LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF100617\
Data File : BF099170.D
Acq On : 7 Oct 2017 1:51
Operator : SJ/JU
Sample : I5542-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : OWL-C6-GW

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:\HPCHEM1\BNA_F\METHODS\8270-BF092317.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	2.146	5	10	21	rVB	369693	479470	14.02%	1.744%
2	5.087	506	510	521	rVB	303763	373131	10.91%	1.357%
3	5.487	573	578	581	rBV	1791745	1775450	51.93%	6.457%
4	6.493	744	749	752	rBV	1223311	1143509	33.45%	4.159%
5	6.634	769	773	776	rBV	2906235	2854788	83.50%	10.382%
6	6.863	808	812	816	rBV	976716	772657	22.60%	2.810%
7	7.022	834	839	842	rBV	2772096	2507500	73.34%	9.119%
8	7.257	876	879	888	rVB3	43558	53879	1.58%	0.196%
9	7.428	903	908	911	rBV	1958779	2007217	58.71%	7.300%
10	7.545	924	928	932	rVB3	48814	58359	1.71%	0.212%
11	7.604	932	938	941	rBV2	29145	39532	1.16%	0.144%
12	7.781	966	968	974	rVB3	38240	44839	1.31%	0.163%
13	8.039	1009	1012	1015	rBV	50253	57517	1.68%	0.209%
14	8.145	1026	1030	1033	rBV	1198940	1032239	30.19%	3.754%
15	8.386	1069	1071	1075	rBV2	48090	48640	1.42%	0.177%
16	8.516	1091	1093	1096	rVB3	56200	50835	1.49%	0.185%
17	8.598	1101	1107	1109	rBV4	35649	64973	1.90%	0.236%
18	8.704	1123	1125	1128	rVB	63012	55313	1.62%	0.201%
19	8.739	1128	1131	1133	rBV3	60113	68986	2.02%	0.251%
20	8.951	1158	1167	1171	rBV9	77118	182796	5.35%	0.665%
21	8.986	1171	1173	1177	rVB3	74021	93523	2.74%	0.340%
22	9.051	1177	1184	1185	rBV7	51263	75465	2.21%	0.274%
23	9.145	1196	1200	1204	rBV	241589	247855	7.25%	0.901%
24	9.222	1208	1213	1216	rBV	3754928	3418967	100.00%	12.434%
25	9.292	1220	1225	1226	rBV4	39445	62442	1.83%	0.227%
26	9.381	1238	1240	1243	rVB2	96571	85550	2.50%	0.311%
27	9.469	1251	1255	1256	rBV4	53311	65228	1.91%	0.237%
28	9.528	1263	1265	1269	rBV3	88116	74534	2.18%	0.271%
29	9.563	1269	1271	1276	rVB	176961	165515	4.84%	0.602%
30	9.610	1276	1279	1282	rBV3	81789	96415	2.82%	0.351%
31	9.645	1282	1285	1286	rBV3	50102	46421	1.36%	0.169%
32	9.710	1294	1296	1299	rVB2	71925	60897	1.78%	0.221%
33	9.763	1300	1305	1308	rVV4	63920	93845	2.74%	0.341%
34	9.869	1320	1323	1324	rBV	92132	75887	2.22%	0.276%

LSC Area Percent Report

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Data File : BF099170.D
Acq On : 7 Oct 2017 1:51
Operator : SJ/JU
Sample : I5542-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : OWL-C6-GW

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:\HPCHEM1\BNA_F\METHODS\8270-BF092317.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	9.898	1325	1328	1334	rVB2	1311335	1210043	35.39%	4.401%
36	9.969	1338	1340	1344	rBV4	51663	68484	2.00%	0.249%
37	10.057	1353	1355	1361	rVB5	54737	77858	2.28%	0.283%
38	10.110	1361	1364	1367	rVB2	69212	66939	1.96%	0.243%
39	10.151	1367	1371	1372	rBV3	34859	46412	1.36%	0.169%
40	10.204	1377	1380	1383	rBV3	66101	112551	3.29%	0.409%
41	10.257	1387	1389	1395	rVB	219640	186739	5.46%	0.679%
42	10.328	1399	1401	1404	rVB2	63829	53863	1.58%	0.196%
43	10.510	1429	1432	1435	rBV	262406	250668	7.33%	0.912%
44	10.633	1451	1453	1455	rBV2	44598	42873	1.25%	0.156%
45	10.692	1458	1463	1468	rVB	1783702	1825044	53.38%	6.637%
46	10.810	1481	1483	1489	rVB3	137784	176593	5.17%	0.642%
47	10.880	1489	1495	1498	rBV5	77181	144525	4.23%	0.526%
48	11.022	1516	1519	1521	rBV	121784	107532	3.15%	0.391%
49	11.128	1535	1537	1539	rBV2	46267	43841	1.28%	0.159%
50	11.186	1545	1547	1553	rVB6	41140	43180	1.26%	0.157%
51	11.363	1574	1577	1578	rBV	59379	59056	1.73%	0.215%
52	11.386	1578	1581	1584	rVV	1085470	888528	25.99%	3.231%
53	11.439	1587	1590	1594	rVV	169734	154857	4.53%	0.563%
54	11.498	1597	1600	1605	rVB	148037	146823	4.29%	0.534%
55	11.904	1667	1669	1673	rVB4	48363	55097	1.61%	0.200%
56	12.969	1845	1850	1853	rBV	2315245	2114098	61.83%	7.689%
57	14.021	2025	2029	2034	rBV	653627	569877	16.67%	2.073%
58	15.492	2274	2279	2290	rBV	514606	659064	19.28%	2.397%
59	17.127	2551	2557	2564	rBV7	24880	57541	1.68%	0.209%

Sum of corrected areas: 27496260

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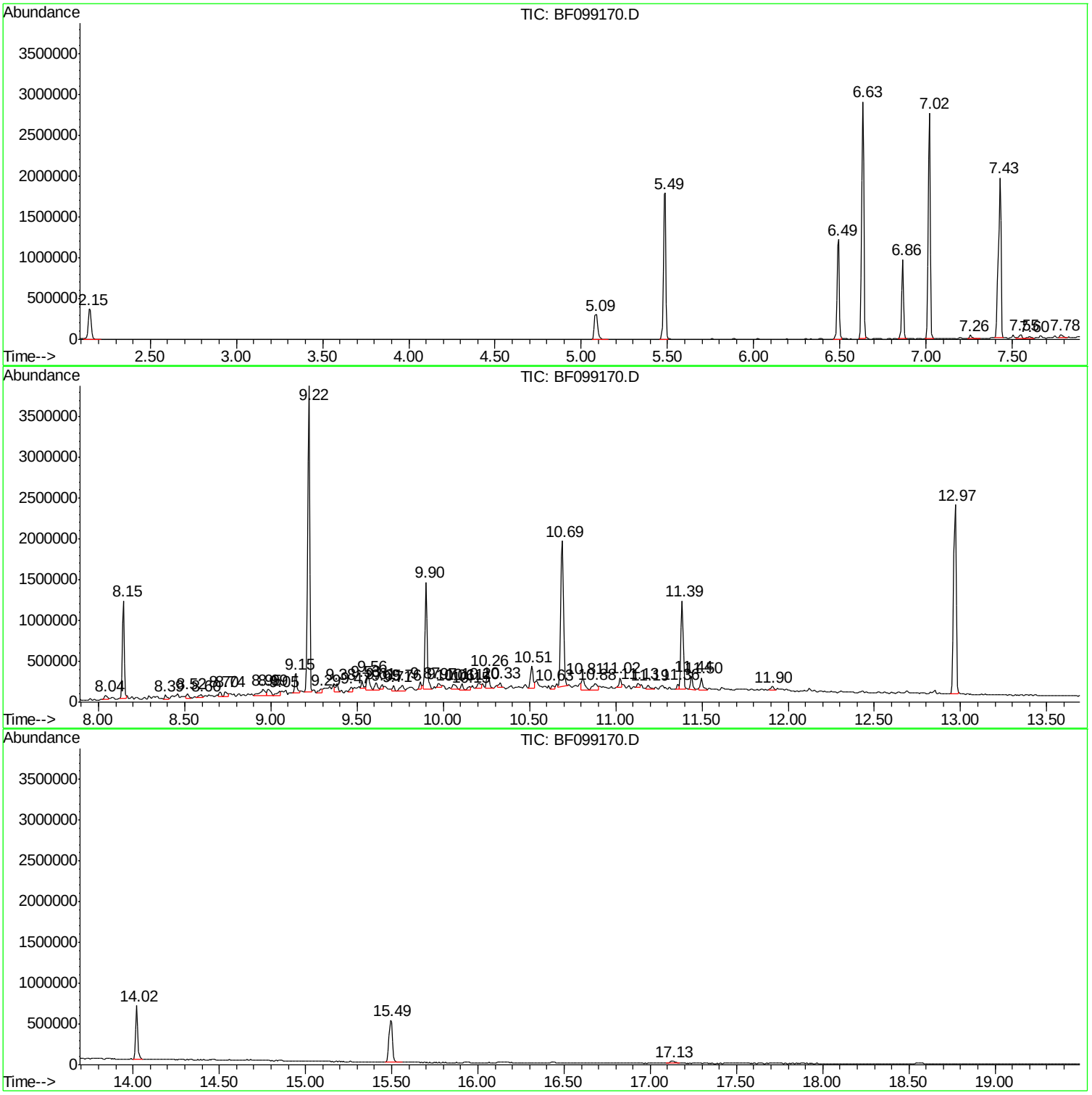
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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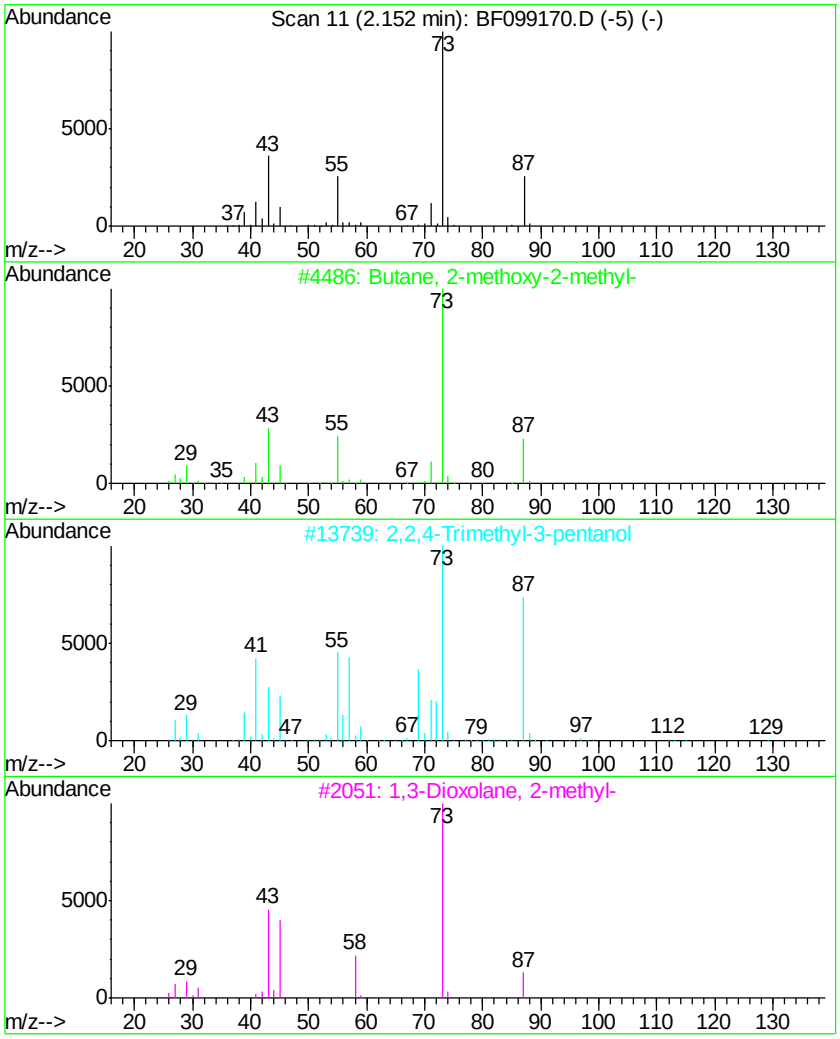
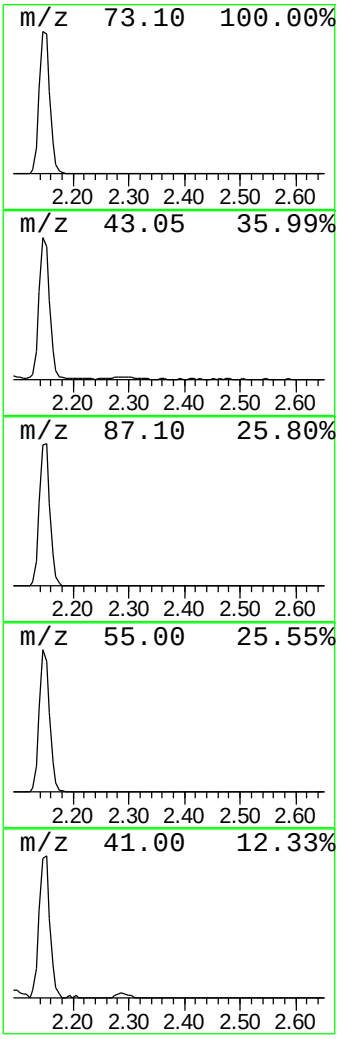
****

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.15	12.41 ng	479470	1,4-Dichlorobenzene-d4	6.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	32
4	1-(2-Methoxyethoxy)-2-methyl-2-p...	162	C8H18O3	1000364-24-4	23
5	Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	14





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Operator : SJ/JU
Sample : I5542-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : OWL-C6-GW

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.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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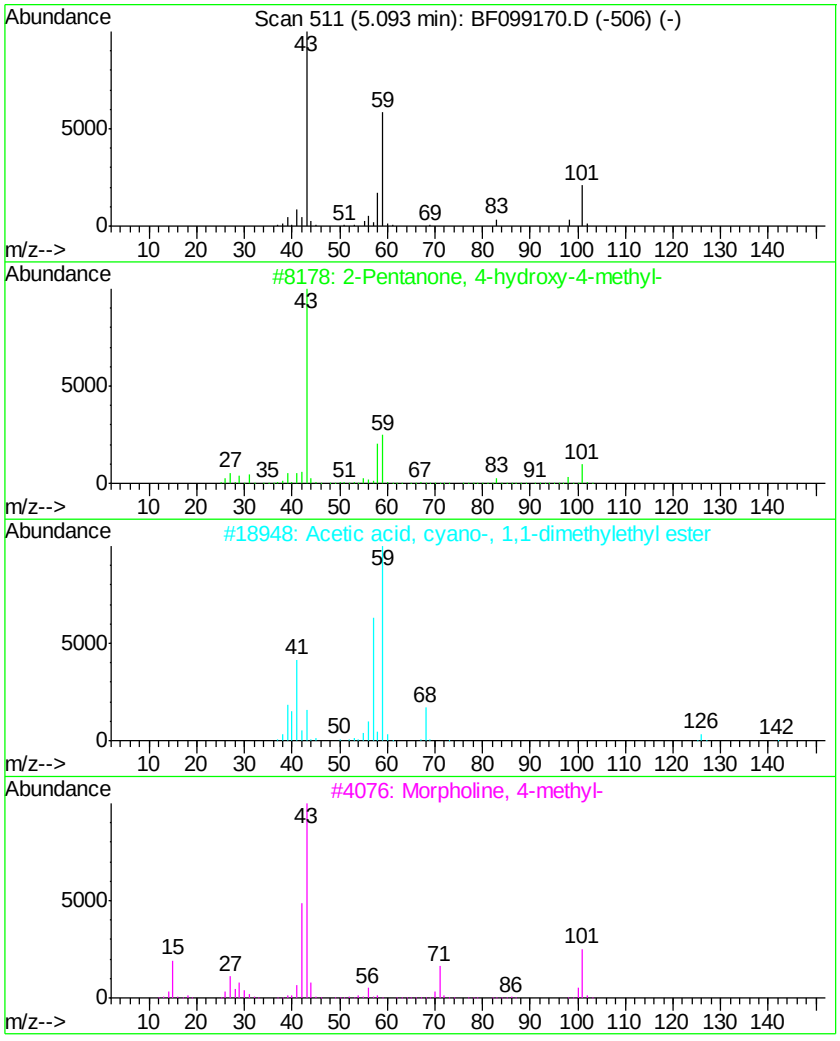
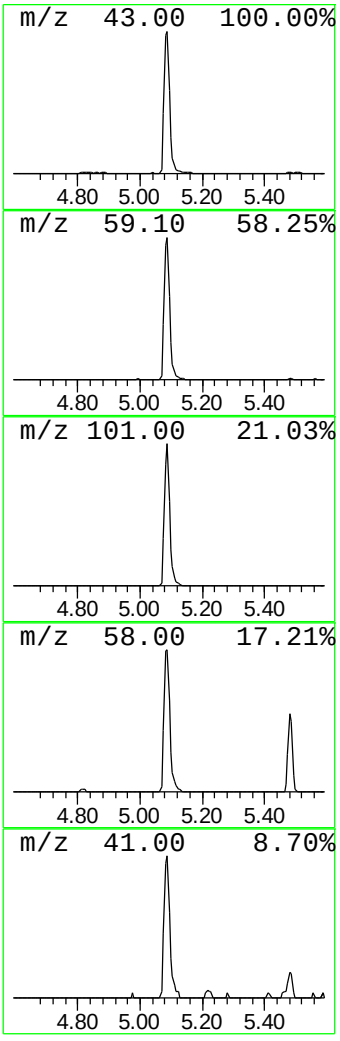
Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.09	9.66 ng	373131	1,4-Dichlorobenzene-d4	6.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17
3	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5	5-Hexen-2-one	98	C6H10O	000109-49-9	9





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Operator : SJ/JU
Sample : I5542-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : OWL-C6-GW

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.M
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TIC Library : C:\DATABASE\NIST11.L
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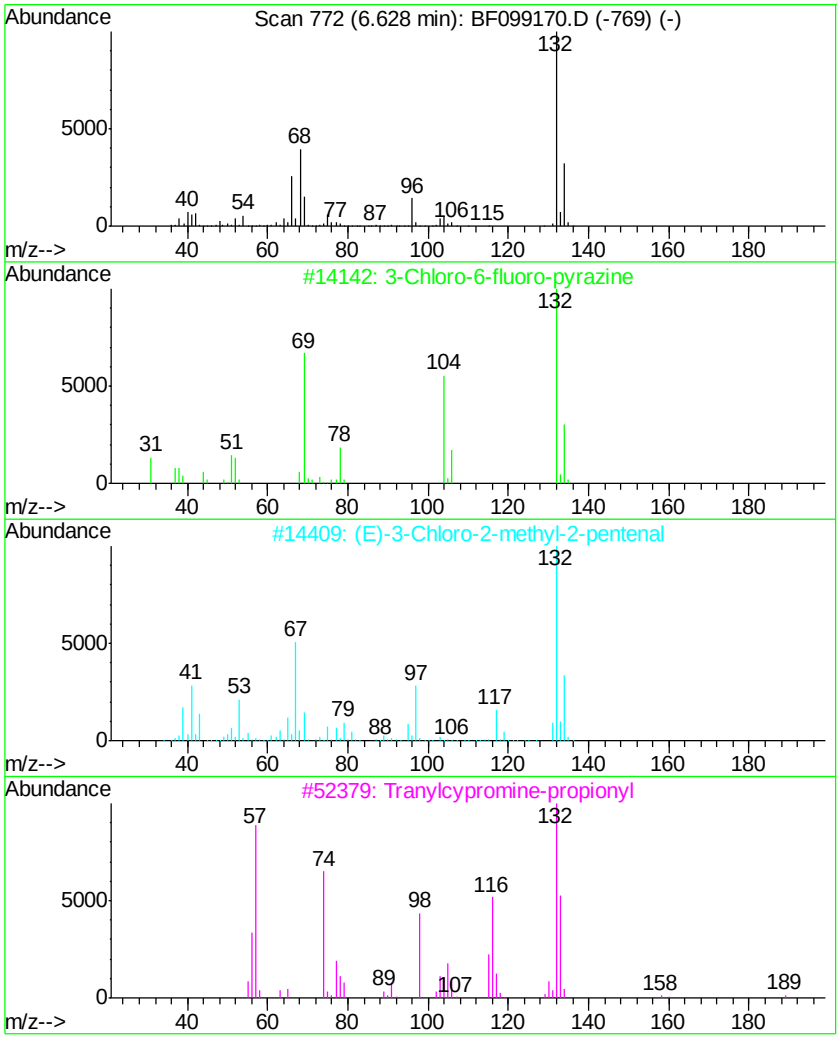
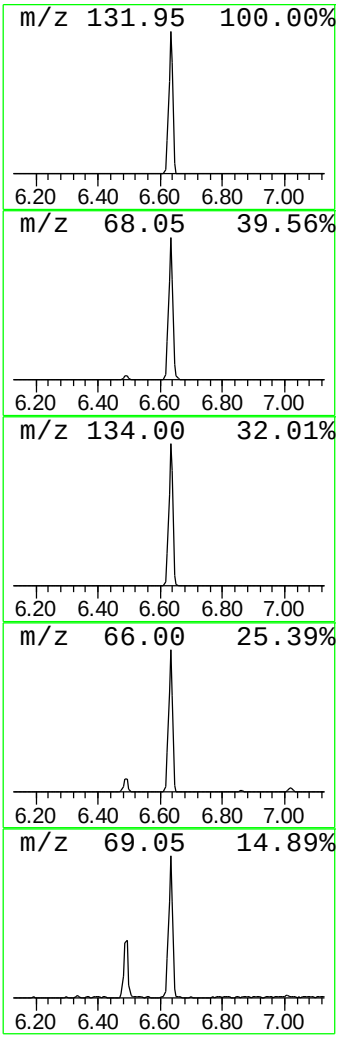
Peak Number 3 unknown6.63 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.63	73.90 ng	2854790	1,4-Dichlorobenzene-d4	6.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
3	Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
4	5-Aminoindole	132	C8H8N2	005192-03-0	12
5	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	10





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ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : OWL-C6-GW

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

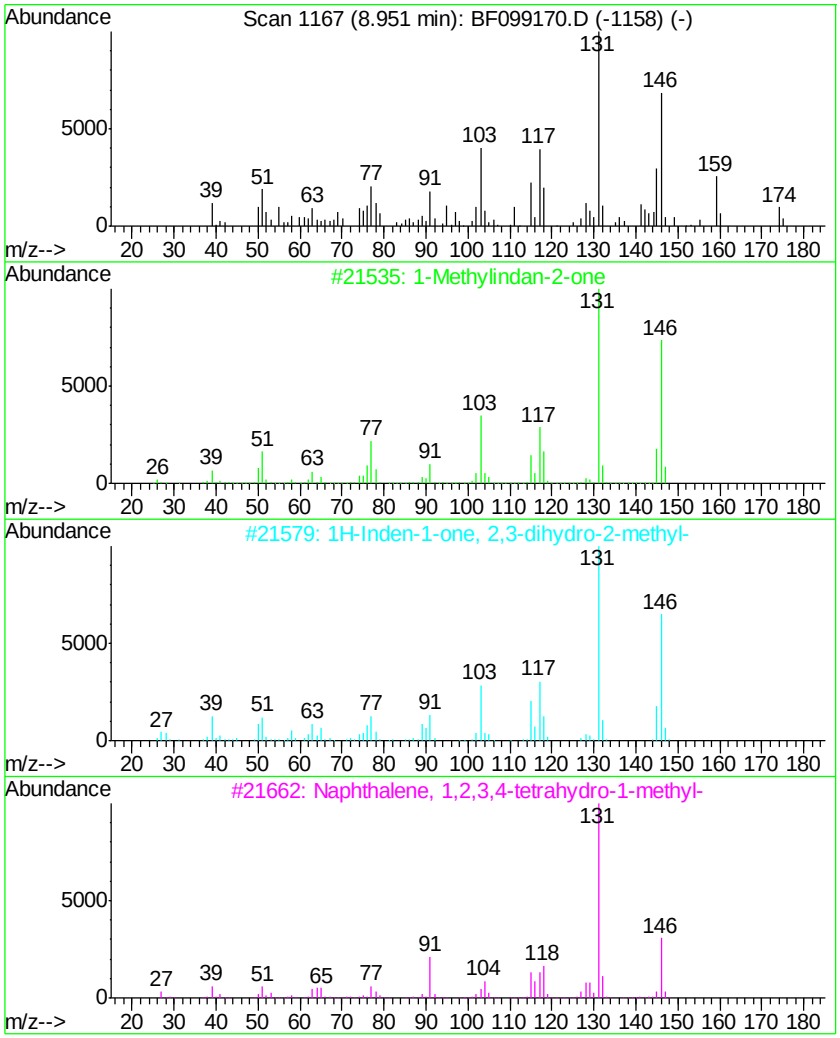
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

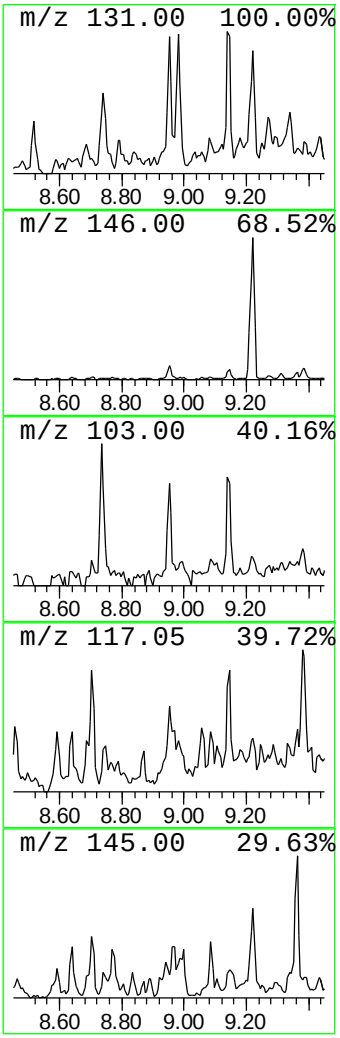
****

Peak Number 5 1-Methylindan-2-one Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.95	3.54 ng	182796	Naphthalene-d8	8.15

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Methylindan-2-one	146	C10H10O	035587-60-1	89
2	1H-Inden-1-one, 2,3-dihydro-2-me...	146	C10H10O	017496-14-9	72
3	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	46
4	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	45
5	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	43





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

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

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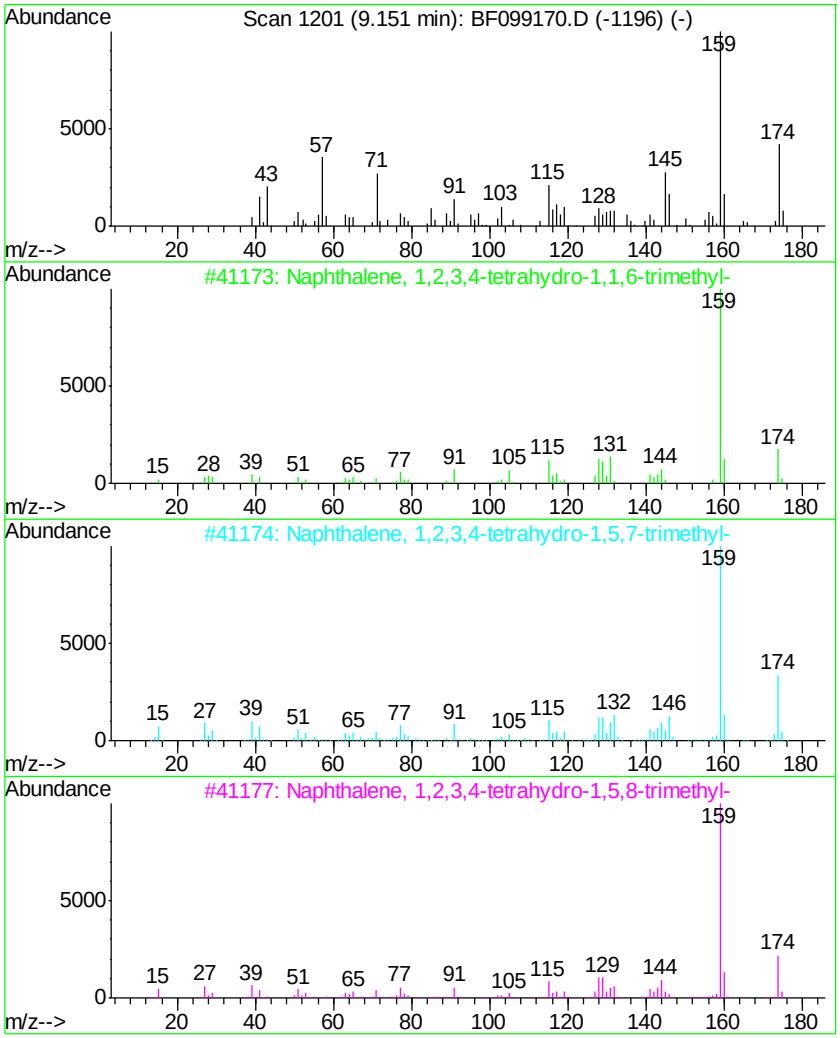
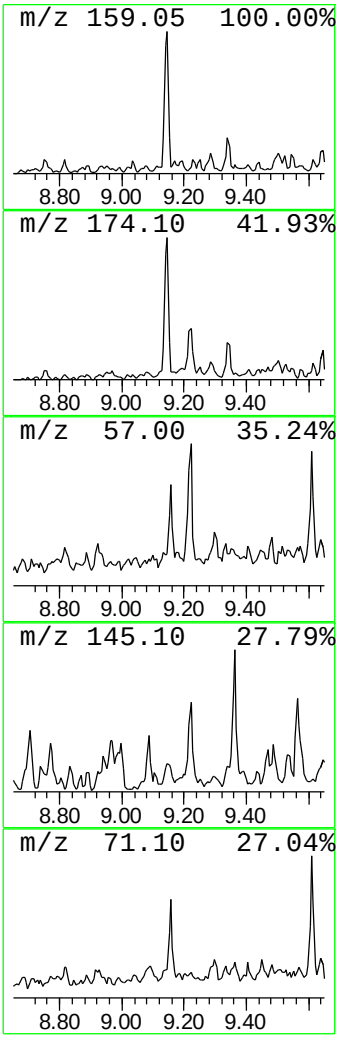
Peak Number 6 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.15	4.10 ng	247855	Acenaphthene-d10	9.90

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	174	C13H18	000475-03-6	93
2	Naphthalene, 1,2,3,4-tetrahydro-...	174	C13H18	021693-55-0	87
3	Naphthalene, 1,2,3,4-tetrahydro-...	174	C13H18	021693-51-6	87
4	1H-Inden-1-one, 2,3-dihydro-3,3,...	174	C12H14O	054484-71-8	83
5	1H-Indene, 2,3-dihydro-1,1,5,6-t...	174	C13H18	000942-43-8	81





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

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

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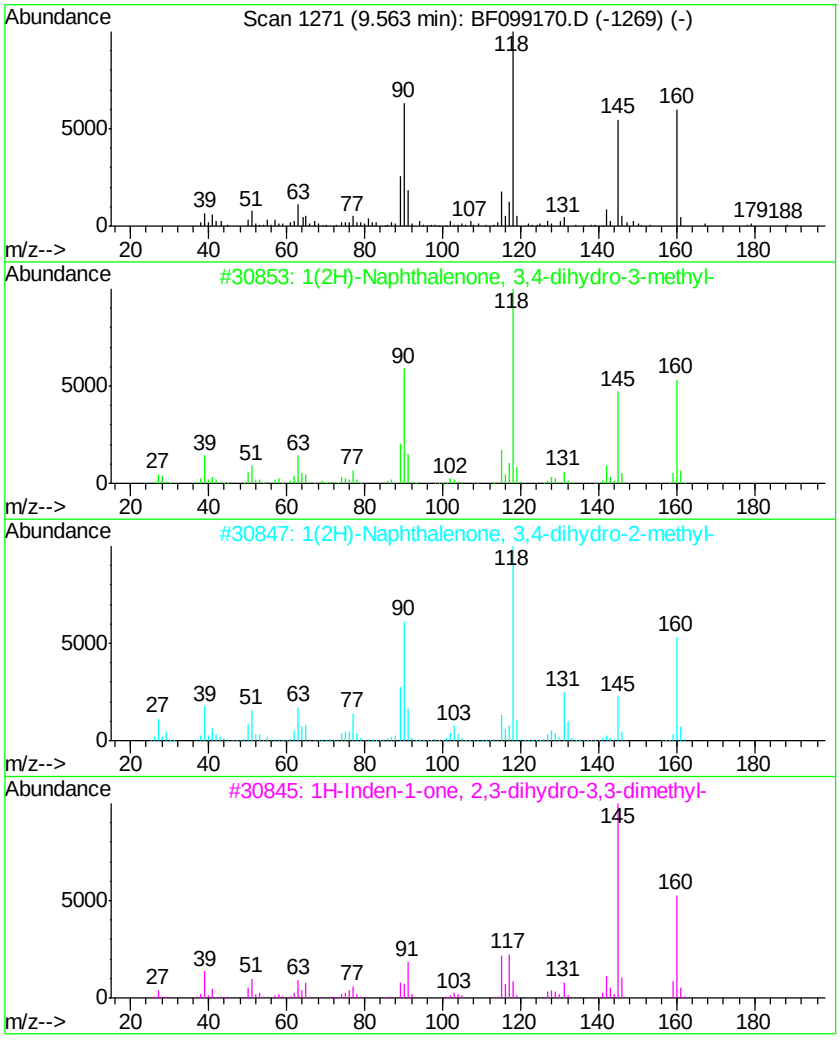
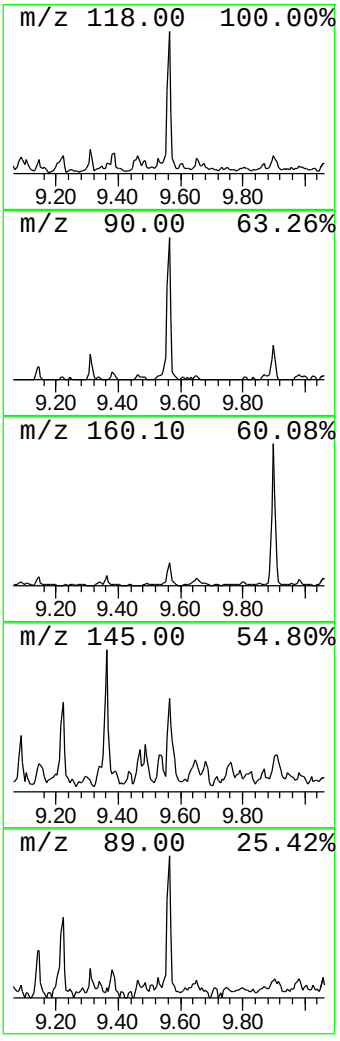
Peak Number 7 1(2H)-Naphthalenone, 3,4-di... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.56	2.74 ng	165515	Acenaphthene-d10	9.90

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1(2H)-Naphthalenone, 3,4-dihydro...	160	C11H12O	014944-23-1	94
2	1(2H)-Naphthalenone, 3,4-dihydro...	160	C11H12O	001590-08-5	80
3	1H-Inden-1-one, 2,3-dihydro-3,3-...	160	C11H12O	026465-81-6	38
4	1,4-Epoxynaphthalene, 1,2,3,4-te...	146	C10H10O	035185-96-7	30
5	Benzene, 4-chloro-2-fluoro-1-met...	160	C7H6ClFO	000452-09-5	27





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Misc :
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Instrument : BNA_F
ClientSampleId : OWL-C6-GW

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

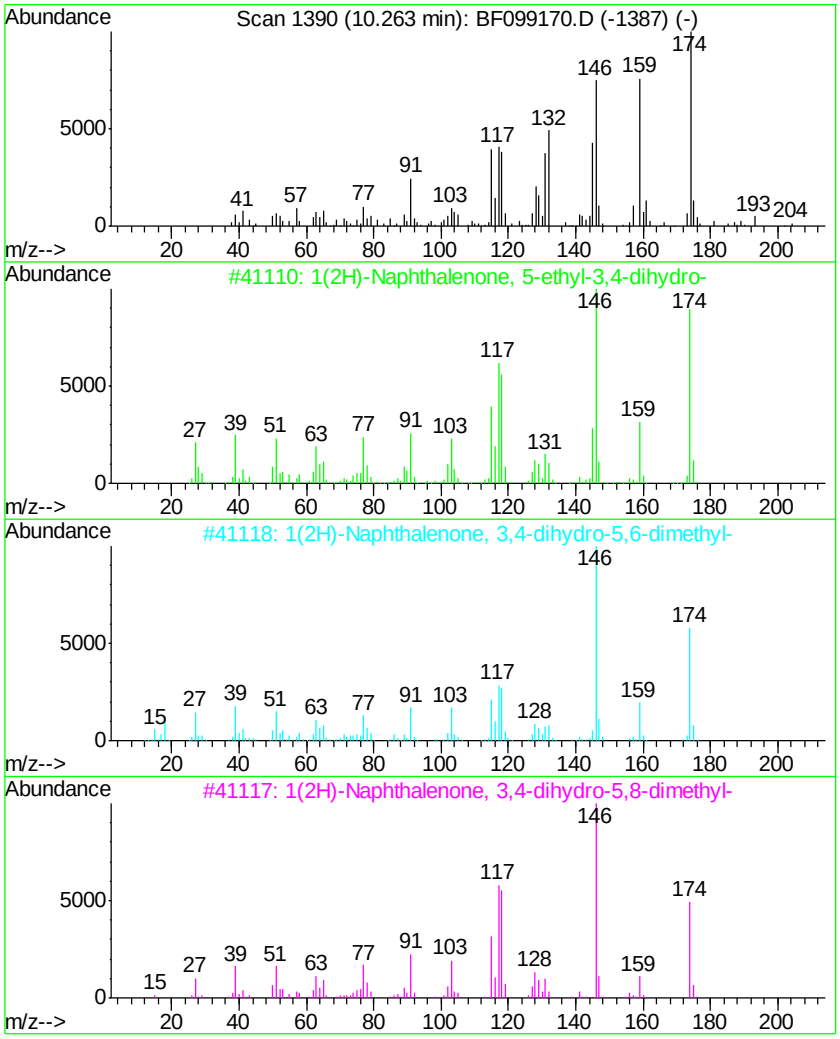
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

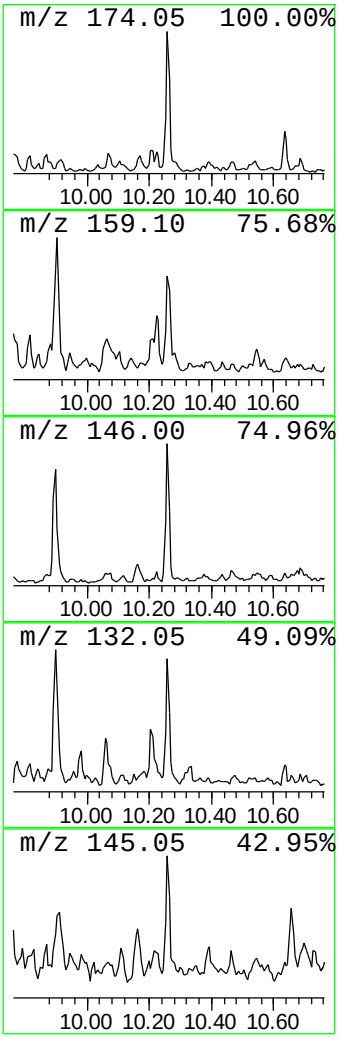
****

Peak Number 8 1(2H)-Naphthalenone, 5-ethy... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.26	3.09 ng	186739	Acenaphthene-d10	9.90

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1(2H)-Naphthalenone, 5-ethyl-3,4...	174	C12H14O	051015-31-7	81
2	1(2H)-Naphthalenone, 3,4-dihydro...	174	C12H14O	032281-65-5	76
3	1(2H)-Naphthalenone, 3,4-dihydro...	174	C12H14O	005037-63-8	64
4	Quinazoline, 4-ethyl-, 3-oxide	174	C10H10N2O	037920-74-4	53
5	1(2H)-Naphthalenone, 3,4-dihydro...	174	C12H14O	030316-30-4	53





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Instrument : BNA_F
ClientSampleId : OWL-C6-GW

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

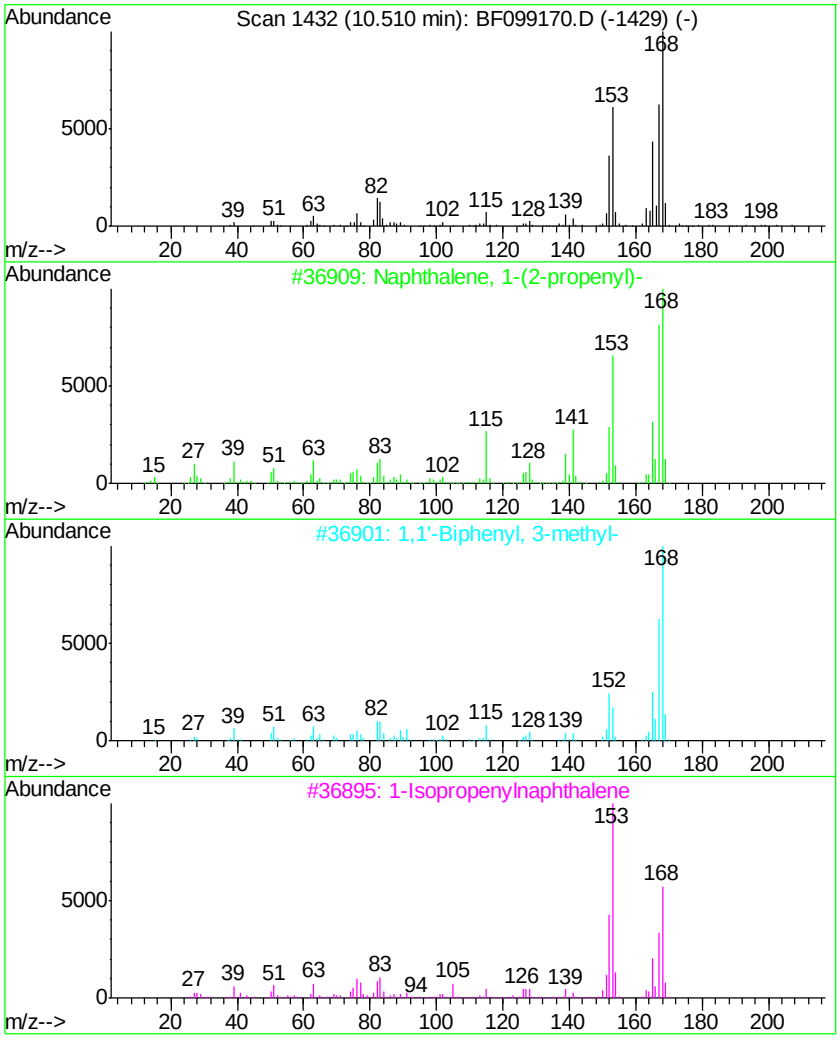
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

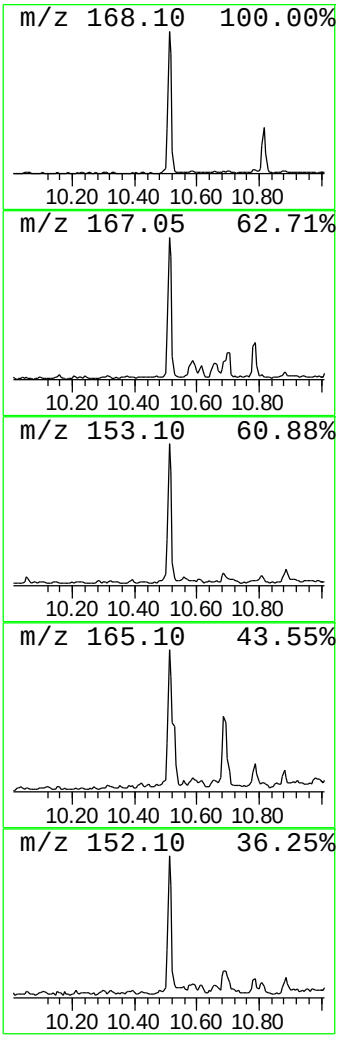
****

Peak Number 9 Naphthalene, 1-(2-propenyl)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.51	4.14 ng	250668	Acenaphthene-d10	9.90

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1-(2-propenyl)-	168	C13H12	002489-86-3	74
2	1,1'-Biphenyl, 3-methyl-	168	C13H12	000643-93-6	64
3	1-Isopropenylnaphthalene	168	C13H12	001855-47-6	64
4	1,1'-Biphenyl, 4-methyl-	168	C13H12	000644-08-6	62
5	1,1'-Biphenyl, 2-methyl-	168	C13H12	000643-58-3	53





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Acq On : 7 Oct 2017 1:51
Operator : SJ/JU
Sample : I5542-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : OWL-C6-GW

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.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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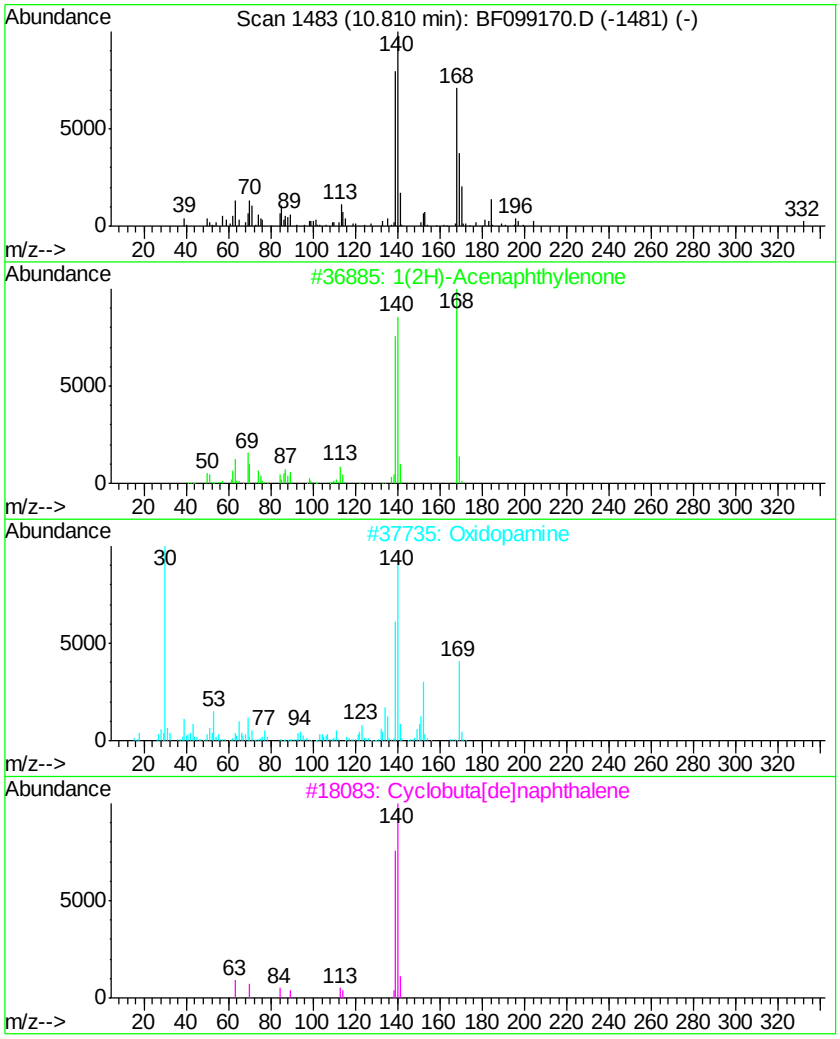
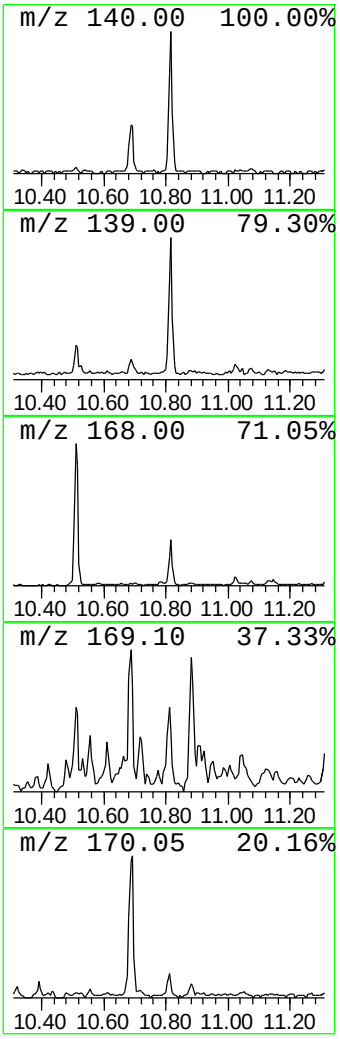
Peak Number 10 1(2H)-Acenaphthylenone Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.81	3.97 ng	176593	Phenanthrene-d10	11.39

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1(2H)-Acenaphthylenone	168	C12H8O	002235-15-6	74
2	Oxidopamine	169	C8H11NO3	001199-18-4	53
3	Cyclobuta[de]naphthalene	140	C11H8	024973-91-9	52
4	4H-Pyran-4-one, 2-ethyl-3-hydroxy-	140	C7H8O3	004940-11-8	47
5	Dibenzofuran	168	C12H8O	000132-64-9	38





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Acq On : 7 Oct 2017 1:51
Operator : SJ/JU
Sample : I5542-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : OWL-C6-GW

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.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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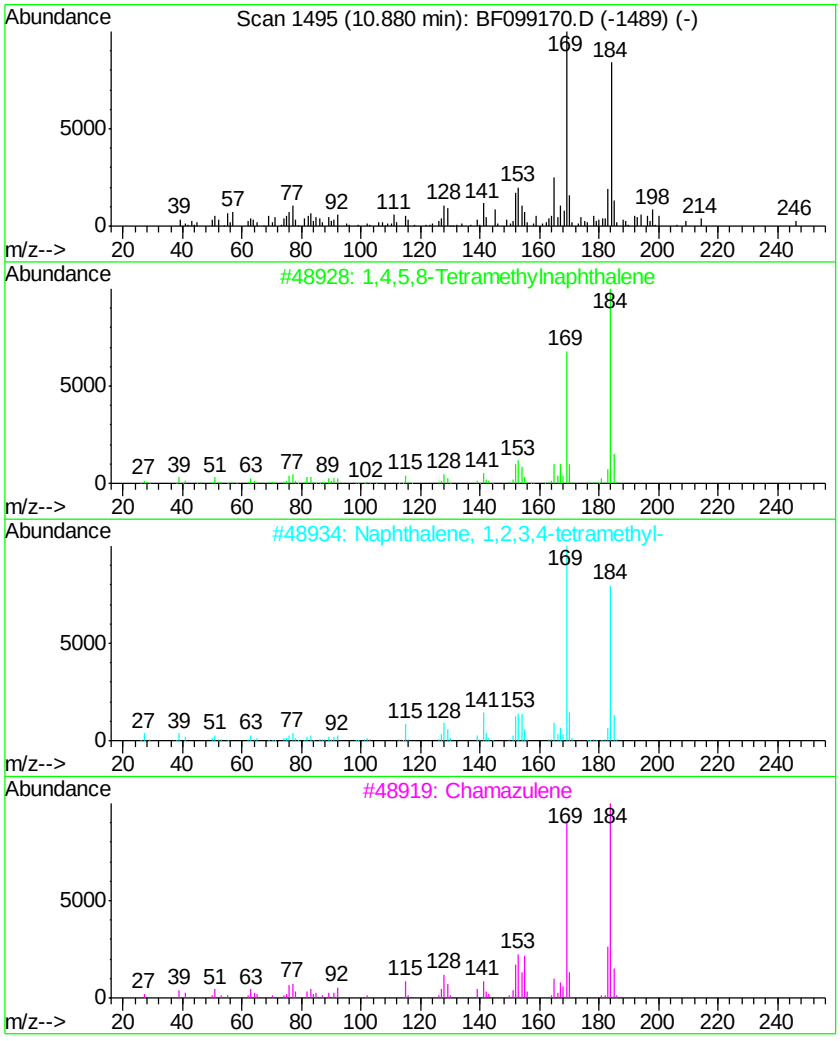
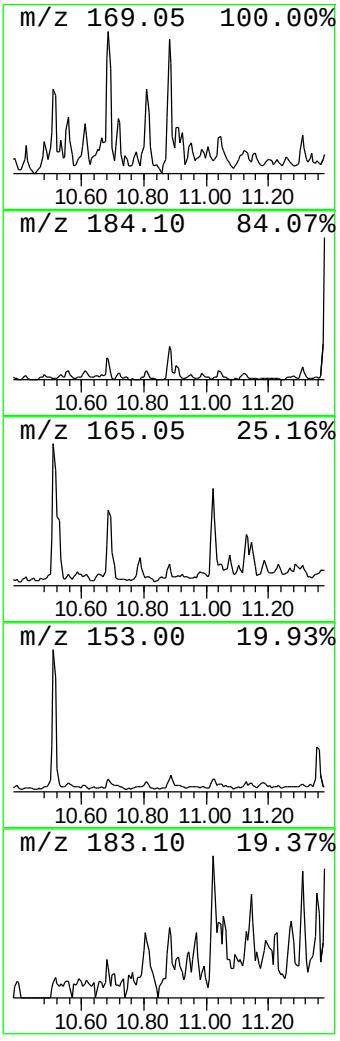
Peak Number 11 1,4,5,8-Tetramethylnaphthalene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.88	3.25 ng	144525	Phenanthrene-d10	11.39

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,4,5,8-Tetramethylnaphthalene	184	C14H16	002717-39-7	94
2	Naphthalene, 1,2,3,4-tetramethyl-	184	C14H16	003031-15-0	94
3	Chamazulene	184	C14H16	000529-05-5	86
4	Cyclopentene, 1,2-dimethyl-4-met...	184	C14H16	1000152-22-4	72
5	Naphthalene, 1-methyl-7-(1-methy...	184	C14H16	000490-65-3	70





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Acq On : 7 Oct 2017 1:51
Operator : SJ/JU
Sample : I5542-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : OWL-C6-GW

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.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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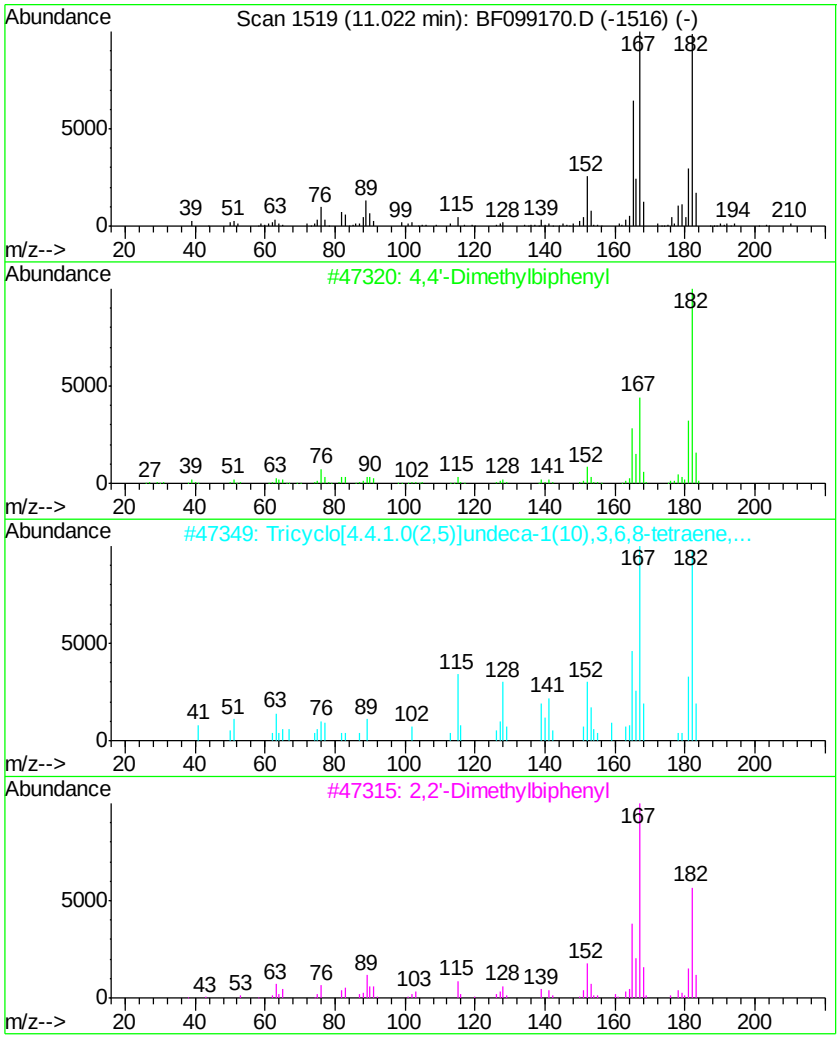
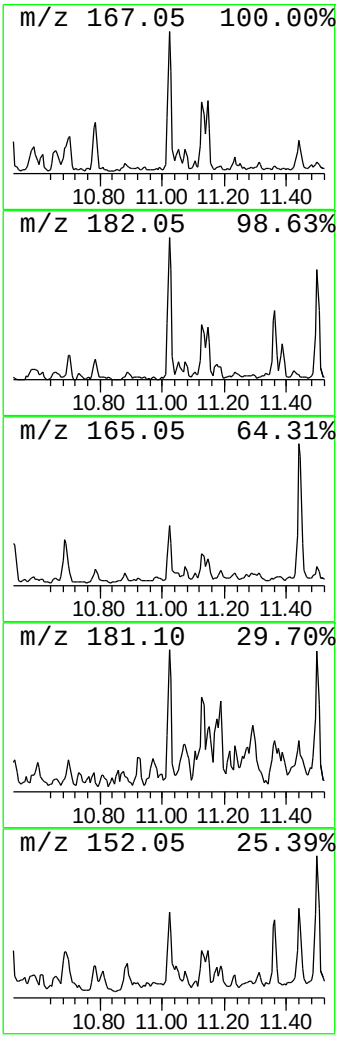
Peak Number 12 4,4'-Dimethylbiphenyl Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.02	2.42 ng	107532	Phenanthrene-d10	11.39

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4,4'-Dimethylbiphenyl	182	C14H14	000613-33-2	95
2	Tricyclo[4.4.1.0(2,5)]undeca-1(1...	182	C14H14	055836-29-8	90
3	2,2'-Dimethylbiphenyl	182	C14H14	000605-39-0	90
4	1,1'-Biphenyl, 2,4'-dimethyl-	182	C14H14	000611-61-0	87
5	1,1'-Biphenyl, 2,3'-dimethyl-	182	C14H14	000611-43-8	87





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Data File : BF099170.D
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Operator : SJ/JU
Sample : I5542-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : OWL-C6-GW

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.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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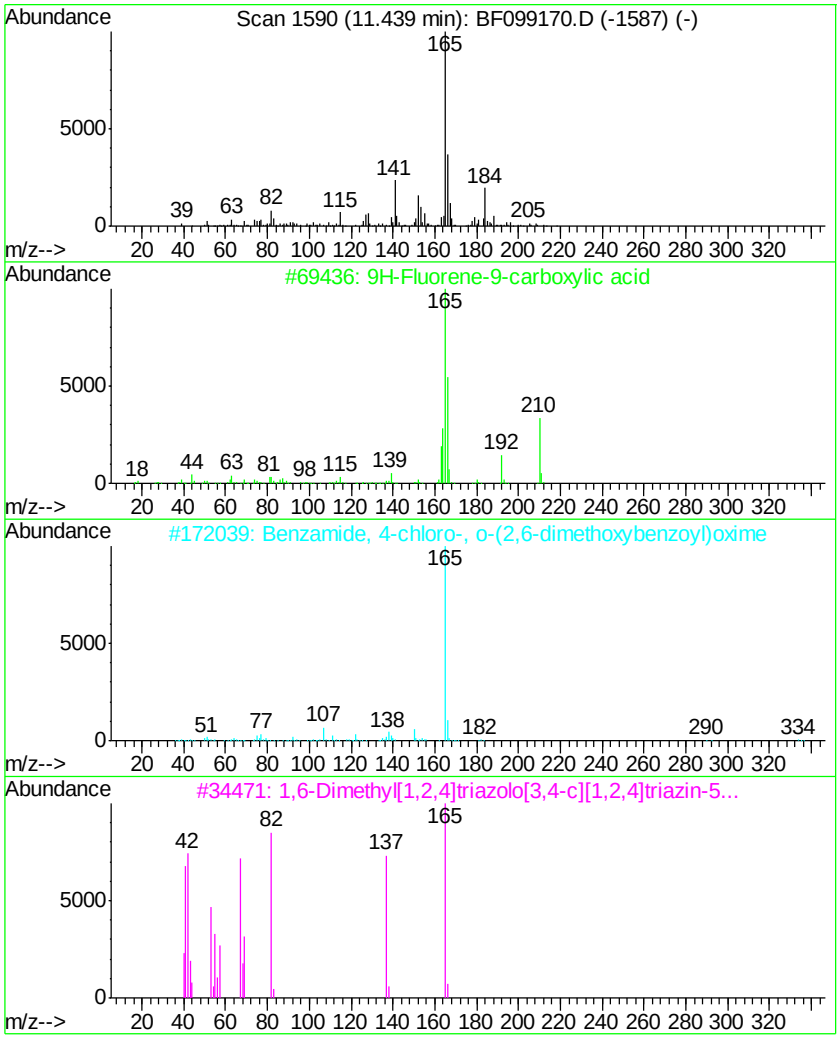
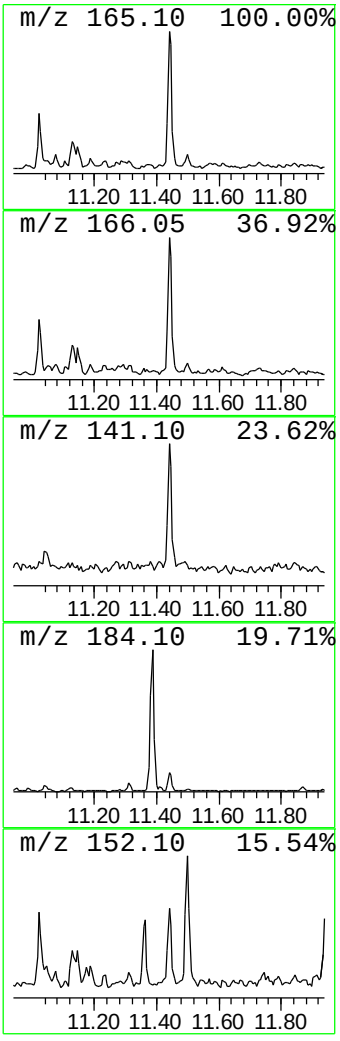
Peak Number 13 unknown11.44 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.44	3.49 ng	154857	Phenanthrene-d10	11.39

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9H-Fluorene-9-carboxylic acid	210	C14H10O2	001989-33-9	38
2	Benzamide, 4-chloro-, o-(2,6-dim...	334	C16H15ClN2O4	1000263-23-0	25
3	1,6-Dimethyl[1,2,4]triazolo[3,4-...	165	C6H7N5O	1000146-89-4	25
4	1H-Phenalene	166	C13H10	000203-80-5	25
5	1-(3,4-methylenedioxyphenyl)-1-m...	208	C11H12O4	1000378-91-3	23





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Data File : BF099170.D
Acq On : 7 Oct 2017 1:51
Operator : SJ/JU
Sample : I5542-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : OWL-C6-GW

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.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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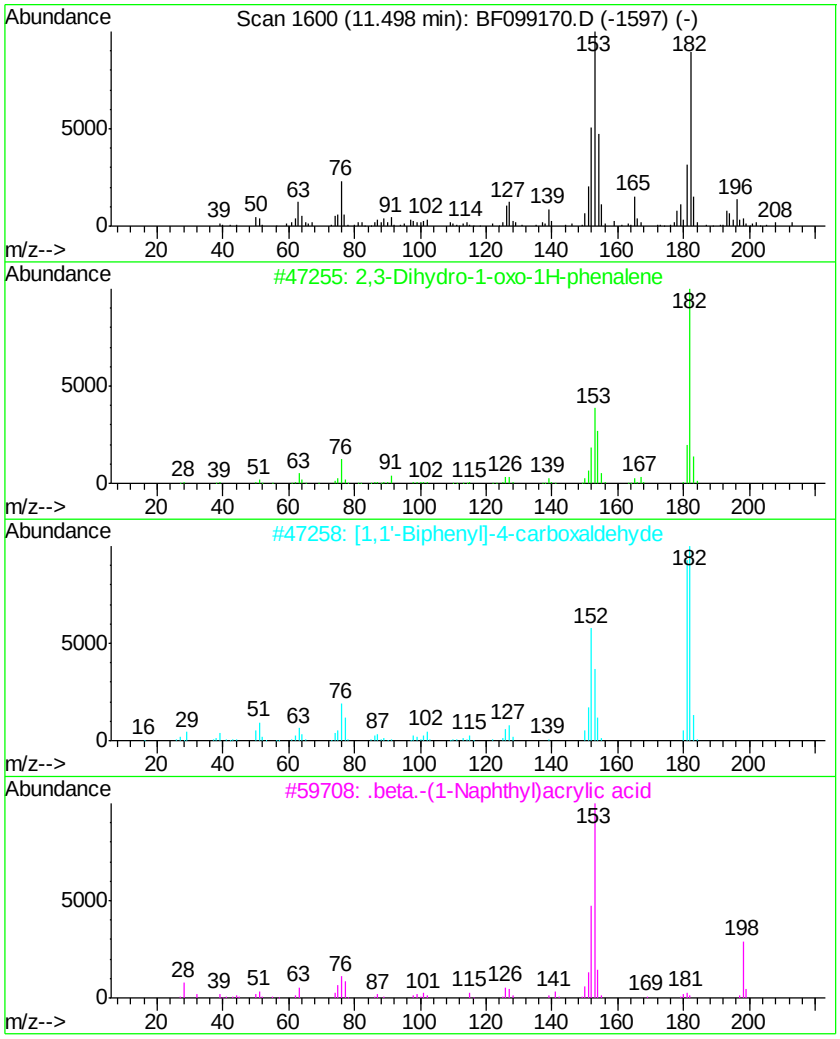
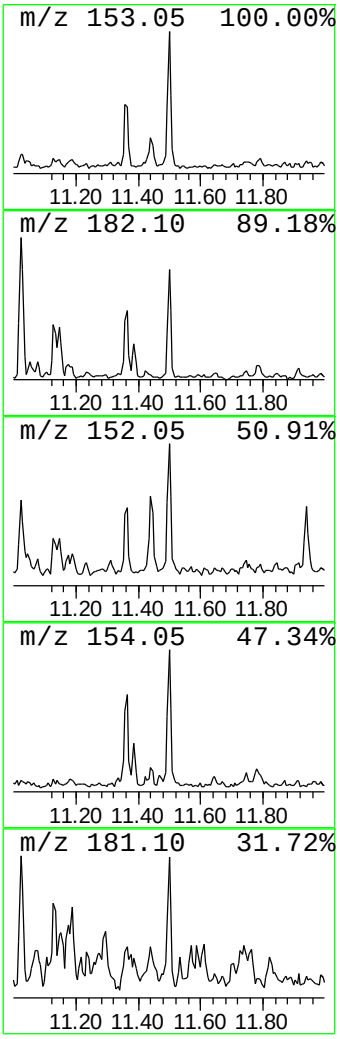
Peak Number 14 2,3-Dihydro-1-oxo-1H-phenalene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.50	3.30 ng	146823	Phenanthrene-d10	11.39

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,3-Dihydro-1-oxo-1H-phenalene	182	C13H10O	000518-85-4	87
2	[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	53
3	.beta.-(1-Naphthyl)acrylic acid	198	C13H10O2	013026-12-5	49
4	Phenanthrene, 1,2,3,4-tetrahydro-	182	C14H14	001013-08-7	43
5	Bicyclo[3.2.1]octane-6-carboxyli...	182	C10H14O3	1000362-16-3	42





Tentatively Identified Compound (LSC) summary

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF100617\
Data File : BF099170.D
Acq On : 7 Oct 2017 1:51
Operator : SJ/JU
Sample : I5542-05
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Instrument : BNA_F
ClientSampleId : OWL-C6-GW

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Butane, 2-methoxy...	2.15	12.4	ng	479470	1	6.86	772657	20.0
2-Pentanone, 4-hy...	5.09	9.7	ng	373131	1	6.86	772657	20.0
unknown6.63	6.63	73.9	ng	2854790	1	6.86	772657	20.0
1-Methylindan-2-one	8.95	3.5	ng	182796	2	8.15	1032240	20.0
Naphthalene, 1,2,...	9.15	4.1	ng	247855	3	9.90	1210040	20.0
1(2H)-Naphthaleno...	9.56	2.7	ng	165515	3	9.90	1210040	20.0
1(2H)-Naphthaleno...	10.26	3.1	ng	186739	3	9.90	1210040	20.0
Naphthalene, 1-(2...	10.51	4.1	ng	250668	3	9.90	1210040	20.0
1(2H)-Acenaphthyl...	10.81	4.0	ng	176593	4	11.39	888528	20.0
1,4,5,8-Tetrameth...	10.88	3.3	ng	144525	4	11.39	888528	20.0
4,4'-Dimethylbiph...	11.02	2.4	ng	107532	4	11.39	888528	20.0
unknown11.44	11.44	3.5	ng	154857	4	11.39	888528	20.0
2,3-Dihydro-1-oxo...	11.50	3.3	ng	146823	4	11.39	888528	20.0