LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092325\
Data File : BG064439.D
Acq On : 23 Sep 2025 14:40
Operator : RC/JU
Sample : Q3155-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
COMP-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-BG082825.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG064439.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.434	392	399	410	rVB	311688	486008	30.11%	2.138%
2	7.014	661	668	677	rBV	336470	562871	34.87%	2.476%
3	7.837	800	808	814	rBV	94704	156569	9.70%	0.689%
4	8.988	994	1004	1016	rBV2	206442	370717	22.97%	1.631%
5	10.628	1275	1283	1288	rBV	119343	215622	13.36%	0.949%
6	10.675	1288	1291	1296	rVB3	15162	24634	1.53%	0.108%
7	13.084	1694	1701	1708	rBV	559727	838403	51.94%	3.688%
8	14.464	1926	1936	1941	rBV	198111	307391	19.04%	1.352%
9	14.523	1941	1946	1953	rVB	49862	78041	4.83%	0.343%
10	14.858	1996	2003	2013	rBV	40081	66105	4.10%	0.291%
11	15.510	2109	2114	2119	rBV	50647	75381	4.67%	0.332%
12	15.816	2159	2166	2173	rVB2	72869	120534	7.47%	0.530%
13	15.951	2183	2189	2197	rBV	550093	816047	50.56%	3.590%
14	16.856	2337	2343	2350	rVB	51854	76857	4.76%	0.338%
15	17.020	2367	2371	2378	rVB2	40885	61866	3.83%	0.272%
16	17.108	2381	2386	2392	rVV3	18790	30091	1.86%	0.132%
17	17.202	2394	2402	2405	rVV2	243582	388252	24.05%	1.708%
18	17.249	2405	2410	2416	rVV	723128	1091111	67.60%	4.800%
19	17.337	2416	2425	2430	rVV	131527	212822	13.18%	0.936%
20	17.384	2430	2433	2438	rVB6	10807	18967	1.18%	0.083%
21	17.602	2461	2470	2476	rBV3	70374	127211	7.88%	0.560%
22	17.719	2486	2490	2496	rBV4	14870	22168	1.37%	0.098%
23	17.784	2496	2501	2507	rVB9	12978	22030	1.36%	0.097%
24	17.978	2529	2534	2541	rBV8	21578	48758	3.02%	0.214%
25	18.042	2541	2545	2547	rVV2	26205	35017	2.17%	0.154%
26	18.078	2547	2551	2552	rVV	68970	91753	5.68%	0.404%
27	18.107	2552	2556	2568	rVV3	189690	416596	25.81%	1.833%
28	18.207	2568	2573	2578	rVV2	30666	49526	3.07%	0.218%
29	18.272	2578	2584	2588	rVV3	100616	183122	11.34%	0.806%
30	18.307	2588	2590	2597	rVB	43753	57599	3.57%	0.253%
31	18.401	2599	2606	2615	rBV10	27706	69920	4.33%	0.308%
32	18.577	2631	2636	2639	rVV	59552	81329	5.04%	0.358%
33	18.612	2639	2642	2648	rVB	73987	106060	6.57%	0.467%
34	18.824	2675	2678	2682	rBV4	16147	19216	1.19%	0.085%
35	18.871	2683	2686	2690	rVV3	15205	17425	1.08%	0.077%
36	19.000	2703	2708	2712	rBV3	31314	47762	2.96%	0.210%

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Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
COMP-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-BG082825.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	19.053	2712	2717	2721	rBV7	10777	18524	1.15%	0.081%
38	19.129	2726	2730	2738	rBV	48330	81961	5.08%	0.361%
39	19.259	2745	2752	2760	rBV	1091186	1614155	100.00%	7.101%
40	19.376	2765	2772	2779	rVB5	38568	85259	5.28%	0.375%
41	19.500	2789	2793	2797	rVB	34170	44765	2.77%	0.197%
42	19.623	2809	2814	2821	rVB	820739	1217780	75.44%	5.357%
43	19.688	2821	2825	2830	rVB2	35369	51637	3.20%	0.227%
44	19.823	2838	2848	2852	rBV	1141623	1523461	94.38%	6.702%
45	19.858	2852	2854	2859	rVV	51070	58256	3.61%	0.256%
46	19.946	2862	2869	2875	rBV3	54721	147595	9.14%	0.649%
47	19.999	2875	2878	2882	rVB	16196	21547	1.33%	0.095%
48	20.075	2886	2891	2893	rVV5	37794	67398	4.18%	0.296%
49	20.111	2893	2897	2903	rVB	83367	129738	8.04%	0.571%
50	20.210	2910	2914	2917	rBV	37379	51814	3.21%	0.228%
51	20.269	2918	2924	2928	rVV3	72076	134808	8.35%	0.593%
52	20.304	2928	2930	2935	rVV	28096	34997	2.17%	0.154%
53	20.352	2935	2938	2944	rVB5	18220	27030	1.67%	0.119%
54	20.410	2944	2948	2951	rBV	38581	52833	3.27%	0.232%
55	20.451	2951	2955	2960	rVV4	37443	56177	3.48%	0.247%
56	20.857	3020	3024	3031	rVB	110229	172057	10.66%	0.757%
57	21.033	3047	3054	3058	rBV3	98868	197275	12.22%	0.868%
58	21.074	3058	3061	3064	rVV	80771	120495	7.46%	0.530%
59	21.115	3064	3068	3074	rVV4	133919	285644	17.70%	1.257%
60	21.174	3074	3078	3086	rVB2	125400	199364	12.35%	0.877%
61	21.339	3102	3106	3110	rVB	76508	89609	5.55%	0.394%
62	21.421	3110	3120	3126	rVV3	602603	1453453	90.04%	6.394%
63	21.480	3126	3130	3142	rVB	451590	739945	45.84%	3.255%
64	21.621	3150	3154	3159	rVB2	64965	106790	6.62%	0.470%
65	21.814	3182	3187	3193	rBV2	63767	130317	8.07%	0.573%
66	22.120	3235	3239	3244	rVB2	57099	97675	6.05%	0.430%
67	22.191	3248	3251	3261	rVB6	54996	114044	7.07%	0.502%
68	22.373	3278	3282	3285	rBV3	34120	57295	3.55%	0.252%
69	22.649	3321	3329	3337	rBV	543846	1041747	64.54%	4.583%
70	23.013	3388	3391	3397	rVB2	43904	74974	4.64%	0.330%
71	23.119	3400	3409	3410	rBV4	31348	67505	4.18%	0.297%
72	23.160	3411	3416	3424	rVB7	87633	196704	12.19%	0.865%
73	23.495	3464	3473	3479	rBV2	327706	1031646	63.91%	4.538%
74	23.636	3493	3497	3505	rVB3	115108	235502	14.59%	1.036%
75	23.724	3509	3512	3520	rVB	64257	109823	6.80%	0.483%
76	23.912	3539	3544	3550	rBV8	31287	77660	4.81%	0.342%

LSC Area Percent Report

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Acq On : 23 Sep 2025 14:40
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ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
COMP-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

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Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG082825.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

77	24.159	3579	3586	3595	rVB	183177	491185	30.43%	2.161%
78	24.294	3600	3609	3621	rVB	279756	722354	44.75%	3.178%
79	24.429	3625	3632	3639	rVV2	178793	495638	30.71%	2.180%
80	24.500	3642	3644	3659	rVB2	88406	204962	12.70%	0.902%
81	24.934	3712	3718	3729	rVB7	57392	154417	9.57%	0.679%
82	25.493	3808	3813	3822	rVB6	34141	85364	5.29%	0.376%
83	25.604	3828	3832	3838	rVB9	20212	48990	3.04%	0.216%
84	27.379	4126	4134	4135	rBV3	23413	42672	2.64%	0.188%
85	27.807	4195	4207	4222	rBV4	123712	585403	36.27%	2.575%
86	28.377	4297	4304	4305	rBV4	21924	42668	2.64%	0.188%
87	28.859	4374	4386	4400	rVB2	89427	376137	23.30%	1.655%
88	29.294	4451	4460	4461	rBV7	30445	69748	4.32%	0.307%

Sum of corrected areas: 22732578

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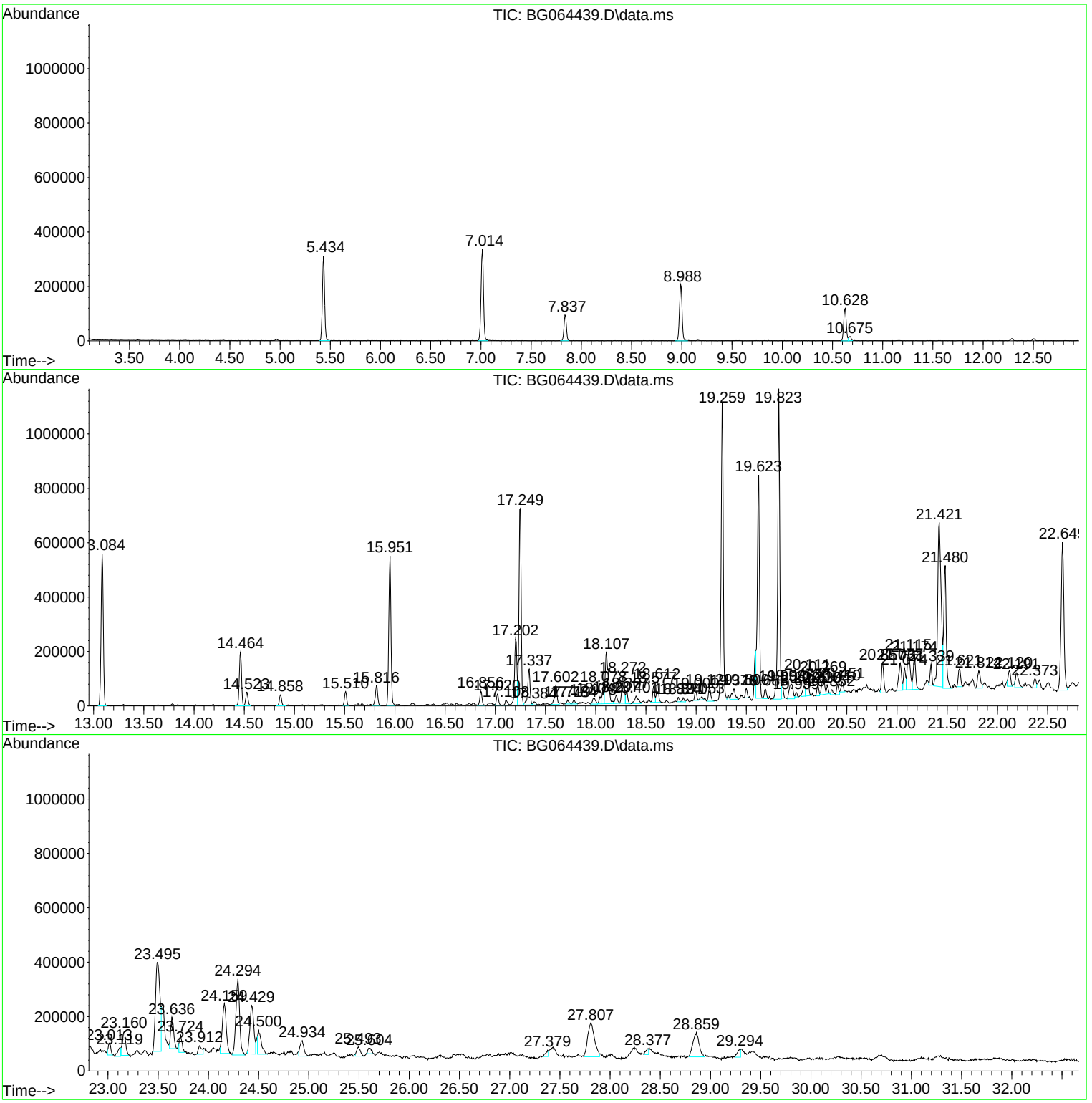
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
COMP-3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG082825.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\BG092325\
Data File : BG064439.D
Acq On : 23 Sep 2025 14:40
Operator : RC/JU
Sample : Q3155-03
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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG082825.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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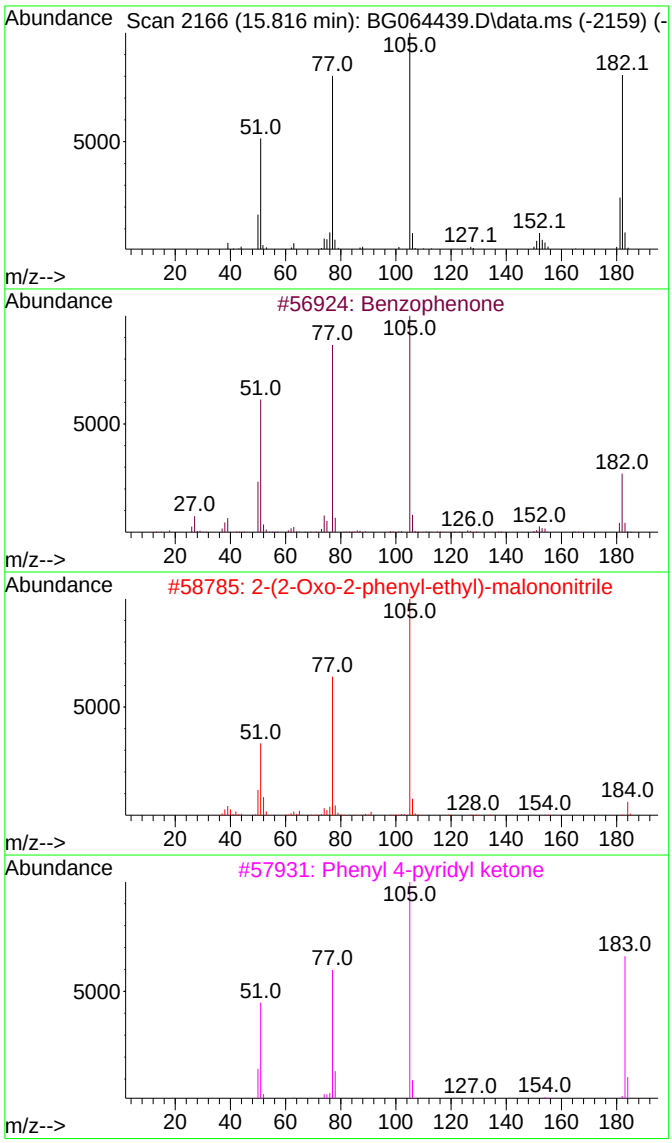
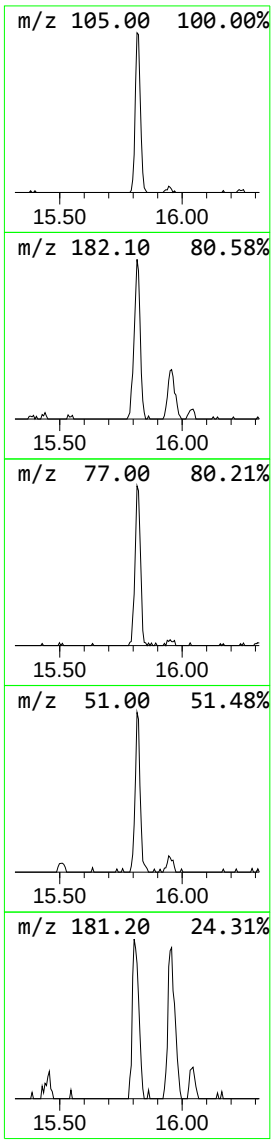
Peak Number 1 Benzophenone Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.816	7.84 ng	120534	Acenaphthene-d10	14.464

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzophenone	182	C13H10O	000119-61-9	93
2	2-(2-Oxo-2-phenyl-ethyl)-malonon...	184	C11H8N2O	1000296-76-9	55
3	Phenyl 4-pyridyl ketone	183	C12H9NO	014548-46-0	49
4	Benzenecarbothioic acid, S-methy...	152	C8H8OS	005925-68-8	49
5	Phenazine, 5,10-dihydro-	182	C12H10N2	000613-32-1	38





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092325\
Data File : BG064439.D
Acq On : 23 Sep 2025 14:40
Operator : RC/JU
Sample : Q3155-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
COMP-3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG082825.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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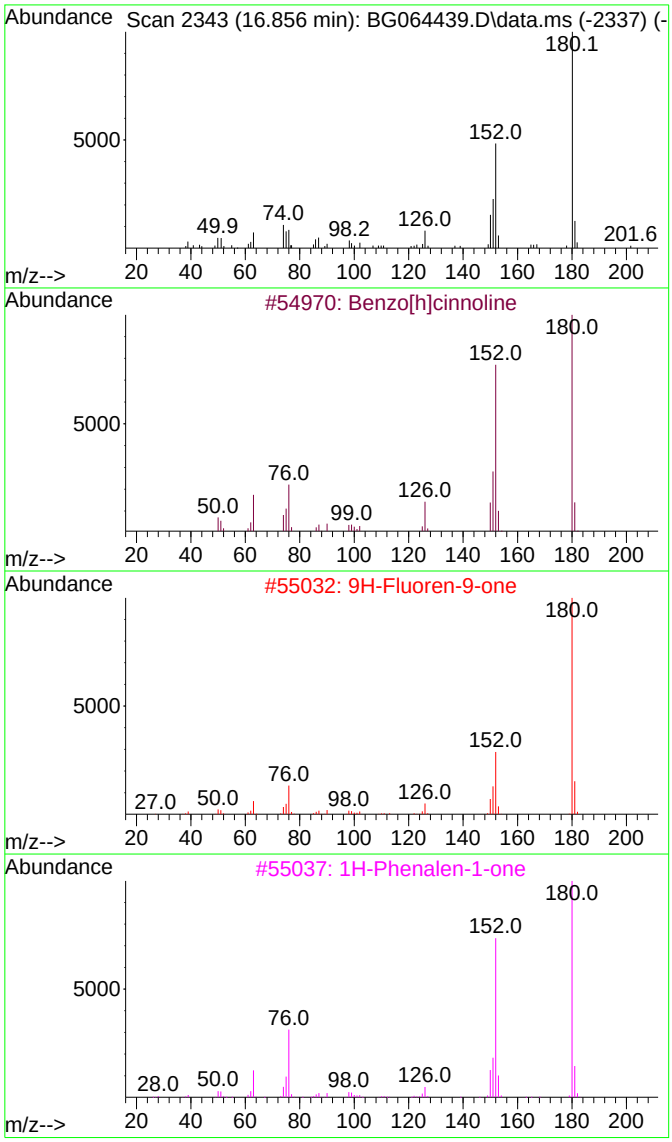
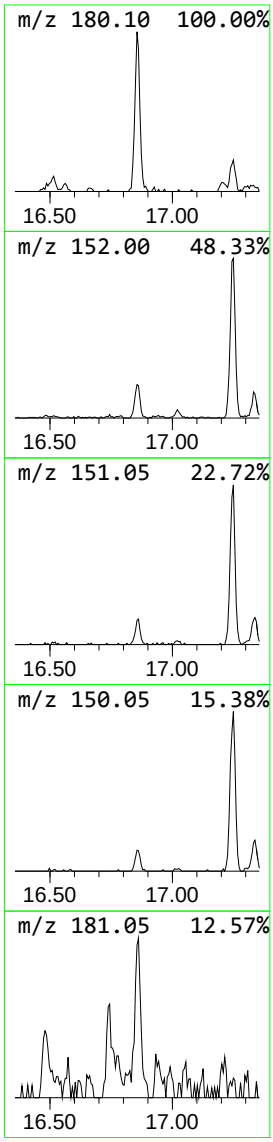
Peak Number 2 Benzo[h]cinnoline Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.856	3.96 ng	76857	Phenanthrene-d10	17.202

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[h]cinnoline	180	C12H8N2	000230-31-9	95
2	9H-Fluoren-9-one	180	C13H8O	000486-25-9	91
3	1H-Phenalen-1-one	180	C13H8O	000548-39-0	64
4	Phenazine	180	C12H8N2	000092-82-0	43
5	2,4-Diazaphenanthrene	180	C12H8N2	000230-28-4	43





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Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG082825.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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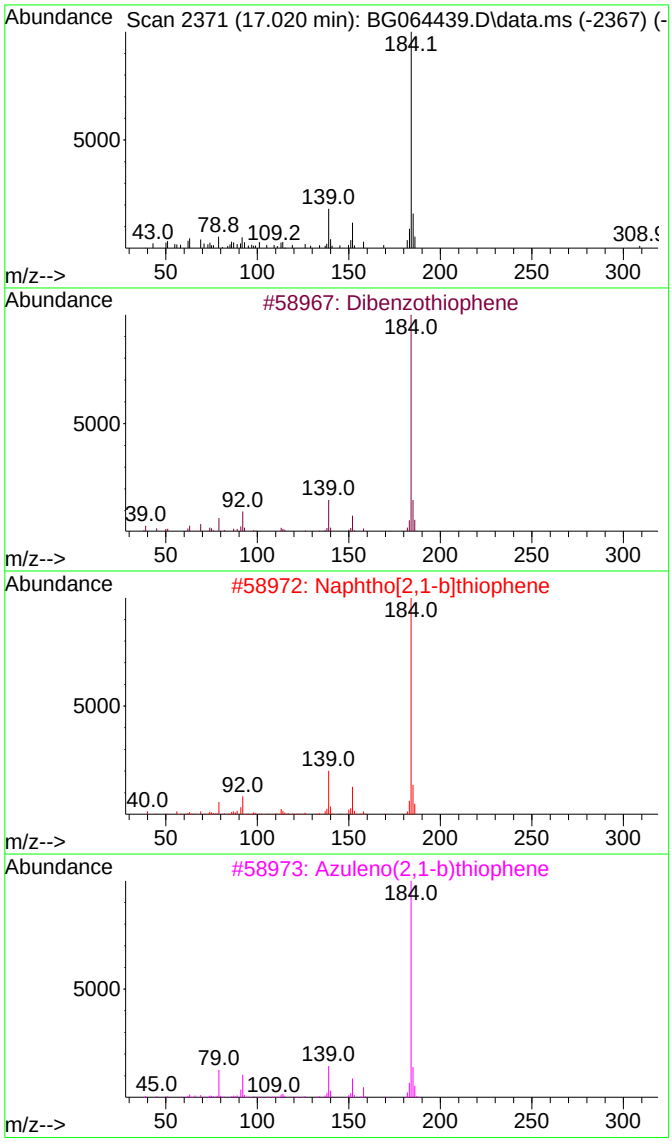
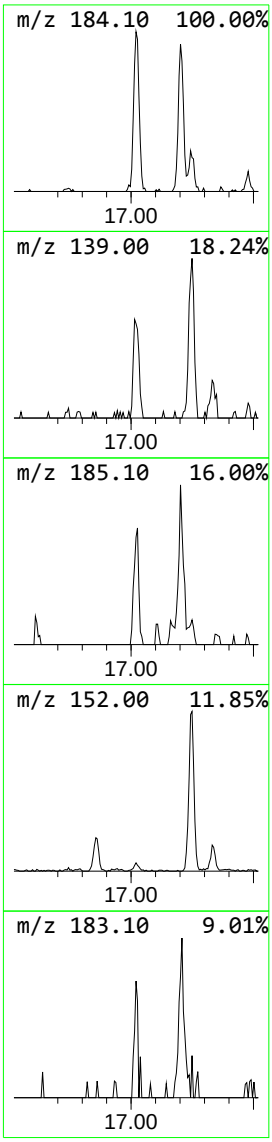
Peak Number 3 Dibenzothiophene Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.020	3.19 ng	61866	Phenanthrene-d10	17.202

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dibenzothiophene	184	C12H8S	000132-65-0	95
2	Naphtho[2,1-b]thiophene	184	C12H8S	000233-02-3	91
3	Azuleno(2,1-b)thiophene	184	C12H8S	000248-13-5	91
4	Naphtho[2,3-b]thiophene	184	C12H8S	000268-77-9	90
5	Naphtho[1,2-b]thiophene	184	C12H8S	000234-41-3	90





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Acq On : 23 Sep 2025 14:40
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Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG082825.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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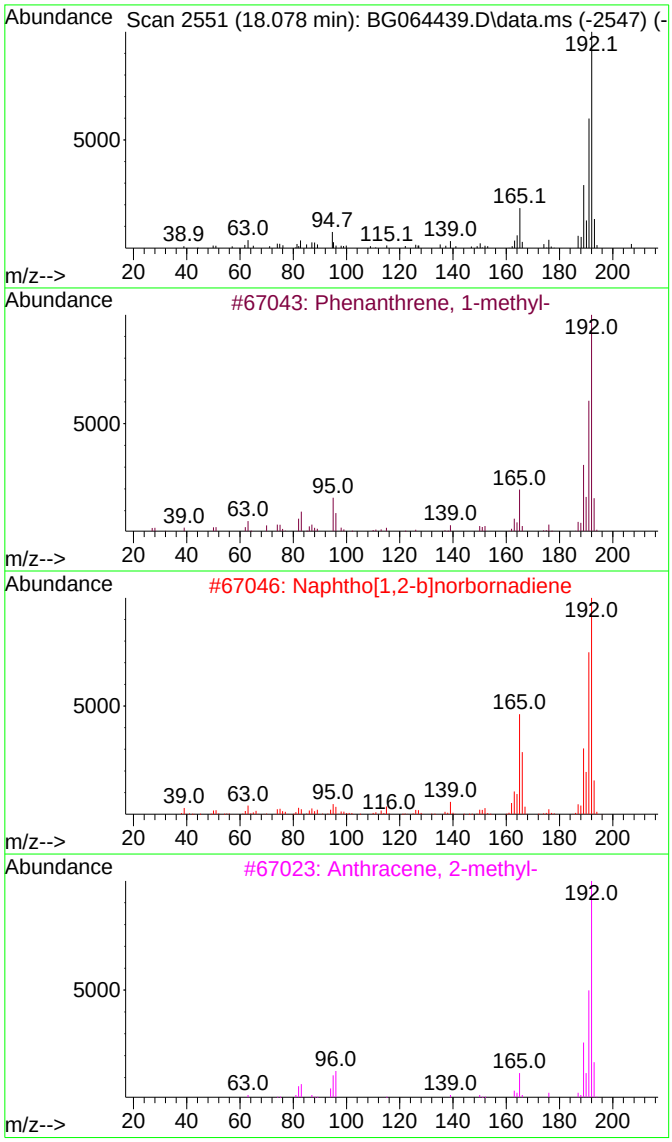
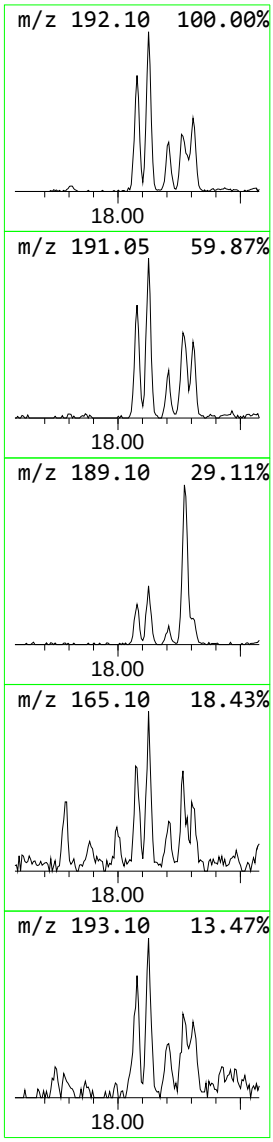
Peak Number 4 Phenanthrene, 1-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.078	4.73 ng	91753	Phenanthrene-d10	17.202

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	87
2	Naphtho[1,2-b]norbornadiene	192	C15H12	1000210-14-8	87
3	Anthracene, 2-methyl-	192	C15H12	000613-12-7	87
4	Anthracene, 1-methyl-	192	C15H12	000610-48-0	87
5	Anthracene, 9-methyl-	192	C15H12	000779-02-2	87





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Data File : BG064439.D
Acq On : 23 Sep 2025 14:40
Operator : RC/JU
Sample : Q3155-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
COMP-3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG082825.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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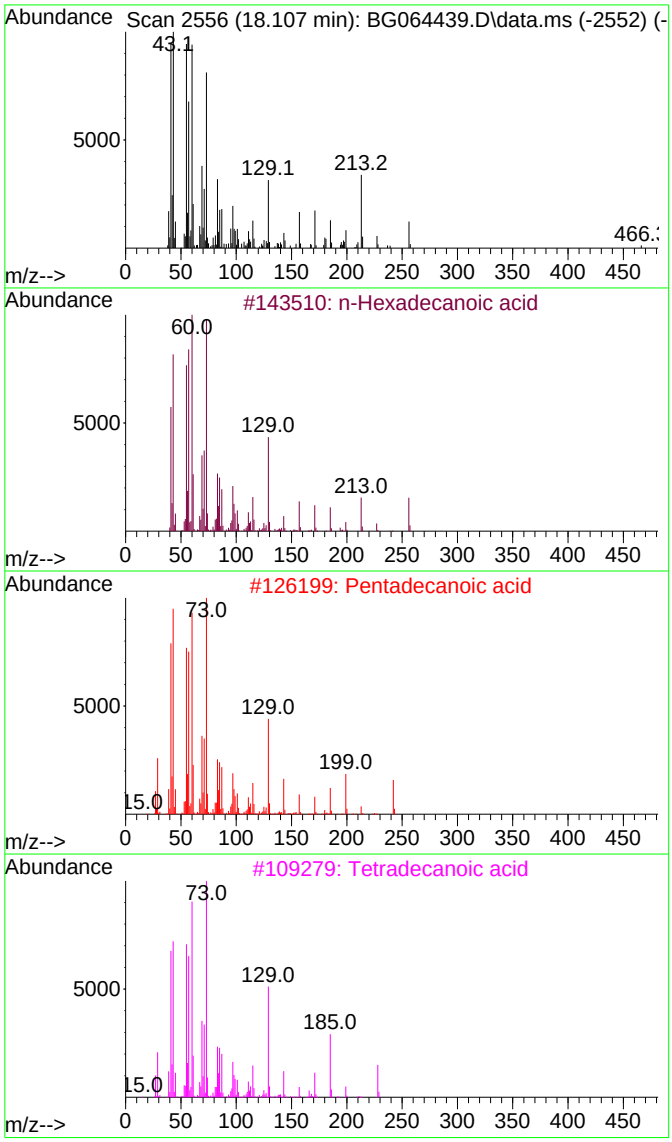
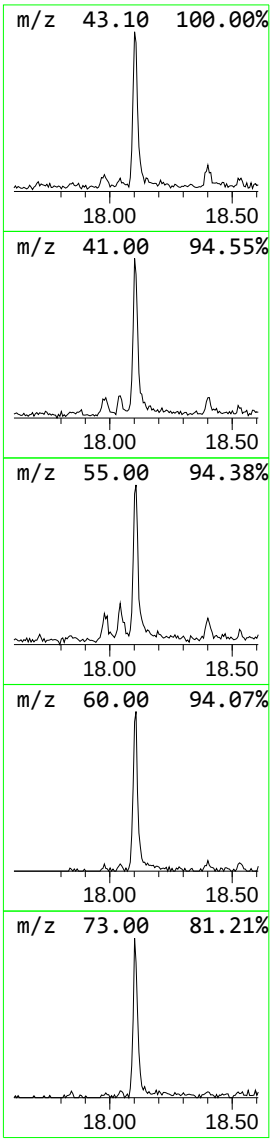
Peak Number 5 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.107	21.46 ng	416596	Phenanthrene-d10	17.202

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	68
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	59
4	Octadecanoic acid	284	C18H36O2	000057-11-4	53
5	Tridecanoic acid	214	C13H26O2	000638-53-9	50





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092325\
Data File : BG064439.D
Acq On : 23 Sep 2025 14:40
Operator : RC/JU
Sample : Q3155-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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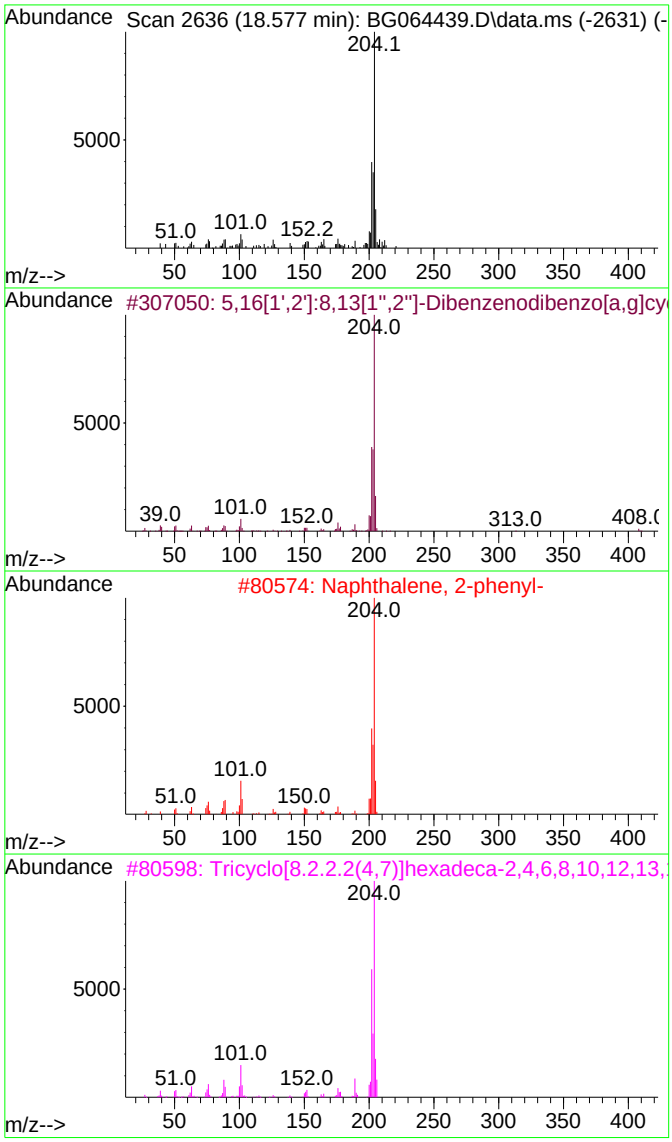
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG082825.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

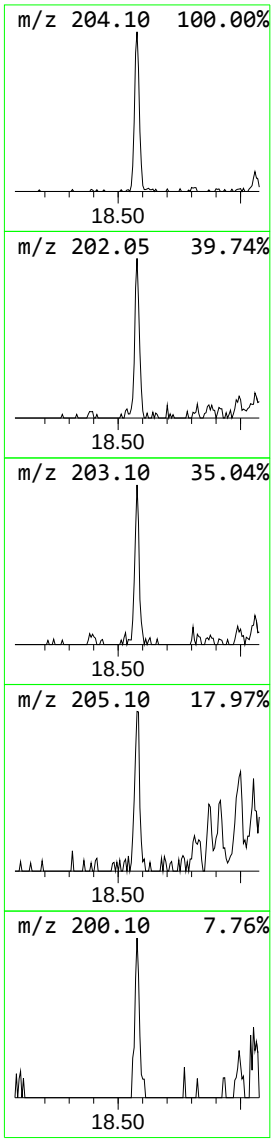
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 5,16[1',2']:8,13[1'',2'']-D... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.577	4.19 ng	81329	Phenanthrene-d10	17.202

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	87
2	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	83
3	Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	58
4	[1,1'-Biphenyl]-2-ol, 5-chloro-	204	C12H9ClO	000607-12-5	49
5	Naphthalene, 1-phenyl-	204	C16H12	000605-02-7	49





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092325\
Data File : BG064439.D
Acq On : 23 Sep 2025 14:40
Operator : RC/JU
Sample : Q3155-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
COMP-3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG082825.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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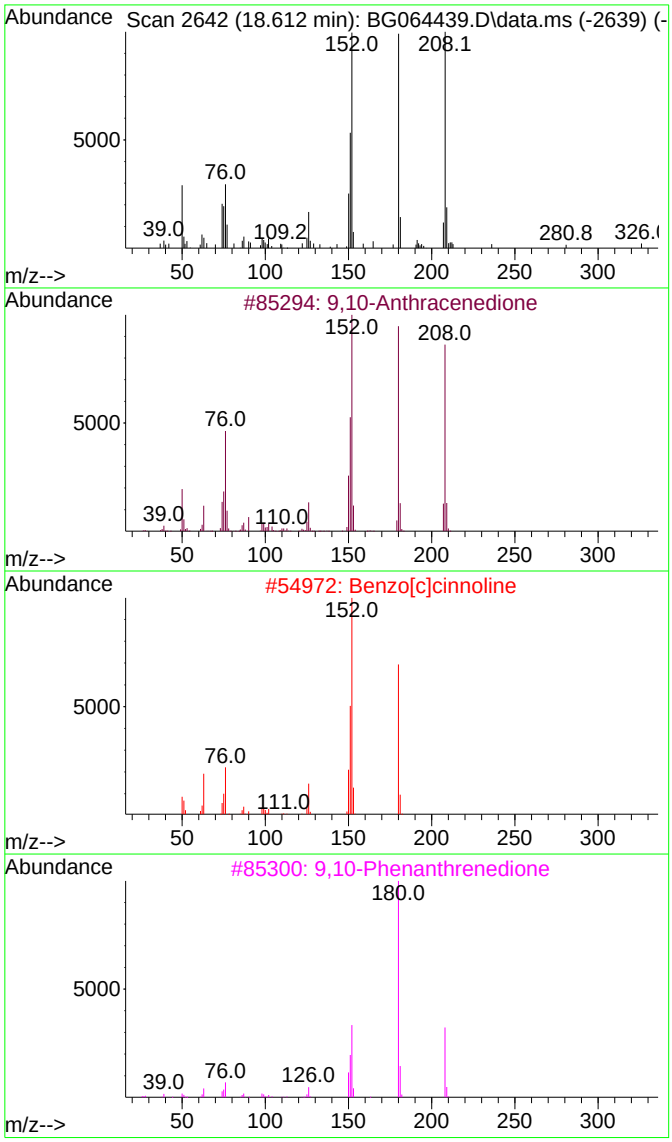
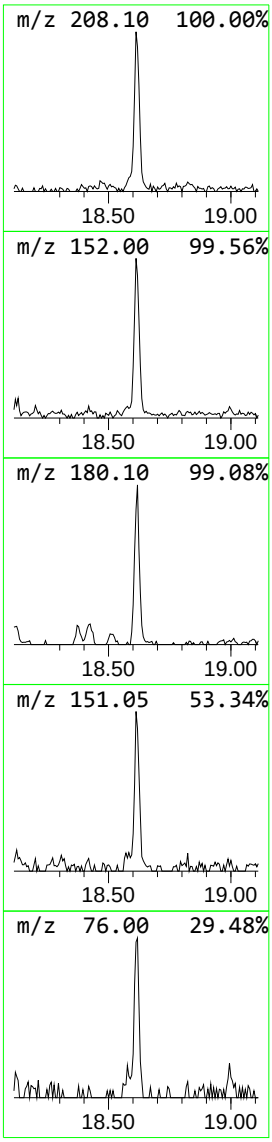
Peak Number 9 9,10-Anthracenedione Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.612	5.46 ng	106060	Phenanthrene-d10	17.202

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Anthracenedione	208	C14H8O2	000084-65-1	93
2	Benzo[c]cinnoline	180	C12H8N2	000230-17-1	86
3	9,10-Phenanthrenedione	208	C14H8O2	000084-11-7	70
4	9H-Fluoren-9-one	180	C13H8O	000486-25-9	64
5	1H-Phenalen-1-one	180	C13H8O	000548-39-0	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092325\
Data File : BG064439.D
Acq On : 23 Sep 2025 14:40
Operator : RC/JU
Sample : Q3155-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
COMP-3

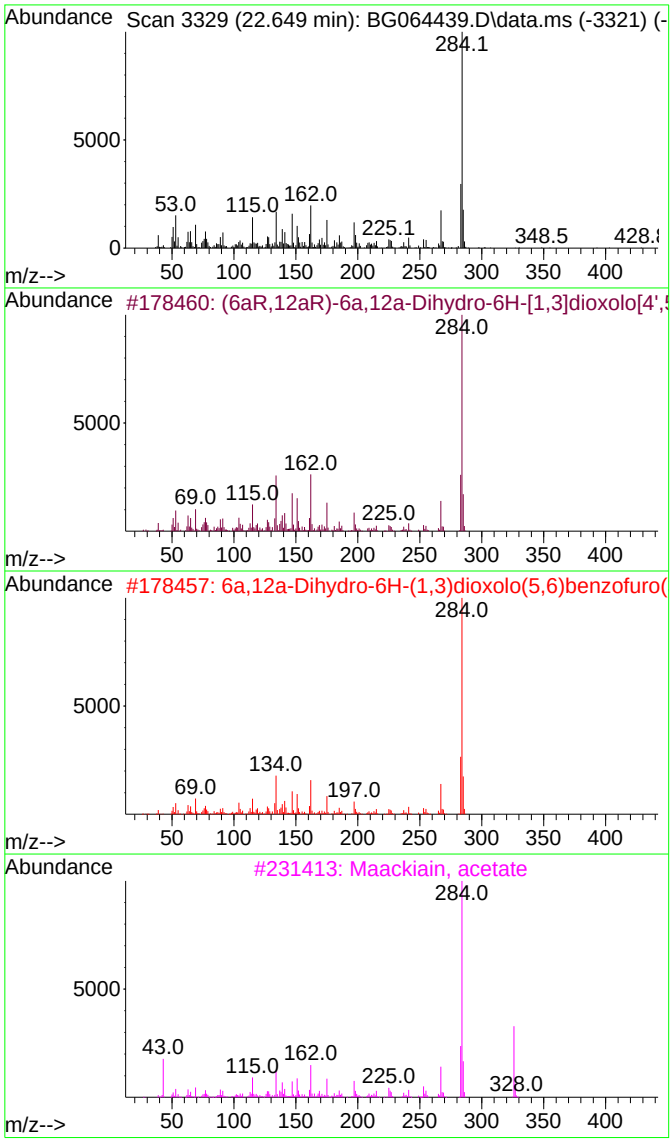
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG082825.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

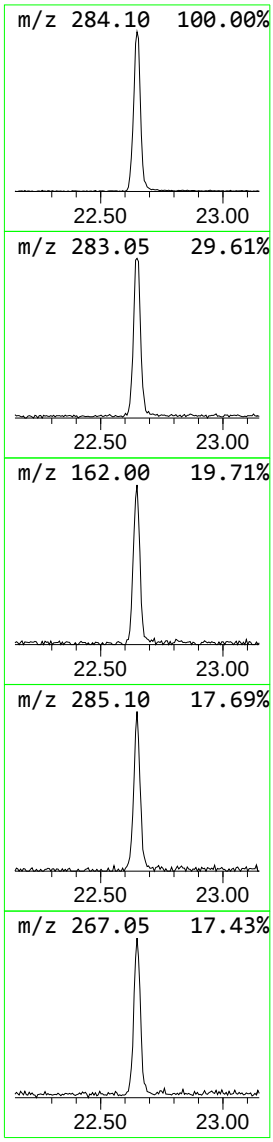
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 (6aR,12aR)-6a,12a-Dihydro-6... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.649	14.33 ng	1041750	Chrysene-d12	21.433

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(6aR,12aR)-6a,12a-Dihydro-6H-[1,...	284	C16H12O5	002035-15-6	96
2	6a,12a-Dihydro-6H-(1,3)dioxolo(5...	284	C16H12O5	076496-57-6	95
3	Maackiain, acetate	326	C18H14O6	002035-16-7	87
4	9,10-Anthracenedione, 1,8-dihydr...	284	C16H12O5	000521-61-9	49
5	3-Hydroxy-2-(4-hydroxy-3-methoxy...	284	C16H12O5	159184-95-9	49





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092325\
Data File : BG064439.D
Acq On : 23 Sep 2025 14:40
Operator : RC/JU
Sample : Q3155-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
COMP-3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG082825.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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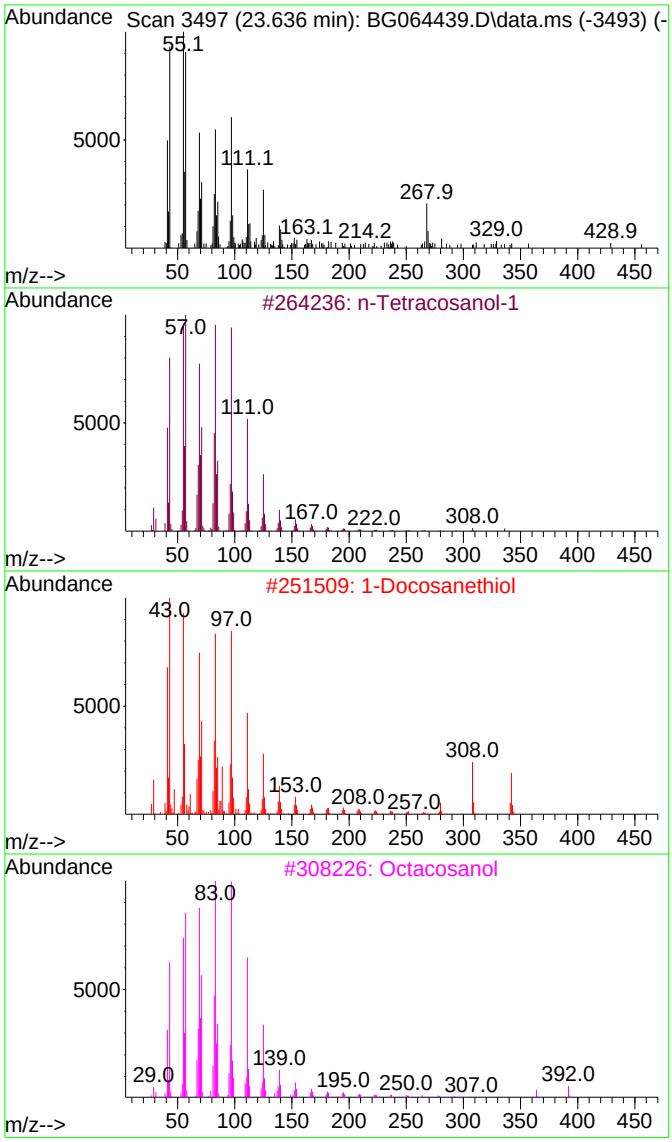
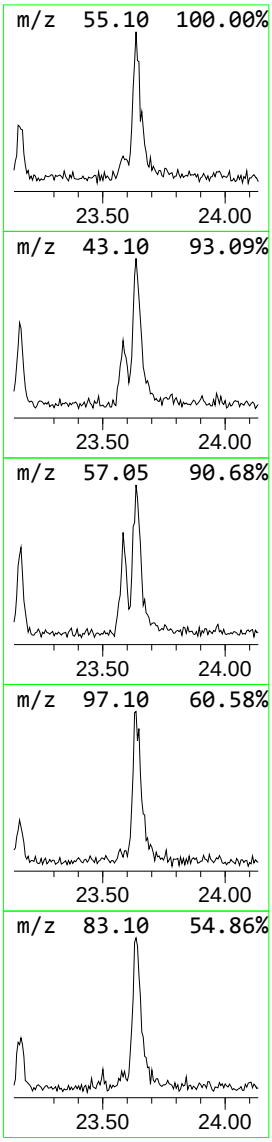
Peak Number 14 n-Tetracosanol-1 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.636	9.50 ng	235502	Perylene-d12	24.435

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Tetracosanol-1	354	C24H50O	000506-51-4	93
2	1-Docosanethiol	342	C22H46S	007773-83-3	89
3	Octacosanol	410	C28H58O	000557-61-9	89
4	1-Heptadecene	238	C17H34	006765-39-5	78
5	Pentafluoropropionic acid, hepta...	402	C20H35F5O2	959218-78-1	76





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092325\
Data File : BG064439.D
Acq On : 23 Sep 2025 14:40
Operator : RC/JU
Sample : Q3155-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
COMP-3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG082825.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
Benzophenone	15.816	7.8	ng	120534	3	14.464	307391	20.0
Benzo[h]cinnoline	16.856	4.0	ng	76857	4	17.202	388252	20.0
Dibenzothiophene	17.020	3.2	ng	61866	4	17.202	388252	20.0
Phenanthrene, 1...	18.078	4.7	ng	91753	4	17.202	388252	20.0
n-Hexadecanoic ...	18.107	21.5	ng	416596	4	17.202	388252	20.0
5,16[1',2']:8,1...	18.577	4.2	ng	81329	4	17.202	388252	20.0
9,10-Anthracene...	18.612	5.5	ng	106060	4	17.202	388252	20.0
(6aR,12aR)-6a,1...	22.649	14.3	ng	1041750	5	21.433	1453450	20.0
n-Tetracosanol-1	23.636	9.5	ng	235502	6	24.435	495638	20.0