LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121922\
Data File : BG056023.D
Acq On : 19 Dec 2022 21:33
Operator : CG/JU
Sample : N6026-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
B-29-SB01

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-BG121322.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG056023.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.400	14	19	26	rBV	30650	54506	8.01%	1.025%
2	5.416	357	362	368	rBB	19270	27967	4.11%	0.526%
3	5.892	436	443	449	rBB	113902	182592	26.83%	3.434%
4	7.501	709	717	724	rBB	121104	210683	30.96%	3.962%
5	8.371	859	865	871	rBB2	38879	63002	9.26%	1.185%
6	9.540	1057	1064	1072	rBB	74763	134305	19.73%	2.526%
7	11.203	1340	1347	1353	rBV	49285	86736	12.74%	1.631%
8	13.606	1749	1756	1763	rBB	271748	404227	59.39%	7.602%
9	14.975	1984	1989	1995	rVB	88983	136068	19.99%	2.559%
10	15.040	1995	2000	2004	rBV2	5200	7986	1.17%	0.150%
11	15.369	2051	2056	2061	rVB2	5132	7851	1.15%	0.148%
12	16.015	2161	2166	2174	rBV2	6362	11398	1.67%	0.214%
13	16.450	2234	2240	2247	rVB	209779	315835	46.41%	5.939%
14	16.673	2274	2278	2284	rVB3	4555	7119	1.05%	0.134%
15	17.360	2389	2395	2400	rBV2	5769	8058	1.18%	0.152%
16	17.437	2402	2408	2413	rVB3	4316	7097	1.04%	0.133%
17	17.531	2419	2424	2431	rBV3	4967	7950	1.17%	0.150%
18	17.713	2449	2455	2459	rBV	122212	182814	26.86%	3.438%
19	17.754	2459	2462	2468	rVB	96082	134858	19.81%	2.536%
20	17.848	2470	2478	2482	rBV	20135	32960	4.84%	0.620%
21	18.107	2516	2522	2526	rBV3	7366	12021	1.77%	0.226%
22	18.171	2529	2533	2538	rVB5	4081	6939	1.02%	0.130%
23	18.559	2595	2599	2607	rBV2	24693	50493	7.42%	0.950%
24	18.630	2607	2611	2616	rVB	22120	31752	4.67%	0.597%
25	18.712	2618	2625	2629	rVB	7621	12786	1.88%	0.240%
26	18.776	2629	2636	2640	rBV2	22197	47034	6.91%	0.884%
27	18.812	2640	2642	2645	rVB	12907	12020	1.77%	0.226%
28	19.064	2680	2685	2689	rBV	17639	25122	3.69%	0.472%
29	19.105	2689	2692	2700	rVB3	11420	20886	3.07%	0.393%
30	19.305	2719	2726	2731	rBV4	10807	22652	3.33%	0.426%
31	19.352	2731	2734	2738	rBV2	8383	12250	1.80%	0.230%
32	19.399	2738	2742	2750	rVB5	6633	11247	1.65%	0.212%
33	19.476	2750	2755	2761	rBV	19119	35530	5.22%	0.668%
34	19.570	2769	2771	2775	rVB3	8254	8728	1.28%	0.164%
35	19.622	2775	2780	2785	rBV4	15470	29563	4.34%	0.556%
36	19.746	2795	2801	2807	rBV	237086	329622	48.43%	6.199%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121922\
Data File : BG056023.D
Acq On : 19 Dec 2022 21:33
Operator : CG/JU
Sample : N6026-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
B-29-SB01

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-BG121322.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	19.805	2807	2811	2812	rBV4	8849	12324	1.81%	0.232%
38	19.981	2839	2841	2845	rBV	5973	7526	1.11%	0.142%
39	20.069	2851	2856	2858	rBV2	39032	61108	8.98%	1.149%
40	20.104	2858	2862	2869	rVV	231224	341207	50.13%	6.416%
41	20.169	2869	2873	2877	rVB	16243	23006	3.38%	0.433%
42	20.286	2886	2893	2898	rBV	498416	680593	100.00%	12.799%
43	20.427	2910	2917	2921	rBV4	20679	54513	8.01%	1.025%
44	20.680	2958	2960	2964	rVB	14604	17670	2.60%	0.332%
45	20.886	2993	2995	2999	rBV2	14165	16400	2.41%	0.308%
46	21.373	3075	3078	3083	rVB2	20659	31267	4.59%	0.588%
47	21.608	3115	3118	3124	rVB2	42690	72287	10.62%	1.359%
48	22.008	3179	3186	3193	rBV3	159919	476147	69.96%	8.954%
49	22.073	3193	3197	3204	rVB	108868	194829	28.63%	3.664%
50	24.411	3587	3595	3603	rBV	78606	270746	39.78%	5.091%
51	25.363	3751	3757	3767	rVB	57695	160394	23.57%	3.016%
52	25.527	3777	3785	3794	rBV2	64620	204982	30.12%	3.855%

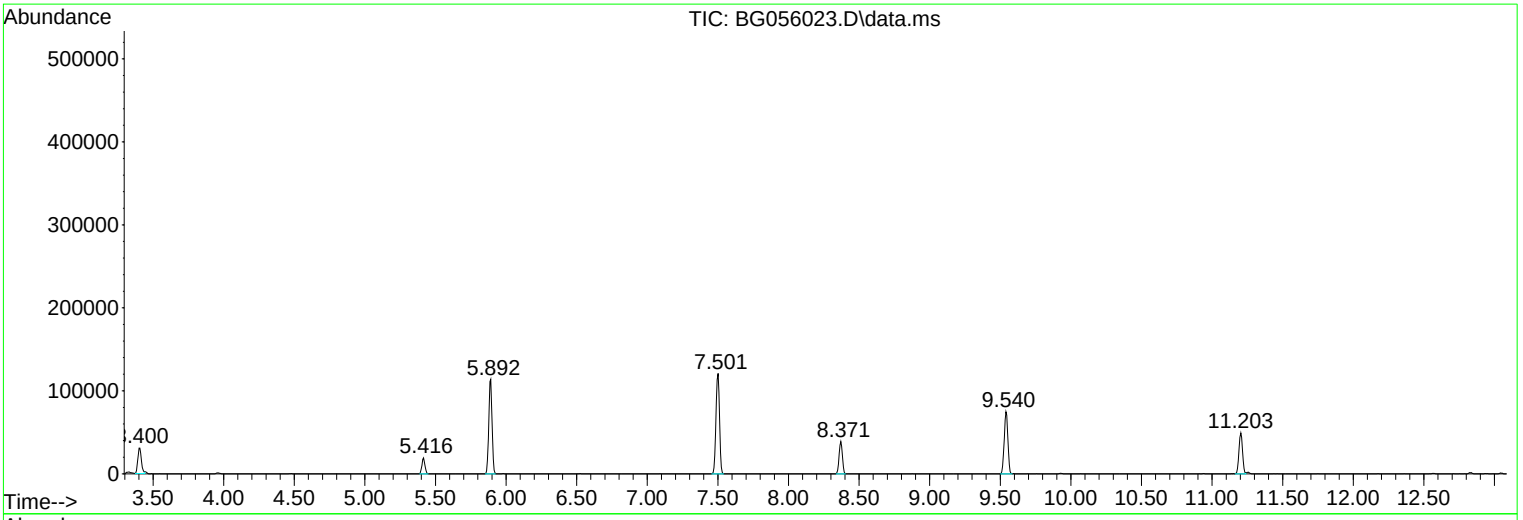
Sum of corrected areas: 5317656

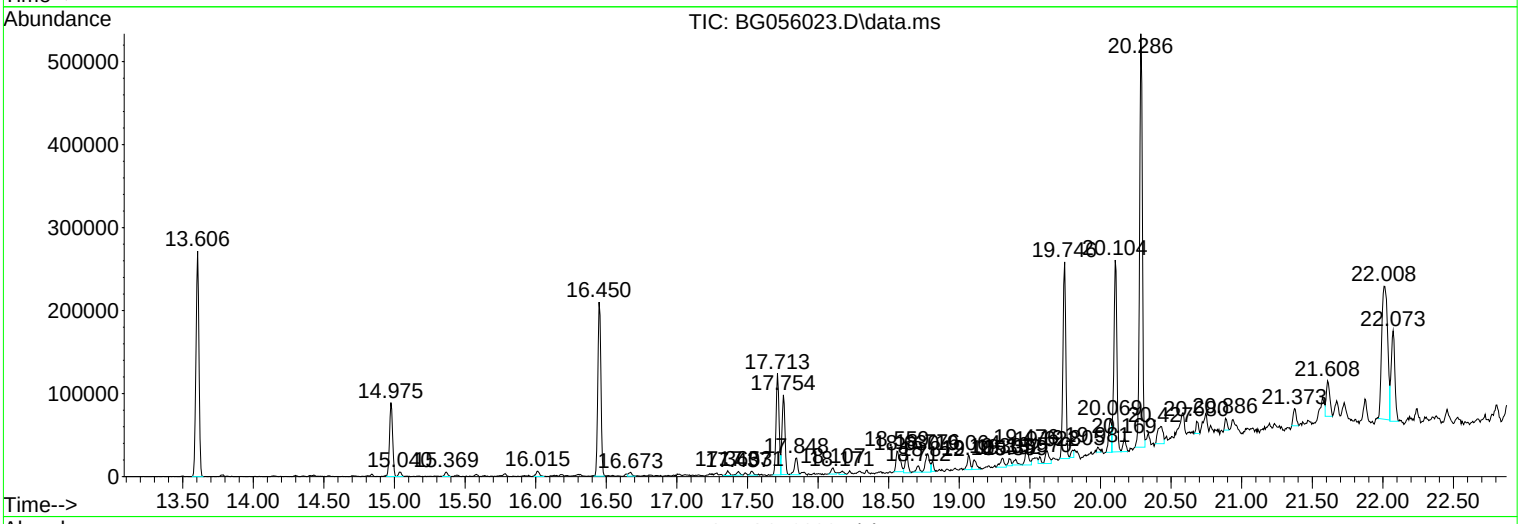
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121922\
Data File : BG056023.D
Acq On : 19 Dec 2022 21:33
Operator : CG/JU
Sample : N6026-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

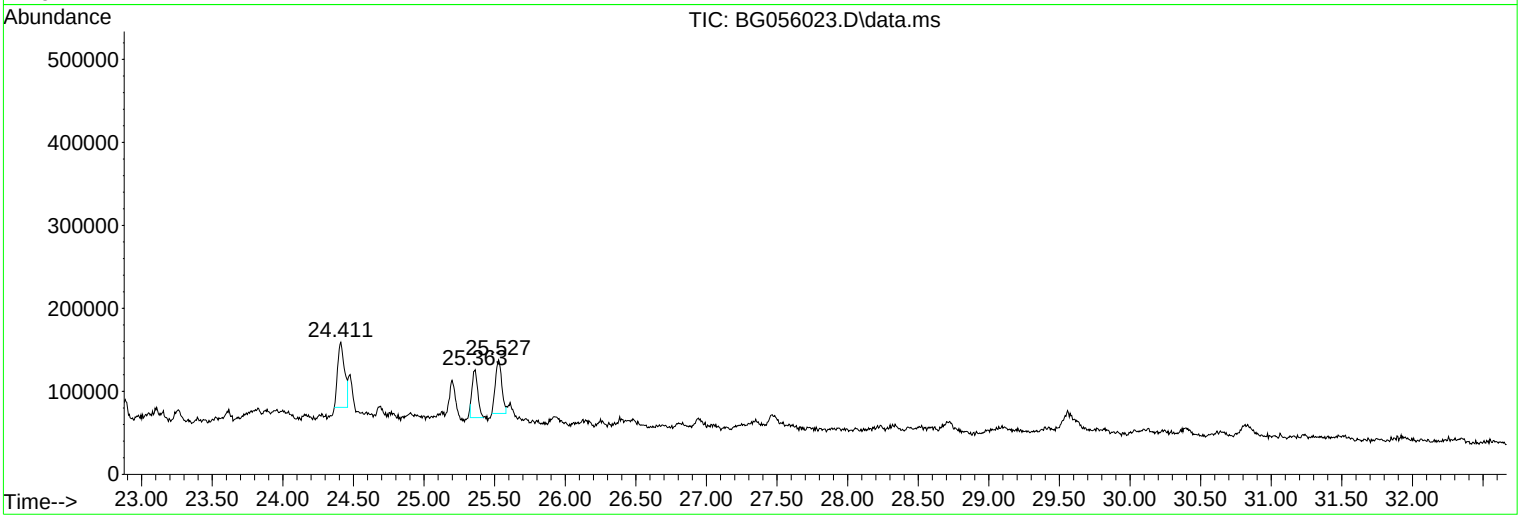
Instrument :
BNA_G
ClientSampleId :
B-29-SB01

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG121322.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\BG121922\
Data File : BG056023.D
Acq On : 19 Dec 2022 21:33
Operator : CG/JU
Sample : N6026-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
B-29-SB01

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG121322.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

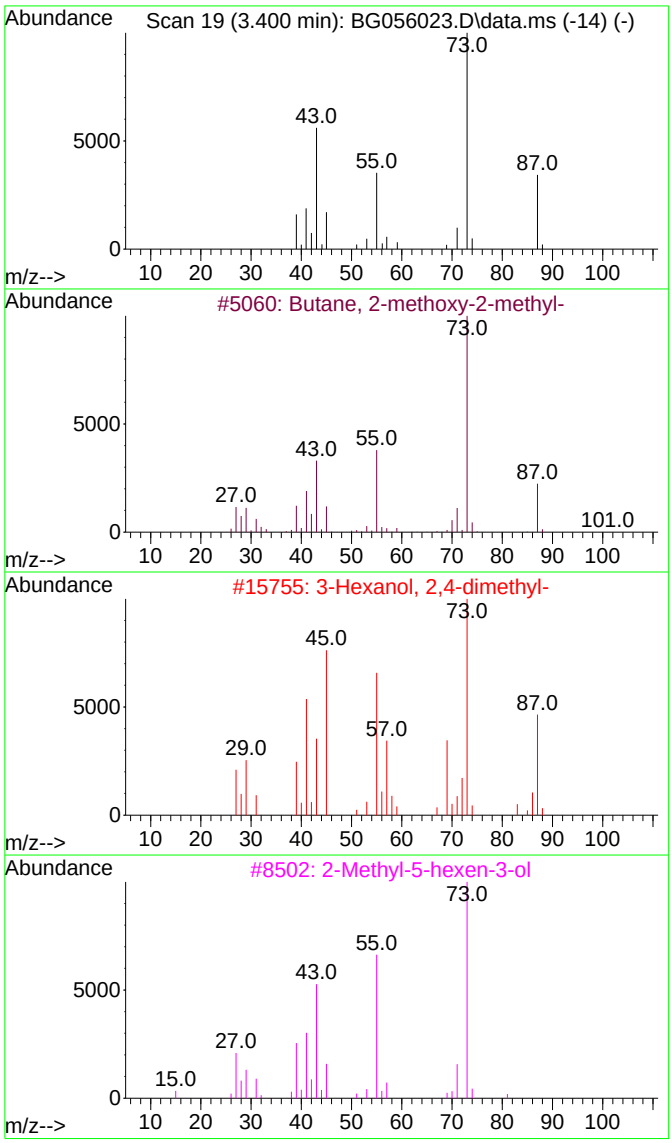
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

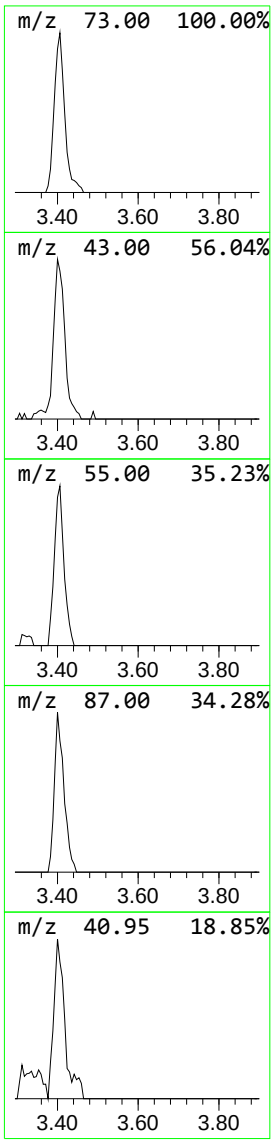
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.400	17.30 ng	54506	1,4-Dichlorobenzene-d4	8.371

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2	3-Hexanol, 2,4-dimethyl-	130	C8H18O	013432-25-2	56
3	2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	43
4	3-Pentanol, 3-methyl-	102	C6H14O	000077-74-7	33
5	Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121922\
Data File : BG056023.D
Acq On : 19 Dec 2022 21:33
Operator : CG/JU
Sample : N6026-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
B-29-SB01

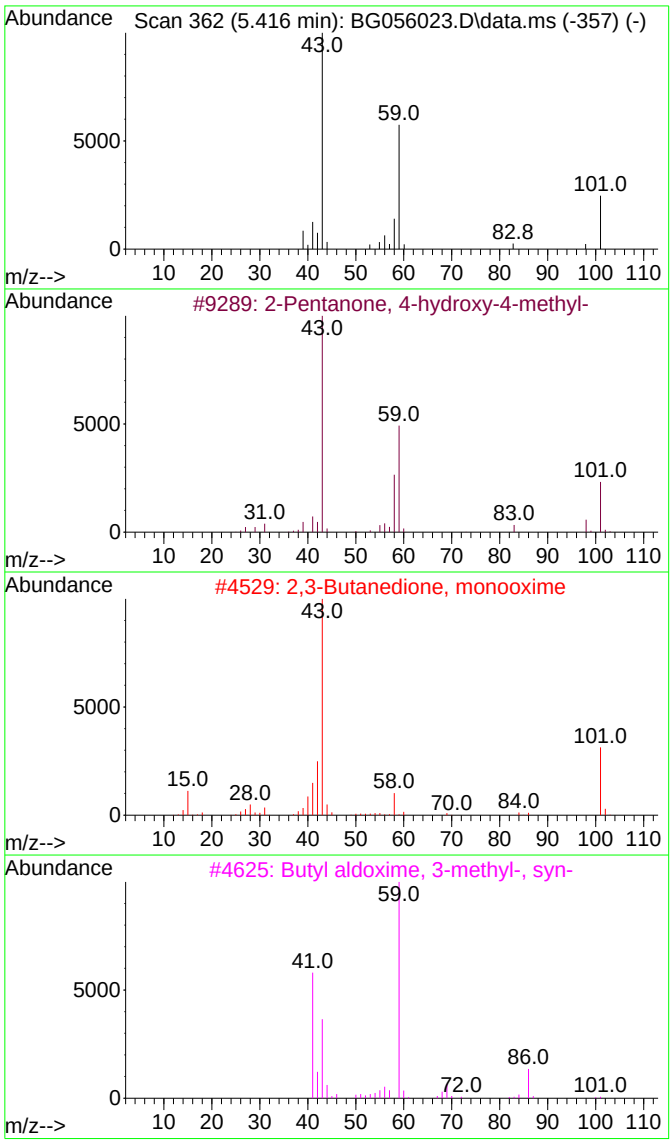
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG121322.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

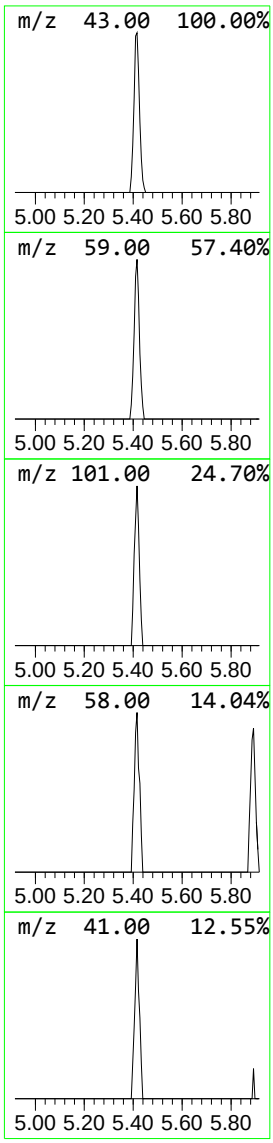
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.416	8.88 ng	27967	1,4-Dichlorobenzene-d4	8.371

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	17
3	Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	17
4	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
5	Propylamine	59	C3H9N	000107-10-8	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121922\
Data File : BG056023.D
Acq On : 19 Dec 2022 21:33
Operator : CG/JU
Sample : N6026-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
B-29-SB01

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG121322.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

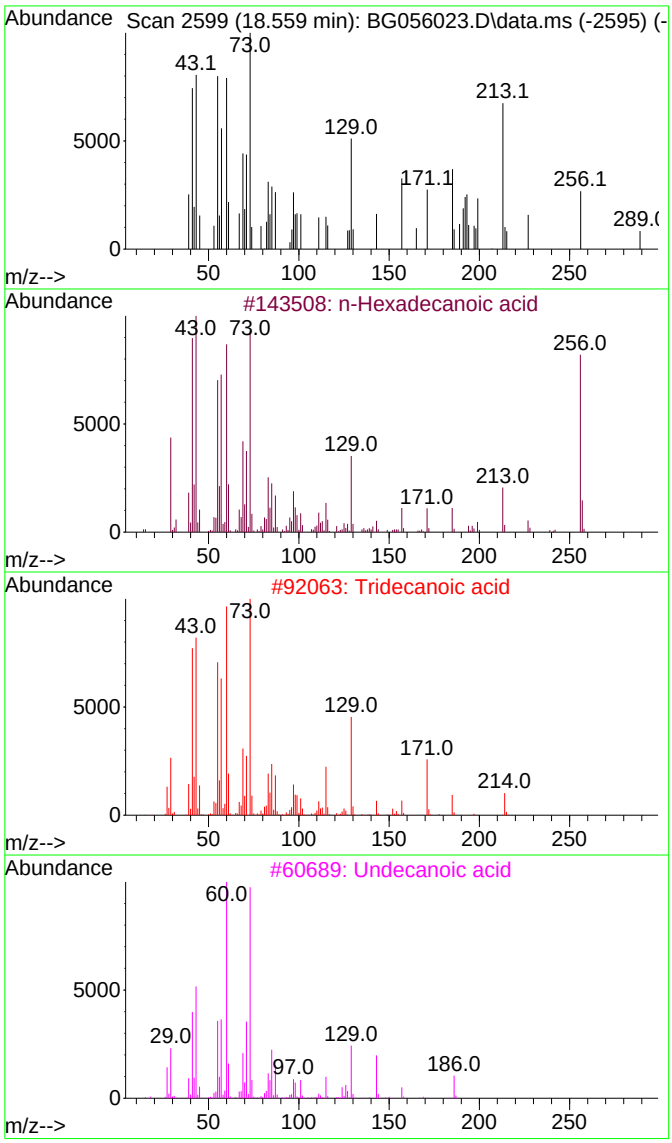
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

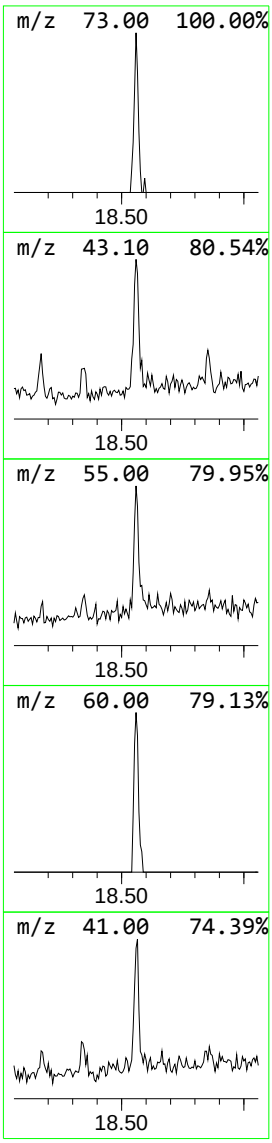
****

Peak Number 3 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.559	5.52 ng	50493	Phenanthrene-d10	17.713

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91
2	Tridecanoic acid	214	C13H26O2	000638-53-9	46
3	Undecanoic acid	186	C11H22O2	000112-37-8	46
4	Pentadecanoic acid	242	C15H30O2	001002-84-2	41
5	Octadecanoic acid	284	C18H36O2	000057-11-4	38





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121922\
Data File : BG056023.D
Acq On : 19 Dec 2022 21:33
Operator : CG/JU
Sample : N6026-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
B-29-SB01

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG121322.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

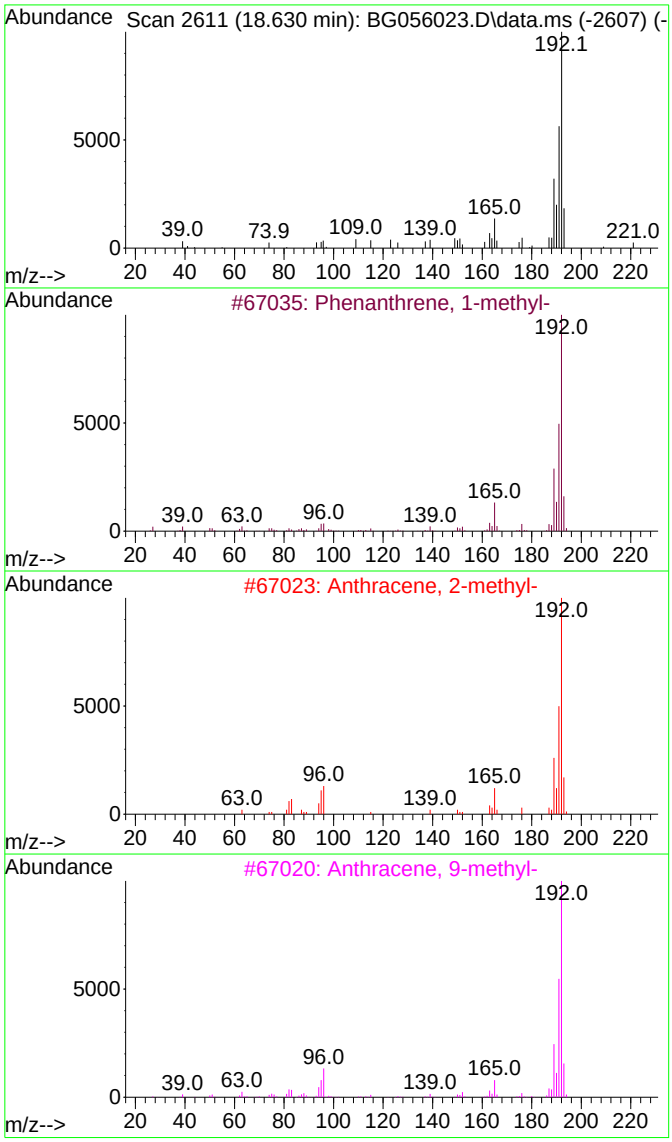
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

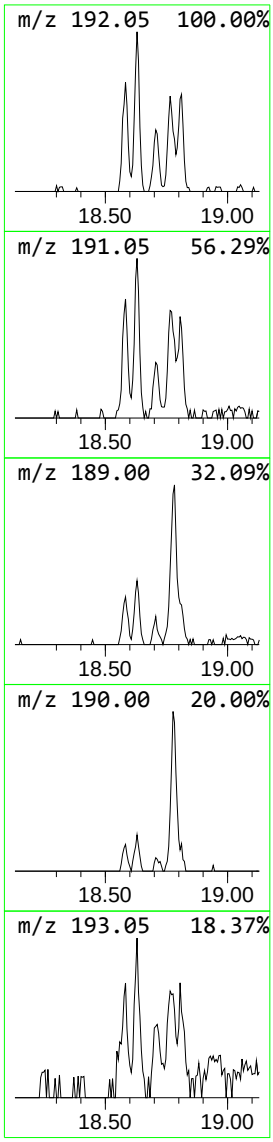
****

Peak Number 4 Phenanthrene, 1-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.630	3.47 ng	31752	Phenanthrene-d10	17.713

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
2	Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
3	Anthracene, 9-methyl-	192	C15H12	000779-02-2	91
4	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	90
5	Phenanthrene, 3-methyl-	192	C15H12	000832-71-3	89





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121922\
Data File : BG056023.D
Acq On : 19 Dec 2022 21:33
Operator : CG/JU
Sample : N6026-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
B-29-SB01

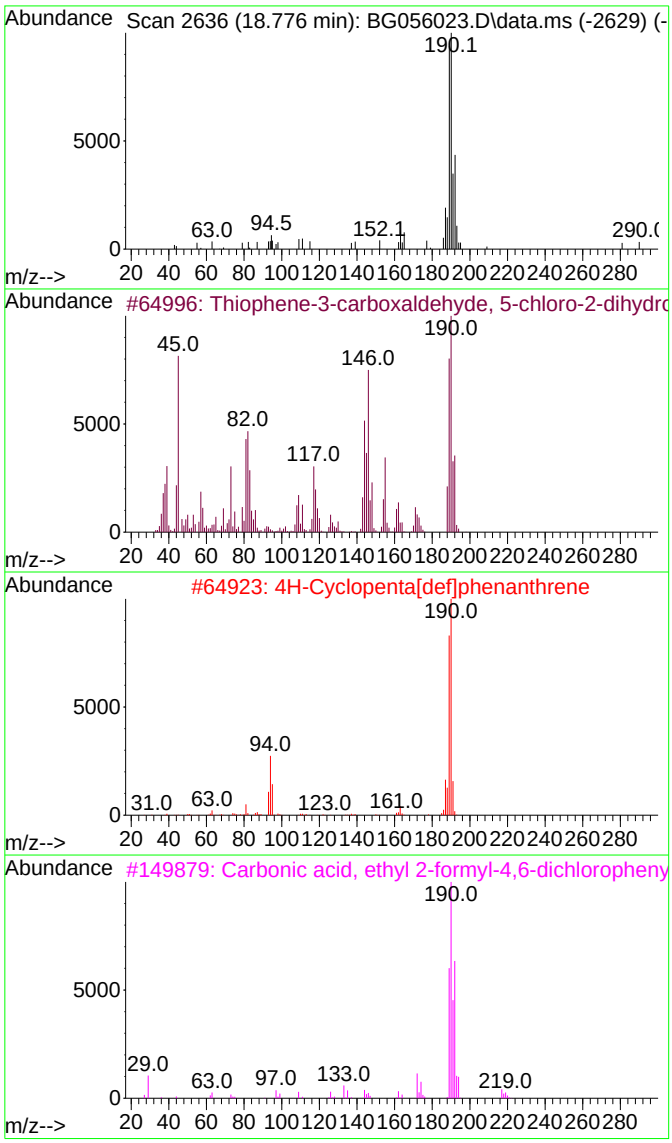
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG121322.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

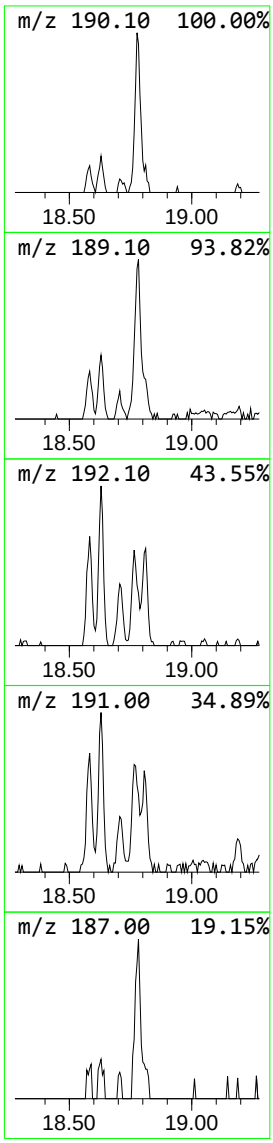
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Thiophene-3-carboxaldehyde,... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.776	5.15 ng	47034	Phenanthrene-d10	17.713

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	72
2	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	64
3	Carbonic acid, ethyl 2-formyl-4,...	262	C10H8Cl2O4	1000331-34-4	50
4	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	38
5	Carbonic acid, 2-formyl-4,6-dich...	276	C11H10Cl2O4	1000331-39-3	25





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121922\
Data File : BG056023.D
Acq On : 19 Dec 2022 21:33
Operator : CG/JU
Sample : N6026-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
B-29-SB01

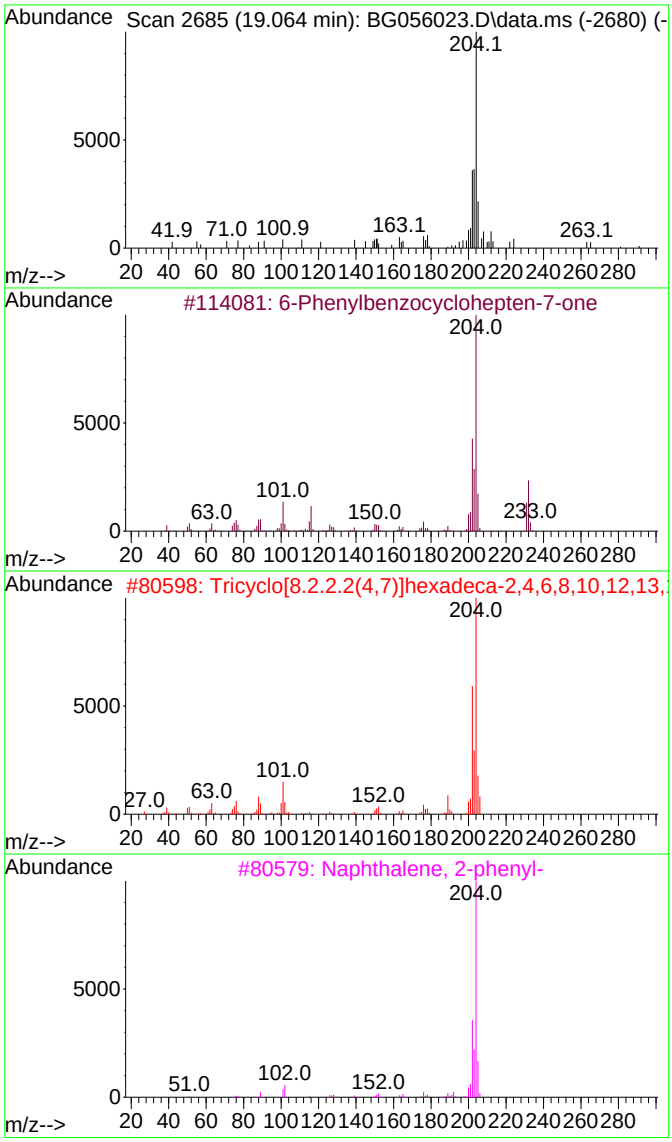
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG121322.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

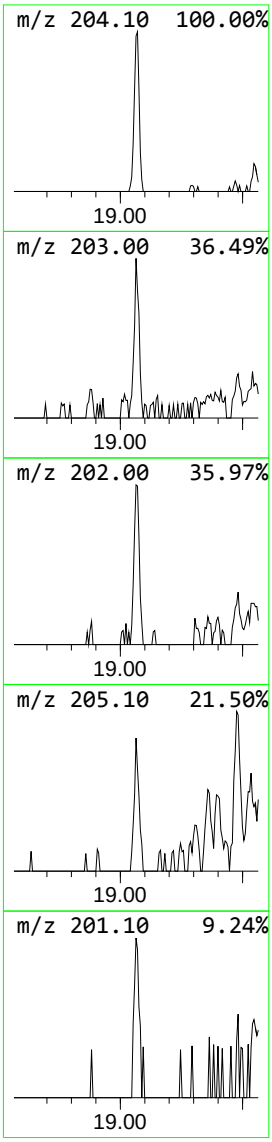
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 6-Phenylbenzocyclohepten-7-one Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.064	2.75 ng	25122	Phenanthrene-d10	17.713

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	6-Phenylbenzocyclohepten-7-one	232	C17H12O	093327-56-1	64
2	Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	64
3	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	64
4	5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	64
5	9,10-di(Chloromethyl)anthracene	274	C16H12Cl2	010387-13-0	56





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121922\
Data File : BG056023.D
Acq On : 19 Dec 2022 21:33
Operator : CG/JU
Sample : N6026-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
B-29-SB01

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG121322.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

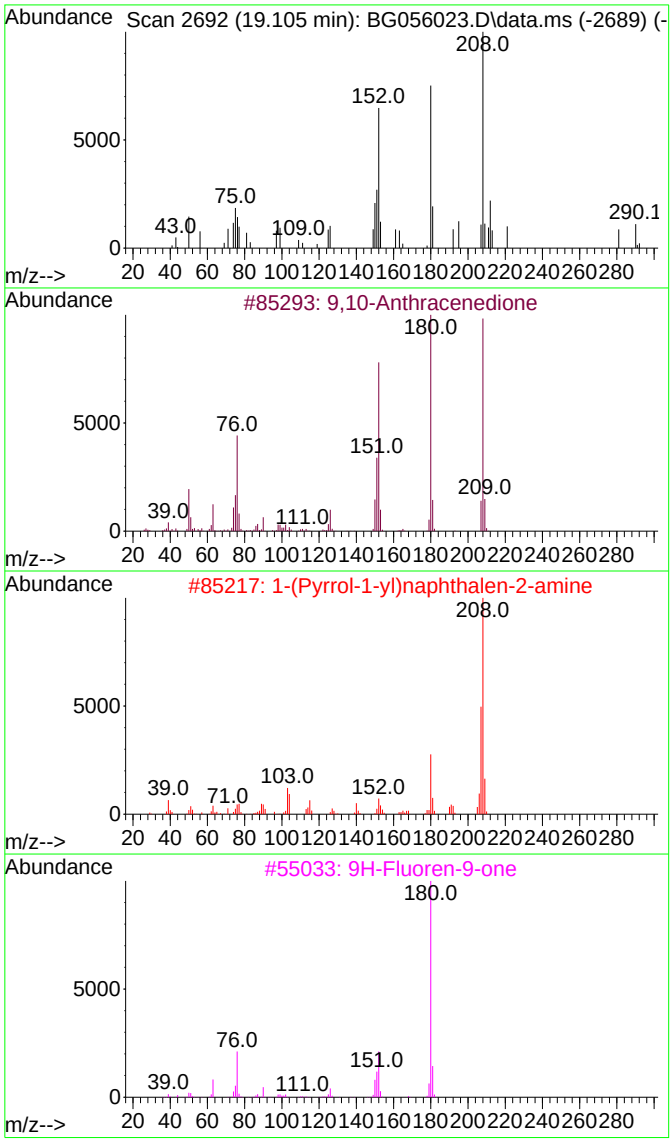
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

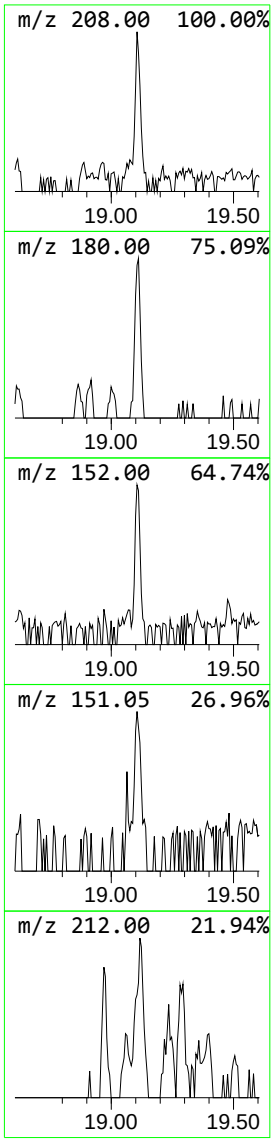
****

Peak Number 7 9,10-Anthracenedione Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.105	2.28 ng	20886	Phenanthrene-d10	17.713

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Anthracenedione	208	C14H8O2	000084-65-1	89
2	1-(Pyrrol-1-yl)naphthalen-2-amine	208	C14H12N2	1000459-83-6	52
3	9H-Fluoren-9-one	180	C13H8O	000486-25-9	49
4	15,19-Dioxatetracyclo[12.5.0.0(2...	266	C17H14O3	1000436-55-0	47
5	Benzo[c]cinnoline	180	C12H8N2	000230-17-1	43





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121922\
Data File : BG056023.D
Acq On : 19 Dec 2022 21:33
Operator : CG/JU
Sample : N6026-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
B-29-SB01

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG121322.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

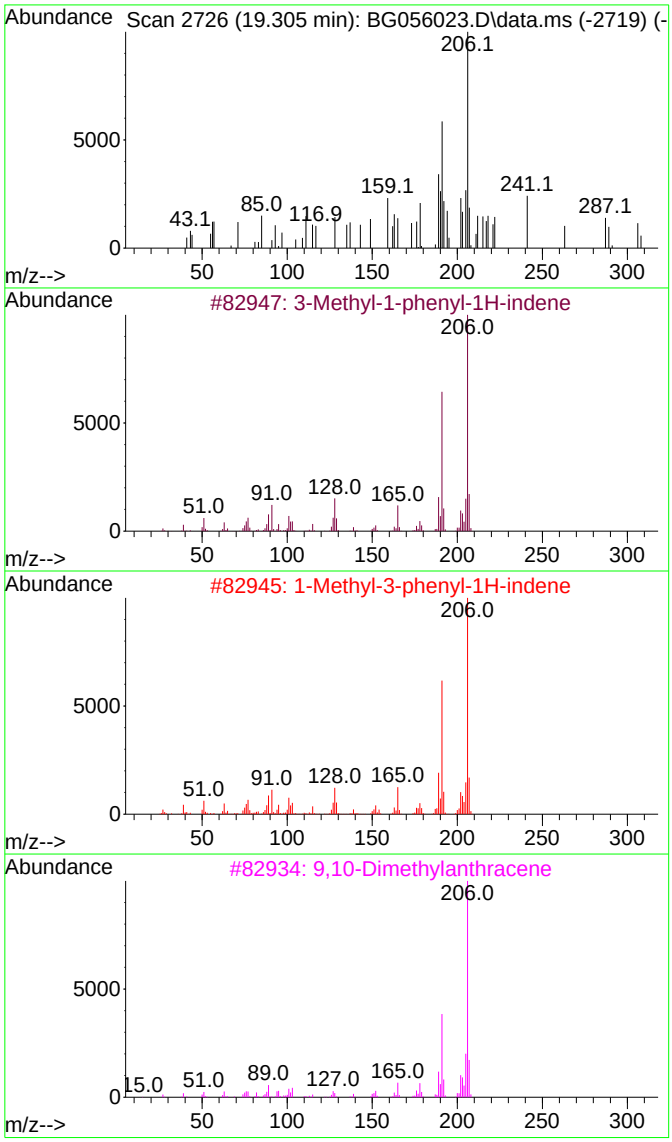
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

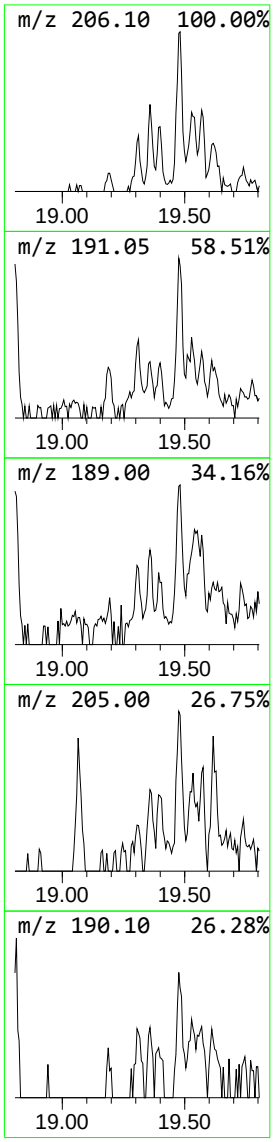
****

Peak Number 8 unknown19.305 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.305	2.48 ng	22652	Phenanthrene-d10	17.713

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Methyl-1-phenyl-1H-indene	206	C16H14	022360-63-0	49
2	1-Methyl-3-phenyl-1H-indene	206	C16H14	022360-62-9	49
3	9,10-Dimethylanthracene	206	C16H14	000781-43-1	49
4	Phenanthrene, 4,5-dimethyl-	206	C16H14	003674-69-9	49
5	1,3-Diphenyl-3-methylcyclopropene	206	C16H14	086544-79-8	49





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121922\
Data File : BG056023.D
Acq On : 19 Dec 2022 21:33
Operator : CG/JU
Sample : N6026-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
B-29-SB01

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG121322.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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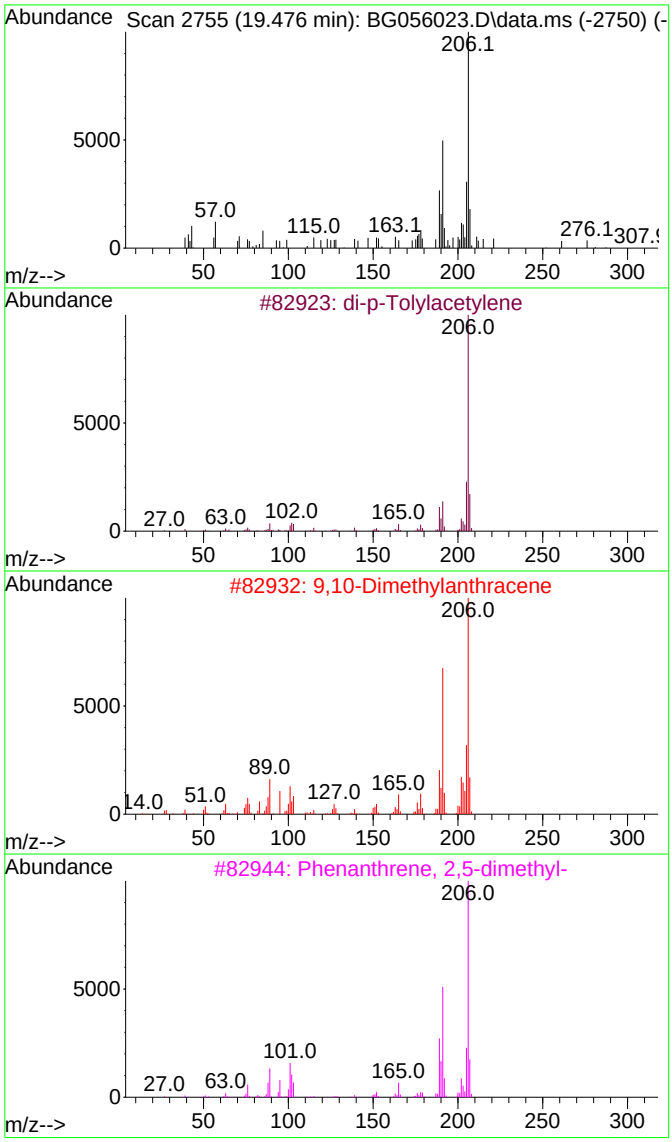
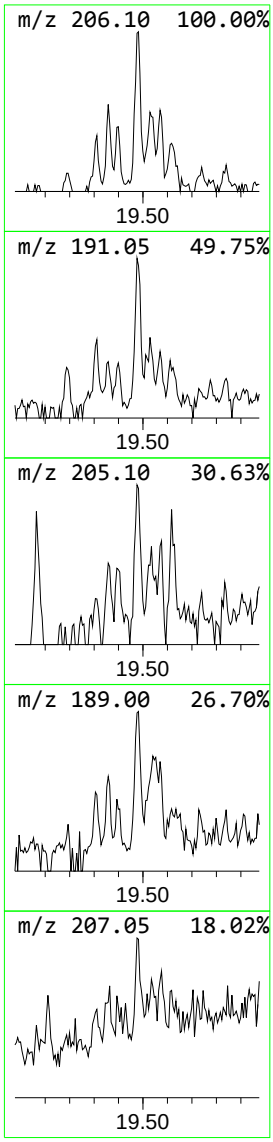
Peak Number 9 di-p-Tolylacetylene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.476	3.89 ng	35530	Phenanthrene-d10	17.713

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	di-p-Tolylacetylene	206	C16H14	002789-88-0	91
2	9,10-Dimethylanthracene	206	C16H14	000781-43-1	90
3	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	87
4	[14]Annulene, 1,6:8,13-bis(metha...	206	C16H14	085385-68-8	83
5	Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	76





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121922\
Data File : BG056023.D
Acq On : 19 Dec 2022 21:33
Operator : CG/JU
Sample : N6026-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
B-29-SB01

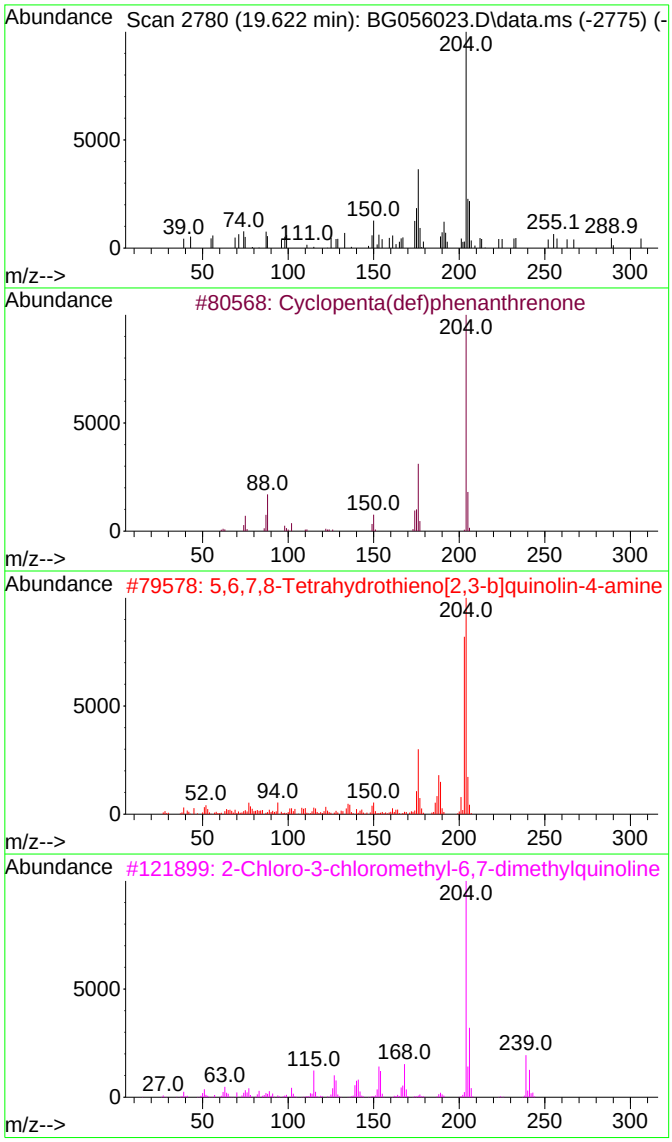
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG121322.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

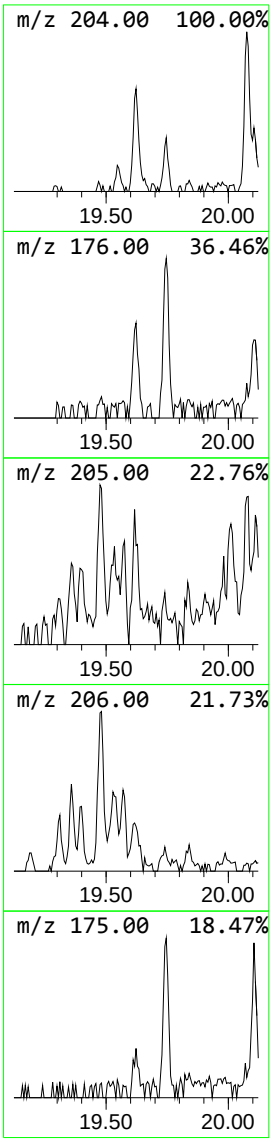
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.622	3.23 ng	29563	Phenanthrene-d10	17.713

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	62
2	5,6,7,8-Tetrahydrothieno[2,3-b]q...	204	C11H12N2S	122914-50-5	53
3	2-Chloro-3-chloromethyl-6,7-dime...	239	C12H11Cl2N	182052-67-1	50
4	Tetracyano-p-quinodimethane	204	C12H4N4	001518-16-7	47
5	Pyrene, 4,5-dihydro-	204	C16H12	006628-98-4	47





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121922\
Data File : BG056023.D
Acq On : 19 Dec 2022 21:33
Operator : CG/JU
Sample : N6026-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
B-29-SB01

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG121322.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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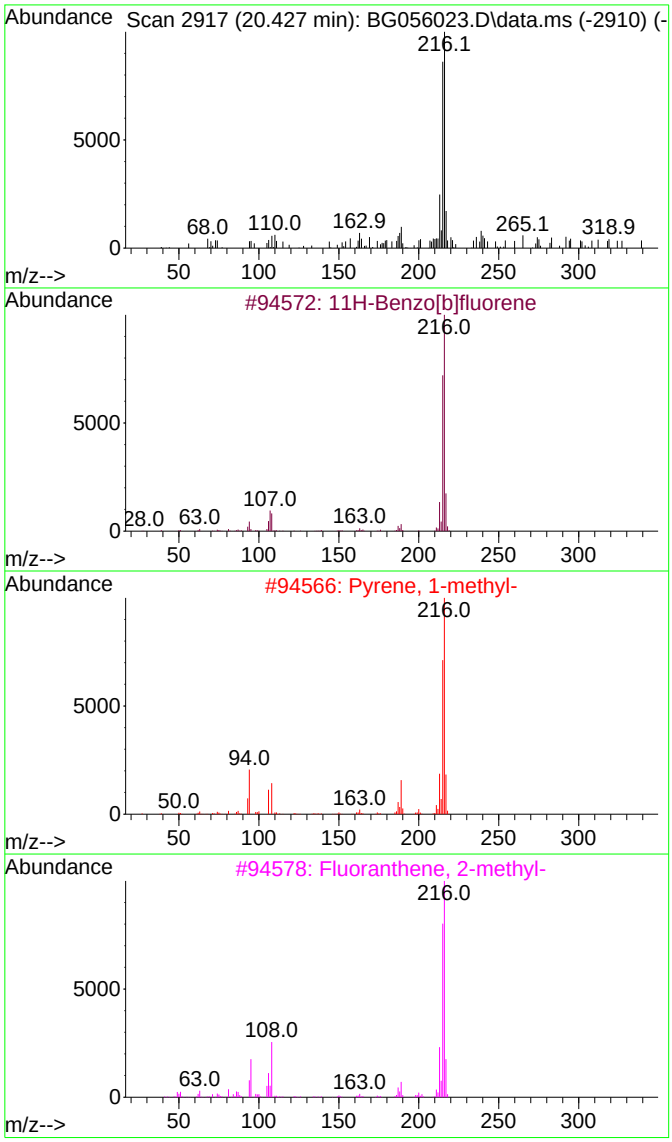
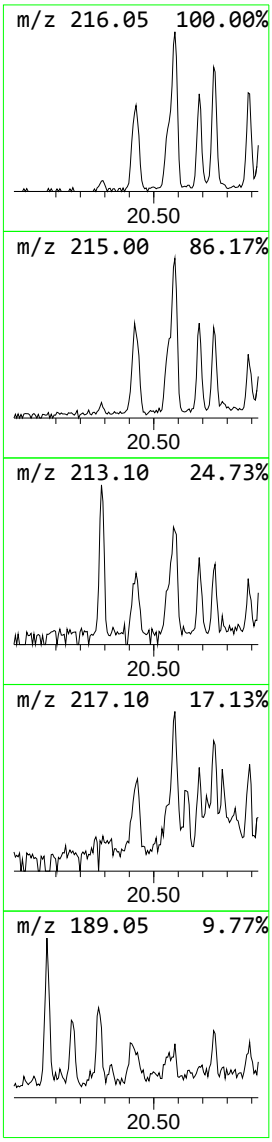
Peak Number 11 11H-Benzo[b]fluorene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.427	2.29 ng	54513	Chrysene-d12	22.026

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	90
2	Pyrene, 1-methyl-	216	C17H12	002381-21-7	76
3	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	76
4	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	76
5	7H-Benzo[c]fluorene	216	C17H12	000205-12-9	59





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121922\
Data File : BG056023.D
Acq On : 19 Dec 2022 21:33
Operator : CG/JU
Sample : N6026-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
B-29-SB01

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG121322.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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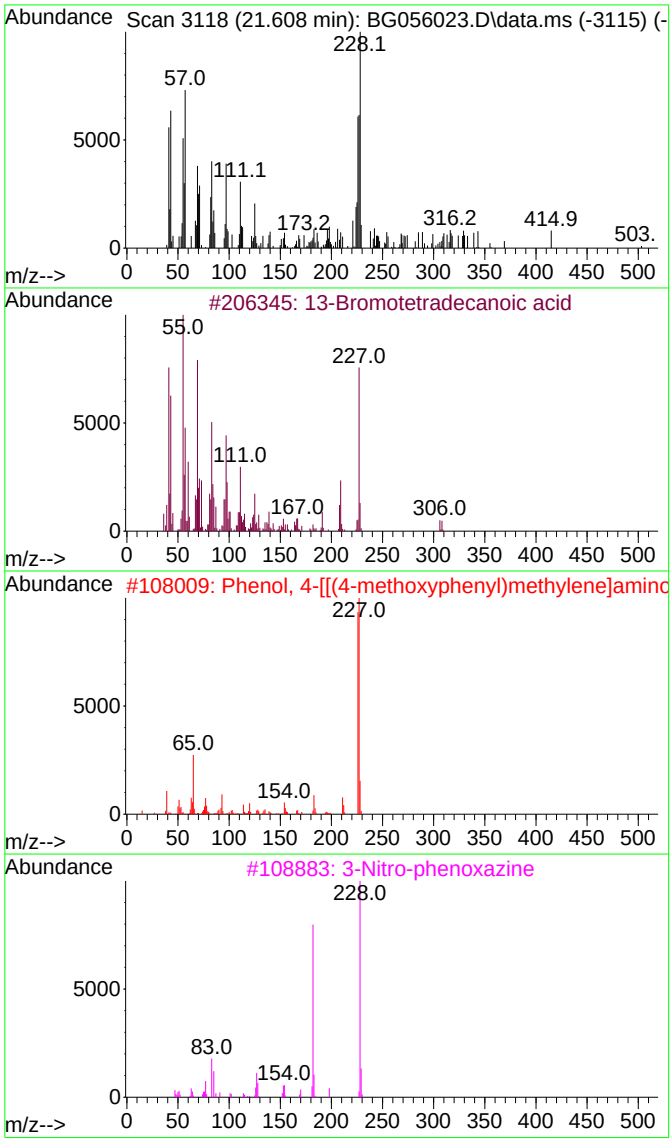
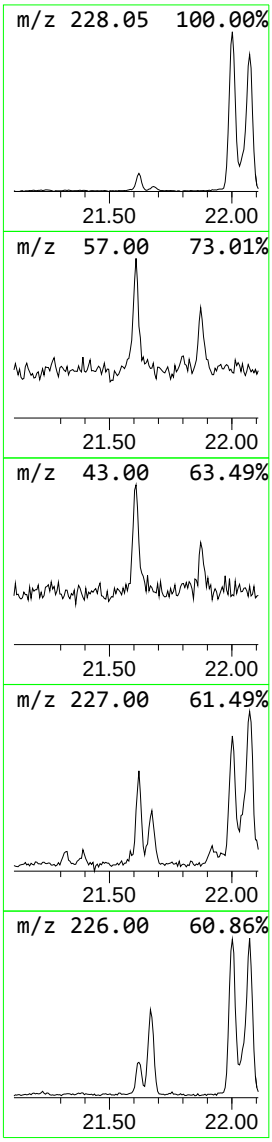
Peak Number 12 unknown21.608 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.608	3.04 ng	72287	Chrysene-d12	22.026

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	13-Bromotetradecanoic acid	306	C14H27BrO2	1000099-12-8	42
2	Phenol, 4-[[(4-methoxyphenyl)met...	227	C14H13NO2	003230-39-5	18
3	3-Nitro-phenoxazine	228	C12H8N2O3	020464-44-2	18
4	3H-Spiro[1,3-benzothiazole-2,3'-...	254	C14H10N2OS	120105-59-1	18
5	Ferrocene, acetyl-	228	C12H12FeO	001271-55-2	18





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121922\
Data File : BG056023.D
Acq On : 19 Dec 2022 21:33
Operator : CG/JU
Sample : N6026-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
B-29-SB01

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG121322.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...	3.400	17.3	ng	54506	1	8.371	63002	20.0
2-Pentanone, 4-...	5.416	8.9	ng	27967	1	8.371	63002	20.0
n-Hexadecanoic ...	18.559	5.5	ng	50493	4	17.713	182814	20.0
Phenanthrene, 1...	18.630	3.5	ng	31752	4	17.713	182814	20.0
Thiophene-3-car...	18.776	5.2	ng	47034	4	17.713	182814	20.0
6-Phenylbenzocy...	19.064	2.8	ng	25122	4	17.713	182814	20.0
9,10-Anthracene...	19.105	2.3	ng	20886	4	17.713	182814	20.0
unknown19.305	19.305	2.5	ng	22652	4	17.713	182814	20.0
di-p-Tolylacety...	19.476	3.9	ng	35530	4	17.713	182814	20.0
Cyclopenta(def)...	19.622	3.2	ng	29563	4	17.713	182814	20.0
11H-Benzo[b]flu...	20.427	2.3	ng	54513	5	22.026	476147	20.0
unknown21.608	21.608	3.0	ng	72287	5	22.026	476147	20.0