LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122018\
Data File : BF111651.D
Acq On : 20 Dec 2018 21:00
Operator : JU/SJ
Sample : J6346-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : DW-5

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-BF121918.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	2.234	21	25	34	rVB	59791	90153	1.83%	0.160%
2	5.128	511	517	523	rVB	181903	241001	4.90%	0.428%
3	5.522	579	584	587	rBV	4491211	4308102	87.59%	7.652%
4	6.522	748	754	759	rBV2	4376889	4918578	100.00%	8.736%
5	6.669	774	779	782	rBV	4646351	4514970	91.79%	8.019%
6	6.898	814	818	821	rBV	1348026	1221188	24.83%	2.169%
7	7.051	840	844	848	rVB	3693152	3538704	71.95%	6.285%
8	7.451	907	912	915	rBV	3053444	2890648	58.77%	5.134%
9	8.175	1031	1035	1039	rBV	1803719	1554174	31.60%	2.760%
10	8.781	1133	1138	1146	rBV3	85319	120185	2.44%	0.213%
11	8.886	1150	1156	1160	rVB	56619	55311	1.12%	0.098%
12	8.986	1170	1173	1178	rBV2	77465	69103	1.40%	0.123%
13	9.251	1213	1218	1221	rBV	5481348	4759307	96.76%	8.453%
14	9.351	1228	1235	1239	rBV2	72328	97285	1.98%	0.173%
15	9.586	1270	1275	1278	rBV	65375	93440	1.90%	0.166%
16	9.639	1281	1284	1287	rBV	537514	405414	8.24%	0.720%
17	9.933	1330	1334	1337	rBV	1666299	1507804	30.66%	2.678%
18	9.963	1337	1339	1342	rVB	454220	332141	6.75%	0.590%
19	10.133	1364	1368	1371	rBV	214286	183829	3.74%	0.327%
20	10.475	1423	1426	1432	rBV	283714	249255	5.07%	0.443%
21	10.586	1443	1445	1448	rVV	97566	87652	1.78%	0.156%
22	10.633	1450	1453	1457	rVB2	105591	103542	2.11%	0.184%
23	10.716	1463	1467	1470	rBV	2915442	2433149	49.47%	4.322%
24	10.851	1485	1490	1495	rBV3	418809	502857	10.22%	0.893%
25	10.898	1495	1498	1501	rBV	386175	342454	6.96%	0.608%
26	10.933	1501	1504	1507	rVV2	346804	398918	8.11%	0.709%
27	10.969	1507	1510	1512	rVV	346046	331583	6.74%	0.589%
28	10.992	1512	1514	1516	rVV	252971	200180	4.07%	0.356%
29	11.033	1516	1521	1522	rVV2	177089	248468	5.05%	0.441%
30	11.080	1526	1529	1532	rVV3	296282	346271	7.04%	0.615%
31	11.116	1532	1535	1537	rVV	380816	321438	6.54%	0.571%
32	11.157	1540	1542	1547	rVB2	235578	197656	4.02%	0.351%
33	11.316	1565	1569	1573	rVB3	269131	343523	6.98%	0.610%
34	11.416	1582	1586	1588	rBV	1337134	1305609	26.54%	2.319%

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122018\
Data File : BF111651.D
Acq On : 20 Dec 2018 21:00
Operator : JU/SJ
Sample : J6346-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : DW-5

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-BF121918.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	11.439	1588	1590	1593	rVV	1258532	918149	18.67%	1.631%
36	11.486	1593	1598	1601	rVB	354811	312525	6.35%	0.555%
37	11.522	1601	1604	1608	rBV	120171	140020	2.85%	0.249%
38	11.633	1620	1623	1626	rBV	227033	218371	4.44%	0.388%
39	11.669	1626	1629	1632	rVB2	88063	79583	1.62%	0.141%
40	11.939	1672	1675	1680	rVB	709114	654943	13.32%	1.163%
41	12.027	1687	1690	1692	rBV2	298324	274025	5.57%	0.487%
42	12.210	1718	1721	1722	rBV	111843	83160	1.69%	0.148%
43	12.463	1762	1764	1767	rVB2	106324	96037	1.95%	0.171%
44	12.627	1788	1792	1795	rBV	2264235	1978172	40.22%	3.514%
45	12.668	1795	1799	1805	rVB2	368687	380332	7.73%	0.676%
46	12.727	1806	1809	1811	rBV	371966	334497	6.80%	0.594%
47	12.857	1825	1831	1836	rBV	1558432	1789746	36.39%	3.179%
48	12.968	1848	1850	1851	rBV2	98042	71163	1.45%	0.126%
49	12.998	1851	1855	1858	rVB	3599006	3184968	64.75%	5.657%
50	13.180	1884	1886	1889	rVB	266430	219899	4.47%	0.391%
51	13.245	1895	1897	1900	rVB	284593	212915	4.33%	0.378%
52	13.510	1940	1942	1945	rBV	187433	165594	3.37%	0.294%
53	13.892	2004	2007	2014	rVB2	635382	652614	13.27%	1.159%
54	14.039	2028	2032	2036	rBV3	1688550	2737518	55.66%	4.862%
55	14.074	2036	2038	2041	rVB	852618	652797	13.27%	1.159%
56	15.092	2208	2211	2214	rBV	601911	863999	17.57%	1.535%
57	15.398	2260	2263	2269	rVV	402669	516363	10.50%	0.917%
58	15.462	2271	2274	2277	rVB	412794	447682	9.10%	0.795%
59	15.527	2282	2285	2288	rVB	697579	779771	15.85%	1.385%
60	17.451	2608	2612	2618	rVB2	130391	223107	4.54%	0.396%

Sum of corrected areas: 56301842

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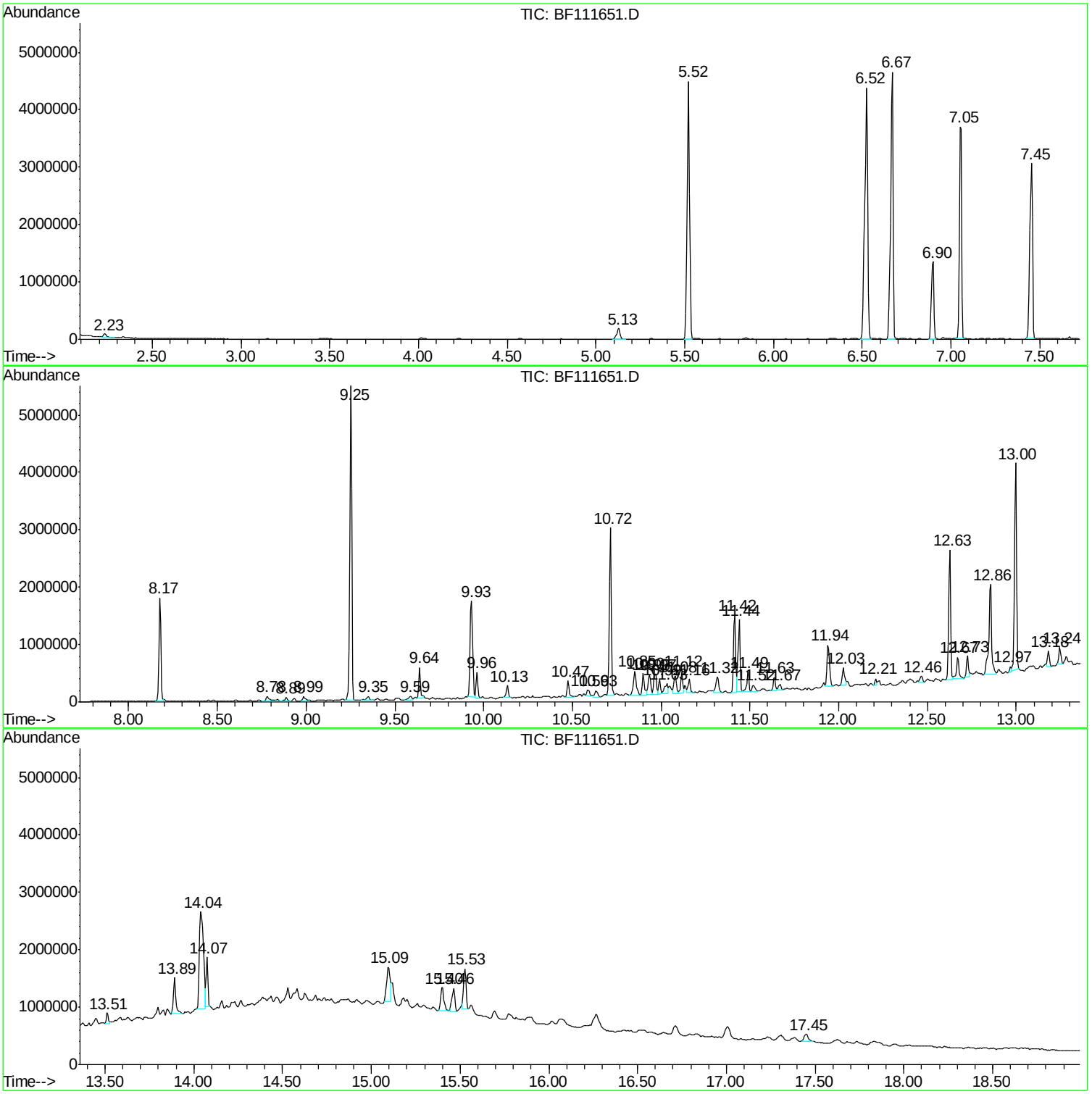
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : DW-5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF121918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122018\
Data File : BF111651.D
Acq On : 20 Dec 2018 21:00
Operator : JU/SJ
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Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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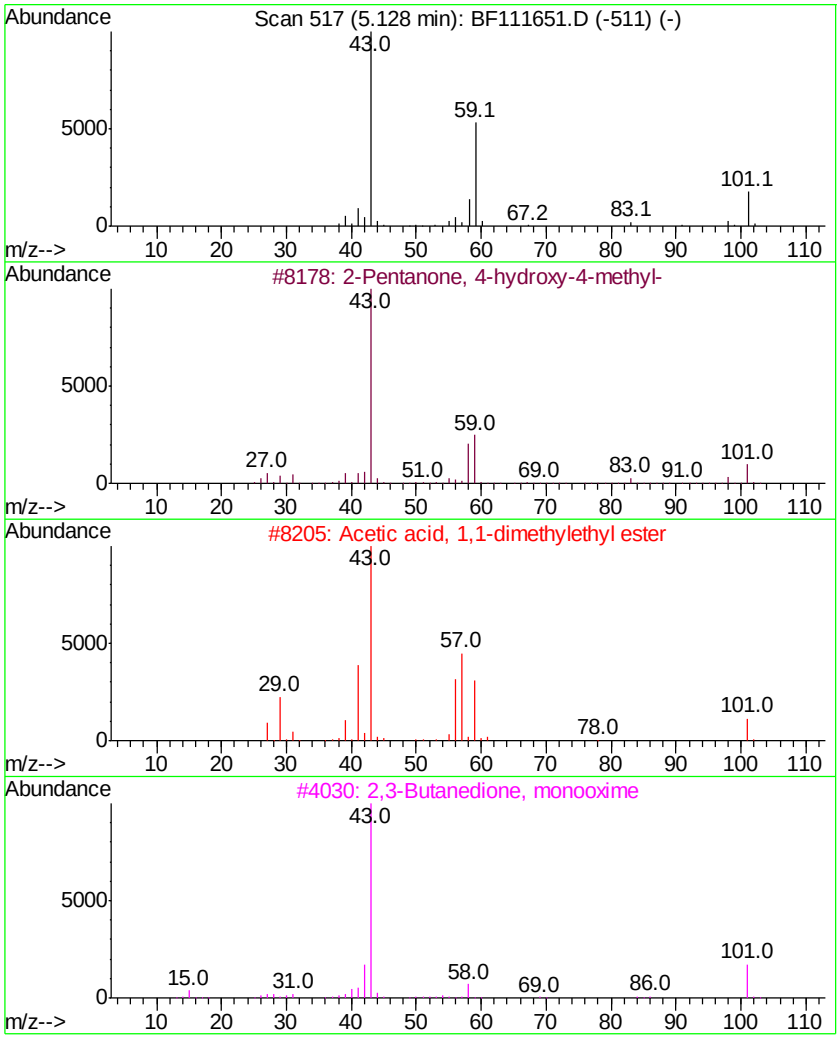
TIC Integration Parameters: LSCINT.P

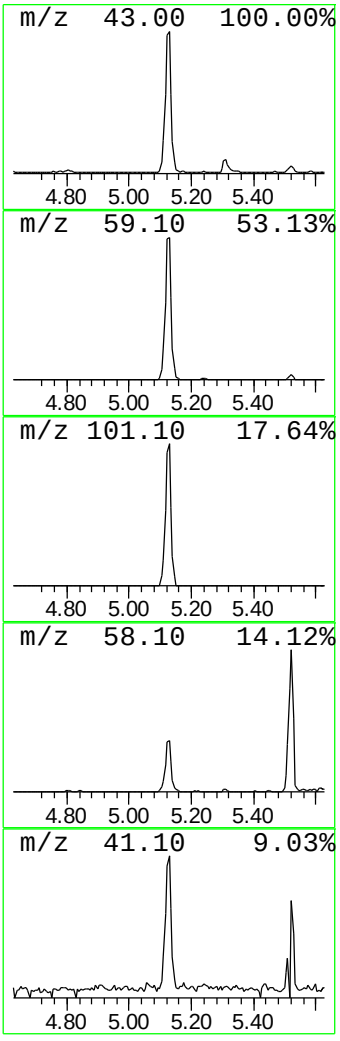
****

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.13	3.95 ng	241001	1,4-Dichlorobenzene-d4	6.90

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	28
3	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	27
4	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
5	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122018\
Data File : BF111651.D
Acq On : 20 Dec 2018 21:00
Operator : JU/SJ
Sample : J6346-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : DW-5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF121918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

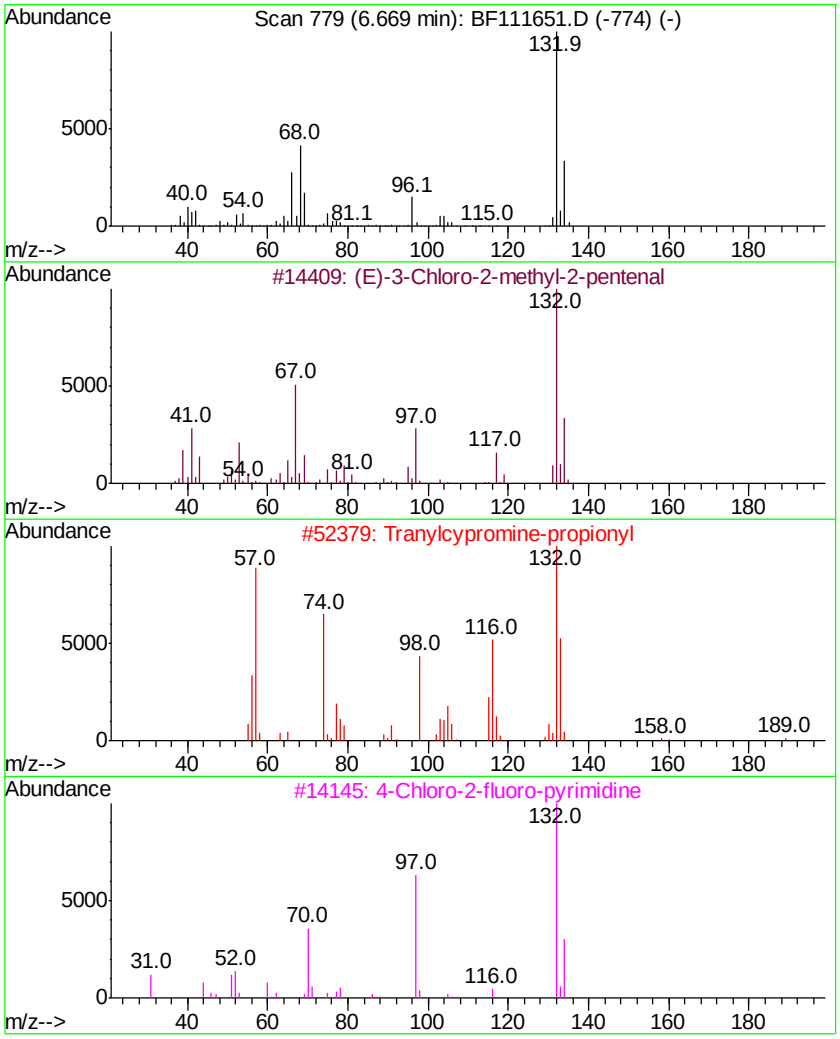
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

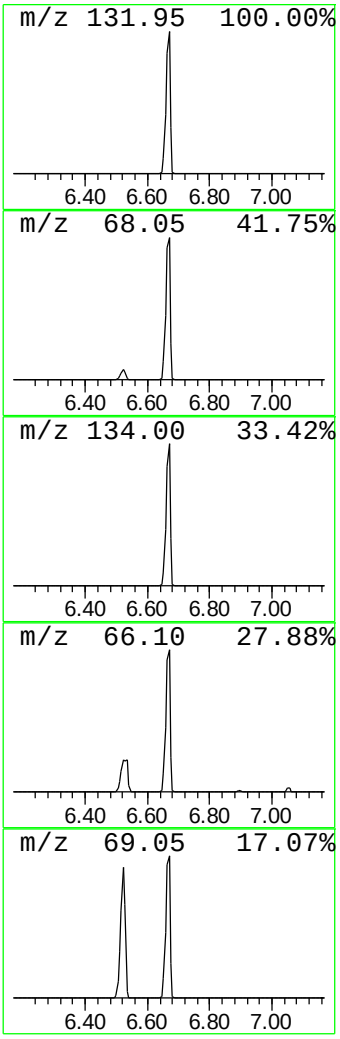
****

Peak Number 2 unknown6.67 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.67	73.94 ng	4514970	1,4-Dichlorobenzene-d4	6.90

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
2	Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
3	4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
4	Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5	Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122018\
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Operator : JU/SJ
Sample : J6346-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : DW-5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF121918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

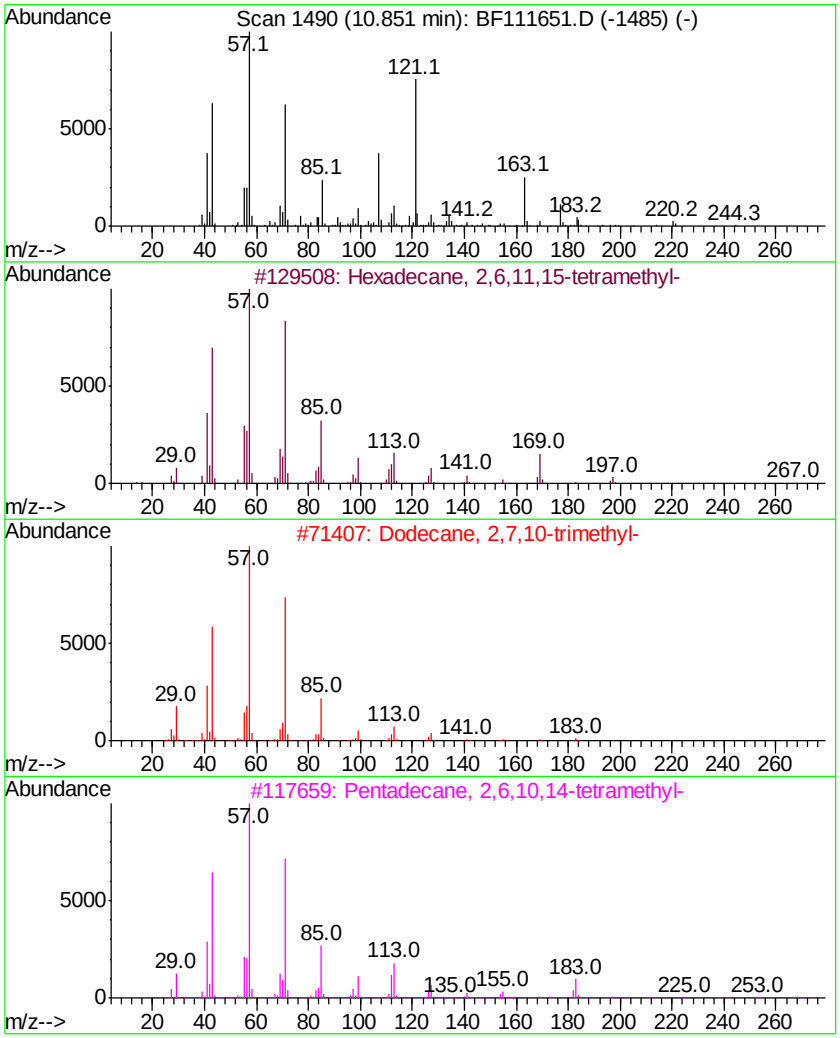
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

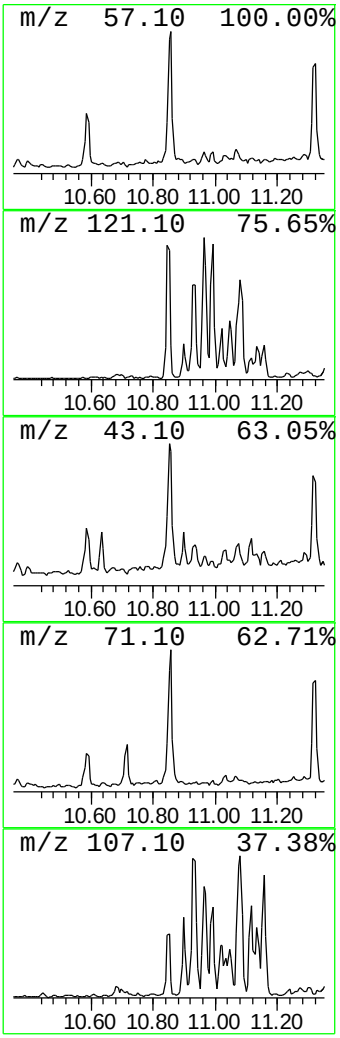
****

Peak Number 5 unknown10.85 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.85	7.70 ng	502857	Phenanthrene-d10	11.42

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane, 2,6,11,15-tetramethyl-	282	C20H42	000504-44-9	42
2	Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	38
3	Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	38
4	Tridecane, 5-propyl-	226	C16H34	055045-11-9	30
5	Nonane, 3,7-dimethyl-	156	C11H24	017302-32-8	30





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122018\
Data File : BF111651.D
Acq On : 20 Dec 2018 21:00
Operator : JU/SJ
Sample : J6346-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : DW-5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF121918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

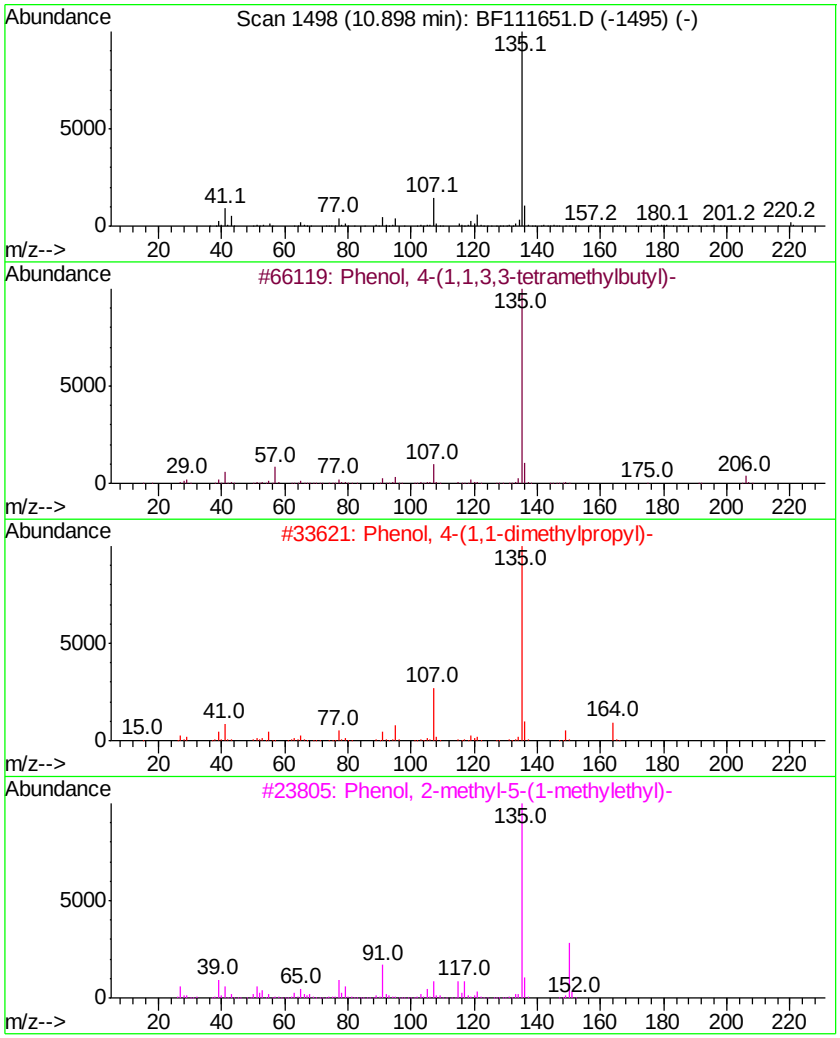
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

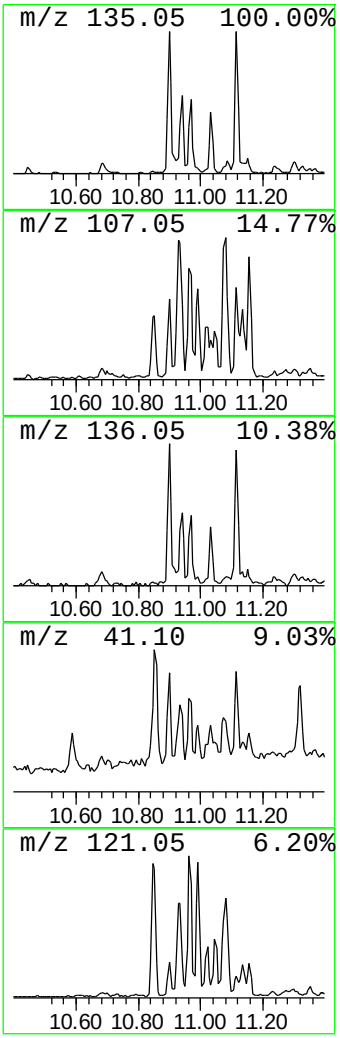
****

Peak Number 6 Phenol, 4-(1,1,3,3-tetramet... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.90	5.25 ng	342454	Phenanthrene-d10	11.42

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	78
2	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	78
3	Phenol, 2-methyl-5-(1-methylethyl)-	150	C10H14O	000499-75-2	72
4	m-Anisic acid, 4-nitrophenyl ester	273	C14H11NO5	1000307-80-4	64
5	1-n-Hexyladamantane	220	C16H28	022458-75-9	64





Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122018\
Data File : BF111651.D
Acq On : 20 Dec 2018 21:00
Operator : JU/SJ
Sample : J6346-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : DW-5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF121918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

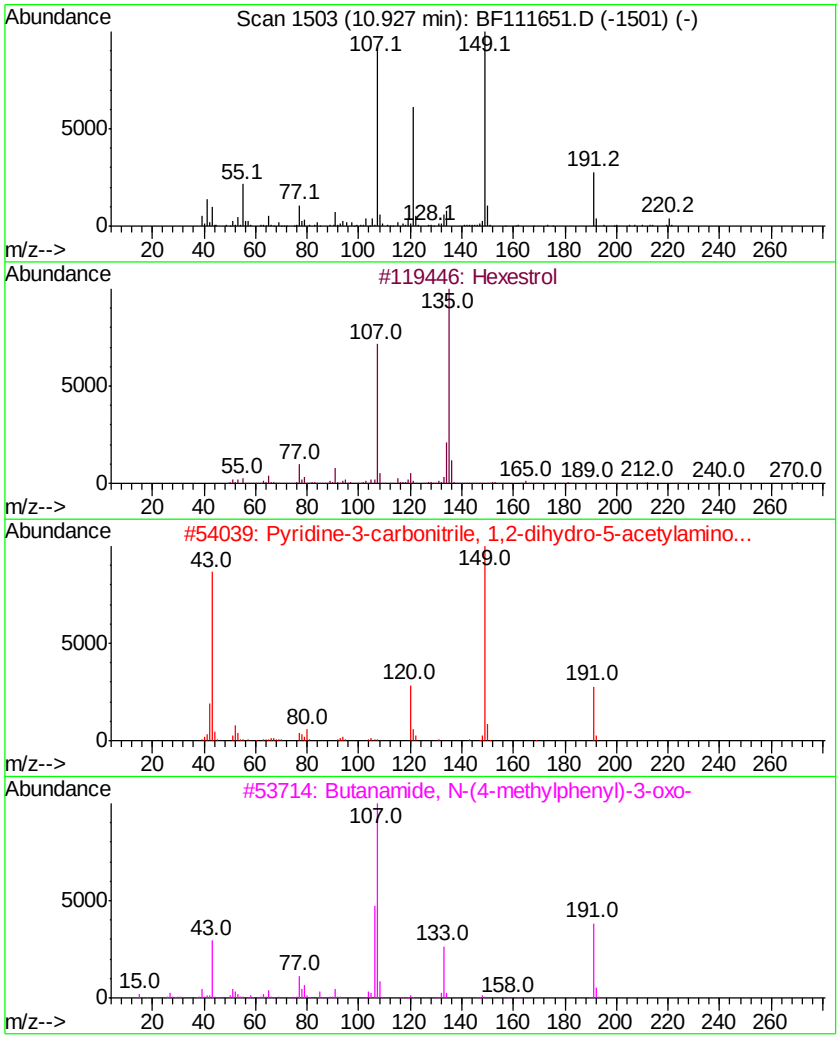
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

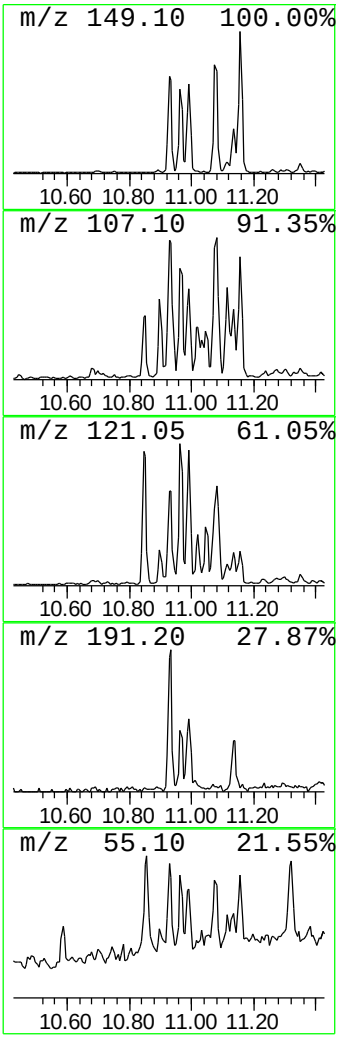
****

Peak Number 7 unknown10.93 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.93	6.11 ng	398918	Phenanthrene-d10	11.42

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexestrol	270	C18H22O2	000084-16-2	43
2	Pyridine-3-carbonitrile, 1,2-dih...	191	C9H9N3O2	220098-92-0	25
3	Butanamide, N-(4-methylphenyl)-3...	191	C11H13NO2	002415-85-2	25
4	Phenol, 2-propyl-	136	C9H12O	000644-35-9	18
5	Silane, trimethylphenyl-	150	C9H14Si	000768-32-1	18





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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF121918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

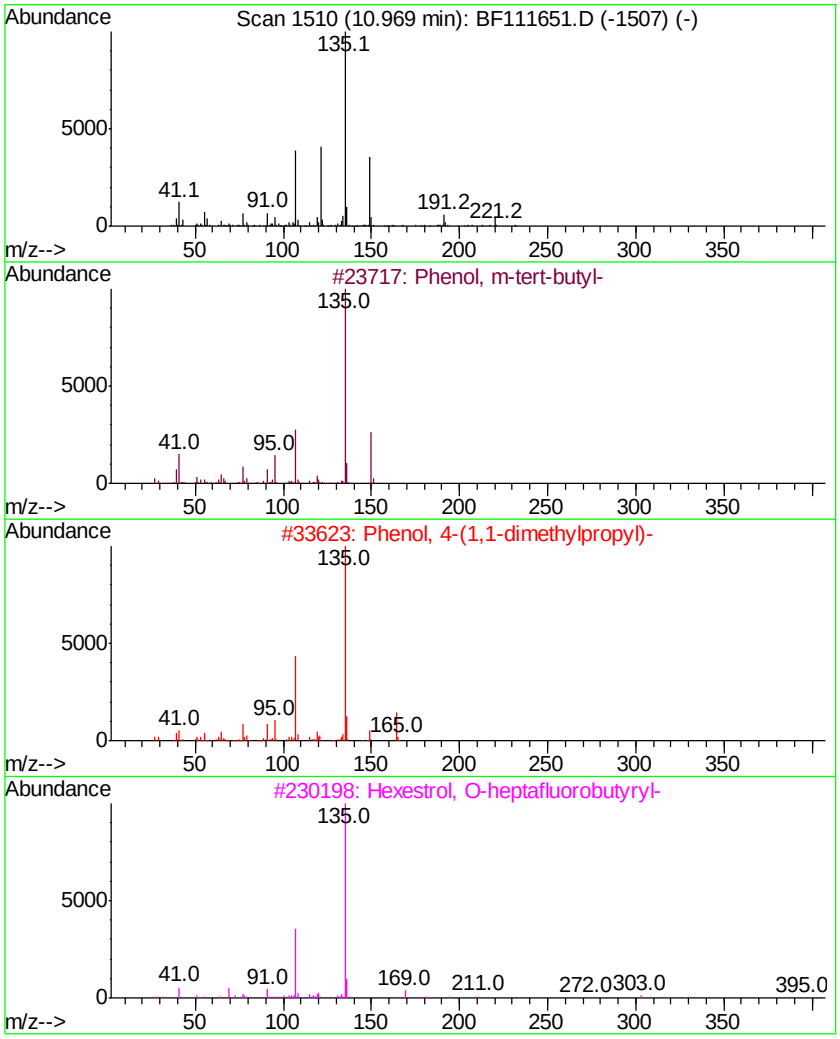
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

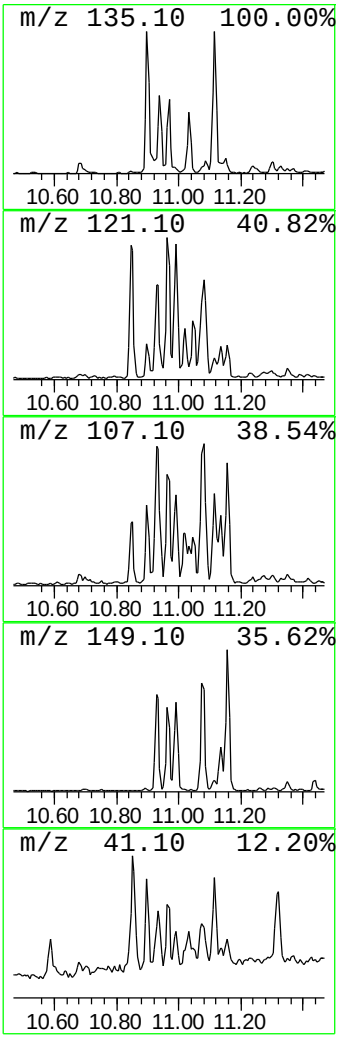
****

Peak Number 8 Phenol, m-tert-butyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.97	5.08 ng	331583	Phenanthrene-d10	11.42

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	64
2	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	59
3	Hexestrol, O-heptafluorobutyryl-	466	C22H21F7O3	1000365-31-9	53
4	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	53
5	Hexestrol, O-acetyl-	312	C20H24O3	1000374-87-8	53





Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122018\
Data File : BF111651.D
Acq On : 20 Dec 2018 21:00
Operator : JU/SJ
Sample : J6346-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : DW-5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF121918.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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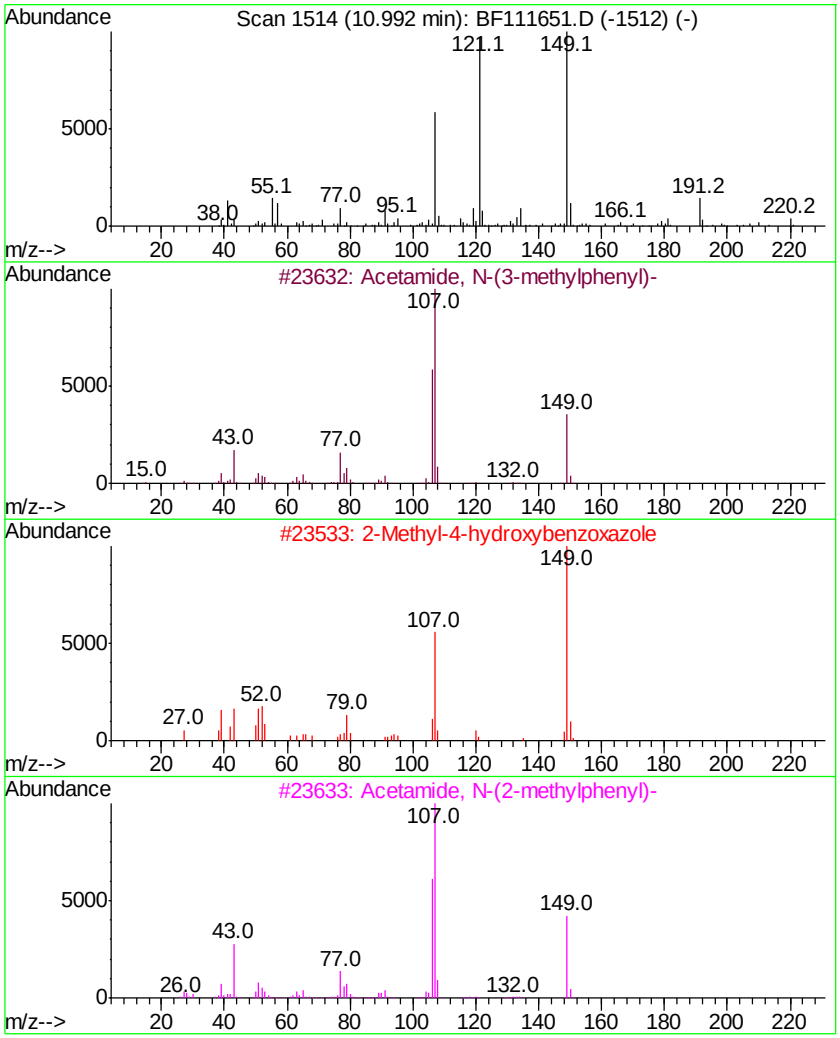
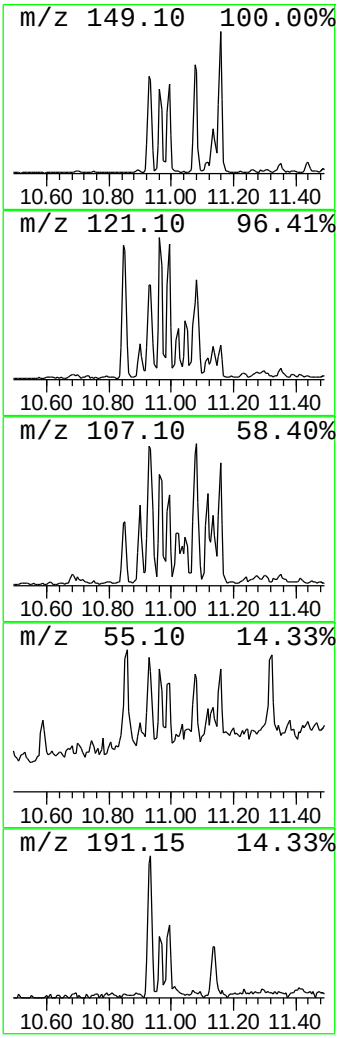
Peak Number 9 unknown10.99 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.99	3.07 ng	200180	Phenanthrene-d10	11.42

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	38
2	2-Methyl-4-hydroxybenzoxazole	149	C8H7NO2	051110-60-2	38
3	Acetamide, N-(2-methylphenyl)-	149	C9H11NO	000120-66-1	30
4	2-Methyl-2-(4'-methoxyphenyl)pen...	206	C13H18O2	185700-49-6	27
5	Phthalic acid, 5-methoxy-3-methy...	322	C18H26O5	1000315-37-5	27





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Data File : BF111651.D
Acq On : 20 Dec 2018 21:00
Operator : JU/SJ
Sample : J6346-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : DW-5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF121918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

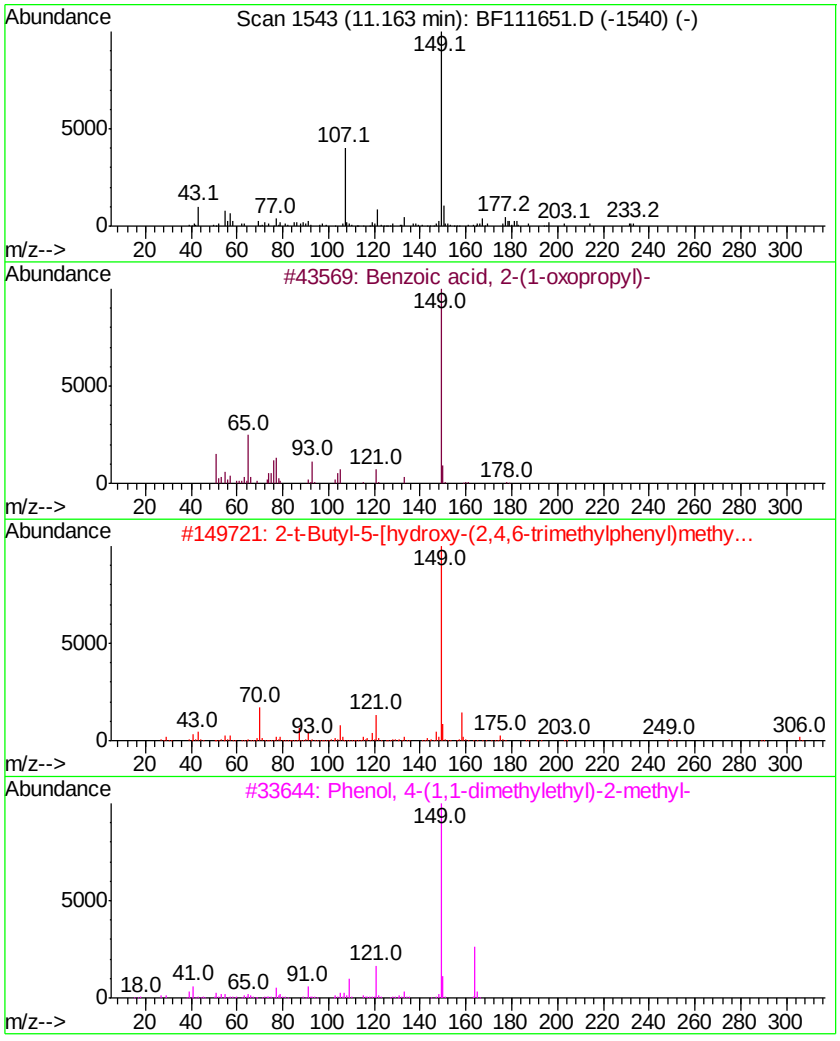
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

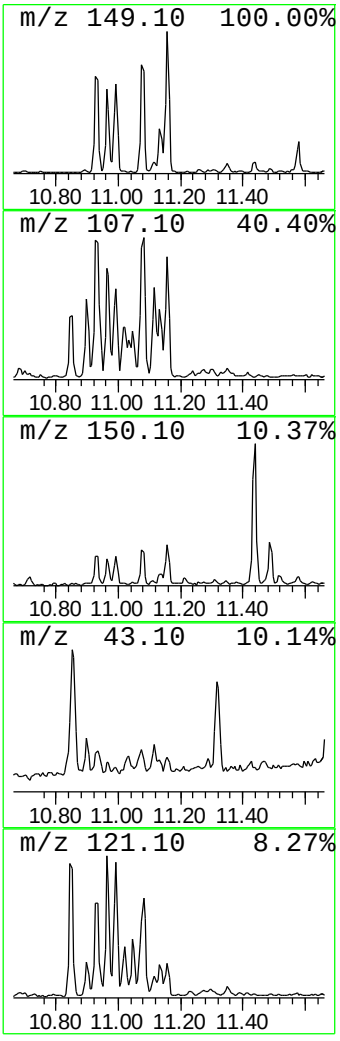
****

Peak Number 13 Benzoic acid, 2-(1-oxopropyl)- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.16	3.03 ng	197656	Phenanthrene-d10	11.42

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzoic acid, 2-(1-oxopropyl)-	178	C10H10O3	002360-45-4	50
2	2-t-Butyl-5-[hydroxy-(2,4,6-trim...	306	C18H26O4	092572-63-9	50
3	Phenol, 4-(1,1-dimethylethyl)-2-...	164	C11H16O	000098-27-1	47
4	Phthalic acid, dodecyl ethyl ester	362	C22H34O4	1000308-93-4	47
5	Phthalic acid, ethyl 3-methylbut...	264	C15H20O4	1000356-80-1	47





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122018\
Data File : BF111651.D
Acq On : 20 Dec 2018 21:00
Operator : JU/SJ
Sample : J6346-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : DW-5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF121918.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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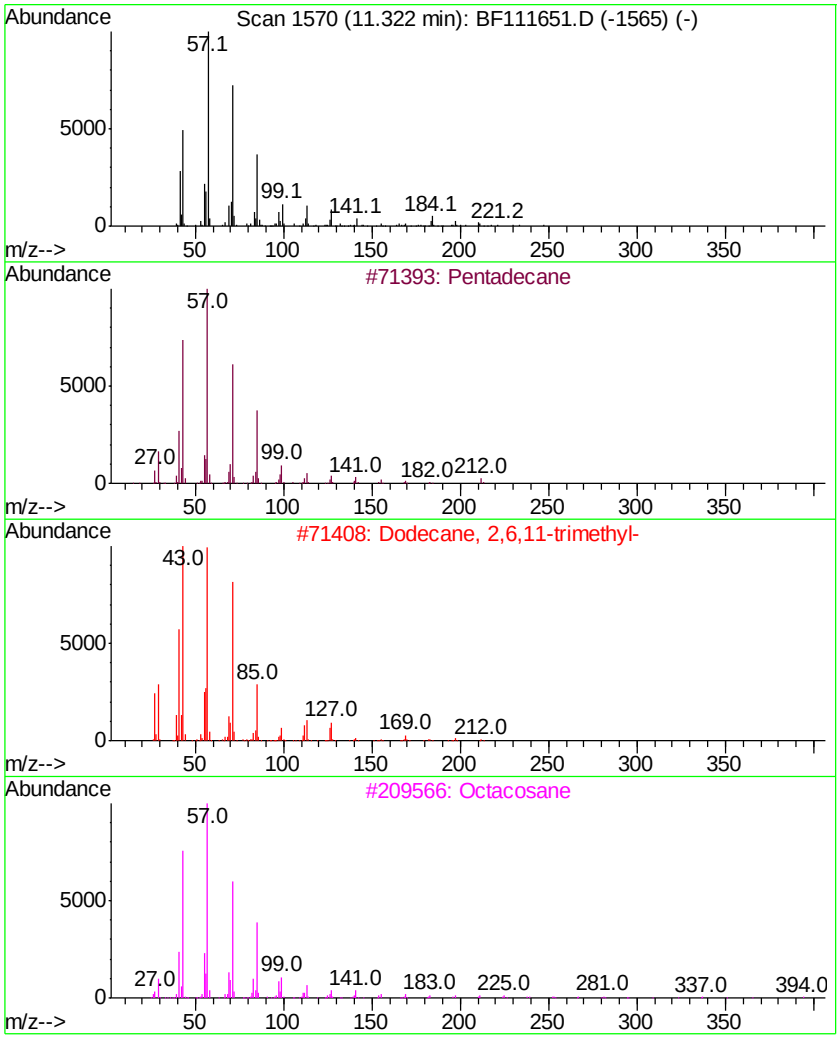
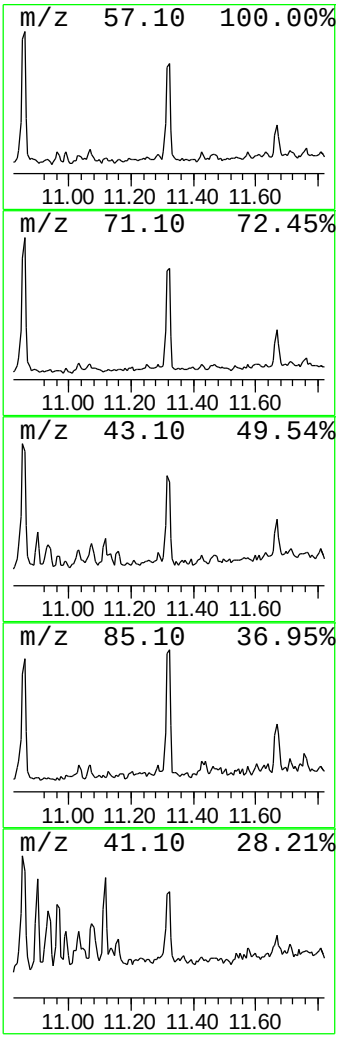
Peak Number 14 Pentadecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.32	5.26 ng	343523	Phenanthrene-d10	11.42

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecane	212	C15H32	000629-62-9	58
2	Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	53
3	Octacosane	394	C28H58	000630-02-4	52
4	Tetracosane	338	C24H50	000646-31-1	52
5	Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	52





Library Search Compound Report

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Operator : JU/SJ
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Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : DW-5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF121918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

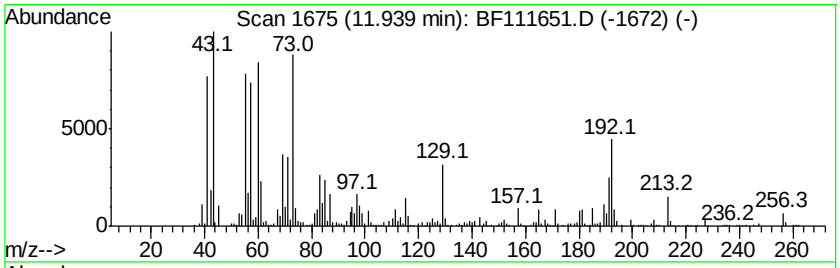
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

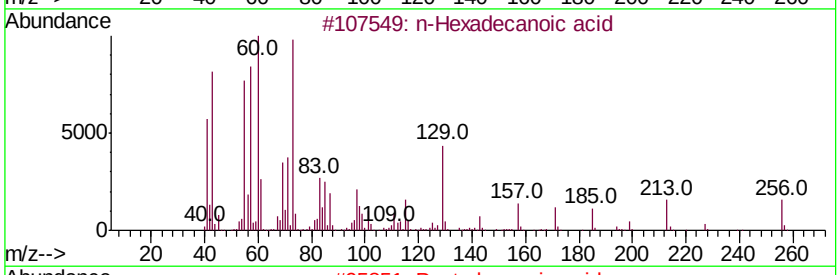
****

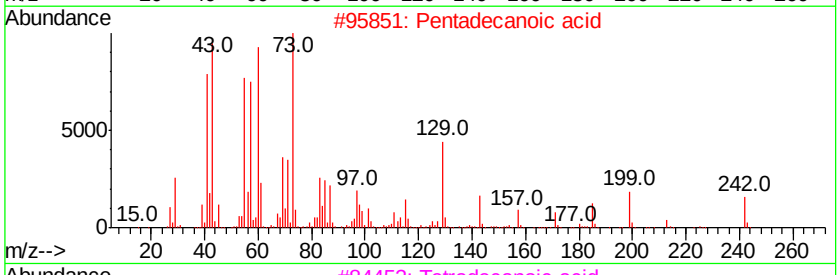
Peak Number 15 n-Hexadecanoic acid Concentration Rank 4

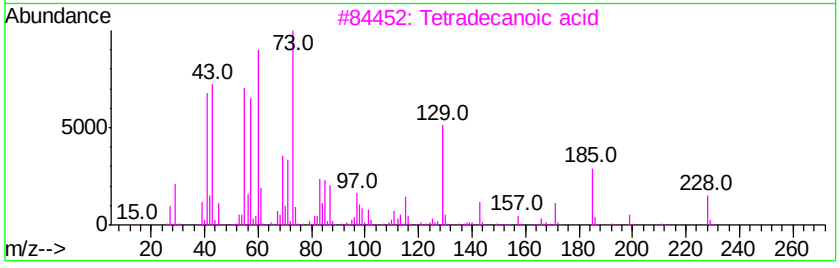
R.T.	EstConc	Area	Relative to ISTD	R.T.
11.94	10.03 ng	654943	Phenanthrene-d10	11.42

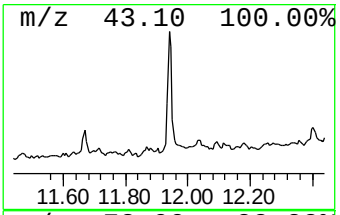
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	93
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	92
4	Tridecanoic acid	214	C13H26O2	000638-53-9	58
5	n-Decanoic acid	172	C10H20O2	000334-48-5	50

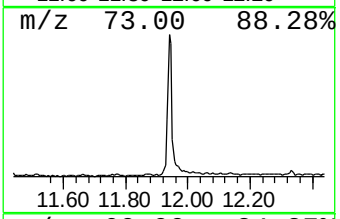


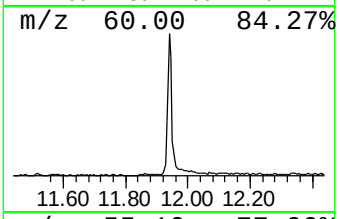


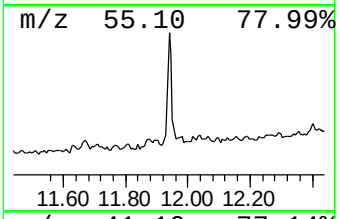


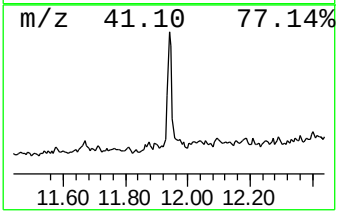












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Acq On : 20 Dec 2018 21:00
Operator : JU/SJ
Sample : J6346-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : DW-5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF121918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

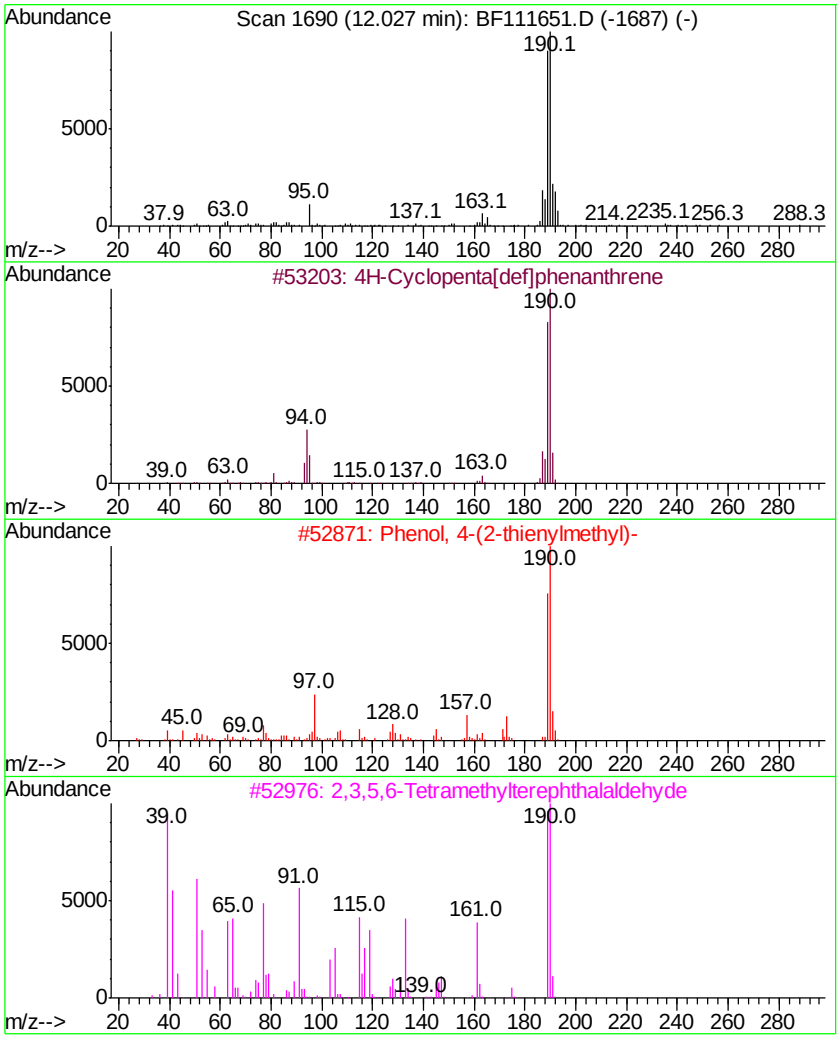
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

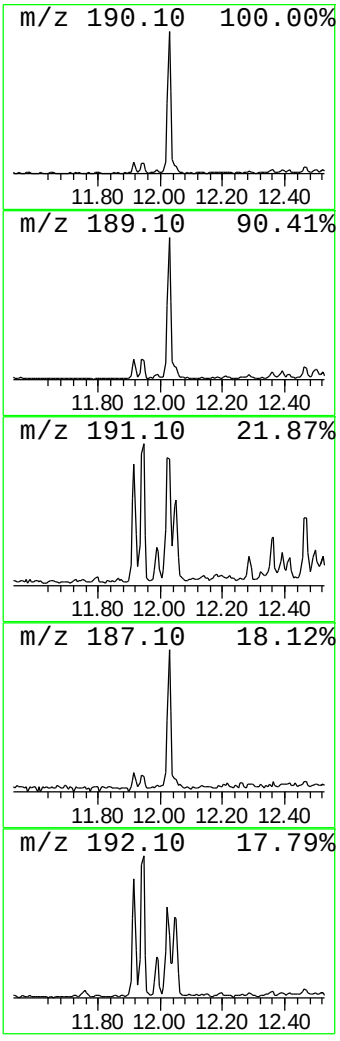
****

Peak Number 16 4H-Cyclopenta[def]phenanthrene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.03	4.20 ng	274025	Phenanthrene-d10	11.42

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	90
2	Phenol, 4-(2-thienylmethyl)-	190	C11H10OS	091680-55-6	64
3	2,3,5,6-Tetramethylterephthalald...	190	C12H14O2	007072-01-7	53
4	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	50
5	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	45





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Misc :
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Instrument : BNA_F
ClientSampleId : DW-5

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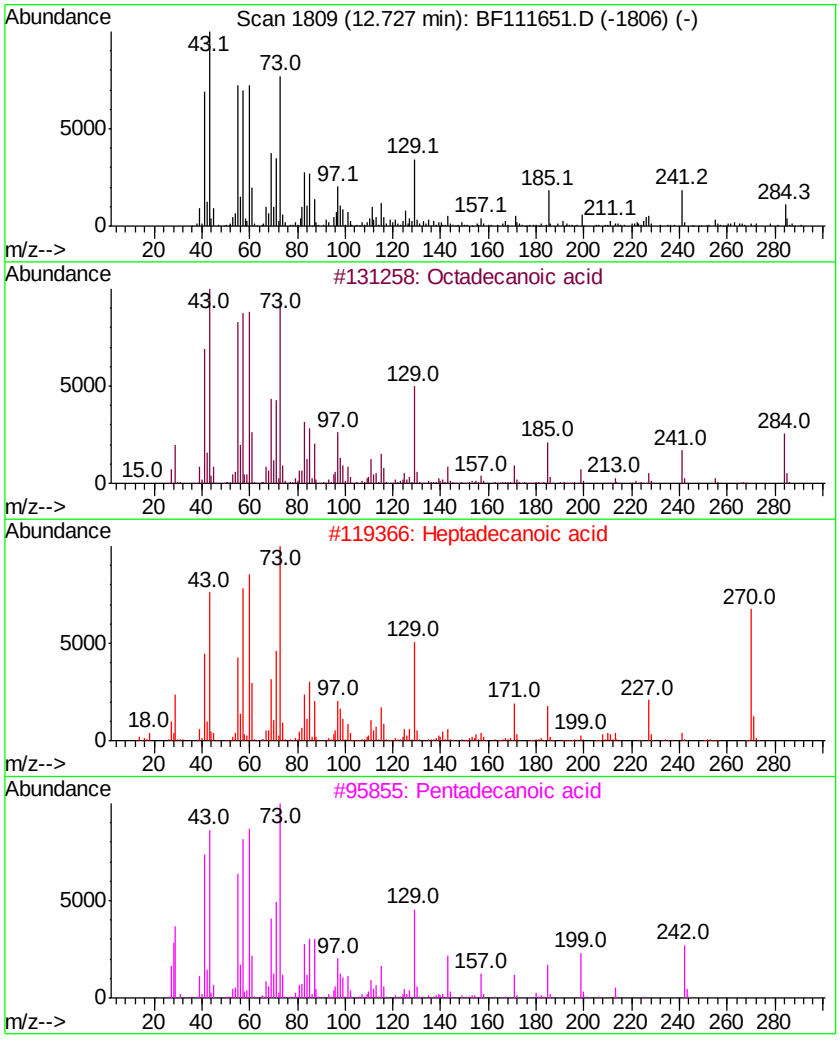
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

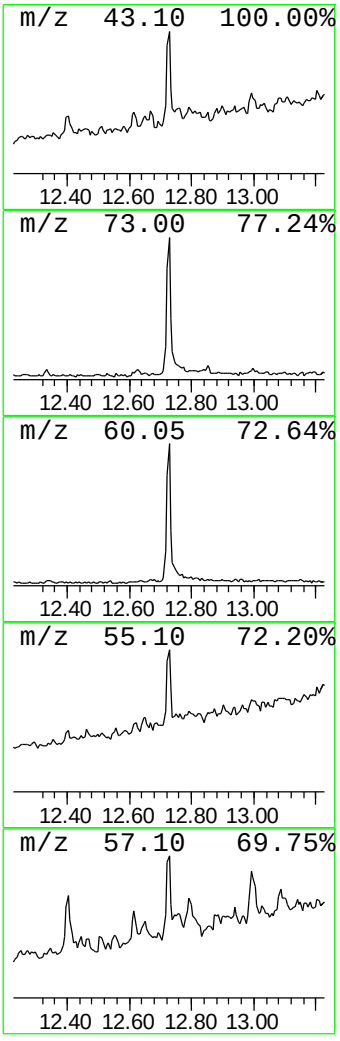
****

Peak Number 17 Octadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.73	5.12 ng	334497	Phenanthrene-d10	11.42

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	Heptadecanoic acid	270	C17H34O2	000506-12-7	93
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	83
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	78
5	Tridecanoic acid	214	C13H26O2	000638-53-9	70





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Instrument : BNA_F
ClientSampleId : DW-5

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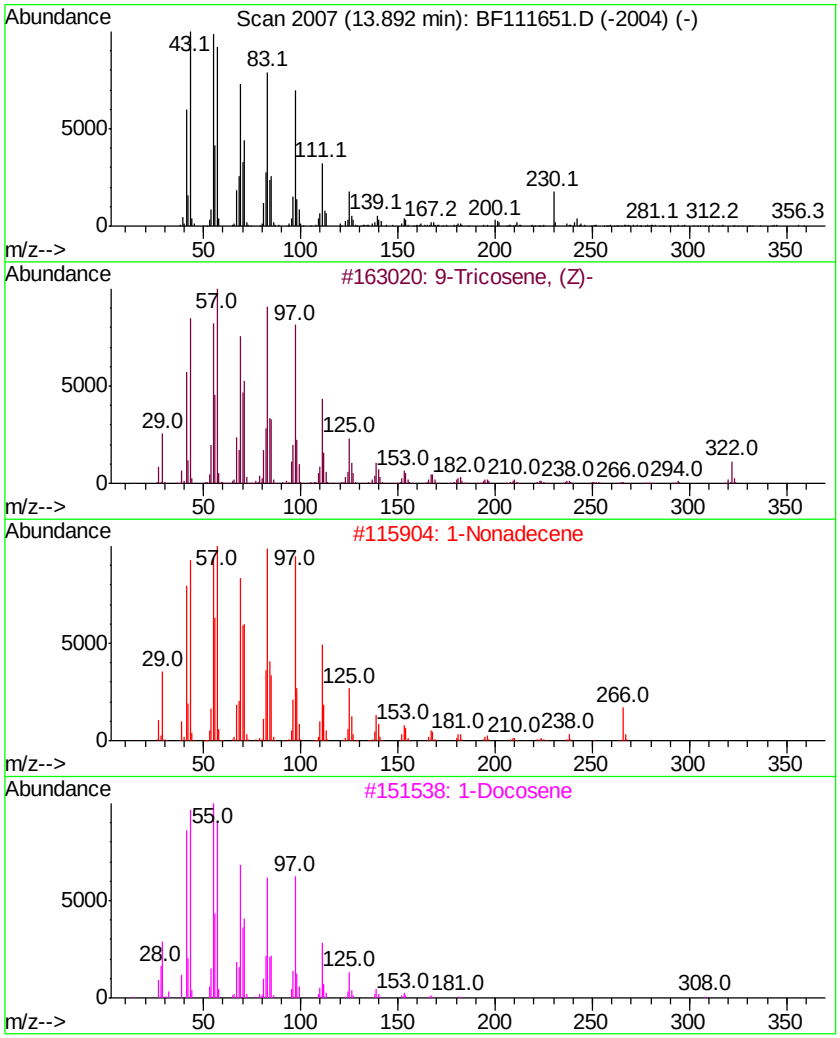
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

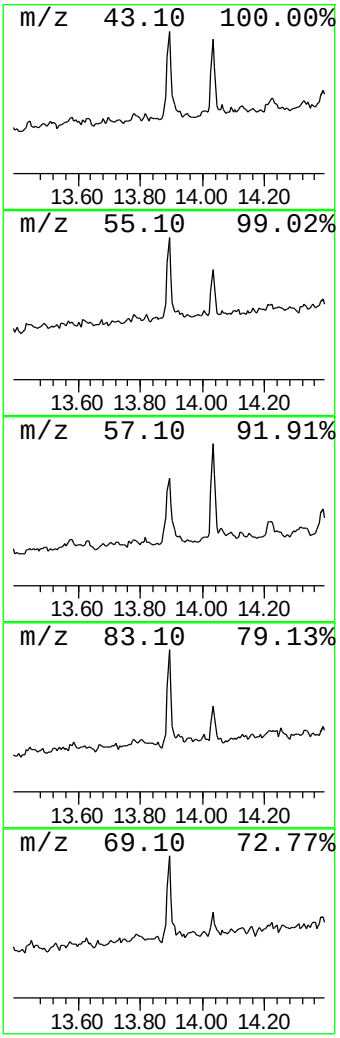
****

Peak Number 18 9-Tricosene, (Z)- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.89	4.77 ng	652614	Chrysene-d12	14.05

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Tricosene, (Z)-	322	C23H46	027519-02-4	95
2	1-Nonadecene	266	C19H38	018435-45-5	95
3	1-Docosene	308	C22H44	001599-67-3	94
4	1-Heneicosanol	312	C21H44O	015594-90-8	93
5	Bromoacetic acid, hexadecyl ester	362	C18H35BrO2	005454-48-8	91





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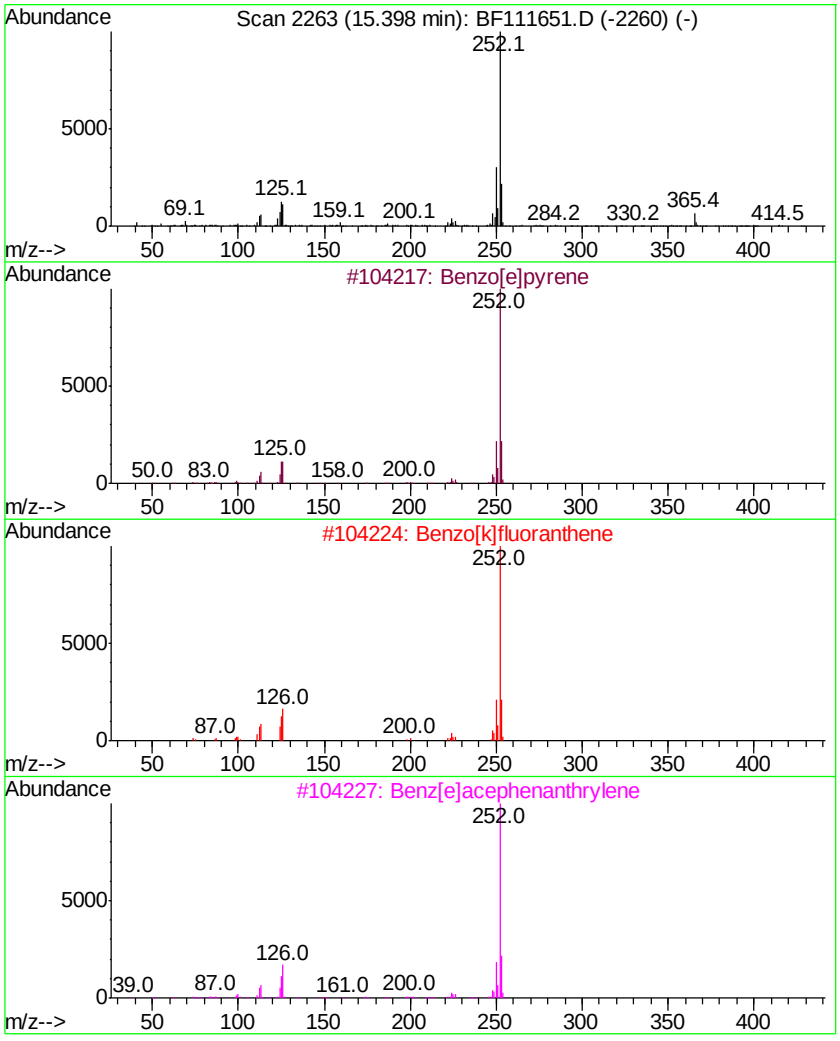
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

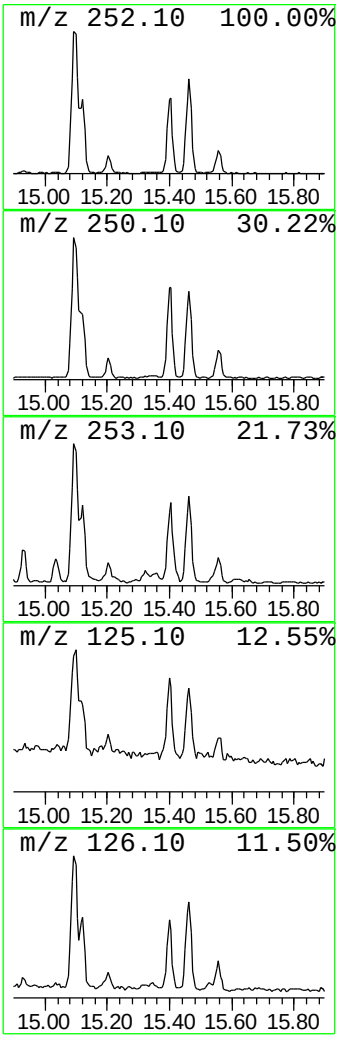
****

Peak Number 19 Benzo[e]pyrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.40	13.24 ng	516363	Perylene-d12	15.53

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[e]pyrene	252	C20H12	000192-97-2	98
2	Benzo[k]fluoranthene	252	C20H12	000207-08-9	96
3	Benz[e]acephenanthrylene	252	C20H12	000205-99-2	95
4	Benzo[a]pyrene	252	C20H12	000050-32-8	91
5	Benzo[j]fluoranthene	252	C20H12	000205-82-3	90





Tentatively Identified Compound (LSC) summary

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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...	5.13	4.0	ng	241001	1	6.90	1221190	20.0
unknown6.67	6.67	73.9	ng	4514970	1	6.90	1221190	20.0
unknown10.85	10.85	7.7	ng	502857	4	11.42	1305610	20.0
Phenol, 4-(1,1,3,...	10.90	5.3	ng	342454	4	11.42	1305610	20.0
unknown10.93	10.93	6.1	ng	398918	4	11.42	1305610	20.0
Phenol, m-tert-bu...	10.97	5.1	ng	331583	4	11.42	1305610	20.0
unknown10.99	10.99	3.1	ng	200180	4	11.42	1305610	20.0
Benzoic acid, 2-(...	11.16	3.0	ng	197656	4	11.42	1305610	20.0
Pentadecane	11.32	5.3	ng	343523	4	11.42	1305610	20.0
n-Hexadecanoic acid	11.94	10.0	ng	654943	4	11.42	1305610	20.0
4H-Cyclopenta[def...	12.03	4.2	ng	274025	4	11.42	1305610	20.0
Octadecanoic acid	12.73	5.1	ng	334497	4	11.42	1305610	20.0
9-Tricosene, (Z)-	13.89	4.8	ng	652614	5	14.05	2737520	20.0
Benzo[e]pyrene	15.40	13.2	ng	516363	6	15.53	779771	20.0