LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082124\
Data File : BF139108.D
Acq On : 21 Aug 2024 18:20
Operator : RC/JU
Sample : P3663-09
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
915-J-SD-0-0.5-081624

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-BF082024.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF139108.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.228	13	23	30	rVB	633324	1020250	52.01%	4.916%
2	4.005	318	325	332	rBV	38298	58900	3.00%	0.284%
3	5.122	505	515	523	rVB	82599	127529	6.50%	0.615%
4	5.516	576	582	587	rBV	1467834	1922138	97.99%	9.262%
5	6.516	745	752	762	rBV	1423694	1961563	100.00%	9.452%
6	6.898	811	817	822	rBV	434687	545317	27.80%	2.628%
7	7.457	906	912	917	rBV	873068	1134516	57.84%	5.467%
8	8.175	1029	1034	1039	rBV	518638	650400	33.16%	3.134%
9	8.487	1082	1087	1094	rVB	27260	34117	1.74%	0.164%
10	9.251	1212	1217	1222	rBV	1263826	1514847	77.23%	7.299%
11	9.928	1327	1332	1336	rBV	464449	592836	30.22%	2.857%
12	9.963	1336	1338	1342	rVB	24074	24093	1.23%	0.116%
13	10.028	1343	1349	1354	rVB	22320	32568	1.66%	0.157%
14	10.475	1421	1425	1431	rVB	28521	35117	1.79%	0.169%
15	10.716	1461	1466	1471	rBV	634660	778767	39.70%	3.753%
16	11.310	1564	1567	1573	rVB	20309	24405	1.24%	0.118%
17	11.416	1580	1585	1587	rBV	377230	502316	25.61%	2.420%
18	11.439	1587	1589	1594	rVV	440357	478344	24.39%	2.305%
19	11.492	1594	1598	1601	rVV	40296	48864	2.49%	0.235%
20	11.639	1617	1623	1632	rVB2	23215	35992	1.83%	0.173%
21	11.916	1665	1670	1672	rBV	27811	38530	1.96%	0.186%
22	11.945	1672	1675	1679	rVB2	61952	77811	3.97%	0.375%
23	12.028	1685	1689	1697	rVB2	76290	119194	6.08%	0.574%
24	12.233	1717	1724	1729	rVB2	56124	103302	5.27%	0.498%
25	12.551	1774	1778	1783	rVB2	17308	23355	1.19%	0.113%
26	12.627	1786	1791	1796	rBV	804894	1035882	52.81%	4.992%
27	12.675	1796	1799	1805	rBV2	20129	29108	1.48%	0.140%
28	12.780	1814	1817	1822	rVB	21126	27299	1.39%	0.132%
29	12.857	1822	1830	1845	rVB2	728839	1408747	71.82%	6.788%
30	13.004	1846	1855	1860	rVB	761411	1000549	51.01%	4.821%
31	13.075	1861	1867	1871	rBV2	34203	68810	3.51%	0.332%
32	13.186	1875	1886	1890	rBV2	90227	157386	8.02%	0.758%
33	13.251	1893	1897	1900	rBV	85196	103733	5.29%	0.500%
34	13.286	1900	1903	1906	rBV	33196	38707	1.97%	0.187%
35	13.380	1916	1919	1922	rBV	17532	22878	1.17%	0.110%
36	13.410	1922	1924	1929	rVB2	16011	20545	1.05%	0.099%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082124\
Data File : BF139108.D
Acq On : 21 Aug 2024 18:20
Operator : RC/JU
Sample : P3663-09
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
915-J-SD-0-0.5-081624

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-BF082024.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	13.533	1942	1945	1948	rBV3	14758	21599	1.10%	0.104%
38	13.686	1968	1971	1978	rVB	41027	63409	3.23%	0.306%
39	13.804	1985	1991	1994	rBV3	95250	161250	8.22%	0.777%
40	13.833	1994	1996	1998	rVV	22639	26374	1.34%	0.127%
41	13.898	2003	2007	2014	rVV4	94144	188792	9.62%	0.910%
42	13.969	2015	2019	2023	rVV2	66273	97623	4.98%	0.470%
43	14.051	2023	2033	2036	rVV3	652415	1172364	59.77%	5.649%
44	14.080	2036	2038	2045	rVB	390543	454203	23.16%	2.189%
45	14.163	2048	2052	2055	rVV	55411	68682	3.50%	0.331%
46	14.274	2068	2071	2075	rVB2	38304	53961	2.75%	0.260%
47	14.439	2096	2099	2102	rVB	30429	35801	1.83%	0.173%
48	14.474	2102	2105	2108	rVB4	35994	38102	1.94%	0.184%
49	14.533	2112	2115	2118	rVV3	48571	69960	3.57%	0.337%
50	14.586	2121	2124	2128	rVB2	91898	121235	6.18%	0.584%
51	14.745	2148	2151	2155	rBV3	25593	39994	2.04%	0.193%
52	14.839	2163	2167	2171	rBV4	25430	47370	2.41%	0.228%
53	15.098	2205	2211	2222	rVB	321891	772219	39.37%	3.721%
54	15.198	2223	2228	2238	rVB4	81553	151028	7.70%	0.728%
55	15.298	2242	2245	2252	rBV3	17389	35583	1.81%	0.171%
56	15.404	2258	2263	2269	rVB	152649	221722	11.30%	1.068%
57	15.468	2269	2274	2279	rVB	189788	282159	14.38%	1.360%
58	15.533	2279	2285	2289	rBV	223971	371077	18.92%	1.788%
59	16.033	2365	2370	2375	rBV	40725	72902	3.72%	0.351%
60	17.004	2529	2535	2544	rVB2	83679	187155	9.54%	0.902%
61	17.251	2572	2577	2582	rBV	14026	30555	1.56%	0.147%
62	17.451	2605	2611	2623	rVB	63560	134278	6.85%	0.647%
63	18.227	2738	2743	2750	rVB9	13200	34744	1.77%	0.167%

Sum of corrected areas: 20752856

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082124\
Data File : BF139108.D
Acq On : 21 Aug 2024 18:20
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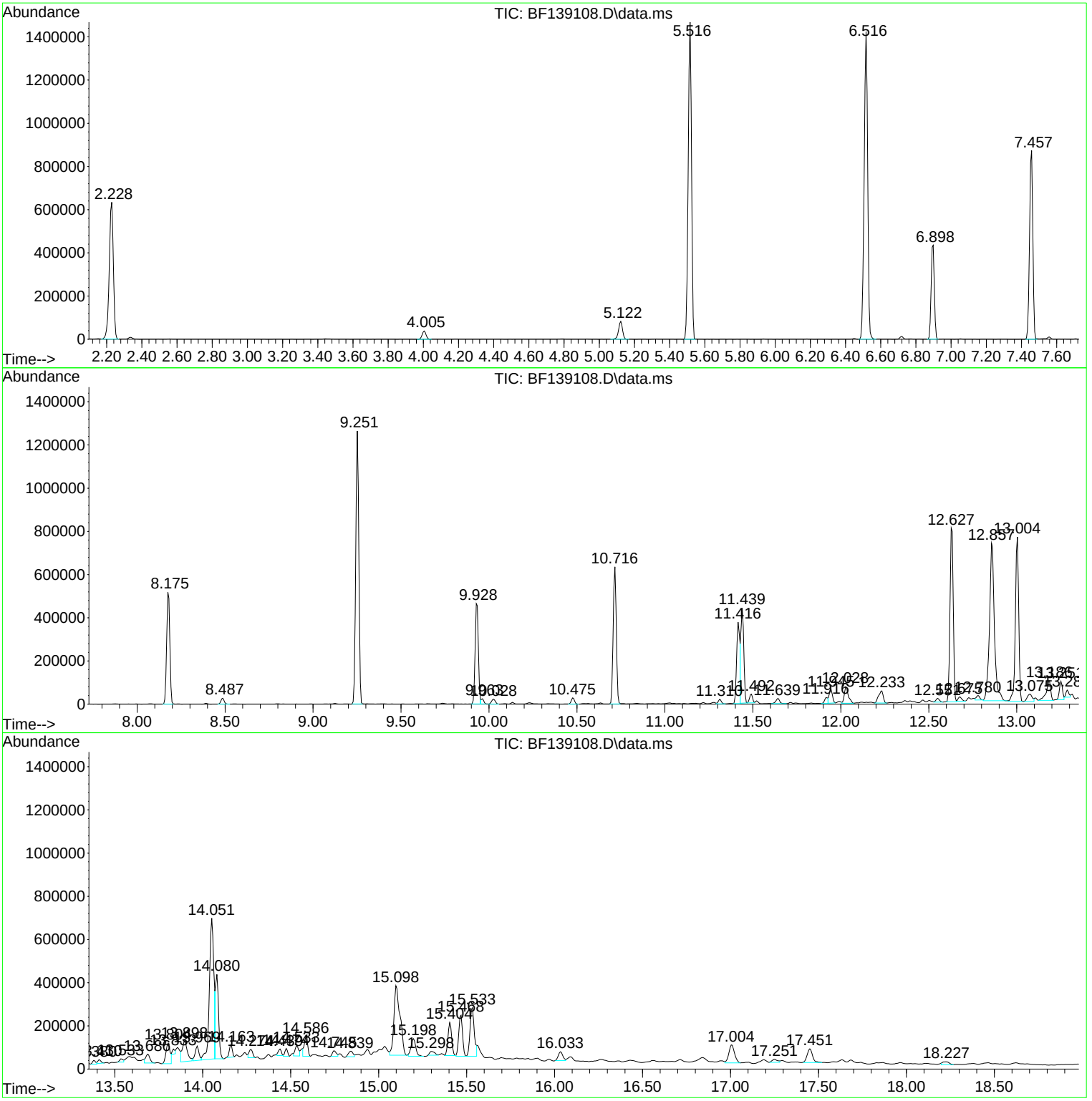
Sample : P3663-09
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
915-J-SD-0-0.5-081624

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF082024.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\BF082124\
Data File : BF139108.D
Acq On : 21 Aug 2024 18:20
Operator : RC/JU
Sample : P3663-09
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
915-J-SD-0-0.5-081624

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF082024.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

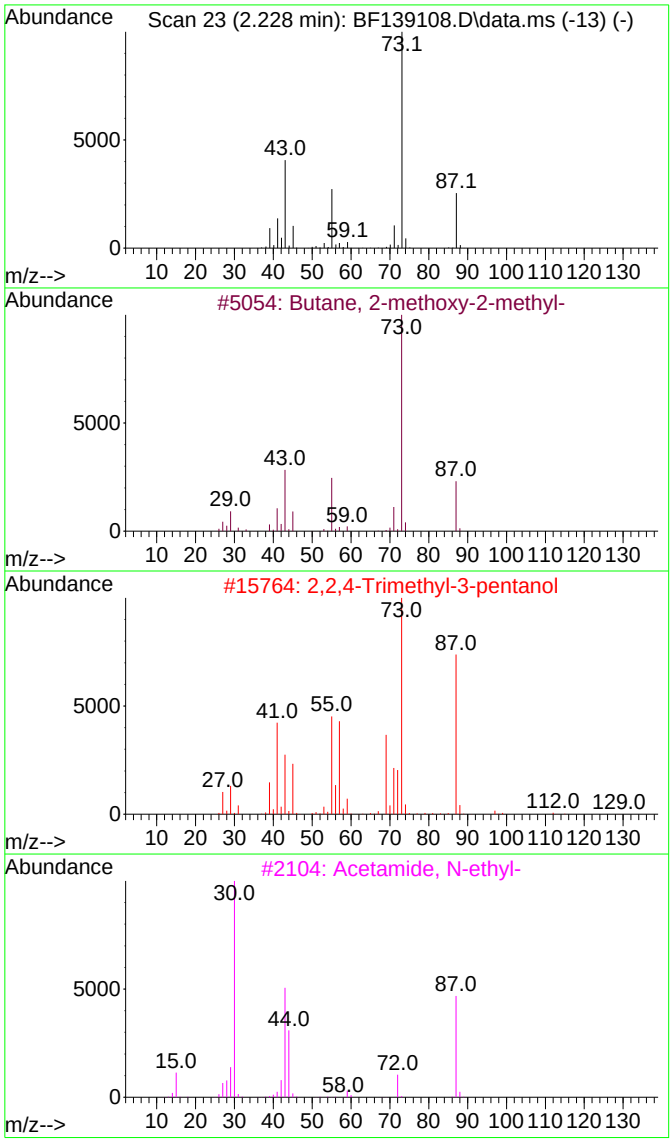
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

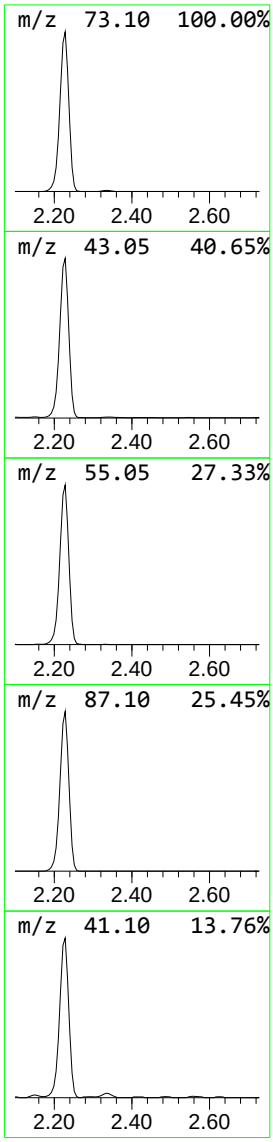
****

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.228	37.42 ng	1020250	1,4-Dichlorobenzene-d4	6.898

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3	Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	27
4	Octanal, 7-methoxy-3,7-dimethyl-	186	C11H22O2	003613-30-7	17
5	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082124\
Data File : BF139108.D
Acq On : 21 Aug 2024 18:20
Operator : RC/JU
Sample : P3663-09
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
915-J-SD-0-0.5-081624

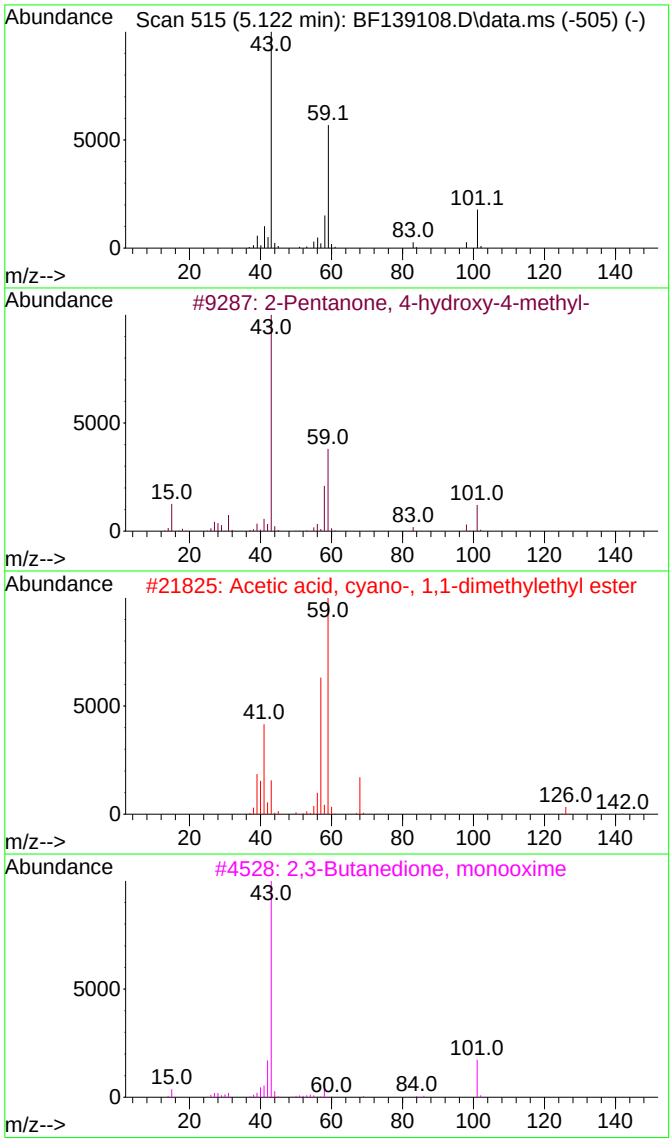
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF082024.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

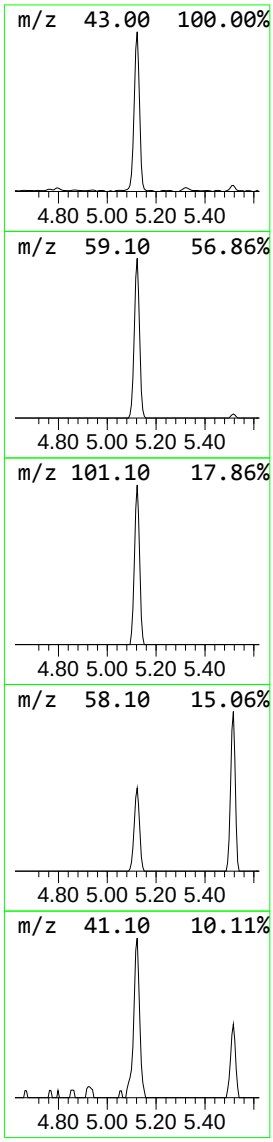
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.122	4.68 ng	127529	1,4-Dichlorobenzene-d4	6.898

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
3	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
4	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	12
5	Acetamide, N-(cyanomethyl)-	98	C4H6N2O	004814-80-6	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082124\
Data File : BF139108.D
Acq On : 21 Aug 2024 18:20
Operator : RC/JU
Sample : P3663-09
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
915-J-SD-0-0.5-081624

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF082024.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

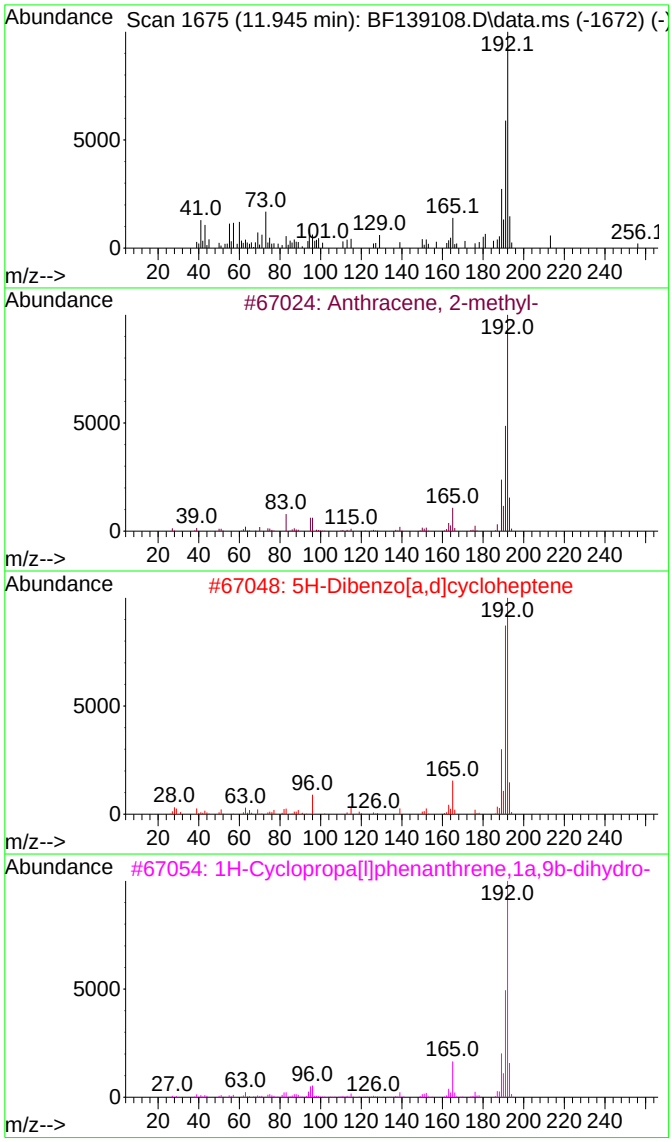
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

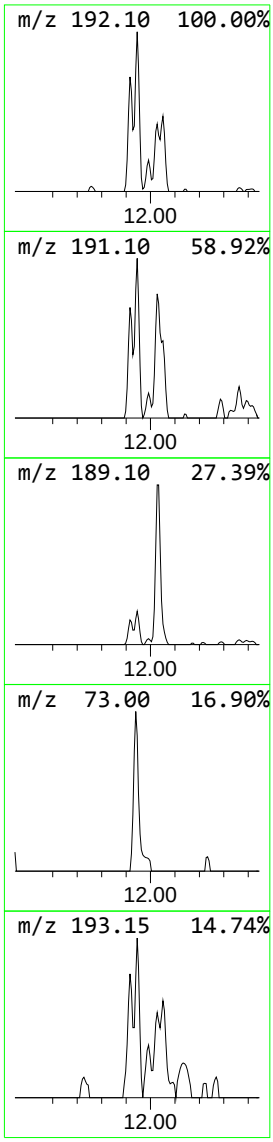
****

Peak Number 4 Anthracene, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.945	3.10 ng	77811	Phenanthrene-d10	11.416

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
2	5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	93
3	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	93
4	Anthracene, 1-methyl-	192	C15H12	000610-48-0	89
5	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	89





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082124\
Data File : BF139108.D
Acq On : 21 Aug 2024 18:20
Operator : RC/JU
Sample : P3663-09
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
915-J-SD-0-0.5-081624

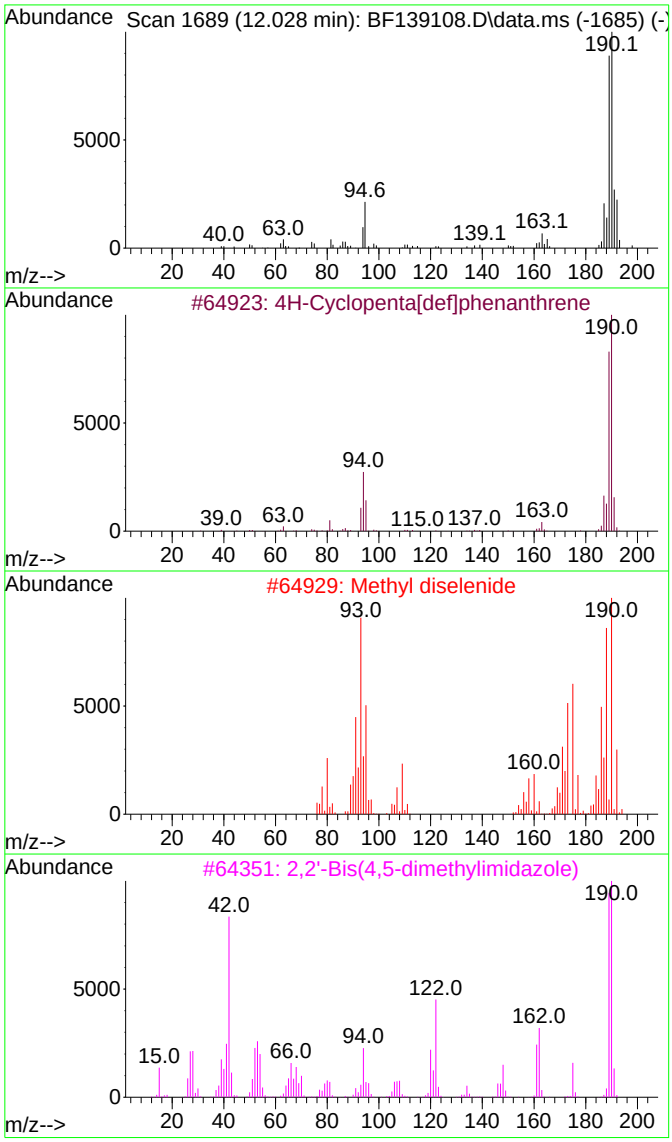
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF082024.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

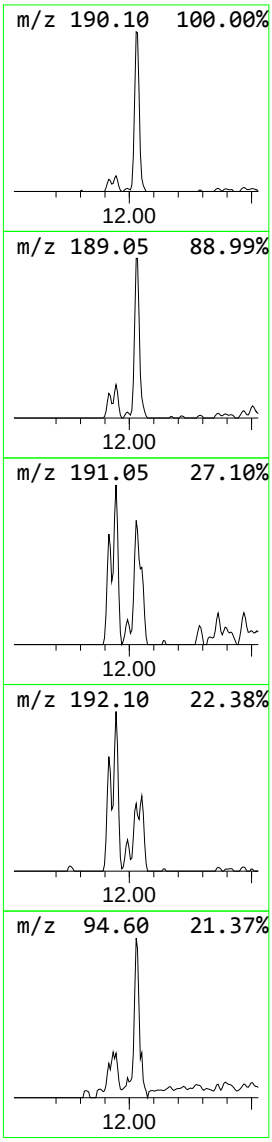
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.028	4.75 ng	119194	Phenanthrene-d10	11.416

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	93
2	Methyl diselenide	190	C2H6Se2	007101-31-7	46
3	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	45
4	6,7-Methylenedioxy-4(3H)-quinazo...	190	C9H6N2O3	028310-11-4	33
5	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	32





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082124\
Data File : BF139108.D
Acq On : 21 Aug 2024 18:20
Operator : RC/JU
Sample : P3663-09
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
915-J-SD-0-0.5-081624

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF082024.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

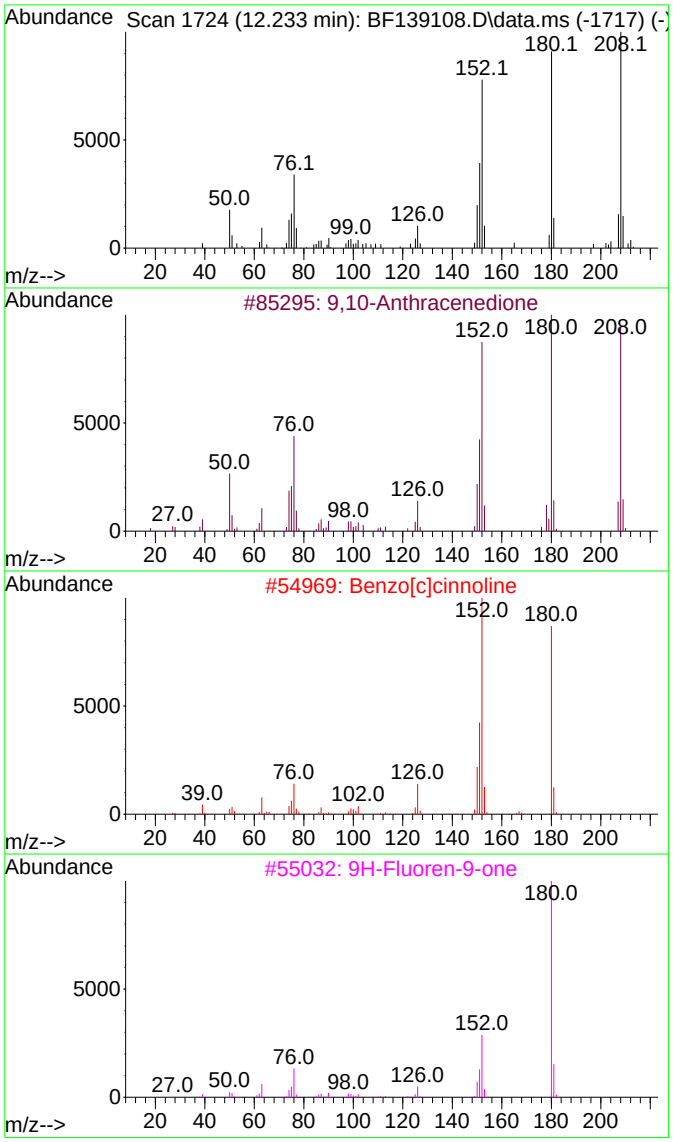
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

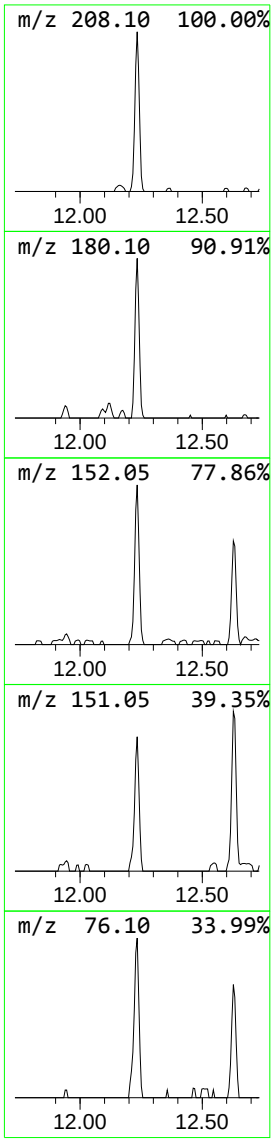
****

Peak Number 6 9,10-Anthracenedione Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.233	4.11 ng	103302	Phenanthrene-d10	11.416

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Anthracenedione	208	C14H8O2	000084-65-1	98
2	Benzo[c]cinnoline	180	C12H8N2	000230-17-1	93
3	9H-Fluoren-9-one	180	C13H8O	000486-25-9	58
4	9,10-Phenanthrenedione	208	C14H8O2	000084-11-7	58
5	1H-Phenalen-1-one	180	C13H8O	000548-39-0	53





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Data File : BF139108.D
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Sample : P3663-09
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
915-J-SD-0-0.5-081624

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF082024.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

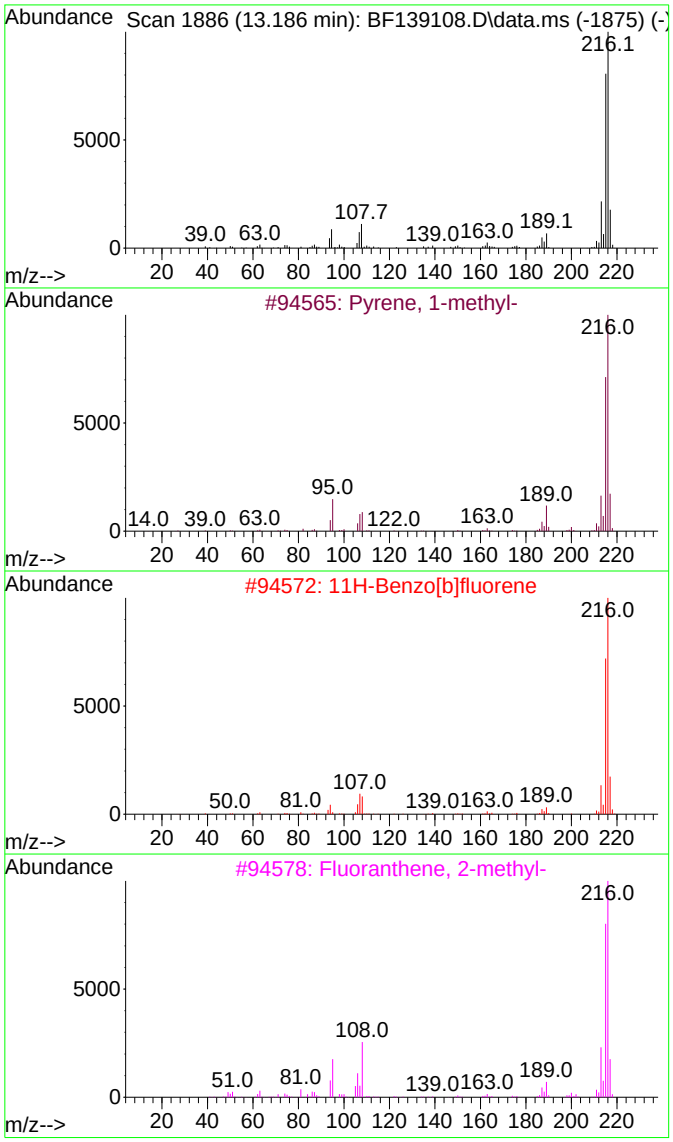
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

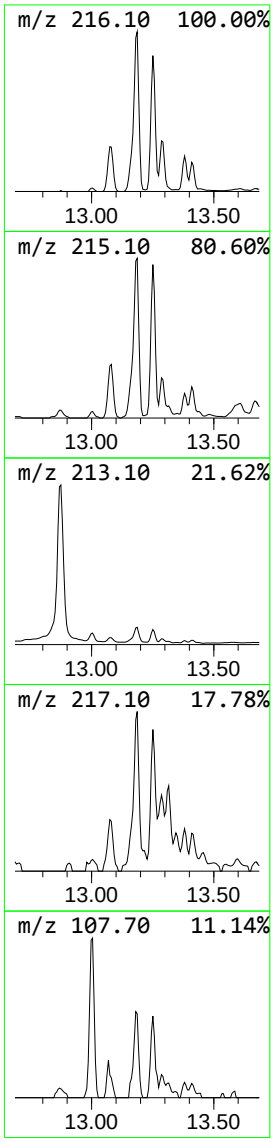
****

Peak Number 7 Pyrene, 1-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.186	2.68 ng	157386	Chrysene-d12	14.057

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-	216	C17H12	002381-21-7	94
2	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	94
3	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	91
4	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	90
5	7H-Benzo[c]fluorene	216	C17H12	000205-12-9	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082124\
Data File : BF139108.D
Acq On : 21 Aug 2024 18:20
Operator : RC/JU
Sample : P3663-09
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
915-J-SD-0-0.5-081624

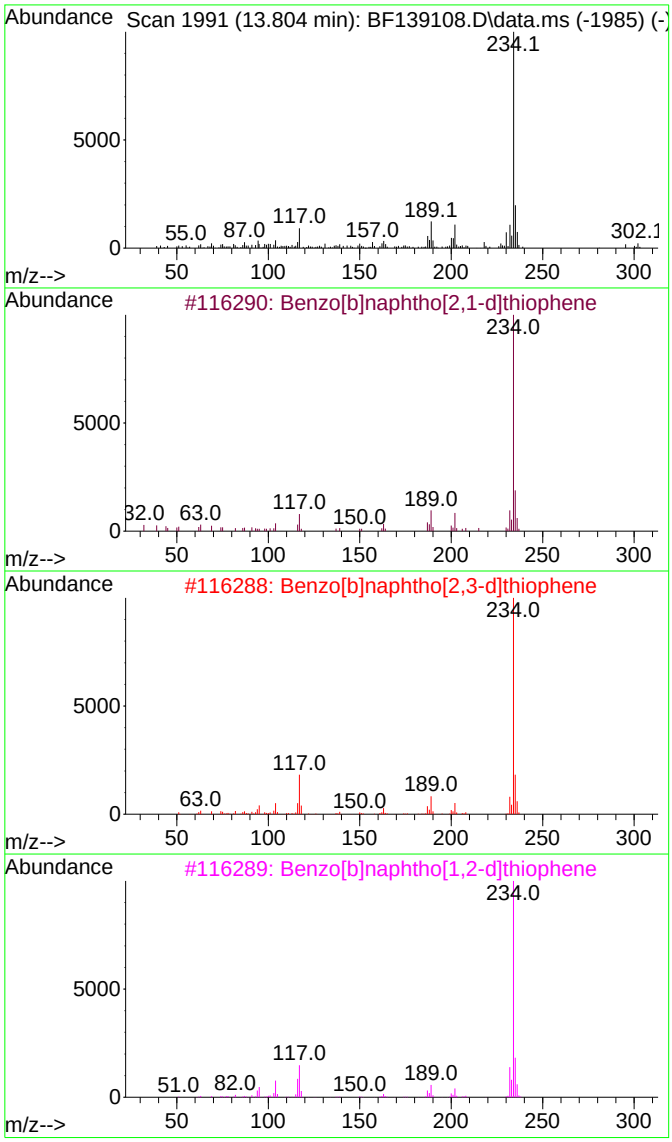
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF082024.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

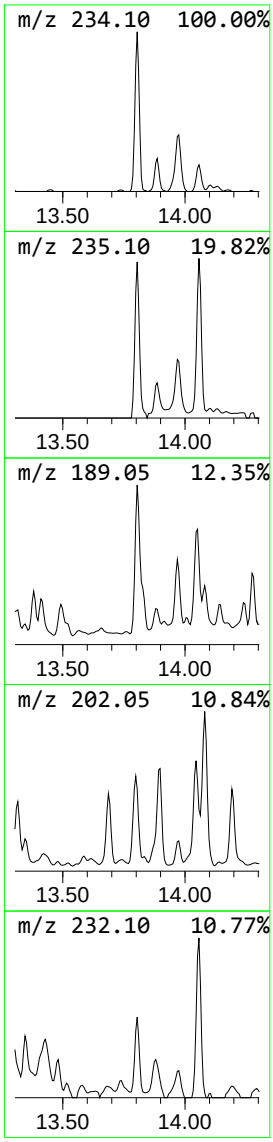
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzo[b]naphtho[2,1-d]thiop... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.804	2.75 ng	161250	Chrysene-d12	14.057

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[b]naphtho[2,1-d]thiophene	234	C16H10S	000239-35-0	96
2	Benzo[b]naphtho[2,3-d]thiophene	234	C16H10S	000243-46-9	94
3	Benzo[b]naphtho[1,2-d]thiophene	234	C16H10S	000205-43-6	83
4	Phenanthro(2,1-b)thiophene	234	C16H10S	000219-25-0	83
5	pyrido[1',2':1,2]imidazo[4,5-b]q...	234	C14H10N4	1000401-06-3	64





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Acq On : 21 Aug 2024 18:20
Operator : RC/JU
Sample : P3663-09
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
915-J-SD-0-0.5-081624

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF082024.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

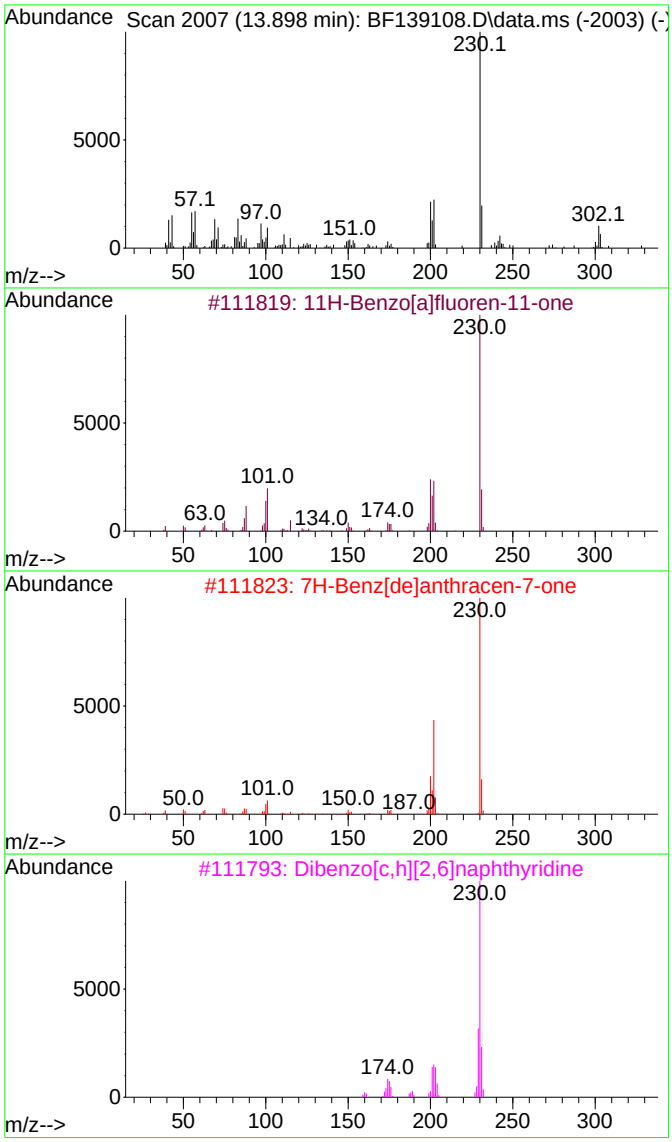
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

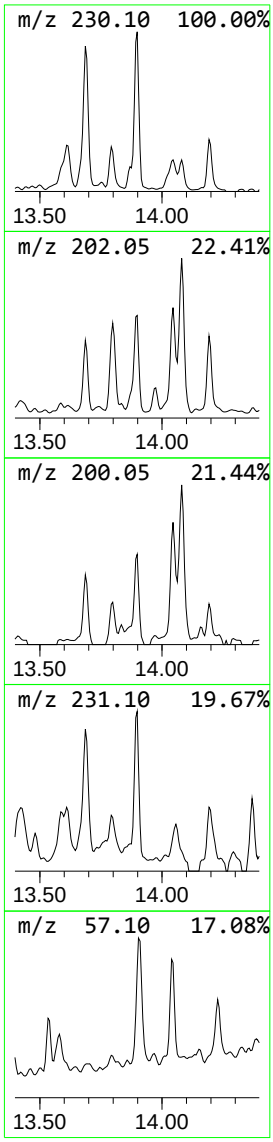
****

Peak Number 9 11H-Benzo[a]fluoren-11-one Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.898	3.22 ng	188792	Chrysene-d12	14.057

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[a]fluoren-11-one	230	C17H10O	000479-79-8	91
2	7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	89
3	Dibenzo[c,h][2,6]naphthyridine	230	C16H10N2	000218-30-4	53
4	2,6-Dimethyl-4-hydroxyquinoline,...	287	C17H25NOSi	1000463-05-5	50
5	5-Amino-7-imino-4,8-diazatricycl...	230	C12H14N4O	1000388-35-6	49





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Instrument :
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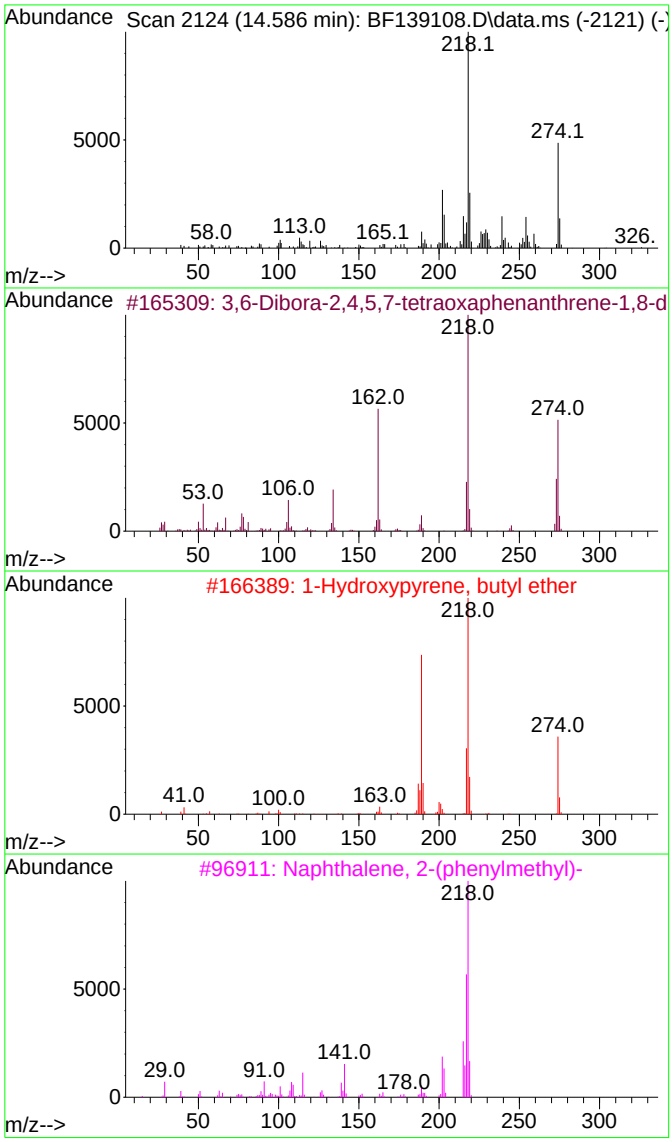
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

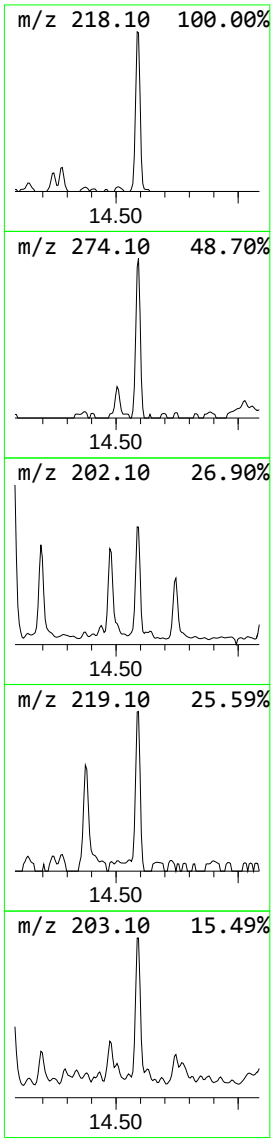
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 3,6-Dibora-2,4,5,7-tetraoxa... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.586	2.07 ng	121235	Chrysene-d12	14.057

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,6-Dibora-2,4,5,7-tetraoxaphena...	274	C12H12B2O6	1000159-66-2	50
2	1-Hydroxypyrene, butyl ether	274	C20H18O	1000453-43-2	42
3	Naphthalene, 2-(phenylmethyl)-	218	C17H14	000613-59-2	38
4	3,4-Dihydro-2-bromo-4,4-dimethyl...	312	C18H17Br	071161-44-9	38
5	1H-Pyrazole-5-carboxylic acid, 3...	274	C15H18N2O3	1000319-72-5	35





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Sample : P3663-09
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915-J-SD-0-0.5-081624

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF082024.M
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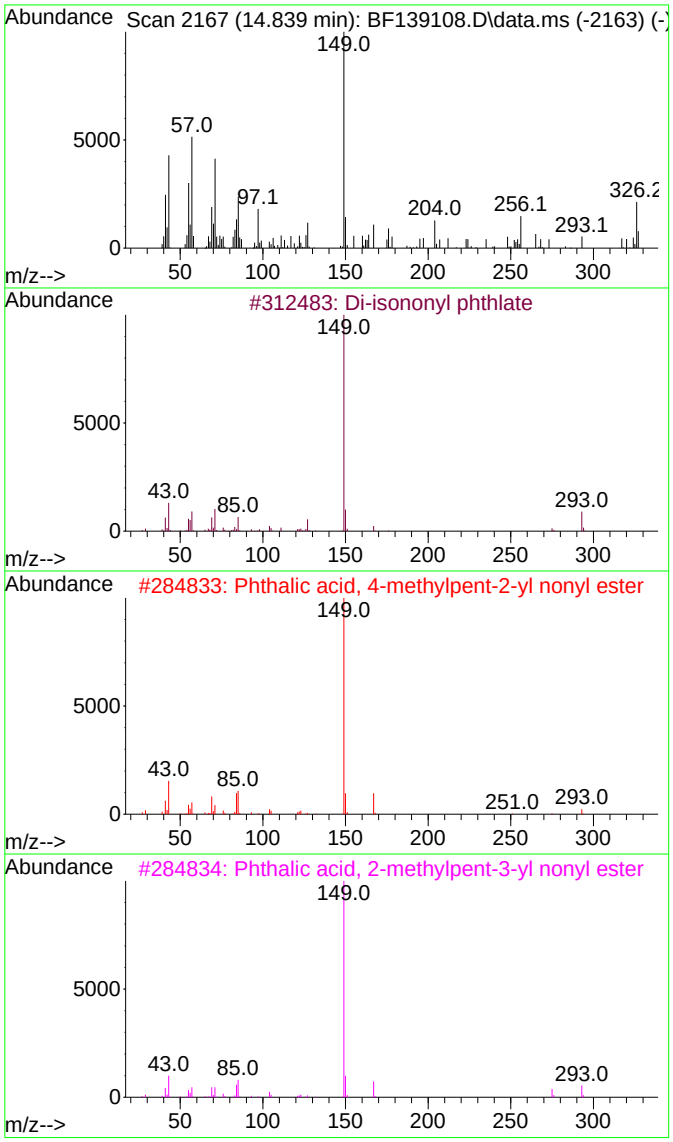
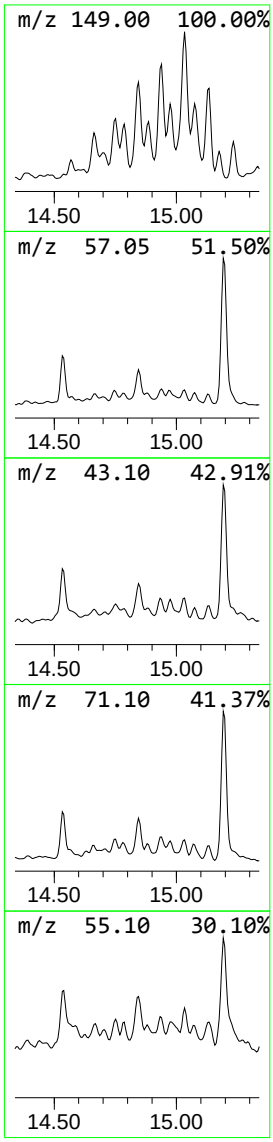
Peak Number 11 Di-isononyl phthlate Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.839	2.55 ng	47370	Perylene-d12	15.533

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Di-isononyl phthlate	418	C26H42O4	020548-62-3	59
2	Phthalic acid, 4-methylpent-2-yl...	376	C23H36O4	1000356-87-9	38
3	Phthalic acid, 2-methylpent-3-yl...	376	C23H36O4	1000356-91-4	38
4	1,2-Benzenedicarboxylic acid, di...	418	C26H42O4	000084-76-4	38
5	1,2-Benzenedicarboxylic acid, di...	474	C30H50O4	003648-20-2	38





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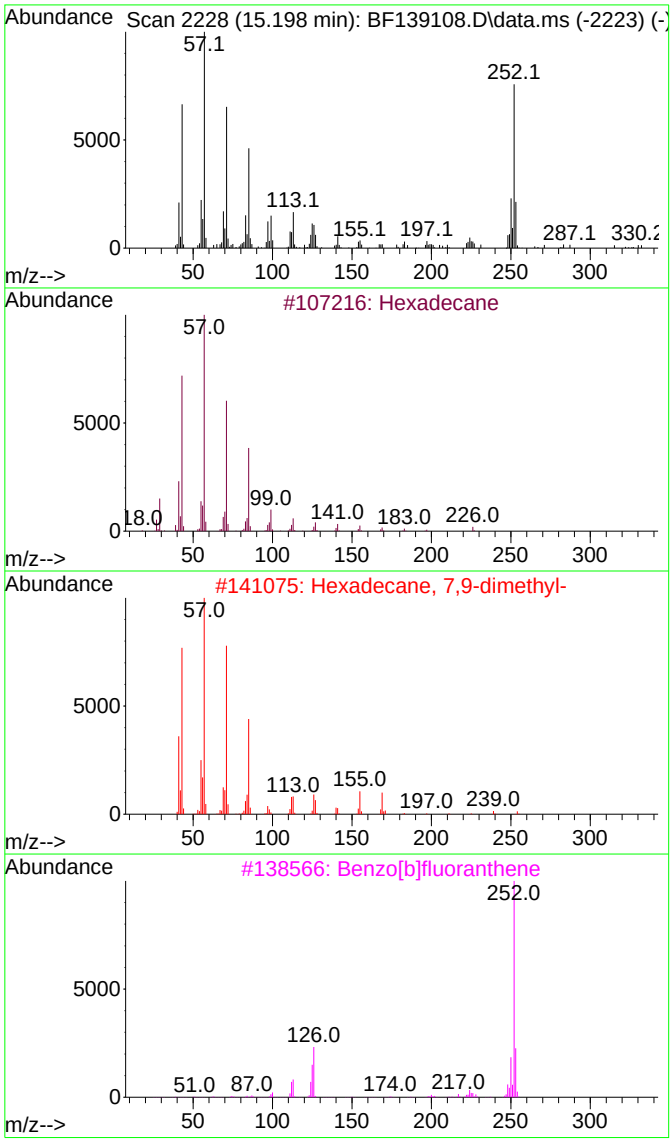
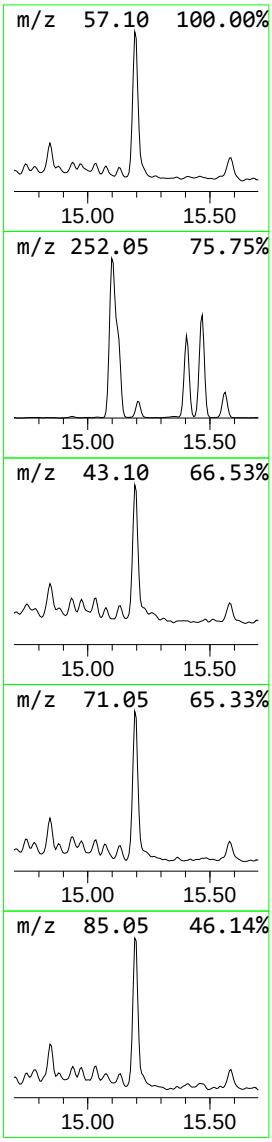
Peak Number 12 Hexadecane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.198	8.14 ng	151028	Perylene-d12	15.533

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	70
2	Hexadecane, 7,9-dimethyl-	254	C18H38	021164-95-4	70
3	Benzo[b]fluoranthene	252	C20H12	000205-99-2	64
4	Benzo[a]pyrene	252	C20H12	000050-32-8	50
5	Perylene	252	C20H12	000198-55-0	45





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Instrument :
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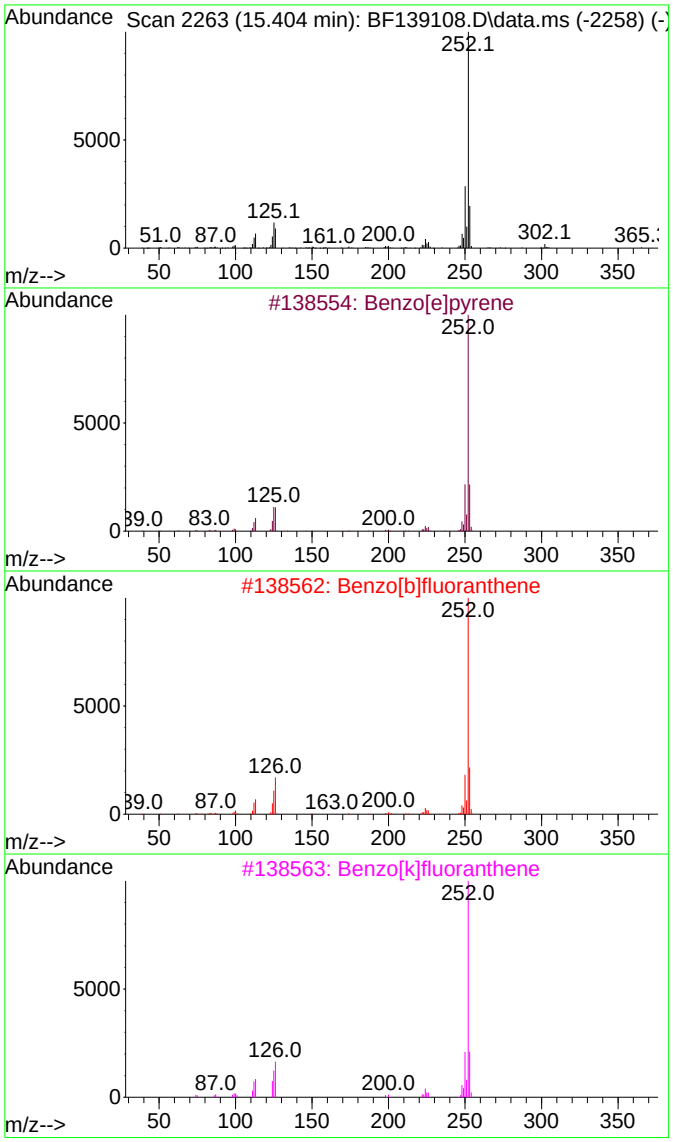
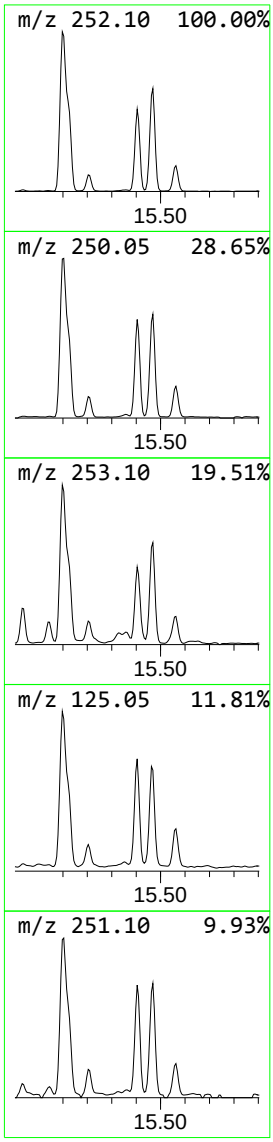
Peak Number 13 Benzo[e]pyrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.404	11.95 ng	221722	Perylene-d12	15.533

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[e]pyrene	252	C20H12	000192-97-2	94
2	Benzo[b]fluoranthene	252	C20H12	000205-99-2	93
3	Benzo[k]fluoranthene	252	C20H12	000207-08-9	93
4	Perylene	252	C20H12	000198-55-0	90
5	Benzo[a]pyrene	252	C20H12	000050-32-8	90





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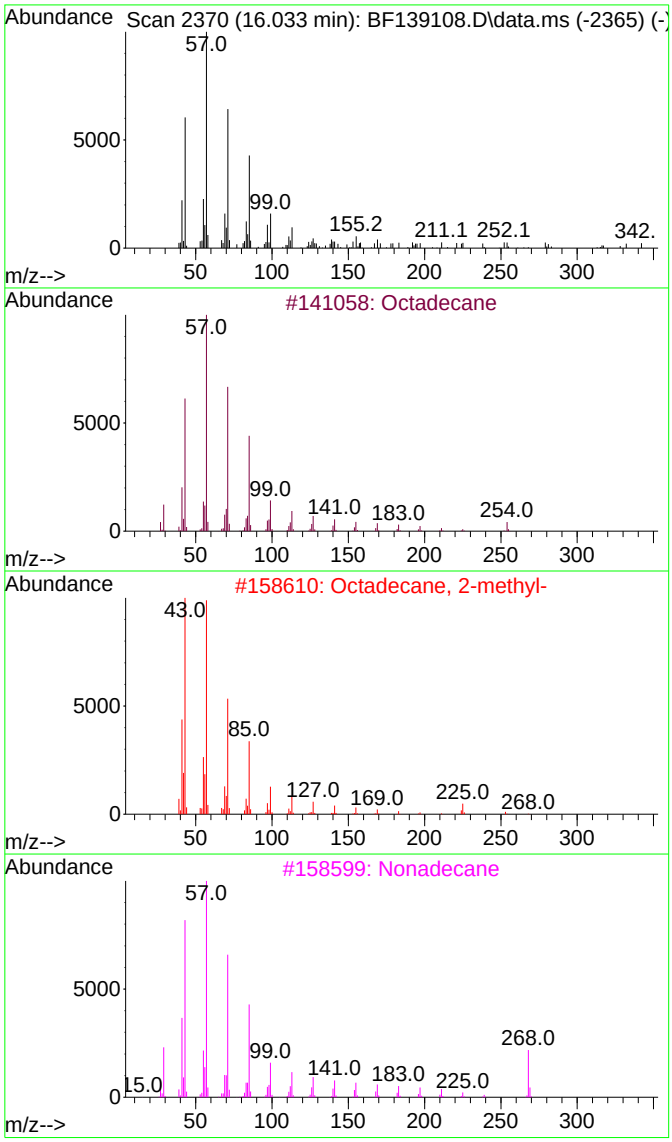
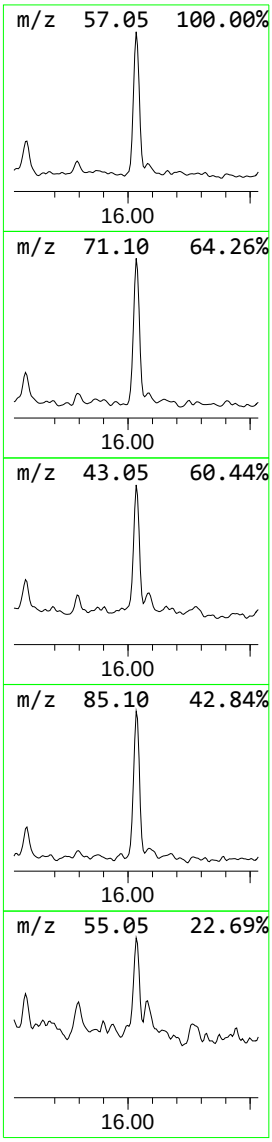
Peak Number 14 Octadecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.033	3.93 ng	72902	Perylene-d12	15.533

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane	254	C18H38	000593-45-3	93
2	Octadecane, 2-methyl-	268	C19H40	001560-88-9	93
3	Nonadecane	268	C19H40	000629-92-5	91
4	Tetracosane	338	C24H50	000646-31-1	87
5	Heptadecane	240	C17H36	000629-78-7	87





Tentatively Identified Compound (LSC) summary

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF082024.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
Butane, 2-metho...	2.228	37.4	ng	1020250	1	6.898	545317	20.0
2-Pentanone, 4-...	5.122	4.7	ng	127529	1	6.898	545317	20.0
Anthracene, 2-m...	11.945	3.1	ng	77811	4	11.416	502316	20.0
4H-Cyclopenta[d...	12.028	4.8	ng	119194	4	11.416	502316	20.0
9,10-Anthracene...	12.233	4.1	ng	103302	4	11.416	502316	20.0
Pyrene, 1-methyl-	13.186	2.7	ng	157386	5	14.057	1172360	20.0
Benzo[b]naphtho...	13.804	2.8	ng	161250	5	14.057	1172360	20.0
11H-Benzo[a]flu...	13.898	3.2	ng	188792	5	14.057	1172360	20.0
3,6-Dibora-2,4,...	14.586	2.1	ng	121235	5	14.057	1172360	20.0
Di-isononyl pht...	14.839	2.5	ng	47370	6	15.533	371077	20.0
Hexadecane	15.198	8.1	ng	151028	6	15.533	371077	20.0
Benzo[e]pyrene	15.404	11.9	ng	221722	6	15.533	371077	20.0
Octadecane	16.033	3.9	ng	72902	6	15.533	371077	20.0