LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100821\
Data File : BG050528.D
Acq On : 9 Oct 2021 3:56
Operator : CG/JU
Sample : M4113-05
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : TP-3

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-BG100621.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG050528.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.798	386	393	403	rBV	768241	1288590	58.46%	5.866%
2	6.926	578	585	592	rBV3	18871	30299	1.37%	0.138%
3	7.396	657	665	676	rBV	798025	1444747	65.55%	6.577%
4	8.259	805	812	823	rVB	259248	459760	20.86%	2.093%
5	9.429	1003	1011	1020	rBV	512295	931374	42.26%	4.240%
6	10.827	1243	1249	1256	rBV	74001	127926	5.80%	0.582%
7	11.085	1283	1293	1310	rBV	372874	694664	31.52%	3.162%
8	13.500	1696	1704	1712	rBV	1143653	1837042	83.35%	8.363%
9	13.747	1742	1746	1753	rVB2	24671	40292	1.83%	0.183%
10	14.200	1816	1823	1828	rVB	162405	246412	11.18%	1.122%
11	14.282	1833	1837	1844	rVB2	14233	22127	1.00%	0.101%
12	14.828	1921	1930	1933	rBV	64277	100378	4.55%	0.457%
13	14.875	1933	1938	1944	rVV	561131	904625	41.04%	4.118%
14	14.940	1944	1949	1957	rVB3	32575	64883	2.94%	0.295%
15	15.110	1974	1978	1985	rBV3	27349	41978	1.90%	0.191%
16	15.269	1998	2005	2011	rBV2	11829	23196	1.05%	0.106%
17	15.921	2110	2116	2124	rVB	24145	40056	1.82%	0.182%
18	16.356	2175	2190	2198	rBV	796491	1347511	61.14%	6.134%
19	16.426	2198	2202	2209	rVB8	10058	22600	1.03%	0.103%
20	17.272	2341	2346	2351	rVB	20536	28837	1.31%	0.131%
21	17.437	2370	2374	2380	rVB2	15083	24075	1.09%	0.110%
22	17.619	2399	2405	2409	rBV	650991	1036049	47.01%	4.716%
23	17.660	2409	2412	2418	rVV	348237	524576	23.80%	2.388%
24	17.754	2418	2428	2433	rVV	39181	72618	3.29%	0.331%
25	18.019	2468	2473	2480	rVB2	29652	51381	2.33%	0.234%
26	18.259	2510	2514	2519	rVB	19754	24334	1.10%	0.111%
27	18.483	2547	2552	2558	rBV2	46811	99761	4.53%	0.454%
28	18.541	2558	2562	2569	rVB	48229	77491	3.52%	0.353%
29	18.688	2580	2587	2591	rBV3	56856	104077	4.72%	0.474%
30	18.982	2632	2637	2640	rBV	31825	48473	2.20%	0.221%
31	19.023	2640	2644	2649	rVB	74862	109677	4.98%	0.499%
32	19.417	2702	2711	2715	rBV5	71671	128490	5.83%	0.585%
33	19.540	2727	2732	2736	rBV3	36933	57405	2.60%	0.261%
34	19.658	2747	2752	2759	rBV	604057	849771	38.55%	3.868%
35	19.987	2802	2808	2809	rBV	93884	128052	5.81%	0.583%
36	20.022	2809	2814	2820	rVV	426209	654736	29.71%	2.980%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100821\
Data File : BG050528.D
Acq On : 9 Oct 2021 3:56
Operator : CG/JU
Sample : M4113-05
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : TP-3

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-BG100621.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	20.081	2821	2824	2831	rVB2	22981	37828	1.72%	0.172%
38	20.210	2839	2846	2851	rBV	1610702	2204071	100.00%	10.033%
39	20.257	2851	2854	2860	rVB3	16879	25607	1.16%	0.117%
40	20.333	2861	2867	2870	rBV3	34347	77200	3.50%	0.351%
41	20.474	2886	2891	2892	rBV6	22875	39156	1.78%	0.178%
42	20.504	2892	2896	2901	rVB	72156	116400	5.28%	0.530%
43	20.604	2909	2913	2917	rBV2	40325	57499	2.61%	0.262%
44	20.845	2951	2954	2958	rBV4	25070	37193	1.69%	0.169%
45	21.285	3025	3029	3034	rBV2	37923	61635	2.80%	0.281%
46	21.344	3034	3039	3045	rVB	503097	690753	31.34%	3.144%
47	21.520	3065	3069	3074	rVB2	132718	196524	8.92%	0.895%
48	21.632	3084	3088	3098	rVB	55496	111899	5.08%	0.509%
49	21.920	3128	3137	3142	rBV2	586386	1312731	59.56%	5.976%
50	21.967	3142	3145	3153	rVB	170563	297809	13.51%	1.356%
51	22.760	3274	3280	3290	rVB7	89583	218634	9.92%	0.995%
52	23.677	3431	3436	3442	rVB	84112	171209	7.77%	0.779%
53	24.247	3525	3533	3540	rBV2	105468	361657	16.41%	1.646%
54	24.329	3542	3547	3554	rVV3	147685	367403	16.67%	1.672%
55	24.399	3555	3559	3571	rVB7	55180	145762	6.61%	0.664%
56	25.016	3657	3664	3674	rVB	65556	198343	9.00%	0.903%
57	25.175	3684	3691	3703	rVB2	87925	262825	11.92%	1.196%
58	25.339	3709	3719	3729	rBV2	324281	987602	44.81%	4.496%
59	26.567	3921	3928	3938	rVB3	97219	299594	13.59%	1.364%

Sum of corrected areas: 21967597

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100821\
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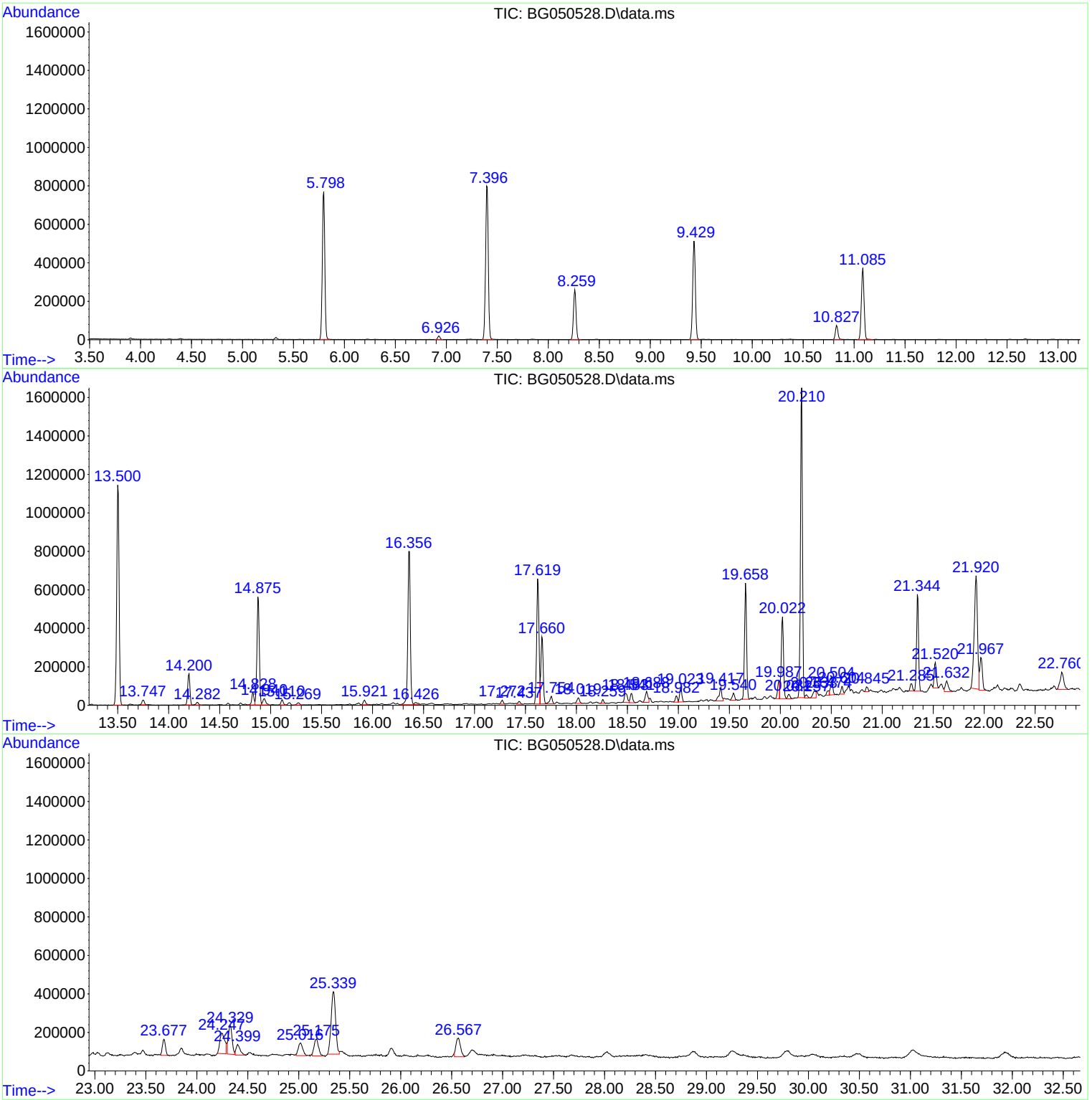
Operator : CG/JU
Sample : M4113-05
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
TP-3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.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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Data File : BG050528.D
Acq On : 9 Oct 2021 3:56
Operator : CG/JU
Sample : M4113-05
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : TP-3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

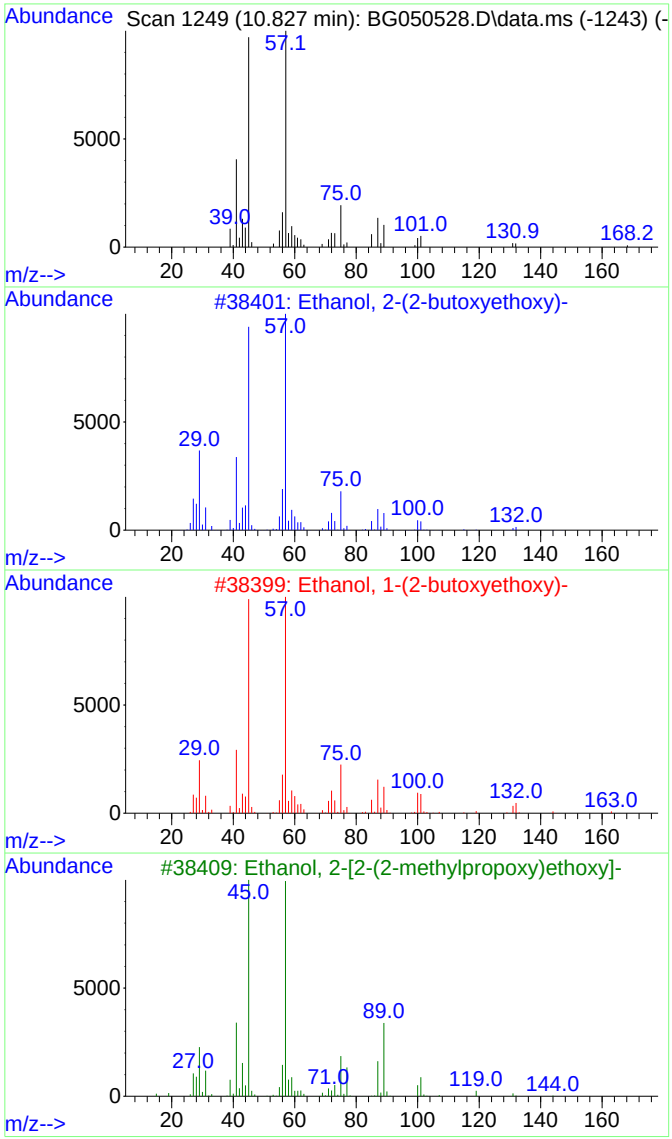
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

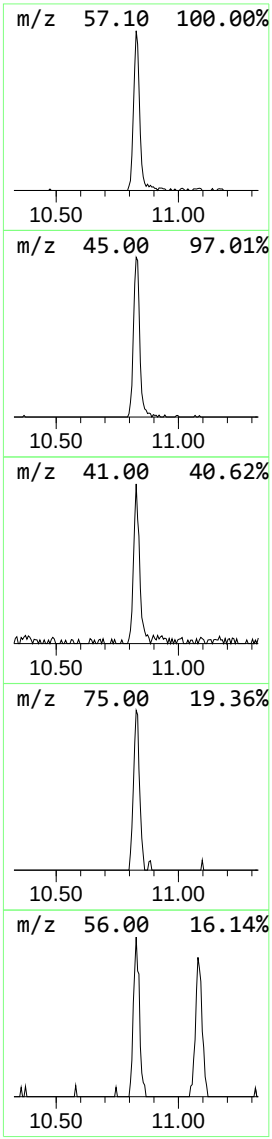
****

Peak Number 1 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.827	3.68 ng	127926	Naphthalene-d8	11.086

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2	Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	83
3	Ethanol, 2-[2-(2-methylpropoxy)e...	162	C8H18O3	018912-80-6	56
4	2-[2-[2-[2-[2-[2-(2-Hydroxyet...	370	C16H34O9	005117-19-1	50
5	Di-sec-Butyl ether	130	C8H18O	006863-58-7	50





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100821\
Data File : BG050528.D
Acq On : 9 Oct 2021 3:56
Operator : CG/JU
Sample : M4113-05
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : TP-3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.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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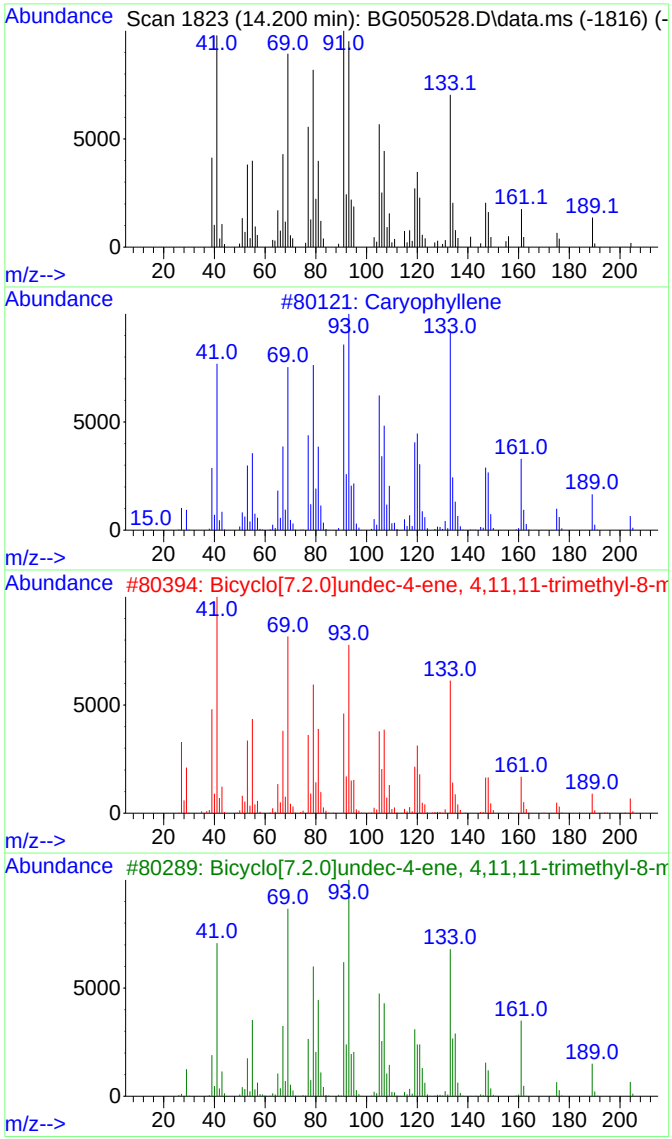
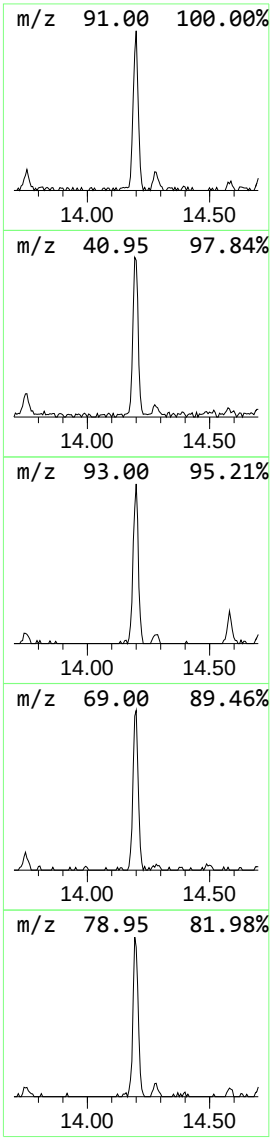
Peak Number 2 Caryophyllene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.200	5.45 ng	246412	Acenaphthene-d10	14.875

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Caryophyllene	204	C15H24	000087-44-5	99
2	Bicyclo[7.2.0]undec-4-ene, 4,11,...	204	C15H24	000118-65-0	92
3	Bicyclo[7.2.0]undec-4-ene, 4,11,...	204	C15H24	013877-93-5	70
4	6,6-Dimethyl-2-vinylidenebicyclo...	148	C11H16	039021-75-5	43
5	trans-.beta.-Ocimene	136	C10H16	003779-61-1	38





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100821\
Data File : BG050528.D
Acq On : 9 Oct 2021 3:56
Operator : CG/JU
Sample : M4113-05
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : TP-3

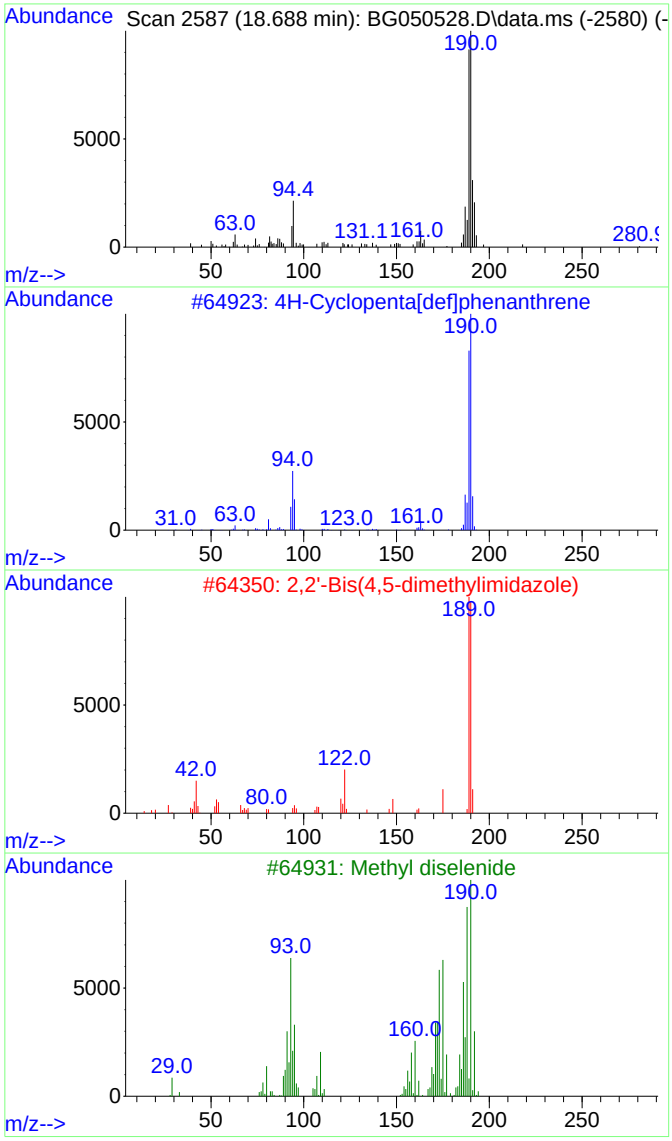
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

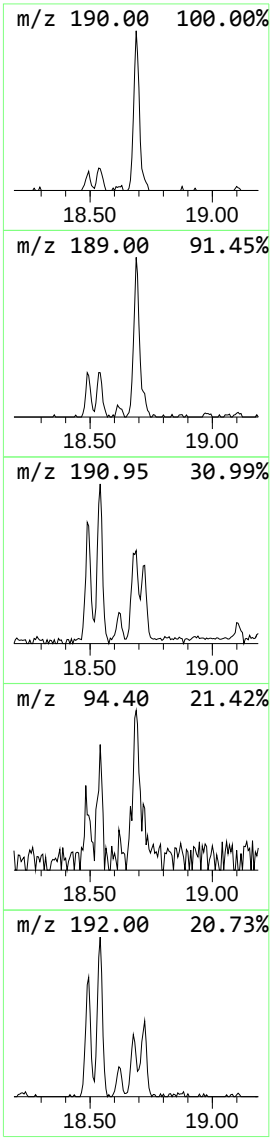
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 4H-Cyclopenta[def]phenanthrene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.688	2.01 ng	104077	Phenanthrene-d10	17.619

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	80
2	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	50
3	Methyl diselenide	190	C2H6Se2	007101-31-7	43
4	1,4-Naphthalenedione, 2,5-dihydr...	190	C10H6O4	004923-55-1	35
5	5,8-Quinoxalinediol, 2,3-dimethyl-	190	C10H10N2O2	007698-00-2	27





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100821\
Data File : BG050528.D
Acq On : 9 Oct 2021 3:56
Operator : CG/JU
Sample : M4113-05
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : TP-3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

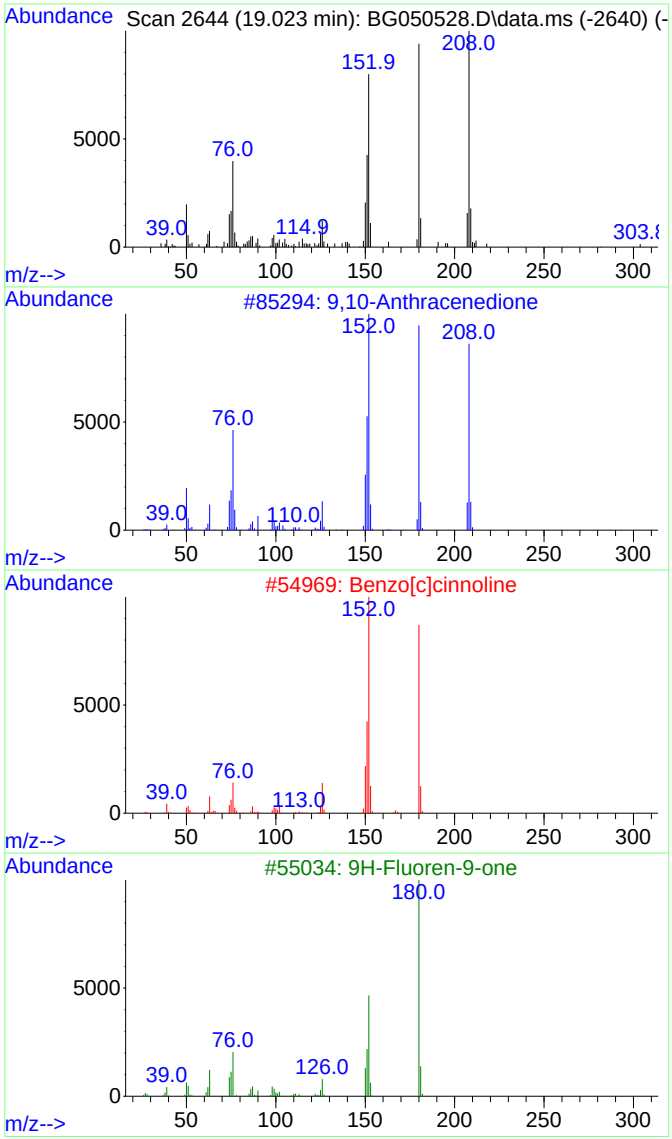
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

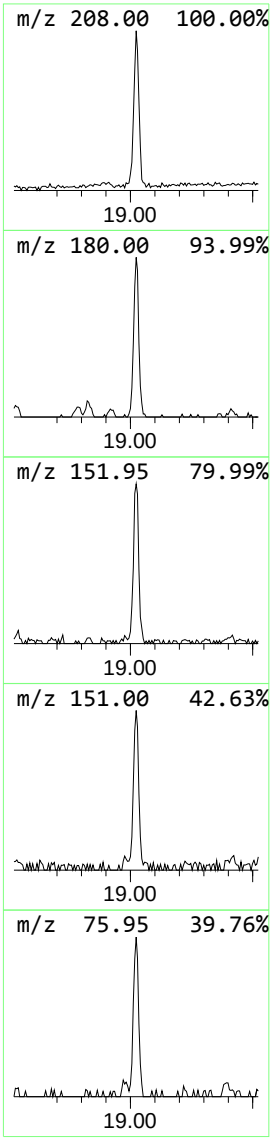
****

Peak Number 4 9,10-Anthracenedione Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.023	2.12 ng	109677	Phenanthrene-d10	17.619

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Anthracenedione	208	C14H8O2	000084-65-1	97
2	Benzo[c]cinnoline	180	C12H8N2	000230-17-1	93
3	9H-Fluoren-9-one	180	C13H8O	000486-25-9	83
4	1H-Phenalen-1-one	180	C13H8O	000548-39-0	64
5	9,10-Phenanthrenedione	208	C14H8O2	000084-11-7	60





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100821\
Data File : BG050528.D
Acq On : 9 Oct 2021 3:56
Operator : CG/JU
Sample : M4113-05
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : TP-3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

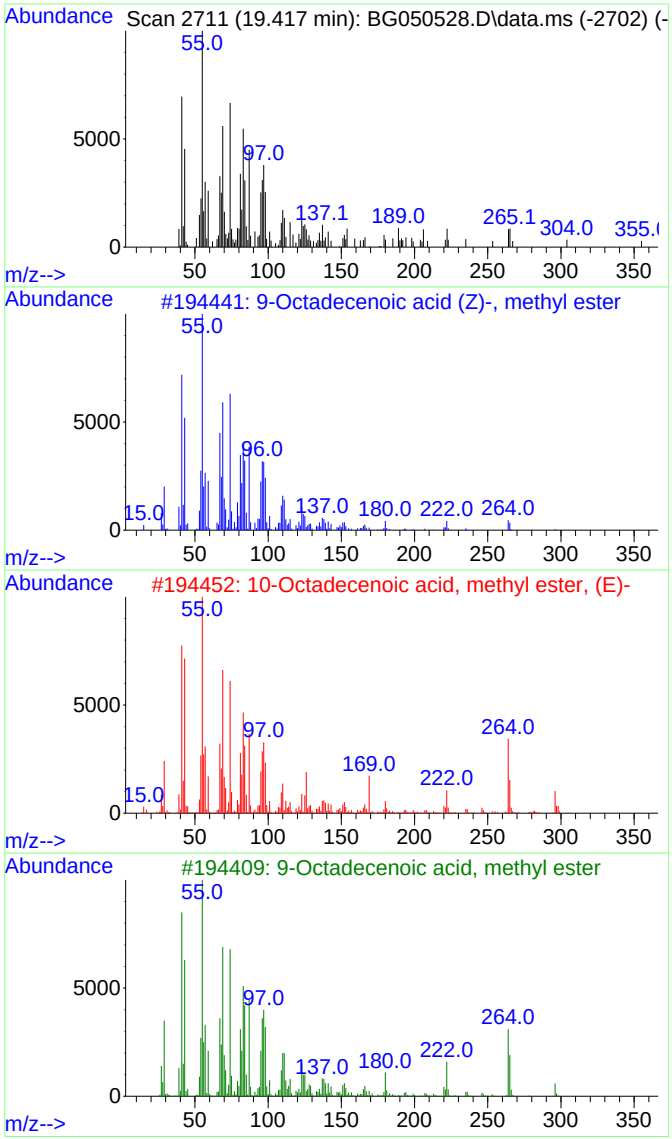
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

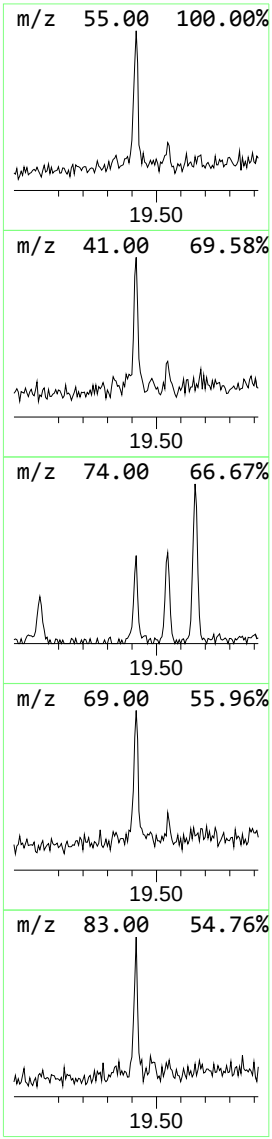
****

Peak Number 5 9-Octadecenoic acid (Z)-, m... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.417	2.48 ng	128490	Phenanthrene-d10	17.619

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	91
2	10-Octadecenoic acid, methyl est...	296	C19H36O2	013038-45-4	90
3	9-Octadecenoic acid, methyl ester	296	C19H36O2	002462-84-2	90
4	12-Octadecenoic acid, methyl ester	296	C19H36O2	056554-46-2	90
5	15-Octadecenoic acid, methyl ester	296	C19H36O2	004764-72-1	87





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Sample : M4113-05
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

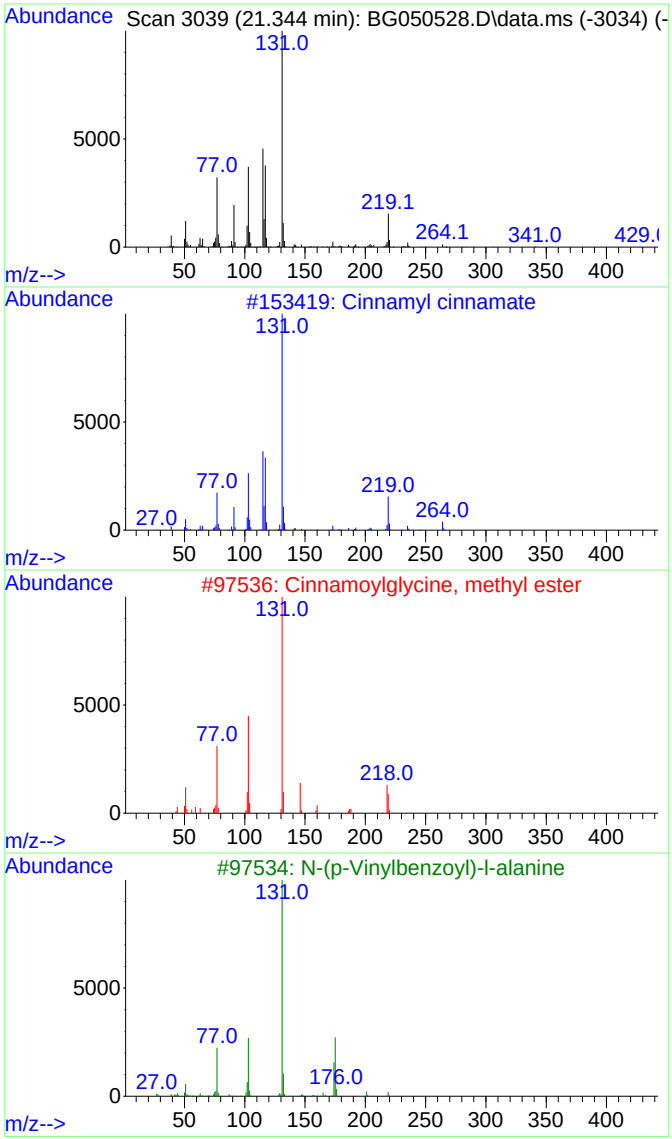
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

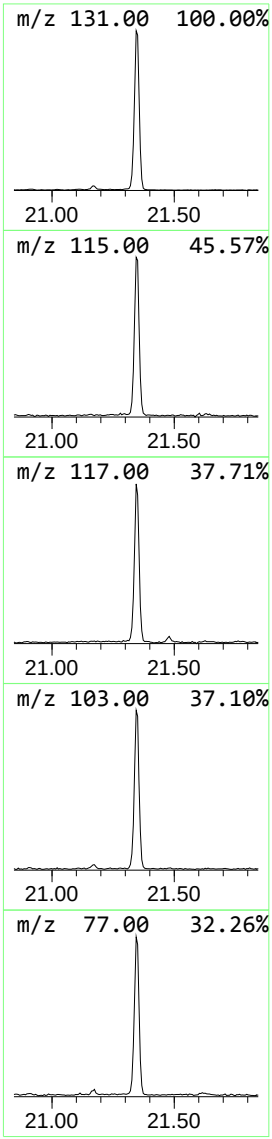
****

Peak Number 6 Cinnamyl cinnamate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.344	10.52 ng	690753	Chrysene-d12	21.920

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cinnamyl cinnamate	264	C18H16O2	000122-69-0	91
2	Cinnamoylglycine, methyl ester	219	C12H13NO3	040778-04-9	64
3	N-(p-Vinylbenzoyl)-l-alanine	219	C12H13NO3	139184-60-4	50
4	trans-1-Cinnamoylimidazole	198	C12H10N2O	001138-15-4	47
5	3-phenylprop-2-enoic anhydride	278	C18H14O3	1010401-87-5	47





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100821\
Data File : BG050528.D
Acq On : 9 Oct 2021 3:56
Operator : CG/JU
Sample : M4113-05
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : TP-3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.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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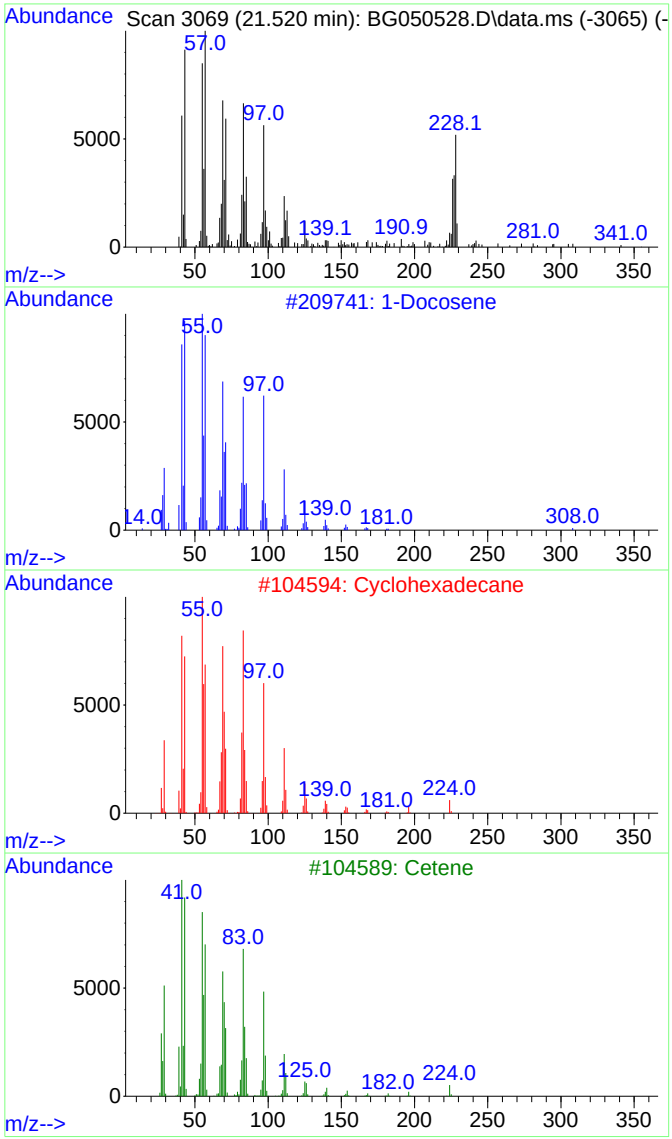
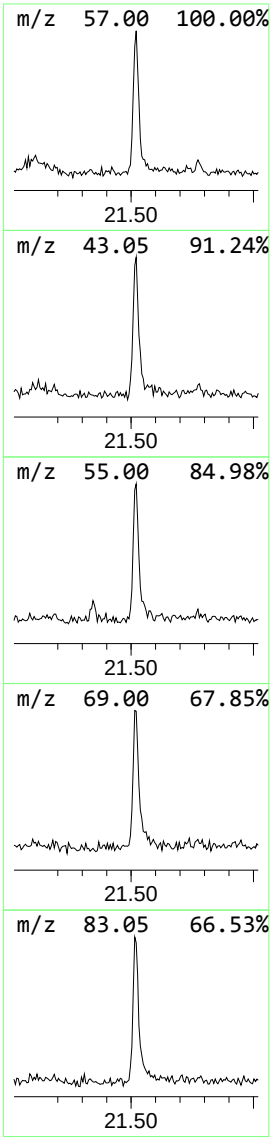
Peak Number 7 1-Docosene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.520	2.99 ng	196524	Chrysene-d12	21.920

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Docosene	308	C22H44	001599-67-3	90
2	Cyclohexadecane	224	C16H32	000295-65-8	87
3	Cetene	224	C16H32	000629-73-2	83
4	Henicos-1-ene	294	C21H42	001599-68-4	83
5	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100821\
Data File : BG050528.D
Acq On : 9 Oct 2021 3:56
Operator : CG/JU
Sample : M4113-05
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : TP-3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.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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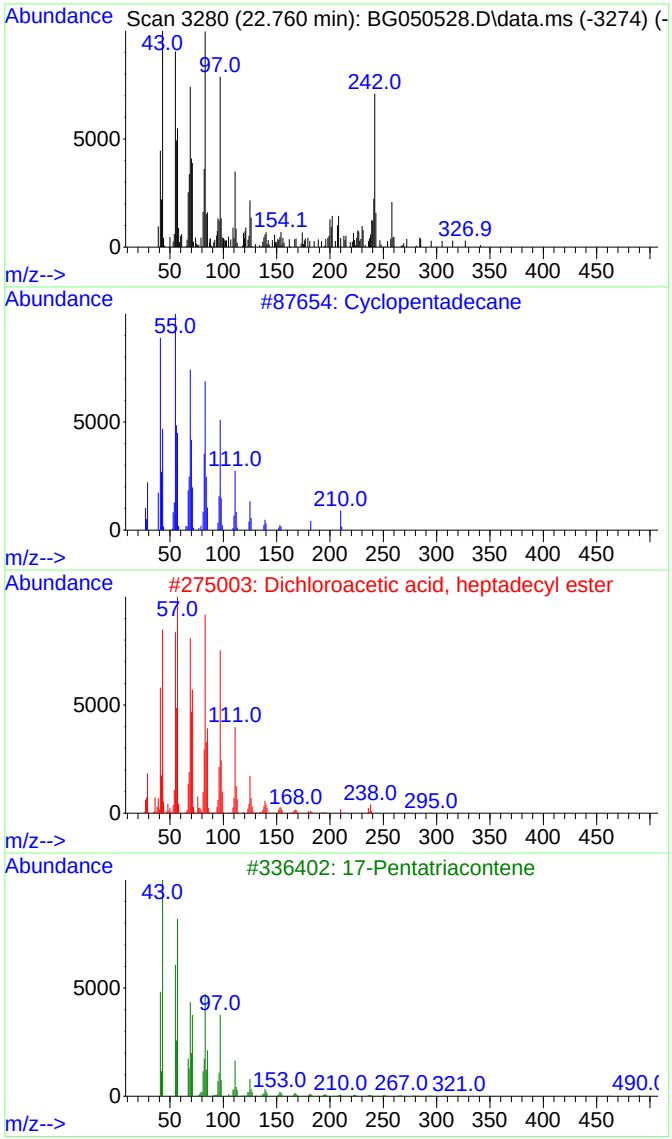
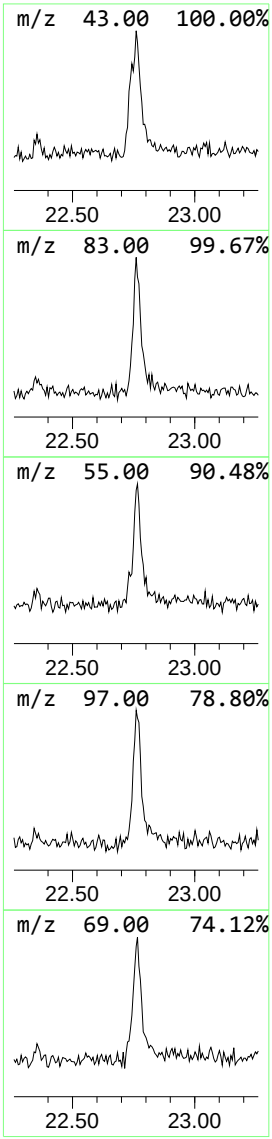
Peak Number 8 Cyclopentadecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.760	3.33 ng	218634	Chrysene-d12	21.920

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentadecane	210	C15H30	000295-48-7	70
2	Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	58
3	17-Pentatriacontene	491	C35H70	006971-40-0	52
4	Cyclohexadecane	224	C16H32	000295-65-8	49
5	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	49





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Data File : BG050528.D
Acq On : 9 Oct 2021 3:56
Operator : CG/JU
Sample : M4113-05
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : TP-3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.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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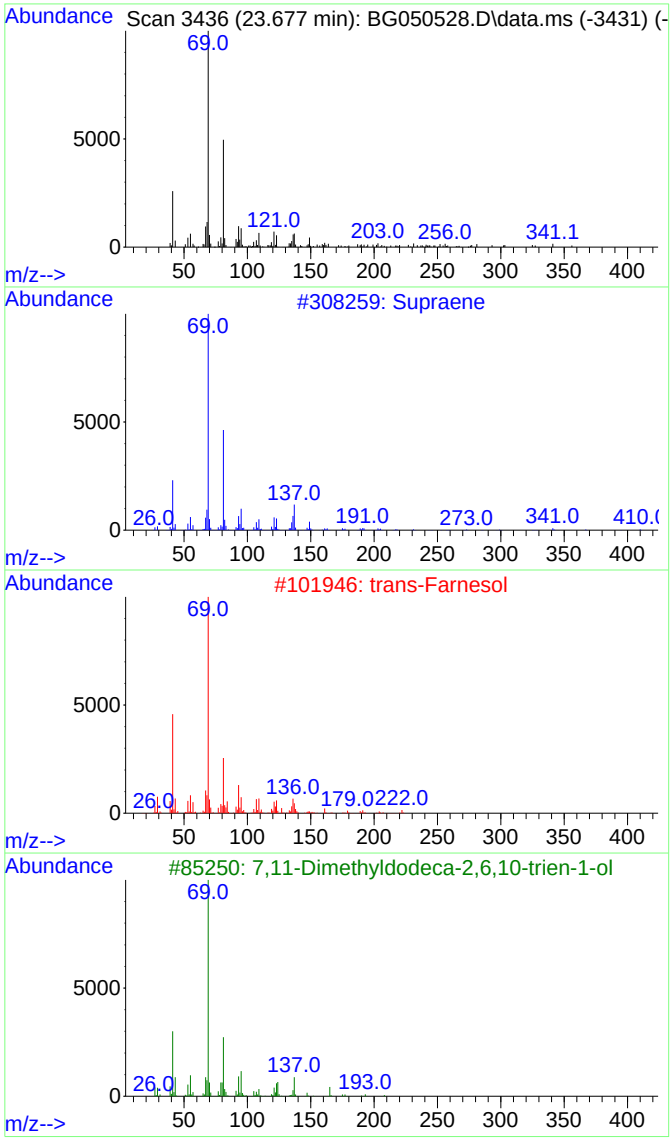
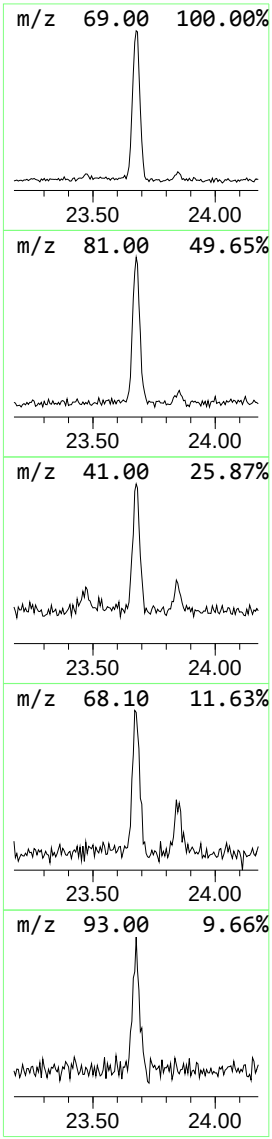
Peak Number 9 Supraene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.677	3.47 ng	171209	Perylene-d12	25.339

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Supraene	410	C30H50	007683-64-9	72
2	trans-Farnesol	222	C15H26O	000106-28-5	72
3	7,11-Dimethyldodeca-2,6,10-trien...	208	C14H24O	125529-10-4	64
4	1,5-Heptadiene, 3,3,6-trimethyl-	138	C10H18	035387-63-4	58
5	4-Methyl-1,5-Heptadiene	110	C8H14	000998-94-7	58





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Operator : CG/JU
Sample : M4113-05
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : TP-3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.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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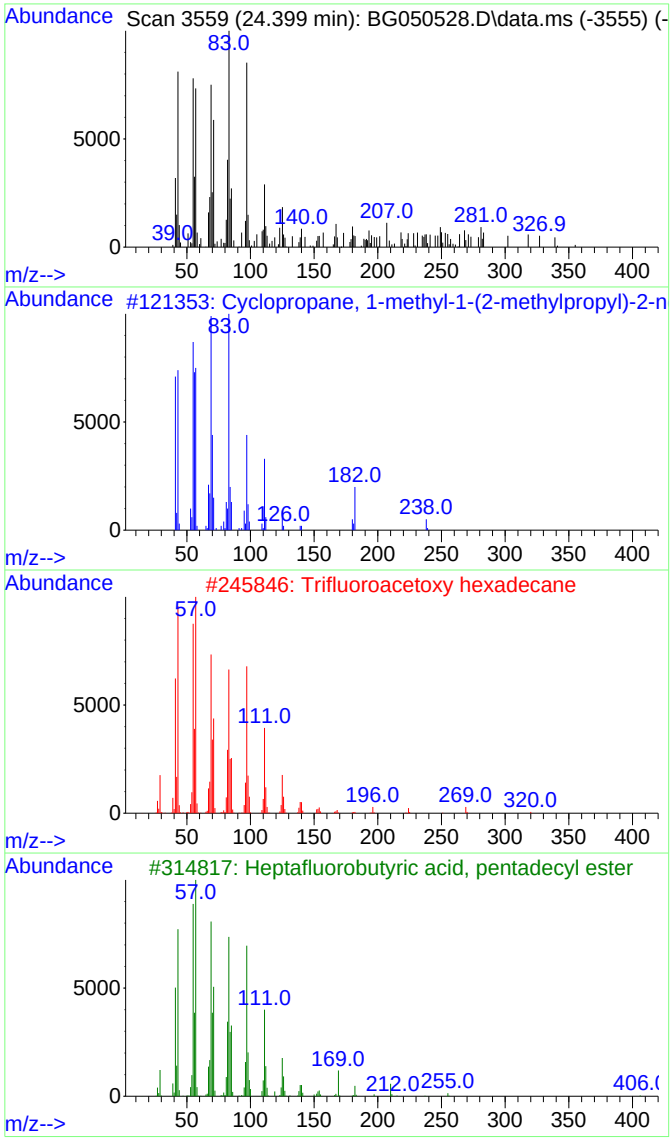
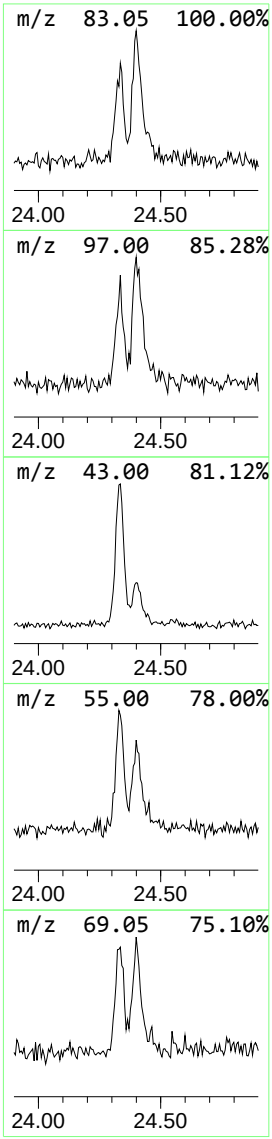
Peak Number 10 Cyclopropane, 1-methyl-1-(2... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.399	2.95 ng	145762	Perylene-d12	25.339

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopropane, 1-methyl-1-(2-meth...	238	C17H34	041977-41-7	70
2	Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	64
3	Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	64
4	Heptafluorobutyric acid, n-tetra...	410	C18H29F7O2	007365-36-8	64
5	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	60





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Acq On : 9 Oct 2021 3:56
Operator : CG/JU
Sample : M4113-05
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : TP-3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.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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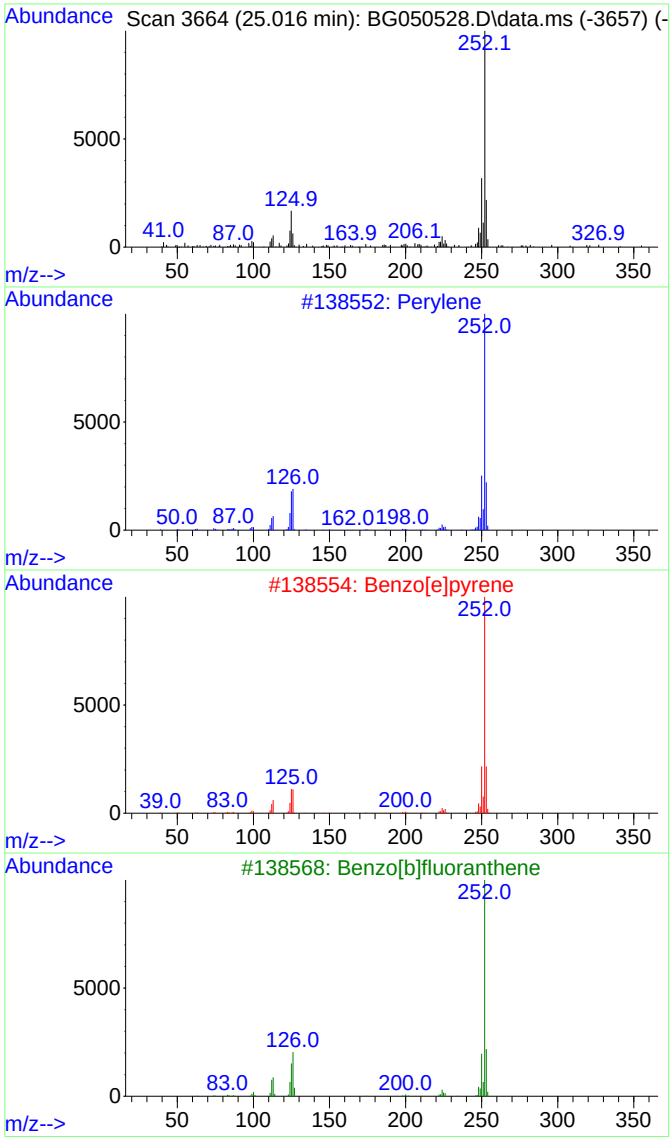
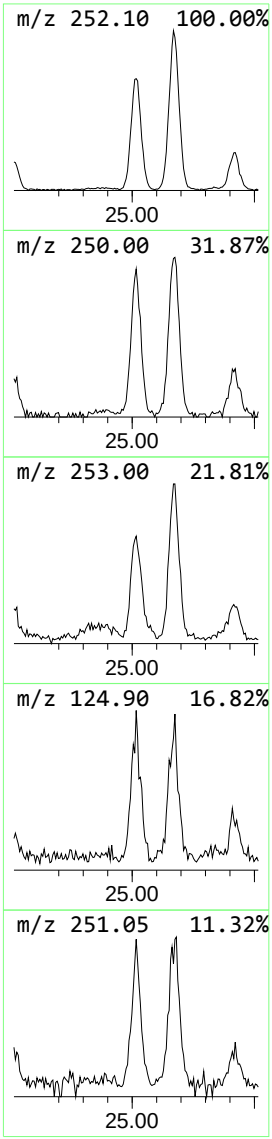
Peak Number 11 Perylene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.016	4.02 ng	198343	Perylene-d12	25.339

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Perylene	252	C20H12	000198-55-0	93
2	Benzo[e]pyrene	252	C20H12	000192-97-2	91
3	Benzo[b]fluoranthene	252	C20H12	000205-99-2	91
4	Benzo[k]fluoranthene	252	C20H12	000207-08-9	90
5	Benzo[a]pyrene	252	C20H12	000050-32-8	90





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Data File : BG050528.D
Acq On : 9 Oct 2021 3:56
Operator : CG/JU
Sample : M4113-05
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : TP-3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.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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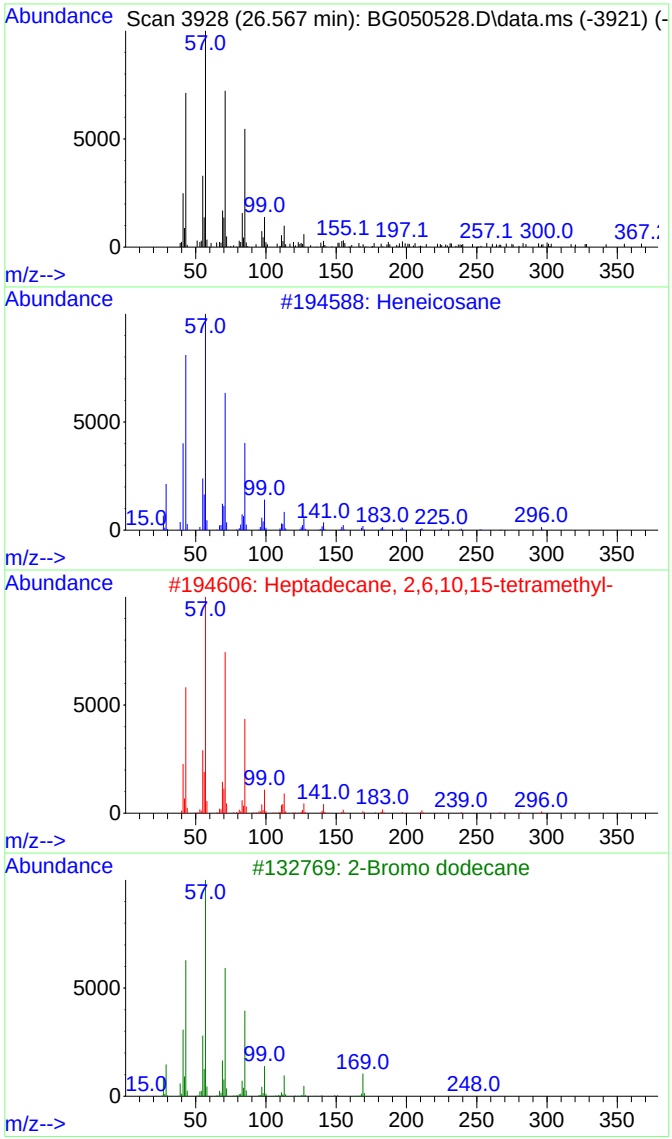
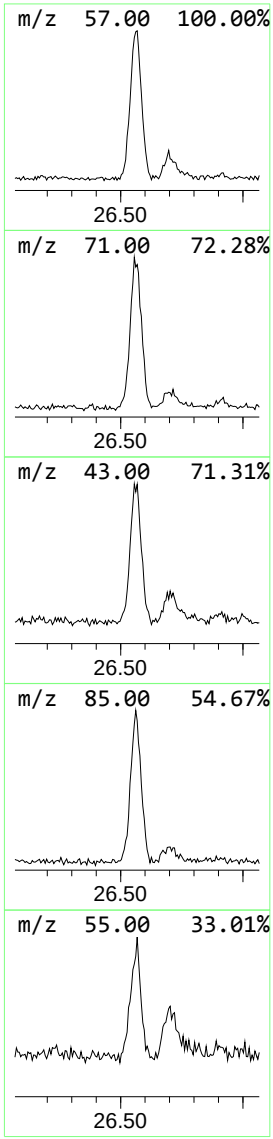
Peak Number 12 Heneicosane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.567	6.07 ng	299594	Perylene-d12	25.339

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	96
2	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	93
3	2-Bromo dodecane	248	C12H25Br	013187-99-0	86
4	Octacosane	394	C28H58	000630-02-4	86
5	Tetracosane	338	C24H50	000646-31-1	86





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100821\
Data File : BG050528.D
Acq On : 9 Oct 2021 3:56
Operator : CG/JU
Sample : M4113-05
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : TP-3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.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
Ethanol, 2-(2-b...	10.827	3.7	ng	127926	2	11.086	694664	20.0
Caryophyllene	14.200	5.5	ng	246412	3	14.875	904625	20.0
4H-Cyclopenta[d...	18.688	2.0	ng	104077	4	17.619	1036050	20.0
9,10-Anthracene...	19.023	2.1	ng	109677	4	17.619	1036050	20.0
9-Octadecenoic ...	19.417	2.5	ng	128490	4	17.619	1036050	20.0
Cinnamyl cinnamate	21.344	10.5	ng	690753	5	21.920	1312730	20.0
1-Docosene	21.520	3.0	ng	196524	5	21.920	1312730	20.0
Cyclopentadecane	22.760	3.3	ng	218634	5	21.920	1312730	20.0
Supraene	23.677	3.5	ng	171209	6	25.339	987602	20.0
Cyclopropane, 1...	24.399	3.0	ng	145762	6	25.339	987602	20.0
Perylene	25.016	4.0	ng	198343	6	25.339	987602	20.0
Heneicosane	26.567	6.1	ng	299594	6	25.339	987602	20.0