LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040121\
Data File : BG048557.D
Acq On : 1 Apr 2021 23:51
Operator : CG/JU
Sample : M1886-07
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : RO-124037

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-BG033021.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG048557.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.168	305	311	317	rVB3	8137	13148	1.06%	0.101%
2	5.644	384	392	400	rBV	415548	667249	53.63%	5.135%
3	7.248	656	665	678	rBV	415343	722978	58.11%	5.564%
4	8.094	802	809	816	rBV2	87657	143859	11.56%	1.107%
5	9.264	1001	1008	1017	rVB	261728	465760	37.43%	3.585%
6	10.920	1284	1290	1295	rBV	109316	187520	15.07%	1.443%
7	10.973	1295	1299	1307	rVB	23890	39209	3.15%	0.302%
8	12.577	1566	1572	1578	rVB2	15052	22162	1.78%	0.171%
9	12.795	1602	1609	1614	rBV3	8463	13414	1.08%	0.103%
10	13.365	1700	1706	1716	rVB	521156	808989	65.02%	6.226%
11	14.064	1820	1825	1829	rBV3	8612	15133	1.22%	0.116%
12	14.187	1840	1846	1854	rBV2	26533	40248	3.23%	0.310%
13	14.745	1930	1941	1947	rBV	164427	263012	21.14%	2.024%
14	14.810	1947	1952	1959	rVB	139918	210512	16.92%	1.620%
15	15.145	2002	2009	2016	rBV	62426	91831	7.38%	0.707%
16	15.797	2113	2120	2127	rBV	104786	163530	13.14%	1.259%
17	15.956	2144	2147	2153	rVB3	12240	16735	1.35%	0.129%
18	16.085	2166	2169	2177	rVB2	19098	27633	2.22%	0.213%
19	16.238	2187	2195	2204	rBV	601005	920816	74.01%	7.087%
20	16.802	2288	2291	2296	rVV3	13551	19961	1.60%	0.154%
21	17.154	2346	2351	2357	rVB4	13586	22816	1.83%	0.176%
22	17.248	2365	2367	2372	rVB4	13173	15136	1.22%	0.116%
23	17.313	2373	2378	2385	rVB	31042	52080	4.19%	0.401%
24	17.501	2403	2410	2413	rBV	224676	342994	27.57%	2.640%
25	17.542	2413	2417	2423	rVV	506676	749613	60.25%	5.769%
26	17.630	2423	2432	2437	rVV	63557	115919	9.32%	0.892%
27	17.689	2437	2442	2447	rVB3	13030	22203	1.78%	0.171%
28	17.901	2469	2478	2482	rBV3	27955	68382	5.50%	0.526%
29	18.018	2494	2498	2504	rBV	14964	19800	1.59%	0.152%
30	18.382	2555	2560	2564	rBV2	118164	172365	13.85%	1.327%
31	18.423	2564	2567	2574	rVV	58419	91820	7.38%	0.707%
32	18.500	2575	2580	2586	rVB2	17386	30947	2.49%	0.238%
33	18.576	2586	2593	2603	rVB3	97134	193731	15.57%	1.491%
34	18.870	2638	2643	2647	rBV	35846	53246	4.28%	0.410%
35	18.911	2647	2650	2655	rVB	29286	38213	3.07%	0.294%
36	19.117	2677	2685	2688	rBV5	10215	19324	1.55%	0.149%

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Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : RO-124037

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-BG033021.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	19.287	2708	2714	2718	rBV3	12742	23306	1.87%	0.179%
38	19.352	2719	2725	2732	rVB9	11668	23569	1.89%	0.181%
39	19.428	2732	2738	2747	rVB	32736	54612	4.39%	0.420%
40	19.557	2753	2760	2765	rVB	724872	1014360	81.53%	7.807%
41	19.616	2765	2770	2772	rBV5	11730	14929	1.20%	0.115%
42	19.651	2772	2776	2785	rVB4	35701	56477	4.54%	0.435%
43	19.740	2785	2791	2794	rBV8	10807	17877	1.44%	0.138%
44	19.792	2794	2800	2804	rVB3	14827	28574	2.30%	0.220%
45	19.881	2809	2815	2817	rVV2	107959	167372	13.45%	1.288%
46	19.916	2817	2821	2828	rVV	532324	741382	59.59%	5.706%
47	19.980	2828	2832	2838	rVB2	22383	34578	2.78%	0.266%
48	20.110	2846	2854	2859	rBV	938003	1244196	100.00%	9.576%
49	20.151	2859	2861	2865	rVB	32013	37601	3.02%	0.289%
50	20.227	2869	2874	2881	rBV5	25459	64646	5.20%	0.498%
51	20.403	2892	2904	2909	rBV2	65879	143326	11.52%	1.103%
52	20.503	2916	2921	2925	rBV2	50752	69439	5.58%	0.534%
53	20.562	2925	2931	2935	rVV2	33917	64418	5.18%	0.496%
54	20.597	2935	2937	2941	rVV	23323	26582	2.14%	0.205%
55	20.638	2941	2944	2949	rVB6	9938	15065	1.21%	0.116%
56	20.703	2951	2955	2958	rBV3	21261	34085	2.74%	0.262%
57	20.744	2958	2962	2967	rVB3	21392	33506	2.69%	0.258%
58	21.173	3031	3035	3040	rVB2	25713	40922	3.29%	0.315%
59	21.361	3061	3067	3072	rBV2	27359	60387	4.85%	0.465%
60	21.408	3072	3075	3080	rVV3	35603	70753	5.69%	0.545%
61	21.455	3080	3083	3088	rVV3	35705	64963	5.22%	0.500%
62	21.514	3088	3093	3098	rVB2	22997	39640	3.19%	0.305%
63	21.737	3127	3131	3132	rBV3	9986	12528	1.01%	0.096%
64	21.796	3132	3141	3146	rVV2	270024	698011	56.10%	5.372%
65	21.843	3146	3149	3157	rVB	115579	188541	15.15%	1.451%
66	21.996	3170	3175	3185	rVB2	25620	53421	4.29%	0.411%
67	22.207	3207	3211	3217	rVB4	12694	21330	1.71%	0.164%
68	22.548	3266	3269	3277	rVB4	13934	25243	2.03%	0.194%
69	22.630	3277	3283	3287	rBV7	14193	31889	2.56%	0.245%
70	22.724	3295	3299	3304	rBV8	7775	14913	1.20%	0.115%
71	22.836	3312	3318	3321	rBV4	12482	21685	1.74%	0.167%
72	22.877	3323	3325	3332	rVB8	10762	21809	1.75%	0.168%
73	23.241	3381	3387	3390	rBV5	7761	17089	1.37%	0.132%
74	24.070	3519	3528	3537	rBV3	67830	227887	18.32%	1.754%
75	24.340	3570	3574	3579	rVB6	7869	13343	1.07%	0.103%
76	24.816	3650	3655	3656	rBV2	25897	33689	2.71%	0.259%

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Misc :
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Instrument : BNA_G
ClientSampleId : RO-124037

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Integrator: RTE
Smoothing : ON Filtering: 5
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Stop Thrs : 0 Peak Location: TOP

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Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG033021.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

77	24.980	3674	3683	3692	rVB2	35233	104966	8.44%	0.808%
78	25.139	3699	3710	3720	rBV3	137661	421720	33.89%	3.246%
79	28.946	4351	4358	4368	rBV6	10738	36519	2.94%	0.281%

Sum of corrected areas: 12993436

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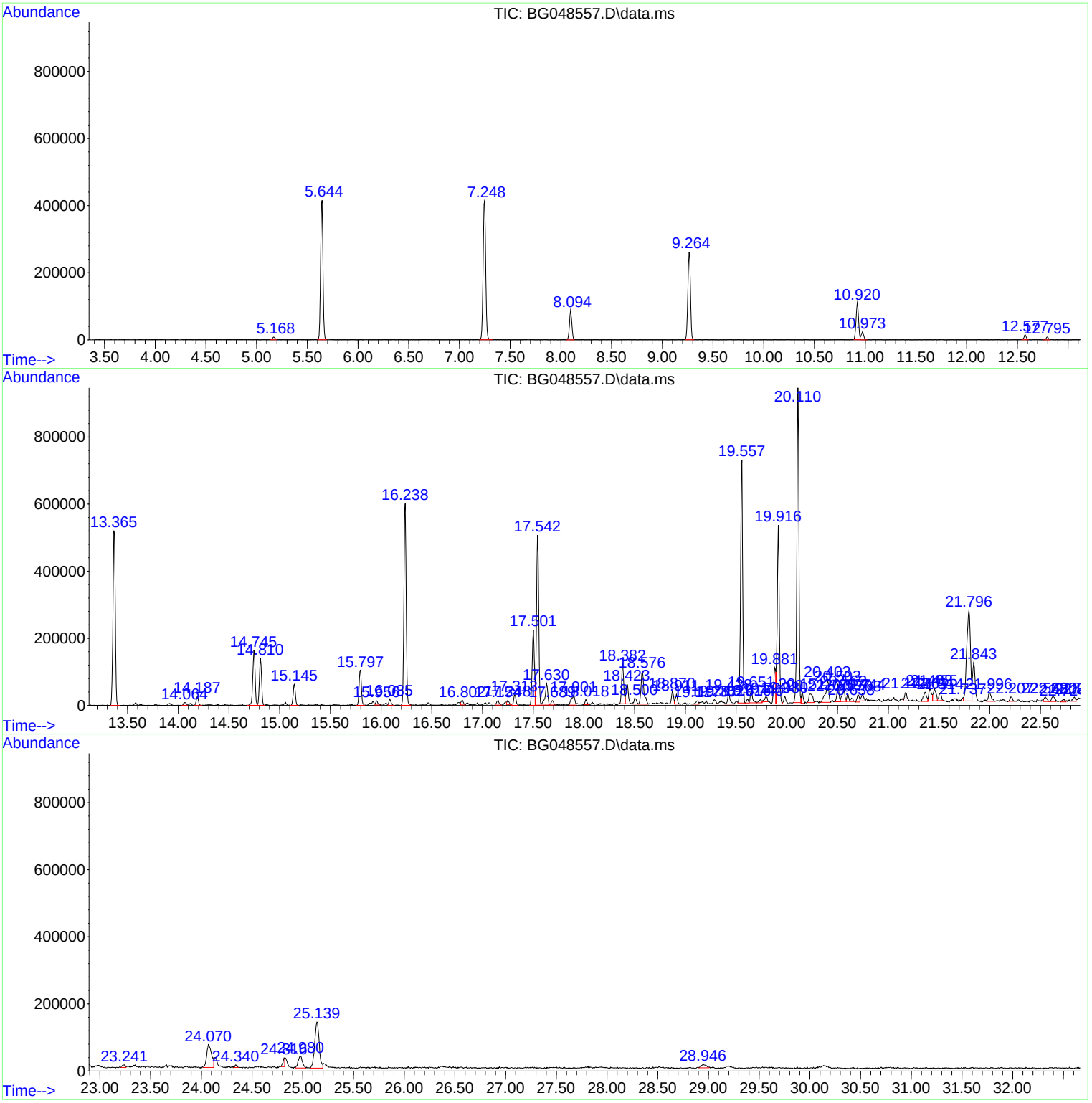
Sample : M1886-07
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
RO-124037

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG033021.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040121\
Data File : BG048557.D
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Sample : M1886-07
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : RO-124037

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG033021.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

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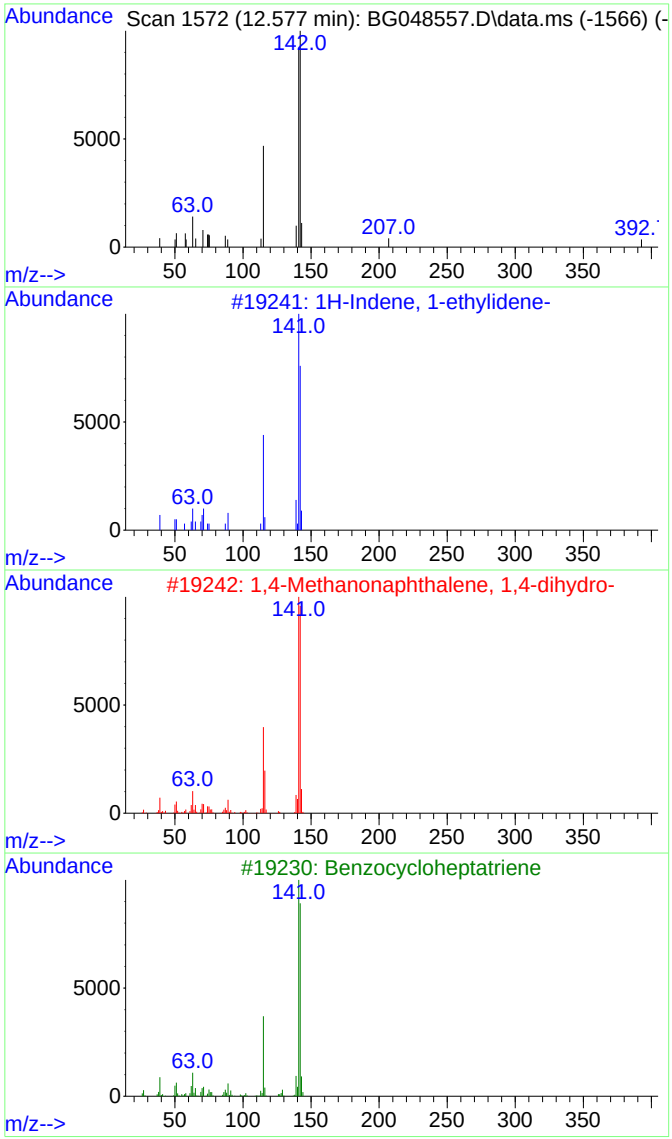
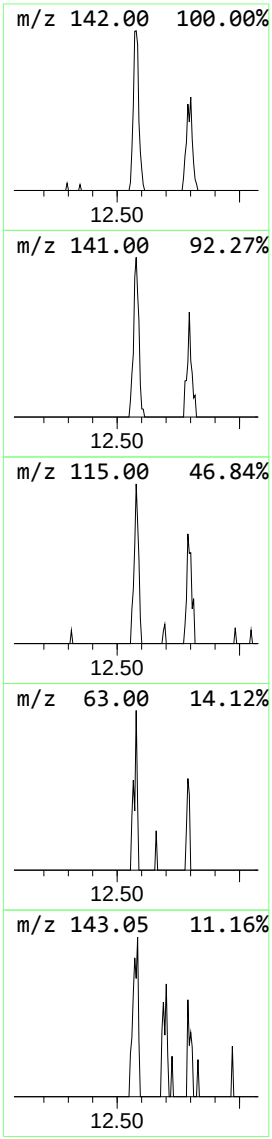
Peak Number 1 1H-Indene, 1-ethylidene- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.577	2.36 ng	22162	Naphthalene-d8	10.921

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	86
2	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	86
3	Benzocycloheptatriene	142	C11H10	000264-09-5	86
4	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	86
5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	80





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040121\
Data File : BG048557.D
Acq On : 1 Apr 2021 23:51
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Sample : M1886-07
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : RO-124037

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG033021.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

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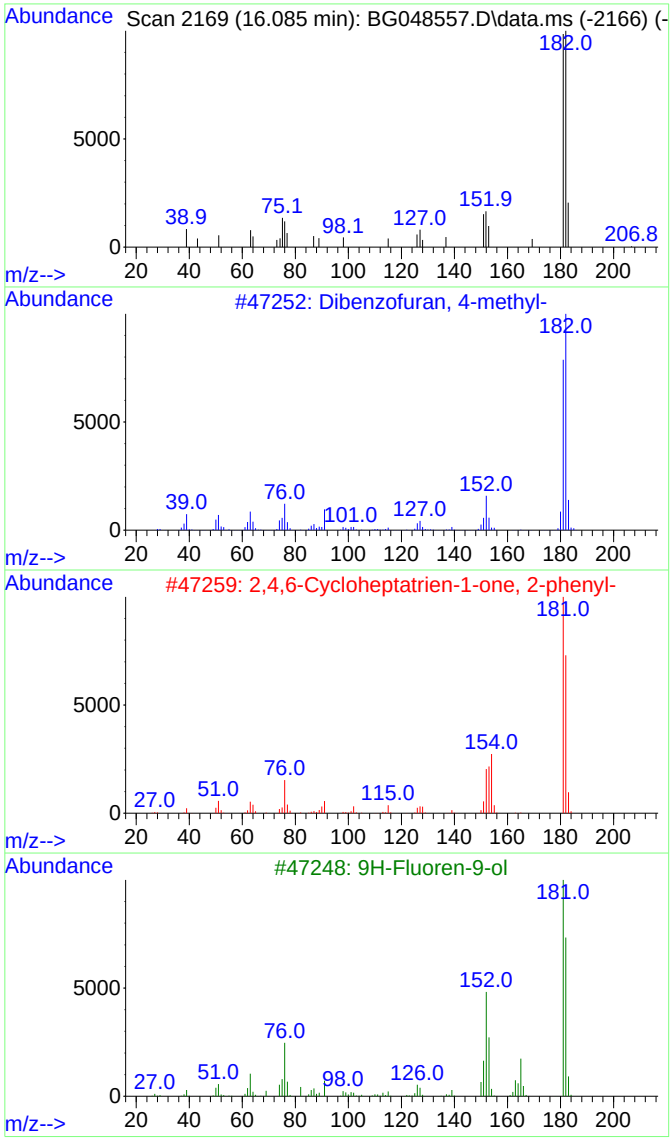
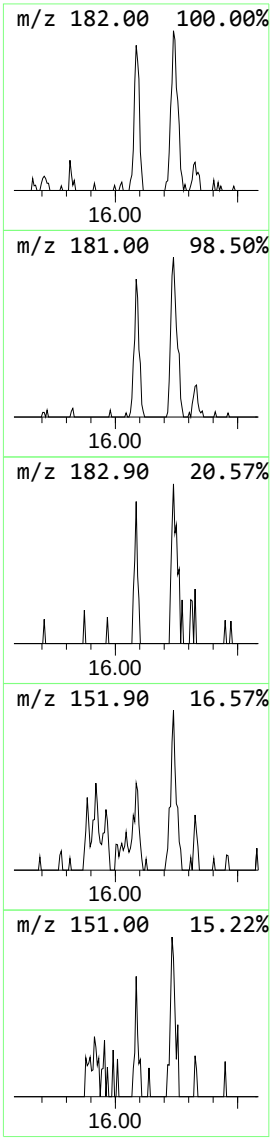
Peak Number 2 Dibenzofuran, 4-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.085	2.10 ng	27633	Acenaphthene-d10	14.745

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dibenzofuran, 4-methyl-	182	C13H10O	007320-53-8	86
2	2,4,6-Cycloheptatrien-1-one, 2-p...	182	C13H10O	014562-09-5	72
3	9H-Fluoren-9-ol	182	C13H10O	001689-64-1	72
4	Pyrido[2,3-b]indole, 6-methyl-	182	C12H10N2	108349-67-3	59
5	[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	59





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040121\
Data File : BG048557.D
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Sample : M1886-07
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : RO-124037

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG033021.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
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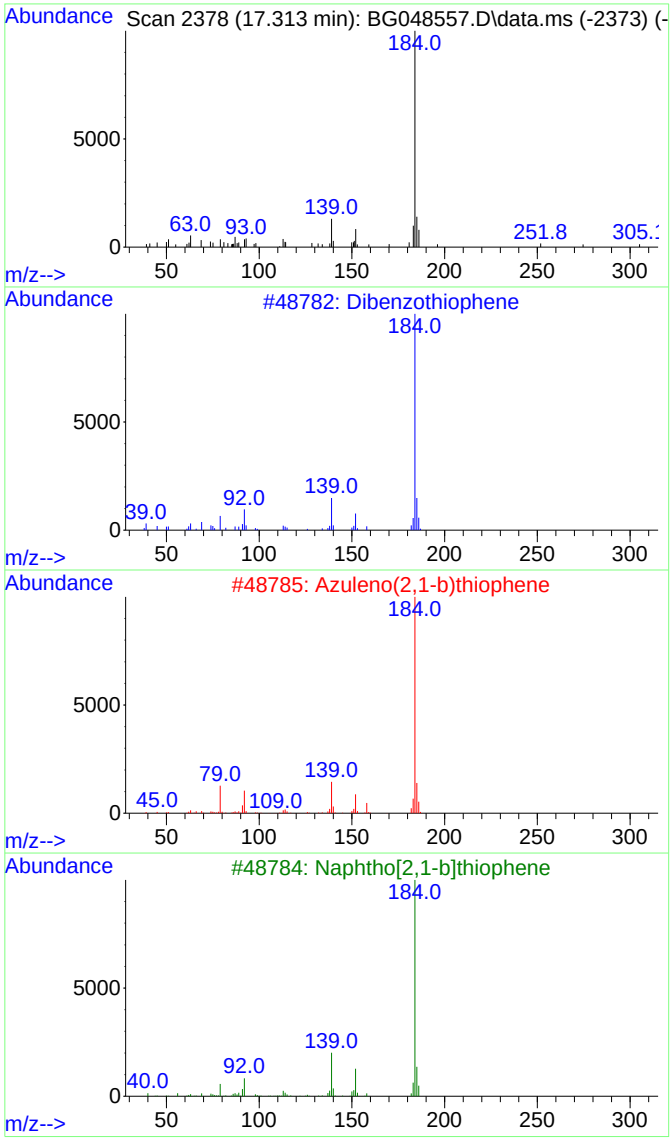
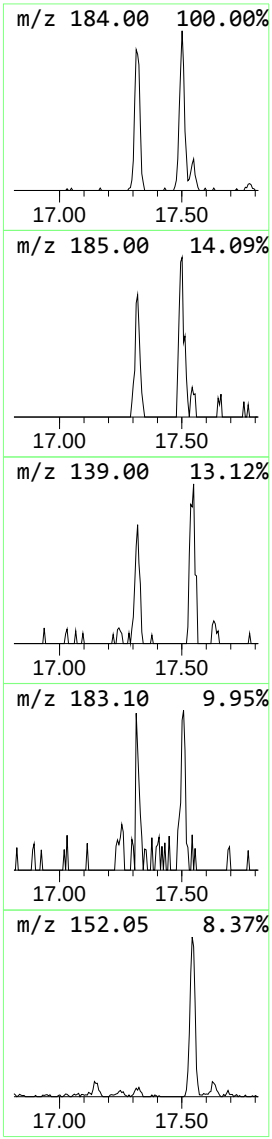
Peak Number 3 Dibenzothiophene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.313	3.04 ng	52080	Phenanthrene-d10	17.501

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dibenzothiophene	184	C12H8S	000132-65-0	90
2	Azuleno(2,1-b)thiophene	184	C12H8S	000248-13-5	86
3	Naphtho[2,1-b]thiophene	184	C12H8S	000233-02-3	83
4	Naphtho[2,3-b]thiophene	184	C12H8S	000268-77-9	74
5	Naphtho[1,2-b]thiophene	184	C12H8S	000234-41-3	72





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Sample : M1886-07
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : RO-124037

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG033021.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

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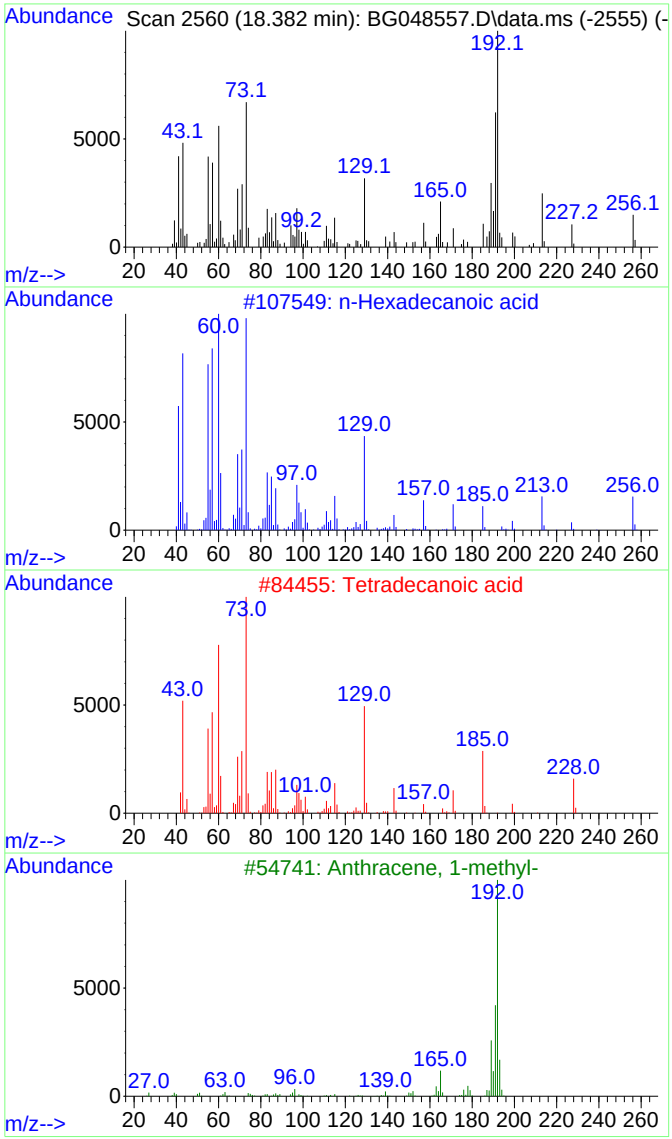
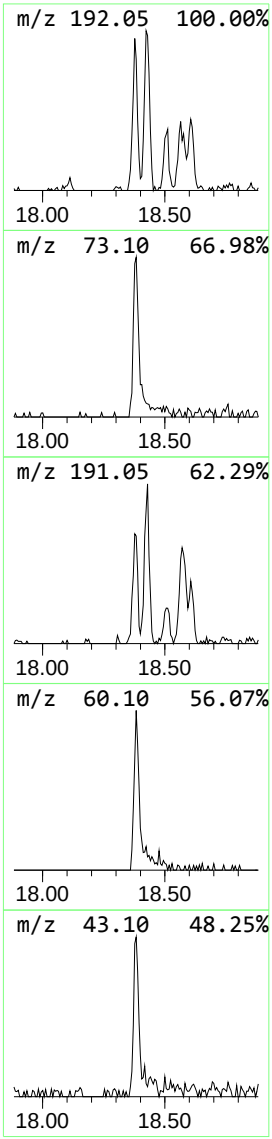
Peak Number 4 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.382	10.05 ng	172365	Phenanthrene-d10	17.501

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	90
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	58
3	Anthracene, 1-methyl-	192	C15H12	000610-48-0	55
4	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	50
5	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	50





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040121\
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Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : RO-124037

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG033021.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

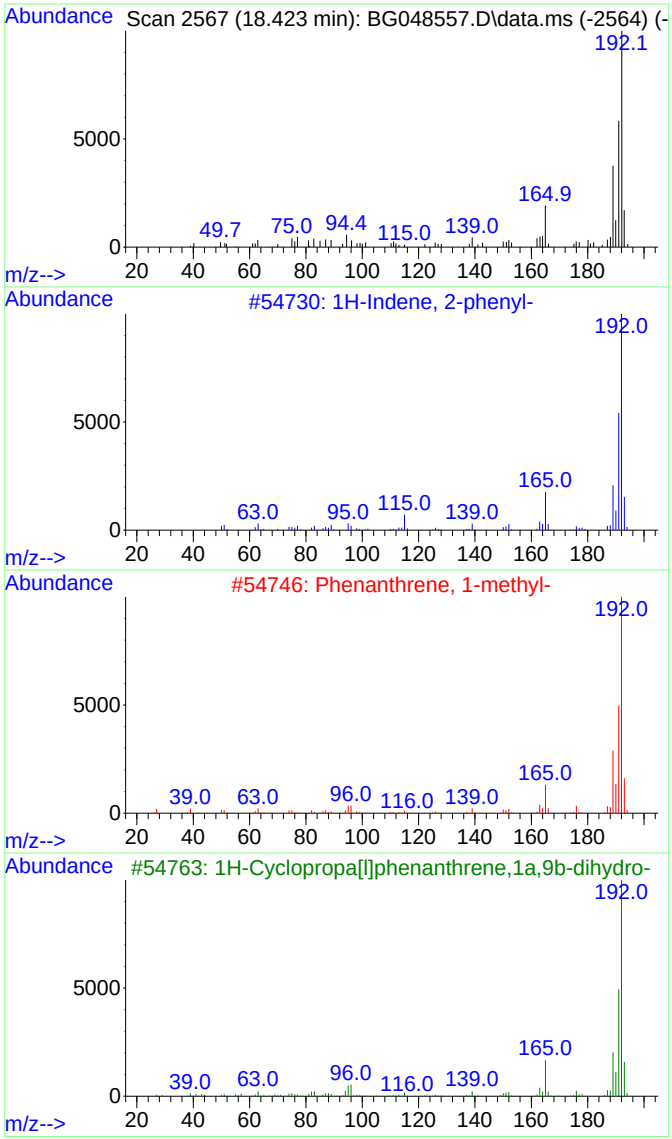
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

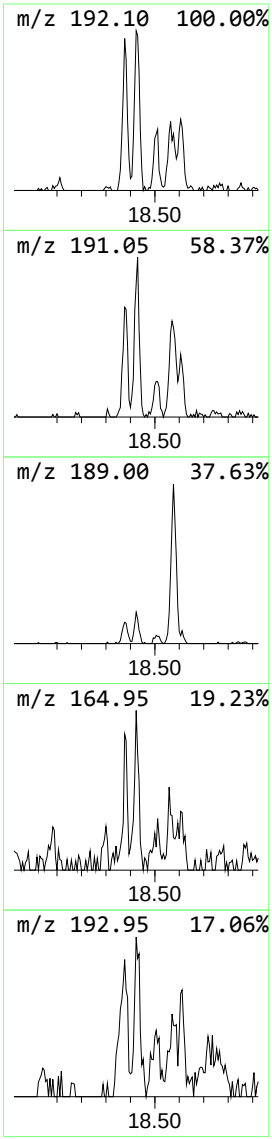
****

Peak Number 5 1H-Indene, 2-phenyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.423	5.35 ng	91820	Phenanthrene-d10	17.501

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	94
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
3	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	91
4	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	91
5	Naphtho[2,3-b]norbornadiene	192	C15H12	107426-38-0	90





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040121\
Data File : BG048557.D
Acq On : 1 Apr 2021 23:51
Operator : CG/JU
Sample : M1886-07
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : RO-124037

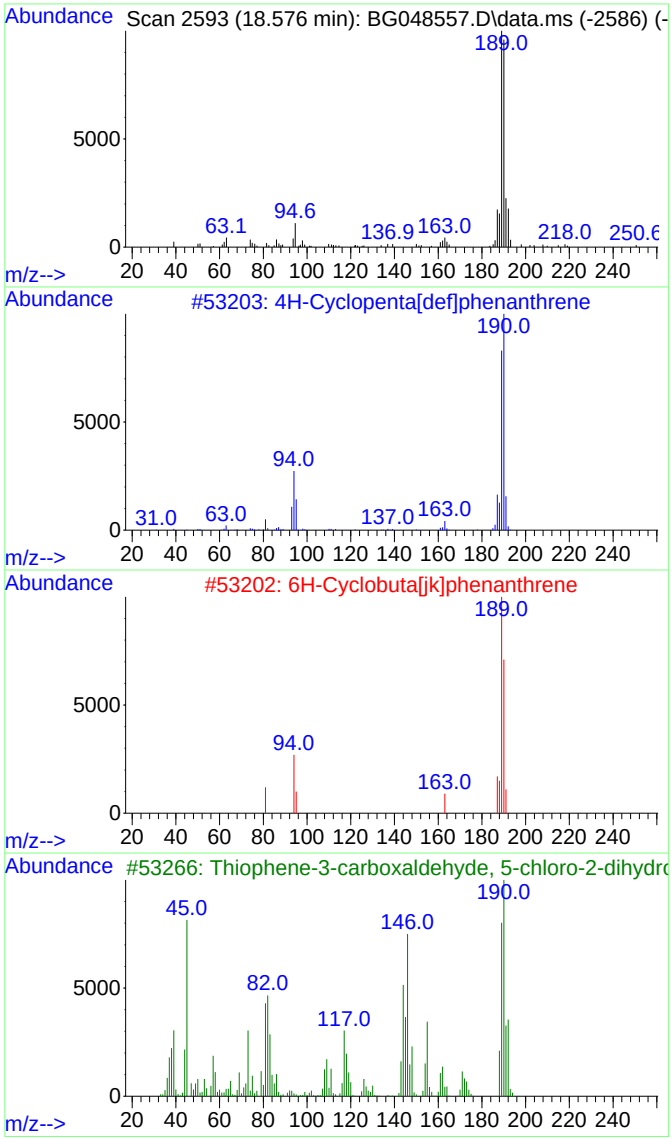
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG033021.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

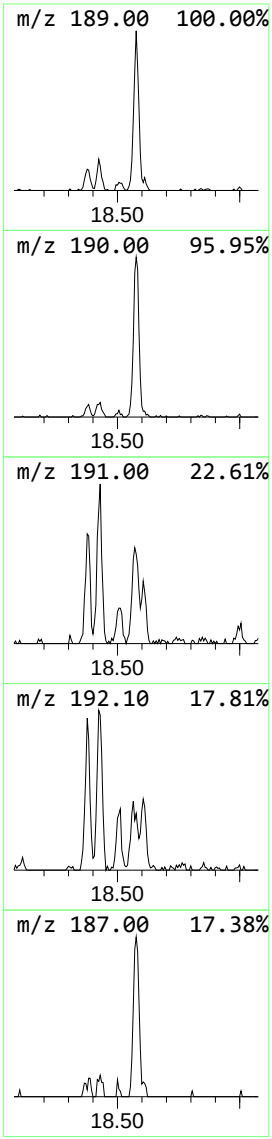
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 4H-Cyclopenta[def]phenanthrene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.576	11.30 ng	193731	Phenanthrene-d10	17.501

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	87
2	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	72
3	Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	45
4	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	45
5	1,4-Naphthalenedione, 2,5-dihydr...	190	C10H6O4	004923-55-1	32





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Acq On : 1 Apr 2021 23:51
Operator : CG/JU
Sample : M1886-07
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : RO-124037

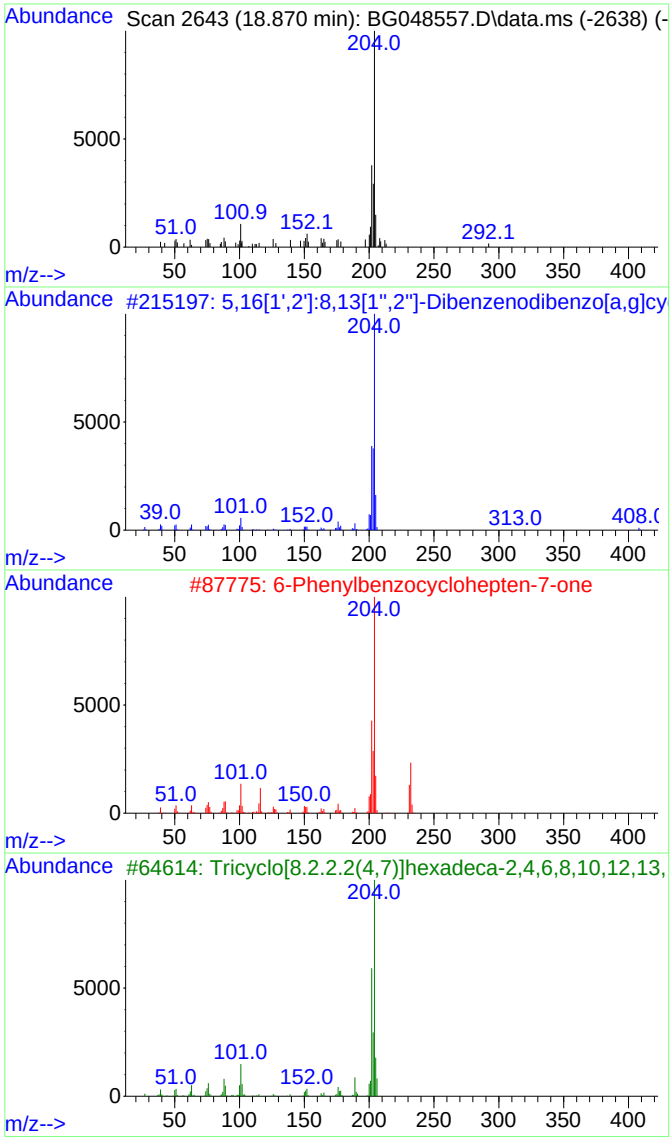
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG033021.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

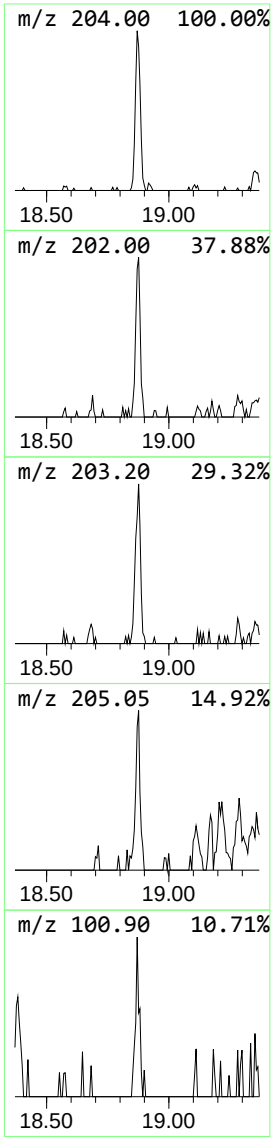
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 5,16[1',2']:8,13[1'',2'']-D... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.870	3.10 ng	53246	Phenanthrene-d10	17.501

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	86
2	6-Phenylbenzocyclohepten-7-one	232	C17H12O	093327-56-1	86
3	Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	81
4	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	68
5	Naphthalene, 1-phenyl-	204	C16H12	000605-02-7	58





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Data File : BG048557.D
Acq On : 1 Apr 2021 23:51
Operator : CG/JU
Sample : M1886-07
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : RO-124037

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG033021.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

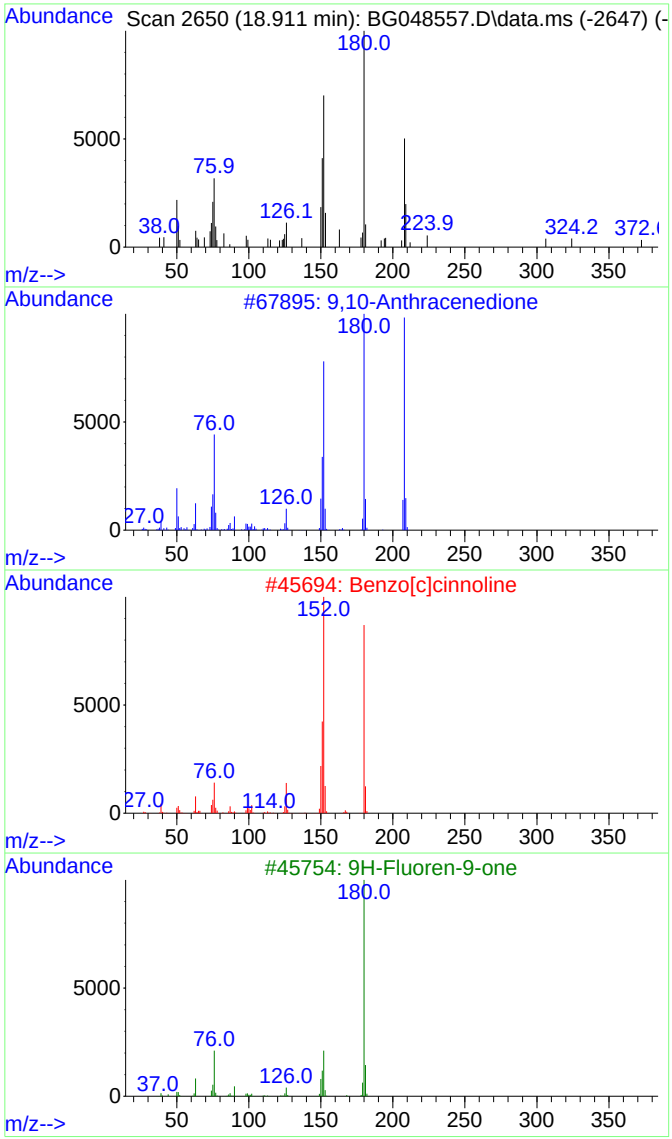
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

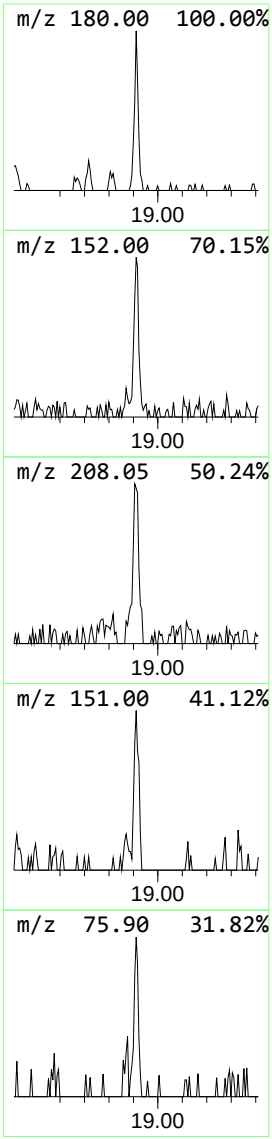
****

Peak Number 8 9,10-Anthracenedione Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.911	2.23 ng	38213	Phenanthrene-d10	17.501

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Anthracenedione	208	C14H8O2	000084-65-1	94
2	Benzo[c]cinnoline	180	C12H8N2	000230-17-1	62
3	9H-Fluoren-9-one	180	C13H8O	000486-25-9	60
4	9,10-Phenanthrenedione	208	C14H8O2	000084-11-7	55
5	Diphenic anhydride	224	C14H8O3	006050-13-1	53





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Acq On : 1 Apr 2021 23:51
Operator : CG/JU
Sample : M1886-07
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : RO-124037

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG033021.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

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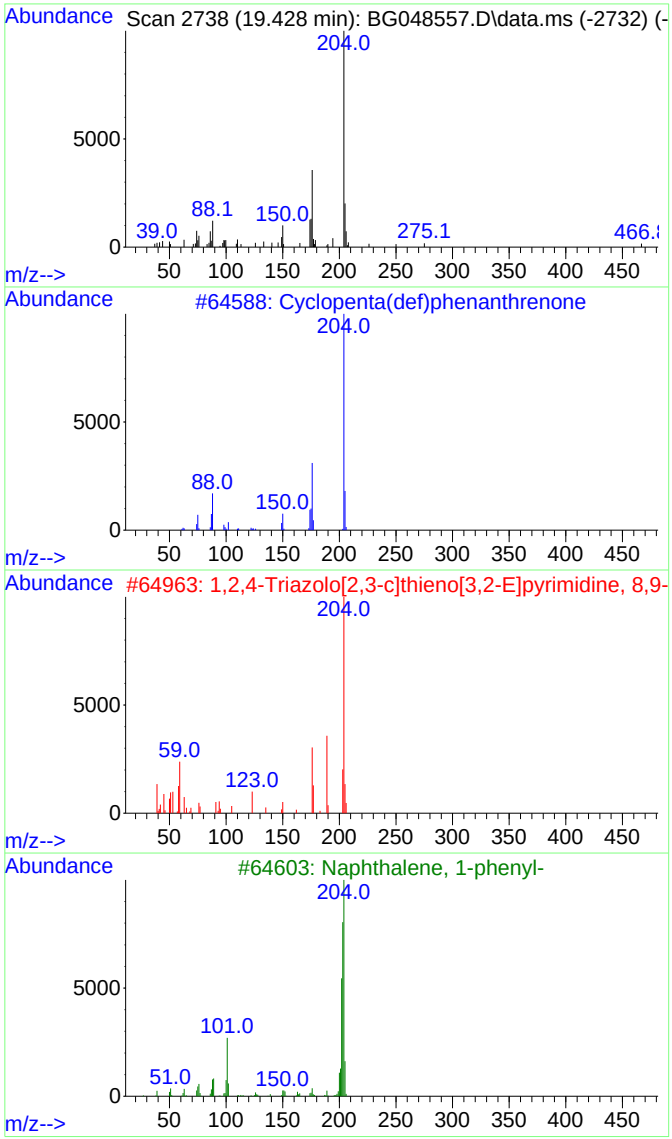
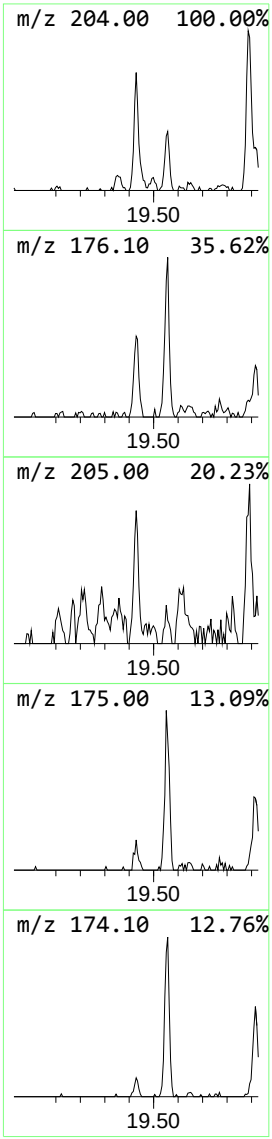
Peak Number 9 Cyclopenta(def)phenanthrenone Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.428	3.18 ng	54612	Phenanthrene-d10	17.501

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	91
2	1,2,4-Triazolo[2,3-c]thieno[3,2-...	204	C9H8N4S	113246-88-1	59
3	Naphthalene, 1-phenyl-	204	C16H12	000605-02-7	53
4	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	50
5	3,4-Biphenyldicarbonitrile	204	C14H8N2	004128-63-6	50





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Operator : CG/JU
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Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : RO-124037

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG033021.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

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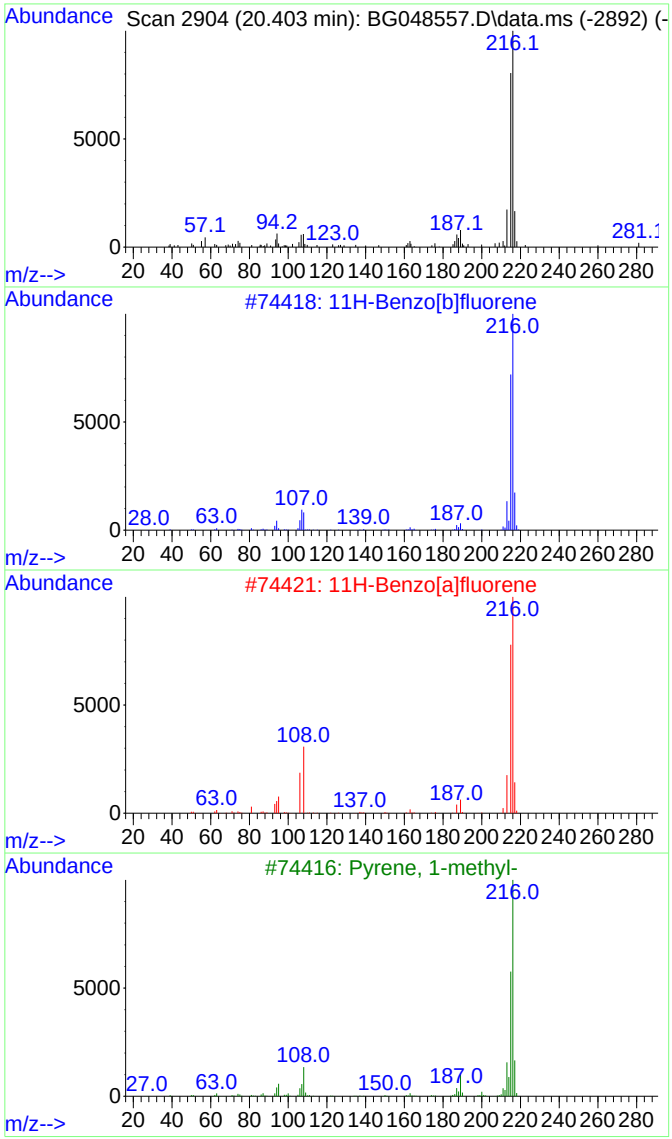
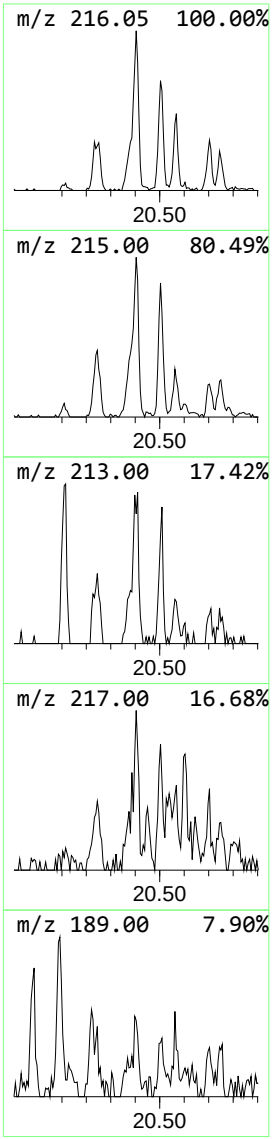
Peak Number 10 11H-Benzo[b]fluorene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.403	4.11 ng	143326	Chrysene-d12	21.796

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	91
2	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	87
3	Pyrene, 1-methyl-	216	C17H12	002381-21-7	86
4	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	86
5	7H-Benzo[c]fluorene	216	C17H12	000205-12-9	78





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Data File : BG048557.D
Acq On : 1 Apr 2021 23:51
Operator : CG/JU
Sample : M1886-07
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
RO-124037

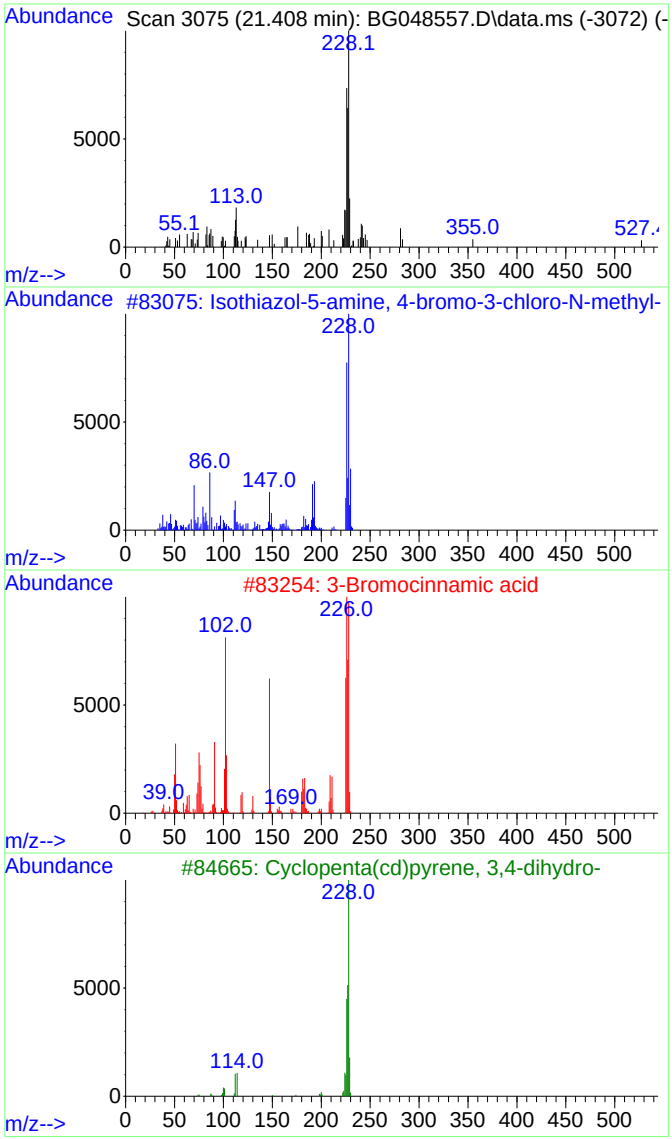
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG033021.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

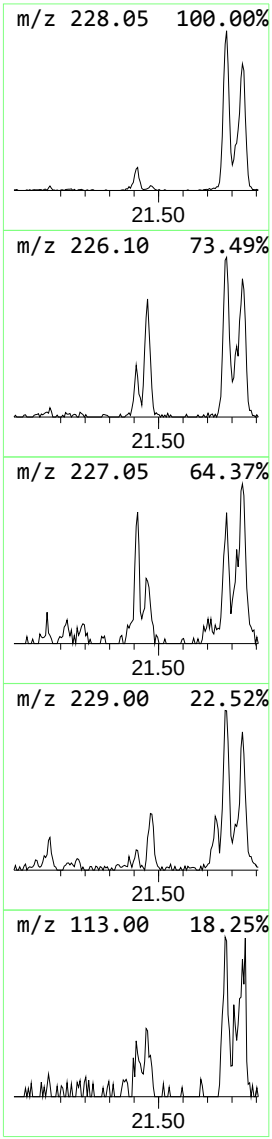
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 Isothiazol-5-amine, 4-bromo... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.408	2.03 ng	70753	Chrysene-d12	21.796

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Isothiazol-5-amine, 4-bromo-3-ch...	226	C4H4BrClN2S	1000272-59-9	59
2	3-Bromocinnamic acid	226	C9H7BrO2	032862-97-8	45
3	Cyclopenta(cd)pyrene, 3,4-dihydro-	228	C18H12	025732-74-5	43
4	Dimethyl-[4-[2-(3-methylisoxazol...	228	C14H16N2O	1000306-39-6	38
5	Benzo[ghi]fluoranthene	226	C18H10	000203-12-3	38





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040121\
Data File : BG048557.D
Acq On : 1 Apr 2021 23:51
Operator : CG/JU
Sample : M1886-07
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : RO-124037

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG033021.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
1H-Indene, 1-et...	12.577	2.4	ng	22162	2	10.921	187520	20.0
Dibenzofuran, 4...	16.085	2.1	ng	27633	3	14.745	263012	20.0
Dibenzothiophene	17.313	3.0	ng	52080	4	17.501	342994	20.0
n-Hexadecanoic ...	18.382	10.1	ng	172365	4	17.501	342994	20.0
1H-Indene, 2-ph...	18.423	5.3	ng	91820	4	17.501	342994	20.0
4H-Cyclopenta[d...	18.576	11.3	ng	193731	4	17.501	342994	20.0
5,16[1',2']:8,1...	18.870	3.1	ng	53246	4	17.501	342994	20.0
9,10-Anthracene...	18.911	2.2	ng	38213	4	17.501	342994	20.0
Cyclopenta(def)...	19.428	3.2	ng	54612	4	17.501	342994	20.0
11H-Benzo[b]flu...	20.403	4.1	ng	143326	5	21.796	698011	20.0
Isothiazol-5-am...	21.408	2.0	ng	70753	5	21.796	698011	20.0