LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062923\
Data File : BG057914.D
Acq On : 29 Jun 2023 21:38
Operator : CG/JU
Sample : 03434-04
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : 350

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-BG062723.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG057914.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.021	323	329	338	rVB	66763	103131	5.08%	0.505%
2	5.491	401	409	419	rBV	647900	1058313	52.09%	5.184%
3	7.077	671	679	692	rBV	645773	1121802	55.22%	5.494%
4	7.911	814	821	831	rBV	196442	328426	16.17%	1.609%
5	9.075	1010	1019	1029	rBV	392880	687782	33.85%	3.369%
6	10.714	1287	1298	1306	rBV	281033	502475	24.73%	2.461%
7	13.170	1709	1716	1724	rBV	1010952	1537632	75.69%	7.531%
8	14.275	1899	1904	1909	rBV	15576	28634	1.41%	0.140%
9	14.545	1941	1950	1957	rBV2	430748	691114	34.02%	3.385%
10	14.950	2011	2019	2026	rBV4	18139	37251	1.83%	0.182%
11	15.591	2123	2128	2133	rBV4	13907	22377	1.10%	0.110%
12	15.902	2172	2181	2187	rBV	136983	204008	10.04%	0.999%
13	16.031	2195	2203	2217	rBV	733467	1190533	58.60%	5.831%
14	16.266	2237	2243	2249	rVB5	15751	30905	1.52%	0.151%
15	16.948	2353	2359	2365	rVB2	21202	35810	1.76%	0.175%
16	17.107	2381	2386	2394	rVB	16425	32701	1.61%	0.160%
17	17.206	2395	2403	2408	rBV	33596	56941	2.80%	0.279%
18	17.294	2412	2418	2422	rBV	538352	837442	41.22%	4.102%
19	17.336	2422	2425	2430	rVB	232875	309944	15.26%	1.518%
20	17.424	2436	2440	2444	rBV	25910	35596	1.75%	0.174%
21	17.477	2444	2449	2455	rVB6	25828	42003	2.07%	0.206%
22	17.659	2476	2480	2483	rBV	56723	83131	4.09%	0.407%
23	17.694	2483	2486	2492	rVB	26028	41103	2.02%	0.201%
24	17.870	2512	2516	2522	rBV	18862	30109	1.48%	0.147%
25	18.011	2537	2540	2547	rBV2	10276	21058	1.04%	0.103%
26	18.182	2561	2569	2572	rBV2	346980	542224	26.69%	2.656%
27	18.211	2572	2574	2580	rVB	103983	147788	7.27%	0.724%
28	18.358	2594	2599	2603	rVV2	86506	157653	7.76%	0.772%
29	18.399	2603	2606	2612	rVB	54574	77147	3.80%	0.378%
30	18.475	2616	2619	2623	rBV5	15148	21732	1.07%	0.106%
31	18.564	2630	2634	2638	rBV4	17549	23933	1.18%	0.117%
32	18.664	2647	2651	2654	rBV	47219	66346	3.27%	0.325%
33	18.705	2654	2658	2667	rVB3	51135	96897	4.77%	0.475%
34	18.910	2689	2693	2698	rVB2	62522	78243	3.85%	0.383%
35	18.963	2698	2702	2705	rVV	34821	46198	2.27%	0.226%
36	18.998	2705	2708	2712	rVB2	21794	30283	1.49%	0.148%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062923\
Data File : BG057914.D
Acq On : 29 Jun 2023 21:38
Operator : CG/JU
Sample : 03434-04
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
350

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-BG062723.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	19.081	2717	2722	2727	rBV	86672	145053	7.14%	0.710%
38	19.139	2728	2732	2736	rVV4	25723	39945	1.97%	0.196%
39	19.222	2742	2746	2752	rBV2	44306	75727	3.73%	0.371%
40	19.345	2758	2767	2773	rBV	638457	1052041	51.78%	5.153%
41	19.457	2781	2786	2792	rBV3	116035	186148	9.16%	0.912%
42	19.586	2806	2808	2818	rVB6	27741	52441	2.58%	0.257%
43	19.709	2824	2829	2835	rVB	514982	761007	37.46%	3.727%
44	19.780	2838	2841	2846	rVB2	41449	59901	2.95%	0.293%
45	19.909	2857	2863	2868	rBV	1636302	2031587	100.00%	9.951%
46	20.027	2879	2883	2884	rBV2	55532	68523	3.37%	0.336%
47	20.197	2903	2912	2918	rVB2	86852	228309	11.24%	1.118%
48	20.297	2927	2929	2934	rVB	34907	37989	1.87%	0.186%
49	20.356	2935	2939	2943	rVV2	73196	104310	5.13%	0.511%
50	20.444	2951	2954	2959	rVB4	41530	52693	2.59%	0.258%
51	20.497	2959	2963	2967	rBV	80558	115412	5.68%	0.565%
52	20.544	2968	2971	2974	rVB	59097	73402	3.61%	0.360%
53	20.949	3037	3040	3046	rVB	65555	95749	4.71%	0.469%
54	21.196	3075	3082	3091	rVB4	184128	398183	19.60%	1.950%
55	21.437	3119	3123	3131	rVB	179271	240716	11.85%	1.179%
56	21.542	3133	3141	3146	rVV3	581543	1278961	62.95%	6.264%
57	21.589	3146	3149	3156	rVB	236396	363003	17.87%	1.778%
58	21.736	3170	3174	3178	rBV2	92795	123149	6.06%	0.603%
59	22.330	3271	3275	3286	rVB4	111389	202132	9.95%	0.990%
60	23.687	3500	3506	3515	rBV	180663	559951	27.56%	2.743%
61	24.398	3622	3627	3637	rVB	150376	363614	17.90%	1.781%
62	24.539	3645	3651	3665	rVB	165701	464255	22.85%	2.274%
63	24.692	3669	3677	3687	rBV	303848	883991	43.51%	4.330%

Sum of corrected areas: 20416857

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062923\
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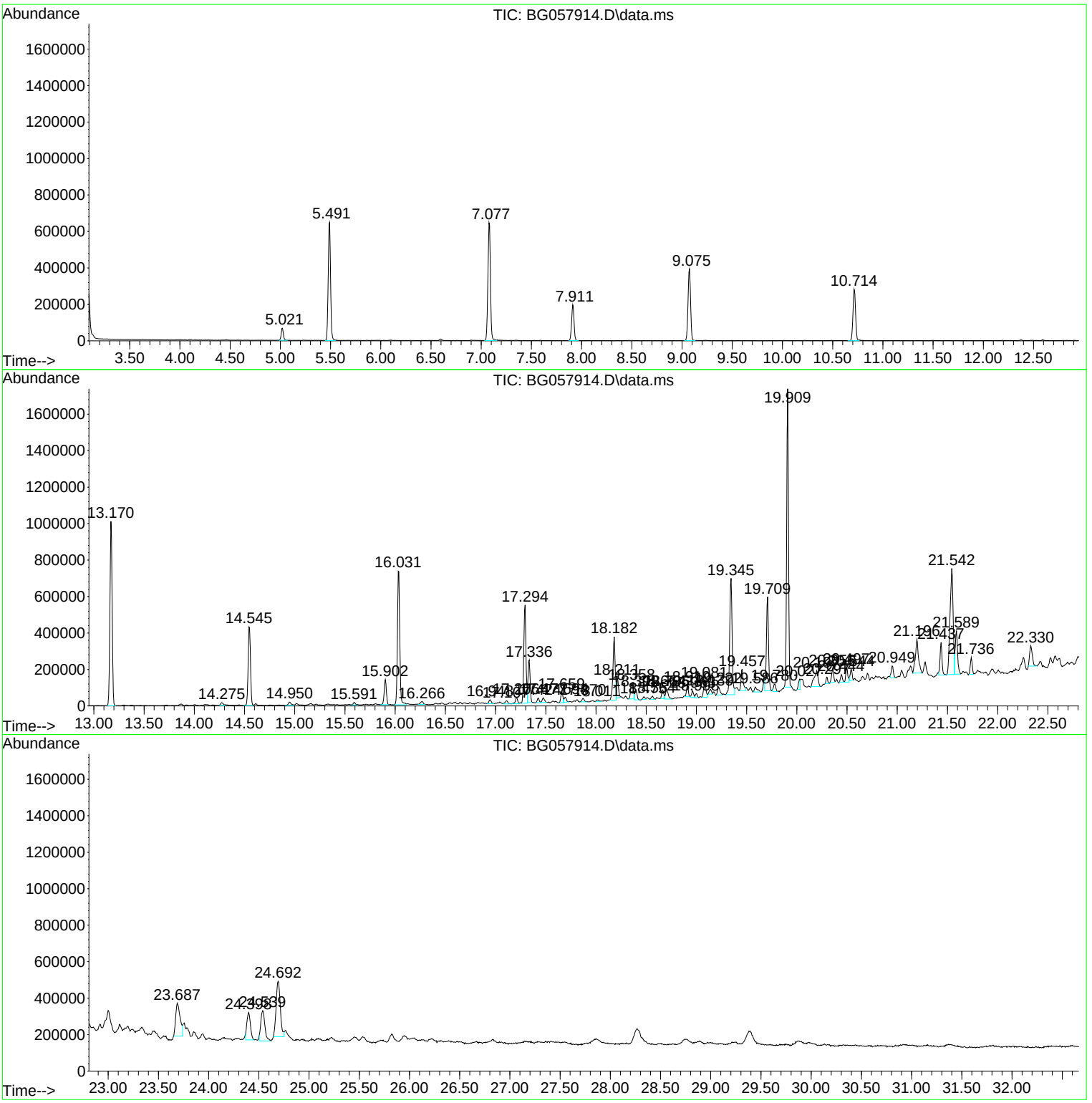
Sample : 03434-04
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
350

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG062723.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\BG062923\
Data File : BG057914.D
Acq On : 29 Jun 2023 21:38
Operator : CG/JU
Sample : 03434-04
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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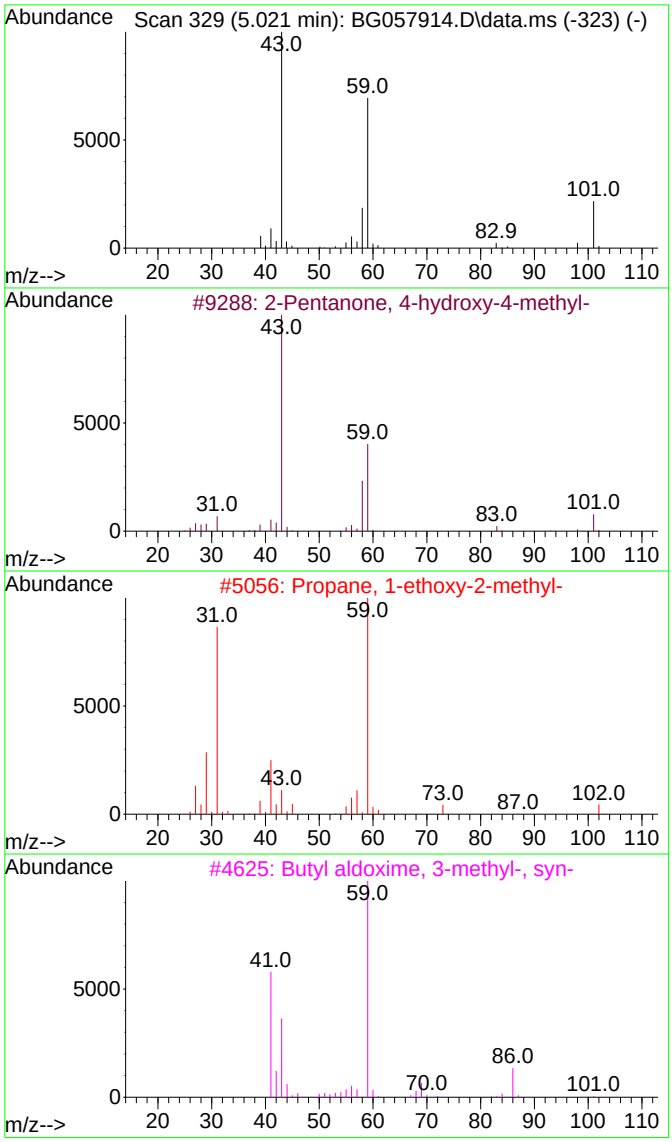
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG062723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

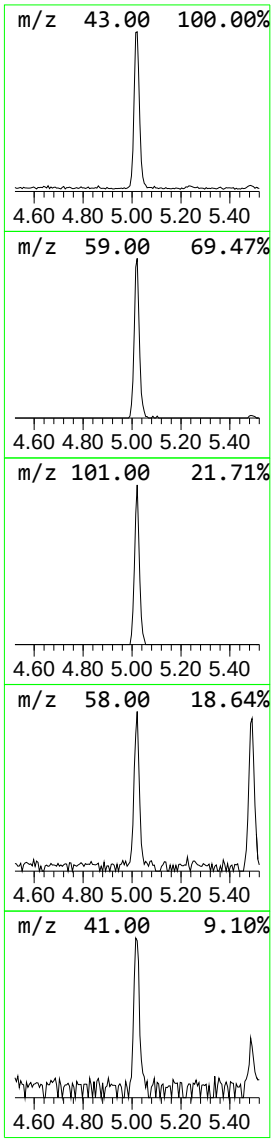
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.021	6.28 ng	103131	1,4-Dichlorobenzene-d4	7.911

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2	Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	35
3	Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	25
4	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	25
5	1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	17





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062923\
Data File : BG057914.D
Acq On : 29 Jun 2023 21:38
Operator : CG/JU
Sample : 03434-04
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
350

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG062723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

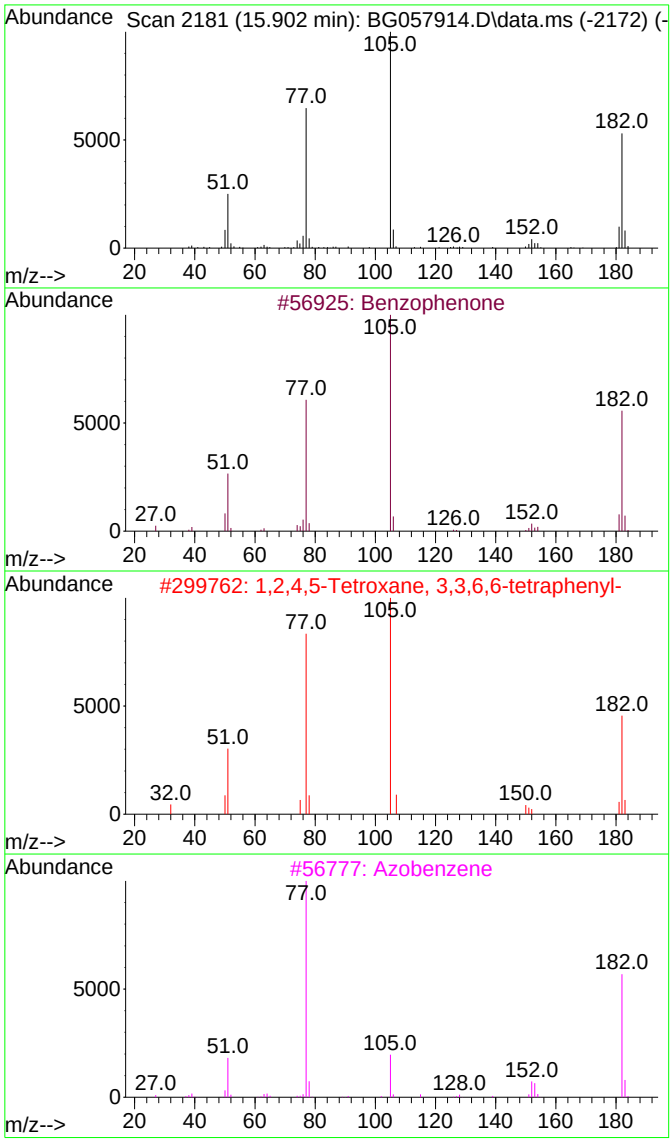
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

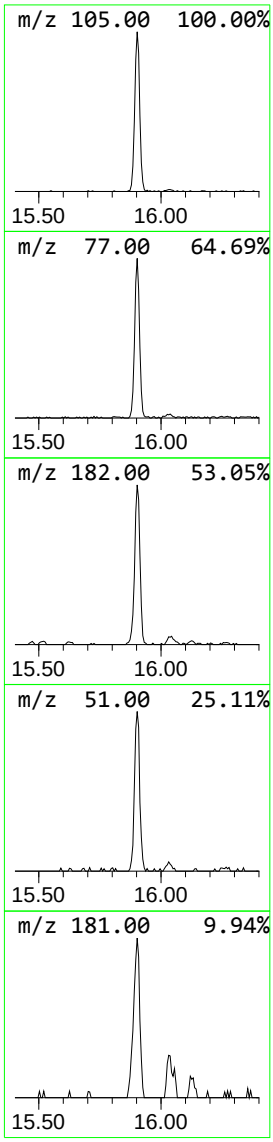
****

Peak Number 2 Benzophenone Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.902	5.90 ng	204008	Acenaphthene-d10	14.545

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzophenone	182	C13H10O	000119-61-9	96
2	1,2,4,5-Tetroxane, 3,3,6,6-tetra...	396	C26H20O4	016204-36-7	78
3	Azobenzene	182	C12H10N2	000103-33-3	59
4	Acetophenone, 2-chloro-	154	C8H7ClO	000532-27-4	49
5	N-Methylbenzamide, N-chlorodiflu...	247	C10H8ClF2NO2	1000446-98-2	47





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062923\
Data File : BG057914.D
Acq On : 29 Jun 2023 21:38
Operator : CG/JU
Sample : 03434-04
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
350

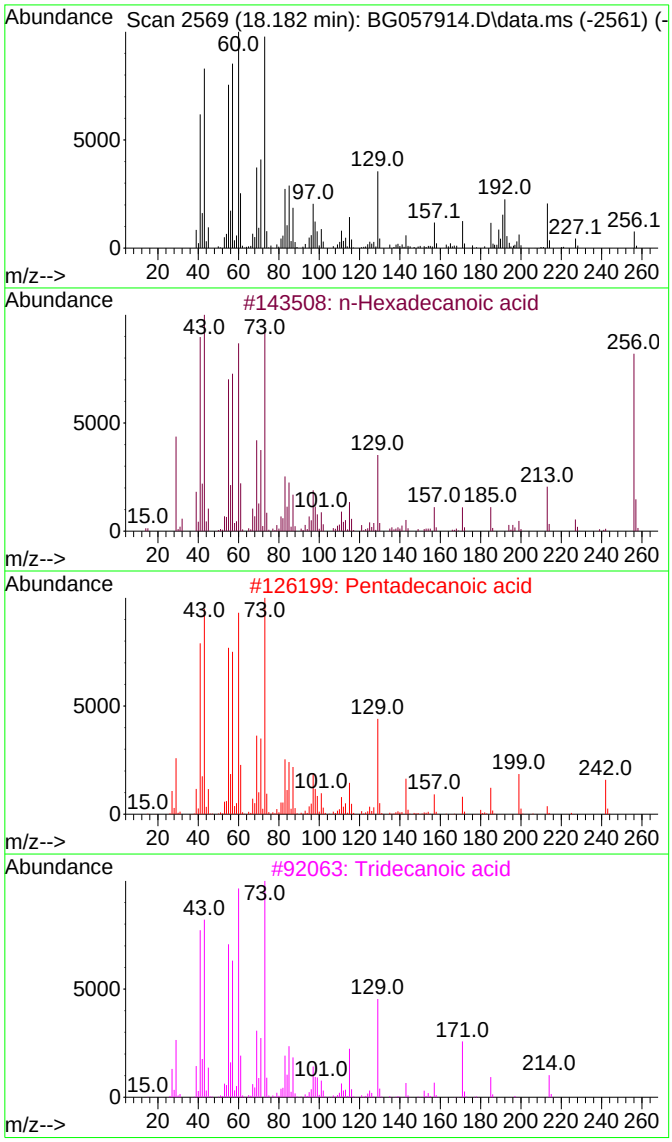
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG062723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

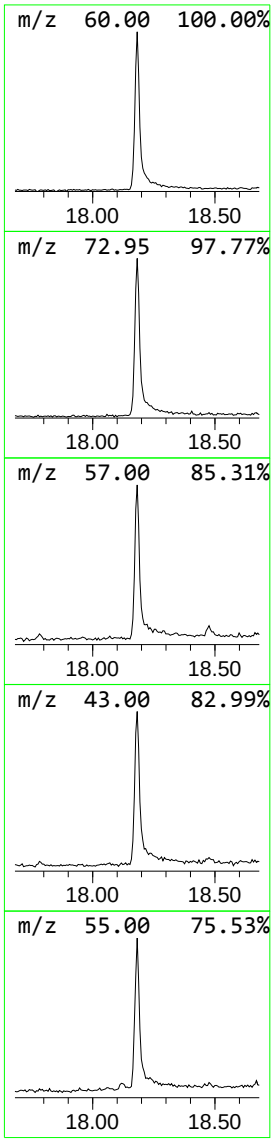
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.182	12.95 ng	542224	Phenanthrene-d10	17.294

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	90
3	Tridecanoic acid	214	C13H26O2	000638-53-9	86
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	72
5	n-Decanoic acid	172	C10H20O2	000334-48-5	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062923\
Data File : BG057914.D
Acq On : 29 Jun 2023 21:38
Operator : CG/JU
Sample : 03434-04
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
350

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG062723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

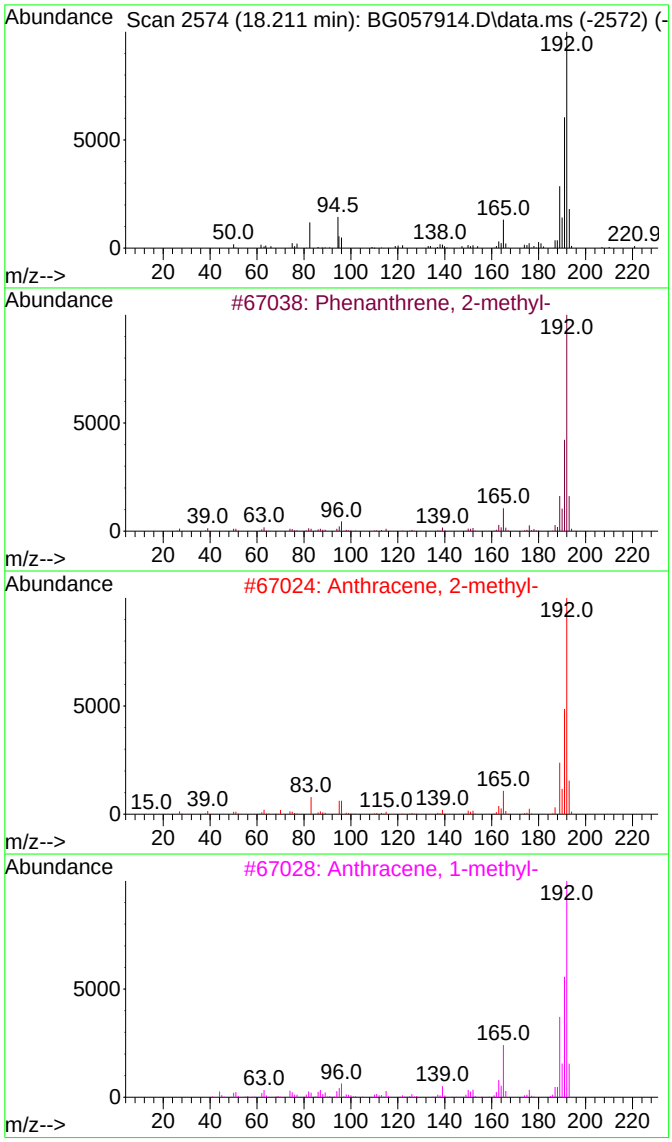
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

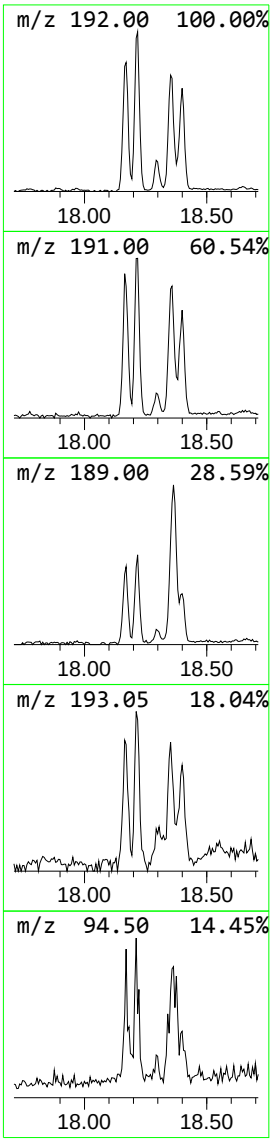
****

Peak Number 4 Phenanthrene, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.211	3.53 ng	147788	Phenanthrene-d10	17.294

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
2	Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
3	Anthracene, 1-methyl-	192	C15H12	000610-48-0	93
4	Anthracene, 9-methyl-	192	C15H12	000779-02-2	91
5	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	90





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062923\
Data File : BG057914.D
Acq On : 29 Jun 2023 21:38
Operator : CG/JU
Sample : 03434-04
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
350

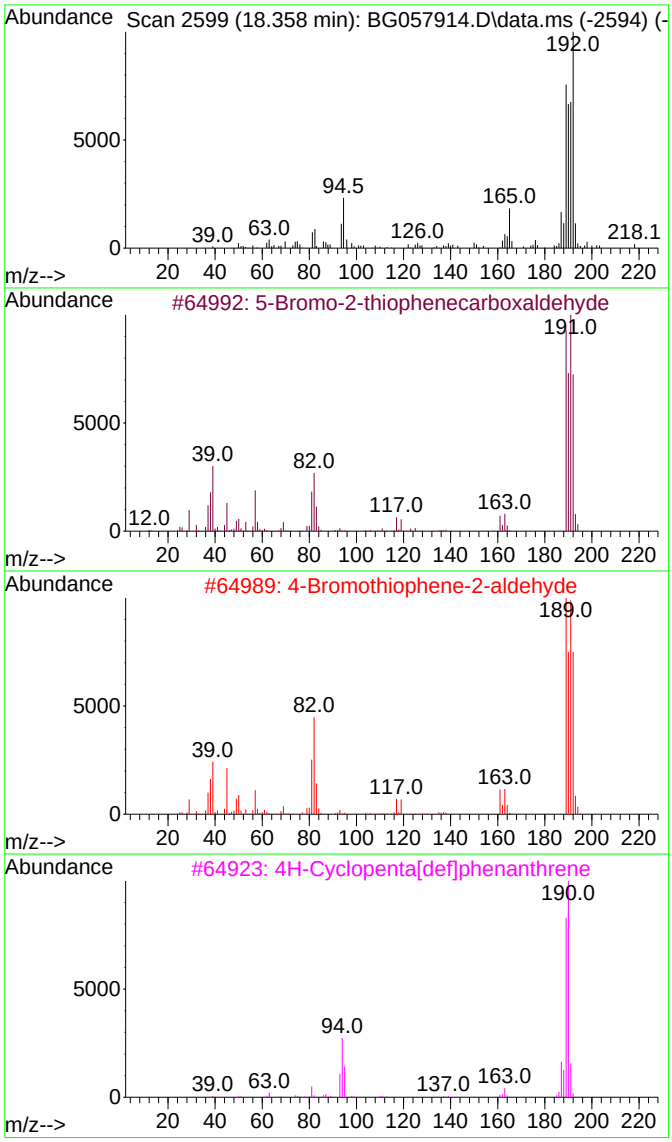
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG062723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

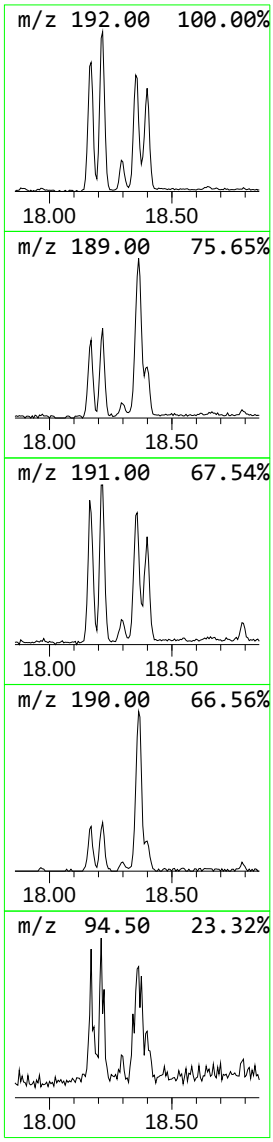
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 5-Bromo-2-thiophenecarboxal... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.358	3.77 ng	157653	Phenanthrene-d10	17.294

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Bromo-2-thiophenecarboxaldehyde	190	C5H3BrOS	004701-17-1	58
2	4-Bromothiophene-2-aldehyde	190	C5H3BrOS	018791-75-8	53
3	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	46
4	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	46
5	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	45





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062923\
Data File : BG057914.D
Acq On : 29 Jun 2023 21:38
Operator : CG/JU
Sample : 03434-04
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
350

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG062723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

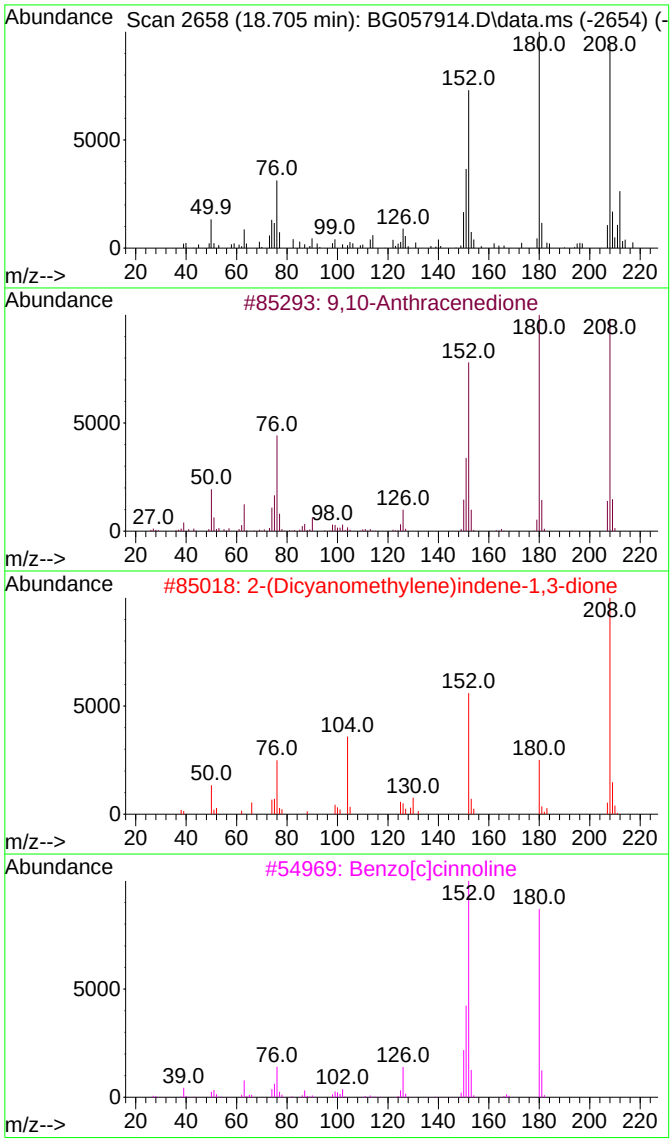
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

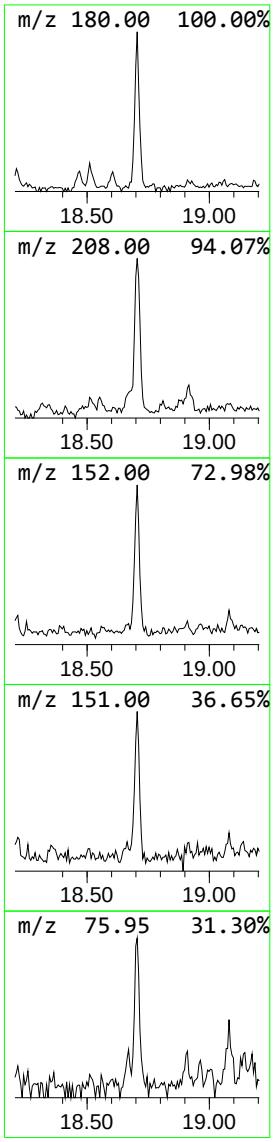
****

Peak Number 6 9,10-Anthracenedione Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.705	2.31 ng	96897	Phenanthrene-d10	17.294

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Anthracenedione	208	C14H8O2	000084-65-1	98
2	2-(Dicyanomethylene)indene-1,3-d...	208	C12H4N2O2	016954-74-8	78
3	Benzo[c]cinnoline	180	C12H8N2	000230-17-1	55
4	1H-Phenalen-1-one	180	C13H8O	000548-39-0	50
5	Benzo[h]cinnoline	180	C12H8N2	000230-31-9	49





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062923\
Data File : BG057914.D
Acq On : 29 Jun 2023 21:38
Operator : CG/JU
Sample : 03434-04
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
350

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG062723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

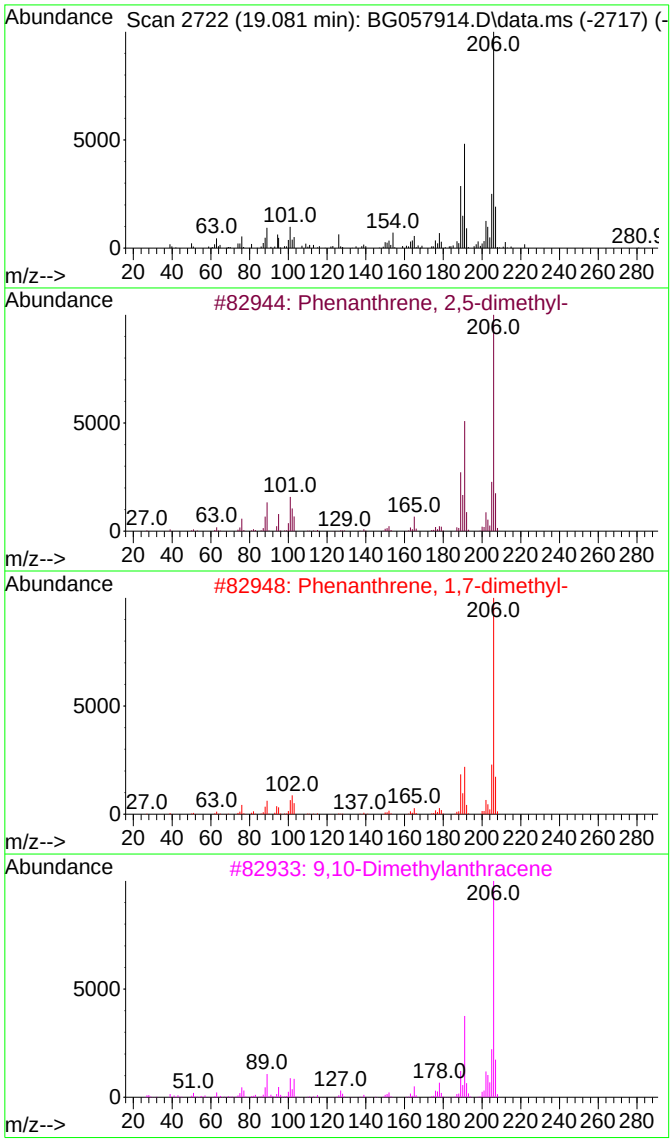
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

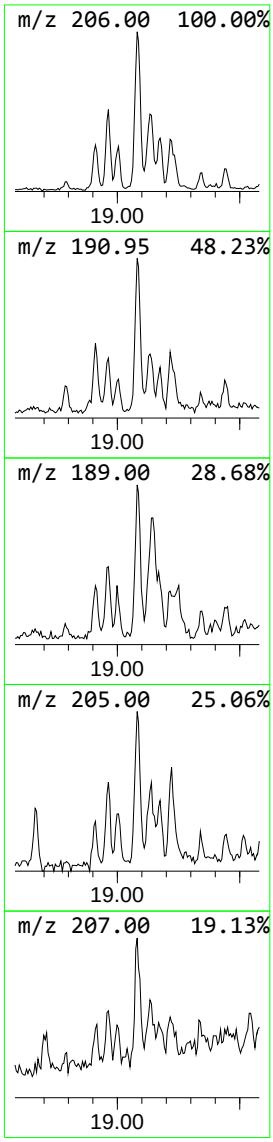
****

Peak Number 7 Phenanthrene, 2,5-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.081	3.46 ng	145053	Phenanthrene-d10	17.294

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	94
2	Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	94
3	9,10-Dimethylanthracene	206	C16H14	000781-43-1	93
4	Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	93
5	1,3-Diphenyl-3-methylcyclopropene	206	C16H14	086544-79-8	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062923\
Data File : BG057914.D
Acq On : 29 Jun 2023 21:38
Operator : CG/JU
Sample : 03434-04
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
350

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG062723.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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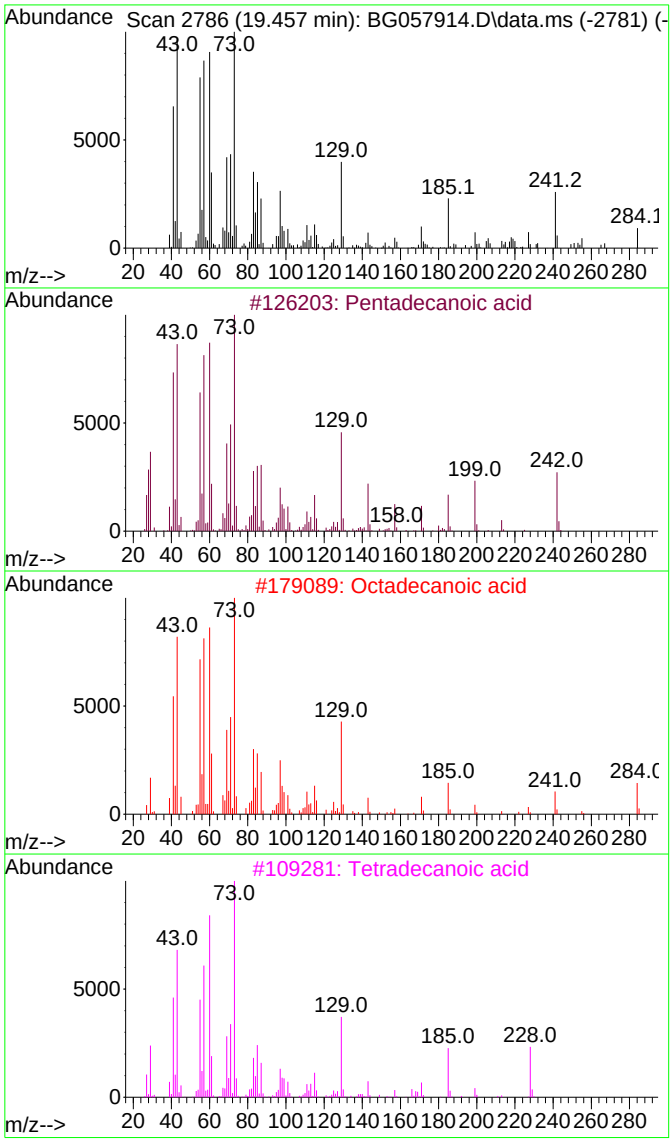
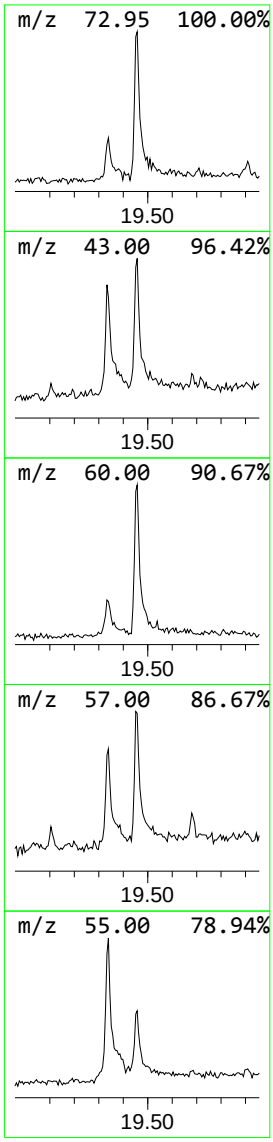
Peak Number 8 Pentadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.457	2.91 ng	186148	Chrysene-d12	21.543

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecanoic acid	242	C15H30O2	001002-84-2	93
2	Octadecanoic acid	284	C18H36O2	000057-11-4	86
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	80
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	64
5	n-Decanoic acid	172	C10H20O2	000334-48-5	62





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Data File : BG057914.D
Acq On : 29 Jun 2023 21:38
Operator : CG/JU
Sample : 03434-04
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
350

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG062723.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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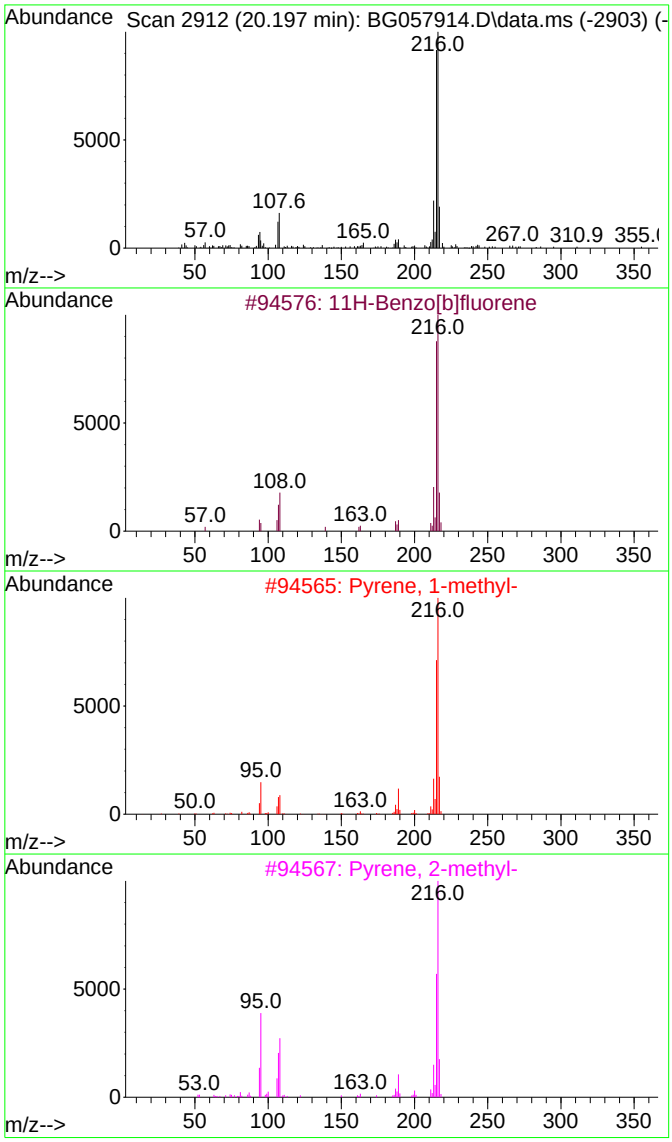
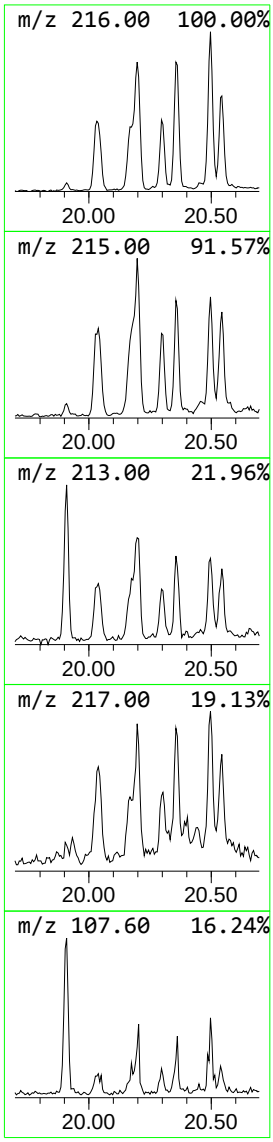
Peak Number 9 11H-Benzo[b]fluorene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.197	3.57 ng	228309	Chrysene-d12	21.543

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	91
2	Pyrene, 1-methyl-	216	C17H12	002381-21-7	87
3	Pyrene, 2-methyl-	216	C17H12	003442-78-2	74
4	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	74
5	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	72





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Data File : BG057914.D
Acq On : 29 Jun 2023 21:38
Operator : CG/JU
Sample : 03434-04
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
350

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG062723.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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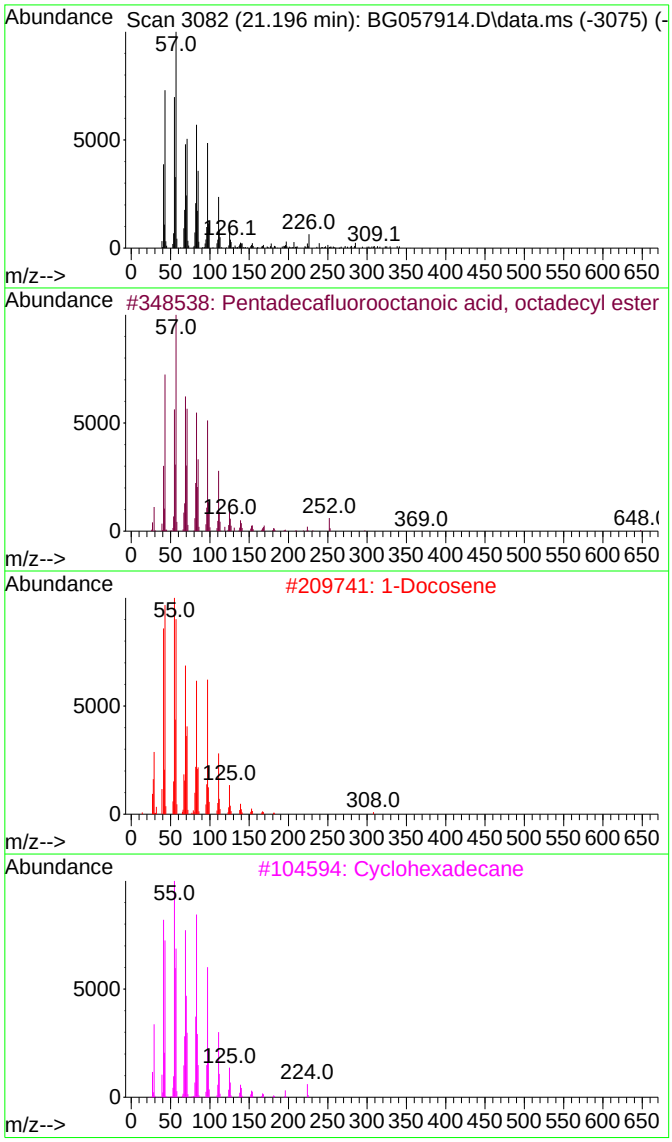
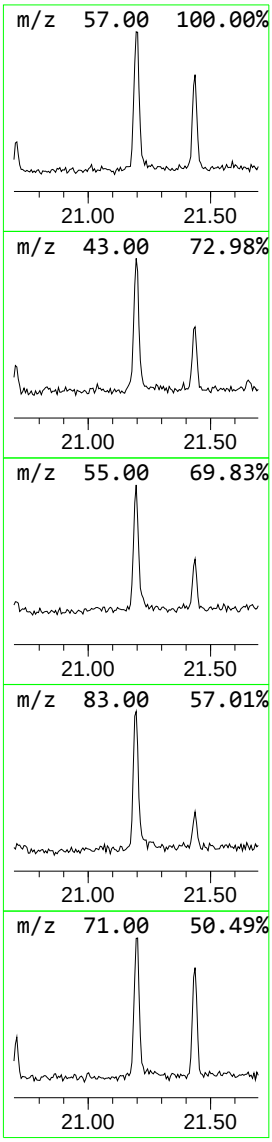
Peak Number 10 Pentadecafluorooctanoic aci... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.196	6.23 ng	398183	Chrysene-d12	21.543

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	94
2	1-Docosene	308	C22H44	001599-67-3	93
3	Cyclohexadecane	224	C16H32	000295-65-8	93
4	Pentafluoropropionic acid, hexad...	388	C19H33F5O2	006222-07-7	91
5	Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062923\
Data File : BG057914.D
Acq On : 29 Jun 2023 21:38
Operator : CG/JU
Sample : 03434-04
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
350

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG062723.M
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TIC Library : C:\Database\NIST20.L
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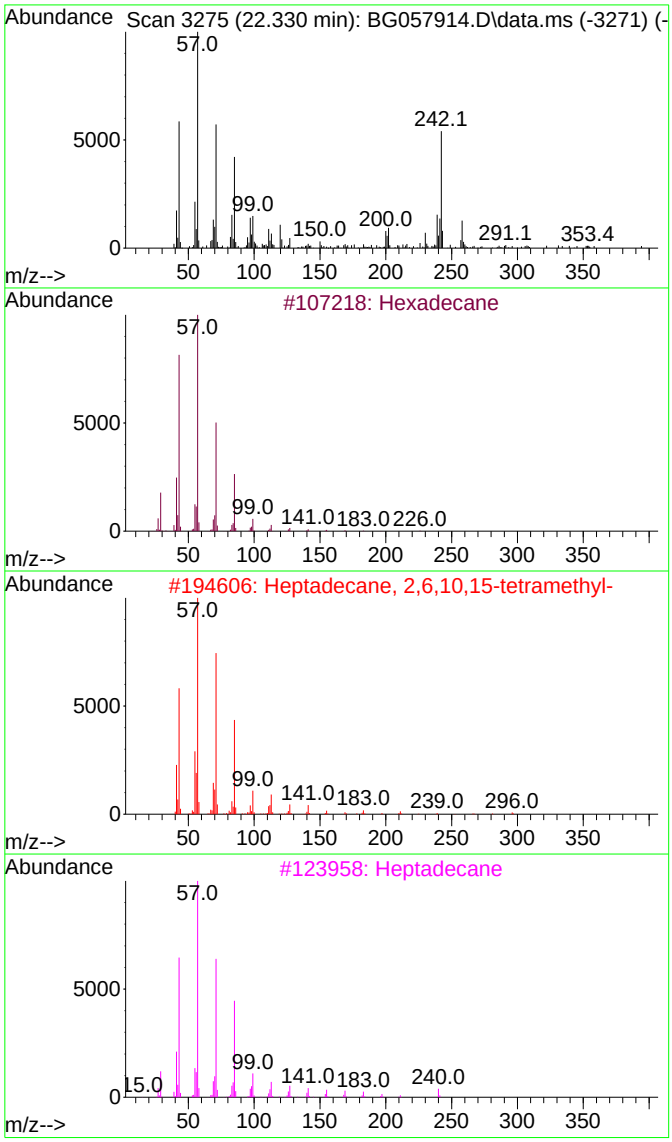
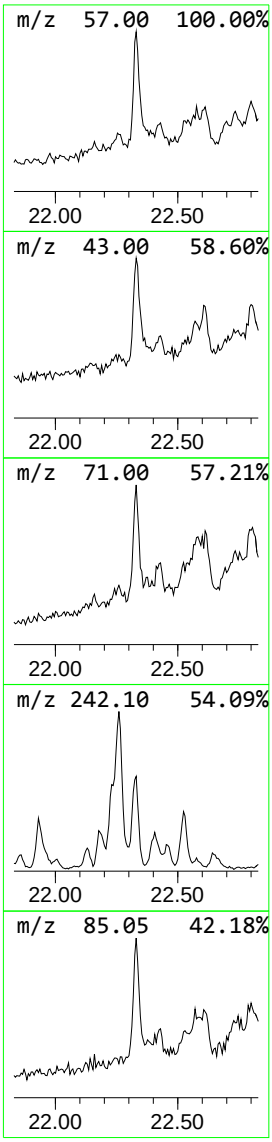
Peak Number 11 Hexadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.330	3.16 ng	202132	Chrysene-d12	21.543

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	89
2	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	56
3	Heptadecane	240	C17H36	000629-78-7	53
4	Heneicosane	296	C21H44	000629-94-7	46
5	Octadecane, 1-iodo-	380	C18H37I	000629-93-6	38





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Operator : CG/JU
Sample : 03434-04
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG062723.M
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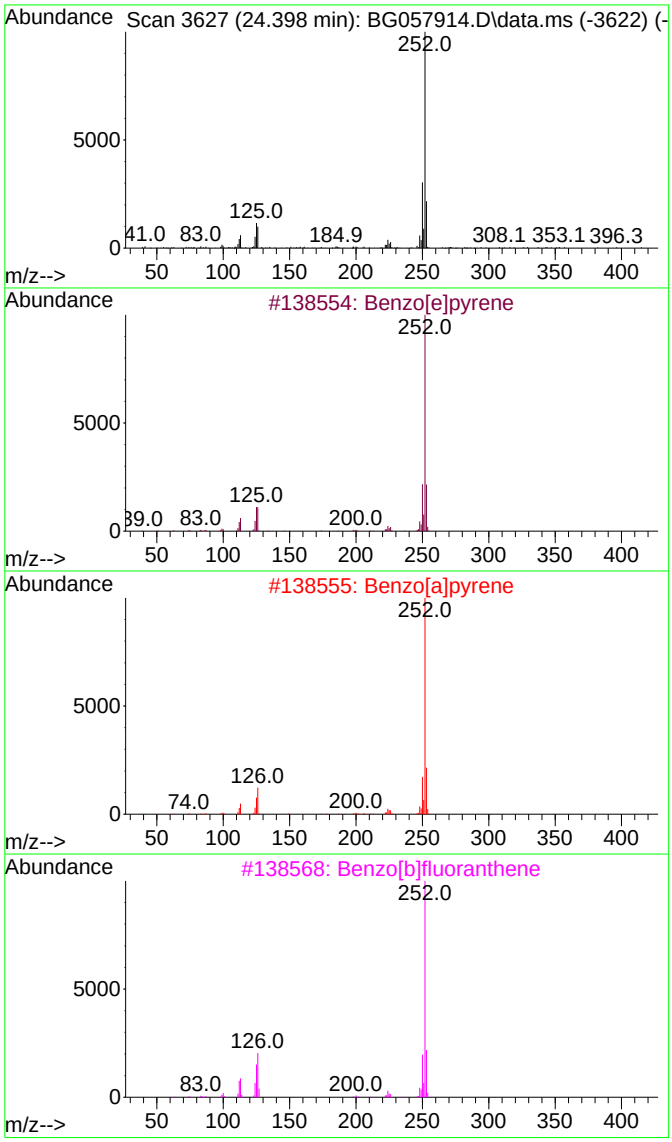
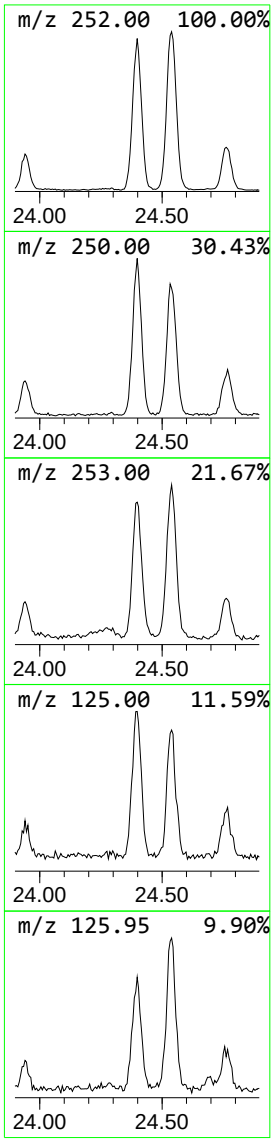
Peak Number 12 Benzo[e]pyrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.398	8.23 ng	363614	Perylene-d12	24.692

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[e]pyrene	252	C20H12	000192-97-2	94
2	Benzo[a]pyrene	252	C20H12	000050-32-8	93
3	Benzo[b]fluoranthene	252	C20H12	000205-99-2	91
4	Benzo[k]fluoranthene	252	C20H12	000207-08-9	91
5	Benzo[j]fluoranthene	252	C20H12	000205-82-3	90





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062923\
Data File : BG057914.D
Acq On : 29 Jun 2023 21:38
Operator : CG/JU
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Misc :
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BNA_G
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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-...	5.021	6.3	ng	103131	1	7.911	328426	20.0
Benzophenone	15.902	5.9	ng	204008	3	14.545	691114	20.0
n-Hexadecanoic ...	18.182	12.9	ng	542224	4	17.294	837442	20.0
Phenanthrene, 2...	18.211	3.5	ng	147788	4	17.294	837442	20.0
5-Bromo-2-thiop...	18.358	3.8	ng	157653	4	17.294	837442	20.0
9,10-Anthracene...	18.705	2.3	ng	96897	4	17.294	837442	20.0
Phenanthrene, 2...	19.081	3.5	ng	145053	4	17.294	837442	20.0
Pentadecanoic acid	19.457	2.9	ng	186148	5	21.543	1278960	20.0
11H-Benzo[b]flu...	20.197	3.6	ng	228309	5	21.543	1278960	20.0
Pentadecafluoro...	21.196	6.2	ng	398183	5	21.543	1278960	20.0
Hexadecane	22.330	3.2	ng	202132	5	21.543	1278960	20.0
Benzo[e]pyrene	24.398	8.2	ng	363614	6	24.692	883991	20.0