LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121025\
Data File : BG064926.D
Acq On : 10 Dec 2025 18:32
Operator : RC/JU
Sample : Q3826-09
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : TP-1

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-BG111225.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG064926.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.123	3	6	15	rVB3	10701	20388	1.58%	0.152%
2	3.199	15	19	25	rBV	25775	41237	3.20%	0.307%
3	5.209	355	361	368	rBV	48479	75762	5.87%	0.565%
4	5.690	436	443	453	rBV	180056	296937	23.03%	2.214%
5	7.300	709	717	731	rBV	167970	314344	24.38%	2.344%
6	8.152	855	862	871	rBV	65075	110810	8.59%	0.826%
7	9.316	1052	1060	1070	rBV	98791	193402	15.00%	1.442%
8	10.967	1333	1341	1350	rBV	77047	150278	11.65%	1.120%
9	13.393	1747	1754	1762	rBV	315575	515358	39.96%	3.843%
10	14.762	1979	1987	1994	rBV	155875	255993	19.85%	1.909%
11	15.802	2158	2164	2172	rVB2	11366	21054	1.63%	0.157%
12	16.102	2209	2215	2222	rBV2	42921	66942	5.19%	0.499%
13	16.237	2231	2238	2252	rBV	418103	680523	52.77%	5.074%
14	16.343	2252	2256	2261	rVB4	30617	42361	3.28%	0.316%
15	17.148	2388	2393	2398	rBV	8266	13164	1.02%	0.098%
16	17.312	2417	2421	2427	rVB2	9884	14795	1.15%	0.110%
17	17.494	2446	2452	2456	rBV	249929	400594	31.06%	2.987%
18	17.535	2456	2459	2466	rVB	232194	330093	25.60%	2.461%
19	17.624	2471	2474	2480	rVB	17823	25865	2.01%	0.193%
20	17.888	2514	2519	2530	rVB	14587	28467	2.21%	0.212%
21	18.070	2547	2550	2559	rVB5	9945	17917	1.39%	0.134%
22	18.305	2583	2590	2594	rBV8	5936	13262	1.03%	0.099%
23	18.364	2594	2600	2604	rVV2	110721	166261	12.89%	1.240%
24	18.411	2604	2608	2616	rVV	67850	110855	8.60%	0.827%
25	18.499	2618	2623	2626	rVV3	9063	14177	1.10%	0.106%
26	18.558	2627	2633	2637	rVV2	58569	123502	9.58%	0.921%
27	18.593	2637	2639	2646	rVB2	36188	46588	3.61%	0.347%
28	18.851	2679	2683	2686	rBV	31000	44144	3.42%	0.329%
29	18.893	2686	2690	2699	rVB2	45599	76550	5.94%	0.571%
30	19.098	2721	2725	2730	rVB3	13066	16529	1.28%	0.123%
31	19.151	2730	2734	2737	rBV	16069	20322	1.58%	0.152%
32	19.186	2737	2740	2748	rVB3	11233	19046	1.48%	0.142%
33	19.269	2749	2754	2759	rBV2	40953	65216	5.06%	0.486%
34	19.322	2759	2763	2767	rVV5	16858	33058	2.56%	0.246%
35	19.363	2767	2770	2773	rVB2	12250	14450	1.12%	0.108%
36	19.404	2773	2777	2788	rBV3	28351	52449	4.07%	0.391%

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121025\
Data File : BG064926.D
Acq On : 10 Dec 2025 18:32
Operator : RC/JU
Sample : Q3826-09
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
TP-1

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-BG111225.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	19.533	2792	2799	2805	rVV	521642	761679	59.06%	5.679%
38	19.586	2805	2808	2811	rVV2	11061	14944	1.16%	0.111%
39	19.621	2811	2814	2819	rVB3	27055	36306	2.82%	0.271%
40	19.721	2826	2831	2834	rBV5	13303	22195	1.72%	0.165%
41	19.774	2835	2840	2845	rVV	17956	25902	2.01%	0.193%
42	19.856	2849	2854	2856	rBV2	77695	119507	9.27%	0.891%
43	19.891	2856	2860	2868	rVB	451775	666685	51.70%	4.971%
44	19.956	2868	2871	2878	rVB3	28726	45523	3.53%	0.339%
45	20.079	2879	2892	2897	rBV	943067	1289591	100.00%	9.615%
46	20.126	2897	2900	2908	rVB2	23018	36974	2.87%	0.276%
47	20.215	2908	2915	2920	rBV3	44159	114620	8.89%	0.855%
48	20.350	2932	2938	2939	rBV4	29977	52473	4.07%	0.391%
49	20.373	2939	2942	2948	rVB	90470	127751	9.91%	0.953%
50	20.479	2954	2960	2963	rBV	52038	74834	5.80%	0.558%
51	20.532	2963	2969	2973	rVV2	64839	102849	7.98%	0.767%
52	20.673	2989	2993	2997	rBV	36961	51495	3.99%	0.384%
53	20.720	2997	3001	3005	rVB	32506	45306	3.51%	0.338%
54	20.832	3016	3020	3023	rVB4	10307	14769	1.15%	0.110%
55	20.926	3033	3036	3039	rBV4	9863	15524	1.20%	0.116%
56	20.967	3039	3043	3045	rVV5	15162	20549	1.59%	0.153%
57	21.090	3062	3064	3067	rVV3	13737	15856	1.23%	0.118%
58	21.131	3067	3071	3078	rVB	51683	86710	6.72%	0.647%
59	21.319	3096	3103	3107	rBV3	56568	125244	9.71%	0.934%
60	21.366	3107	3111	3116	rVV2	89167	159222	12.35%	1.187%
61	21.413	3116	3119	3124	rVV2	56843	92129	7.14%	0.687%
62	21.466	3124	3128	3139	rVB4	65712	151346	11.74%	1.128%
63	21.742	3161	3175	3180	rBV2	410932	1078926	83.66%	8.044%
64	21.789	3180	3183	3191	rVB	236078	386140	29.94%	2.879%
65	21.942	3205	3209	3215	rVB2	34503	64469	5.00%	0.481%
66	22.148	3240	3244	3253	rBV2	33801	75772	5.88%	0.565%
67	22.218	3253	3256	3266	rVB2	16448	41773	3.24%	0.311%
68	22.465	3285	3298	3306	rBV3	86364	323814	25.11%	2.414%
69	22.547	3309	3312	3319	rVB4	48424	98308	7.62%	0.733%
70	22.753	3343	3347	3352	rBV2	23890	46376	3.60%	0.346%
71	22.906	3366	3373	3379	rVB5	19862	45838	3.55%	0.342%
72	23.969	3545	3554	3561	rBV	166636	579506	44.94%	4.321%
73	24.222	3592	3597	3605	rVB2	24997	58106	4.51%	0.433%
74	24.427	3627	3632	3637	rBV8	13551	32317	2.51%	0.241%
75	24.698	3670	3678	3692	rVB2	100321	309331	23.99%	2.306%
76	24.839	3694	3702	3716	rVB	141424	405110	31.41%	3.020%

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121025\
Data File : BG064926.D
Acq On : 10 Dec 2025 18:32
Operator : RC/JU
Sample : Q3826-09
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
TP-1

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-BG111225.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

77	25.003	3719	3730	3738	rBV	200801	615612	47.74%	4.590%
78	28.711	4348	4361	4378	rBV4	62418	319913	24.81%	2.385%
79	29.868	4546	4558	4572	rVB2	49072	228039	17.68%	1.700%

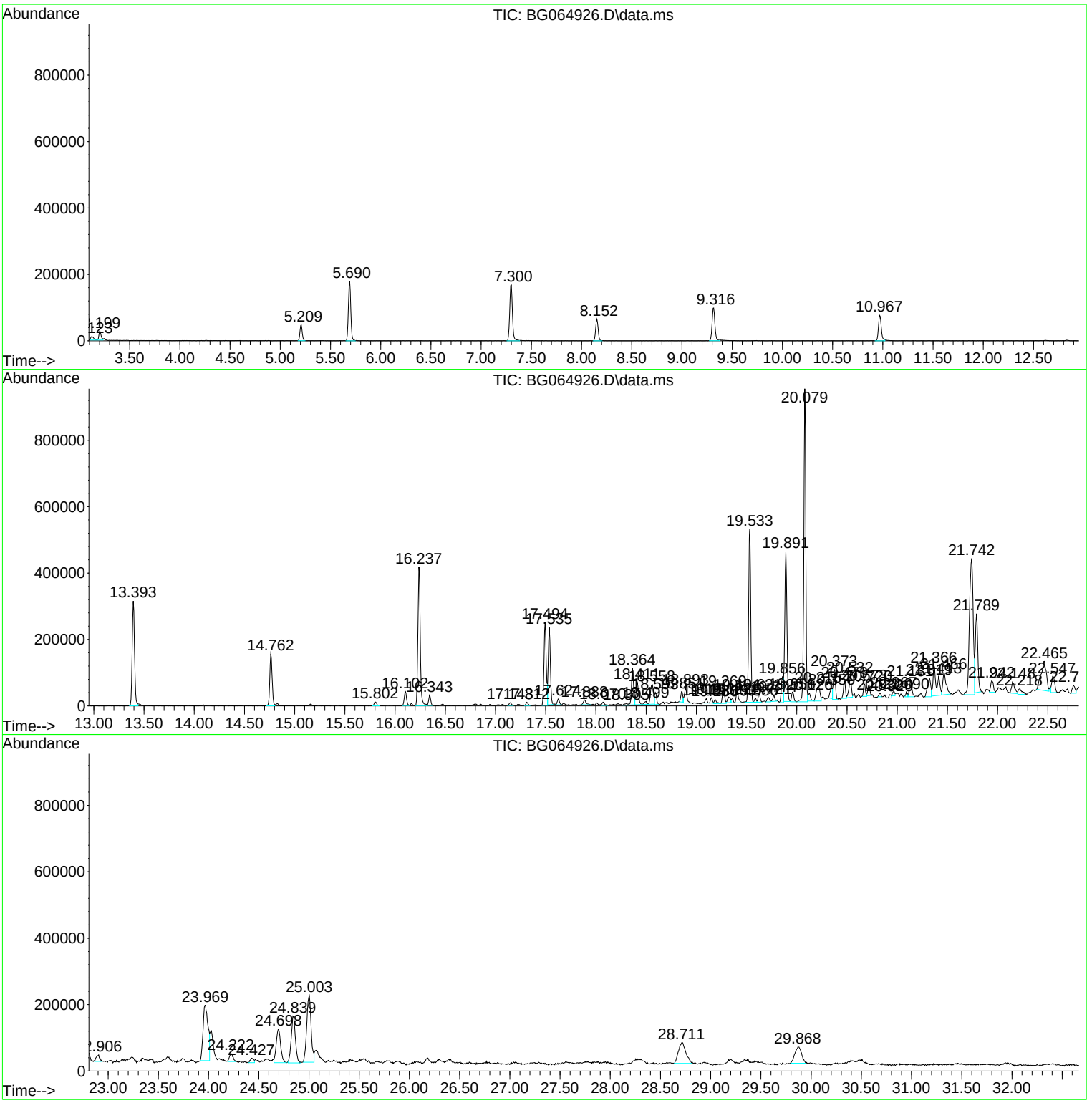
Sum of corrected areas: 13412026

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121025\
Data File : BG064926.D
Acq On : 10 Dec 2025 18:32
Operator : RC/JU
Sample : Q3826-09
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
TP-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111225.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\BG121025\
Data File : BG064926.D
Acq On : 10 Dec 2025 18:32
Operator : RC/JU
Sample : Q3826-09
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
TP-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111225.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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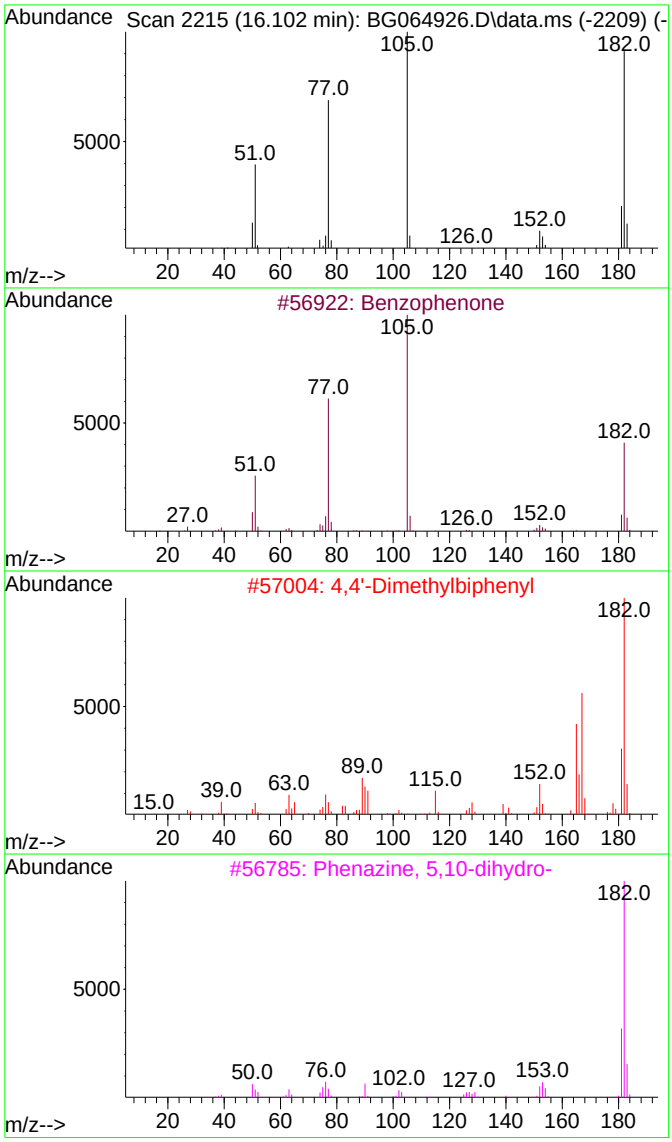
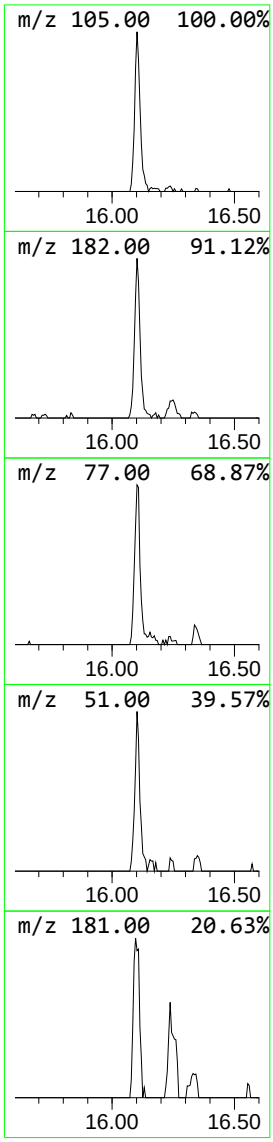
Peak Number 4 Benzophenone Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.102	5.23 ng	66942	Acenaphthene-d10	14.762

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzophenone	182	C13H10O	000119-61-9	95
2	4,4'-Dimethylbiphenyl	182	C14H14	000613-33-2	46
3	Phenazine, 5,10-dihydro-	182	C12H10N2	000613-32-1	43
4	Benzenecarbothioic acid, S-methy...	152	C8H8OS	005925-68-8	43
5	1,1'-Biphenyl, 3,4'-dimethyl-	182	C14H14	007383-90-6	43





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121025\
Data File : BG064926.D
Acq On : 10 Dec 2025 18:32
Operator : RC/JU
Sample : Q3826-09
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
TP-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111225.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

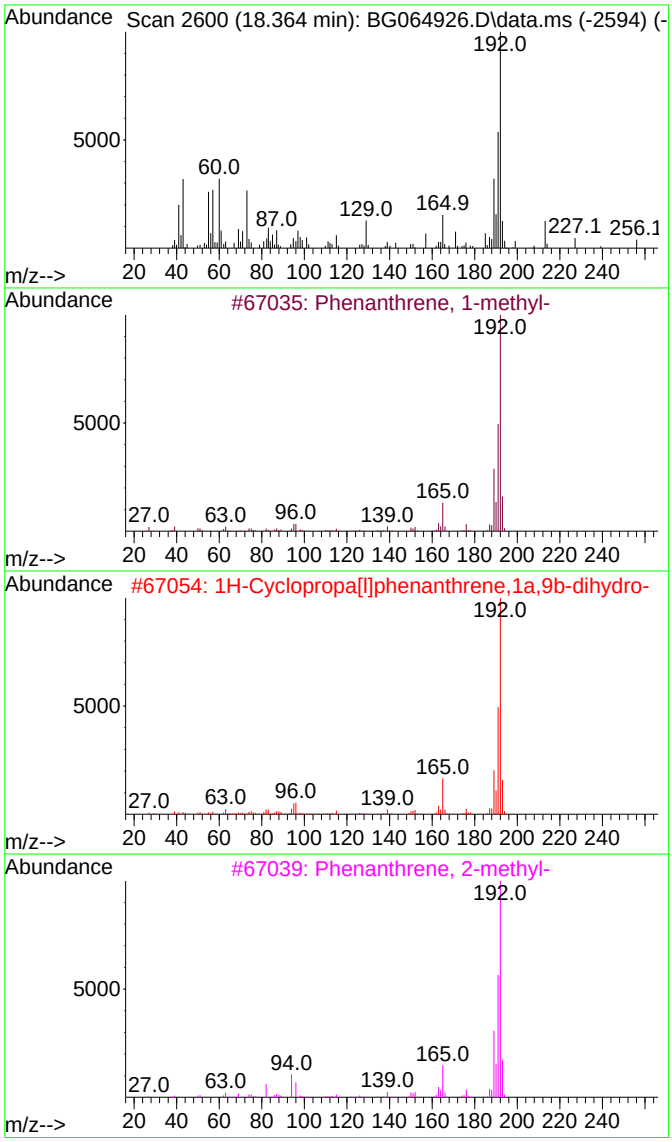
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

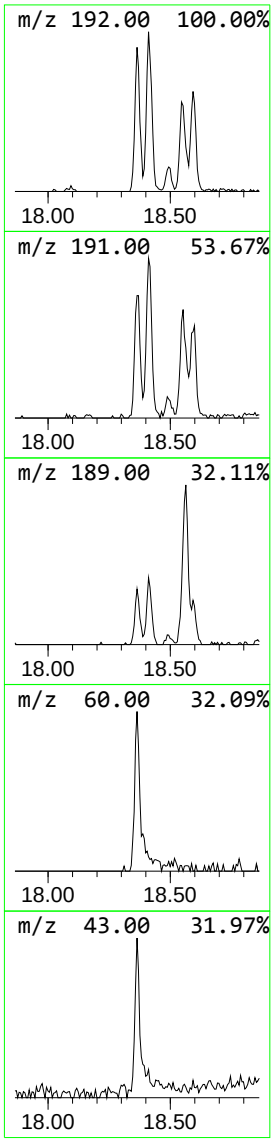
****

Peak Number 6 Phenanthrene, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.364	8.30 ng	166261	Phenanthrene-d10	17.494

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	91
2	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	83
3	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	83
4	Anthracene, 2-methyl-	192	C15H12	000613-12-7	70
5	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121025\
Data File : BG064926.D
Acq On : 10 Dec 2025 18:32
Operator : RC/JU
Sample : Q3826-09
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
TP-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111225.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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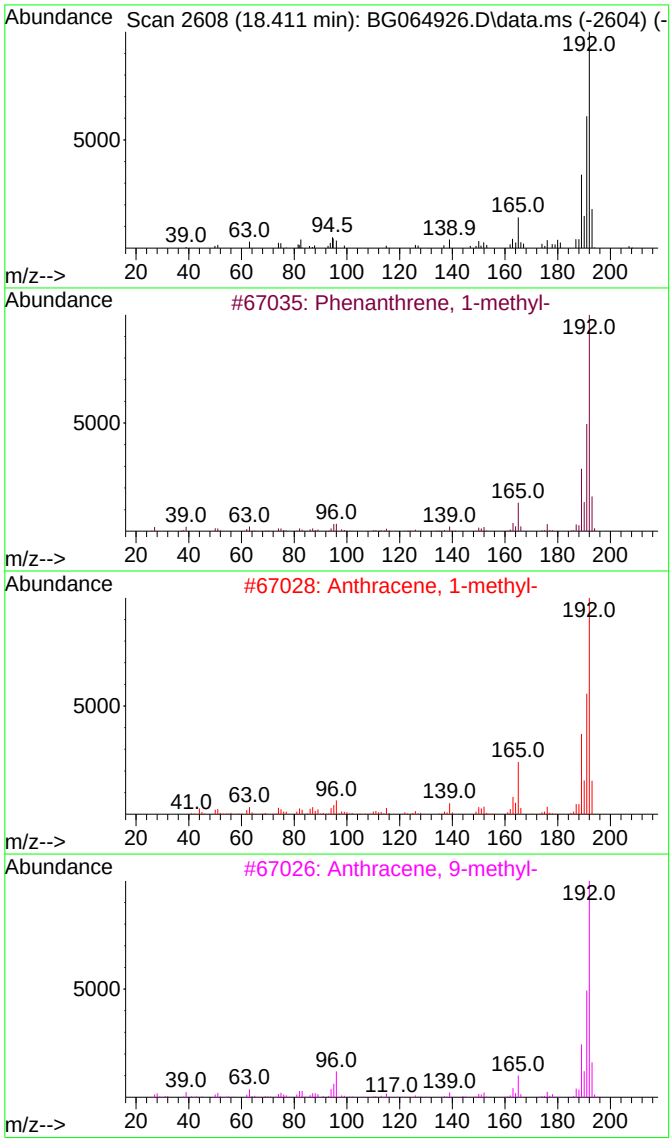
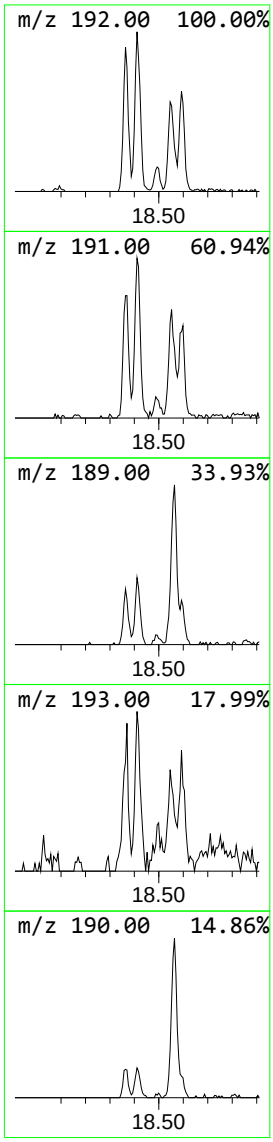
Peak Number 7 Anthracene, 1-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.411	5.53 ng	110855	Phenanthrene-d10	17.494

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





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121025\
Data File : BG064926.D
Acq On : 10 Dec 2025 18:32
Operator : RC/JU
Sample : Q3826-09
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
TP-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111225.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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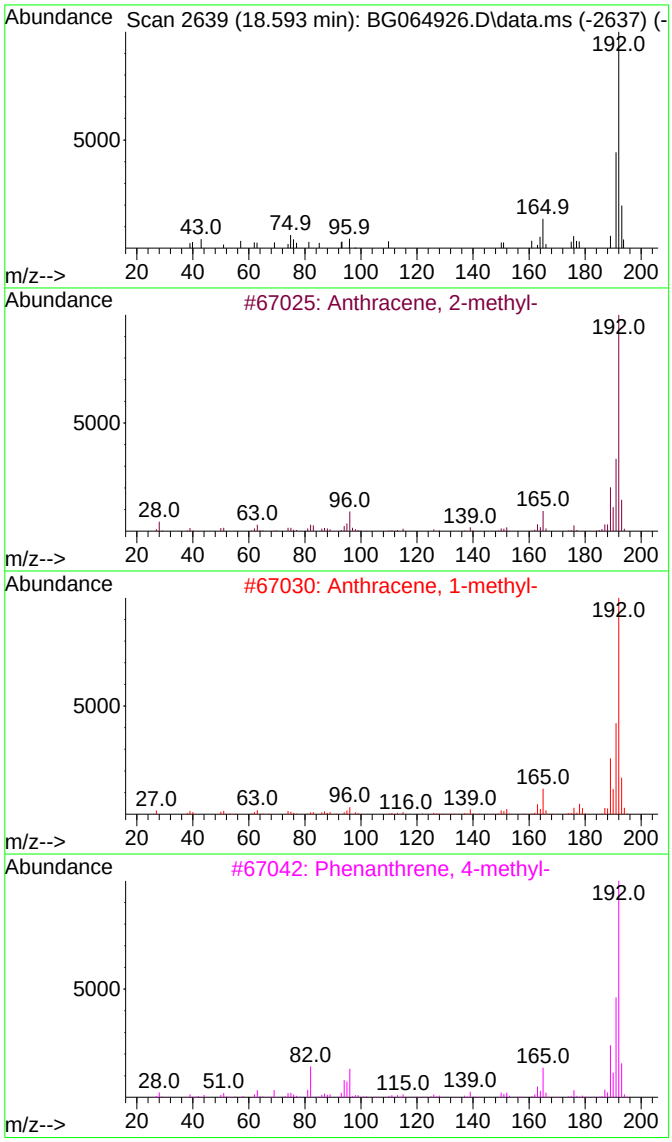
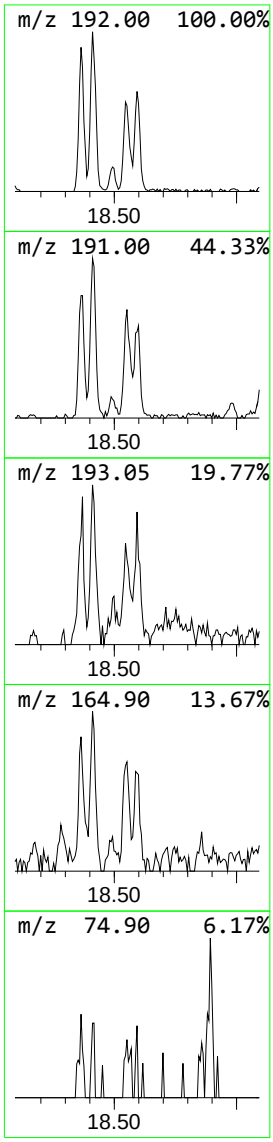
Peak Number 9 Anthracene, 2-methyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.593	2.33 ng	46588	Phenanthrene-d10	17.494

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 2-methyl-	192	C15H12	000613-12-7	87
2	Anthracene, 1-methyl-	192	C15H12	000610-48-0	87
3	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	86
4	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	80
5	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	80





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121025\
Data File : BG064926.D
Acq On : 10 Dec 2025 18:32
Operator : RC/JU
Sample : Q3826-09
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
TP-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111225.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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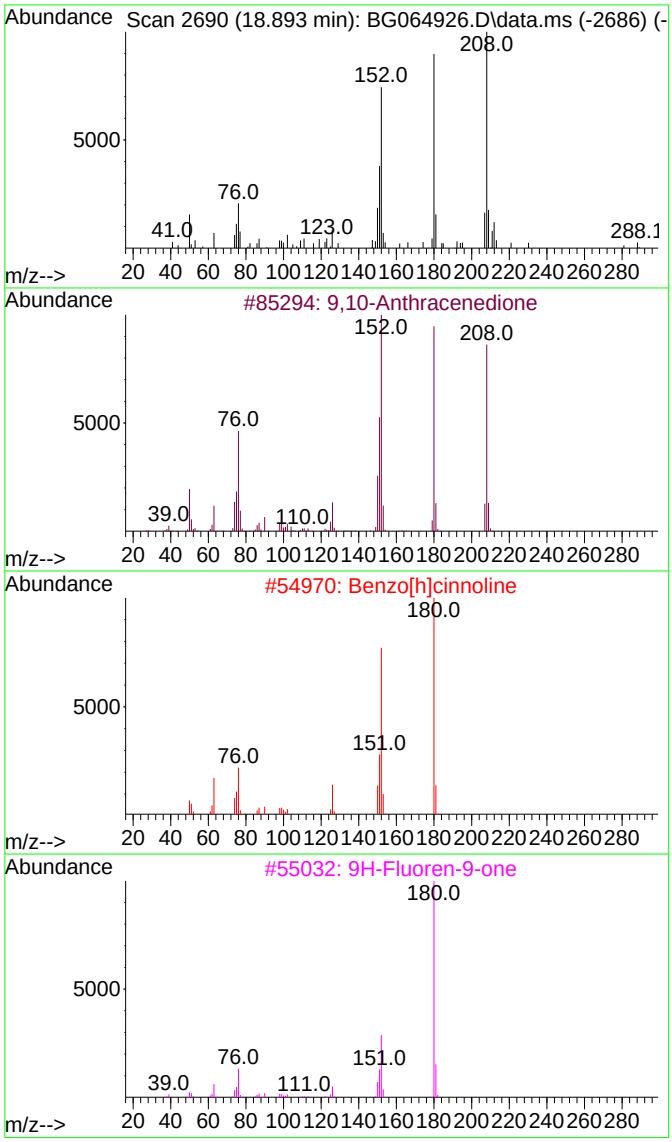
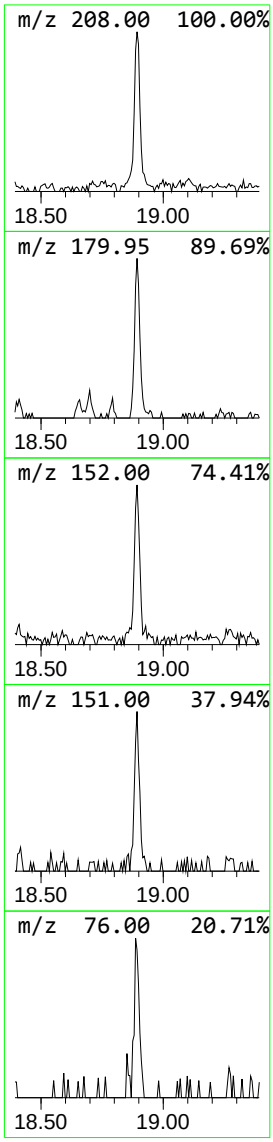
Peak Number 11 9,10-Anthracenedione Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.893	3.82 ng	76550	Phenanthrene-d10	17.494

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Anthracenedione	208	C14H8O2	000084-65-1	97
2	Benzo[h]cinnoline	180	C12H8N2	000230-31-9	64
3	9H-Fluoren-9-one	180	C13H8O	000486-25-9	58
4	Benzo[c]cinnoline	180	C12H8N2	000230-17-1	58
5	1H-Phenalen-1-one	180	C13H8O	000548-39-0	50





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121025\
Data File : BG064926.D
Acq On : 10 Dec 2025 18:32
Operator : RC/JU
Sample : Q3826-09
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
TP-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111225.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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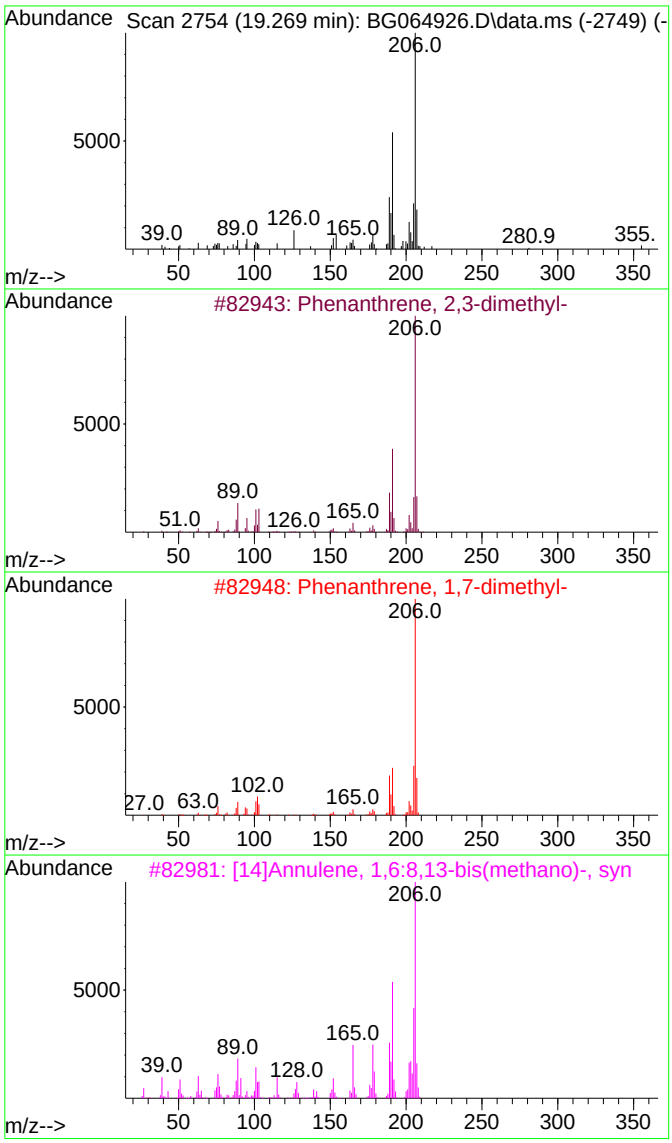
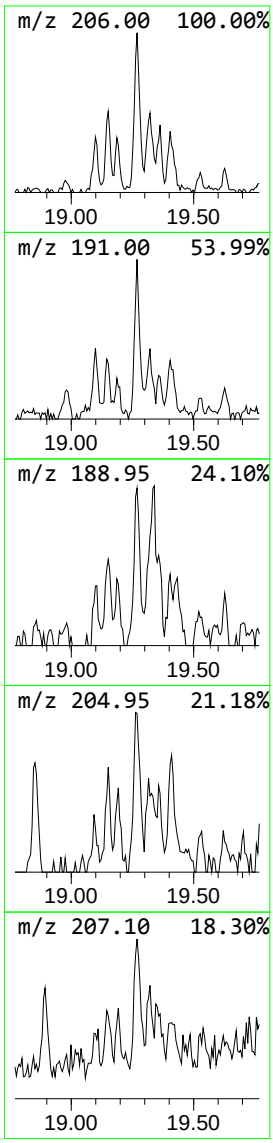
Peak Number 12 Phenanthrene, 2,3-dimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.269	3.26 ng	65216	Phenanthrene-d10	17.494

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	90
2	Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	81
3	[14]Annulene, 1,6:8,13-bis(metha...	206	C16H14	085385-68-8	81
4	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	80
5	9,10-Dimethylanthracene	206	C16H14	000781-43-1	76





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121025\
Data File : BG064926.D
Acq On : 10 Dec 2025 18:32
Operator : RC/JU
Sample : Q3826-09
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
TP-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111225.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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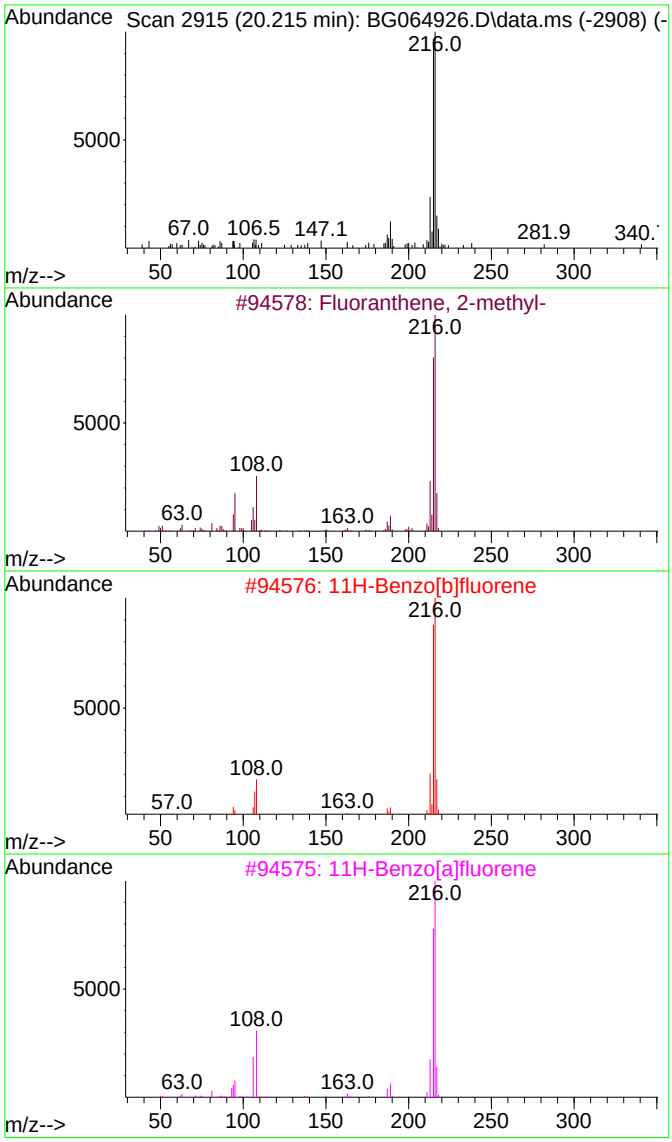
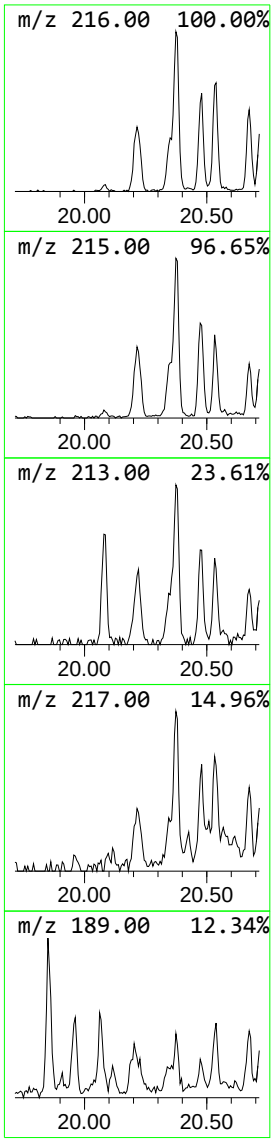
Peak Number 14 Fluoranthene, 2-methyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.215	2.12 ng	114620	Chrysene-d12	21.742

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	91
2	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	91
3	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	90
4	7H-Benzo[c]fluorene	216	C17H12	000205-12-9	80
5	Pyrene, 1-methyl-	216	C17H12	002381-21-7	70





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Acq On : 10 Dec 2025 18:32
Operator : RC/JU
Sample : Q3826-09
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
TP-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111225.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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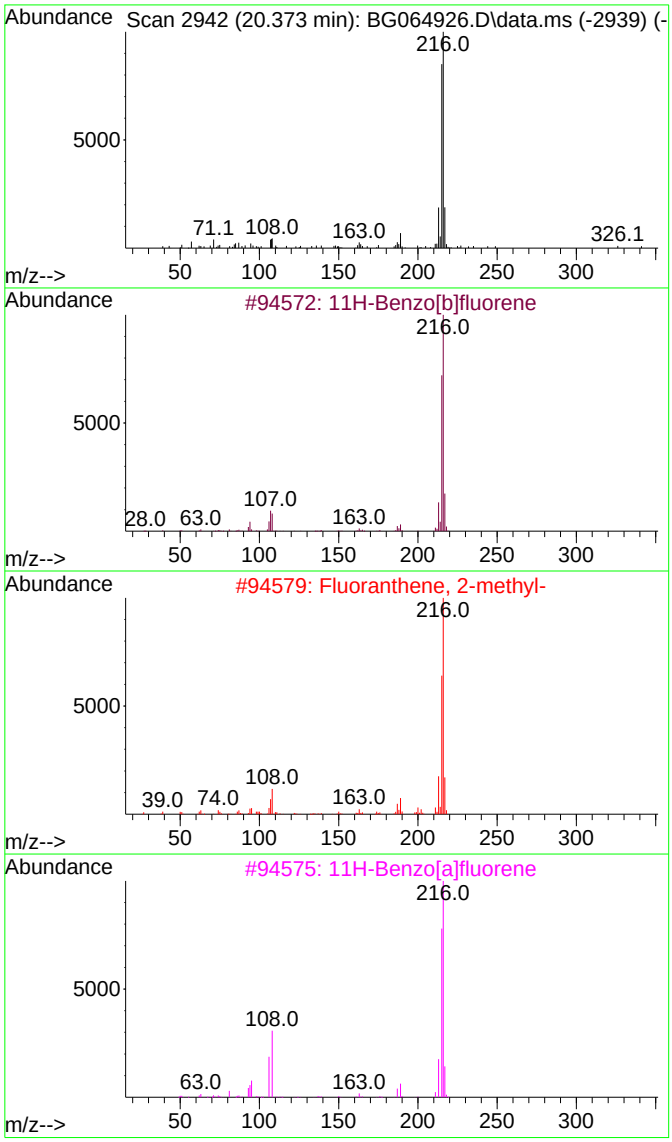
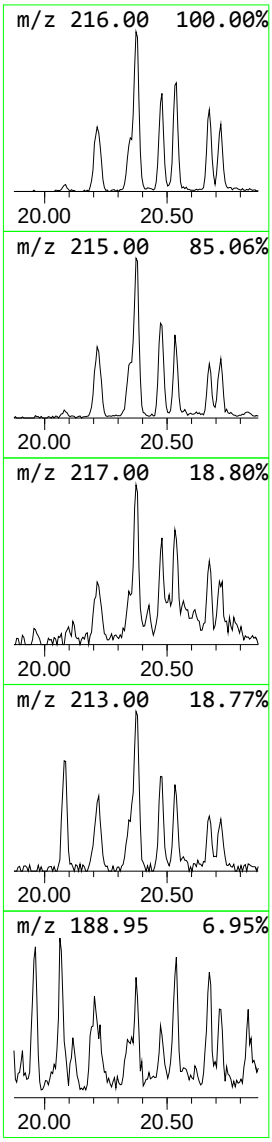
Peak Number 15 11H-Benzo[b]fluorene Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.373	2.37 ng	127751	Chrysene-d12	21.742

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	91
2	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	87
3	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	86
4	Pyrene, 1-methyl-	216	C17H12	002381-21-7	80
5	Pyrene, 2-methyl-	216	C17H12	003442-78-2	64





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Data File : BG064926.D
Acq On : 10 Dec 2025 18:32
Operator : RC/JU
Sample : Q3826-09
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
TP-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111225.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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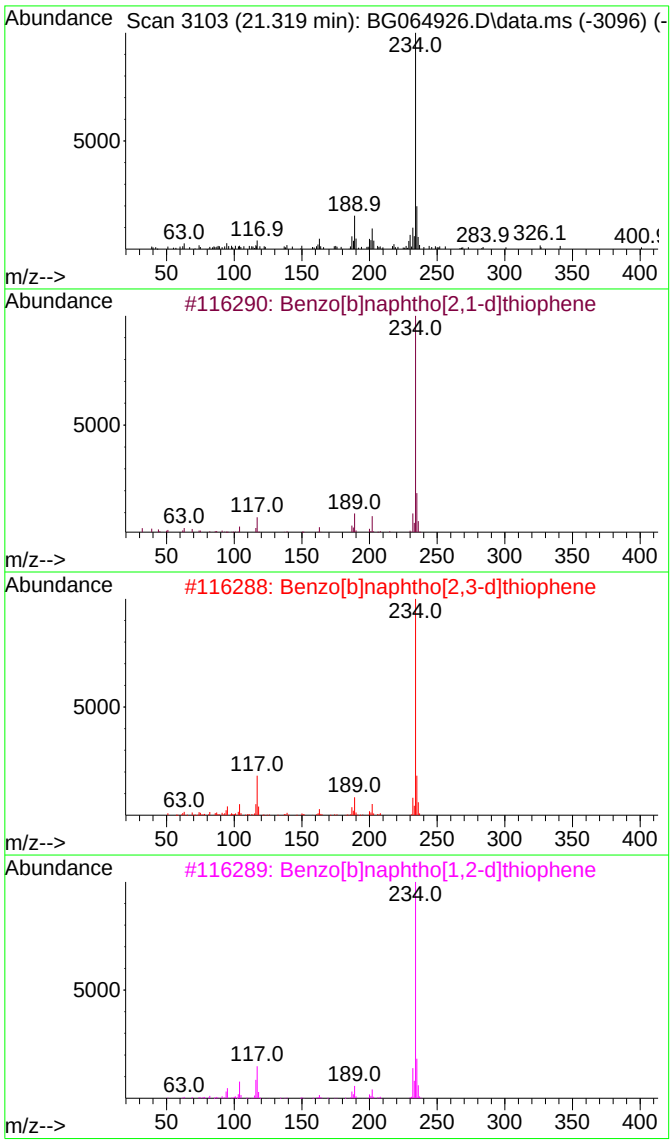
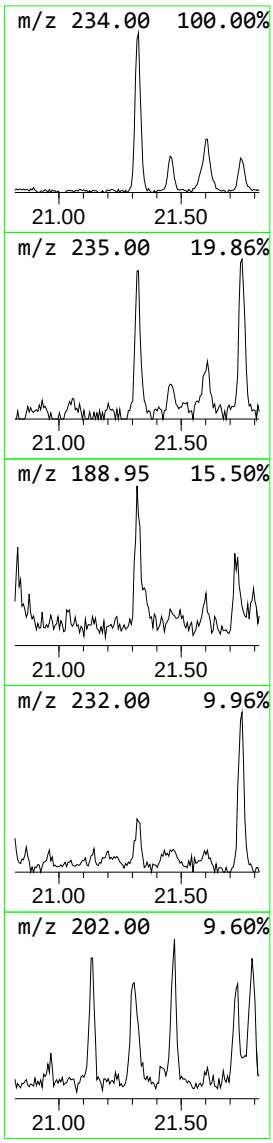
Peak Number 16 Benzo[b]naphtho[2,1-d]thiop... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.319	2.32 ng	125244	Chrysene-d12	21.742

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[b]naphtho[2,1-d]thiophene	234	C16H10S	000239-35-0	94
2	Benzo[b]naphtho[2,3-d]thiophene	234	C16H10S	000243-46-9	91
3	Benzo[b]naphtho[1,2-d]thiophene	234	C16H10S	000205-43-6	87
4	Phenanthro(2,1-b)thiophene	234	C16H10S	000219-25-0	81
5	Anthra(2,3-b)thiophene	234	C16H10S	022108-55-0	80





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121025\
Data File : BG064926.D
Acq On : 10 Dec 2025 18:32
Operator : RC/JU
Sample : Q3826-09
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
TP-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111225.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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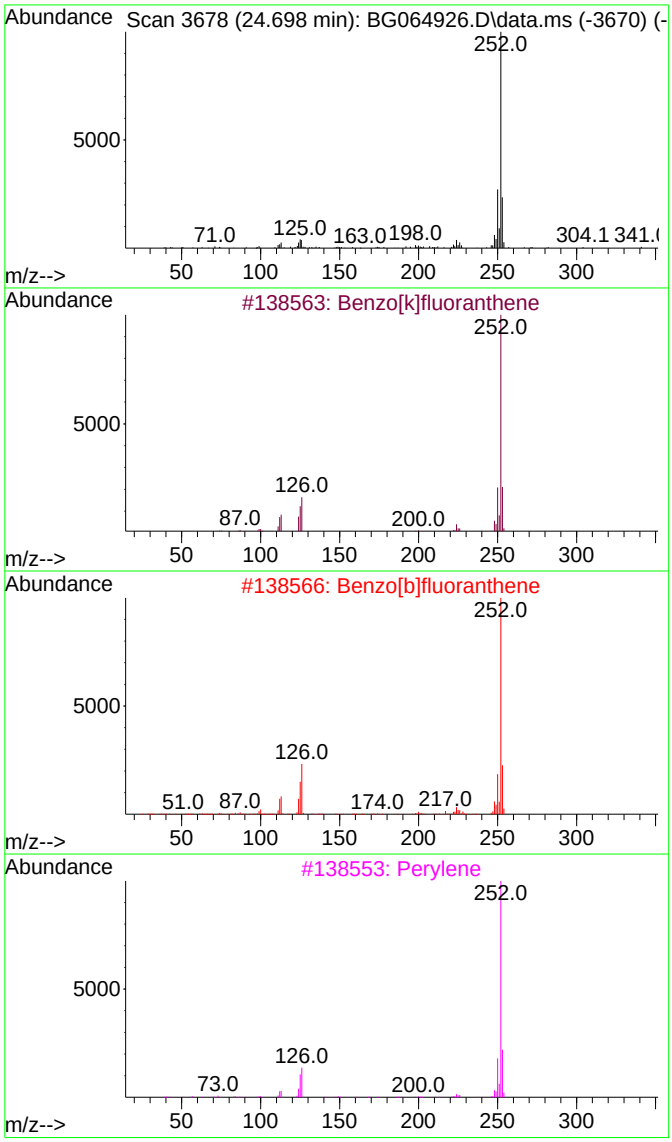
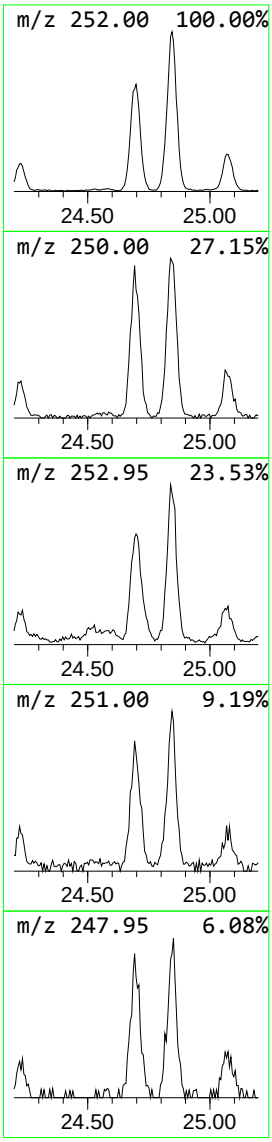
Peak Number 20 Perylene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.698	10.05 ng	309331	Perylene-d12	25.003

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[k]fluoranthene	252	C20H12	000207-08-9	94
2	Benzo[b]fluoranthene	252	C20H12	000205-99-2	90
3	Perylene	252	C20H12	000198-55-0	87
4	Benzo[e]pyrene	252	C20H12	000192-97-2	87
5	Benzo[a]pyrene	252	C20H12	000050-32-8	87





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121025\
Data File : BG064926.D
Acq On : 10 Dec 2025 18:32
Operator : RC/JU
Sample : Q3826-09
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
TP-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111225.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	16.102	5.2	ng	66942	3	14.762	255993	20.0
Phenanthrene, 1...	18.364	8.3	ng	166261	4	17.494	400594	20.0
Anthracene, 1-m...	18.411	5.5	ng	110855	4	17.494	400594	20.0
Anthracene, 2-m...	18.593	2.3	ng	46588	4	17.494	400594	20.0
9,10-Anthracene...	18.893	3.8	ng	76550	4	17.494	400594	20.0
Phenanthrene, 2...	19.269	3.3	ng	65216	4	17.494	400594	20.0
Fluoranthene, 2...	20.215	2.1	ng	114620	5	21.742	1078930	20.0
11H-Benzo[b]flu...	20.373	2.4	ng	127751	5	21.742	1078930	20.0
Benzo[b]naphtho...	21.319	2.3	ng	125244	5	21.742	1078930	20.0
Perylene	24.698	10.1	ng	309331	6	25.003	615612	20.0