

Data Path : Z:\HPCHEM1\BNA G\DATA\BG042215\
 Data File : BG016640.D
 Acq On : 23 Apr 2015 5:31
 Operator : TP/IZ
 Sample : G1976-02
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 S8-0435

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-BG042115.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	2.721	3	5	15	rVB	305902	518247	1.70%	0.376%
2	2.803	15	19	26	rVB	404503	533509	1.75%	0.387%
3	3.332	99	109	118	rBV2	385605	758429	2.48%	0.551%
4	3.820	183	192	203	rBV	451600	690615	2.26%	0.501%
5	5.153	411	419	427	rBV	5603081	9390280	30.74%	6.817%
6	5.247	427	435	436	rBV	600550	788467	2.58%	0.572%
7	5.300	436	444	452	rVB	9913992	30551601	100.00%	22.180%
8	5.653	496	504	514	rVB	5013667	7642606	25.02%	5.549%
9	6.123	573	584	599	rBV	933342	1506280	4.93%	1.094%
10	6.616	660	668	674	rBV	965312	1438067	4.71%	1.044%
11	6.728	679	687	692	rBV	3602667	5458045	17.87%	3.963%
12	6.787	692	697	702	rVB	1728227	2406273	7.88%	1.747%
13	6.863	702	710	715	rBV2	2404832	3858341	12.63%	2.801%
14	7.027	731	738	745	rBV	2143121	3212961	10.52%	2.333%
15	7.180	758	764	771	rVV	1023452	1720251	5.63%	1.249%
16	7.292	775	783	789	rVB	5333591	8476467	27.74%	6.154%
17	7.633	835	841	845	rBV2	250506	457022	1.50%	0.332%
18	7.674	845	848	855	rVB	249842	368283	1.21%	0.267%
19	7.750	855	861	871	rVB	2948620	4724259	15.46%	3.430%
20	7.956	885	896	900	rBV	1056450	1875244	6.14%	1.361%
21	8.003	900	904	908	rBV	665987	1002907	3.28%	0.728%
22	8.179	929	934	940	rVB	330130	504350	1.65%	0.366%
23	8.273	944	950	953	rBV	587975	974137	3.19%	0.707%
24	8.432	971	977	982	rBV	287176	533697	1.75%	0.387%
25	8.526	984	993	997	rVV6	175730	524271	1.72%	0.381%
26	8.585	998	1003	1007	rVV	341074	588632	1.93%	0.427%
27	8.637	1007	1012	1020	rVB2	643707	1427181	4.67%	1.036%
28	8.737	1023	1029	1033	rBV	469721	756260	2.48%	0.549%
29	8.802	1033	1040	1047	rBV3	948925	2606891	8.53%	1.893%
30	8.978	1065	1070	1076	rVB4	1244732	2911158	9.53%	2.114%
31	9.066	1080	1085	1093	rBV3	490469	1005209	3.29%	0.730%
32	9.178	1100	1104	1112	rBV6	148784	328309	1.07%	0.238%
33	9.260	1112	1118	1123	rVB2	260806	489682	1.60%	0.356%
34	9.319	1123	1128	1132	rBV	475094	934676	3.06%	0.679%

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Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG042115.M
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35	9.501	1151	1159	1172	rBV4	486457	1273664	4.17%	0.925%
36	9.624	1172	1180	1184	rBV5	184953	397756	1.30%	0.289%
37	9.777	1201	1206	1209	rBV2	182684	312231	1.02%	0.227%
38	9.830	1209	1215	1222	rVV2	913900	1646073	5.39%	1.195%
39	10.030	1244	1249	1250	rVV2	210732	320013	1.05%	0.232%
40	10.053	1250	1253	1258	rVB	291526	427197	1.40%	0.310%
41	10.424	1305	1316	1319	rBV2	589403	1237243	4.05%	0.898%
42	10.471	1319	1324	1331	rVV	2115983	3708748	12.14%	2.693%
43	10.958	1398	1407	1412	rBV2	304820	790464	2.59%	0.574%
44	11.064	1419	1425	1434	rVB5	286357	749547	2.45%	0.544%
45	11.211	1443	1450	1457	rBV7	287702	804364	2.63%	0.584%
46	11.281	1458	1462	1465	rVV4	255632	493699	1.62%	0.358%
47	11.334	1466	1471	1476	rVV4	404624	945988	3.10%	0.687%
48	11.393	1477	1481	1486	rVB4	180776	355382	1.16%	0.258%
49	11.781	1535	1547	1551	rBV6	228946	770553	2.52%	0.559%
50	11.822	1551	1554	1562	rVB5	124911	340488	1.11%	0.247%
51	11.922	1567	1571	1585	rVV7	322729	920341	3.01%	0.668%
52	12.039	1585	1591	1595	rVV2	760471	1399673	4.58%	1.016%
53	12.092	1595	1600	1606	rVB	1196480	1997053	6.54%	1.450%
54	12.239	1617	1625	1629	rBV2	235752	524423	1.72%	0.381%
55	12.310	1630	1637	1644	rVB	850140	1514404	4.96%	1.099%
56	12.492	1665	1668	1673	rVB3	219084	347504	1.14%	0.252%
57	12.621	1685	1690	1693	rBV3	238352	401246	1.31%	0.291%
58	12.762	1706	1714	1727	rVB5	671113	1664860	5.45%	1.209%
59	12.903	1732	1738	1742	rVV	1908730	2748573	9.00%	1.995%
60	12.950	1742	1746	1754	rVB2	275090	489393	1.60%	0.355%
61	13.432	1822	1828	1830	rBV4	166095	328291	1.07%	0.238%
62	13.596	1853	1856	1862	rVV	198167	354989	1.16%	0.258%
63	13.655	1862	1866	1878	rVB2	201911	441057	1.44%	0.320%
64	13.908	1904	1909	1917	rVB7	140736	350088	1.15%	0.254%
65	14.284	1968	1973	1983	rVB2	598087	999150	3.27%	0.725%
66	14.419	1991	1996	2000	rBV	236293	387374	1.27%	0.281%
67	15.776	2222	2227	2232	rBV	1062609	1496385	4.90%	1.086%
68	17.022	2435	2439	2444	rBV2	636888	896216	2.93%	0.651%
69	17.938	2591	2595	2603	rBV	278286	437432	1.43%	0.318%
70	19.654	2883	2887	2893	rVB	2075129	2620517	8.58%	1.902%
71	21.211	3148	3152	3159	rVB	805342	971374	3.18%	0.705%

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72	23.432	3525	3530	3540	rVB2	493699	914658	2.99%	0.664%
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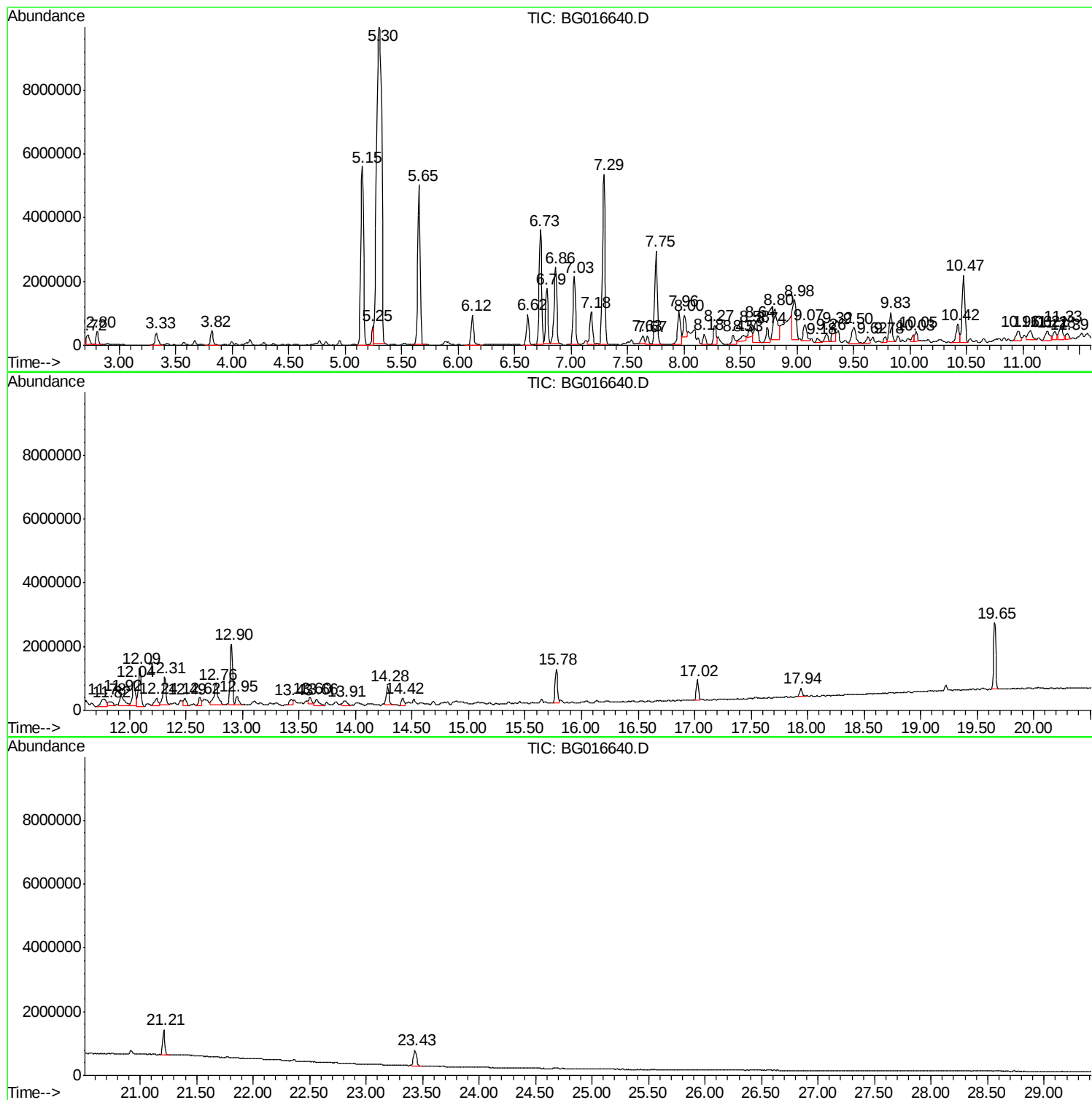
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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG042115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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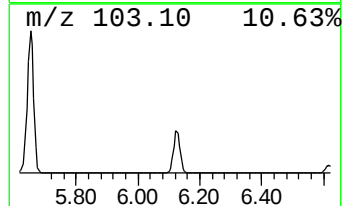
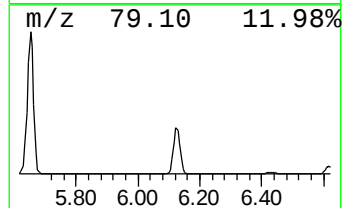
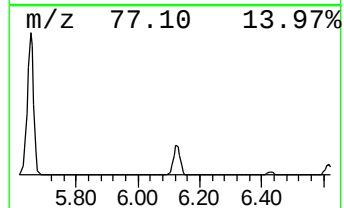
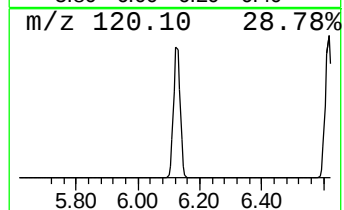
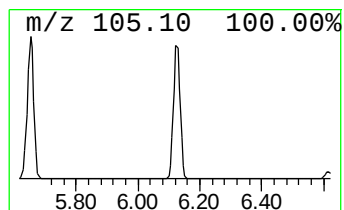
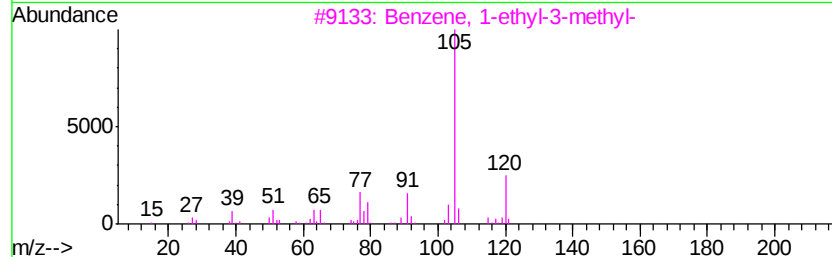
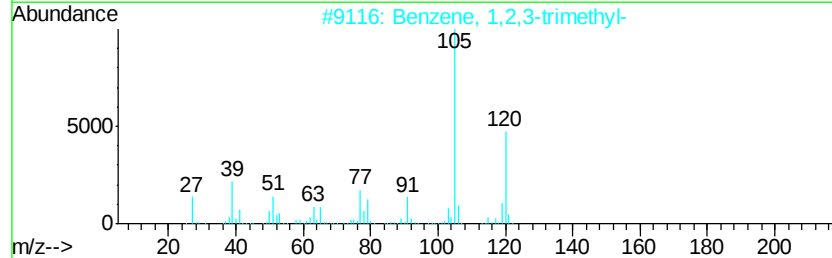
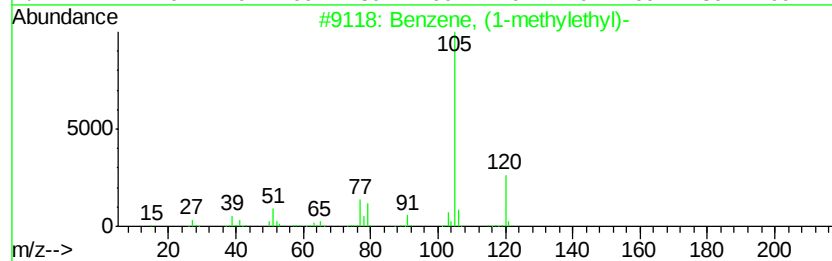
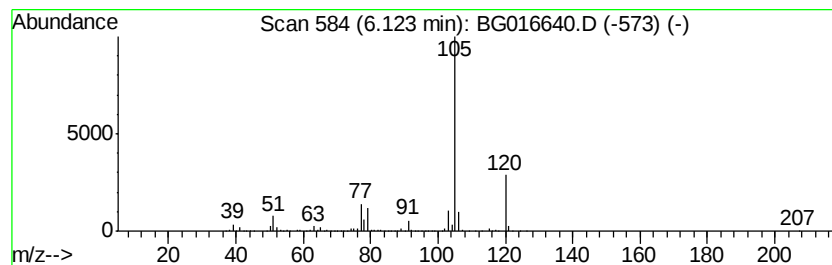
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG042115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Benzene, (1-methylethyl)- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.12	65.92 ng	1506280	1,4-Dichlorobenzene-d4	7.64

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	94
2		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
3		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	86
4		Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	83
5		2,4-Nonadiyne	120	C9H12	063621-15-8	83



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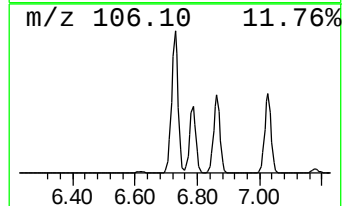
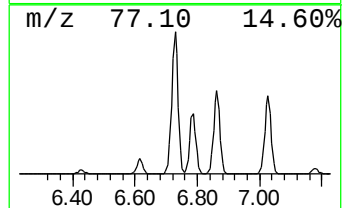
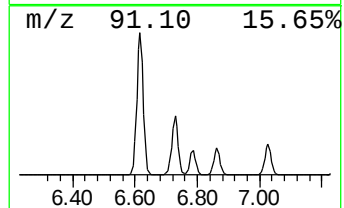
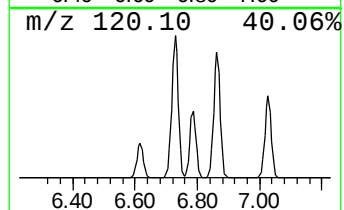
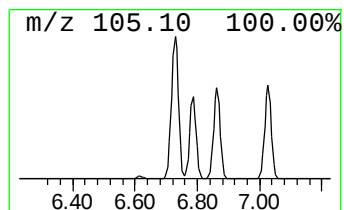
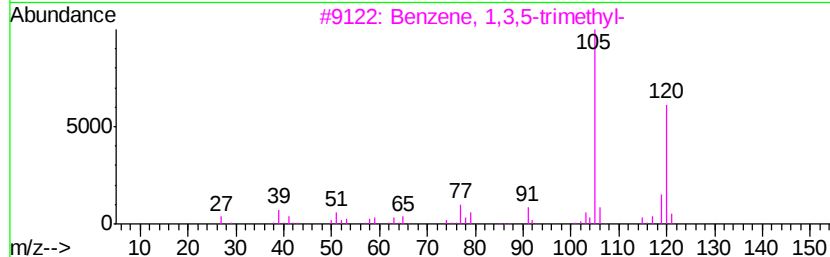
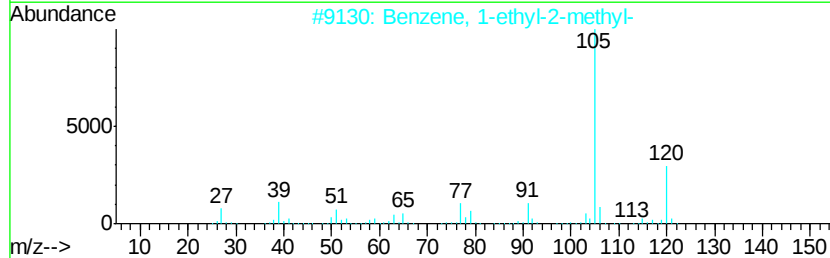
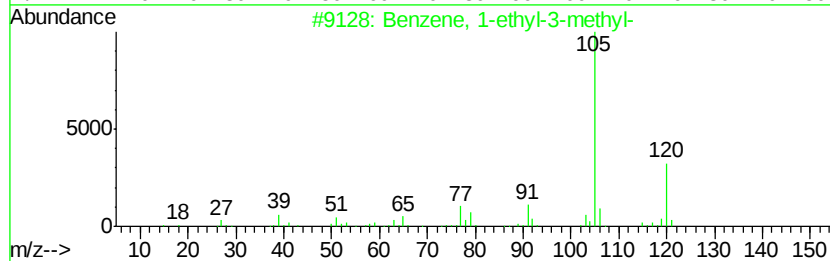
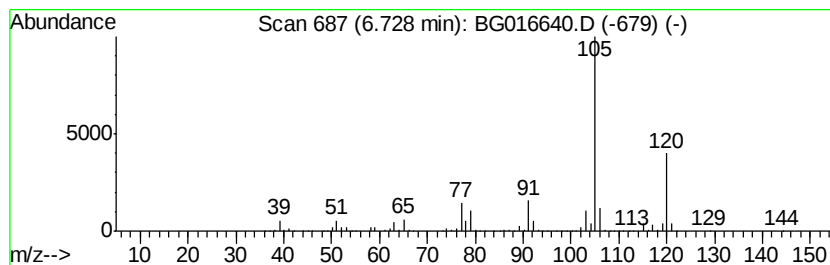
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG042115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Benzene, 1-ethyl-3-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.73	238.85 ng	5458050	1,4-Dichlorobenzene-d4	7.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	97
2		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3		Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	91
4		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	90



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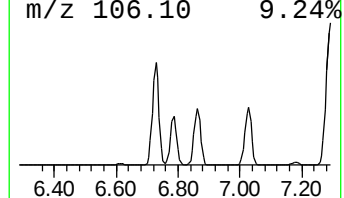
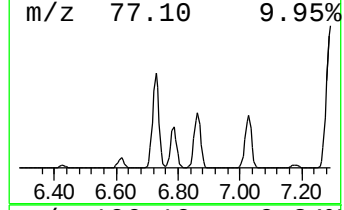
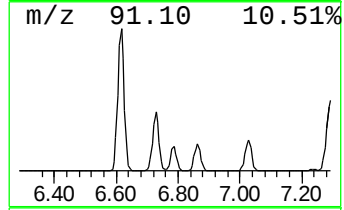
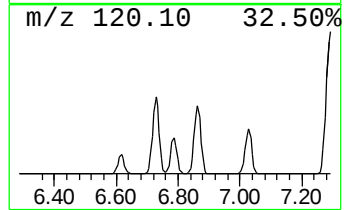
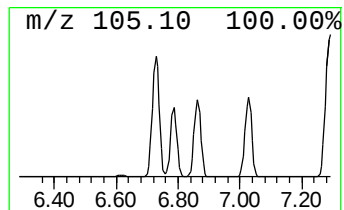
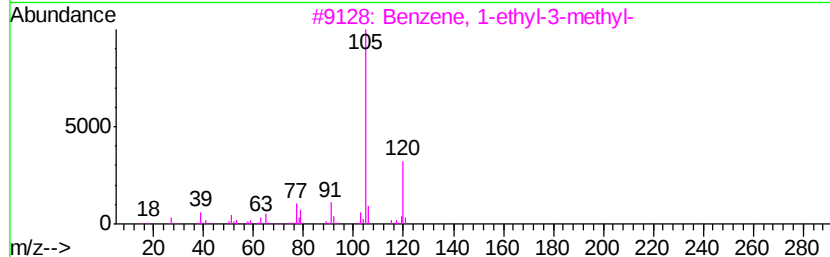
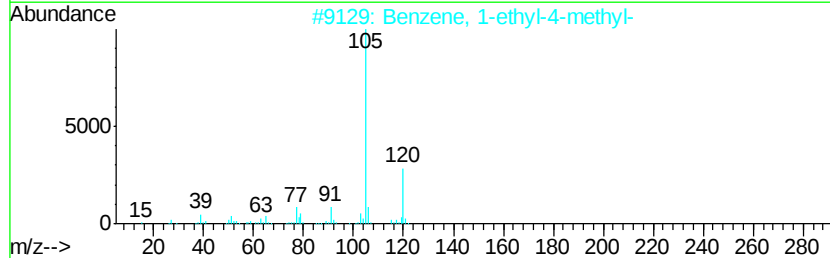
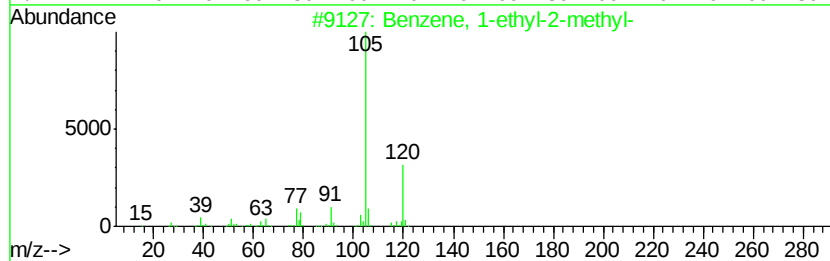
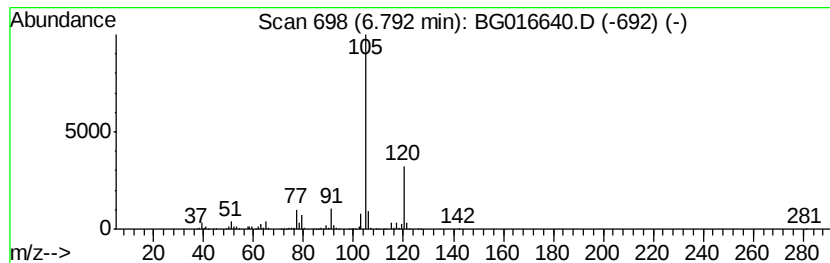
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Benzene, 1-ethyl-2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.79	105.30 ng	2406270	1,4-Dichlorobenzene-d4	7.64

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	94
3			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



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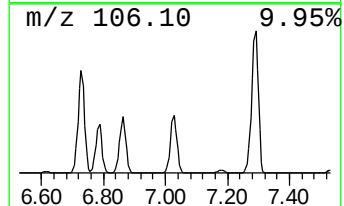
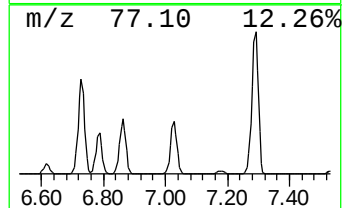
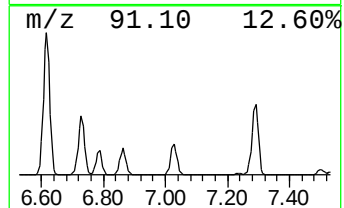
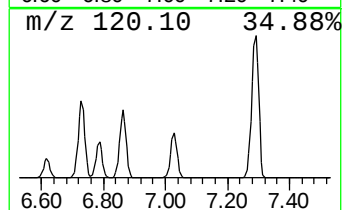
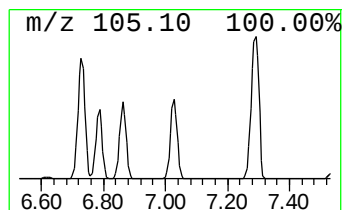
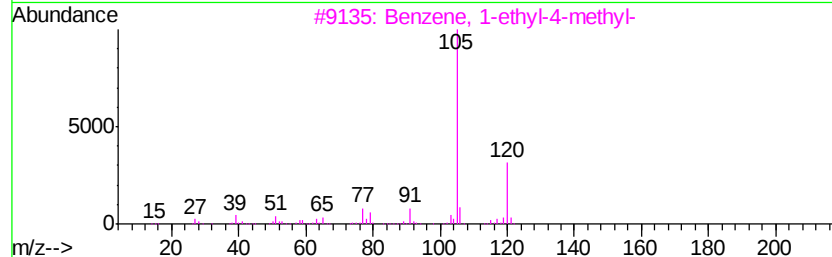
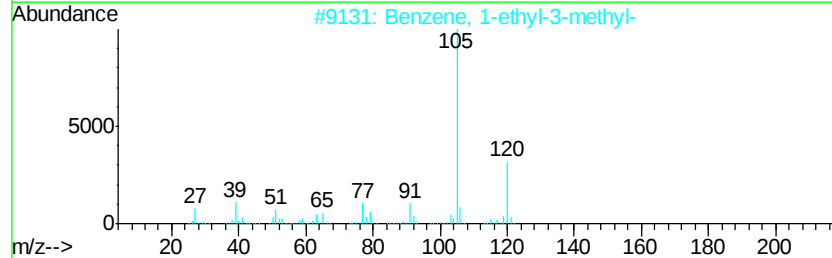
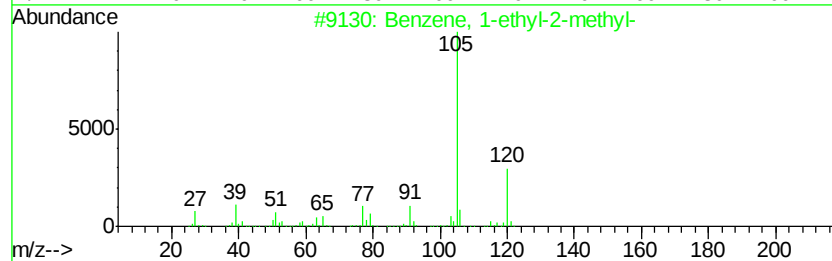
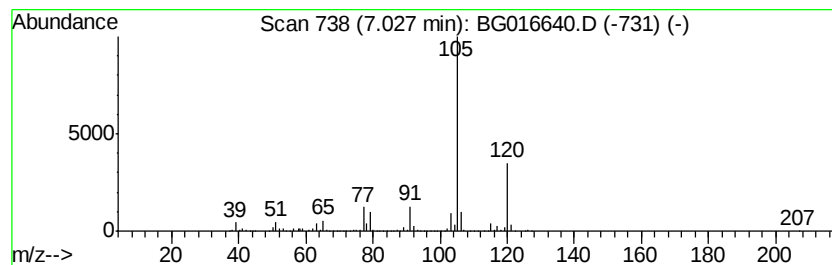
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Benzene, 1-ethyl-4-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.03	140.60 ng	3212960	1,4-Dichlorobenzene-d4	7.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
3		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5		Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	91



Data Path : Z:\HPCHEM1\BNA G\DATA\BG042215\
Data File : BG016640.D
Acq On : 23 Apr 2015 5:31
Operator : TP/IZ
Sample : G1976-02
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
S8-0435

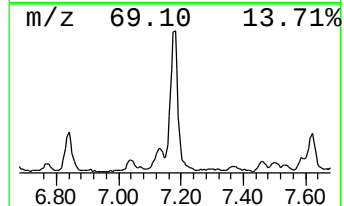
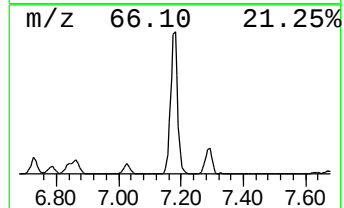
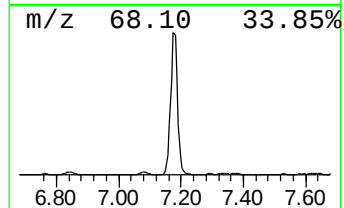
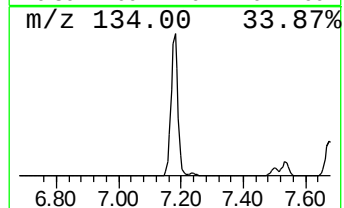
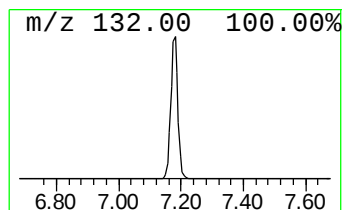
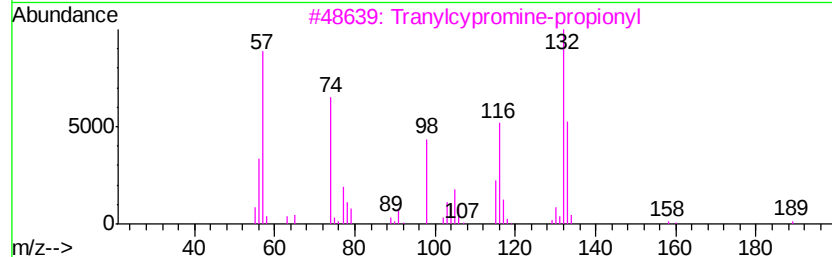
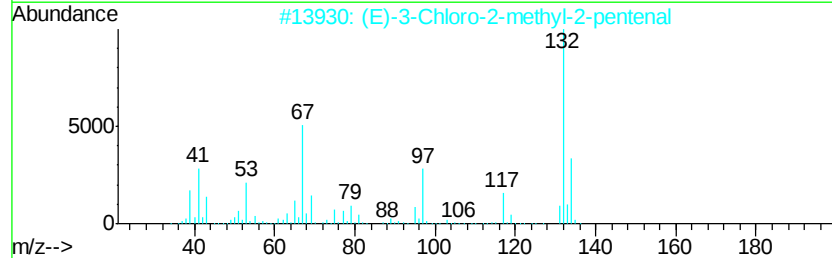
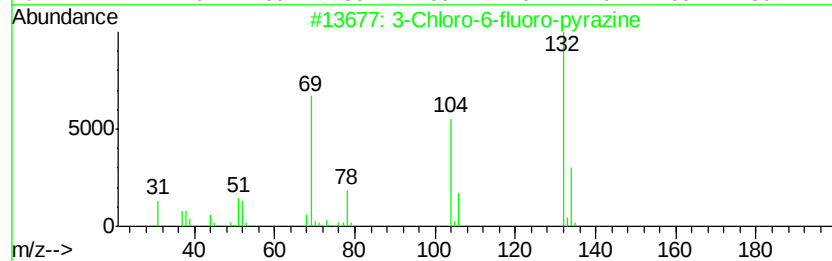
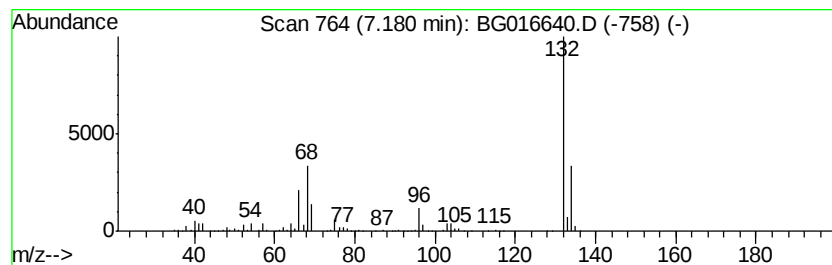
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown7.18 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.18	75.28 ng	1720250	1,4-Dichlorobenzene-d4	7.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9
5		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9



Data Path : Z:\HPCHEM1\BNA G\DATA\BG042215\
Data File : BG016640.D
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Instrument :
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ClientSampleId :
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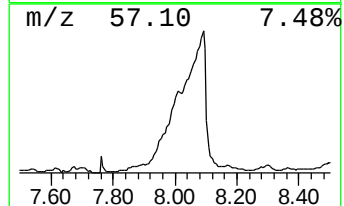
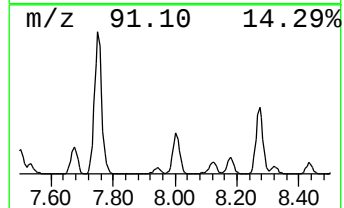
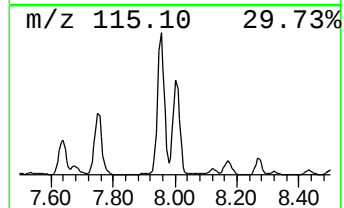
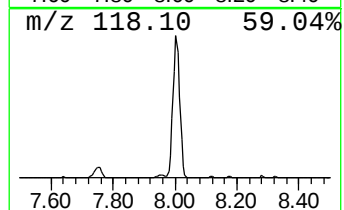
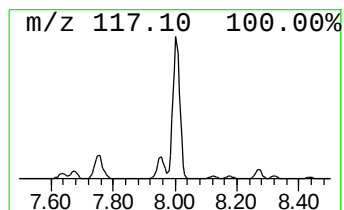
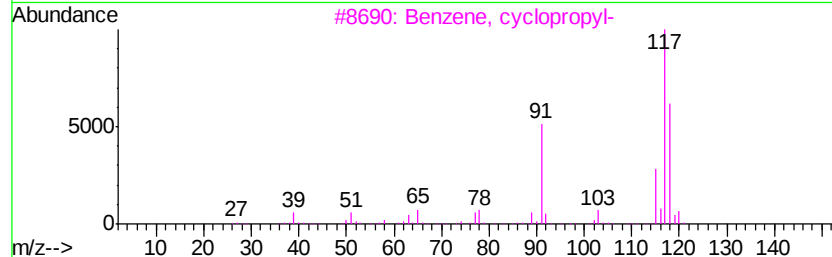
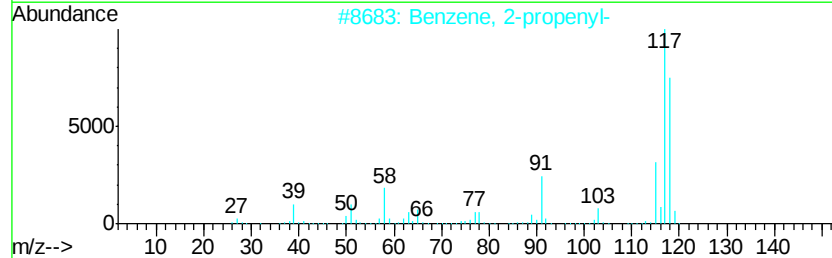
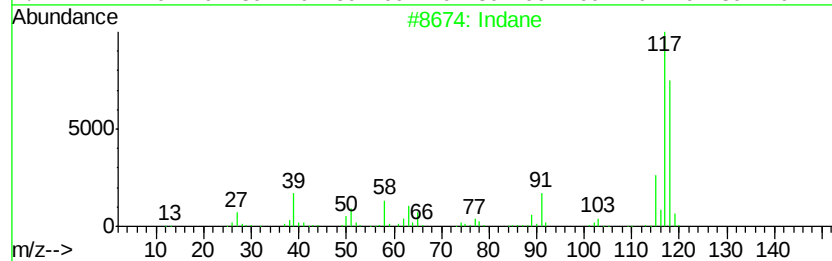
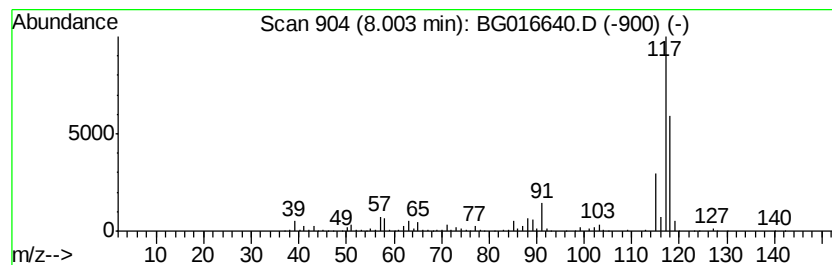
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Indane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.00	43.89 ng	1002910	1,4-Dichlorobenzene-d4	7.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	90
2	Benzene, 2-propenyl-	118	C9H10	000300-57-2	81
3	Benzene, cyclopropyl-	118	C9H10	000873-49-4	74
4	Deltacyclene	118	C9H10	007785-10-6	72
5	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	64



Data Path : Z:\HPCHEM1\BNA G\DATA\BG042215\
Data File : BG016640.D
Acq On : 23 Apr 2015 5:31
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Misc :
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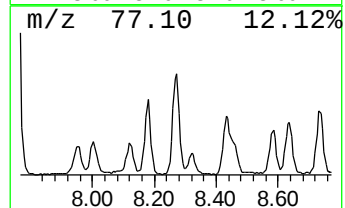
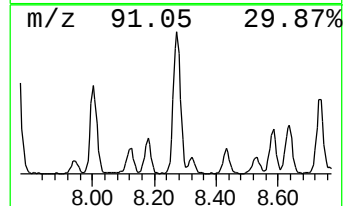
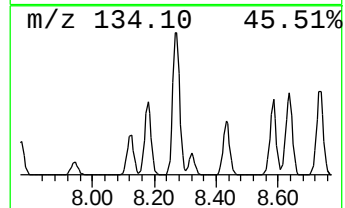
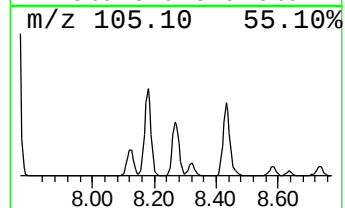
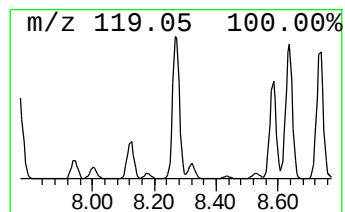
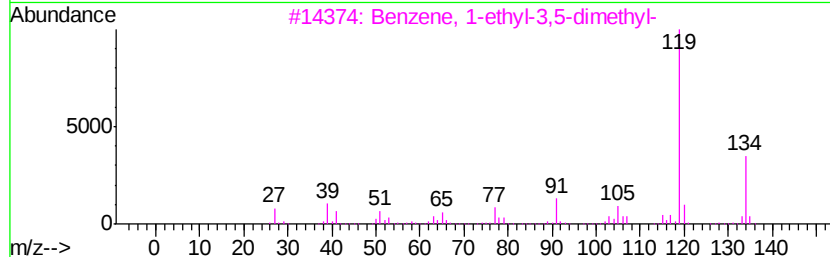
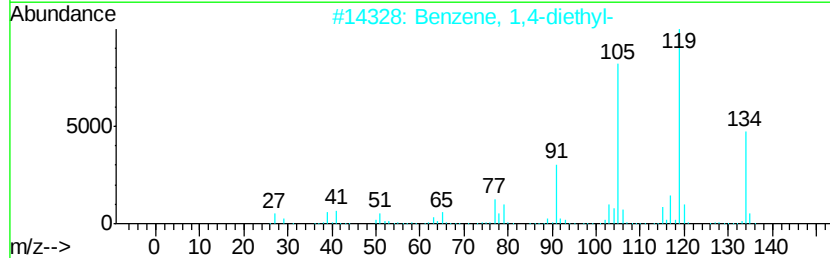
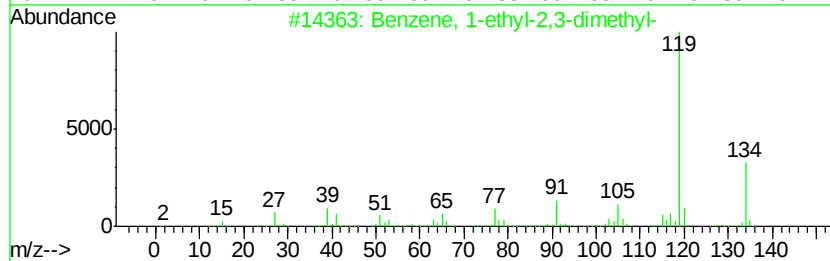
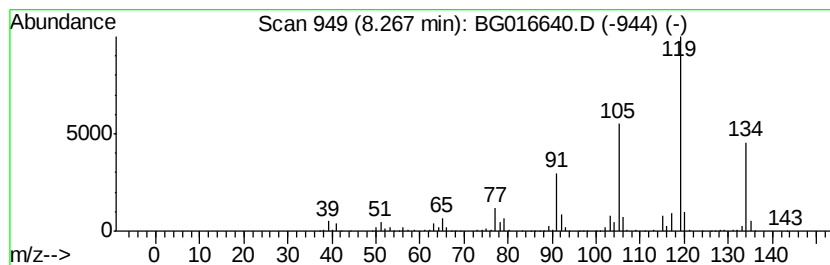
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.27	42.63 ng	974137	1,4-Dichlorobenzene-d4	7.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91
2		Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	91
3		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91
4		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91
5		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	91



Data Path : Z:\HPCHEM1\BNA G\DATA\BG042215\
Data File : BG016640.D
Acq On : 23 Apr 2015 5:31
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Sample : G1976-02
Misc :
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Instrument :
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ClientSampleId :
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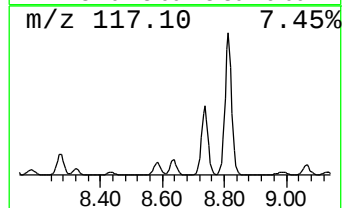
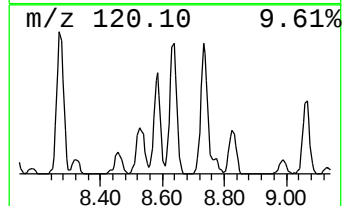
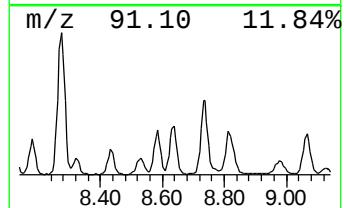
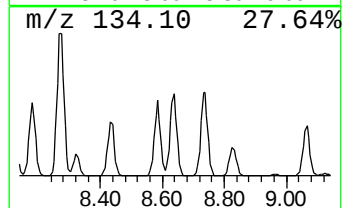
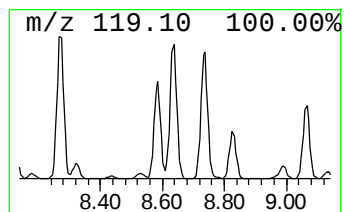
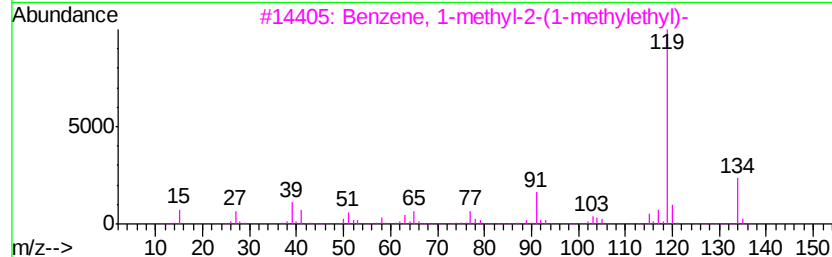
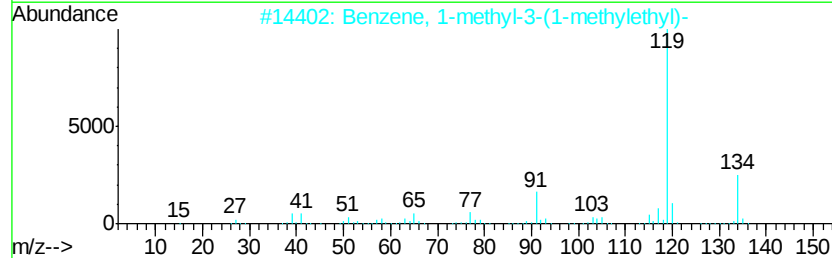
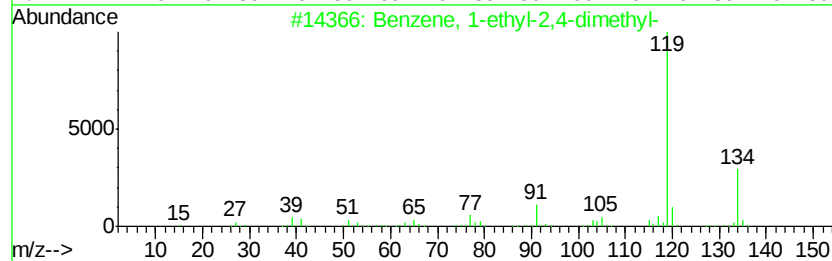
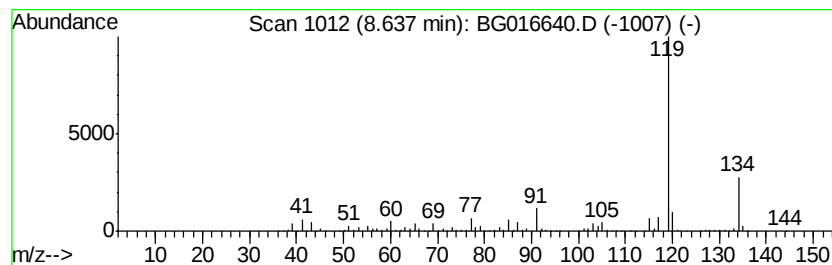
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.64	62.46 ng	1427180	1,4-Dichlorobenzene-d4	7.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	93
2		Benzene, 1-methyl-3-(1-methylethyl-)	134	C10H14	000535-77-3	93
3		Benzene, 1-methyl-2-(1-methylethyl-)	134	C10H14	000527-84-4	90
4		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	90
5		Benzene, 1-methyl-4-(1-methylethyl-)	134	C10H14	000099-87-6	90



Data Path : Z:\HPCHEM1\BNA G\DATA\BG042215\
Data File : BG016640.D
Acq On : 23 Apr 2015 5:31
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Sample : G1976-02
Misc :
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Instrument :
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ClientSampleId :
S8-0435

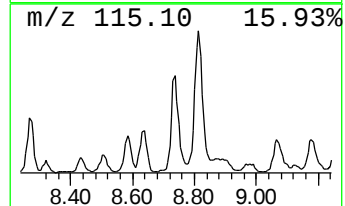
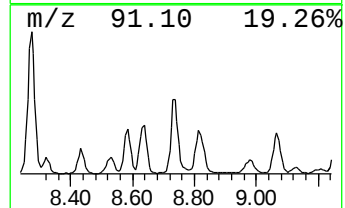
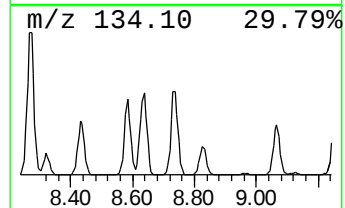
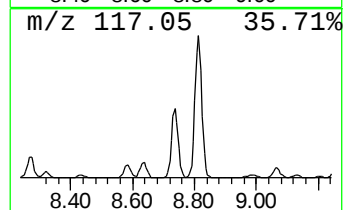
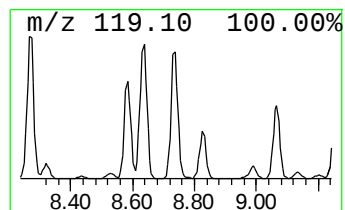
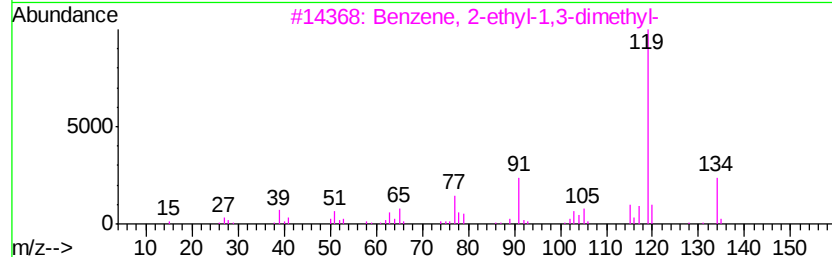
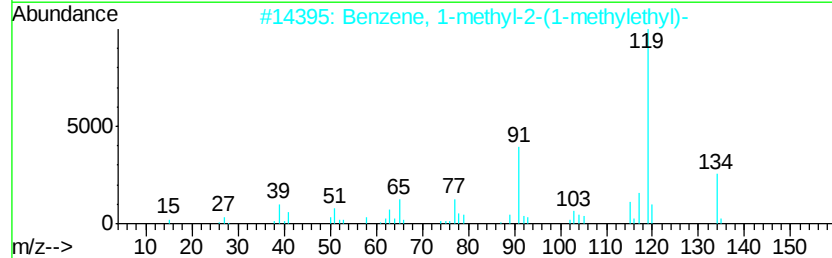
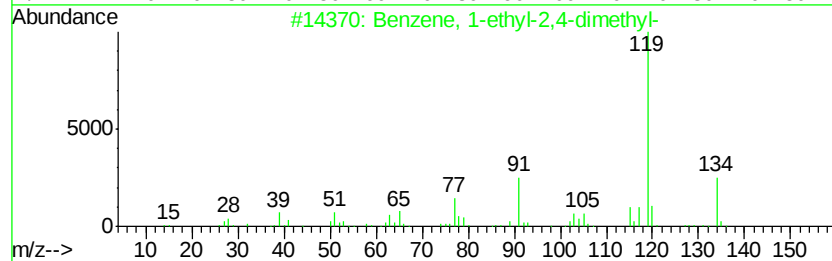
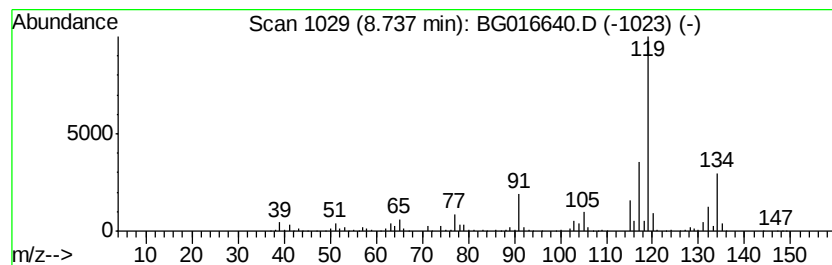
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Benzene, 1-methyl-2-(1-meth... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.74	33.10 ng	756260	1,4-Dichlorobenzene-d4	7.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	81
2		Benzene, 1-methyl-2-(1-methyleth...	134	C10H14	000527-84-4	81
3		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	81
4		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	76
5		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	76



Data Path : Z:\HPCHEM1\BNA G\DATA\BG042215\
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Acq On : 23 Apr 2015 5:31
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Sample : G1976-02
Misc :
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ClientSampled :
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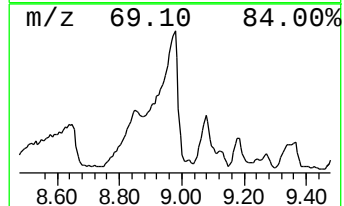
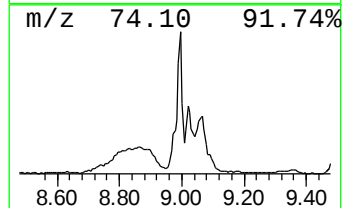
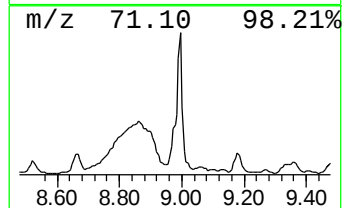
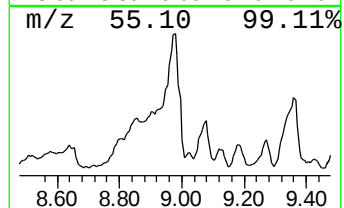
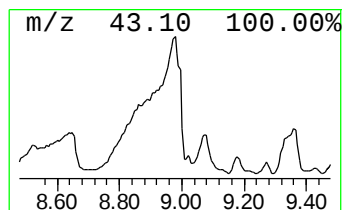
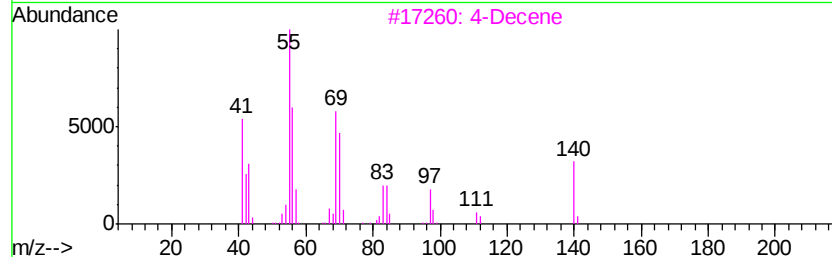
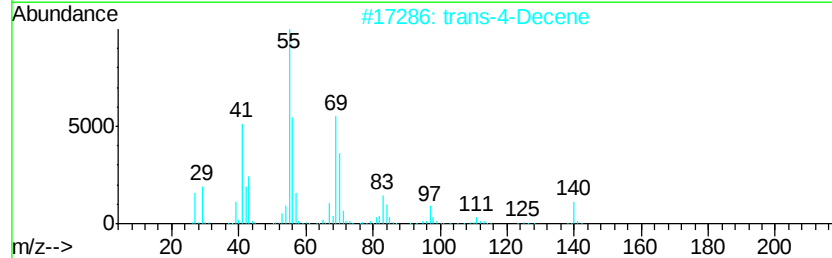
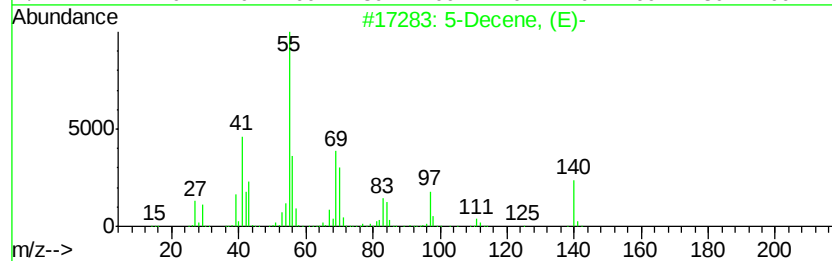
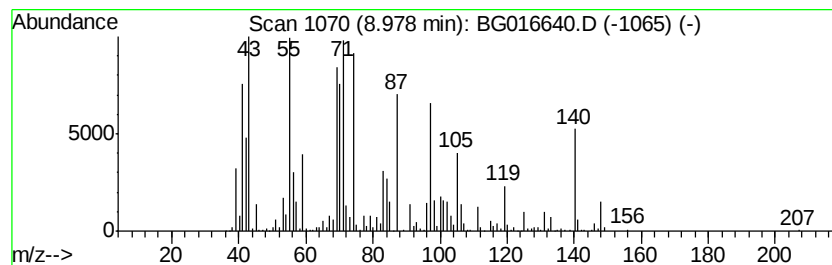
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 unknown8.98 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.98	127.40 ng	2911160	1,4-Dichlorobenzene-d4	7.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Decene, (E)-		140	C10H20	007433-56-9	45
2	trans-4-Decene		140	C10H20	019398-89-1	41
3	4-Decene		140	C10H20	019689-18-0	41
4	5-Decene		140	C10H20	019689-19-1	38
5	3-Decene		140	C10H20	019398-37-9	38



Data Path : Z:\HPCHEM1\BNA G\DATA\BG042215\
Data File : BG016640.D
Acq On : 23 Apr 2015 5:31
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ClientSampleId :
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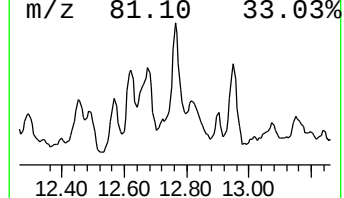
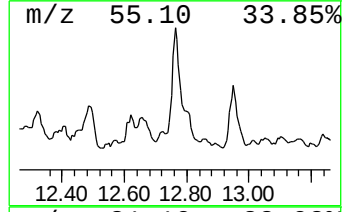
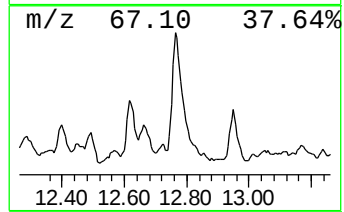
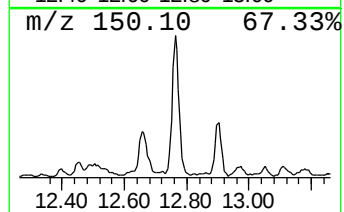
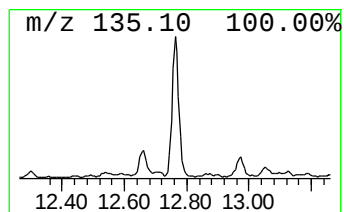
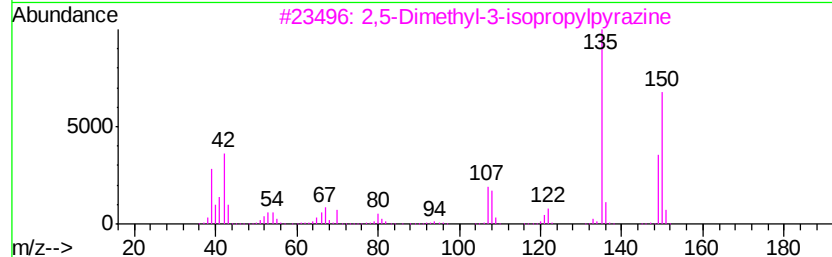
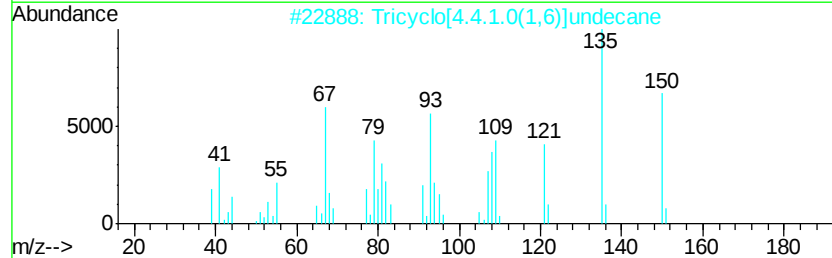
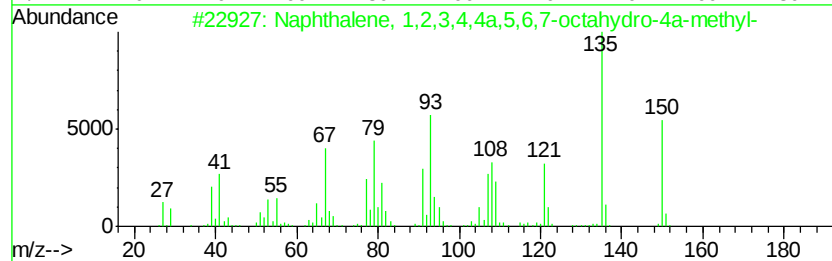
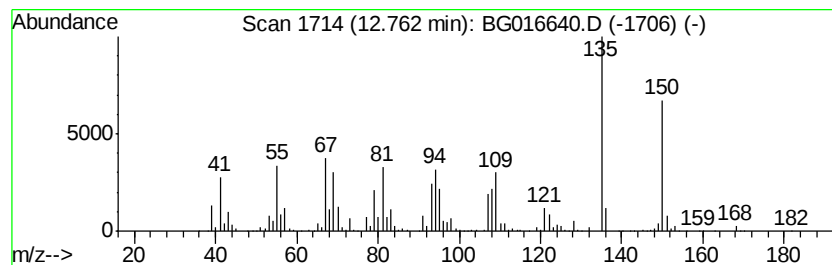
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG042115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 Naphthalene, 1,2,3,4,4a,5,6... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.76	33.33 ng	1664860	Acenaphthene-d10	14.28

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4,4a,5,6,7-oc...	150	C11H18	013943-77-6	74
2		Tricyclo[4.4.1.0(1,6)]undecane	150	C11H18	006571-73-9	59
3		2,5-Dimethyl-3-isopropylpyrazine	150	C9H14N2	013610-20-3	58
4		Cyclohexanone, 2-(2-butenyl)-	150	C10H14O	054166-48-2	49
5		Cyclohexene, 2-ethenyl-1,3,3-tri...	150	C11H18	005293-90-3	47



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG042215\
Data File : BG016640.D
Acq On : 23 Apr 2015 5:31
Operator : TP/IZ
Sample : G1976-02
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
S8-0435

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG042115.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
Benzene, (1-methy...	6.12	65.9	ng	1506280	1	7.64	457022	20.0
Benzene, 1-ethyl-...	6.73	238.8	ng	5458050	1	7.64	457022	20.0
Benzene, 1-ethyl-...	6.79	105.3	ng	2406270	1	7.64	457022	20.0
Benzene, 1-ethyl-...	7.03	140.6	ng	3212960	1	7.64	457022	20.0
unknown7.18	7.18	75.3	ng	1720250	1	7.64	457022	20.0
Indane	8.00	43.9	ng	1002910	1	7.64	457022	20.0
Benzene, 1-ethyl-...	8.27	42.6	ng	974137	1	7.64	457022	20.0
Benzene, 1-ethyl-...	8.64	62.5	ng	1427180	1	7.64	457022	20.0
Benzene, 1-methyl...	8.74	33.1	ng	756260	1	7.64	457022	20.0
unknown8.98	8.98	127.4	ng	2911160	1	7.64	457022	20.0
Naphthalene, 1,2,...	12.76	33.3	ng	1664860	3	14.28	999150	20.0