LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF121217\
Data File : BF101334.D
Acq On : 12 Dec 2017 17:16
Operator : SJ/JU
Sample : I6872-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 293-311

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-BF113017.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	5.040	499	502	513	rBV	70310	100062	8.25%	0.926%
2	5.445	567	571	588	rBV	1020552	959084	79.10%	8.874%
3	6.463	739	744	754	rBV	1075539	1096004	90.39%	10.140%
4	6.592	762	766	770	rBV	1123151	1068375	88.11%	9.885%
5	6.822	801	805	811	rBV	349934	274793	22.66%	2.542%
6	6.981	828	832	835	rBV	1010908	878608	72.46%	8.129%
7	7.222	865	873	881	rBV2	19321	34032	2.81%	0.315%
8	7.392	897	902	905	rBV	717446	663897	54.75%	6.142%
9	7.845	976	979	983	rBV2	11371	14731	1.21%	0.136%
10	8.075	1016	1018	1020	rBV	32512	28265	2.33%	0.262%
11	8.104	1020	1023	1029	rVV	431188	376394	31.04%	3.482%
12	8.151	1029	1031	1034	rVB	33258	28268	2.33%	0.262%
13	8.386	1065	1071	1076	rBV5	25015	36094	2.98%	0.334%
14	8.516	1089	1093	1095	rBV	54190	45075	3.72%	0.417%
15	8.686	1119	1122	1124	rBV	49293	34484	2.84%	0.319%
16	8.775	1135	1137	1142	rVB3	14743	20551	1.69%	0.190%
17	8.969	1167	1170	1174	rBV4	10201	16396	1.35%	0.152%
18	9.051	1180	1184	1186	rBV3	12136	15050	1.24%	0.139%
19	9.116	1192	1195	1198	rBV	56935	44519	3.67%	0.412%
20	9.180	1201	1206	1209	rBV	1506241	1212509	100.00%	11.218%
21	9.251	1215	1218	1221	rBV	78480	63576	5.24%	0.588%
22	9.328	1229	1231	1234	rVB3	18045	14884	1.23%	0.138%
23	9.575	1268	1273	1275	rBV	168965	171401	14.14%	1.586%
24	9.722	1294	1298	1301	rBV5	15706	17871	1.47%	0.165%
25	9.780	1302	1308	1311	rVV	73448	64343	5.31%	0.595%
26	9.863	1318	1322	1325	rVB	418734	333713	27.52%	3.088%
27	10.004	1342	1346	1350	rBV6	13298	21556	1.78%	0.199%
28	10.104	1360	1363	1367	rVB2	20551	21132	1.74%	0.196%
29	10.280	1386	1393	1396	rBV	79025	72068	5.94%	0.667%
30	10.410	1409	1415	1418	rBV4	10570	15075	1.24%	0.139%
31	10.498	1425	1430	1433	rBV	40269	49453	4.08%	0.458%
32	10.657	1453	1457	1465	rBV	539623	469210	38.70%	4.341%
33	10.757	1471	1474	1482	rVB2	69410	104974	8.66%	0.971%
34	11.204	1546	1550	1552	rBV	36239	29523	2.43%	0.273%

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF121217\
Data File : BF101334.D
Acq On : 12 Dec 2017 17:16
Operator : SJ/JU
Sample : I6872-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 293-311

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-BF113017.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	11.233	1552	1555	1557	rVV2	22670	20817	1.72%	0.193%
36	11.351	1572	1575	1578	rBV	339660	322281	26.58%	2.982%
37	11.374	1578	1579	1584	rVB	56968	45499	3.75%	0.421%
38	11.627	1619	1622	1626	rVB	20157	17911	1.48%	0.166%
39	11.868	1658	1663	1665	rBV3	19490	30431	2.51%	0.282%
40	12.039	1688	1692	1694	rBV2	21477	18808	1.55%	0.174%
41	12.427	1756	1758	1763	rVB3	16272	16279	1.34%	0.151%
42	12.574	1779	1783	1786	rBV	88224	87063	7.18%	0.806%
43	12.804	1816	1822	1828	rBV	107456	140480	11.59%	1.300%
44	12.857	1828	1831	1833	rVV4	17774	19240	1.59%	0.178%
45	12.939	1841	1845	1849	rBV	1007702	933414	76.98%	8.636%
46	13.157	1880	1882	1884	rVB2	21336	15856	1.31%	0.147%
47	13.327	1909	1911	1914	rBV2	21918	20615	1.70%	0.191%
48	13.357	1914	1916	1918	rBV2	21023	22369	1.84%	0.207%
49	13.827	1993	1996	1999	rBV4	74996	75428	6.22%	0.698%
50	14.010	2022	2027	2029	rBV	315454	343726	28.35%	3.180%
51	14.439	2097	2100	2106	rBV3	50502	94918	7.83%	0.878%
52	15.486	2275	2278	2282	rVB	167707	187212	15.44%	1.732%

Sum of corrected areas: 10808317

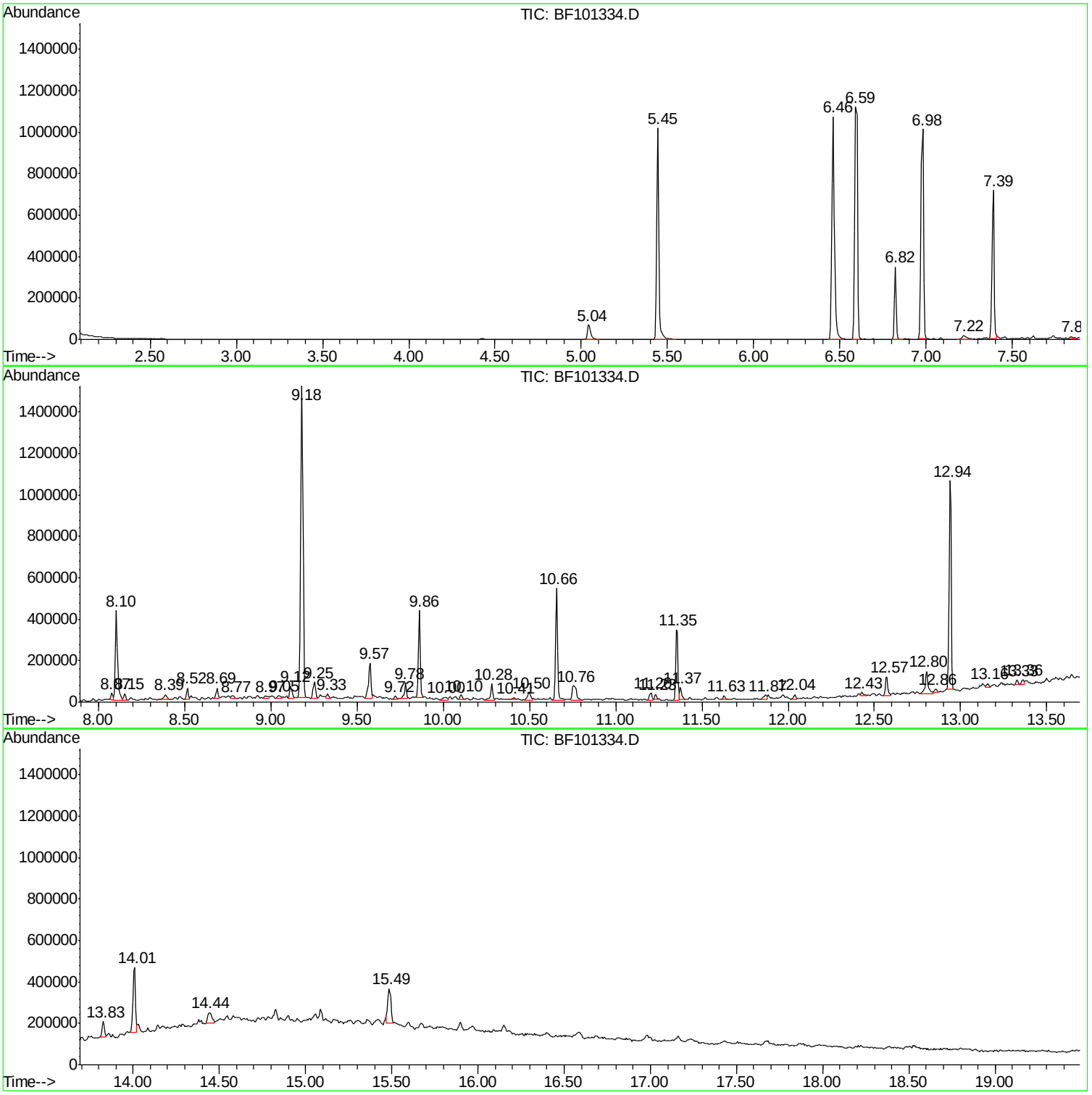
LSC Report - Integrated Chromatogram

Data Path : Z:\HPCHEM1\BNA F\DATA\BF121217\
Data File : BF101334.D
Acq On : 12 Dec 2017 17:16
Operator : SJ/JU
Sample : I6872-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 293-311

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF113017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\HPCHEM1\BNA F\DATA\BF121217\
Data File : BF101334.D
Acq On : 12 Dec 2017 17:16
Operator : SJ/JU
Sample : I6872-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 293-311

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF113017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

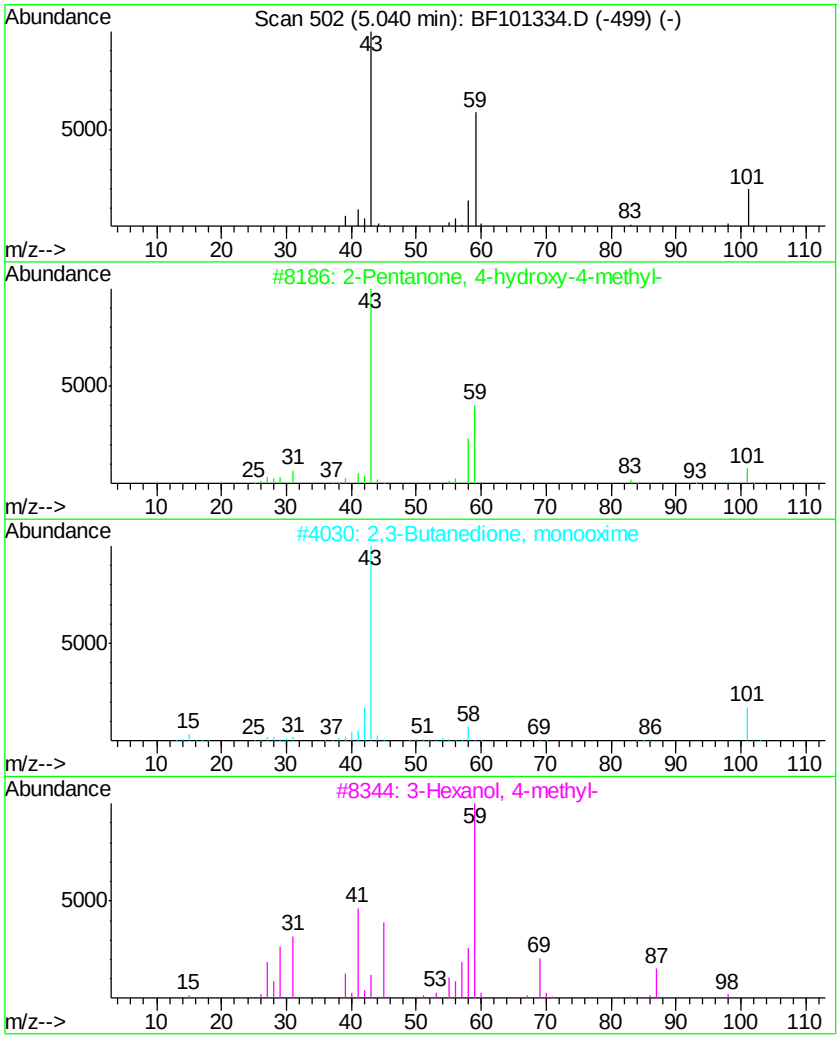
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

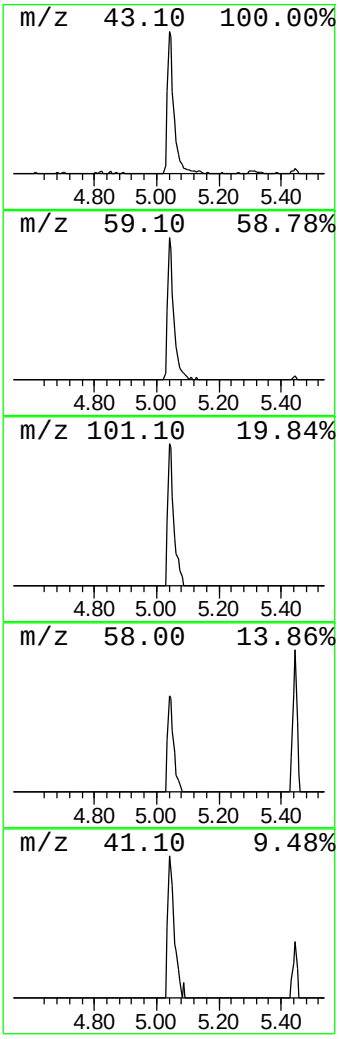
****

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.04	7.28 ng	100062	1,4-Dichlorobenzene-d4	6.82

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	35
3	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
4	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
5	4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	9





Data Path : Z:\HPCHEM1\BNA F\DATA\BF121217\
Data File : BF101334.D
Acq On : 12 Dec 2017 17:16
Operator : SJ/JU
Sample : I6872-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 293-311

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF113017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

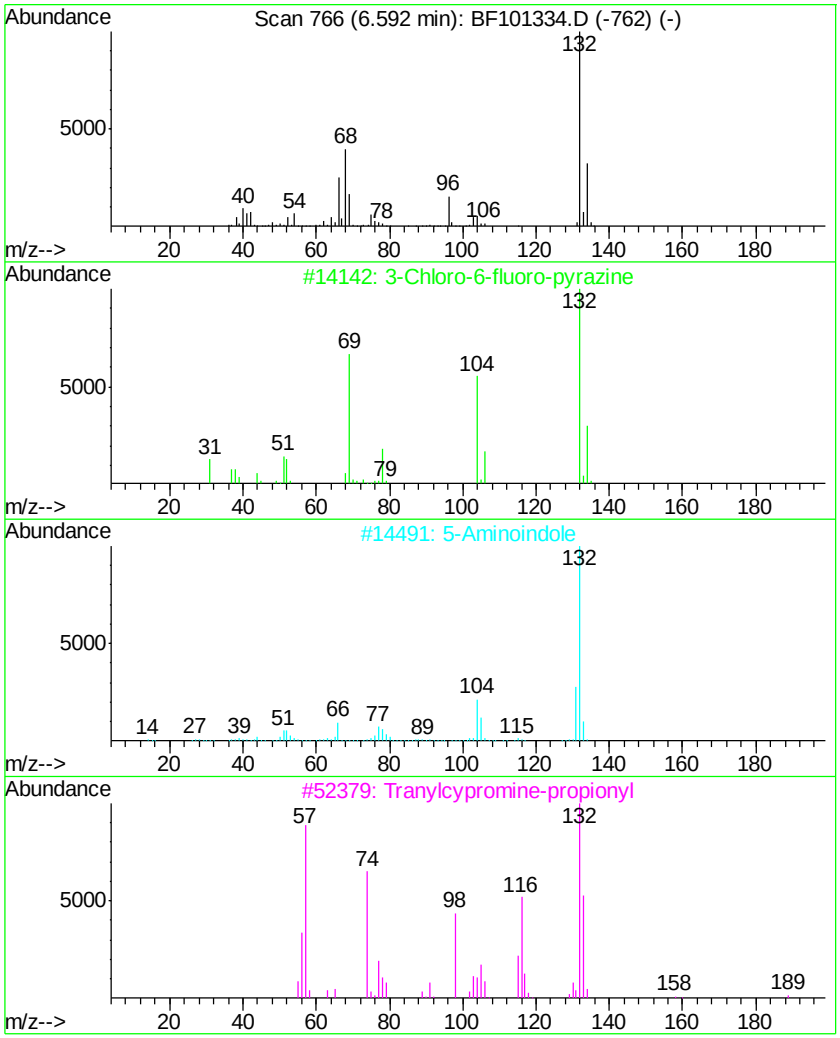
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

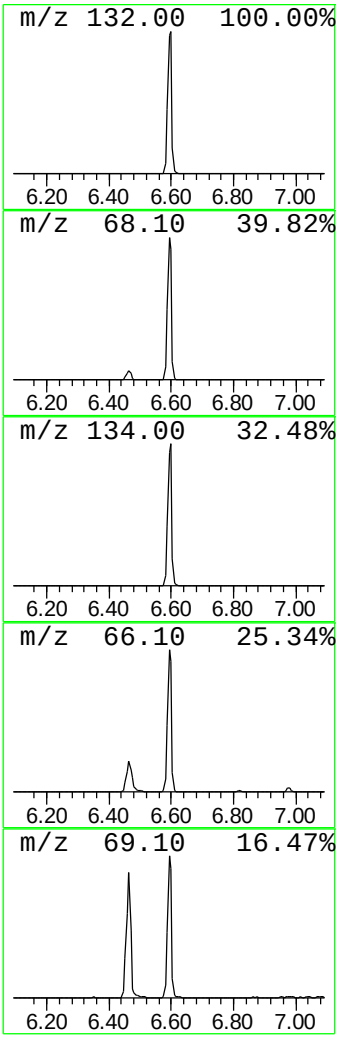
****

Peak Number 2 unknown6.59 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.59	77.76 ng	1068380	1,4-Dichlorobenzene-d4	6.82

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	Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4	4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9
5	2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9





Data Path : Z:\HPCHEM1\BNA F\DATA\BF121217\
Data File : BF101334.D
Acq On : 12 Dec 2017 17:16
Operator : SJ/JU
Sample : I6872-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 293-311

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF113017.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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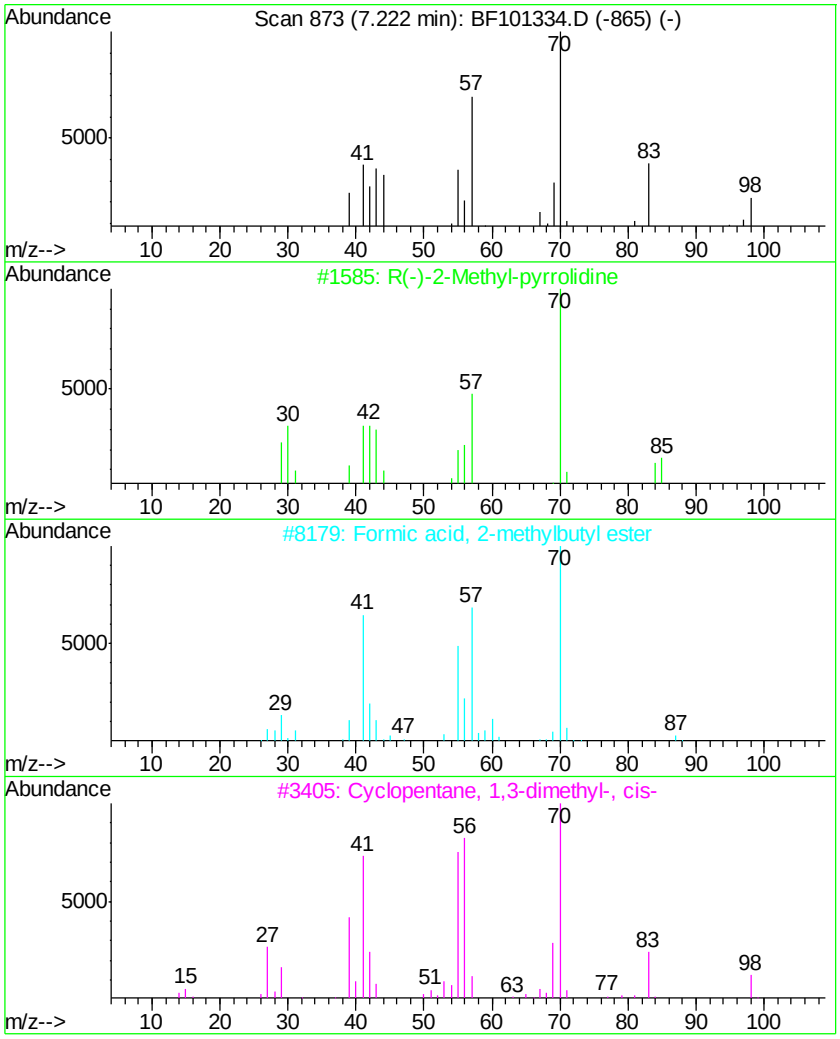
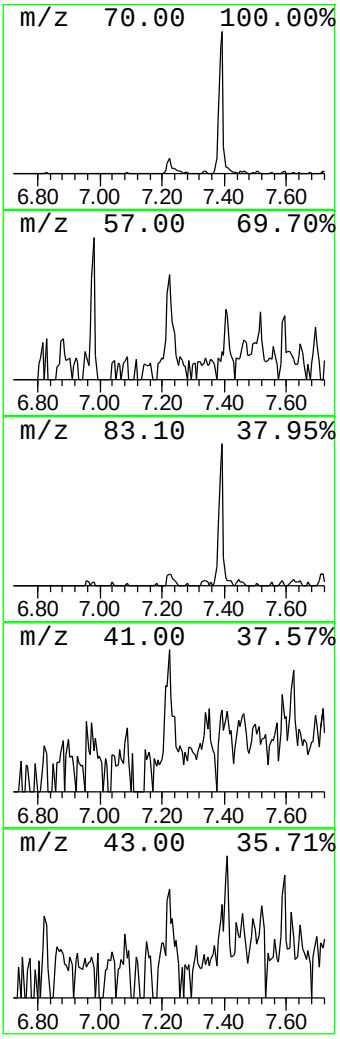
Peak Number 4 unknown7.22 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.22	2.48 ng	34032	1,4-Dichlorobenzene-d4	6.82

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	R(-)-2-Methyl-pyrrolidine	85	C5H11N	1000144-17-1	47
2	Formic acid, 2-methylbutyl ester	116	C6H12O2	1000367-91-1	43
3	Cyclopentane, 1,3-dimethyl-, cis-	98	C7H14	002532-58-3	38
4	2-Cyclohexen-1-ol	98	C6H10O	000822-67-3	38
5	2-Pyrrolidinemethanamine, N-meth...	114	C6H14N2	066411-54-9	28





Data Path : Z:\HPCHEM1\BNA F\DATA\BF121217\
Data File : BF101334.D
Acq On : 12 Dec 2017 17:16
Operator : SJ/JU
Sample : I6872-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 293-311

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF113017.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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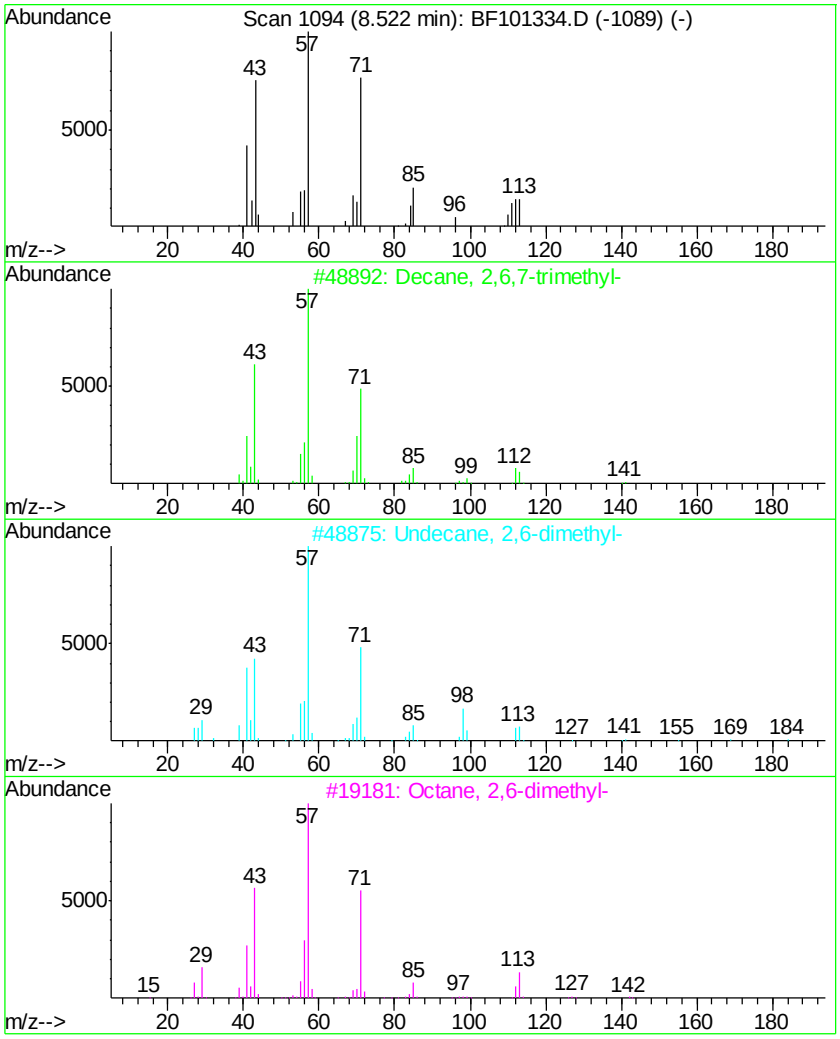
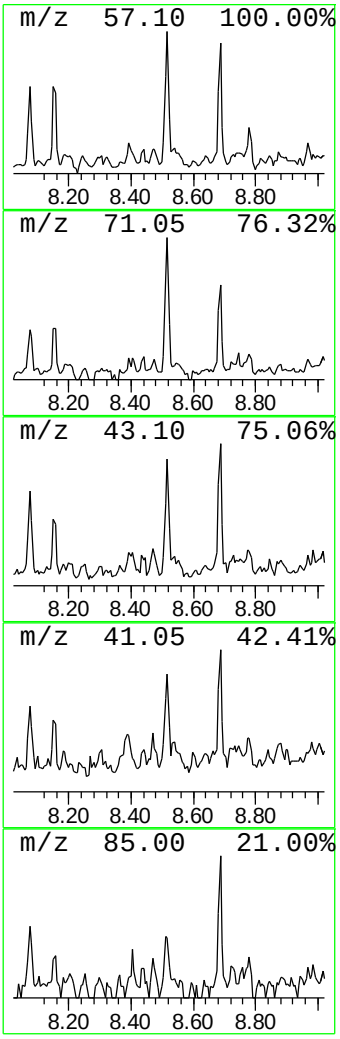
Peak Number 5 Decane, 2,6,7-trimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.52	2.40 ng	45075	Naphthalene-d8	8.10

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane, 2,6,7-trimethyl-	184	C13H28	062108-25-2	83
2	Undecane, 2,6-dimethyl-	184	C13H28	017301-23-4	83
3	Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	72
4	Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	72
5	Undecane, 4,5-dimethyl-	184	C13H28	017312-79-7	59





Data Path : Z:\HPCHEM1\BNA F\DATA\BF121217\
Data File : BF101334.D
Acq On : 12 Dec 2017 17:16
Operator : SJ/JU
Sample : I6872-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 293-311

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF113017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

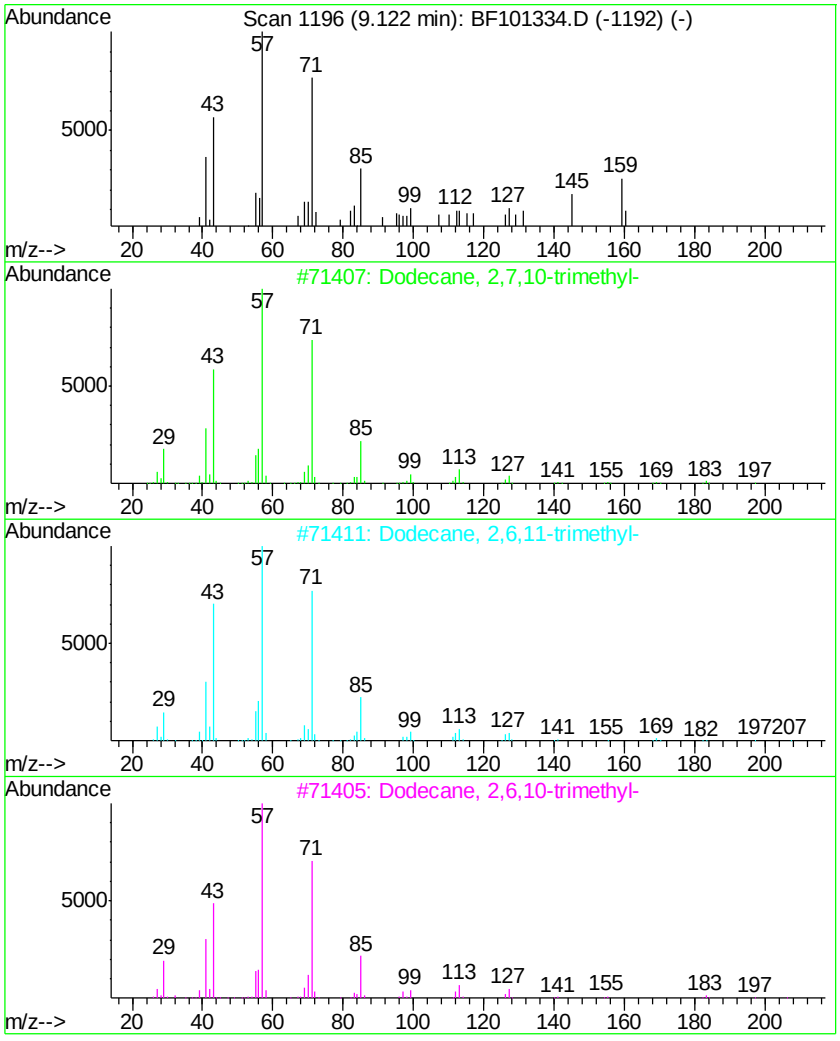
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

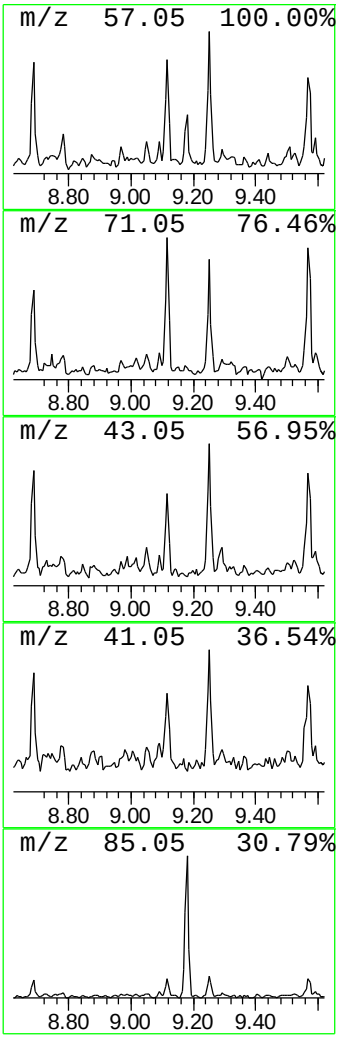
****

Peak Number 6 Dodecane, 2,7,10-trimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.12	2.67 ng	44519	Acenaphthene-d10	9.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	64
2	Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	64
3	Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	64
4	Nonane, 3-methyl-5-propyl-	184	C13H28	031081-18-2	53
5	Nonane, 2-methyl-5-propyl-	184	C13H28	031081-17-1	53





Data Path : Z:\HPCHEM1\BNA F\DATA\BF121217\
Data File : BF101334.D
Acq On : 12 Dec 2017 17:16
Operator : SJ/JU
Sample : I6872-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 293-311

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF113017.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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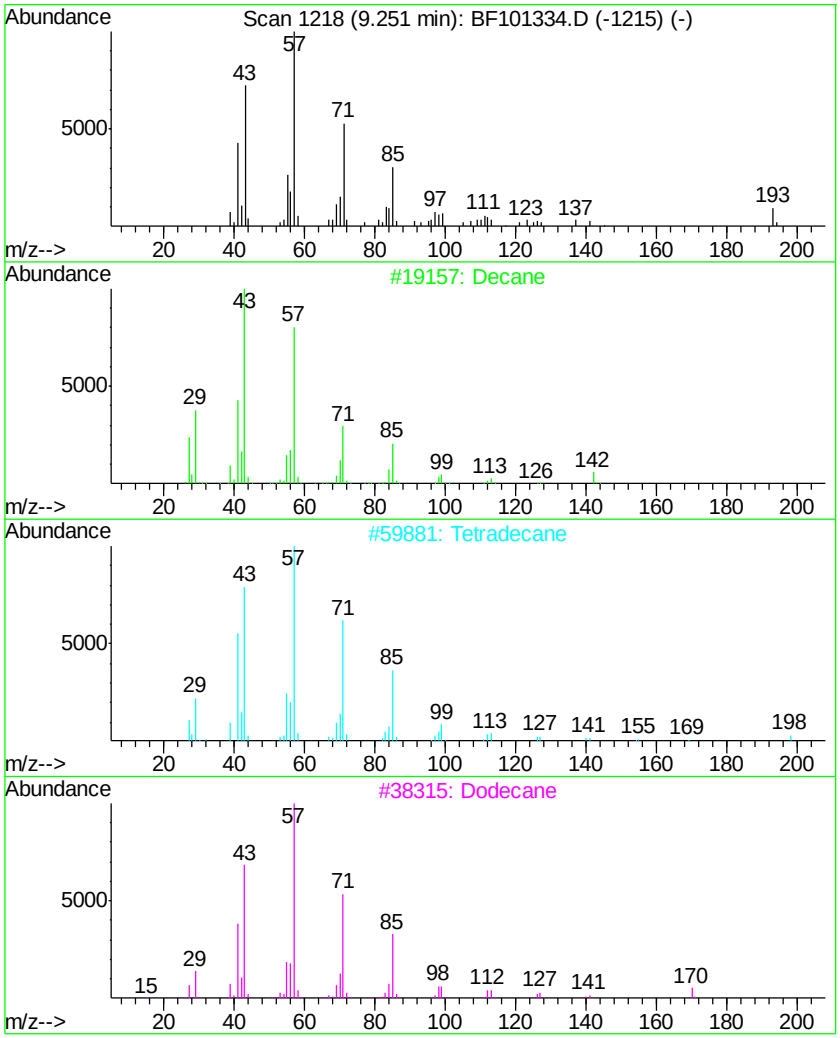
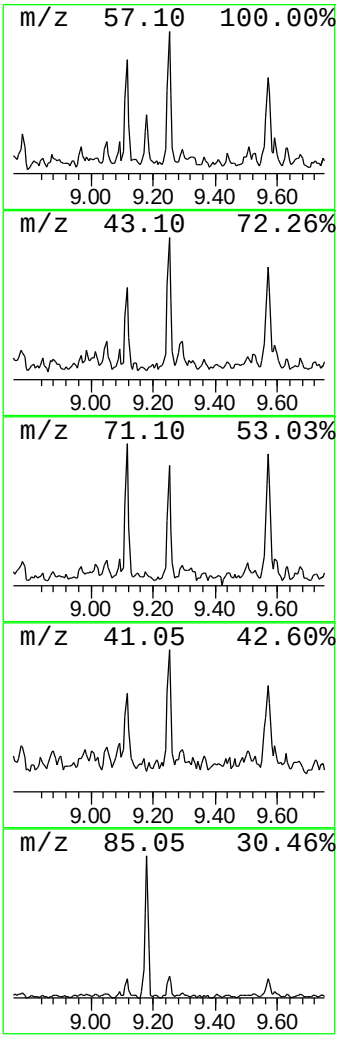
Peak Number 7 Decane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.25	3.81 ng	63576	Acenaphthene-d10	9.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane	142	C10H22	000124-18-5	72
2	Tetradecane	198	C14H30	000629-59-4	72
3	Dodecane	170	C12H26	000112-40-3	72
4	Hexadecane, 7,9-dimethyl-	254	C18H38	021164-95-4	72
5	Nonadecane	268	C19H40	000629-92-5	72





Data Path : Z:\HPCHEM1\BNA F\DATA\BF121217\
Data File : BF101334.D
Acq On : 12 Dec 2017 17:16
Operator : SJ/JU
Sample : I6872-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 293-311

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF113017.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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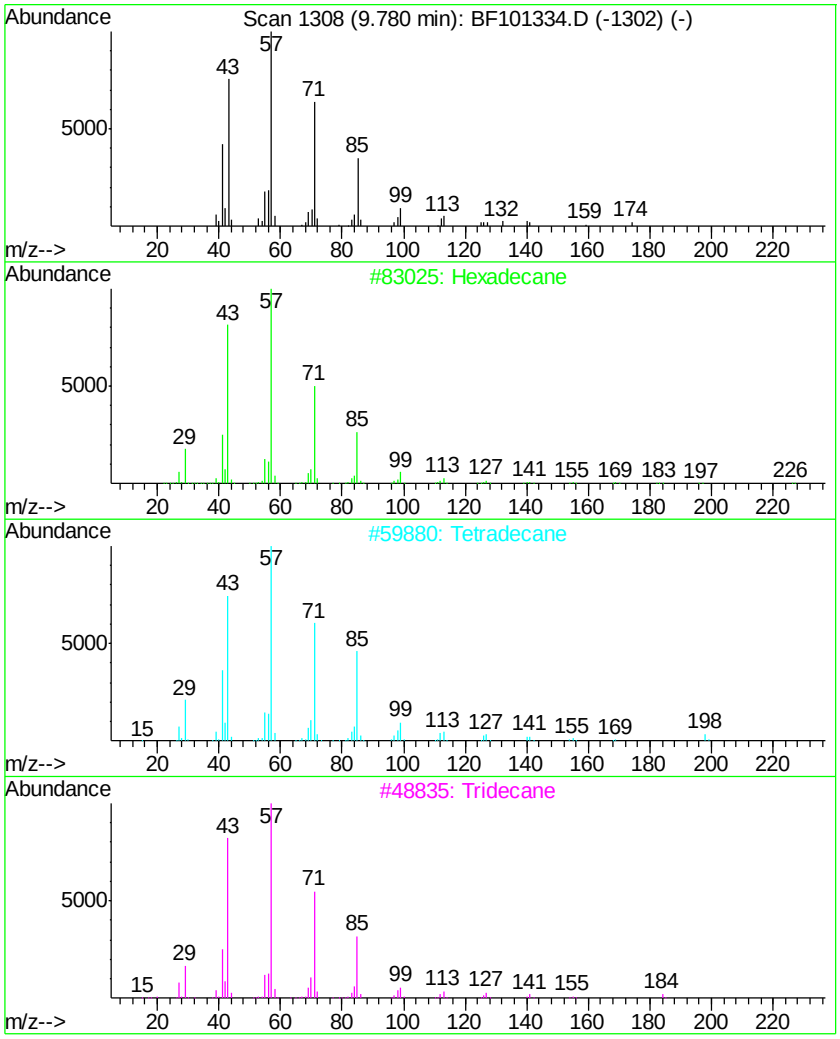
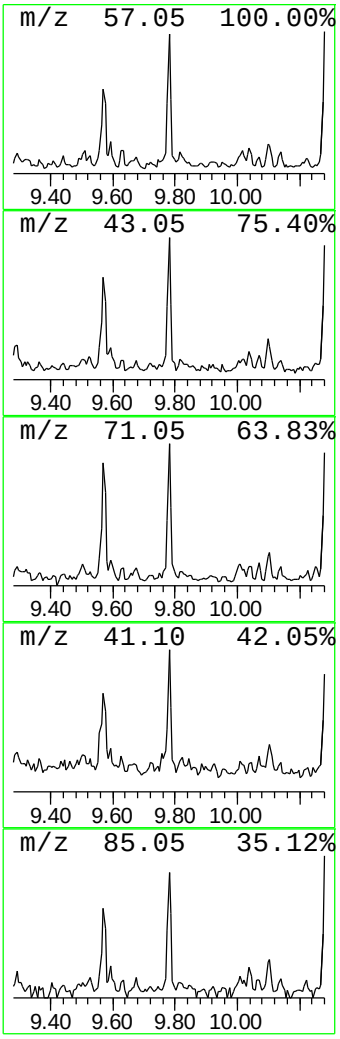
Peak Number 8 Hexadecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.78	3.86 ng	64343	Acenaphthene-d10	9.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	90
2	Tetradecane	198	C14H30	000629-59-4	83
3	Tridecane	184	C13H28	000629-50-5	83
4	Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	78
5	Sulfurous acid, butyl heptadecyl...	376	C21H44O3S	1000309-18-4	78





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF121217\
Data File : BF101334.D
Acq On : 12 Dec 2017 17:16
Operator : SJ/JU
Sample : I6872-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 293-311

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF113017.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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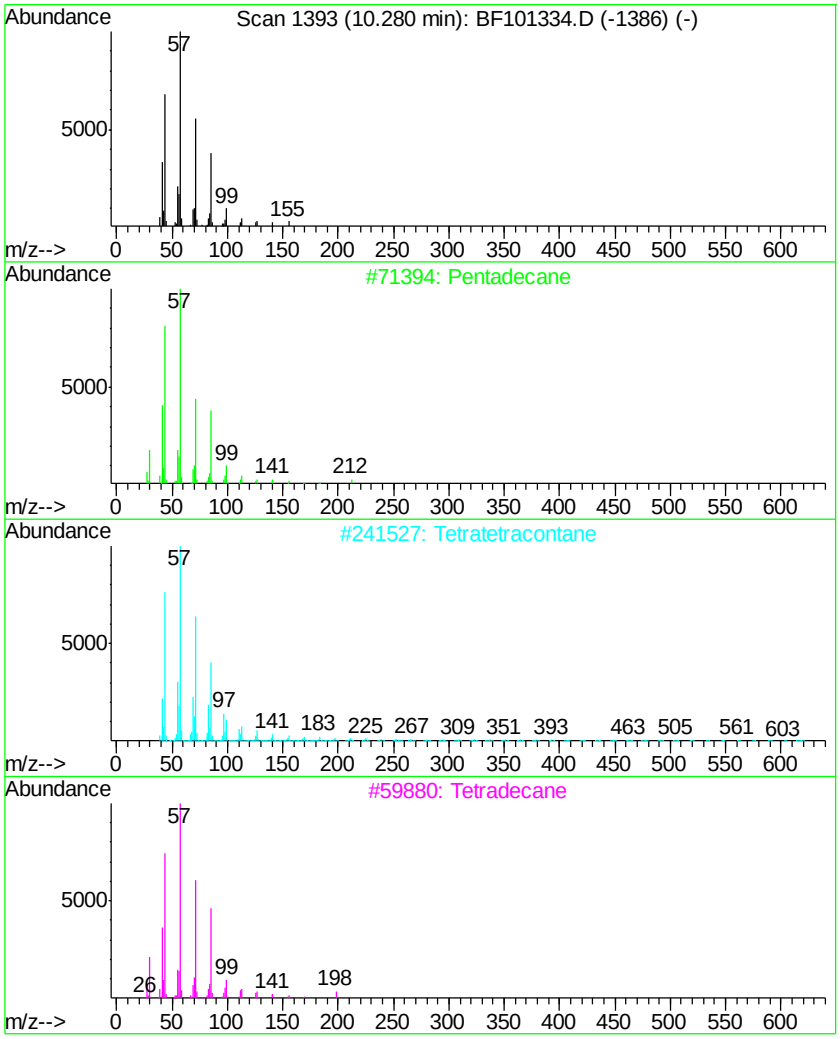
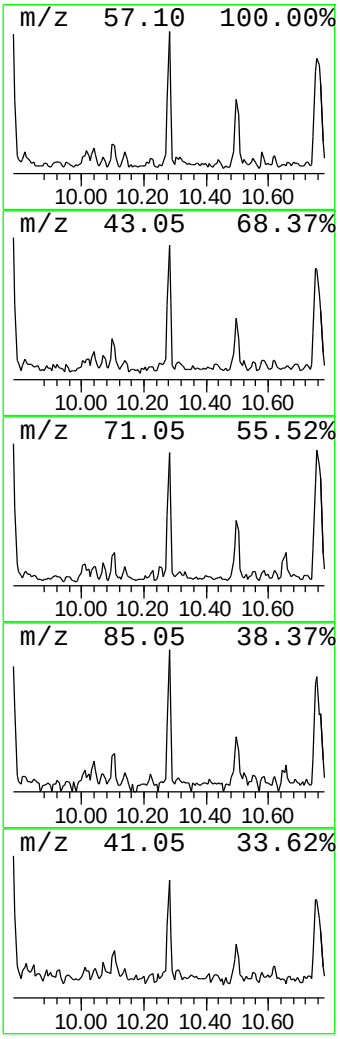
Peak Number 9 Pentadecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.28	4.32 ng	72068	Acenaphthene-d10	9.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecane	212	C15H32	000629-62-9	91
2	Tetratetracontane	619	C44H90	007098-22-8	90
3	Tetradecane	198	C14H30	000629-59-4	90
4	Tridecane	184	C13H28	000629-50-5	90
5	Tetracosane	338	C24H50	000646-31-1	90





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Data File : BF101334.D
Acq On : 12 Dec 2017 17:16
Operator : SJ/JU
Sample : I6872-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 293-311

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF113017.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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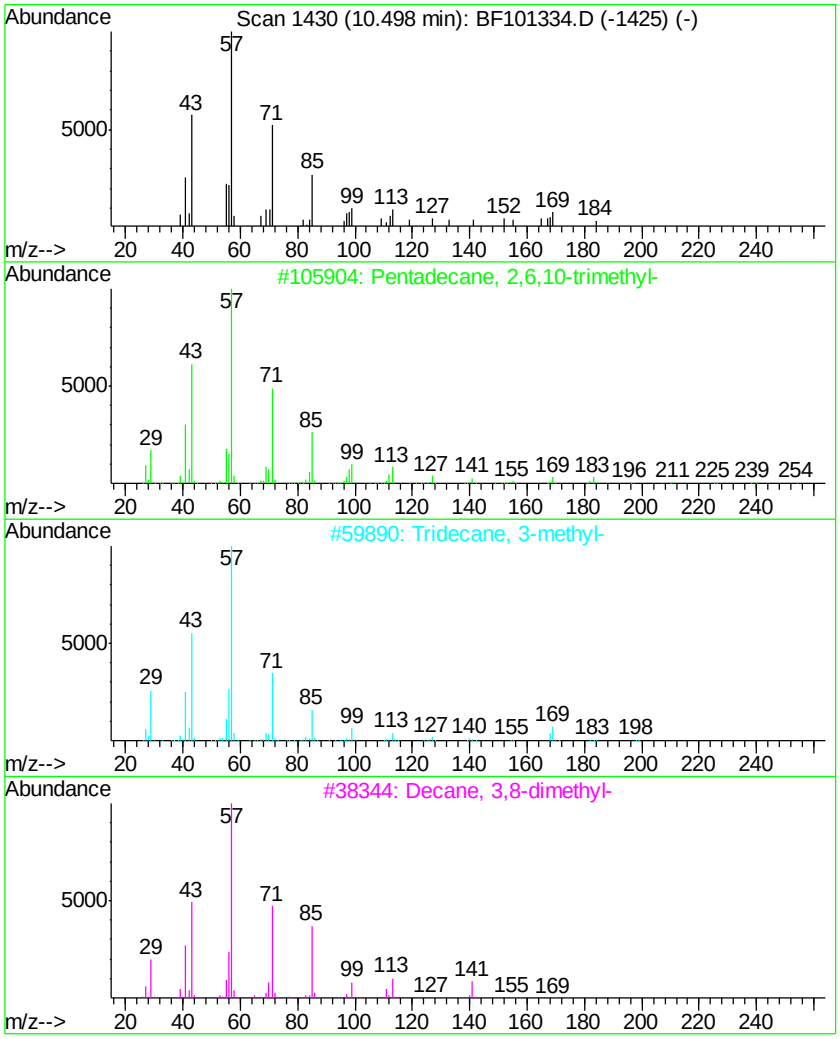
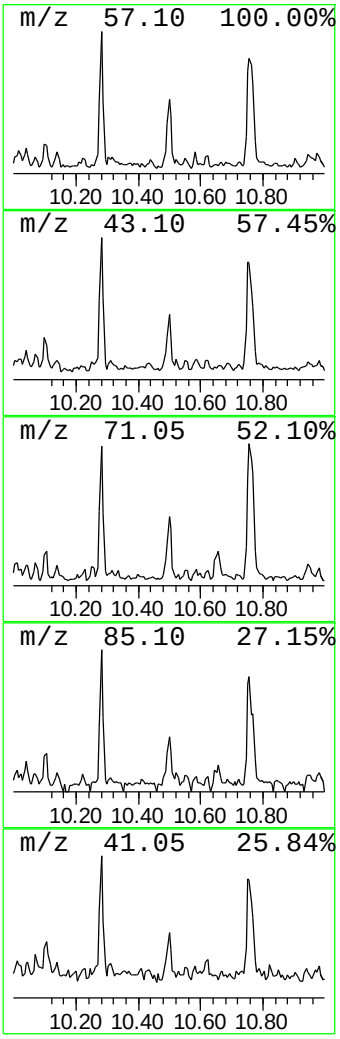
Peak Number 10 Pentadecane, 2,6,10-trimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.50	2.96 ng	49453	Acenaphthene-d10	9.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	80
2	Tridecane, 3-methyl-	198	C14H30	006418-41-3	78
3	Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	72
4	Heneicosane	296	C21H44	000629-94-7	72
5	Eicosane, 10-methyl-	296	C21H44	054833-23-7	64





Library Search Compound Report

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ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 293-311

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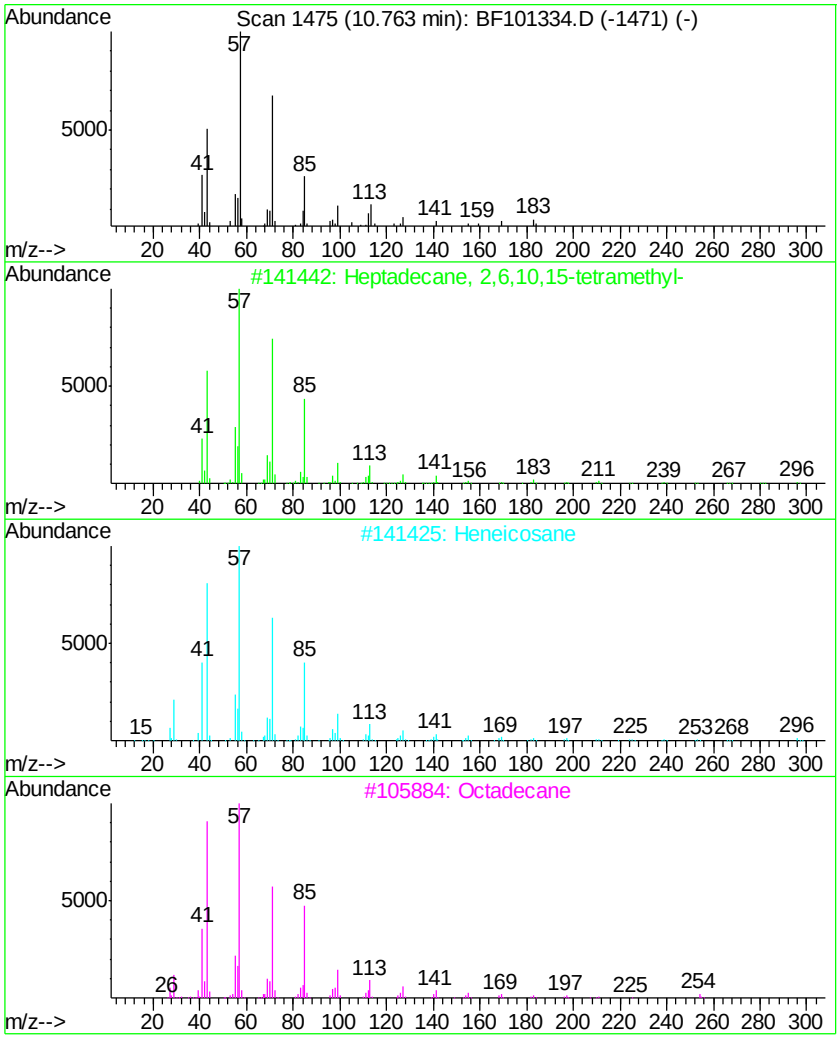
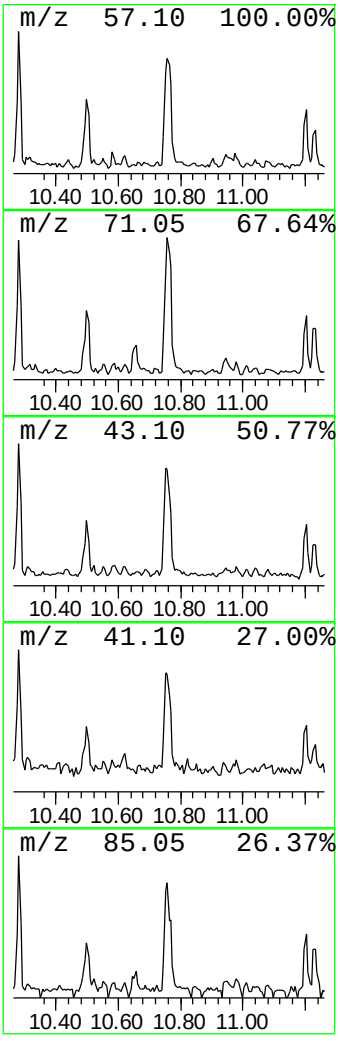
Peak Number 11 Heptadecane, 2,6,10,15-tetr... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.76	6.51 ng	104974	Phenanthrene-d10	11.35

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	91
2	Heneicosane	296	C21H44	000629-94-7	91
3	Octadecane	254	C18H38	000593-45-3	91
4	Tetracosane	338	C24H50	000646-31-1	91
5	Hexadecane	226	C16H34	000544-76-3	90





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Data File : BF101334.D
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Operator : SJ/JU
Sample : I6872-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 293-311

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF113017.M
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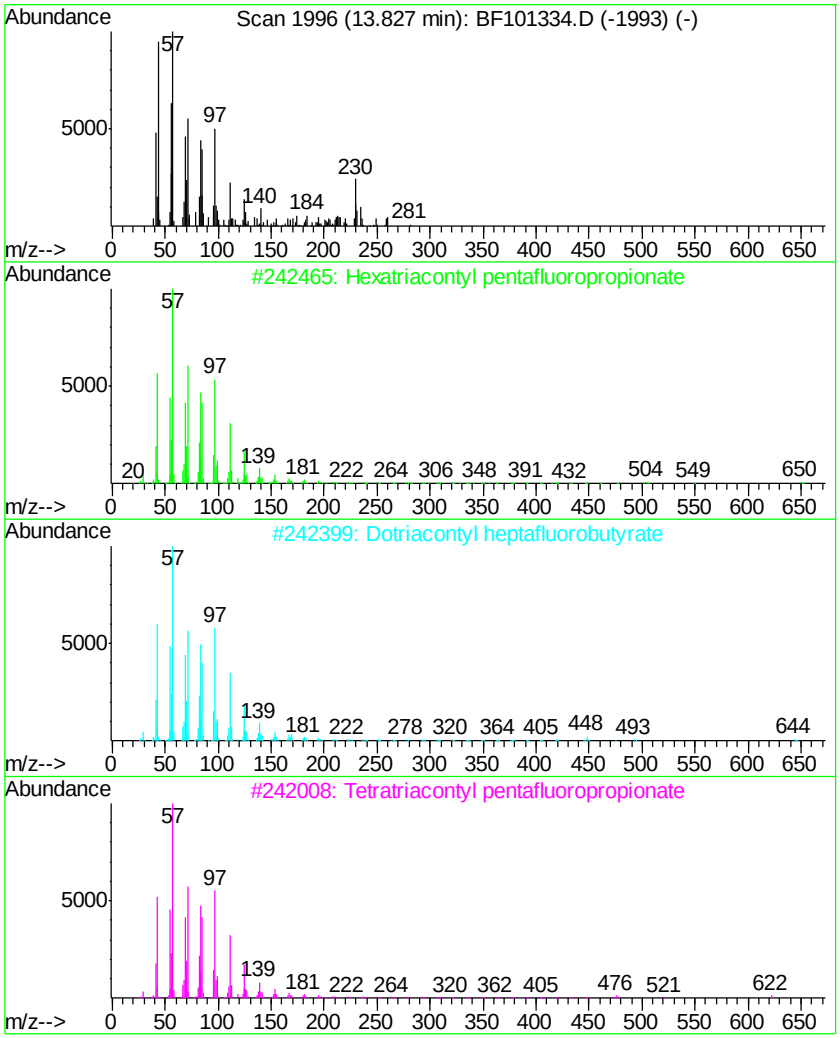
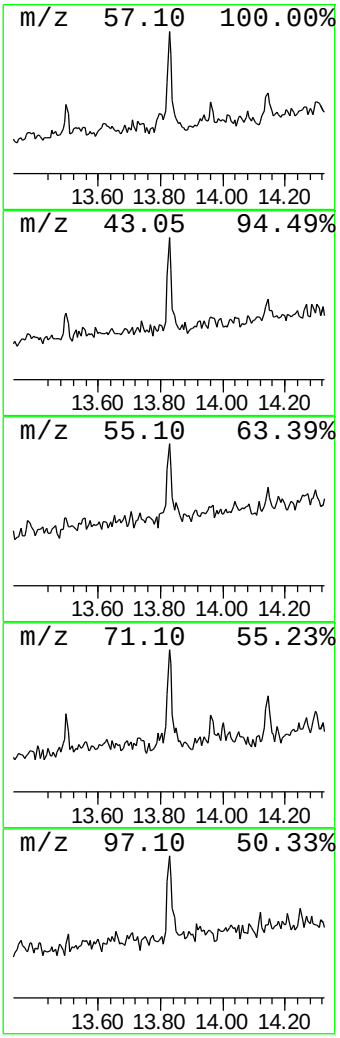
Peak Number 12 Hexatriacontyl pentafluorop... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.83	4.39 ng	75428	Chrysene-d12	14.01

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexatriacontyl pentafluoropropio...	669	C39H73F5O2	1000351-89-0	74
2	Dotriacontyl heptafluorobutyrate	662	C36H65F7O2	1000351-84-2	72
3	Tetratriacontyl pentafluoropropi...	641	C37H69F5O2	1000351-81-5	72
4	1-Decanol, 2-hexyl-	242	C16H34O	002425-77-6	62
5	1-Decanol, 2-octyl-	270	C18H38O	045235-48-1	62





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Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 293-311

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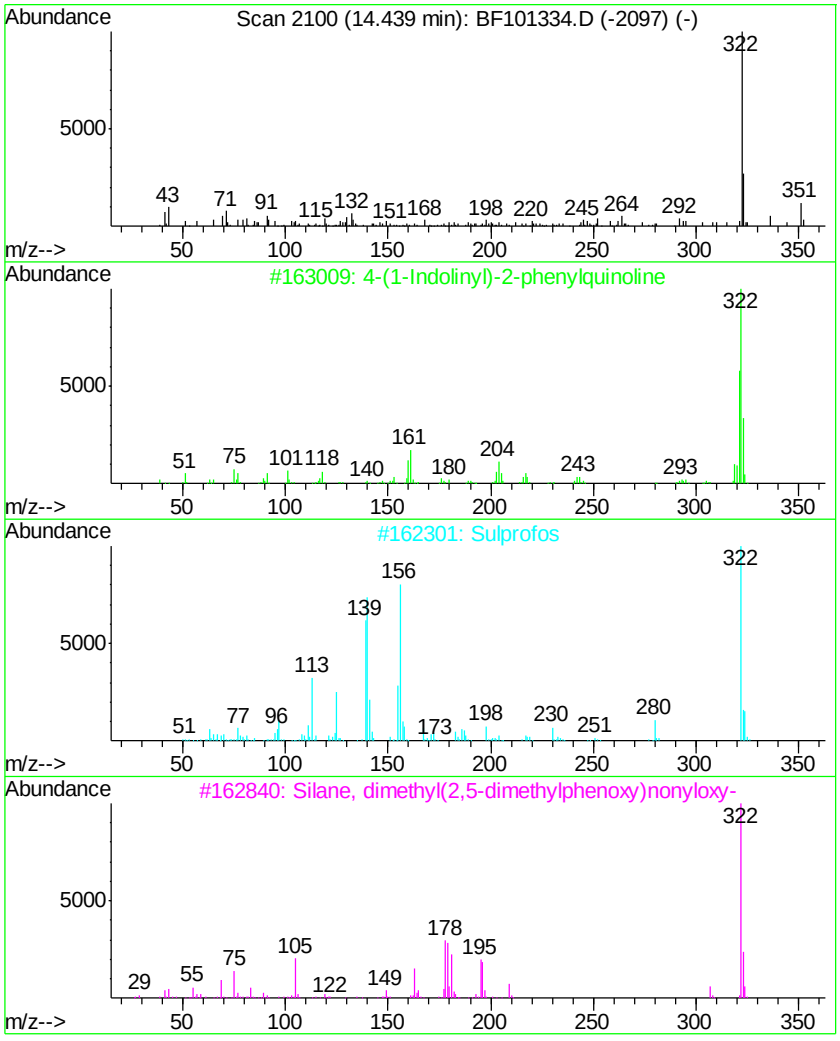
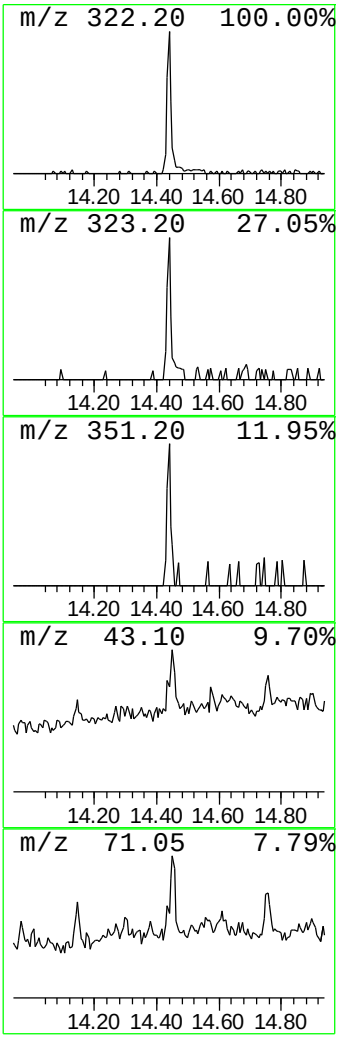
Peak Number 13 4-(1-Indolinyl)-2-phenylqui... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.44	5.52 ng	94918	Chrysene-d12	14.01

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-(1-Indolinyl)-2-phenylquinoline	322	C23H18N2	1000326-78-8	59
2	Sulprofos	322	C12H19O2PS3	035400-43-2	46
3	Silane, dimethyl(2,5-dimethylphe...	322	C19H34O2Si	1000347-29-5	42
4	Pyrazole-4-carboxylic acid, 3-(1...	322	C20H22N2O2	1000272-86-5	42
5	m-Bis(p-methoxyphenoxy)benzene	322	C20H18O4	013118-91-7	42





Tentatively Identified Compound (LSC) summary

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Operator : SJ/JU
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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...	5.04	7.3	ng	100062	1	6.82	274793	20.0
unknown6.59	6.59	77.8	ng	1068380	1	6.82	274793	20.0
unknown7.22	7.22	2.5	ng	34032	1	6.82	274793	20.0
Decane, 2,6,7-tri...	8.52	2.4	ng	45075	2	8.10	376394	20.0
Dodecane, 2,7,10-...	9.12	2.7	ng	44519	3	9.86	333713	20.0
Decane	9.25	3.8	ng	63576	3	9.86	333713	20.0
Hexadecane	9.78	3.9	ng	64343	3	9.86	333713	20.0
Pentadecane	10.28	4.3	ng	72068	3	9.86	333713	20.0
Pentadecane, 2,6,...	10.50	3.0	ng	49453	3	9.86	333713	20.0
Heptadecane, 2,6,...	10.76	6.5	ng	104974	4	11.35	322281	20.0
Hexatriacontyl pe...	13.83	4.4	ng	75428	5	14.01	343726	20.0
4-(1-Indolinyl)-2...	14.44	5.5	ng	94918	5	14.01	343726	20.0