LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062218\
Data File : BF106752.D
Acq On : 23 Jun 2018 11:16
Operator : JU/SJ
Sample : J3597-03
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : W-44

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-BF061918.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	5.190	482	485	492	rBV	27254	28211	2.31%	0.303%
2	5.284	497	501	505	rBV	234117	217917	17.82%	2.340%
3	5.601	551	555	566	rBB	559318	479608	39.22%	5.149%
4	6.601	721	725	733	rBV	304277	295544	24.17%	3.173%
5	6.737	744	748	751	rBV	959827	816986	66.80%	8.772%
6	6.890	771	774	779	rVB2	19950	19478	1.59%	0.209%
7	6.960	782	786	791	rBV	291362	260691	21.32%	2.799%
8	7.113	808	812	816	rBV	834677	804696	65.80%	8.640%
9	7.525	877	882	885	rBV	672371	607717	49.69%	6.525%
10	7.595	892	894	900	rBV4	23029	27379	2.24%	0.294%
11	7.695	909	911	914	rVB2	16436	13432	1.10%	0.144%
12	8.242	999	1004	1008	rBV	300607	261924	21.42%	2.812%
13	8.389	1023	1029	1031	rBV2	15642	18318	1.50%	0.197%
14	8.607	1059	1066	1073	rBV5	16007	33277	2.72%	0.357%
15	8.689	1075	1080	1083	rBV4	16014	24491	2.00%	0.263%
16	8.795	1096	1098	1103	rVB	59471	52672	4.31%	0.566%
17	8.842	1103	1106	1108	rBV2	15874	16690	1.36%	0.179%
18	8.972	1125	1128	1131	rBV3	21591	20602	1.68%	0.221%
19	9.001	1131	1133	1137	rVB3	15056	17693	1.45%	0.190%
20	9.054	1137	1142	1144	rBV4	11942	20656	1.69%	0.222%
21	9.078	1144	1146	1152	rVB5	24890	28650	2.34%	0.308%
22	9.131	1152	1155	1156	rBV2	13749	13477	1.10%	0.145%
23	9.160	1156	1160	1162	rVV2	13324	19336	1.58%	0.208%
24	9.184	1162	1164	1167	rVB4	19231	19559	1.60%	0.210%
25	9.219	1167	1170	1173	rBV4	16381	15985	1.31%	0.172%
26	9.272	1173	1179	1182	rBV7	16051	24361	1.99%	0.262%
27	9.313	1182	1186	1190	rBV	1109934	992895	81.19%	10.660%
28	9.660	1243	1245	1247	rBV3	27758	23487	1.92%	0.252%
29	9.689	1247	1250	1251	rVV	43182	38549	3.15%	0.414%
30	9.707	1251	1253	1258	rVB	45108	42909	3.51%	0.461%
31	9.789	1264	1267	1269	rBV	38536	26177	2.14%	0.281%
32	9.954	1289	1295	1299	rBV5	6781	16053	1.31%	0.172%
33	10.001	1299	1303	1309	rVB	287255	255326	20.88%	2.741%
34	10.307	1353	1355	1358	rBV4	17319	16931	1.38%	0.182%

LSC Area Percent Report

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Acq On : 23 Jun 2018 11:16
Operator : JU/SJ
Sample : J3597-03
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : W-44

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-BF061918.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	10.383	1366	1368	1371	rBV3	16631	15423	1.26%	0.166%
36	10.489	1383	1386	1389	rBV3	12154	14538	1.19%	0.156%
37	10.654	1411	1414	1417	rVB3	24759	22338	1.83%	0.240%
38	10.713	1422	1424	1431	rVB5	13102	17109	1.40%	0.184%
39	10.795	1434	1438	1441	rBV	710321	627995	51.35%	6.743%
40	10.883	1449	1453	1456	rBV	12560	13289	1.09%	0.143%
41	11.036	1474	1479	1482	rVB6	11782	15360	1.26%	0.165%
42	11.078	1484	1486	1491	rBV3	11616	12705	1.04%	0.136%
43	11.489	1553	1556	1560	rBV	290926	267231	21.85%	2.869%
44	12.877	1789	1792	1796	rVB2	25837	24202	1.98%	0.260%
45	12.960	1803	1806	1814	rVB	131148	116560	9.53%	1.251%
46	13.077	1822	1826	1829	rBV	1257338	1222949	100.00%	13.130%
47	13.583	1910	1912	1915	rVB2	16705	14867	1.22%	0.160%
48	13.624	1915	1919	1922	rBV	14747	15638	1.28%	0.168%
49	13.960	1973	1976	1983	rBV	26757	30257	2.47%	0.325%
50	14.148	2004	2008	2012	rBV	354231	339054	27.72%	3.640%
51	14.283	2027	2031	2035	rVB	38740	34928	2.86%	0.375%
52	14.595	2080	2084	2089	rBV	82721	83114	6.80%	0.892%
53	14.918	2135	2139	2143	rVB	116628	121596	9.94%	1.306%
54	15.271	2195	2199	2204	rBV	132577	145848	11.93%	1.566%
55	15.683	2263	2269	2275	rVB2	226764	373863	30.57%	4.014%
56	16.136	2341	2346	2353	rVB	75843	117160	9.58%	1.258%
57	16.671	2432	2437	2443	rVB	35285	59034	4.83%	0.634%
58	17.307	2541	2545	2554	rVB4	11984	24015	1.96%	0.258%
59	18.042	2666	2670	2676	rBV8	5628	13072	1.07%	0.140%

Sum of corrected areas: 9313822

LSC Report - Integrated Chromatogram

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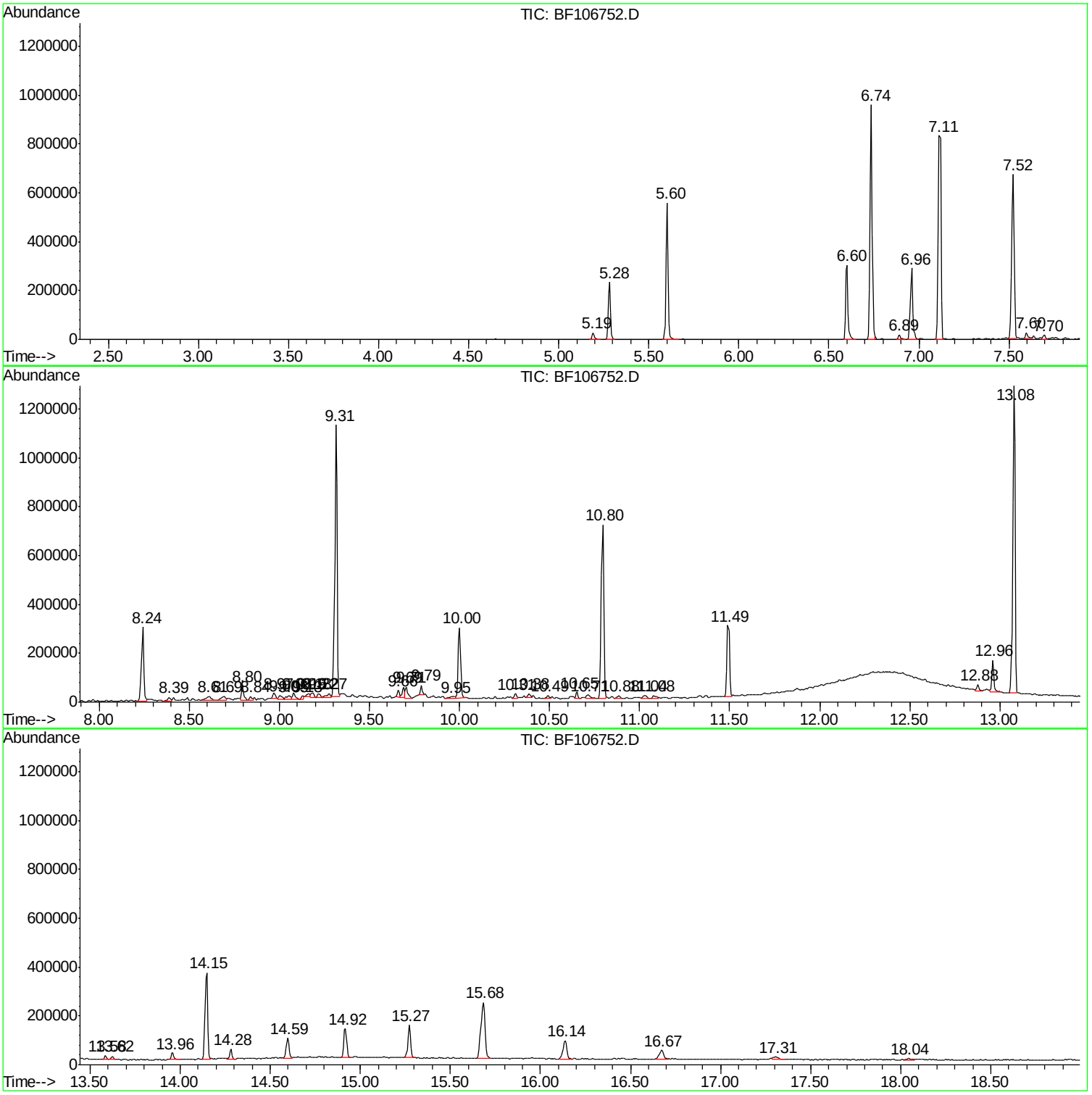
Acq On : 23 Jun 2018 11:16
Operator : JU/SJ
Sample : J3597-03
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : W-44

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.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

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062218\
Data File : BF106752.D
Acq On : 23 Jun 2018 11:16
Operator : JU/SJ
Sample : J3597-03
Misc :
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Instrument : BNA_F
ClientSampleId : W-44

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

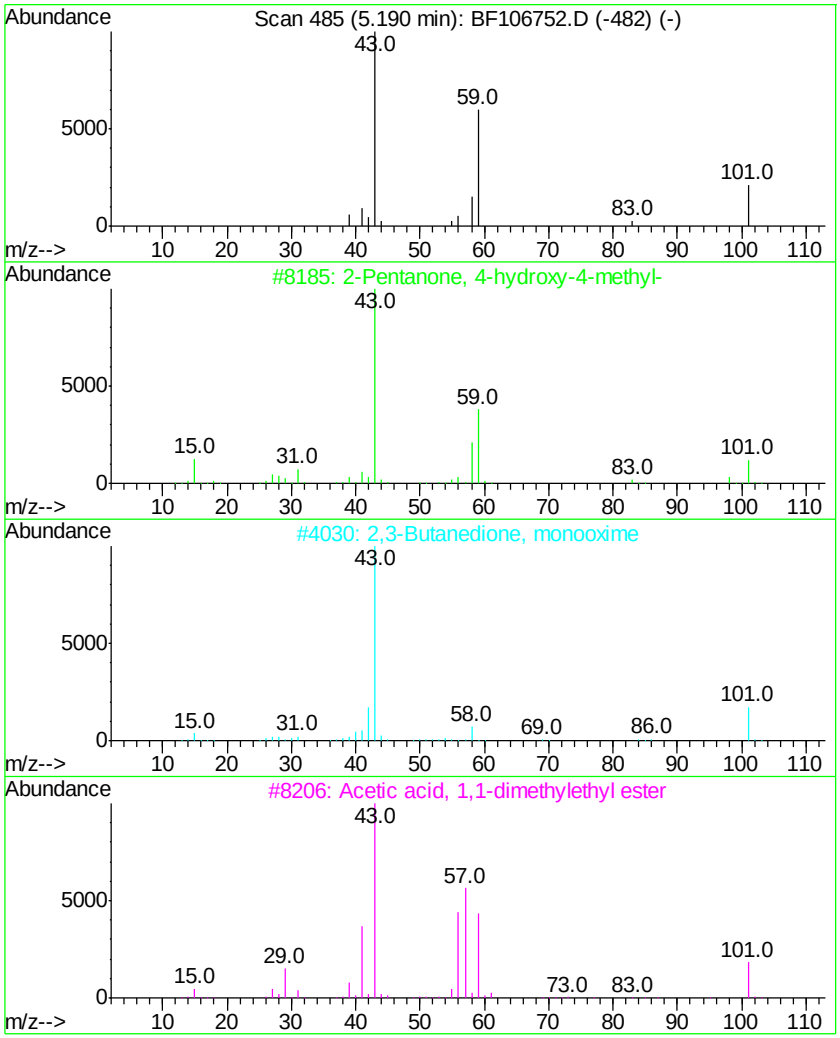
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

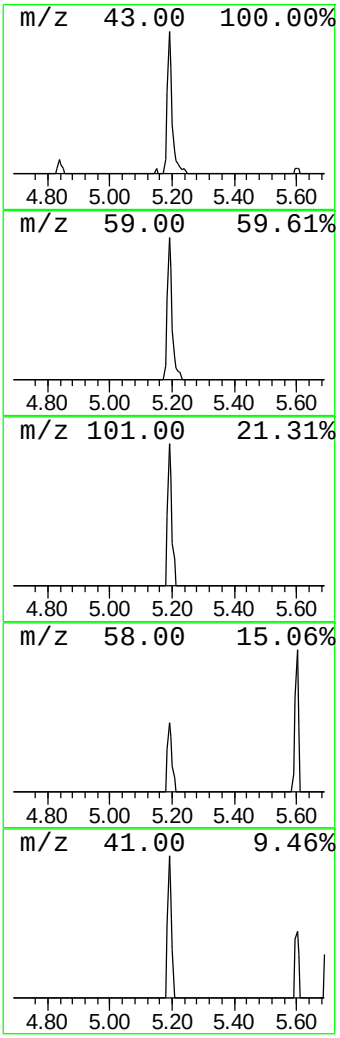
****

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.19	2.16 ng	28211	1,4-Dichlorobenzene-d4	6.96

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
3	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	9
4	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9
5	3-Hexanol	102	C6H14O	000623-37-0	9





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062218\
Data File : BF106752.D
Acq On : 23 Jun 2018 11:16
Operator : JU/SJ
Sample : J3597-03
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : W-44

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

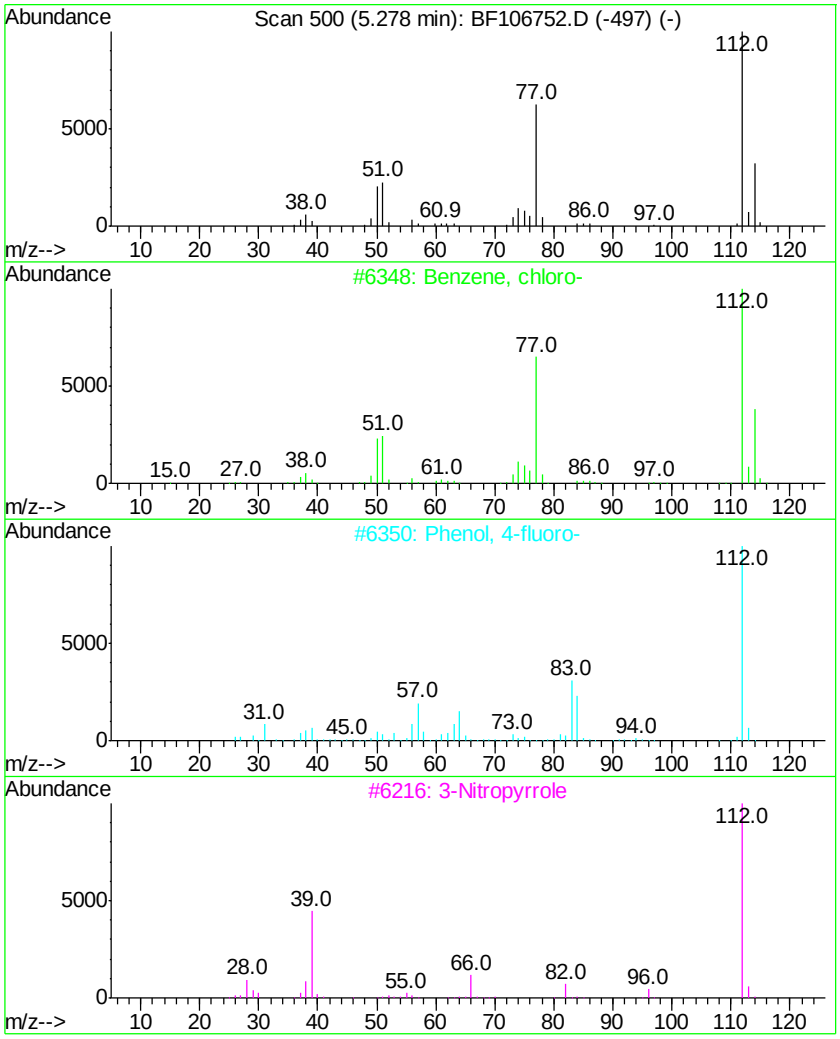
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

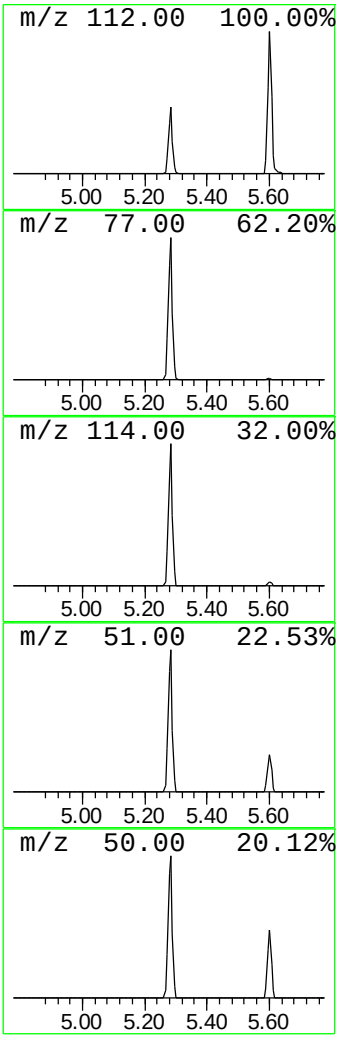
****

Peak Number 2 Benzene, chloro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.28	16.72 ng	217917	1,4-Dichlorobenzene-d4	6.96

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, chloro-	112	C6H5Cl	000108-90-7	96
2	Phenol, 4-fluoro-	112	C6H5FO	000371-41-5	17
3	3-Nitropyrrole	112	C4H4N2O2	005930-94-9	9
4	Cevane-3,4,6,7,14,15,16,20-octol...	793	C41H63NO14	000143-57-7	9
5	Dipipanone	349	C24H31NO	000467-83-4	9





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062218\
Data File : BF106752.D
Acq On : 23 Jun 2018 11:16
Operator : JU/SJ
Sample : J3597-03
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : W-44

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

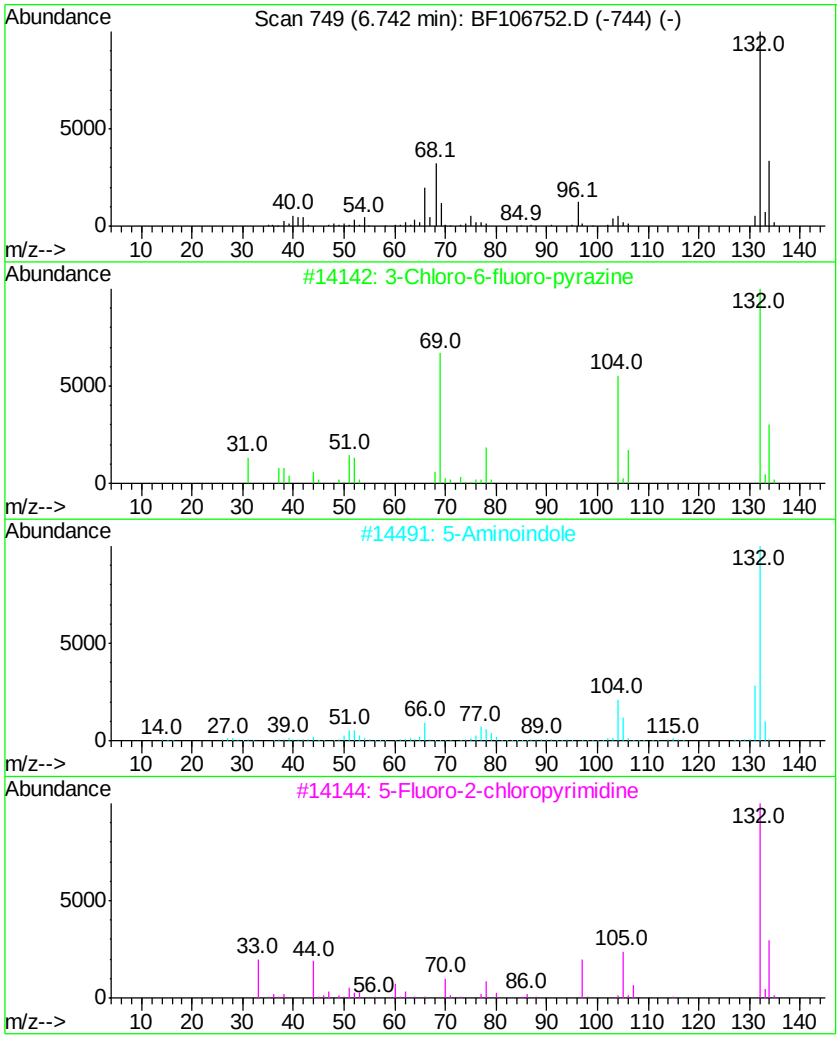
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

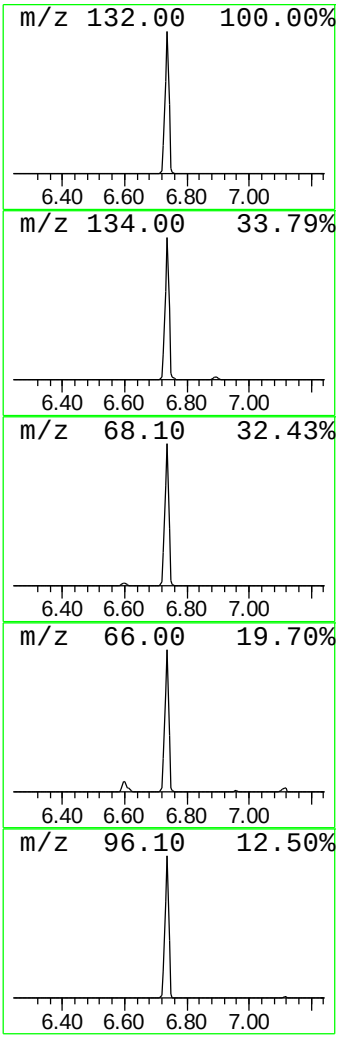
****

Peak Number 3 unknown6.74 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.74	62.68 ng	816986	1,4-Dichlorobenzene-d4	6.96

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	12
3	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
4	2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9
5	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	9





Library Search Compound Report

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Sample : J3597-03
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : W-44

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

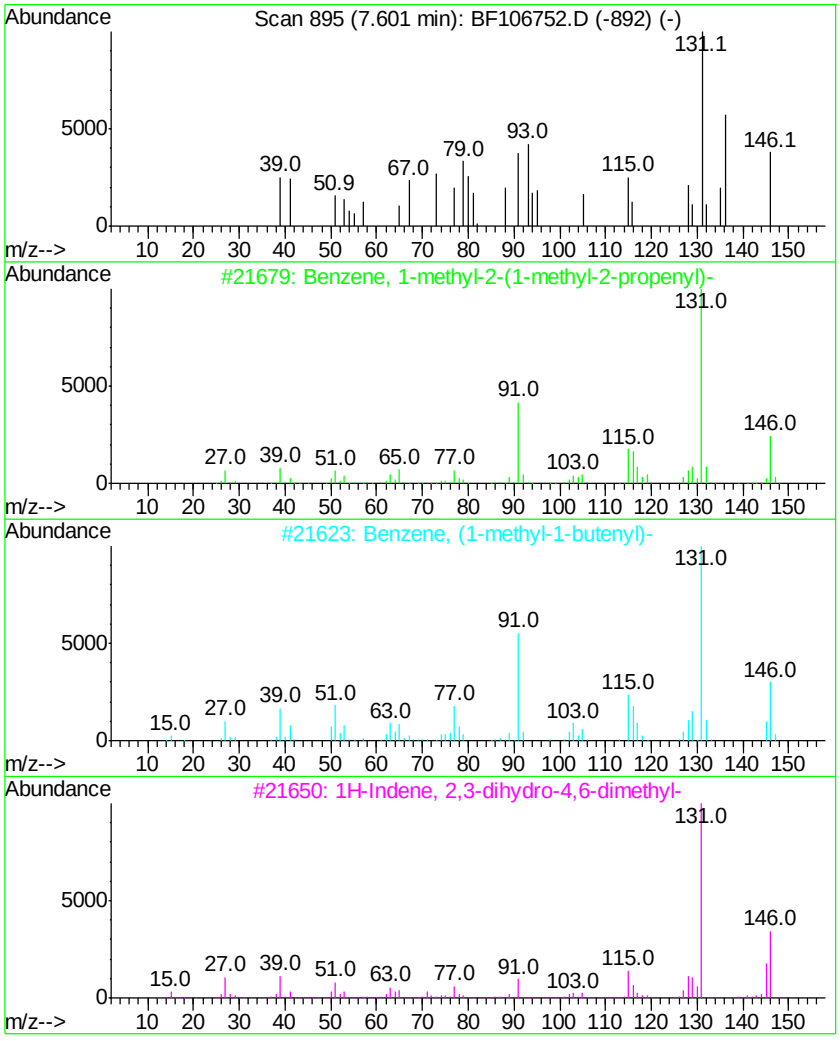
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

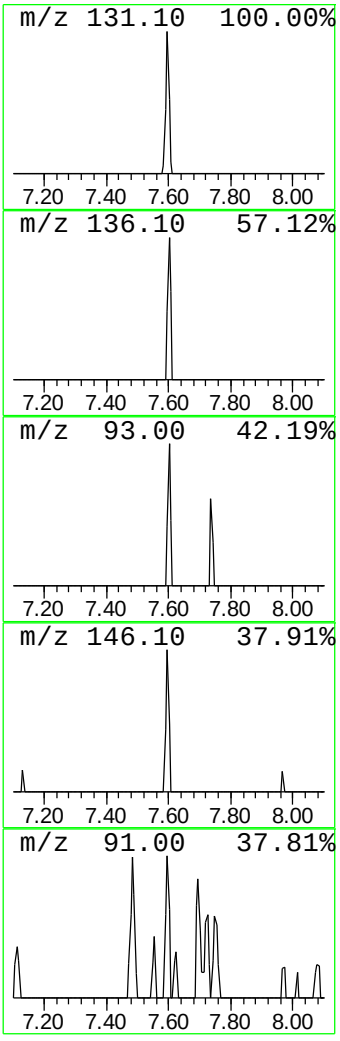
****

Peak Number 5 Benzene, 1-methyl-2-(1-meth... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.60	2.10 ng	27379	1,4-Dichlorobenzene-d4	6.96

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-2-(1-methyl-2-...	146	C11H14	097664-19-2	74
2	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	64
3	1H-Indene, 2,3-dihydro-4,6-dimet...	146	C11H14	001685-82-1	64
4	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	53
5	1,3-Cyclopentadiene, 5-(trans-2-...	146	C11H14	079209-36-2	50





Library Search Compound Report

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Misc :
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Instrument : BNA_F
ClientSampleId : W-44

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

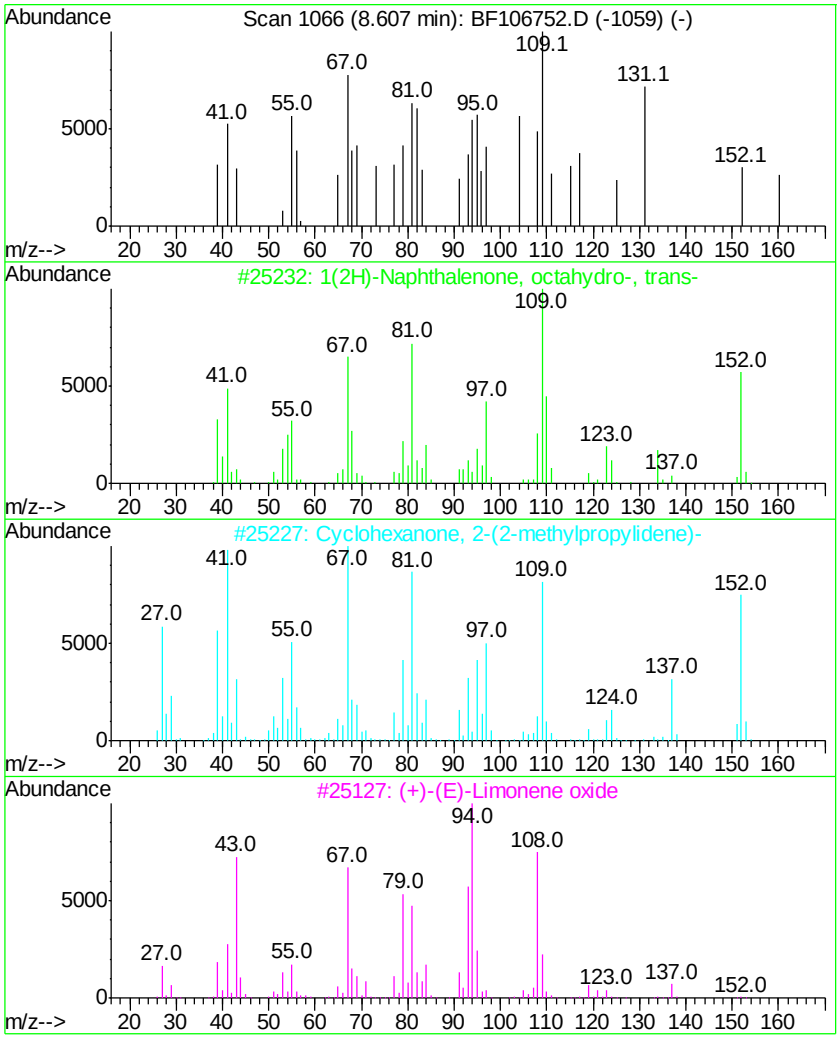
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

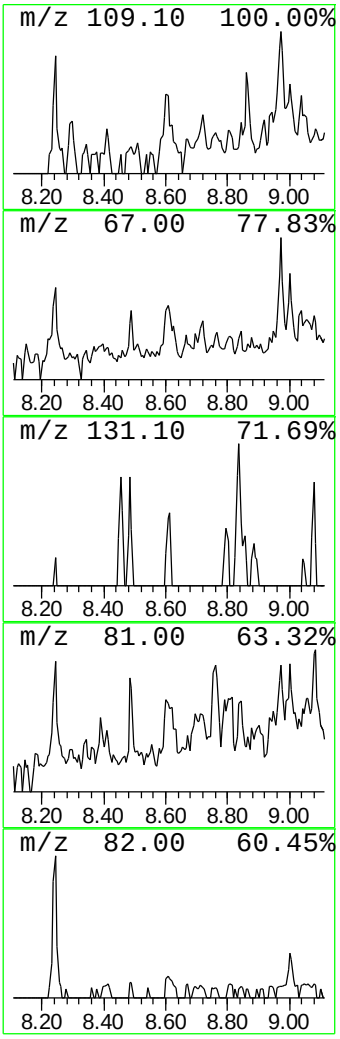
****

Peak Number 6 unknown8.61 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.61	2.54 ng	33277	Naphthalene-d8	8.24

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1(2H)-Naphthalenone, octahydro-,...	152	C10H16O	021370-71-8	43
2	Cyclohexanone, 2-(2-methylpropyl...	152	C10H16O	043108-69-6	35
3	(+)-(E)-Limonene oxide	152	C10H16O	006909-30-4	27
4	2,4-Bis(hydroxylamino)pyrimidine	142	C4H6N4O2	154845-38-2	27
5	Naphth[2,3-b]oxirene, decahydro-	152	C10H16O	021399-51-9	27





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ClientSampleId : W-44

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

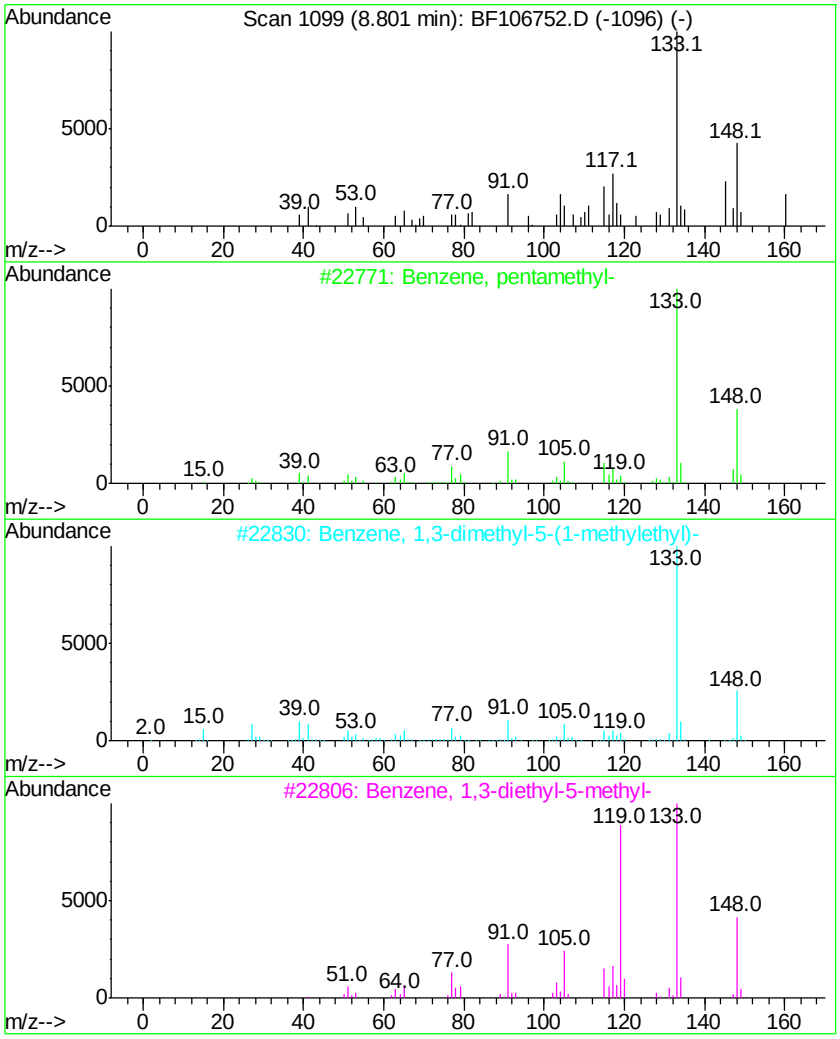
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

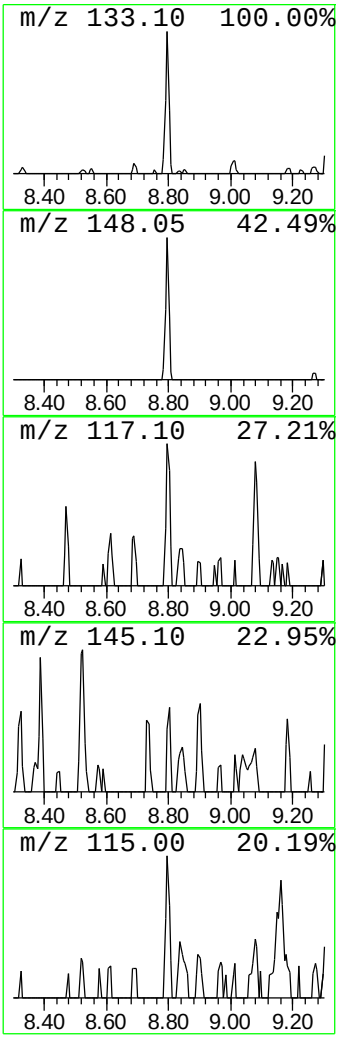
****

Peak Number 7 Benzene, pentamethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.80	4.02 ng	52672	Naphthalene-d8	8.24

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, pentamethyl-	148	C11H16	000700-12-9	95
2	Benzene, 1,3-dimethyl-5-(1-methy...	148	C11H16	004706-90-5	87
3	Benzene, 1,3-diethyl-5-methyl-	148	C11H16	002050-24-0	86
4	3,4-Dimethylcumene	148	C11H16	1000370-34-1	83
5	Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	80





Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062218\
Data File : BF106752.D
Acq On : 23 Jun 2018 11:16
Operator : JU/SJ
Sample : J3597-03
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : W-44

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

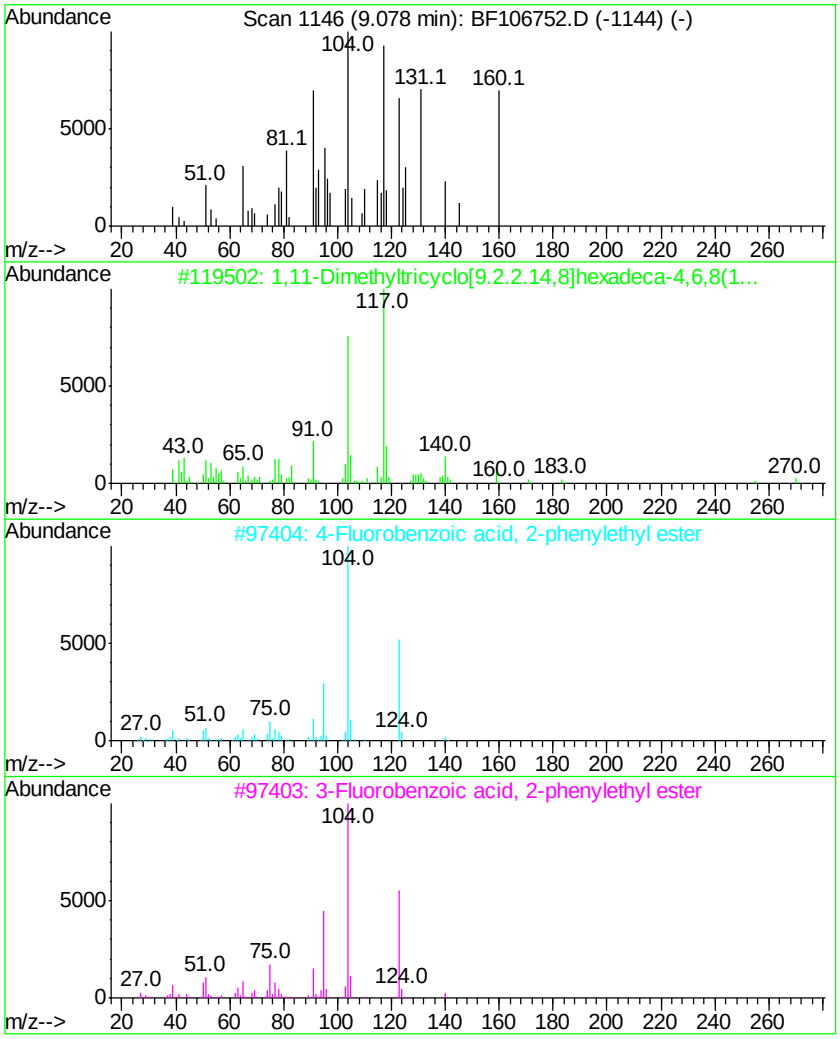
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

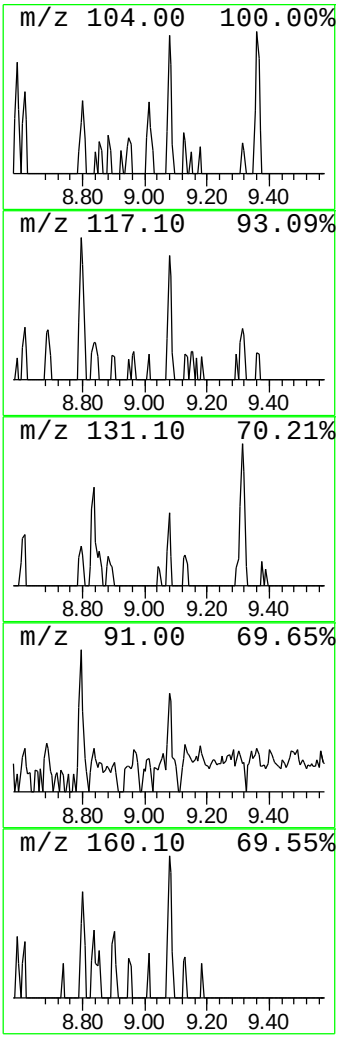
****

Peak Number 8 unknown9.08 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.08	2.19 ng	28650	Naphthalene-d8	8.24

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,11-Dimethyltricyclo[9.2.2.14,8...	270	C18H22O2	1000311-68-2	16
2	4-Fluorobenzoic acid, 2-phenylet...	244	C15H13FO2	090172-19-3	16
3	3-Fluorobenzoic acid, 2-phenylet...	244	C15H13FO2	1000279-04-8	16
4	2-Fluorobenzoic acid, 2-phenylet...	244	C15H13FO2	1000278-98-3	16
5	Benzene, (3-methyl-4-pentenyl)-	160	C12H16	042524-30-1	10





Library Search Compound Report

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Sample : J3597-03
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : W-44

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.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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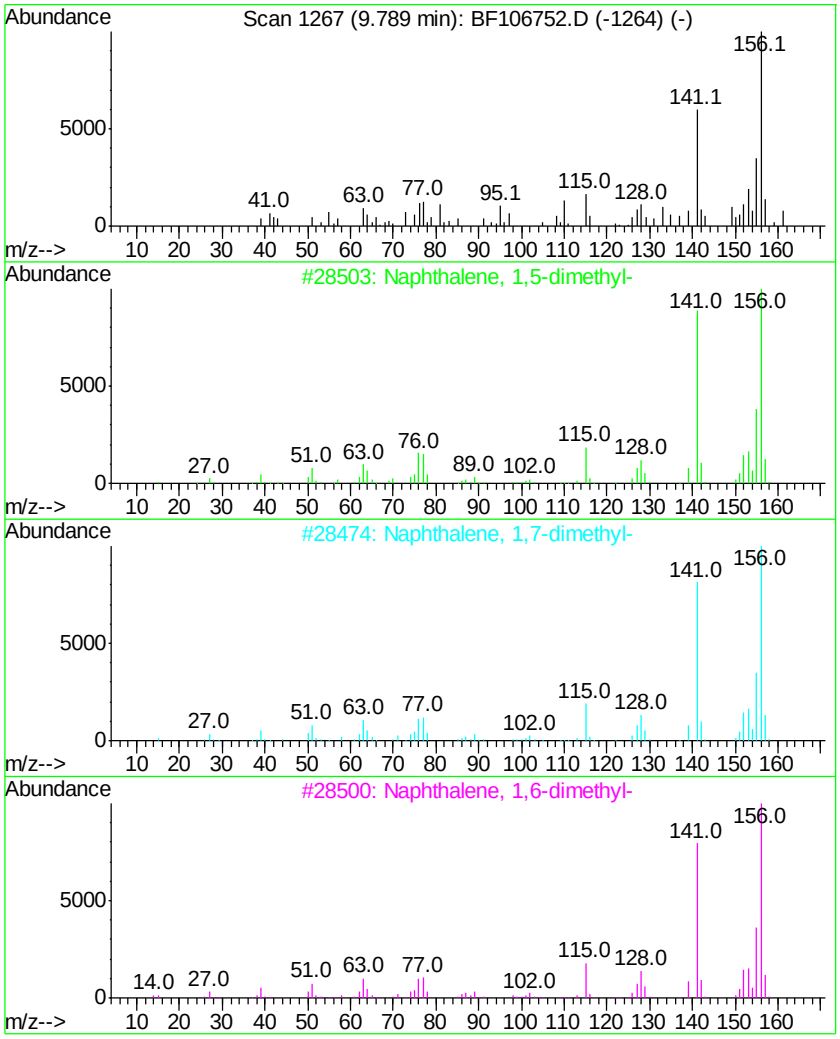
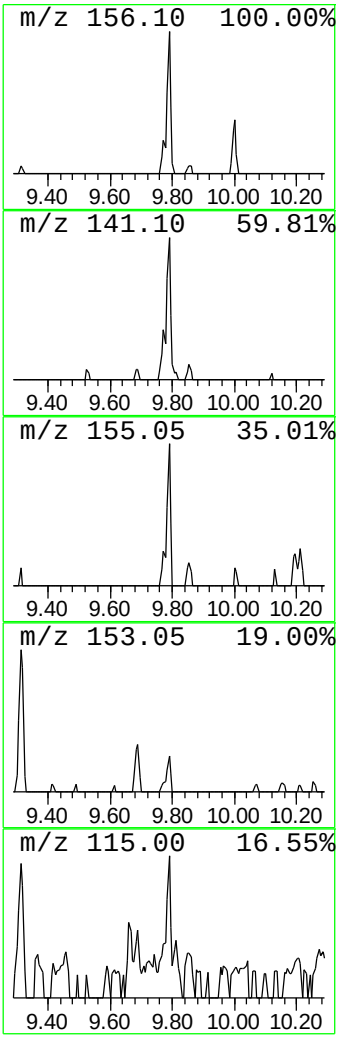
Peak Number 9 Naphthalene, 1,5-dimethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.79	2.05 ng	26177	Acenaphthene-d10	10.00

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	98
2	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	98
3	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
4	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
5	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96





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Acq On : 23 Jun 2018 11:16
Operator : JU/SJ
Sample : J3597-03
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : W-44

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.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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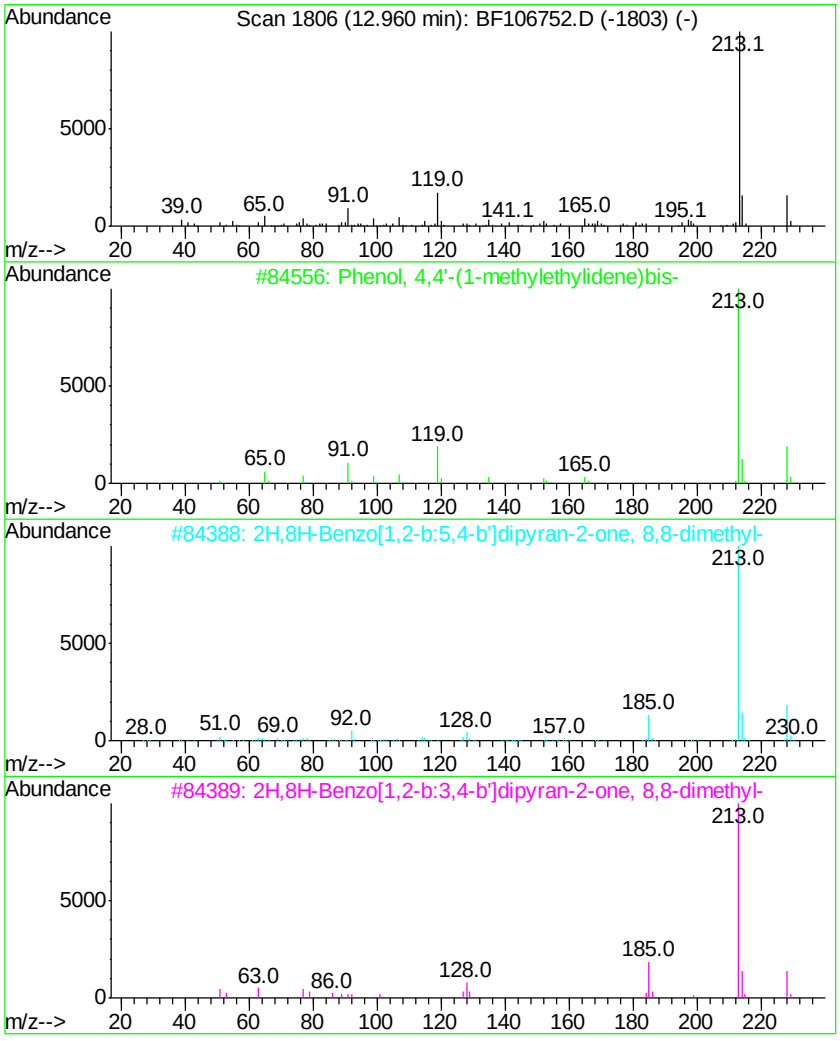
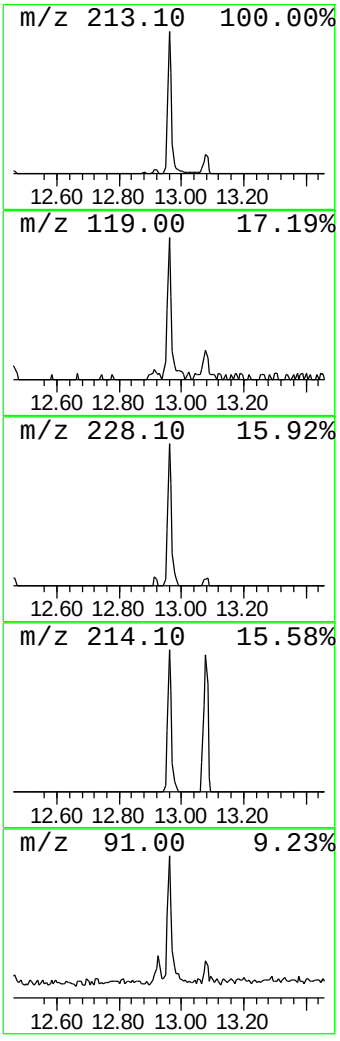
Peak Number 10 Phenol, 4,4'-(1-methylethyl... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.96	6.88 ng	116560	Chrysene-d12	14.15

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	98
2	2H,8H-Benzo[1,2-b:5,4-b']dipyran...	228	C14H12O3	000553-19-5	64
3	2H,8H-Benzo[1,2-b:3,4-b']dipyran...	228	C14H12O3	000523-59-1	59
4	4H-Pyrazolo[5,1-c]-1,4-benzodiaz...	213	C11H7N3O2	1000277-20-1	59
5	Decanedioic acid, diethyl ester	258	C14H26O4	000110-40-7	53





Library Search Compound Report

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Data File : BF106752.D
Acq On : 23 Jun 2018 11:16
Operator : JU/SJ
Sample : J3597-03
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : W-44

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TIC Integration Parameters: LSCINT.P

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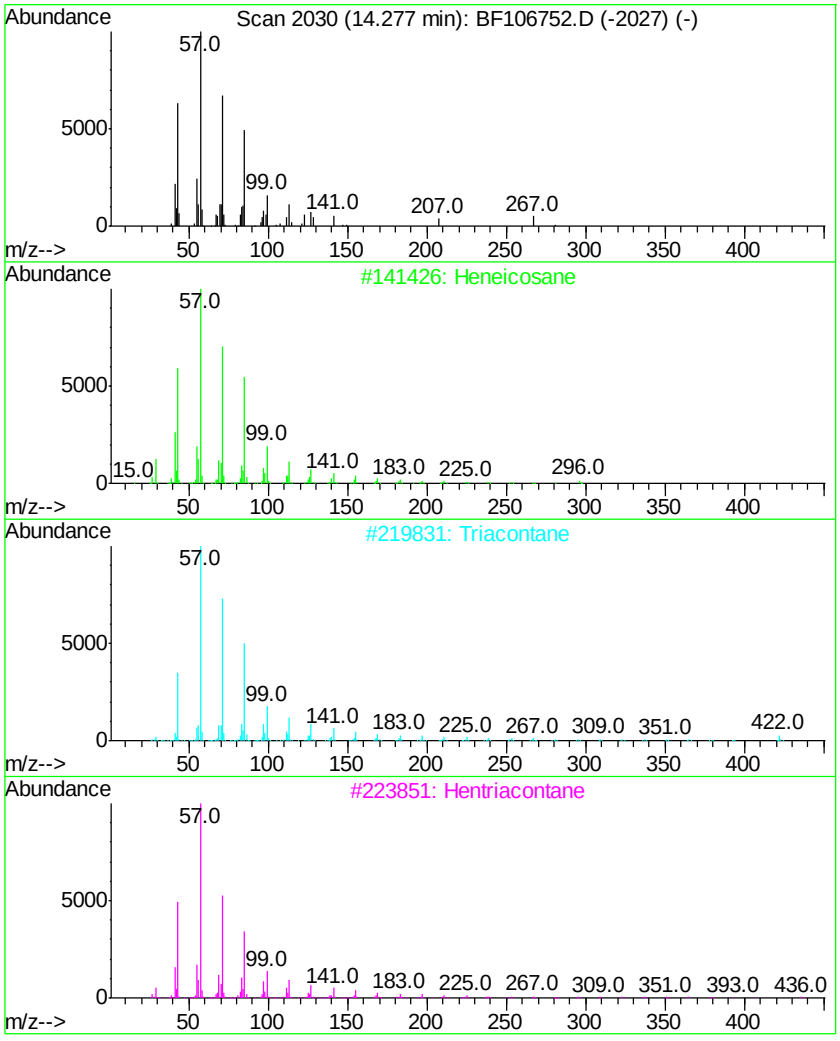
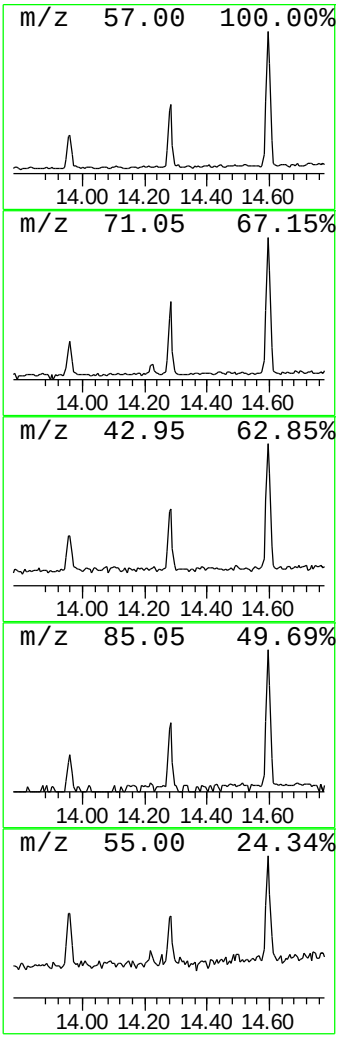
Peak Number 11 Heneicosane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.28	2.06 ng	34928	Chrysene-d12	14.15

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	90
2	Triacontane	422	C30H62	000638-68-6	86
3	Hentriacontane	437	C31H64	000630-04-6	86
4	Tetracosane	338	C24H50	000646-31-1	86
5	Octadecane	254	C18H38	000593-45-3	86





Library Search Compound Report

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Acq On : 23 Jun 2018 11:16
Operator : JU/SJ
Sample : J3597-03
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : W-44

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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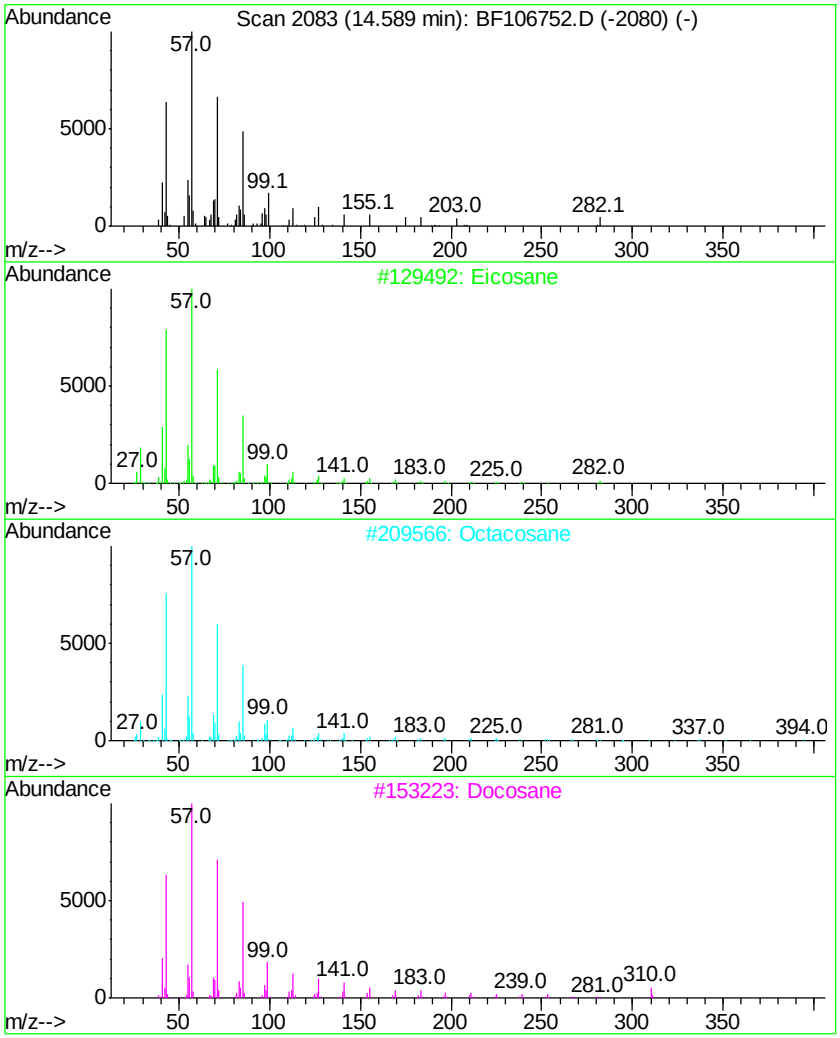
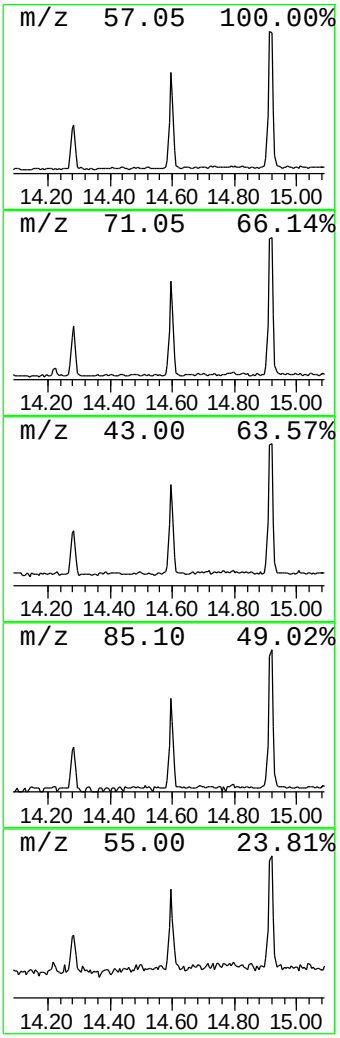
Peak Number 12 Eicosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.59	4.90 ng	83114	Chrysene-d12	14.15

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	97
2	Octacosane	394	C28H58	000630-02-4	91
3	Docosane	310	C22H46	000629-97-0	91
4	2-methyloctacosane	408	C29H60	1000376-72-8	91
5	Heneicosane	296	C21H44	000629-94-7	91





Library Search Compound Report

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Acq On : 23 Jun 2018 11:16
Operator : JU/SJ
Sample : J3597-03
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : W-44

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.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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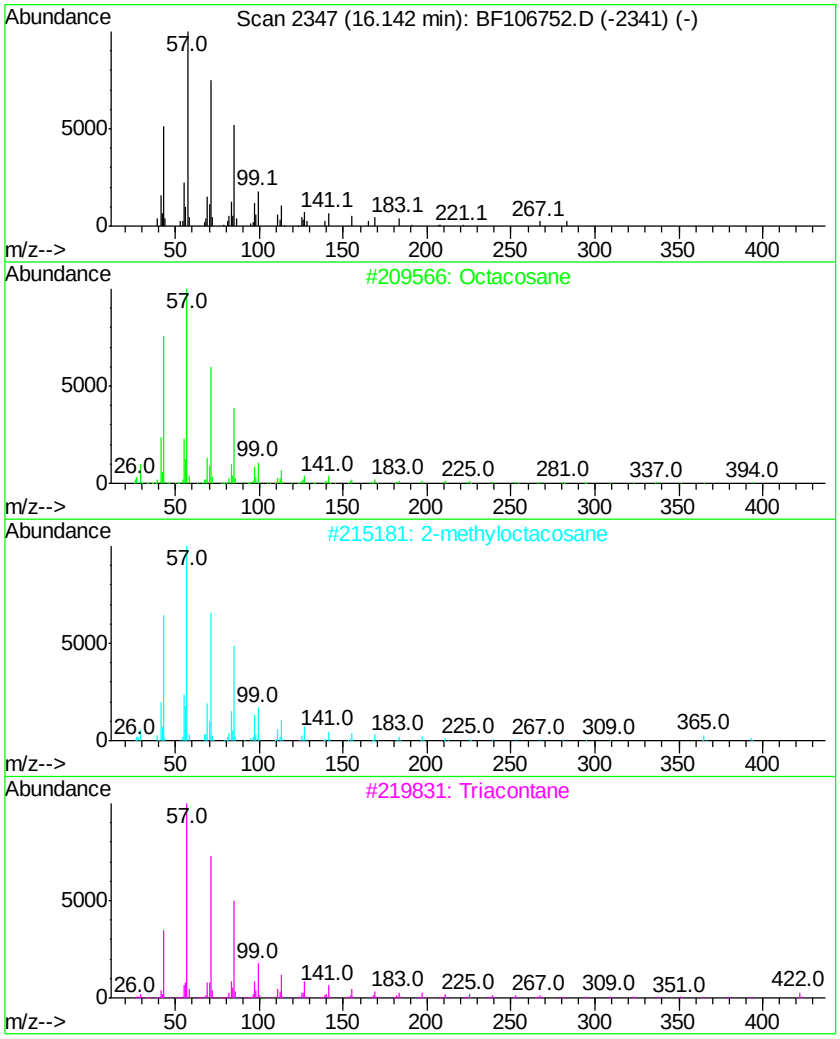
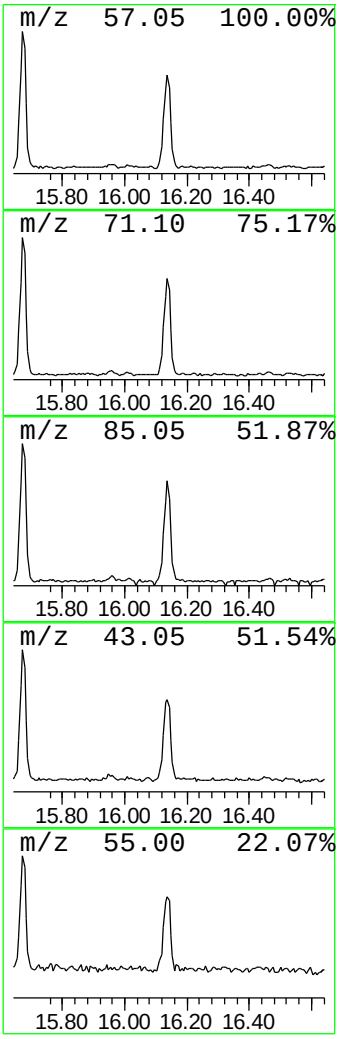
Peak Number 15 Octacosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.14	6.27 ng	117160	Perylene-d12	15.69

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octacosane	394	C28H58	000630-02-4	91
2	2-methyloctacosane	408	C29H60	1000376-72-8	91
3	Triacontane	422	C30H62	000638-68-6	91
4	Heneicosane	296	C21H44	000629-94-7	91
5	Octadecane	254	C18H38	000593-45-3	90





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Operator : JU/SJ
Sample : J3597-03
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : W-44

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.M
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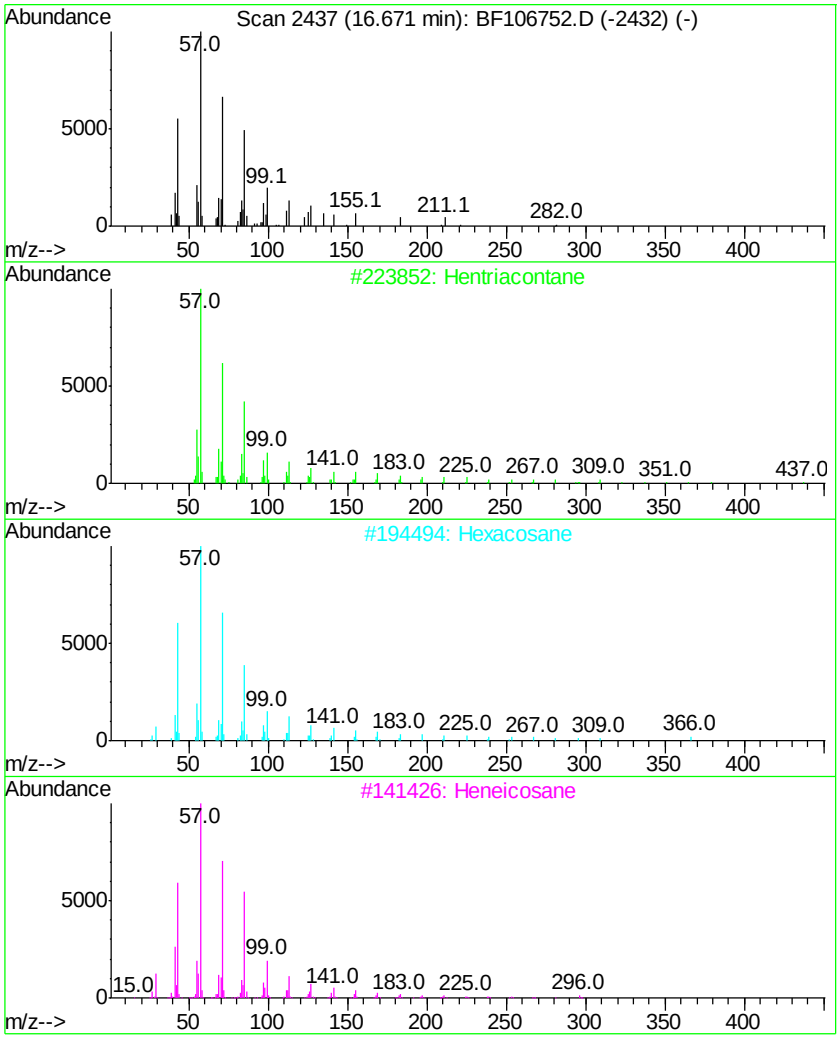
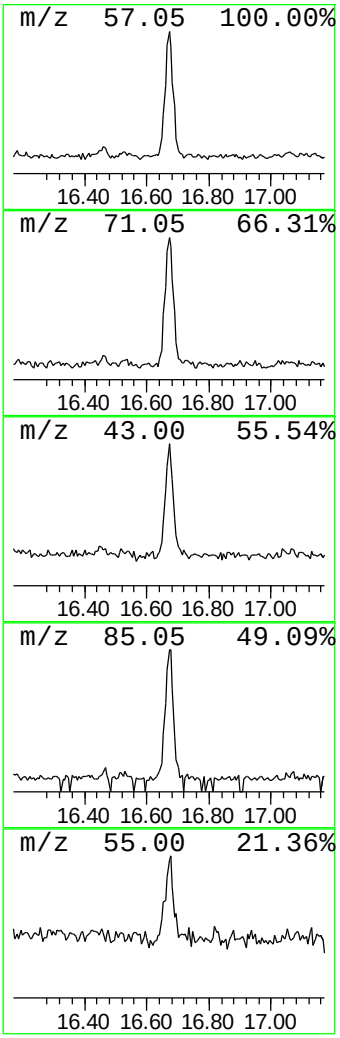
Peak Number 16 Hentriacontane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.67	3.16 ng	59034	Perylene-d12	15.69

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hentriacontane	437	C31H64	000630-04-6	91
2	Hexacosane	366	C26H54	000630-01-3	91
3	Heneicosane	296	C21H44	000629-94-7	91
4	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	90
5	Nonadecane, 2-methyl-	282	C20H42	001560-86-7	90





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF062218\
Data File : BF106752.D
Acq On : 23 Jun 2018 11:16
Operator : JU/SJ
Sample : J3597-03
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : W-44

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.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...	5.19	2.2	ng	28211	1	6.96	260691	20.0
Benzene, chloro-	5.28	16.7	ng	217917	1	6.96	260691	20.0
unknown6.74	6.74	62.7	ng	816986	1	6.96	260691	20.0
Benzene, 1-methyl...	7.60	2.1	ng	27379	1	6.96	260691	20.0
unknown8.61	8.61	2.5	ng	33277	2	8.24	261924	20.0
Benzene, pentamet...	8.80	4.0	ng	52672	2	8.24	261924	20.0
unknown9.08	9.08	2.2	ng	28650	2	8.24	261924	20.0
Naphthalene, 1,5-...	9.79	2.0	ng	26177	3	10.00	255326	20.0
Phenol, 4,4'-(1-m...	12.96	6.9	ng	116560	5	14.15	339054	20.0
Heneicosane	14.28	2.1	ng	34928	5	14.15	339054	20.0
Eicosane	14.59	4.9	ng	83114	5	14.15	339054	20.0
Octacosane	16.14	6.3	ng	117160	6	15.69	373863	20.0
Hentriacontane	16.67	3.2	ng	59034	6	15.69	373863	20.0