LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061422\
Data File : BF128809.D
Acq On : 14 Jun 2022 19:44
Operator : CG\JU
Sample : N3336-01
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
HD-02-061322

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-BF060222.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF128809.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.416	529	532	547	rBV	2197330	1961411	34.65%	6.805%
2	6.051	636	640	643	rVB	86260	75971	1.34%	0.264%
3	6.433	701	705	710	rBV	2232003	1969847	34.80%	6.835%
4	6.774	760	763	767	rVB	576796	506860	8.95%	1.759%
5	7.345	855	860	863	rBV	1500364	1326654	23.44%	4.603%
6	8.063	975	982	985	rBV	750659	611605	10.80%	2.122%
7	9.139	1160	1165	1168	rBV	2733160	2195792	38.79%	7.619%
8	9.662	1251	1254	1260	rBV2	63668	78726	1.39%	0.273%
9	9.815	1277	1280	1283	rVB	763492	597376	10.55%	2.073%
10	10.609	1411	1415	1418	rBV	1190961	961440	16.98%	3.336%
11	11.304	1530	1533	1535	rBV	626552	505104	8.92%	1.753%
12	11.327	1535	1537	1541	rVV	196705	151234	2.67%	0.525%
13	11.380	1541	1546	1549	rVB	63327	58327	1.03%	0.202%
14	11.704	1597	1601	1604	rBV	1459098	1258903	22.24%	4.368%
15	11.839	1621	1624	1629	rBV2	118246	137706	2.43%	0.478%
16	11.915	1634	1637	1640	rBV2	81665	91859	1.62%	0.319%
17	12.356	1707	1712	1714	rBV2	65604	83195	1.47%	0.289%
18	12.398	1714	1719	1720	rVV	5723921	5660909	100.00%	19.641%
19	12.415	1720	1722	1728	rVV2	4050769	3392662	59.93%	11.771%
20	12.498	1732	1736	1738	rBV	612291	575468	10.17%	1.997%
21	12.521	1738	1740	1743	rVB	499690	381478	6.74%	1.324%
22	12.621	1753	1757	1761	rBV3	58320	91168	1.61%	0.316%
23	12.751	1773	1779	1782	rBV	542259	536369	9.47%	1.861%
24	12.898	1800	1804	1807	rVB	1710450	1465404	25.89%	5.084%
25	12.974	1812	1817	1820	rBV2	41550	72826	1.29%	0.253%
26	13.080	1832	1835	1839	rVB	103192	114112	2.02%	0.396%
27	13.145	1844	1846	1849	rVB	79040	65496	1.16%	0.227%
28	13.180	1849	1852	1857	rBV2	65761	89576	1.58%	0.311%
29	13.274	1866	1868	1871	rBV2	71084	86671	1.53%	0.301%
30	13.398	1884	1889	1893	rVB2	55983	79450	1.40%	0.276%
31	13.797	1954	1957	1962	rVB	101967	83863	1.48%	0.291%
32	13.892	1971	1973	1976	rVB2	70354	61894	1.09%	0.215%
33	13.950	1978	1983	1985	rVV2	688706	859864	15.19%	2.983%
34	13.974	1985	1987	1993	rVB	305301	283119	5.00%	0.982%
35	14.050	1995	2000	2004	rBV2	342172	401942	7.10%	1.395%
36	14.421	2060	2063	2067	rBV2	74847	83298	1.47%	0.289%

LSC Area Percent Report

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Acq On : 14 Jun 2022 19:44
Operator : CG\JU
Sample : N3336-01
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
HD-02-061322

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-BF060222.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	14.868	2136	2139	2143	rVB2	71096	69536	1.23%	0.241%
38	14.986	2155	2159	2162	rBV	284680	434914	7.68%	1.509%
39	15.056	2168	2171	2174	rVB	88845	87240	1.54%	0.303%
40	15.280	2206	2209	2215	rVB	175284	197460	3.49%	0.685%
41	15.344	2216	2220	2224	rVB	271555	321670	5.68%	1.116%
42	15.403	2226	2230	2234	rBV	353956	452612	8.00%	1.570%
43	16.838	2470	2474	2483	rVB	80846	180668	3.19%	0.627%
44	17.274	2544	2548	2556	rVB2	66752	119768	2.12%	0.416%

Sum of corrected areas: 28821447

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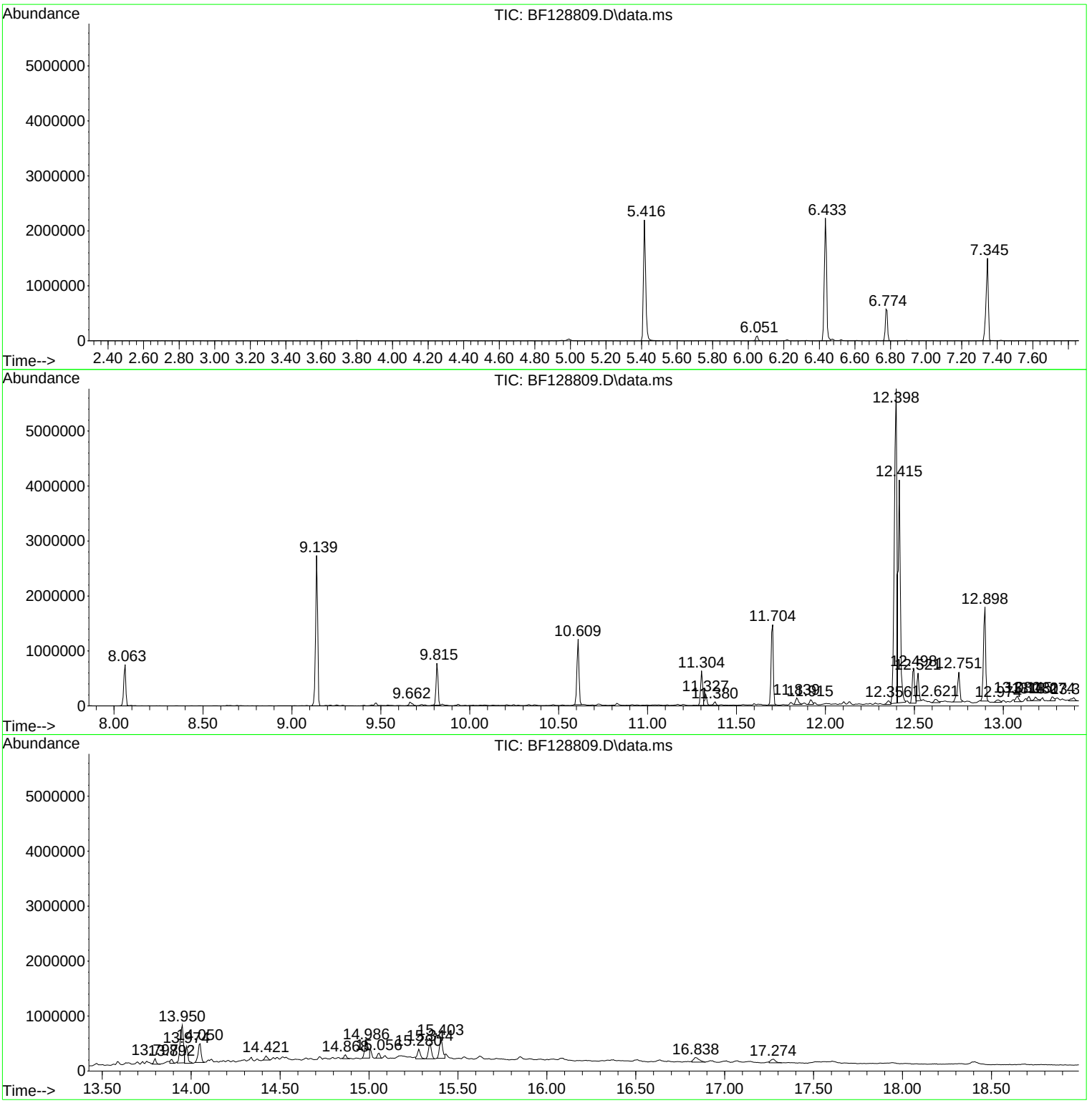
BNA_F
ClientSampleId :
HD-02-061322

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060222.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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Data File : BF128809.D
Acq On : 14 Jun 2022 19:44
Operator : CG\JU
Sample : N3336-01
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Instrument :
BNA_F
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060222.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

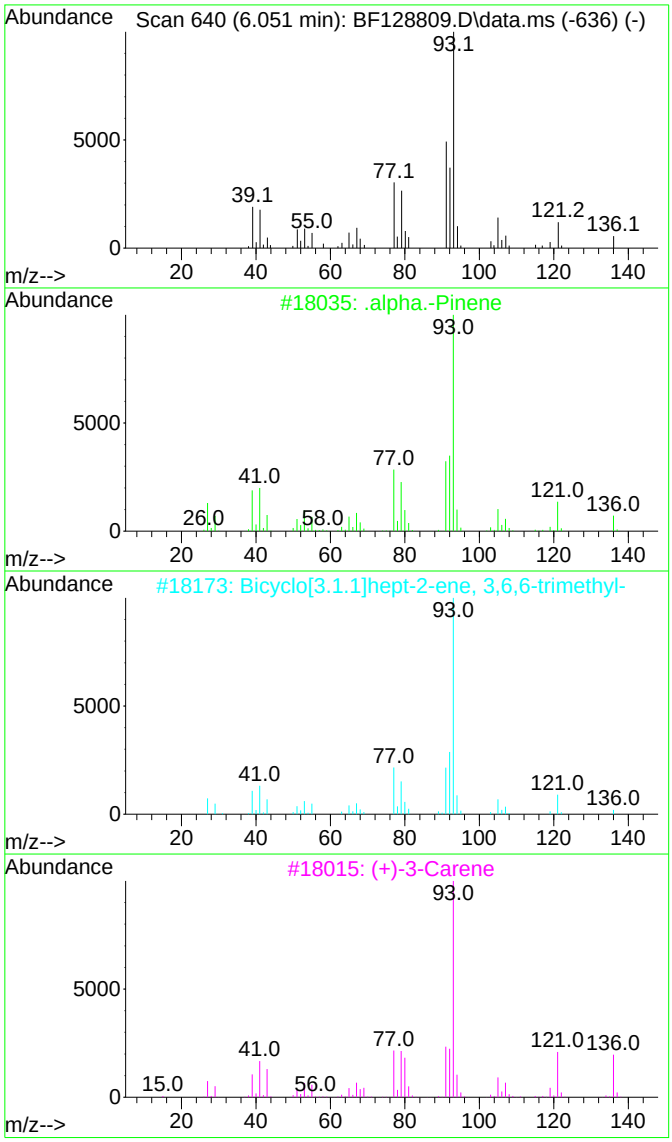
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

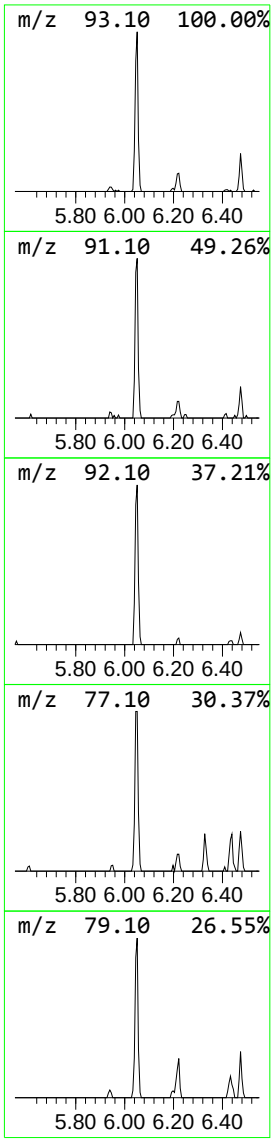
****

Peak Number 1 .alpha.-Pinene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.051	3.00 ng	75971	1,4-Dichlorobenzene-d4	6.780

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.alpha.-Pinene	136	C10H16	000080-56-8	94
2	Bicyclo[3.1.1]hept-2-ene, 3,6,6-...	136	C10H16	004889-83-2	91
3	(+)-3-Carene	136	C10H16	000498-15-7	91
4	Tricyclo[2.2.1.0(2,6)]heptane, 1...	136	C10H16	000508-32-7	91
5	3-Carene	136	C10H16	013466-78-9	91





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Operator : CG\JU
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Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060222.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

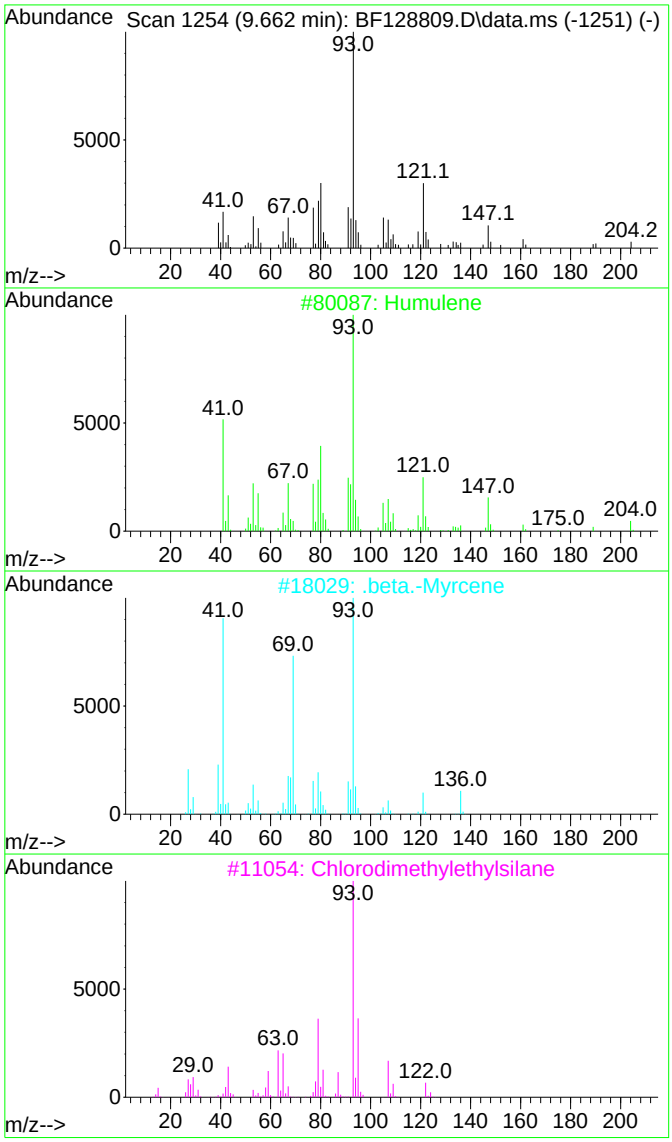
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

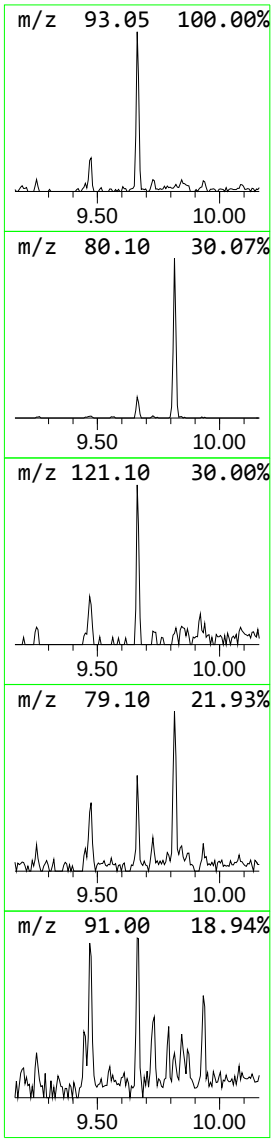
****

Peak Number 2 Humulene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.662	2.64 ng	78726	Acenaphthene-d10	9.815

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Humulene	204	C15H24	006753-98-6	98
2	.beta.-Myrcene	136	C10H16	000123-35-3	64
3	Chlorodimethylethylsilane	122	C4H11ClSi	006917-76-6	53
4	Tricyclo[2.2.1.0(2,6)]heptane, 1...	136	C10H16	000488-97-1	53
5	1,3,7-Octatriene, 3,7-dimethyl-	136	C10H16	000502-99-8	46





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Data File : BF128809.D
Acq On : 14 Jun 2022 19:44
Operator : CG\JU
Sample : N3336-01
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
HD-02-061322

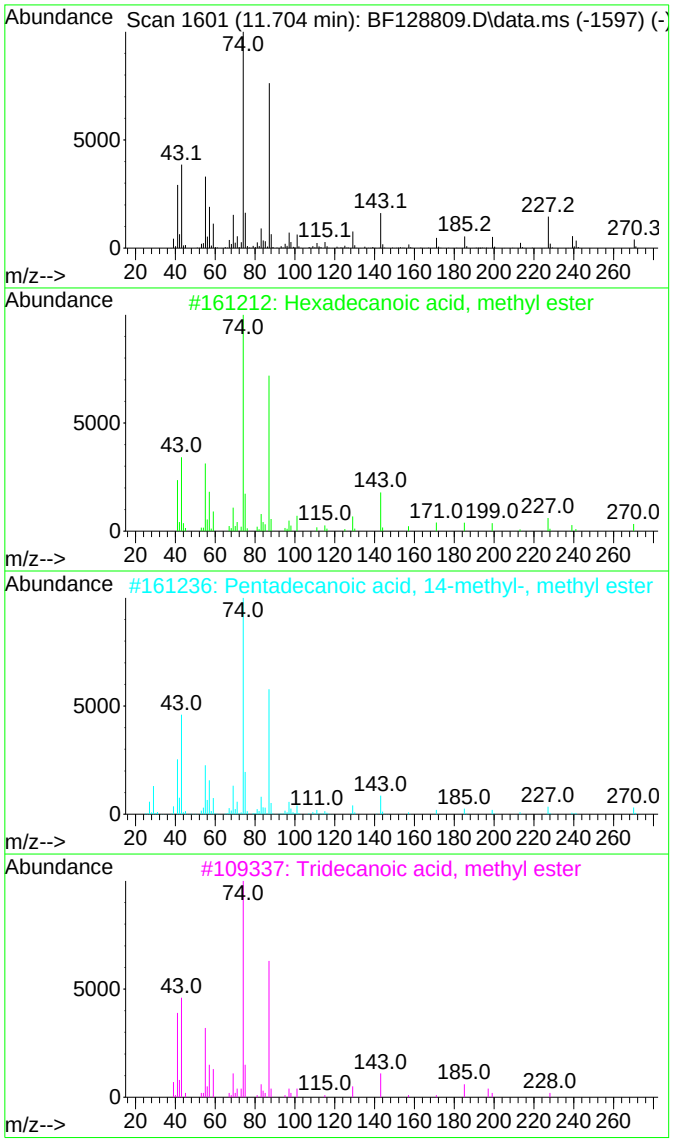
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060222.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

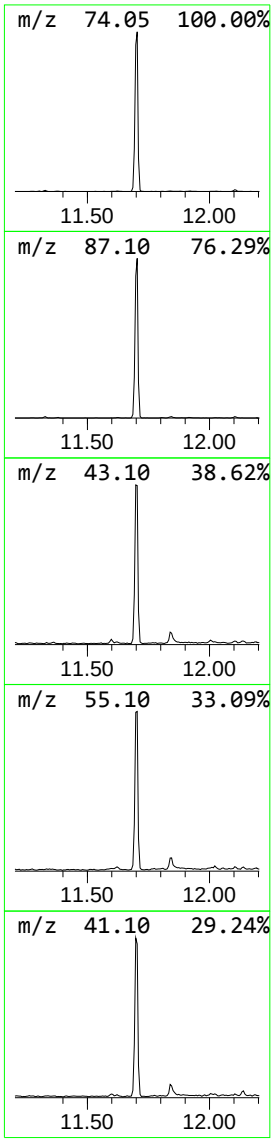
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Hexadecanoic acid, methyl e... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.704	49.85 ng	1258900	Phenanthrene-d10	11.304

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	98
2	Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	93
3	Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	91
4	Methyl tetradecanoate	242	C15H30O2	000124-10-7	87
5	Decanoic acid, methyl ester	186	C11H22O2	000110-42-9	81





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Acq On : 14 Jun 2022 19:44
Operator : CG\JU
Sample : N3336-01
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060222.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

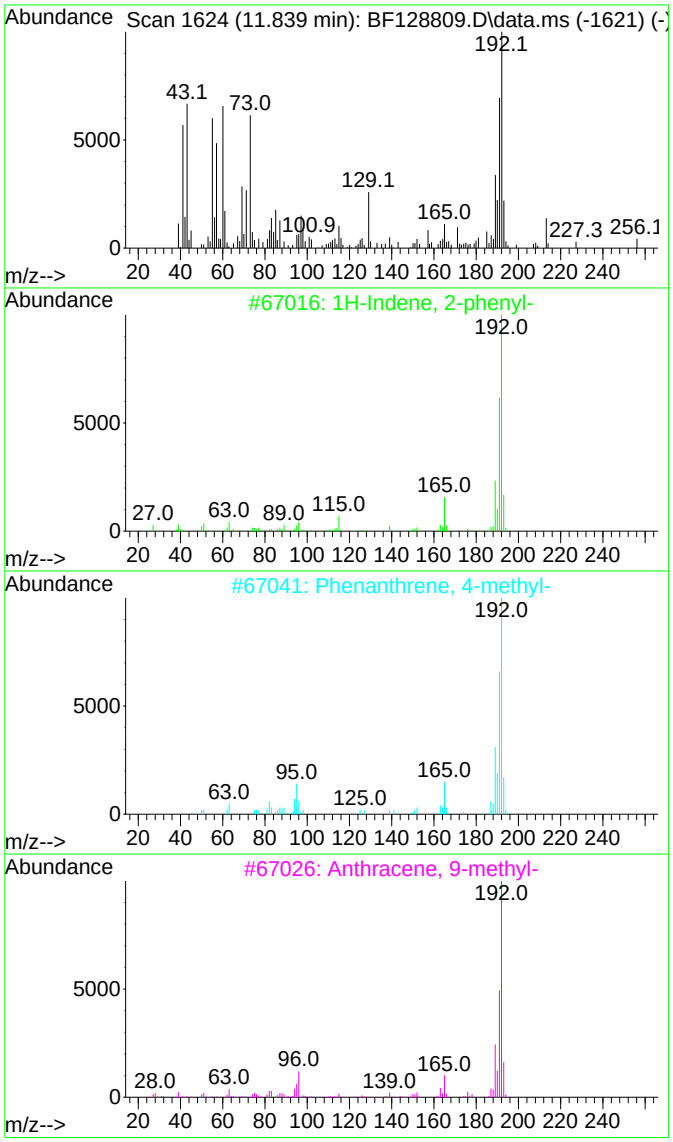
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

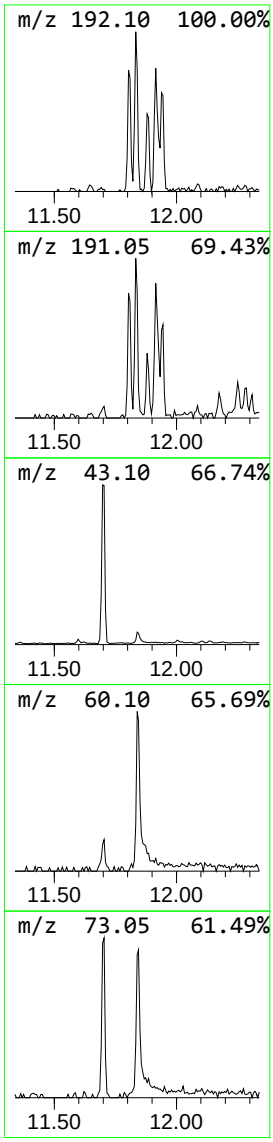
****

Peak Number 4 1H-Indene, 2-phenyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.839	5.45 ng	137706	Phenanthrene-d10	11.304

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	70
2	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	70
3	Anthracene, 9-methyl-	192	C15H12	000779-02-2	55
4	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	55
5	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	55





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Sample : N3336-01
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ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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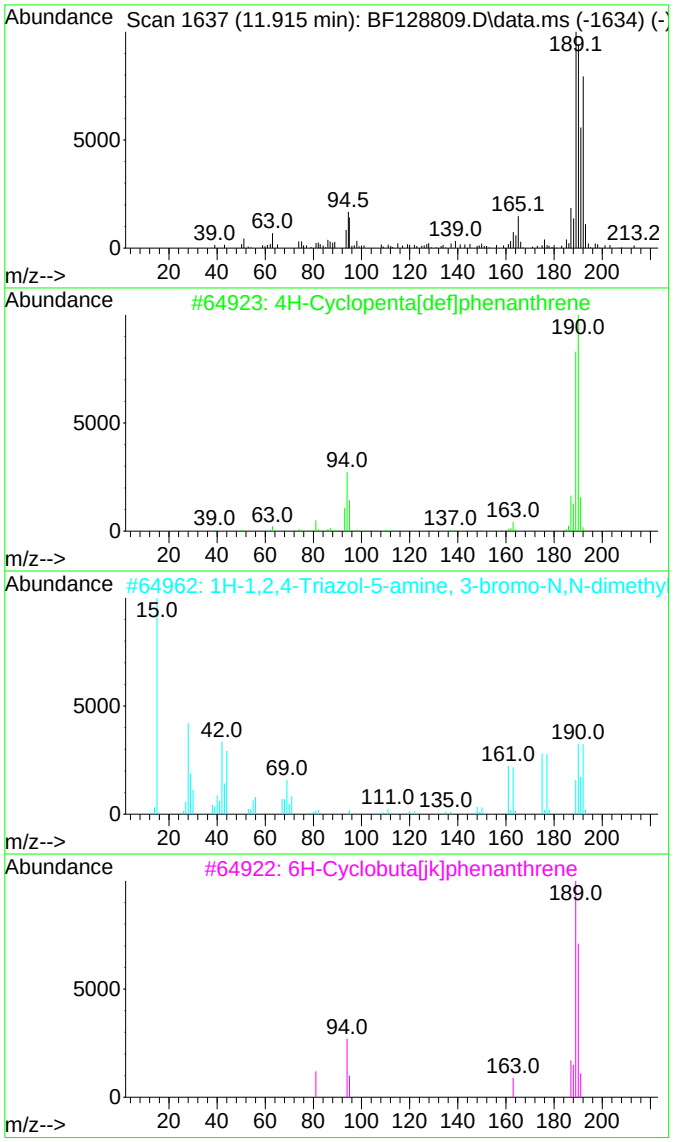
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060222.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

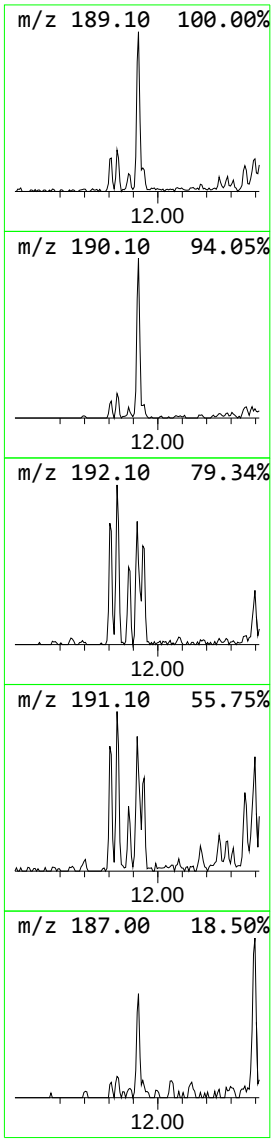
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 4H-Cyclopenta[def]phenanthrene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.915	3.64 ng	91859	Phenanthrene-d10	11.304

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	53
2	1H-1,2,4-Triazol-5-amine, 3-brom...	190	C4H7BrN4	1000460-85-7	53
3	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	47
4	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	46
5	8,9-Dihydrocyclopenta[def]phenan...	192	C15H12	027410-55-5	44





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Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060222.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

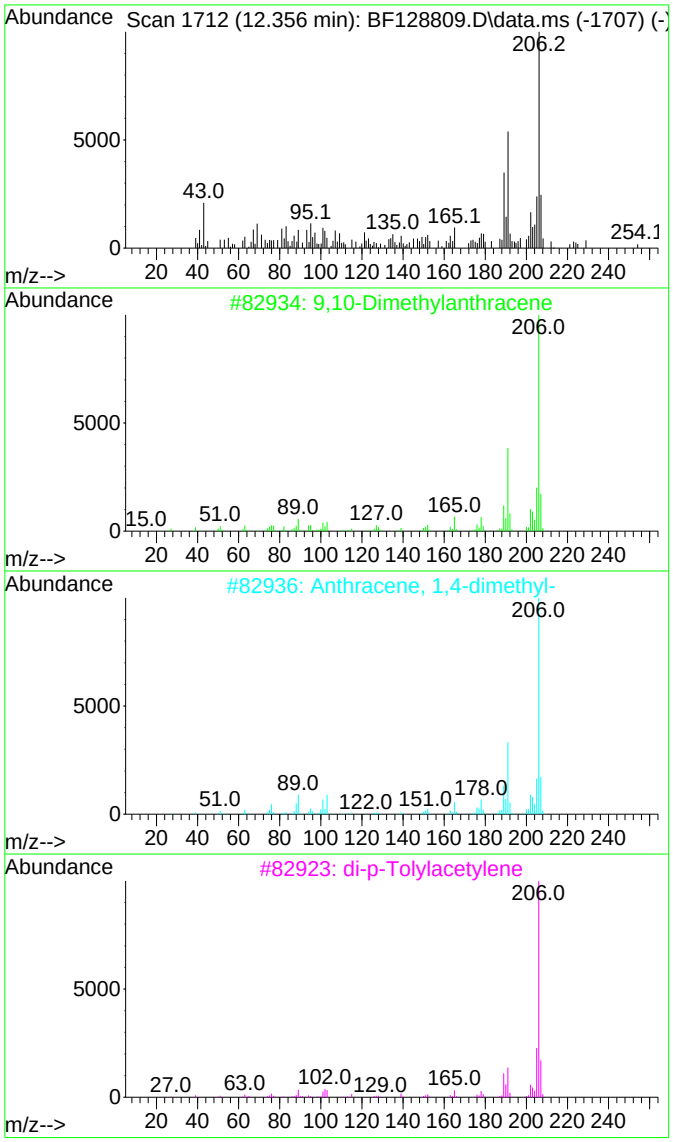
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

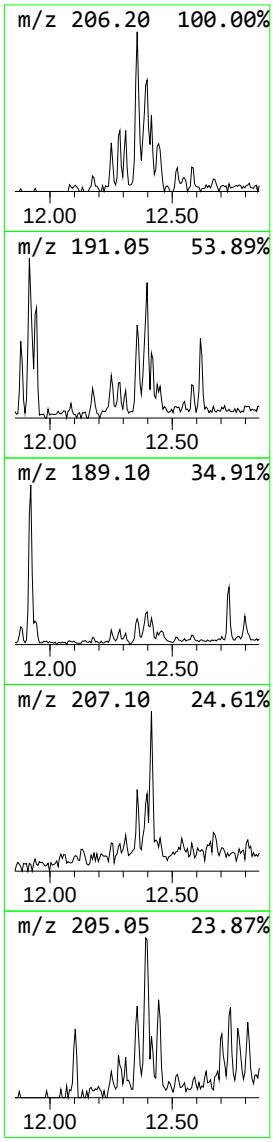
****

Peak Number 6 9,10-Dimethylanthracene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.356	3.29 ng	83195	Phenanthrene-d10	11.304

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Dimethylanthracene	206	C16H14	000781-43-1	91
2	Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	90
3	di-p-Tolylacetylene	206	C16H14	002789-88-0	90
4	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	81
5	3,4-Dimethylphenanthrene	206	C16H14	1000452-24-5	81





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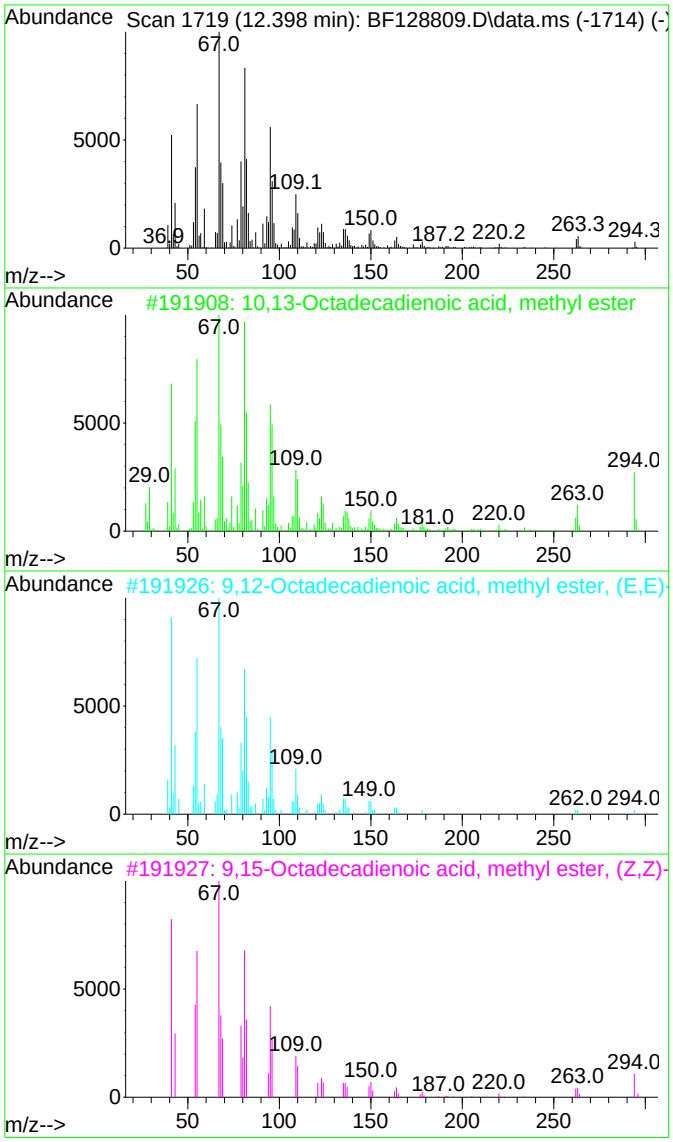
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

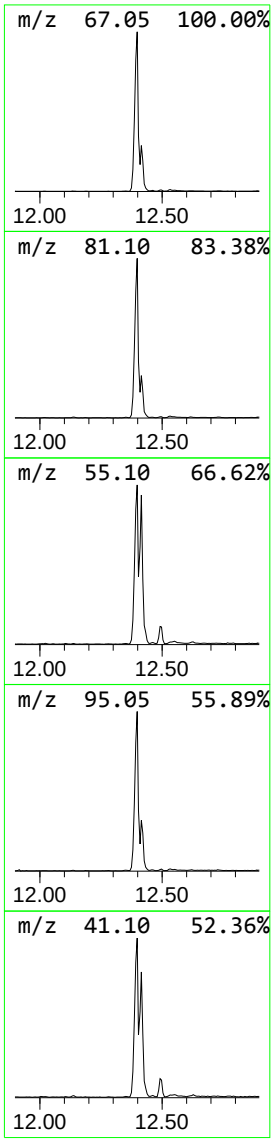
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 10,13-Octadecadienoic acid,... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.398	224.15 ng	5660910	Phenanthrene-d10	11.304

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	10,13-Octadecadienoic acid, meth...	294	C19H34O2	056554-62-2	99
2	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002566-97-4	99
3	9,15-Octadecadienoic acid, methy...	294	C19H34O2	017309-05-6	99
4	Methyl 9-cis,11-trans-octadecadi...	294	C19H34O2	013058-52-1	99
5	9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	99





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Sample : N3336-01
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
HD-02-061322

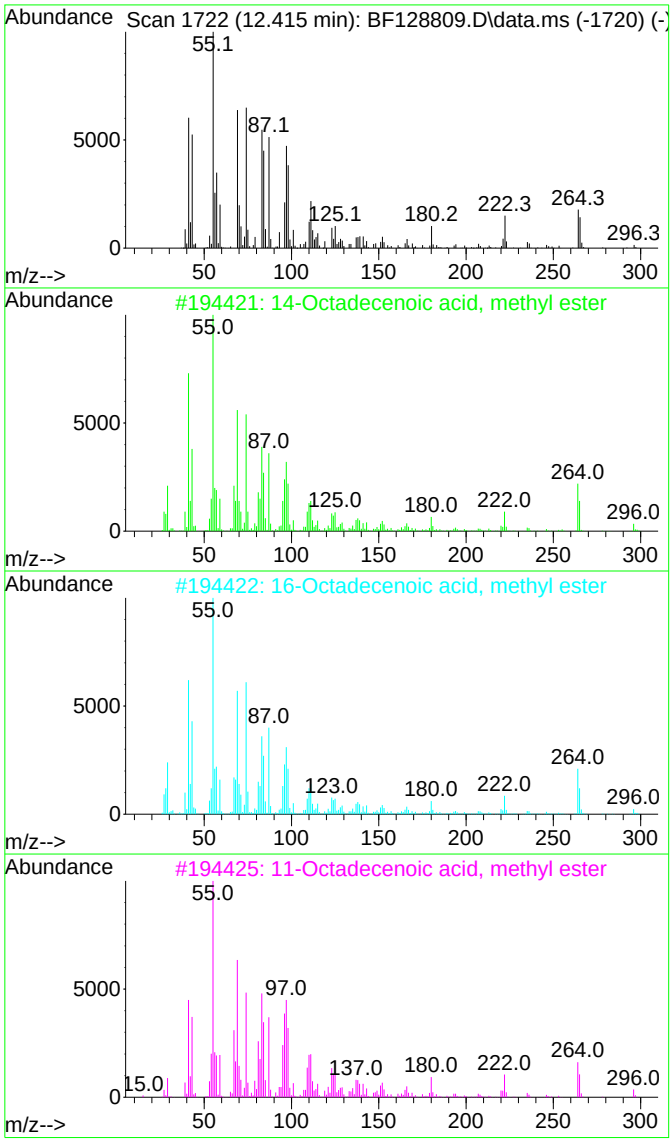
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060222.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

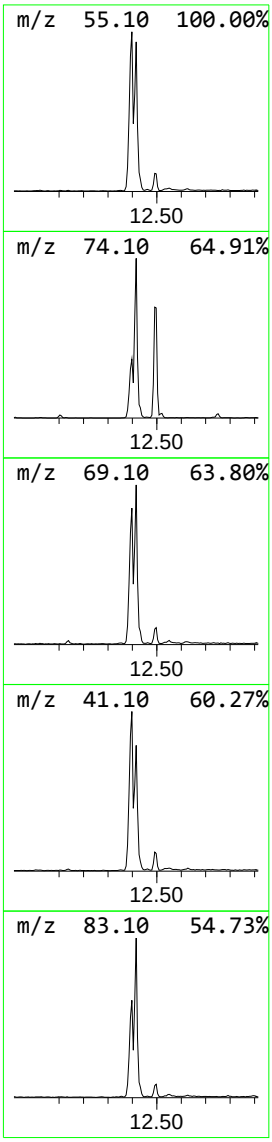
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 14-Octadecenoic acid, methy... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.415	134.34 ng	3392660	Phenanthrene-d10	11.304

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	14-Octadecenoic acid, methyl ester	296	C19H36O2	056554-48-4	91
2	16-Octadecenoic acid, methyl ester	296	C19H36O2	056554-49-5	91
3	11-Octadecenoic acid, methyl ester	296	C19H36O2	052380-33-3	87
4	13-Octadecenoic acid, methyl ester	296	C19H36O2	056554-47-3	87
5	trans-13-Octadecenoic acid, meth...	296	C19H36O2	1000333-61-3	86





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Data File : BF128809.D
Acq On : 14 Jun 2022 19:44
Operator : CG\JU
Sample : N3336-01
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
HD-02-061322

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060222.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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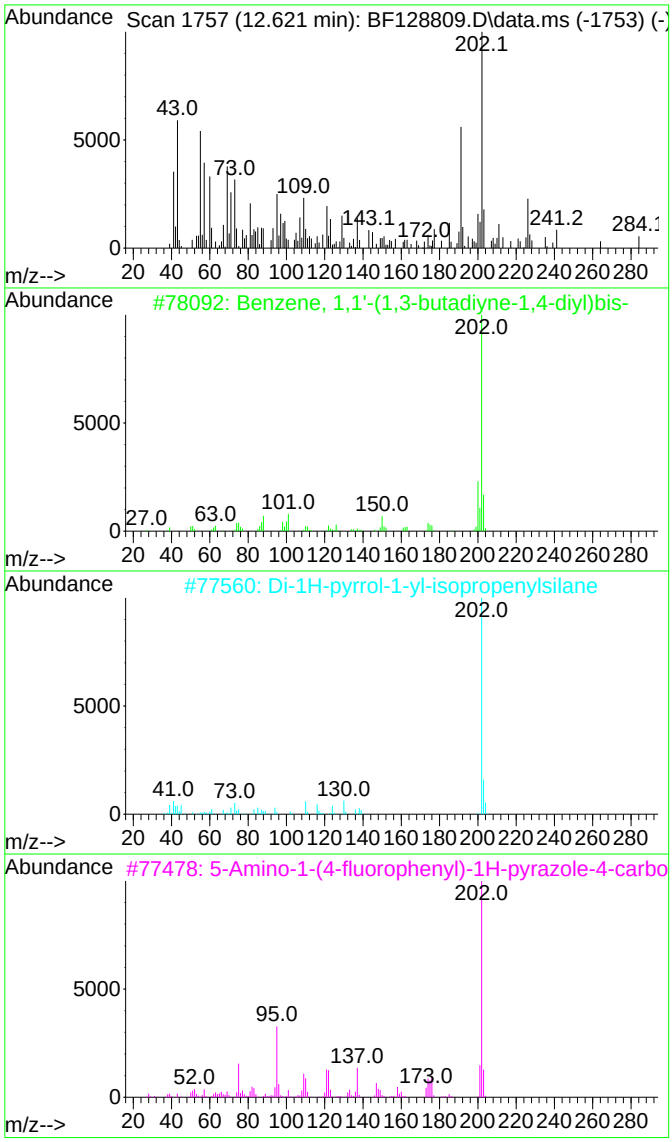
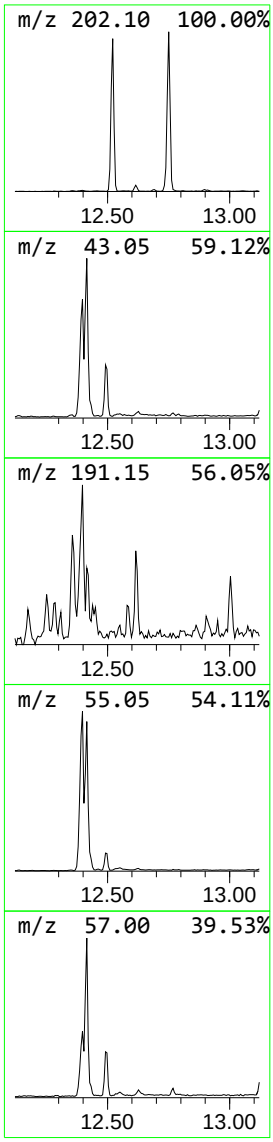
Peak Number 9 unknown12.621 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.621	3.61 ng	91168	Phenanthrene-d10	11.304

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,1'-(1,3-butadiyne-1,4...	202	C16H10	000886-66-8	35
2	Di-1H-pyrrol-1-yl-isopropenylsilane	202	C11H14N2Si	1000306-57-7	30
3	5-Amino-1-(4-fluorophenyl)-1H-py...	202	C10H7FN4	051516-70-2	27
4	Phenol, 4,4'-oxybis-	202	C12H10O3	001965-09-9	22
5	Fluoranthene	202	C16H10	000206-44-0	18





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Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
HD-02-061322

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060222.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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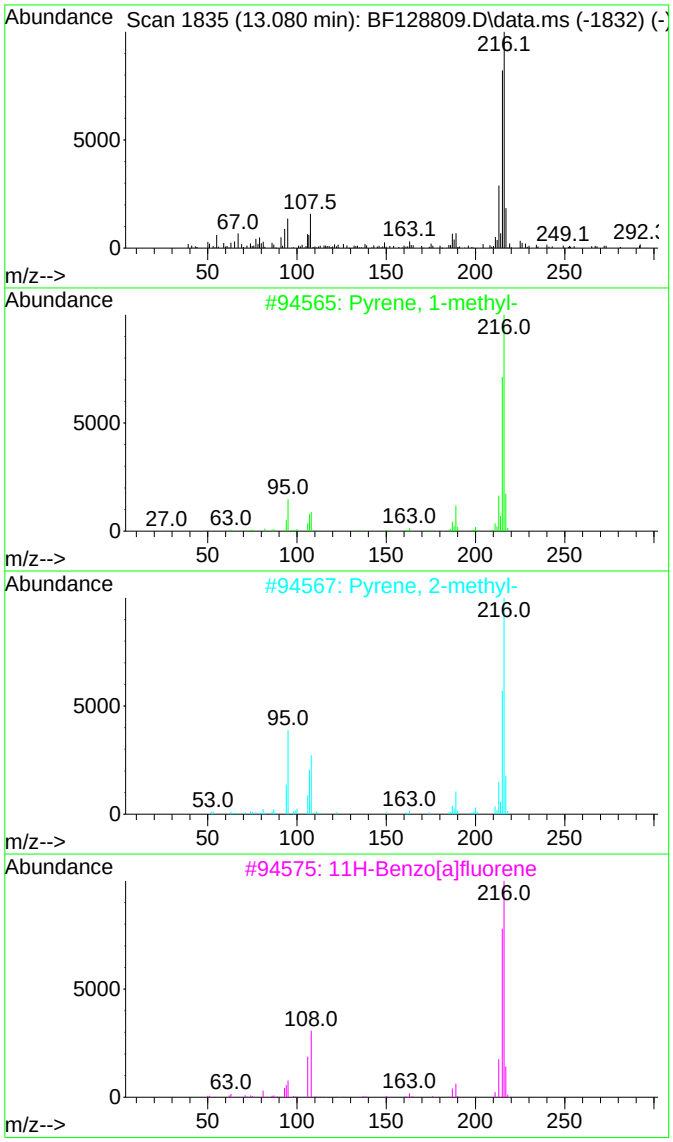
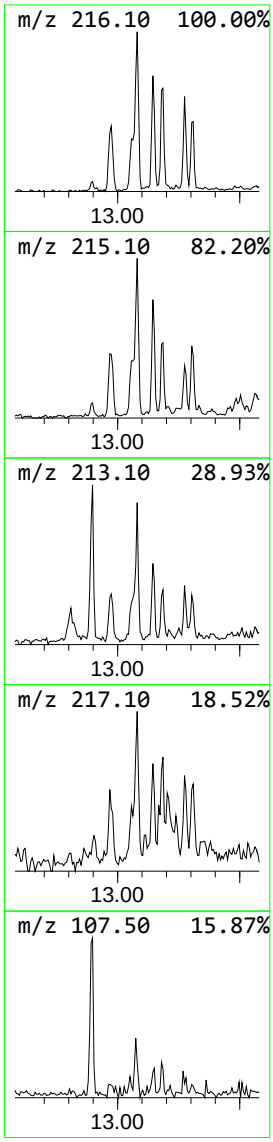
Peak Number 10 Pyrene, 1-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.080	2.65 ng	114112	Chrysene-d12	13.950

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-	216	C17H12	002381-21-7	91
2	Pyrene, 2-methyl-	216	C17H12	003442-78-2	90
3	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	81
4	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	81
5	7H-Benzo[c]fluorene	216	C17H12	000205-12-9	74





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Sample : N3336-01
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
HD-02-061322

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060222.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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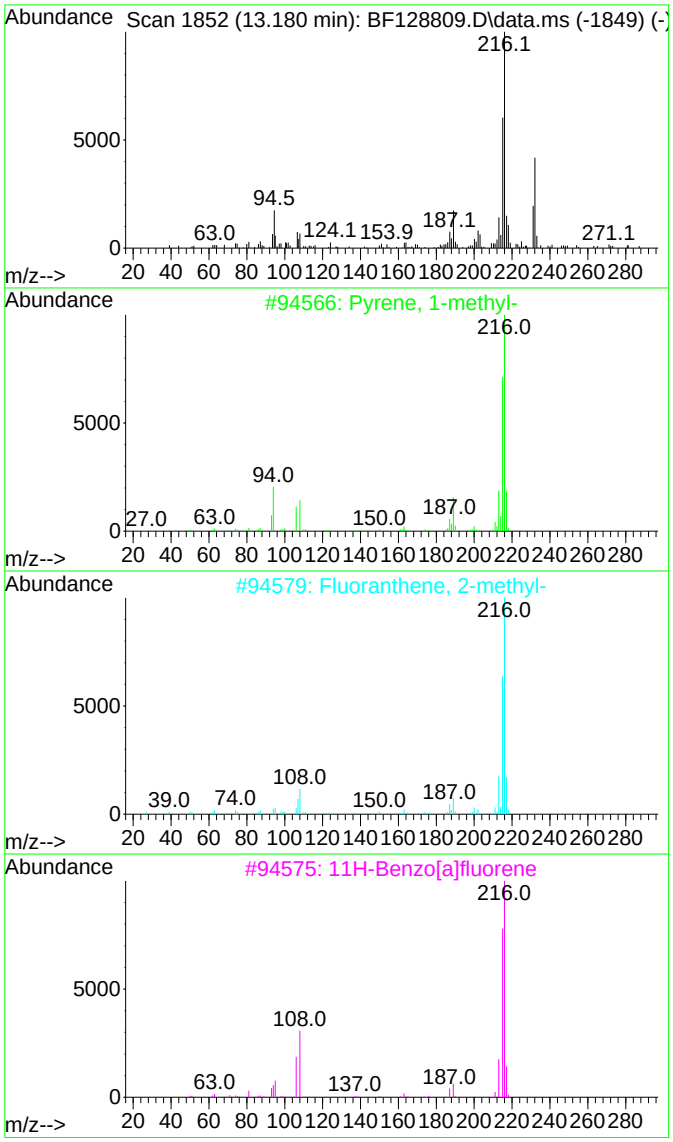
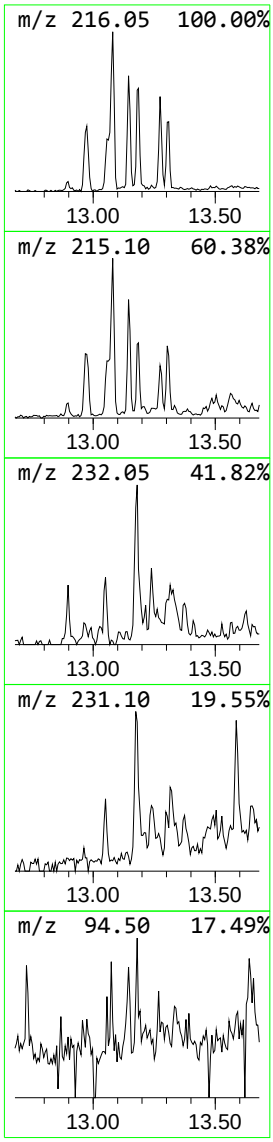
Peak Number 11 Fluoranthene, 2-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.180	2.08 ng	89576	Chrysene-d12	13.950

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-	216	C17H12	002381-21-7	78
2	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	78
3	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	60
4	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	58
5	Pyrene, 2-methyl-	216	C17H12	003442-78-2	49





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Operator : CG\JU
Sample : N3336-01
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
HD-02-061322

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060222.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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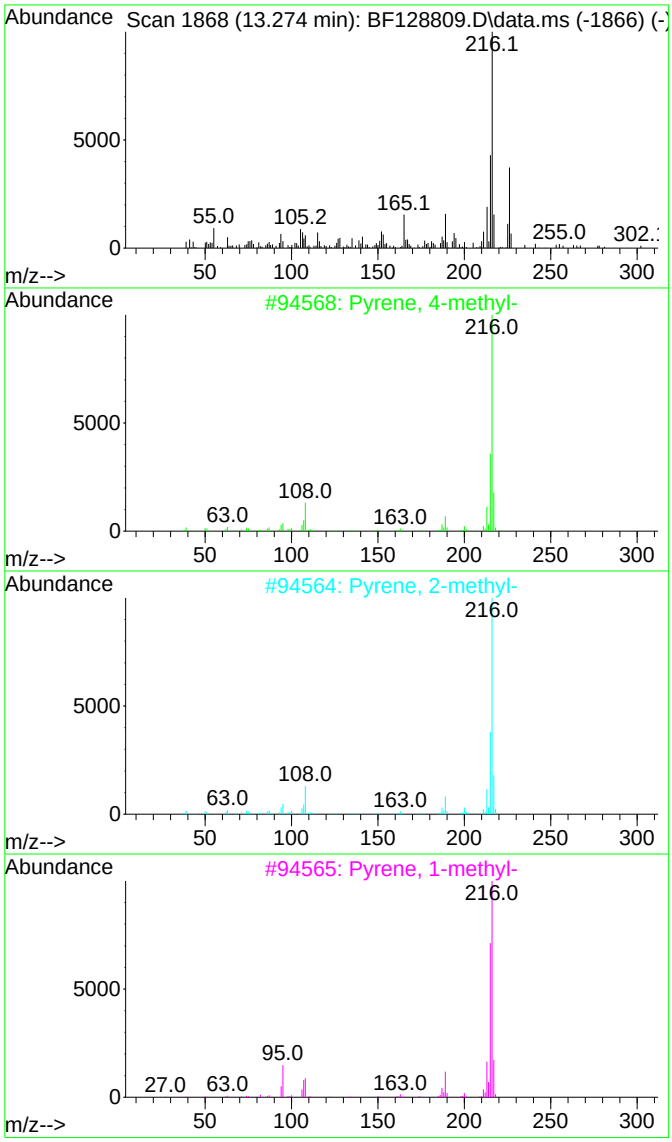
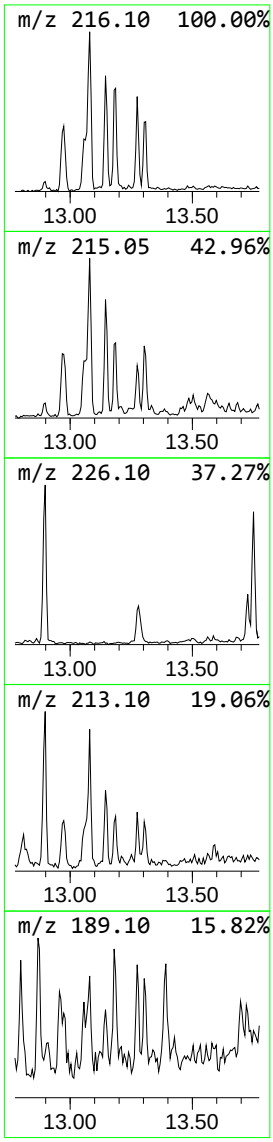
Peak Number 12 Pyrene, 4-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.274	2.02 ng	86671	Chrysene-d12	13.950

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 4-methyl-	216	C17H12	003353-12-6	62
2	Pyrene, 2-methyl-	216	C17H12	003442-78-2	62
3	Pyrene, 1-methyl-	216	C17H12	002381-21-7	53
4	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	52
5	1,2-Difluorostilbene	216	C14H10F2	000643-76-5	50





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Acq On : 14 Jun 2022 19:44
Operator : CG\JU
Sample : N3336-01
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
HD-02-061322

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060222.M
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TIC Library : C:\Database\NIST20.L
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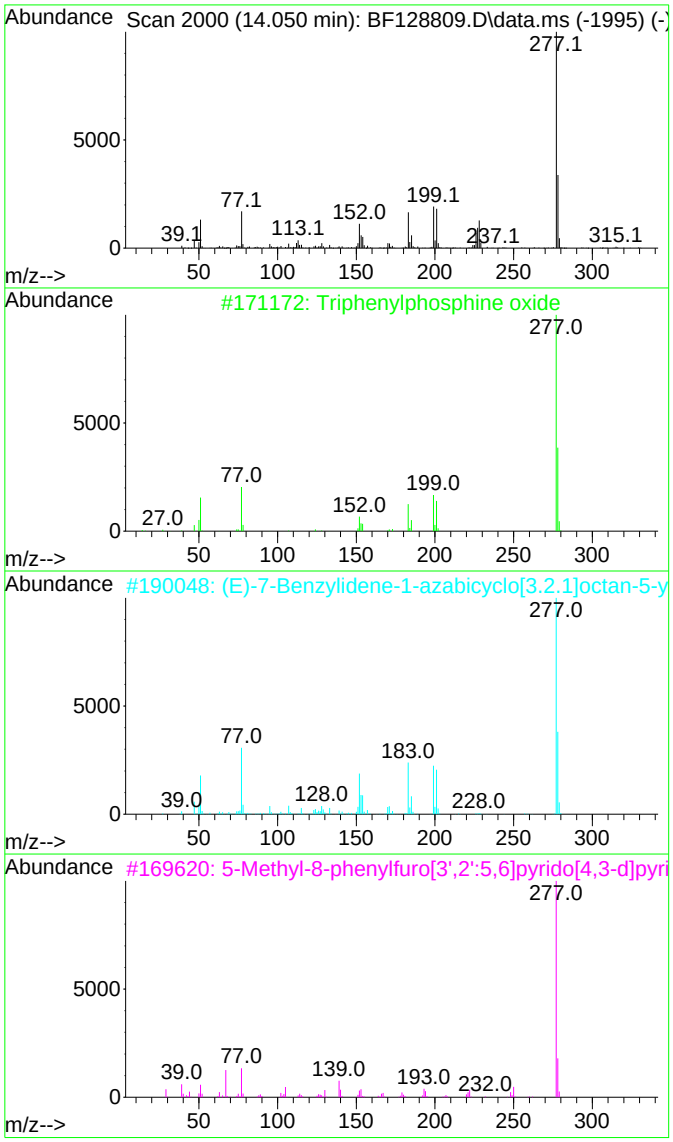
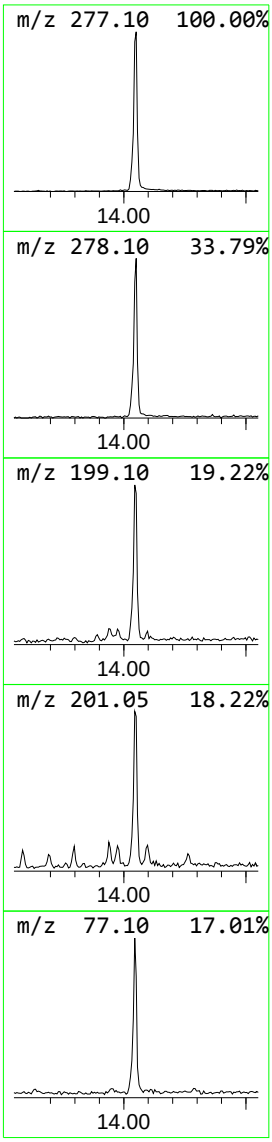
Peak Number 13 Triphenylphosphine oxide Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.050	9.35 ng	401942	Chrysene-d12	13.950

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Triphenylphosphine oxide	278	C18H15OP	000791-28-6	95
2	(E)-7-Benzylidene-1-azabicyclo[3...	293	C15H19NO3S	1000483-83-4	46
3	5-Methyl-8-phenylfuro[3',2':5,6]...	277	C16H11N3O2	1000460-07-9	43
4	2-Phenyl-6-(trifluoromethyl)-1H-...	277	C15H10F3NO	1000387-59-3	38
5	Cobalt,(.eta.<5>-2,4-cyclopentad...	277	C16H15BCo	036682-12-9	37





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Sample : N3336-01
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Instrument :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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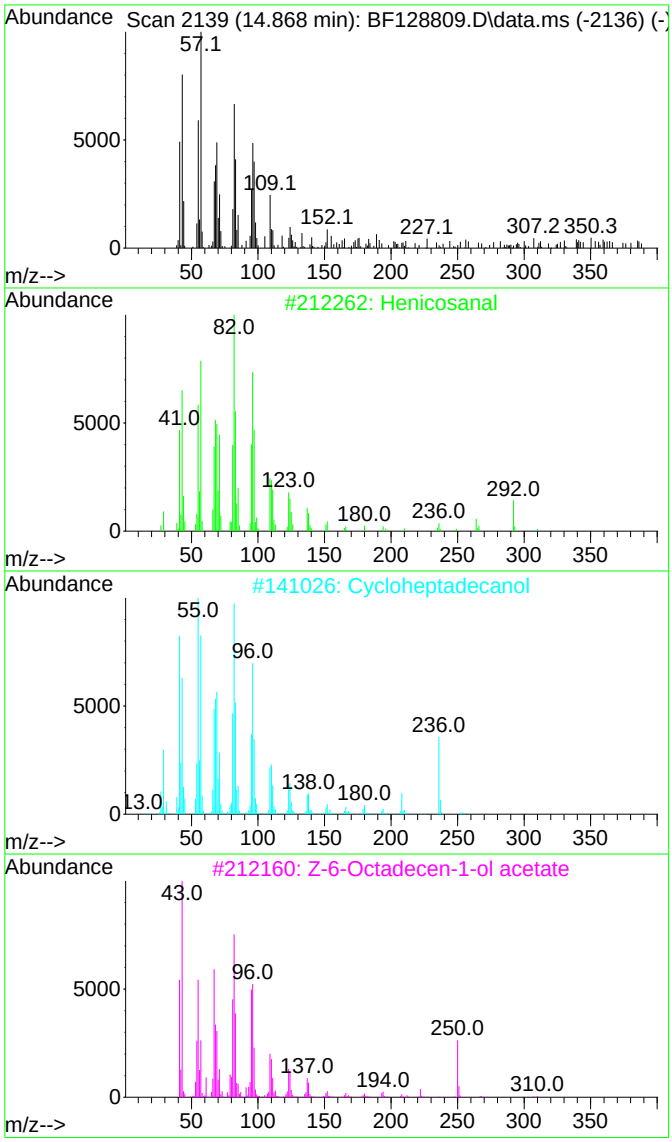
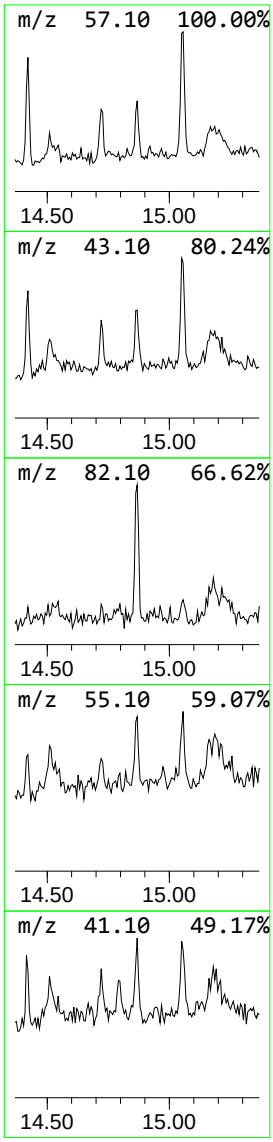
Peak Number 14 Henicosanal Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.868	3.07 ng	69536	Perylene-d12	15.403

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Henicosanal	310	C21H42O	051227-32-8	93
2	Cycloheptadecanol	254	C17H34O	004429-77-0	72
3	Z-6-Octadecen-1-ol acetate	310	C20H38O2	1000131-09-9	72
4	Pentadecanal-	226	C15H30O	002765-11-9	70
5	Z-14-Octadecen-1-ol acetate	310	C20H38O2	1010131-07-6	64





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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
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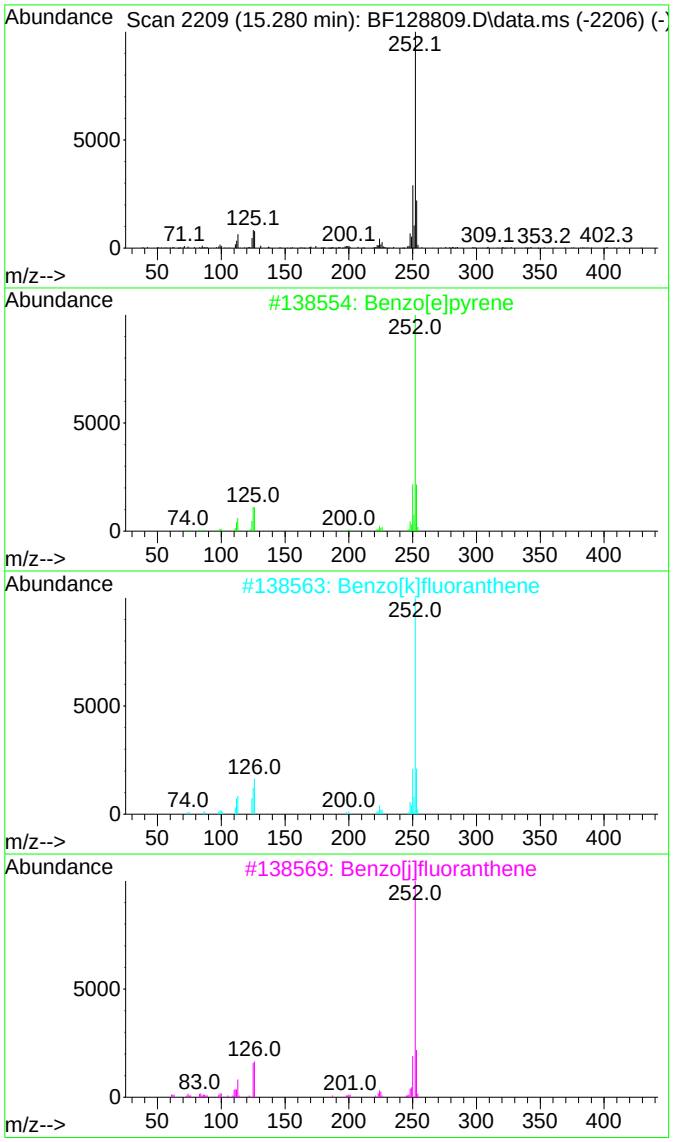
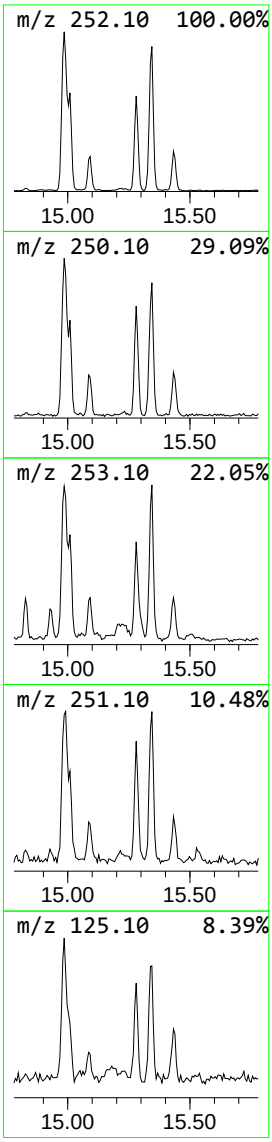
Peak Number 15 Benzo[e]pyrene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.280	8.73 ng	197460	Perylene-d12	15.403

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[e]pyrene	252	C20H12	000192-97-2	96
2	Benzo[k]fluoranthene	252	C20H12	000207-08-9	96
3	Benzo[j]fluoranthene	252	C20H12	000205-82-3	94
4	Benzo[b]fluoranthene	252	C20H12	000205-99-2	90
5	Benzo[a]pyrene	252	C20H12	000050-32-8	90





Tentatively Identified Compound (LSC) summary

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Data File : BF128809.D
Acq On : 14 Jun 2022 19:44
Operator : CG\JU
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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
.alpha.-Pinene	6.051	3.0	ng	75971	1	6.780	506860	20.0
Humulene	9.662	2.6	ng	78726	3	9.815	597376	20.0
Hexadecanoic ac...	11.704	49.9	ng	1258900	4	11.304	505104	20.0
1H-Indene, 2-ph...	11.839	5.5	ng	137706	4	11.304	505104	20.0
4H-Cyclopenta[d...	11.915	3.6	ng	91859	4	11.304	505104	20.0
9,10-Dimethylan...	12.356	3.3	ng	83195	4	11.304	505104	20.0
10,13-Octadecad...	12.398	224.2	ng	5660910	4	11.304	505104	20.0
14-Octadecenoic...	12.415	134.3	ng	3392660	4	11.304	505104	20.0
unknown12.621	12.621	3.6	ng	91168	4	11.304	505104	20.0
Pyrene, 1-methyl-	13.080	2.6	ng	114112	5	13.950	859864	20.0
Fluoranthene, 2...	13.180	2.1	ng	89576	5	13.950	859864	20.0
Pyrene, 4-methyl-	13.274	2.0	ng	86671	5	13.950	859864	20.0
Triphenylphosph...	14.050	9.3	ng	401942	5	13.950	859864	20.0
Henicosanal	14.868	3.1	ng	69536	6	15.403	452612	20.0
Benzo[e]pyrene	15.280	8.7	ng	197460	6	15.403	452612	20.0