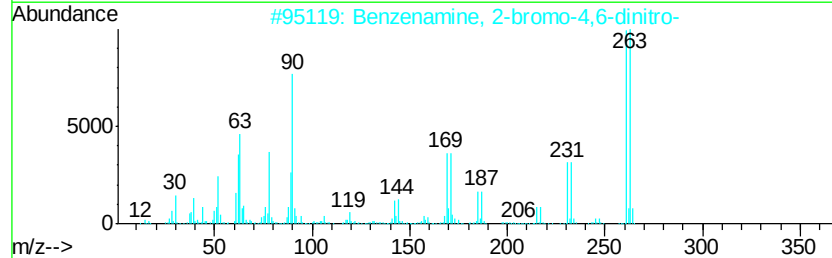
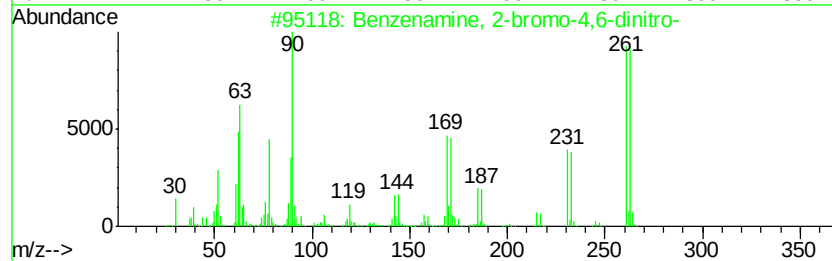
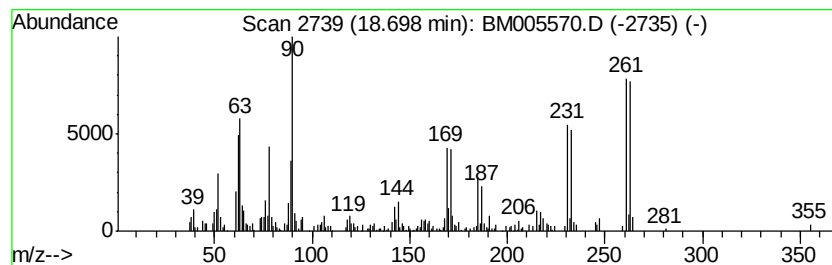
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Benzenamine, 2-bromo-4,6-di... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.70	2.72 ng/ul	236174	Phenanthrene-d10	17.20

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenamine, 2-bromo-4,6-dinitro-	261	C6H4BrN3O4	001817-73-8	99
2			Benzenamine, 2-bromo-4,6-dinitro-	261	C6H4BrN3O4	001817-73-8	99
3			2-Bromo-3,5-dinitrobenzenamine	261	C6H4BrN3O4	116529-41-0	70
4			1,3,5-Norcaratriene	90	C7H6	004646-69-9	50
5			Benzenamine, 2-bromo-4,6-dinitro-	261	C6H4BrN3O4	001817-73-8	42



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM052116\
Data File : BM005570.D
Acq On : 22 May 2016 03:01
Operator : UM/SJ
Sample : H2890-16
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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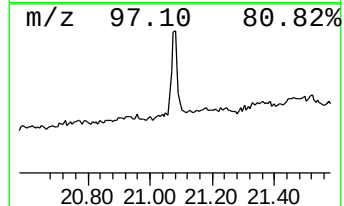
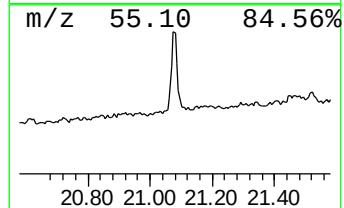
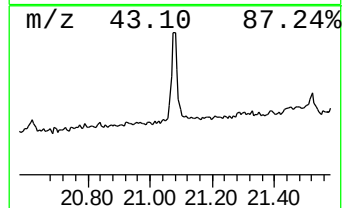
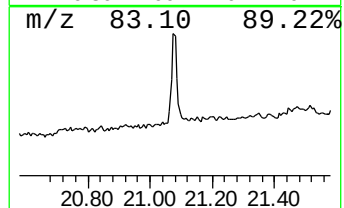
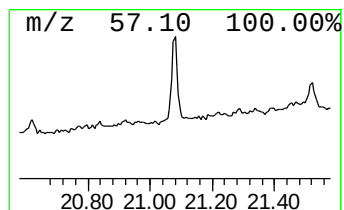
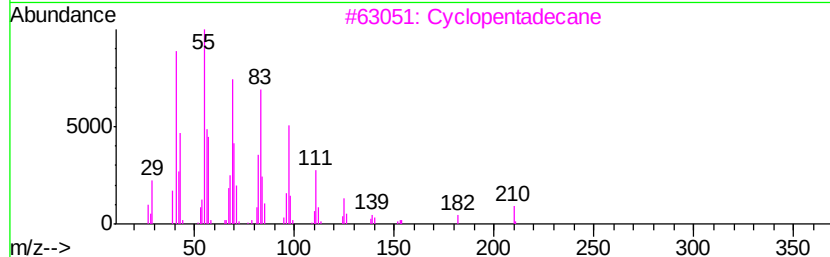
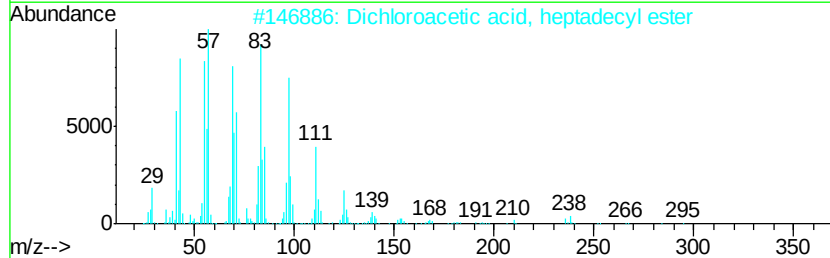
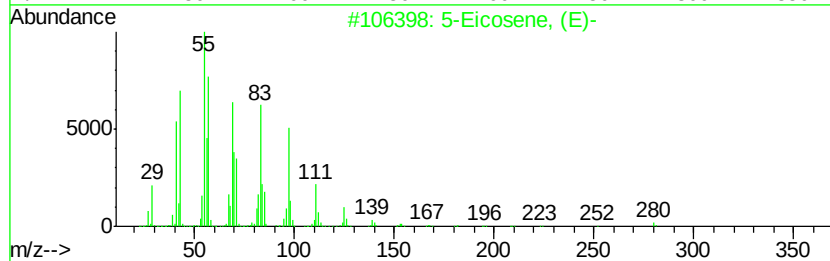
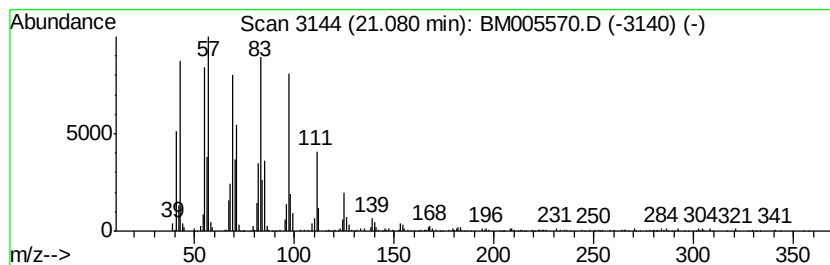
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Cyclic21.08 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.08	4.24 ng/ul	401839	Chrysene-d12	21.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Eicosene, (E)-	280	C20H40	074685-30-6	98
2		Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	94
3		Cyclopentadecane	210	C15H30	000295-48-7	93
4		17-Pentatriacontene	491	C35H70	006971-40-0	91
5		Trifluoroacetic acid, n-octadecy...	366	C20H37F3O2	079392-43-1	91



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM052116\
Data File : BM005570.D
Acq On : 22 May 2016 03:01
Operator : UM/SJ
Sample : H2890-16
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM052016.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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Benzene, 1,2,3-tr...	7.47	3.4	ng/ul	127151	1	7.82	759893	20.0
Benzene, pentamet...	11.94	2.1	ng/ul	117433	2	10.61	1128650	20.0
Naphthalene, 2,6-...	13.62	2.9	ng/ul	228412	3	14.46	1551270	20.0
Naphthalene, 2,3-...	13.77	5.1	ng/ul	392104	3	14.46	1551270	20.0
Naphthalene, 1,3-...	14.01	2.3	ng/ul	174405	3	14.46	1551270	20.0
Naphthalene, 2,3,...	14.85	2.1	ng/ul	159761	3	14.46	1551270	20.0
Benzenamine, 2-br...	18.70	2.7	ng/ul	236174	4	17.20	1738400	20.0
(DEL) Alkane: Cyc...	21.08	4.2	ng/ul	401839	5	21.37	1894520	20.0