LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092618\
Data File : BF109355.D
Acq On : 26 Sep 2018 23:26
Operator : JU/SJ
Sample : J5055-20
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : OILY-WATER

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-BF092218.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	4.887	473	476	484	rBV	55914	62485	2.61%	0.327%
2	5.316	545	549	561	rBV	1190488	1079081	45.01%	5.653%
3	6.340	720	723	733	rVB	845713	709639	29.60%	3.717%
4	6.475	742	746	749	rBV	2063186	1831807	76.42%	9.596%
5	6.640	770	774	782	rBV	75489	70242	2.93%	0.368%
6	6.704	782	785	789	rBV	645216	589163	24.58%	3.086%
7	6.863	808	812	816	rVB	2095831	1820646	75.95%	9.537%
8	7.269	876	881	884	rBV	1550643	1391905	58.06%	7.291%
9	7.569	928	932	939	rBV3	22563	46350	1.93%	0.243%
10	7.816	967	974	977	rBV	33783	46179	1.93%	0.242%
11	7.987	997	1003	1007	rBV	840852	708421	29.55%	3.711%
12	8.087	1015	1020	1023	rBV	73597	65966	2.75%	0.346%
13	8.187	1035	1037	1046	rVB6	23469	39486	1.65%	0.207%
14	8.351	1059	1065	1066	rBV3	26972	34747	1.45%	0.182%
15	8.369	1066	1068	1073	rVB3	33486	36938	1.54%	0.193%
16	8.598	1105	1107	1109	rBV2	37707	30957	1.29%	0.162%
17	8.704	1122	1125	1131	rBV	168724	219215	9.14%	1.148%
18	8.798	1131	1141	1144	rVB2	94454	220048	9.18%	1.153%
19	9.063	1182	1186	1190	rBV	2695122	2397175	100.00%	12.557%
20	9.228	1211	1214	1218	rBV5	32941	41200	1.72%	0.216%
21	9.304	1224	1227	1230	rBV4	41087	45126	1.88%	0.236%
22	9.404	1241	1244	1248	rBV2	62840	75019	3.13%	0.393%
23	9.457	1250	1253	1257	rVB	134718	113186	4.72%	0.593%
24	9.504	1258	1261	1264	rVB4	25873	30462	1.27%	0.160%
25	9.545	1264	1268	1269	rBV3	24516	32880	1.37%	0.172%
26	9.651	1284	1286	1289	rVB2	46949	43740	1.82%	0.229%
27	9.739	1297	1301	1305	rVB	789097	694397	28.97%	3.638%
28	9.975	1338	1341	1343	rBV2	88989	83307	3.48%	0.436%
29	10.151	1368	1371	1374	rBV	69479	63074	2.63%	0.330%
30	10.522	1430	1434	1438	rBV	1360997	1275063	53.19%	6.679%
31	11.216	1549	1552	1555	rBV	734189	615289	25.67%	3.223%
32	11.775	1644	1647	1652	rVB	243373	246506	10.28%	1.291%
33	12.498	1768	1770	1774	rVB	164535	135089	5.64%	0.708%
34	12.804	1818	1822	1825	rVB	2214309	2049978	85.52%	10.739%

LSC Area Percent Report

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Acq On : 26 Sep 2018 23:26
Operator : JU/SJ
Sample : J5055-20
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : OILY-WATER

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-BF092218.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	13.710	1974	1976	1985	rVB	107449	132501	5.53%	0.694%
36	13.845	1995	1999	2004	rVB	875498	779458	32.52%	4.083%
37	14.692	2140	2143	2149	rVB	196440	189652	7.91%	0.993%
38	14.933	2182	2184	2188	rVB	54961	55524	2.32%	0.291%
39	15.245	2233	2237	2241	rVB	569743	639321	26.67%	3.349%
40	15.686	2310	2312	2317	rVB2	45484	52919	2.21%	0.277%
41	16.874	2510	2514	2521	rVB	56015	101990	4.25%	0.534%
42	17.509	2616	2622	2633	rVB	78377	193543	8.07%	1.014%

Sum of corrected areas: 19089674

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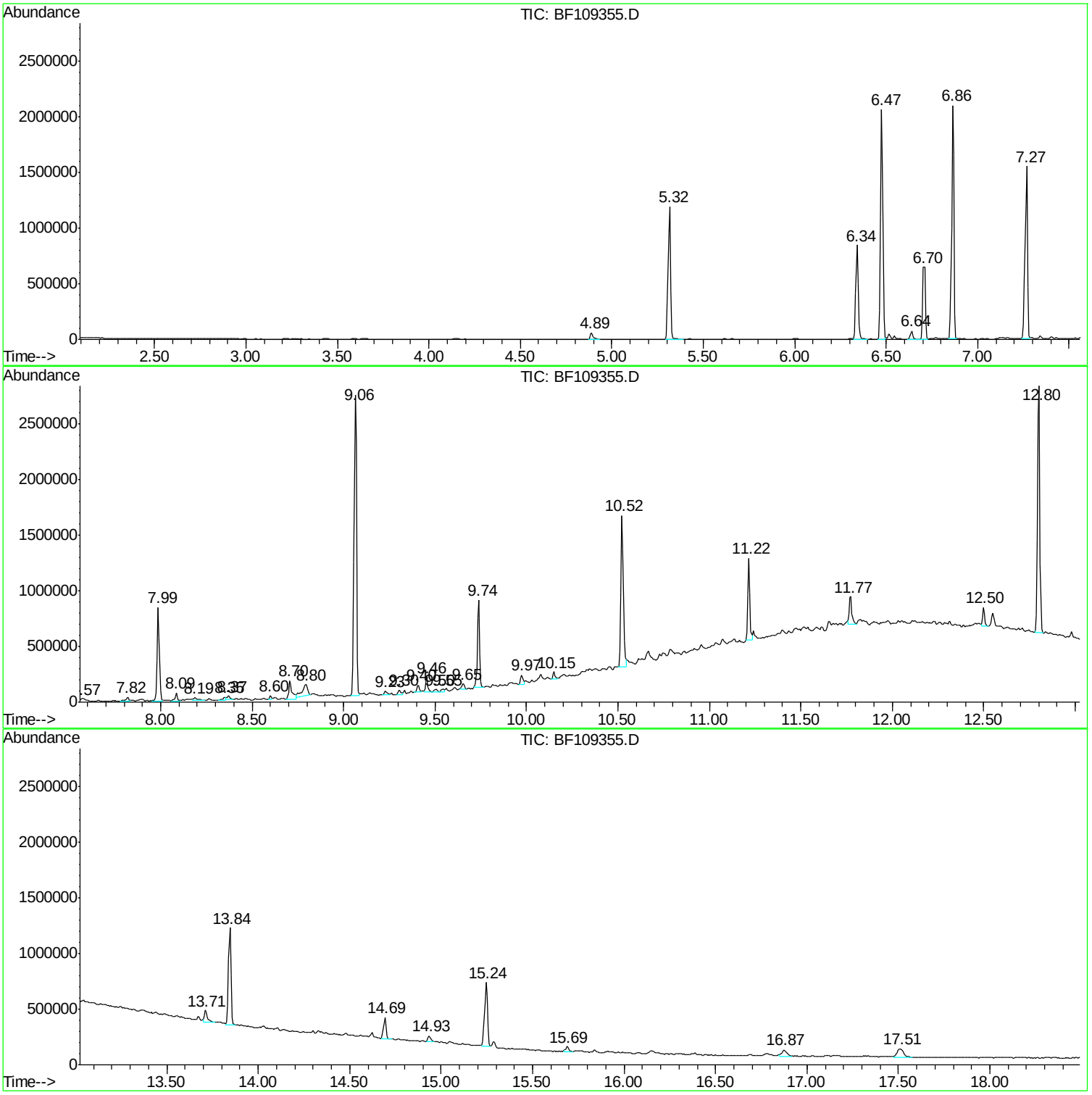
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
OILY-WATER

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF092218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Library Search Compound Report

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Sample : J5055-20
Misc :
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Instrument : BNA_F
ClientSampleId : OILY-WATER

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF092218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

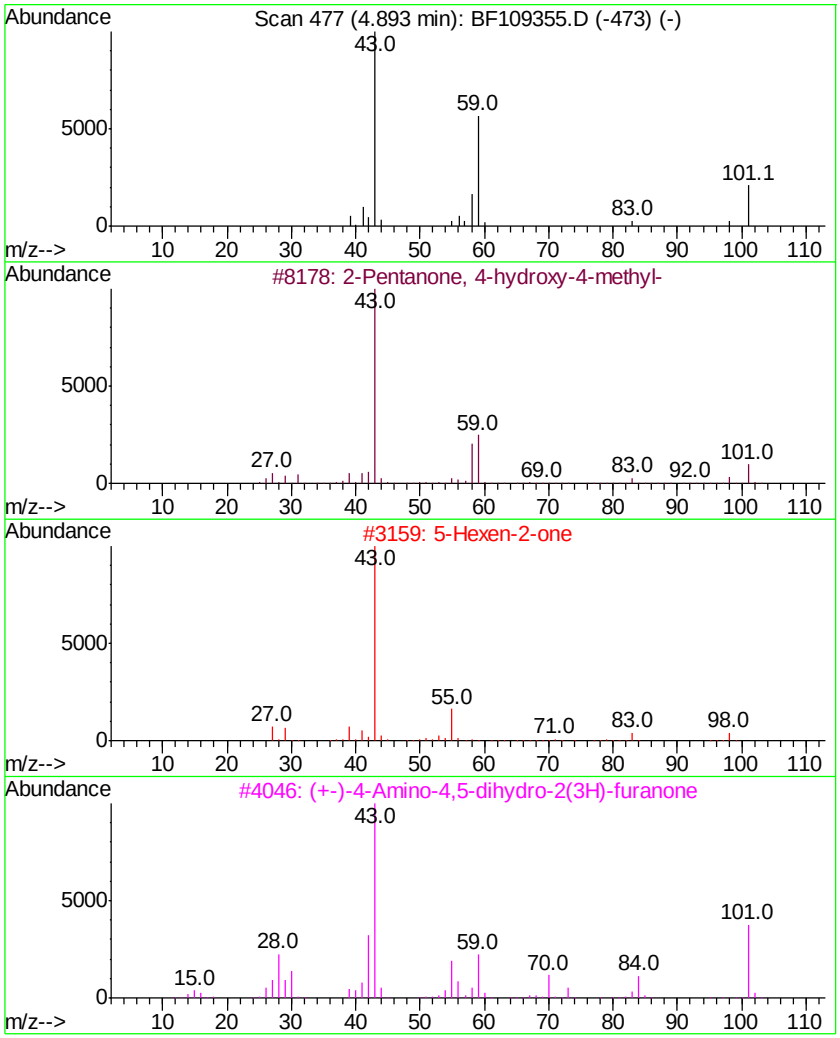
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

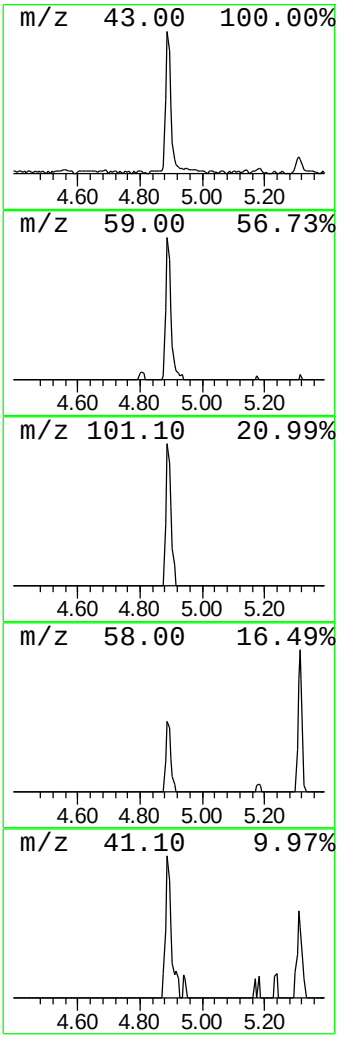
****

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.89	2.12 ng	62485	1,4-Dichlorobenzene-d4	6.71

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2	5-Hexen-2-one	98	C6H10O	000109-49-9	9
3	(+-)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9
4	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5	Acetone	58	C3H6O	000067-64-1	9





Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092618\
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Sample : J5055-20
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : OILY-WATER

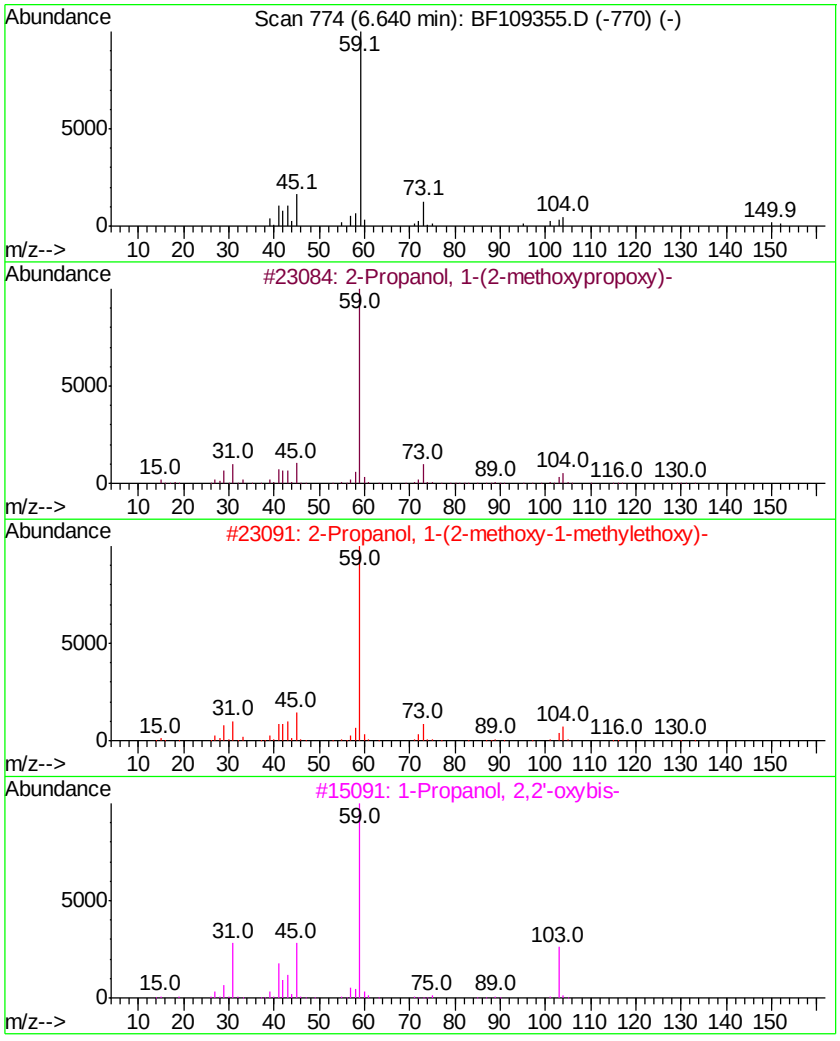
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF092218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

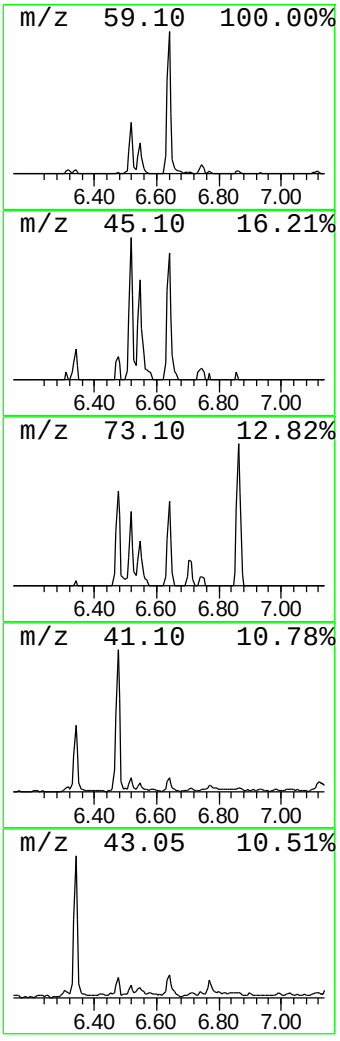
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 2-Propanol, 1-(2-methoxypro... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.64	2.38 ng	70242	1,4-Dichlorobenzene-d4	6.71

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-(2-methoxypropoxy)-	148	C7H16O3	013429-07-7	78
2	2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	78
3	1-Propanol, 2,2'-oxybis-	134	C6H14O3	000108-61-2	59
4	1-Propanol, 2-(2-methoxypropoxy)-	148	C7H16O3	013588-28-8	59
5	1-(1-Methoxypropan-2-yloxy)propa...	232	C12H24O4	1000367-13-2	50





Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092618\
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Operator : JU/SJ
Sample : J5055-20
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : OILY-WATER

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF092218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

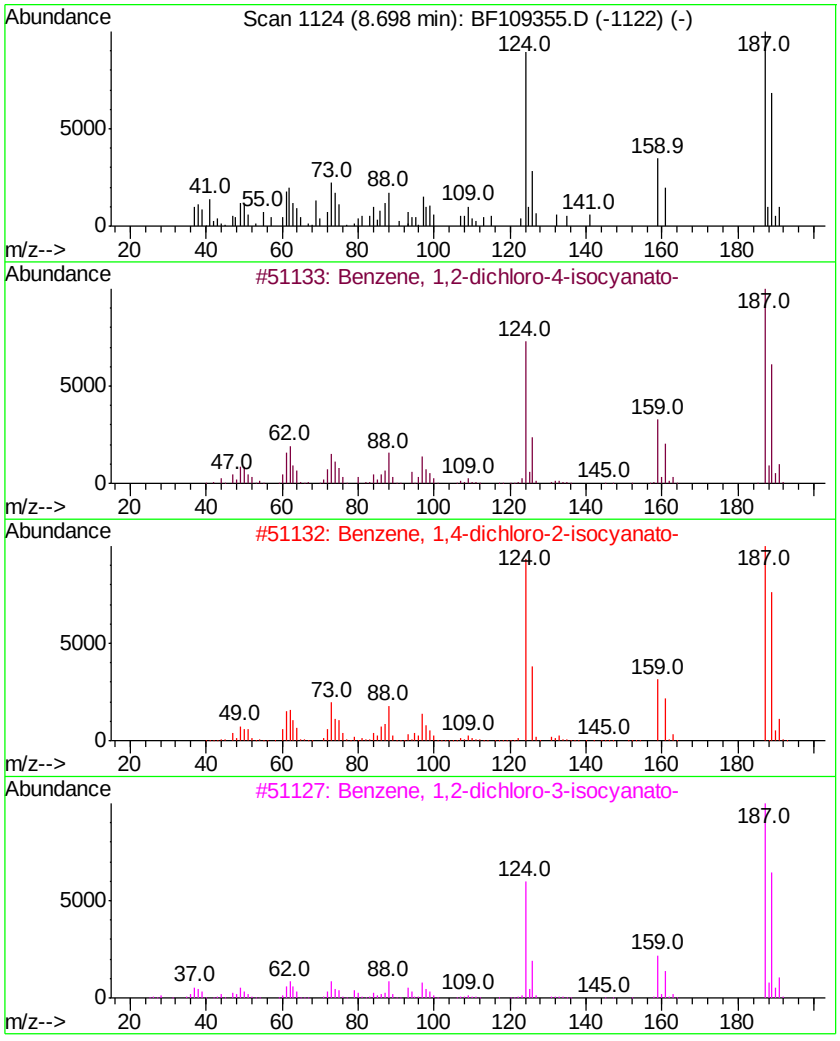
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

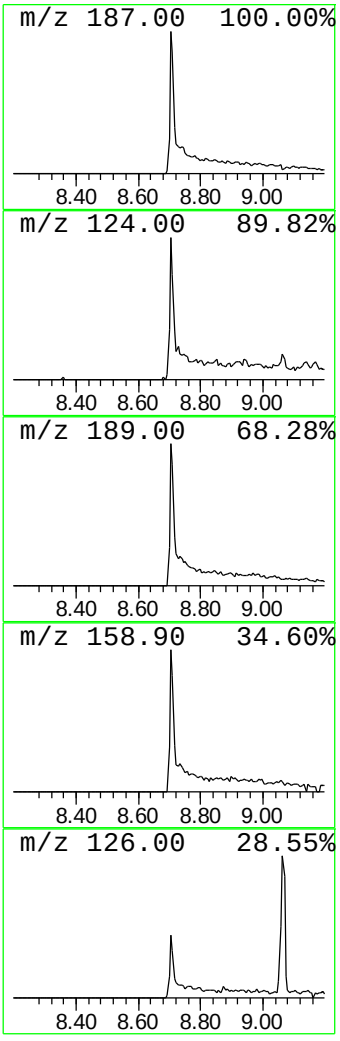
****

Peak Number 5 Benzene, 1,2-dichloro-4-iso... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.70	6.19 ng	219215	Naphthalene-d8	7.99

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2-dichloro-4-isocyanato-	187	C7H3Cl2NO	000102-36-3	99
2	Benzene, 1,4-dichloro-2-isocyanato-	187	C7H3Cl2NO	005392-82-5	98
3	Benzene, 1,2-dichloro-3-isocyanato-	187	C7H3Cl2NO	041195-90-8	98
4	3,5-Dichlorophenyl isocyanate	187	C7H3Cl2NO	034893-92-0	97
5	Benzene, 1,3-dichloro-2-isocyanato-	187	C7H3Cl2NO	039920-37-1	96





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Sample : J5055-20
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : OILY-WATER

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF092218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

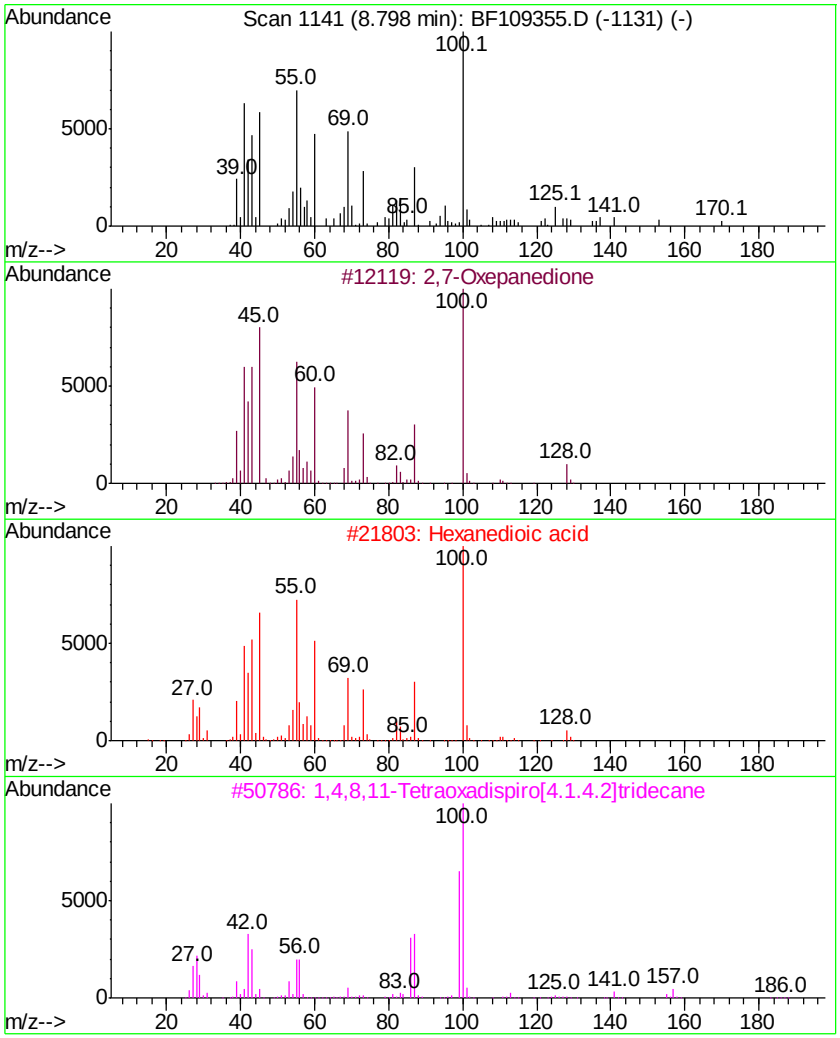
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

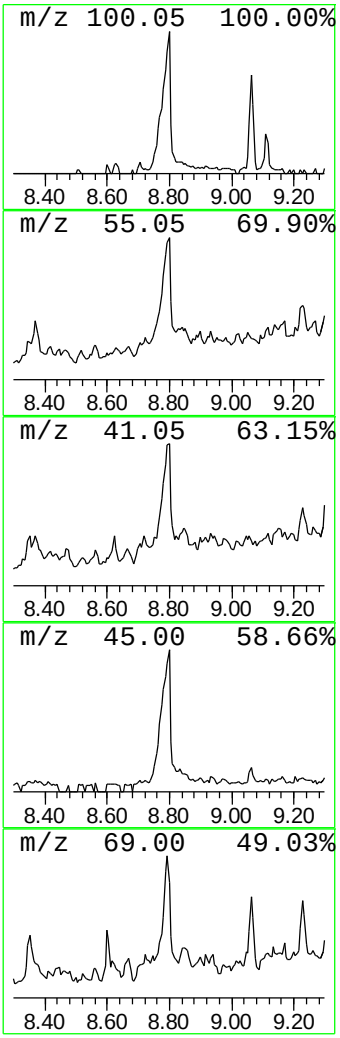
****

Peak Number 6 2,7-Oxepanedione Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.80	6.21 ng	220048	Naphthalene-d8	7.99

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,7-Oxepanedione	128	C6H8O3	002035-75-8	93
2	Hexanedioic acid	146	C6H10O4	000124-04-9	86
3	1,4,8,11-Tetraoxadispiro[4.1.4.2...	186	C9H14O4	004086-01-5	43
4	Ethane-1,2-diamine, N,N'-dinitro...	348	C12H24N6O6	1000271-06-8	35
5	2-Pentenoic acid	100	C5H8O2	000626-98-2	27





Library Search Compound Report

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Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : OILY-WATER

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF092218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
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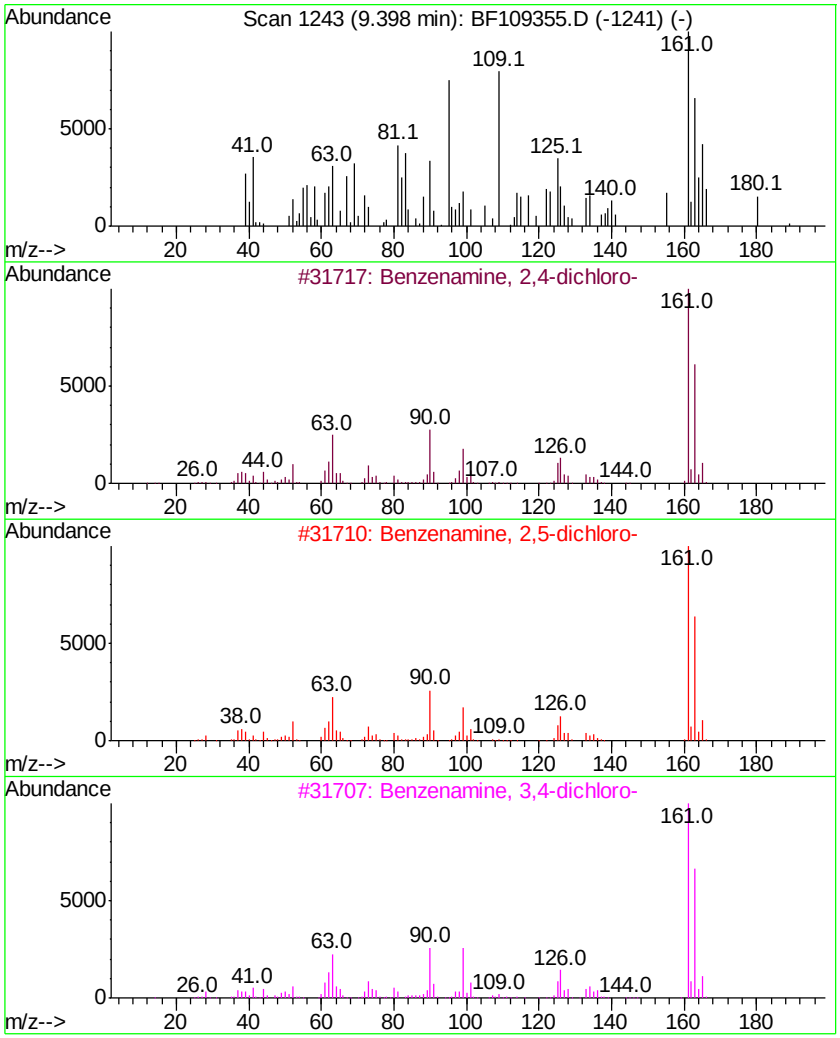
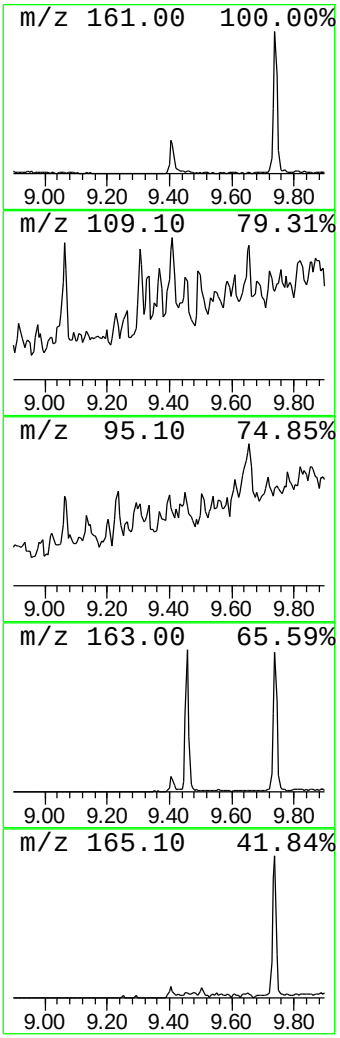
****

Peak Number 7 Benzenamine, 2,4-dichloro- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.40	2.16 ng	75019	Acenaphthene-d10	9.74

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenamine, 2,4-dichloro-	161	C6H5Cl2N	000554-00-7	96
2	Benzenamine, 2,5-dichloro-	161	C6H5Cl2N	000095-82-9	96
3	Benzenamine, 3,4-dichloro-	161	C6H5Cl2N	000095-76-1	95
4	Benzenamine, 3,5-dichloro-	161	C6H5Cl2N	000626-43-7	91
5	Benzenamine, 2,3-dichloro-	161	C6H5Cl2N	000608-27-5	89





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Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : OILY-WATER

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF092218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
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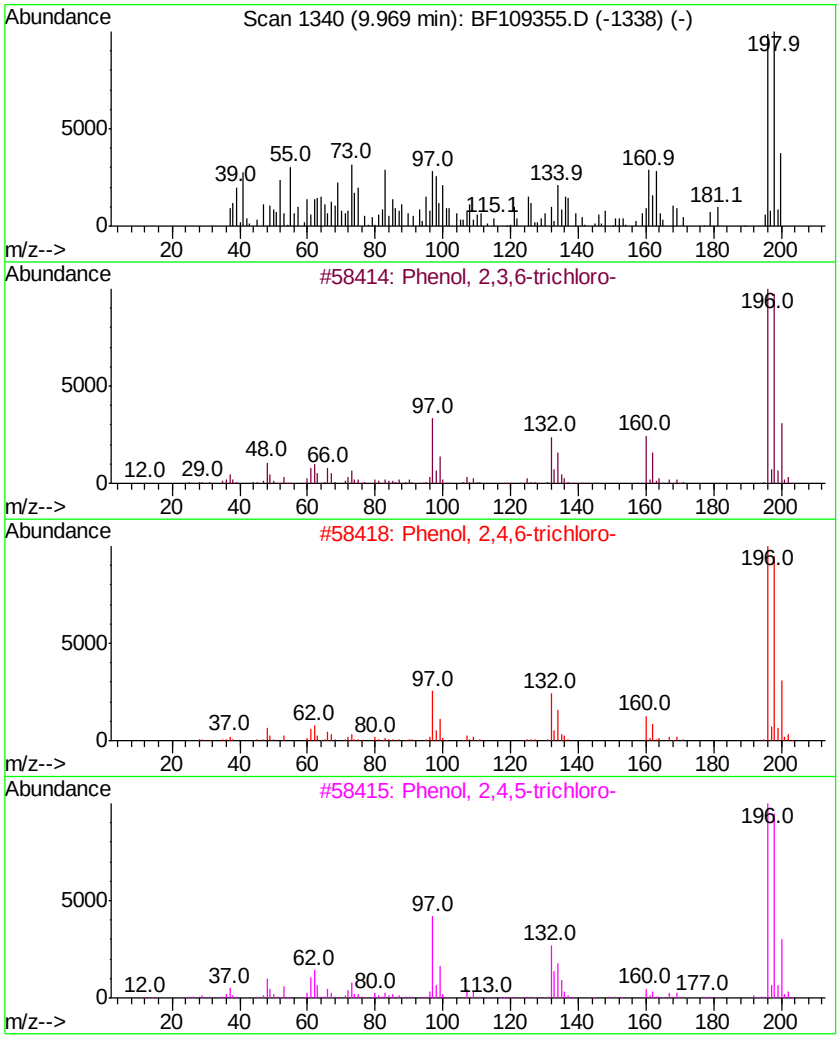
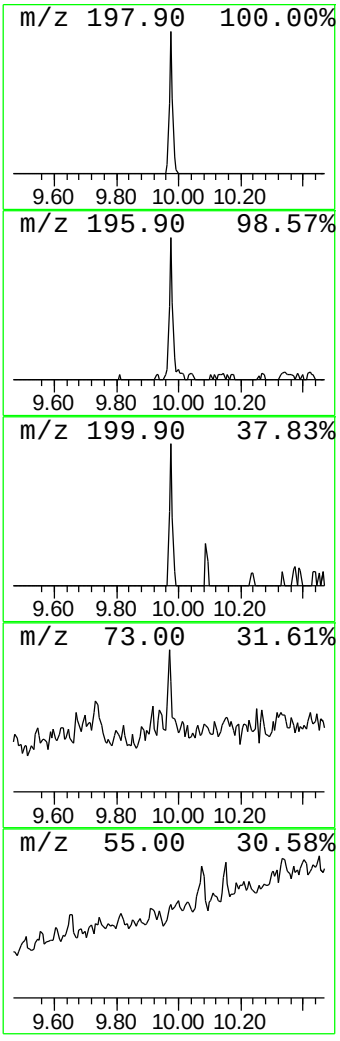
****

Peak Number 8 Phenol, 2,3,6-trichloro- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.97	2.40 ng	83307	Acenaphthene-d10	9.74

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,3,6-trichloro-	196	C6H3Cl3O	000933-75-5	86
2	Phenol, 2,4,6-trichloro-	196	C6H3Cl3O	000088-06-2	70
3	Phenol, 2,4,5-trichloro-	196	C6H3Cl3O	000095-95-4	64
4	Phenol, 2,3,5-trichloro-	196	C6H3Cl3O	000933-78-8	64
5	Phenol, 2,3,4-trichloro-	196	C6H3Cl3O	015950-66-0	62





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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
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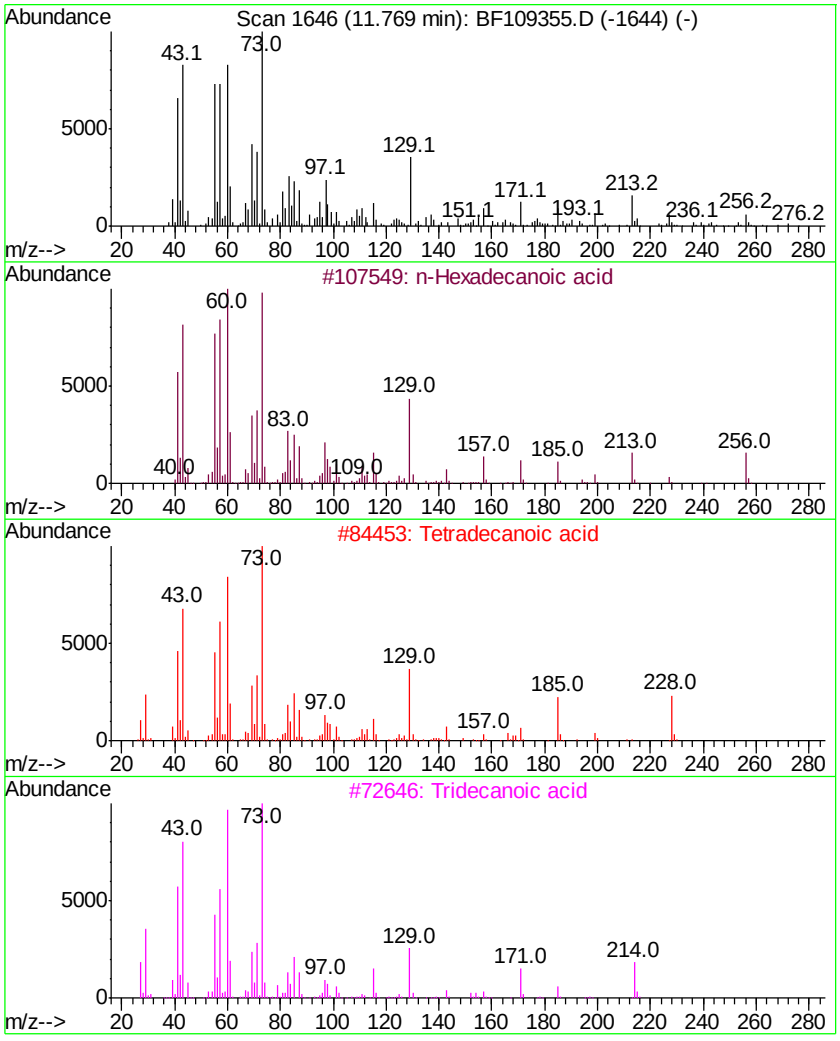
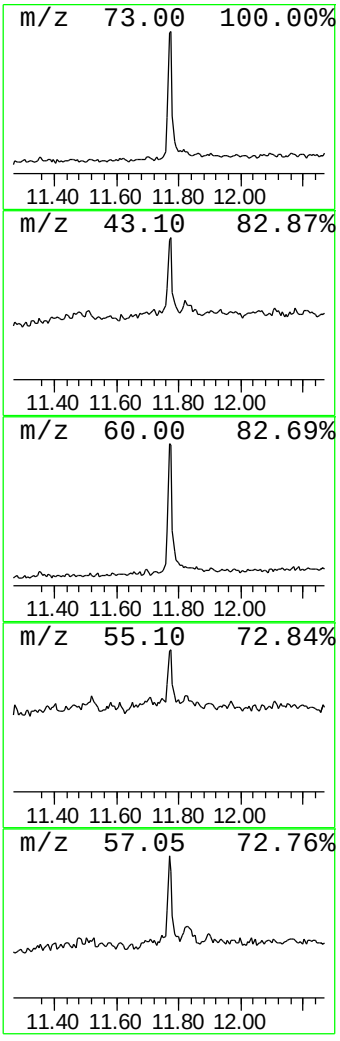
****

Peak Number 9 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.77	8.01 ng	246506	Phenanthrene-d10	11.22

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	91
3	Tridecanoic acid	214	C13H26O2	000638-53-9	89
4	Pentadecanoic acid	242	C15H30O2	001002-84-2	83
5	Dodecanoic acid	200	C12H24O2	000143-07-7	58





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Acq On : 26 Sep 2018 23:26
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Sample : J5055-20
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : OILY-WATER

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF092218.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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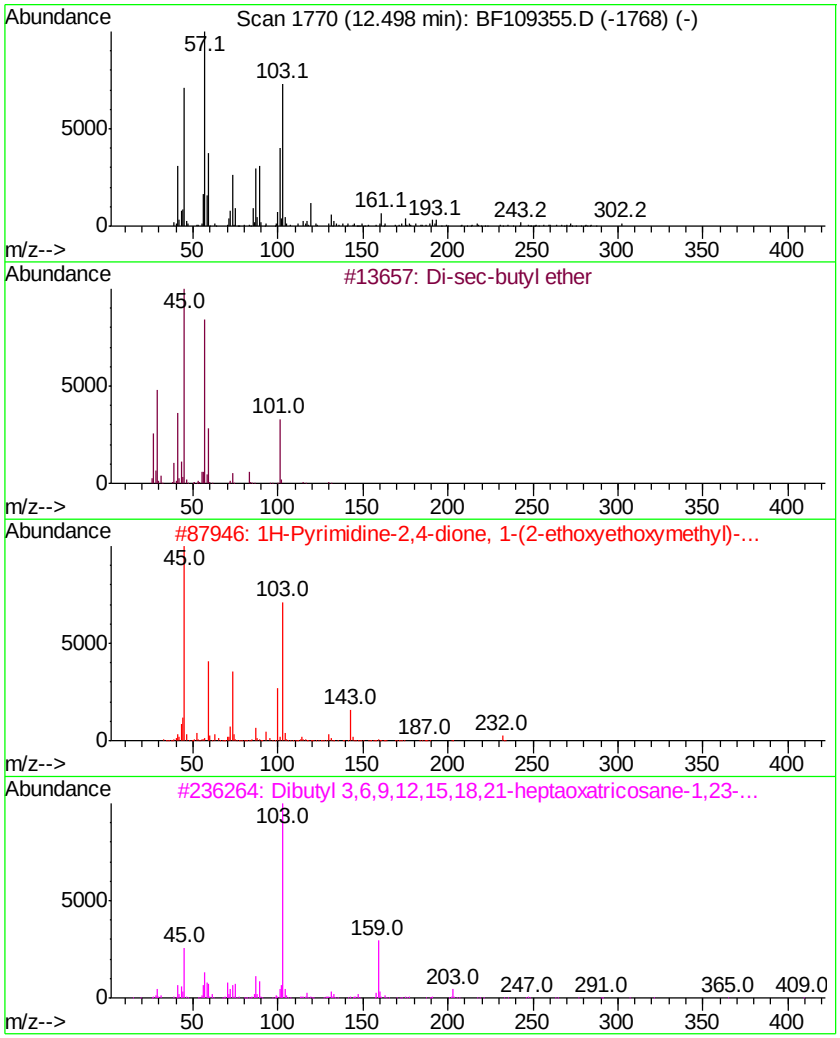
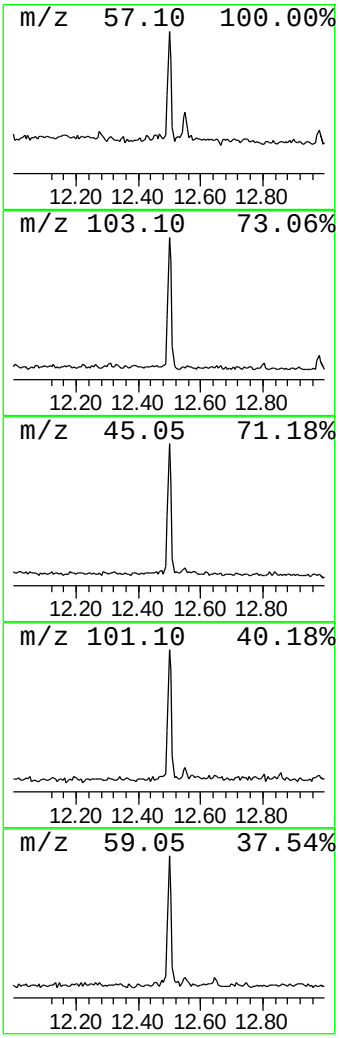
Peak Number 10 unknown12.50 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.50	4.39 ng	135089	Phenanthrene-d10	11.22

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Di-sec-butyl ether	130	C8H18O	006863-58-7	38
2	1H-Pyrimidine-2,4-dione, 1-(2-et...	232	C9H13FN2O4	1000310-13-7	38
3	Dibutyl 3,6,9,12,15,18,21-heptao...	510	C24H46O11	1000366-79-9	35
4	2-[2-[2-[2-[2-(2-Methoxyethox...	340	C15H32O8	1000352-04-1	35
5	1,4,7,10,13,16-Hexaoxacyclooctad...	264	C12H24O6	017455-13-9	27





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092618\
Data File : BF109355.D
Acq On : 26 Sep 2018 23:26
Operator : JU/SJ
Sample : J5055-20
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : OILY-WATER

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF092218.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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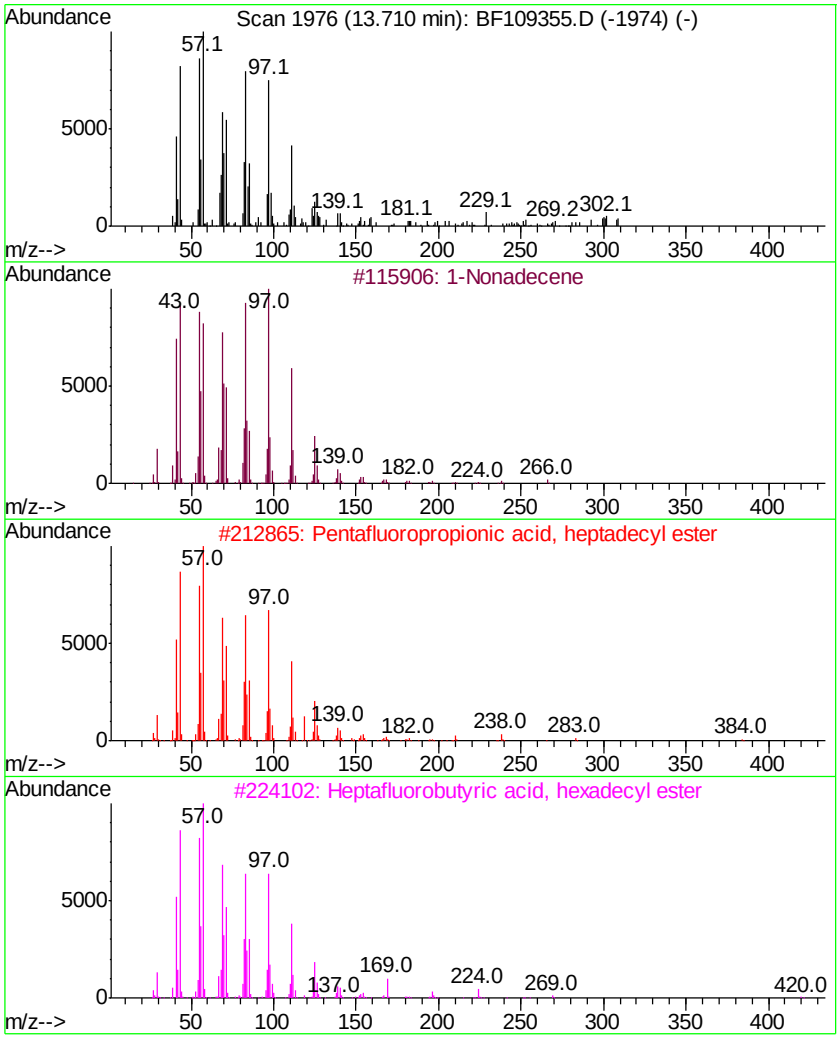
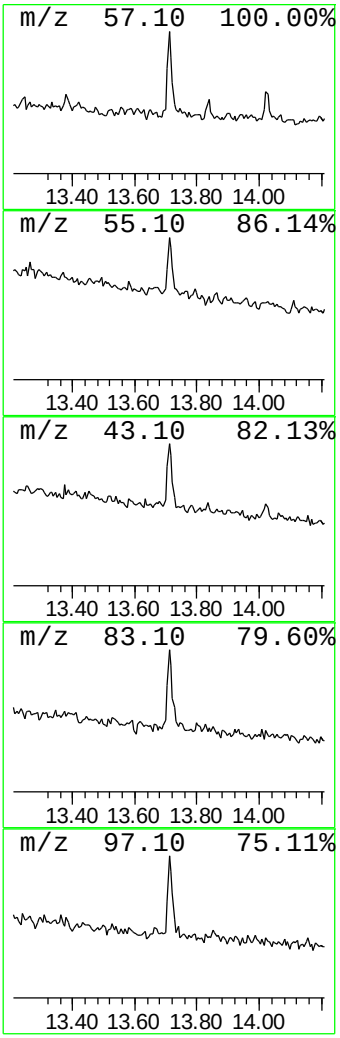
Peak Number 11 1-Nonadecene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.71	3.40 ng	132501	Chrysene-d12	13.85

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Nonadecene	266	C19H38	018435-45-5	97
2	Pentafluoropropionic acid, hepta...	402	C20H35F5O2	959218-78-1	93
3	Heptafluorobutyric acid, hexadec...	438	C20H33F7O2	006385-15-5	93
4	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	93
5	n-Tetracosanol-1	354	C24H50O	000506-51-4	91





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Operator : JU/SJ
Sample : J5055-20
Misc :
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Instrument : BNA_F
ClientSampleId : OILY-WATER

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TIC Integration Parameters: LSCINT.P

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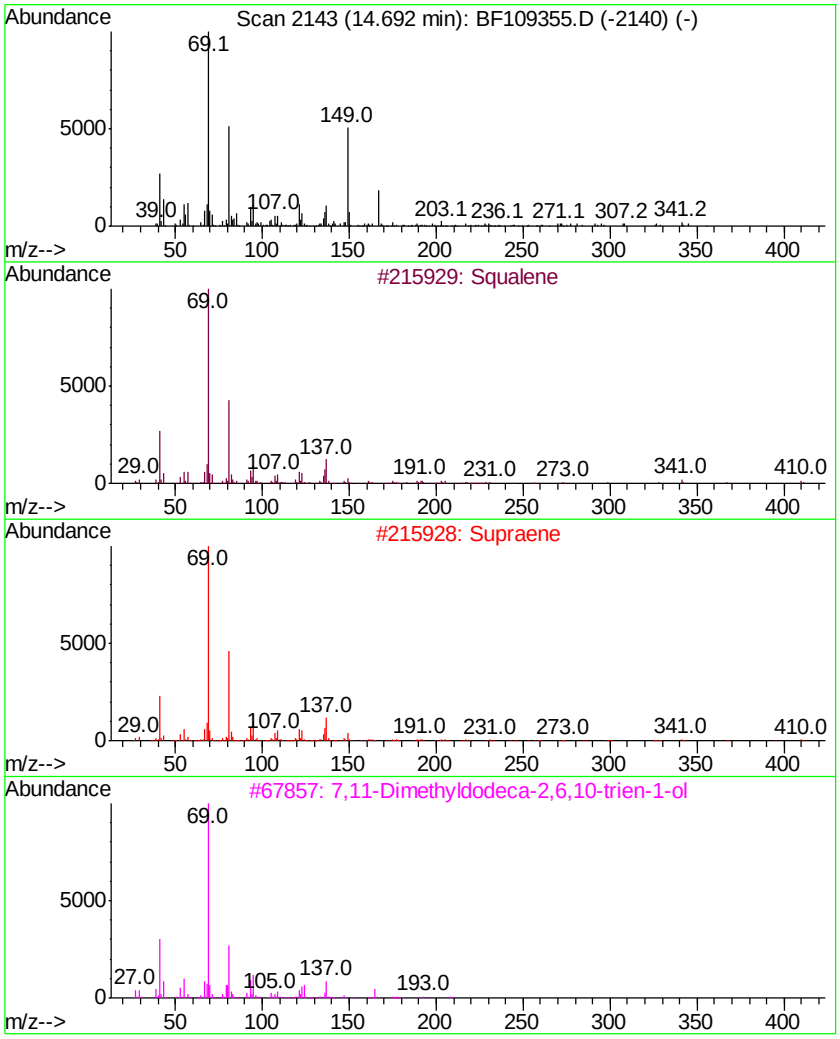
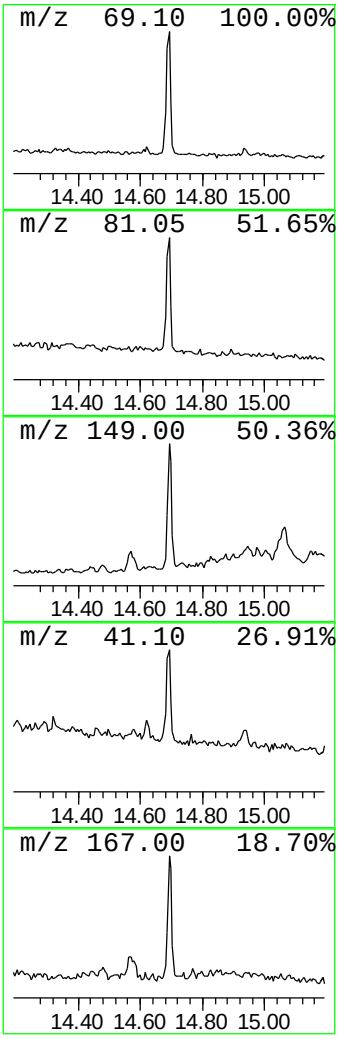
Peak Number 12 Squalene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.69	5.93 ng	189652	Perylene-d12	15.24

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Squalene	410	C30H50	000111-02-4	58
2	Supraene	410	C30H50	007683-64-9	49
3	7,11-Dimethyldodeca-2,6,10-trien...	208	C14H24O	125529-10-4	47
4	2,6-Octadiene, 4-methyl-	124	C9H16	074498-94-5	43
5	Trideca-1,7,11-triene-1,1-dicarb...	270	C18H26N2	1000161-46-0	38





Library Search Compound Report

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Data File : BF109355.D
Acq On : 26 Sep 2018 23:26
Operator : JU/SJ
Sample : J5055-20
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : OILY-WATER

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF092218.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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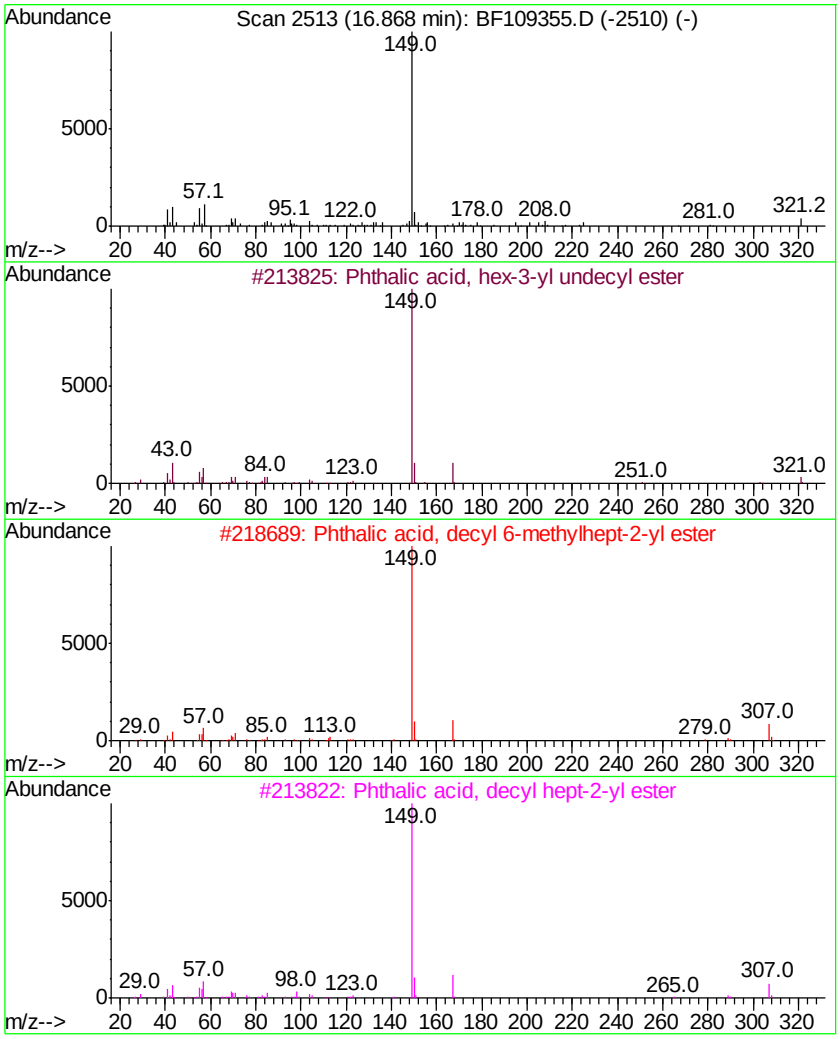
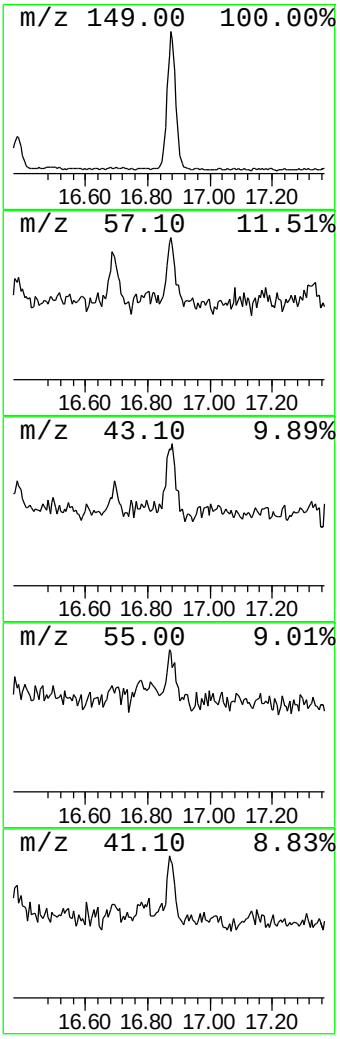
Peak Number 13 Phthalic acid, hex-3-yl und... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.87	3.19 ng	101990	Perylene-d12	15.24

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phthalic acid, hex-3-yl undecyl ...	404	C25H40O4	1000356-96-2	86
2	Phthalic acid, decyl 6-methylhep...	418	C26H42O4	1000377-96-4	72
3	Phthalic acid, decyl hept-2-yl e...	404	C25H40O4	1000356-97-7	72
4	Phthalic acid, 3-fluorophenyl tr...	442	C27H35FO4	1000356-66-7	72
5	Phthalic acid, monoamide, N-meth...	423	C27H37NO3	1000322-84-9	72





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Sample : J5055-20
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : OILY-WATER

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF092218.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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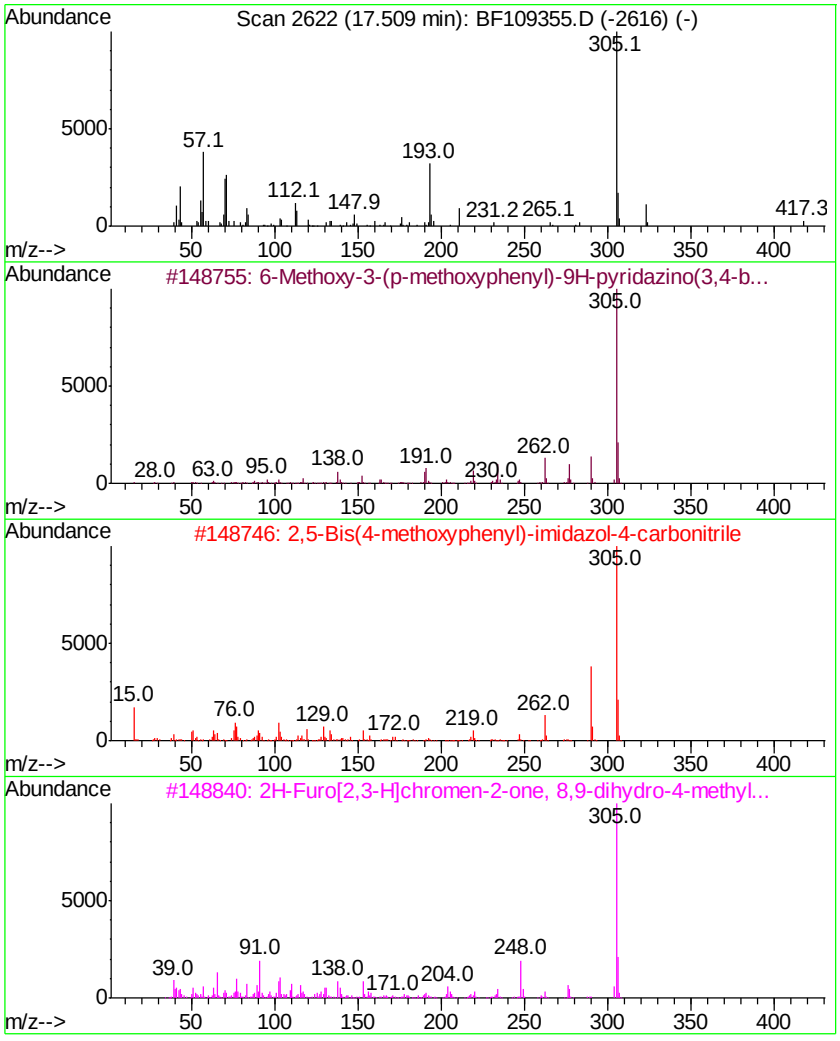
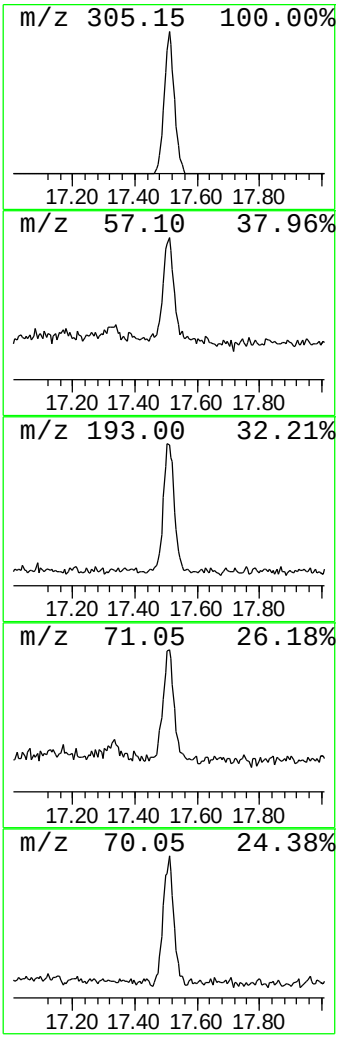
Peak Number 14 unknown17.51 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.51	6.05 ng	193543	Perylene-d12	15.24

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	6-Methoxy-3-(p-methoxyphenyl)-9H...	305	C18H15N3O2	004460-54-2	37
2	2,5-Bis(4-methoxyphenyl)-imidazo...	305	C18H15N3O2	189453-70-1	37
3	2H-Furo[2,3-H]chromen-2-one, 8,9...	305	C19H15NO3	1000271-14-0	25
4	6-Hydroxy-3-[2-methyl-2-[2-hydro...	595	C31H49N3O3Si3	1000292-78-8	25
5	tert-Butyldimethylsilyl 3-iodobe...	362	C13H19IO2Si	1000373-18-3	25





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF092618\
Data File : BF109355.D
Acq On : 26 Sep 2018 23:26
Operator : JU/SJ
Sample : J5055-20
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : OILY-WATER

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF092218.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
2-Pentanone, 4-hy...	4.89	2.1	ng	62485	1	6.71	589163	20.0
2-Propanol, 1-(2-...	6.64	2.4	ng	70242	1	6.71	589163	20.0
Benzene, 1,2-dich...	8.70	6.2	ng	219215	2	7.99	708421	20.0
2,7-Oxepanedione	8.80	6.2	ng	220048	2	7.99	708421	20.0
Benzenamine, 2,4-...	9.40	2.2	ng	75019	3	9.74	694397	20.0
Phenol, 2,3,6-tri...	9.97	2.4	ng	83307	3	9.74	694397	20.0
n-Hexadecanoic acid	11.77	8.0	ng	246506	4	11.22	615289	20.0
unknown12.50	12.50	4.4	ng	135089	4	11.22	615289	20.0
1-Nonadecene	13.71	3.4	ng	132501	5	13.85	779458	20.0
Squalene	14.69	5.9	ng	189652	6	15.24	639321	20.0
Phthalic acid, he...	16.87	3.2	ng	101990	6	15.24	639321	20.0
unknown17.51	17.51	6.0	ng	193543	6	15.24	639321	20.0