

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG111618\
 Data File : BG038055.D
 Acq On : 17 Nov 2018 5:09
 Operator : JU/SJ
 Sample : J5968-02
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MW-2-111418

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:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG111318.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	4.008	118	122	127	rVB3	11117	19483	1.28%	0.089%
2	4.237	155	161	165	rBV	42363	70173	4.61%	0.319%
3	4.725	240	244	249	rVV	15703	23861	1.57%	0.109%
4	4.772	249	252	262	rVB6	6928	16241	1.07%	0.074%
5	5.177	315	321	328	rVB6	10647	24437	1.61%	0.111%
6	5.565	379	387	393	rBV	426024	759065	49.92%	3.453%
7	5.630	393	398	404	rVV	163206	292261	19.22%	1.330%
8	5.718	404	413	428	rVB	308844	899486	59.16%	4.092%
9	6.076	467	474	490	rVB	196136	367436	24.16%	1.672%
10	6.552	548	555	565	rVB	99903	175011	11.51%	0.796%
11	7.046	632	639	650	rVB	224103	393225	25.86%	1.789%
12	7.152	650	657	663	rVV	139703	249878	16.43%	1.137%
13	7.216	663	668	675	rVV2	177204	415136	27.30%	1.889%
14	7.287	675	680	688	rVB	160125	295849	19.46%	1.346%
15	7.457	702	709	720	rVB	157084	272673	17.93%	1.240%
16	7.610	728	735	744	rBV	304570	558530	36.73%	2.541%
17	7.716	746	753	762	rVB	663108	1194594	78.56%	5.435%
18	7.939	785	791	793	rBV2	20035	32093	2.11%	0.146%
19	7.968	793	796	805	rVB2	25850	43666	2.87%	0.199%
20	8.086	809	816	825	rBV	152623	288928	19.00%	1.314%
21	8.192	827	834	843	rVB2	178730	325629	21.42%	1.481%
22	8.409	859	871	875	rBV	287919	554557	36.47%	2.523%
23	8.456	875	879	888	rVB	300191	532751	35.04%	2.424%
24	8.562	890	897	902	rBV	112504	192757	12.68%	0.877%
25	8.620	902	907	915	rVB2	45737	83095	5.46%	0.378%
26	8.709	915	922	928	rBV2	133858	258807	17.02%	1.177%
27	8.761	928	931	937	rVB2	26478	44949	2.96%	0.204%
28	8.879	944	951	955	rBV	26885	49860	3.28%	0.227%
29	9.026	969	976	981	rBV	66191	113876	7.49%	0.518%
30	9.079	981	985	992	rVB2	48546	79557	5.23%	0.362%
31	9.179	994	1002	1010	rBV2	348141	659621	43.38%	3.001%
32	9.267	1010	1017	1025	rVB2	429274	829896	54.58%	3.775%
33	9.425	1037	1044	1050	rVB4	15915	33893	2.23%	0.154%
34	9.514	1054	1059	1065	rBV2	21266	41930	2.76%	0.191%

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Integration Parameters: rteint.p
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 Sampling : 1
 Start Thrs: 0.2
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Filtering: 5
 Min Area: 3 % of largest Peak
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 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG111318.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	9.707	1085	1092	1097	rBV	168102	289178	19.02%	1.316%
36	9.766	1097	1102	1108	rVB	90505	157211	10.34%	0.715%
37	9.960	1130	1135	1139	rBV	22423	37338	2.46%	0.170%
38	10.078	1149	1155	1159	rBV4	22061	43966	2.89%	0.200%
39	10.136	1159	1165	1175	rVV	199364	363966	23.94%	1.656%
40	10.289	1178	1191	1198	rVV	369493	739308	48.62%	3.363%
41	10.342	1198	1200	1205	rVV4	25685	43856	2.88%	0.200%
42	10.407	1205	1211	1213	rVV5	15768	34662	2.28%	0.158%
43	10.465	1217	1221	1226	rVV2	24931	46769	3.08%	0.213%
44	10.530	1226	1232	1237	rVV2	32440	66864	4.40%	0.304%
45	10.577	1237	1240	1252	rVB	29061	50170	3.30%	0.228%
46	10.806	1274	1279	1281	rBV4	15927	26959	1.77%	0.123%
47	10.871	1281	1290	1291	rVV3	61700	139584	9.18%	0.635%
48	10.906	1291	1296	1300	rVV	230012	483758	31.81%	2.201%
49	10.953	1300	1304	1310	rVV	368218	697430	45.87%	3.173%
50	11.000	1310	1312	1319	rVV	41375	59419	3.91%	0.270%
51	11.088	1321	1327	1335	rVB4	26089	59372	3.90%	0.270%
52	11.552	1401	1406	1414	rVB2	28245	53843	3.54%	0.245%
53	11.805	1443	1449	1456	rBV	36007	57630	3.79%	0.262%
54	11.922	1462	1469	1473	rBV9	8037	20753	1.36%	0.094%
55	11.999	1475	1482	1491	rVV2	26241	61140	4.02%	0.278%
56	12.081	1491	1496	1509	rVB6	17713	41373	2.72%	0.188%
57	12.204	1509	1517	1521	rBV3	13913	24518	1.61%	0.112%
58	12.275	1521	1529	1537	rVB5	24547	58198	3.83%	0.265%
59	12.445	1553	1558	1564	rVB4	9041	17237	1.13%	0.078%
60	12.551	1570	1576	1583	rBV	251395	429838	28.27%	1.955%
61	12.616	1583	1587	1593	rVB6	11147	19078	1.25%	0.087%
62	12.769	1605	1613	1622	rVB	160449	286482	18.84%	1.303%
63	13.186	1679	1684	1690	rVB5	10810	16831	1.11%	0.077%
64	13.338	1702	1710	1719	rBV	715653	1142179	75.12%	5.196%
65	13.744	1777	1779	1791	rVB5	9995	21228	1.40%	0.097%
66	13.891	1791	1804	1810	rBV7	24454	66566	4.38%	0.303%
67	14.038	1822	1829	1833	rBV2	27367	50275	3.31%	0.229%
68	14.090	1833	1838	1846	rVB4	17836	29645	1.95%	0.135%
69	14.161	1846	1850	1854	rBV	11160	15955	1.05%	0.073%
70	14.719	1938	1945	1953	rBV2	358127	592042	38.94%	2.693%
71	14.884	1967	1973	1981	rBV6	21923	41820	2.75%	0.190%

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Sampling : 1 Min Area: 3 % of largest Peak
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Stop Thrs : 0 Peak Location: TOP

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72	15.107	2005	2011	2020	rBV	56713	97036	6.38%	0.441%
73	15.589	2088	2093	2107	rVB4	11488	26674	1.75%	0.121%
74	16.206	2191	2198	2210	rBV2	496731	777102	51.11%	3.535%
75	17.469	2406	2413	2423	rBV2	422985	689985	45.38%	3.139%
76	18.321	2552	2558	2566	rBV2	50587	83350	5.48%	0.379%
77	19.590	2769	2774	2786	rBV3	29296	53118	3.49%	0.242%
78	20.066	2849	2855	2861	rBV	1159893	1520547	100.00%	6.917%
79	20.336	2897	2901	2906	rVB2	17175	20962	1.38%	0.095%
80	20.836	2981	2986	2990	rVB	30392	34327	2.26%	0.156%
81	21.341	3068	3072	3077	rVB	37546	51227	3.37%	0.233%
82	21.746	3134	3141	3151	rBV	418899	732223	48.16%	3.331%
83	21.899	3161	3167	3171	rBV	49102	74268	4.88%	0.338%
84	22.516	3266	3272	3278	rVB	47766	81054	5.33%	0.369%
85	23.221	3384	3392	3398	rBV2	53634	104634	6.88%	0.476%
86	23.421	3418	3426	3432	rBV3	34598	69500	4.57%	0.316%
87	24.038	3525	3531	3541	rVB	46871	107561	7.07%	0.489%
88	25.066	3687	3706	3720	rBV2	228407	809180	53.22%	3.681%
89	26.159	3885	3892	3893	rBV6	21799	29916	1.97%	0.136%
90	27.575	4122	4133	4137	rBV9	11514	38937	2.56%	0.177%

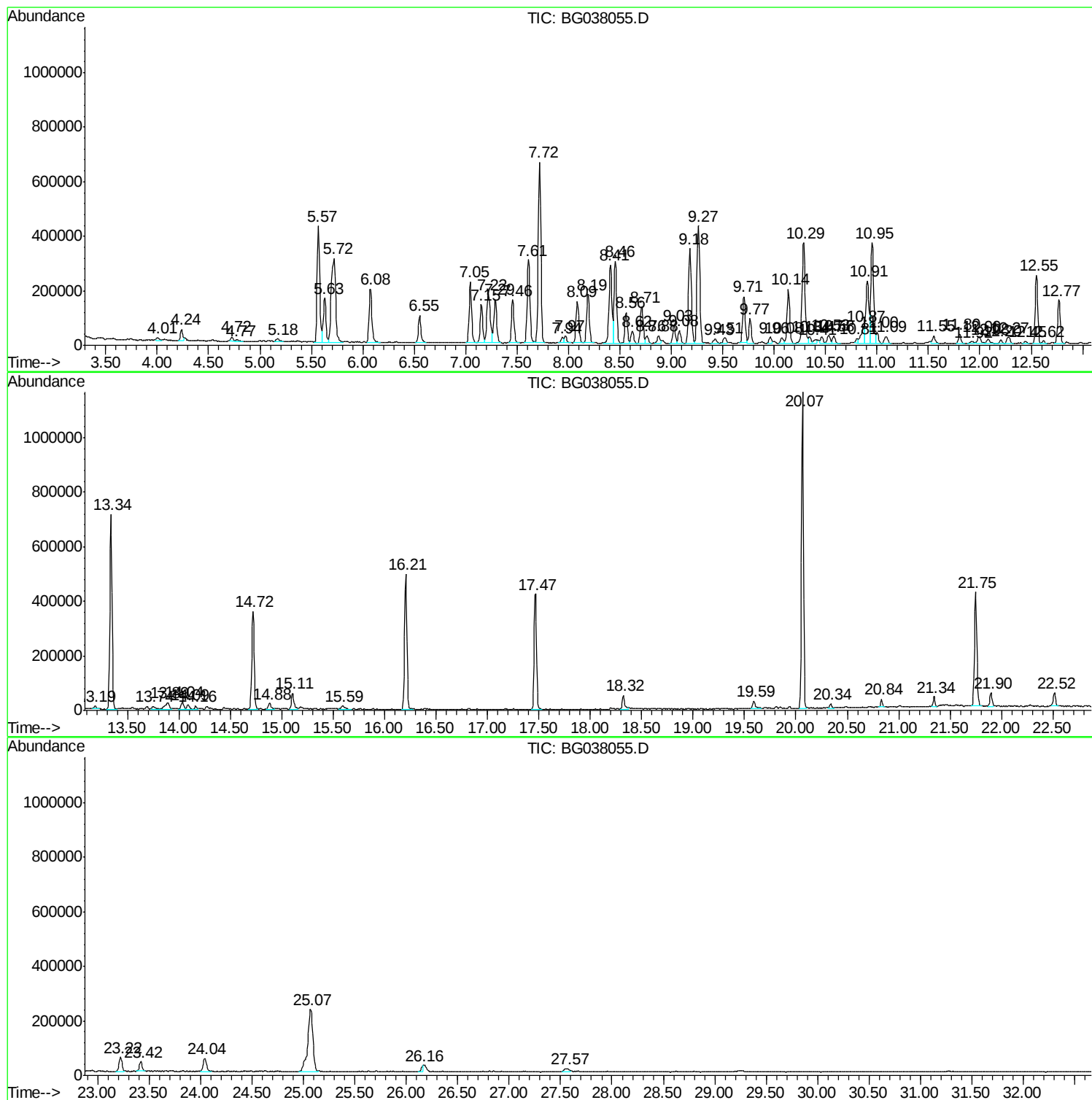
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG111318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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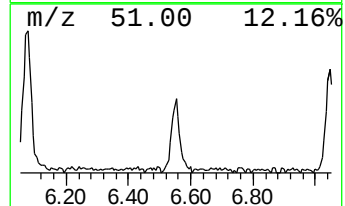
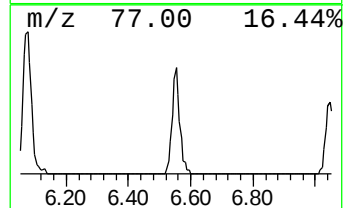
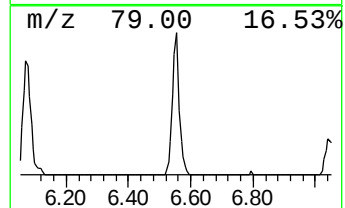
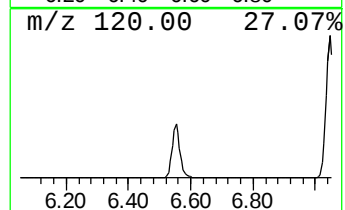
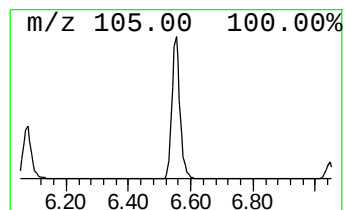
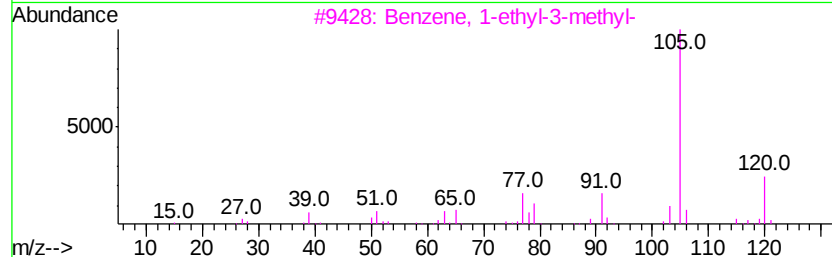
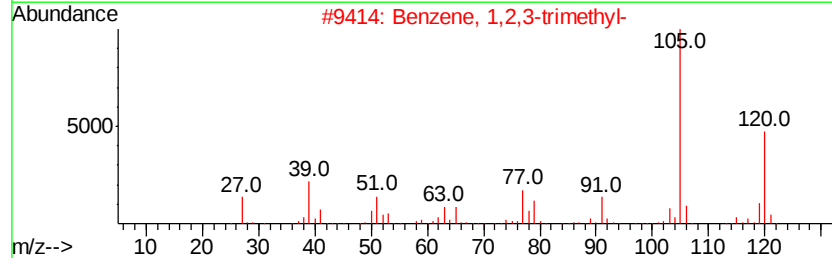
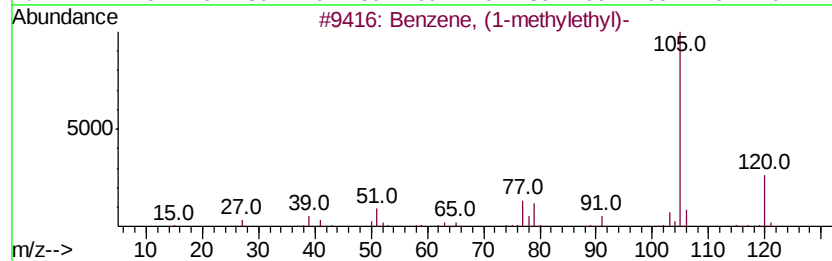
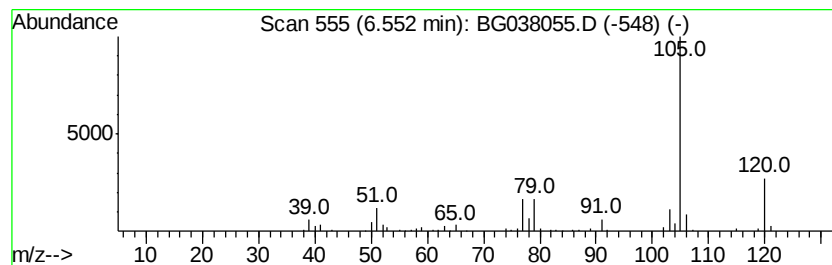
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG111318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzene, (1-methylethyl)- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.55	12.11 ng	175011	1,4-Dichlorobenzene-d4	8.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	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	74
4		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	72
5		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	72



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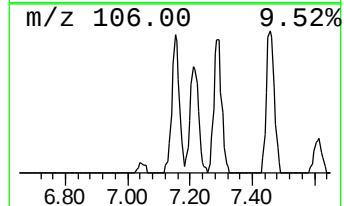
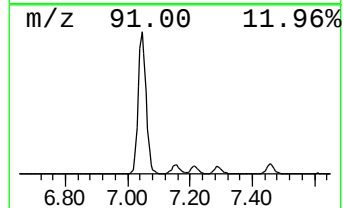
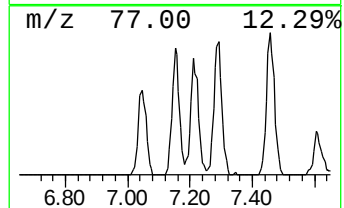
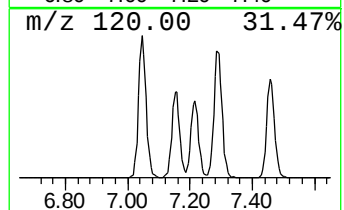
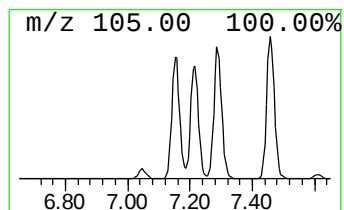
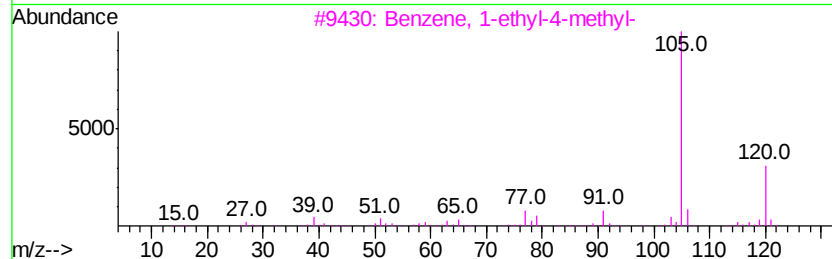
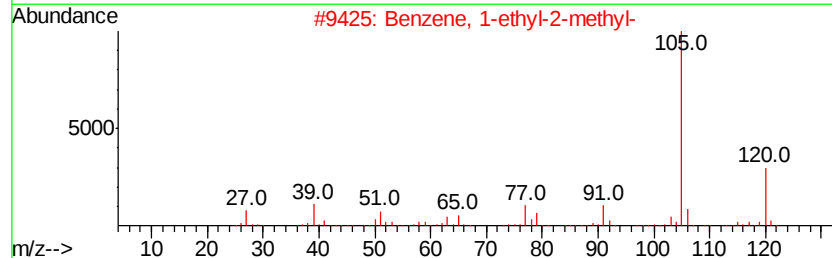
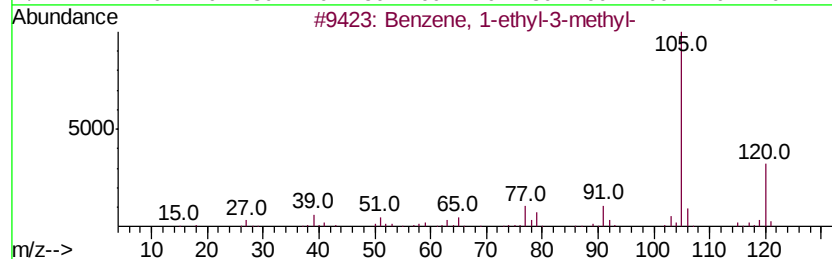
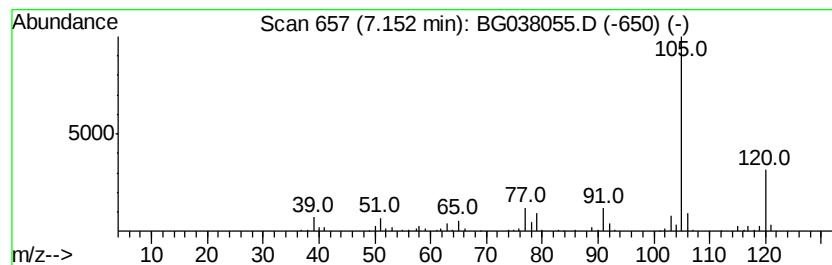
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG111318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzene, 1-ethyl-3-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.15	17.30 ng	249878	1,4-Dichlorobenzene-d4	8.09

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



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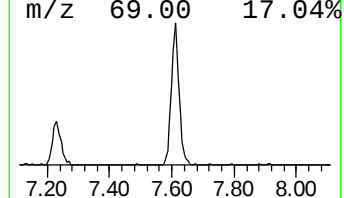
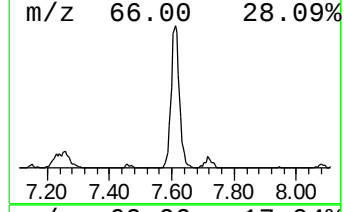
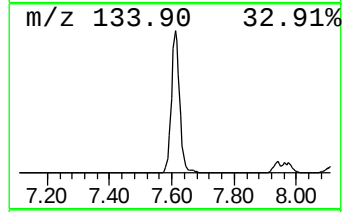
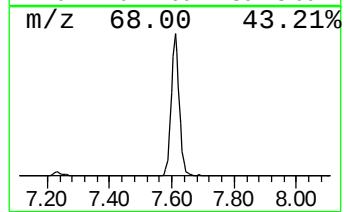
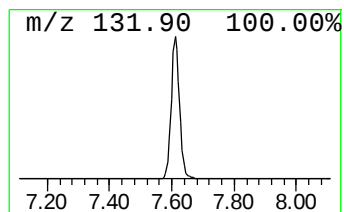
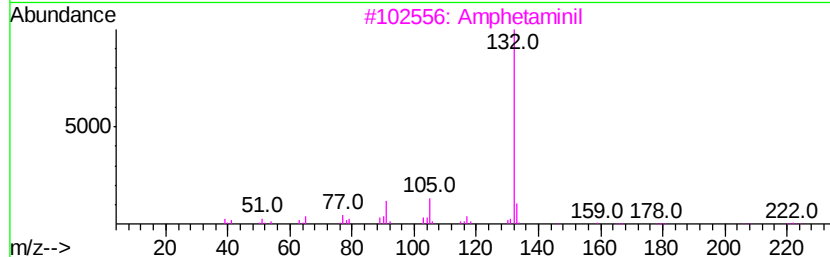
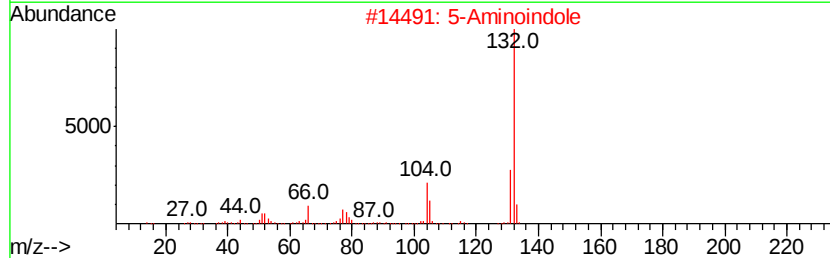
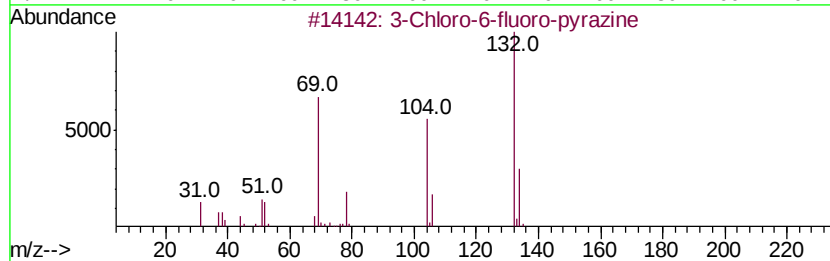
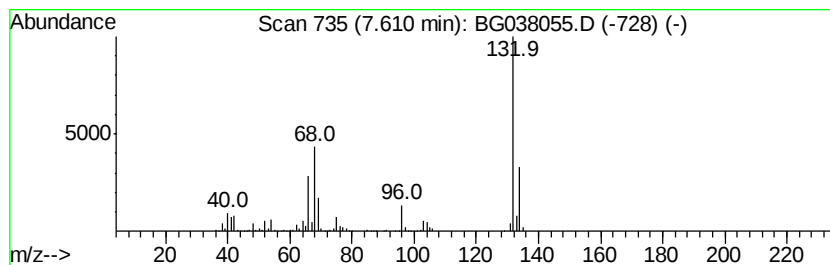
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG111318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 unknown7.61 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.61	38.66 ng	558530	1,4-Dichlorobenzene-d4	8.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Aminoindole	132	C8H8N2	005192-03-0	22
3		Amphetaminil	250	C17H18N2	017590-01-1	16
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
5		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12



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MW-2-111418

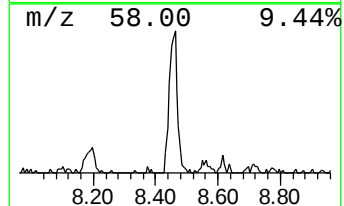
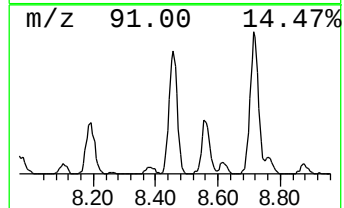
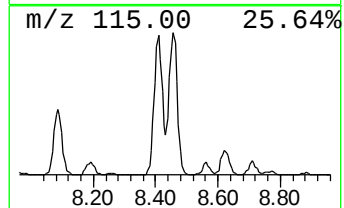
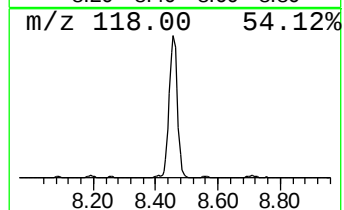
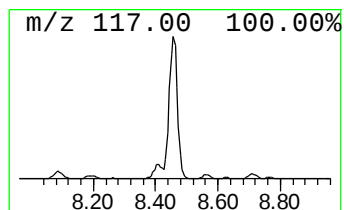
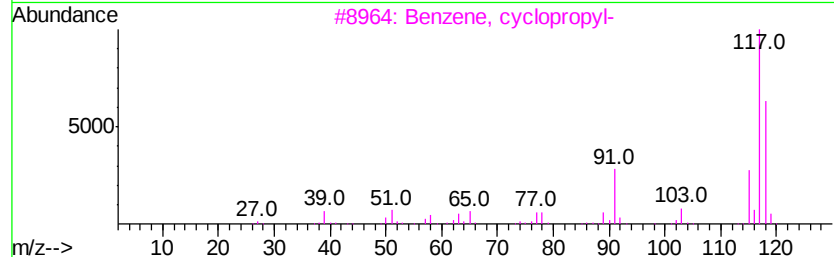
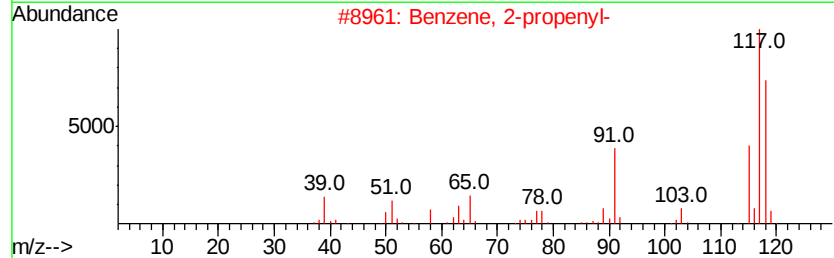
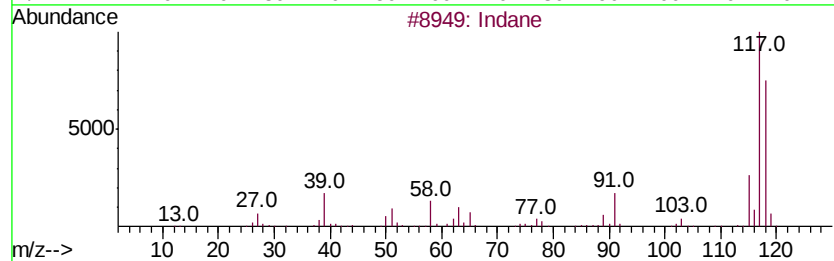
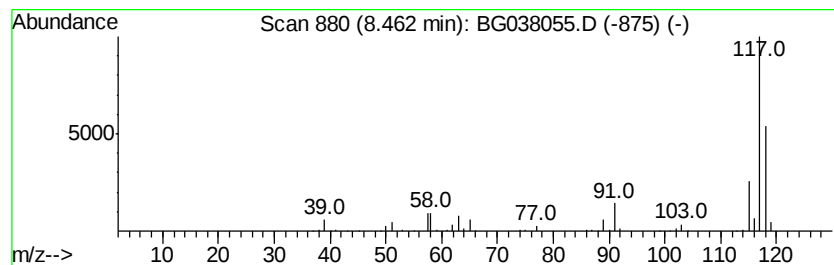
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Indane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.46	36.88 ng	532751	1,4-Dichlorobenzene-d4	8.09

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	91
2			Benzene, 2-propenyl-	118	C9H10	000300-57-2	74
3			Benzene, cyclopropyl-	118	C9H10	000873-49-4	72
4			Tetracyclo[3.3.1.0(2,8).0(4,6)]-	118	C9H10	1000191-13-7	72
5			Benzene, 1-propenyl-	118	C9H10	000637-50-3	55



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG111618\
Data File : BG038055.D
Acq On : 17 Nov 2018 5:09
Operator : JU/SJ
Sample : J5968-02
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
MW-2-111418

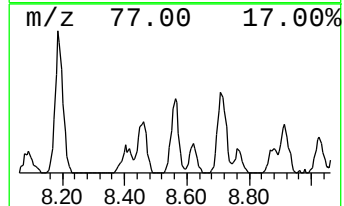
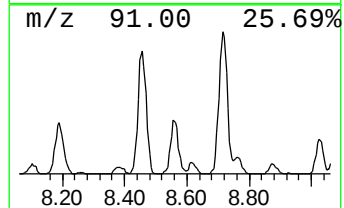
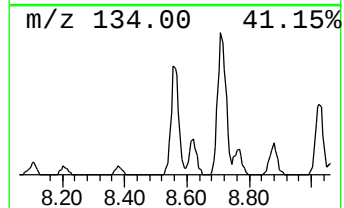
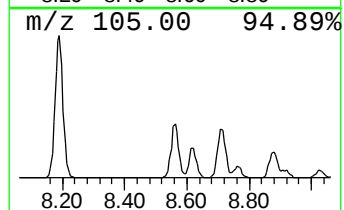
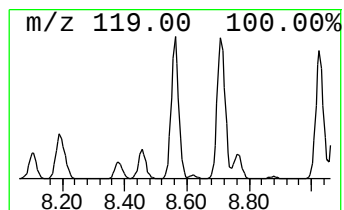
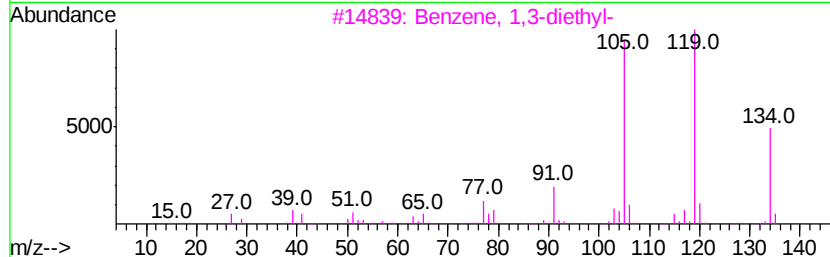
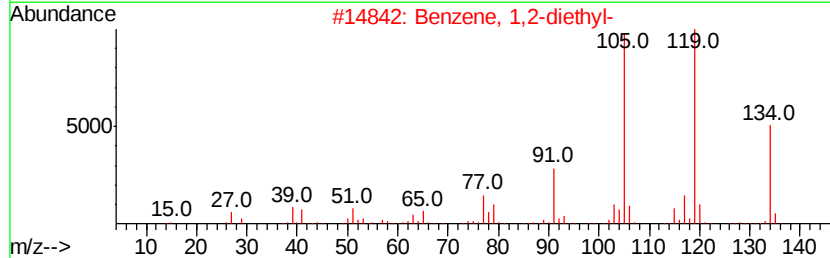
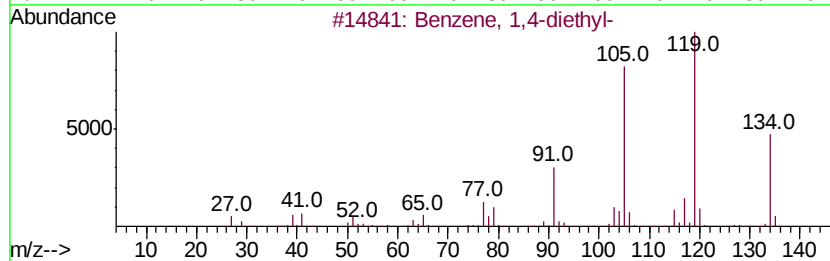
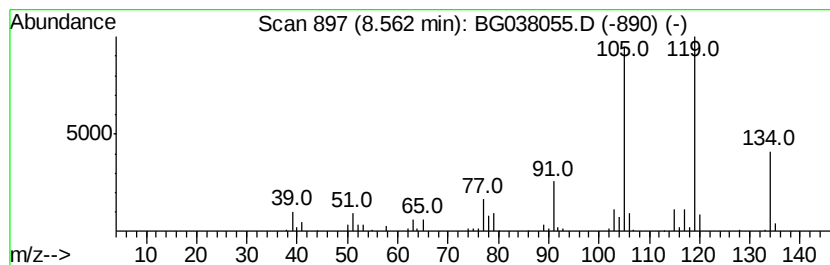
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Benzene, 1,4-diethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.56	13.34 ng	192757	1,4-Dichlorobenzene-d4	8.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	97
2		Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	96
3		Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	94
4		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	93
5		o-Cymene	134	C10H14	000527-84-4	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG111618\
Data File : BG038055.D
Acq On : 17 Nov 2018 5:09
Operator : JU/SJ
Sample : J5968-02
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
MW-2-111418

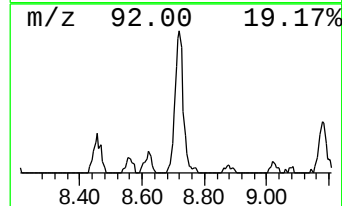
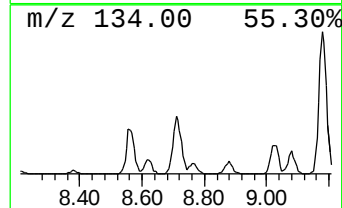
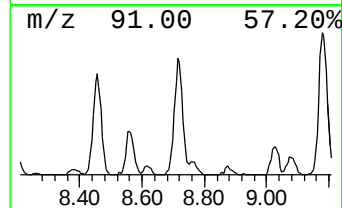
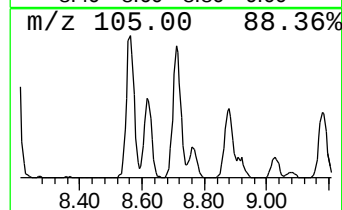
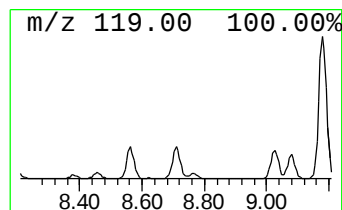
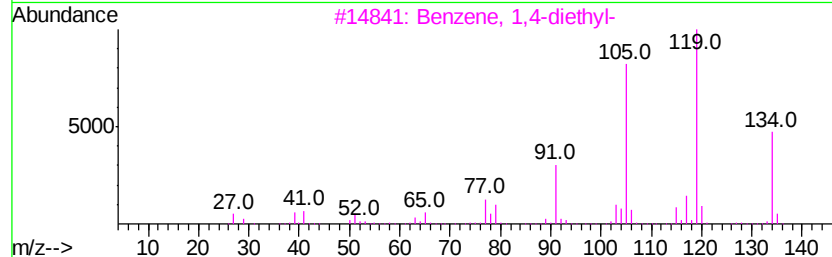
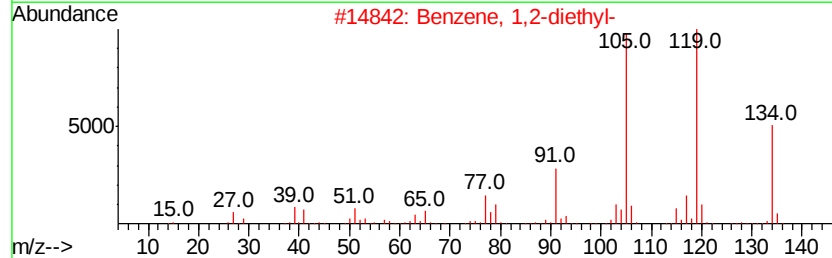
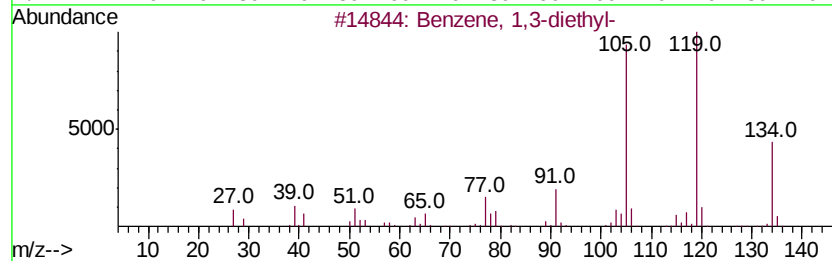
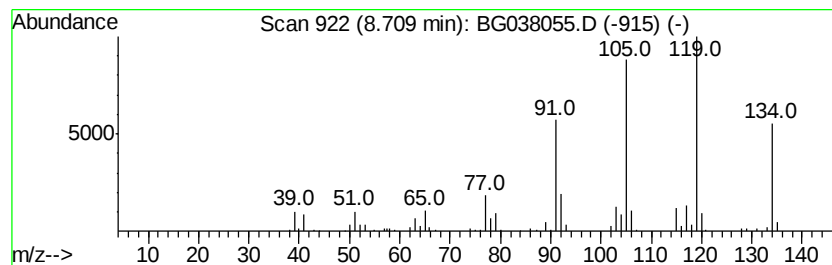
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 Benzene, 1,3-diethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.71	17.92 ng	258807	1,4-Dichlorobenzene-d4	8.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	96
2		Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	96
3		Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	95
4		1,3,8-p-Menthatriene	134	C10H14	018368-95-1	83
5		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	81



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG111618\
Data File : BG038055.D
Acq On : 17 Nov 2018 5:09
Operator : JU/SJ
Sample : J5968-02
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
MW-2-111418

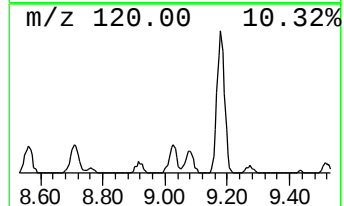
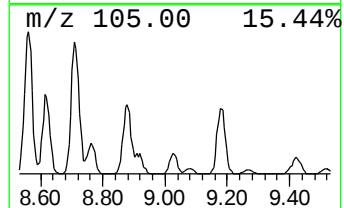
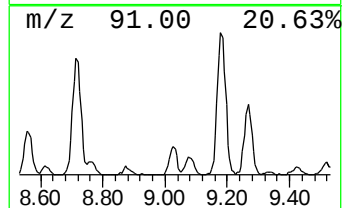
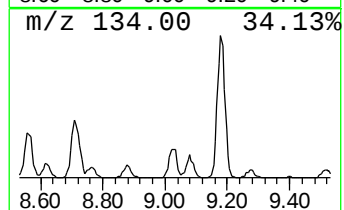
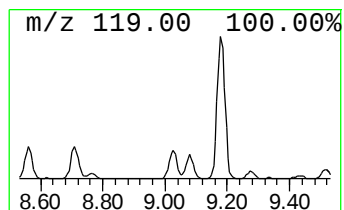
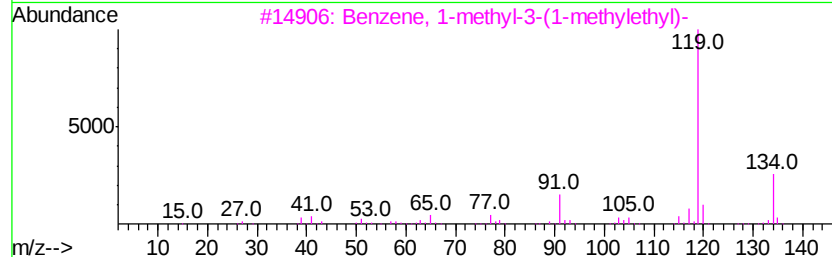
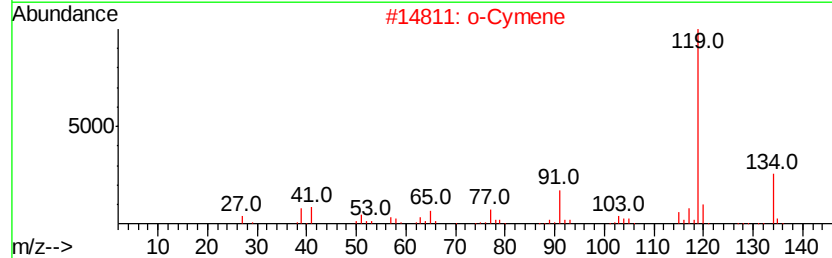
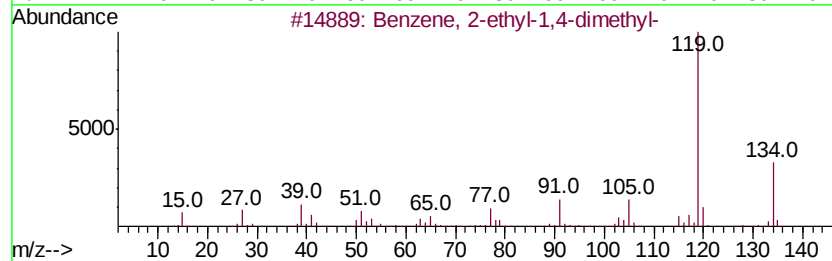
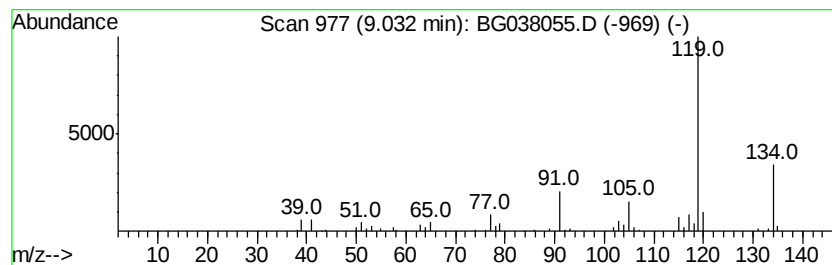
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.03	7.88 ng	113876	1,4-Dichlorobenzene-d4	8.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
2		o-Cymene	134	C10H14	000527-84-4	95
3		Benzene, 1-methyl-3-(1-methylethyl-)	134	C10H14	000535-77-3	94
4		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	94
5		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG111618\
Data File : BG038055.D
Acq On : 17 Nov 2018 5:09
Operator : JU/SJ
Sample : J5968-02
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
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ClientSampleId :
MW-2-111418

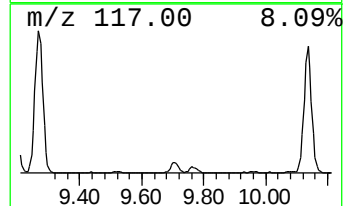
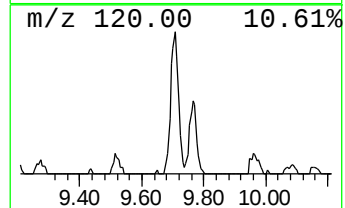
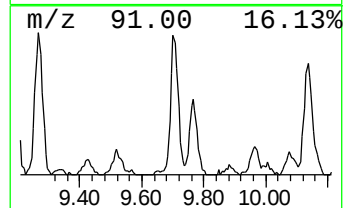
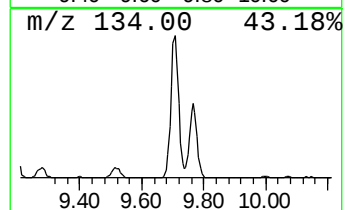
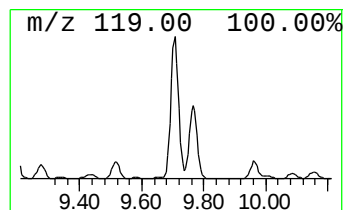
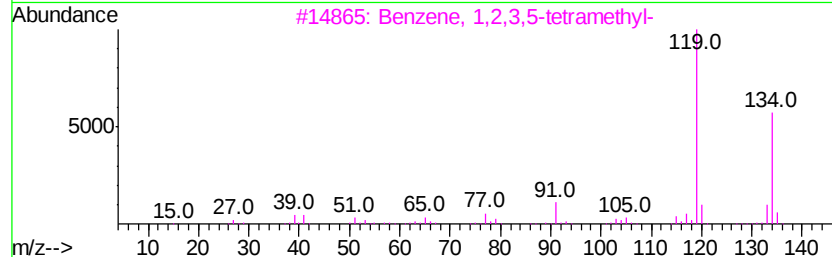
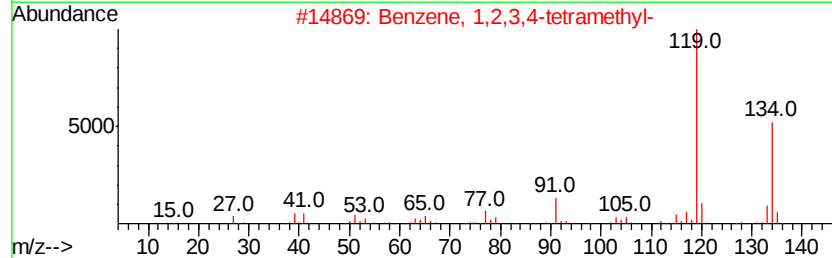
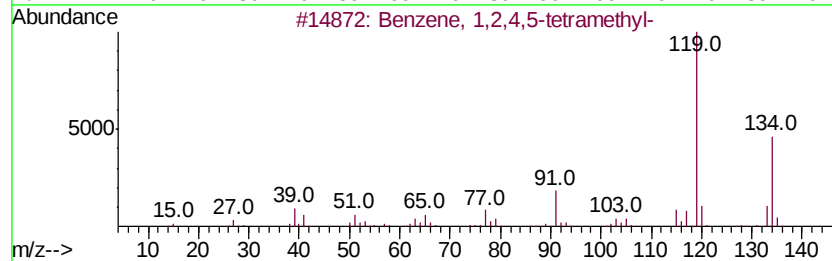
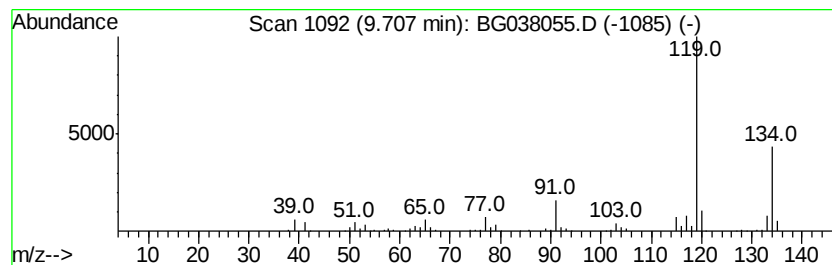
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 17 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.71	11.96 ng	289178	Naphthalene-d8	10.91

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
2		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
3		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
4		o-Cymene	134	C10H14	000527-84-4	95
5		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG111618\
Data File : BG038055.D
Acq On : 17 Nov 2018 5:09
Operator : JU/SJ
Sample : J5968-02
Misc :
ALS Vial : 25 Sample Multiplier: 1

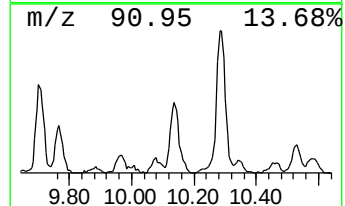
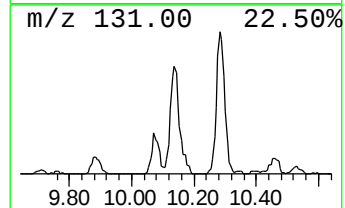
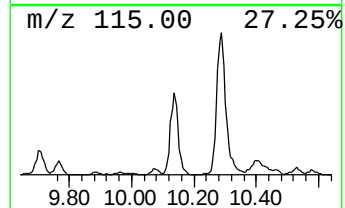
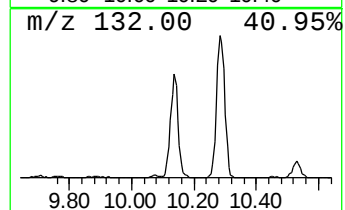
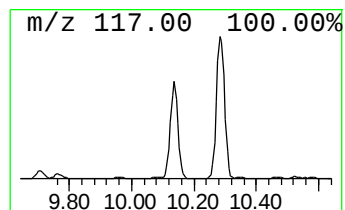
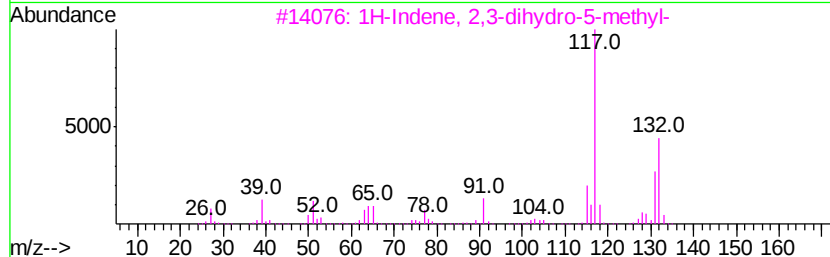
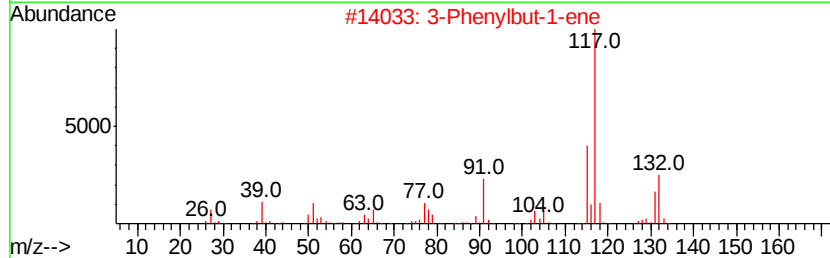
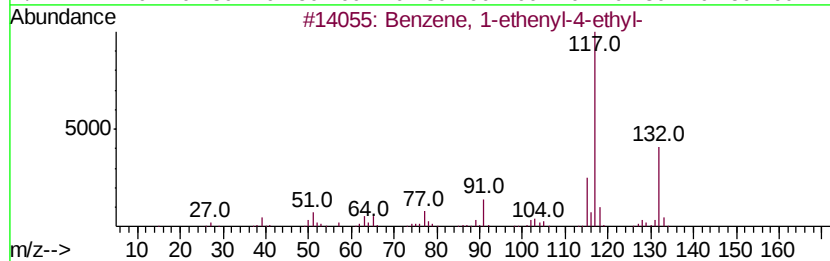
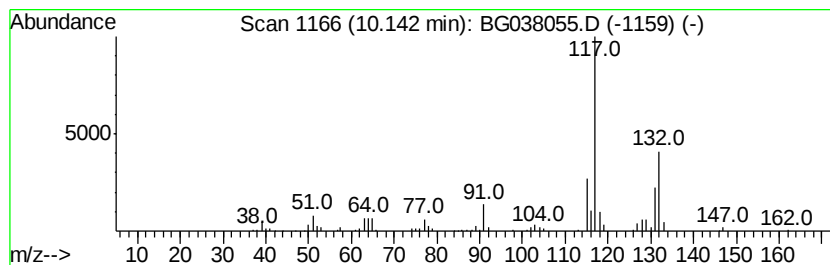
Instrument :
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ClientSampleId :
MW-2-111418

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG111318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 19 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.14	15.05 ng	363966	Napthalene-d8	10.91
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-ethenyl-4-ethyl-	132 C10H12	003454-07-7 93
2		3-Phenylbut-1-ene	132 C10H12	000934-10-1 93
3		1H-Indene, 2,3-dihydro-5-methyl-	132 C10H12	000874-35-1 92
4		Benzene, 2-ethenyl-1,4-dimethyl-	132 C10H12	002039-89-6 92
5		Benzene, 1-ethenyl-3-ethyl-	132 C10H12	007525-62-4 90



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG111618\
Data File : BG038055.D
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Misc :
ALS Vial : 25 Sample Multiplier: 1

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ClientSampleId :
MW-2-111418

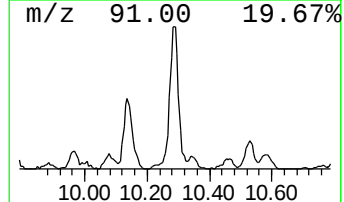
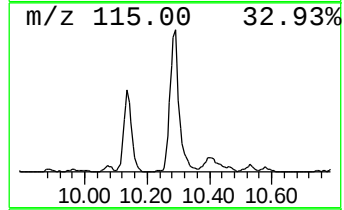
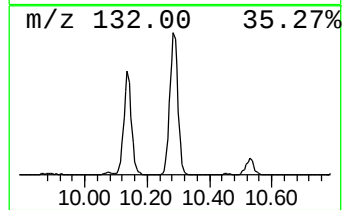
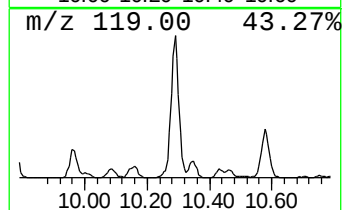
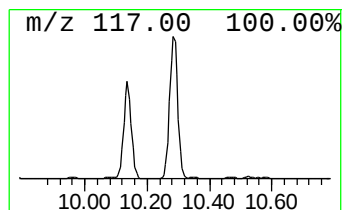
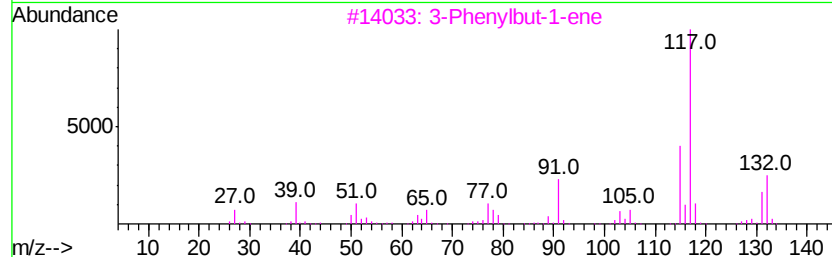
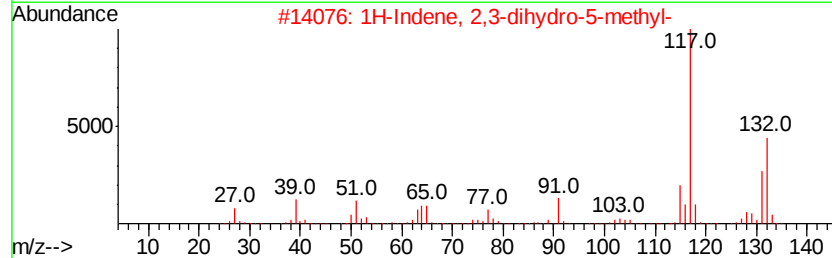
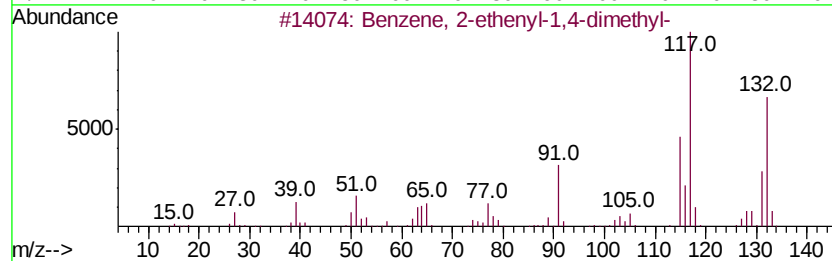
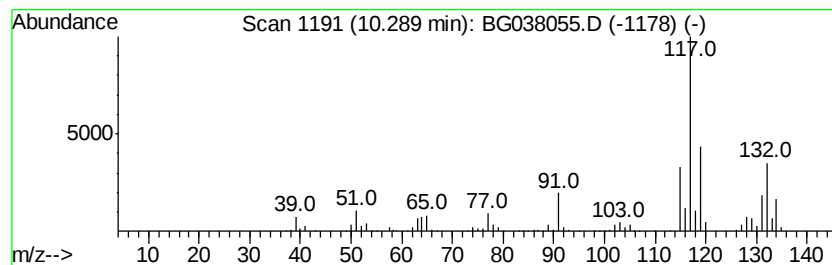
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 20 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.29	30.57 ng	739308	Napthalene-d8	10.91

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
2		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	92
3		3-Phenylbut-1-ene	132	C10H12	000934-10-1	76
4		1-Phenyl-1-butene	132	C10H12	000824-90-8	76
5		Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	76



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG111618\
Data File : BG038055.D
Acq On : 17 Nov 2018 5:09
Operator : JU/SJ
Sample : J5968-02
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
MW-2-111418

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG111318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, (1-methy...	6.55	12.1	ng	175011	1	8.09	288928	20.0
Benzene, 1-ethyl-...	7.15	17.3	ng	249878	1	8.09	288928	20.0
unknown7.61	7.61	38.7	ng	558530	1	8.09	288928	20.0
Indane	8.46	36.9	ng	532751	1	8.09	288928	20.0
Benzene, 1,4-diet...	8.56	13.3	ng	192757	1	8.09	288928	20.0
Benzene, 1,3-diet...	8.71	17.9	ng	258807	1	8.09	288928	20.0
Benzene, 2-ethyl-...	9.03	7.9	ng	113876	1	8.09	288928	20.0
Benzene, 1,2,4,5-...	9.71	12.0	ng	289178	2	10.91	483758	20.0
Benzene, 1-etheny...	10.14	15.1	ng	363966	2	10.91	483758	20.0
Benzene, 2-etheny...	10.29	30.6	ng	739308	2	10.91	483758	20.0