LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF022618\
Data File : BF103212.D
Acq On : 26 Feb 2018 12:21
Operator : SJ/JU
Sample : J1648-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RO-124031-RO124035

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-BF022218.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.146	4	10	19	rVB	1083502	1469113	40.51%	3.865%
2	5.104	510	513	524	rVB	153220	196287	5.41%	0.516%
3	5.492	575	579	593	rBV	3526916	3626771	100.00%	9.542%
4	6.504	746	751	754	rBV	3184822	3343168	92.18%	8.796%
5	6.639	770	774	778	rBV	3581048	3469237	95.66%	9.127%
6	6.869	809	813	818	rBV	1199119	959038	26.44%	2.523%
7	7.028	835	840	843	rBV	2803577	2637391	72.72%	6.939%
8	7.433	904	909	912	rBV	2355516	2288557	63.10%	6.021%
9	8.151	1027	1031	1051	rVB	1392795	1316577	36.30%	3.464%
10	9.222	1209	1213	1223	rBV	3041744	3131162	86.33%	8.238%
11	9.298	1223	1226	1229	rVB	69187	63136	1.74%	0.166%
12	9.563	1265	1271	1274	rBV4	32286	52737	1.45%	0.139%
13	9.616	1277	1280	1288	rVB	268318	265563	7.32%	0.699%
14	9.769	1303	1306	1311	rBV2	53610	54871	1.51%	0.144%
15	9.827	1311	1316	1319	rVV	177069	145133	4.00%	0.382%
16	9.910	1324	1330	1333	rVV	1234535	1217703	33.58%	3.204%
17	9.939	1333	1335	1338	rVB	70005	58777	1.62%	0.155%
18	10.057	1351	1355	1358	rBV2	29484	42341	1.17%	0.111%
19	10.110	1361	1364	1368	rVV2	44868	52032	1.43%	0.137%
20	10.145	1368	1370	1374	rVB3	53996	53278	1.47%	0.140%
21	10.327	1394	1401	1404	rBV	202283	201822	5.56%	0.531%
22	10.457	1420	1423	1427	rVB	54598	54481	1.50%	0.143%
23	10.545	1432	1438	1440	rBV2	75469	89295	2.46%	0.235%
24	10.698	1460	1464	1474	rBV	1727285	1705316	47.02%	4.487%
25	10.798	1478	1481	1487	rVV	164940	202251	5.58%	0.532%
26	11.004	1511	1516	1518	rBV5	32482	52162	1.44%	0.137%
27	11.245	1554	1557	1560	rBV2	75302	71133	1.96%	0.187%
28	11.274	1560	1562	1564	rBV	46863	41216	1.14%	0.108%
29	11.398	1579	1583	1585	rBV	1064048	974713	26.88%	2.564%
30	11.421	1585	1587	1592	rVV	559370	456896	12.60%	1.202%
31	11.474	1593	1596	1599	rVB	99519	89167	2.46%	0.235%
32	11.633	1618	1623	1628	rBV3	52922	95221	2.63%	0.251%
33	11.898	1666	1668	1671	rBV2	76487	102028	2.81%	0.268%
34	11.927	1671	1673	1675	rVV	107135	81984	2.26%	0.216%

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF022618\
Data File : BF103212.D
Acq On : 26 Feb 2018 12:21
Operator : SJ/JU
Sample : J1648-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RO-124031-RO124035

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-BF022218.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	11.974	1679	1681	1684	rBV	50267	39662	1.09%	0.104%
36	12.010	1684	1687	1690	rBV	131057	155881	4.30%	0.410%
37	12.080	1696	1699	1702	rBV2	55576	53471	1.47%	0.141%
38	12.192	1716	1718	1721	rVV	58520	46324	1.28%	0.122%
39	12.227	1721	1724	1729	rVB5	54768	63998	1.76%	0.168%
40	12.339	1738	1743	1746	rBV3	64043	94104	2.59%	0.248%
41	12.445	1759	1761	1764	rBV	64485	56234	1.55%	0.148%
42	12.539	1774	1777	1780	rVB	49178	54819	1.51%	0.144%
43	12.610	1786	1789	1793	rBV	759121	697115	19.22%	1.834%
44	12.792	1817	1820	1823	rVB	164877	147105	4.06%	0.387%
45	12.845	1823	1829	1831	rBV	834077	900372	24.83%	2.369%
46	12.892	1835	1837	1840	rVB2	48571	50716	1.40%	0.133%
47	12.980	1844	1852	1855	rBV	2270868	2089620	57.62%	5.498%
48	13.004	1855	1856	1860	rVB	66455	44373	1.22%	0.117%
49	13.057	1862	1865	1869	rVV3	60250	93573	2.58%	0.246%
50	13.168	1882	1884	1887	rVB	93722	84106	2.32%	0.221%
51	13.233	1893	1895	1898	rVB	63316	53100	1.46%	0.140%
52	13.274	1899	1902	1904	rVV2	88054	78924	2.18%	0.208%
53	13.368	1915	1918	1920	rBV	58230	51021	1.41%	0.134%
54	13.398	1920	1923	1925	rBV2	72568	66016	1.82%	0.174%
55	13.498	1937	1940	1943	rVB	101352	96428	2.66%	0.254%
56	13.680	1969	1971	1975	rVB	74508	68403	1.89%	0.180%
57	13.815	1992	1994	1996	rBV	51313	41942	1.16%	0.110%
58	13.862	2000	2002	2010	rVB4	406020	532977	14.70%	1.402%
59	14.004	2024	2026	2028	rBV	60242	70100	1.93%	0.184%
60	14.045	2028	2033	2035	rVV2	951883	1272882	35.10%	3.349%
61	14.068	2035	2037	2045	rVB	414657	384168	10.59%	1.011%
62	15.098	2209	2212	2216	rBV	308492	479714	13.23%	1.262%
63	15.409	2262	2265	2271	rVB	222702	279650	7.71%	0.736%
64	15.474	2272	2276	2281	rBV	293977	385208	10.62%	1.013%
65	15.539	2282	2287	2290	rBV	604560	817325	22.54%	2.150%

Sum of corrected areas: 38009888

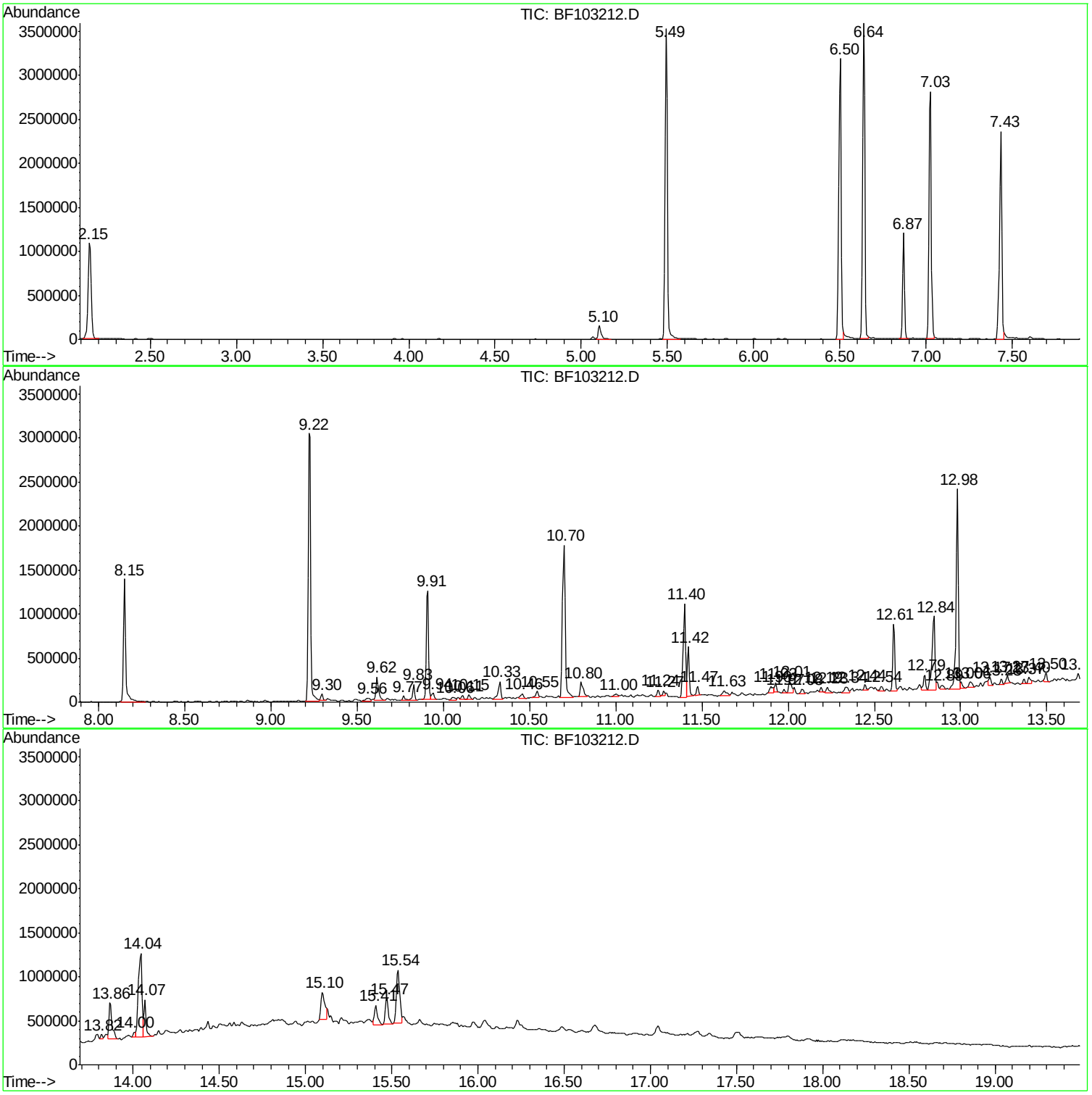
LSC Report - Integrated Chromatogram

Data Path : Z:\HPCHEM1\BNA F\DATA\BF022618\
Data File : BF103212.D
Acq On : 26 Feb 2018 12:21
Operator : SJ/JU
Sample : J1648-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RO-124031-RO124035

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.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\BF022618\
Data File : BF103212.D
Acq On : 26 Feb 2018 12:21
Operator : SJ/JU
Sample : J1648-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RO-124031-RO124035

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

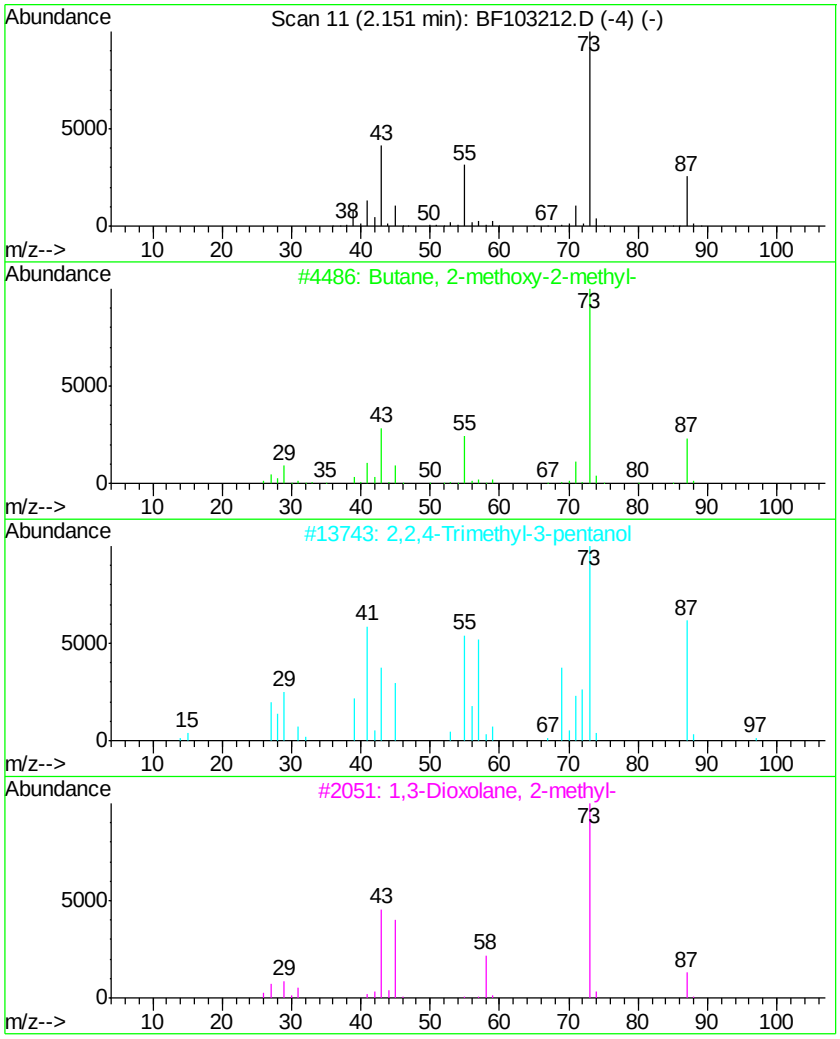
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

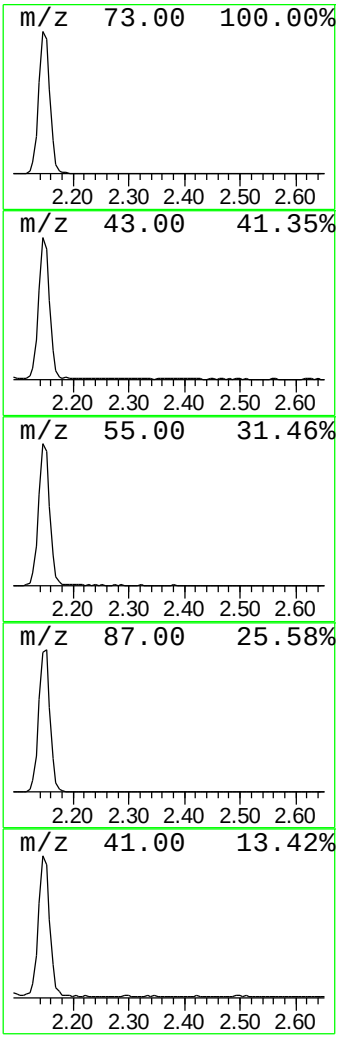
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.15	30.64 ng	1469110	1,4-Dichlorobenzene-d4	6.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	23
4	3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	23
5	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF022618\
Data File : BF103212.D
Acq On : 26 Feb 2018 12:21
Operator : SJ/JU
Sample : J1648-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RO-124031-RO124035

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

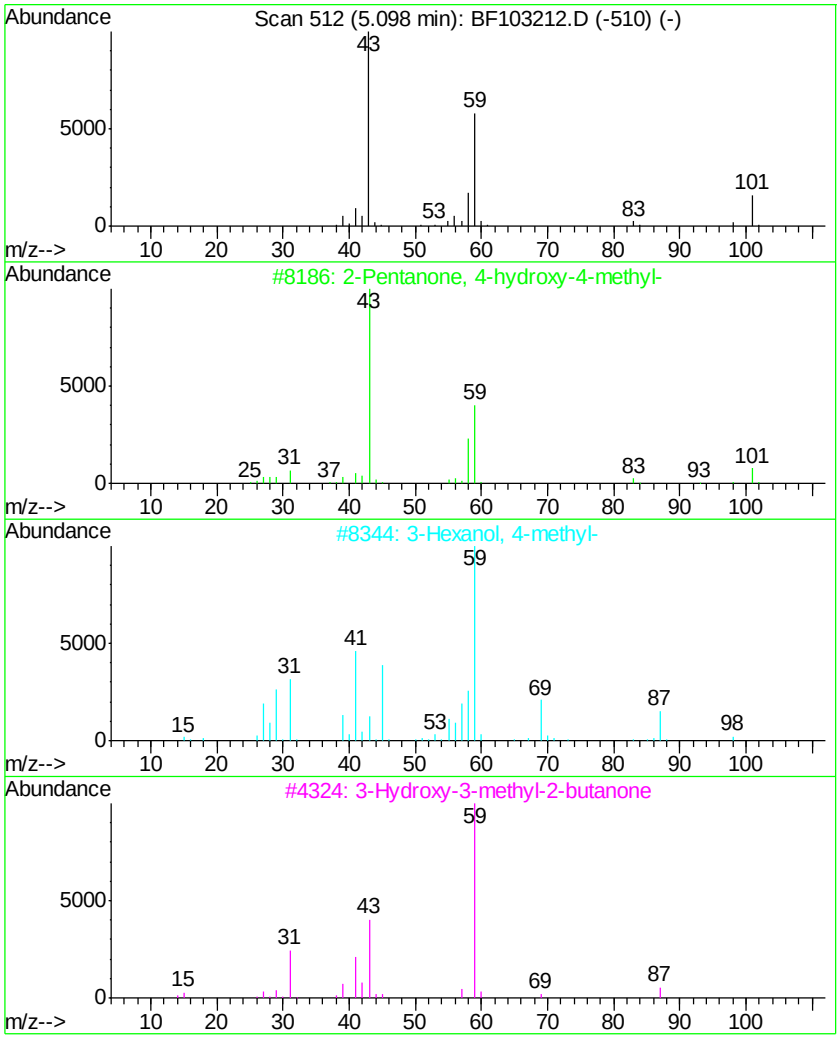
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

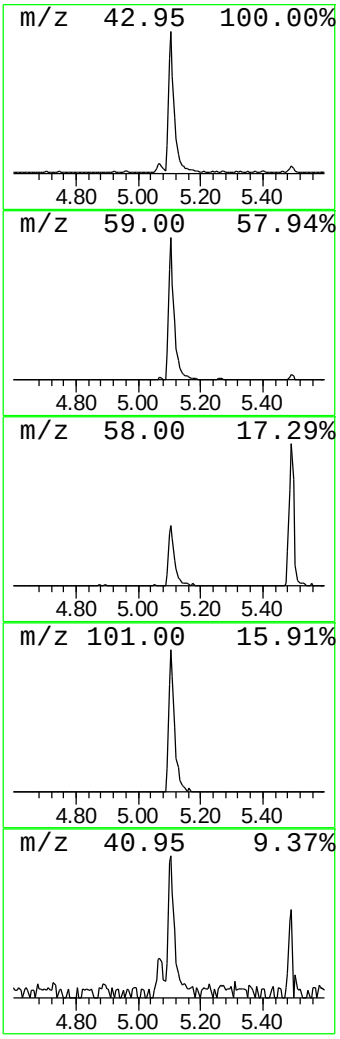
****

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.10	4.09 ng	196287	1,4-Dichlorobenzene-d4	6.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
3	3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	23
4	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
5	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16





Data Path : Z:\HPCHEM1\BNA F\DATA\BF022618\
Data File : BF103212.D
Acq On : 26 Feb 2018 12:21
Operator : SJ/JU
Sample : J1648-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RO-124031-RO124035

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.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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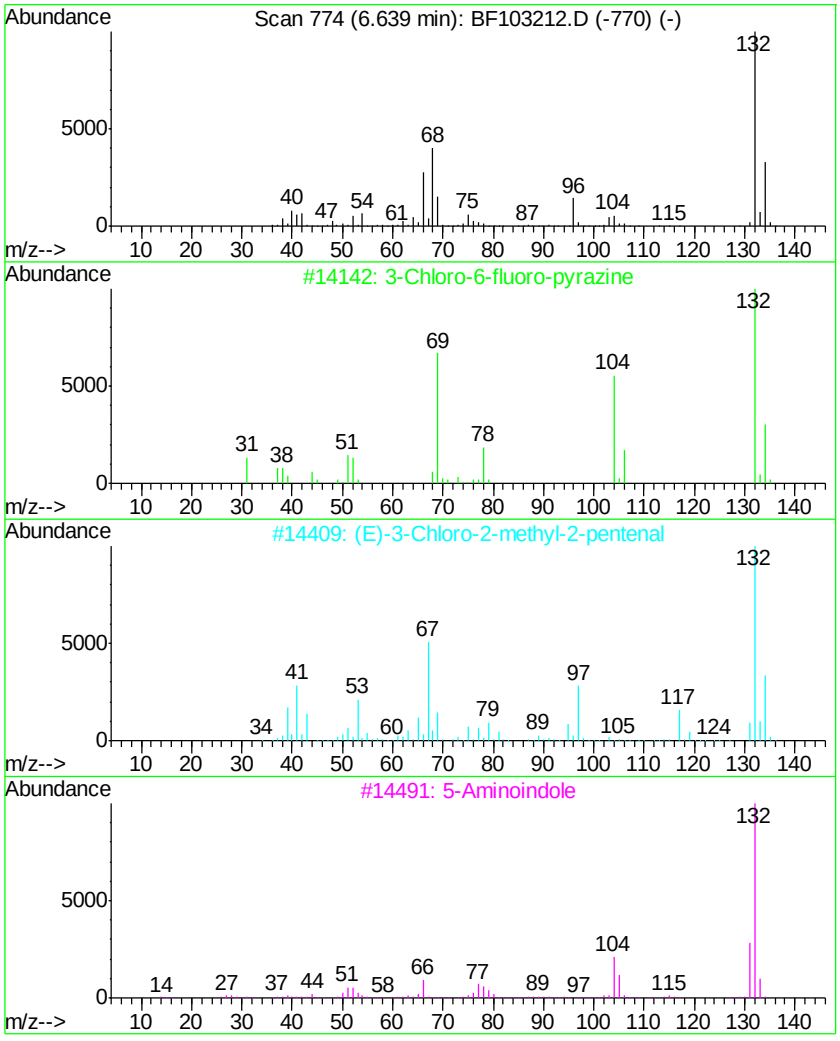
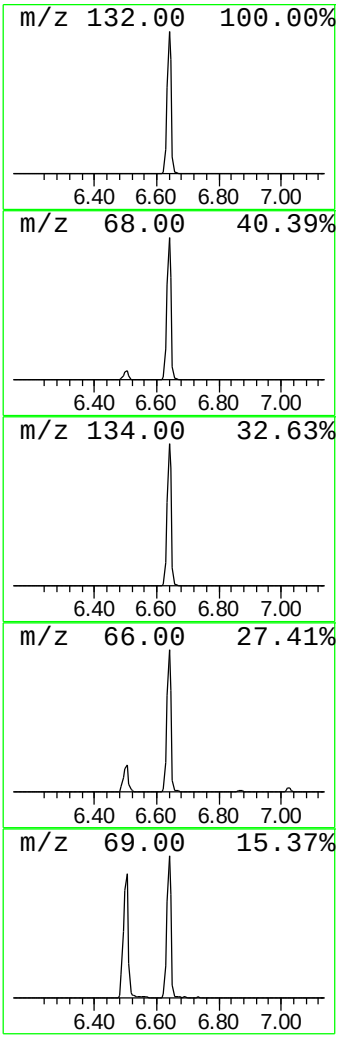
Peak Number 3 unknown6.64 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.64	72.35 ng	3469240	1,4-Dichlorobenzene-d4	6.87

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	1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9





Data Path : Z:\HPCHEM1\BNA F\DATA\BF022618\
Data File : BF103212.D
Acq On : 26 Feb 2018 12:21
Operator : SJ/JU
Sample : J1648-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RO-124031-RO124035

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.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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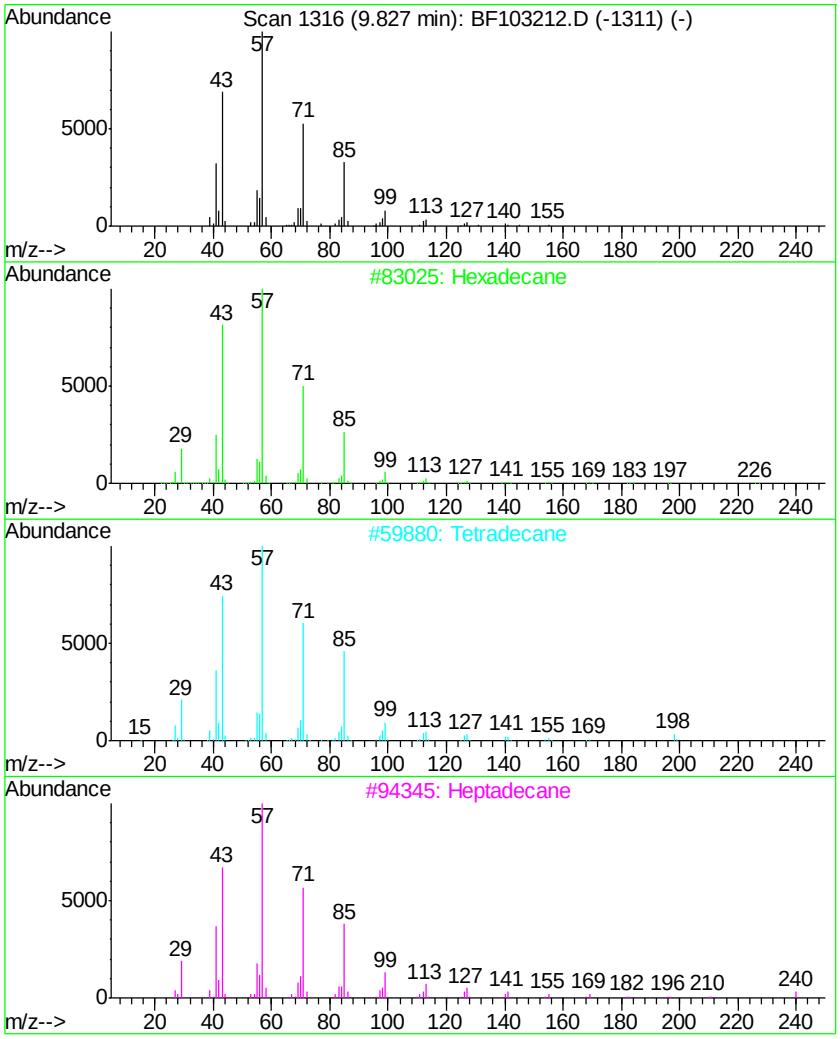
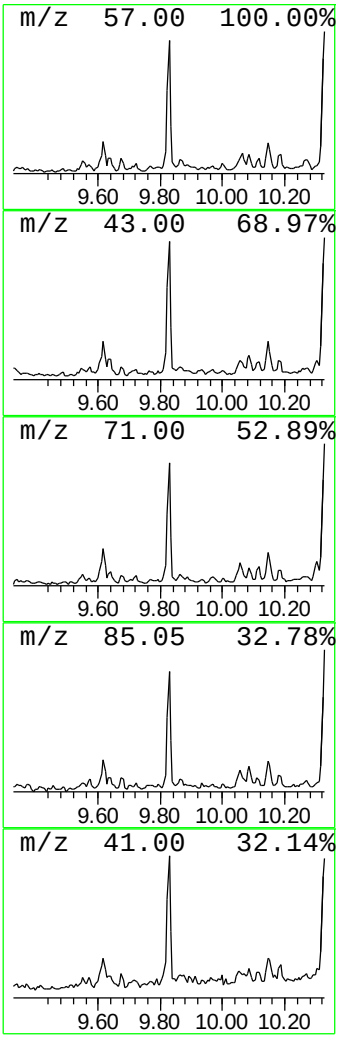
Peak Number 5 Hexadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.83	2.38 ng	145133	Acenaphthene-d10	9.91

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	91
2	Tetradecane	198	C14H30	000629-59-4	86
3	Heptadecane	240	C17H36	000629-78-7	86
4	Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	83
5	Heptadecane, 2,6,10,14-tetramethyl-	296	C21H44	018344-37-1	80





Data Path : Z:\HPCHEM1\BNA F\DATA\BF022618\
Data File : BF103212.D
Acq On : 26 Feb 2018 12:21
Operator : SJ/JU
Sample : J1648-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RO-124031-RO124035

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.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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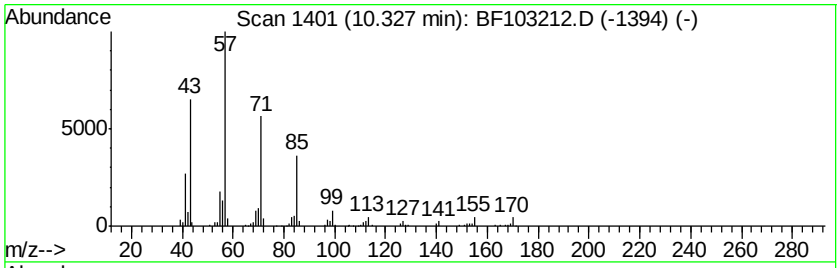
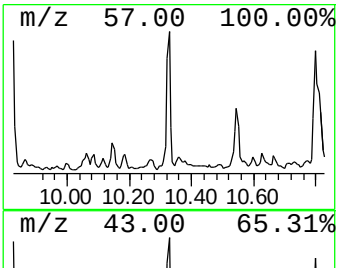
Peak Number 6 Eicosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.33	3.31 ng	201822	Acenaphthene-d10	9.91

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	86
2	Hexadecane	226	C16H34	000544-76-3	86
3	Sulfurous acid, 2-ethylhexyl hex...	278	C14H30O3S	1000309-20-2	80
4	Tetradecane	198	C14H30	000629-59-4	80
5	Dodecane	170	C12H26	000112-40-3	68





Data Path : Z:\HPCHEM1\BNA F\DATA\BF022618\
Data File : BF103212.D
Acq On : 26 Feb 2018 12:21
Operator : SJ/JU
Sample : J1648-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RO-124031-RO124035

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

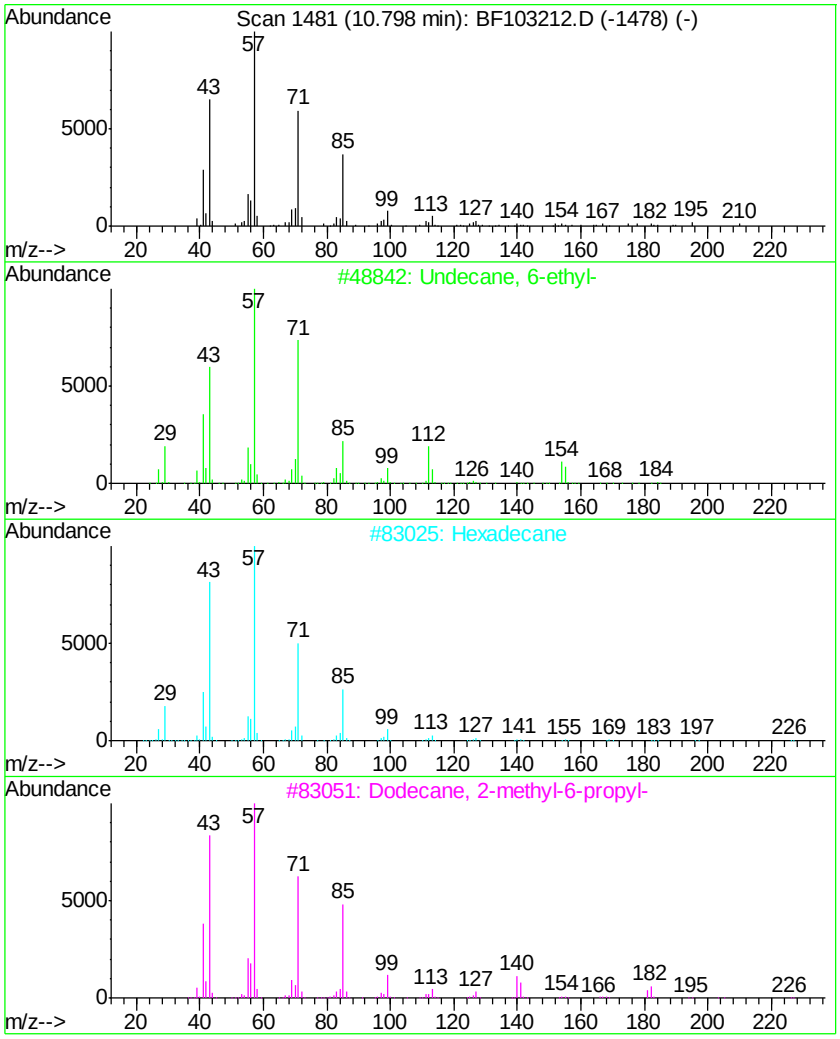
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

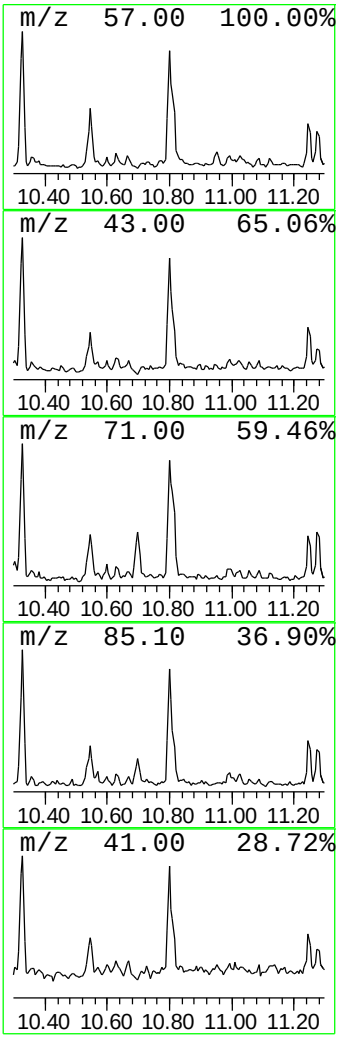
****

Peak Number 7 Undecane, 6-ethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.80	4.15 ng	202251	Phenanthrene-d10	11.40

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane, 6-ethyl-	184	C13H28	017312-60-6	90
2	Hexadecane	226	C16H34	000544-76-3	90
3	Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	80
4	Heptacosane	380	C27H56	000593-49-7	80
5	Pentadecane	212	C15H32	000629-62-9	78





Data Path : Z:\HPCHEM1\BNA F\DATA\BF022618\
Data File : BF103212.D
Acq On : 26 Feb 2018 12:21
Operator : SJ/JU
Sample : J1648-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RO-124031-RO124035

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

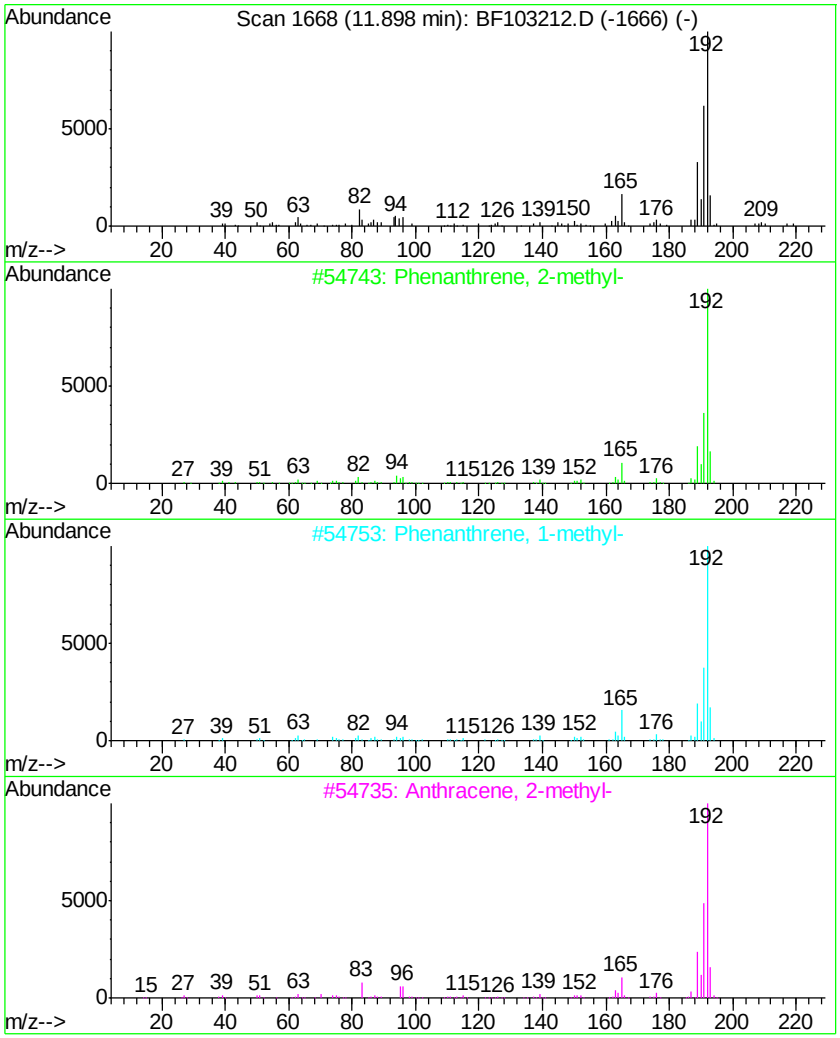
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

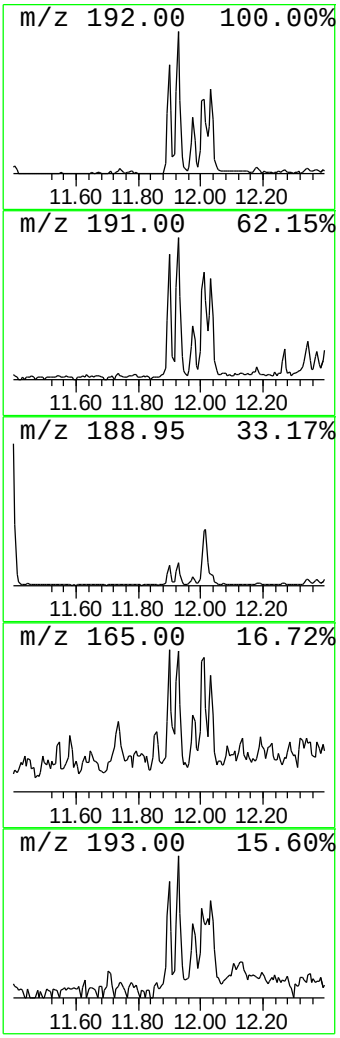
****

Peak Number 8 Phenanthrene, 2-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.90	2.09 ng	102028	Phenanthrene-d10	11.40

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
3	Anthracene, 2-methyl-	192	C15H12	000613-12-7	92
4	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	91
5	Anthracene, 9-methyl-	192	C15H12	000779-02-2	91





Data Path : Z:\HPCHEM1\BNA F\DATA\BF022618\
Data File : BF103212.D
Acq On : 26 Feb 2018 12:21
Operator : SJ/JU
Sample : J1648-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RO-124031-RO124035

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.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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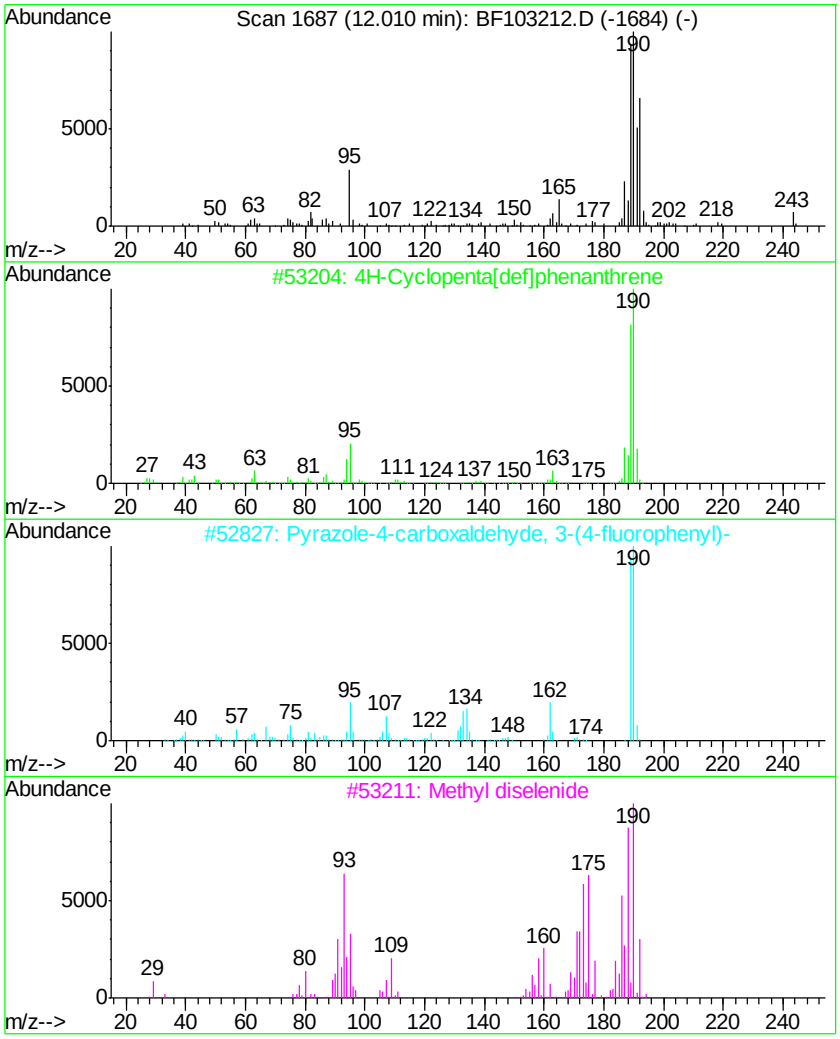
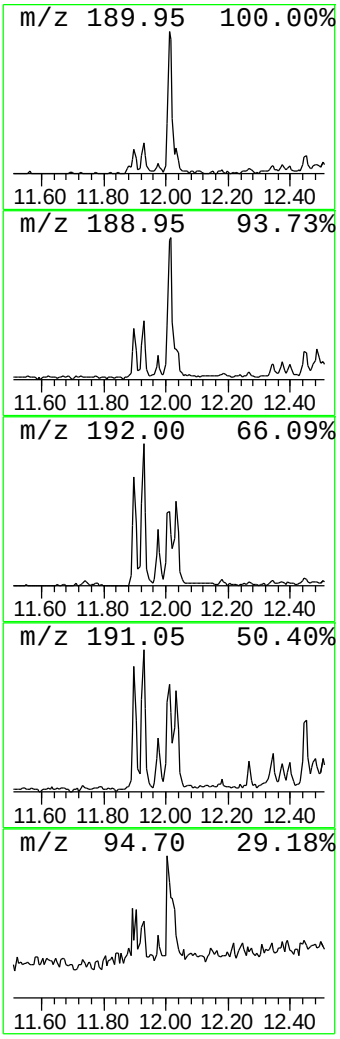
Peak Number 9 4H-Cyclopenta[def]phenanthrene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.01	3.20 ng	155881	Phenanthrene-d10	11.40

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	64
2	Pyrazole-4-carboxaldehyde, 3-(4-...	190	C10H7FN2O	306936-57-2	47
3	Methyl diselenide	190	C2H6Se2	007101-31-7	43
4	Carbonic acid, ethyl 2-formyl-4,...	262	C10H8Cl2O4	1000331-34-4	40
5	9-(Chloromethyl)anthracene	226	C15H11Cl	024463-19-2	38





Data Path : Z:\HPCHEM1\BNA F\DATA\BF022618\
Data File : BF103212.D
Acq On : 26 Feb 2018 12:21
Operator : SJ/JU
Sample : J1648-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RO-124031-RO124035

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.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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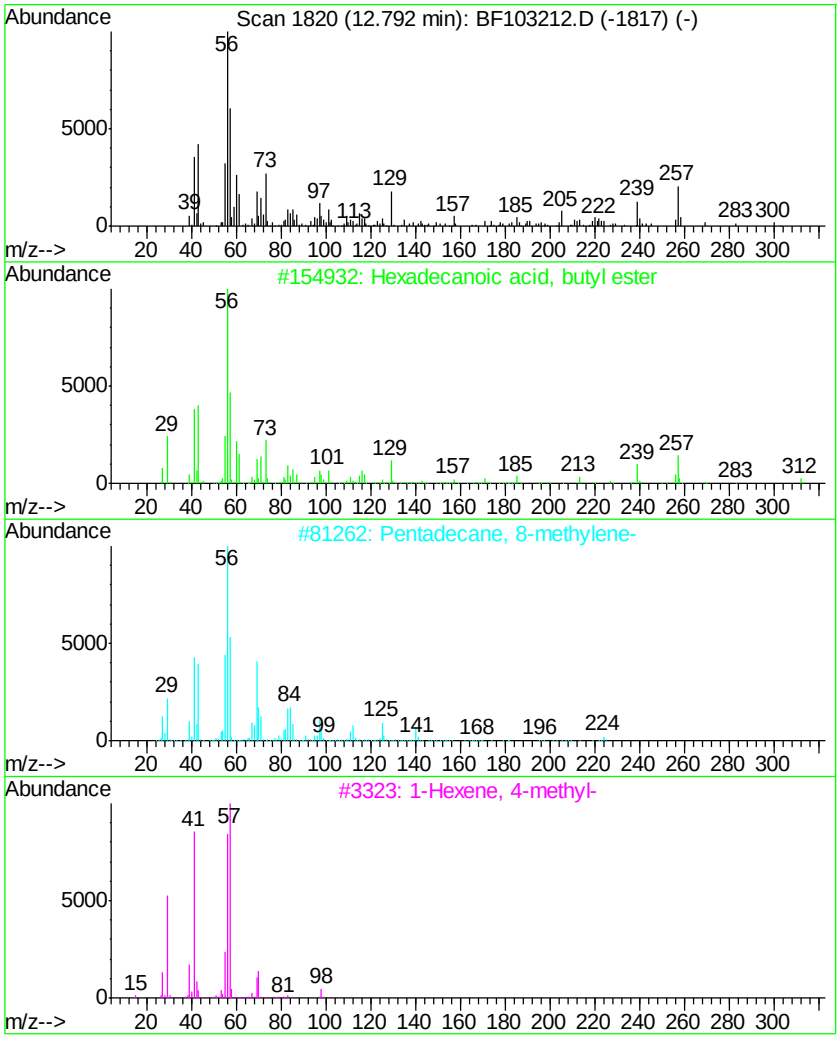
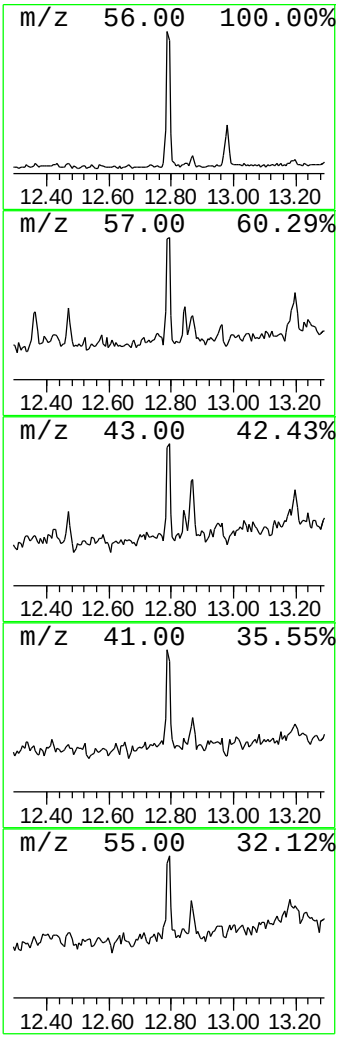
Peak Number 10 Hexadecanoic acid, butyl ester Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.79	2.31 ng	147105	Chrysene-d12	14.04

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecanoic acid, butyl ester	312	C20H40O2	000111-06-8	90
2	Pentadecane, 8-methylene-	224	C16H32	055668-09-2	38
3	1-Hexene, 4-methyl-	98	C7H14	003769-23-1	38
4	2,4-Dipropyl-5-ethyl-1,3-dioxane	200	C12H24O2	006413-83-8	37
5	Pentanal, 3-methyl-	100	C6H12O	015877-57-3	35





Data Path : Z:\HPCHEM1\BNA F\DATA\BF022618\
Data File : BF103212.D
Acq On : 26 Feb 2018 12:21
Operator : SJ/JU
Sample : J1648-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RO-124031-RO124035

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.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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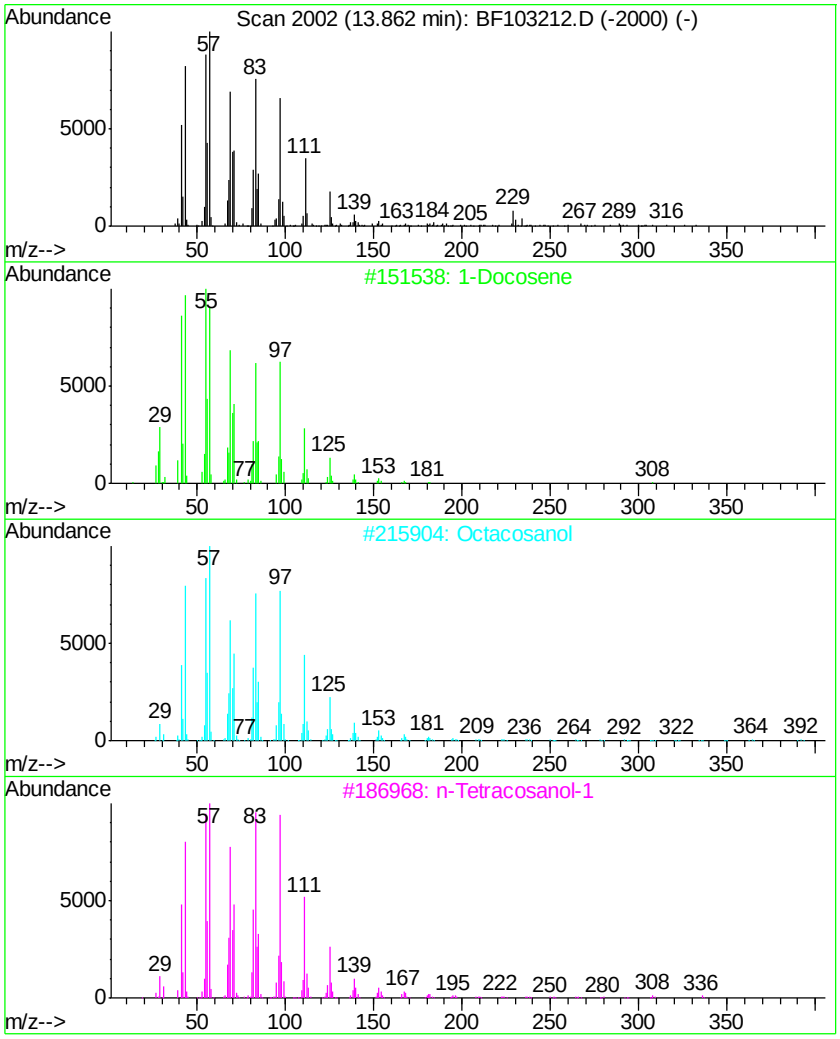
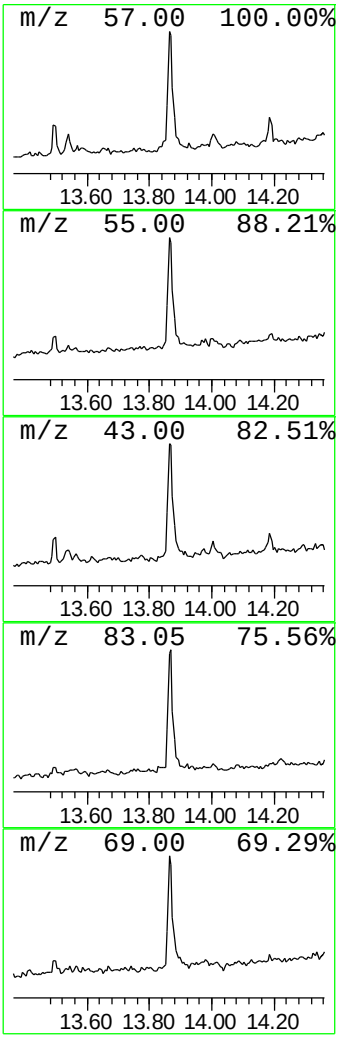
Peak Number 11 1-Docosene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.86	8.37 ng	532977	Chrysene-d12	14.04

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Docosene	308	C22H44	001599-67-3	98
2	Octacosanol	410	C28H58O	000557-61-9	91
3	n-Tetracosanol-1	354	C24H50O	000506-51-4	91
4	Behenic alcohol	326	C22H46O	000661-19-8	91
5	5-Eicosene, (E)-	280	C20H40	074685-30-6	90





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF022618\
Data File : BF103212.D
Acq On : 26 Feb 2018 12:21
Operator : SJ/JU
Sample : J1648-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RO-124031-RO124035

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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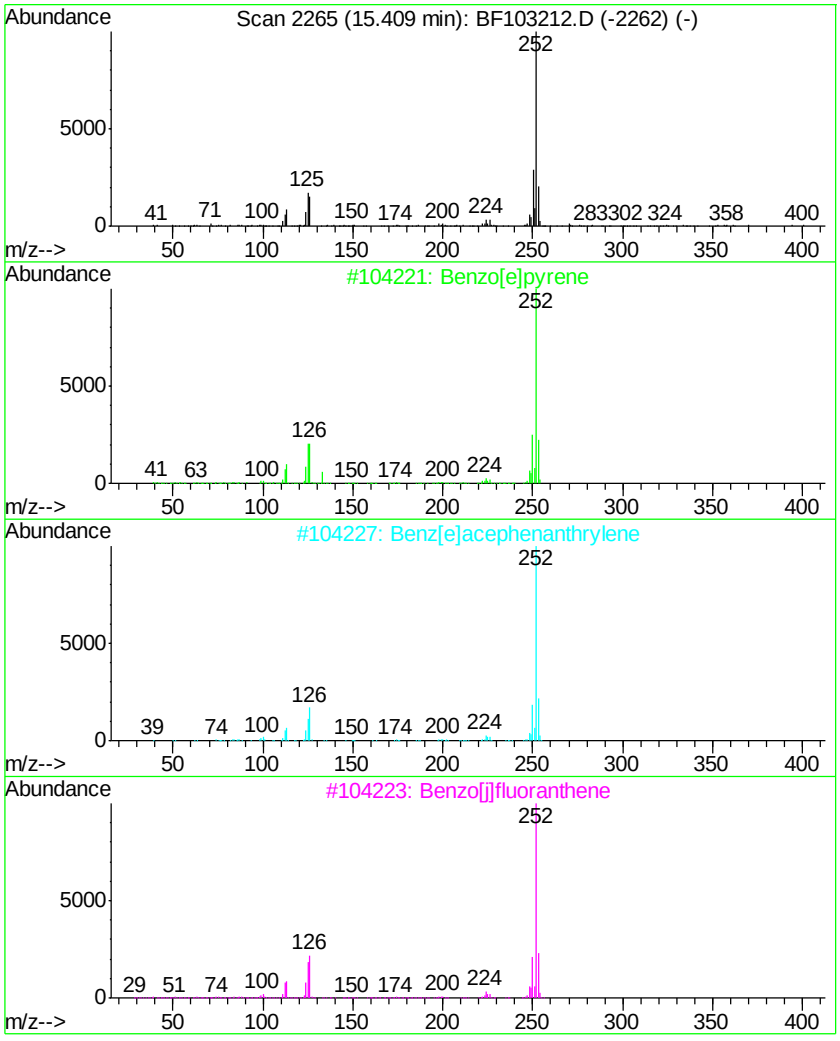
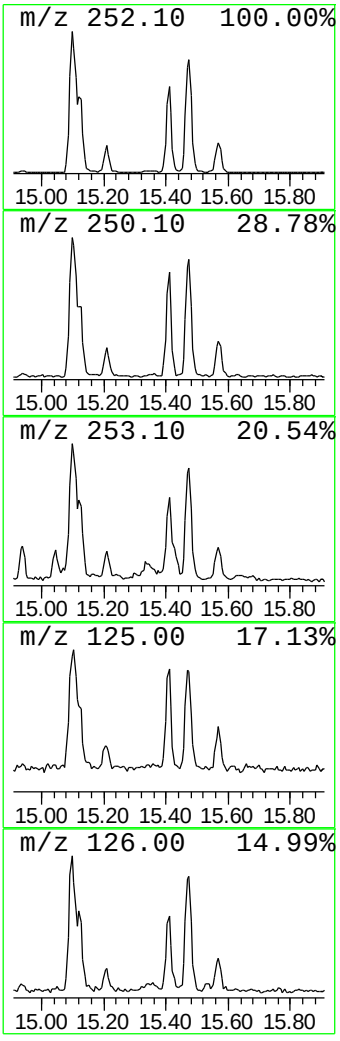
Peak Number 12 Benzo[e]pyrene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.41	6.84 ng	279650	Perylene-d12	15.54

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[e]pyrene	252	C20H12	000192-97-2	98
2	Benz[e]acephenanthrylene	252	C20H12	000205-99-2	98
3	Benzo[j]fluoranthene	252	C20H12	000205-82-3	96
4	Benzo[a]pyrene	252	C20H12	000050-32-8	95
5	Benzo[k]fluoranthene	252	C20H12	000207-08-9	94





Tentatively Identified Compound (LSC) summary

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF022618\
Data File : BF103212.D
Acq On : 26 Feb 2018 12:21
Operator : SJ/JU
Sample : J1648-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RO-124031-RO124035

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Butane, 2-methoxy...	2.15	30.6	ng	1469110	1	6.87	959038	20.0
2-Pentanone, 4-hy...	5.10	4.1	ng	196287	1	6.87	959038	20.0
unknown6.64	6.64	72.3	ng	3469240	1	6.87	959038	20.0
Hexadecane	9.83	2.4	ng	145133	3	9.91	1217700	20.0
Eicosane	10.33	3.3	ng	201822	3	9.91	1217700	20.0
Undecane, 6-ethyl-	10.80	4.2	ng	202251	4	11.40	974713	20.0
Phenanthrene, 2-m...	11.90	2.1	ng	102028	4	11.40	974713	20.0
4H-Cyclopenta[def...	12.01	3.2	ng	155881	4	11.40	974713	20.0
Hexadecanoic acid...	12.79	2.3	ng	147105	5	14.04	1272880	20.0
1-Docosene	13.86	8.4	ng	532977	5	14.04	1272880	20.0
Benzo[e]pyrene	15.41	6.8	ng	279650	6	15.54	817325	20.0