LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072522\
Data File : BG054387.D
Acq On : 25 Jul 2022 20:08
Operator : CG/JU
Sample : N3855-01 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EO-01-072222

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_G\Methods\8270-BG071522.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG054387.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.038	3	8	17	rVV3	6840	13916	1.65%	0.210%
2	3.120	17	22	30	rVB2	10379	19503	2.31%	0.294%
3	5.570	433	439	447	rBV	26075	46585	5.52%	0.702%
4	7.174	706	712	721	rVB	28256	51637	6.12%	0.779%
5	7.991	843	851	858	rBB	82383	144541	17.13%	2.179%
6	9.154	1043	1049	1057	rBV	18160	32701	3.87%	0.493%
7	10.799	1320	1329	1335	rBV	110082	203089	24.06%	3.062%
8	12.456	1605	1611	1616	rVB2	6175	10986	1.30%	0.166%
9	12.673	1641	1648	1657	rVB2	6245	12494	1.48%	0.188%
10	13.161	1728	1731	1736	rVB2	5621	8729	1.03%	0.132%
11	13.243	1740	1745	1753	rVV	62411	102265	12.12%	1.542%
12	13.449	1774	1780	1786	rVB6	4664	9609	1.14%	0.145%
13	13.948	1860	1865	1870	rBV4	7315	11893	1.41%	0.179%
14	14.101	1885	1891	1895	rBV6	4851	8779	1.04%	0.132%
15	14.348	1929	1933	1942	rBV2	11008	20661	2.45%	0.312%
16	14.454	1945	1951	1957	rVB6	3760	8683	1.03%	0.131%
17	14.512	1957	1961	1967	rBV7	6327	10908	1.29%	0.164%
18	14.624	1971	1980	1987	rVV	189792	307950	36.49%	4.643%
19	14.689	1987	1991	1997	rVB3	13609	22795	2.70%	0.344%
20	15.024	2042	2048	2053	rVB3	15445	26198	3.10%	0.395%
21	15.100	2053	2061	2065	rVV3	5866	9585	1.14%	0.145%
22	15.241	2080	2085	2090	rBV5	5975	11996	1.42%	0.181%
23	15.406	2108	2113	2117	rBV6	4555	9741	1.15%	0.147%
24	15.464	2117	2123	2127	rVB3	6785	11861	1.41%	0.179%
25	15.670	2151	2158	2164	rBV	32445	52649	6.24%	0.794%
26	15.876	2190	2193	2198	rVB5	8238	12551	1.49%	0.189%
27	15.964	2203	2208	2215	rVB5	9155	16438	1.95%	0.248%
28	16.116	2224	2234	2242	rBV	62959	107601	12.75%	1.622%
29	16.351	2263	2274	2279	rBV6	16846	42355	5.02%	0.639%
30	16.675	2322	2329	2333	rBV6	8953	18008	2.13%	0.272%
31	16.727	2334	2338	2343	rVB4	10292	14104	1.67%	0.213%
32	16.822	2349	2354	2360	rVB7	6571	12926	1.53%	0.195%
33	16.904	2362	2368	2370	rBV6	6782	10787	1.28%	0.163%
34	17.027	2384	2389	2395	rBV4	7610	12461	1.48%	0.188%
35	17.115	2398	2404	2409	rVV6	15064	27503	3.26%	0.415%
36	17.186	2410	2416	2424	rVB2	17369	44750	5.30%	0.675%

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072522\
Data File : BG054387.D
Acq On : 25 Jul 2022 20:08
Operator : CG/JU
Sample : N3855-01 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EO-01-072222

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_G\Methods\8270-BG071522.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	17.374	2441	2448	2451	rBV	230506	376414	44.60%	5.676%
38	17.415	2451	2455	2461	rVB	220353	325669	38.59%	4.910%
39	17.503	2465	2470	2475	rVB	54396	79100	9.37%	1.193%
40	17.773	2512	2516	2525	rVB	18972	31351	3.71%	0.473%
41	17.856	2527	2530	2532	rVV3	9225	8837	1.05%	0.133%
42	17.950	2543	2546	2554	rVV2	10911	18264	2.16%	0.275%
43	18.026	2555	2559	2565	rVB	36456	52562	6.23%	0.793%
44	18.249	2592	2597	2601	rBV	44065	68452	8.11%	1.032%
45	18.296	2601	2605	2611	rVB	50744	80900	9.59%	1.220%
46	18.379	2613	2619	2625	rBV	20445	40486	4.80%	0.610%
47	18.443	2625	2630	2634	rBV3	66622	127783	15.14%	1.927%
48	18.549	2645	2648	2652	rBV5	12434	15483	1.83%	0.233%
49	18.743	2677	2681	2685	rBV	30513	39012	4.62%	0.588%
50	18.984	2720	2722	2726	rBV4	8557	11834	1.40%	0.178%
51	19.042	2729	2732	2735	rBV2	12190	15396	1.82%	0.232%
52	19.166	2748	2753	2755	rBV3	143016	200284	23.73%	3.020%
53	19.195	2755	2758	2771	rVB4	123804	220095	26.08%	3.319%
54	19.424	2792	2797	2804	rVB	460170	657467	77.90%	9.913%
55	19.477	2804	2806	2809	rBV3	14621	17467	2.07%	0.263%
56	19.759	2849	2854	2855	rBV2	74760	112532	13.33%	1.697%
57	19.789	2855	2859	2866	rVB	380244	557669	66.07%	8.408%
58	19.977	2887	2891	2896	rVB2	136848	183509	21.74%	2.767%
59	20.276	2940	2942	2947	rVB	80375	99002	11.73%	1.493%
60	20.376	2956	2959	2963	rVB	51841	62704	7.43%	0.945%
61	21.628	3166	3172	3178	rBV2	337977	843994	100.00%	12.726%
62	21.681	3178	3181	3193	rVB	196253	346087	41.01%	5.218%
63	23.831	3542	3547	3554	rBV	94232	248411	29.43%	3.746%
64	24.842	3714	3719	3729	rVB	117834	302229	35.81%	4.557%

Sum of corrected areas: 6632231

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072522\
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Operator : CG/JU
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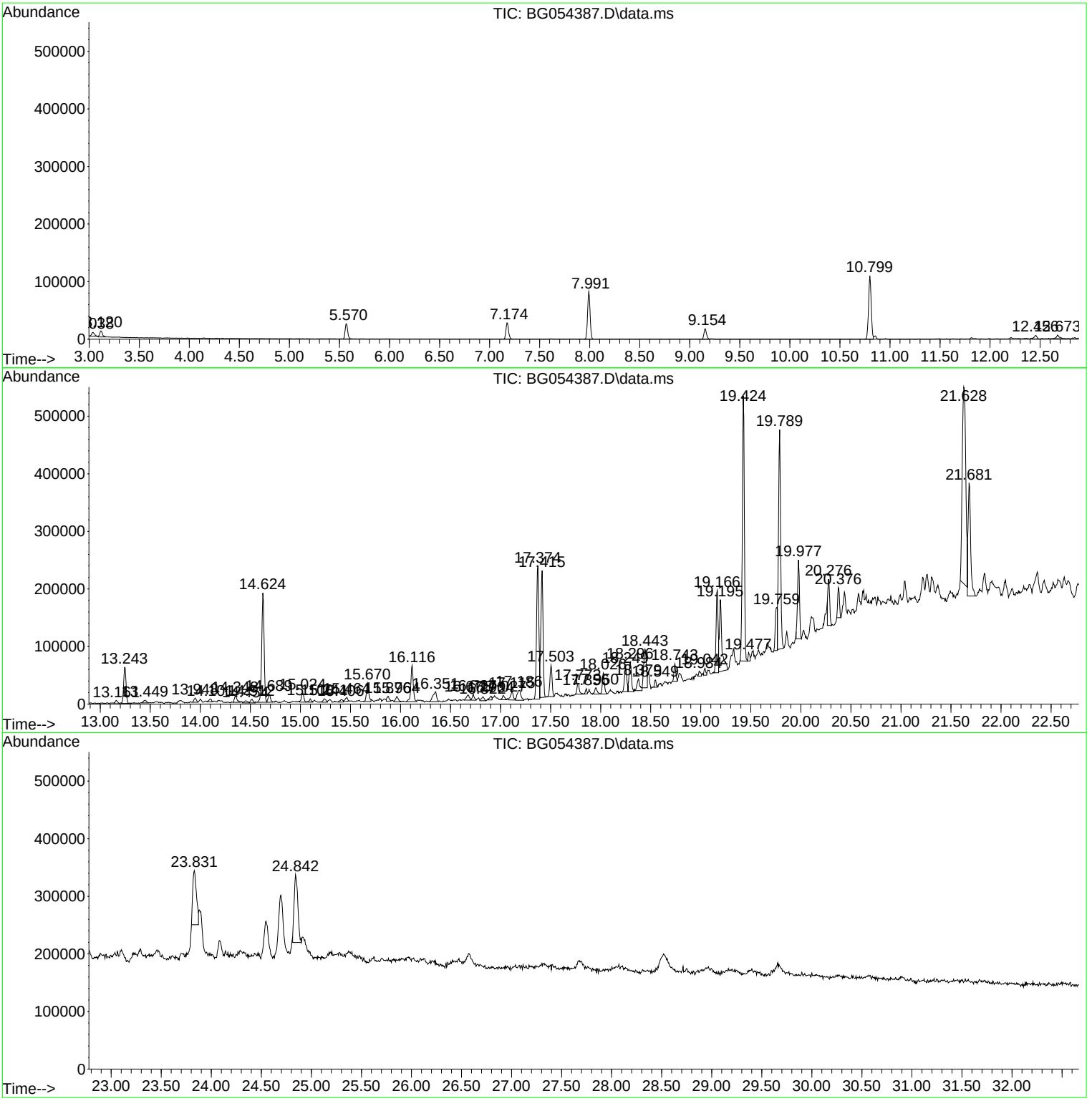
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EO-01-072222

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG071522.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_G\Data\BG072522\
Data File : BG054387.D
Acq On : 25 Jul 2022 20:08
Operator : CG/JU
Sample : N3855-01 5X
Misc :
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Instrument :
BNA_G
ClientSampleId :
EO-01-072222

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG071522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

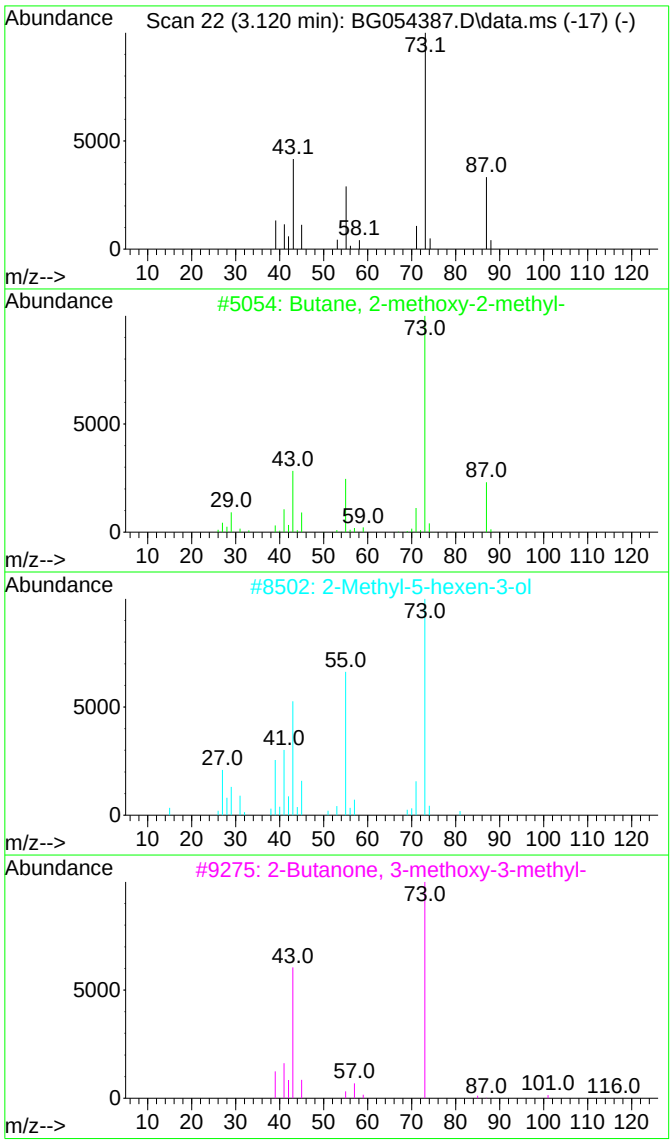
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

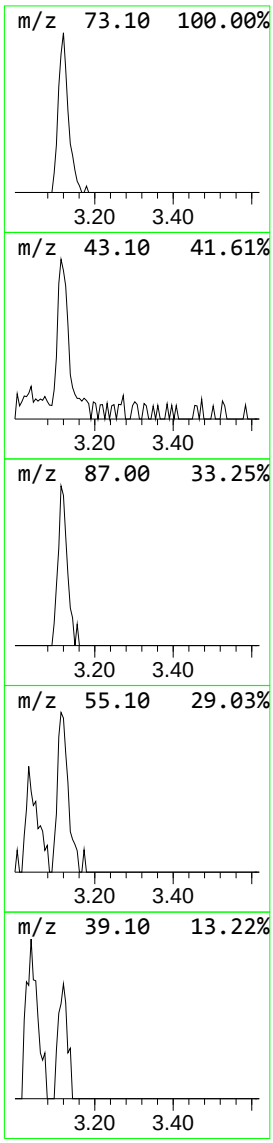
****

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.120	2.70 ng	19503	1,4-Dichlorobenzene-d4	7.991

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	56
2	2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	38
3	2-Butanone, 3-methoxy-3-methyl-	116	C6H12O2	036687-98-6	9
4	Silane, ethyltrimethyl-	102	C5H14Si	003439-38-1	9
5	3-Hexanol, 2,4-dimethyl-	130	C8H18O	013432-25-2	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072522\
Data File : BG054387.D
Acq On : 25 Jul 2022 20:08
Operator : CG/JU
Sample : N3855-01 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EO-01-072222

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG071522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

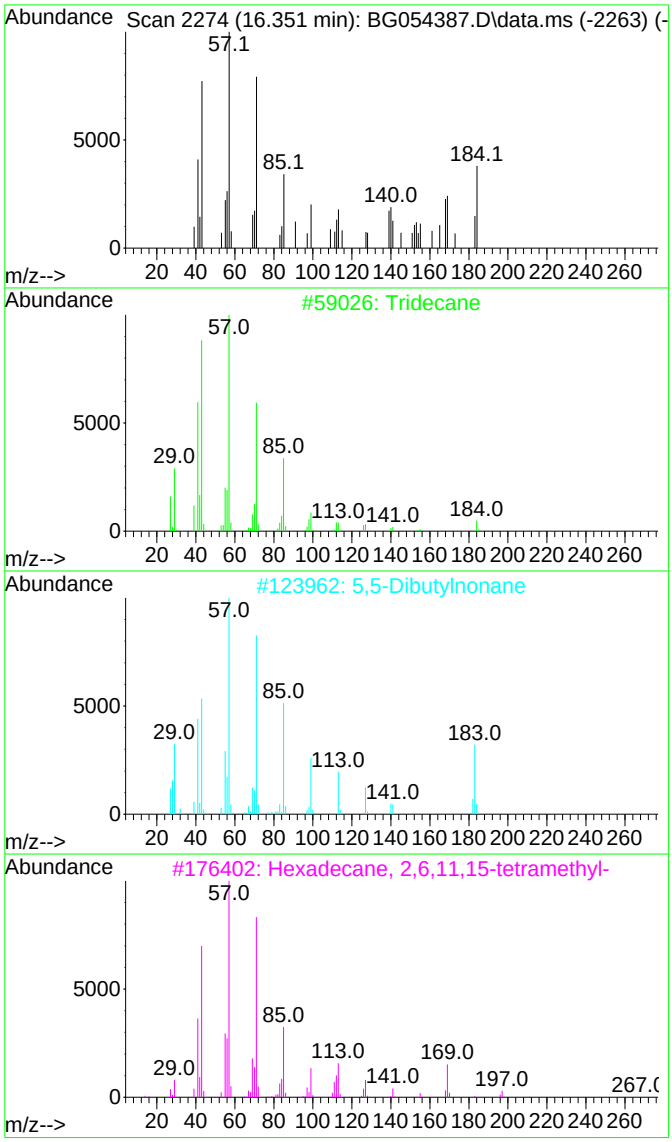
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

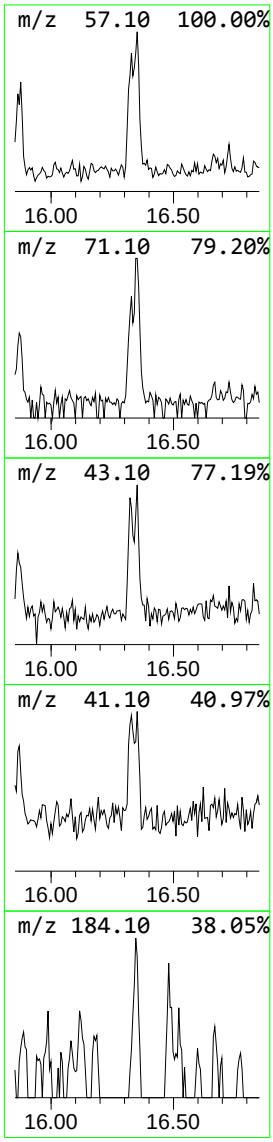
****

Peak Number 2 unknown16.352 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.352	2.25 ng	42355	Phenanthrene-d10	17.374

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane	184	C13H28	000629-50-5	49
2	5,5-Dibutylnonane	240	C17H36	006008-17-9	43
3	Hexadecane, 2,6,11,15-tetramethyl-	282	C20H42	000504-44-9	43
4	Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	38
5	Tridecane, 1-iodo-	310	C13H27I	035599-77-0	38





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072522\
Data File : BG054387.D
Acq On : 25 Jul 2022 20:08
Operator : CG/JU
Sample : N3855-01 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EO-01-072222

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG071522.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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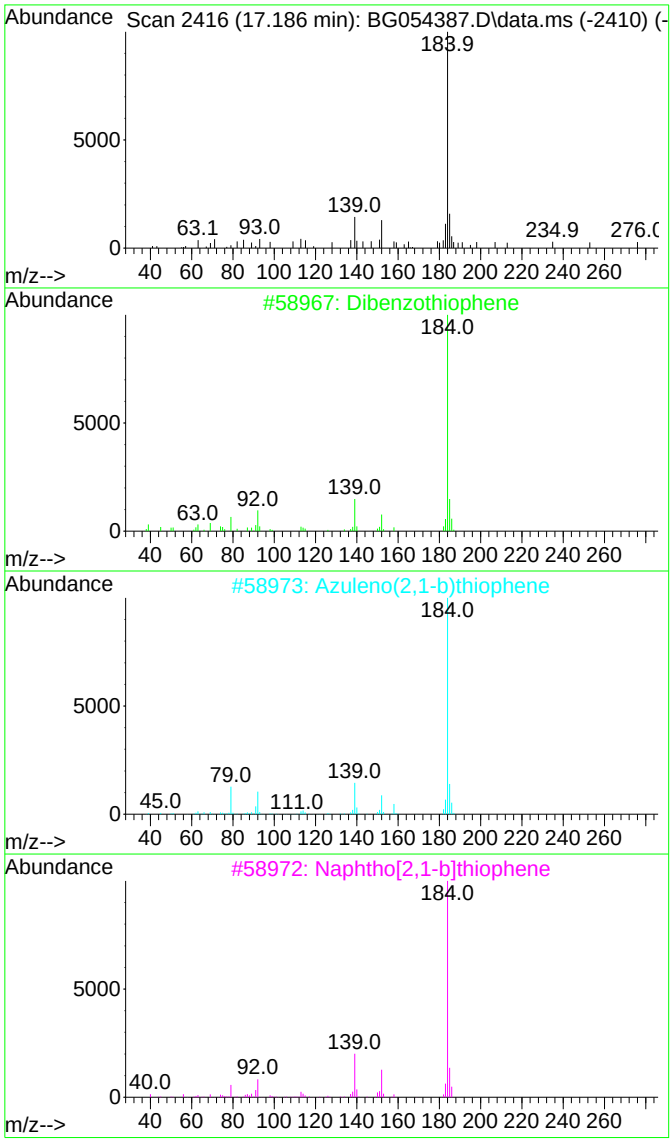
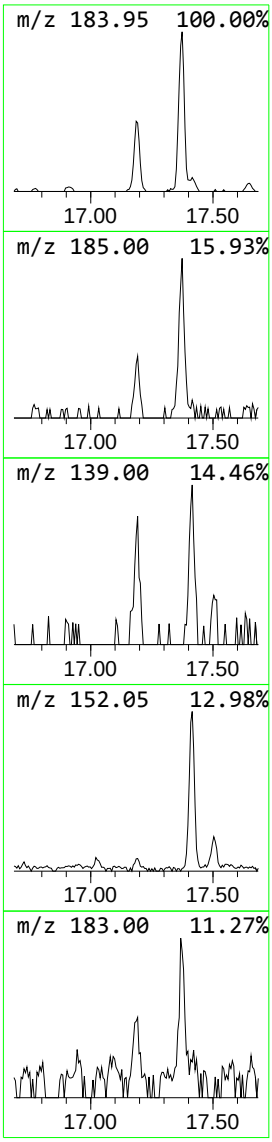
Peak Number 3 Dibenzothiophene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.186	2.38 ng	44750	Phenanthrene-d10	17.374

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dibenzothiophene	184	C12H8S	000132-65-0	91
2	Azuleno(2,1-b)thiophene	184	C12H8S	000248-13-5	91
3	Naphtho[2,1-b]thiophene	184	C12H8S	000233-02-3	80
4	Naphtho[2,3-b]thiophene	184	C12H8S	000268-77-9	74
5	Naphtho[1,2-b]thiophene	184	C12H8S	000234-41-3	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072522\
Data File : BG054387.D
Acq On : 25 Jul 2022 20:08
Operator : CG/JU
Sample : N3855-01 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EO-01-072222

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG071522.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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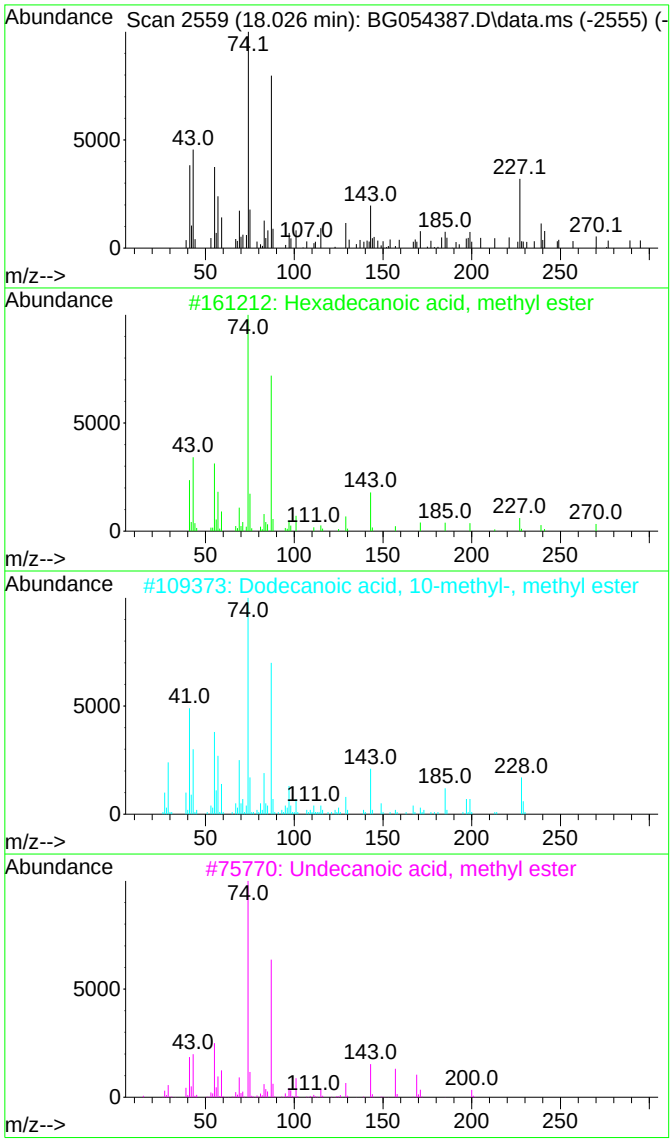
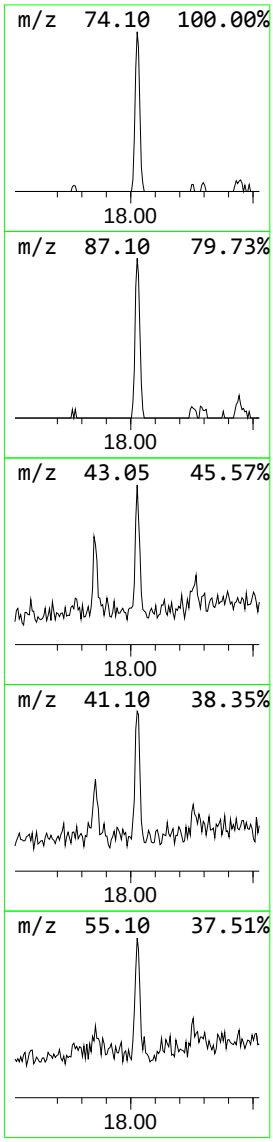
Peak Number 4 Hexadecanoic acid, methyl e... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.026	2.79 ng	52562	Phenanthrene-d10	17.374

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	95
2	Dodecanoic acid, 10-methyl-, met...	228	C14H28O2	005129-65-7	76
3	Undecanoic acid, methyl ester	200	C12H24O2	001731-86-8	76
4	Hexadecanoic acid, 2-methyl-	270	C17H34O2	027147-71-3	76
5	Undecanoic acid, 10-methyl-, met...	214	C13H26O2	005129-56-6	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072522\
Data File : BG054387.D
Acq On : 25 Jul 2022 20:08
Operator : CG/JU
Sample : N3855-01 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EO-01-072222

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG071522.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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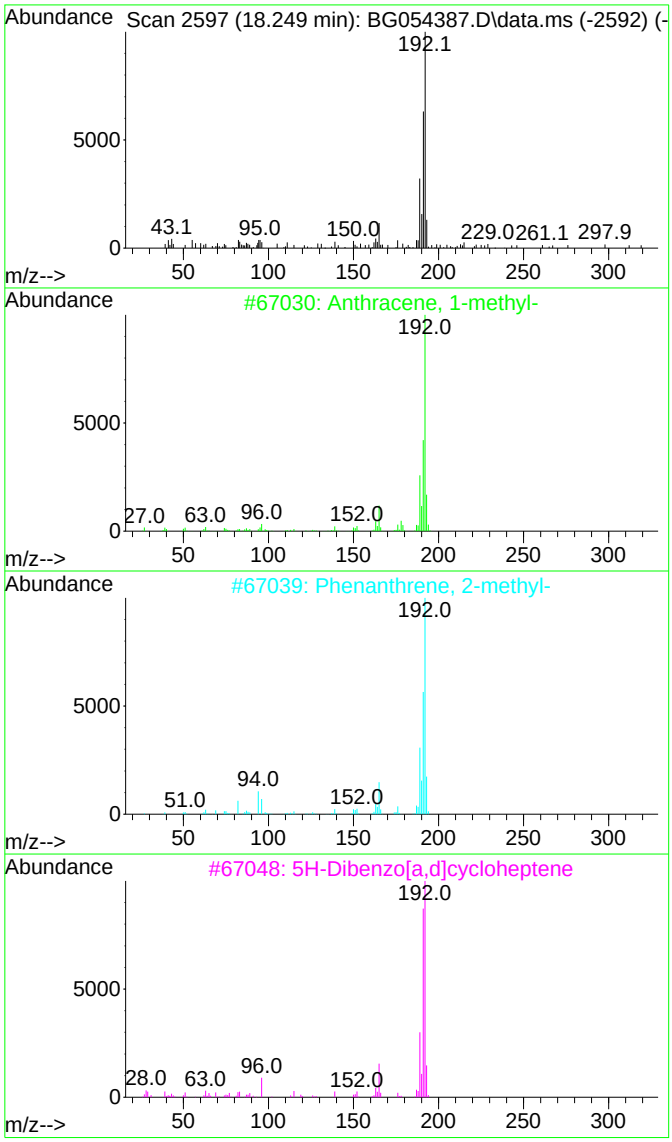
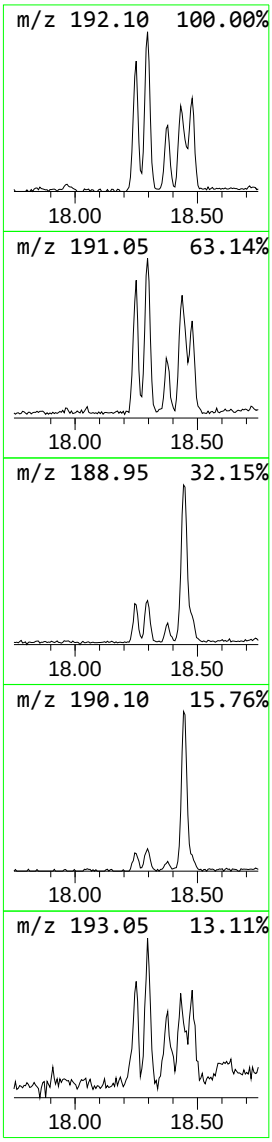
Peak Number 5 Anthracene, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.249	3.64 ng	68452	Phenanthrene-d10	17.374

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 1-methyl-	192	C15H12	000610-48-0	92
2	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	91
3	5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	91
4	Anthracene, 2-methyl-	192	C15H12	000613-12-7	90
5	Anthracene, 9-methyl-	192	C15H12	000779-02-2	87





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072522\
Data File : BG054387.D
Acq On : 25 Jul 2022 20:08
Operator : CG/JU
Sample : N3855-01 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG071522.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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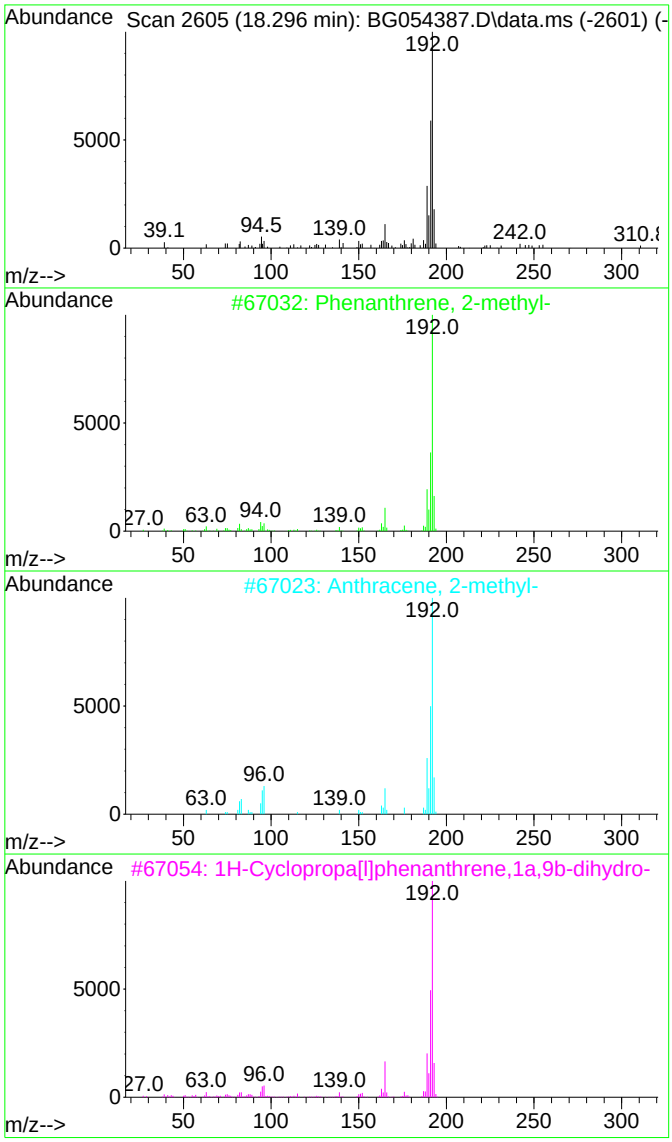
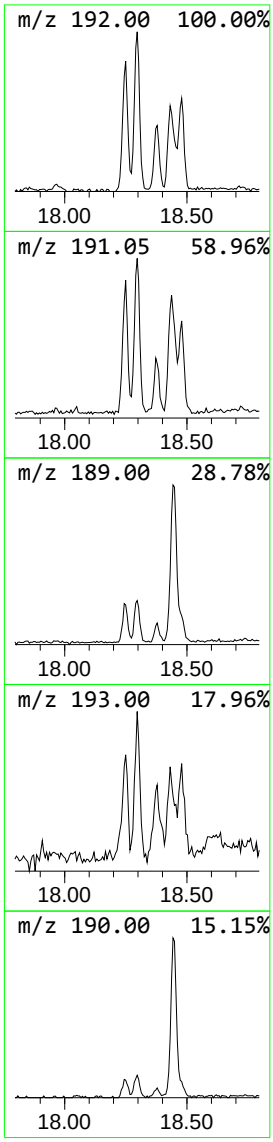
Peak Number 6 Phenanthrene, 2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.296	4.30 ng	80900	Phenanthrene-d10	17.374

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
2	Anthracene, 2-methyl-	192	C15H12	000613-12-7	90
3	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	90
4	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	89
5	Anthracene, 1-methyl-	192	C15H12	000610-48-0	89





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072522\
Data File : BG054387.D
Acq On : 25 Jul 2022 20:08
Operator : CG/JU
Sample : N3855-01 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EO-01-072222

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG071522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

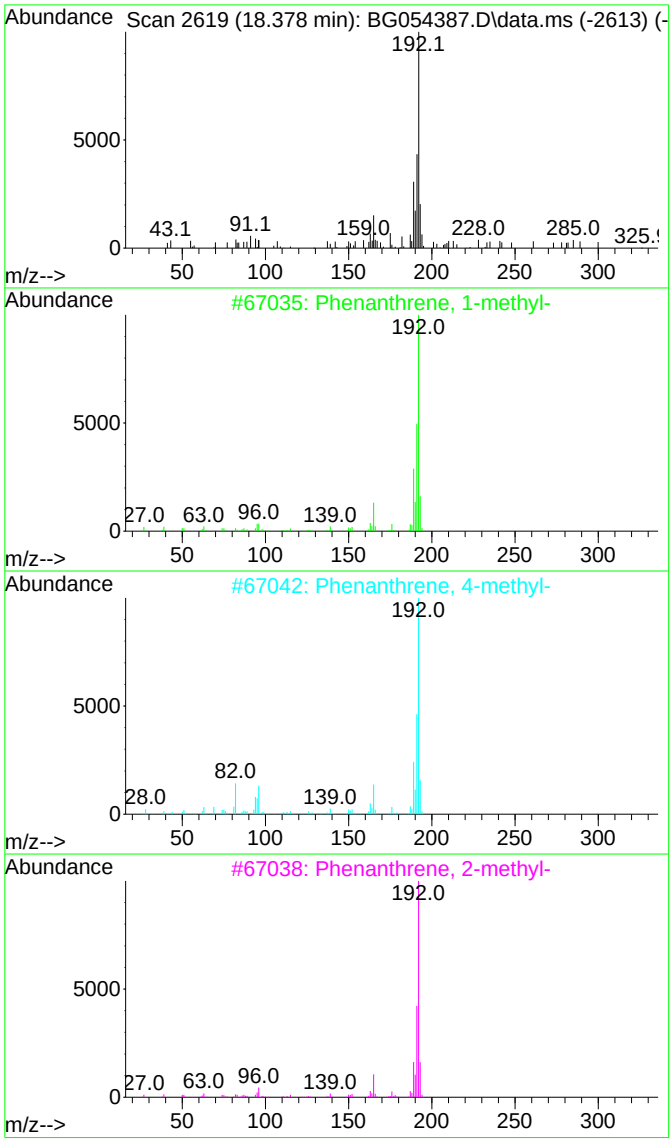
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

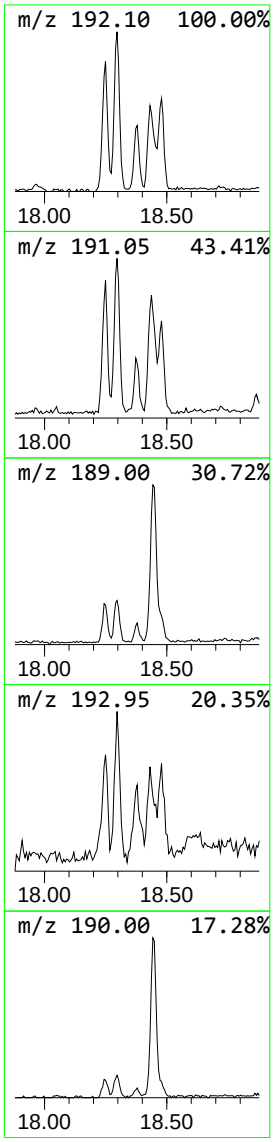
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.378	2.15 ng	40486	Phenanthrene-d10	17.374

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
2	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	81
3	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	81
4	Naphtho[2,3-b]norbornadiene	192	C15H12	107426-38-0	81
5	Anthracene, 1-methyl-	192	C15H12	000610-48-0	81





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Data File : BG054387.D
Acq On : 25 Jul 2022 20:08
Operator : CG/JU
Sample : N3855-01 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EO-01-072222

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG071522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

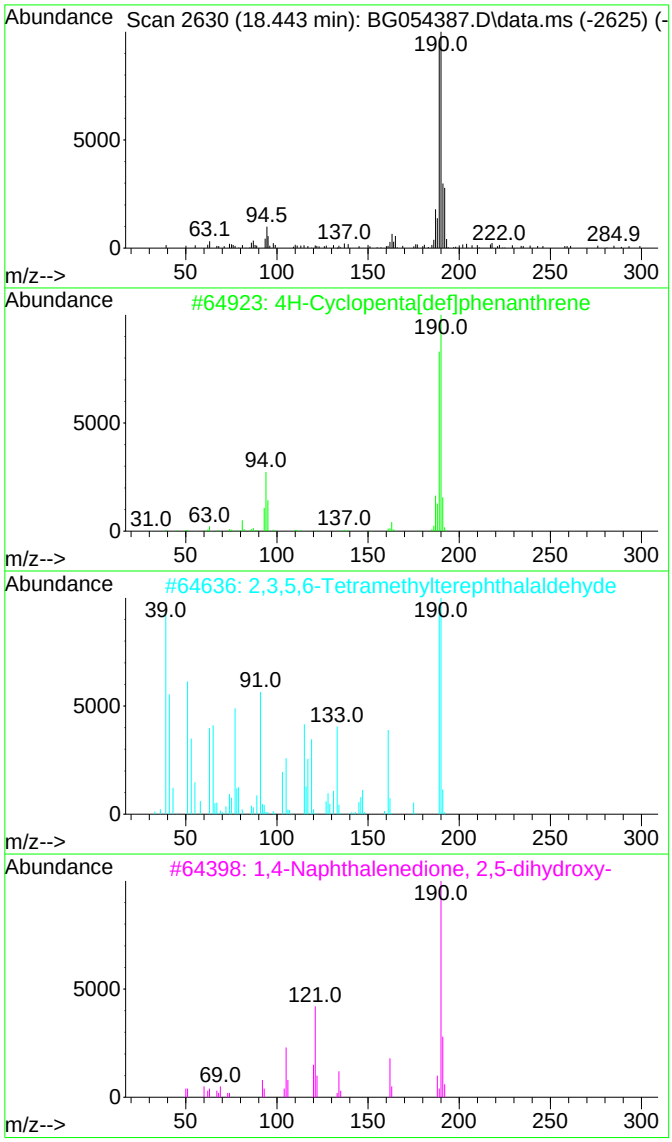
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

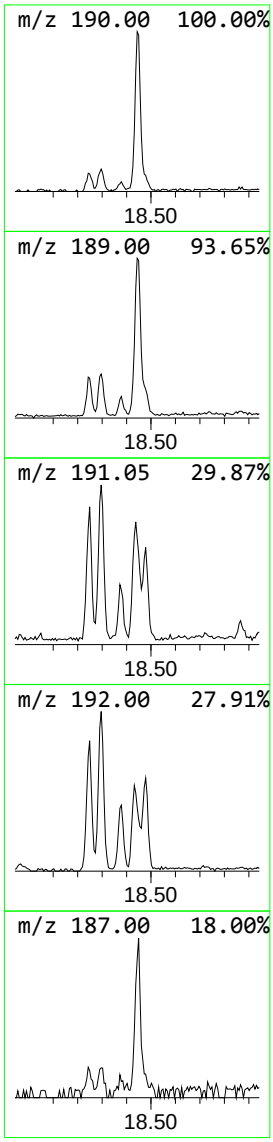
****

Peak Number 8 4H-Cyclopenta[def]phenanthrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.443	6.79 ng	127783	Phenanthrene-d10	17.374

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	87
2	2,3,5,6-Tetramethylterephthalald...	190	C12H14O2	007072-01-7	38
3	1,4-Naphthalenedione, 2,5-dihydr...	190	C10H6O4	004923-55-1	38
4	3,3',5,5'-Tetramethyl-2,2'-bifuryl	190	C12H14O2	222187-59-9	27
5	Epilupeol; 20(29)-Lupen-3alpha-o...	468	C32H52O2	1010513-01-8	25





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Data File : BG054387.D
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Operator : CG/JU
Sample : N3855-01 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EO-01-072222

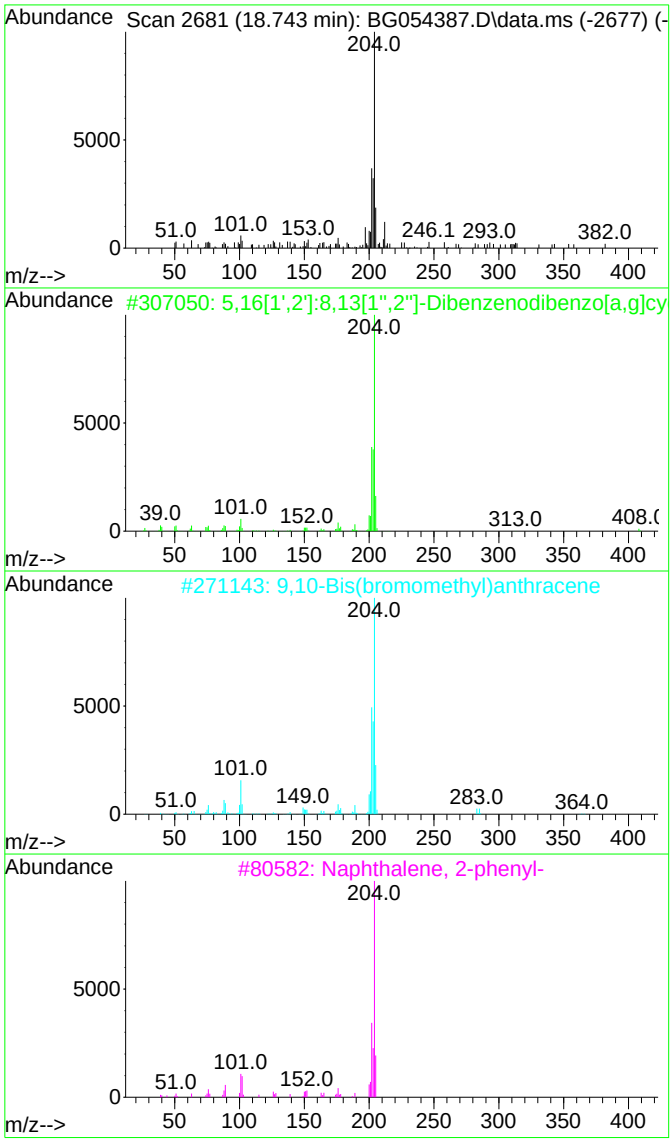
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG071522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

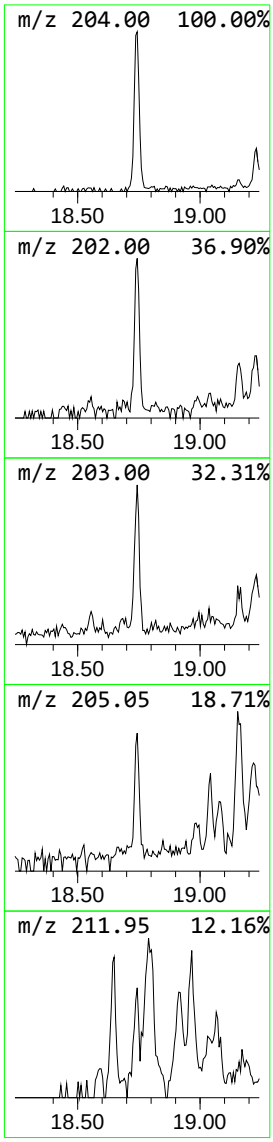
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 5,16[1',2']:8,13[1'',2'']-D... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.743	2.07 ng	39012	Phenanthrene-d10	17.374

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	72
2	9,10-Bis(bromomethyl)anthracene	362	C16H12Br2	034373-96-1	64
3	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	64
4	Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	58
5	[1,1'-Biphenyl]-4,4'-dicarbonitrile	204	C14H8N2	001591-30-6	52





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Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EO-01-072222

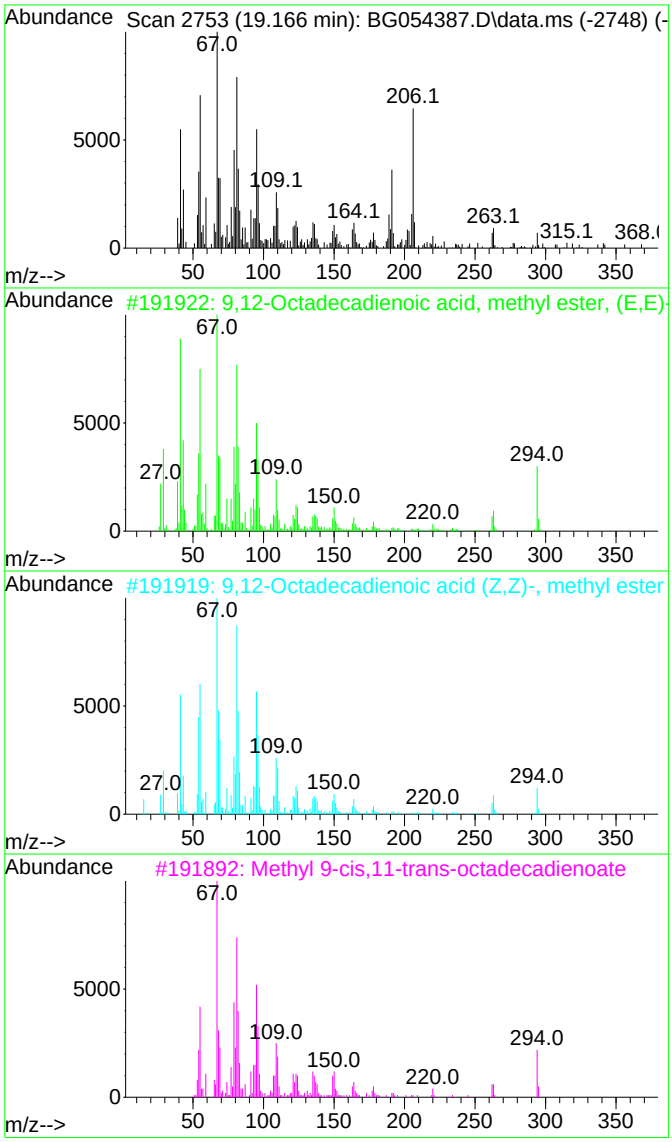
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG071522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

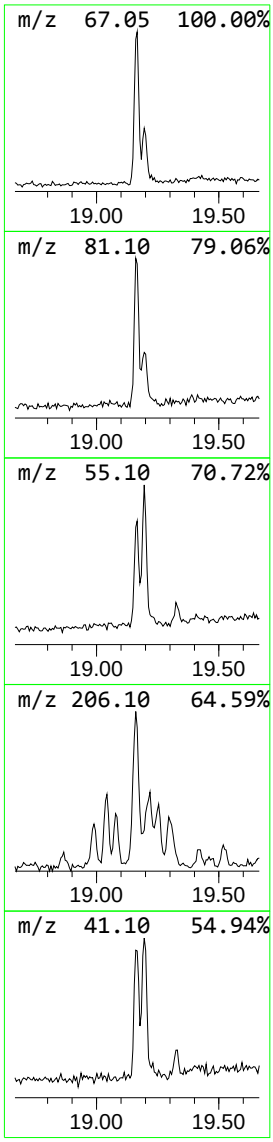
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 9,12-Octadecadienoic acid, ... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.166	10.64 ng	200284	Phenanthrene-d10	17.374

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002566-97-4	99
2	9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	99
3	Methyl 9-cis,11-trans-octadecadi...	294	C19H34O2	013058-52-1	98
4	9,11-Octadecadienoic acid, methy...	294	C19H34O2	013038-47-6	94
5	8,11-Octadecadienoic acid, methy...	294	C19H34O2	056599-58-7	94





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Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EO-01-072222

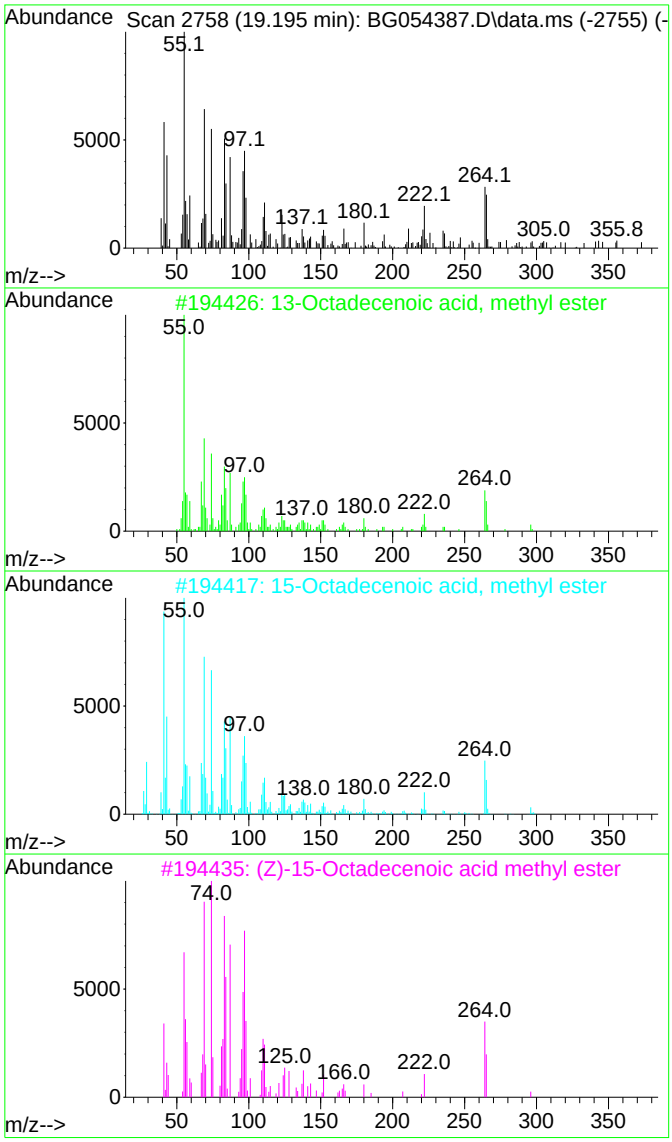
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG071522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

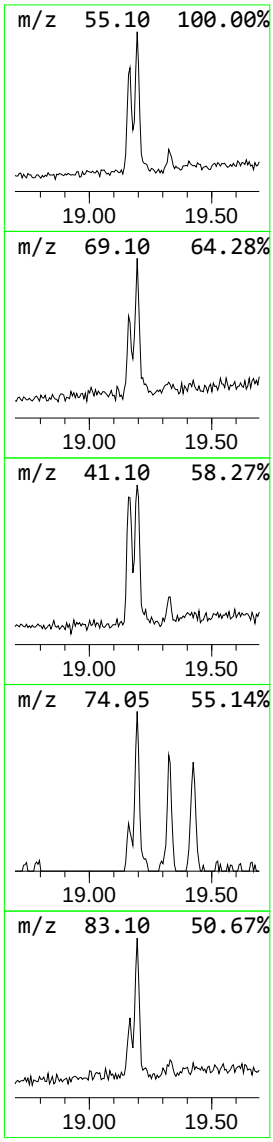
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 13-Octadecenoic acid, methy... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.195	11.69 ng	220095	Phenanthrene-d10	17.374

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	13-Octadecenoic acid, methyl ester	296	C19H36O2	056554-47-3	94
2	15-Octadecenoic acid, methyl ester	296	C19H36O2	004764-72-1	93
3	(Z)-15-Octadecenoic acid methyl ...	296	C19H36O2	1000427-69-3	90
4	3-Octadecenoic acid, methyl ester	296	C19H36O2	056599-33-8	89
5	16-Octadecenoic acid, methyl ester	296	C19H36O2	056554-49-5	86





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Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EO-01-072222

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG071522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

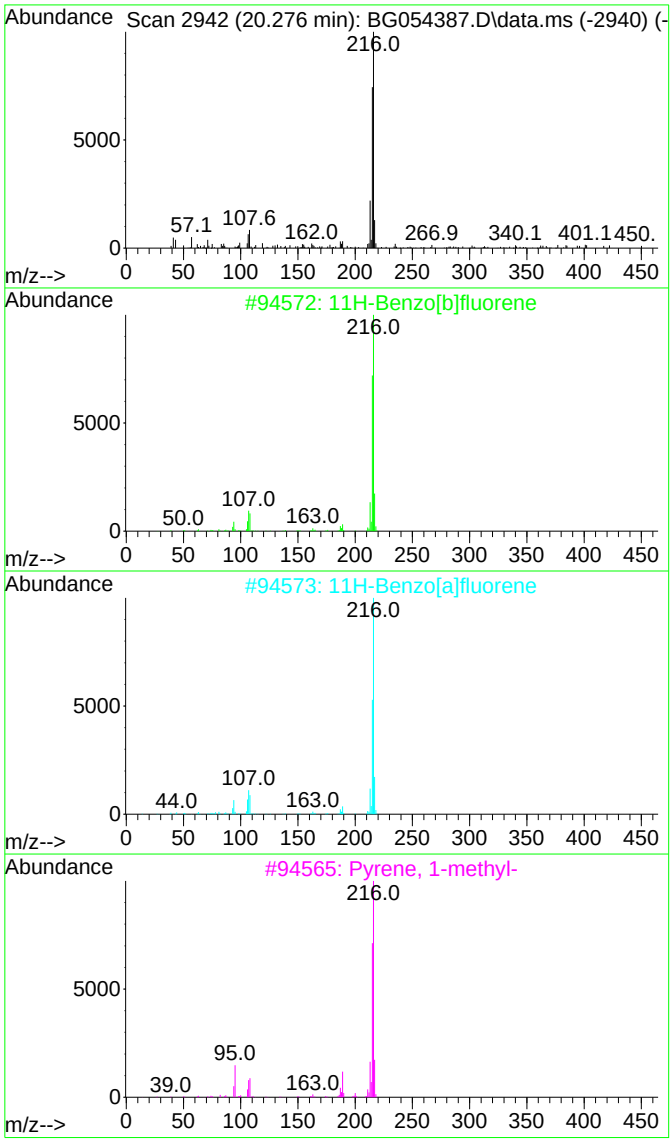
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

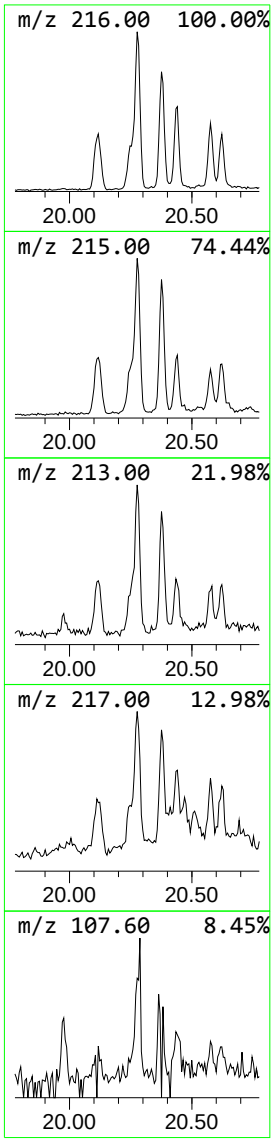
****

Peak Number 12 11H-Benzo[b]fluorene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.276	2.35 ng	99002	Chrysene-d12	21.639

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	87
2	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	80
3	Pyrene, 1-methyl-	216	C17H12	002381-21-7	78
4	3-Trifluoromethylcinnamic acid	216	C10H7F3O2	000779-89-5	72
5	trans-p-(Trifluoromethyl)cinnami...	216	C10H7F3O2	016642-92-5	64





Tentatively Identified Compound (LSC) summary

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Operator : CG/JU
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Misc :
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Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG071522.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...	3.120	2.7	ng	19503	1	7.991	144541	20.0
unknown16.352	16.352	2.3	ng	42355	4	17.374	376414	20.0
Dibenzothiophene	17.186	2.4	ng	44750	4	17.374	376414	20.0
Hexadecanoic ac...	18.026	2.8	ng	52562	4	17.374	376414	20.0
Anthracene, 1-m...	18.249	3.6	ng	68452	4	17.374	376414	20.0
Phenanthrene, 2...	18.296	4.3	ng	80900	4	17.374	376414	20.0
Phenanthrene, 1...	18.378	2.1	ng	40486	4	17.374	376414	20.0
4H-Cyclopenta[d...	18.443	6.8	ng	127783	4	17.374	376414	20.0
5,16[1',2']:8,1...	18.743	2.1	ng	39012	4	17.374	376414	20.0
9,12-Octadecadi...	19.166	10.6	ng	200284	4	17.374	376414	20.0
13-Octadecenoic...	19.195	11.7	ng	220095	4	17.374	376414	20.0
11H-Benzo[b]flu...	20.276	2.4	ng	99002	5	21.639	843994	20.0