

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012418\
 Data File : BF102390.D
 Acq On : 24 Jan 2018 15:35
 Operator : SJ/JU
 Sample : J1235-07 5X
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 30-2

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_F\METHODS\8270-BF012218.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	2.699	99	104	111	rVB3	12830	21357	1.47%	0.126%
2	3.657	260	267	274	rBV	33445	48061	3.30%	0.284%
3	4.887	472	476	481	rBV	56934	63435	4.35%	0.375%
4	5.093	506	511	514	rBV2	21034	25276	1.73%	0.149%
5	5.163	520	523	527	rVB	19004	18299	1.26%	0.108%
6	5.275	538	542	545	rBV	86146	101038	6.93%	0.597%
7	5.310	545	548	557	rVB	678359	665938	45.69%	3.933%
8	5.463	570	574	578	rVB4	12090	16290	1.12%	0.096%
9	5.540	584	587	591	rBV2	22791	26339	1.81%	0.156%
10	5.716	612	617	621	rBV3	17920	22579	1.55%	0.133%
11	5.869	637	643	646	rBV	115151	107288	7.36%	0.634%
12	5.951	654	657	660	rBV2	19783	19910	1.37%	0.118%
13	6.163	690	693	698	rVB	195228	156775	10.76%	0.926%
14	6.234	702	705	708	rBV	40809	41562	2.85%	0.245%
15	6.263	708	710	714	rVB	18456	14619	1.00%	0.086%
16	6.310	714	718	721	rBV	32928	32751	2.25%	0.193%
17	6.345	721	724	730	rVV	728440	703310	48.25%	4.154%
18	6.392	730	732	737	rVB	27852	28431	1.95%	0.168%
19	6.475	743	746	752	rBV	779680	738569	50.67%	4.362%
20	6.534	752	756	760	rBV	177312	154441	10.60%	0.912%
21	6.710	782	786	792	rVB	930708	768579	52.73%	4.539%
22	6.763	792	795	798	rVB	41742	38054	2.61%	0.225%
23	6.863	808	812	815	rBV	637541	567462	38.93%	3.352%
24	6.887	815	816	820	rVB	89755	59699	4.10%	0.353%
25	6.957	825	828	831	rBV	208841	160609	11.02%	0.949%
26	7.028	836	840	843	rBV	90783	88049	6.04%	0.520%
27	7.104	850	853	855	rBV3	21153	22304	1.53%	0.132%
28	7.245	872	877	879	rBV	329561	267013	18.32%	1.577%
29	7.275	879	882	886	rVB	492337	414543	28.44%	2.448%
30	7.351	891	895	898	rVB4	19414	24899	1.71%	0.147%
31	7.481	911	917	920	rBV	134918	122494	8.40%	0.723%
32	7.510	920	922	925	rVB	48424	38655	2.65%	0.228%
33	7.598	931	937	941	rBV4	25848	42491	2.92%	0.251%
34	7.639	941	944	946	rVV3	19340	30927	2.12%	0.183%

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35	7.663	946	948	953	rVV4	36893	43143	2.96%	0.255%
36	7.734	956	960	963	rVV2	118205	125290	8.60%	0.740%
37	7.763	963	965	968	rVV	23957	20691	1.42%	0.122%
38	7.816	968	974	976	rVV4	19773	29152	2.00%	0.172%
39	7.863	978	982	985	rVB	36680	30300	2.08%	0.179%
40	7.992	998	1004	1010	rBV	1172349	1109690	76.13%	6.554%
41	8.081	1017	1019	1022	rVB	28933	19701	1.35%	0.116%
42	8.751	1127	1133	1134	rBV	36819	35432	2.43%	0.209%
43	8.910	1156	1160	1168	rVB9	9686	22161	1.52%	0.131%
44	9.069	1180	1187	1190	rBV	920736	744434	51.07%	4.397%
45	9.463	1249	1254	1257	rBV	93513	87033	5.97%	0.514%
46	9.551	1265	1269	1276	rVB3	39521	68094	4.67%	0.402%
47	9.604	1276	1278	1282	rBV	14643	15942	1.09%	0.094%
48	9.675	1288	1290	1297	rBV3	28565	37773	2.59%	0.223%
49	9.745	1297	1302	1310	rVB2	1323645	1187795	81.49%	7.015%
50	9.845	1315	1319	1323	rVB2	20542	29001	1.99%	0.171%
51	9.951	1332	1337	1342	rBV5	29645	46645	3.20%	0.275%
52	10.045	1349	1353	1357	rBV	31287	64923	4.45%	0.383%
53	10.163	1368	1373	1378	rVB	169479	255589	17.54%	1.510%
54	10.445	1419	1421	1425	rVB4	16465	20173	1.38%	0.119%
55	10.533	1432	1436	1443	rVB	592844	517219	35.49%	3.055%
56	10.645	1453	1455	1460	rBV2	26706	31622	2.17%	0.187%
57	10.739	1468	1471	1476	rBV6	14103	22568	1.55%	0.133%
58	10.898	1494	1498	1502	rBV6	26391	34114	2.34%	0.201%
59	11.075	1525	1528	1530	rBV2	16710	23788	1.63%	0.140%
60	11.127	1535	1537	1541	rVB4	22934	20311	1.39%	0.120%
61	11.233	1551	1555	1557	rBV	1261781	1143956	78.49%	6.756%
62	11.251	1557	1558	1561	rVB	142553	100754	6.91%	0.595%
63	11.616	1618	1620	1627	rVB4	25159	31642	2.17%	0.187%
64	11.733	1638	1640	1641	rBV2	17726	15525	1.07%	0.092%
65	11.763	1642	1645	1647	rVV2	101495	100461	6.89%	0.593%
66	11.792	1647	1650	1653	rVV	65851	74393	5.10%	0.439%
67	11.845	1656	1659	1661	rVV2	25831	27075	1.86%	0.160%
68	12.445	1758	1761	1764	rVB	249151	189571	13.01%	1.120%
69	12.545	1776	1778	1785	rVB7	39214	37913	2.60%	0.224%
70	12.669	1796	1799	1805	rVB	191816	190755	13.09%	1.127%
71	12.816	1820	1824	1831	rVB	846178	795095	54.55%	4.696%

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 Start Thrs: 0.2 Max Peaks: 100
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72	13.357	1912	1916	1918	rBV4	19006	33508	2.30%	0.198%
73	13.710	1973	1976	1980	rBV3	77565	86707	5.95%	0.512%
74	13.868	1995	2003	2006	rBV	1269202	1457543	100.00%	8.609%
75	14.033	2026	2031	2035	rBV6	28187	48472	3.33%	0.286%
76	14.263	2067	2070	2071	rBV2	26600	30747	2.11%	0.182%
77	14.492	2105	2109	2116	rBV3	74638	137505	9.43%	0.812%
78	14.633	2131	2133	2136	rVB2	31944	30591	2.10%	0.181%
79	14.880	2171	2175	2179	rBV	117024	189661	13.01%	1.120%
80	14.951	2184	2187	2190	rVV2	36032	42750	2.93%	0.252%
81	14.992	2190	2194	2196	rVV3	17430	23621	1.62%	0.140%
82	15.021	2196	2199	2203	rVV6	20760	24233	1.66%	0.143%
83	15.168	2220	2224	2231	rVB	94168	124673	8.55%	0.736%
84	15.227	2231	2234	2239	rBV	74425	89229	6.12%	0.527%
85	15.292	2240	2245	2254	rVB	812152	979365	67.19%	5.784%
86	15.480	2272	2277	2284	rBV9	62821	149427	10.25%	0.883%
87	16.180	2390	2396	2403	rVB7	35626	84624	5.81%	0.500%
88	16.668	2475	2479	2486	rVB3	46998	96238	6.60%	0.568%
89	17.086	2547	2550	2557	rVB2	49745	89924	6.17%	0.531%
90	18.456	2773	2783	2794	rVB8	60237	210848	14.47%	1.245%

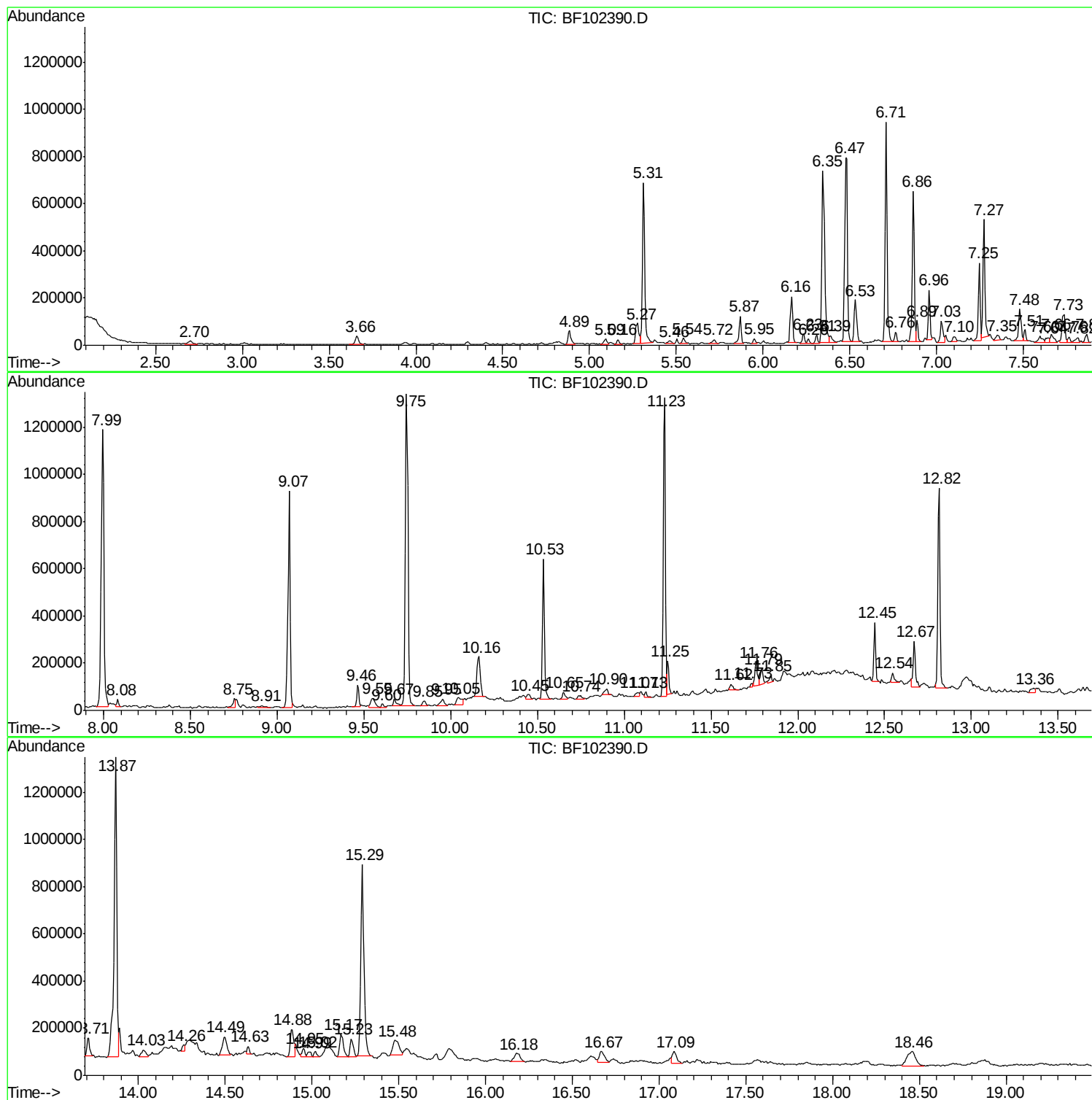
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ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampled :
30-2

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.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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Instrument :
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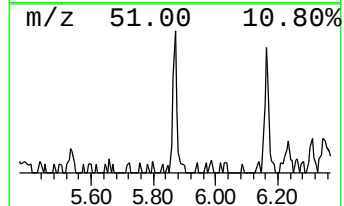
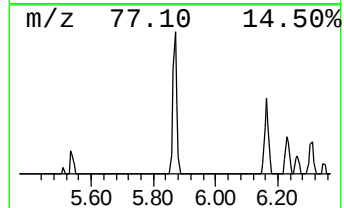
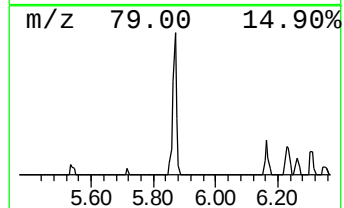
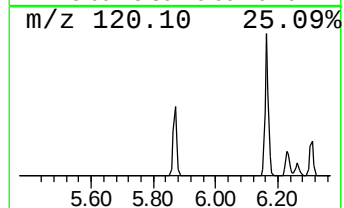
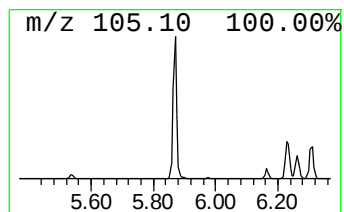
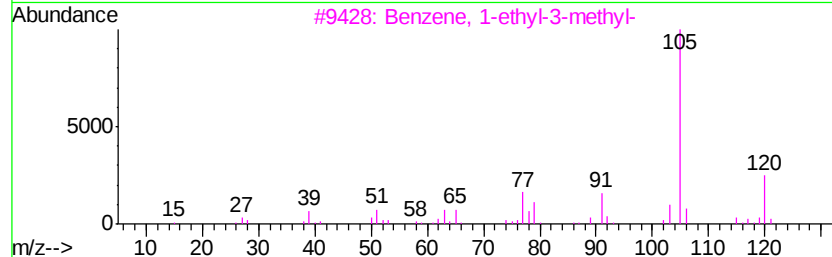
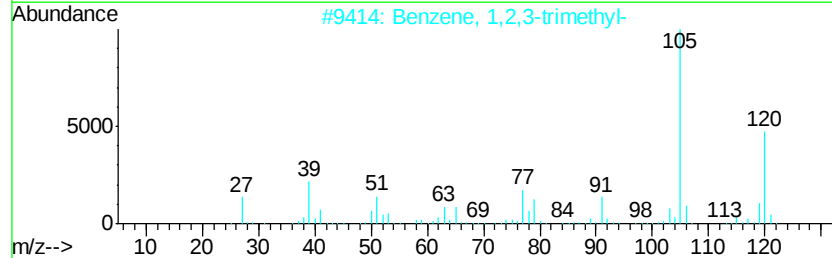
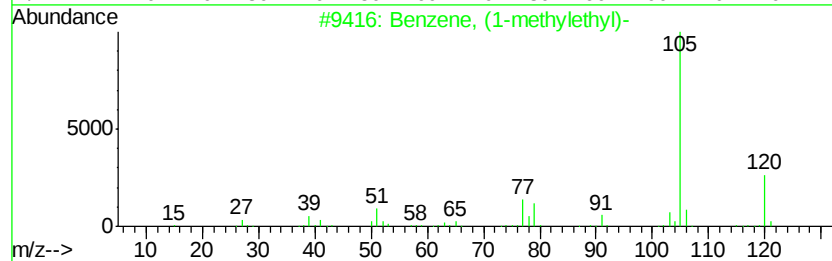
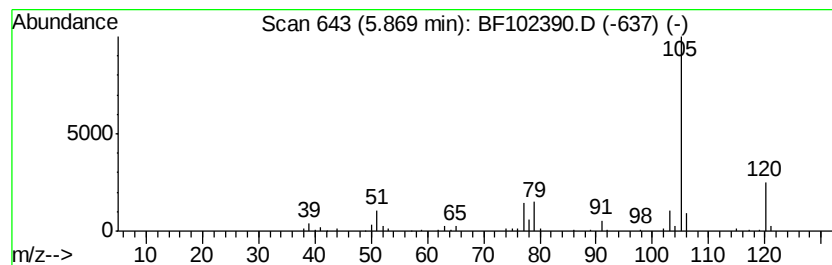
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Benzene, (1-methylethyl)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.87	2.79 ng	107288	1,4-Dichlorobenzene-d4	6.71

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	83
4		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	80
5		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	80



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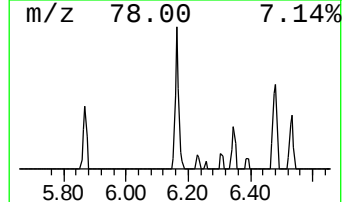
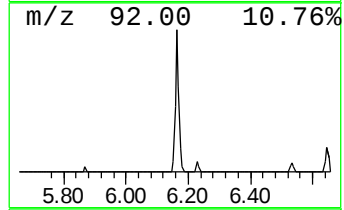
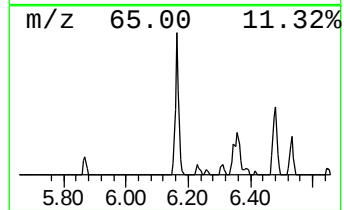
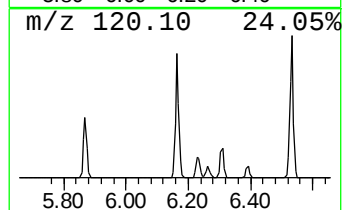
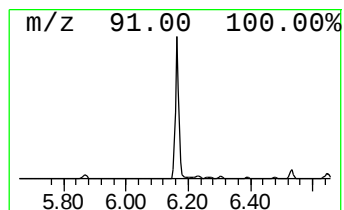
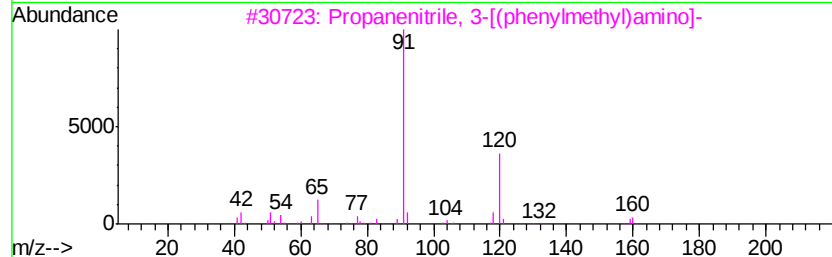
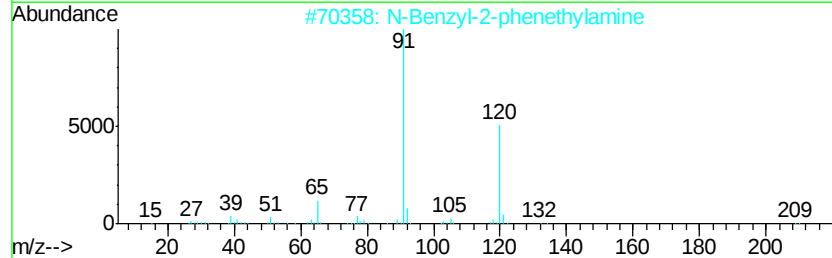
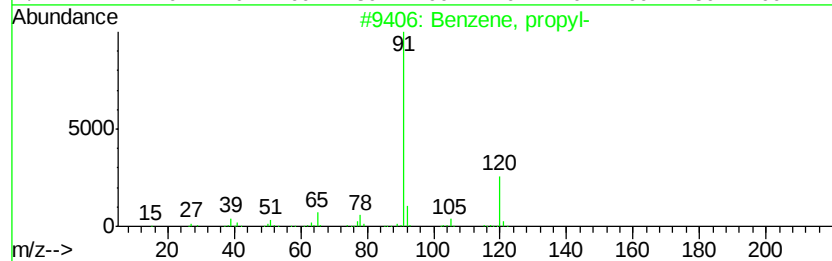
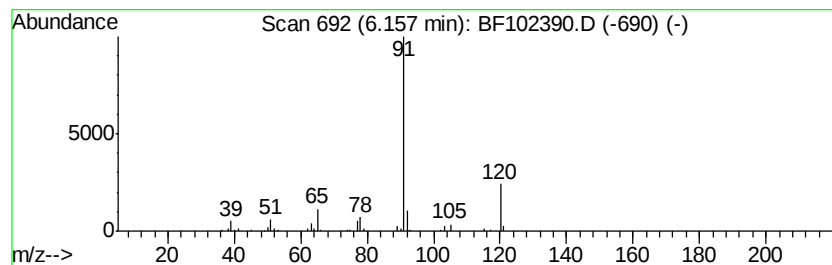
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Benzene, propyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.16	4.08 ng	156775	1,4-Dichlorobenzene-d4	6.71

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, propyl-	120	C9H12	000103-65-1	91
2		N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	64
3		Propanenitrile, 3-[(phenylmethyl)...	160	C10H12N2	000706-03-6	59
4		Ethanol, 2-[(phenylmethyl)amino]-	151	C9H13NO	000104-63-2	56
5		Hydroxylamine, O-(phenylmethyl)-	123	C7H9NO	000622-33-3	47



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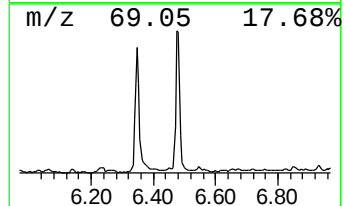
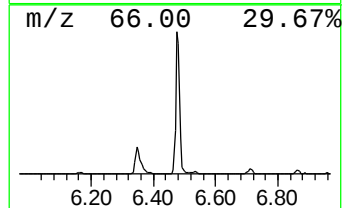
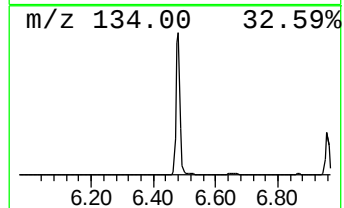
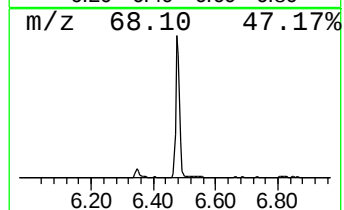
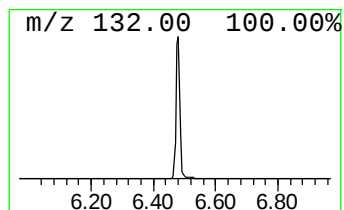
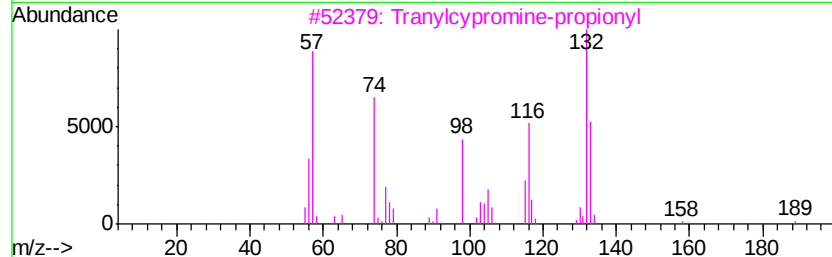
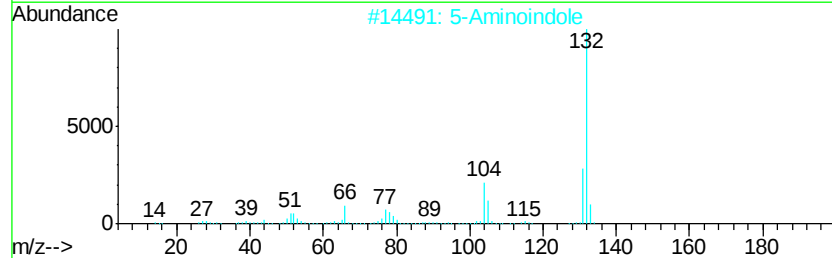
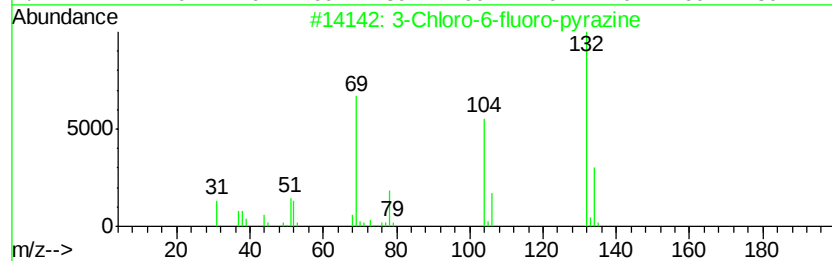
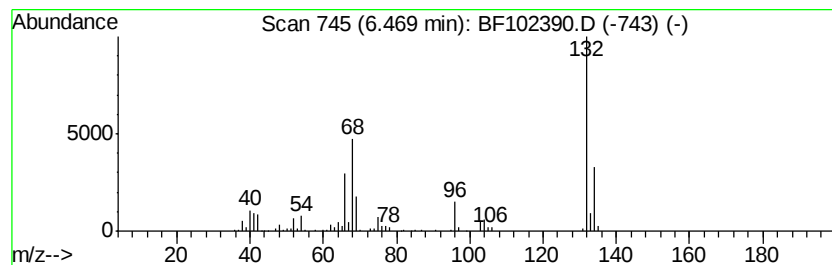
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown6.47 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.47	19.22 ng	738569	1,4-Dichlorobenzene-d4	6.71

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	12
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
5		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



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30-2

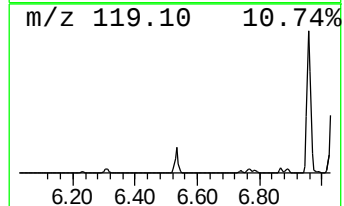
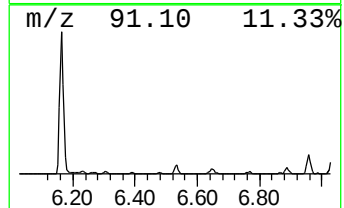
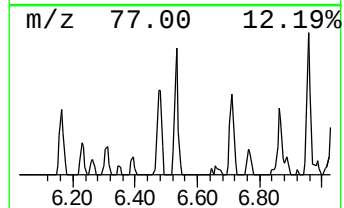
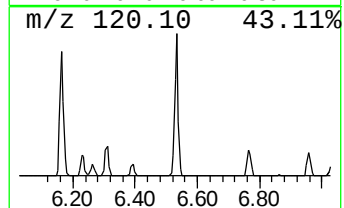
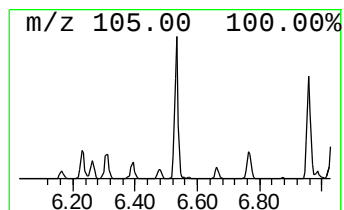
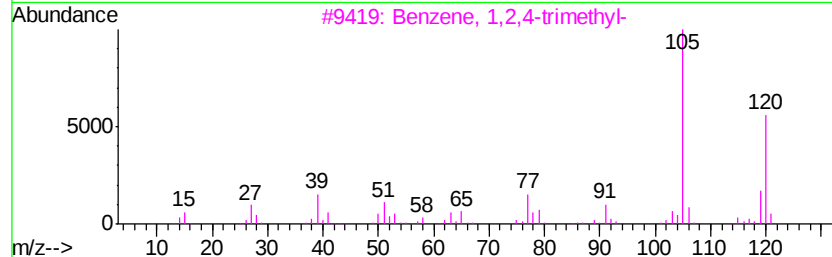
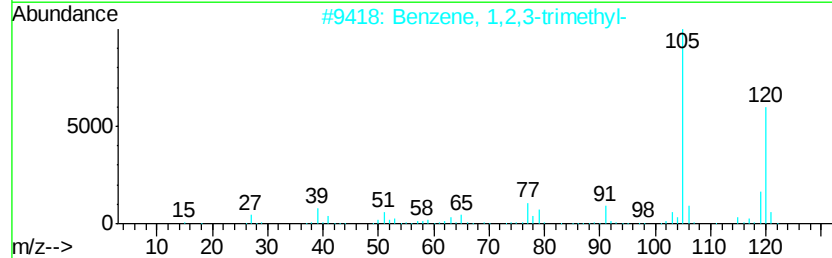
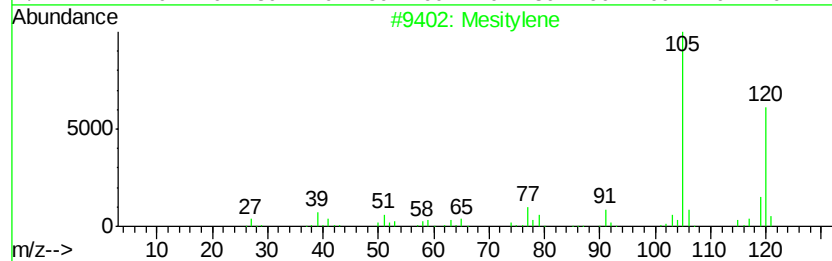
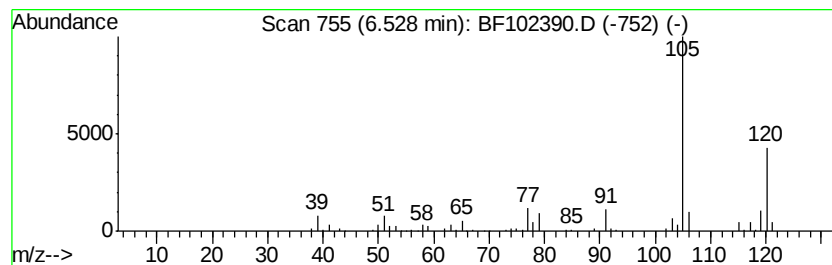
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Mesitylene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.53	4.02 ng	154441	1,4-Dichlorobenzene-d4	6.71

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Mesitylene	120	C9H12	000108-67-8	95
2	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
3	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90
5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012418\
Data File : BF102390.D
Acq On : 24 Jan 2018 15:35
Operator : SJ/JU
Sample : J1235-07 5X
Misc :
ALS Vial : 3 Sample Multiplier: 1

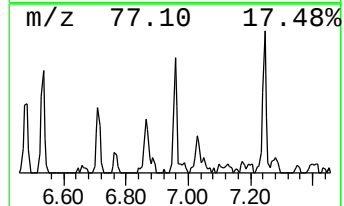
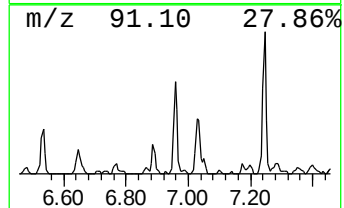
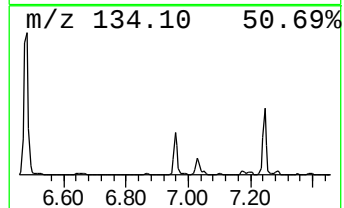
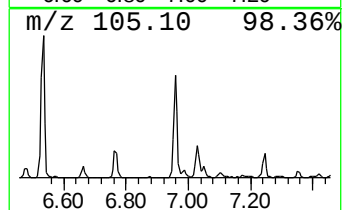
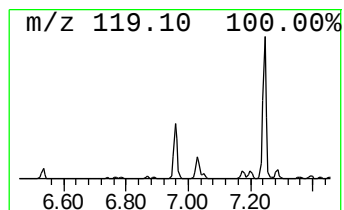
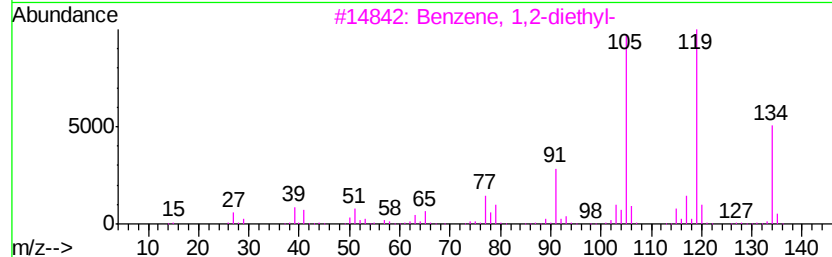
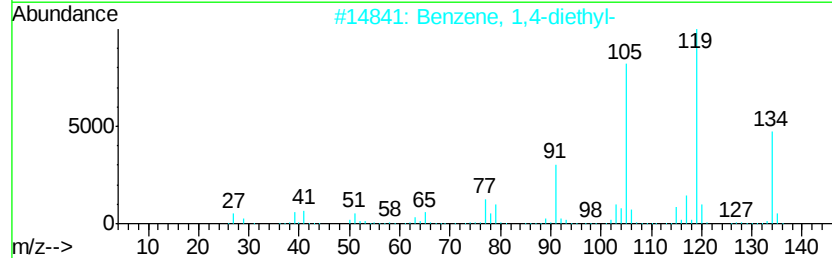
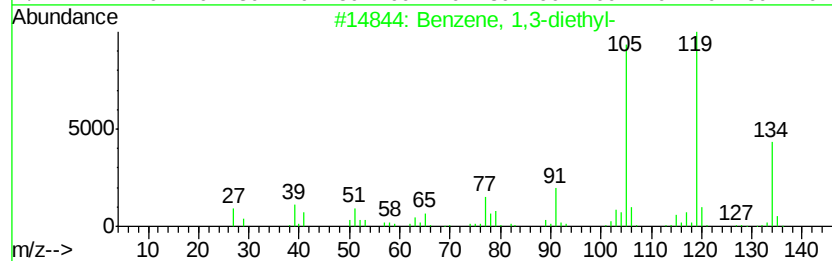
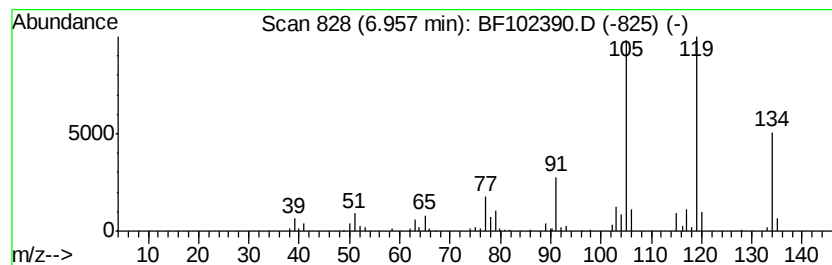
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Benzene, 1,3-diethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
6.96	4.18 ng	160609	1,4-Dichlorobenzene-d4	6.71	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	96
2	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	95
3	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	95
4	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91
5	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012418\
Data File : BF102390.D
Acq On : 24 Jan 2018 15:35
Operator : SJ/JU
Sample : J1235-07 5X
Misc :
ALS Vial : 3 Sample Multiplier: 1

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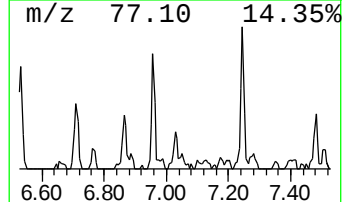
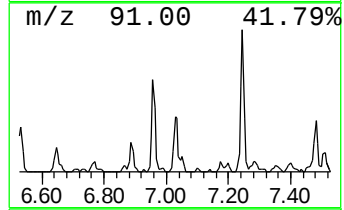
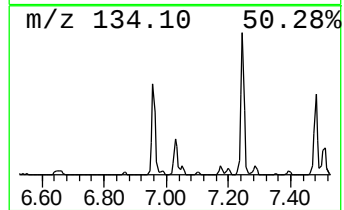
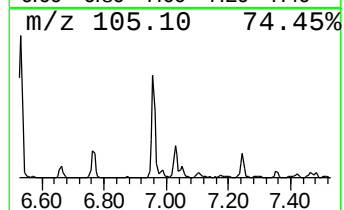
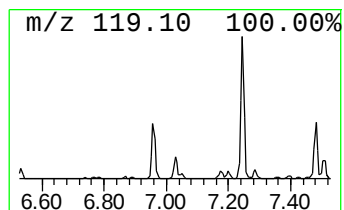
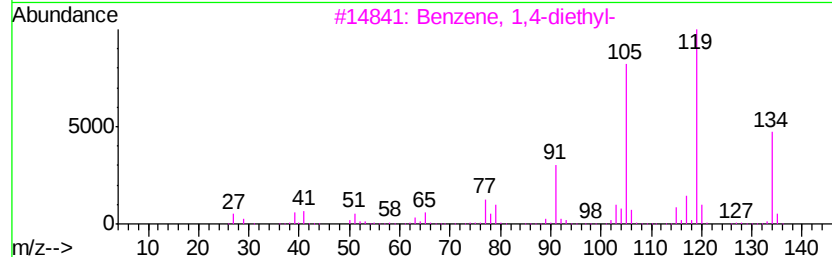
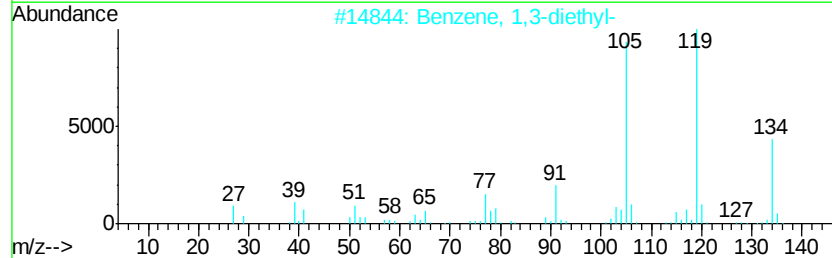
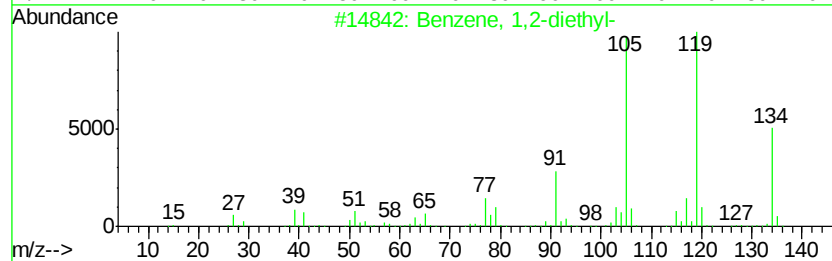
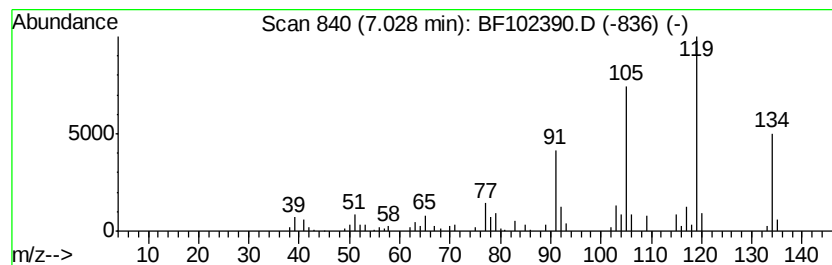
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Benzene, 1,2-diethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.03	2.29 ng	88049	1,4-Dichlorobenzene-d4	6.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	96
2		Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	96
3		Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	96
4		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91
5		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012418\
Data File : BF102390.D
Acq On : 24 Jan 2018 15:35
Operator : SJ/JU
Sample : J1235-07 5X
Misc :
ALS Vial : 3 Sample Multiplier: 1

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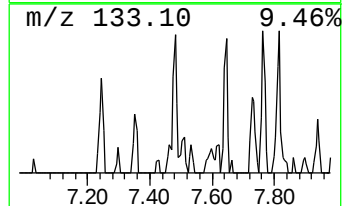
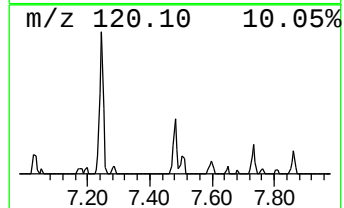
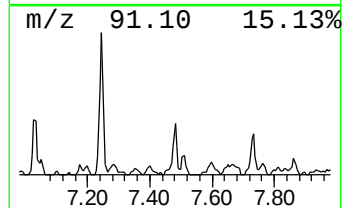
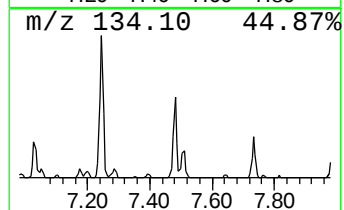
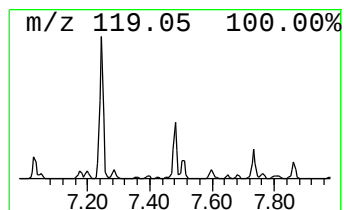
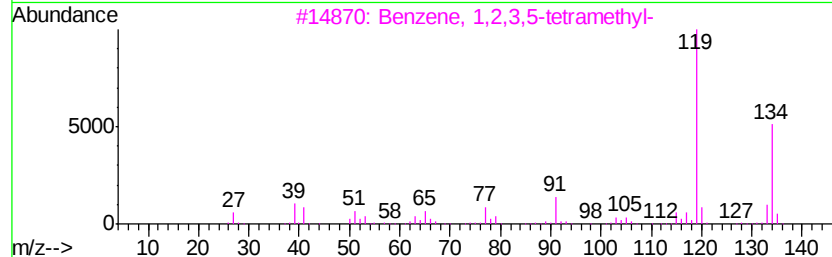
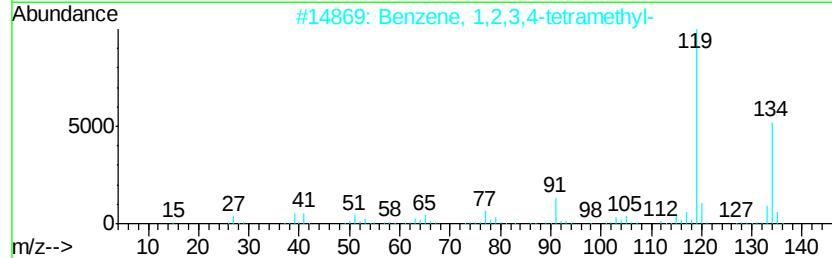
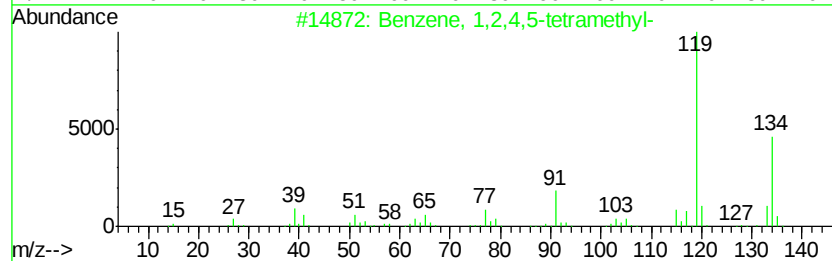
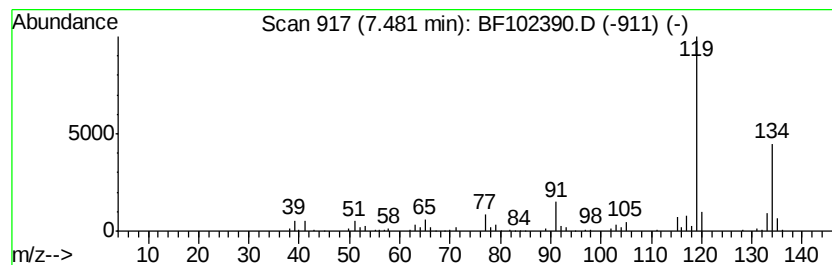
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.48	2.21 ng	122494	Naphthalene-d8	7.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
3		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
4		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	93
5		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	93



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012418\
Data File : BF102390.D
Acq On : 24 Jan 2018 15:35
Operator : SJ/JU
Sample : J1235-07 5X
Misc :
ALS Vial : 3 Sample Multiplier: 1

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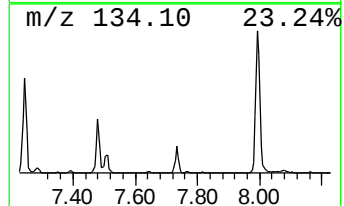
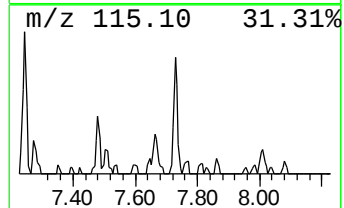
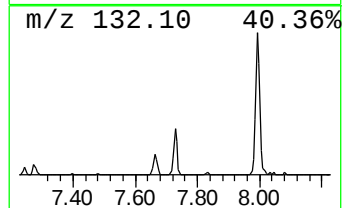
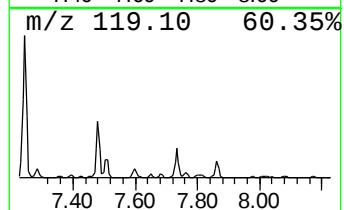
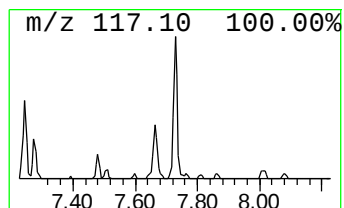
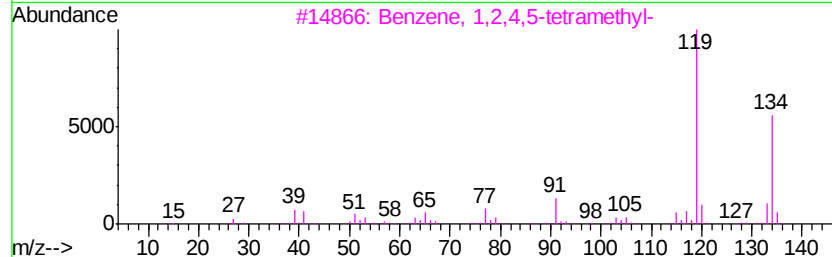
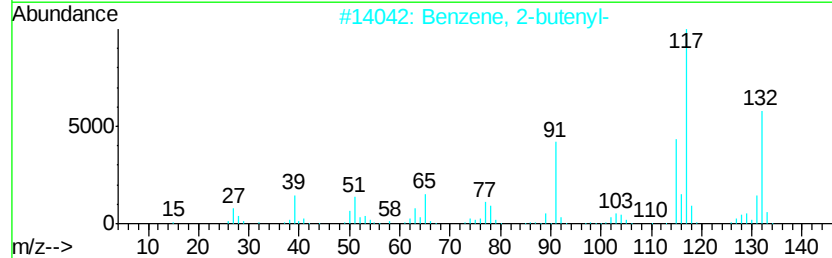
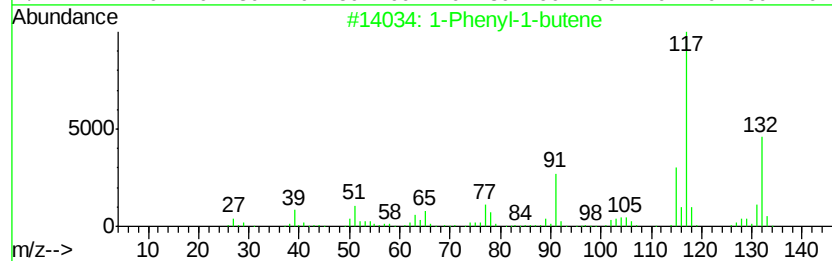
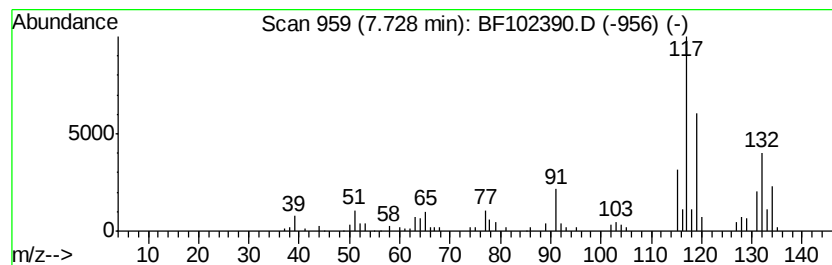
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 1-Phenyl-1-butene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.73	2.26 ng	125290	Naphthalene-d8	7.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Phenyl-1-butene	132	C10H12	000824-90-8	83
2		Benzene, 2-butenyl-	132	C10H12	001560-06-1	80
3		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	64
4		o-Cymene	134	C10H14	000527-84-4	60
5		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	58



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012418\
Data File : BF102390.D
Acq On : 24 Jan 2018 15:35
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Instrument :
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ClientSampleId :
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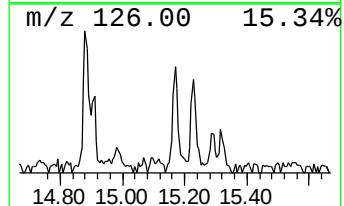
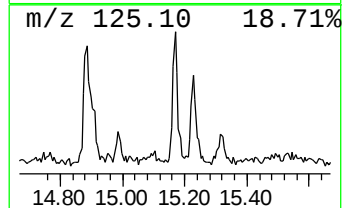
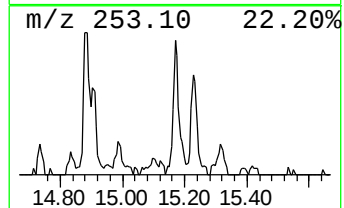
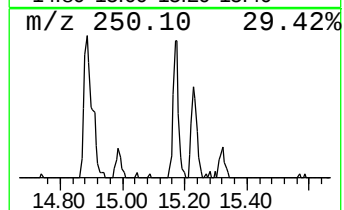
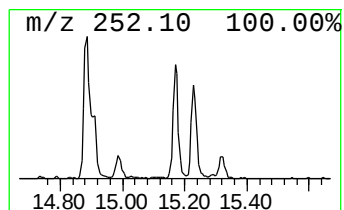
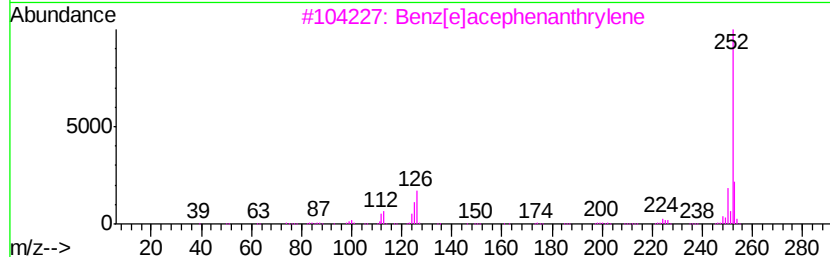
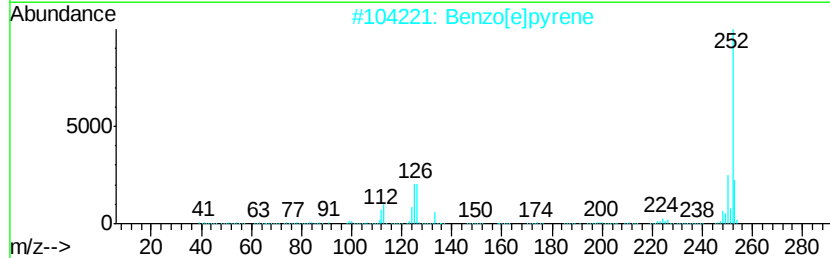
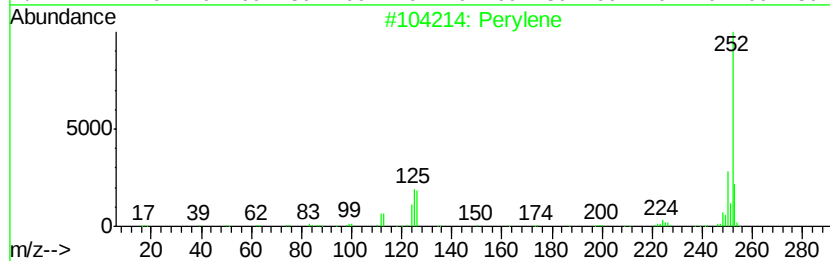
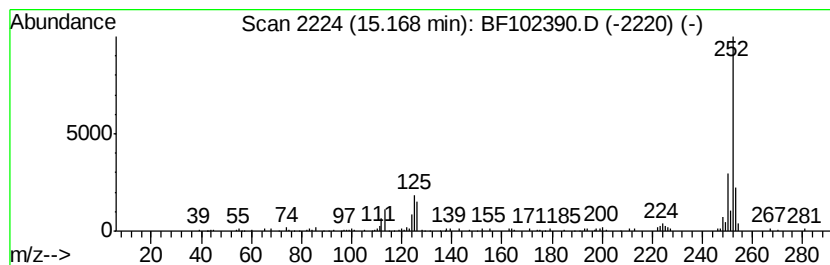
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Perylene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.17	2.55 ng	124673	Perylene-d12	15.29

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Perylene		252	C20H12	000198-55-0	98
2	Benzo[e]pyrene		252	C20H12	000192-97-2	97
3	Benz[e]acephenanthrylene		252	C20H12	000205-99-2	96
4	Benzo[k]fluoranthene		252	C20H12	000207-08-9	93
5	Benzo[a]pyrene		252	C20H12	000050-32-8	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012418\
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Misc :
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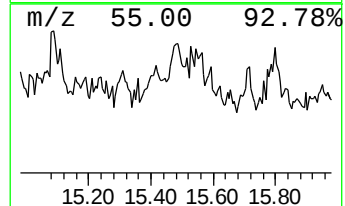
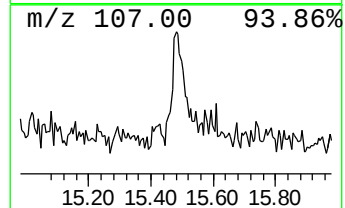
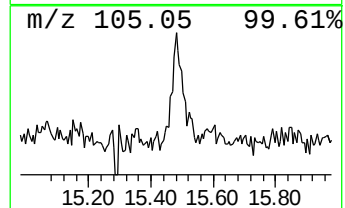
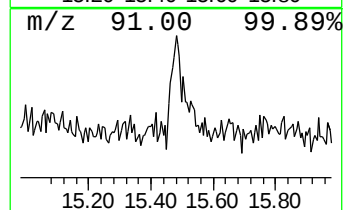
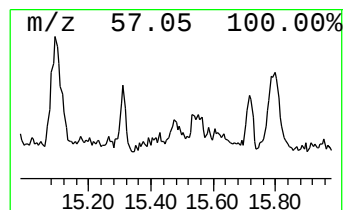
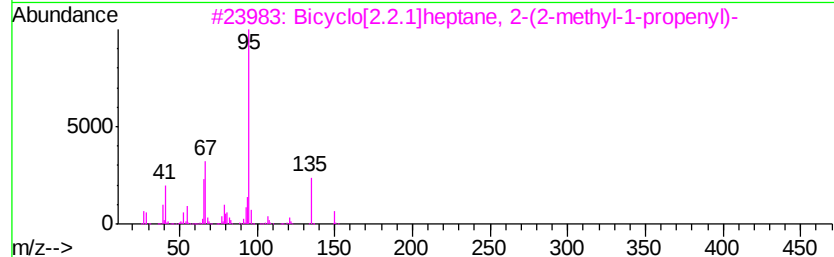
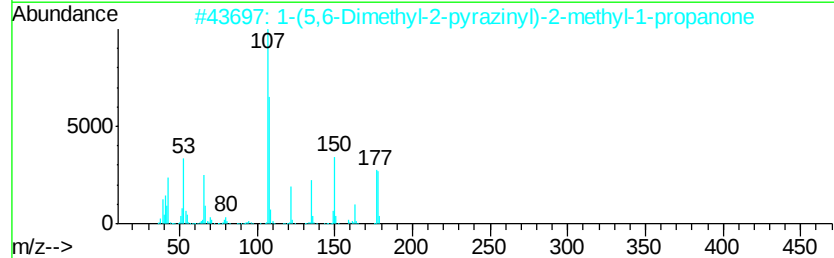
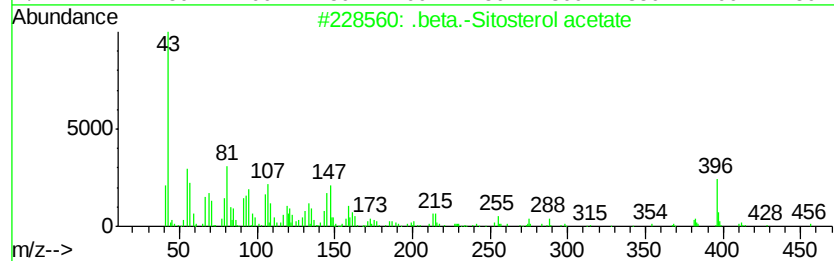
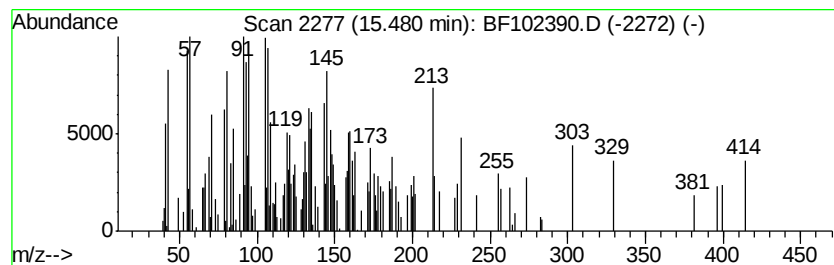
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 unknown15.48 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.48	3.05 ng	149427	Perylene-d12	15.29
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		.beta.-Sitosterol acetate	456 C31H52O2	000915-05-9 35
2		1-(5,6-Dimethyl-2-pyrazinyl)-2-m...	178 C10H14N2O	1000109-57-7 18
3		Bicyclo[2.2.1]heptane, 2-(2-meth...	150 C11H18	061142-27-6 14
4		Cholest-5-ene, 3-ethoxy-, (3.bet...	414 C29H50O	000986-19-6 14
5		3-Pyridinecarboxylic acid, 1,2-d...	139 C6H5NO3	000609-71-2 14



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012418\
Data File : BF102390.D
Acq On : 24 Jan 2018 15:35
Operator : SJ/JU
Sample : J1235-07 5X
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampled :
30-2

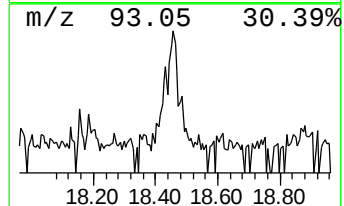
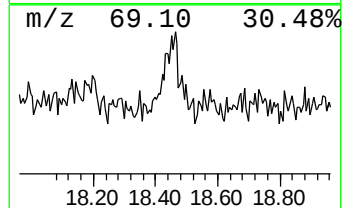
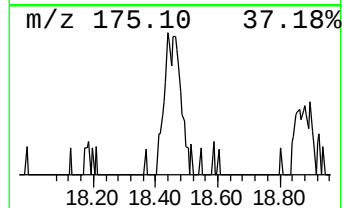
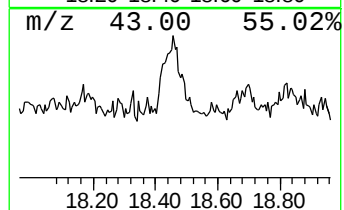
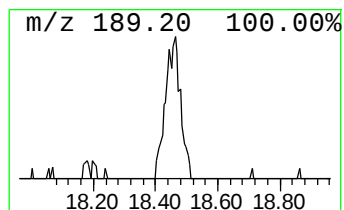
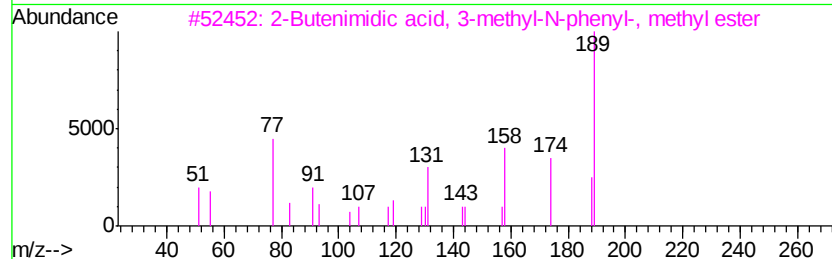
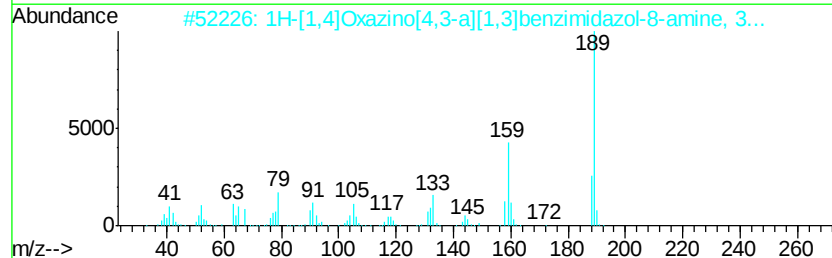
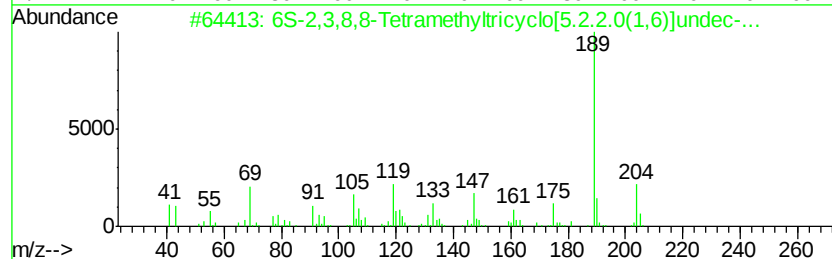
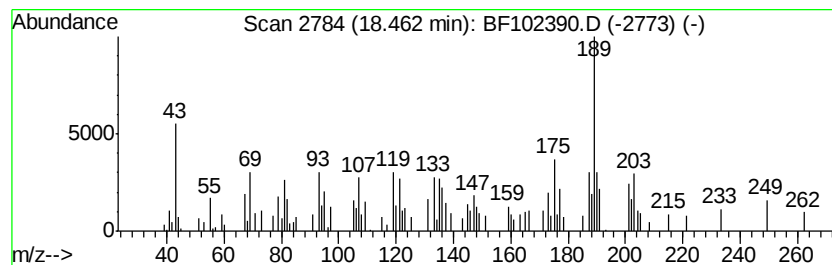
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown18.46 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.46	4.31 ng	210848	Perylene-d12	15.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	6S-2,3,8,8-Tetramethyltricyclo[5...	204	C15H24	137235-48-4	30
2		1H-[1,4]Oxazino[4,3-a][1,3]benzi...	189	C10H11N3O	1000362-54-6	27
3		2-Butenimidic acid, 3-methyl-N-p...	189	C12H15NO	056830-02-5	27
4		2-Fluoro-4-(trifluoromethyl)benz...	189	C8H3F4N	146070-34-0	27
5		9-Methyltetracyclo[7.3.1.0(2,7)....	204	C15H24	1000215-30-5	25



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012418\
 Data File : BF102390.D
 Acq On : 24 Jan 2018 15:35
 Operator : SJ/JU
 Sample : J1235-07 5X
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 30-2

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.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...	5.87	2.8	ng	107288	1	6.71	768579	20.0
Benzene, propyl-	6.16	4.1	ng	156775	1	6.71	768579	20.0
unknown6.47	6.47	19.2	ng	738569	1	6.71	768579	20.0
Mesitylene	6.53	4.0	ng	154441	1	6.71	768579	20.0
Benzene, 1,3-diet...	6.96	4.2	ng	160609	1	6.71	768579	20.0
Benzene, 1,2-diet...	7.03	2.3	ng	88049	1	6.71	768579	20.0
Benzene, 1,2,4,5-...	7.48	2.2	ng	122494	2	7.99	1109690	20.0
1-Phenyl-1-butene	7.73	2.3	ng	125290	2	7.99	1109690	20.0
Perylene	15.17	2.5	ng	124673	6	15.29	979365	20.0
unknown15.48	15.48	3.0	ng	149427	6	15.29	979365	20.0
unknown18.46	18.46	4.3	ng	210848	6	15.29	979365	20.0