LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111822\
Data File : BG055698.D
Acq On : 18 Nov 2022 20:16
Operator : CG/JU
Sample : N5691-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
ETGI-358

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-BG111622.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG055698.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.297	332	342	350	rVB	99713	154578	6.87%	1.014%
2	5.779	417	424	433	rBV	553133	884346	39.28%	5.801%
3	7.395	691	699	717	rVB	578079	1016055	45.13%	6.665%
4	8.247	837	844	851	rBV	114024	192454	8.55%	1.262%
5	9.422	1036	1044	1057	rVB	313957	576091	25.59%	3.779%
6	11.079	1318	1326	1331	rBV	169991	312667	13.89%	2.051%
7	11.126	1331	1334	1340	rVB	31131	49456	2.20%	0.324%
8	12.712	1599	1604	1609	rBV	14220	23738	1.05%	0.156%
9	13.494	1729	1737	1745	rVB	953586	1477215	65.61%	9.690%
10	14.193	1850	1856	1861	rBV2	14653	24691	1.10%	0.162%
11	14.869	1961	1971	1979	rVB	326425	529512	23.52%	3.473%
12	15.909	2142	2148	2155	rVB4	16510	32135	1.43%	0.211%
13	16.214	2191	2200	2207	rVB2	24253	44869	1.99%	0.294%
14	16.355	2217	2224	2234	rBV	915355	1500801	66.66%	9.845%
15	16.537	2248	2255	2259	rBV2	33418	55288	2.46%	0.363%
16	17.260	2374	2378	2386	rVB2	13792	27329	1.21%	0.179%
17	17.430	2404	2407	2414	rVB	16871	26033	1.16%	0.171%
18	17.612	2432	2438	2442	rBV	463035	694788	30.86%	4.558%
19	17.654	2442	2445	2452	rVB	187824	263090	11.68%	1.726%
20	17.742	2452	2460	2464	rBV	48711	86733	3.85%	0.569%
21	18.194	2533	2537	2542	rVB	21395	34060	1.51%	0.223%
22	18.482	2581	2586	2590	rBV	94445	157599	7.00%	1.034%
23	18.529	2590	2594	2600	rVV	91601	157418	6.99%	1.033%
24	18.611	2604	2608	2612	rVV2	15483	25574	1.14%	0.168%
25	18.670	2613	2618	2622	rVV2	70349	142447	6.33%	0.934%
26	18.711	2622	2625	2630	rVB	59209	84637	3.76%	0.555%
27	18.864	2647	2651	2657	rVB4	22758	37634	1.67%	0.247%
28	18.970	2665	2669	2673	rBV	48777	67149	2.98%	0.440%
29	19.017	2673	2677	2682	rVB6	21551	37936	1.68%	0.249%
30	19.216	2708	2711	2715	rVB4	20199	26700	1.19%	0.175%
31	19.269	2715	2720	2723	rVV	29665	44879	1.99%	0.294%
32	19.305	2723	2726	2730	rVB3	21291	29068	1.29%	0.191%
33	19.387	2735	2740	2745	rBV	73307	120143	5.34%	0.788%
34	19.440	2745	2749	2754	rBV5	22498	48782	2.17%	0.320%
35	19.528	2759	2764	2771	rBV4	36137	79784	3.54%	0.523%
36	19.651	2779	2785	2790	rBV	250829	376365	16.72%	2.469%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111822\
Data File : BG055698.D
Acq On : 18 Nov 2022 20:16
Operator : CG/JU
Sample : N5691-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
ETGI-358

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-BG111622.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	19.710	2790	2795	2796	rBV3	17741	25870	1.15%	0.170%
38	19.739	2796	2800	2807	rVV3	55753	115639	5.14%	0.759%
39	20.016	2842	2847	2853	rVB	351354	518620	23.03%	3.402%
40	20.074	2854	2857	2863	rVB	31368	47766	2.12%	0.313%
41	20.198	2872	2878	2883	rBV	1732938	2251569	100.00%	14.769%
42	20.333	2895	2901	2908	rBV6	38975	96805	4.30%	0.635%
43	20.491	2920	2928	2933	rVB2	62668	150227	6.67%	0.985%
44	20.597	2943	2946	2950	rVB	29245	34341	1.53%	0.225%
45	20.656	2953	2956	2960	rVB	61917	72899	3.24%	0.478%
46	20.797	2976	2980	2984	rBV	57795	84926	3.77%	0.557%
47	20.844	2985	2988	2997	rVB2	54270	97377	4.32%	0.639%
48	21.273	3059	3061	3066	rVB	38814	52667	2.34%	0.345%
49	21.514	3098	3102	3108	rBV	102592	179520	7.97%	1.178%
50	21.907	3161	3169	3174	rBV2	422543	939881	41.74%	6.165%
51	21.960	3174	3178	3190	rVB	172788	316049	14.04%	2.073%
52	24.234	3560	3565	3571	rBV2	35447	93835	4.17%	0.616%
53	25.327	3742	3751	3761	rVB2	253512	724840	32.19%	4.755%

Sum of corrected areas: 15244875

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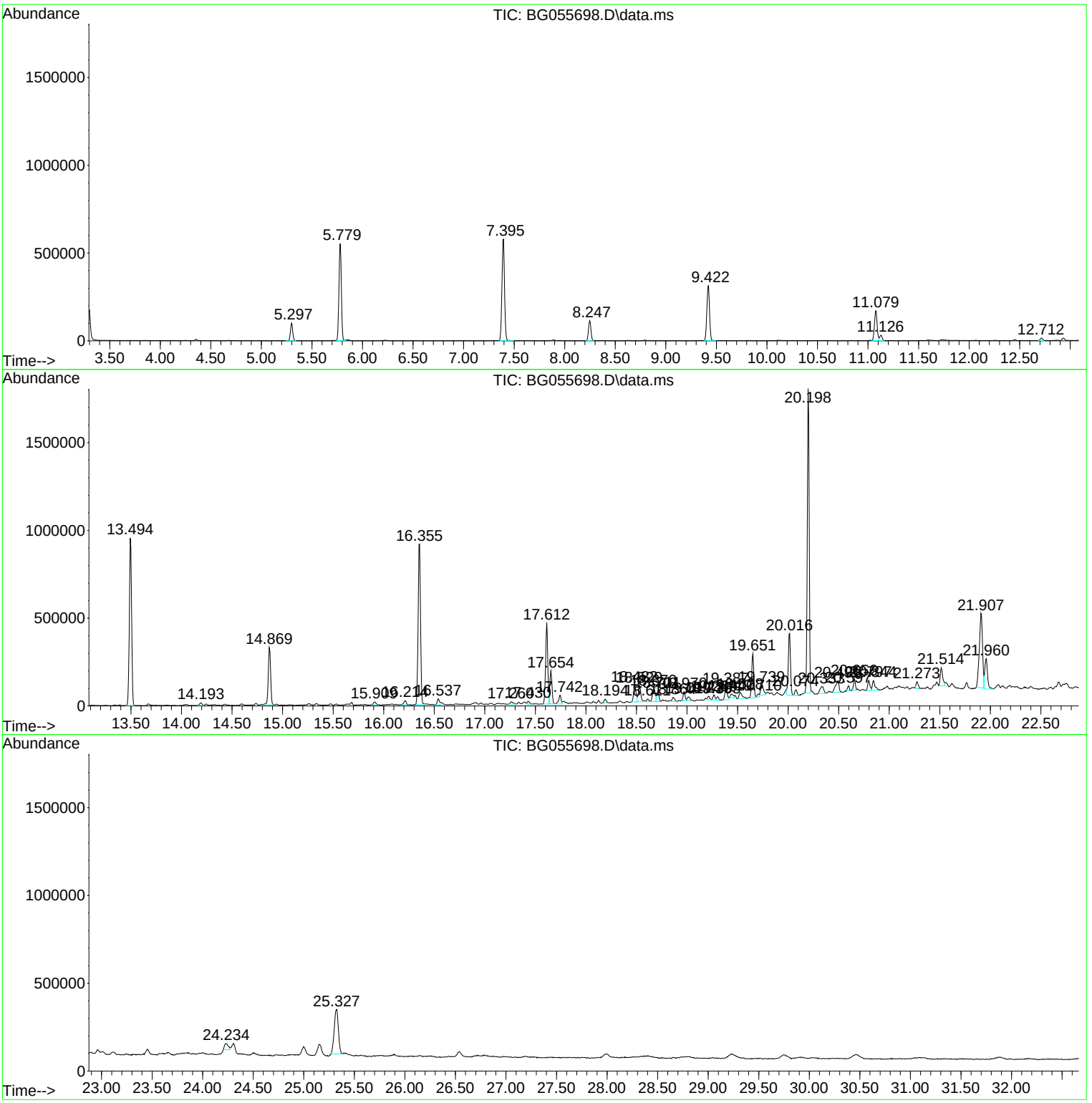
Operator : CG/JU
Sample : N5691-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
ETGI-358

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111622.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\BG111822\
Data File : BG055698.D
Acq On : 18 Nov 2022 20:16
Operator : CG/JU
Sample : N5691-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
ETGI-358

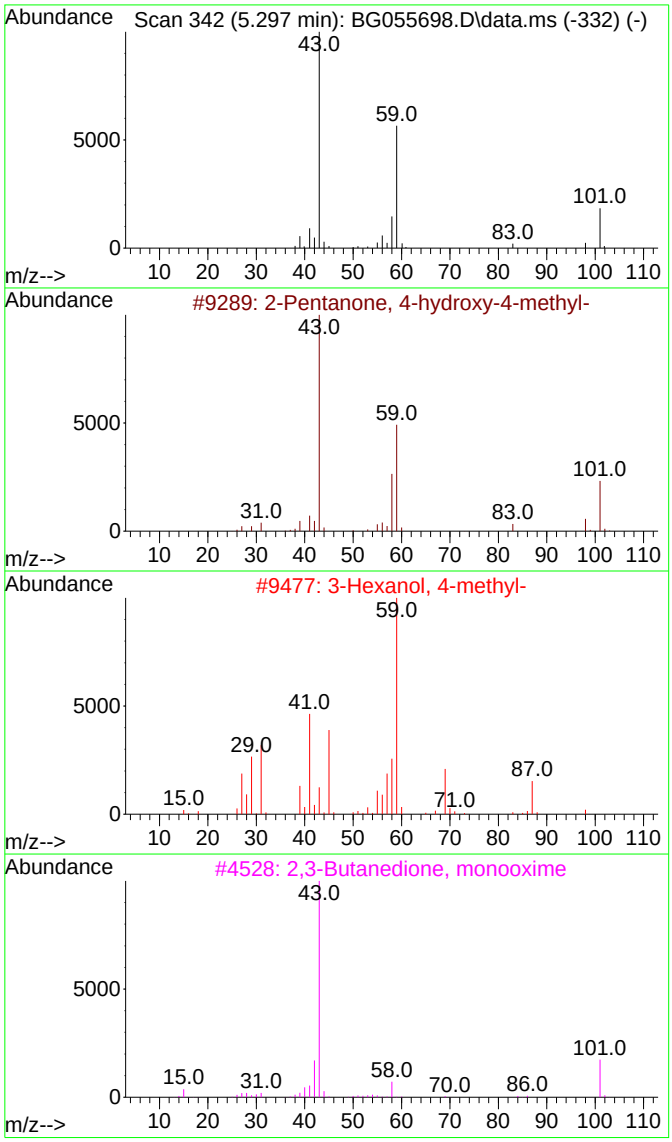
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

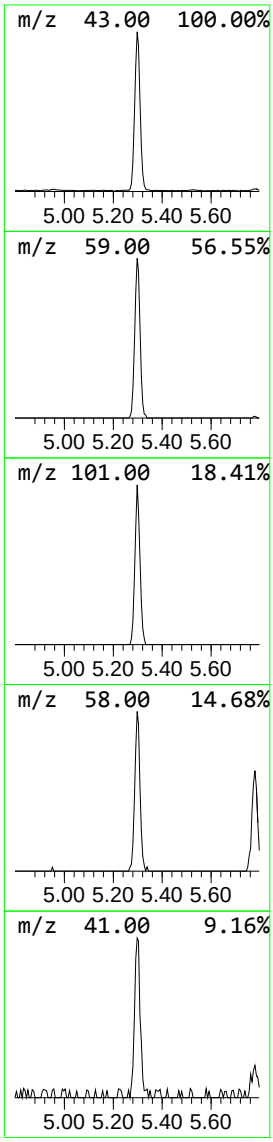
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.297	16.06 ng	154578	1,4-Dichlorobenzene-d4	8.247

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
3	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	27
4	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
5	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	12





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111822\
Data File : BG055698.D
Acq On : 18 Nov 2022 20:16
Operator : CG/JU
Sample : N5691-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
ETGI-358

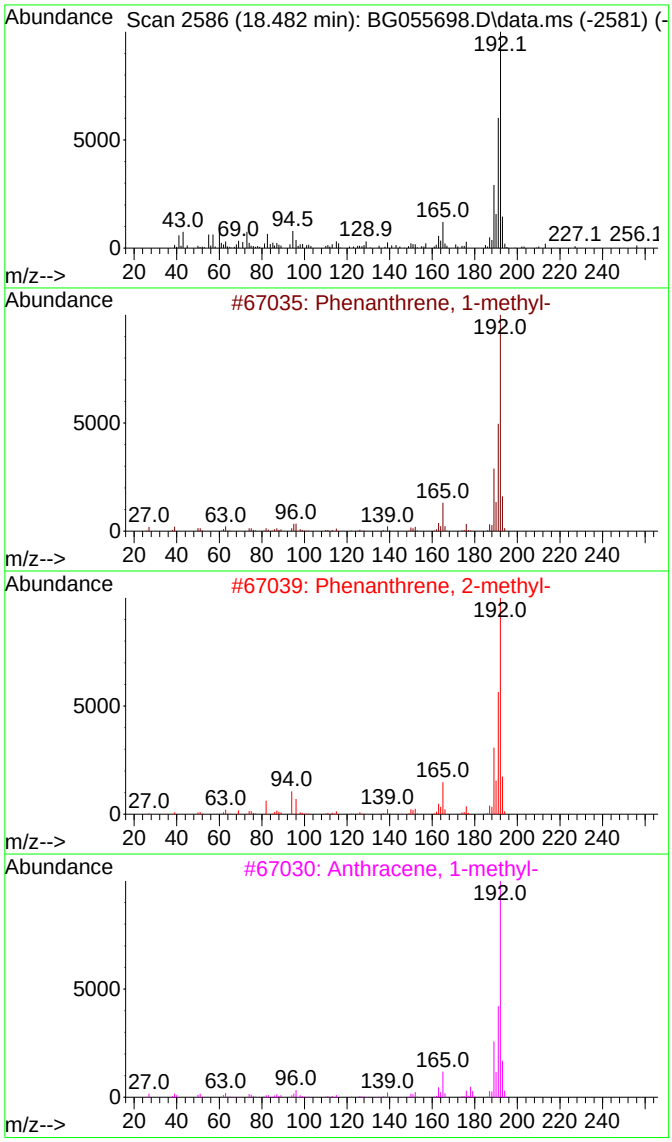
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

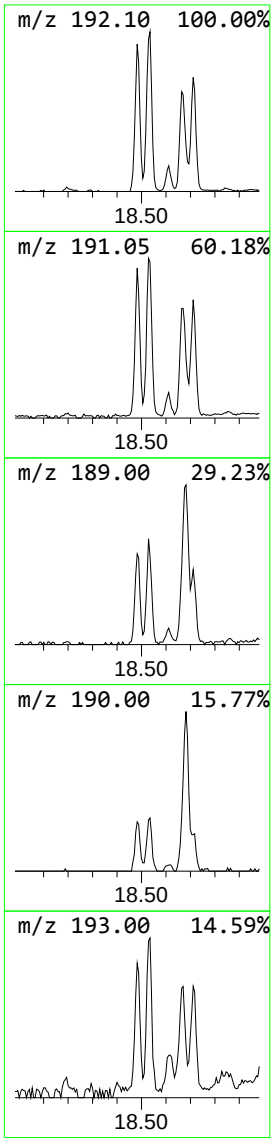
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Phenanthrene, 1-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.482	4.54 ng	157599	Phenanthrene-d10	17.613

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	96
3	Anthracene, 1-methyl-	192	C15H12	000610-48-0	94
4	Anthracene, 2-methyl-	192	C15H12	000613-12-7	94
5	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	93





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111822\
Data File : BG055698.D
Acq On : 18 Nov 2022 20:16
Operator : CG/JU
Sample : N5691-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
ETGI-358

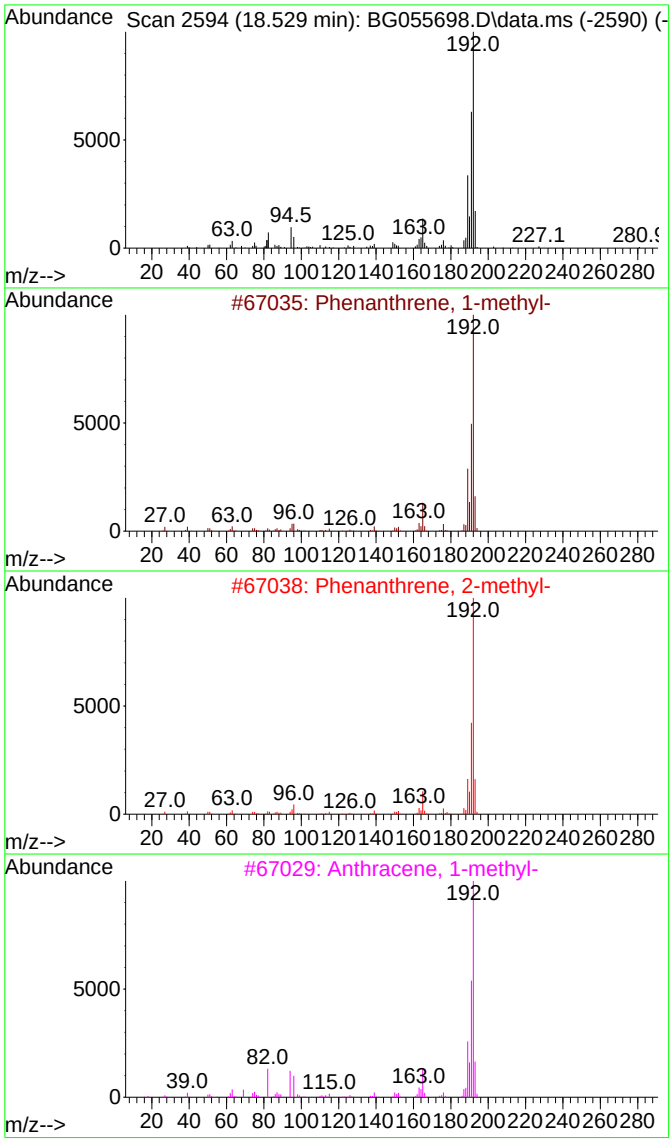
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

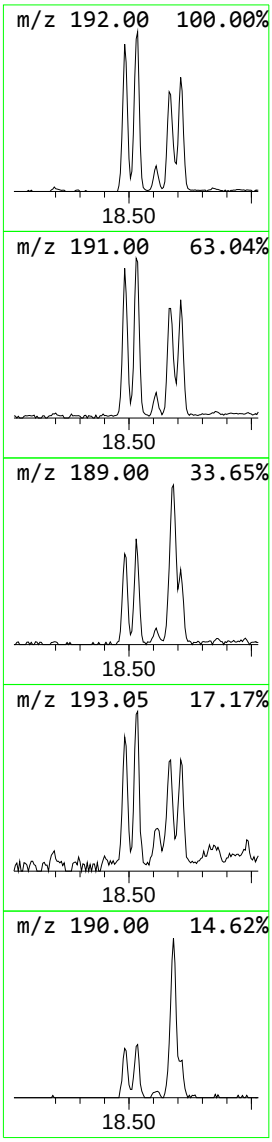
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Phenanthrene, 2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.529	4.53 ng	157418	Phenanthrene-d10	17.613

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
2	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	94
3	Anthracene, 1-methyl-	192	C15H12	000610-48-0	93
4	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	93
5	Anthracene, 2-methyl-	192	C15H12	000613-12-7	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111822\
Data File : BG055698.D
Acq On : 18 Nov 2022 20:16
Operator : CG/JU
Sample : N5691-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
ETGI-358

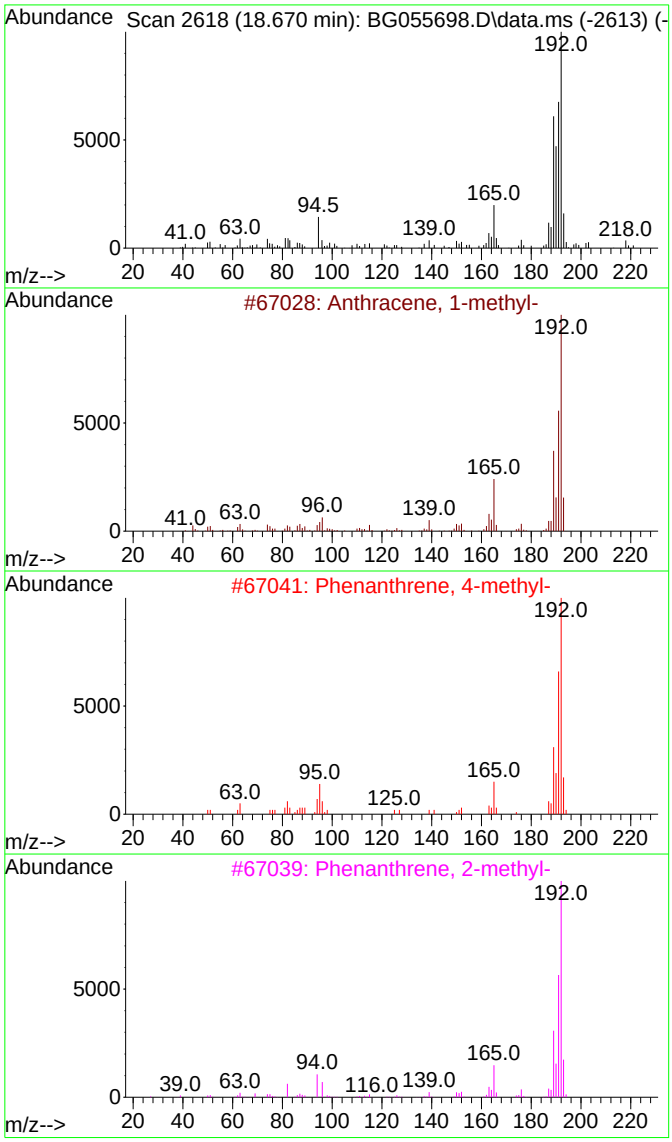
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

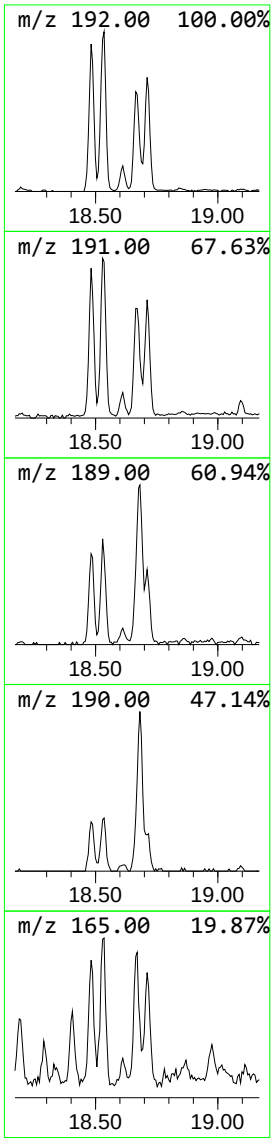
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Anthracene, 1-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.670	4.10 ng	142447	Phenanthrene-d10	17.613

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 1-methyl-	192	C15H12	000610-48-0	89
2	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	78
3	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	64
4	8,9-Dihydrocyclopenta[def]phenan...	192	C15H12	027410-55-5	62
5	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	60





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111822\
Data File : BG055698.D
Acq On : 18 Nov 2022 20:16
Operator : CG/JU
Sample : N5691-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
ETGI-358

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

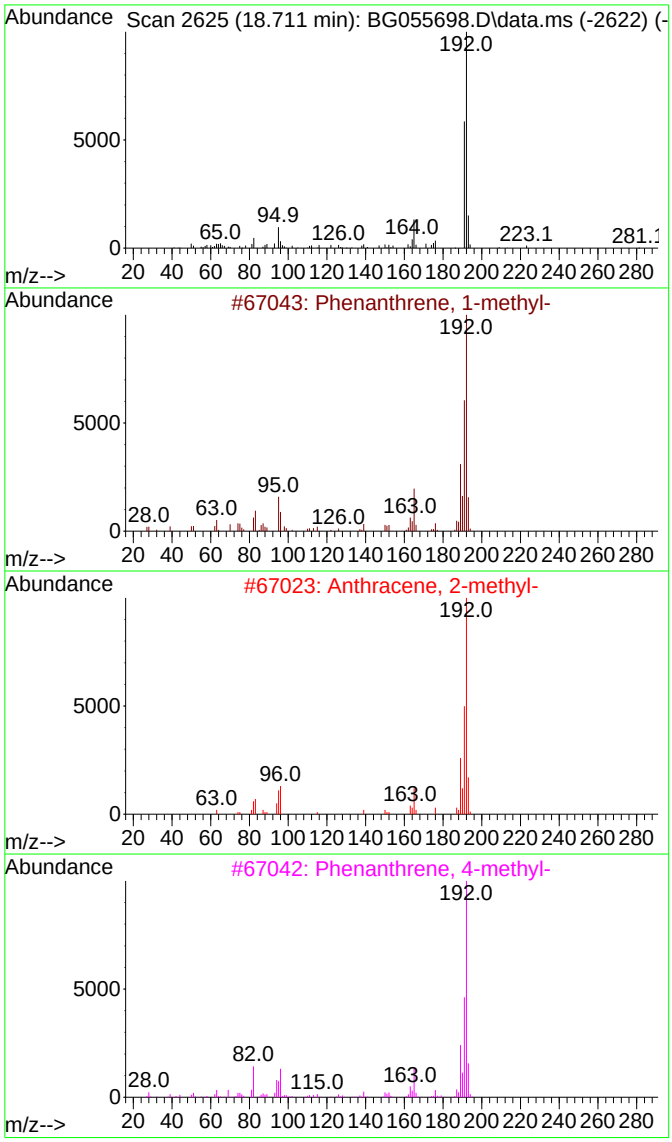
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

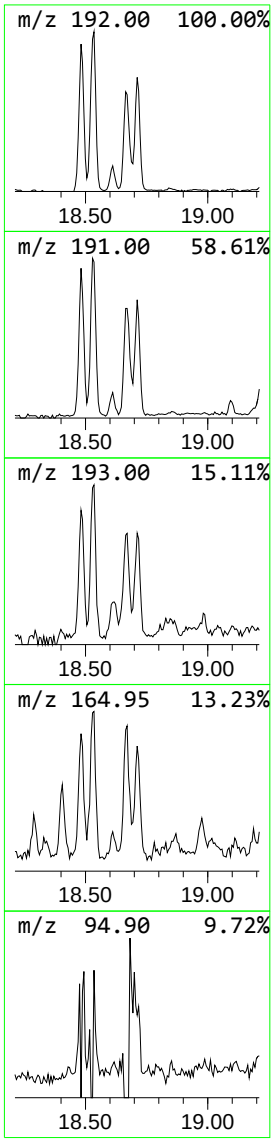
****

Peak Number 5 Anthracene, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.711	2.44 ng	84637	Phenanthrene-d10	17.613

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	91
2	Anthracene, 2-methyl-	192	C15H12	000613-12-7	91
3	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	91
4	Anthracene, 9-methyl-	192	C15H12	000779-02-2	91
5	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	90





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111822\
Data File : BG055698.D
Acq On : 18 Nov 2022 20:16
Operator : CG/JU
Sample : N5691-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
ETGI-358

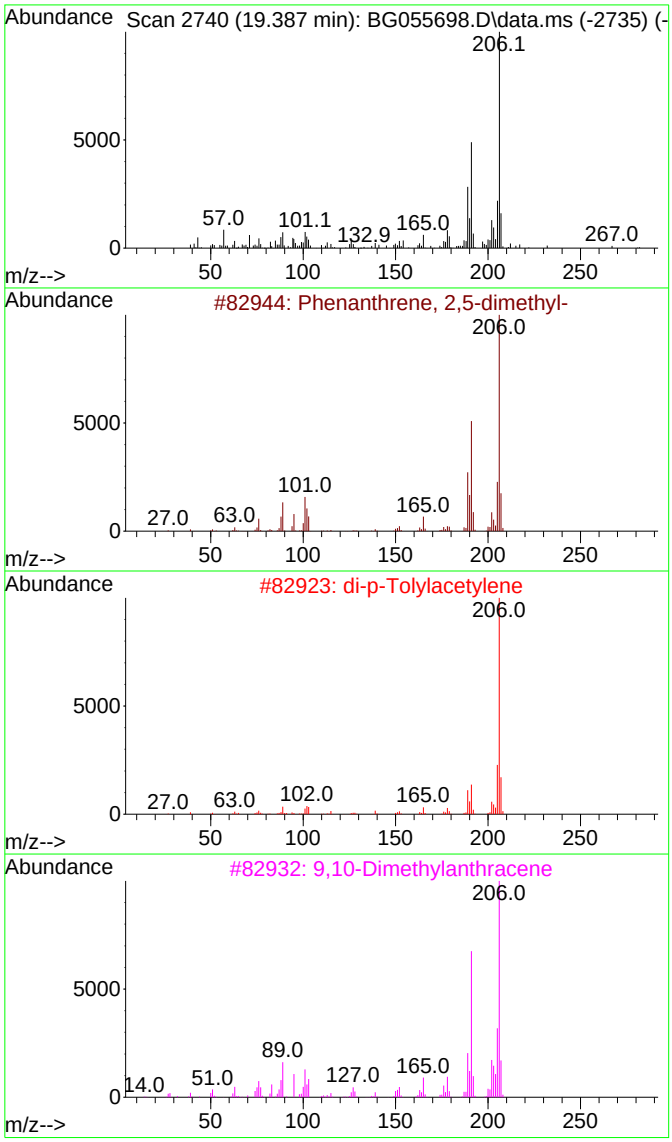
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

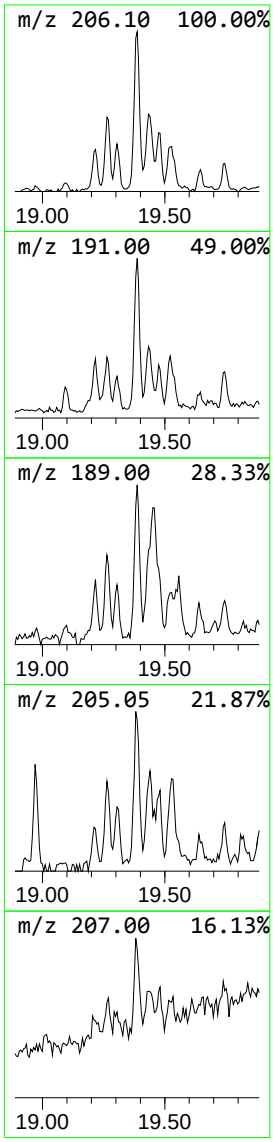
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Phenanthrene, 2,5-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.387	3.46 ng	120143	Phenanthrene-d10	17.613

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	93
2	di-p-Tolylacetylene	206	C16H14	002789-88-0	93
3	9,10-Dimethylanthracene	206	C16H14	000781-43-1	91
4	3,4-Dimethylphenanthrene	206	C16H14	1000452-24-5	90
5	Phenanthrene, 2,7-dimethyl-	206	C16H14	001576-69-8	87





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111822\
Data File : BG055698.D
Acq On : 18 Nov 2022 20:16
Operator : CG/JU
Sample : N5691-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
ETGI-358

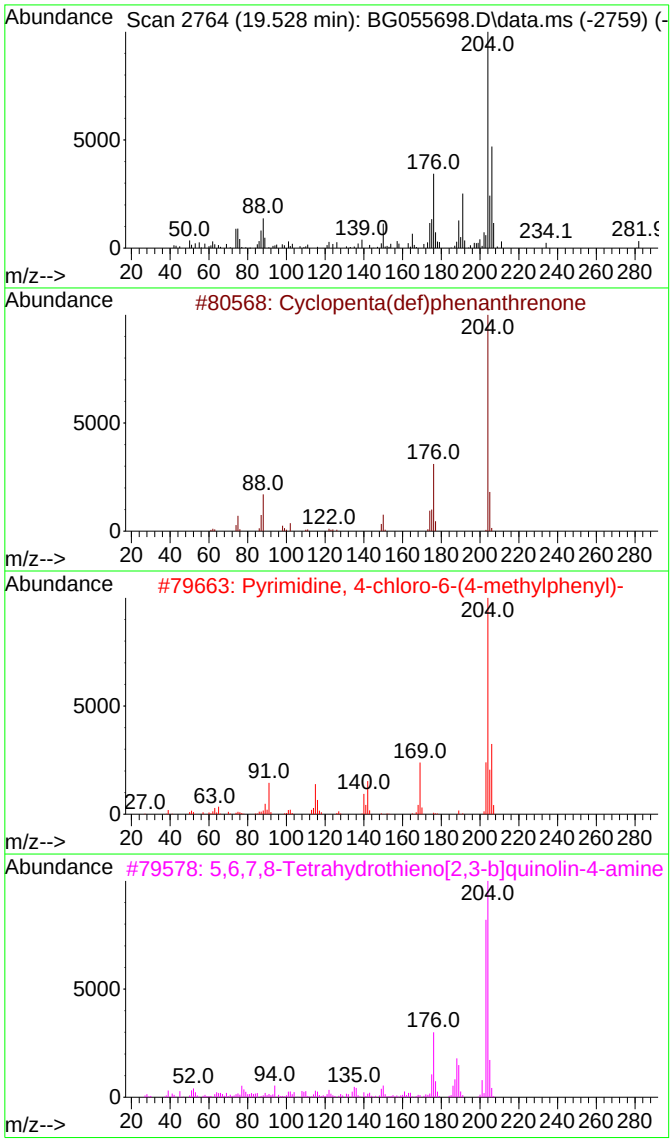
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

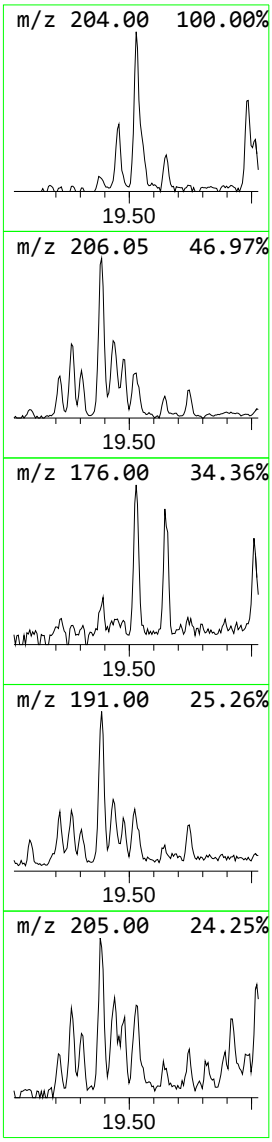
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Cyclopenta(def)phenanthrenone Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.528	2.30 ng	79784	Phenanthrene-d10	17.613

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	60
2	Pyrimidine, 4-chloro-6-(4-methyl...	204	C11H9ClN2	1000380-55-8	50
3	5,6,7,8-Tetrahydrothieno[2,3-b]q...	204	C11H12N2S	122914-50-5	47
4	Pyrimido[5,4-b]benzofurane, 4-ch...	204	C10H5ClN2O	039876-88-5	43
5	2,4-Dichloro-3-ethyl-5-methylphe...	246	C11H12Cl2O2	1000467-41-1	40





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111822\
Data File : BG055698.D
Acq On : 18 Nov 2022 20:16
Operator : CG/JU
Sample : N5691-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
ETGI-358

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

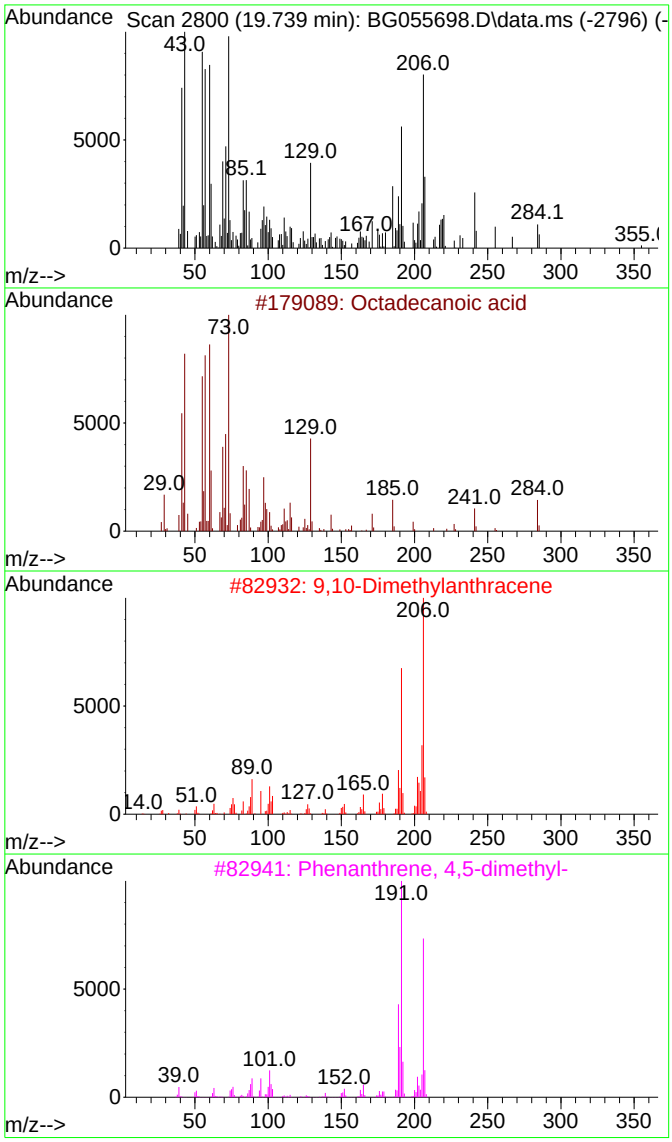
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

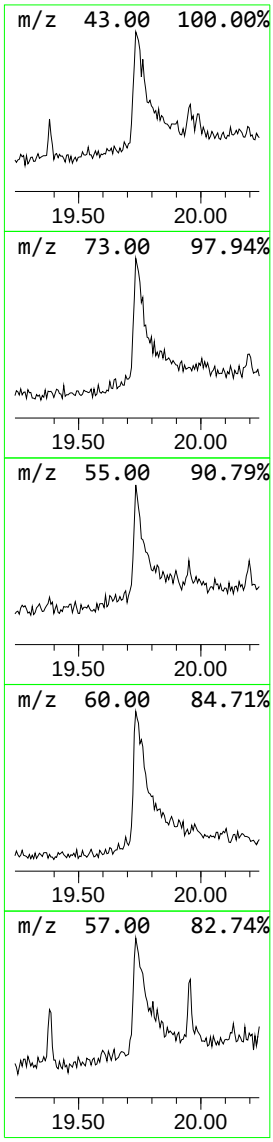
****

Peak Number 8 Octadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.739	3.33 ng	115639	Phenanthrene-d10	17.613

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	83
2	9,10-Dimethylanthracene	206	C16H14	000781-43-1	55
3	Phenanthrene, 4,5-dimethyl-	206	C16H14	003674-69-9	46
4	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	42
5	Anthracene, 2-ethyl-	206	C16H14	052251-71-5	42





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111822\
Data File : BG055698.D
Acq On : 18 Nov 2022 20:16
Operator : CG/JU
Sample : N5691-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
ETGI-358

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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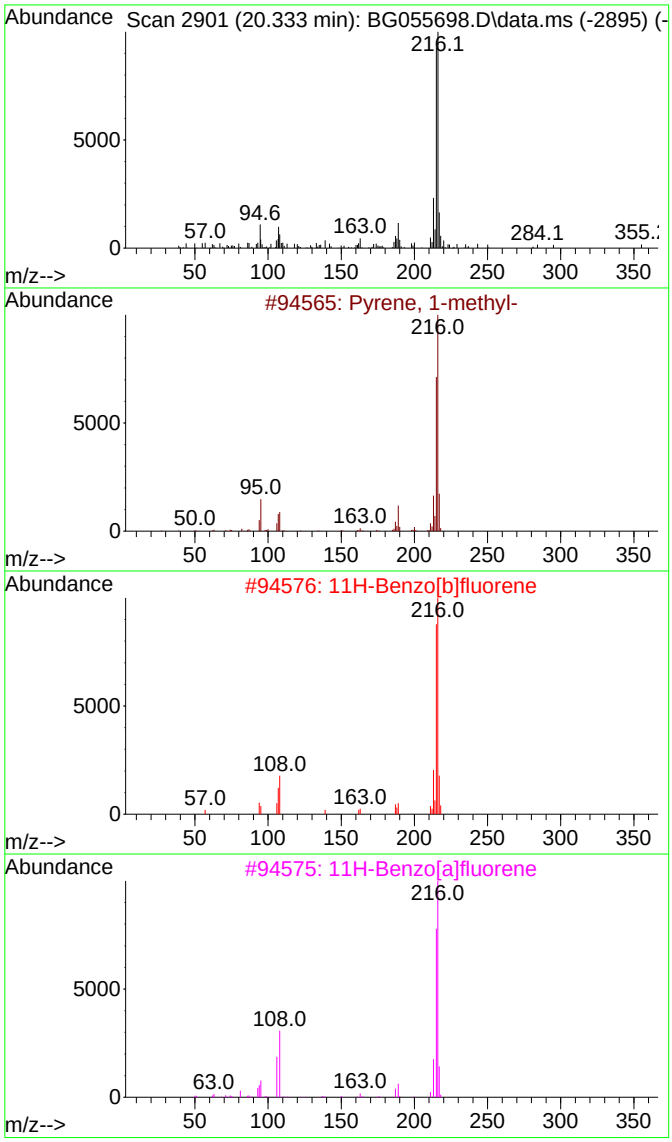
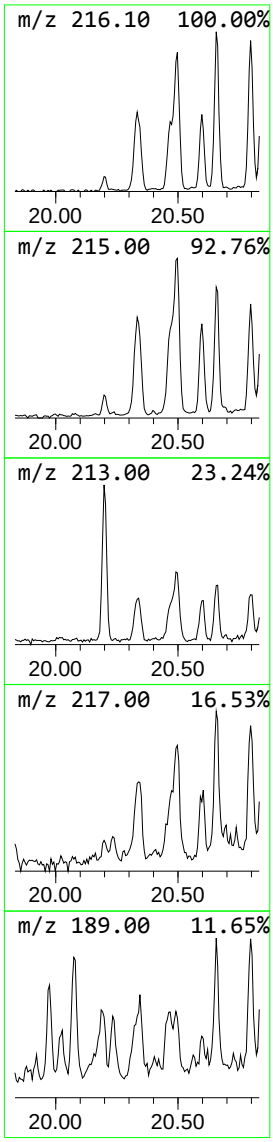
Peak Number 9 Pyrene, 1-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.333	2.06 ng	96805	Chrysene-d12	21.907

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	93
3	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	90
4	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	90
5	7H-Benzo[c]fluorene	216	C17H12	000205-12-9	80





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Data File : BG055698.D
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Operator : CG/JU
Sample : N5691-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
ETGI-358

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111622.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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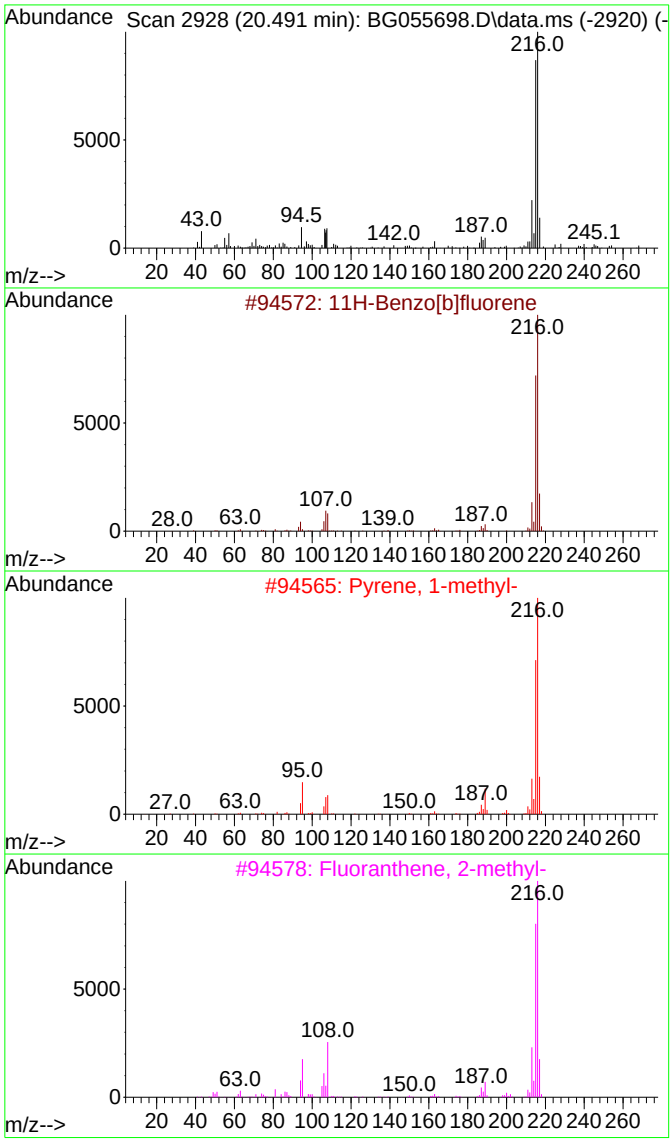
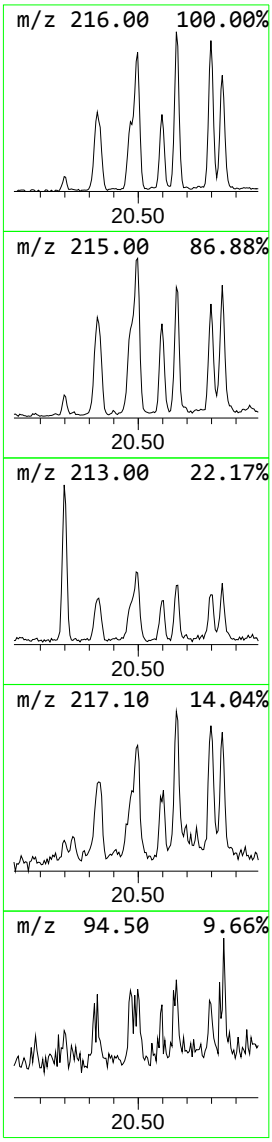
Peak Number 10 11H-Benzo[b]fluorene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.491	3.20 ng	150227	Chrysene-d12	21.907

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	83
2	Pyrene, 1-methyl-	216	C17H12	002381-21-7	83
3	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	81
4	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	62
5	3-Trifluoromethylcinnamic acid	216	C10H7F3O2	000779-89-5	59





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111822\
Data File : BG055698.D
Acq On : 18 Nov 2022 20:16
Operator : CG/JU
Sample : N5691-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
ETGI-358

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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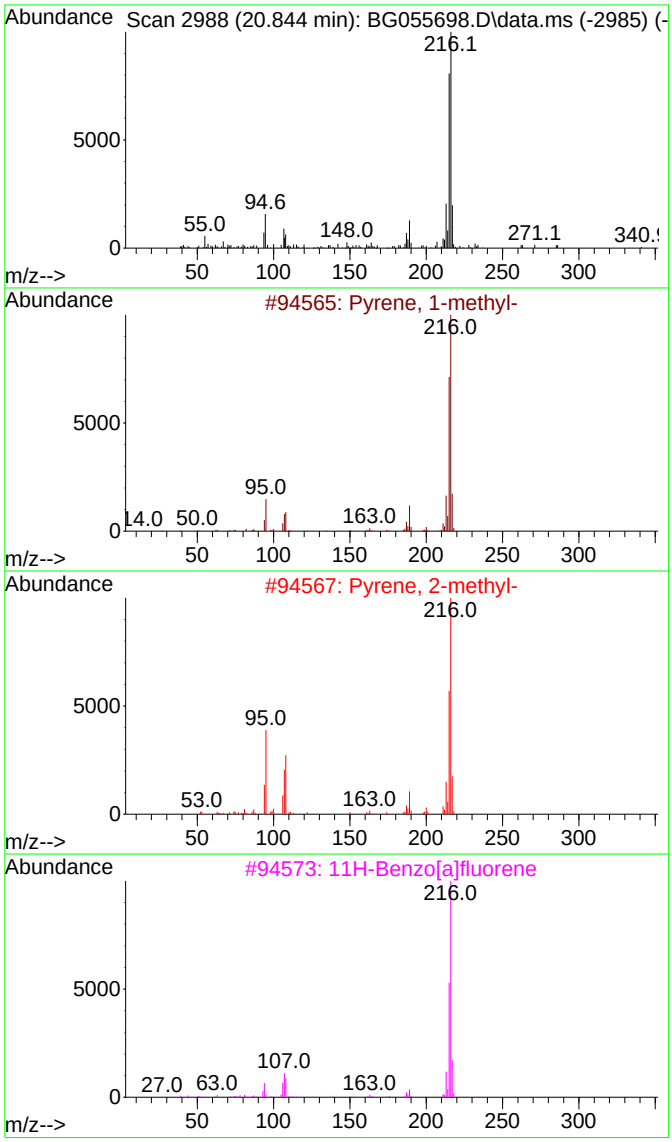
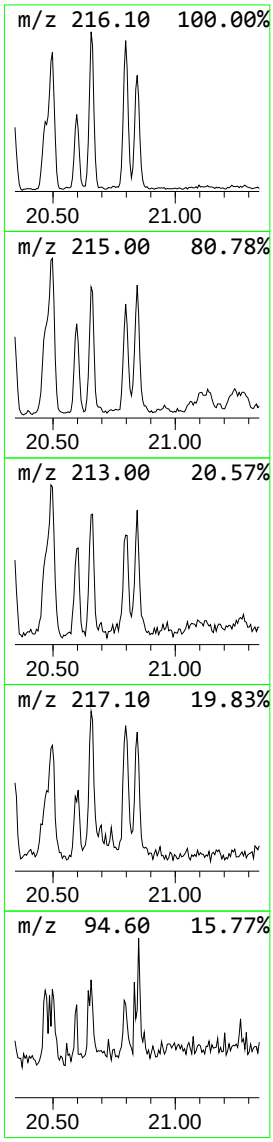
Peak Number 11 Pyrene, 2-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.844	2.07 ng	97377	Chrysene-d12	21.907

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-	216	C17H12	002381-21-7	95
2	Pyrene, 2-methyl-	216	C17H12	003442-78-2	94
3	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	94
4	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	90
5	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	87





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111822\
Data File : BG055698.D
Acq On : 18 Nov 2022 20:16
Operator : CG/JU
Sample : N5691-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
ETGI-358

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111622.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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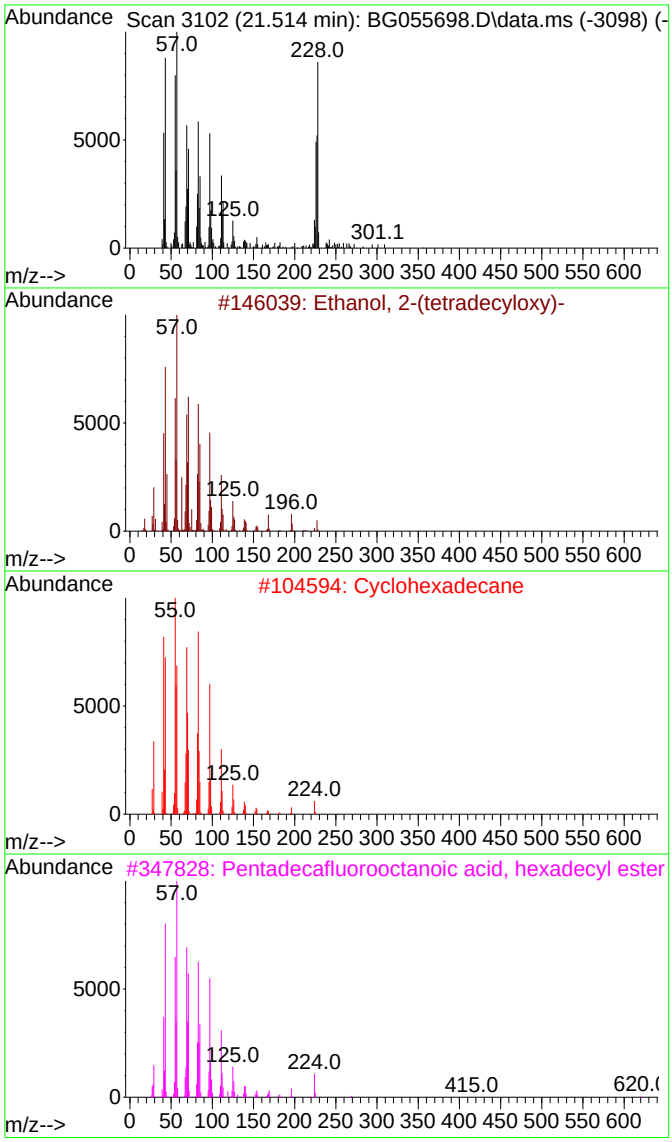
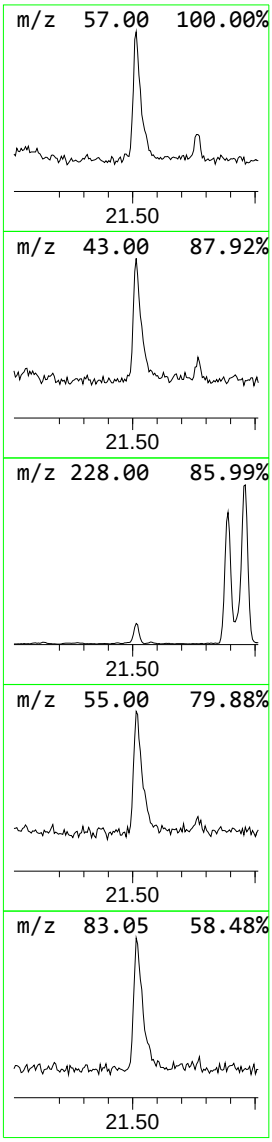
Peak Number 12 Ethanol, 2-(tetradecyloxy)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.514	3.82 ng	179520	Chrysene-d12	21.907

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	94
2	Cyclohexadecane	224	C16H32	000295-65-8	93
3	Pentadecafluorooctanoic acid, he...	638	C24H33F15O2	1000406-04-6	87
4	1-Octadecene	252	C18H36	000112-88-9	70
5	1-Hexadecanol, 2-methyl-	256	C17H36O	002490-48-4	66





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111822\
Data File : BG055698.D
Acq On : 18 Nov 2022 20:16
Operator : CG/JU
Sample : N5691-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
ETGI-358

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111622.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-...	5.297	16.1	ng	154578	1	8.247	192454	20.0
Phenanthrene, 1...	18.482	4.5	ng	157599	4	17.613	694788	20.0
Phenanthrene, 2...	18.529	4.5	ng	157418	4	17.613	694788	20.0
Anthracene, 1-m...	18.670	4.1	ng	142447	4	17.613	694788	20.0
Anthracene, 2-m...	18.711	2.4	ng	84637	4	17.613	694788	20.0
Phenanthrene, 2...	19.387	3.5	ng	120143	4	17.613	694788	20.0
Cyclopenta(def)...	19.528	2.3	ng	79784	4	17.613	694788	20.0
Octadecanoic acid	19.739	3.3	ng	115639	4	17.613	694788	20.0
Pyrene, 1-methyl-	20.333	2.1	ng	96805	5	21.907	939881	20.0
11H-Benzo[b]flu...	20.491	3.2	ng	150227	5	21.907	939881	20.0
Pyrene, 2-methyl-	20.844	2.1	ng	97377	5	21.907	939881	20.0
Ethanol, 2-(tet...	21.514	3.8	ng	179520	5	21.907	939881	20.0