

Data Path : Z:\HPCHEM1\BNA F\DATA\BF062817\
 Data File : BF096263.D
 Acq On : 28 Jun 2017 21:36
 Operator : SJ/MA
 Sample : I3852-16 5X
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 G-38-1.5-3

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-BF062717.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	2.851	124	130	134	rVB4	14740	25445	1.98%	0.164%
2	3.792	283	290	297	rBV	36600	60130	4.69%	0.387%
3	4.410	391	395	403	rVB5	20206	31608	2.47%	0.204%
4	4.934	480	484	487	rBV2	11532	16849	1.31%	0.109%
5	4.963	487	489	498	rVB	124487	144362	11.26%	0.930%
6	5.386	558	561	565	rVB	847093	761778	59.42%	4.909%
7	5.628	600	602	612	rVB2	33997	38705	3.02%	0.249%
8	5.704	612	615	619	rBV2	28274	29059	2.27%	0.187%
9	5.951	650	657	661	rVB5	20495	23616	1.84%	0.152%
10	6.045	669	673	680	rVB5	19328	23248	1.81%	0.150%
11	6.239	700	706	710	rBV5	13770	26037	2.03%	0.168%
12	6.328	717	721	723	rBV3	19558	23936	1.87%	0.154%
13	6.404	729	734	737	rBV	955408	734061	57.26%	4.730%
14	6.545	755	758	762	rBV	836072	727241	56.73%	4.686%
15	6.616	766	770	771	rBV	30849	29064	2.27%	0.187%
16	6.633	771	773	775	rVB	29953	19749	1.54%	0.127%
17	6.781	795	798	802	rBV	991068	897824	70.03%	5.786%
18	6.845	807	809	814	rVB3	23638	24294	1.90%	0.157%
19	6.939	821	825	828	rBV	705537	537473	41.93%	3.464%
20	7.110	850	854	856	rBV3	19748	18909	1.47%	0.122%
21	7.186	862	867	869	rBV3	25320	26237	2.05%	0.169%
22	7.333	886	892	896	rBV	504843	416202	32.47%	2.682%
23	7.392	900	902	905	rVB	43962	35131	2.74%	0.226%
24	7.428	905	908	914	rBV5	23221	23906	1.86%	0.154%
25	7.751	961	963	967	rBV4	23843	24444	1.91%	0.158%
26	8.063	1013	1016	1019	rBV	1389312	1204328	93.94%	7.761%
27	8.086	1019	1020	1023	rVV	110470	60313	4.70%	0.389%
28	8.139	1027	1029	1031	rVB	24014	16330	1.27%	0.105%
29	8.498	1085	1090	1095	rBV	37369	38295	2.99%	0.247%
30	8.669	1116	1119	1122	rVB	43757	32261	2.52%	0.208%
31	8.774	1134	1137	1140	rVB	115899	90402	7.05%	0.583%
32	8.874	1151	1154	1157	rBV	73614	55688	4.34%	0.359%
33	9.098	1190	1192	1195	rVB	27863	21935	1.71%	0.141%
34	9.139	1195	1199	1202	rBV	877333	679397	53.00%	4.378%

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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-BF062717.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	9.233	1211	1215	1220	rBV2	70541	79086	6.17%	0.510%
36	9.333	1230	1232	1237	rBV2	16757	18878	1.47%	0.122%
37	9.398	1240	1243	1251	rBV2	33098	54841	4.28%	0.353%
38	9.474	1254	1256	1259	rVV	49738	47721	3.72%	0.308%
39	9.504	1259	1261	1263	rVB	46468	32753	2.55%	0.211%
40	9.527	1263	1265	1268	rBV	87594	81997	6.40%	0.528%
41	9.551	1268	1269	1271	rVV	35942	24009	1.87%	0.155%
42	9.592	1274	1276	1278	rVV	25361	16178	1.26%	0.104%
43	9.674	1287	1290	1293	rBV	87038	68216	5.32%	0.440%
44	9.763	1302	1305	1309	rVB	55036	46742	3.65%	0.301%
45	9.816	1309	1314	1318	rBV	1322718	1141206	89.02%	7.354%
46	9.845	1318	1319	1323	rVB3	28217	22811	1.78%	0.147%
47	9.933	1328	1334	1337	rBV6	18154	35715	2.79%	0.230%
48	10.021	1345	1349	1352	rBV	69377	65067	5.08%	0.419%
49	10.057	1352	1355	1358	rBV	16531	17277	1.35%	0.111%
50	10.098	1358	1362	1365	rBV	64003	67403	5.26%	0.434%
51	10.163	1370	1373	1376	rVB4	31319	30343	2.37%	0.196%
52	10.227	1380	1384	1387	rBV4	33198	50888	3.97%	0.328%
53	10.263	1387	1390	1392	rVB	52392	45599	3.56%	0.294%
54	10.357	1403	1406	1408	rBV2	51843	41064	3.20%	0.265%
55	10.480	1424	1427	1430	rBV2	51212	52898	4.13%	0.341%
56	10.521	1431	1434	1439	rVB5	29216	32471	2.53%	0.209%
57	10.598	1443	1447	1451	rBV	455496	405887	31.66%	2.616%
58	10.639	1451	1454	1458	rVB6	25994	30892	2.41%	0.199%
59	10.698	1461	1464	1466	rBV	45118	40420	3.15%	0.260%
60	10.745	1466	1472	1476	rVB	180999	227305	17.73%	1.465%
61	10.804	1480	1482	1484	rBV2	16615	19783	1.54%	0.127%
62	11.098	1529	1532	1537	rVB3	31718	43135	3.36%	0.278%
63	11.180	1543	1546	1548	rBV	66918	66600	5.20%	0.429%
64	11.210	1548	1551	1554	rVB	99051	98645	7.69%	0.636%
65	11.298	1562	1566	1568	rBV	1100718	1006345	78.50%	6.485%
66	11.316	1568	1569	1573	rVV	227989	172503	13.46%	1.112%
67	11.368	1575	1578	1581	rVB	148152	119668	9.33%	0.771%
68	11.404	1581	1584	1591	rBV9	23547	52365	4.08%	0.337%
69	11.521	1602	1604	1607	rVB2	33840	28876	2.25%	0.186%
70	11.604	1615	1618	1621	rBV2	65202	59376	4.63%	0.383%
71	11.798	1648	1651	1653	rBV2	50182	51308	4.00%	0.331%

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Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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72	11.827	1653	1656	1660	rVB	174478	150484	11.74%	0.970%
73	11.904	1666	1669	1672	rBV2	90601	102326	7.98%	0.659%
74	12.010	1685	1687	1689	rVB	48941	32773	2.56%	0.211%
75	12.092	1699	1701	1702	rBV	33285	23920	1.87%	0.154%
76	12.345	1742	1744	1748	rBV5	53200	53366	4.16%	0.344%
77	12.398	1751	1753	1756	rVB2	65097	53413	4.17%	0.344%
78	12.427	1756	1758	1764	rBV6	43785	47826	3.73%	0.308%
79	12.504	1768	1771	1775	rBV	478264	394850	30.80%	2.544%
80	12.733	1804	1810	1812	rBV	395608	404985	31.59%	2.610%
81	12.851	1828	1830	1832	rBV2	45435	41323	3.22%	0.266%
82	12.874	1832	1834	1841	rVB	657172	619820	48.35%	3.994%
83	13.127	1874	1877	1880	rBV	88952	72942	5.69%	0.470%
84	13.927	2009	2013	2016	rBV2	1149014	1281972	100.00%	8.261%
85	14.951	2184	2187	2189	rBV	177173	227369	17.74%	1.465%
86	15.356	2253	2256	2260	rVB	513442	596265	46.51%	3.842%

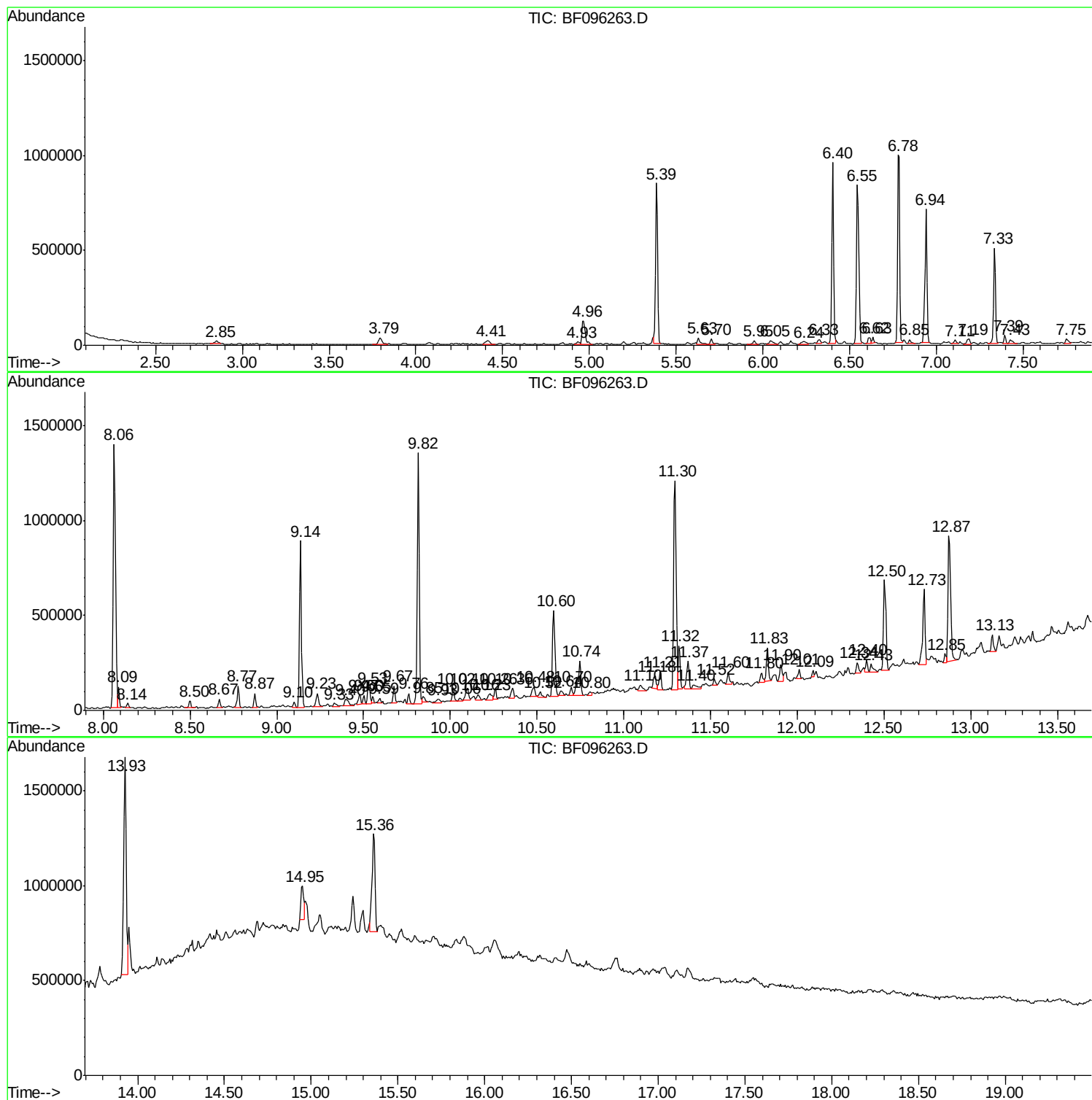
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Instrument :
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ClientSampled :
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF062717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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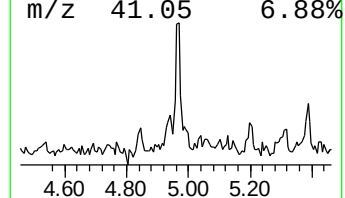
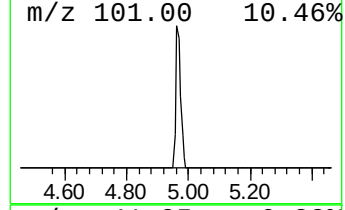
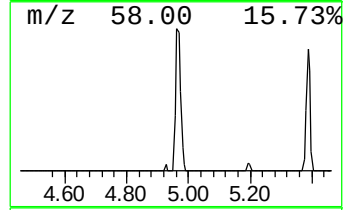
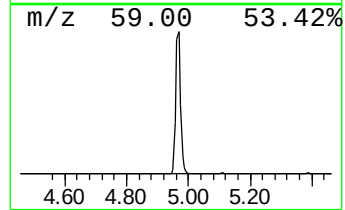
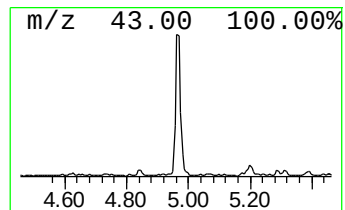
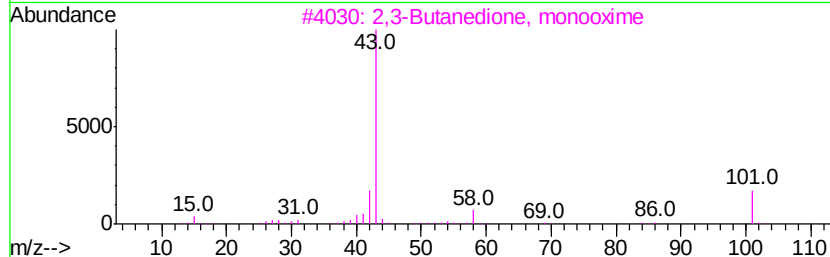
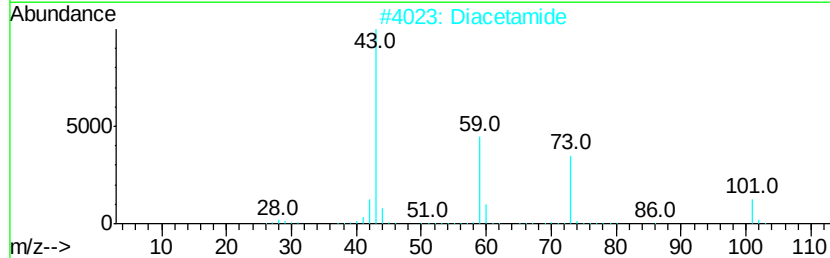
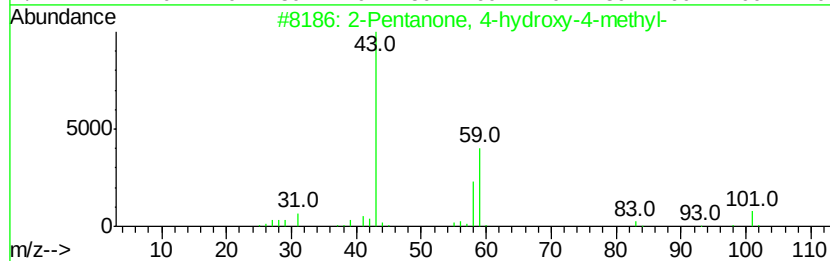
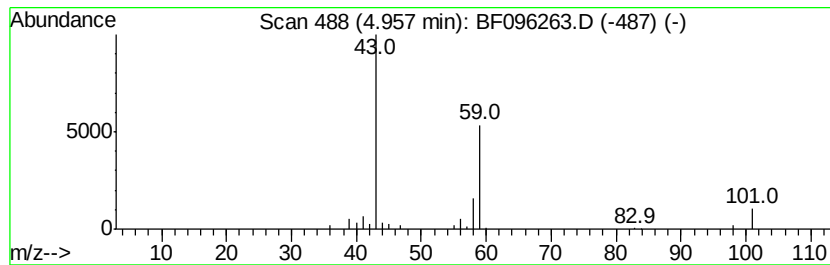
Instrument :
BNA_F
ClientSampleId :
G-38-1.5-3

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF062717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.		
4.96	3.22 ng	144362	1,4-Dichlorobenzene-d4	6.79		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50	
2	Diacetamide	101	C4H7NO2	000625-77-4	9	
3	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9	
4	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9	
5	Formamide, N,N-diethyl-	101	C5H11NO	000617-84-5	7	



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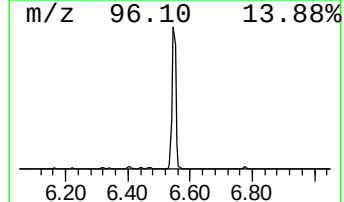
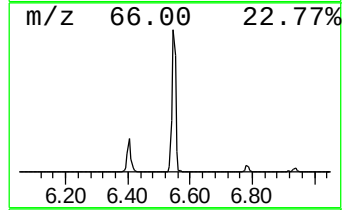
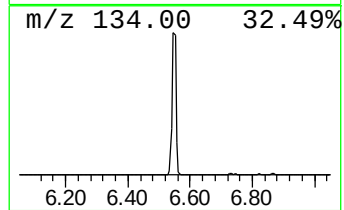
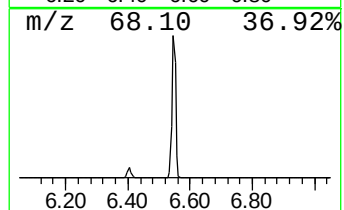
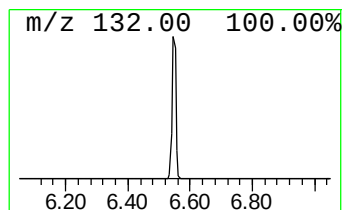
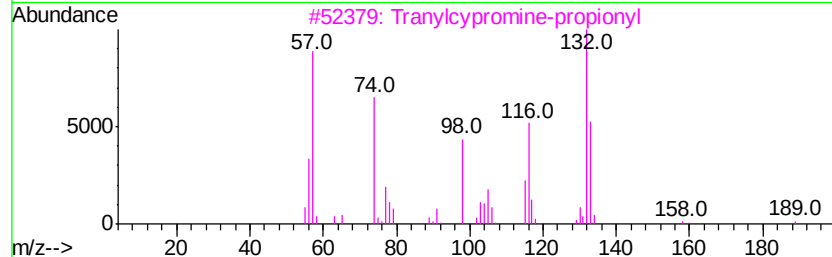
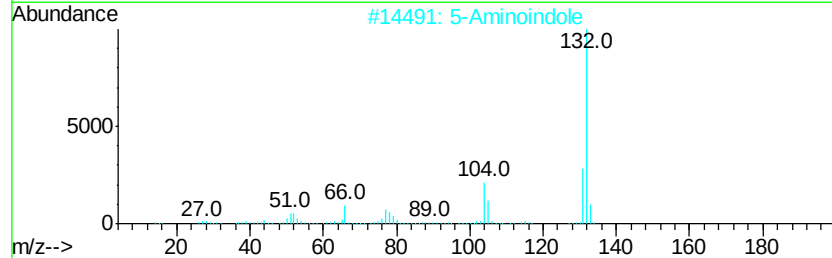
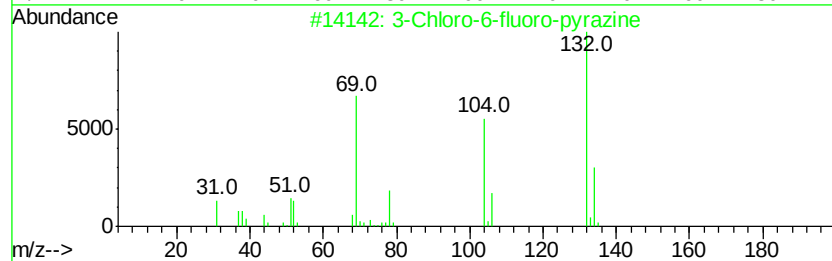
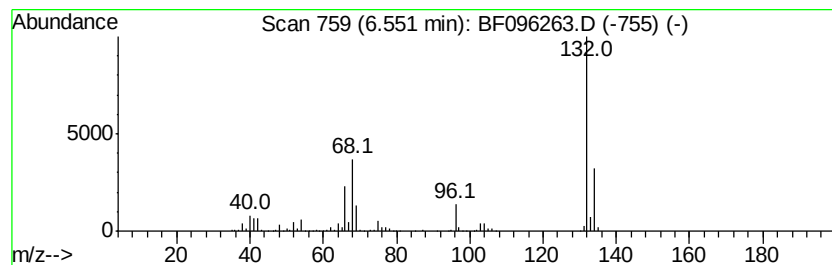
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF062717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 unknown6.55 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.55	16.20 ng	727241	1,4-Dichlorobenzene-d4	6.79

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2			5-Aminoindole	132	C8H8N2	005192-03-0	10
3			Tranvlcypromine-propionvl	189	C12H15NO	1000123-86-3	10
4			1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9
5			4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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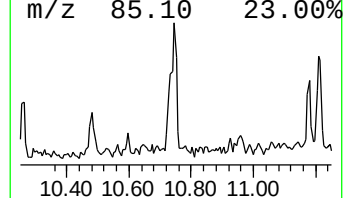
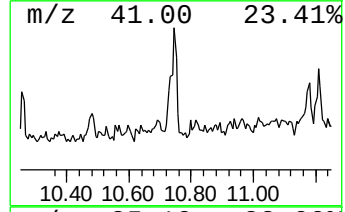
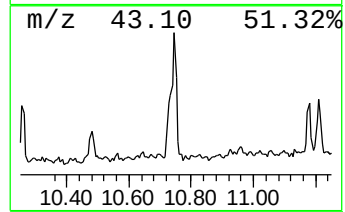
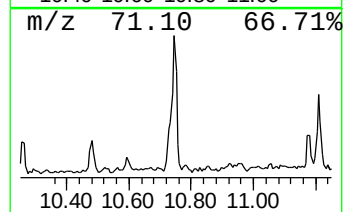
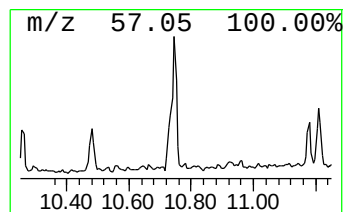
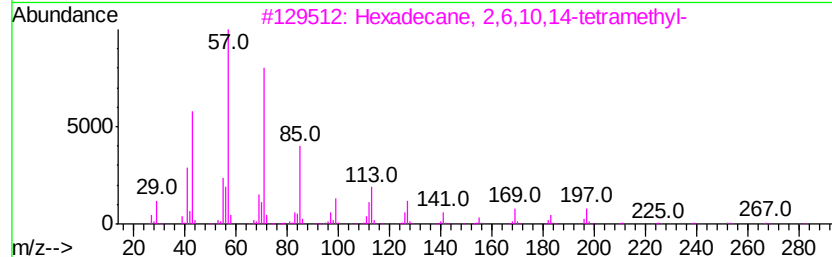
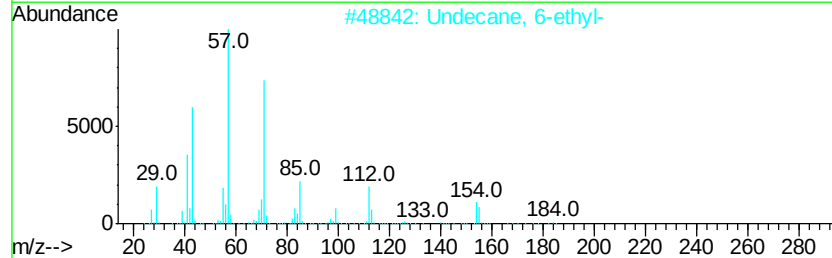
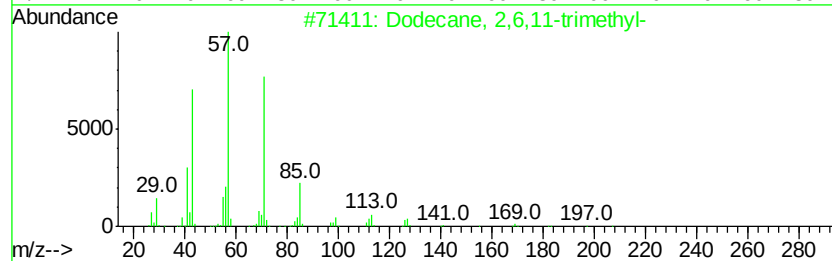
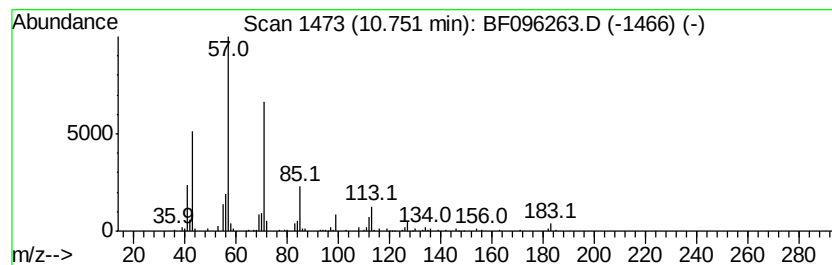
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF062717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Dodecane, 2,6,11-trimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.75	4.52 ng	227305	Phenanthrene-d10	11.30

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	90	
2	Undecane, 6-ethyl-	184	C13H28	017312-60-6	83	
3	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	78	
4	Heptane, 3-ethyl-5-methyl-	142	C10H22	052896-90-9	64	
5	Octane, 3,5-dimethyl-	142	C10H22	015869-93-9	64	



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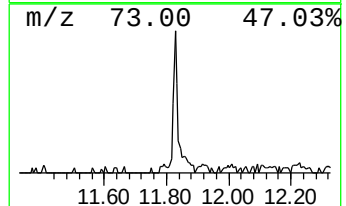
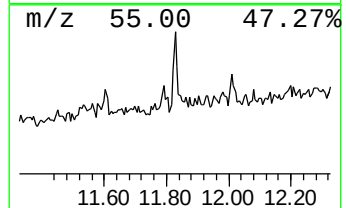
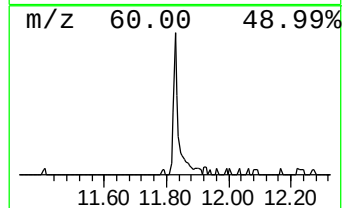
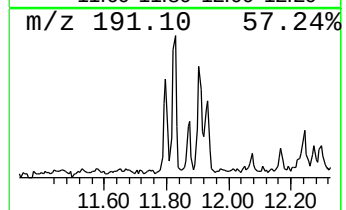
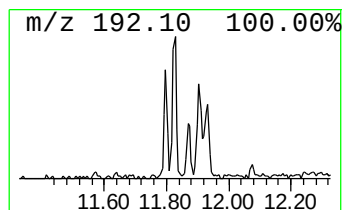
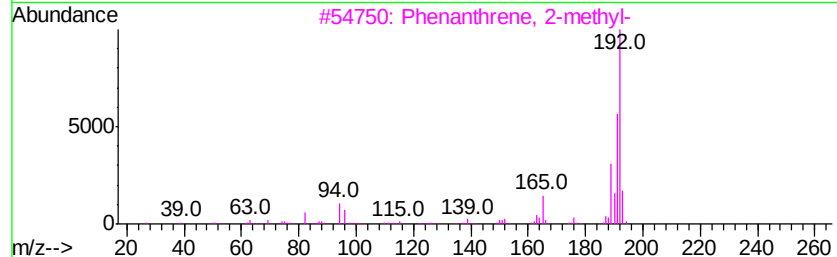
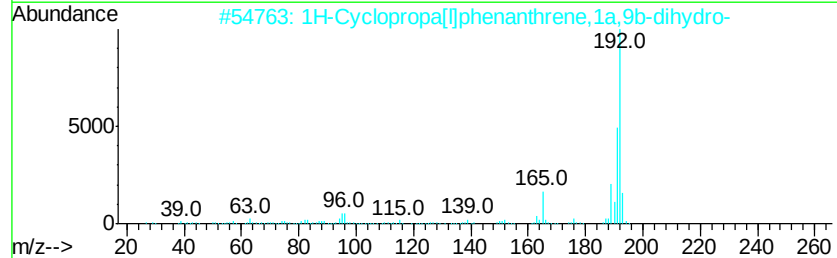
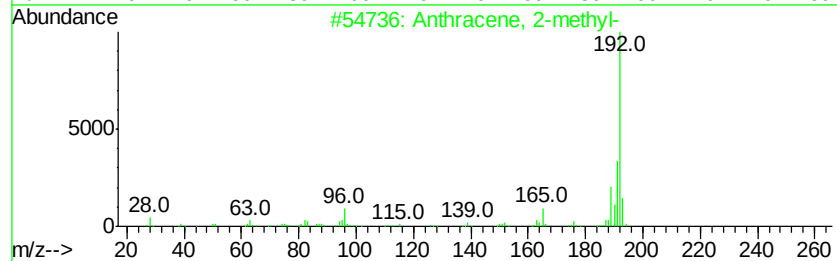
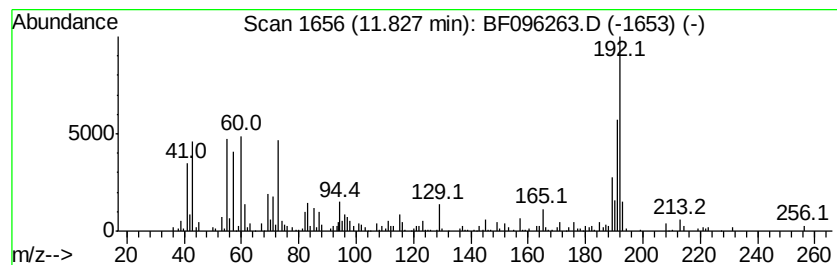
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Anthracene, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.83	2.99 ng	150484	Phenanthrene-d10	11.30

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Anthracene, 2-methyl-	192	C15H12	000613-12-7	86
2		1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	78
3		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	62
4		1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	60
5		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	60



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062817\
Data File : BF096263.D
Acq On : 28 Jun 2017 21:36
Operator : SJ/MA
Sample : I3852-16 5X
Misc :
ALS Vial : 22 Sample Multiplier: 1

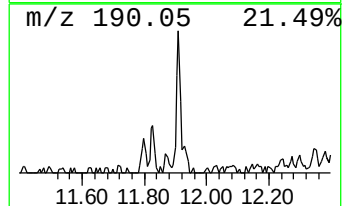
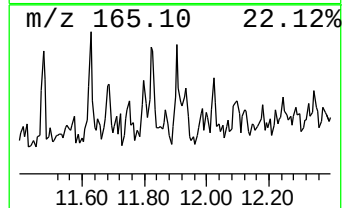
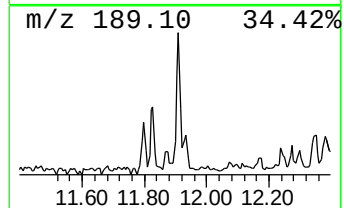
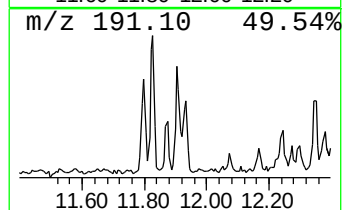
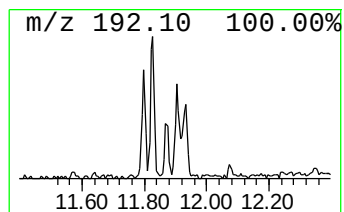
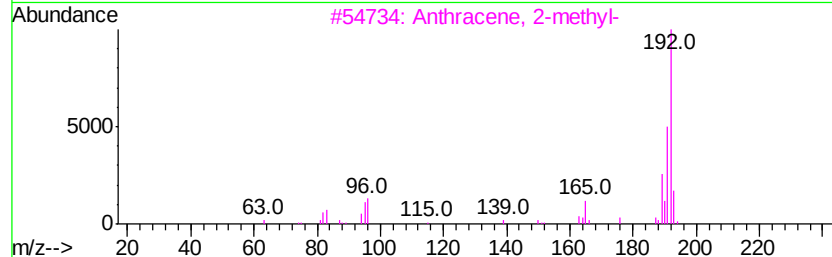
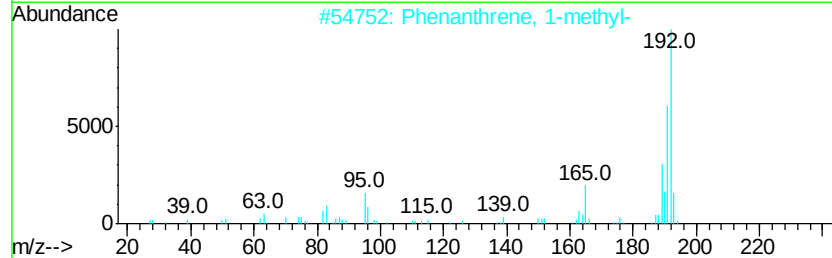
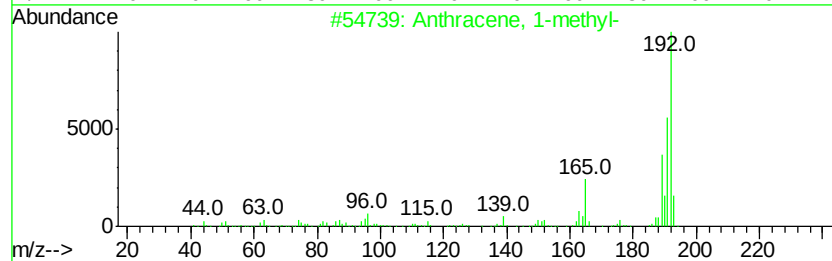
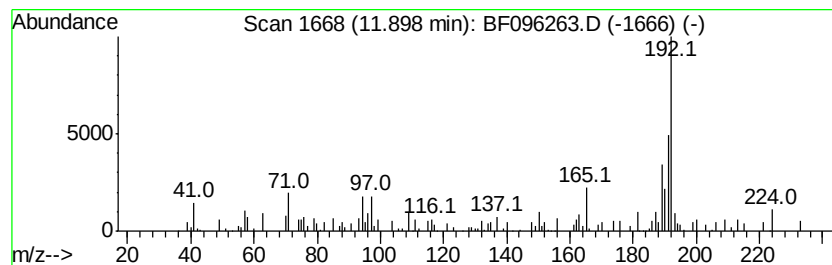
Instrument :
BNA_F
ClientSampleId :
G-38-1.5-3

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF062717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Anthracene, 1-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.90	2.03 ng	102326	Phenanthrene-d10		11.30	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Anthracene, 1-methyl-	192	C15H12	000610-48-0	55
2		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	55
3		Anthracene, 2-methyl-	192	C15H12	000613-12-7	55
4		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	53
5		1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	49



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062817\
Data File : BF096263.D
Acq On : 28 Jun 2017 21:36
Operator : SJ/MA
Sample : I3852-16 5X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
G-38-1.5-3

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF062717.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
2-Pentanone, 4-hy...	4.96	3.2	ng	144362	1	6.79	897824	20.0
unknown6.55	6.55	16.2	ng	727241	1	6.79	897824	20.0
Dodecane, 2,6,11-...	10.75	4.5	ng	227305	4	11.30	1006350	20.0
Anthracene, 2-met...	11.83	3.0	ng	150484	4	11.30	1006350	20.0
Anthracene, 1-met...	11.90	2.0	ng	102326	4	11.30	1006350	20.0