

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG012615\
 Data File : BG015931.D
 Acq On : 27 Jan 2015 2:40
 Operator : TP/IZ
 Sample : G1139-09
 Misc : S
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 121B2B

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 3 % 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\8270-BG012015.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.826	17	23	31	rBV2	481784	814180	7.76%	1.020%
2	2.902	31	36	48	rVB	858277	1064574	10.15%	1.334%
3	4.812	356	361	367	rVB2	105941	154603	1.47%	0.194%
4	5.212	423	429	433	rVB	124697	170718	1.63%	0.214%
5	5.282	433	441	446	rBV	769148	1095092	10.44%	1.372%
6	5.341	446	451	459	rVB	259538	541283	5.16%	0.678%
7	5.705	508	513	519	rBV	256646	372773	3.55%	0.467%
8	6.780	688	696	701	rBV	303946	447500	4.27%	0.561%
9	6.839	701	706	707	rVV2	134569	176658	1.68%	0.221%
10	6.868	707	711	715	rVV2	494911	768733	7.33%	0.963%
11	6.910	715	718	724	rVB2	208013	291290	2.78%	0.365%
12	7.074	740	746	753	rVB	224485	365950	3.49%	0.458%
13	7.221	765	771	780	rVB	1485243	2202482	21.00%	2.759%
14	7.333	783	790	798	rVB	738980	1086783	10.36%	1.361%
15	7.685	842	850	854	rBV2	455890	744922	7.10%	0.933%
16	7.797	863	869	877	rVB	396815	624079	5.95%	0.782%
17	8.002	897	904	910	rBV	1656787	2702131	25.77%	3.385%
18	8.055	910	913	923	rVB	225502	310746	2.96%	0.389%
19	8.231	939	943	948	rVB2	120271	178701	1.70%	0.224%
20	8.320	952	958	963	rBV2	223785	342150	3.26%	0.429%
21	8.484	980	986	993	rBV2	94575	182762	1.74%	0.229%
22	8.631	1005	1011	1016	rBV	144784	221127	2.11%	0.277%
23	8.684	1016	1020	1027	rVB2	143467	209556	2.00%	0.263%
24	8.784	1029	1037	1041	rBV	248666	387449	3.69%	0.485%
25	8.843	1041	1047	1057	rVB	1571109	2734950	26.08%	3.426%
26	9.119	1087	1094	1099	rBV3	103126	197933	1.89%	0.248%
27	9.307	1121	1126	1130	rBV	122558	168128	1.60%	0.211%
28	9.365	1130	1136	1143	rVB2	221235	335367	3.20%	0.420%
29	9.677	1182	1189	1192	rBV8	63240	141431	1.35%	0.177%
30	9.724	1192	1197	1206	rVB2	209439	344373	3.28%	0.431%
31	9.871	1216	1222	1230	rVB2	558968	994395	9.48%	1.246%
32	10.106	1257	1262	1268	rVB2	321675	501116	4.78%	0.628%
33	10.470	1313	1324	1328	rBV2	626685	1231914	11.75%	1.543%
34	10.517	1328	1332	1340	rVV	826656	1436741	13.70%	1.800%

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Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

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Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG012015.M
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35	10.588	1340	1344	1348	rVB3	106904	167862	1.60%	0.210%
36	10.934	1397	1403	1412	rVB3	126724	221664	2.11%	0.278%
37	11.046	1412	1422	1427	rBV3	98293	190853	1.82%	0.239%
38	11.381	1474	1479	1487	rVB2	144803	262046	2.50%	0.328%
39	11.581	1505	1513	1518	rVB4	139439	301459	2.87%	0.378%
40	11.663	1518	1527	1536	rBV2	295036	564190	5.38%	0.707%
41	11.827	1546	1555	1558	rBV5	121241	272749	2.60%	0.342%
42	12.027	1582	1589	1594	rVB3	229319	418803	3.99%	0.525%
43	12.139	1598	1608	1615	rVV	1579182	2519130	24.02%	3.156%
44	12.362	1637	1646	1657	rBV	1193638	2126659	20.28%	2.664%
45	12.579	1677	1683	1688	rBV4	131471	195318	1.86%	0.245%
46	12.873	1727	1733	1738	rBV8	70221	135894	1.30%	0.170%
47	12.949	1738	1746	1752	rBV	3976484	5799050	55.30%	7.264%
48	13.161	1776	1782	1787	rVB2	259965	403088	3.84%	0.505%
49	13.267	1794	1800	1805	rBV5	114964	216679	2.07%	0.271%
50	13.331	1805	1811	1813	rBV4	258838	389706	3.72%	0.488%
51	13.490	1831	1838	1839	rBV	426965	582066	5.55%	0.729%
52	13.649	1854	1865	1870	rBV2	833404	1585376	15.12%	1.986%
53	13.702	1870	1874	1884	rVB2	601678	928797	8.86%	1.164%
54	13.884	1900	1905	1908	rBV2	310218	481464	4.59%	0.603%
55	13.919	1908	1911	1917	rVB6	205515	310628	2.96%	0.389%
56	14.054	1929	1934	1939	rVB	252143	354694	3.38%	0.444%
57	14.236	1959	1965	1970	rVB2	127735	187542	1.79%	0.235%
58	14.330	1971	1981	1987	rVV3	1119777	1965430	18.74%	2.462%
59	14.389	1987	1991	1995	rVV5	144412	301693	2.88%	0.378%
60	14.548	2011	2018	2026	rVB4	133814	373423	3.56%	0.468%
61	14.630	2026	2032	2036	rBV8	69838	146002	1.39%	0.183%
62	14.730	2043	2049	2055	rBV2	253266	423240	4.04%	0.530%
63	14.806	2057	2062	2066	rVB2	171107	221840	2.12%	0.278%
64	14.953	2080	2087	2091	rBV4	169403	313685	2.99%	0.393%
65	15.006	2091	2096	2100	rVB	149417	202256	1.93%	0.253%
66	15.188	2124	2127	2131	rVB2	147049	189732	1.81%	0.238%
67	15.376	2154	2159	2164	rBV2	264149	433759	4.14%	0.543%
68	15.417	2164	2166	2172	rVB5	151428	207596	1.98%	0.260%
69	15.511	2177	2182	2183	rVV3	327498	456383	4.35%	0.572%
70	15.535	2183	2186	2192	rVV	715494	1081950	10.32%	1.355%
71	15.593	2192	2196	2200	rVB5	117945	153161	1.46%	0.192%

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72	15.670	2207	2209	2216	rVB8	88729	174724	1.67%	0.219%
73	15.823	2227	2235	2244	rBV	2975009	4423010	42.18%	5.541%
74	16.052	2269	2274	2275	rBV4	171317	228024	2.17%	0.286%
75	16.440	2335	2340	2346	rBV4	155716	262928	2.51%	0.329%
76	16.839	2404	2408	2413	rBV3	115228	162764	1.55%	0.204%
77	16.892	2413	2417	2426	rVB5	126806	284403	2.71%	0.356%
78	17.080	2443	2449	2453	rBV	1539240	2211626	21.09%	2.771%
79	17.121	2453	2456	2460	rVB	318773	396219	3.78%	0.496%
80	17.280	2478	2483	2487	rBV2	230685	350560	3.34%	0.439%
81	17.485	2514	2518	2523	rBV2	179133	281250	2.68%	0.352%
82	17.573	2530	2533	2540	rBV4	110458	157440	1.50%	0.197%
83	17.750	2556	2563	2568	rBV2	1210217	1555471	14.83%	1.949%
84	17.950	2592	2597	2601	rBV3	159763	229537	2.19%	0.288%
85	18.002	2601	2606	2611	rVB	481223	685250	6.53%	0.858%
86	18.138	2624	2629	2633	rVB5	125572	173050	1.65%	0.217%
87	18.261	2646	2650	2654	rBV3	197897	303206	2.89%	0.380%
88	18.302	2654	2657	2663	rVB2	260474	334407	3.19%	0.419%
89	18.402	2668	2674	2681	rVB3	308167	660947	6.30%	0.828%
90	18.707	2723	2726	2729	rBV	156455	191376	1.83%	0.240%
91	18.743	2729	2732	2736	rVV3	148338	207261	1.98%	0.260%
92	18.790	2736	2740	2745	rVB	217740	303502	2.89%	0.380%
93	19.025	2777	2780	2785	rVB2	182904	279615	2.67%	0.350%
94	19.095	2785	2792	2795	rVV2	188784	291832	2.78%	0.366%
95	19.125	2795	2797	2803	rVB4	108880	159853	1.52%	0.200%
96	19.219	2809	2813	2822	rVB3	113122	174668	1.67%	0.219%
97	19.530	2862	2866	2873	rBV3	101008	189019	1.80%	0.237%
98	19.712	2891	2897	2908	rVB	8713159	10485943	100.00%	13.136%
99	21.263	3156	3161	3169	rBV	2502905	3177304	30.30%	3.980%
100	23.519	3538	3545	3556	rBV	1452068	2863857	27.31%	3.588%

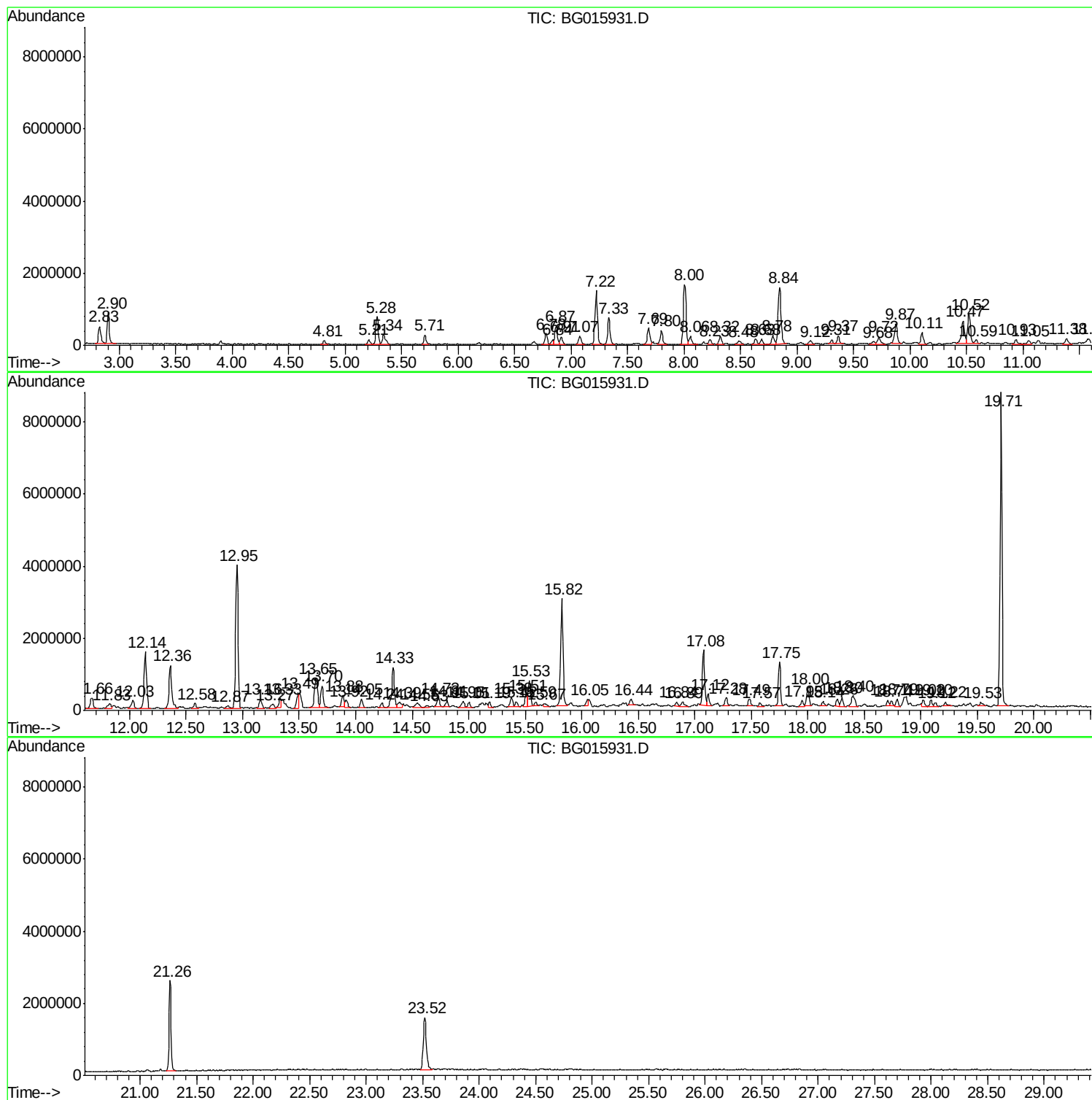
Sum of corrected areas: 79827503

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Instrument :
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG012015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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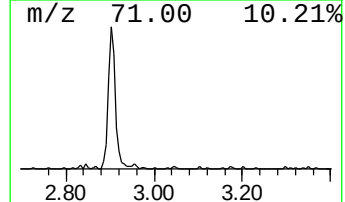
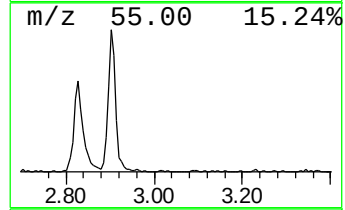
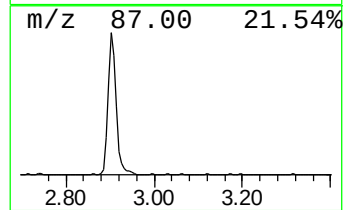
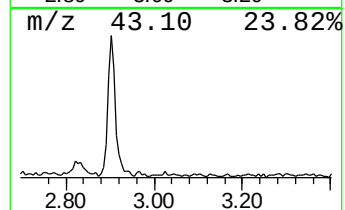
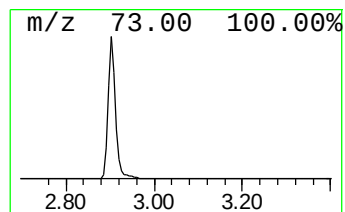
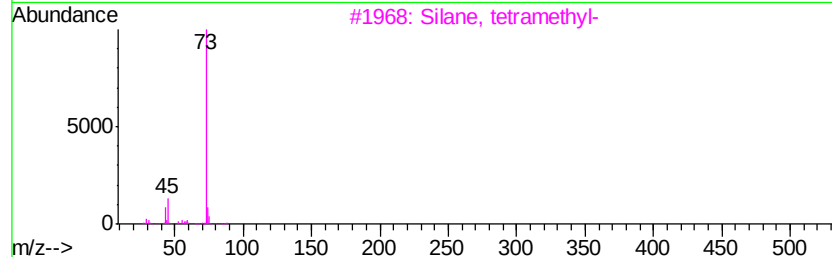
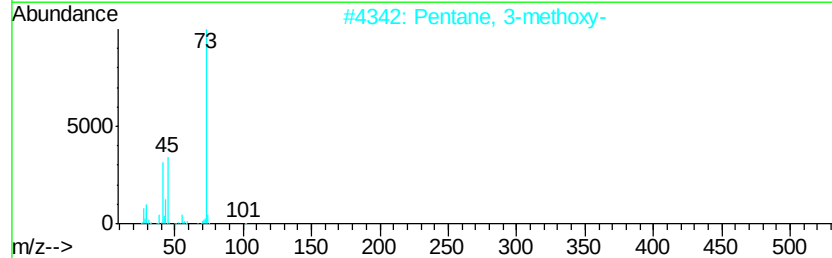
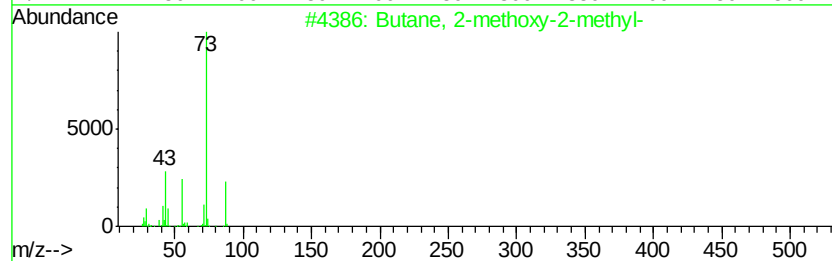
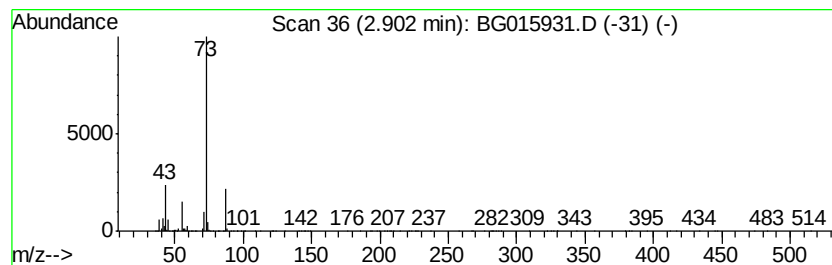
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG012015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.90	28.58 ng	1064570	1,4-Dichlorobenzene-d4	7.69

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
3		Silane, tetramethyl-	88	C4H12Si	000075-76-3	9
4		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9
5		N-Ethylformamide	73	C3H7NO	000627-45-2	9



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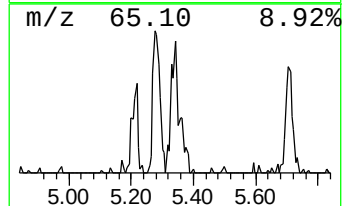
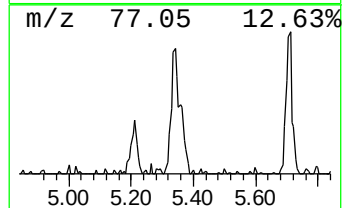
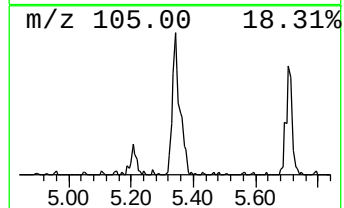
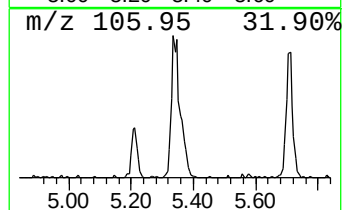
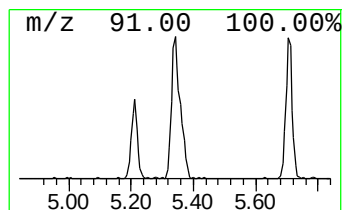
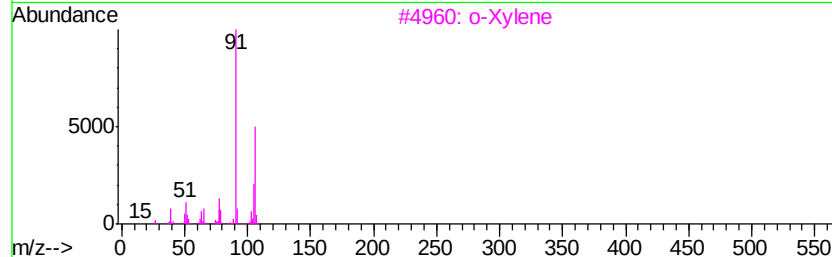
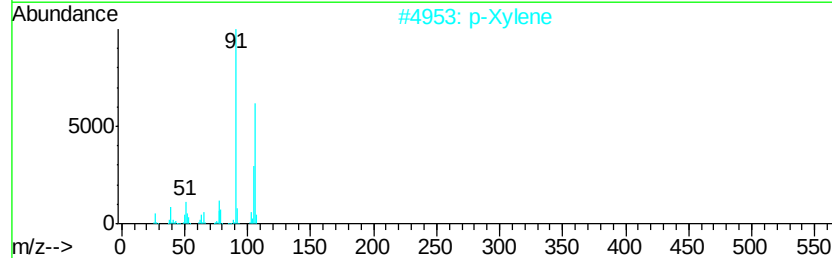
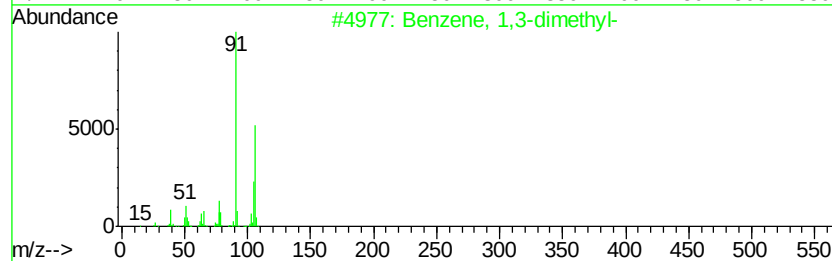
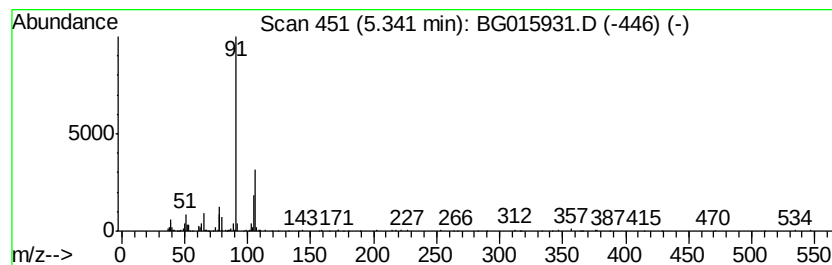
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG012015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Benzene, 1,3-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.34	14.53 ng	541283	1,4-Dichlorobenzene-d4	7.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	68
2		p-Xylene	106	C8H10	000106-42-3	68
3		o-Xylene	106	C8H10	000095-47-6	64
4		Ethylbenzene	106	C8H10	000100-41-4	52
5		Cyclopentene, 1-ethenyl-3-methyl...	106	C8H10	061142-07-2	49



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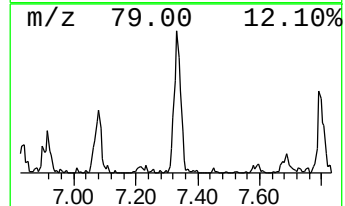
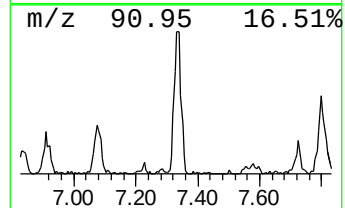
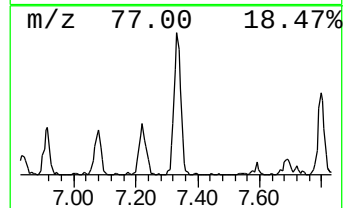
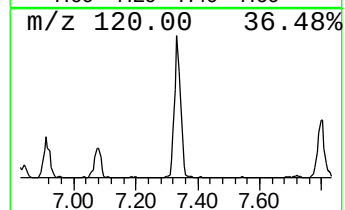
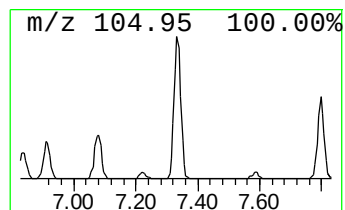
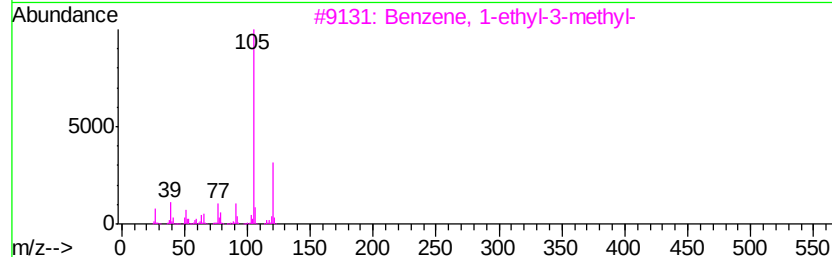
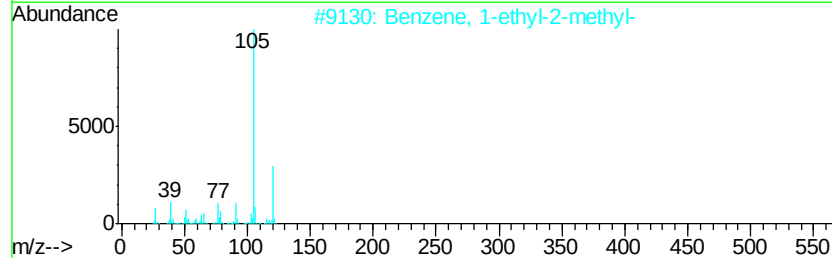
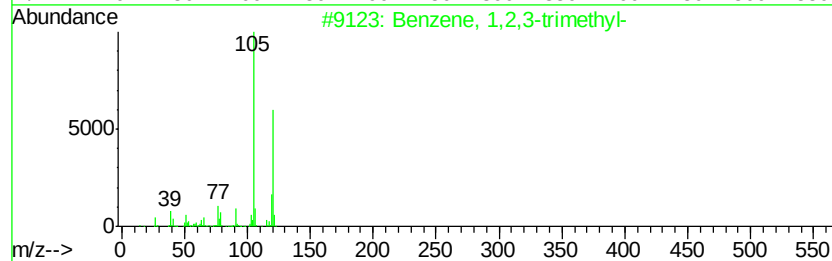
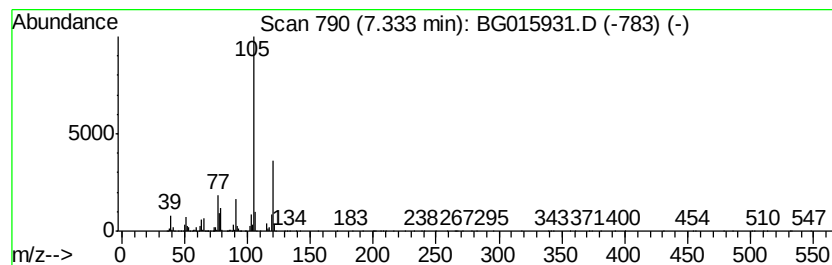
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Peak Number 5 Benzene, 1,2,3-trimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.33	29.18 ng	1086780	1,4-Dichlorobenzene-d4	7.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	93
2		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91
3		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4		Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	90
5		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	87



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG012615\
Data File : BG015931.D
Acq On : 27 Jan 2015 2:40
Operator : TP/IZ
Sample : G1139-09
Misc : S
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
121B2B

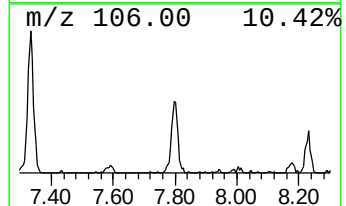
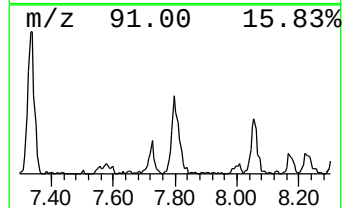
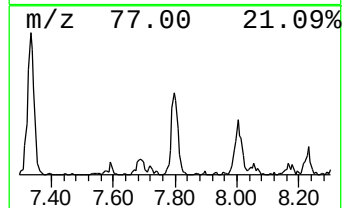
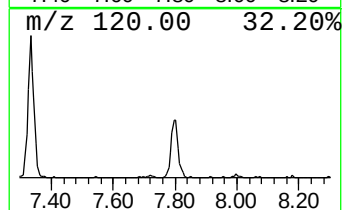
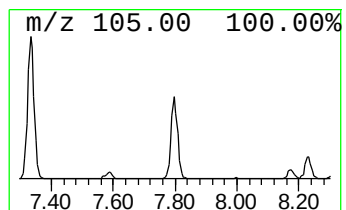
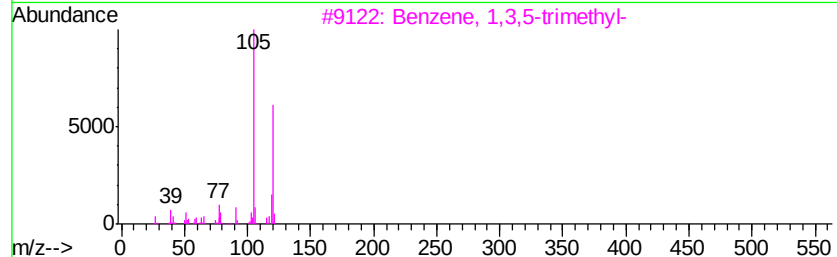
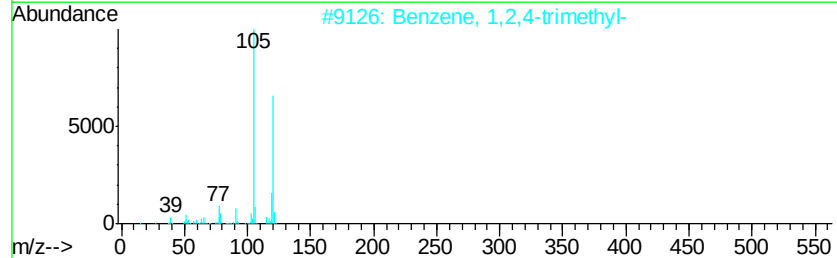
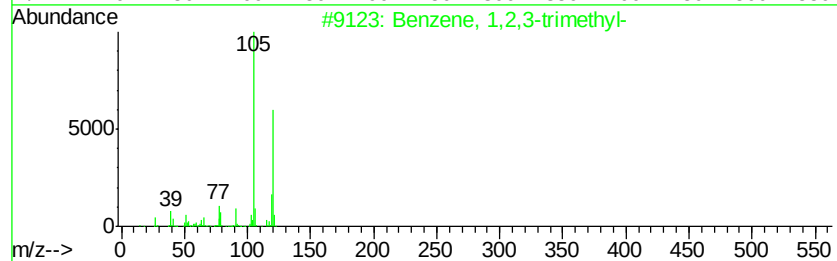
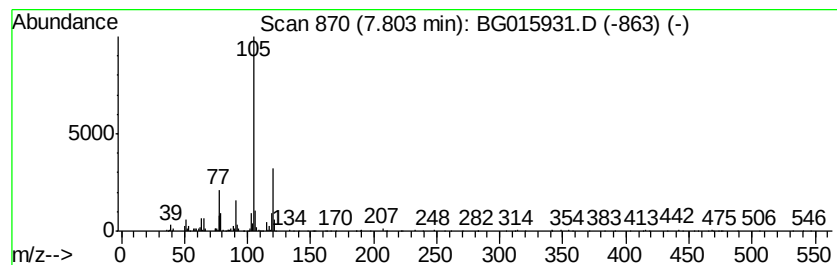
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Benzene, 1,2,4-trimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.80	16.76 ng	624079	1,4-Dichlorobenzene-d4	7.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	90
2		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	90
3		Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	90
4		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	81
5		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	76



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG012615\
Data File : BG015931.D
Acq On : 27 Jan 2015 2:40
Operator : TP/IZ
Sample : G1139-09
Misc : S
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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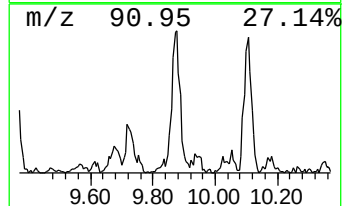
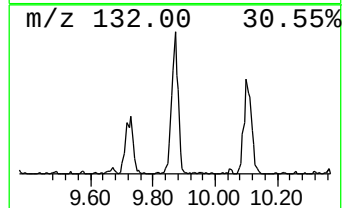
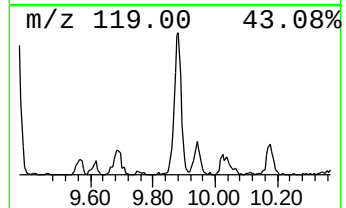
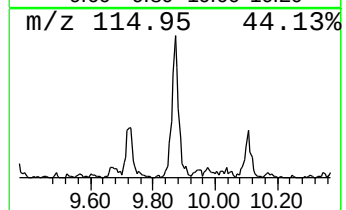
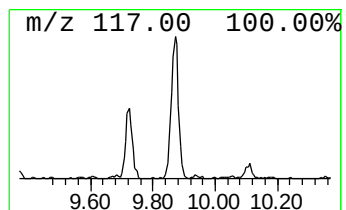
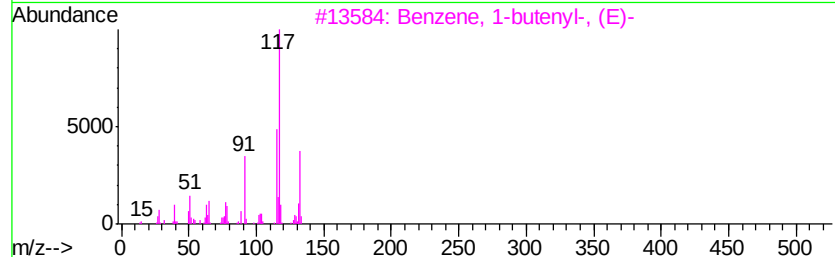
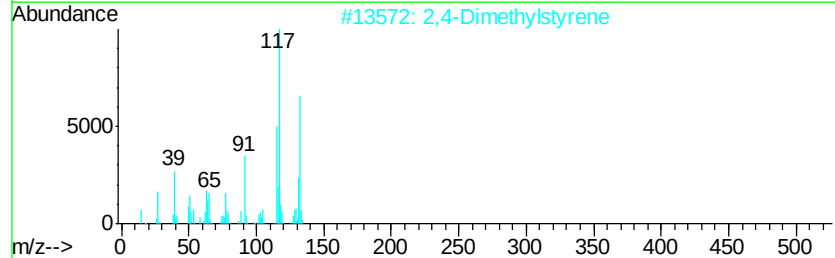
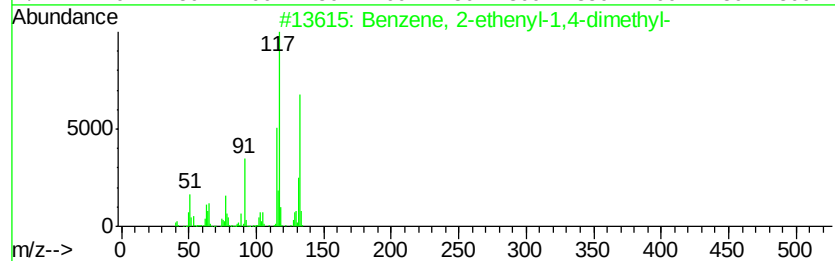
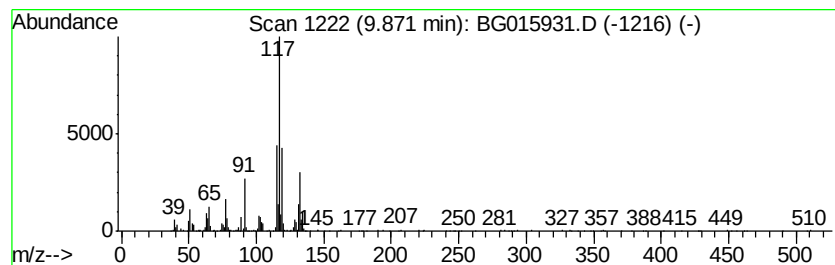
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.87	16.14 ng	994395	Naphthalene-d8	10.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	86
2		2,4-Dimethylstyrene	132	C10H12	002234-20-0	86
3		Benzene, 1-butenyl-, (E)-	132	C10H12	001005-64-7	86
4		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	86
5		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	83



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG012615\
Data File : BG015931.D
Acq On : 27 Jan 2015 2:40
Operator : TP/IZ
Sample : G1139-09
Misc : S
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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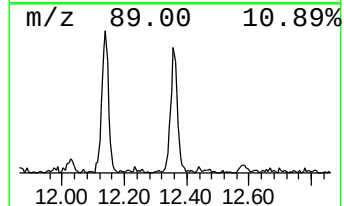
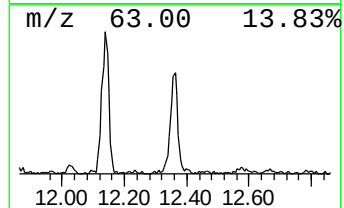
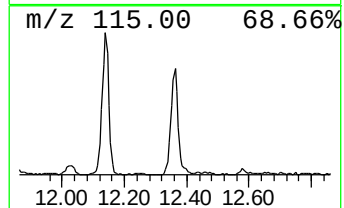
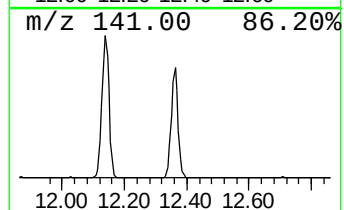
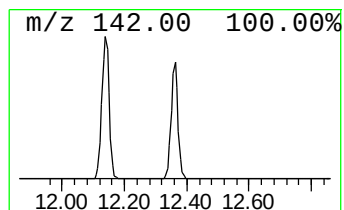
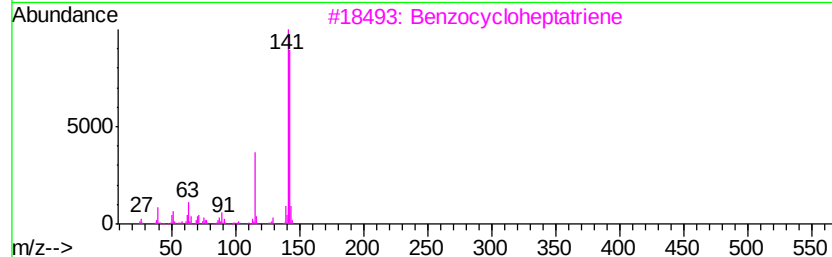
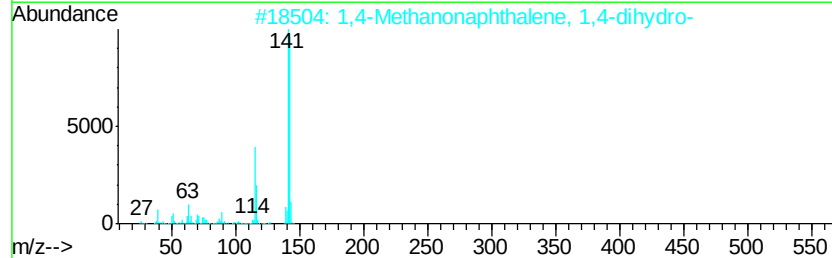
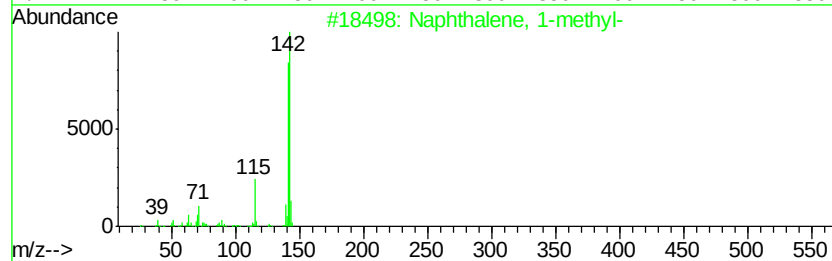
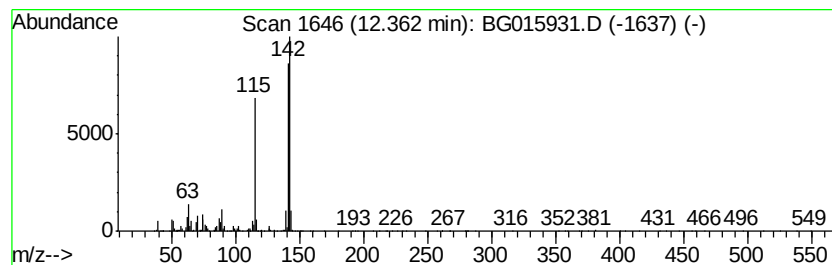
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG012015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Naphthalene, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.36	34.53 ng	2126660	Naphthalene-d8	10.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94
2		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	94
3		Benzocycloheptatriene	142	C11H10	000264-09-5	91
4		Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	90
5		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	80



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG012615\
Data File : BG015931.D
Acq On : 27 Jan 2015 2:40
Operator : TP/IZ
Sample : G1139-09
Misc : S
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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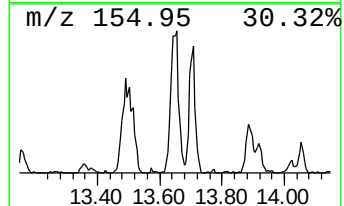
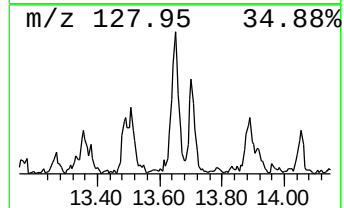
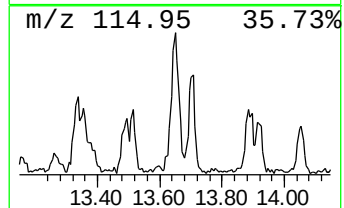
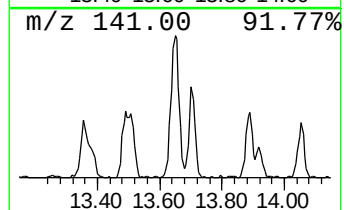
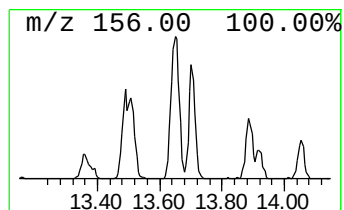
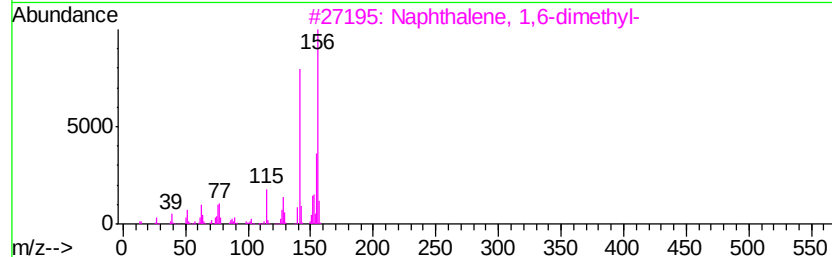
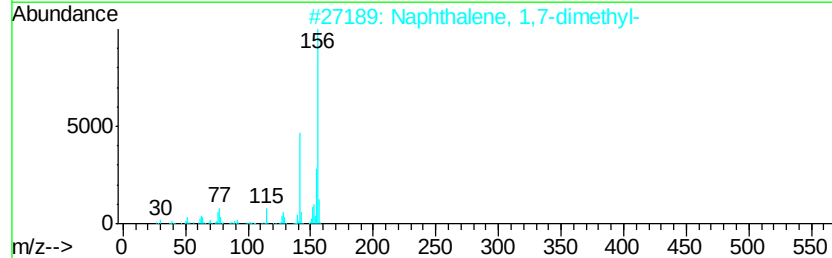
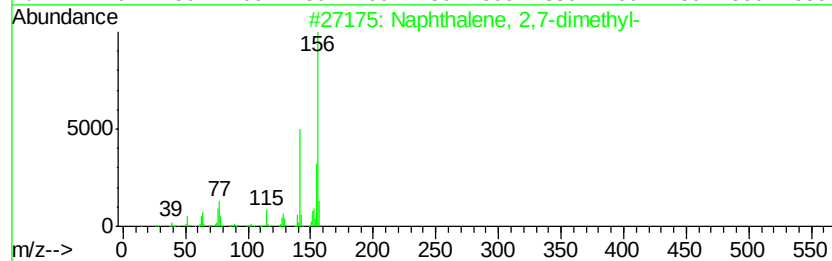
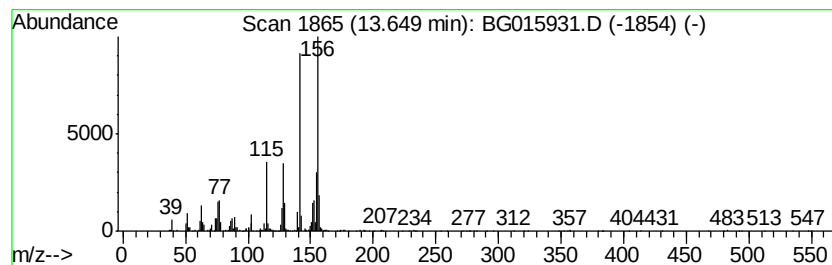
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG012015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Naphthalene, 2,7-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.65	16.13 ng	1585380	Acenaphthene-d10	14.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	95
2		Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	93
3		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	93
4		Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	93
5		Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	90



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG012615\
Data File : BG015931.D
Acq On : 27 Jan 2015 2:40
Operator : TP/IZ
Sample : G1139-09
Misc : S
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG012015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	2.90	28.6	ng	1064570	1	7.69	744922	20.0
Benzene, 1,3-dime...	5.34	14.5	ng	541283	1	7.69	744922	20.0
Benzene, 1,2,3-tr...	7.33	29.2	ng	1086780	1	7.69	744922	20.0
Benzene, 1,2,4-tr...	7.80	16.8	ng	624079	1	7.69	744922	20.0
Benzene, 2-etheny...	9.87	16.1	ng	994395	2	10.47	1231910	20.0
Naphthalene, 1-me...	12.36	34.5	ng	2126660	2	10.47	1231910	20.0
Naphthalene, 2,7-...	13.65	16.1	ng	1585380	3	14.33	1965430	20.0