

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061721\
 Data File : BF124364.D
 Acq On : 17 Jun 2021 15:45
 Operator : JU/CG
 Sample : M2736-01 2X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PL-01-061621

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

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF124364.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.404	562	564	572	rBV	114803	131895	42.71%	3.578%
2	6.428	736	738	752	rBB	135998	156848	50.79%	4.254%
3	6.781	795	798	803	rVB	262470	212257	68.73%	5.757%
4	7.345	891	894	898	rBV	92812	72496	23.47%	1.966%
5	7.381	898	900	903	rVB	6107	4704	1.52%	0.128%
6	7.816	971	974	979	rBV3	3277	4186	1.36%	0.114%
7	8.004	998	1006	1007	rBV3	6150	10241	3.32%	0.278%
8	8.069	1011	1017	1020	rVV	341158	283862	91.91%	7.700%
9	8.128	1026	1027	1030	rVB	14565	9100	2.95%	0.247%
10	8.275	1048	1052	1057	rVB4	6343	7866	2.55%	0.213%
11	8.357	1064	1066	1069	rBV3	9037	9526	3.08%	0.258%
12	8.410	1073	1075	1079	rBV3	7550	7364	2.38%	0.200%
13	8.445	1079	1081	1085	rVB2	7990	9289	3.01%	0.252%
14	8.492	1085	1089	1091	rBV2	19726	22222	7.20%	0.603%
15	8.575	1100	1103	1106	rBV3	6735	6924	2.24%	0.188%
16	8.616	1106	1110	1111	rBV3	4883	5717	1.85%	0.155%
17	8.663	1115	1118	1120	rVV	26838	23988	7.77%	0.651%
18	8.722	1126	1128	1130	rBV3	6553	6552	2.12%	0.178%
19	8.757	1130	1134	1137	rVB3	10010	13164	4.26%	0.357%
20	8.822	1141	1145	1146	rBV3	3929	5036	1.63%	0.137%
21	8.875	1150	1154	1155	rBV4	3820	4421	1.43%	0.120%
22	8.886	1155	1156	1160	rVB4	9115	7874	2.55%	0.214%
23	8.939	1163	1165	1168	rBV3	7560	8867	2.87%	0.241%
24	9.022	1176	1179	1183	rVB4	9517	10357	3.35%	0.281%
25	9.057	1183	1185	1188	rVB4	8747	8011	2.59%	0.217%
26	9.086	1188	1190	1195	rBV3	26455	26465	8.57%	0.718%
27	9.139	1196	1199	1203	rVB	226769	189959	61.51%	5.153%
28	9.228	1208	1214	1217	rVV	36712	38181	12.36%	1.036%
29	9.269	1217	1221	1223	rVV3	7739	10481	3.39%	0.284%
30	9.292	1223	1225	1228	rVV2	12020	12124	3.93%	0.329%
31	9.339	1230	1233	1236	rBV3	7243	7372	2.39%	0.200%
32	9.416	1244	1246	1249	rBV3	5358	4417	1.43%	0.120%
33	9.463	1249	1254	1255	rBV5	8012	9046	2.93%	0.245%
34	9.498	1258	1260	1263	rVV3	9410	10940	3.54%	0.297%
35	9.545	1263	1268	1270	rVV4	33547	33668	10.90%	0.913%
36	9.569	1270	1272	1276	rVB4	16299	12625	4.09%	0.342%

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Instrument :
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 ClientSampleId :
 PL-01-061621

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

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	9.604	1276	1278	1280	rBV3	14357	11738	3.80%	0.318%
38	9.686	1289	1292	1297	rBV5	8424	10594	3.43%	0.287%
39	9.728	1297	1299	1301	rVV2	7915	7557	2.45%	0.205%
40	9.751	1301	1303	1308	rVV2	33177	32733	10.60%	0.888%

41	9.792	1308	1310	1312	rVV2	8274	8407	2.72%	0.228%
42	9.822	1312	1315	1319	rVV	372774	308841	100.00%	8.377%
43	9.857	1319	1321	1323	rVB3	12498	9143	2.96%	0.248%
44	9.933	1329	1334	1336	rVB5	5489	5649	1.83%	0.153%
45	9.980	1340	1342	1345	rVV4	4938	5575	1.81%	0.151%

46	10.016	1345	1348	1354	rVV7	5771	14594	4.73%	0.396%
47	10.075	1355	1358	1362	rVB5	7379	11122	3.60%	0.302%
48	10.251	1385	1388	1391	rVB	30382	28419	9.20%	0.771%
49	10.286	1391	1394	1396	rBV3	5143	6139	1.99%	0.167%
50	10.310	1396	1398	1403	rVV4	6227	7687	2.49%	0.209%

51	10.375	1403	1409	1415	rVV7	8700	15697	5.08%	0.426%
52	10.475	1421	1426	1432	rVB3	16797	24760	8.02%	0.672%
53	10.527	1432	1435	1438	rBV3	6276	5145	1.67%	0.140%
54	10.557	1438	1440	1444	rBV4	5549	6092	1.97%	0.165%
55	10.592	1444	1446	1447	rBV	9148	5909	1.91%	0.160%

56	10.616	1447	1450	1453	rVB2	79409	72389	23.44%	1.964%
57	10.727	1466	1469	1474	rBV2	23406	38692	12.53%	1.050%
58	10.869	1488	1493	1495	rVB4	5156	7449	2.41%	0.202%
59	10.922	1498	1502	1504	rVV4	6570	9783	3.17%	0.265%
60	10.951	1505	1507	1511	rVB3	8598	6641	2.15%	0.180%

61	11.174	1540	1545	1547	rBV3	24525	22208	7.19%	0.602%
62	11.204	1547	1550	1553	rVB	18886	18060	5.85%	0.490%
63	11.263	1557	1560	1565	rBV7	10755	17456	5.65%	0.473%
64	11.310	1565	1568	1571	rBV	324620	268143	86.82%	7.273%
65	11.333	1571	1572	1575	rVV	45186	31627	10.24%	0.858%

66	11.363	1575	1577	1579	rVB3	5131	4620	1.50%	0.125%
67	11.386	1579	1581	1584	rBV2	16767	15970	5.17%	0.433%
68	11.533	1600	1606	1608	rBV7	7425	12889	4.17%	0.350%
69	11.551	1608	1609	1614	rVB4	7355	9064	2.93%	0.246%
70	11.604	1614	1618	1621	rBV3	21861	20312	6.58%	0.551%

71	11.704	1629	1635	1637	rBV5	10157	14574	4.72%	0.395%
72	11.757	1643	1644	1646	rBV2	5624	4574	1.48%	0.124%
73	11.774	1646	1647	1651	rVV3	6406	6263	2.03%	0.170%
74	11.816	1651	1654	1656	rBV3	11900	13921	4.51%	0.378%
75	11.839	1656	1658	1661	rVV4	12577	12471	4.04%	0.338%

76	11.892	1664	1667	1669	rBV4	8941	9502	3.08%	0.258%
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Instrument :
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 ClientSampleId :
 PL-01-061621

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

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

77	11.927	1669	1673	1675	rBV3	22599	22499	7.28%	0.610%
78	12.010	1683	1687	1689	rBV4	24011	23447	7.59%	0.636%
79	12.145	1708	1710	1713	rVB3	10143	11052	3.58%	0.300%
80	12.174	1713	1715	1717	rBV3	8445	8009	2.59%	0.217%
81	12.210	1719	1721	1724	rVB4	5551	5210	1.69%	0.141%
82	12.239	1724	1726	1728	rBV2	10104	8147	2.64%	0.221%
83	12.292	1733	1735	1737	rVB3	10706	8054	2.61%	0.218%
84	12.363	1743	1747	1749	rBV5	21751	27690	8.97%	0.751%
85	12.392	1749	1752	1753	rVV2	41682	42724	13.83%	1.159%
86	12.410	1753	1755	1758	rVB3	54205	48635	15.75%	1.319%
87	12.527	1771	1775	1778	rBV	121538	100151	32.43%	2.717%
88	12.563	1780	1781	1783	rBV2	9035	7235	2.34%	0.196%
89	12.733	1808	1810	1811	rBV2	22567	18158	5.88%	0.493%
90	12.757	1811	1814	1820	rVB	106052	128879	41.73%	3.496%
91	12.839	1826	1828	1829	rVB2	9170	4962	1.61%	0.135%
92	12.892	1834	1837	1842	rVV	180178	179156	58.01%	4.860%
93	12.974	1849	1851	1852	rBV2	13322	9277	3.00%	0.252%
94	13.127	1875	1877	1879	rBV3	20244	16475	5.33%	0.447%
95	13.151	1879	1881	1883	rVB3	19807	15179	4.91%	0.412%
96	13.192	1883	1888	1889	rBV5	20740	28010	9.07%	0.760%
97	13.280	1901	1903	1905	rBV3	15846	9138	2.96%	0.248%
98	13.957	2014	2018	2021	rBV2	254487	231886	75.08%	6.290%
99	15.415	2263	2266	2273	rVB2	135894	184043	59.59%	4.992%
100	18.603	2806	2808	2809	rVB2	13395	8222	2.66%	0.223%

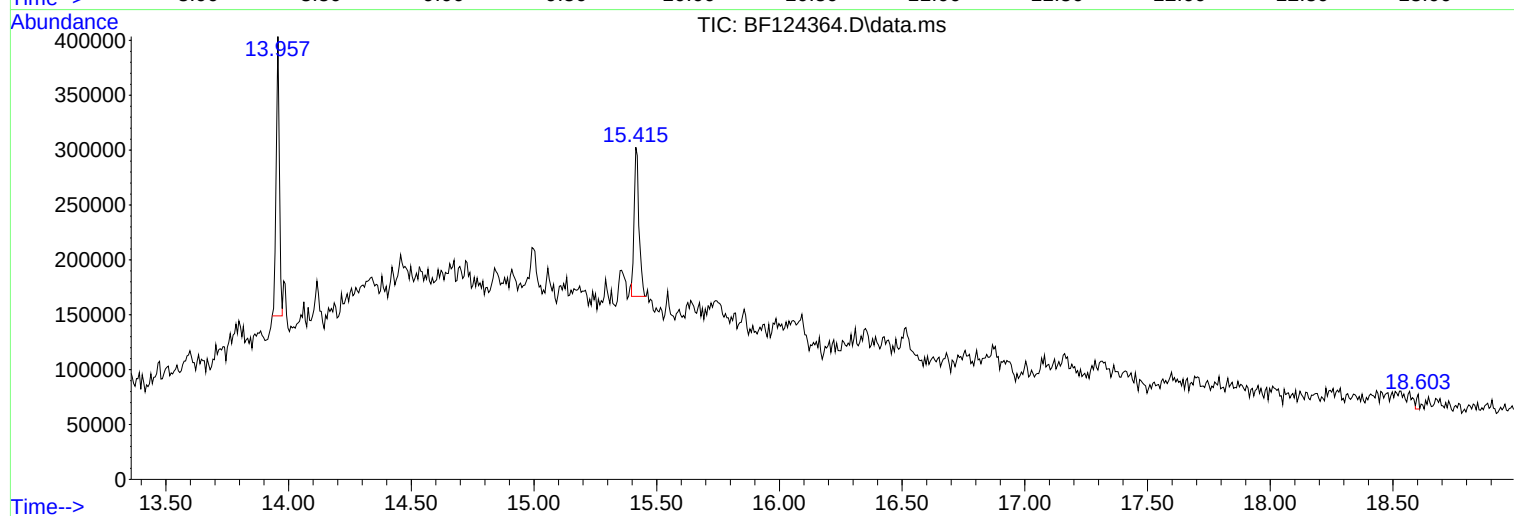
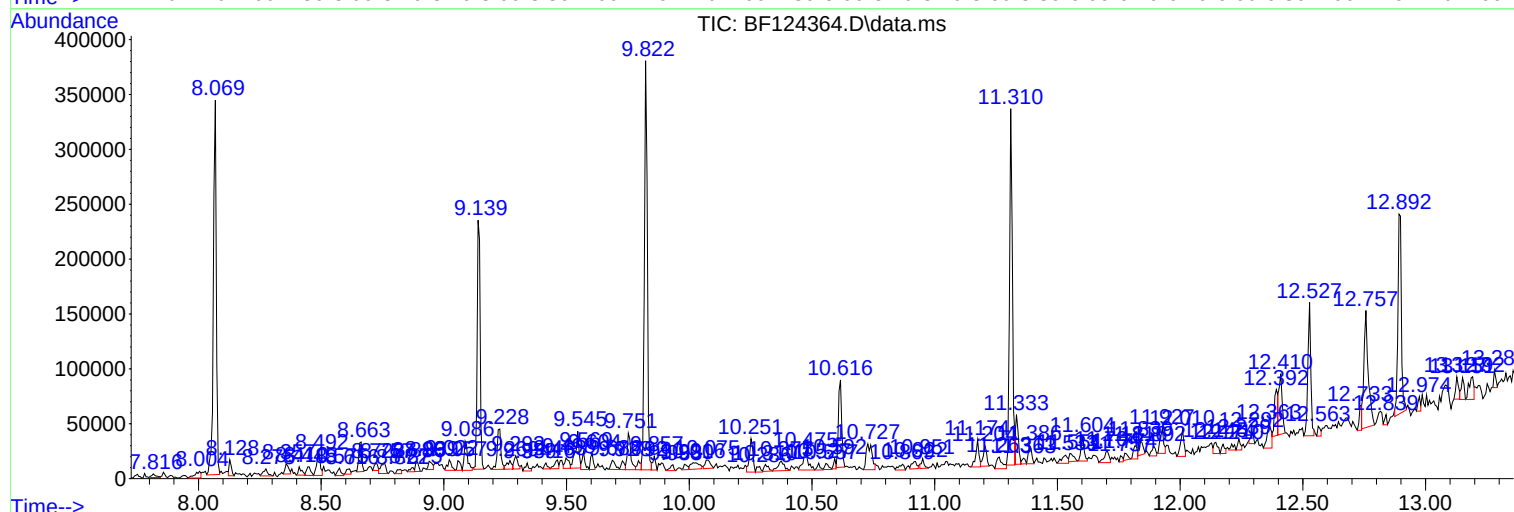
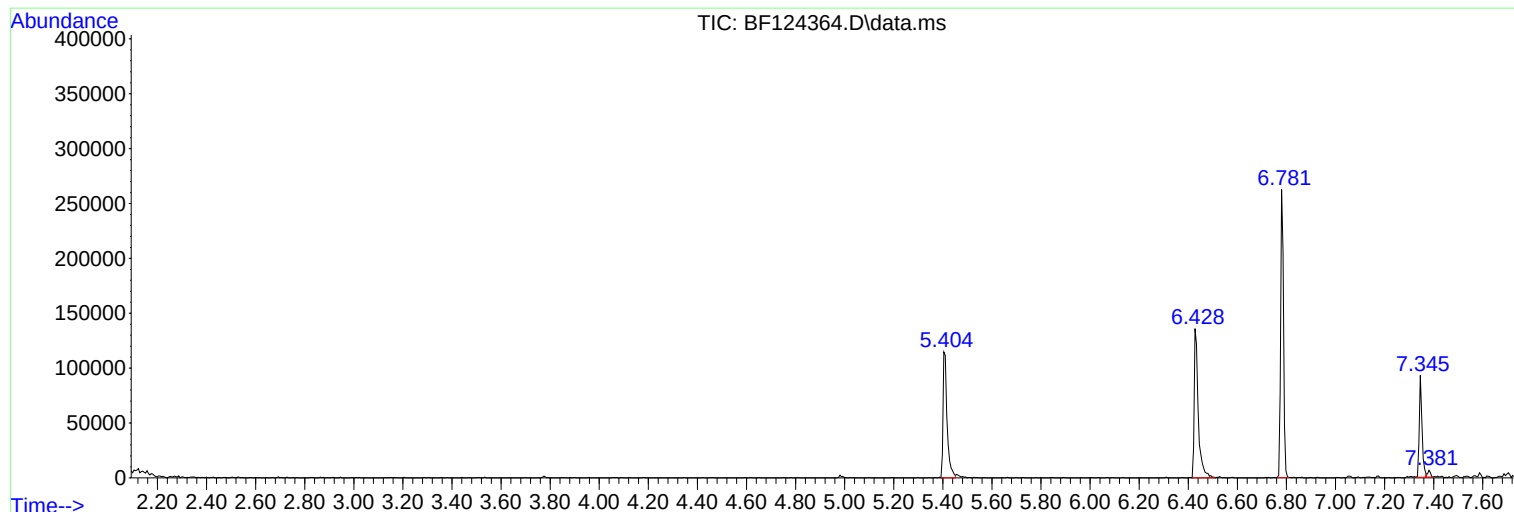
Sum of corrected areas: 3686692

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Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PL-01-061621

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.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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PL-01-061621

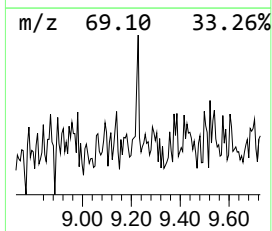
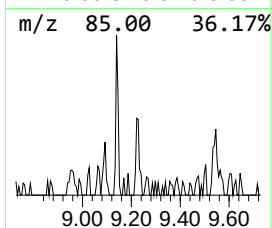
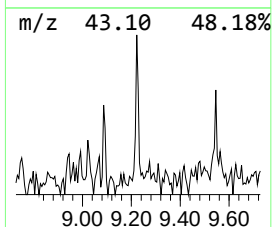
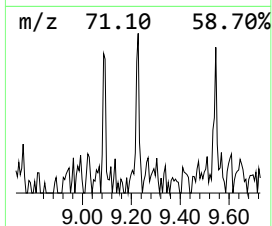
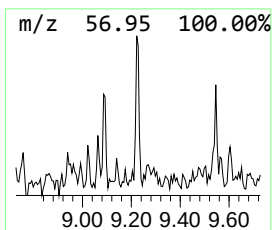
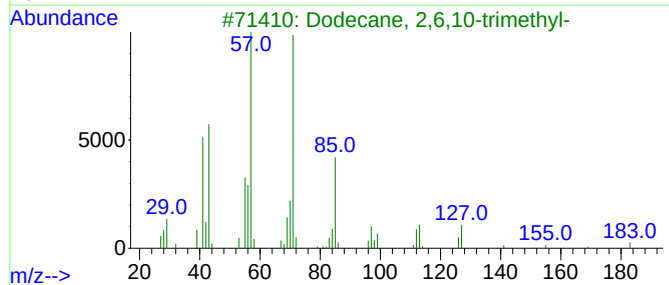
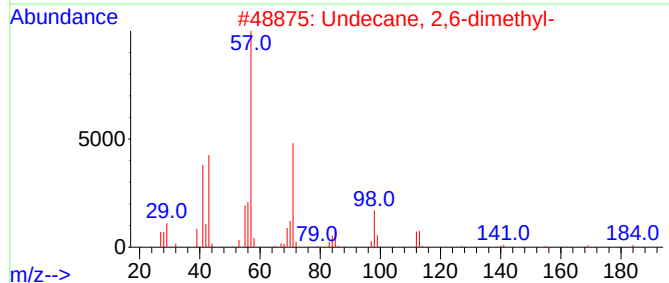
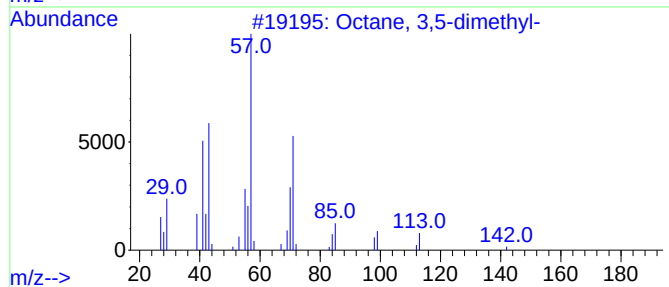
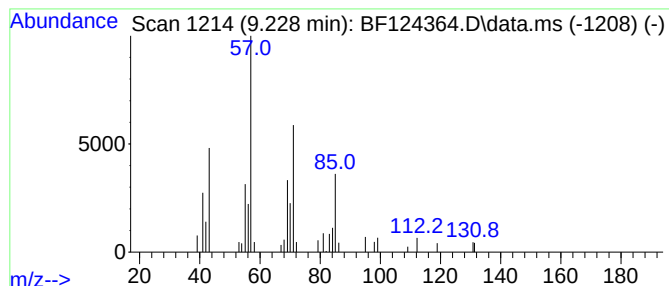
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Octane, 3,5-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.228	2.47 ng	38181	Acenaphthene-d10	9.822

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octane, 3,5-dimethyl-	142	C10H22	015869-93-9	72
2			Undecane, 2,6-dimethyl-	184	C13H28	017301-23-4	64
3			Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	64
4			Hexadecane	226	C16H34	000544-76-3	53
5			Nonane, 3,7-dimethyl-	156	C11H24	017302-32-8	53



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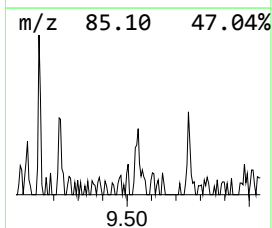
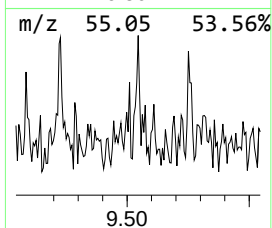
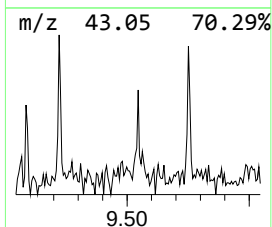
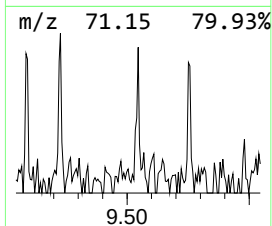
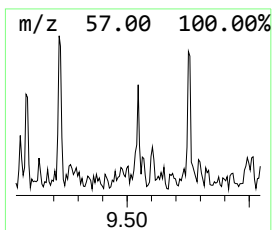
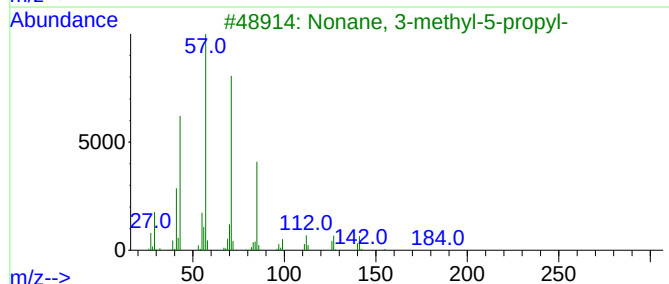
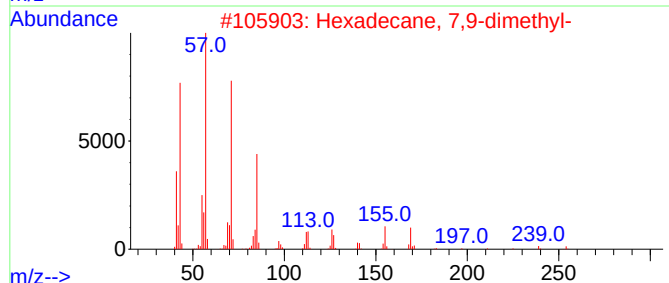
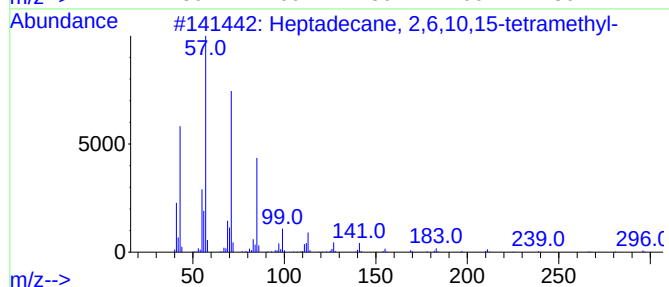
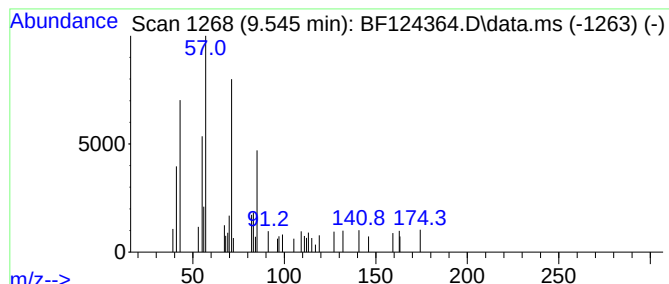
Instrument :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Heptadecane, 2,6,10,15-tetr... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
9.545	2.18 ng	33668	Acenaphthene-d10	9.822	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	64
2	Hexadecane, 7,9-dimethyl-	254	C18H38	021164-95-4	64
3	Nonane, 3-methyl-5-propyl-	184	C13H28	031081-18-2	59
4	Pentadecane	212	C15H32	000629-62-9	59
5	Tetratetracontane	619	C44H90	007098-22-8	59



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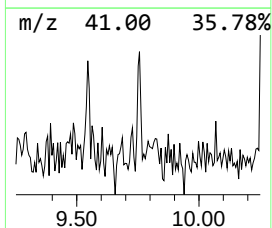
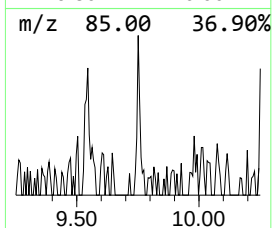
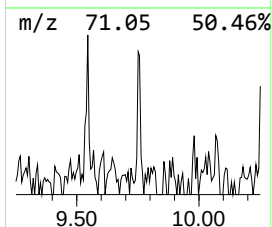
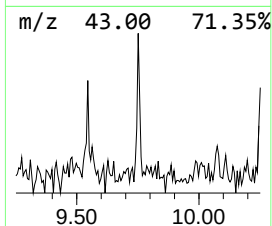
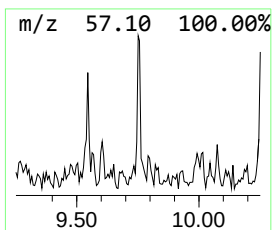
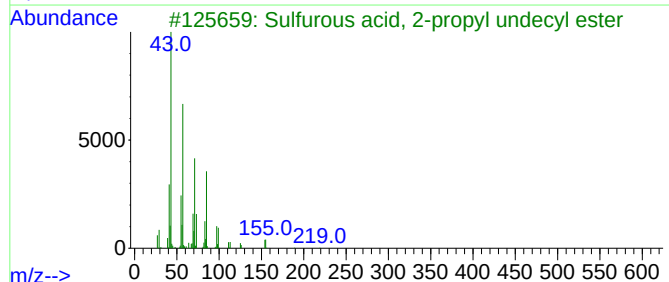
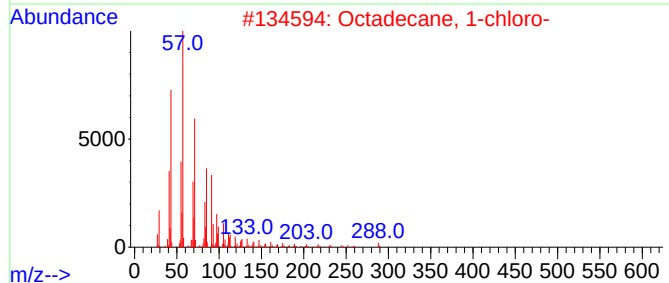
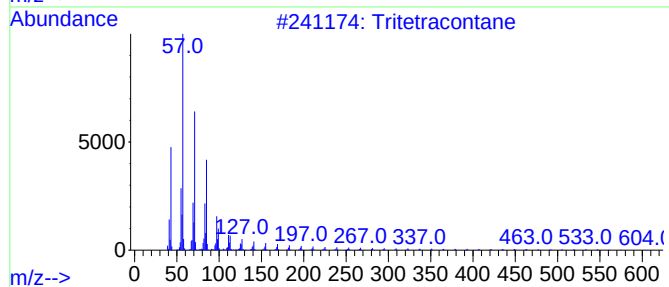
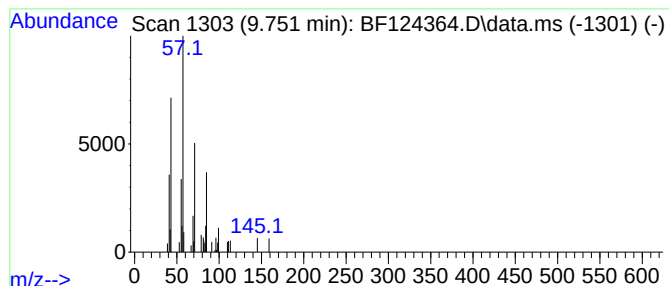
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Tritetracontane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.751	2.12 ng	32733	Acenaphthene-d10	9.822

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tritetracontane	605	C43H88	007098-21-7	72
2			Octadecane, 1-chloro-	288	C18H37Cl	003386-33-2	72
3			Sulfurous acid, 2-propyl undecyl...	278	C14H30O3S	1000309-12-2	64
4			Sulfurous acid, butyl heptadecyl...	376	C21H44O3S	1000309-18-4	64
5			Pentadecane	212	C15H32	000629-62-9	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061721\
Data File : BF124364.D
Acq On : 17 Jun 2021 15:45
Operator : JU/CG
Sample : M2736-01 2X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PL-01-061621

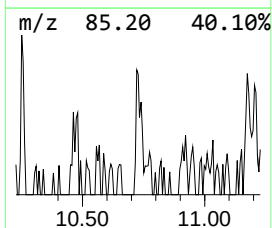
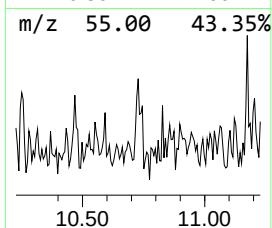
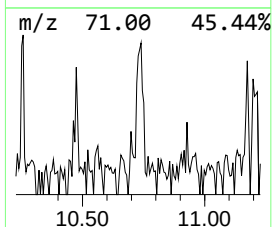
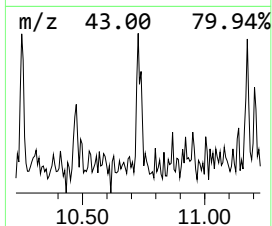
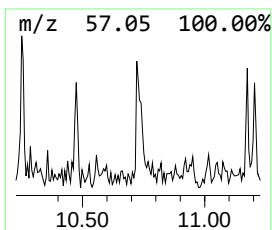
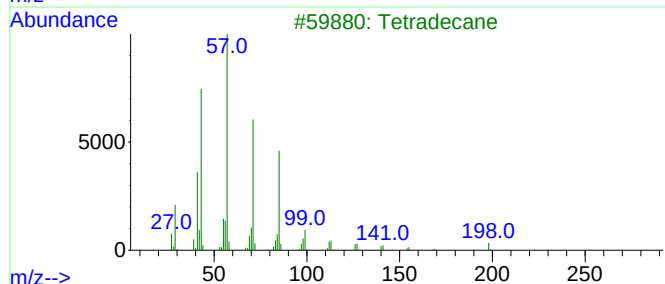
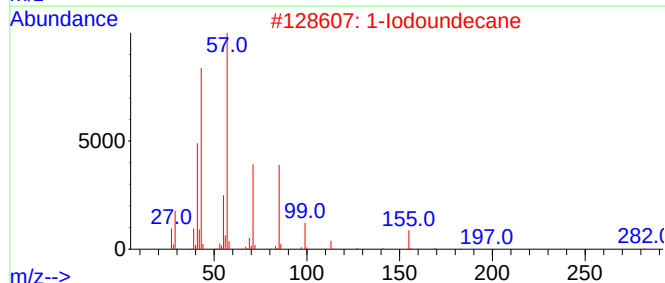
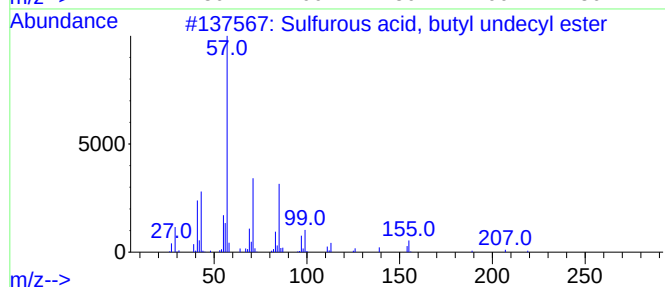
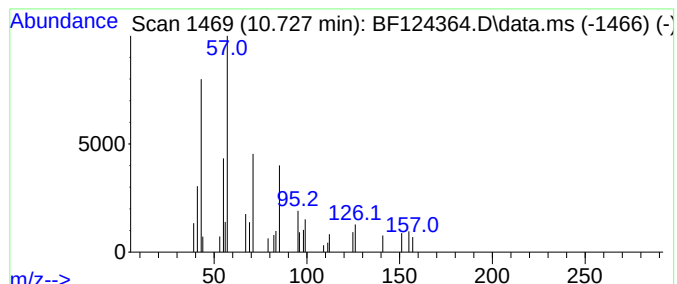
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Sulfurous acid, butyl undec... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.727	2.89 ng	38692	Phenanthrene-d10	11.310

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Sulfurous acid, butyl undecyl ester	292	C15H32O3S	1000309-17-8	53
2			1-Iodoundecane	282	C11H23I	004282-44-4	50
3			Tetradecane	198	C14H30	000629-59-4	50
4			Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	47
5			Heptadecane	240	C17H36	000629-78-7	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061721\
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Misc :
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ClientSampleId :
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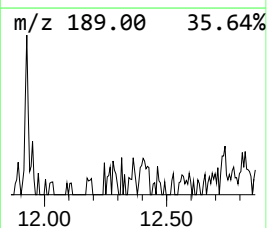
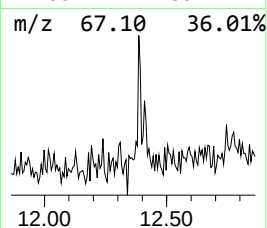
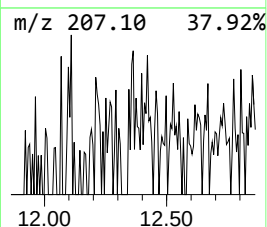
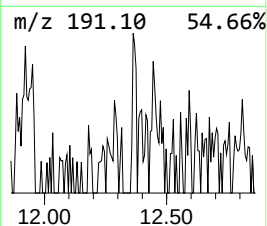
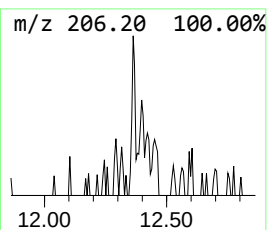
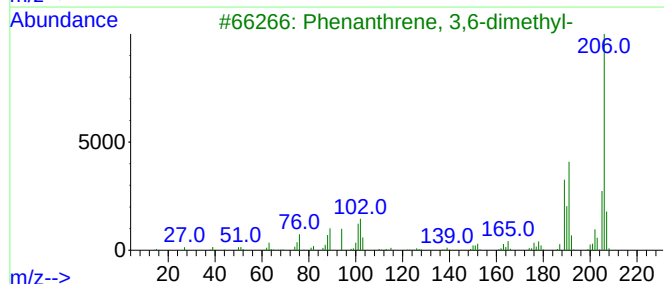
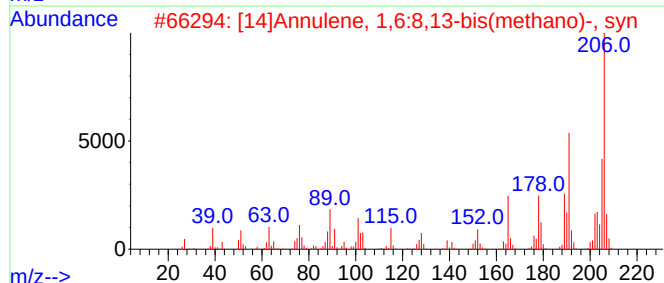
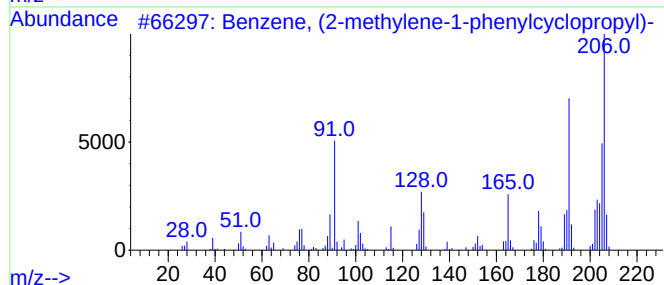
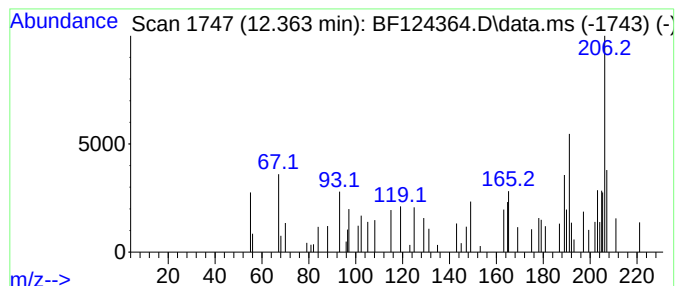
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Benzene, (2-methylene-1-phe... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.363	2.07 ng	27690	Phenanthrene-d10	11.310

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (2-methylene-1-phenylcy...	206	C16H14	025152-47-0	76
2			[14]Annulene, 1,6:8,13-bis(metha...	206	C16H14	085385-68-8	68
3			Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	64
4			1,4-Diphenyl-1,3-butadiene	206	C16H14	000886-65-7	62
5			Naphthalene, 1,2-dihydro-4-phenyl-	206	C16H14	007469-40-1	60



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061721\
Data File : BF124364.D
Acq On : 17 Jun 2021 15:45
Operator : JU/CG
Sample : M2736-01 2X
Misc :
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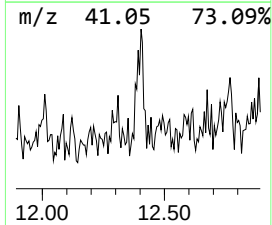
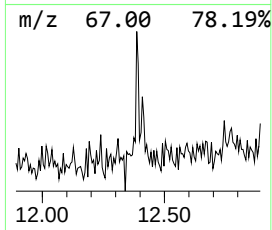
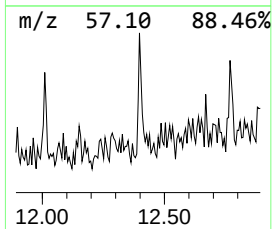
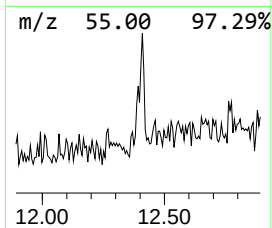
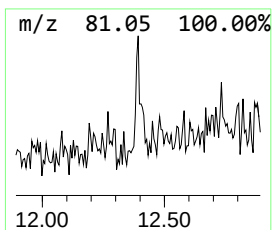
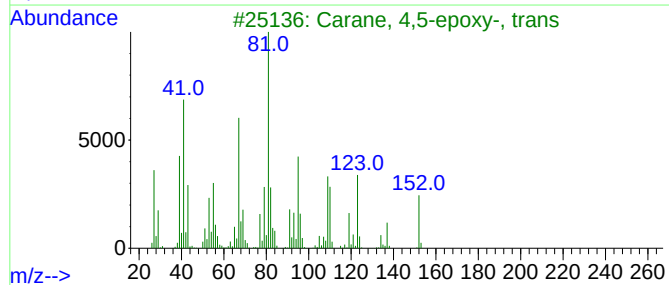
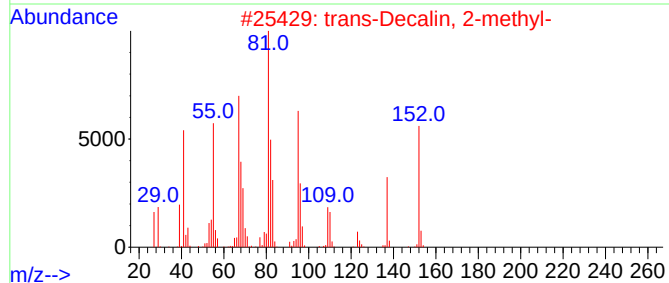
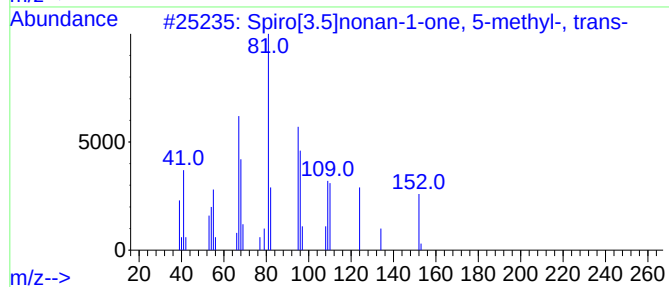
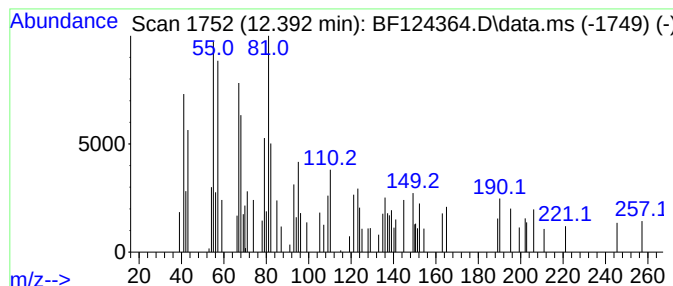
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Spiro[3.5]nonan-1-one, 5-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.392	3.19 ng	42724	Phenanthrene-d10	11.310

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Spiro[3.5]nonan-1-one, 5-methyl-...	152	C10H16O	065147-56-0	58
2			trans-Decalin, 2-methyl-	152	C11H20	1000152-47-3	50
3			Carane, 4,5-epoxy-, trans	152	C10H16O	006909-20-2	46
4			Cyclopentane, 1-methyl-3-(2-meth...	138	C10H18	075873-01-7	46
5			13-Tetradecene-11-yn-1-ol	208	C14H24O	1000131-00-4	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061721\
Data File : BF124364.D
Acq On : 17 Jun 2021 15:45
Operator : JU/CG
Sample : M2736-01 2X
Misc :
ALS Vial : 12 Sample Multiplier: 1

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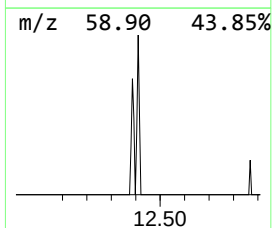
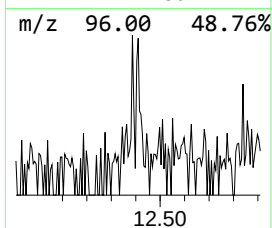
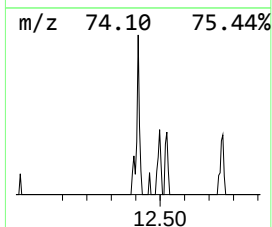
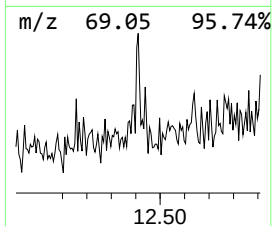
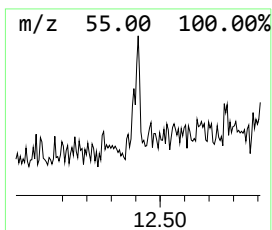
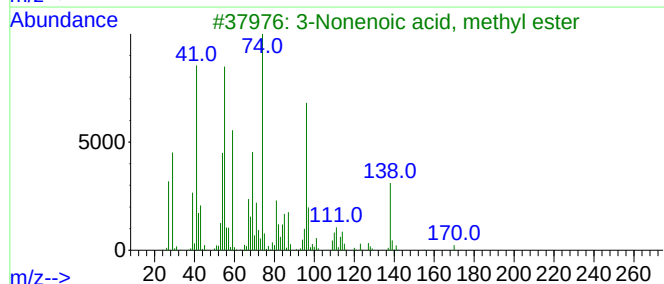
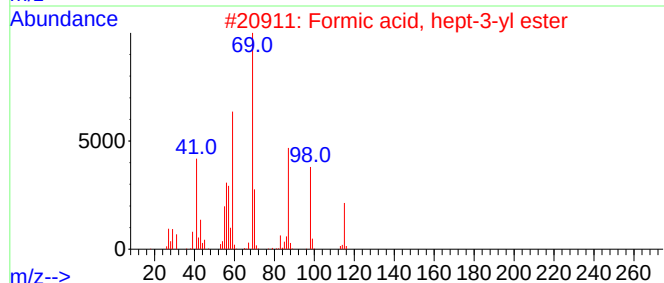
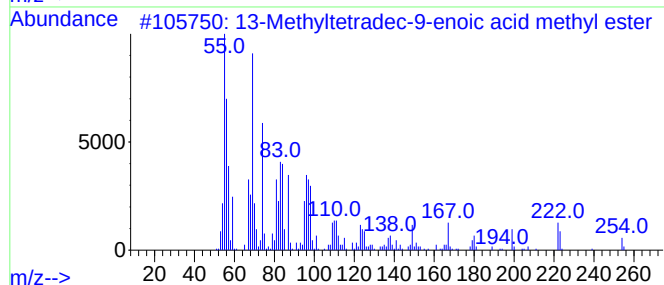
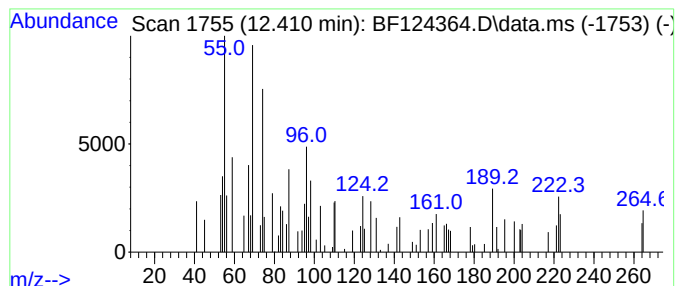
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 13-Methyltetradec-9-enoic a... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.410	3.63 ng	48635	Phenanthrene-d10	11.310

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			13-Methyltetradec-9-enoic acid m...	254	C16H30O2	1000365-89-8	50
2			Formic acid, hept-3-yl ester	144	C8H16O2	1000368-76-5	27
3			3-Nonenoic acid, methyl ester	170	C10H18O2	013481-87-3	27
4			9-Hexadecenoic acid, methyl este...	268	C17H32O2	001120-25-8	27
5			Heptadecanoic acid, 16-methyl-, ...	298	C19H38O2	005129-61-3	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061721\
Data File : BF124364.D
Acq On : 17 Jun 2021 15:45
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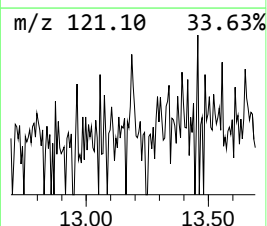
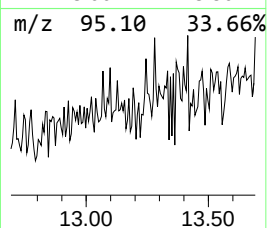
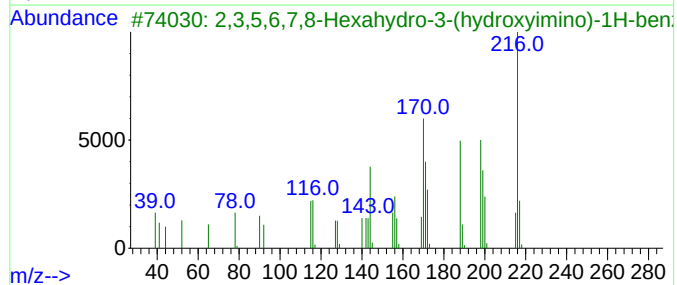
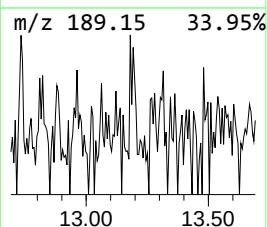
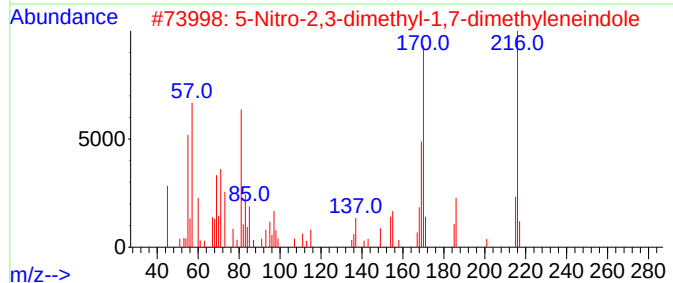
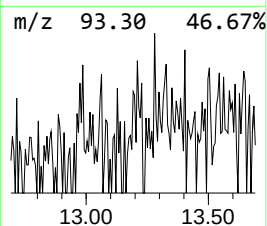
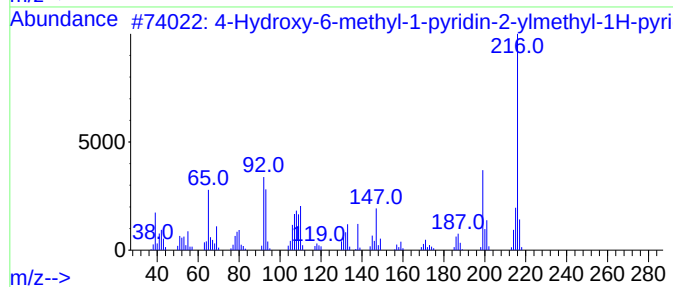
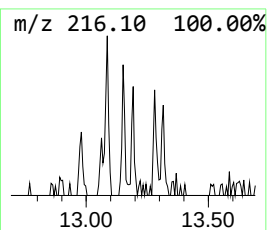
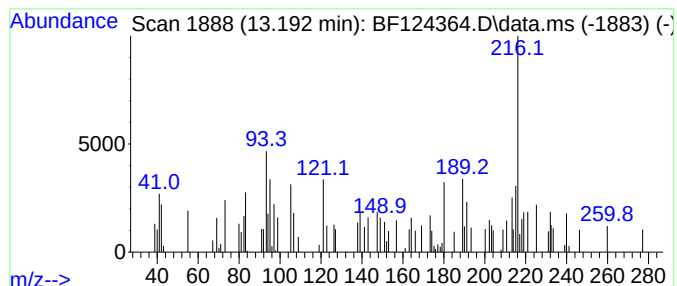
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown13.192 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.192	2.42 ng	28010	Chrysene-d12	13.957

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Hydroxy-6-methyl-1-pyridin-2-y...	216	C12H12N2O2	1000318-25-4	43
2			5-Nitro-2,3-dimethyl-1,7-dimethy...	216	C12H12N2O2	056147-50-3	38
3			2,3,5,6,7,8-Hexahydro-3-(hydroxy...	216	C12H12N2O2	107976-77-2	35
4			9,10-Dihydro-9-oxa-10-phosphaphe...	216	C12H9O2P	035948-25-5	32
5			Hydrazidiamantane	216	C14H20N2	1000138-39-6	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061721\
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Acq On : 17 Jun 2021 15:45
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Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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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
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Heptadecane, 2,...	9.545	2.2	ng	33668	3	9.822	308841	20.0
Tritetracontane	9.751	2.1	ng	32733	3	9.822	308841	20.0
Sulfurous acid,...	10.727	2.9	ng	38692	4	11.310	268143	20.0
Benzene, (2-met...	12.363	2.1	ng	27690	4	11.310	268143	20.0
Spiro[3.5]nonan...	12.392	3.2	ng	42724	4	11.310	268143	20.0
13-Methyltetrad...	12.410	3.6	ng	48635	4	11.310	268143	20.0
unknown13.192	13.192	2.4	ng	28010	5	13.957	231886	20.0