LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF011725\
Data File : BF141196.D
Acq On : 17 Jan 2025 16:48
Operator : RC/JU
Sample : Q1103-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : VNJ-241

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-BF011025.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF141196.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.193	9	17	29	rVB	72589	119499	2.72%	0.281%
2	5.022	485	498	507	rBV	246379	394864	8.99%	0.928%
3	5.422	560	566	572	rBV	2520685	3452209	78.57%	8.116%
4	6.434	732	738	750	rBV	2523034	3558383	80.99%	8.366%
5	6.810	796	802	807	rBV	966627	1198218	27.27%	2.817%
6	7.369	890	897	902	rBV	1746955	2336053	53.17%	5.492%
7	7.628	935	941	947	rVB	76688	99104	2.26%	0.233%
8	8.092	1014	1020	1037	rVB	1232495	1633697	37.18%	3.841%
9	8.310	1051	1057	1061	rBV	33349	44957	1.02%	0.106%
10	9.163	1196	1202	1207	rBV	3390146	4393639	100.00%	10.330%
11	9.704	1290	1294	1299	rBV	43042	55072	1.25%	0.129%
12	9.839	1311	1317	1322	rBV	1402033	1813574	41.28%	4.264%
13	10.557	1433	1439	1444	rBV	714048	876746	19.95%	2.061%
14	10.628	1445	1451	1456	rBV	2016009	2596559	59.10%	6.105%
15	10.745	1466	1471	1479	rVB4	67155	114372	2.60%	0.269%
16	10.863	1487	1491	1497	rVB	101778	128710	2.93%	0.303%
17	10.992	1509	1513	1520	rBV	78821	103909	2.36%	0.244%
18	11.328	1565	1570	1578	rVV2	1259428	1898549	43.21%	4.464%
19	11.398	1578	1582	1585	rVV	80959	115903	2.64%	0.272%
20	11.557	1603	1609	1617	rBV3	92894	190502	4.34%	0.448%
21	11.857	1650	1660	1666	rBV	690591	1098045	24.99%	2.582%
22	11.939	1670	1674	1683	rVV2	104179	205332	4.67%	0.483%
23	12.122	1700	1705	1707	rBV2	100291	166969	3.80%	0.393%
24	12.275	1727	1731	1734	rBV2	50614	70449	1.60%	0.166%
25	12.380	1746	1749	1752	rVB	51787	60024	1.37%	0.141%
26	12.463	1759	1763	1769	rVB	114184	154003	3.51%	0.362%
27	12.539	1771	1776	1789	rBV	1380610	1976684	44.99%	4.647%
28	12.645	1789	1794	1805	rVB5	264931	493372	11.23%	1.160%
29	12.769	1808	1815	1821	rBV2	1071634	1587477	36.13%	3.732%
30	12.916	1831	1840	1847	rVB	2481989	3259645	74.19%	7.664%
31	12.986	1847	1852	1860	rBV4	116124	238778	5.43%	0.561%
32	13.080	1863	1868	1874	rVB3	126262	268717	6.12%	0.632%
33	13.198	1884	1888	1892	rBV	135128	185276	4.22%	0.436%
34	13.292	1901	1904	1907	rBV	55403	69364	1.58%	0.163%
35	13.327	1907	1910	1914	rVV4	35467	53336	1.21%	0.125%
36	13.604	1952	1957	1962	rBV	174558	253023	5.76%	0.595%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF011725\
Data File : BF141196.D
Acq On : 17 Jan 2025 16:48
Operator : RC/JU
Sample : Q1103-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
VNJ-241

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-BF011025.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	13.716	1971	1976	1979	rBV2	155509	261127	5.94%	0.614%
38	13.810	1988	1992	1997	rVV2	204414	300094	6.83%	0.706%
39	13.857	1998	2000	2003	rVB	62758	62180	1.42%	0.146%
40	13.963	2011	2018	2030	rBV2	1060445	2426250	55.22%	5.704%
41	14.351	2082	2084	2087	rVB2	60290	63124	1.44%	0.148%
42	14.392	2087	2091	2095	rVB2	95906	123940	2.82%	0.291%
43	14.451	2097	2101	2104	rBV3	84873	138113	3.14%	0.325%
44	14.604	2124	2127	2131	rBV	57526	77157	1.76%	0.181%
45	14.657	2132	2136	2145	rVB2	72359	139544	3.18%	0.328%
46	14.998	2189	2194	2204	rBV	558905	1204416	27.41%	2.832%
47	15.086	2205	2209	2211	rBV2	106956	172108	3.92%	0.405%
48	15.298	2240	2245	2250	rVB	265305	403422	9.18%	0.948%
49	15.357	2250	2255	2260	rBV	219679	310291	7.06%	0.730%
50	15.421	2260	2266	2271	rBV	711952	1122112	25.54%	2.638%
51	15.898	2342	2347	2351	rBV	98823	170141	3.87%	0.400%
52	16.851	2505	2509	2517	rVB2	72821	146337	3.33%	0.344%
53	17.286	2579	2583	2598	rVB	59640	148174	3.37%	0.348%

Sum of corrected areas: 42533543

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF011725\
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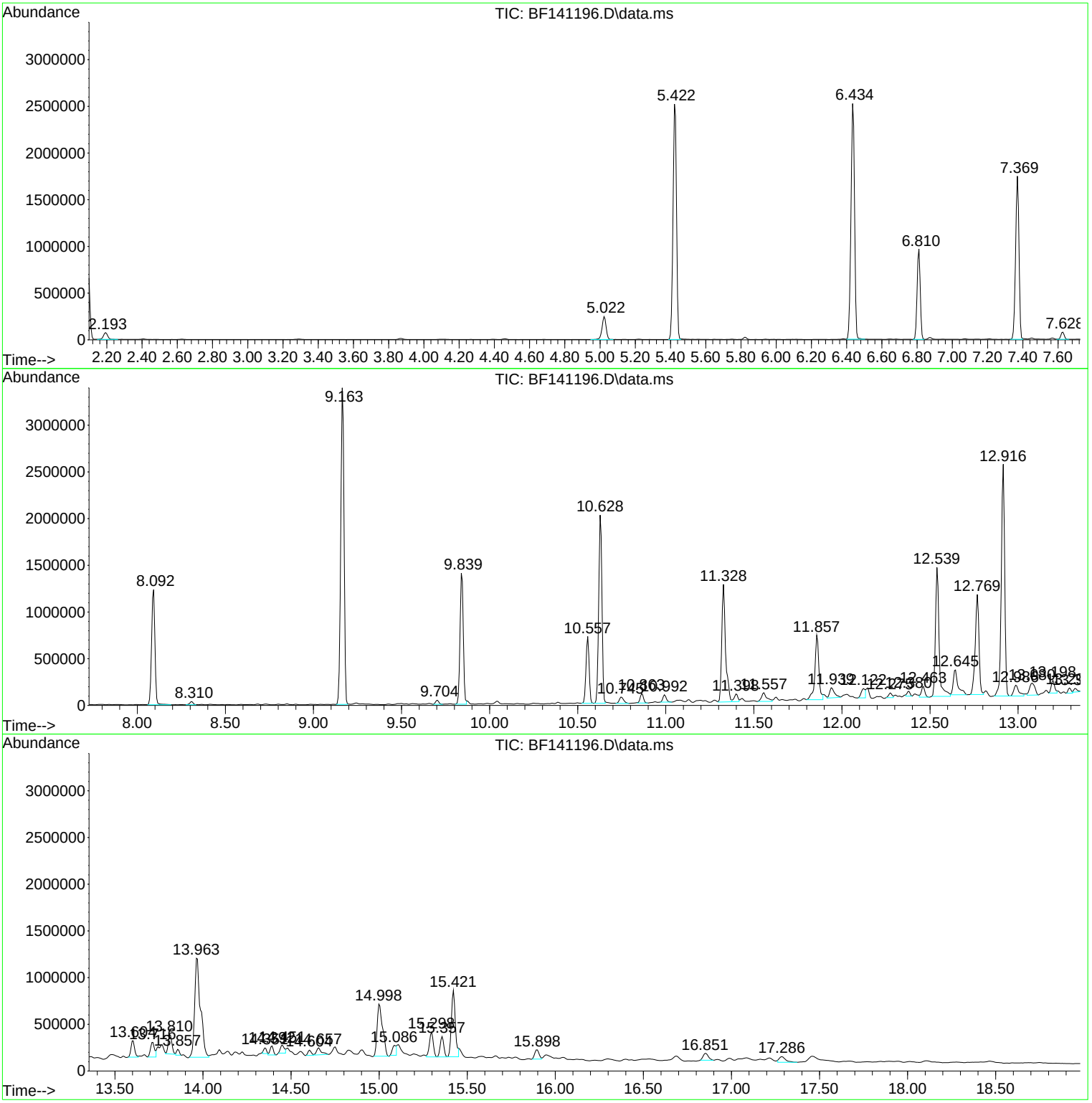
Acq On : 17 Jan 2025 16:48
Operator : RC/JU
Sample : Q1103-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
VNJ-241

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF011025.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\BF011725\
Data File : BF141196.D
Acq On : 17 Jan 2025 16:48
Operator : RC/JU
Sample : Q1103-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
VNJ-241

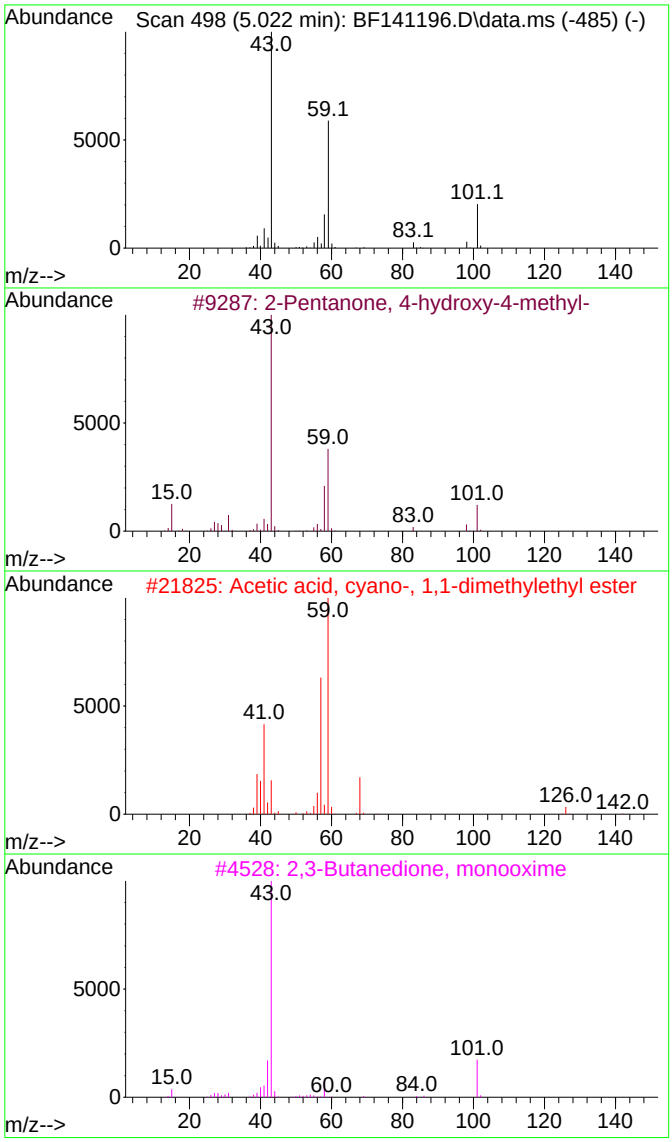
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF011025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

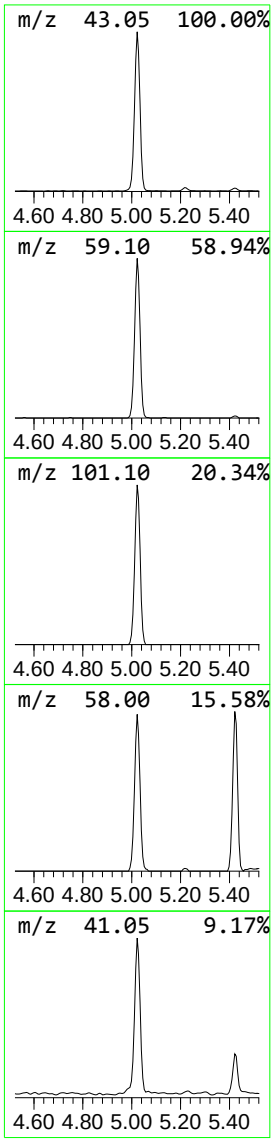
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.022	6.59 ng	394864	1,4-Dichlorobenzene-d4	6.810

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, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	32
3	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	12
4	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
5	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF011725\
Data File : BF141196.D
Acq On : 17 Jan 2025 16:48
Operator : RC/JU
Sample : Q1103-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
VNJ-241

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF011025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

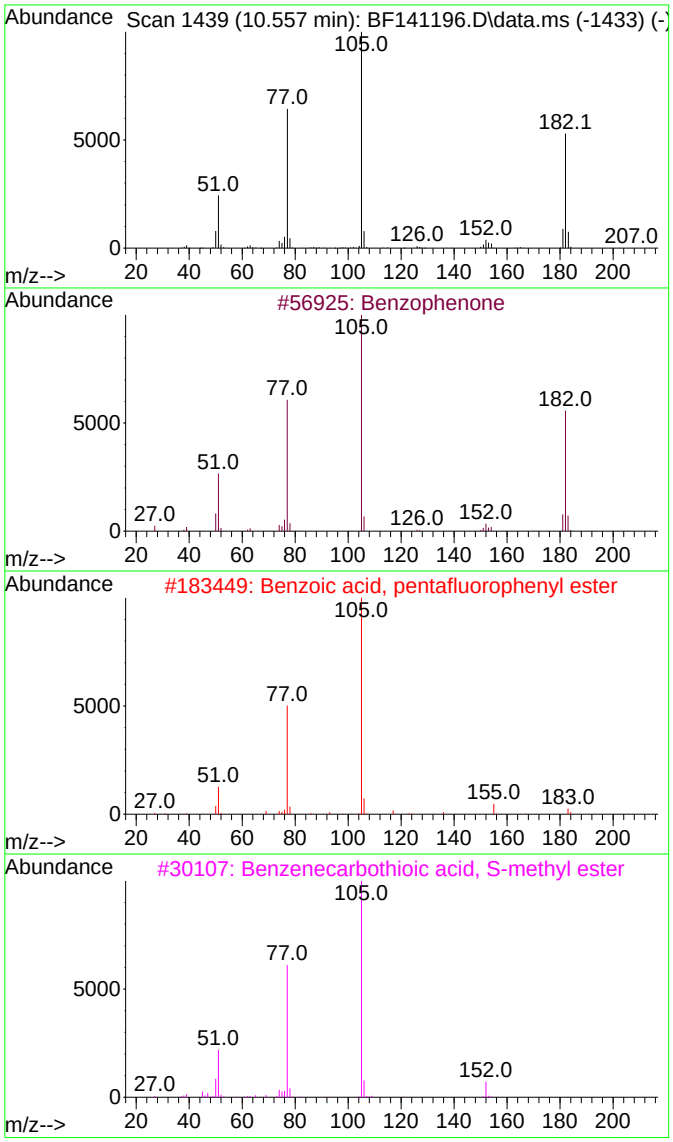
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

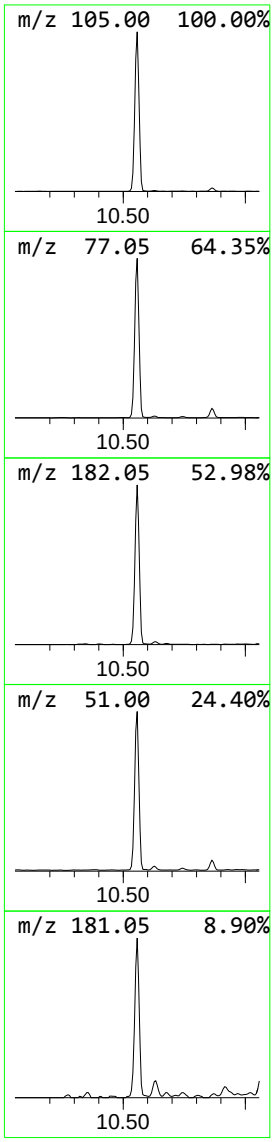
****

Peak Number 2 Benzophenone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.557	9.67 ng	876746	Acenaphthene-d10	9.839

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzophenone	182	C13H10O	000119-61-9	96
2	Benzoic acid, pentafluorophenyl ...	288	C13H5F5O2	1000360-67-9	50
3	Benzenecarbothioic acid, S-methy...	152	C8H8OS	005925-68-8	50
4	Acetophenone, 2-chloro-	154	C8H7ClO	000532-27-4	49
5	N-Methylbenzamide, N-trifluoroac...	231	C10H8F3NO2	1000446-91-5	47





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF011725\
Data File : BF141196.D
Acq On : 17 Jan 2025 16:48
Operator : RC/JU
Sample : Q1103-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
VNJ-241

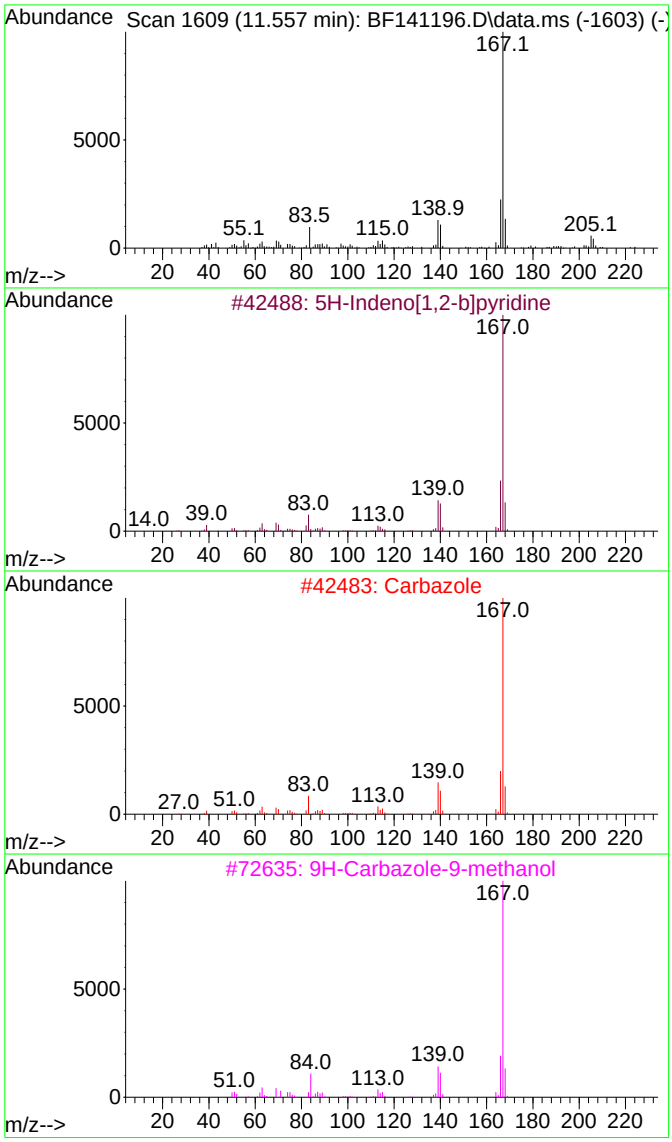
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF011025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

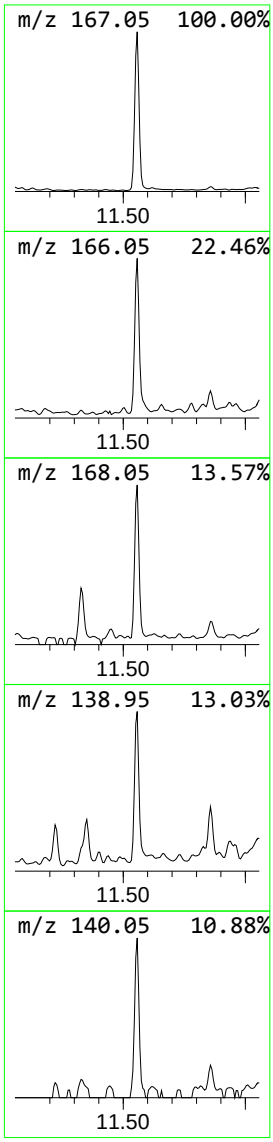
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 5H-Indeno[1,2-b]pyridine Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.557	2.01 ng	190502	Phenanthrene-d10	11.328

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	87
3	9H-Carbazole-9-methanol	197	C13H11NO	002409-36-1	74
4	2-Azafluorene	167	C12H9N	000244-40-6	64
5	Benzene, 1,1'-(chloromethylene)bis-	202	C13H11Cl	000090-99-3	58





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF011725\
Data File : BF141196.D
Acq On : 17 Jan 2025 16:48
Operator : RC/JU
Sample : Q1103-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
VNJ-241

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF011025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

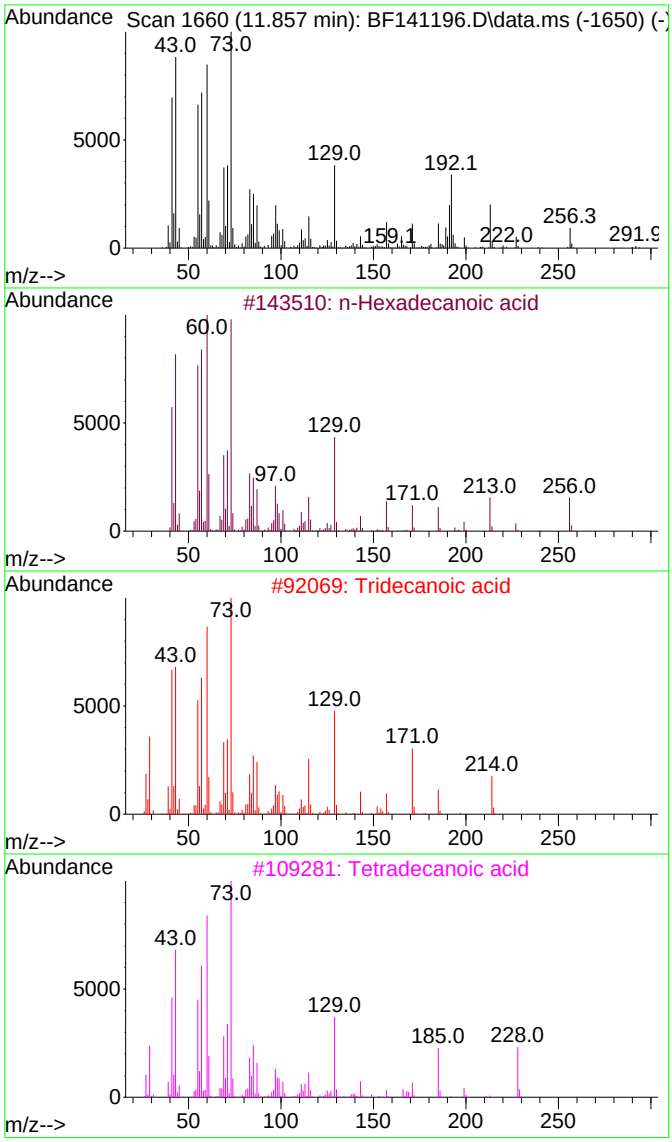
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

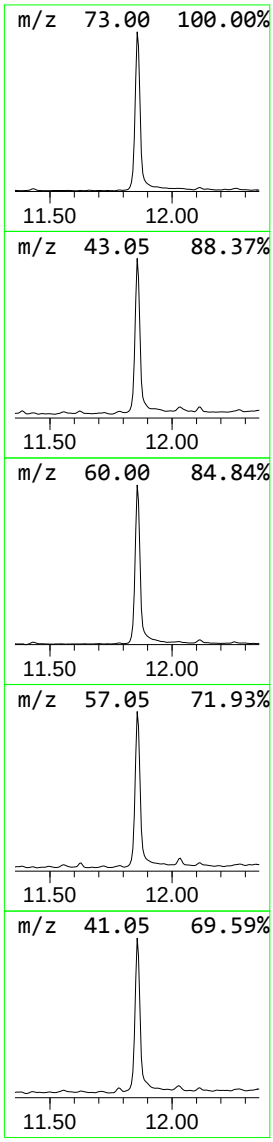
****

Peak Number 4 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.857	11.57 ng	1098050	Phenanthrene-d10	11.328

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	Tridecanoic acid	214	C13H26O2	000638-53-9	91
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	90
4	Pentadecanoic acid	242	C15H30O2	001002-84-2	90
5	Octadecanoic acid	284	C18H36O2	000057-11-4	81





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF011725\
Data File : BF141196.D
Acq On : 17 Jan 2025 16:48
Operator : RC/JU
Sample : Q1103-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
VNJ-241

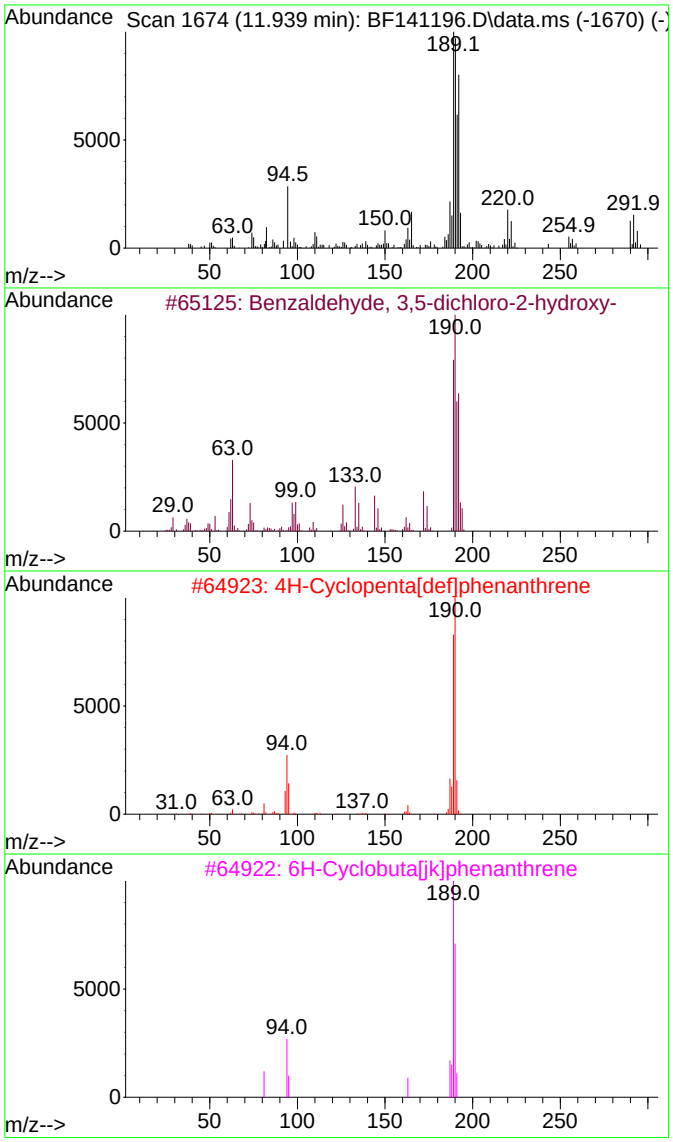
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF011025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

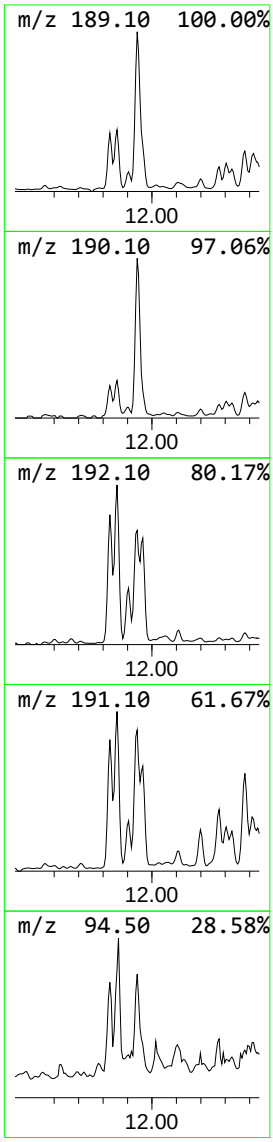
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Benzaldehyde, 3,5-dichloro-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.939	2.16 ng	205332	Phenanthrene-d10	11.328

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	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	46
4	Methyl 3,5-dichloro-4-hydroxyben...	220	C8H6Cl2O3	003337-59-5	38
5	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	30





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF011725\
Data File : BF141196.D
Acq On : 17 Jan 2025 16:48
Operator : RC/JU
Sample : Q1103-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
VNJ-241

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF011025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

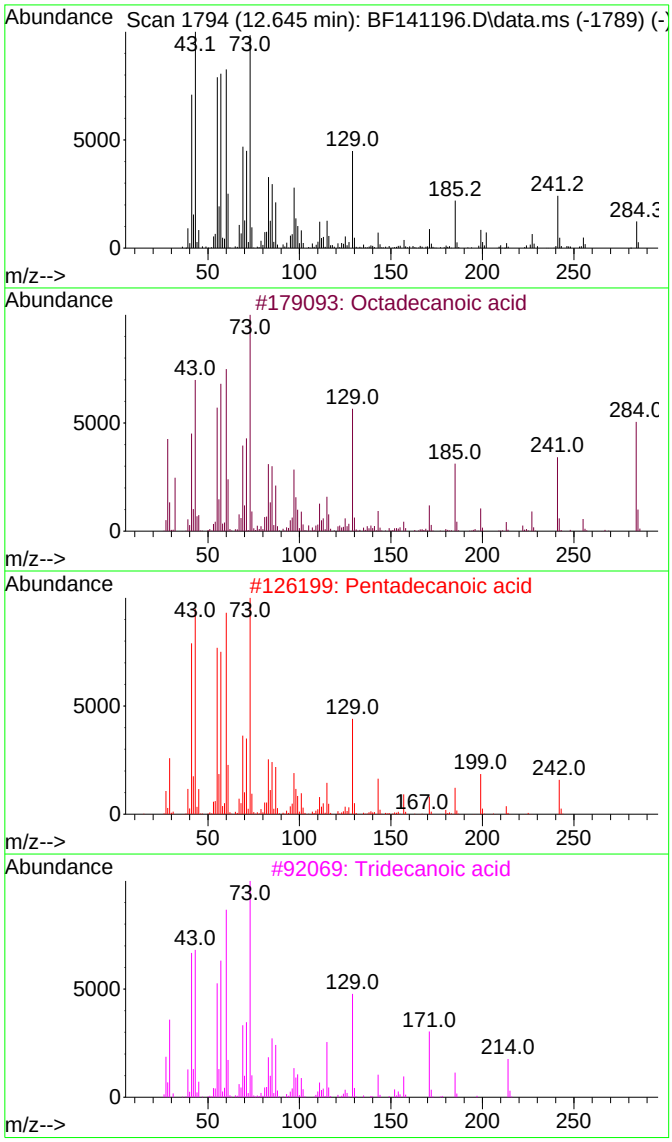
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

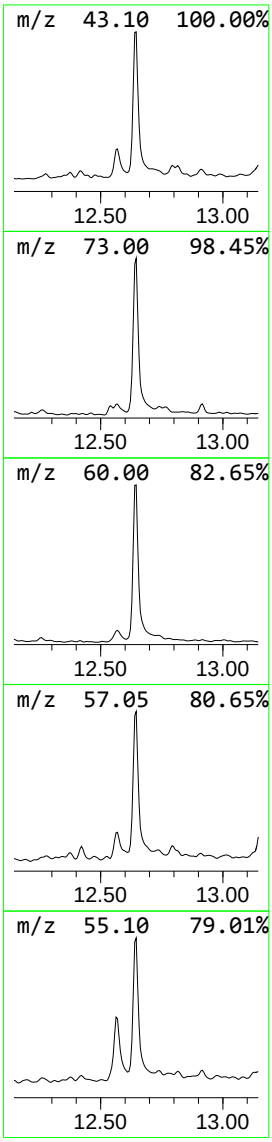
****

Peak Number 6 Octadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.645	5.20 ng	493372	Phenanthrene-d10	11.328

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	90
3	Tridecanoic acid	214	C13H26O2	000638-53-9	74
4	n-Decanoic acid	172	C10H20O2	000334-48-5	70
5	Tetradecanoic acid	228	C14H28O2	000544-63-8	64





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Data File : BF141196.D
Acq On : 17 Jan 2025 16:48
Operator : RC/JU
Sample : Q1103-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
VNJ-241

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF011025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

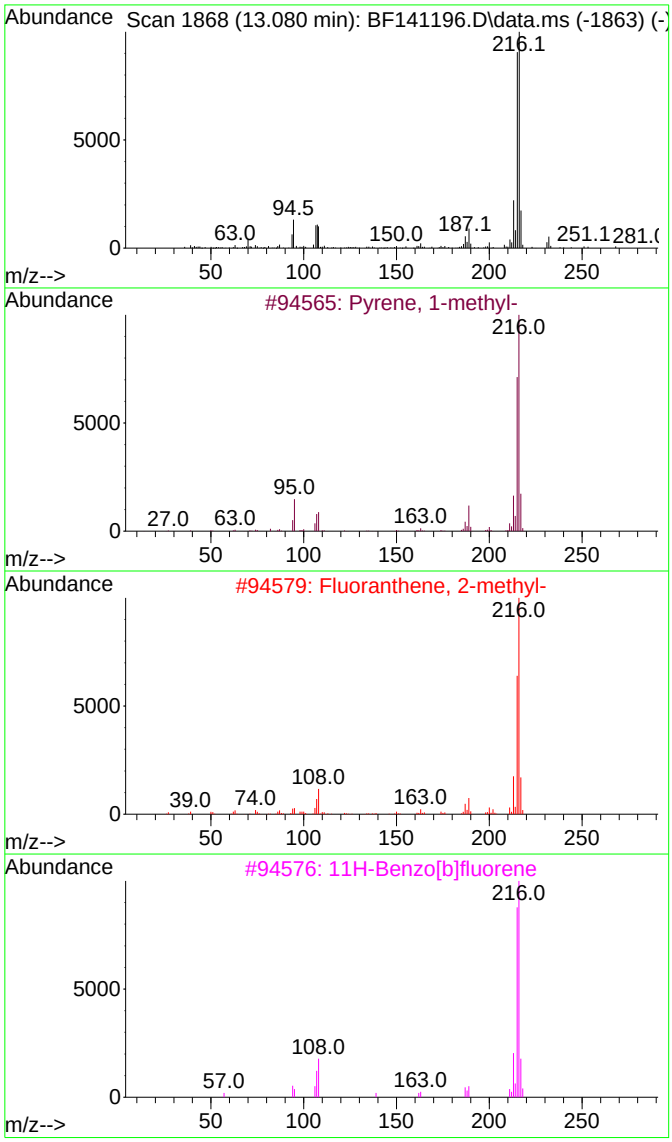
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

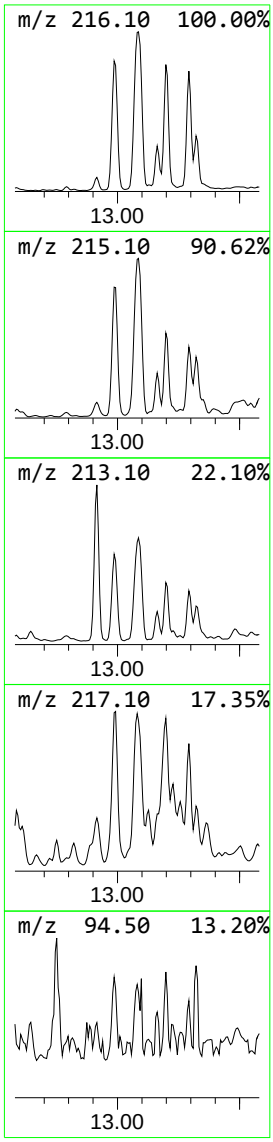
****

Peak Number 7 Pyrene, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.080	2.22 ng	268717	Chrysene-d12	13.969

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-	216	C17H12	002381-21-7	93
2	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	93
3	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	90
4	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	89
5	Pyrene, 2-methyl-	216	C17H12	003442-78-2	83





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Data File : BF141196.D
Acq On : 17 Jan 2025 16:48
Operator : RC/JU
Sample : Q1103-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
VNJ-241

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF011025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

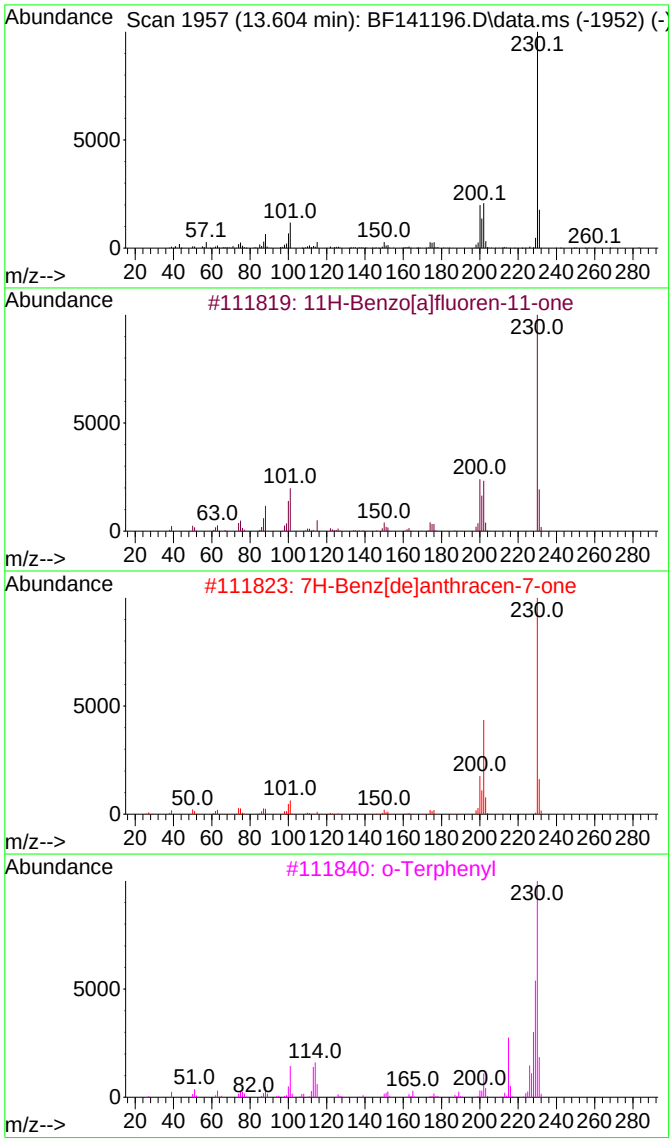
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

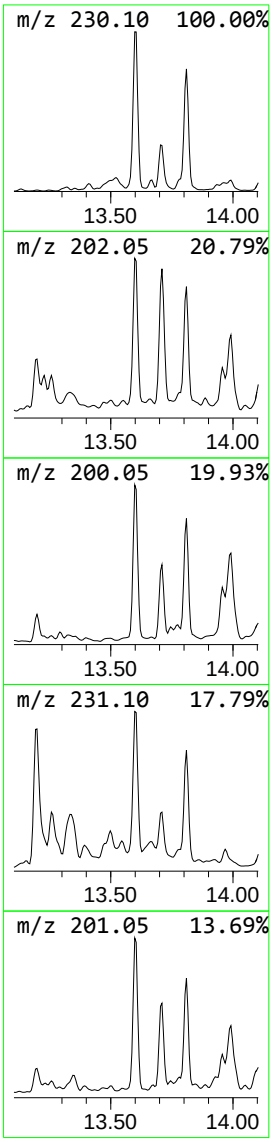
****

Peak Number 8 11H-Benzo[a]fluoren-11-one Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.604	2.09 ng	253023	Chrysene-d12	13.969

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[a]fluoren-11-one	230	C17H10O	000479-79-8	97
2	7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	90
3	o-Terphenyl	230	C18H14	000084-15-1	58
4	5,6-Dihydrochrysene	230	C18H14	002091-92-1	50
5	5-Amino-7-imino-4,8-diazatricycl...	230	C12H14N4O	1000388-35-6	50





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Acq On : 17 Jan 2025 16:48
Operator : RC/JU
Sample : Q1103-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
VNJ-241

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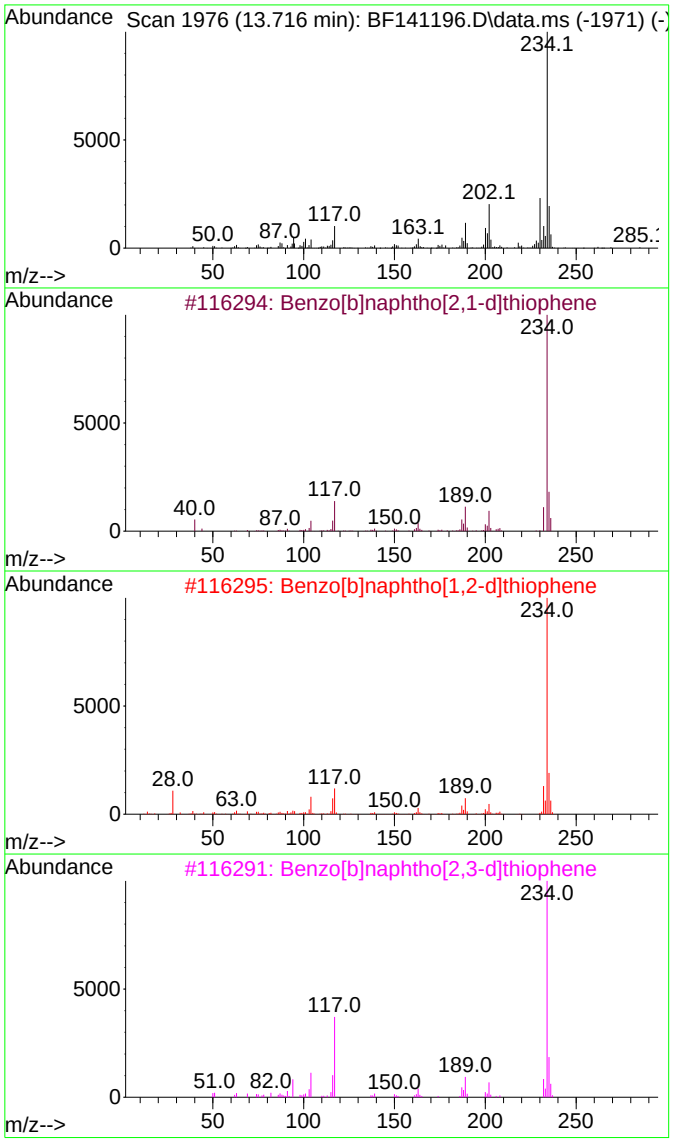
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

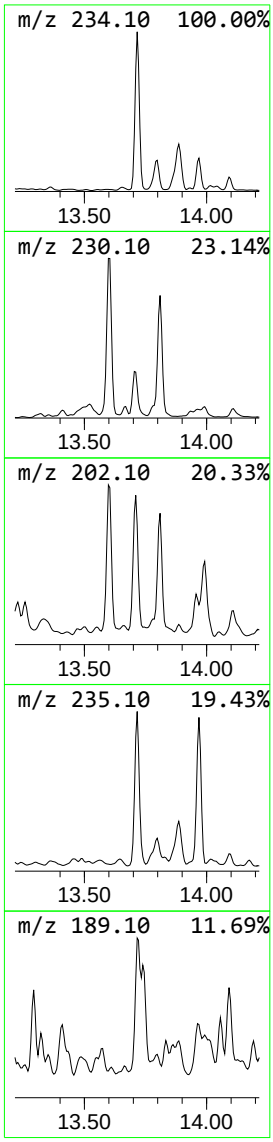
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzo[b]naphtho[2,1-d]thiop... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.716	2.15 ng	261127	Chrysene-d12	13.969

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[b]naphtho[2,1-d]thiophene	234	C16H10S	000239-35-0	95
2	Benzo[b]naphtho[1,2-d]thiophene	234	C16H10S	000205-43-6	86
3	Benzo[b]naphtho[2,3-d]thiophene	234	C16H10S	000243-46-9	70
4	Phenanthro(2,1-b)thiophene	234	C16H10S	000219-25-0	58
5	Phenanthro[4,3-b]thiophene	234	C16H10S	000195-68-6	52





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Sample : Q1103-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
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ClientSampleId :
VNJ-241

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF011025.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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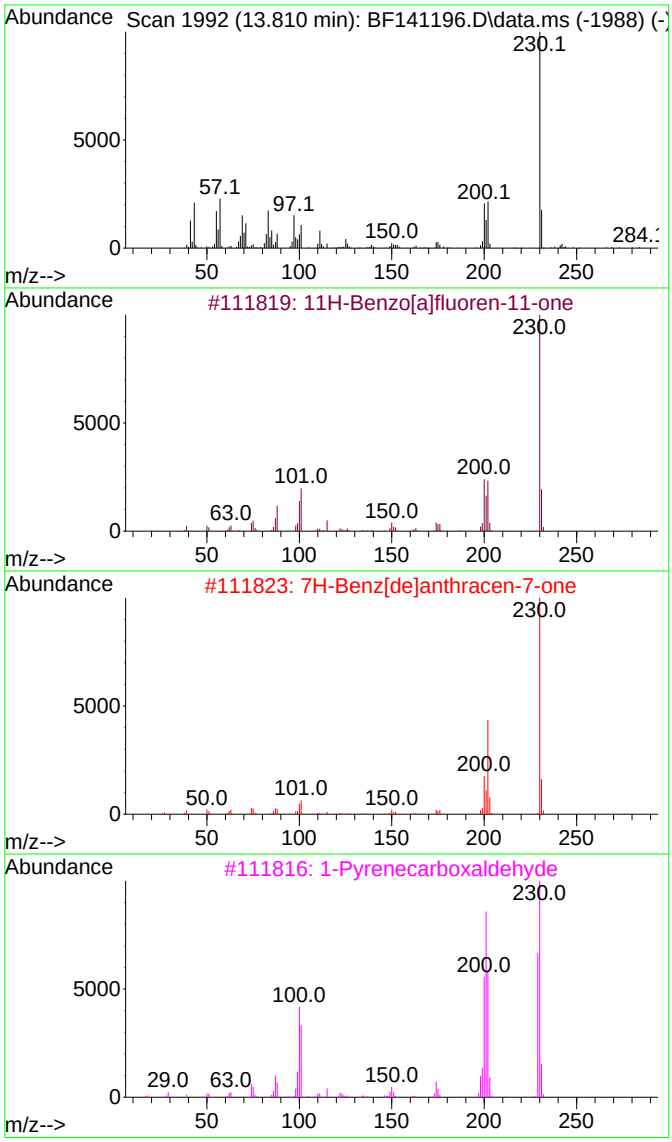
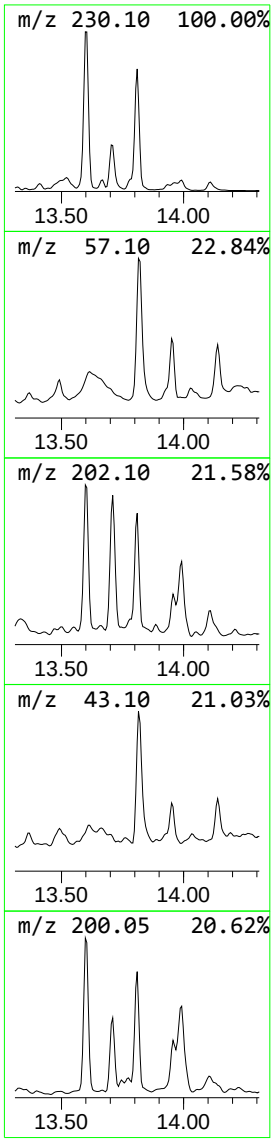
Peak Number 10 7H-Benz[de]anthracen-7-one Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.810	2.47 ng	300094	Chrysene-d12	13.969

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[a]fluoren-11-one	230	C17H10O	000479-79-8	80
2	7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	76
3	1-Pyrenecarboxaldehyde	230	C17H10O	003029-19-4	64
4	benzonitrile, 2-(3-isoquinolinyl)-	230	C16H10N2	1000401-00-3	53
5	Pyrene, 1,3-dimethyl-	230	C18H14	064401-21-4	50





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Sample : Q1103-01
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Instrument :
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ClientSampleId :
VNJ-241

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TIC Library : C:\Database\NIST20.L
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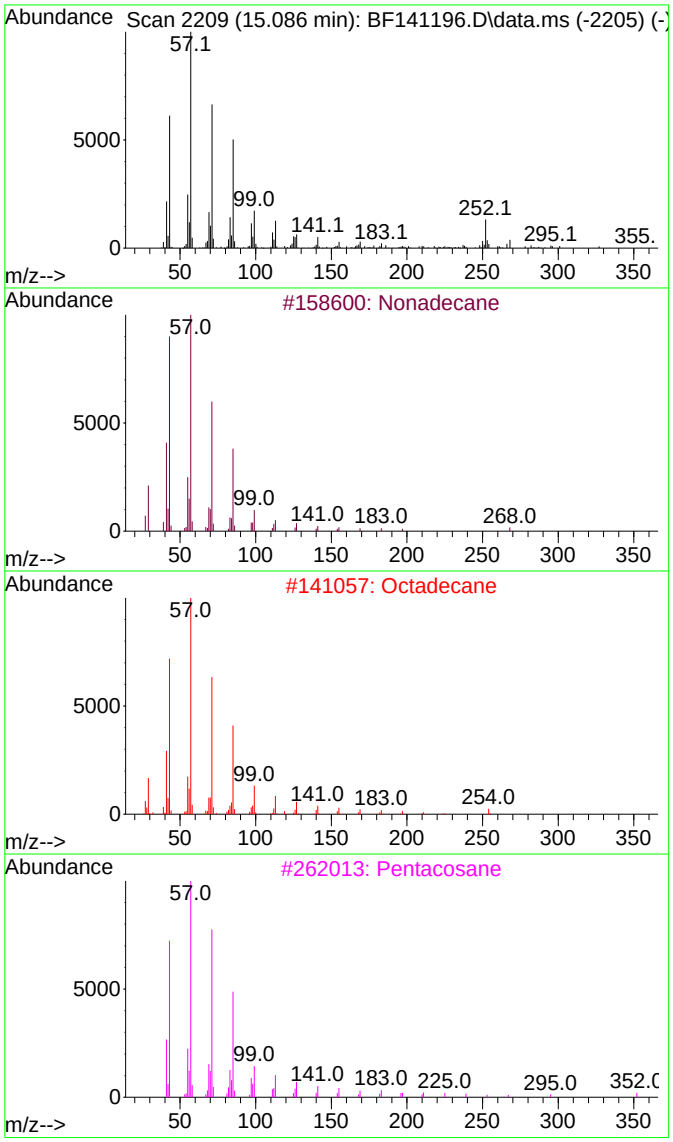
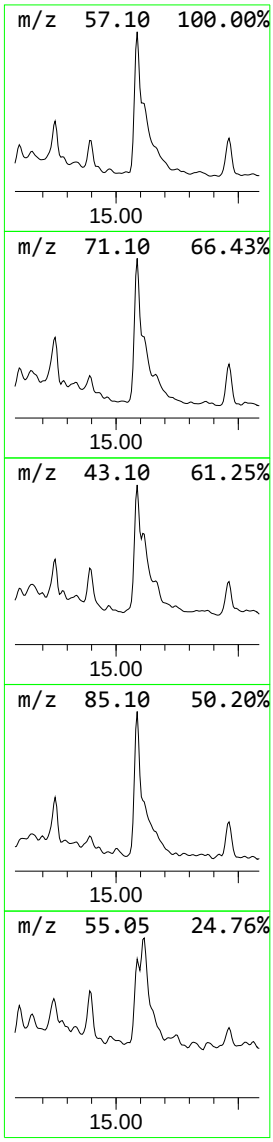
Peak Number 11 Nonadecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.086	3.07 ng	172108	Perylene-d12	15.421

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane	268	C19H40	000629-92-5	95
2	Octadecane	254	C18H38	000593-45-3	93
3	Pentacosane	352	C25H52	000629-99-2	87
4	Hentriacontane	437	C31H64	000630-04-6	83
5	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	83





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Acq On : 17 Jan 2025 16:48
Operator : RC/JU
Sample : Q1103-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
VNJ-241

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF011025.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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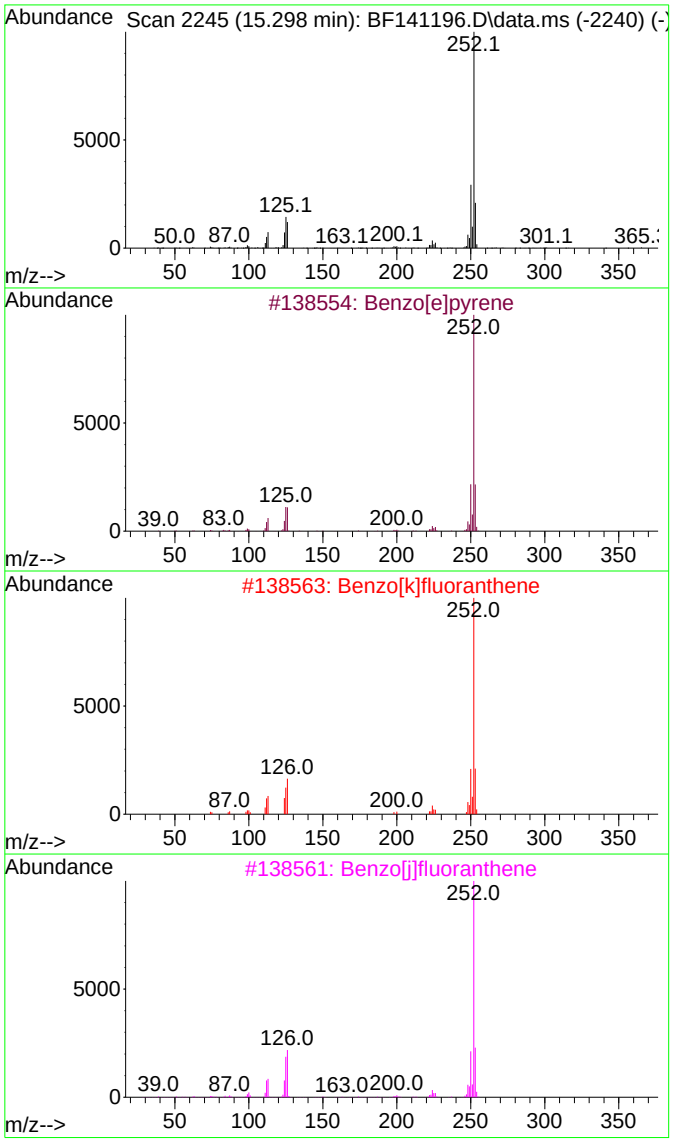
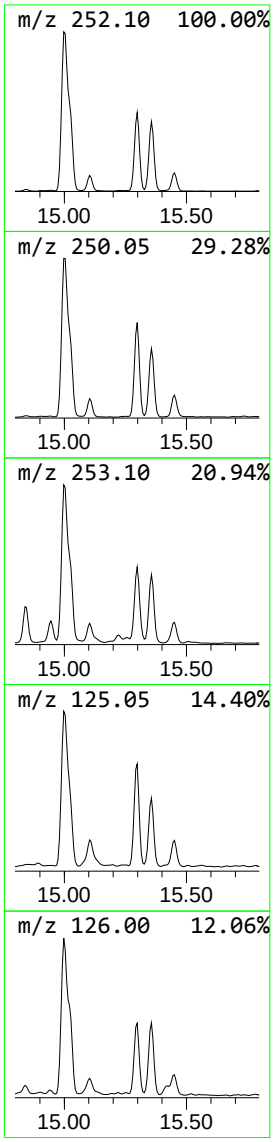
Peak Number 12 Benzo[e]pyrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.298	7.19 ng	403422	Perylene-d12	15.421

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	94
3	Benzo[j]fluoranthene	252	C20H12	000205-82-3	93
4	Benzo[b]fluoranthene	252	C20H12	000205-99-2	93
5	Benzo[a]pyrene	252	C20H12	000050-32-8	90





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Data File : BF141196.D
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Operator : RC/JU
Sample : Q1103-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
VNJ-241

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF011025.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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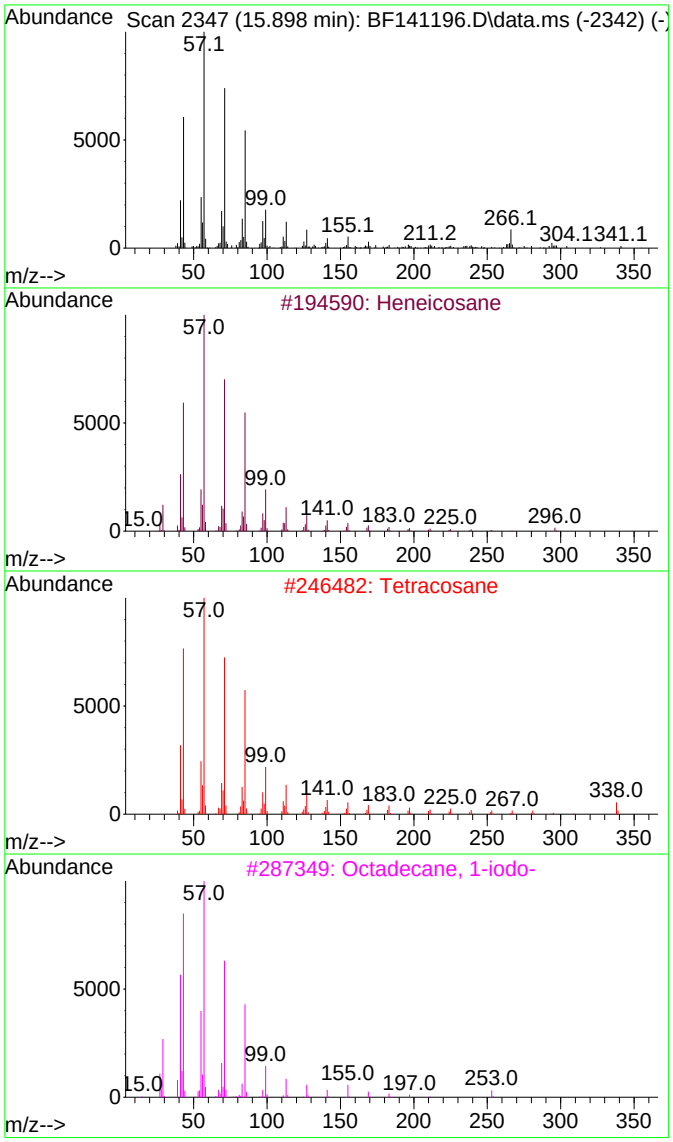
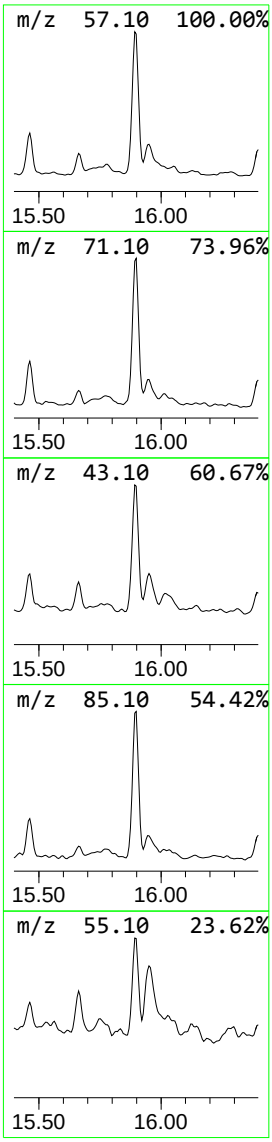
Peak Number 13 Heneicosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.898	3.03 ng	170141	Perylene-d12	15.421

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	99
2	Tetracosane	338	C24H50	000646-31-1	87
3	Octadecane, 1-iodo-	380	C18H37I	000629-93-6	87
4	Eicosane, 1-iodo-	408	C20H41I	1000406-31-8	87
5	Docosane, 1-iodo-	436	C22H45I	1000406-31-9	87





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF011725\
Data File : BF141196.D
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BNA_F
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF011025.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-...	5.022	6.6	ng	394864	1	6.810	1198220	20.0
Benzophenone	10.557	9.7	ng	876746	3	9.839	1813570	20.0
5H-Indeno[1,2-b...	11.557	2.0	ng	190502	4	11.328	1898550	20.0
n-Hexadecanoic ...	11.857	11.6	ng	1098050	4	11.328	1898550	20.0
Benzaldehyde, 3...	11.939	2.2	ng	205332	4	11.328	1898550	20.0
Octadecanoic acid	12.645	5.2	ng	493372	4	11.328	1898550	20.0
Pyrene, 1-methyl-	13.080	2.2	ng	268717	5	13.969	2426250	20.0
11H-Benzo[a]flu...	13.604	2.1	ng	253023	5	13.969	2426250	20.0
Benzo[b]naphtho...	13.716	2.1	ng	261127	5	13.969	2426250	20.0
7H-Benz[de]anth...	13.810	2.5	ng	300094	5	13.969	2426250	20.0
Nonadecane	15.086	3.1	ng	172108	6	15.421	1122110	20.0
Benzo[e]pyrene	15.298	7.2	ng	403422	6	15.421	1122110	20.0
Heneicosane	15.898	3.0	ng	170141	6	15.421	1122110	20.0