LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF121117\
Data File : BF101293.D
Acq On : 11 Dec 2017 16:56
Operator : SJ/JU
Sample : I6744-06
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : PROS-2-E(0-1)

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	5.057	501	505	516	rBV	48936	73590	5.83%	0.590%
2	5.451	568	572	589	rBV	1105687	1147117	90.91%	9.192%
3	6.469	740	745	759	rBV	1098122	1200191	95.11%	9.617%
4	6.604	763	768	771	rBV	1402505	1212630	96.10%	9.716%
5	6.828	803	806	813	rBV	378330	305655	24.22%	2.449%
6	6.986	829	833	836	rBV	1133366	922303	73.09%	7.390%
7	7.398	898	903	906	rBV	834825	719860	57.05%	5.768%
8	8.110	1021	1024	1032	rBV	459439	395341	31.33%	3.168%
9	9.186	1203	1207	1210	rBV	1523383	1261847	100.00%	10.111%
10	9.257	1216	1219	1222	rVB	20919	15935	1.26%	0.128%
11	9.580	1268	1274	1281	rVB	105321	93613	7.42%	0.750%
12	9.786	1306	1309	1314	rBV	56969	47912	3.80%	0.384%
13	9.869	1320	1323	1327	rBV2	510707	408265	32.35%	3.271%
14	10.263	1387	1390	1392	rBV	14383	14969	1.19%	0.120%
15	10.286	1392	1394	1397	rVB	62570	44319	3.51%	0.355%
16	10.522	1424	1434	1438	rBV2	32732	49584	3.93%	0.397%
17	10.663	1454	1458	1462	rBV	790729	708215	56.13%	5.675%
18	10.763	1472	1475	1479	rBV	52767	54740	4.34%	0.439%
19	10.921	1498	1502	1505	rBV	75152	60667	4.81%	0.486%
20	11.210	1548	1551	1553	rBV	33594	25688	2.04%	0.206%
21	11.280	1560	1563	1565	rBV	78530	64000	5.07%	0.513%
22	11.304	1565	1567	1570	rVB	32195	23737	1.88%	0.190%
23	11.363	1573	1577	1579	rBV	480778	411598	32.62%	3.298%
24	11.386	1579	1581	1585	rVB	107795	87133	6.91%	0.698%
25	11.451	1586	1592	1595	rVB2	45237	59890	4.75%	0.480%
26	11.633	1621	1623	1628	rVB2	23252	21369	1.69%	0.171%
27	11.698	1630	1634	1638	rVB2	85360	111420	8.83%	0.893%
28	11.768	1643	1646	1649	rBV	50241	46015	3.65%	0.369%
29	11.845	1657	1659	1661	rBV	21216	17860	1.42%	0.143%
30	11.874	1661	1664	1666	rBV2	27619	26710	2.12%	0.214%
31	11.968	1678	1680	1683	rBV2	44141	44794	3.55%	0.359%
32	12.004	1683	1686	1688	rVV3	24980	22084	1.75%	0.177%
33	12.033	1688	1691	1696	rVB6	14842	20640	1.64%	0.165%
34	12.139	1706	1709	1711	rBV	31824	24381	1.93%	0.195%

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF121117\
Data File : BF101293.D
Acq On : 11 Dec 2017 16:56
Operator : SJ/JU
Sample : I6744-06
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : PROS-2-E(0-1)

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	12.580	1780	1784	1787	rBV	196459	188321	14.92%	1.509%
36	12.810	1816	1823	1829	rBV	177764	204864	16.24%	1.642%
37	12.945	1839	1846	1849	rBV	1056479	889983	70.53%	7.131%
38	13.033	1856	1861	1864	rVB2	14173	24534	1.94%	0.197%
39	13.121	1871	1876	1877	rBV4	19962	28197	2.23%	0.226%
40	13.139	1877	1879	1882	rVB	22263	21600	1.71%	0.173%
41	13.239	1893	1896	1903	rBV4	15433	27822	2.20%	0.223%
42	13.651	1964	1966	1970	rBV	17331	17680	1.40%	0.142%
43	13.757	1982	1984	1986	rBV2	17225	16200	1.28%	0.130%
44	13.827	1994	1996	2000	rVB2	49581	49390	3.91%	0.396%
45	13.862	2000	2002	2004	rVB	21022	17429	1.38%	0.140%
46	13.968	2017	2020	2023	rBV	124471	110697	8.77%	0.887%
47	14.010	2023	2027	2030	rVV2	378414	415050	32.89%	3.326%
48	14.033	2030	2031	2035	rVB	86741	67089	5.32%	0.538%
49	14.115	2043	2045	2049	rBV2	18439	16583	1.31%	0.133%
50	15.057	2202	2205	2208	rBV	74540	97068	7.69%	0.778%
51	15.362	2254	2257	2263	rVB	48163	61111	4.84%	0.490%
52	15.427	2265	2268	2272	rVB	77257	96823	7.67%	0.776%
53	15.492	2273	2279	2282	rBV	217325	300086	23.78%	2.405%
54	16.986	2529	2533	2539	rVB2	43408	85560	6.78%	0.686%

Sum of corrected areas: 12480159

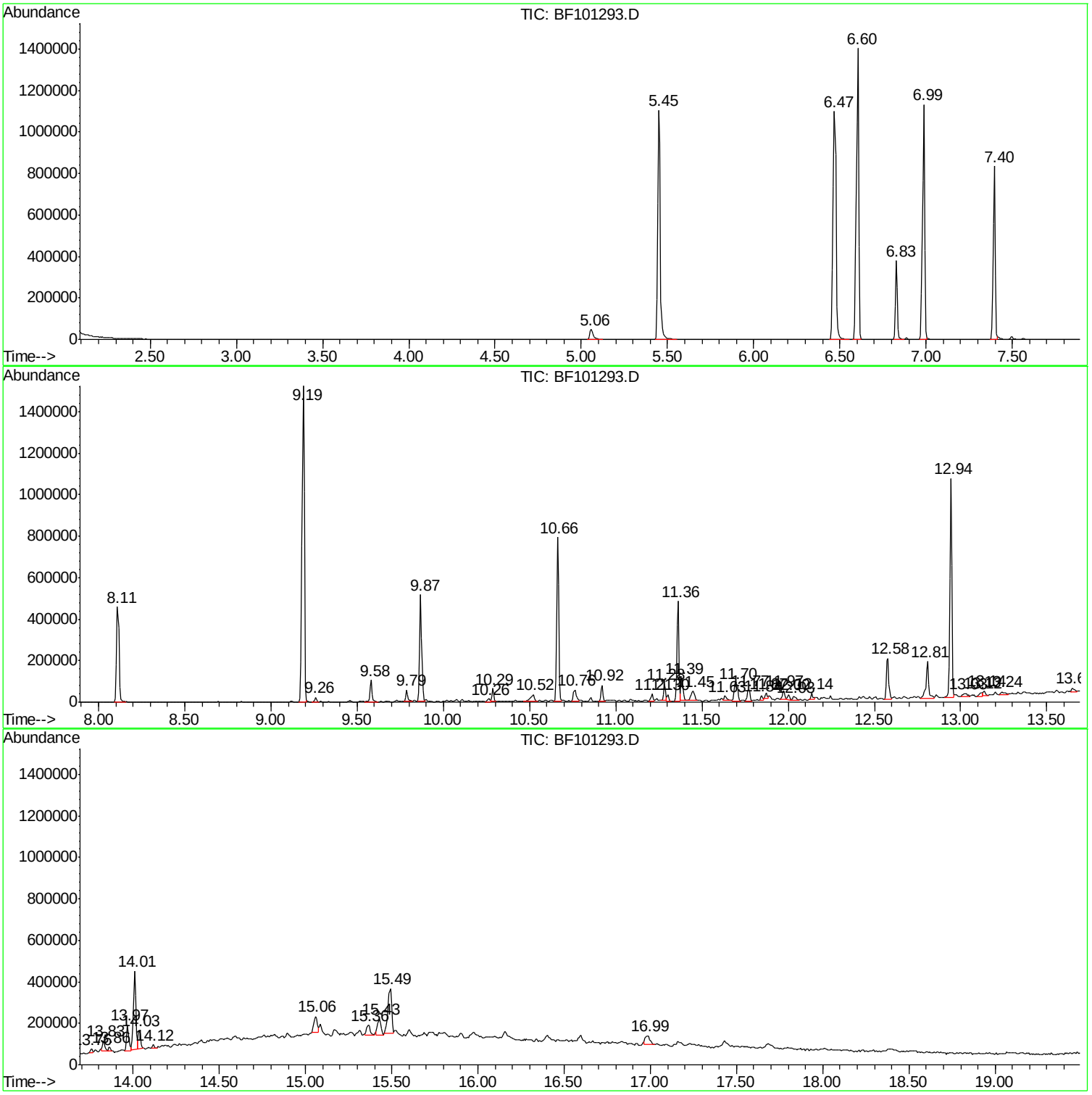
LSC Report - Integrated Chromatogram

Data Path : Z:\HPCHEM1\BNA F\DATA\BF121117\
Data File : BF101293.D
Acq On : 11 Dec 2017 16:56
Operator : SJ/JU
Sample : I6744-06
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PROS-2-E(0-1)

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



Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF121117\
Data File : BF101293.D
Acq On : 11 Dec 2017 16:56
Operator : SJ/JU
Sample : I6744-06
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : PROS-2-E(0-1)

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

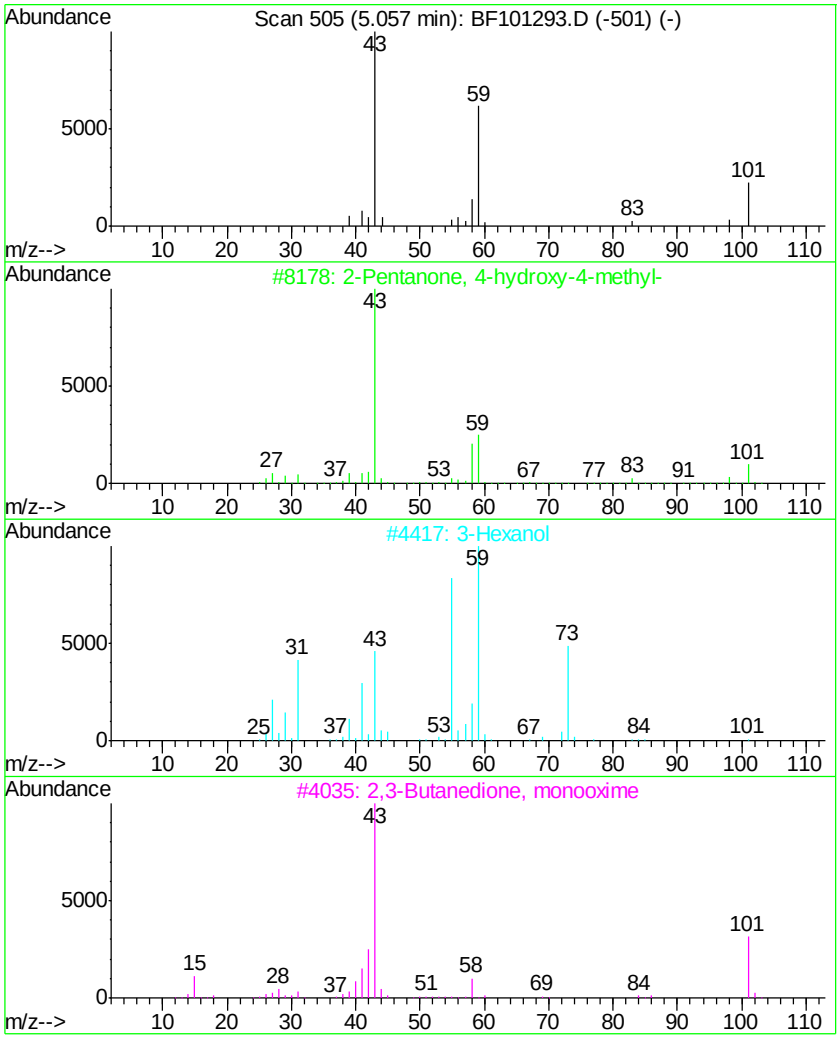
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

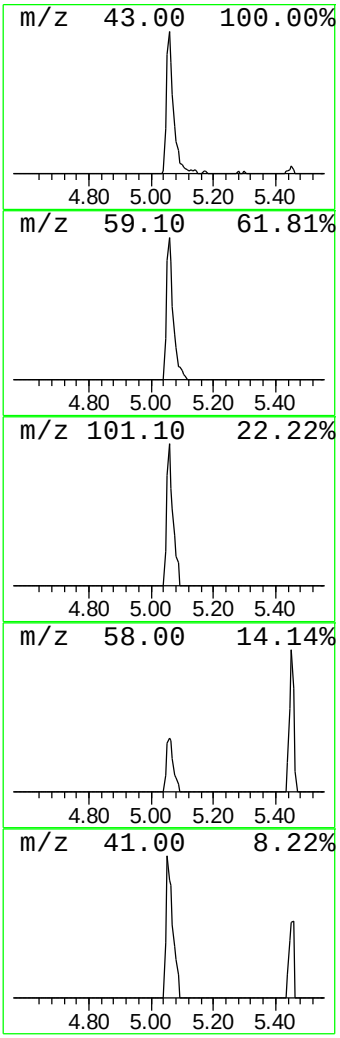
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.06	4.82 ng	73590	1,4-Dichlorobenzene-d4	6.83

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2	3-Hexanol	102	C6H14O	000623-37-0	33
3	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	32
4	Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	23
5	1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	23





Data Path : Z:\HPCHEM1\BNA F\DATA\BF121117\
Data File : BF101293.D
Acq On : 11 Dec 2017 16:56
Operator : SJ/JU
Sample : I6744-06
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : PROS-2-E(0-1)

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

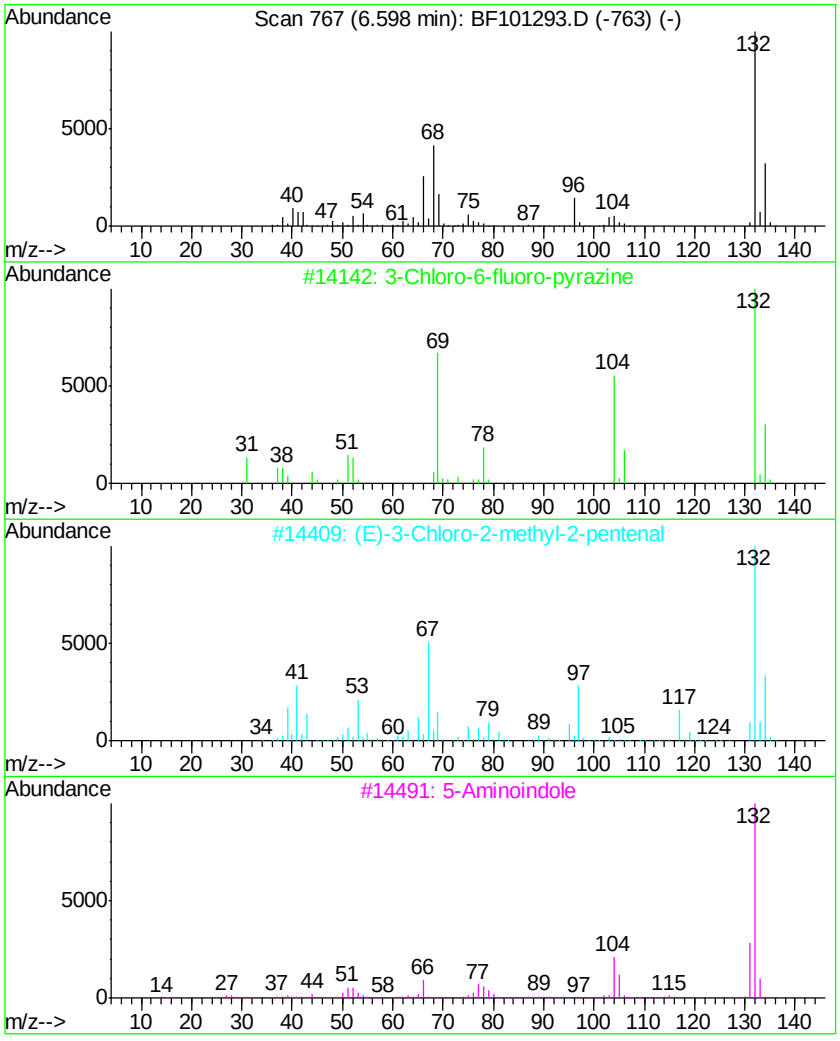
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

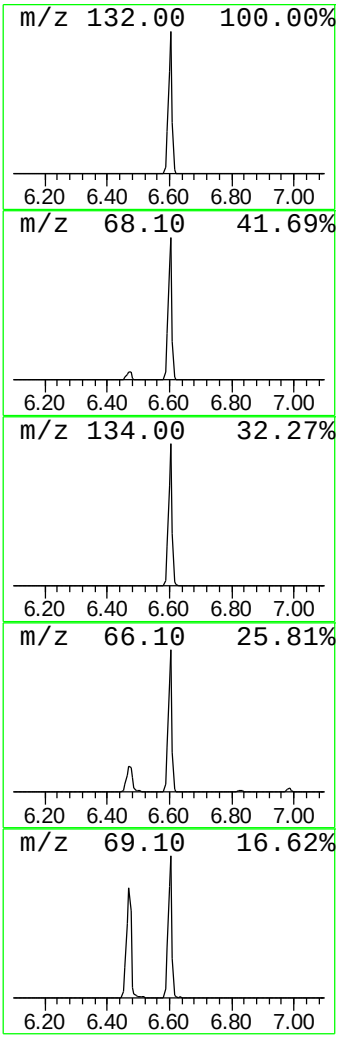
****

Peak Number 2 unknown6.60 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.60	79.35 ng	1212630	1,4-Dichlorobenzene-d4	6.83

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3	5-Aminoindole	132	C8H8N2	005192-03-0	12
4	Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5	4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9





Data Path : Z:\HPCHEM1\BNA F\DATA\BF121117\
Data File : BF101293.D
Acq On : 11 Dec 2017 16:56
Operator : SJ/JU
Sample : I6744-06
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : PROS-2-E(0-1)

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

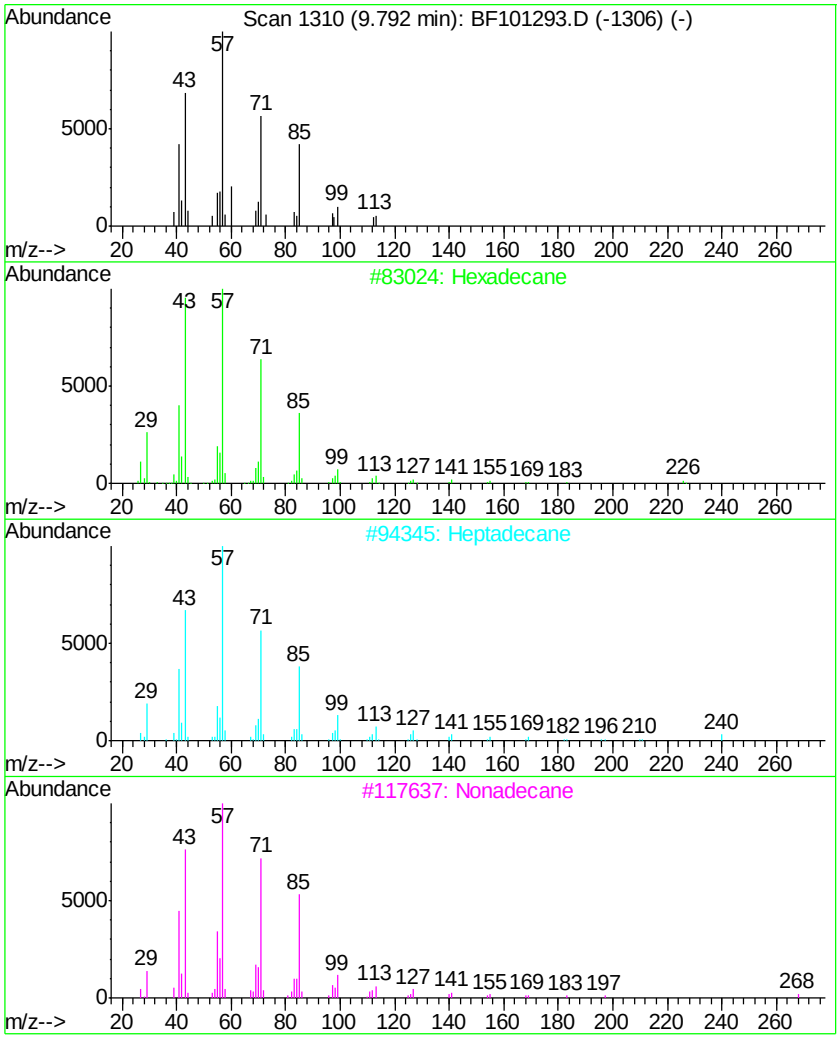
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

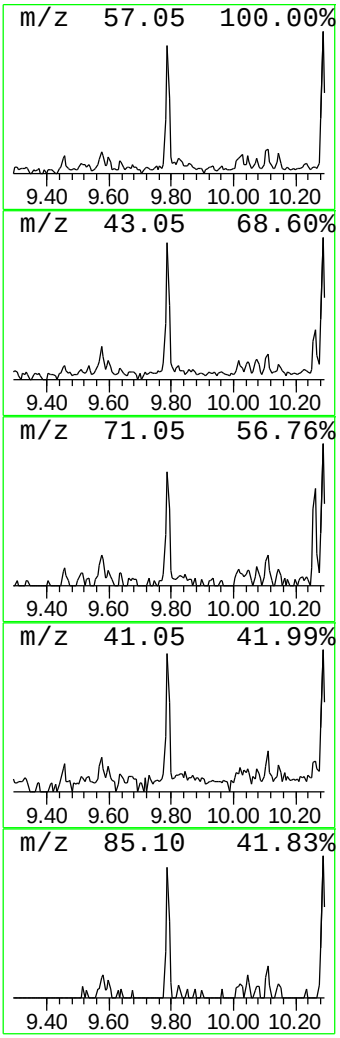
****

Peak Number 4 Hexadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.79	2.35 ng	47912	Acenaphthene-d10	9.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	86
2	Heptadecane	240	C17H36	000629-78-7	78
3	Nonadecane	268	C19H40	000629-92-5	78
4	Tridecane	184	C13H28	000629-50-5	78
5	Pentadecane	212	C15H32	000629-62-9	64





Data Path : Z:\HPCHEM1\BNA F\DATA\BF121117\
Data File : BF101293.D
Acq On : 11 Dec 2017 16:56
Operator : SJ/JU
Sample : I6744-06
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : PROS-2-E(0-1)

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

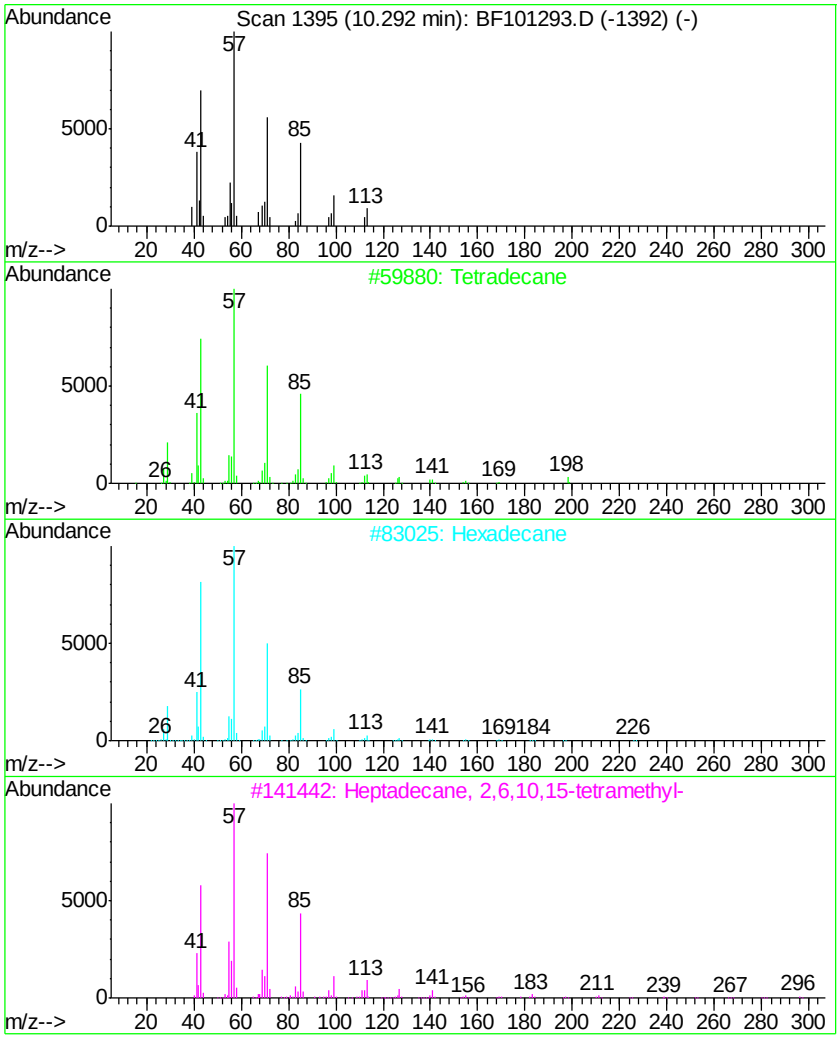
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

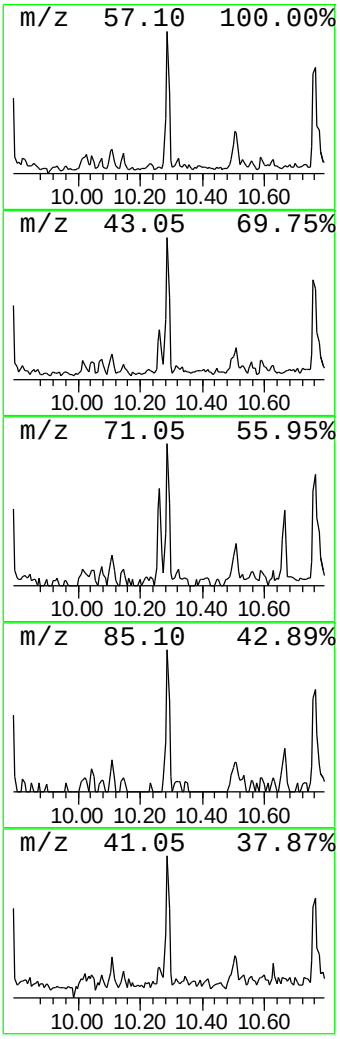
****

Peak Number 5 Tetradecane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.29	2.17 ng	44319	Acenaphthene-d10	9.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecane	198	C14H30	000629-59-4	90
2	Hexadecane	226	C16H34	000544-76-3	86
3	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	86
4	Tridecane	184	C13H28	000629-50-5	83
5	Octadecane	254	C18H38	000593-45-3	83





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Sample : I6744-06
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : PROS-2-E(0-1)

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

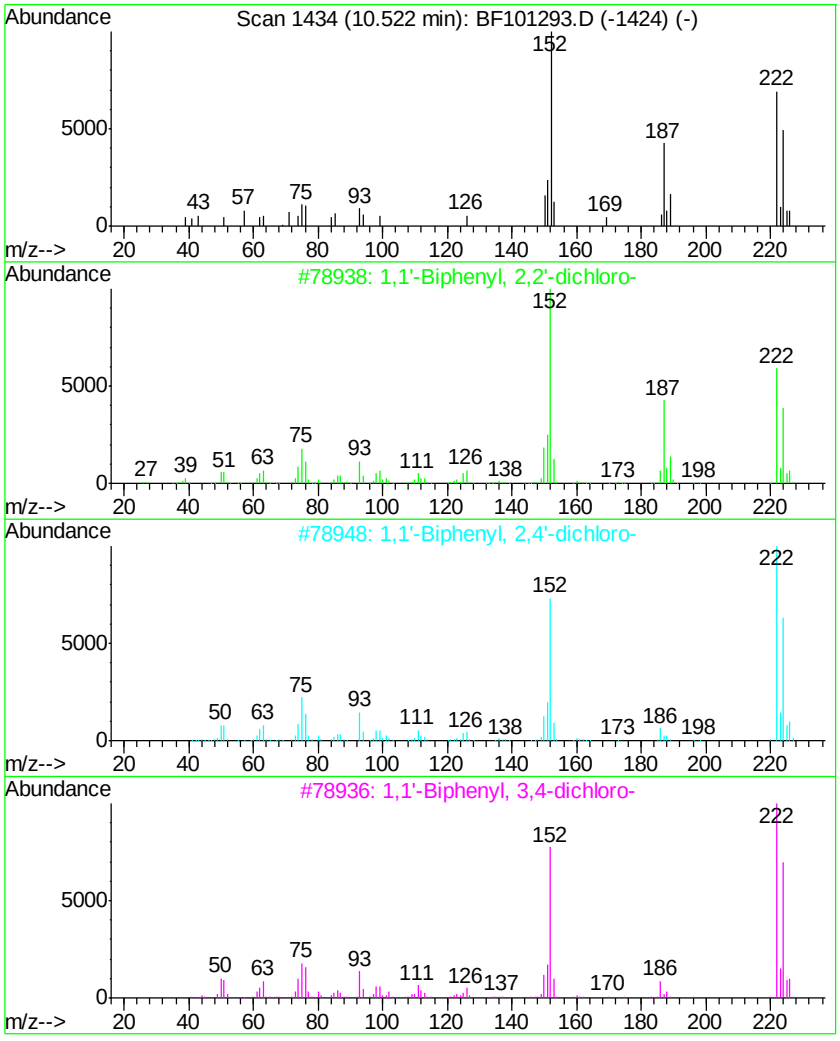
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

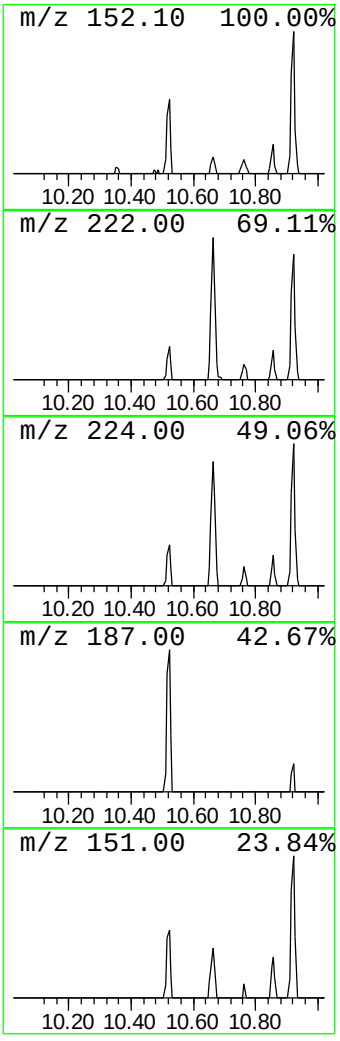
****

Peak Number 6 1,1'-Biphenyl, 2,2'-dichloro- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.52	2.43 ng	49584	Acenaphthene-d10	9.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2'-dichloro-	222	C12H8Cl2	013029-08-8	98
2	1,1'-Biphenyl, 2,4'-dichloro-	222	C12H8Cl2	034883-43-7	95
3	1,1'-Biphenyl, 3,4-dichloro-	222	C12H8Cl2	002974-92-7	95
4	1,1'-Biphenyl 2,3'-dichloro-	222	C12H8Cl2	025569-80-6	94
5	1,1'-Biphenyl, 3,4'-dichloro-	222	C12H8Cl2	002974-90-5	93





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Operator : SJ/JU
Sample : I6744-06
Misc :
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Instrument : BNA_F
ClientSampleId : PROS-2-E(0-1)

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

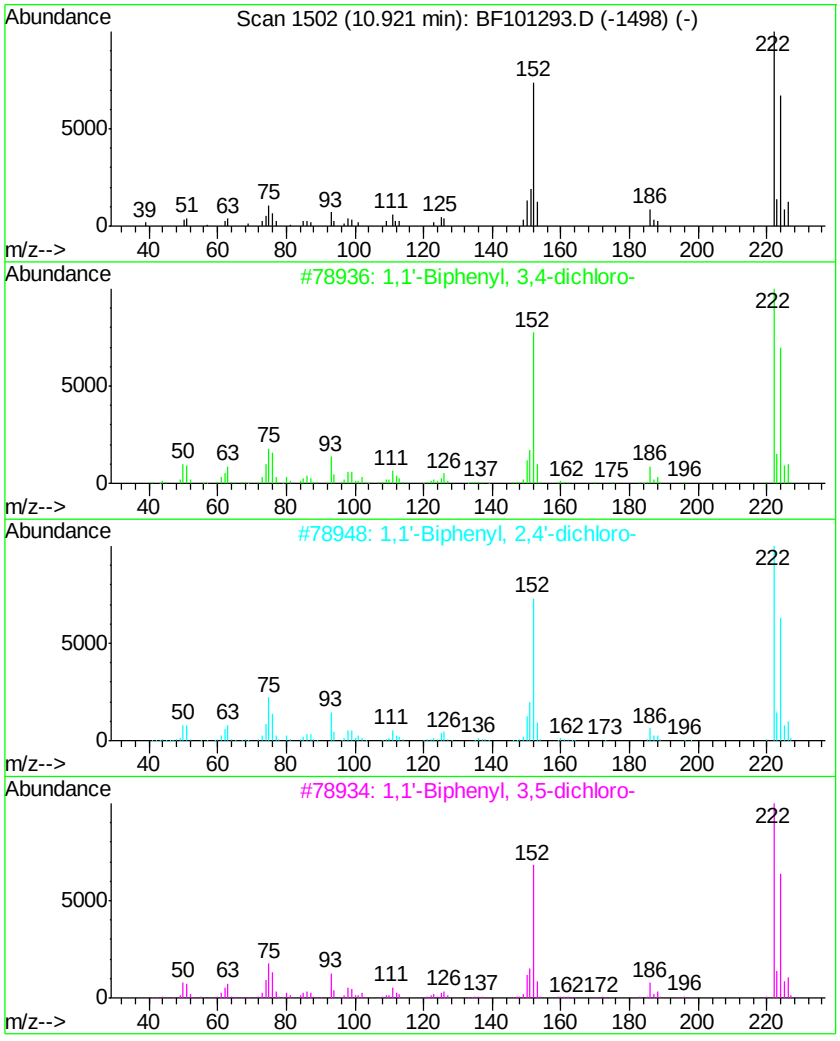
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

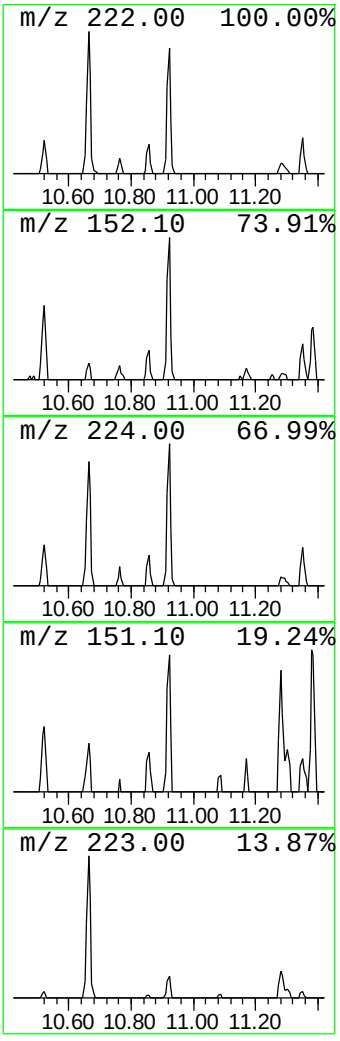
****

Peak Number 8 1,1'-Biphenyl, 3,4-dichloro- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.92	2.95 ng	60667	Phenanthrene-d10	11.36

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 3,4-dichloro-	222	C12H8Cl2	002974-92-7	98
2	1,1'-Biphenyl, 2,4'-dichloro-	222	C12H8Cl2	034883-43-7	98
3	1,1'-Biphenyl, 3,5-dichloro-	222	C12H8Cl2	034883-41-5	98
4	1,1'-Biphenyl, 3,3'-dichloro-	222	C12H8Cl2	002050-67-1	98
5	1,1'-Biphenyl, 2,3-dichloro-	222	C12H8Cl2	016605-91-7	97





Data Path : Z:\HPCHEM1\BNA F\DATA\BF121117\
Data File : BF101293.D
Acq On : 11 Dec 2017 16:56
Operator : SJ/JU
Sample : I6744-06
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : PROS-2-E(0-1)

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

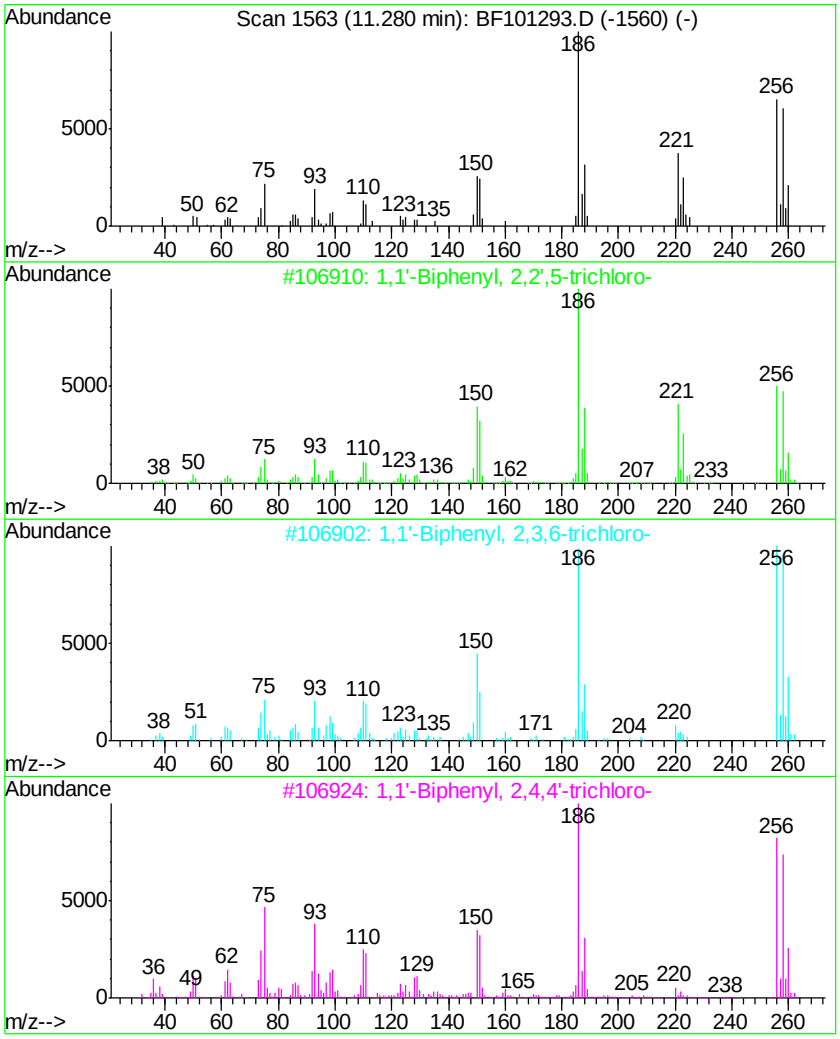
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

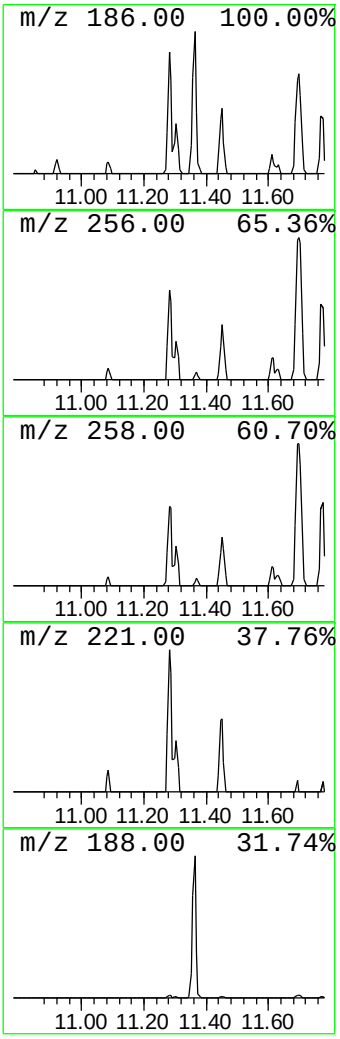
****

Peak Number 9 1,1'-Biphenyl, 2,2',5-trich... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.28	3.11 ng	64000	Phenanthrene-d10	11.36

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	99
2	1,1'-Biphenyl, 2,3,6-trichloro-	256	C12H7Cl3	055702-45-9	95
3	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	93
4	1,1'-Biphenyl, 3,4,4'-Trichloro-	256	C12H7Cl3	038444-90-5	70
5	1,1'-Biphenyl, 2,4,6-trichloro-	256	C12H7Cl3	035693-92-6	60





Data Path : Z:\HPCHEM1\BNA F\DATA\BF121117\
Data File : BF101293.D
Acq On : 11 Dec 2017 16:56
Operator : SJ/JU
Sample : I6744-06
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : PROS-2-E(0-1)

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

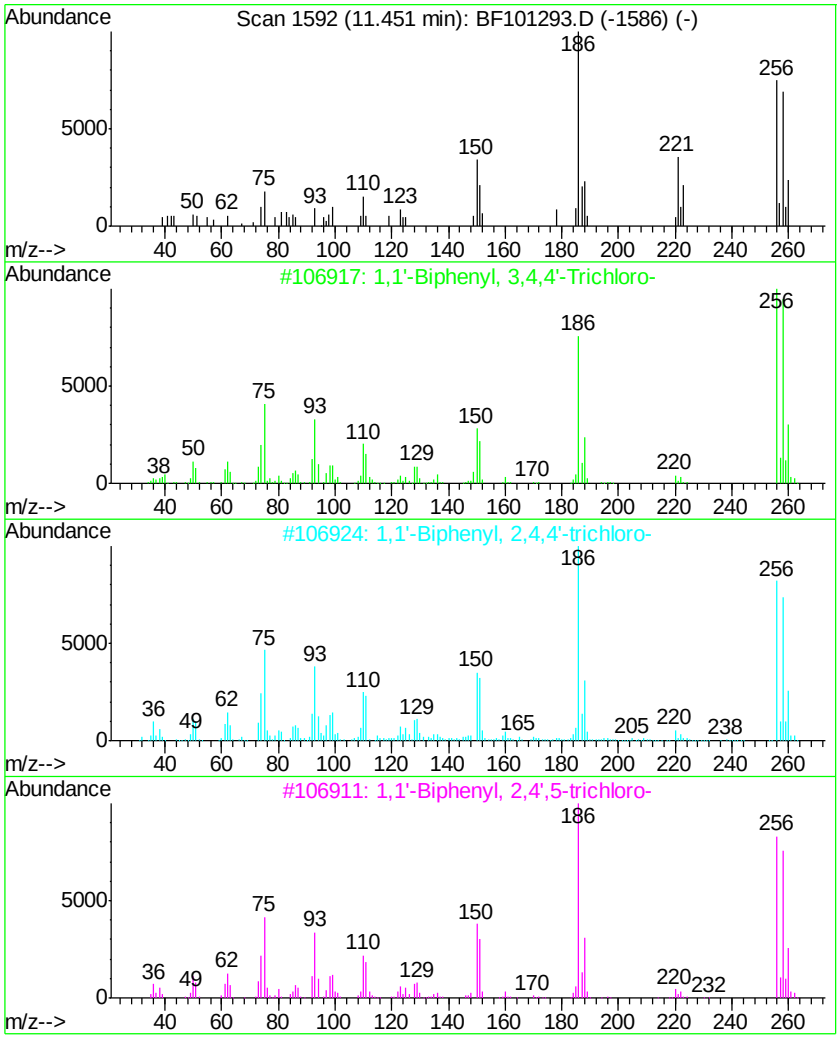
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

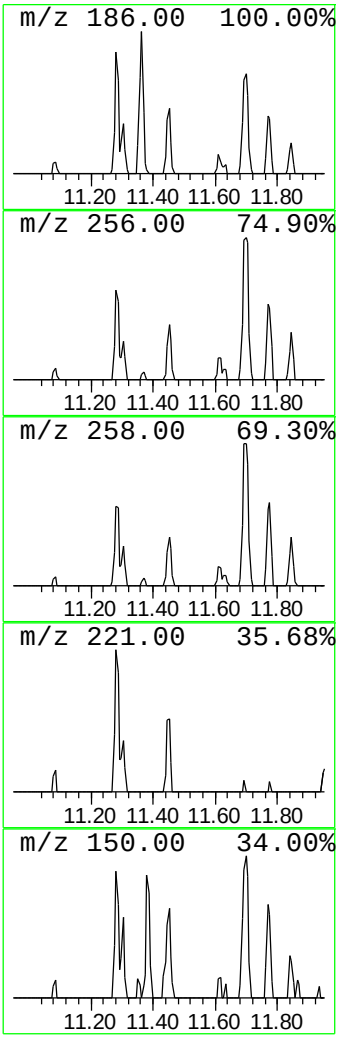
****

Peak Number 10 1,1'-Biphenyl, 3,4,4'-Trich... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.45	2.91 ng	59890	Phenanthrene-d10	11.36

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 3,4,4'-Trichloro-	256	C12H7Cl3	038444-90-5	93
2	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	90
3	1,1'-Biphenyl, 2,4',5-trichloro-	256	C12H7Cl3	016606-02-3	90
4	1,1'-Biphenyl, 2,4,6-trichloro-	256	C12H7Cl3	035693-92-6	90
5	1,1'-Biphenyl, 2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	89





Data Path : Z:\HPCHEM1\BNA F\DATA\BF121117\
Data File : BF101293.D
Acq On : 11 Dec 2017 16:56
Operator : SJ/JU
Sample : I6744-06
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : PROS-2-E(0-1)

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

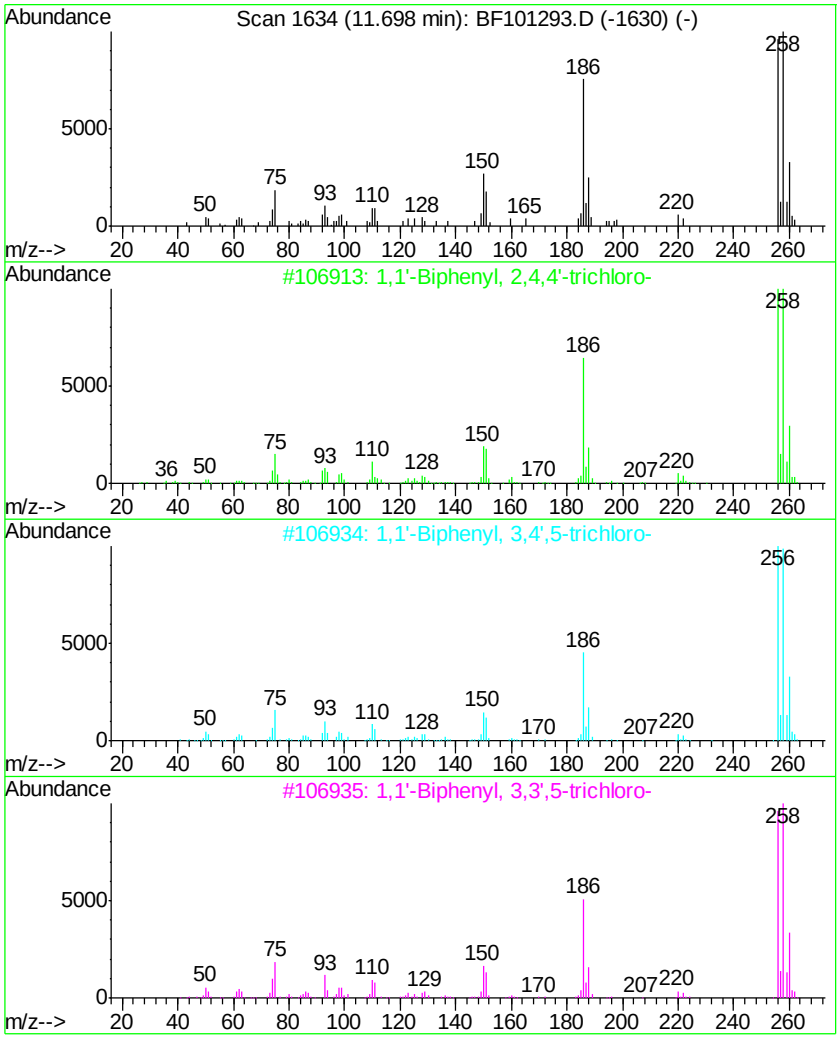
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

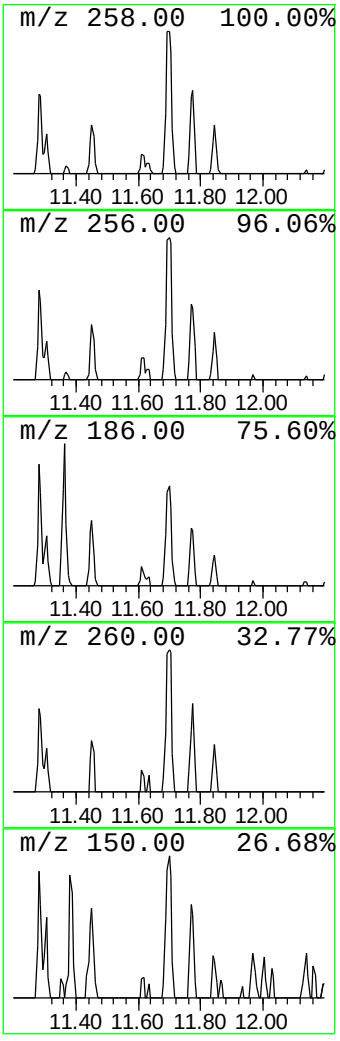
****

Peak Number 11 1,1'-Biphenyl, 2,4,4'-trich... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.70	5.41 ng	111420	Phenanthrene-d10	11.36

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	99
2	1,1'-Biphenyl, 3,4',5-trichloro-	256	C12H7Cl3	038444-88-1	99
3	1,1'-Biphenyl, 3,3',5-trichloro-	256	C12H7Cl3	038444-87-0	99
4	1,1'-Biphenyl, 2,3,4-trichloro-	256	C12H7Cl3	055702-46-0	99
5	1,1'-Biphenyl, 2',3,4-trichloro-	256	C12H7Cl3	038444-86-9	99





Data Path : Z:\HPCHEM1\BNA F\DATA\BF121117\
Data File : BF101293.D
Acq On : 11 Dec 2017 16:56
Operator : SJ/JU
Sample : I6744-06
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : PROS-2-E(0-1)

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

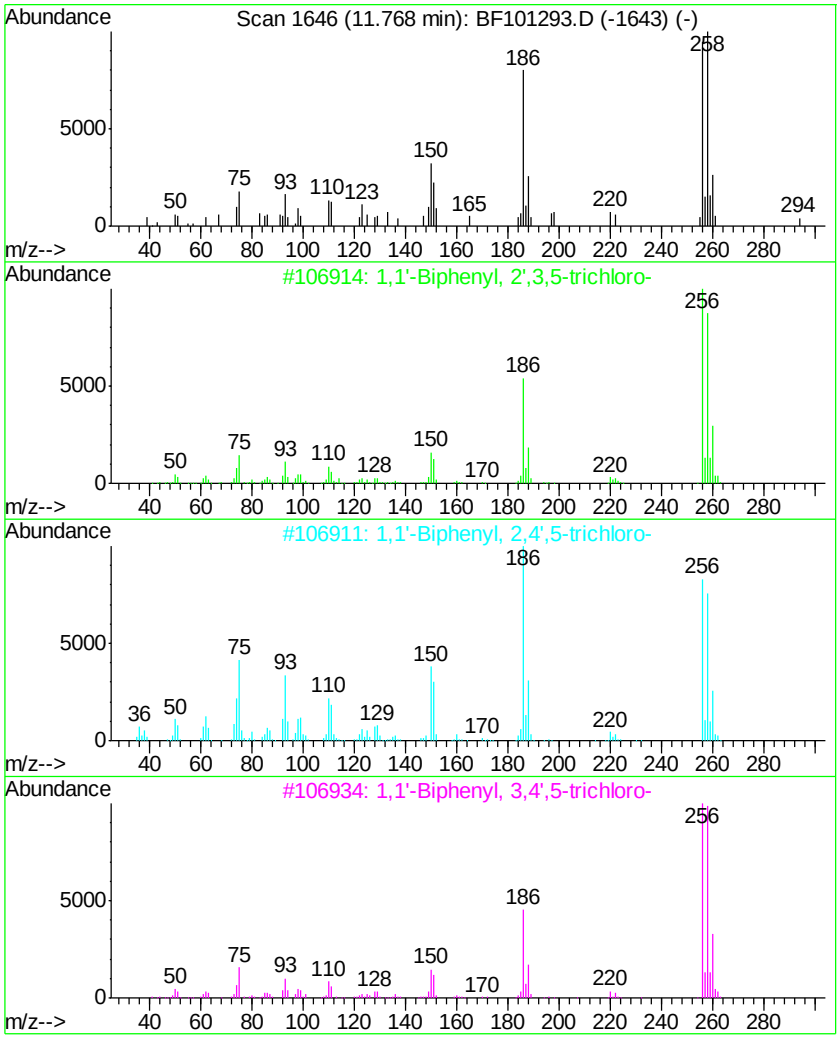
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

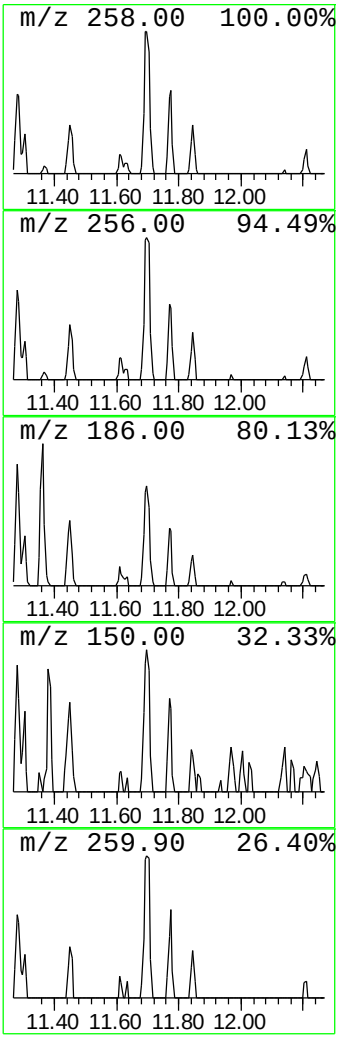
****

Peak Number 12 1,1'-Biphenyl, 2',3,5-trich... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.77	2.24 ng	46015	Phenanthrene-d10	11.36

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2',3,5-trichloro-	256	C12H7Cl3	037680-68-5	99
2	1,1'-Biphenyl, 2,4',5-trichloro-	256	C12H7Cl3	016606-02-3	99
3	1,1'-Biphenyl, 3,4',5-trichloro-	256	C12H7Cl3	038444-88-1	99
4	1,1'-Biphenyl, 2',3,4-trichloro-	256	C12H7Cl3	038444-86-9	99
5	1,1'-Biphenyl, 2,3,6-trichloro-	256	C12H7Cl3	055702-45-9	99





Data Path : Z:\HPCHEM1\BNA F\DATA\BF121117\
Data File : BF101293.D
Acq On : 11 Dec 2017 16:56
Operator : SJ/JU
Sample : I6744-06
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : PROS-2-E(0-1)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF113017.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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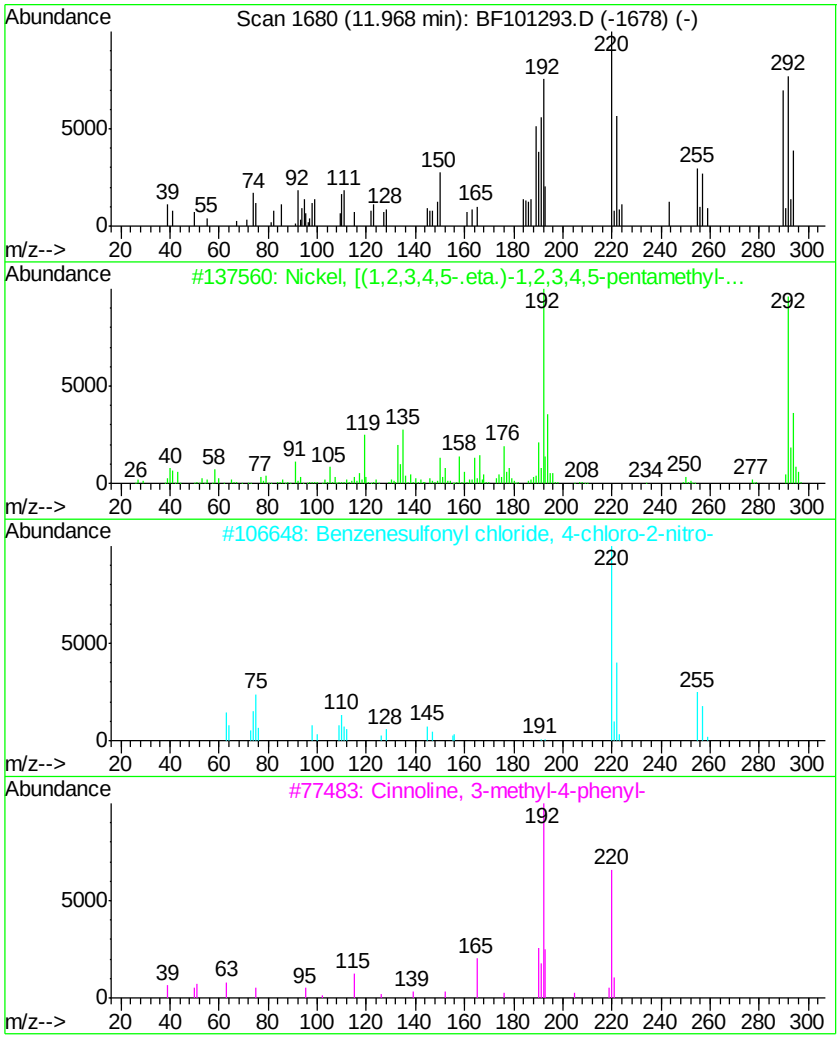
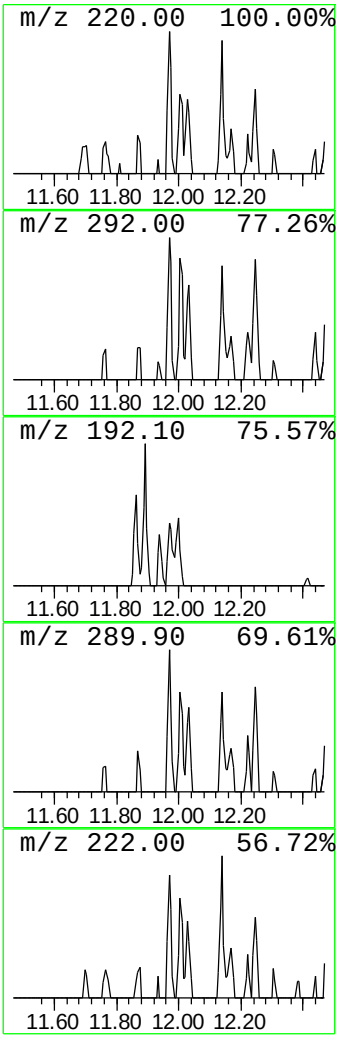
Peak Number 13 unknown11.97 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.97	2.18 ng	44794	Phenanthrene-d10	11.36

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nickel, [(1,2,3,4,5-.eta.)-1,2,3...	292	C15H22NiO2	097210-29-2	35
2	Benzenesulfonyl chloride, 4-chlo...	255	C6H3Cl2NO4S	004533-96-4	25
3	Cinnoline, 3-methyl-4-phenyl-	220	C15H12N2	021039-71-4	22
4	6H-Indolo[3,2,1-de][1,5]naphthyr...	220	C14H8N2O	000479-43-6	22
5	2,2'-Bifuran, 3,3'-dibromo-	290	C8H4Br2O2	066721-51-5	18





Data Path : Z:\HPCHEM1\BNA F\DATA\BF121117\
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Operator : SJ/JU
Sample : I6744-06
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : PROS-2-E(0-1)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF113017.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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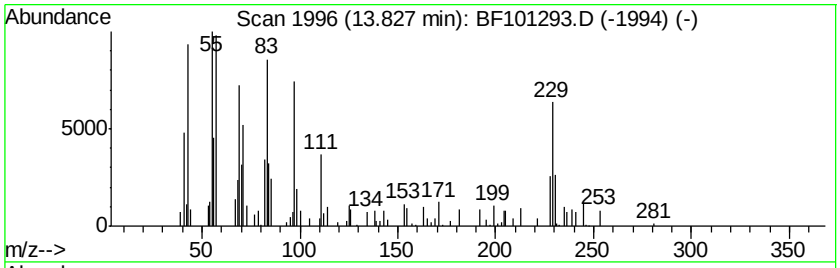
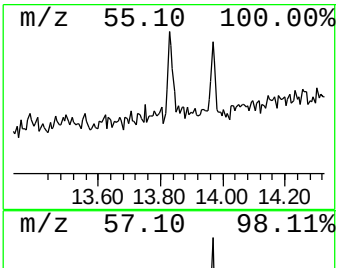
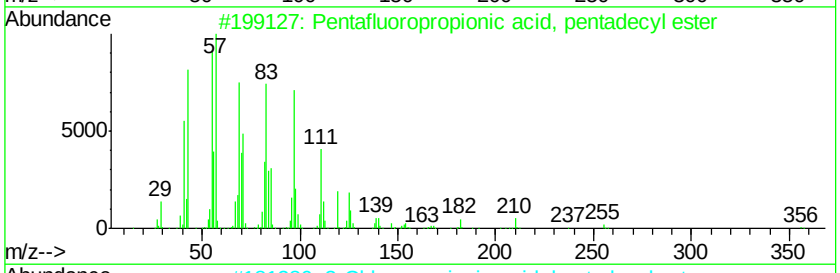
Peak Number 14 Pentafluoropropionic acid, ... Concentration Rank 11

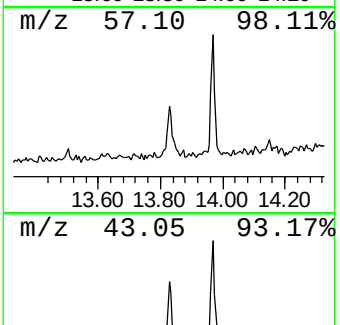
R.T.	EstConc	Area	Relative to ISTD	R.T.
13.83	2.38 ng	49390	Chrysene-d12	14.01

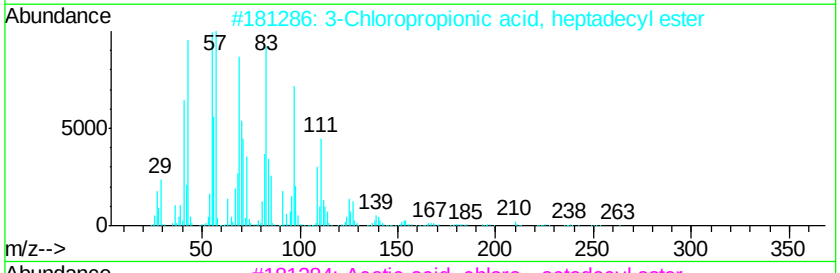
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	58
2	3-Chloropropionic acid, heptadec...	346	C20H39ClO2	1000283-05-1	52
3	Acetic acid, chloro-, octadecyl ...	346	C20H39ClO2	005348-82-3	52
4	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	52
5	Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	52

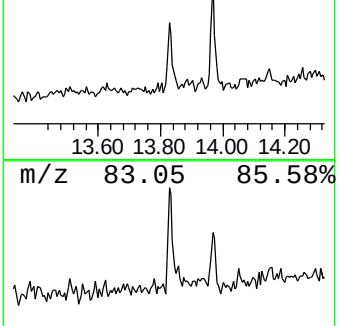


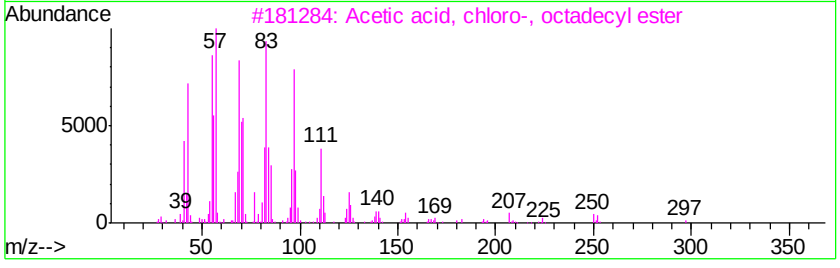


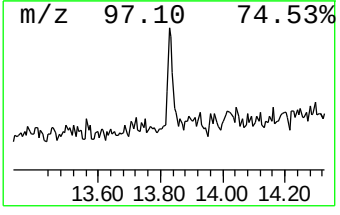












Tentatively Identified Compound (LSC) summary

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121117\
Data File : BF101293.D
Acq On : 11 Dec 2017 16:56
Operator : SJ/JU
Sample : I6744-06
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : PROS-2-E(0-1)

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

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.06	4.8	ng	73590	1	6.83	305655	20.0
unknown6.60	6.60	79.3	ng	1212630	1	6.83	305655	20.0
Hexadecane	9.79	2.4	ng	47912	3	9.87	408265	20.0
Tetradecane	10.29	2.2	ng	44319	3	9.87	408265	20.0
1,1'-Biphenyl, 2,...	10.52	2.4	ng	49584	3	9.87	408265	20.0
1,1'-Biphenyl, 3,...	10.92	3.0	ng	60667	4	11.36	411598	20.0
1,1'-Biphenyl, 2,...	11.28	3.1	ng	64000	4	11.36	411598	20.0
1,1'-Biphenyl, 3,...	11.45	2.9	ng	59890	4	11.36	411598	20.0
1,1'-Biphenyl, 2,...	11.70	5.4	ng	111420	4	11.36	411598	20.0
1,1'-Biphenyl, 2'...	11.77	2.2	ng	46015	4	11.36	411598	20.0
unknown11.97	11.97	2.2	ng	44794	4	11.36	411598	20.0
Pentafluoropropio...	13.83	2.4	ng	49390	5	14.01	415050	20.0