LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF033120\
Data File : BF119663.D
Acq On : 31 Mar 2020 20:22
Operator : MA/SJ
Sample : L2055-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-7-(1-3)

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-BF031820.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	2.193	14	18	24	rVB3	114714	163515	2.90%	0.293%
2	5.040	494	502	510	rBV	160199	195149	3.46%	0.350%
3	5.445	566	571	574	rBV	5912972	5639336	100.00%	10.107%
4	6.457	738	743	746	rBV	5157776	5594223	99.20%	10.026%
5	6.828	803	806	810	rBV	1914764	1636787	29.02%	2.933%
6	6.986	829	833	836	rBV	5139774	4493989	79.69%	8.054%
7	7.392	897	902	905	rBV	3832420	3455824	61.28%	6.193%
8	8.116	1020	1025	1028	rBV	2560511	2000617	35.48%	3.585%
9	9.198	1205	1209	1212	rBV	6767957	5487877	97.31%	9.835%
10	9.280	1220	1223	1227	rBV2	88863	105355	1.87%	0.189%
11	9.586	1272	1275	1280	rBV	530387	502971	8.92%	0.901%
12	9.739	1297	1301	1305	rBV	120085	120728	2.14%	0.216%
13	9.810	1311	1313	1318	rVB2	80644	88467	1.57%	0.159%
14	9.880	1321	1325	1328	rVB	2054521	1880006	33.34%	3.369%
15	10.310	1396	1398	1401	rVB	102967	83950	1.49%	0.150%
16	10.422	1414	1417	1424	rBV3	55405	80767	1.43%	0.145%
17	10.533	1431	1436	1439	rBV	85557	106660	1.89%	0.191%
18	10.669	1454	1459	1462	rBV	2802353	2553927	45.29%	4.577%
19	10.769	1473	1476	1478	rBV	305108	285307	5.06%	0.511%
20	10.786	1478	1479	1484	rVB2	205253	256419	4.55%	0.460%
21	11.239	1553	1556	1558	rBV	132190	106800	1.89%	0.191%
22	11.269	1558	1561	1565	rVB2	126493	134259	2.38%	0.241%
23	11.316	1566	1569	1574	rBV3	54356	70026	1.24%	0.125%
24	11.369	1574	1578	1580	rVV	1878223	1538593	27.28%	2.757%
25	11.392	1580	1582	1585	rVV	511509	403575	7.16%	0.723%
26	11.439	1585	1590	1594	rVB	339173	353468	6.27%	0.633%
27	11.598	1612	1617	1619	rBV	58773	77941	1.38%	0.140%
28	11.663	1626	1628	1631	rVB	151430	122795	2.18%	0.220%
29	11.869	1659	1663	1665	rBV2	113873	131285	2.33%	0.235%
30	11.898	1665	1668	1674	rVV	910579	793001	14.06%	1.421%
31	11.980	1679	1682	1685	rBV	193256	194289	3.45%	0.348%
32	12.074	1695	1698	1701	rBV	178017	128547	2.28%	0.230%
33	12.169	1711	1714	1716	rBV	77582	69413	1.23%	0.124%
34	12.351	1742	1745	1751	rVB3	56969	85934	1.52%	0.154%

LSC Area Percent Report

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Data File : BF119663.D
Acq On : 31 Mar 2020 20:22
Operator : MA/SJ
Sample : L2055-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-7-(1-3)

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-BF031820.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	12.421	1754	1757	1760	rBV	105581	102161	1.81%	0.183%
36	12.463	1760	1764	1769	rVV3	152677	176522	3.13%	0.316%
37	12.504	1769	1771	1776	rBV2	75697	85955	1.52%	0.154%
38	12.586	1781	1785	1788	rBV	1693416	1478769	26.22%	2.650%
39	12.680	1798	1801	1809	rBV2	296770	422353	7.49%	0.757%
40	12.816	1818	1824	1826	rBV	1333488	1367074	24.24%	2.450%
41	12.839	1826	1828	1830	rVB	144599	105190	1.87%	0.189%
42	12.933	1841	1844	1845	rBV	93871	78822	1.40%	0.141%
43	12.963	1845	1849	1852	rVB	4349143	3946413	69.98%	7.073%
44	13.027	1857	1860	1865	rBV2	95433	188370	3.34%	0.338%
45	13.145	1877	1880	1883	rVB2	292571	266982	4.73%	0.478%
46	13.210	1886	1891	1894	rVV2	260799	337346	5.98%	0.605%
47	13.245	1894	1897	1901	rVV2	142623	150809	2.67%	0.270%
48	13.368	1916	1918	1922	rBV2	76369	80904	1.43%	0.145%
49	13.539	1945	1947	1950	rBV	152802	123999	2.20%	0.222%
50	13.792	1988	1990	1992	rBV	113369	83570	1.48%	0.150%
51	13.815	1992	1994	1997	rVV2	101030	114334	2.03%	0.205%
52	13.874	2001	2004	2012	rVB	674413	966646	17.14%	1.732%
53	14.015	2023	2028	2030	rBV2	1721580	2243535	39.78%	4.021%
54	14.039	2030	2032	2038	rVB	763334	694047	12.31%	1.244%
55	14.121	2044	2046	2049	rBV	184459	135881	2.41%	0.244%
56	14.186	2055	2057	2059	rBV	156933	127686	2.26%	0.229%
57	14.492	2106	2109	2112	rBV2	341964	346001	6.14%	0.620%
58	14.798	2159	2161	2164	rBV	169765	156856	2.78%	0.281%
59	15.057	2202	2205	2213	rVB	523449	1005277	17.83%	1.802%
60	15.139	2217	2219	2222	rBV2	205792	210944	3.74%	0.378%
61	15.357	2254	2256	2262	rVB	299684	379890	6.74%	0.681%
62	15.421	2264	2267	2272	rBV	429742	473115	8.39%	0.848%
63	15.486	2274	2278	2281	rBV	851297	1007358	17.86%	1.805%

Sum of corrected areas: 55798609

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Sample : L2055-08
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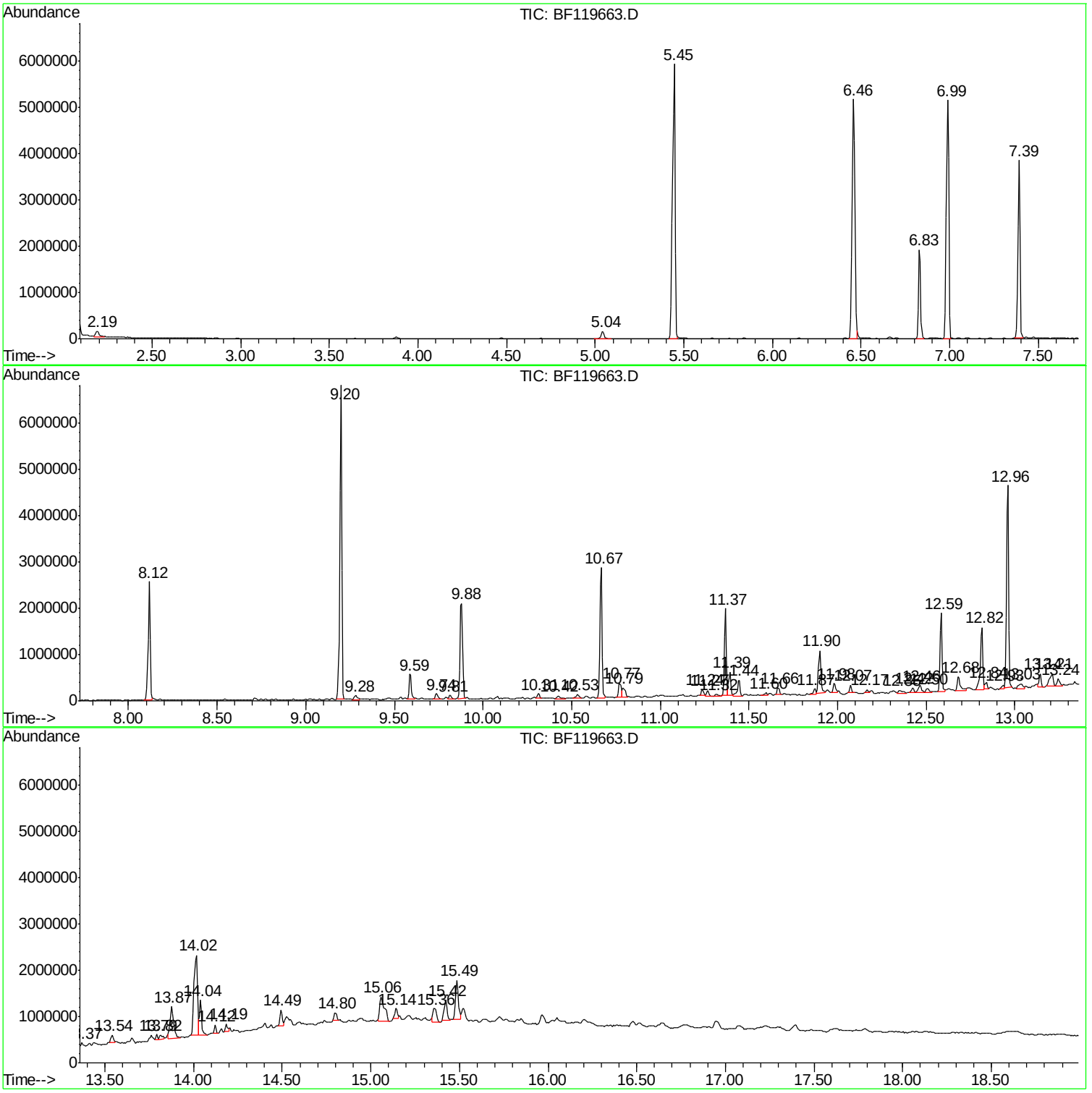
Instrument :
BNA_F
ClientSampleId :
SB-7-(1-3)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF031820.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

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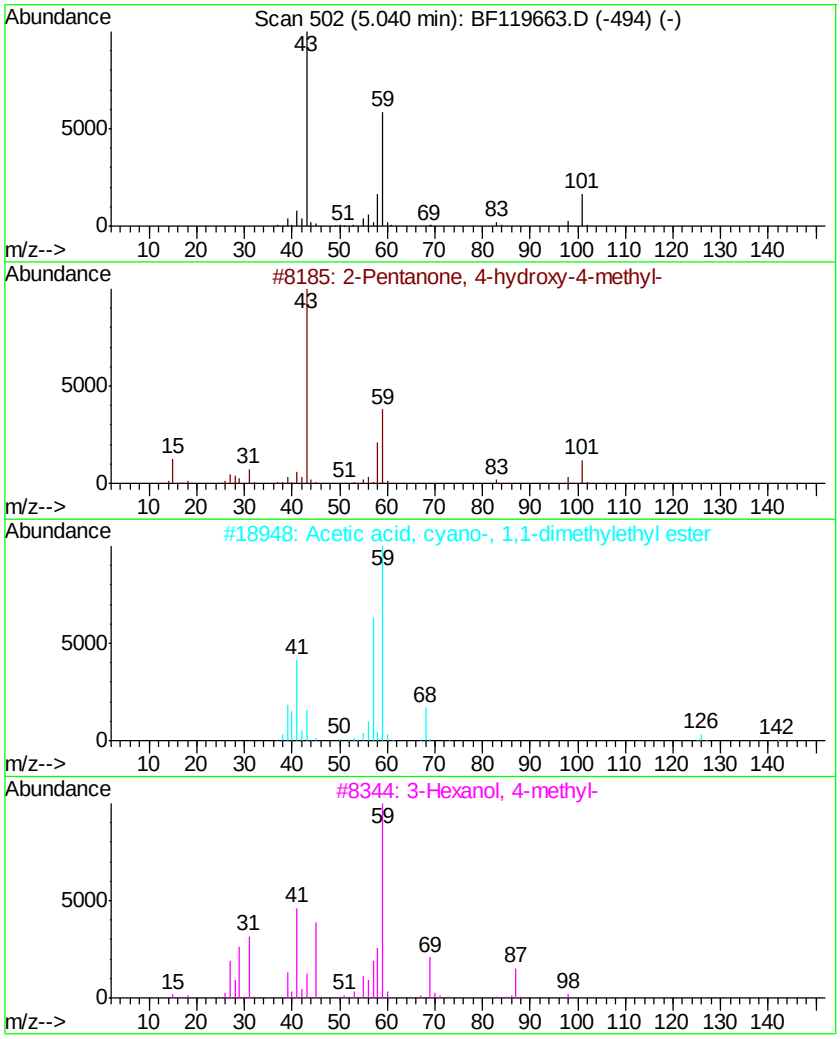
TIC Integration Parameters: LSCINT.P

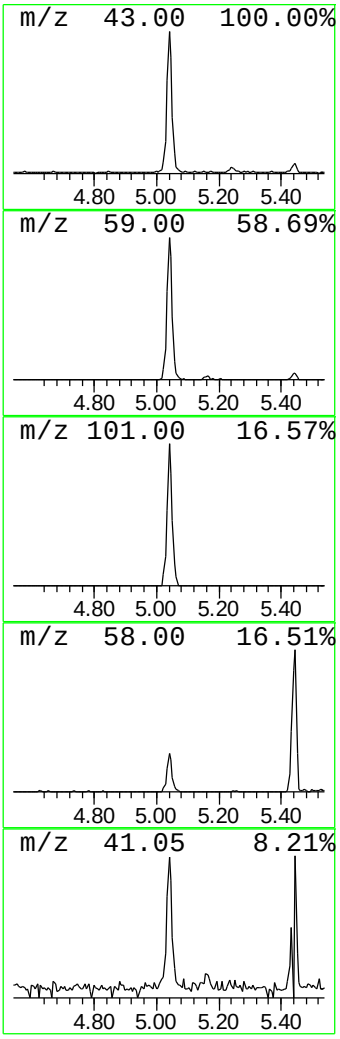
****

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.04	2.38 ng	195149	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	72
2	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	37
3	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
4	3-Hexanol, 4-ethyl-	130	C8H18O	019780-44-0	28
5	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28





Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF033120\
Data File : BF119663.D
Acq On : 31 Mar 2020 20:22
Operator : MA/SJ
Sample : L2055-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-7-(1-3)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF031820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

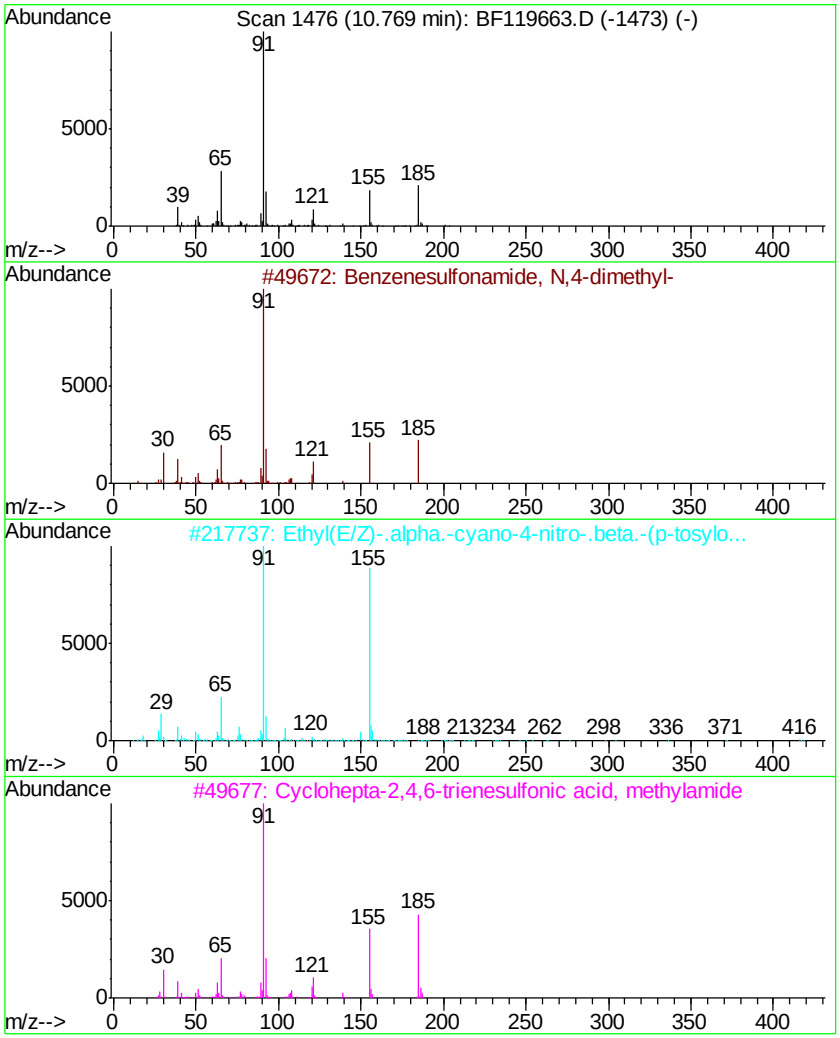
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

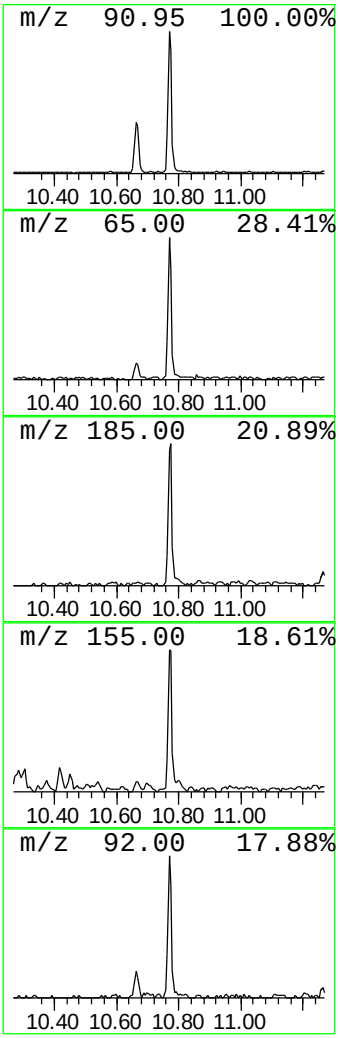
****

Peak Number 3 Benzenesulfonamide, N,4-dim... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.77	3.71 ng	285307	Phenanthrene-d10	11.37

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenesulfonamide, N,4-dimethyl-	185	C8H11NO2S	000640-61-9	87
2	Ethyl(E/Z)-.alpha.-cyano-4-nitro...	416	C19H16N2O7S	1000287-26-8	53
3	Cyclohepta-2,4,6-trienesulfonic ...	185	C8H11NO2S	1000187-28-7	46
4	Benzene, 1,1'-[sulfinylbis(methy...	230	C14H14OS	000621-08-9	43
5	Toluene-4-sulfonic acid, 4-benzy...	354	C20H18O4S	1000295-51-0	43





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Operator : MA/SJ
Sample : L2055-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-7-(1-3)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF031820.M
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TIC Library : C:\DATABASE\NIST11.L
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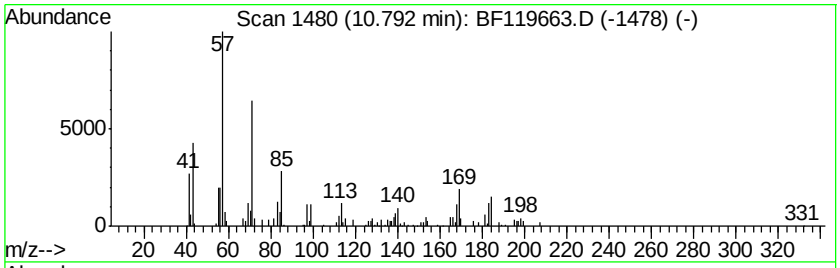
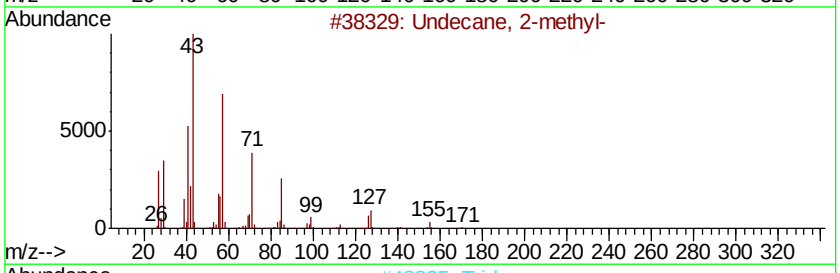
****

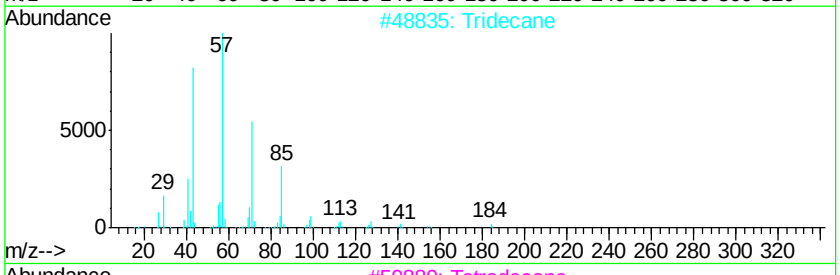
Peak Number 4 Undecane, 2-methyl- Concentration Rank 8

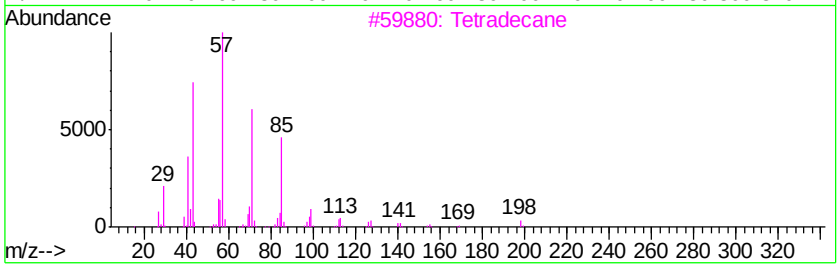
R.T.	EstConc	Area	Relative to ISTD	R.T.
10.79	3.33 ng	256419	Phenanthrene-d10	11.37

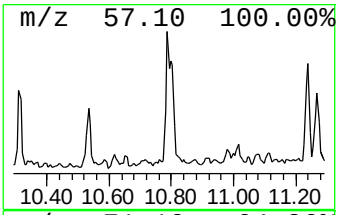
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane, 2-methyl-	170	C12H26	007045-71-8	72
2	Tridecane	184	C13H28	000629-50-5	70
3	Tetradecane	198	C14H30	000629-59-4	70
4	Dodecane, 2-methyl-	184	C13H28	001560-97-0	68
5	Tetracosane	338	C24H50	000646-31-1	64

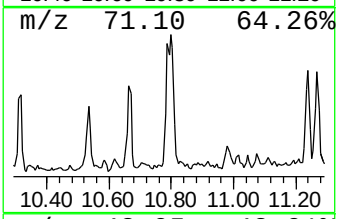


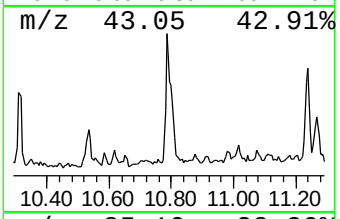


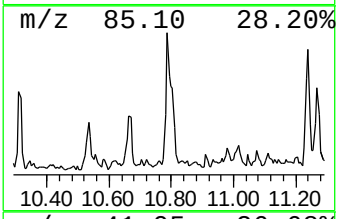


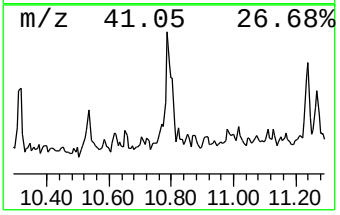












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Instrument : BNA_F
ClientSampleId : SB-7-(1-3)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF031820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

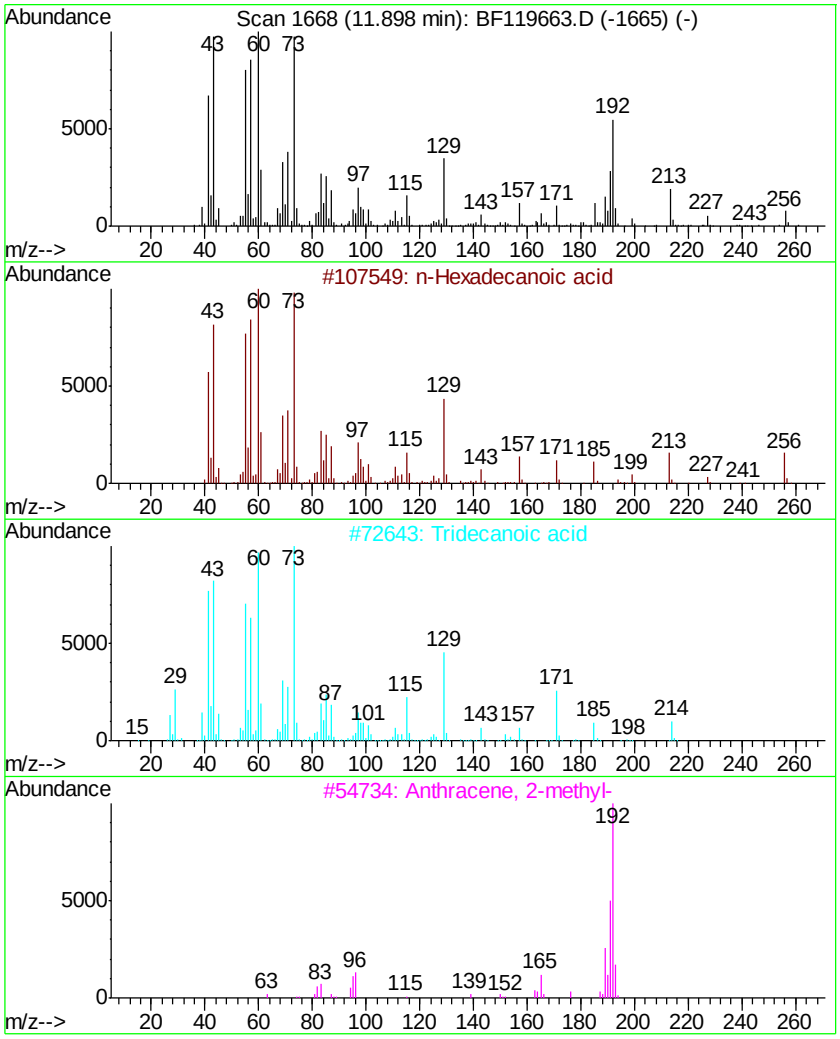
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

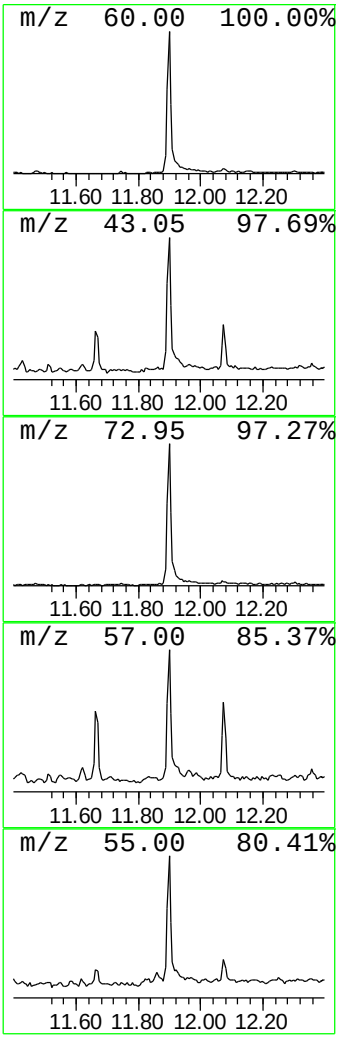
****

Peak Number 5 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.90	10.31 ng	793001	Phenanthrene-d10	11.37

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	86
2	Tridecanoic acid	214	C13H26O2	000638-53-9	78
3	Anthracene, 2-methyl-	192	C15H12	000613-12-7	74
4	Pentadecanoic acid	242	C15H30O2	001002-84-2	60
5	Tetradecanoic acid	228	C14H28O2	000544-63-8	58





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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

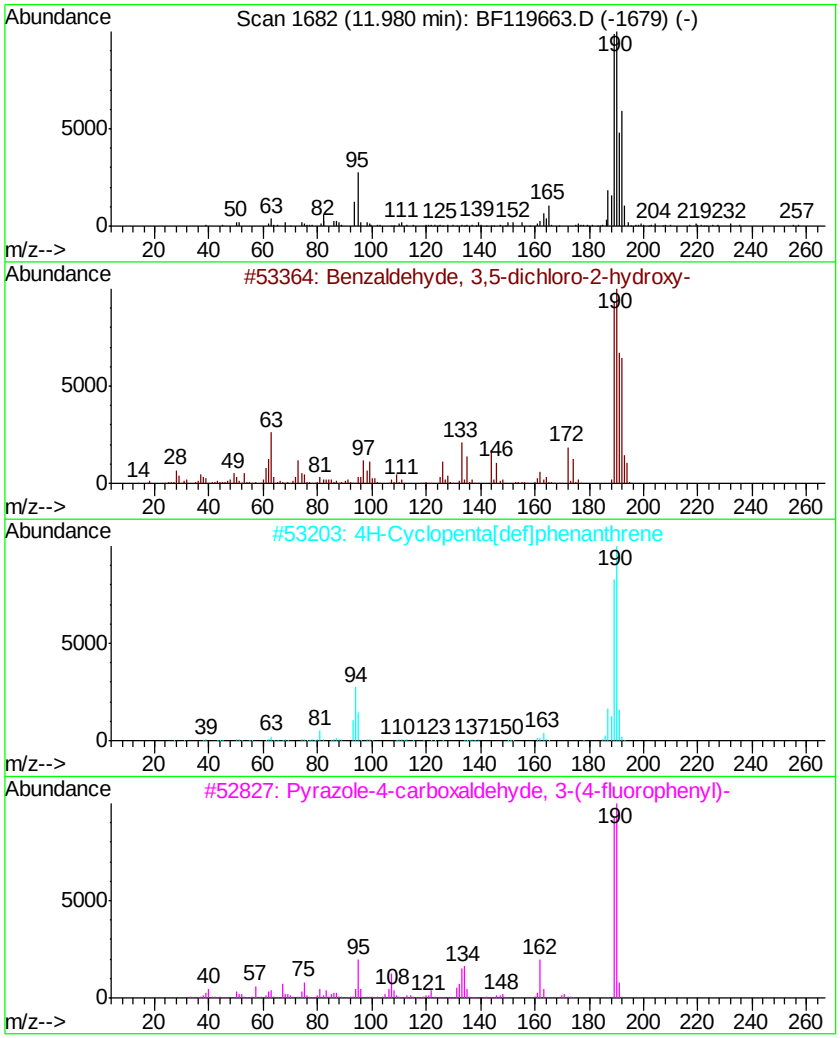
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

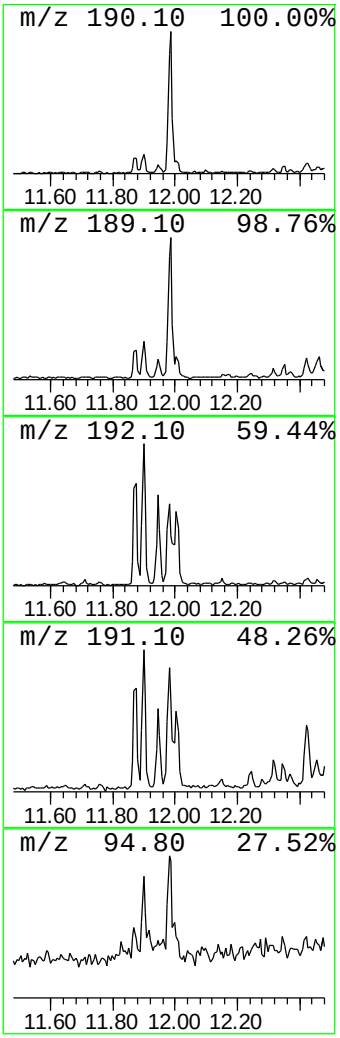
****

Peak Number 6 Benzaldehyde, 3,5-dichloro-... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.98	2.53 ng	194289	Phenanthrene-d10	11.37

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	72
2	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	64
3	Pyrazole-4-carboxaldehyde, 3-(4-...	190	C10H7FN2O	306936-57-2	49
4	2,3,5,6-Tetramethylterephthalald...	190	C12H14O2	007072-01-7	43
5	Naphtho[1,2-b]norbornadiene	192	C15H12	1000210-14-8	38





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Instrument : BNA_F
ClientSampleId : SB-7-(1-3)

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

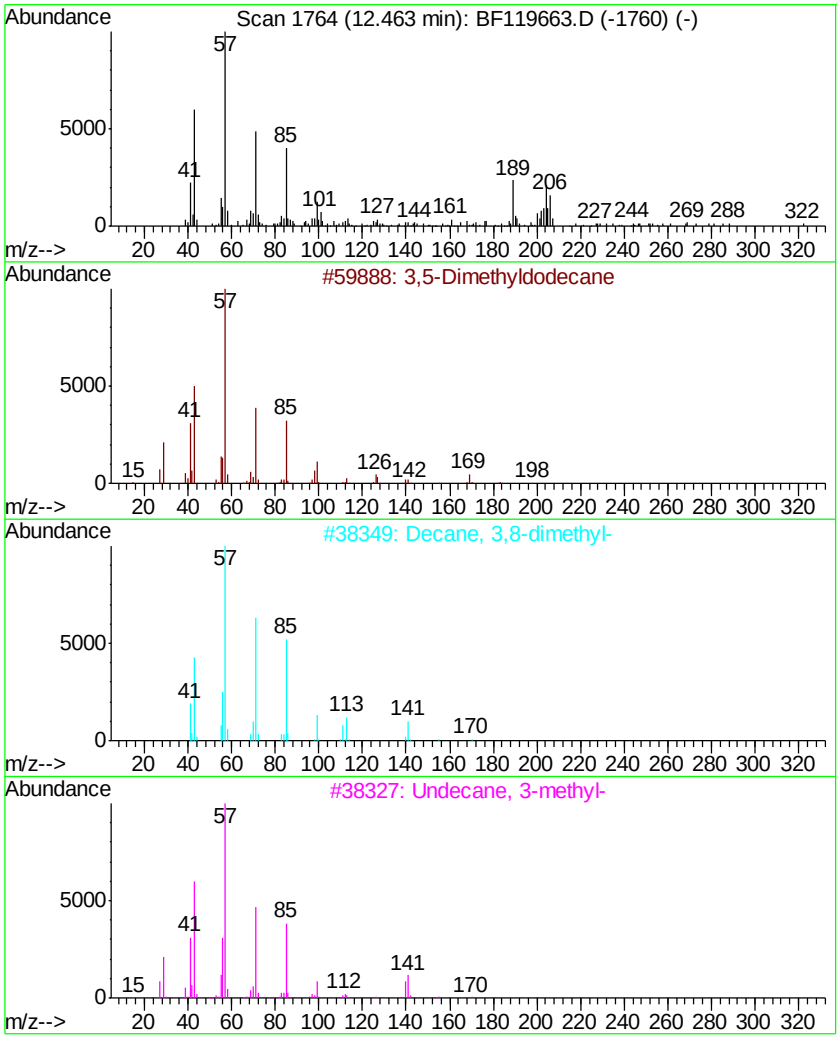
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

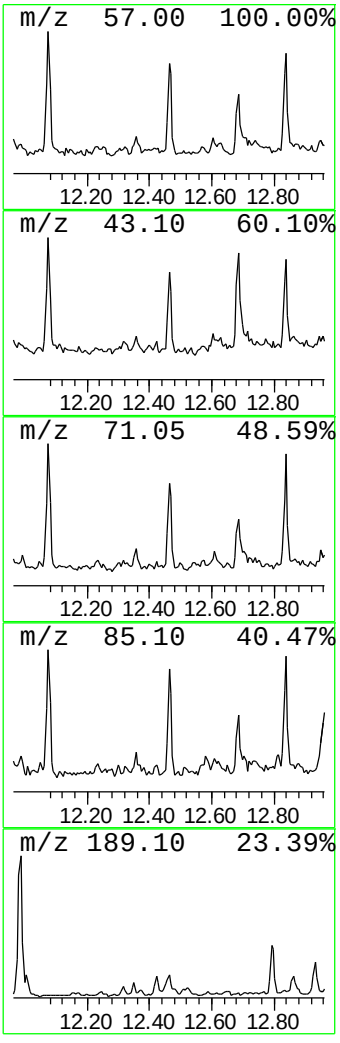
****

Peak Number 7 3,5-Dimethyldodecane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.46	2.29 ng	176522	Phenanthrene-d10	11.37

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,5-Dimethyldodecane	198	C14H30	107770-99-0	60
2	Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	60
3	Undecane, 3-methyl-	170	C12H26	001002-43-3	49
4	Nonadecane	268	C19H40	000629-92-5	43
5	Tridecane	184	C13H28	000629-50-5	42





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Data File : BF119663.D
Acq On : 31 Mar 2020 20:22
Operator : MA/SJ
Sample : L2055-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-7-(1-3)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF031820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

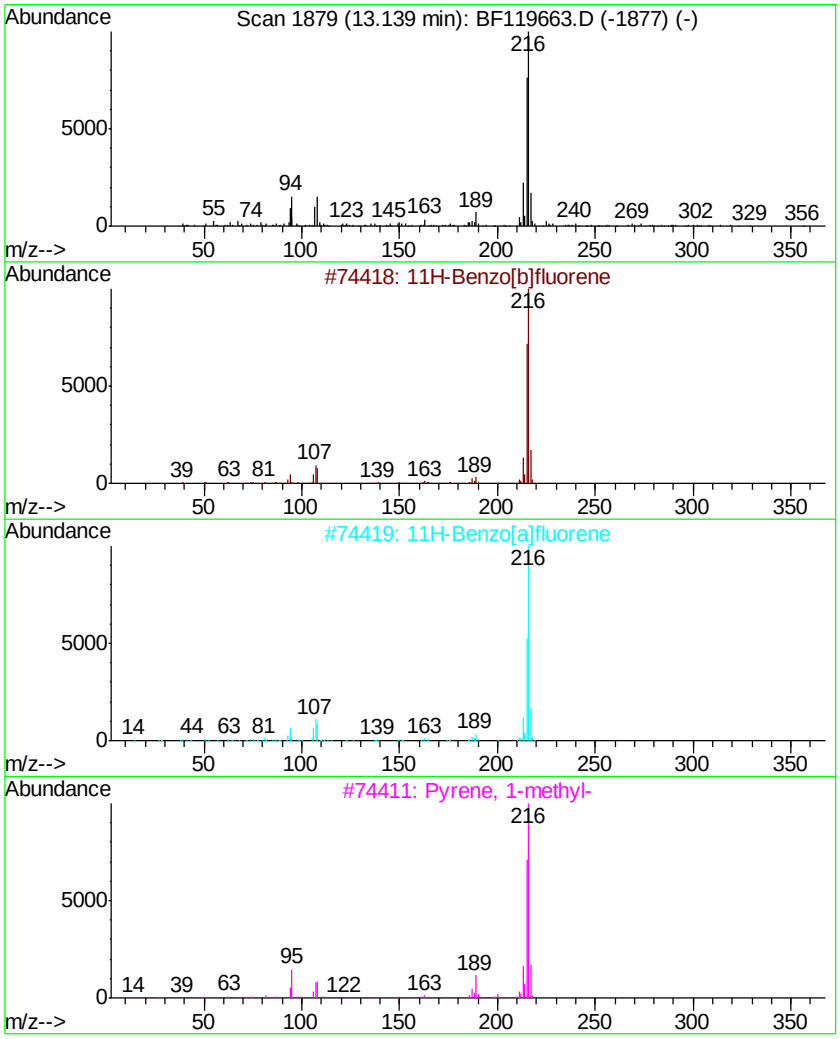
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

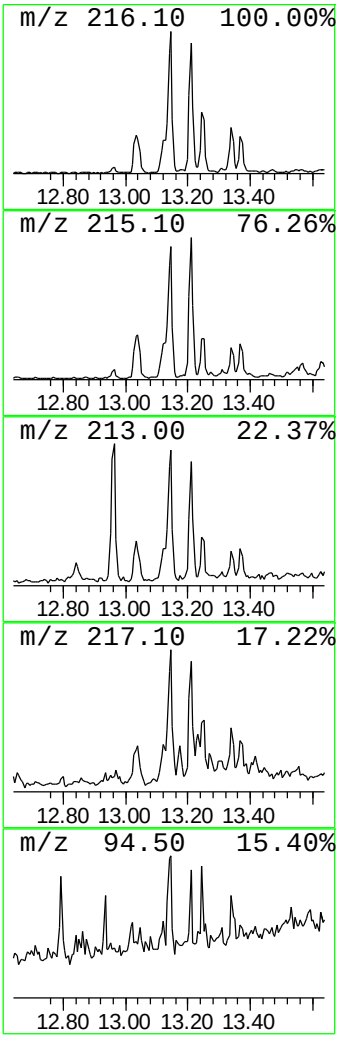
****

Peak Number 9 11H-Benzo[b]fluorene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.14	2.38 ng	266982	Chrysene-d12	14.02

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	94
2	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	92
3	Pyrene, 1-methyl-	216	C17H12	002381-21-7	86
4	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	74
5	Pyrene, 2-methyl-	216	C17H12	003442-78-2	60





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Data File : BF119663.D
Acq On : 31 Mar 2020 20:22
Operator : MA/SJ
Sample : L2055-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-7-(1-3)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF031820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

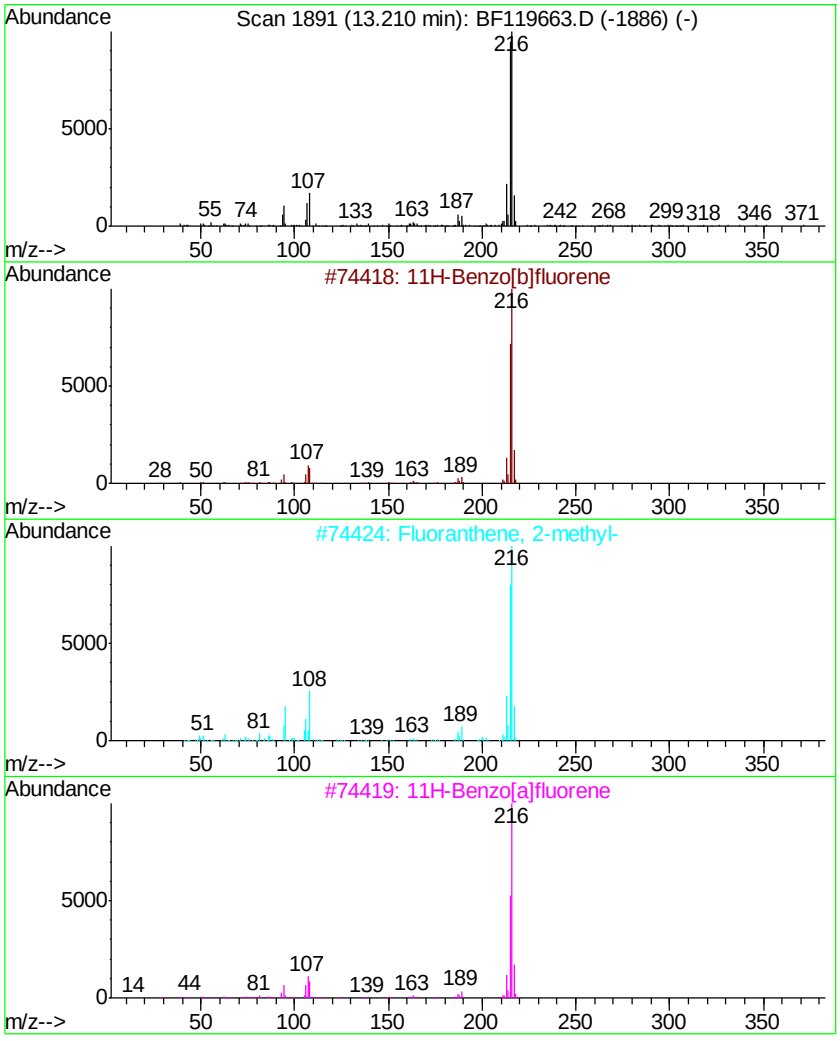
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

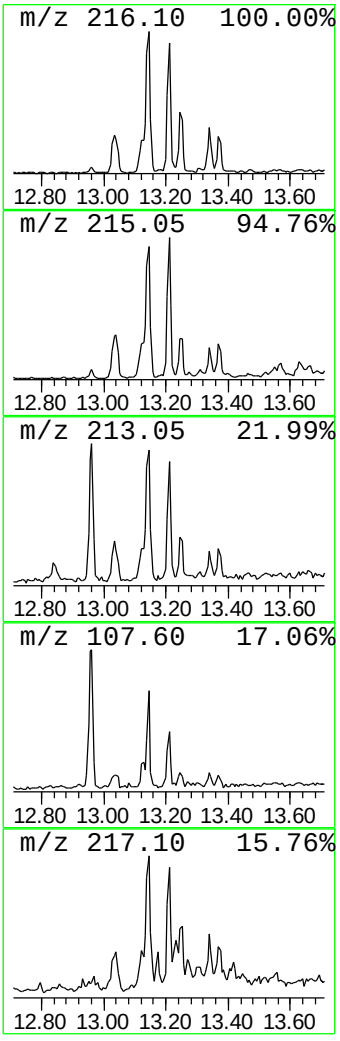
****

Peak Number 10 Fluoranthene, 2-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.21	3.01 ng	337346	Chrysene-d12	14.02

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	83
2	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	64
3	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	62
4	Benzoic acid, 3-(3,5-dimethyl-1-...	216	C12H12N2O2	312531-88-7	58
5	Pyrene, 2-methyl-	216	C17H12	003442-78-2	58





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF033120\
Data File : BF119663.D
Acq On : 31 Mar 2020 20:22
Operator : MA/SJ
Sample : L2055-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-7-(1-3)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF031820.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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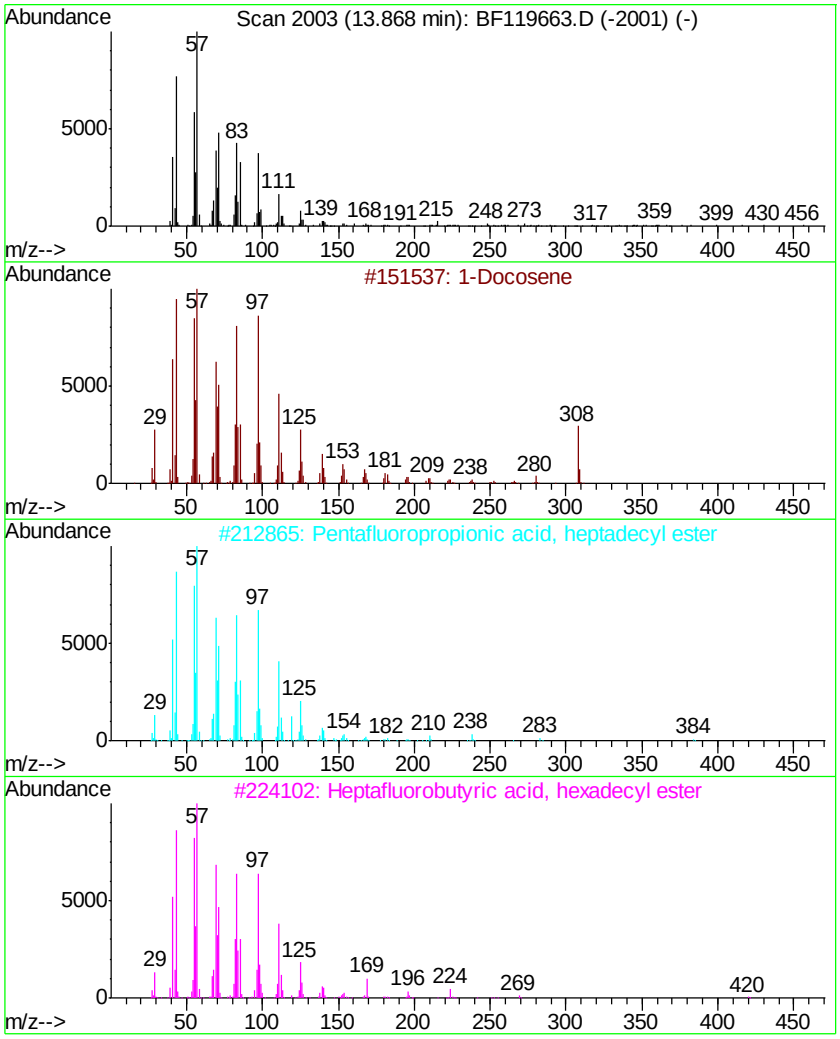
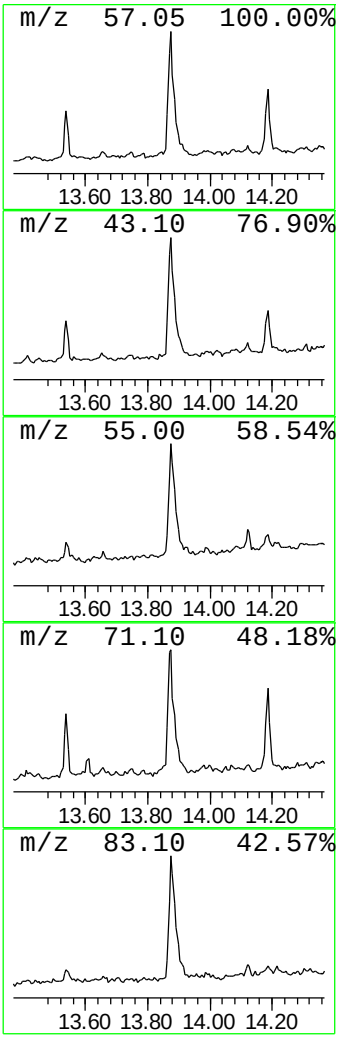
Peak Number 11 1-Docosene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.87	8.62 ng	966646	Chrysene-d12	14.02

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Docosene	308	C22H44	001599-67-3	95
2	Pentafluoropropionic acid, hepta...	402	C20H35F5O2	959218-78-1	94
3	Heptafluorobutyric acid, hexadec...	438	C20H33F7O2	006385-15-5	91
4	Docosyl pentafluoropropionate	472	C25H45F5O2	1000351-80-9	91
5	Octadecyl trifluoroacetate	366	C20H37F3O2	079392-43-1	91





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Data File : BF119663.D
Acq On : 31 Mar 2020 20:22
Operator : MA/SJ
Sample : L2055-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-7-(1-3)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF031820.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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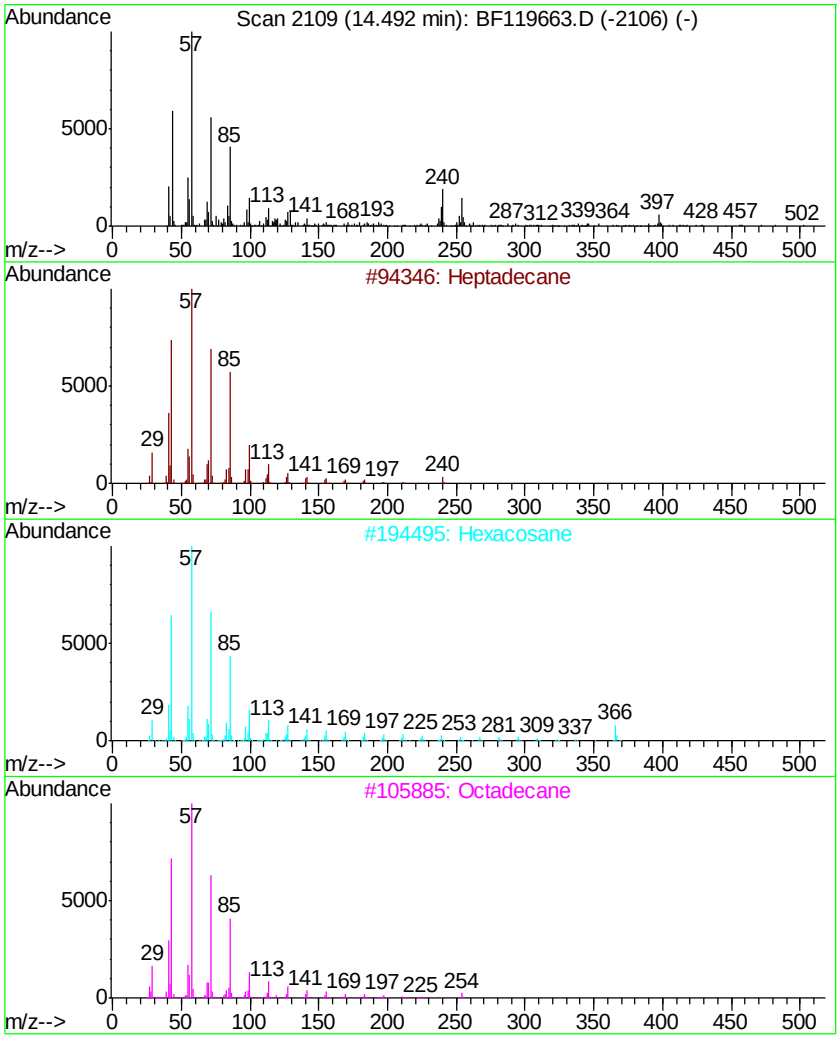
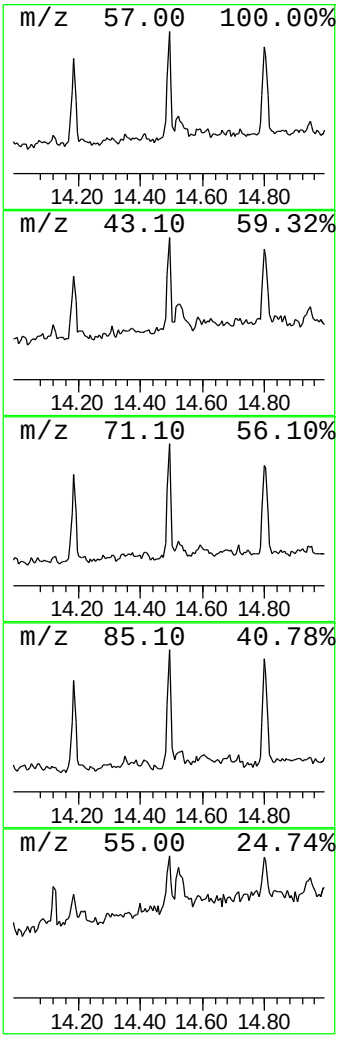
Peak Number 12 Heptadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.49	3.08 ng	346001	Chrysene-d12	14.02

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	86
2	Hexacosane	366	C26H54	000630-01-3	64
3	Octadecane	254	C18H38	000593-45-3	62
4	Octacosane	394	C28H58	000630-02-4	60
5	Heptacosane	380	C27H56	000593-49-7	58





Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF033120\
Data File : BF119663.D
Acq On : 31 Mar 2020 20:22
Operator : MA/SJ
Sample : L2055-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-7-(1-3)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF031820.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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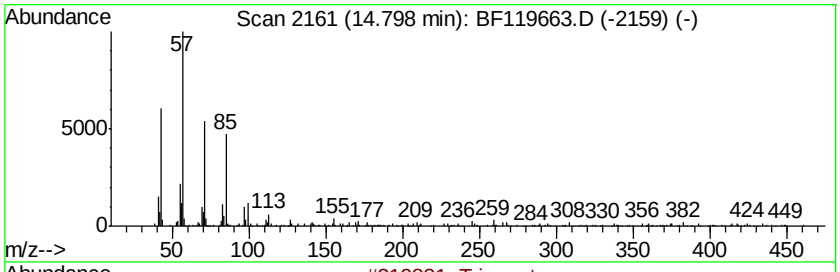
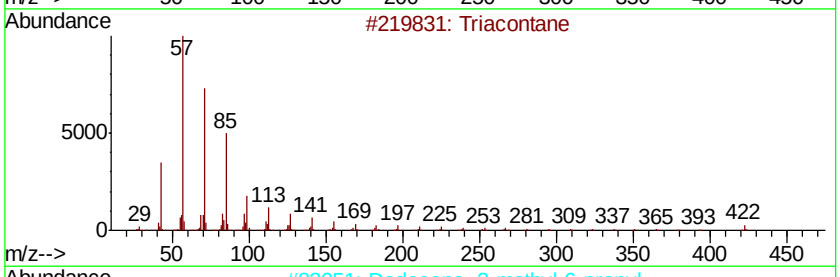
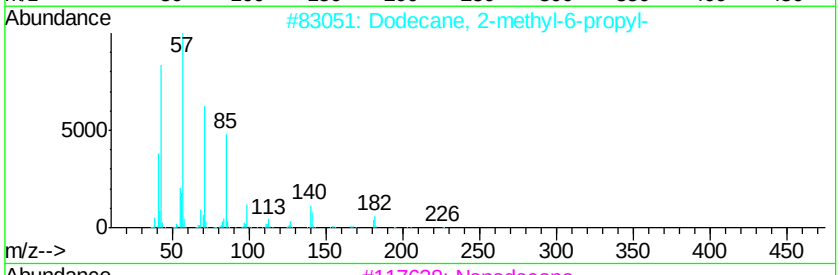
Peak Number 13 Triacontane Concentration Rank 9

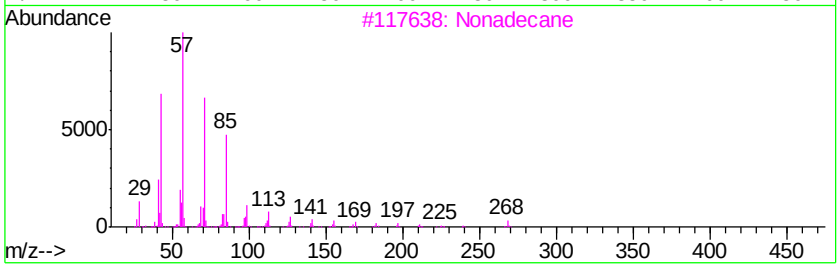
R.T.	EstConc	Area	Relative to ISTD	R.T.
14.80	3.11 ng	156856	Perylene-d12	15.49

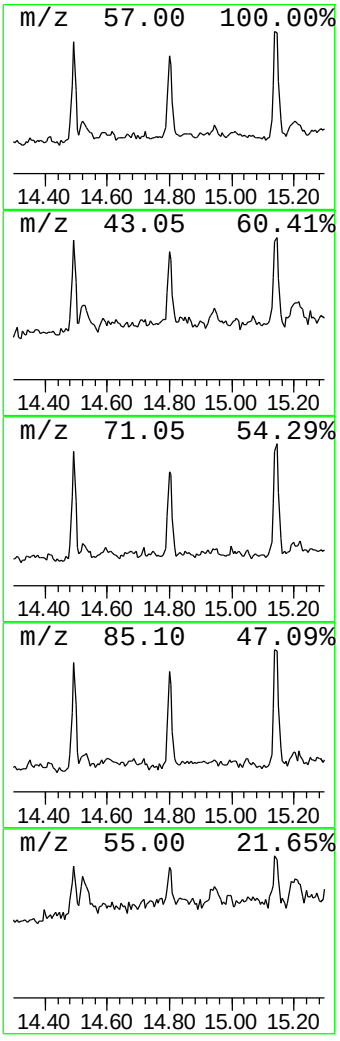
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Triacontane	422	C30H62	000638-68-6	74
2	Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	72
3	Nonadecane	268	C19H40	000629-92-5	64
4	Heneicosane	296	C21H44	000629-94-7	58
5	Octacosane	394	C28H58	000630-02-4	58











Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF033120\
Data File : BF119663.D
Acq On : 31 Mar 2020 20:22
Operator : MA/SJ
Sample : L2055-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-7-(1-3)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF031820.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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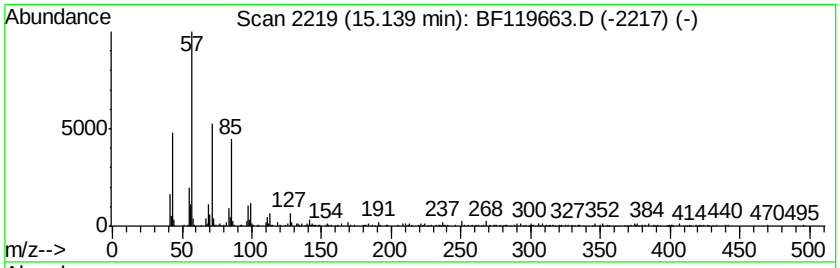
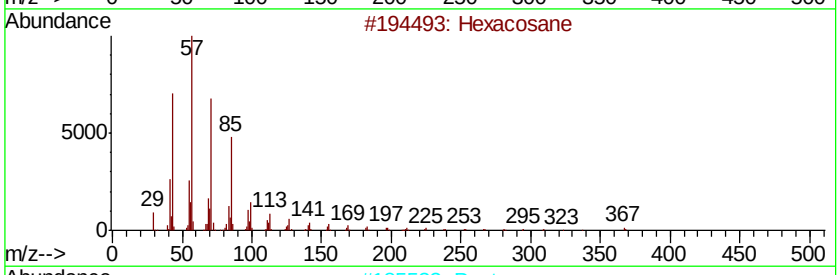
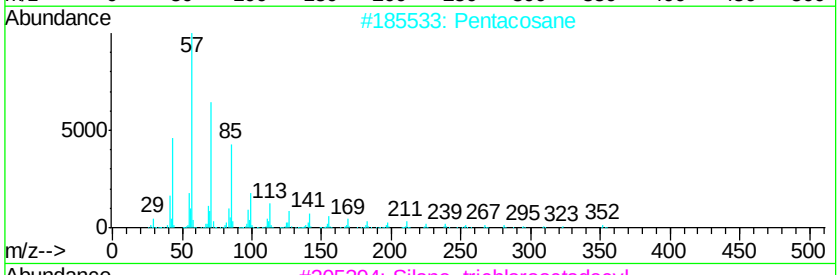
Peak Number 14 Hexacosane Concentration Rank 6

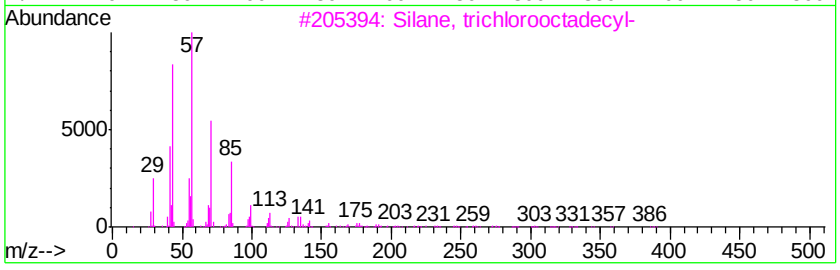
R.T.	EstConc	Area	Relative to ISTD	R.T.
15.14	4.19 ng	210944	Perylene-d12	15.49

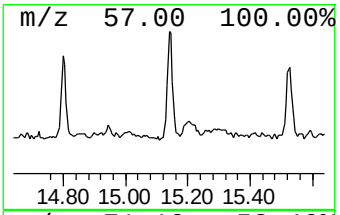
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexacosane	366	C26H54	000630-01-3	87
2	Pentacosane	352	C25H52	000629-99-2	83
3	Silane, trichlorooctadecyl-	386	C18H37Cl3Si	000112-04-9	83
4	Octadecane	254	C18H38	000593-45-3	80
5	1-Iodoundecane	282	C11H23I	004282-44-4	74

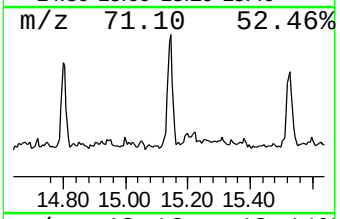


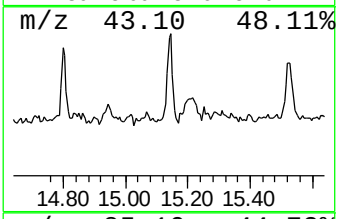


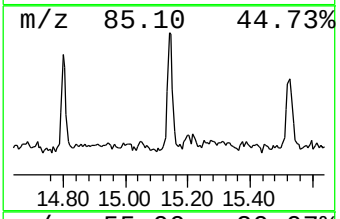


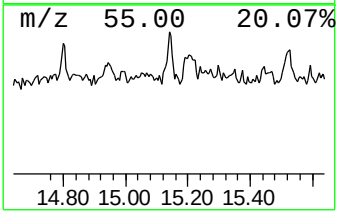












Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF033120\
Data File : BF119663.D
Acq On : 31 Mar 2020 20:22
Operator : MA/SJ
Sample : L2055-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-7-(1-3)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF031820.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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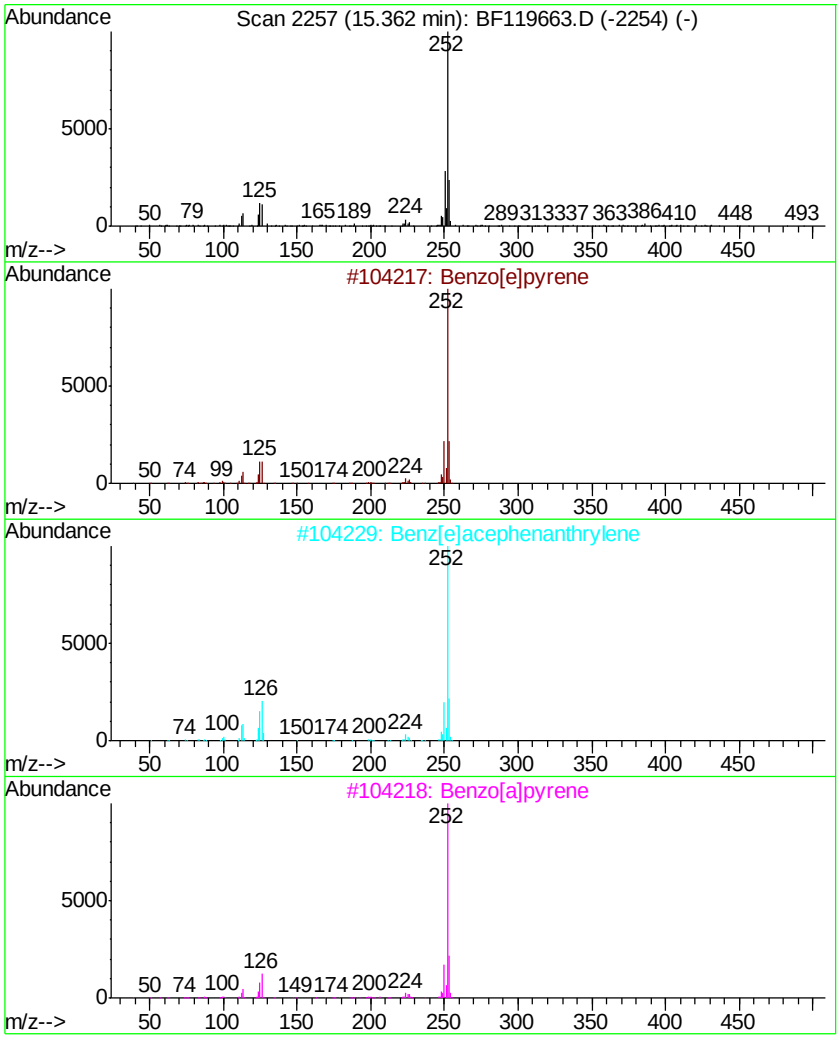
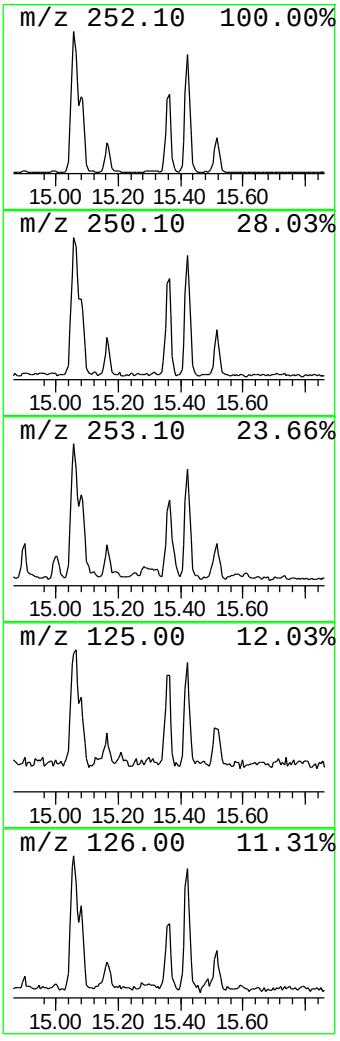
Peak Number 15 Benzo[e]pyrene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.36	7.54 ng	379890	Perylene-d12	15.49

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	94
3	Benzo[a]pyrene	252	C20H12	000050-32-8	94
4	Benzo[j]fluoranthene	252	C20H12	000205-82-3	94
5	Benzo[k]fluoranthene	252	C20H12	000207-08-9	93





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF033120\
Data File : BF119663.D
Acq On : 31 Mar 2020 20:22
Operator : MA/SJ
Sample : L2055-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-7-(1-3)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF031820.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...	5.04	2.4	ng	195149	1	6.83	1636790	20.0
Benzenesulfonamid...	10.77	3.7	ng	285307	4	11.37	1538590	20.0
Undecane, 2-methyl-	10.79	3.3	ng	256419	4	11.37	1538590	20.0
n-Hexadecanoic acid	11.90	10.3	ng	793001	4	11.37	1538590	20.0
Benzaldehyde, 3,5...	11.98	2.5	ng	194289	4	11.37	1538590	20.0
3,5-Dimethyldodecane	12.46	2.3	ng	176522	4	11.37	1538590	20.0
11H-Benzo[b]fluorene	13.14	2.4	ng	266982	5	14.02	2243540	20.0
Fluoranthene, 2-m...	13.21	3.0	ng	337346	5	14.02	2243540	20.0
1-Docosene	13.87	8.6	ng	966646	5	14.02	2243540	20.0
Heptadecane	14.49	3.1	ng	346001	5	14.02	2243540	20.0
Triacontane	14.80	3.1	ng	156856	6	15.49	1007360	20.0
Hexacosane	15.14	4.2	ng	210944	6	15.49	1007360	20.0
Benzo[e]pyrene	15.36	7.5	ng	379890	6	15.49	1007360	20.0