LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111121\
Data File : BF126136.D
Acq On : 11 Nov 2021 20:55
Operator : JU/CG
Sample : M4630-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-13N-5.0

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-BF110321.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF126136.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.416	392	396	403	rVB	87332	99121	2.88%	0.252%
2	5.045	499	503	510	rVB	135649	159044	4.63%	0.405%
3	5.457	568	573	576	rBV	2880689	2849545	82.94%	7.255%
4	6.469	740	745	748	rBV	3287695	3068905	89.32%	7.814%
5	6.839	805	808	811	rBV	1139850	857434	24.96%	2.183%
6	7.398	899	903	906	rBV	2140447	1977402	57.55%	5.035%
7	8.027	1006	1010	1022	rBV	256185	281401	8.19%	0.716%
8	8.122	1022	1026	1029	rBV	1275525	1096458	31.91%	2.792%
9	9.198	1205	1209	1212	rBV	4317239	3435840	100.00%	8.748%
10	9.739	1297	1301	1304	rVB	53995	45734	1.33%	0.116%
11	9.880	1321	1325	1328	rBV	1488971	1271264	37.00%	3.237%
12	9.910	1328	1330	1333	rVB	124889	91311	2.66%	0.232%
13	10.080	1356	1359	1363	rBV	105749	84029	2.45%	0.214%
14	10.421	1414	1417	1423	rVB	97278	86443	2.52%	0.220%
15	10.663	1454	1458	1462	rBV	2696933	2281798	66.41%	5.810%
16	10.974	1504	1511	1513	rBV4	27083	48029	1.40%	0.122%
17	11.163	1541	1543	1548	rVB	73040	67559	1.97%	0.172%
18	11.210	1548	1551	1557	rVV5	34210	63829	1.86%	0.163%
19	11.257	1557	1559	1564	rVB	72040	62883	1.83%	0.160%
20	11.363	1573	1577	1579	rBV	1708798	1367371	39.80%	3.481%
21	11.386	1579	1581	1584	rVV	1379923	993735	28.92%	2.530%
22	11.433	1584	1589	1593	rVB	264420	272698	7.94%	0.694%
23	11.592	1610	1616	1618	rBV2	146805	150914	4.39%	0.384%
24	11.757	1641	1644	1652	rVB6	33579	56221	1.64%	0.143%
25	11.863	1657	1662	1664	rVV	172324	150277	4.37%	0.383%
26	11.892	1664	1667	1671	rVB	257154	213106	6.20%	0.543%
27	11.939	1671	1675	1677	rBV	87364	73914	2.15%	0.188%
28	11.974	1677	1681	1683	rVV	227398	228606	6.65%	0.582%
29	11.998	1683	1685	1688	rVB	137226	103557	3.01%	0.264%
30	12.163	1710	1713	1714	rBV	122348	110507	3.22%	0.281%
31	12.180	1714	1716	1720	rVB	179125	135683	3.95%	0.345%
32	12.310	1735	1738	1741	rVB2	58389	41480	1.21%	0.106%
33	12.415	1750	1756	1759	rBV	120128	134560	3.92%	0.343%
34	12.451	1759	1762	1764	rVV3	63250	82590	2.40%	0.210%
35	12.498	1767	1770	1776	rVB2	140482	140959	4.10%	0.359%
36	12.574	1780	1783	1787	rVB	1874426	1586401	46.17%	4.039%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111121\
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Acq On : 11 Nov 2021 20:55
Operator : JU/CG
Sample : M4630-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-13N-5.0

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-BF110321.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	12.639	1787	1794	1797	rBV3	43749	66001	1.92%	0.168%
38	12.668	1797	1799	1801	rVB	59424	43285	1.26%	0.110%
39	12.704	1801	1805	1807	rBV2	55415	70813	2.06%	0.180%
40	12.727	1807	1809	1812	rVB	54279	42827	1.25%	0.109%
41	12.804	1816	1822	1828	rBV	1626703	1643594	47.84%	4.185%
42	12.851	1828	1830	1836	rVB2	92027	101483	2.95%	0.258%
43	12.921	1836	1842	1843	rBV	113075	99228	2.89%	0.253%
44	12.951	1843	1847	1850	rVV	3941771	3171778	92.31%	8.076%
45	13.021	1854	1859	1862	rBV2	118434	195994	5.70%	0.499%
46	13.110	1871	1874	1875	rBV2	86016	89345	2.60%	0.227%
47	13.133	1875	1878	1880	rVB	131693	140820	4.10%	0.359%
48	13.198	1885	1889	1892	rBV	66471	59177	1.72%	0.151%
49	13.233	1892	1895	1899	rBV	163249	168625	4.91%	0.429%
50	13.327	1908	1911	1914	rBV	94475	74712	2.17%	0.190%
51	13.357	1914	1916	1920	rBV2	91346	98222	2.86%	0.250%
52	13.639	1960	1964	1968	rVB	183807	186362	5.42%	0.474%
53	13.704	1968	1975	1977	rBV5	25901	52350	1.52%	0.133%
54	13.745	1978	1982	1986	rBV2	116620	168307	4.90%	0.429%
55	13.780	1986	1988	1990	rVV	121065	95177	2.77%	0.242%
56	13.804	1990	1992	1996	rVV2	82983	105499	3.07%	0.269%
57	13.845	1996	1999	2002	rVV2	137249	125726	3.66%	0.320%
58	13.880	2002	2005	2010	rVB6	53831	90420	2.63%	0.230%
59	14.004	2017	2026	2028	rBV2	1393427	1913741	55.70%	4.873%
60	14.027	2028	2030	2033	rVB	653524	482099	14.03%	1.227%
61	14.145	2047	2050	2052	rBV	38218	41812	1.22%	0.106%
62	14.227	2060	2064	2067	rBV2	67056	97164	2.83%	0.247%
63	14.351	2083	2085	2088	rBV3	39986	48373	1.41%	0.123%
64	14.386	2088	2091	2095	rVV	150990	174477	5.08%	0.444%
65	14.421	2095	2097	2100	rVV2	103482	101651	2.96%	0.259%
66	14.480	2104	2107	2110	rVV3	73076	91919	2.68%	0.234%
67	14.515	2110	2113	2115	rVV2	90613	108912	3.17%	0.277%
68	14.533	2115	2116	2120	rVV3	70579	75048	2.18%	0.191%
69	14.586	2123	2125	2129	rVB2	47265	53486	1.56%	0.136%
70	14.692	2139	2143	2146	rBV3	61099	75193	2.19%	0.191%
71	14.792	2155	2160	2162	rBV6	35015	52790	1.54%	0.134%
72	14.862	2168	2172	2179	rBV2	200229	258959	7.54%	0.659%
73	14.945	2184	2186	2190	rVB3	40627	45585	1.33%	0.116%
74	15.039	2198	2202	2206	rBV	598526	994407	28.94%	2.532%
75	15.121	2213	2216	2218	rBV2	34999	47392	1.38%	0.121%
76	15.145	2218	2220	2223	rVB	111492	105994	3.08%	0.270%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111121\
Data File : BF126136.D
Acq On : 11 Nov 2021 20:55
Operator : JU/CG
Sample : M4630-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-13N-5.0

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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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

77	15.239	2233	2236	2237	rBV2	40764	42377	1.23%	0.108%
78	15.345	2249	2254	2259	rVB	353885	483635	14.08%	1.231%
79	15.403	2259	2264	2270	rVB	526129	639675	18.62%	1.629%
80	15.468	2270	2275	2278	rBV	930315	1128188	32.84%	2.872%
81	15.498	2278	2280	2287	rVB2	159374	215871	6.28%	0.550%
82	15.945	2352	2356	2360	rVB2	55832	73338	2.13%	0.187%
83	16.374	2426	2429	2436	rVB8	29054	55265	1.61%	0.141%
84	16.450	2438	2442	2445	rBV5	34301	43552	1.27%	0.111%
85	16.739	2485	2491	2492	rBV3	45143	77751	2.26%	0.198%
86	16.915	2516	2521	2531	rVB2	237878	465129	13.54%	1.184%
87	17.092	2547	2551	2556	rVB2	44129	71501	2.08%	0.182%
88	17.150	2557	2561	2566	rBV2	37732	68027	1.98%	0.173%
89	17.356	2589	2596	2604	rVB	171921	362569	10.55%	0.923%
90	17.592	2631	2636	2640	rVB3	37314	74878	2.18%	0.191%
91	17.739	2657	2661	2667	rVB8	20703	45230	1.32%	0.115%

Sum of corrected areas: 39276341

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111121\
Data File : BF126136.D
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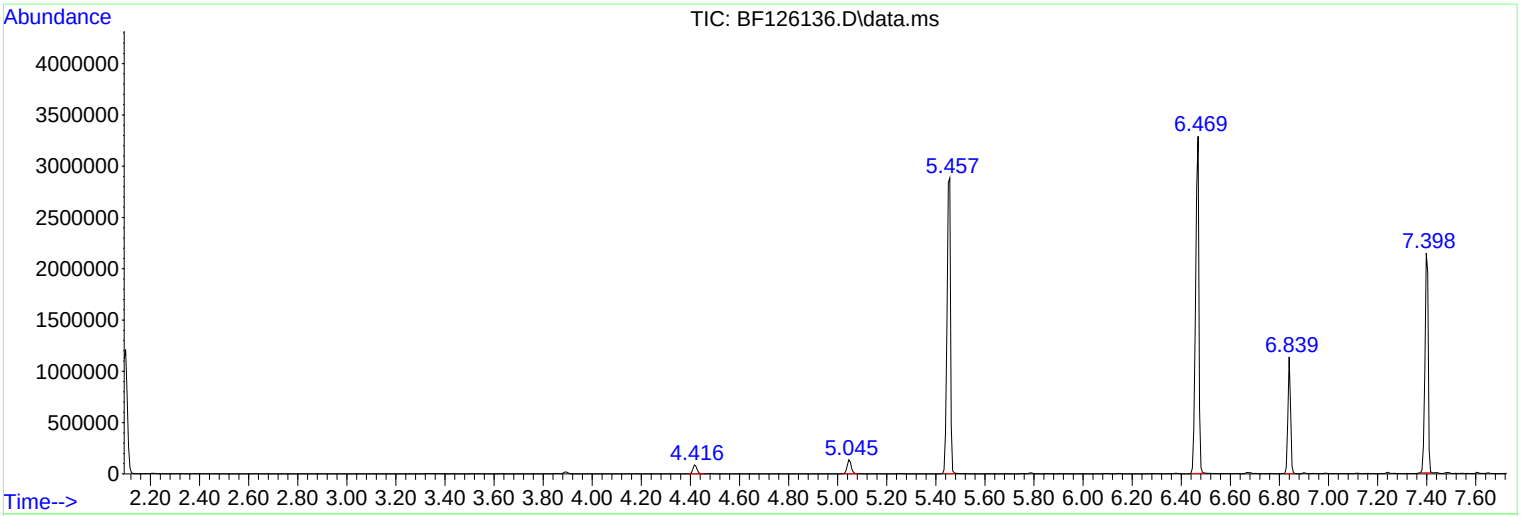
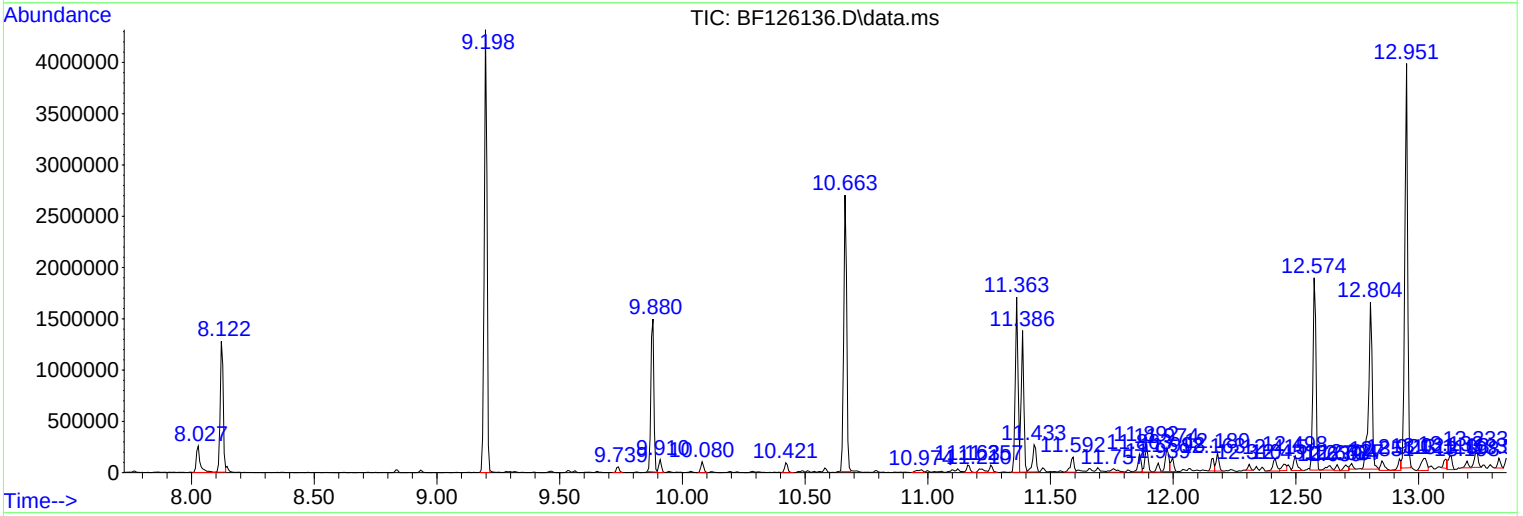
Sample : M4630-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

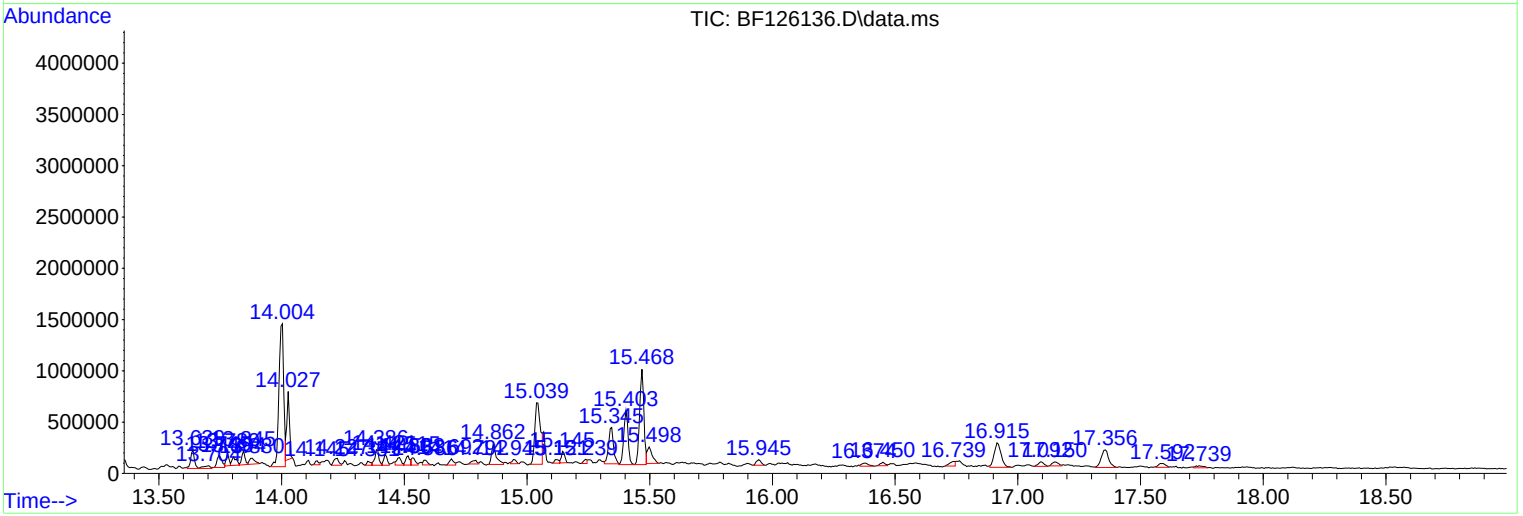
Instrument :
BNA_F
ClientSampleId :
SB-13N-5.0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110321.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\BF111121\
Data File : BF126136.D
Acq On : 11 Nov 2021 20:55
Operator : JU/CG
Sample : M4630-06
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ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-13N-5.0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

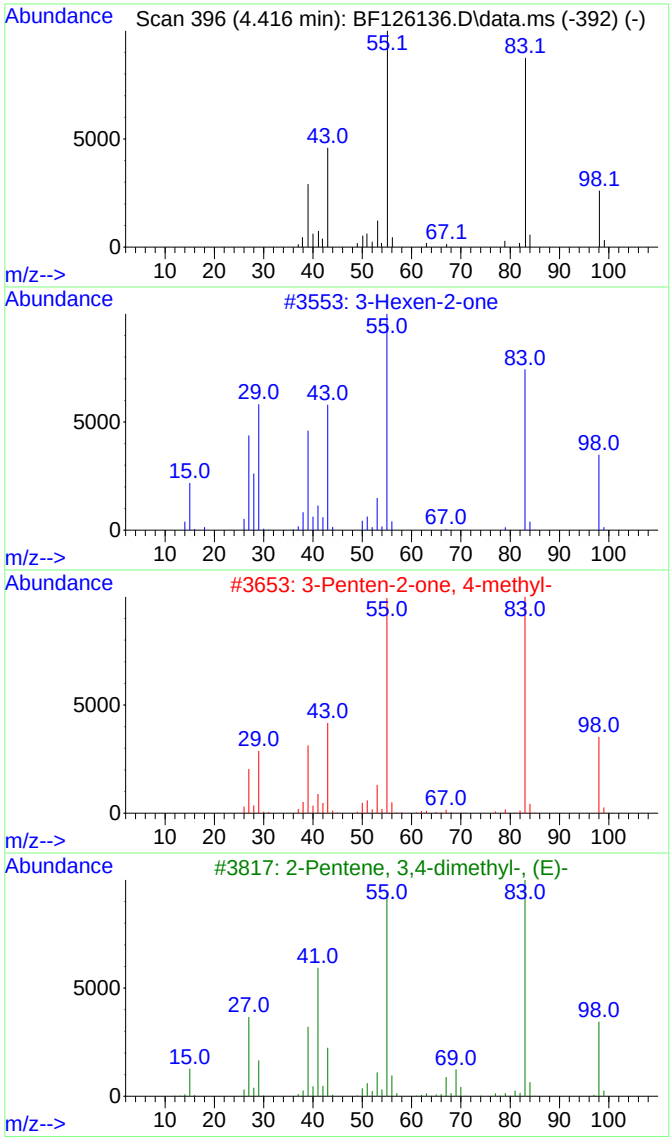
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

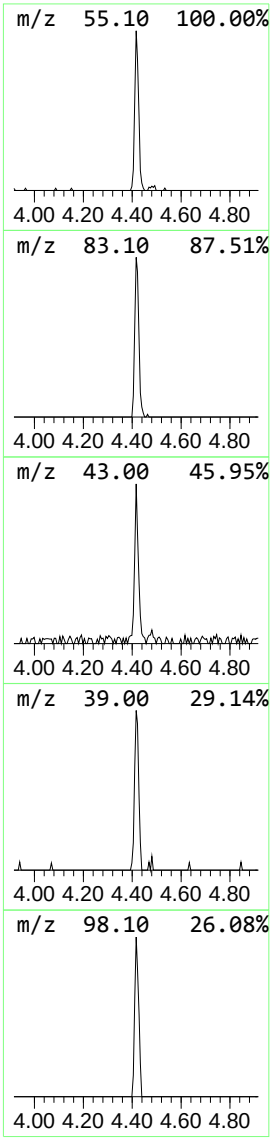
****

Peak Number 1 3-Hexen-2-one Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.416	2.31 ng	99121	1,4-Dichlorobenzene-d4	6.839

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Hexen-2-one	98	C6H10O	000763-93-9	91
2	3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	91
3	2-Pentene, 3,4-dimethyl-, (E)-	98	C7H14	004914-92-5	90
4	2-Pentene, 4,4-dimethyl-, (Z)-	98	C7H14	000762-63-0	86
5	Cyclopropane, 1,1,2,2-tetramethyl-	98	C7H14	004127-47-3	86





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111121\
Data File : BF126136.D
Acq On : 11 Nov 2021 20:55
Operator : JU/CG
Sample : M4630-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-13N-5.0

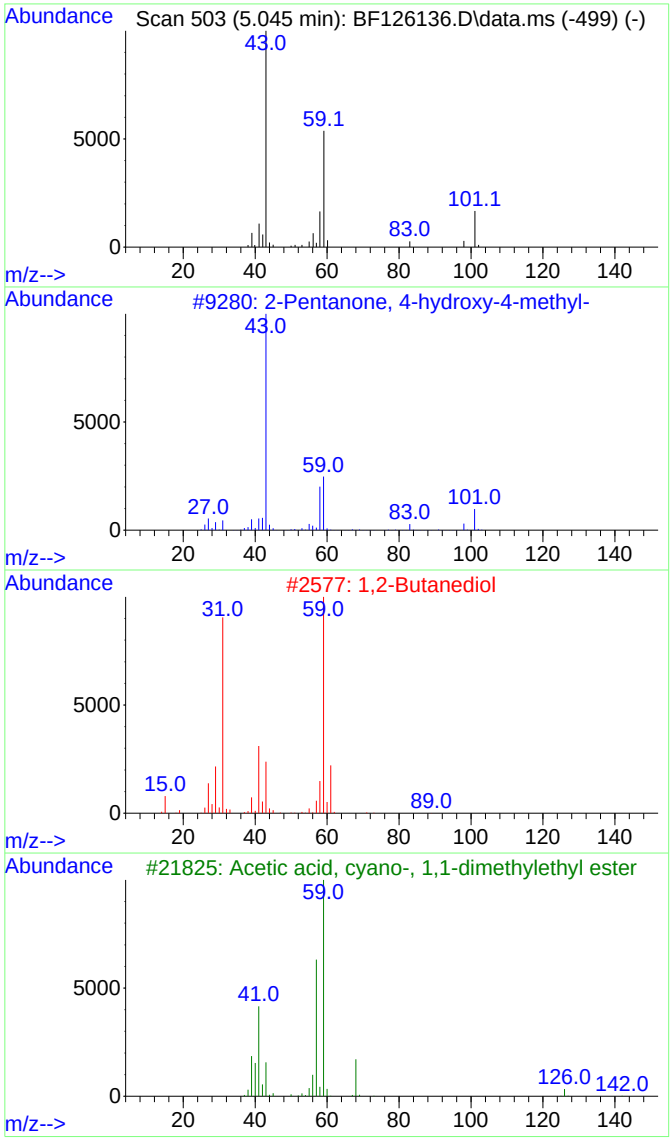
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

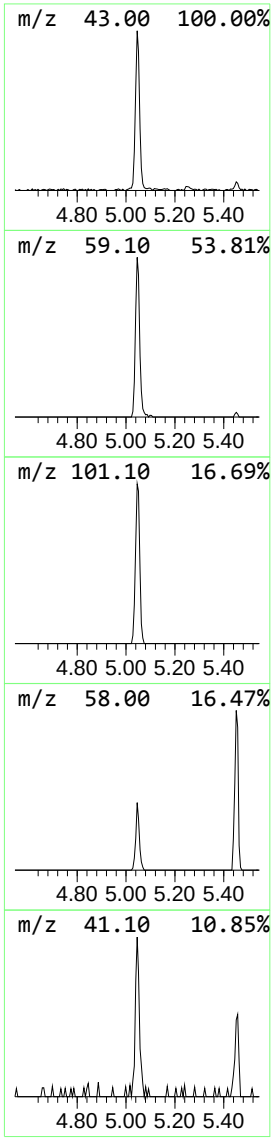
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.045	3.71 ng	159044	1,4-Dichlorobenzene-d4	6.839

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2	1,2-Butanediol	90	C4H10O2	000584-03-2	28
3	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
4	Acetone	58	C3H6O	000067-64-1	12
5	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	12





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111121\
Data File : BF126136.D
Acq On : 11 Nov 2021 20:55
Operator : JU/CG
Sample : M4630-06
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ALS Vial : 12 Sample Multiplier: 1

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ClientSampleId :
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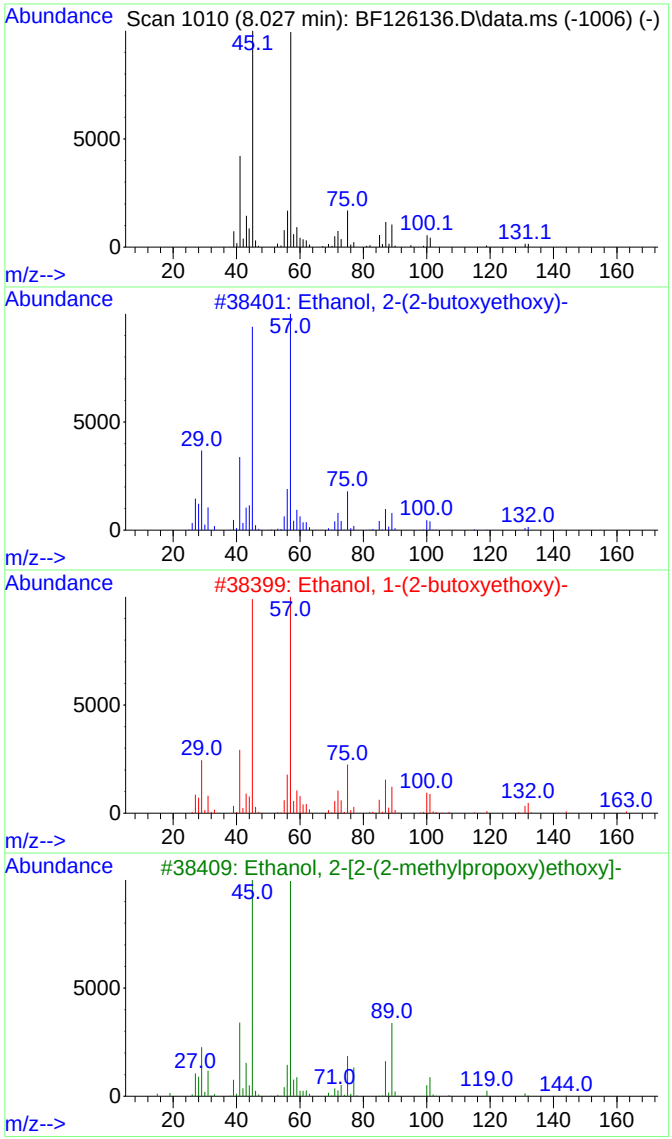
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

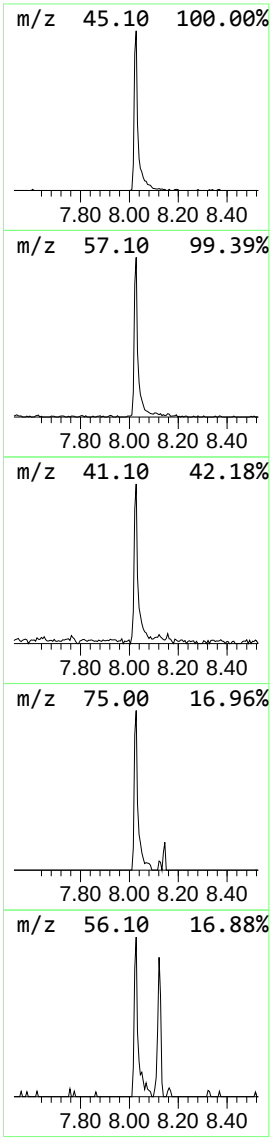
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.027	5.13 ng	281401	Naphthalene-d8	8.122

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2	Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	90
3	Ethanol, 2-[2-(2-methylpropoxy)e...	162	C8H18O3	018912-80-6	78
4	Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	59
5	2-Heptanol, 3-methyl-	130	C8H18O	031367-46-1	50





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111121\
Data File : BF126136.D
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ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-13N-5.0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

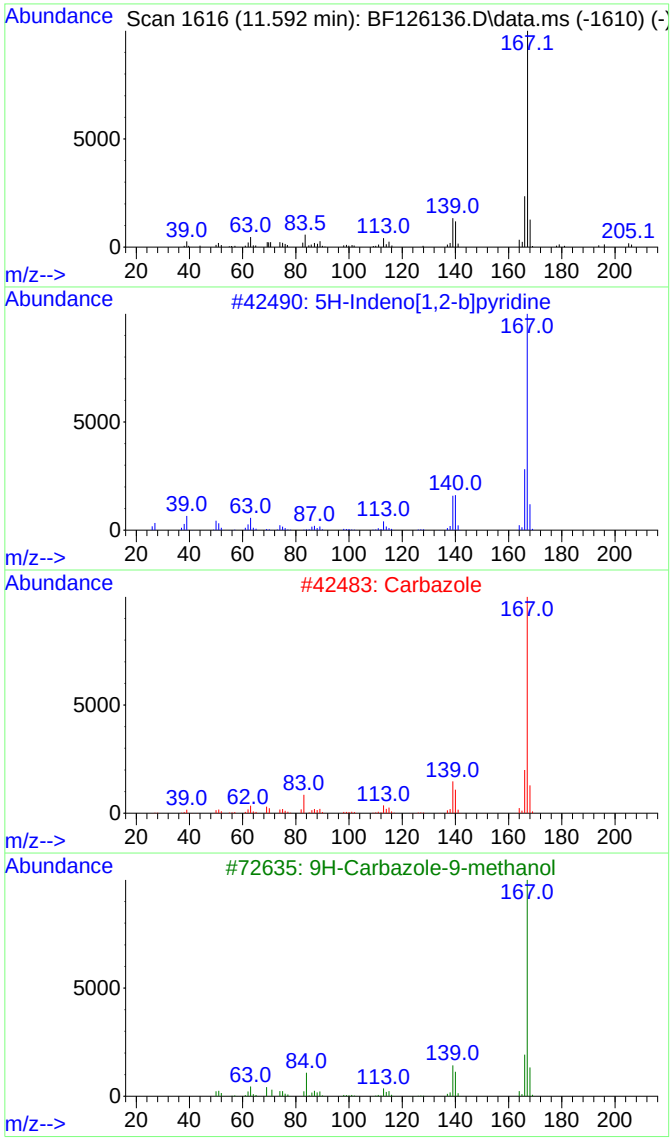
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

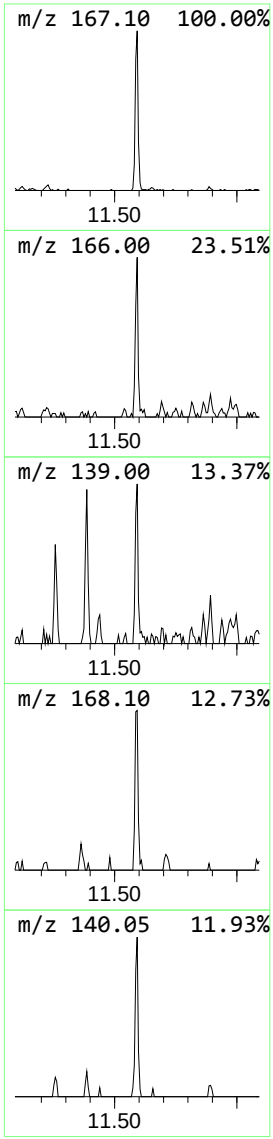
****

Peak Number 4 5H-Indeno[1,2-b]pyridine Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.592	2.21 ng	150914	Phenanthrene-d10	11.363

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5H-Indeno[1,2-b]pyridine	167	C12H9N	000244-99-5	94
2	Carbazole	167	C12H9N	000086-74-8	94
3	9H-Carbazole-9-methanol	197	C13H11NO	002409-36-1	83
4	9H-Carbazole, 9-nitroso-	196	C12H8N2O	002788-23-0	83
5	1-Naphthaleneacetonitrile	167	C12H9N	000132-75-2	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111121\
Data File : BF126136.D
Acq On : 11 Nov 2021 20:55
Operator : JU/CG
Sample : M4630-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-13N-5.0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

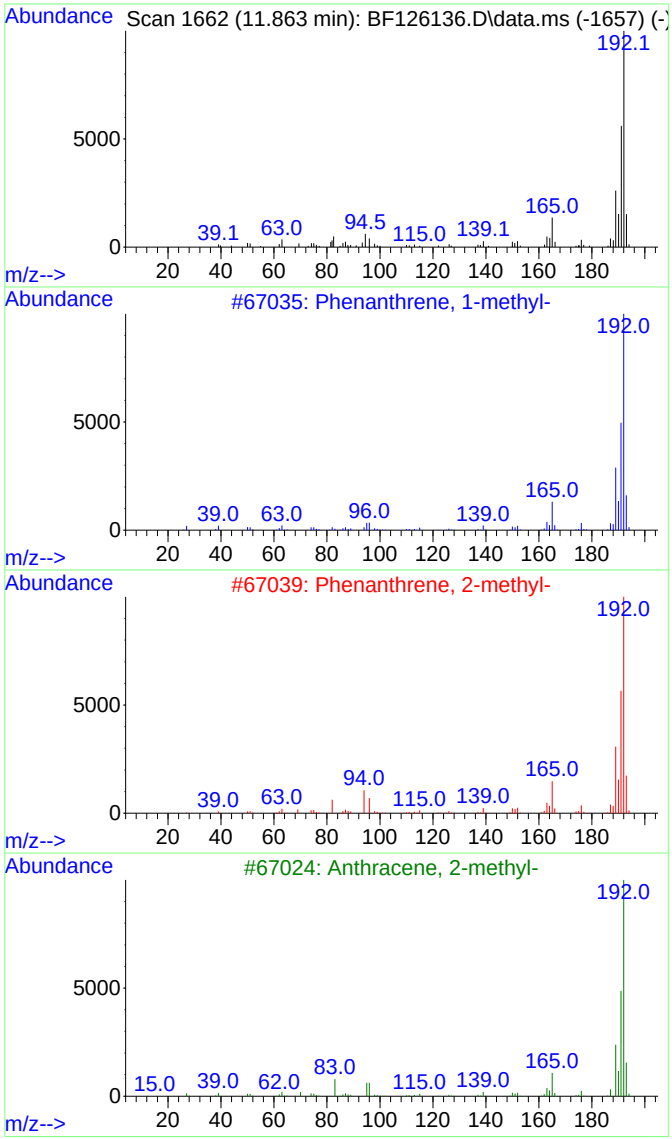
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

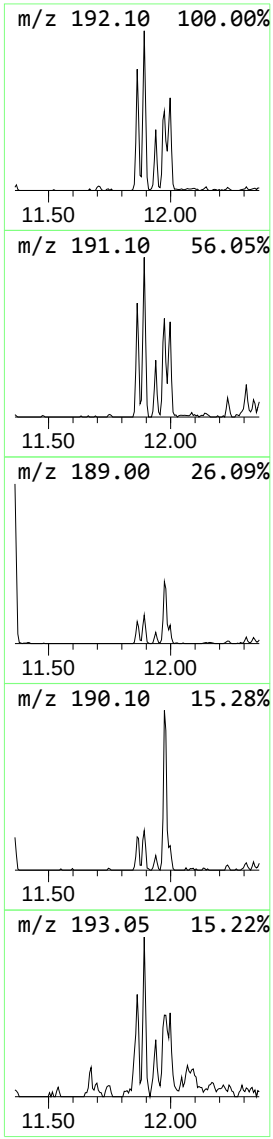
****

Peak Number 5 Phenanthrene, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.863	2.20 ng	150277	Phenanthrene-d10	11.363

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
2	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
3	Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
4	Anthracene, 9-methyl-	192	C15H12	000779-02-2	95
5	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	94





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Acq On : 11 Nov 2021 20:55
Operator : JU/CG
Sample : M4630-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-13N-5.0

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

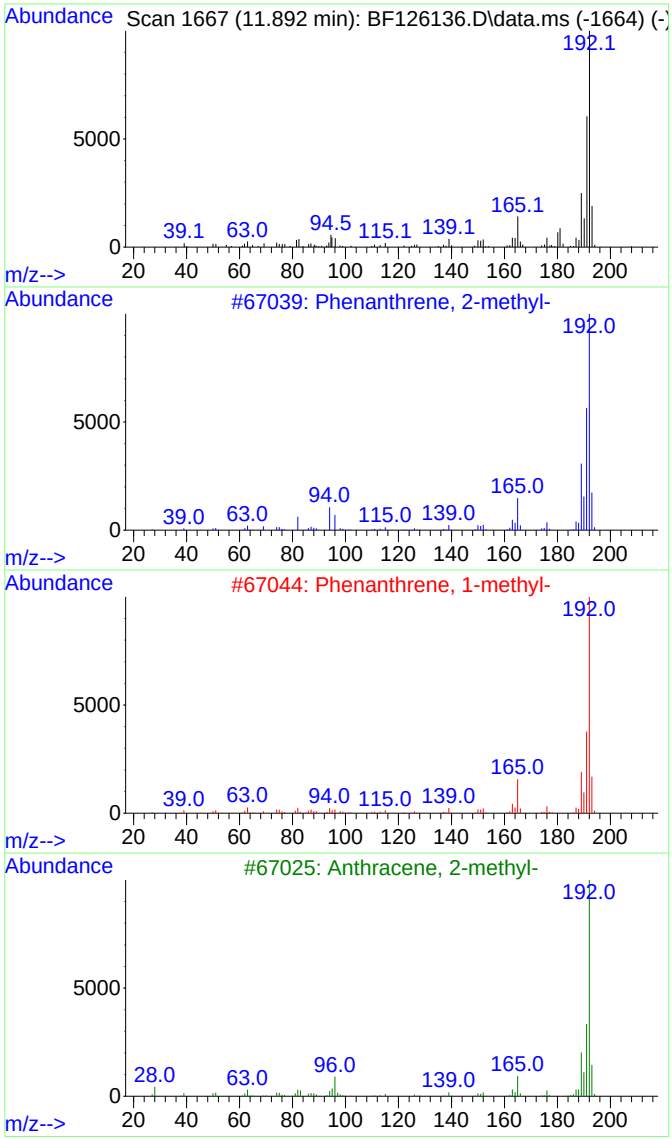
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

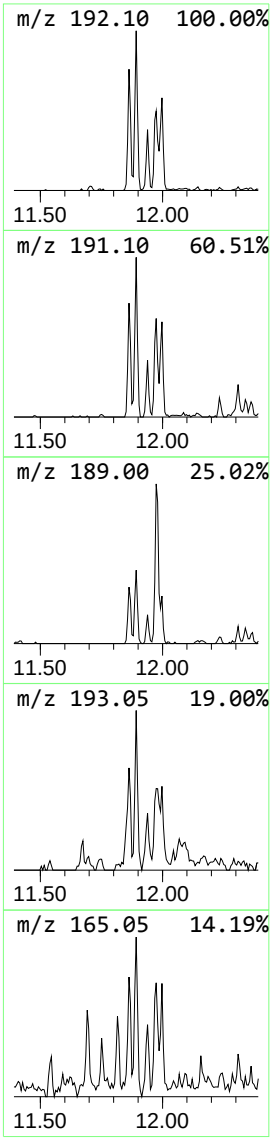
****

Peak Number 6 Phenanthrene, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.892	3.12 ng	213106	Phenanthrene-d10	11.363

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
3	Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
4	Anthracene, 9-methyl-	192	C15H12	000779-02-2	91
5	Anthracene, 1-methyl-	192	C15H12	000610-48-0	91





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Acq On : 11 Nov 2021 20:55
Operator : JU/CG
Sample : M4630-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-13N-5.0

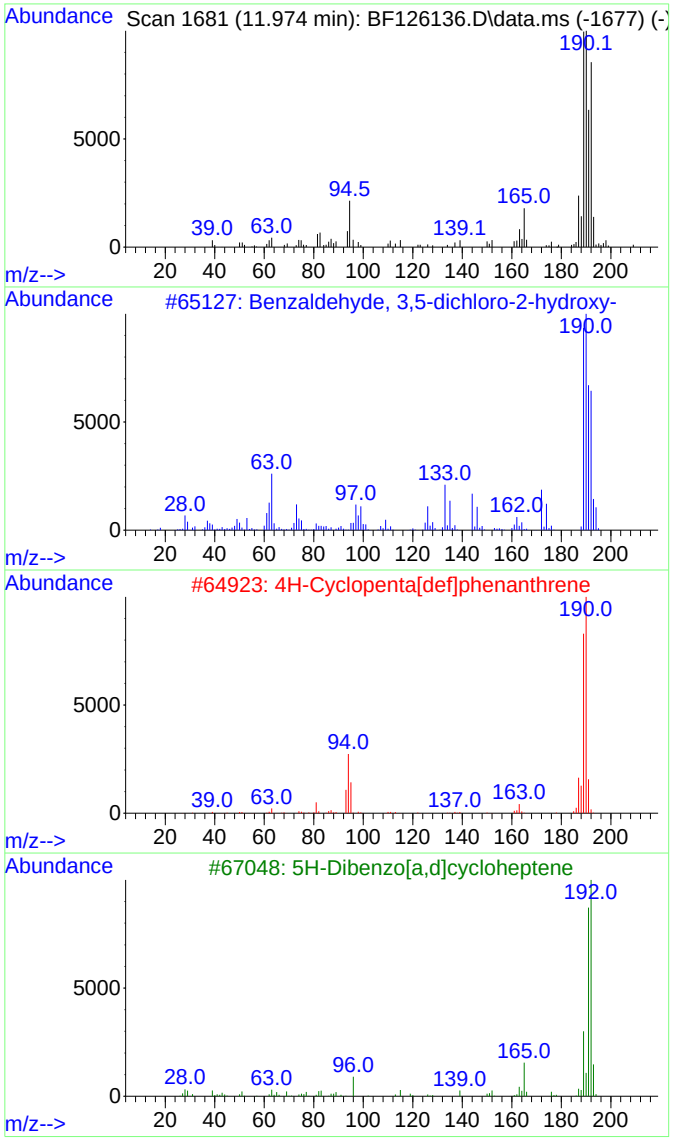
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

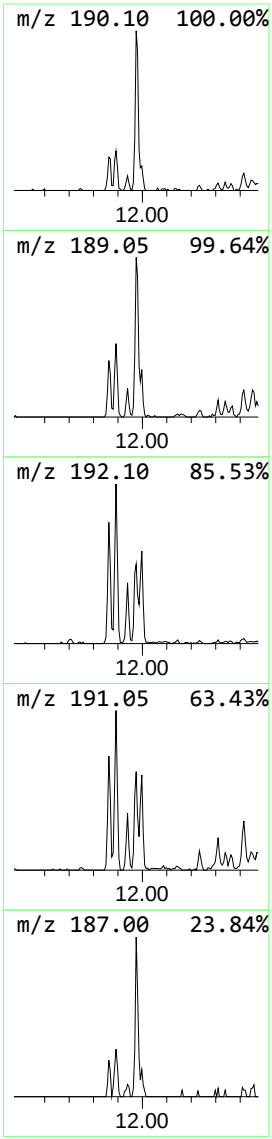
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzaldehyde, 3,5-dichloro-... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.974	3.34 ng	228606	Phenanthrene-d10	11.363

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	46
3	5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	42
4	Anthracene, 1-methyl-	192	C15H12	000610-48-0	42
5	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	42





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Acq On : 11 Nov 2021 20:55
Operator : JU/CG
Sample : M4630-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-13N-5.0

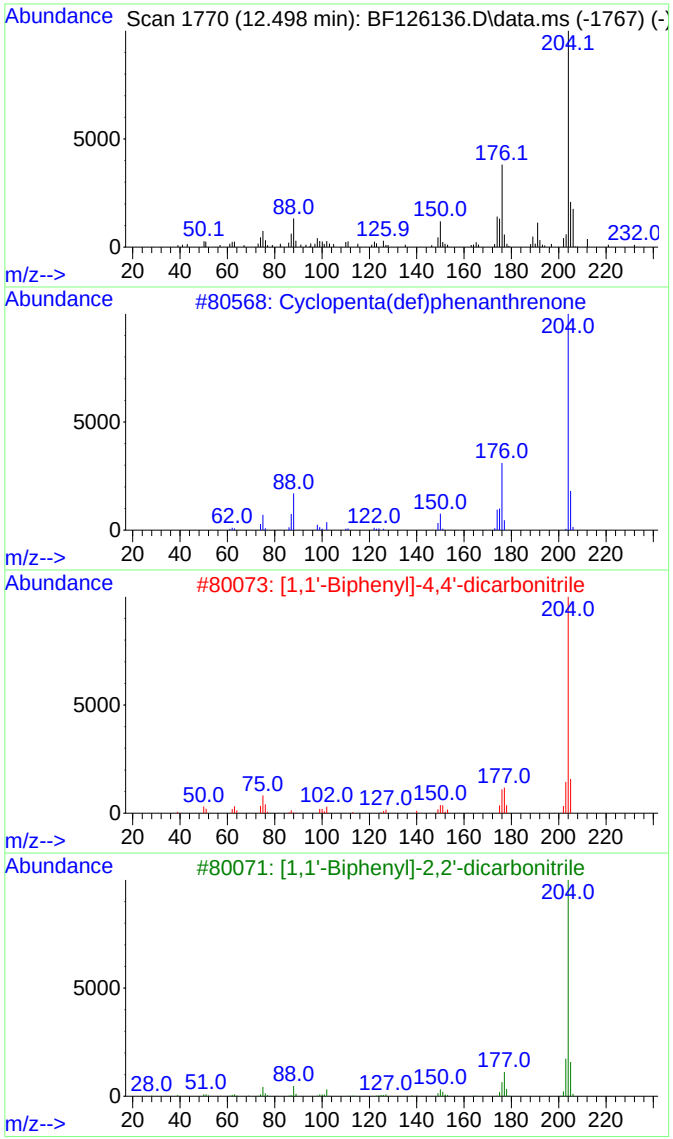
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

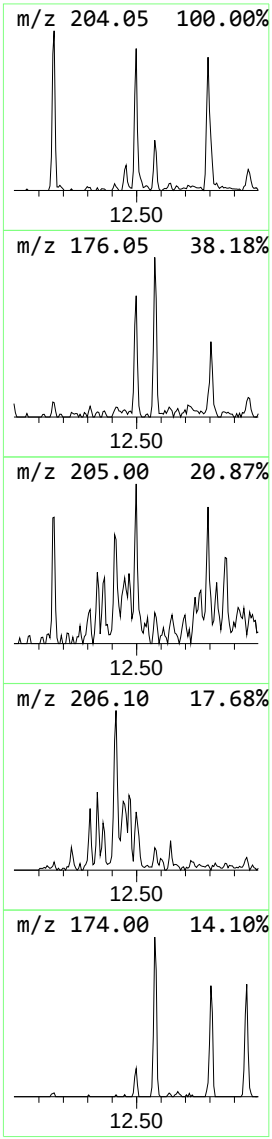
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Cyclopenta(def)phenanthrenone Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.498	2.06 ng	140959	Phenanthrene-d10	11.363

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	93
2	[1,1'-Biphenyl]-4,4'-dicarbonitrile	204	C14H8N2	001591-30-6	50
3	[1,1'-Biphenyl]-2,2'-dicarbonitrile	204	C14H8N2	004341-02-0	47
4	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	47
5	5,6,7,8-Tetrahydrothieno[2,3-b]q...	204	C11H12N2S	122914-50-5	45





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Sample : M4630-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-13N-5.0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

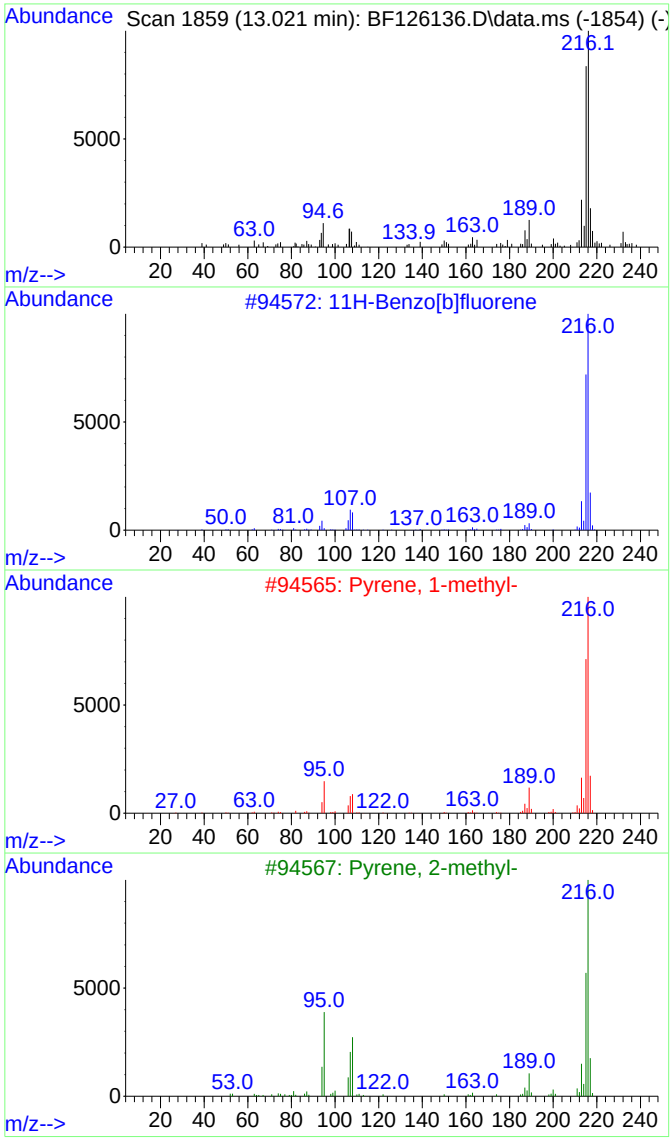
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

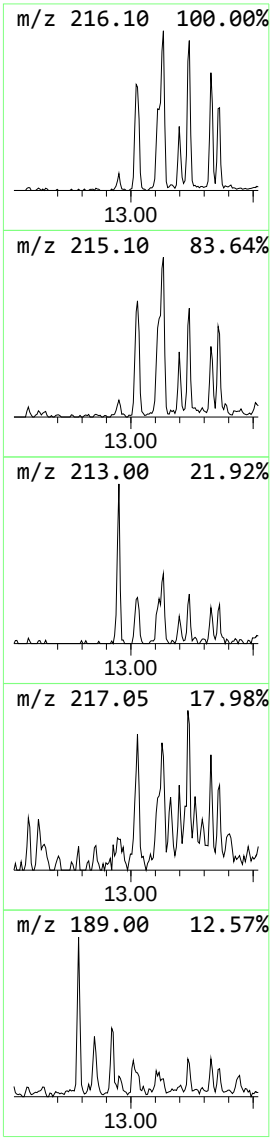
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.021	2.05 ng	195994	Chrysene-d12	14.004

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	95
2	Pyrene, 1-methyl-	216	C17H12	002381-21-7	94
3	Pyrene, 2-methyl-	216	C17H12	003442-78-2	93
4	7H-Benzo[c]fluorene	216	C17H12	000205-12-9	83
5	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	81





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Data File : BF126136.D
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Operator : JU/CG
Sample : M4630-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-13N-5.0

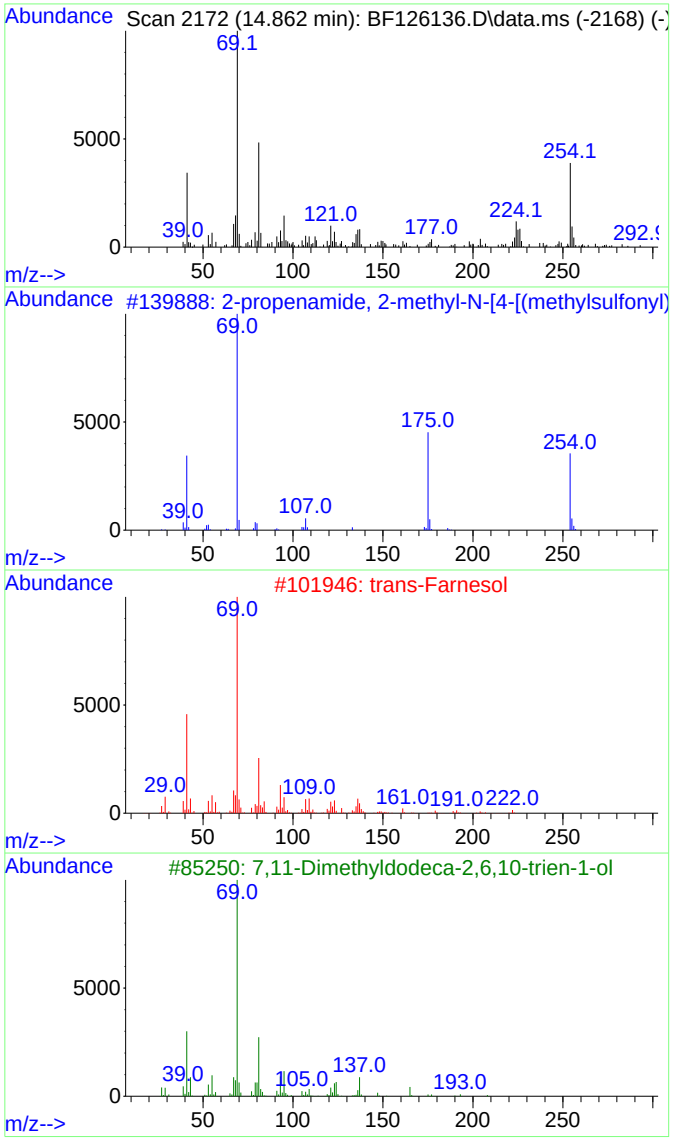
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

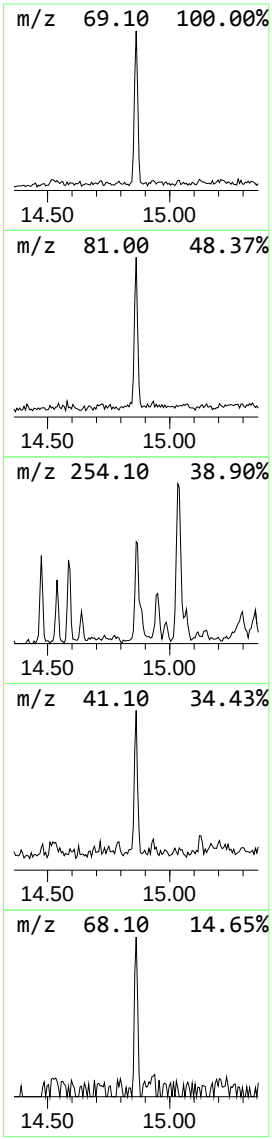
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 2-propenamide, 2-methyl-N-[... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.862	4.59 ng	258959	Perylene-d12	15.468

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-propenamide, 2-methyl-N-[4-[(m...	254	C11H14N2O3S	1000401-43-5	76
2	trans-Farnesol	222	C15H26O	000106-28-5	55
3	7,11-Dimethyldodeca-2,6,10-trien...	208	C14H24O	125529-10-4	46
4	1,3-Dioxolan-2-one, 3-methyl-3-(...	264	C16H24O3	336879-76-6	46
5	Farnesol isomer a	222	C15H26O	1000108-92-4	43





Tentatively Identified Compound (LSC) summary

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BNA_F
ClientSampleId :
SB-13N-5.0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110321.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
3-Hexen-2-one	4.416	2.3	ng	99121	1	6.839	857434	20.0
2-Pentanone, 4-...	5.045	3.7	ng	159044	1	6.839	857434	20.0
Ethanol, 2-(2-b...	8.027	5.1	ng	281401	2	8.122	1096460	20.0
5H-Indeno[1,2-b...	11.592	2.2	ng	150914	4	11.363	1367370	20.0
Phenanthrene, 1...	11.863	2.2	ng	150277	4	11.363	1367370	20.0
Phenanthrene, 2...	11.892	3.1	ng	213106	4	11.363	1367370	20.0
Benzaldehyde, 3...	11.974	3.3	ng	228606	4	11.363	1367370	20.0
Cyclopenta(def)...	12.498	2.1	ng	140959	4	11.363	1367370	20.0
11H-Benzo[b]flu...	13.021	2.0	ng	195994	5	14.004	1913740	20.0
2-propenamide, ...	14.862	4.6	ng	258959	6	15.468	1128190	20.0