LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF052422\
Data File : BF128436.D
Acq On : 24 May 2022 15:48
Operator : CG\JU
Sample : N2981-02
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-13(3-3.5)

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-BF052022.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF128436.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.569	384	388	397	rBV	132146	153658	7.53%	0.710%
2	5.081	472	475	490	rVB	43482	66455	3.26%	0.307%
3	5.469	538	541	561	rBV	1737567	1815526	88.95%	8.391%
4	6.481	709	713	726	rBV	2003598	1906020	93.39%	8.810%
5	6.587	729	731	742	rVB	27666	38876	1.90%	0.180%
6	6.845	771	775	782	rBV	640092	575776	28.21%	2.661%
7	7.410	867	871	879	rBV	1413397	1333748	65.35%	6.165%
8	8.128	989	993	996	rBV	851543	765910	37.53%	3.540%
9	8.845	1111	1115	1121	rBV	25573	24412	1.20%	0.113%
10	9.198	1171	1175	1179	rBV	2099506	2040971	100.00%	9.433%
11	9.269	1183	1187	1191	rBV	28344	24855	1.22%	0.115%
12	9.539	1230	1233	1236	rBV	22216	22268	1.09%	0.103%
13	9.745	1264	1268	1272	rBV	49158	45460	2.23%	0.210%
14	9.881	1286	1291	1295	rBV	936069	868179	42.54%	4.013%
15	10.092	1323	1327	1330	rVV3	21048	25510	1.25%	0.118%
16	10.298	1355	1362	1366	rBV3	23974	39314	1.93%	0.182%
17	10.439	1381	1386	1391	rVB4	18697	31523	1.54%	0.146%
18	10.675	1423	1426	1433	rBV	1307996	1256612	61.57%	5.808%
19	10.775	1440	1443	1449	rBV4	20467	33633	1.65%	0.155%
20	10.986	1472	1479	1481	rBV3	19713	26335	1.29%	0.122%
21	11.110	1496	1500	1502	rBV2	20897	27011	1.32%	0.125%
22	11.186	1508	1513	1516	rBV2	29356	34290	1.68%	0.158%
23	11.222	1516	1519	1522	rBV	31226	27503	1.35%	0.127%
24	11.375	1542	1545	1548	rBV	1100024	942132	46.16%	4.355%
25	11.398	1548	1549	1554	rVB	320217	213962	10.48%	0.989%
26	11.451	1554	1558	1563	rBV	58534	73481	3.60%	0.340%
27	11.616	1581	1586	1589	rBV3	39608	54633	2.68%	0.253%
28	11.710	1599	1602	1607	rVB2	15889	24906	1.22%	0.115%
29	11.751	1607	1609	1612	rBV	25567	20505	1.00%	0.095%
30	11.880	1628	1631	1633	rBV	85102	93304	4.57%	0.431%
31	11.904	1633	1635	1641	rVV	133688	138629	6.79%	0.641%
32	11.957	1641	1644	1646	rVV	39560	41213	2.02%	0.190%
33	11.992	1646	1650	1652	rVV2	103345	123850	6.07%	0.572%
34	12.016	1652	1654	1658	rVB	60383	49870	2.44%	0.231%
35	12.174	1677	1681	1685	rBV	64082	61142	3.00%	0.283%
36	12.216	1685	1688	1691	rVB2	31913	30527	1.50%	0.141%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF052422\
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Operator : CG\JU
Sample : N2981-02
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-13(3-3.5)

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-BF052022.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	12.322	1701	1706	1709	rBV2	29298	39409	1.93%	0.182%
38	12.351	1709	1711	1714	rVB	35986	31337	1.54%	0.145%
39	12.427	1720	1724	1726	rBV	86916	84957	4.16%	0.393%
40	12.486	1732	1734	1736	rVB	38716	28311	1.39%	0.131%
41	12.527	1736	1741	1748	rVB3	40476	70184	3.44%	0.324%
42	12.592	1748	1752	1756	rBV	749215	641094	31.41%	2.963%
43	12.633	1756	1759	1761	rVV2	30455	27314	1.34%	0.126%
44	12.657	1761	1763	1766	rVB2	24845	22646	1.11%	0.105%
45	12.692	1766	1769	1772	rBV	41222	42584	2.09%	0.197%
46	12.745	1774	1778	1780	rBV2	29915	30449	1.49%	0.141%
47	12.804	1785	1788	1789	rBV2	116676	99915	4.90%	0.462%
48	12.822	1789	1791	1797	rVB	558539	510096	24.99%	2.358%
49	12.874	1797	1800	1803	rVB	56647	56064	2.75%	0.259%
50	12.963	1807	1815	1818	rBV	2073323	1948135	95.45%	9.004%
51	13.033	1823	1827	1834	rBV4	61090	113423	5.56%	0.524%
52	13.092	1834	1837	1839	rBV3	24029	23401	1.15%	0.108%
53	13.127	1839	1843	1845	rBV3	44683	64605	3.17%	0.299%
54	13.151	1845	1847	1849	rVB	92803	70978	3.48%	0.328%
55	13.216	1856	1858	1861	rBV	57966	49327	2.42%	0.228%
56	13.257	1861	1865	1868	rVB2	72714	82607	4.05%	0.382%
57	13.351	1878	1881	1883	rBV	41005	36871	1.81%	0.170%
58	13.380	1883	1886	1890	rVV3	38720	40285	1.97%	0.186%
59	13.516	1906	1909	1913	rBV4	37510	44280	2.17%	0.205%
60	13.574	1917	1919	1926	rVB6	27409	44135	2.16%	0.204%
61	13.633	1926	1929	1932	rBV4	21125	26400	1.29%	0.122%
62	13.669	1932	1935	1938	rVV	73909	68020	3.33%	0.314%
63	13.774	1948	1953	1955	rBV	78243	78792	3.86%	0.364%
64	13.798	1955	1957	1960	rVB	51271	39591	1.94%	0.183%
65	13.845	1960	1965	1967	rBV2	55229	77437	3.79%	0.358%
66	13.880	1967	1971	1980	rVV3	124745	214762	10.52%	0.993%
67	14.021	1990	1995	1998	rBV2	721814	925240	45.33%	4.276%
68	14.051	1998	2000	2004	rVV	309228	254874	12.49%	1.178%
69	14.116	2007	2011	2013	rBV	47194	54824	2.69%	0.253%
70	14.163	2016	2019	2020	rBV	36312	30357	1.49%	0.140%
71	14.186	2020	2023	2027	rVB3	40249	46635	2.28%	0.216%
72	14.374	2053	2055	2057	rBV	24363	26182	1.28%	0.121%
73	14.416	2057	2062	2064	rVB	79917	94496	4.63%	0.437%
74	14.445	2064	2067	2069	rBV	38967	35834	1.76%	0.166%
75	14.539	2080	2083	2085	rBV	41023	38502	1.89%	0.178%
76	14.745	2115	2118	2119	rBV2	20757	22571	1.11%	0.104%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF052422\
Data File : BF128436.D
Acq On : 24 May 2022 15:48
Operator : CG\JU
Sample : N2981-02
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-13(3-3.5)

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

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Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052022.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

77	14.768	2120	2122	2126	rVB3	37605	36110	1.77%	0.167%
78	14.915	2144	2147	2150	rBV	45265	44632	2.19%	0.206%
79	15.074	2170	2174	2177	rBV	280682	394809	19.34%	1.825%
80	15.186	2190	2193	2197	rBV	68387	73332	3.59%	0.339%
81	15.274	2205	2208	2209	rBV3	22580	24181	1.18%	0.112%
82	15.380	2222	2226	2233	rVB	161236	201864	9.89%	0.933%
83	15.445	2233	2237	2242	rBV	231836	288944	14.16%	1.336%
84	15.510	2243	2248	2251	rVV	448518	582248	28.53%	2.691%
85	15.539	2251	2253	2260	rVB2	60057	96325	4.72%	0.445%
86	15.686	2275	2278	2282	rBV4	17017	23641	1.16%	0.109%
87	16.080	2343	2345	2350	rVB2	15370	21010	1.03%	0.097%
88	17.009	2497	2503	2512	rVB	125528	257265	12.61%	1.189%
89	17.192	2529	2534	2540	rBV	29759	62146	3.04%	0.287%
90	17.468	2575	2581	2592	rBV	81234	175257	8.59%	0.810%
91	17.727	2621	2625	2635	rVB2	25795	64820	3.18%	0.300%

Sum of corrected areas: 21635507

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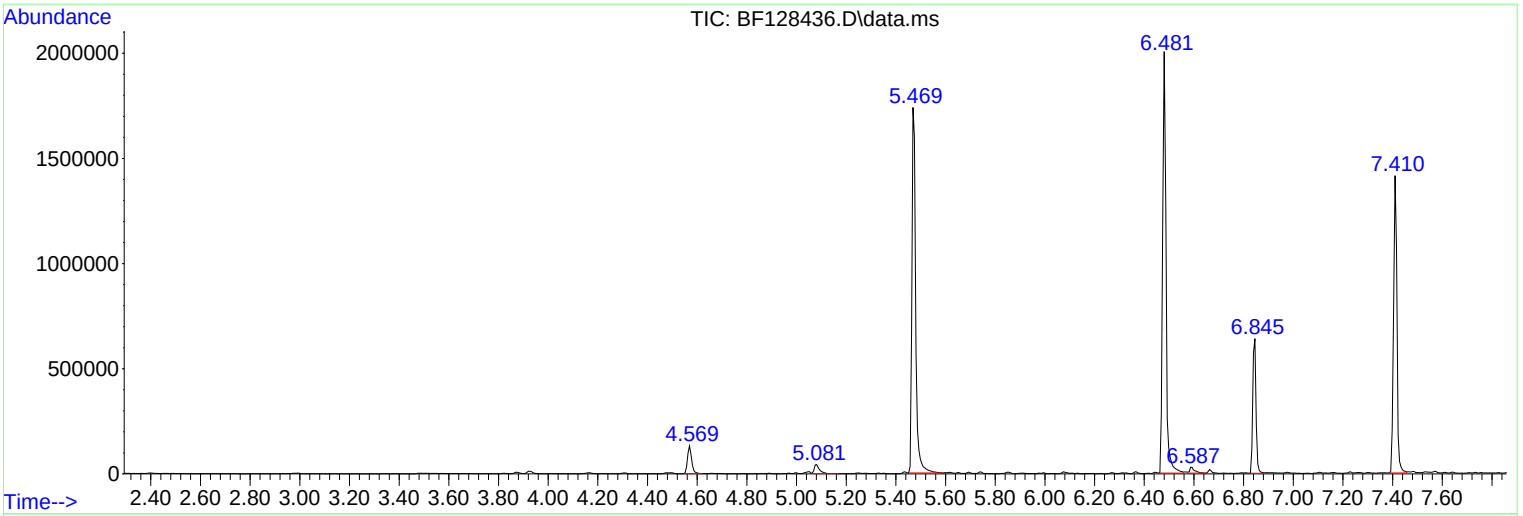
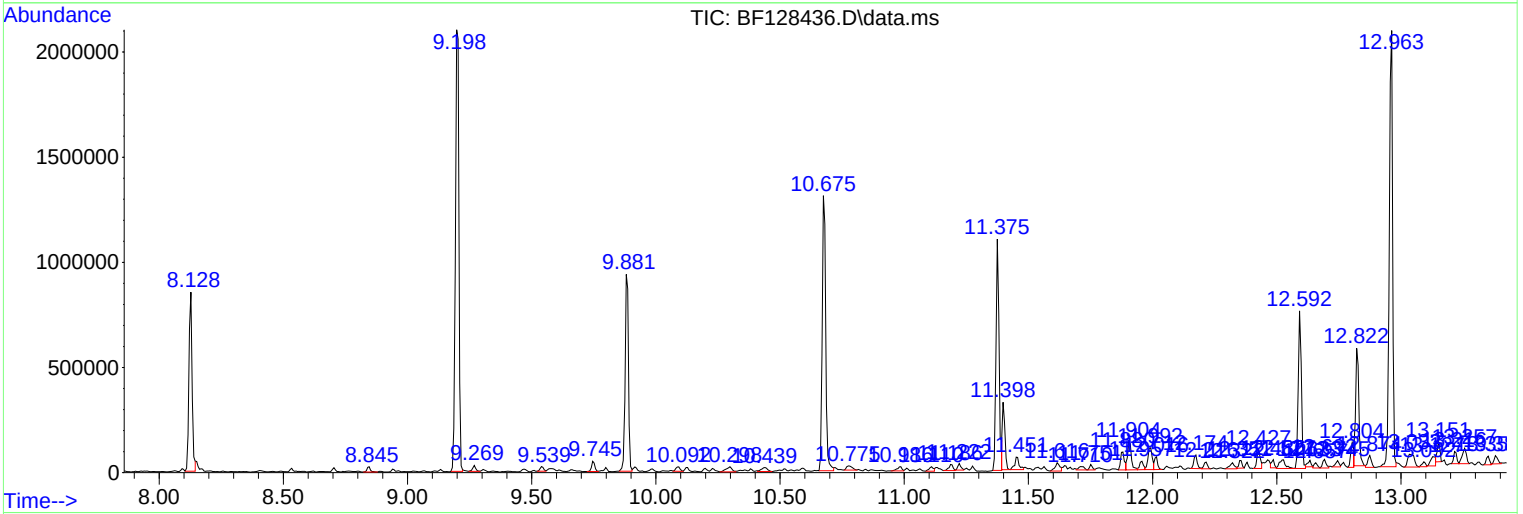
Operator : CG\JU
Sample : N2981-02
Misc :
ALS Vial : 9 Sample Multiplier: 1

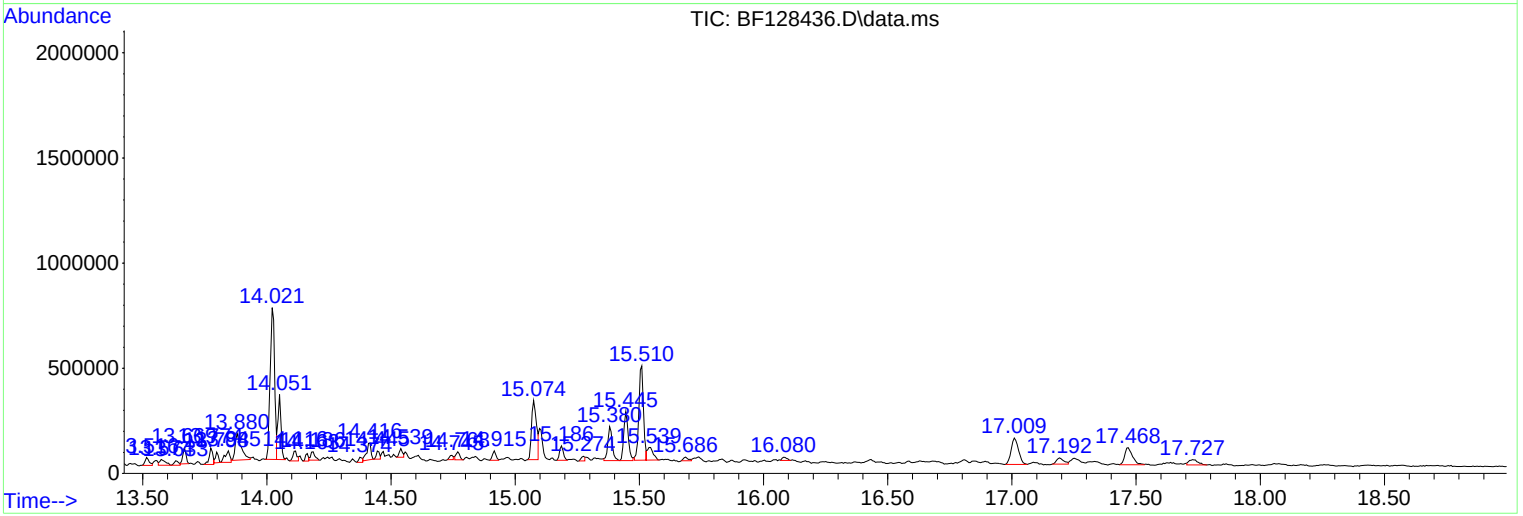
Instrument :
BNA_F
ClientSampleId :
SB-13(3-3.5)

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052022.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\BF052422\
Data File : BF128436.D
Acq On : 24 May 2022 15:48
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Sample : N2981-02
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Instrument :
BNA_F
ClientSampleId :
SB-13(3-3.5)

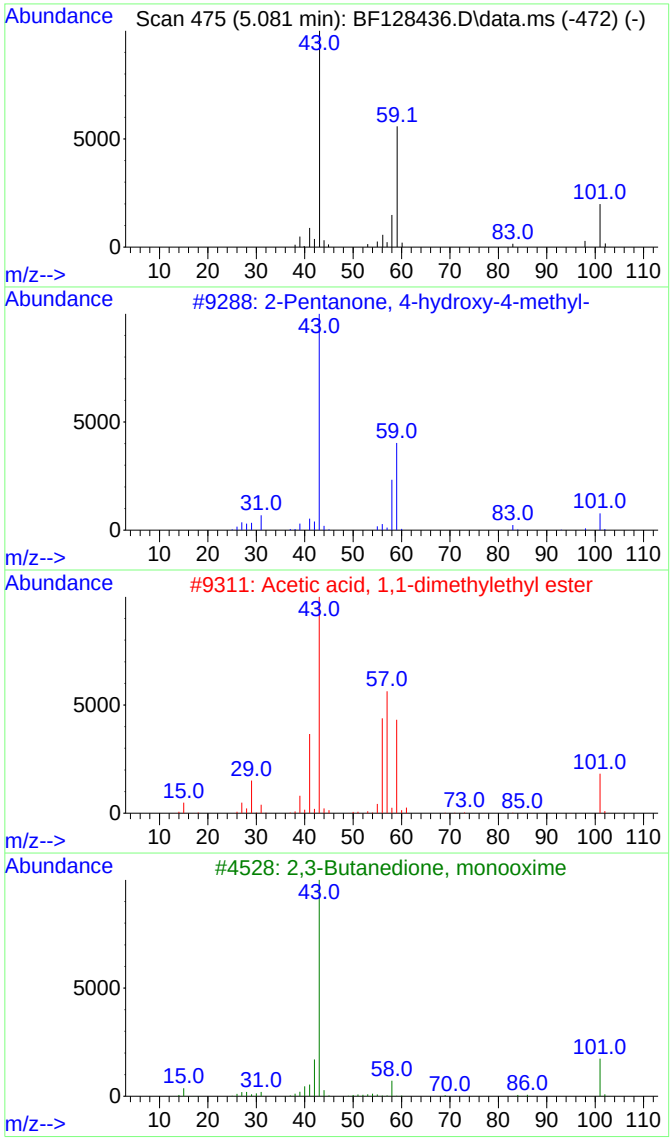
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052022.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

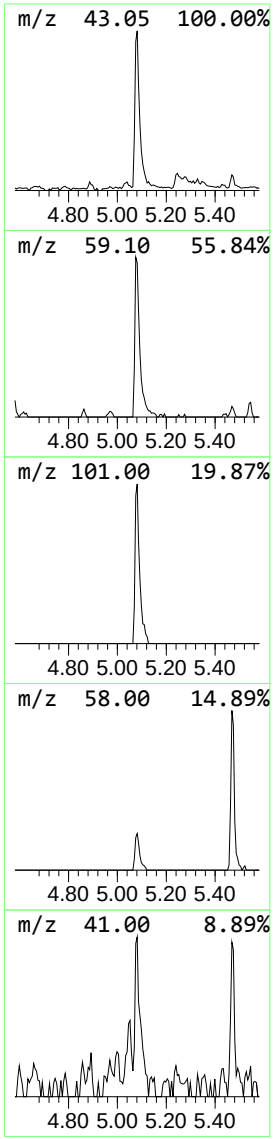
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.081	2.31 ng	66455	1,4-Dichlorobenzene-d4	6.845

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	35
4	3-Hexanol	102	C6H14O	000623-37-0	28
5	Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	23





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF052422\
Data File : BF128436.D
Acq On : 24 May 2022 15:48
Operator : CG\JU
Sample : N2981-02
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-13(3-3.5)

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052022.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

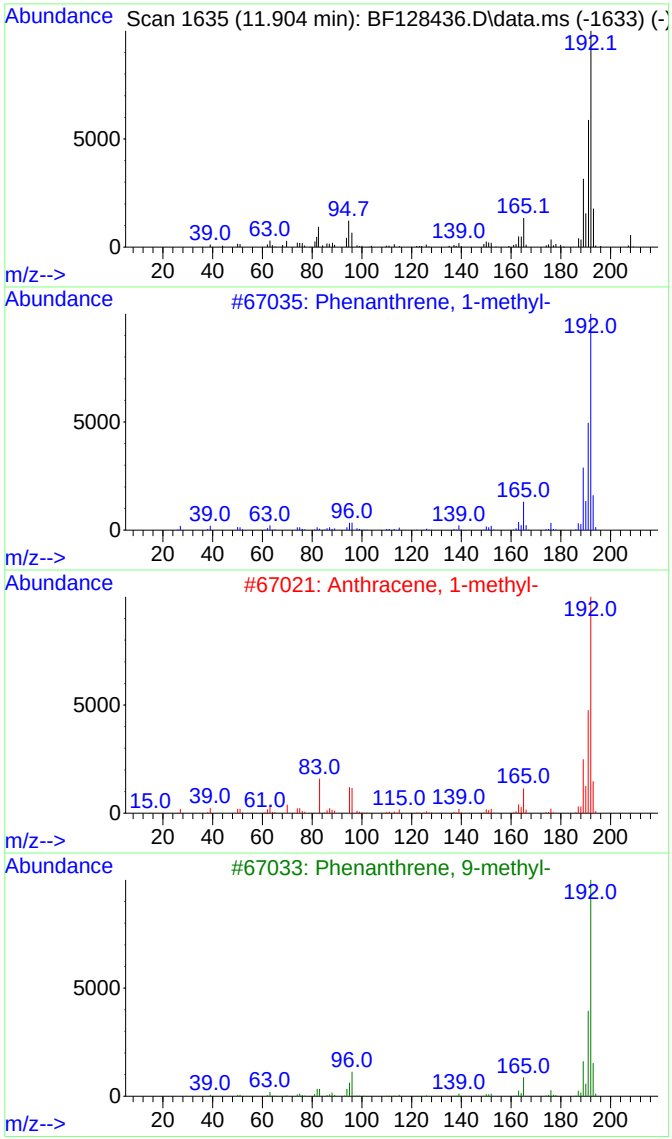
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

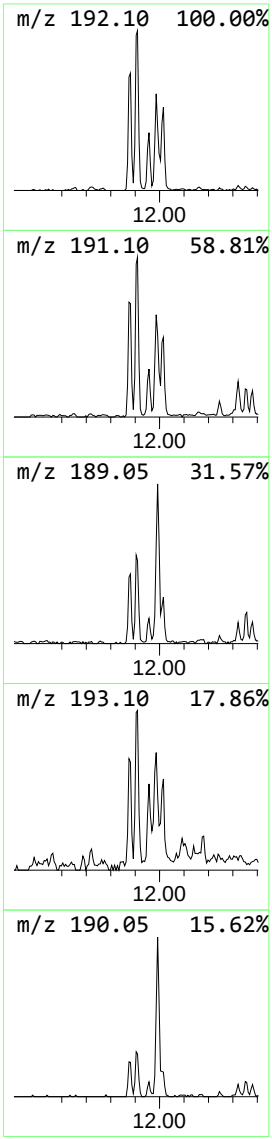
****

Peak Number 3 Phenanthrene, 1-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.904	2.94 ng	138629	Phenanthrene-d10	11.374

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
2	Anthracene, 1-methyl-	192	C15H12	000610-48-0	93
3	Phenanthrene, 9-methyl-	192	C15H12	000883-20-5	91
4	Anthracene, 2-methyl-	192	C15H12	000613-12-7	91
5	Anthracene, 9-methyl-	192	C15H12	000779-02-2	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF052422\
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Operator : CG\JU
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ClientSampleId :
SB-13(3-3.5)

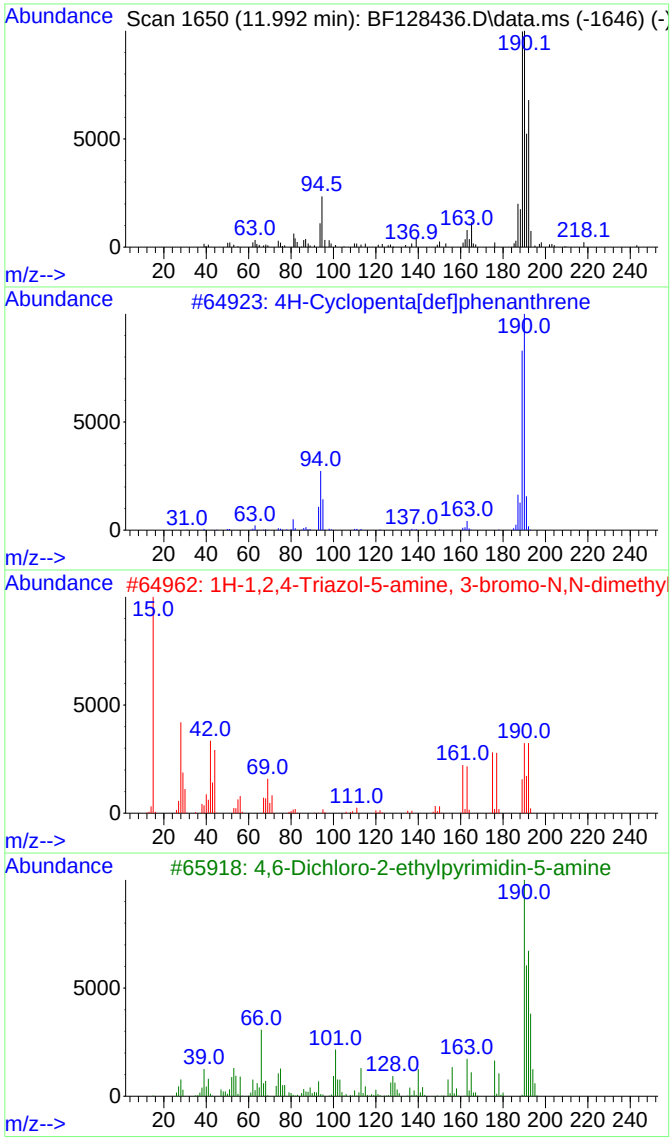
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052022.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

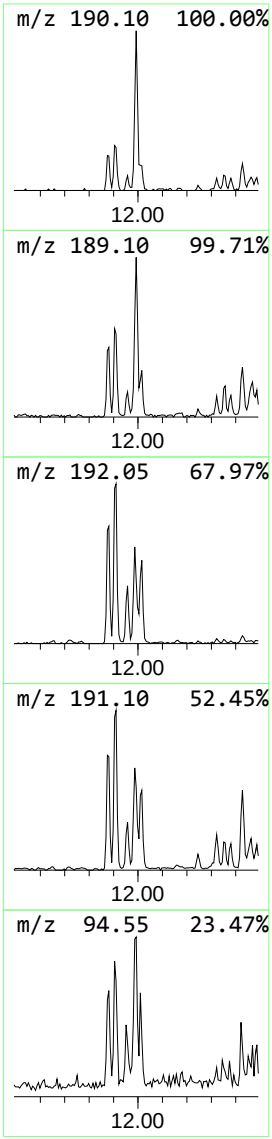
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.992	2.63 ng	123850	Phenanthrene-d10	11.374

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	70
2	1H-1,2,4-Triazol-5-amine, 3-brom...	190	C4H7BrN4	1000460-85-7	47
3	4,6-Dichloro-2-ethylpyrimidin-5-...	191	C6H7Cl2N3	006237-96-3	43
4	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	42
5	Methyl diselenide	190	C2H6Se2	007101-31-7	38





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF052422\
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Instrument :
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SB-13(3-3.5)

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052022.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

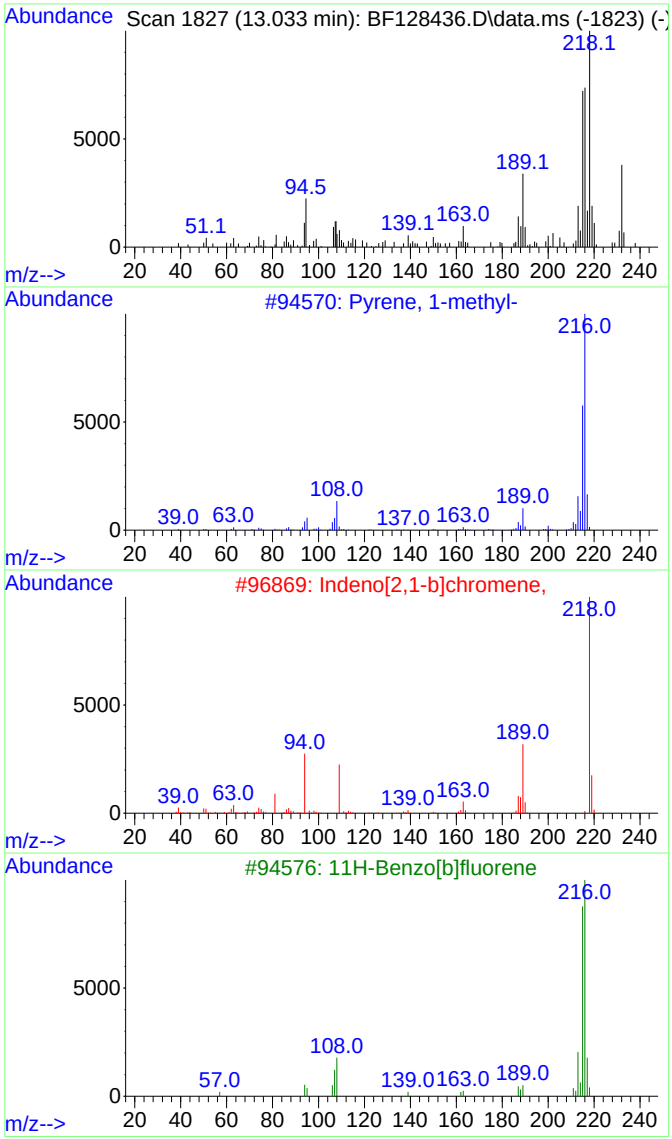
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

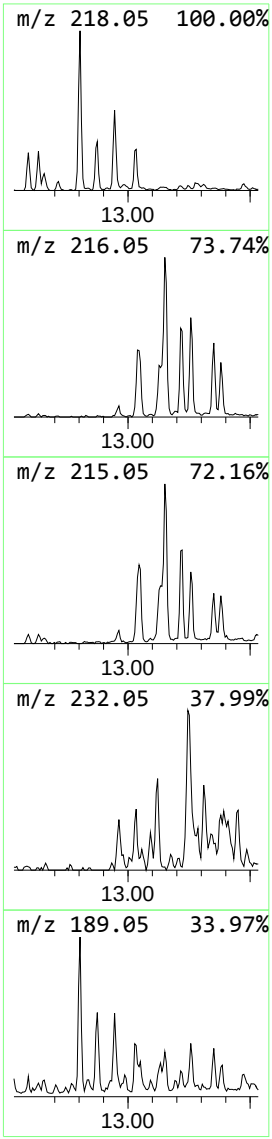
****

Peak Number 5 Pyrene, 1-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.033	2.45 ng	113423	Chrysene-d12	14.027

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-	216	C17H12	002381-21-7	70
2	Indeno[2,1-b]chromene,	218	C16H10O	000243-24-3	55
3	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	50
4	Pyrene, 2-methyl-	216	C17H12	003442-78-2	50
5	Benzo[b]naphtho[2,3-d]furan	218	C16H10O	000243-42-5	42





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF052422\
Data File : BF128436.D
Acq On : 24 May 2022 15:48
Operator : CG\JU
Sample : N2981-02
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-13(3-3.5)

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052022.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

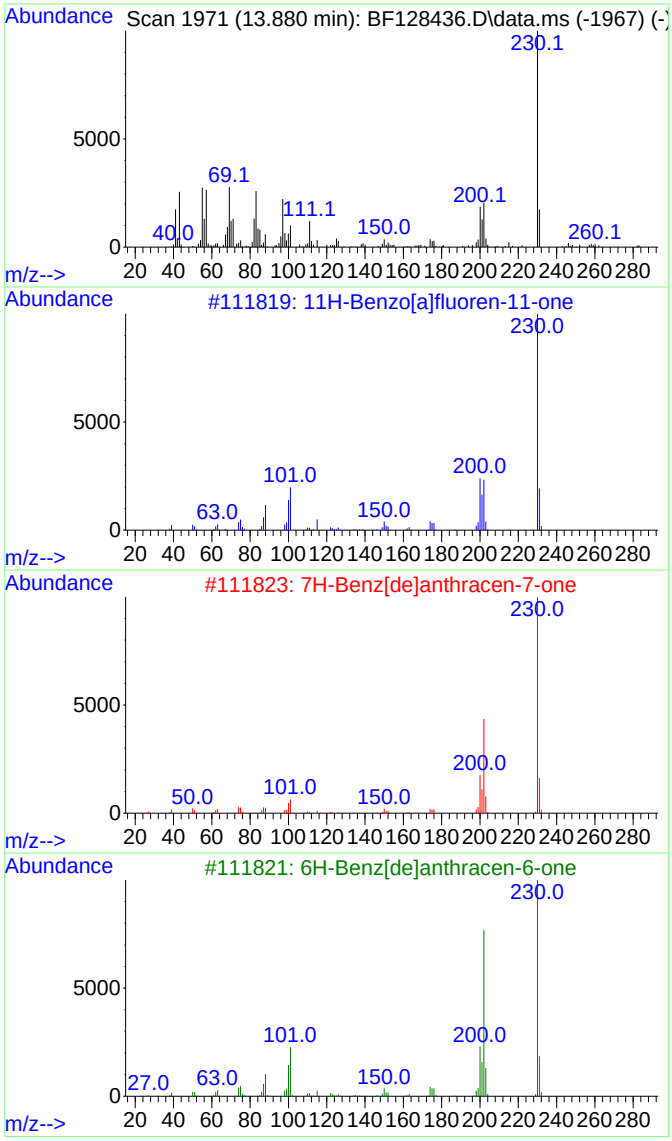
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

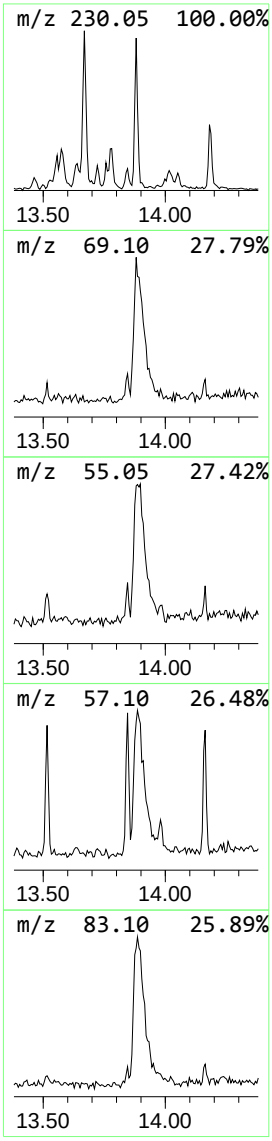
****

Peak Number 6 11H-Benzo[a]fluoren-11-one Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.880	4.64 ng	214762	Chrysene-d12	14.027

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[a]fluoren-11-one	230	C17H10O	000479-79-8	95
2	7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	78
3	6H-Benz[de]anthracen-6-one	230	C17H10O	080252-14-8	55
4	4-Pyrenecarbaldehyde	230	C17H10O	022245-51-8	50
5	.delta.(sup1),.alpha.-Acenaphthe...	230	C15H6N2O	014619-86-4	47





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Acq On : 24 May 2022 15:48
Operator : CG\JU
Sample : N2981-02
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-13(3-3.5)

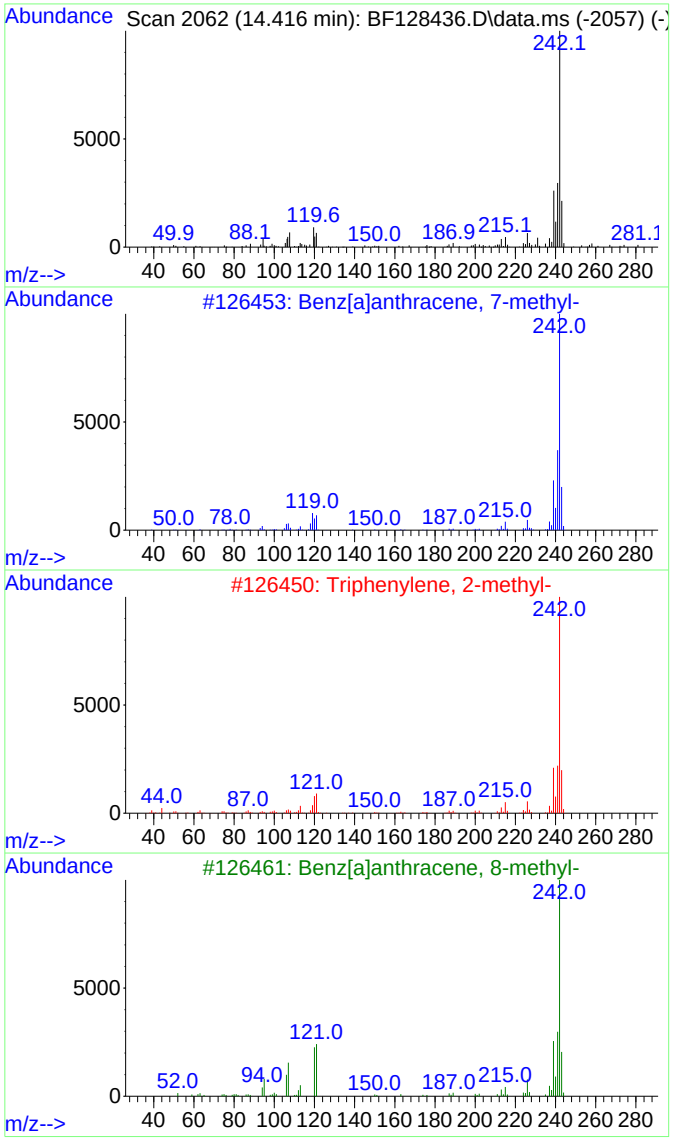
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052022.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

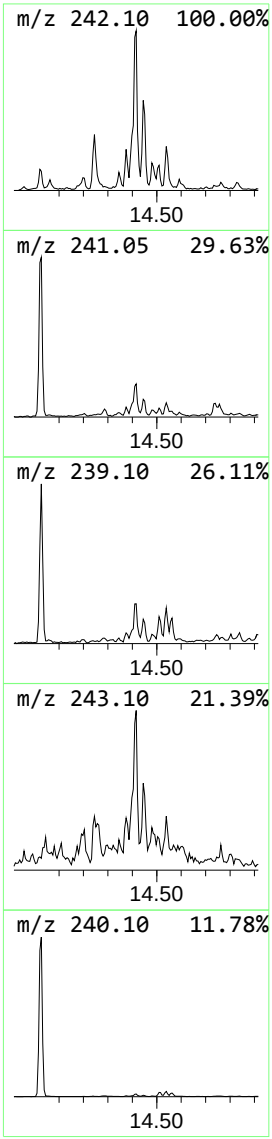
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benz[a]anthracene, 7-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.416	2.04 ng	94496	Chrysene-d12	14.027

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benz[a]anthracene, 7-methyl-	242	C19H14	002541-69-7	96
2	Triphenylene, 2-methyl-	242	C19H14	001705-84-6	93
3	Benz[a]anthracene, 8-methyl-	242	C19H14	002381-31-9	93
4	Chrysene, 3-methyl-	242	C19H14	003351-31-3	93
5	Benz[a]anthracene, 1-methyl-	242	C19H14	002498-77-3	92





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Data File : BF128436.D
Acq On : 24 May 2022 15:48
Operator : CG\JU
Sample : N2981-02
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-13(3-3.5)

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052022.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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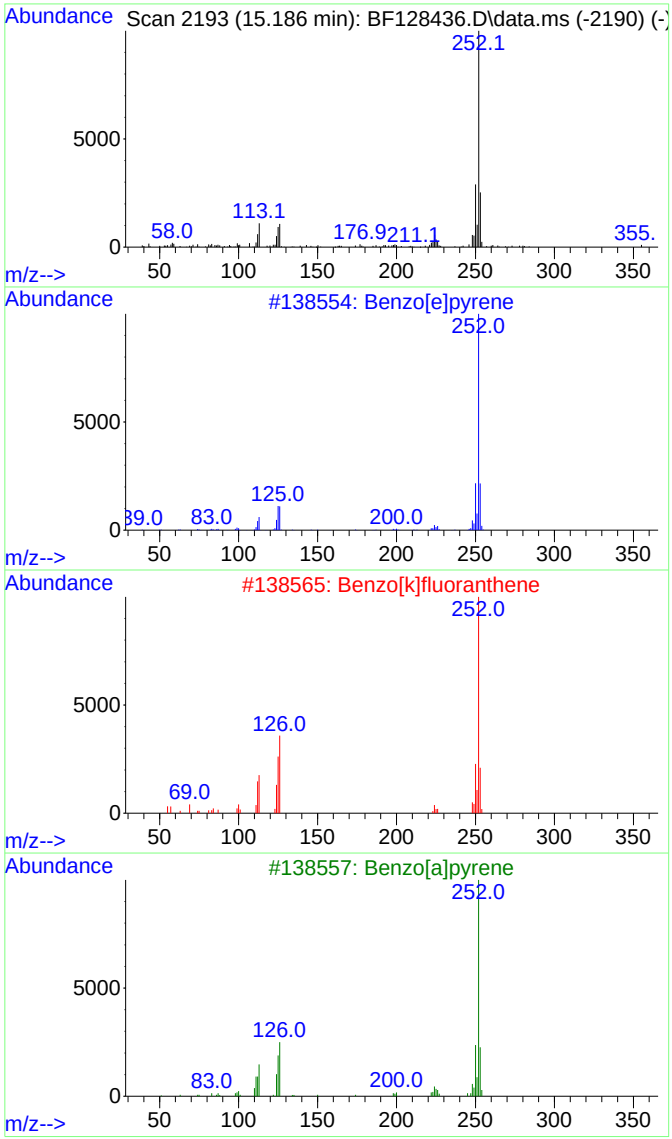
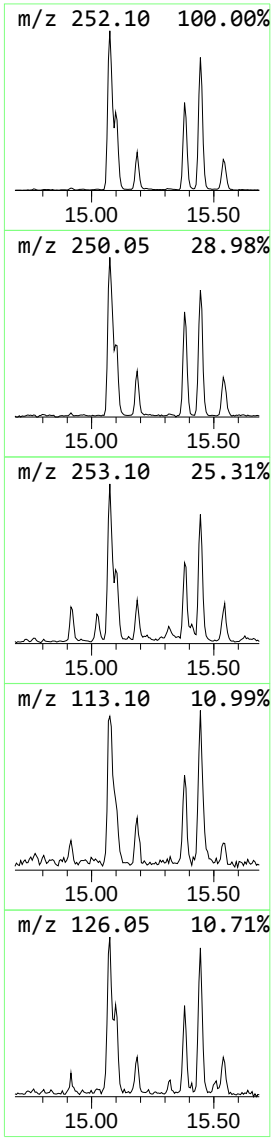
Peak Number 8 Benzo[e]pyrene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.186	2.52 ng	73332	Perylene-d12	15.510

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	94
3	Benzo[a]pyrene	252	C20H12	000050-32-8	91
4	Benzo[b]fluoranthene	252	C20H12	000205-99-2	90
5	Perylene	252	C20H12	000198-55-0	90





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Operator : CG\JU
Sample : N2981-02
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-13(3-3.5)

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052022.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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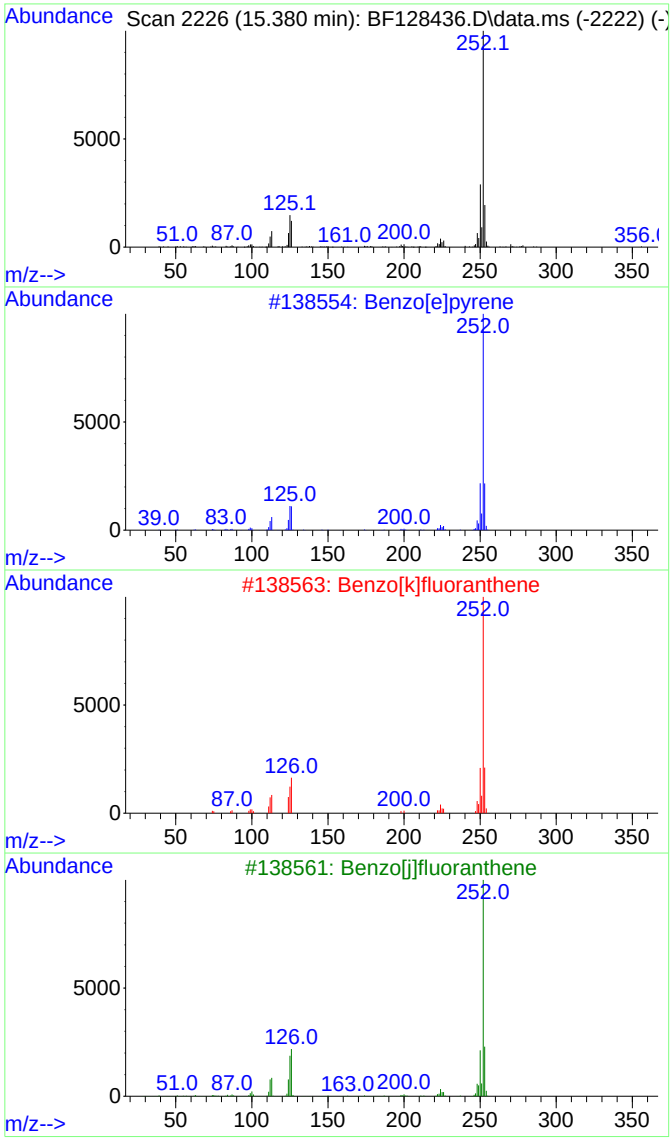
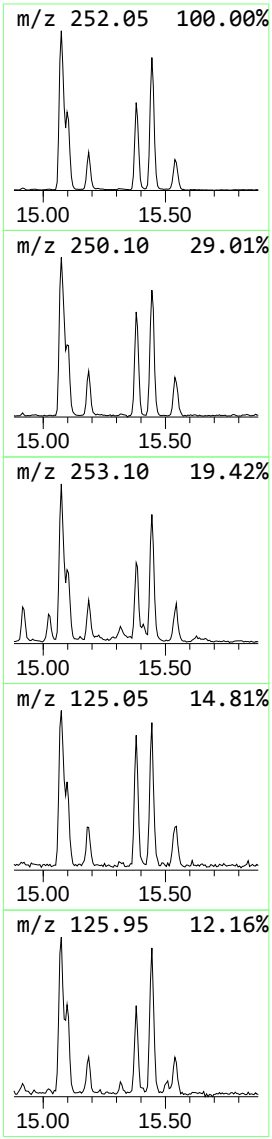
Peak Number 9 Benzo[j]fluoranthene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.380	6.93 ng	201864	Perylene-d12	15.510

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[e]pyrene	252	C20H12	000192-97-2	99
2	Benzo[k]fluoranthene	252	C20H12	000207-08-9	97
3	Benzo[j]fluoranthene	252	C20H12	000205-82-3	97
4	Perylene	252	C20H12	000198-55-0	96
5	Benzo[b]fluoranthene	252	C20H12	000205-99-2	96





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Data File : BF128436.D
Acq On : 24 May 2022 15:48
Operator : CG\JU
Sample : N2981-02
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-13(3-3.5)

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052022.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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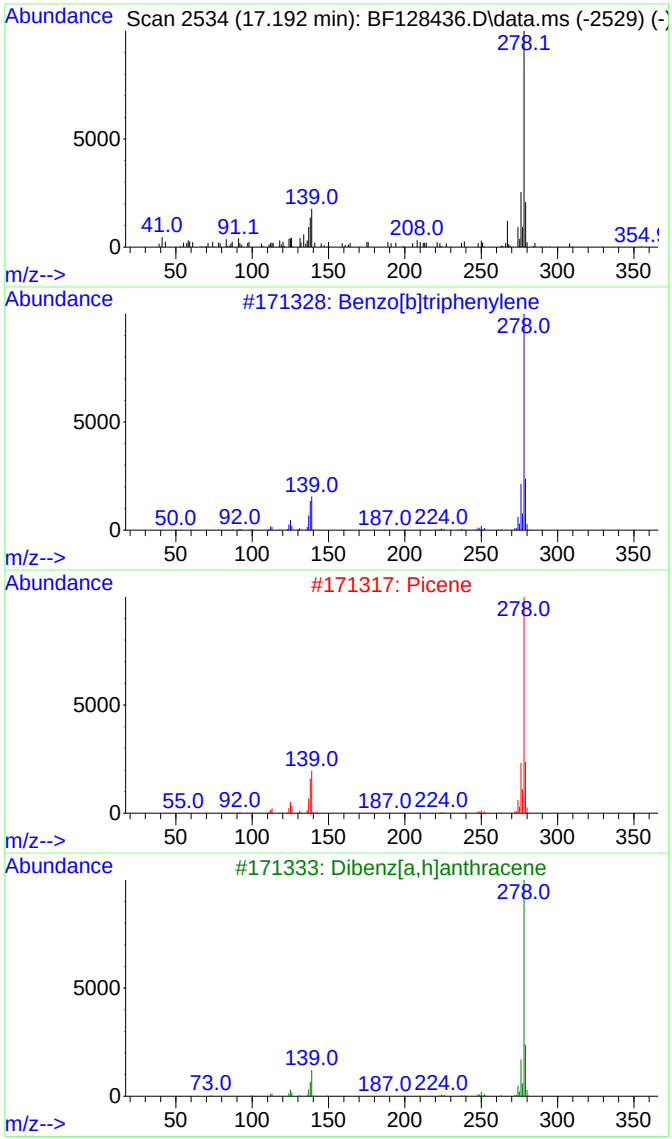
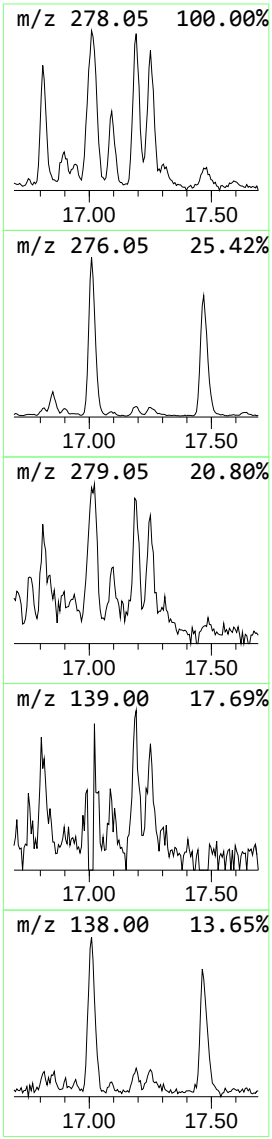
Peak Number 10 Benzo[b]triphenylene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.192	2.13 ng	62146	Perylene-d12	15.510

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[b]triphenylene	278	C22H14	000215-58-7	96
2	Picene	278	C22H14	000213-46-7	95
3	Dibenz[a,h]anthracene	278	C22H14	000053-70-3	90
4	Benzo[b]chrysene	278	C22H14	000214-17-5	89
5	Benzo[a]naphthacene	278	C22H14	000226-88-0	89





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Data File : BF128436.D
Acq On : 24 May 2022 15:48
Operator : CG\JU
Sample : N2981-02
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-13(3-3.5)

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052022.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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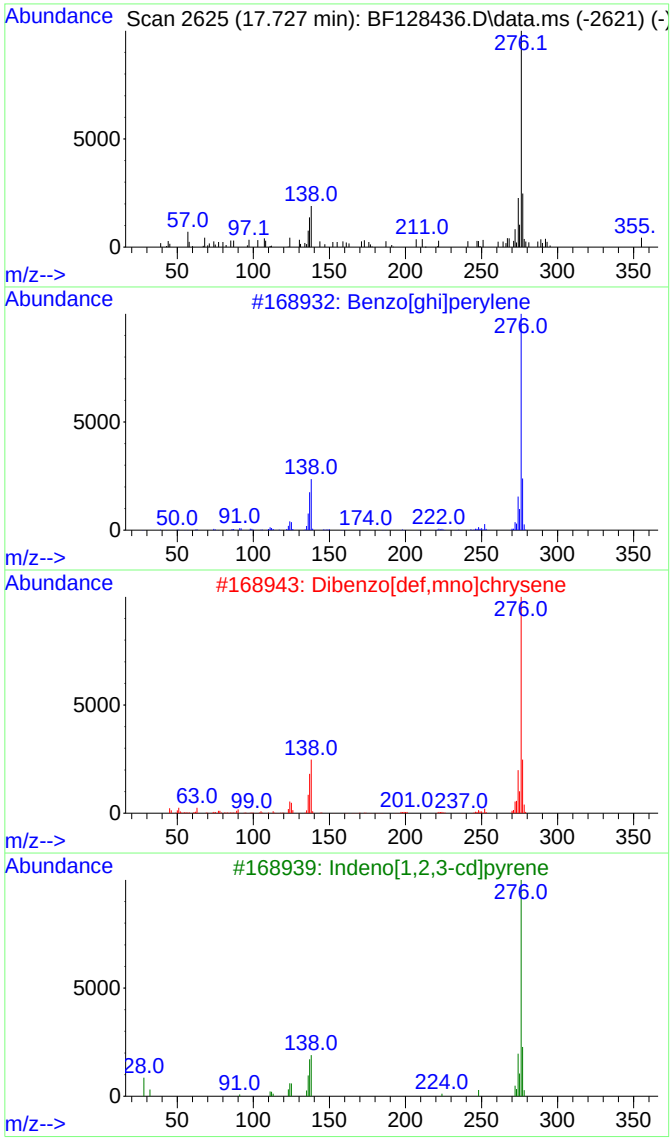
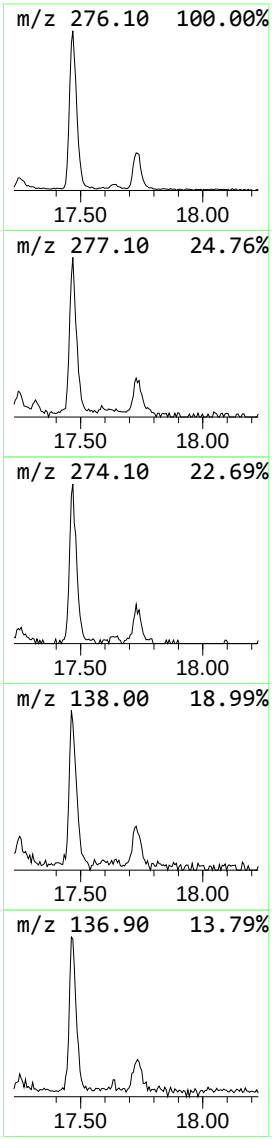
Peak Number 11 Dibenzo[def,mno]chrysene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.727	2.23 ng	64820	Perylene-d12	15.510

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[ghi]perylene	276	C22H12	000191-24-2	92
2	Dibenzo[def,mno]chrysene	276	C22H12	000191-26-4	76
3	Indeno[1,2,3-cd]pyrene	276	C22H12	000193-39-5	76
4	Indeno[1,2,3-cd]fluoranthene	276	C22H12	000193-43-1	62
5	Naphthalen-1-yl-(3-nitro-benzyli...	276	C17H12N2O2	200421-53-0	53





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF052422\
Data File : BF128436.D
Acq On : 24 May 2022 15:48
Operator : CG\JU
Sample : N2981-02
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-13(3-3.5)

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052022.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
2-Pentanone, 4-...	5.081	2.3	ng	66455	1	6.845	575776	20.0
Phenanthrene, 1...	11.904	2.9	ng	138629	4	11.374	942132	20.0
4H-Cyclopenta[d...	11.992	2.6	ng	123850	4	11.374	942132	20.0
Pyrene, 1-methyl-	13.033	2.5	ng	113423	5	14.027	925240	20.0
11H-Benzo[a]flu...	13.880	4.6	ng	214762	5	14.027	925240	20.0
Benz[a]anthrace...	14.416	2.0	ng	94496	5	14.027	925240	20.0
Benzo[e]pyrene	15.186	2.5	ng	73332	6	15.510	582248	20.0
Benzo[j]fluoran...	15.380	6.9	ng	201864	6	15.510	582248	20.0
Benzo[b]triphen...	17.192	2.1	ng	62146	6	15.510	582248	20.0
Dibenzo[def,mno...	17.727	2.2	ng	64820	6	15.510	582248	20.0