LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF103122\
Data File : BF130892.D
Acq On : 31 Oct 2022 22:03
Operator : CG\JU
Sample : N5374-01 2X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : OK-2-102822

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-BF102722.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF130892.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.204	15	20	31	rVB	175796	233970	34.63%	2.633%
2	2.316	35	39	46	rVB	15257	20826	3.08%	0.234%
3	4.040	328	332	337	rVB	9354	12147	1.80%	0.137%
4	5.157	518	522	528	rBV	90479	97756	14.47%	1.100%
5	5.363	553	557	560	rBV4	9252	11860	1.76%	0.133%
6	5.551	585	589	601	rVB	799288	675608	100.00%	7.604%
7	6.557	756	760	764	rBV	746203	639466	94.65%	7.197%
8	6.769	793	796	800	rBV	11341	12259	1.81%	0.138%
9	6.945	822	826	829	rBV	622514	485986	71.93%	5.470%
10	7.504	917	921	924	rBV	438745	355500	52.62%	4.001%
11	7.534	924	926	930	rVB2	18645	16752	2.48%	0.189%
12	7.581	930	934	938	rBV5	9385	15083	2.23%	0.170%
13	8.139	1024	1029	1033	rBV5	11919	25532	3.78%	0.287%
14	8.198	1037	1039	1041	rBV	14224	12008	1.78%	0.135%
15	8.228	1041	1044	1047	rBV	668507	535438	79.25%	6.026%
16	8.516	1089	1093	1097	rBV6	10868	14000	2.07%	0.158%
17	8.810	1141	1143	1148	rBV3	24009	23699	3.51%	0.267%
18	9.045	1178	1183	1186	rBV3	20619	24792	3.67%	0.279%
19	9.175	1202	1205	1210	rBV6	9856	14626	2.16%	0.165%
20	9.304	1223	1227	1230	rBV	840737	627512	92.88%	7.062%
21	9.380	1237	1240	1243	rBV	35174	26728	3.96%	0.301%
22	9.451	1248	1252	1256	rVB4	14295	15746	2.33%	0.177%
23	9.569	1269	1272	1276	rBV5	14952	20923	3.10%	0.235%
24	9.645	1282	1285	1287	rBV	19601	17606	2.61%	0.198%
25	9.698	1291	1294	1297	rVB4	15677	15729	2.33%	0.177%
26	9.763	1302	1305	1308	rBV4	18180	18222	2.70%	0.205%
27	9.851	1314	1320	1323	rBV	26712	29945	4.43%	0.337%
28	9.910	1327	1330	1334	rVB	36645	31022	4.59%	0.349%
29	9.992	1340	1344	1347	rVB	694373	562773	83.30%	6.334%
30	10.022	1347	1349	1352	rVB	30959	24453	3.62%	0.275%
31	10.092	1358	1361	1365	rVB6	13463	18270	2.70%	0.206%
32	10.145	1365	1370	1372	rBV5	14817	25417	3.76%	0.286%
33	10.192	1376	1378	1383	rVB3	16567	18030	2.67%	0.203%
34	10.233	1383	1385	1389	rBV4	18588	17392	2.57%	0.196%
35	10.310	1394	1398	1399	rBV3	15884	16032	2.37%	0.180%
36	10.410	1409	1415	1419	rBV2	45435	55982	8.29%	0.630%

LSC Area Percent Report

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Acq On : 31 Oct 2022 22:03
Operator : CG\JU
Sample : N5374-01 2X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
OK-2-102822

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-BF102722.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	10.539	1434	1437	1440	rBV	26166	24988	3.70%	0.281%
38	10.627	1450	1452	1455	rBV	20863	22089	3.27%	0.249%
39	10.780	1475	1478	1482	rBV	466692	384295	56.88%	4.325%
40	10.886	1493	1496	1502	rBV	137787	149325	22.10%	1.681%
41	11.022	1516	1519	1522	rBV	24819	26237	3.88%	0.295%
42	11.116	1533	1535	1539	rBV4	13263	18316	2.71%	0.206%
43	11.186	1543	1547	1548	rBV4	17421	19623	2.90%	0.221%
44	11.210	1548	1551	1552	rBV2	18713	19874	2.94%	0.224%
45	11.274	1560	1562	1566	rBV4	15641	24549	3.63%	0.276%
46	11.339	1570	1573	1576	rVV2	99011	100738	14.91%	1.134%
47	11.369	1576	1578	1584	rVB2	36925	62582	9.26%	0.704%
48	11.486	1594	1598	1600	rBV	676352	606842	89.82%	6.830%
49	11.510	1600	1602	1605	rVV	248312	196688	29.11%	2.214%
50	11.557	1608	1610	1612	rVV	39204	33275	4.93%	0.374%
51	11.586	1612	1615	1619	rVV2	33129	37555	5.56%	0.423%
52	11.710	1634	1636	1643	rVB2	38919	42567	6.30%	0.479%
53	11.769	1643	1646	1651	rBV2	54742	64304	9.52%	0.724%
54	11.816	1652	1654	1659	rVB2	37485	49665	7.35%	0.559%
55	11.986	1681	1683	1686	rBV2	47269	62605	9.27%	0.705%
56	12.016	1686	1688	1690	rVB	54449	43134	6.38%	0.485%
57	12.098	1699	1702	1704	rBV2	77512	80046	11.85%	0.901%
58	12.174	1713	1715	1720	rBV2	65132	61060	9.04%	0.687%
59	12.310	1735	1738	1742	rVB3	48268	58800	8.70%	0.662%
60	12.539	1774	1777	1779	rBV2	66072	67853	10.04%	0.764%
61	12.569	1779	1782	1785	rVV3	110031	111344	16.48%	1.253%
62	12.704	1802	1805	1808	rBV	369428	327689	48.50%	3.688%
63	12.933	1841	1844	1851	rVB2	352290	360891	53.42%	4.062%
64	13.074	1865	1868	1871	rBV	658681	547185	80.99%	6.158%
65	13.298	1904	1906	1909	rBV	114941	105028	15.55%	1.182%
66	14.133	2045	2048	2051	rBV2	409361	406949	60.23%	4.580%

Sum of corrected areas: 8885186

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF103122\
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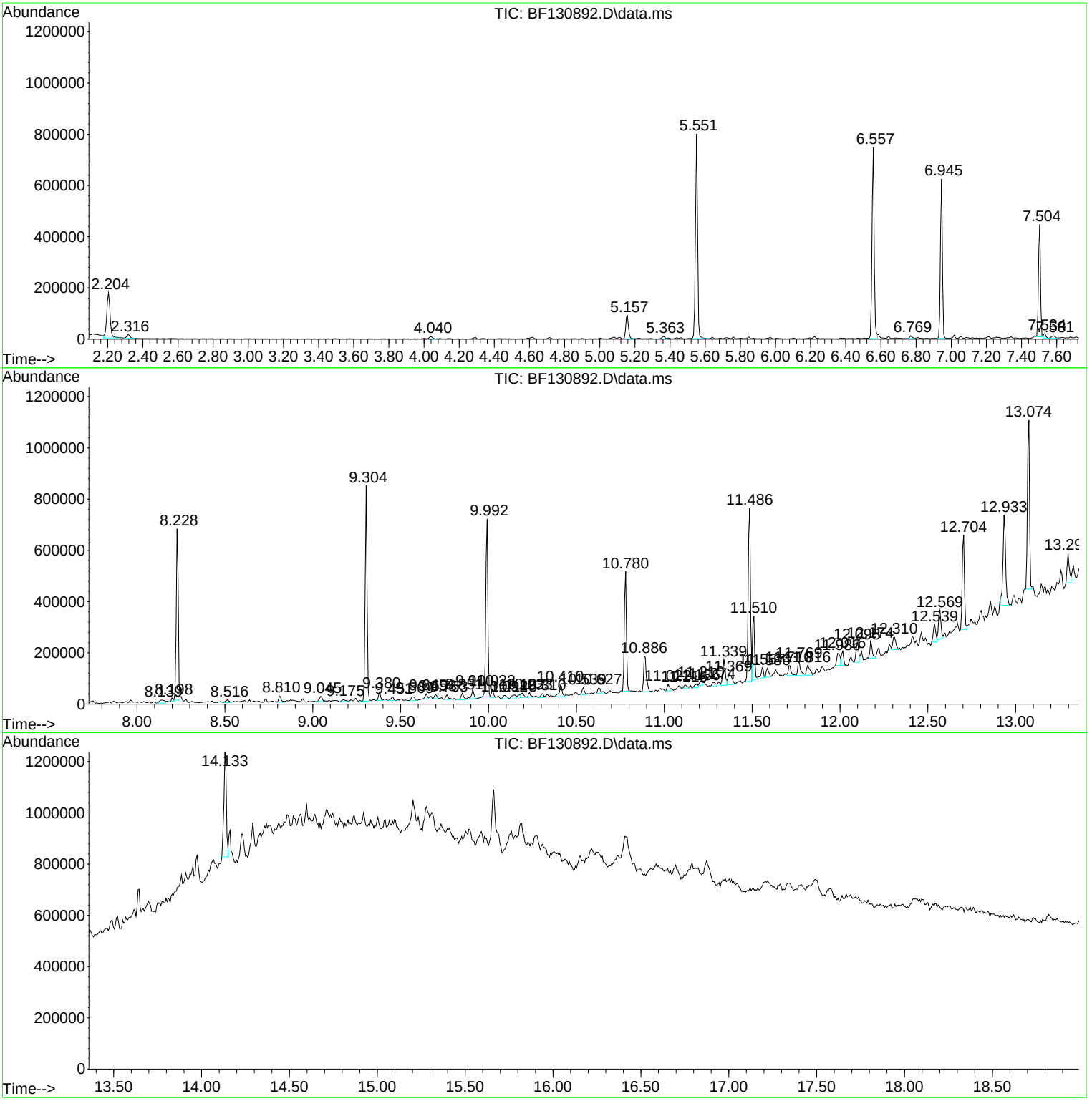
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
OK-2-102822

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF102722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF103122\
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

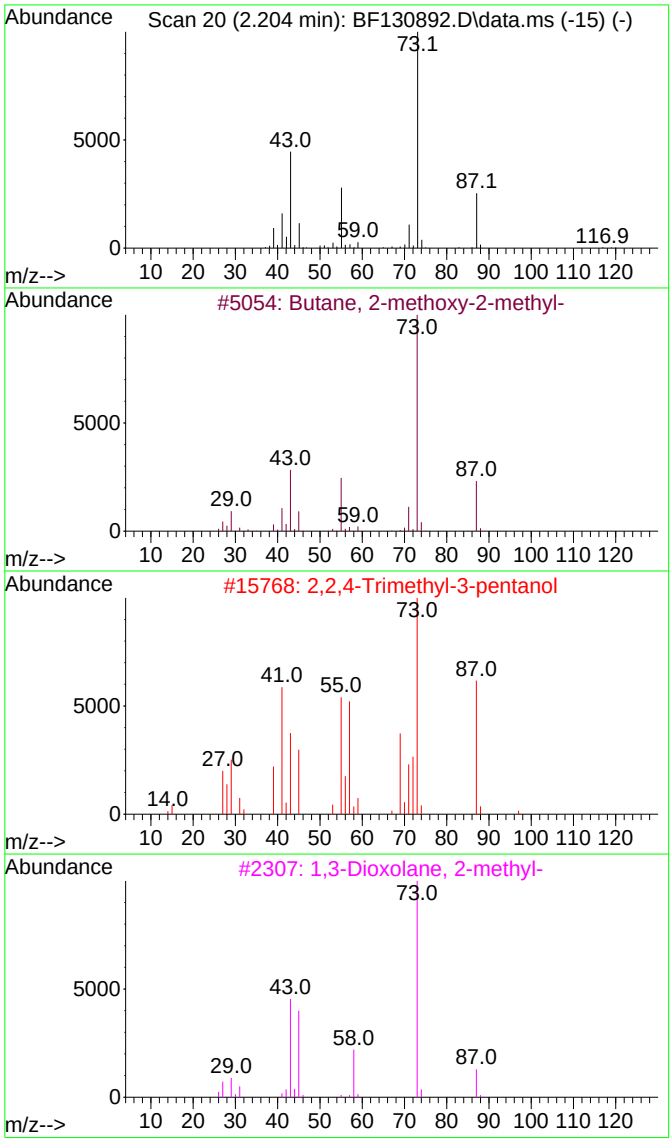
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

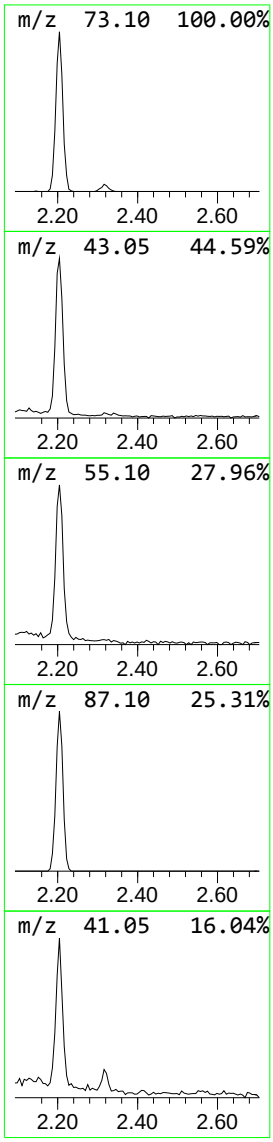
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.204	9.63 ng	233970	1,4-Dichlorobenzene-d4	6.945

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	Silane, tetramethyl-	88	C4H12Si	000075-76-3	23
5	1,3-Dioxolane, 2-propyl-	116	C6H12O2	003390-13-4	17





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF103122\
Data File : BF130892.D
Acq On : 31 Oct 2022 22:03
Operator : CG\JU
Sample : N5374-01 2X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
OK-2-102822

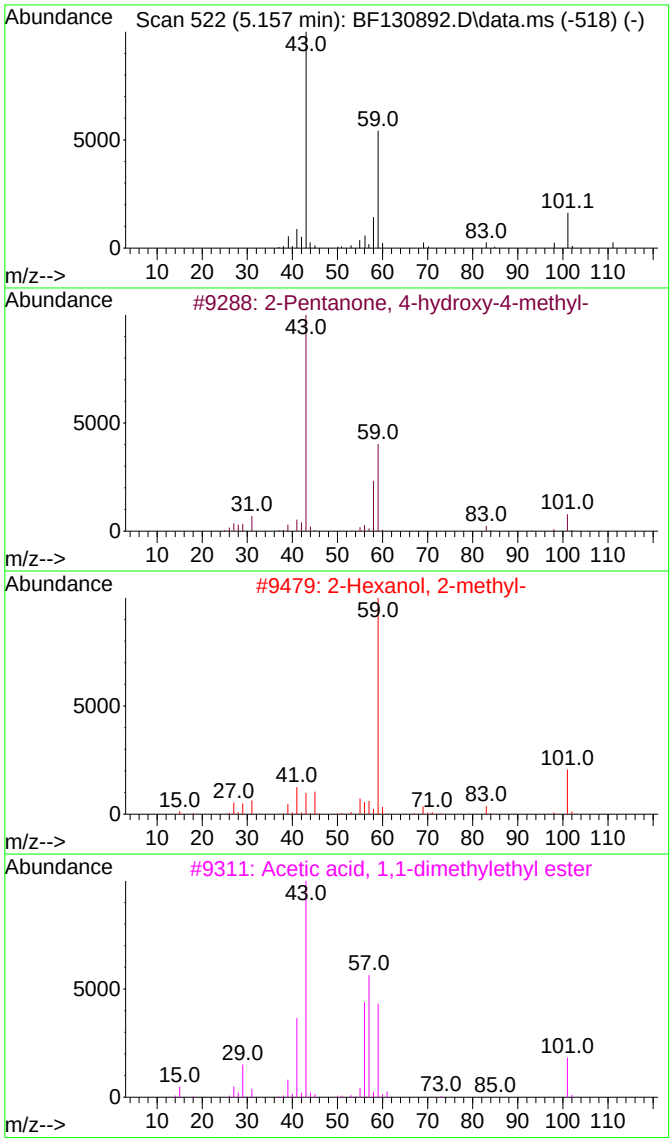
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF102722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

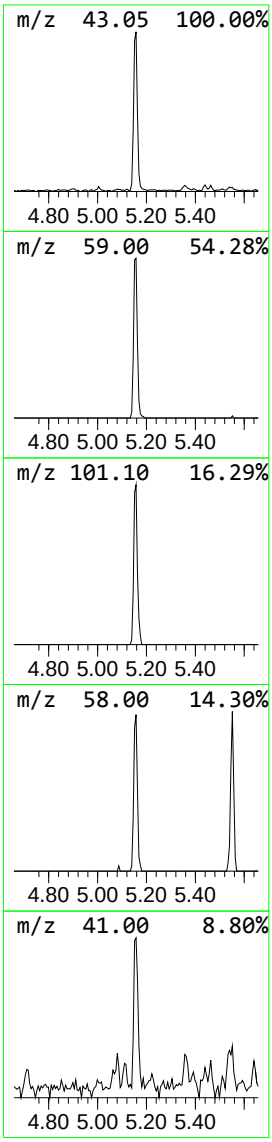
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.157	4.02 ng	97756	1,4-Dichlorobenzene-d4	6.945

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	40
3	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
4	Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	36
5	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25





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Misc :
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Instrument :
BNA_F
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF102722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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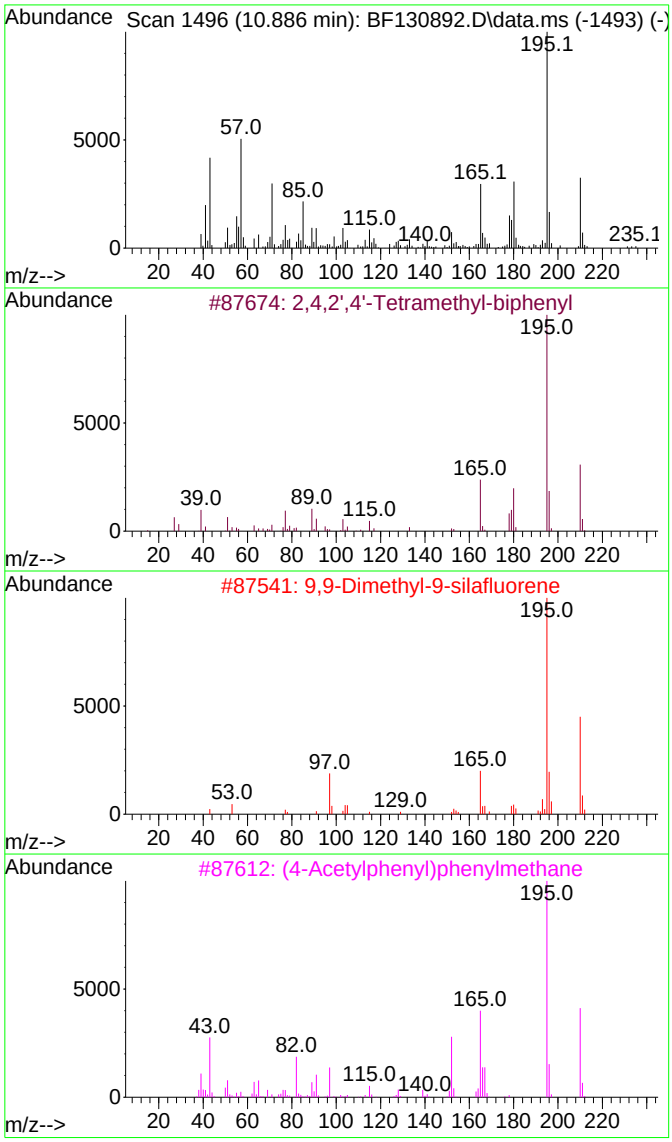
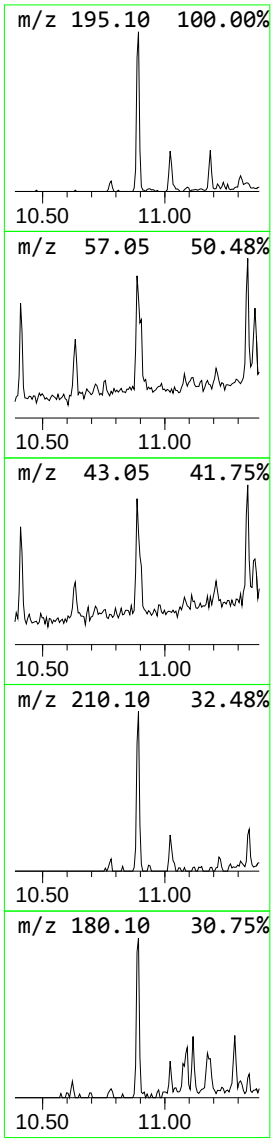
Peak Number 3 2,4,2',4'-Tetramethyl-biphenyl Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.886	4.92 ng	149325	Phenanthrene-d10	11.486

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,2',4'-Tetramethyl-biphenyl	210	C16H18	1000486-97-7	93
2	9,9-Dimethyl-9-silafluorene	210	C14H14Si	013688-68-1	83
3	(4-Acetylphenyl)phenylmethane	210	C15H14O	000782-92-3	58
4	Benzoic acid, 4-hydroxy-, trimet...	210	C10H14O3Si	1000395-53-3	55
5	Carbazole, 4,5-dimethyl-	195	C14H13N	018992-66-0	53





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF103122\
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ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF102722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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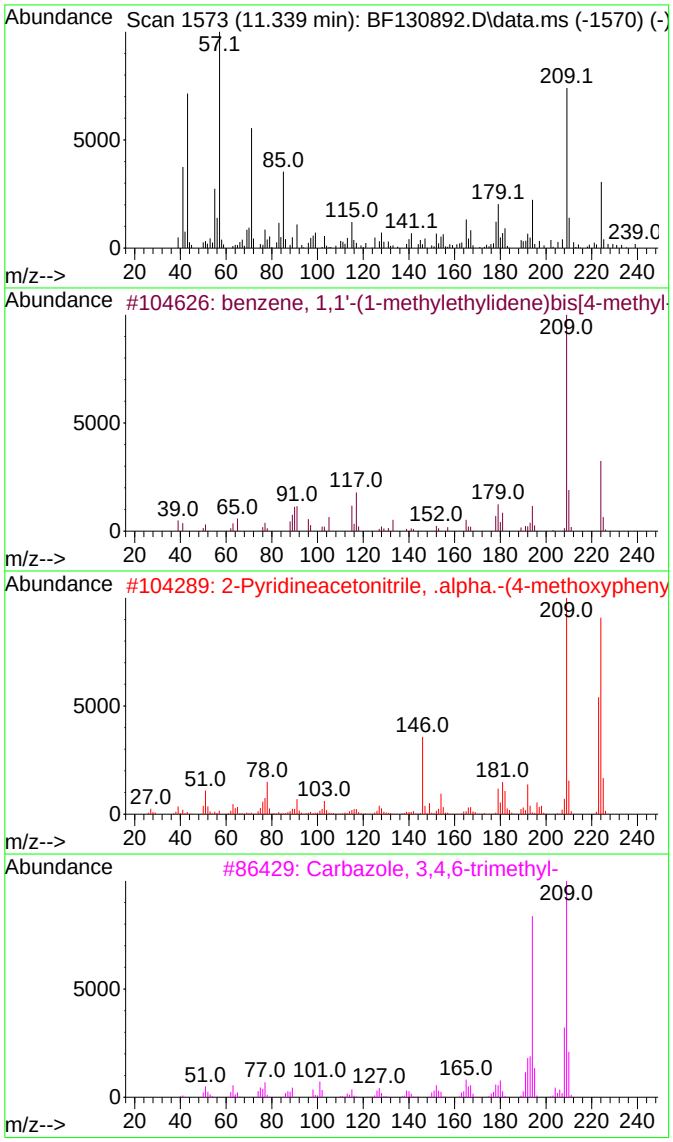
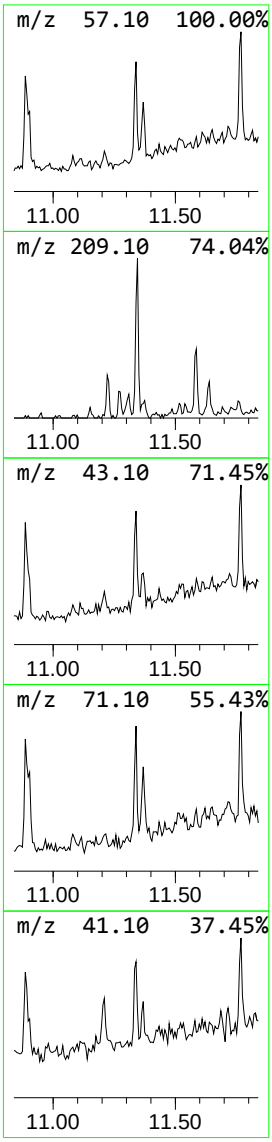
Peak Number 4 unknown11.339 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.339	3.32 ng	100738	Phenanthrene-d10	11.486

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	benzene, 1,1'-(1-methylethyliden...	224	C17H20	1000400-89-0	46
2	2-Pyridineacetonitrile, .alpha.-...	224	C14H12N2O	038236-51-0	46
3	Carbazole, 3,4,6-trimethyl-	209	C15H15N	078787-90-3	43
4	1,4,7-Trimethyl-2-azafluorene	209	C15H15N	062736-78-1	38
5	1,5,6,7-Tetramethyl-3-phenylbicy...	224	C17H20	126584-00-7	38





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Misc :
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BNA_F
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF102722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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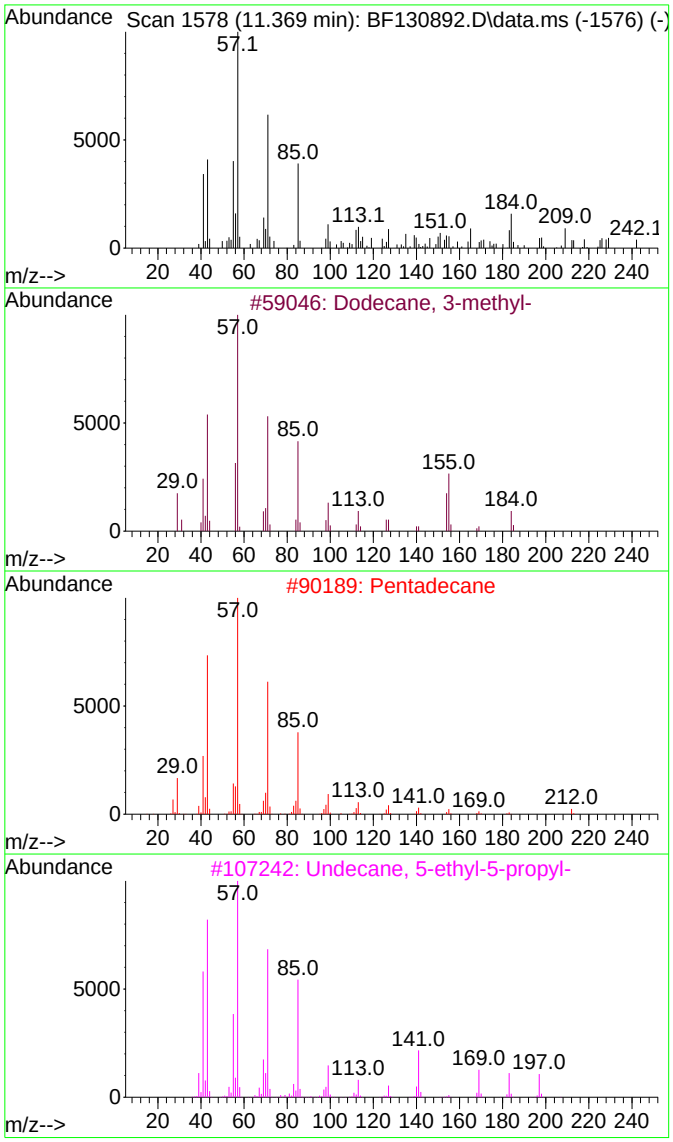
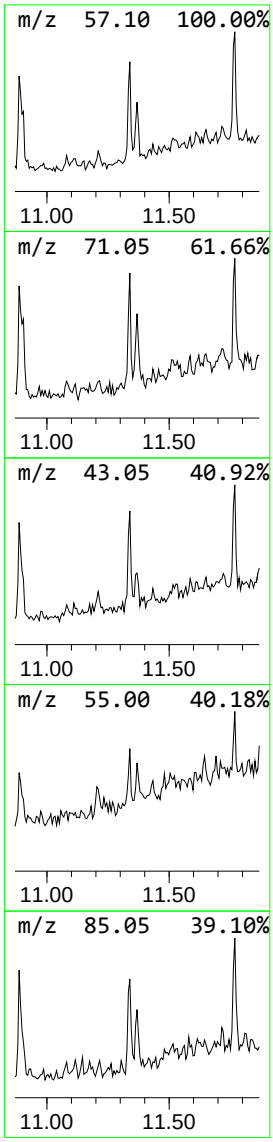
Peak Number 5 Dodecane, 3-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.369	2.06 ng	62582	Phenanthrene-d10	11.486

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane, 3-methyl-	184	C13H28	017312-57-1	76
2	Pentadecane	212	C15H32	000629-62-9	70
3	Undecane, 5-ethyl-5-propyl-	226	C16H34	002755-07-9	59
4	Dodecane	170	C12H26	000112-40-3	58
5	Tetradecane	198	C14H30	000629-59-4	58





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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF102722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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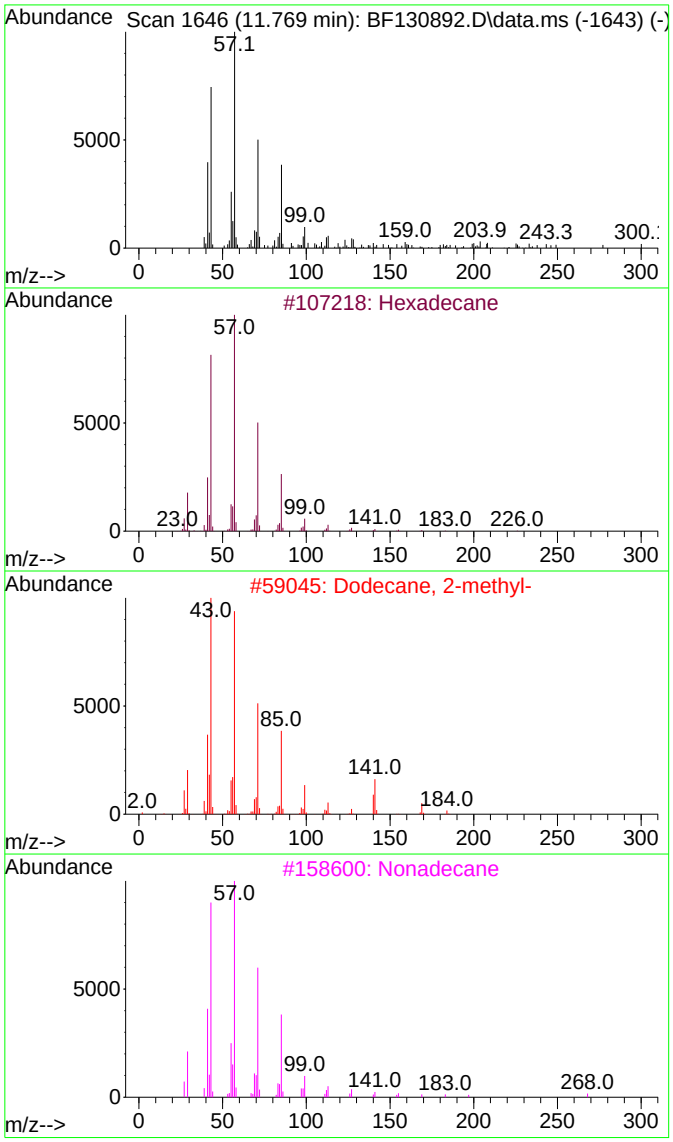
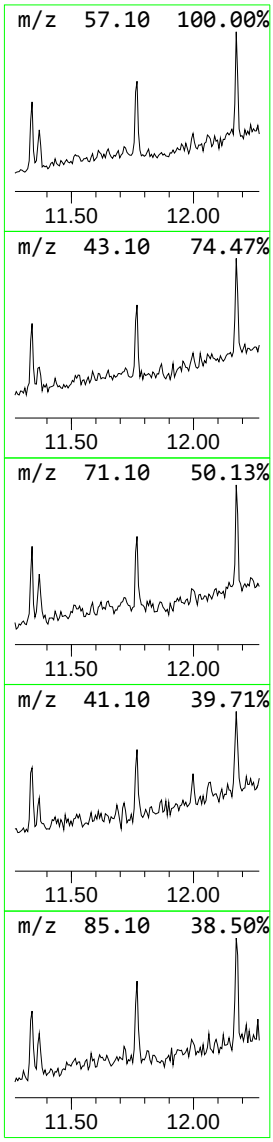
Peak Number 6 Hexadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.769	2.12 ng	64304	Phenanthrene-d10	11.486

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	95
2	Dodecane, 2-methyl-	184	C13H28	001560-97-0	87
3	Nonadecane	268	C19H40	000629-92-5	83
4	Dodecane, 1-iodo-	296	C12H25I	004292-19-7	80
5	Tetradecane	198	C14H30	000629-59-4	80





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF103122\
Data File : BF130892.D
Acq On : 31 Oct 2022 22:03
Operator : CG\JU
Sample : N5374-01 2X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
OK-2-102822

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF102722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

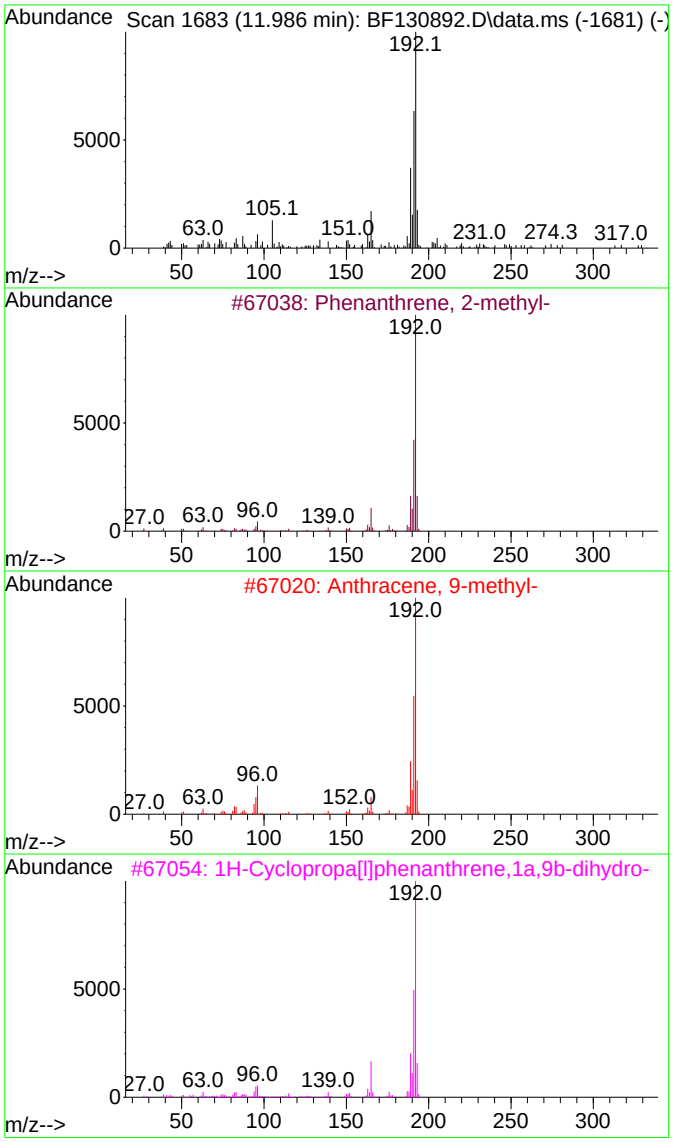
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

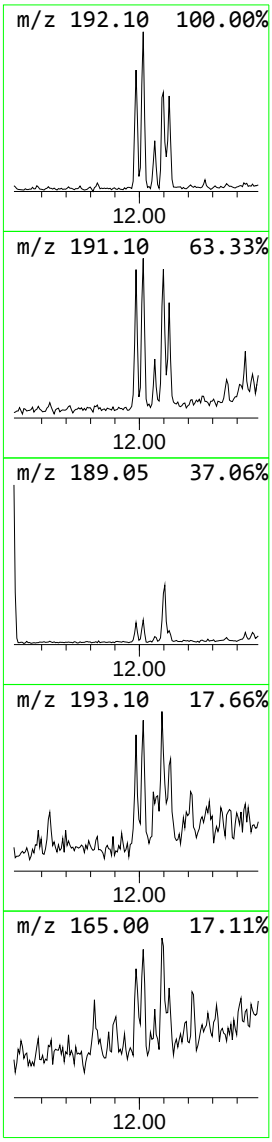
****

Peak Number 7 Phenanthrene, 2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.986	2.06 ng	62605	Phenanthrene-d10	11.486

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	94
2	Anthracene, 9-methyl-	192	C15H12	000779-02-2	90
3	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	90
4	Anthracene, 2-methyl-	192	C15H12	000613-12-7	90
5	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	90





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Acq On : 31 Oct 2022 22:03
Operator : CG\JU
Sample : N5374-01 2X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
OK-2-102822

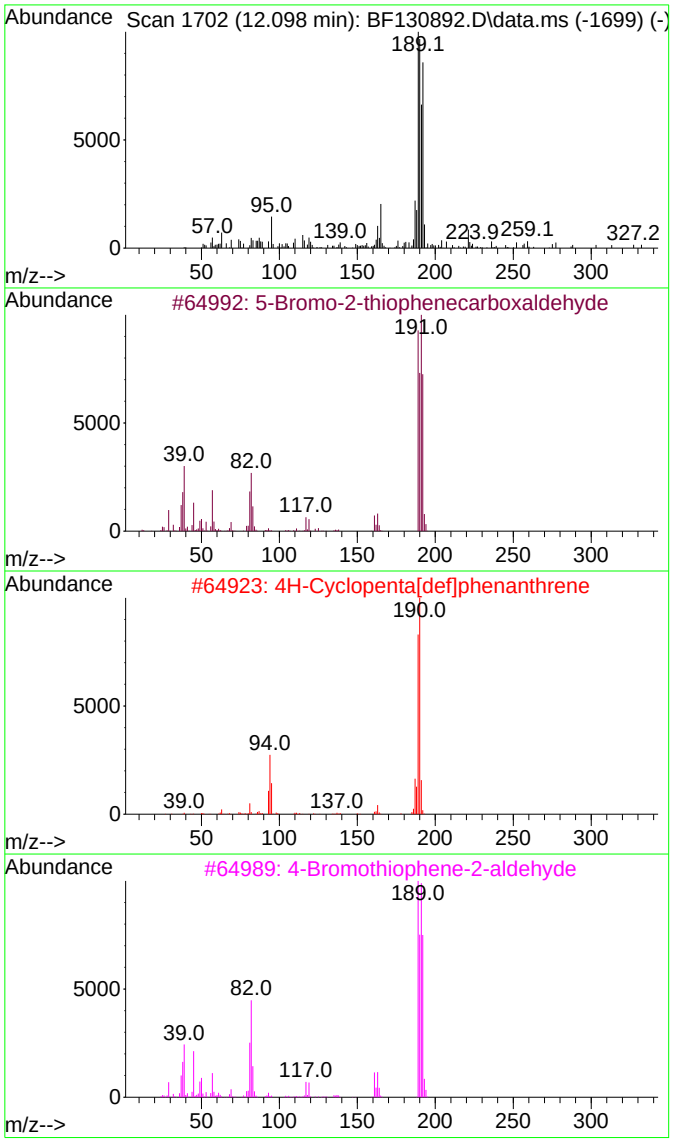
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF102722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

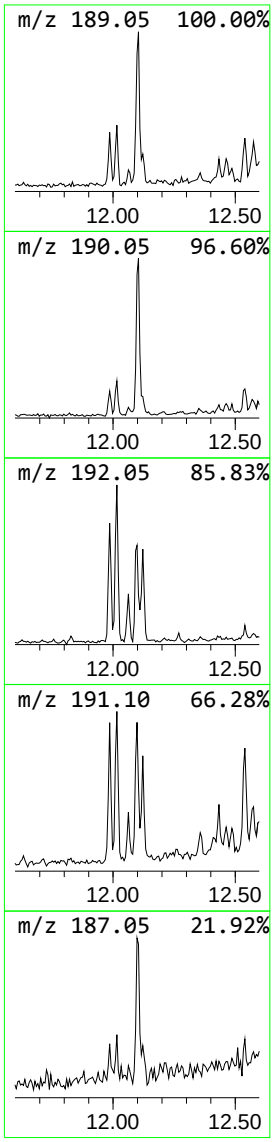
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 5-Bromo-2-thiophenecarboxal... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.098	2.64 ng	80046	Phenanthrene-d10	11.486

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Bromo-2-thiophenecarboxaldehyde	190	C5H3BrOS	004701-17-1	58
2	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	53
3	4-Bromothiophene-2-aldehyde	190	C5H3BrOS	018791-75-8	53
4	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	49
5	5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	46





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Sample : N5374-01 2X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
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ClientSampleId :
OK-2-102822

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF102722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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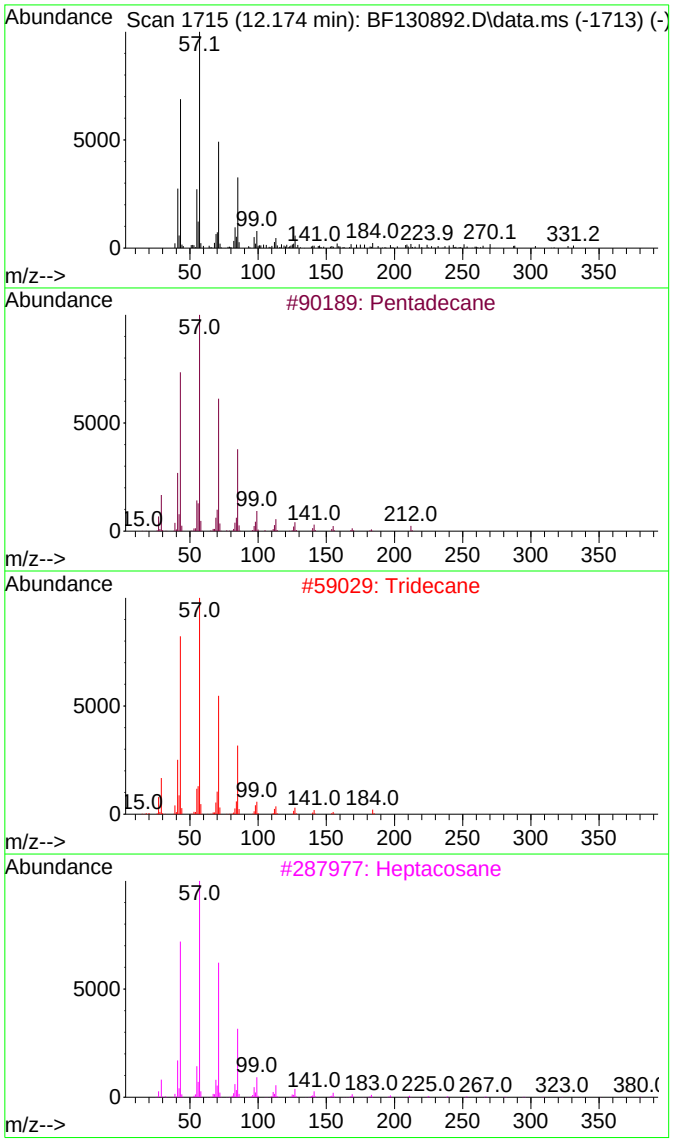
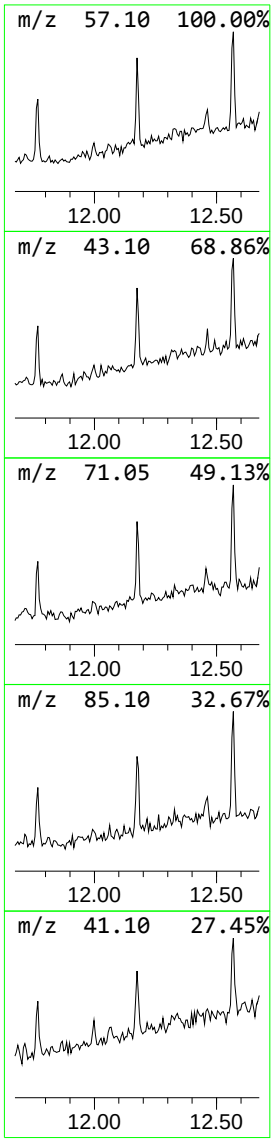
Peak Number 9 Pentadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.174	2.01 ng	61060	Phenanthrene-d10	11.486

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecane	212	C15H32	000629-62-9	91
2	Tridecane	184	C13H28	000629-50-5	81
3	Heptacosane	380	C27H56	000593-49-7	72
4	Eicosane	282	C20H42	000112-95-8	72
5	Tridecane, 2-methyl-	198	C14H30	001560-96-9	58





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Acq On : 31 Oct 2022 22:03
Operator : CG\JU
Sample : N5374-01 2X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
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ClientSampleId :
OK-2-102822

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF102722.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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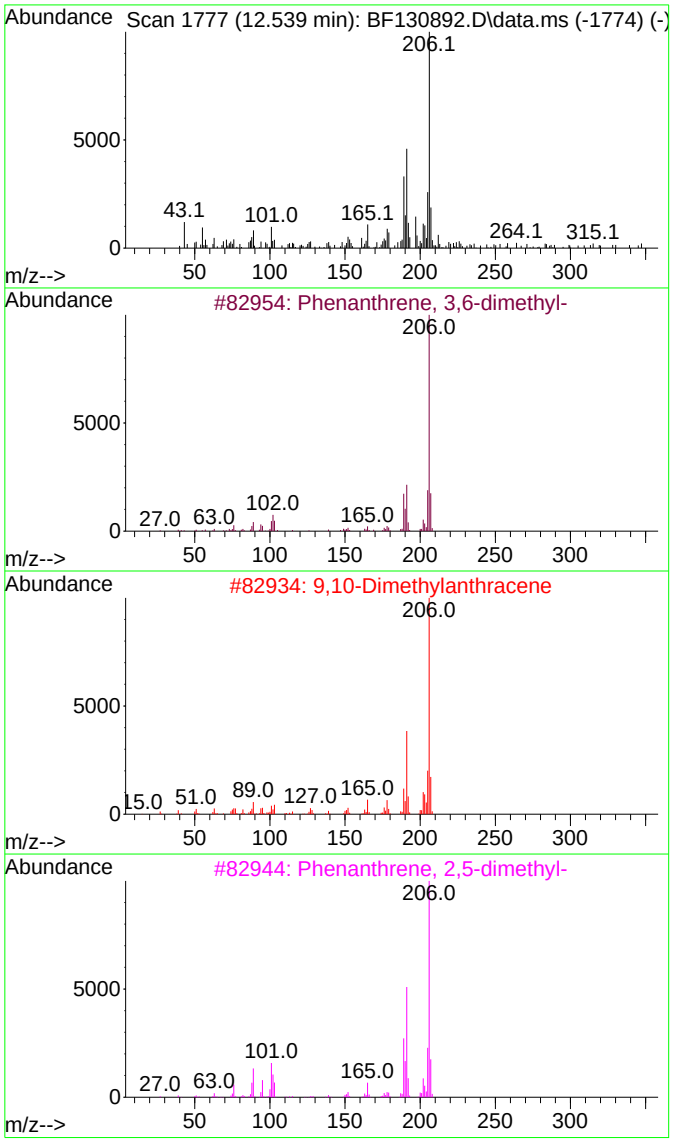
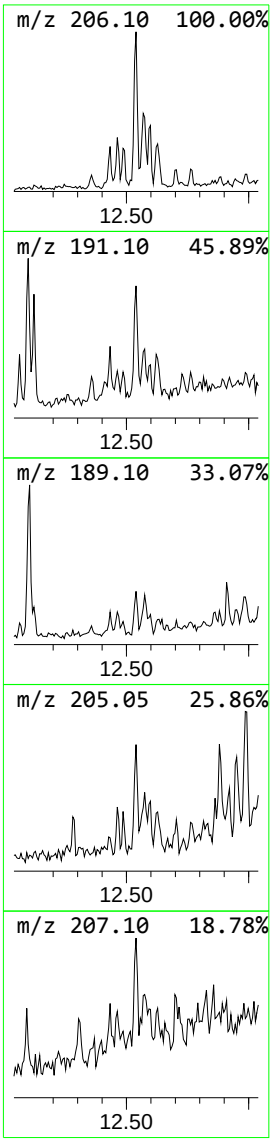
Peak Number 10 Phenanthrene, 3,6-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.539	2.24 ng	67853	Phenanthrene-d10	11.486

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	94
2	9,10-Dimethylanthracene	206	C16H14	000781-43-1	93
3	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	93
4	Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	93
5	di-p-Tolylacetylene	206	C16H14	002789-88-0	93





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Data File : BF130892.D
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Operator : CG\JU
Sample : N5374-01 2X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
OK-2-102822

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF102722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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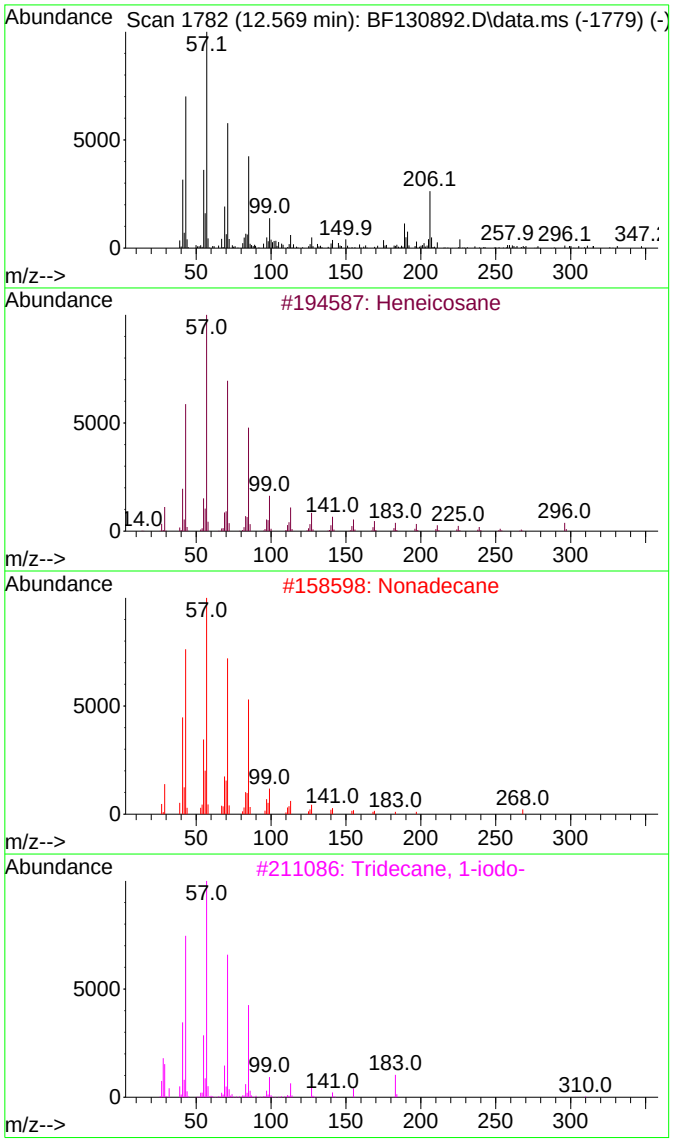
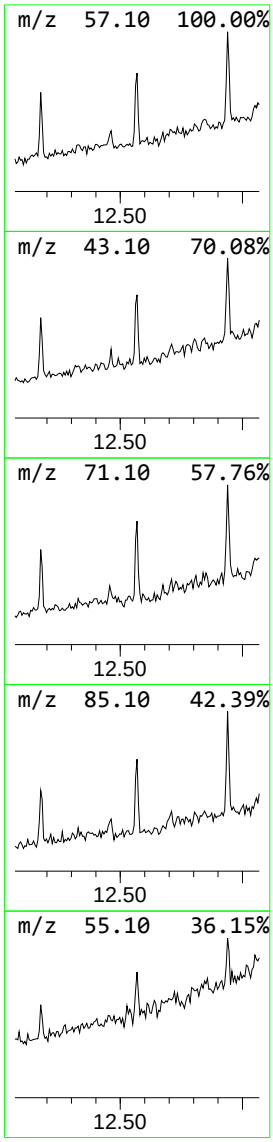
Peak Number 11 Heneicosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.569	3.67 ng	111344	Phenanthrene-d10	11.486

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	76
2	Nonadecane	268	C19H40	000629-92-5	70
3	Tridecane, 1-iodo-	310	C13H27I	035599-77-0	62
4	Heptacosane	380	C27H56	000593-49-7	62
5	Tetracosane	338	C24H50	000646-31-1	58





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Operator : CG\JU
Sample : N5374-01 2X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
OK-2-102822

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF102722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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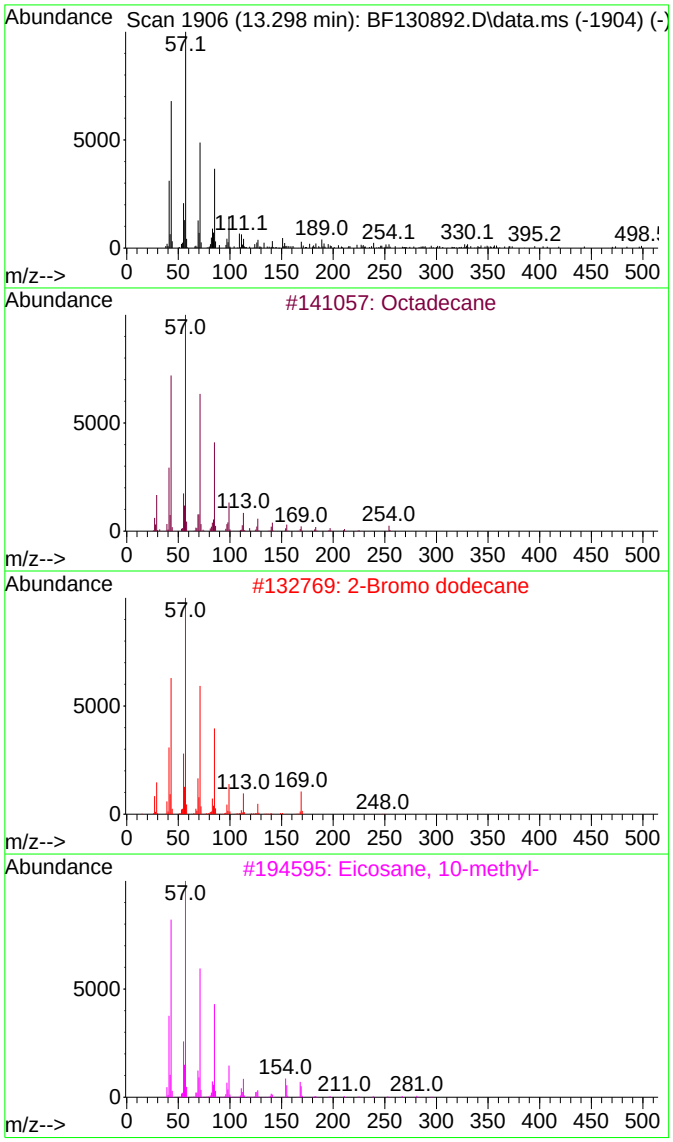
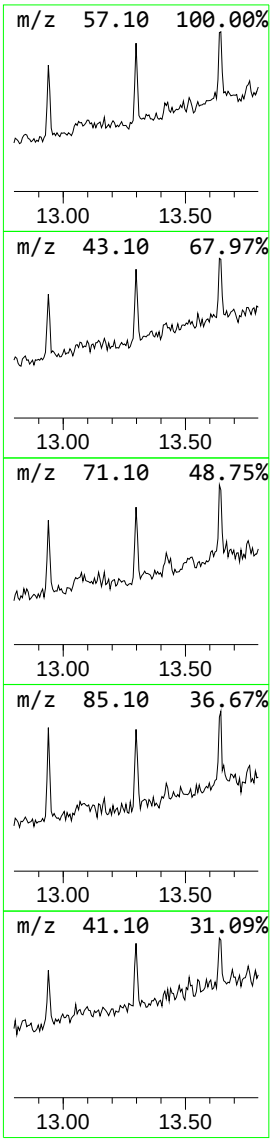
Peak Number 12 Octadecane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.298	5.16 ng	105028	Chrysene-d12	14.133

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane	254	C18H38	000593-45-3	91
2	2-Bromo dodecane	248	C12H25Br	013187-99-0	87
3	Eicosane, 10-methyl-	296	C21H44	054833-23-7	87
4	Octacosane	394	C28H58	000630-02-4	83
5	Eicosane	282	C20H42	000112-95-8	83





Tentatively Identified Compound (LSC) summary

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Misc :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF102722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Butane, 2-metho...	2.204	9.6	ng	233970	1	6.945	485986	20.0
2-Pentanone, 4-...	5.157	4.0	ng	97756	1	6.945	485986	20.0
2,4,2',4'-Tetra...	10.886	4.9	ng	149325	4	11.486	606842	20.0
unknown11.339	11.339	3.3	ng	100738	4	11.486	606842	20.0
Dodecane, 3-met...	11.369	2.1	ng	62582	4	11.486	606842	20.0
Hexadecane	11.769	2.1	ng	64304	4	11.486	606842	20.0
Phenanthrene, 2...	11.986	2.1	ng	62605	4	11.486	606842	20.0
5-Bromo-2-thiop...	12.098	2.6	ng	80046	4	11.486	606842	20.0
Pentadecane	12.174	2.0	ng	61060	4	11.486	606842	20.0
Phenanthrene, 3...	12.539	2.2	ng	67853	4	11.486	606842	20.0
Heneicosane	12.569	3.7	ng	111344	4	11.486	606842	20.0
Octadecane	13.298	5.2	ng	105028	5	14.133	406949	20.0