

Data Path : Z:\HPCHEM1\BNA F\DATA\BF043017\
 Data File : BF094738.D
 Acq On : 1 May 2017 3:19
 Operator : SJ/MA
 Sample : I2902-05
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-7BS

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-BF041717.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	3.919	333	337	348	rBV	112729	117181	3.25%	0.401%
2	4.166	376	379	384	rVB	71123	64120	1.78%	0.219%
3	4.290	393	400	403	rVB	25967	41429	1.15%	0.142%
4	4.325	403	406	419	rBV	1358343	1359073	37.75%	4.645%
5	4.525	436	440	448	rVB	36553	38967	1.08%	0.133%
6	4.643	456	460	465	rBV	68092	72318	2.01%	0.247%
7	5.395	585	588	605	rBV	1053027	1087836	30.22%	3.718%
8	5.519	605	609	615	rVV	2450134	2481608	68.93%	8.482%
9	5.572	615	618	625	rVB2	54227	59109	1.64%	0.202%
10	5.766	647	651	655	rBV	647777	630479	17.51%	2.155%
11	5.831	660	662	665	rVB	92246	75615	2.10%	0.258%
12	5.942	676	681	684	rBV	2483440	2313488	64.26%	7.907%
13	5.966	684	685	691	rVB	74433	50545	1.40%	0.173%
14	6.054	697	700	704	rBV	183069	170277	4.73%	0.582%
15	6.419	751	762	765	rBV	1883071	1929874	53.61%	6.596%
16	6.513	775	778	783	rVB	114076	110957	3.08%	0.379%
17	6.695	806	809	814	rBV	48830	42718	1.19%	0.146%
18	6.848	830	835	838	rBV2	60773	83382	2.32%	0.285%
19	6.960	849	854	859	rBV2	119555	153134	4.25%	0.523%
20	7.001	859	861	863	rVV2	65823	62000	1.72%	0.212%
21	7.019	863	864	870	rVB2	35708	39523	1.10%	0.135%
22	7.254	900	904	907	rBV	945558	924113	25.67%	3.159%
23	7.284	907	909	913	rVB	329132	273706	7.60%	0.936%
24	7.495	938	945	950	rBV3	30355	44095	1.22%	0.151%
25	7.601	960	963	967	rVB2	35207	36158	1.00%	0.124%
26	7.801	983	997	1000	rBV	629109	1311533	36.43%	4.483%
27	7.878	1008	1010	1016	rVB4	25713	41423	1.15%	0.142%
28	8.142	1047	1055	1060	rVB2	39014	63007	1.75%	0.215%
29	8.242	1067	1072	1080	rBV	453426	446532	12.40%	1.526%
30	8.583	1125	1130	1133	rBV	3519470	3600167	100.00%	12.305%
31	8.695	1143	1149	1153	rBV	201918	181288	5.04%	0.620%
32	8.789	1158	1165	1168	rBV4	23640	41219	1.14%	0.141%
33	8.819	1168	1170	1175	rVB	57362	56055	1.56%	0.192%
34	8.901	1179	1184	1186	rBV3	44439	63643	1.77%	0.218%

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35	8.983	1192	1198	1201	rBV	114651	126158	3.50%	0.431%
36	9.083	1209	1215	1218	rBV	94915	95788	2.66%	0.327%
37	9.125	1218	1222	1224	rBV	47706	57662	1.60%	0.197%
38	9.219	1231	1238	1244	rBV3	218304	277282	7.70%	0.948%
39	9.389	1260	1267	1271	rBV2	1023838	1041216	28.92%	3.559%
40	9.430	1271	1274	1277	rVB	680587	600196	16.67%	2.051%
41	9.536	1288	1292	1297	rBV	39012	61801	1.72%	0.211%
42	9.648	1304	1311	1317	rBV	58530	86189	2.39%	0.295%
43	9.789	1329	1335	1338	rBV3	38753	56361	1.57%	0.193%
44	9.901	1350	1354	1359	rBV	61328	80459	2.23%	0.275%
45	9.954	1359	1363	1366	rVB	76369	75314	2.09%	0.257%
46	10.007	1369	1372	1374	rVV	122006	114948	3.19%	0.393%
47	10.030	1374	1376	1379	rVB	53691	44653	1.24%	0.153%
48	10.066	1379	1382	1388	rVB	222871	213615	5.93%	0.730%
49	10.130	1390	1393	1398	rBV2	48940	69845	1.94%	0.239%
50	10.342	1423	1429	1430	rBV	49674	54957	1.53%	0.188%
51	10.372	1430	1434	1441	rVB	1616228	1761053	48.92%	6.019%
52	10.572	1464	1468	1471	rBV	47326	53659	1.49%	0.183%
53	10.854	1512	1516	1519	rVB2	30465	37542	1.04%	0.128%
54	10.995	1537	1540	1545	rVB2	73817	76275	2.12%	0.261%
55	11.213	1573	1577	1580	rBV	899015	924491	25.68%	3.160%
56	11.242	1580	1582	1587	rVV	602129	525736	14.60%	1.797%
57	11.307	1587	1593	1596	rVB	62017	65127	1.81%	0.223%
58	11.871	1686	1689	1694	rVB	42904	45667	1.27%	0.156%
59	11.948	1696	1702	1703	rBV2	41234	45894	1.27%	0.157%
60	11.971	1703	1706	1708	rVV2	65052	75660	2.10%	0.259%
61	12.001	1708	1711	1715	rVB	76603	79172	2.20%	0.271%
62	12.707	1828	1831	1835	rVB	41578	38358	1.07%	0.131%
63	12.983	1874	1878	1881	rVB	44611	46866	1.30%	0.160%
64	13.024	1881	1885	1889	rBV	64568	62973	1.75%	0.215%
65	13.195	1908	1914	1917	rBV	2557554	2810366	78.06%	9.606%
66	13.901	2030	2034	2040	rBV	43338	44755	1.24%	0.153%
67	14.360	2106	2112	2118	rBV3	26422	45247	1.26%	0.155%
68	14.465	2125	2130	2137	rBV	579723	613064	17.03%	2.095%
69	16.095	2402	2407	2416	rBV	534624	611575	16.99%	2.090%
70	16.906	2536	2545	2557	rBV	39595	170818	4.74%	0.584%

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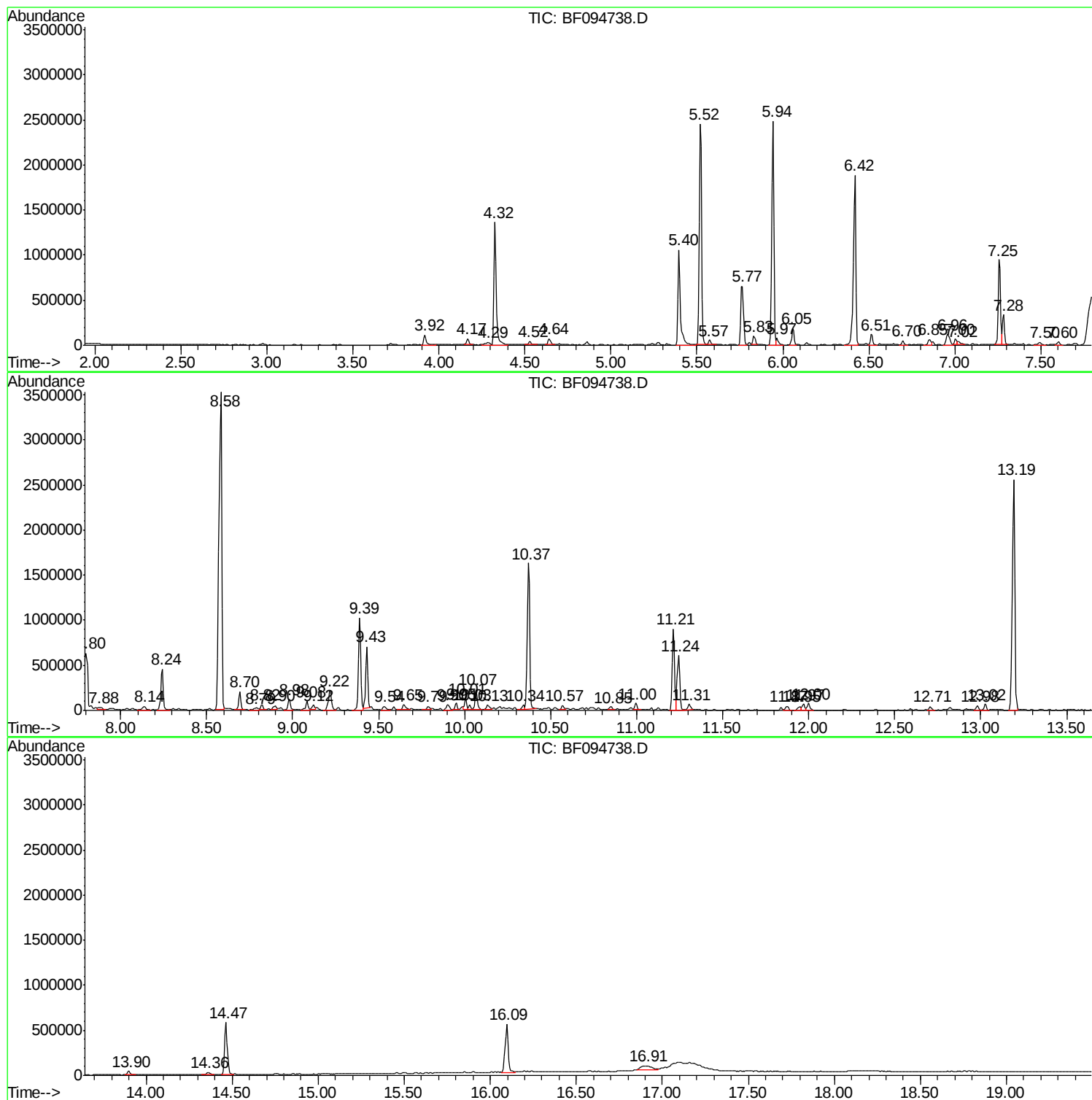
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF041717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Instrument :
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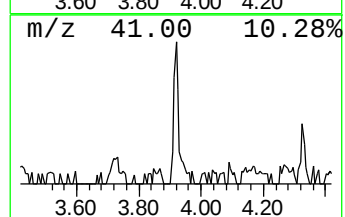
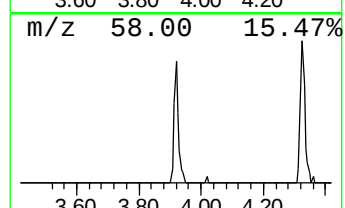
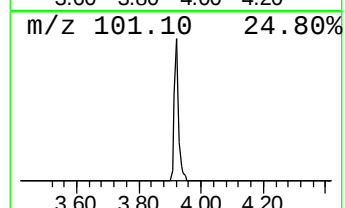
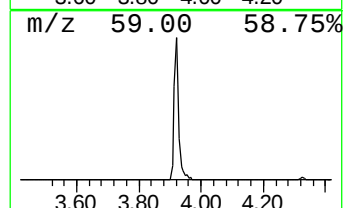
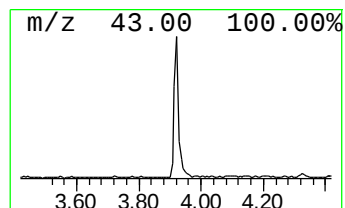
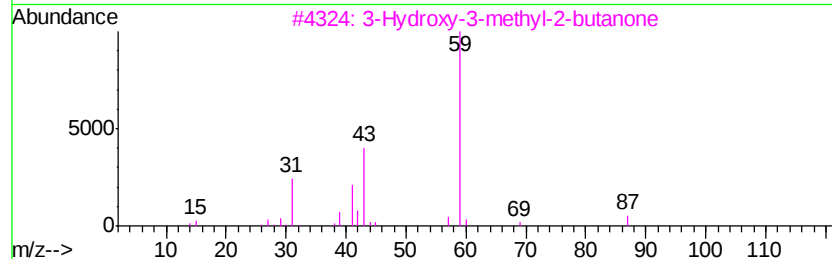
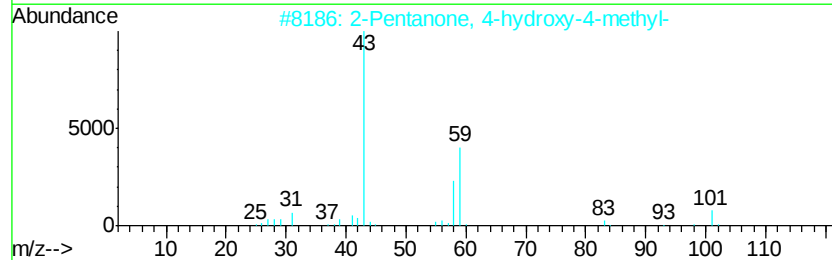
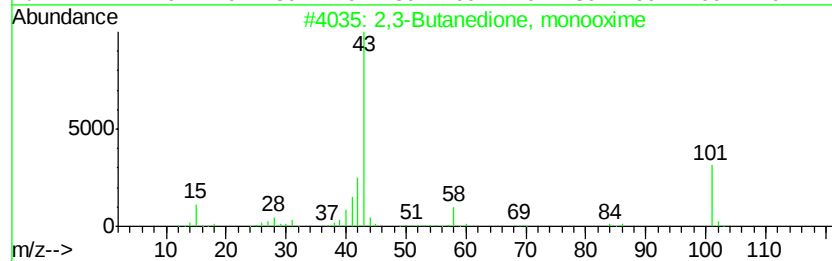
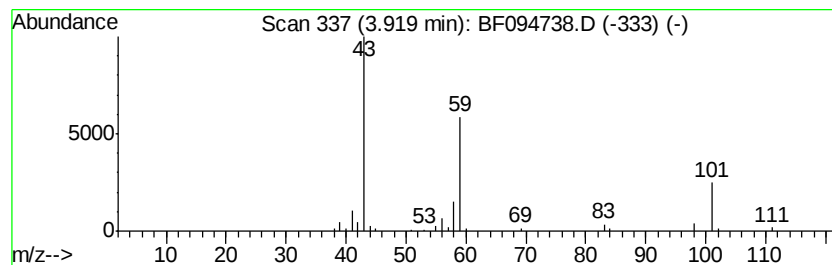
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF041717.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown3.92 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.92	3.72 ng	117181	1,4-Dichlorobenzene-d4	5.77

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	38
2		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	28
3		3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	12
4		5-Hexen-2-one	98	C6H10O	000109-49-9	9
5		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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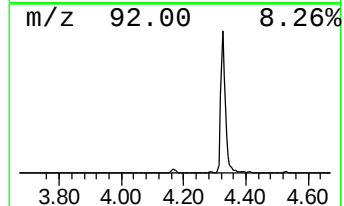
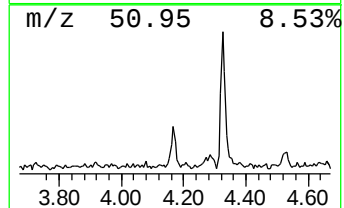
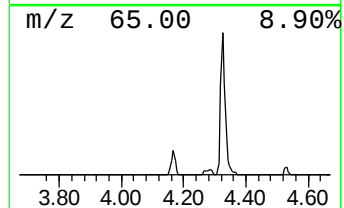
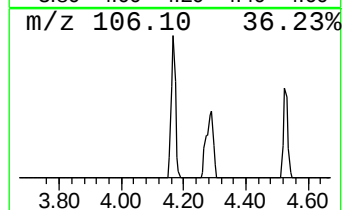
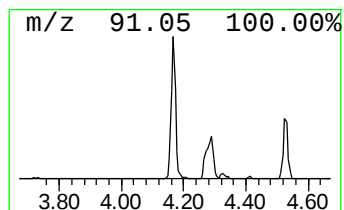
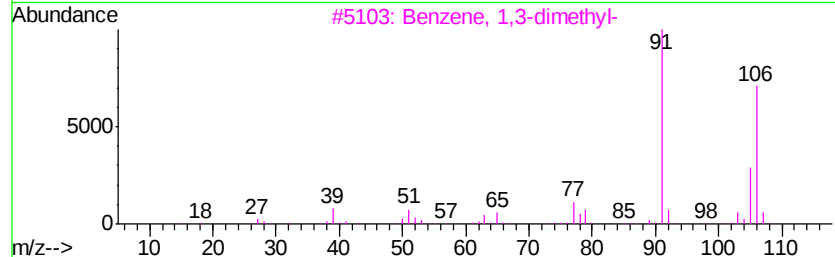
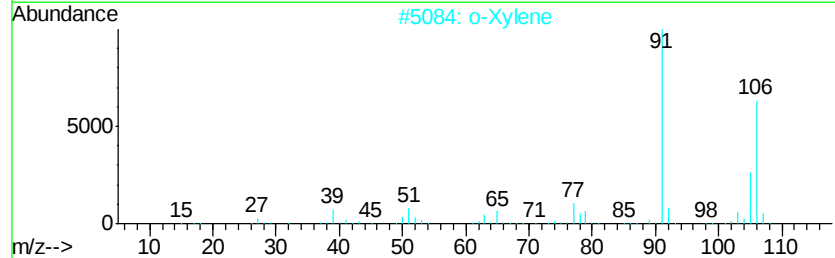
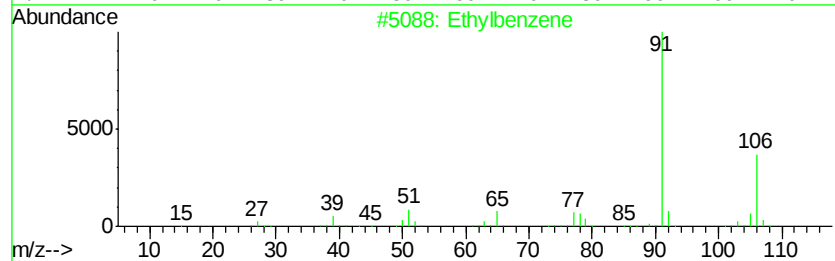
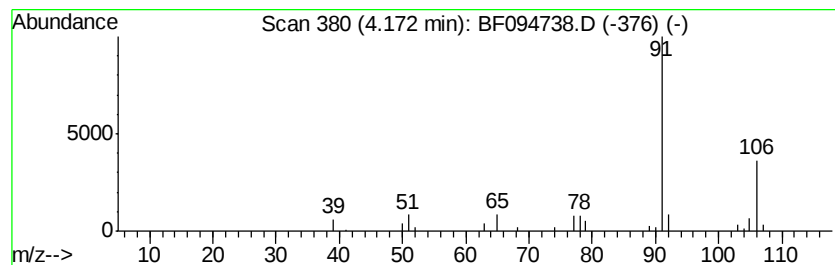
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Ethylbenzene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.17	2.03 ng	64120	1,4-Dichlorobenzene-d4	5.77

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethylbenzene	106	C8H10	000100-41-4	91
2			o-Xylene	106	C8H10	000095-47-6	90
3			Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	87
4			p-Xylene	106	C8H10	000106-42-3	83
5			1,3-Cyclopentadiene, 5-(1-methyl...	106	C8H10	002175-91-9	80



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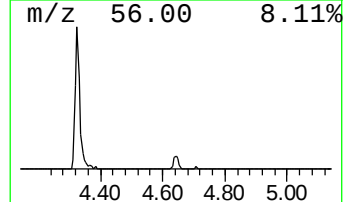
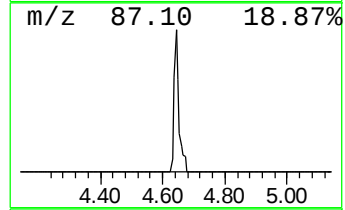
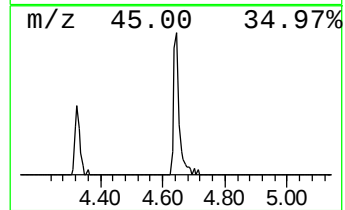
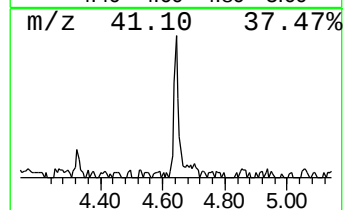
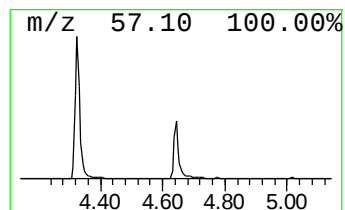
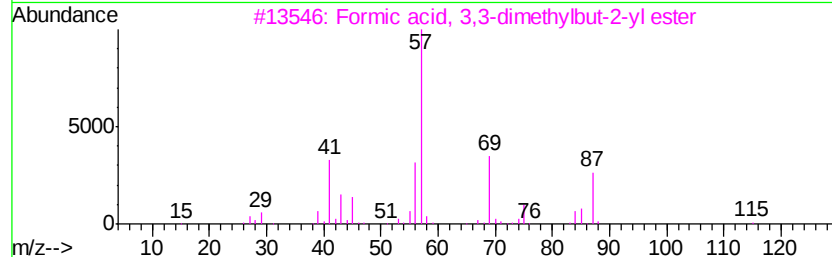
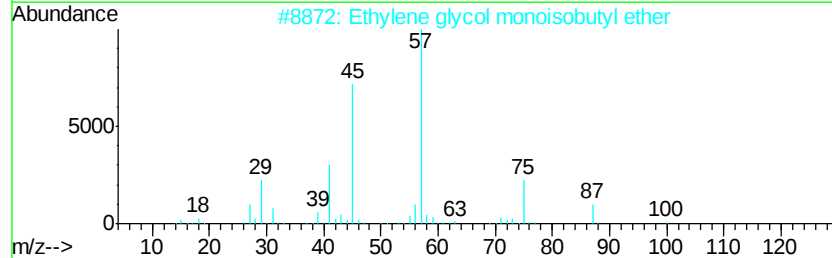
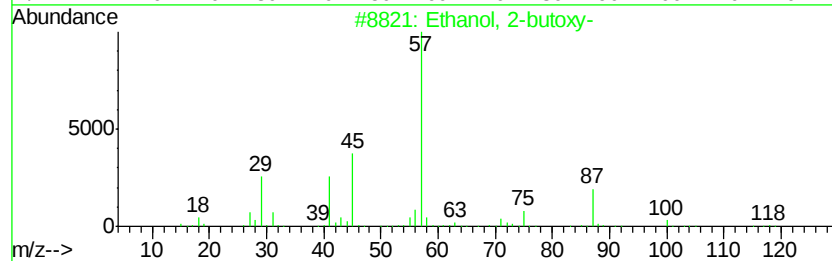
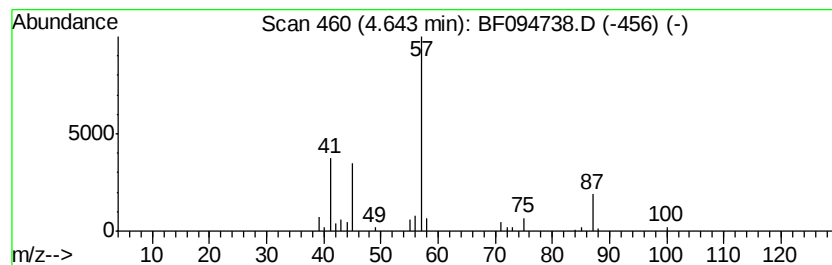
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 Ethanol, 2-butoxy- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.64	2.29 ng	72318	1,4-Dichlorobenzene-d4	5.77

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	78
2		Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	42
3		Formic acid, 3,3-dimethylbut-2-yl...	130	C7H14O2	1000368-20-5	33
4		Propanoic acid, 2-hydroxyethyl e...	118	C5H10O3	024567-27-9	33
5		3-Buten-2-ol	72	C4H8O	000598-32-3	28



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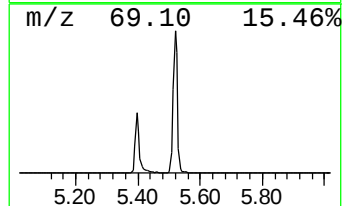
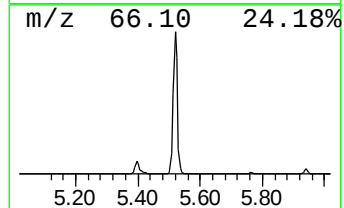
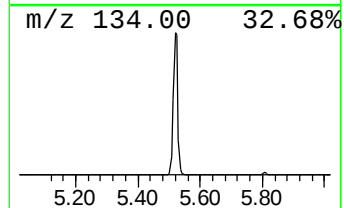
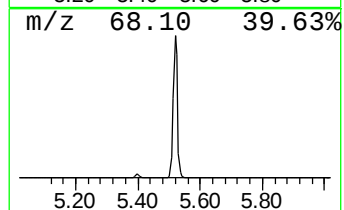
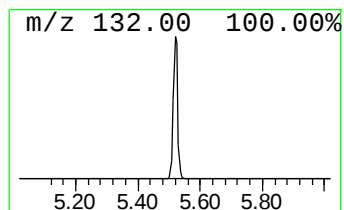
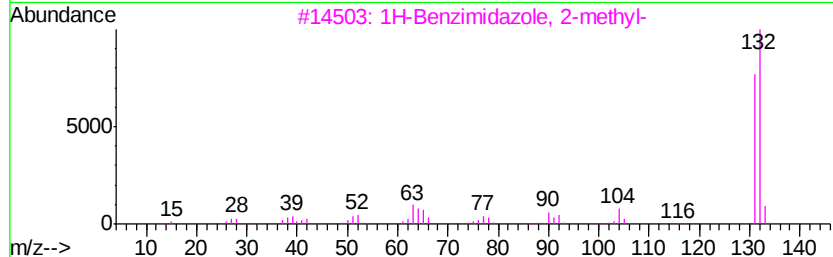
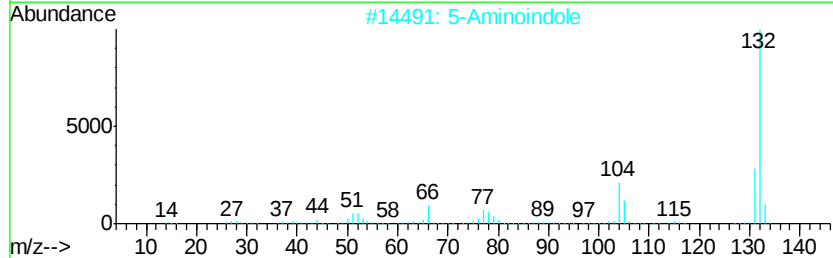
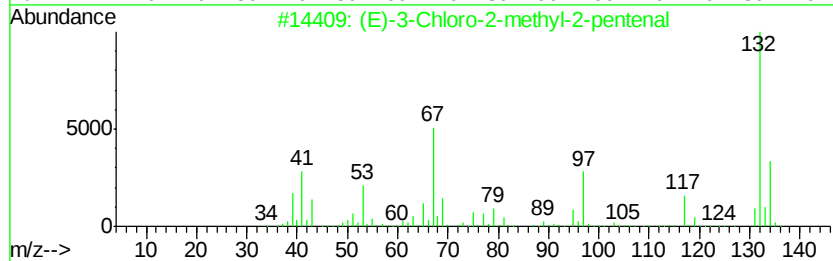
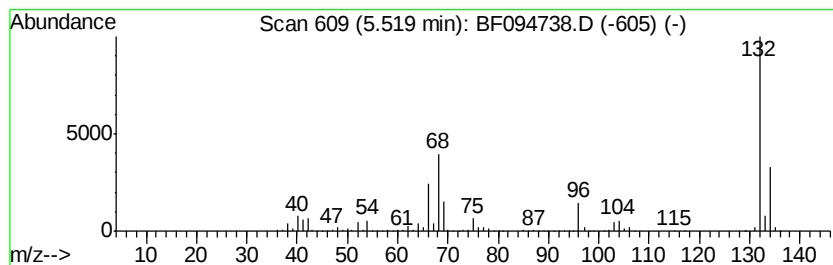
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 4 unknown5.52 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.52	78.72 ng	2481610	1,4-Dichlorobenzene-d4	5.77

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
2		5-Aminoindole	132	C8H8N2	005192-03-0	12
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	10
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF043017\
Data File : BF094738.D
Acq On : 1 May 2017 3:19
Operator : SJ/MA
Sample : I2902-05
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-7BS

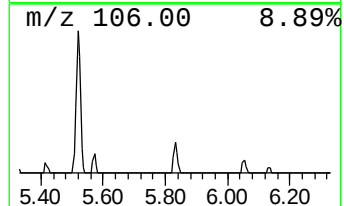
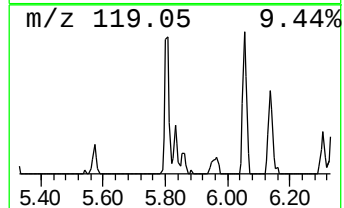
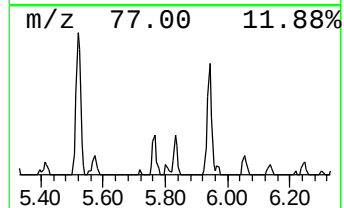
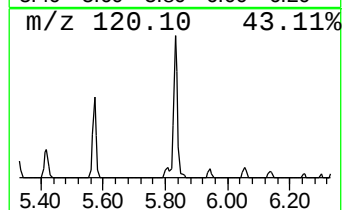
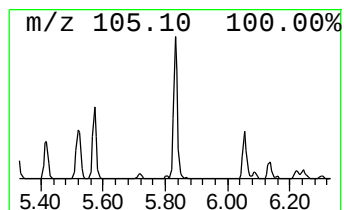
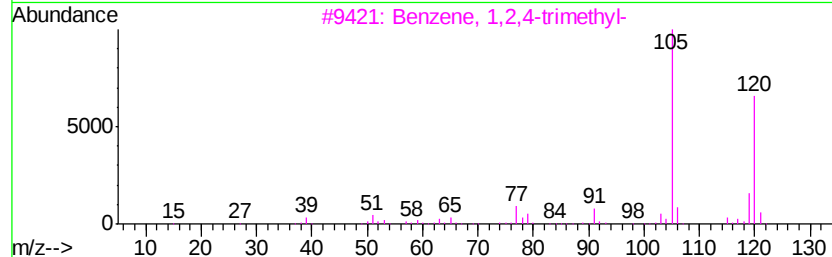
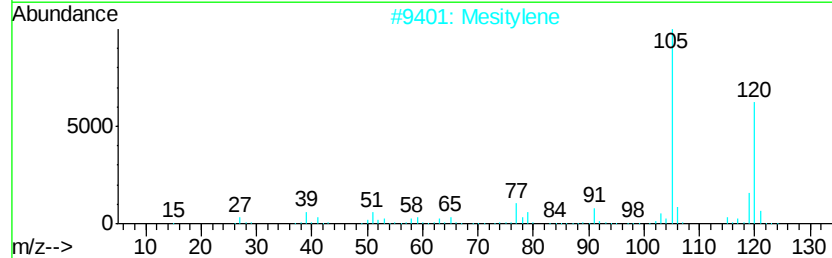
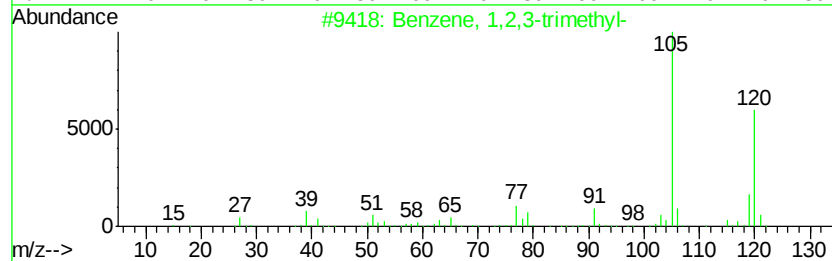
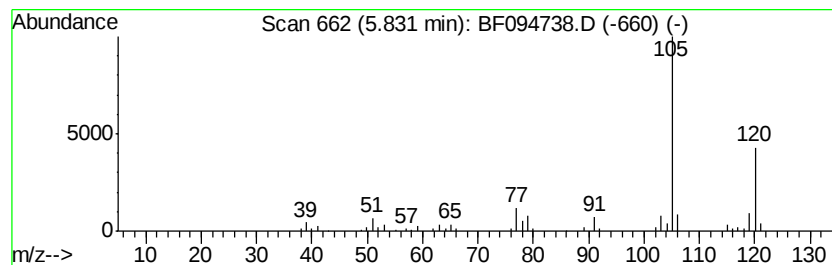
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Benzene, 1,2,3-trimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.83	2.40 ng	75615	1,4-Dichlorobenzene-d4	5.77

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
2		Mesitylene	120	C9H12	000108-67-8	91
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\BF043017\
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Acq On : 1 May 2017 3:19
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ClientSampleId :
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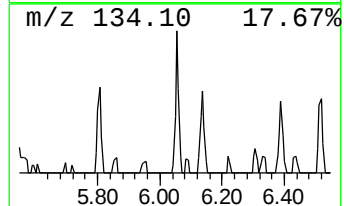
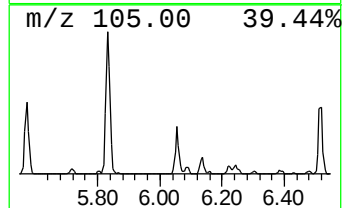
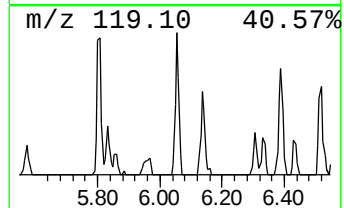
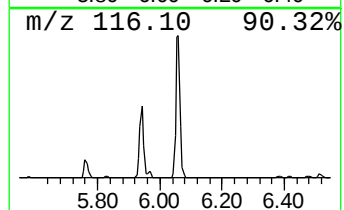
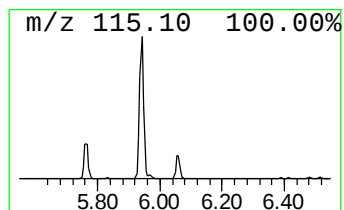
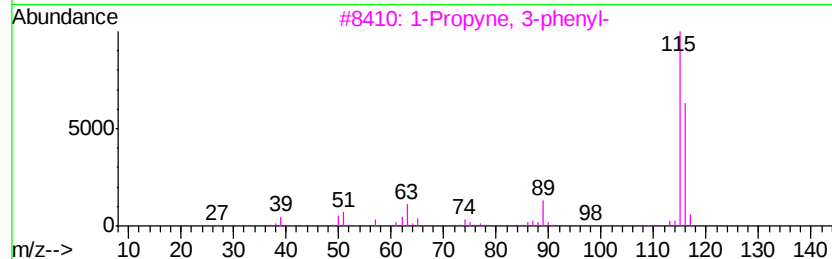
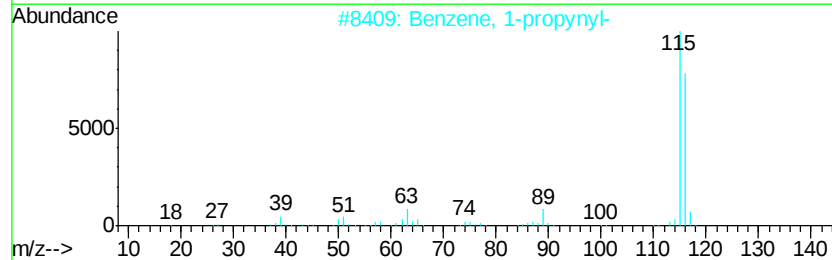
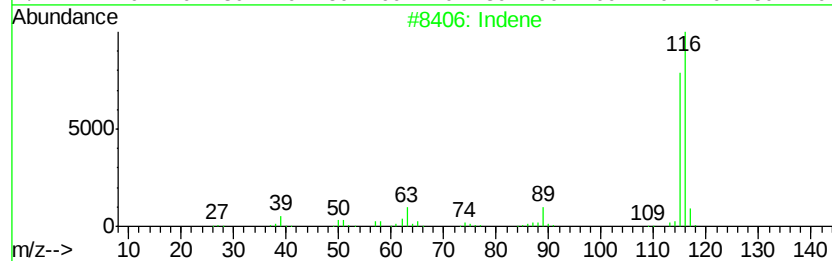
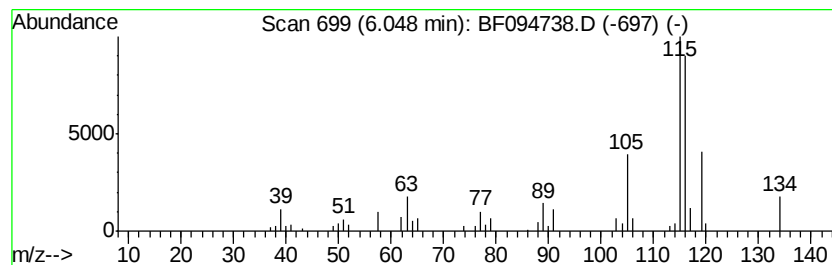
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Indene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.05	5.40 ng	170277	1,4-Dichlorobenzene-d4	5.77

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indene	116	C9H8	000095-13-6	95
2	Benzene, 1-propynyl-	116	C9H8	000673-32-5	94
3	1-Propyne, 3-phenyl-	116	C9H8	010147-11-2	90
4	3-Methylphenylacetylene	116	C9H8	000766-82-5	90
5	Benzene, 1-ethynyl-4-methyl-	116	C9H8	000766-97-2	76



Data Path : Z:\HPCHEM1\BNA F\DATA\BF043017\
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Acq On : 1 May 2017 3:19
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Sample : I2902-05
Misc :
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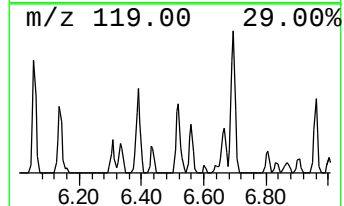
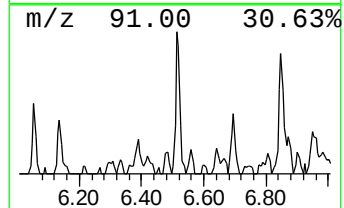
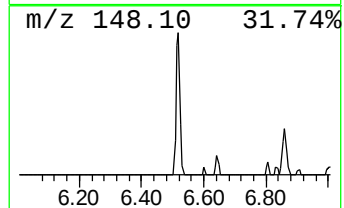
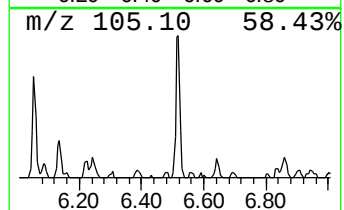
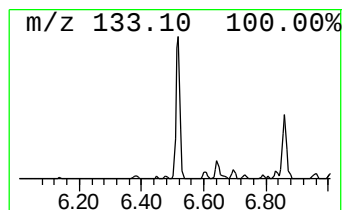
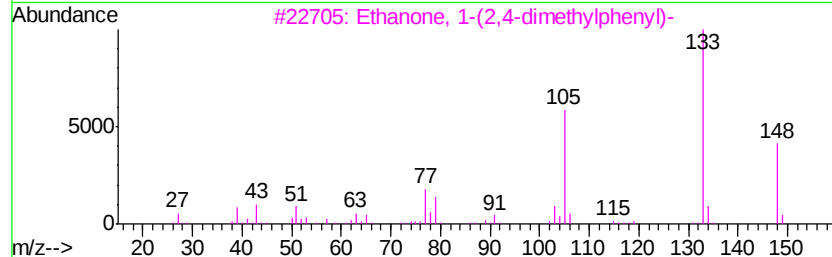
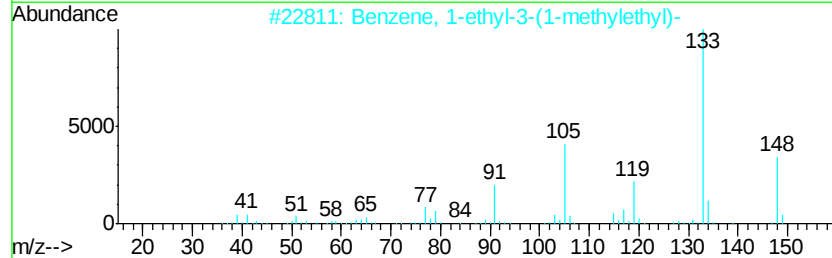
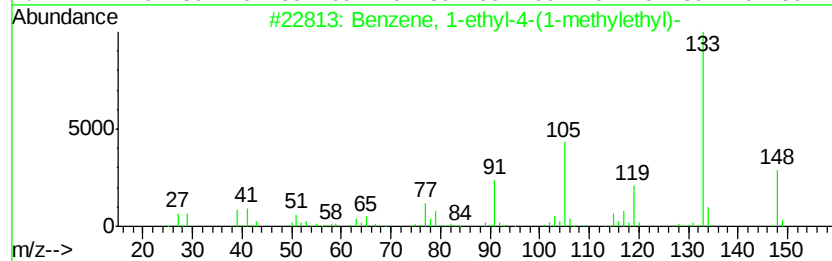
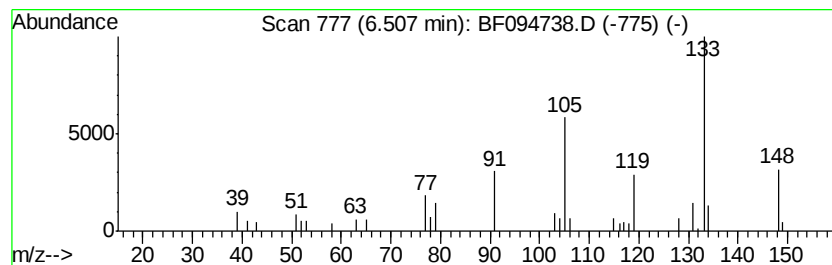
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Benzene, 1-ethyl-4-(1-methylethyl)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.51	2.40 ng	110957	Naphthalene-d8	7.25

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	94
2		Benzene, 1-ethyl-3-(1-methylethyl)-	148	C11H16	004920-99-4	94
3		Ethanone, 1-(2,4-dimethylphenyl)-	148	C10H12O	000089-74-7	92
4		4-tert-Butyltoluene	148	C11H16	000098-51-1	81
5		Benzene, 1-(1,1-dimethylethyl)-3...	148	C11H16	001075-38-3	76



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF043017\
Data File : BF094738.D
Acq On : 1 May 2017 3:19
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Misc :
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Instrument :
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ClientSampleId :
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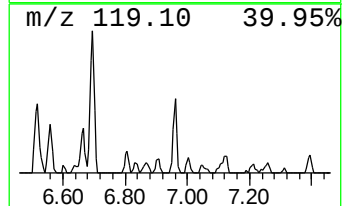
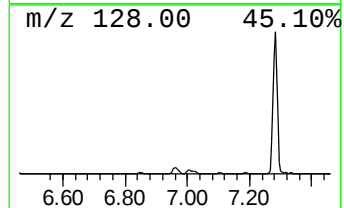
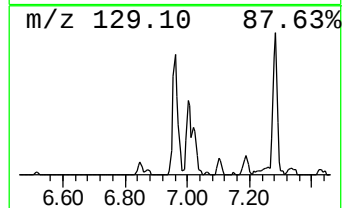
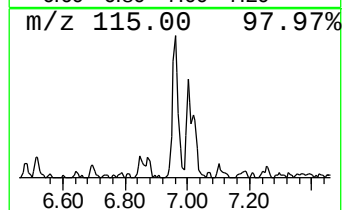
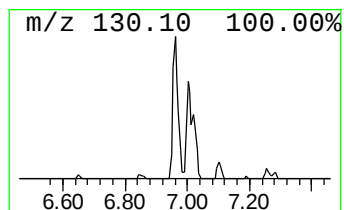
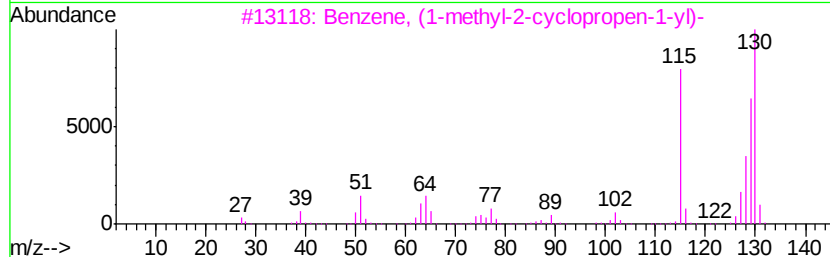
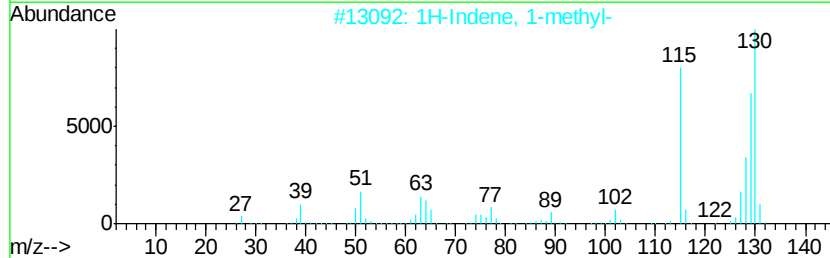
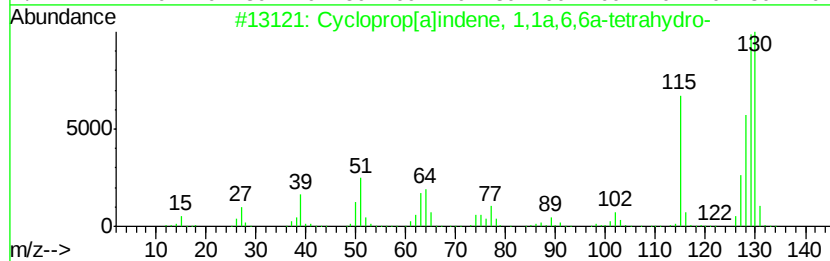
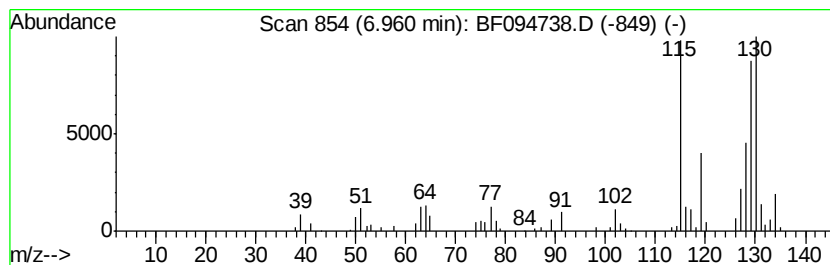
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Cycloprop[a]indene, 1,1a,6,6a,... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.96	3.31 ng	153134	Napthalene-d8	7.25

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cycloprop[a]indene, 1,1a,6,6a-te...	130	C10H10	015677-15-3	97
2			1H-Indene, 1-methyl-	130	C10H10	000767-59-9	93
3			Benzene, (1-methyl-2-cyclopropen...	130	C10H10	065051-83-4	93
4			1,4-Dihydronaphthalene	130	C10H10	000612-17-9	91
5			Benzene, 1-butynyl-	130	C10H10	000622-76-4	87



Data Path : Z:\HPCHEM1\BNA F\DATA\BF043017\
Data File : BF094738.D
Acq On : 1 May 2017 3:19
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Misc :
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ClientSampleId :
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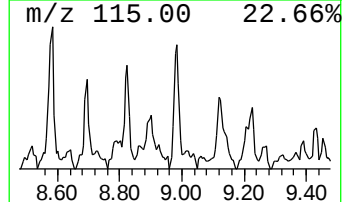
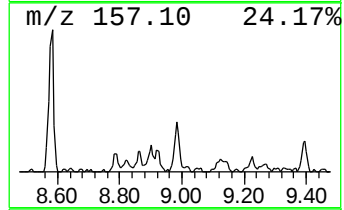
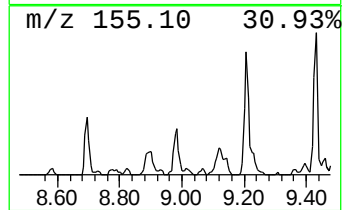
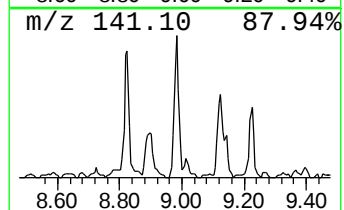
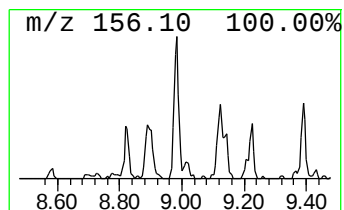
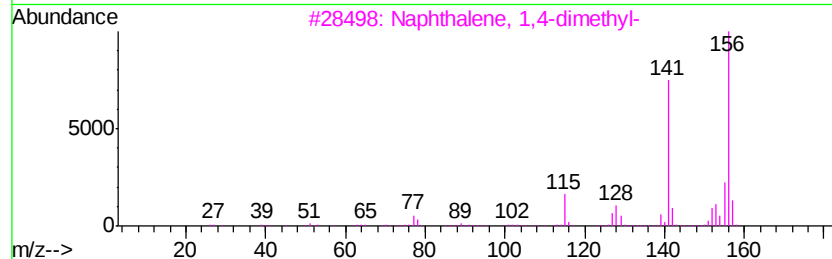
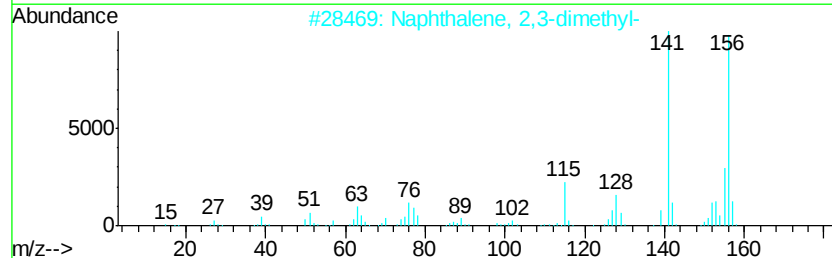
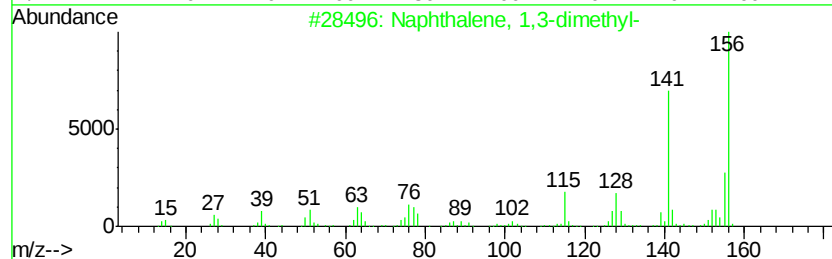
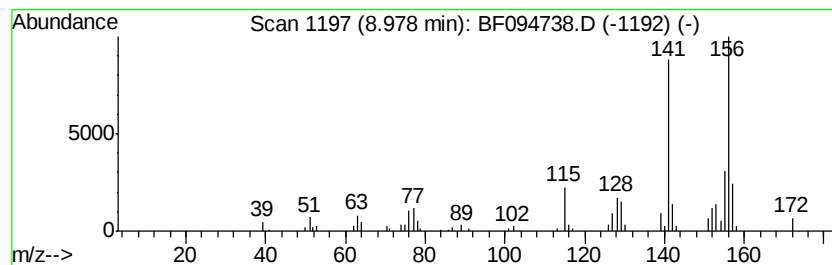
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Naphthalene, 1,3-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.98	2.42 ng	126158	Acenaphthene-d10	9.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	97
2		Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	95
3		Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	94
4		Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	94
5		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	93



Data Path : Z:\HPCHEM1\BNA F\DATA\BF043017\
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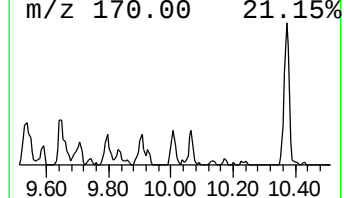
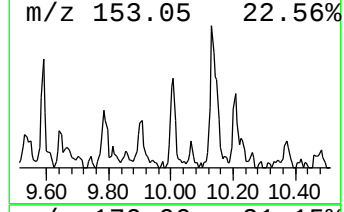
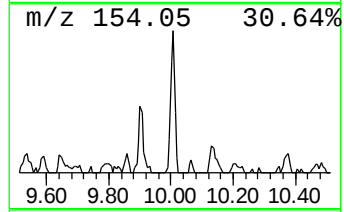
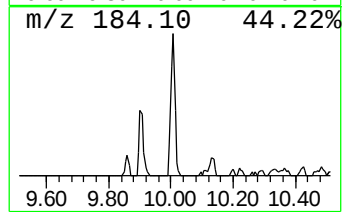
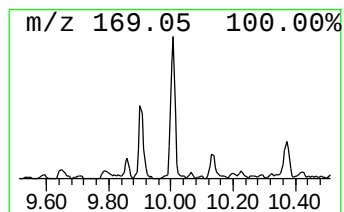
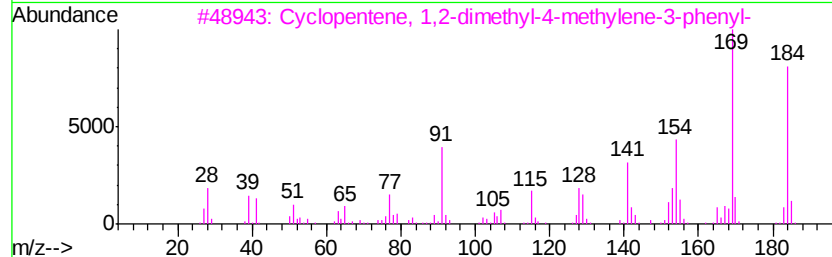
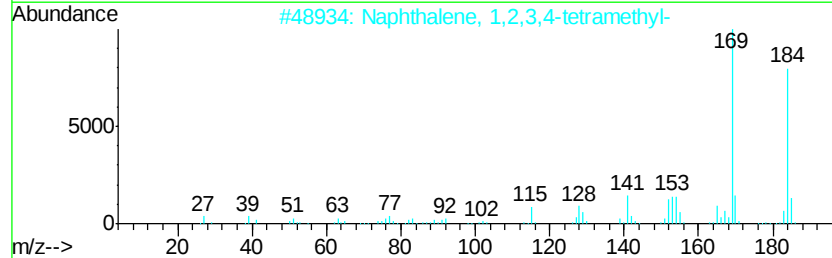
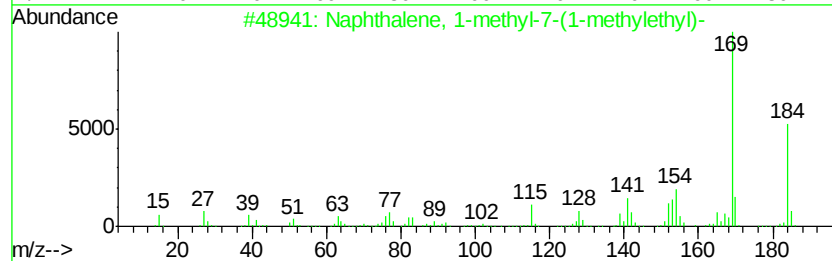
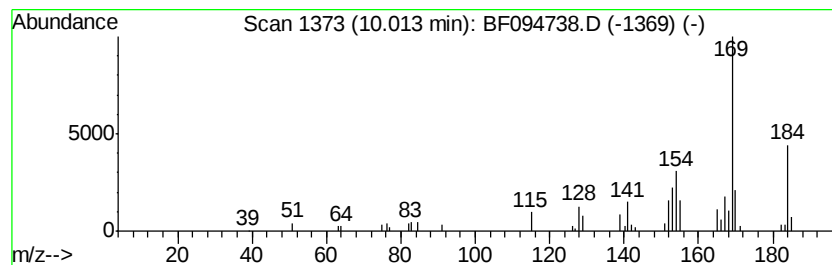
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Naphthalene, 1-methyl-7-(1-... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.01	2.21 ng	114948	Acenaphthene-d10	9.39

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1-methyl-7-(1-methyl-...	184	C14H16	000490-65-3	87
2			Naphthalene, 1,2,3,4-tetramethyl-	184	C14H16	003031-15-0	87
3			Cyclopentene, 1,2-dimethyl-4-met...	184	C14H16	1000152-22-4	76
4			1-Acenaphthylenol, 1,2-dihydro-1...	184	C13H12O	041002-74-8	58
5			Naphthalene, 1-(1,1-dimethylethyl)-	184	C14H16	017085-91-5	53



Data Path : Z:\HPCHEM1\BNA F\DATA\BF043017\
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ALS Vial : 31 Sample Multiplier: 1

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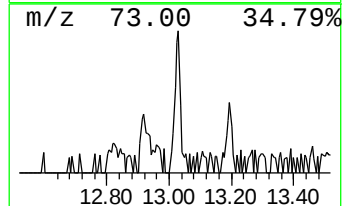
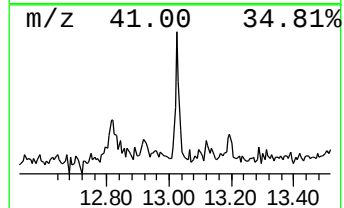
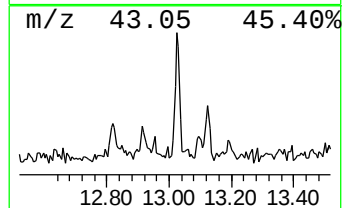
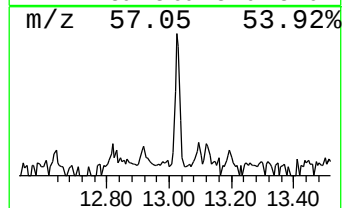
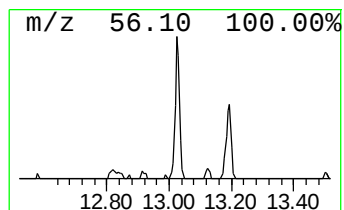
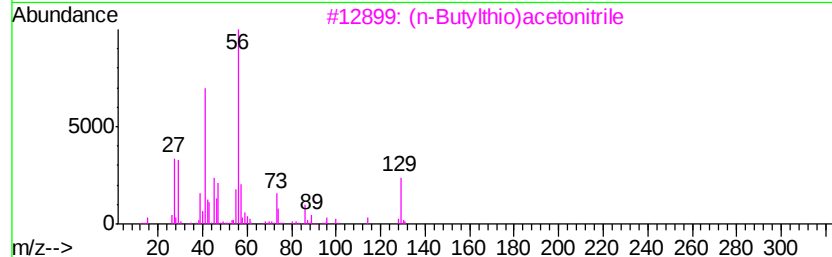
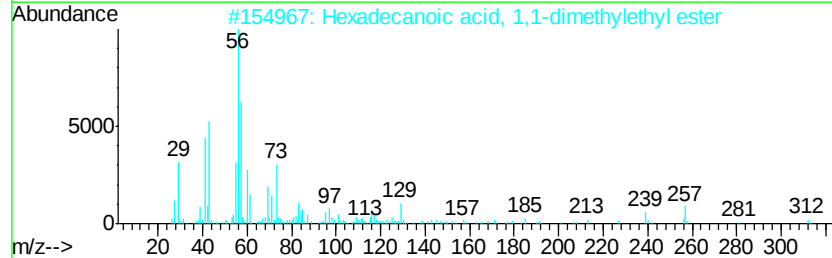
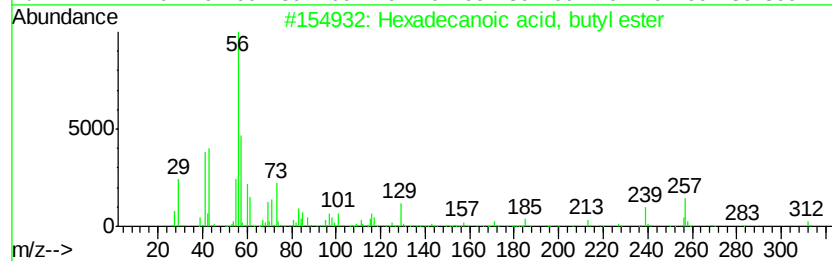
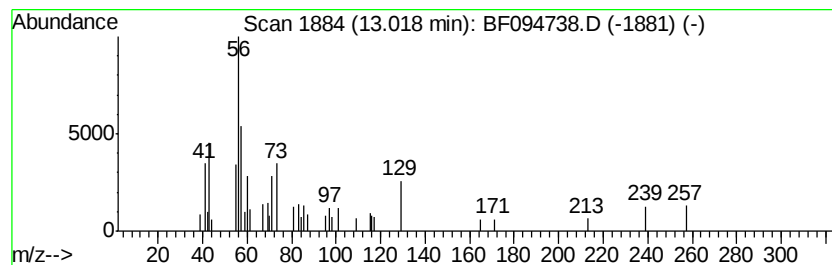
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Hexadecanoic acid, butyl ester Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.02	2.05 ng	62973	Chrysene-d12	14.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecanoic acid, butyl ester	312	C20H40O2	000111-06-8	87
2		Hexadecanoic acid, 1,1-dimethyle...	312	C20H40O2	031158-91-5	64
3		(n-Butylthio)acetonitrile	129	C6H11NS	071037-08-6	47
4		1-Pentene, 2,4-dimethyl-	98	C7H14	002213-32-3	27
5		Cyclobutanone, 3,3-dimethyl-	98	C6H10O	001192-33-2	27



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF043017\
Data File : BF094738.D
Acq On : 1 May 2017 3:19
Operator : SJ/MA
Sample : I2902-05
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-7BS

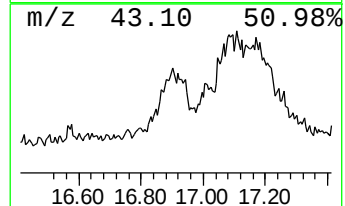
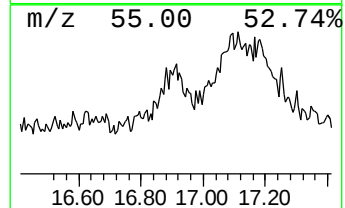
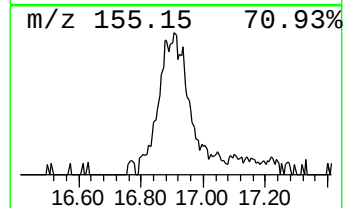
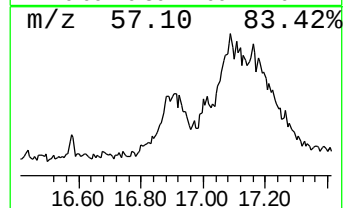
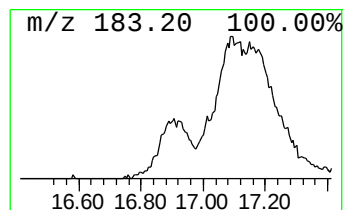
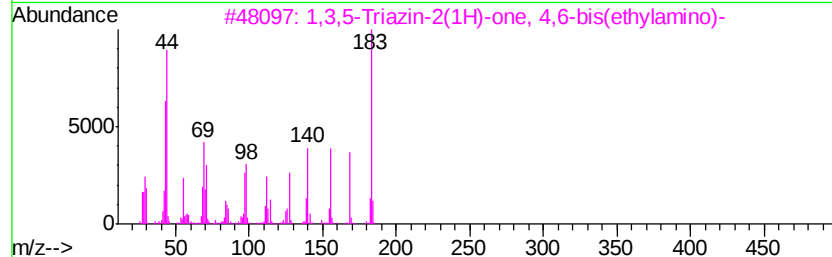
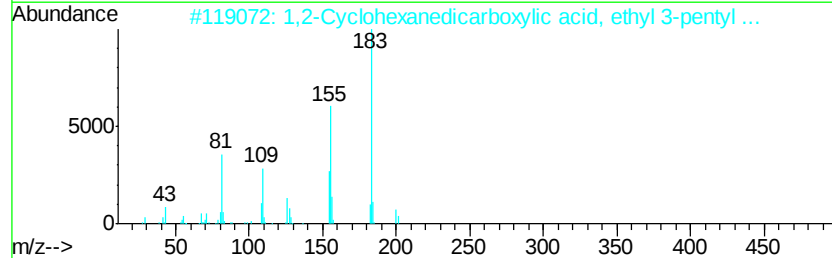
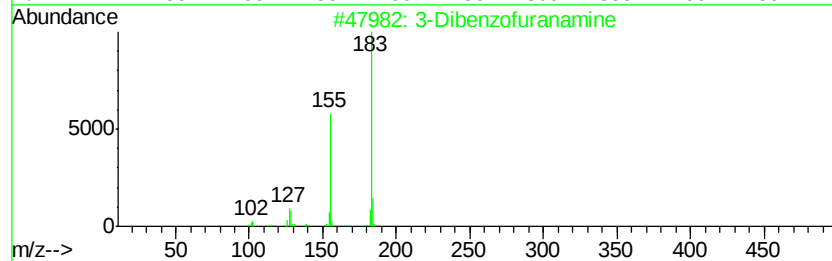
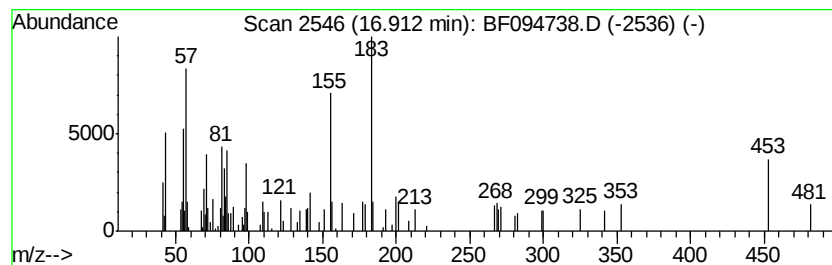
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown16.91 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.91	5.59 ng	170818	Perylene-d12	16.09

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Dibenzofuranamine	183	C12H9NO	004106-66-5	43
2			1,2-Cyclohexanedicarboxylic acid...	270	C15H26O4	1000339-50-2	37
3			1,3,5-Triazin-2(1H)-one, 4,6-bis...	183	C7H13N5O	002599-11-3	35
4			1H-Indazole, 6-(pyrrol-1-yl)-	183	C11H9N3	1000316-64-8	35
5			(Z)-4,6-Dioxohept-2-enoic acid, ...	300	C13H24O4Si2	1000099-84-6	27



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF043017\
Data File : BF094738.D
Acq On : 1 May 2017 3:19
Operator : SJ/MA
Sample : I2902-05
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-7BS

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF041717.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
unknown3.92	3.92	3.7	ng	117181	1	5.77	630479	20.0
Ethylbenzene	4.17	2.0	ng	64120	1	5.77	630479	20.0
Ethanol, 2-butoxy-	4.64	2.3	ng	72318	1	5.77	630479	20.0
unknown5.52	5.52	78.7	ng	2481610	1	5.77	630479	20.0
Benzene, 1,2,3-tr...	5.83	2.4	ng	75615	1	5.77	630479	20.0
Indene	6.05	5.4	ng	170277	1	5.77	630479	20.0
Benzene, 1-ethyl-...	6.51	2.4	ng	110957	2	7.25	924113	20.0
Cycloprop[a]inden...	6.96	3.3	ng	153134	2	7.25	924113	20.0
Naphthalene, 1,3-...	8.98	2.4	ng	126158	3	9.39	1041220	20.0
Naphthalene, 1-me...	10.01	2.2	ng	114948	3	9.39	1041220	20.0
Hexadecanoic acid...	13.02	2.0	ng	62973	5	14.47	613064	20.0
unknown16.91	16.91	5.6	ng	170818	6	16.09	611575	20.0