LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG101625\
Data File : BG064521.D
Acq On : 16 Oct 2025 18:20
Operator : RC/JU
Sample : Q3348-05
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
WCS-TP2

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-BG100225.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG064521.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.342	206	213	219	rVB	15039	20868	1.30%	0.096%
2	5.411	389	395	403	rBV	185250	289415	18.06%	1.329%
3	6.992	658	664	673	rBV	198114	330319	20.61%	1.517%
4	7.814	798	804	812	rVB2	57167	92519	5.77%	0.425%
5	8.972	994	1001	1009	rBV	117412	211275	13.18%	0.970%
6	10.605	1272	1279	1283	rBV	73094	130380	8.14%	0.599%
7	10.652	1283	1287	1294	rVB	126750	214599	13.39%	0.985%
8	12.262	1555	1561	1569	rVB	34933	57986	3.62%	0.266%
9	12.485	1593	1599	1609	rVB2	47786	78869	4.92%	0.362%
10	13.067	1692	1698	1706	rVB	330169	489221	30.53%	2.246%
11	14.166	1879	1885	1894	rBV4	28014	53649	3.35%	0.246%
12	14.448	1925	1933	1938	rBV	125398	194705	12.15%	0.894%
13	14.506	1938	1943	1953	rVB	53174	80454	5.02%	0.369%
14	14.841	1994	2000	2005	rBV	15212	23404	1.46%	0.107%
15	15.493	2105	2111	2117	rBV	31612	46931	2.93%	0.215%
16	15.805	2157	2164	2170	rBV	60007	97014	6.05%	0.445%
17	15.934	2179	2186	2196	rVB	383842	563714	35.18%	2.588%
18	16.363	2254	2259	2263	rBV2	11833	19414	1.21%	0.089%
19	17.009	2364	2369	2377	rVB2	12918	21134	1.32%	0.097%
20	17.186	2394	2399	2403	rBV2	185353	287209	17.92%	1.319%
21	17.227	2403	2406	2414	rVV	233307	328366	20.49%	1.508%
22	17.321	2416	2422	2427	rVV	59885	90490	5.65%	0.416%
23	17.368	2427	2430	2438	rVB6	12128	23583	1.47%	0.108%
24	17.591	2457	2468	2473	rBV	36648	68505	4.27%	0.315%
25	18.061	2539	2548	2549	rBV	40281	58147	3.63%	0.267%
26	18.090	2549	2553	2555	rBV2	62770	85283	5.32%	0.392%
27	18.184	2564	2569	2574	rVB2	28866	43661	2.72%	0.200%
28	18.255	2576	2581	2585	rVV3	91953	161865	10.10%	0.743%
29	18.290	2585	2587	2594	rVB	35590	44334	2.77%	0.204%
30	18.560	2629	2633	2636	rVV	53981	70039	4.37%	0.322%
31	18.602	2636	2640	2646	rVB2	36633	52490	3.28%	0.241%
32	18.807	2672	2675	2679	rVB6	14526	18439	1.15%	0.085%
33	18.936	2694	2697	2700	rBV4	10073	16085	1.00%	0.074%
34	18.978	2700	2704	2710	rBV2	40014	64019	3.99%	0.294%
35	19.042	2710	2715	2718	rBV5	21057	36132	2.25%	0.166%
36	19.119	2723	2728	2737	rVB3	45769	89117	5.56%	0.409%

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Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
WCS-TP2

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-BG100225.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	19.242	2743	2749	2755	rBV	926614	1272969	79.43%	5.845%
38	19.307	2755	2760	2762	rBV5	24200	32928	2.05%	0.151%
39	19.336	2762	2765	2766	rVV	22650	20593	1.29%	0.095%
40	19.359	2766	2769	2772	rVV4	34035	55453	3.46%	0.255%
41	19.389	2772	2774	2778	rVB	40679	45820	2.86%	0.210%
42	19.436	2779	2782	2785	rVB4	15342	16964	1.06%	0.078%
43	19.483	2785	2790	2794	rBV2	30255	44087	2.75%	0.202%
44	19.571	2800	2805	2806	rBV	151228	172959	10.79%	0.794%
45	19.606	2806	2811	2819	rVB	876284	1257251	78.45%	5.773%
46	19.677	2819	2823	2829	rVB2	52739	81165	5.06%	0.373%
47	19.806	2829	2845	2849	rBV	748640	1045709	65.25%	4.802%
48	19.929	2860	2866	2872	rBV2	75478	192933	12.04%	0.886%
49	19.982	2873	2875	2879	rVV4	18715	28178	1.76%	0.129%
50	20.065	2883	2889	2890	rVV4	75549	128108	7.99%	0.588%
51	20.094	2890	2894	2900	rVB	221286	343030	21.41%	1.575%
52	20.194	2906	2911	2915	rBV2	198841	272842	17.03%	1.253%
53	20.253	2915	2921	2926	rVV2	117442	200107	12.49%	0.919%
54	20.335	2932	2935	2941	rBV5	28086	47332	2.95%	0.217%
55	20.394	2941	2945	2948	rBV	65876	90977	5.68%	0.418%
56	20.435	2948	2952	2957	rVB2	73619	107257	6.69%	0.492%
57	20.652	2985	2989	2991	rBV5	24114	32521	2.03%	0.149%
58	20.687	2991	2995	2997	rVV3	45291	71672	4.47%	0.329%
59	20.711	2997	2999	3004	rVV4	38802	57156	3.57%	0.262%
60	20.752	3004	3006	3009	rVV4	16629	17495	1.09%	0.080%
61	20.805	3010	3015	3018	rVV3	44126	83905	5.24%	0.385%
62	20.846	3018	3022	3028	rVB	78766	136256	8.50%	0.626%
63	21.016	3045	3051	3054	rBV3	90613	154516	9.64%	0.710%
64	21.057	3054	3058	3061	rVV	114327	166079	10.36%	0.763%
65	21.099	3061	3065	3071	rVV2	121603	258395	16.12%	1.186%
66	21.157	3071	3075	3082	rVB3	65369	135328	8.44%	0.621%
67	21.281	3093	3096	3102	rVB2	19134	30586	1.91%	0.140%
68	21.398	3110	3116	3123	rBV2	718869	1545191	96.42%	7.095%
69	21.463	3123	3127	3139	rVB	629551	999803	62.39%	4.591%
70	21.545	3139	3141	3145	rBV4	14837	22393	1.40%	0.103%
71	21.604	3146	3151	3155	rBV	84601	136789	8.54%	0.628%
72	21.663	3156	3161	3167	rBV	56937	90091	5.62%	0.414%
73	21.798	3178	3184	3191	rBV3	69972	168300	10.50%	0.773%
74	21.857	3191	3194	3198	rVB5	25496	36884	2.30%	0.169%
75	22.027	3220	3223	3226	rVV3	24200	30590	1.91%	0.140%
76	22.097	3227	3235	3240	rVV	118898	258634	16.14%	1.188%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG101625\
Data File : BG064521.D
Acq On : 16 Oct 2025 18:20
Operator : RC/JU
Sample : Q3348-05
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
WCS-TP2

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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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100225.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

77	22.168	3243	3247	3254	rVB2	75567	138104	8.62%	0.634%
78	22.250	3257	3261	3264	rVV4	32360	56863	3.55%	0.261%
79	22.291	3265	3268	3274	rVV3	68543	120024	7.49%	0.551%
80	22.356	3274	3279	3282	rVV	72973	136293	8.50%	0.626%
81	22.391	3282	3285	3293	rVB2	93071	173866	10.85%	0.798%
82	22.485	3295	3301	3306	rBV5	43938	77387	4.83%	0.355%
83	22.791	3349	3353	3362	rVB5	57039	140168	8.75%	0.644%
84	23.472	3458	3469	3475	rBV	501613	1602549	100.00%	7.359%
85	23.696	3501	3507	3516	rBV2	105895	225229	14.05%	1.034%
86	23.889	3534	3540	3542	rBV3	34983	71950	4.49%	0.330%
87	24.130	3573	3581	3593	rVB	282512	798847	49.85%	3.668%
88	24.271	3593	3605	3617	rBV	452504	1213283	75.71%	5.571%
89	24.401	3618	3627	3633	rVV2	151873	433334	27.04%	1.990%
90	24.471	3634	3639	3652	rVB3	129125	371020	23.15%	1.704%
91	25.464	3802	3808	3811	rBV8	15448	35751	2.23%	0.164%
92	27.767	4185	4200	4217	rVB3	203325	917564	57.26%	4.213%
93	27.955	4222	4232	4241	rBV2	19701	67136	4.19%	0.308%
94	28.179	4259	4270	4271	rBV3	44976	109960	6.86%	0.505%
95	28.325	4284	4295	4297	rBV5	41076	118440	7.39%	0.544%
96	28.801	4364	4376	4377	rBV5	137179	339586	21.19%	1.559%
97	29.360	4462	4471	4485	rVB2	43835	187877	11.72%	0.863%

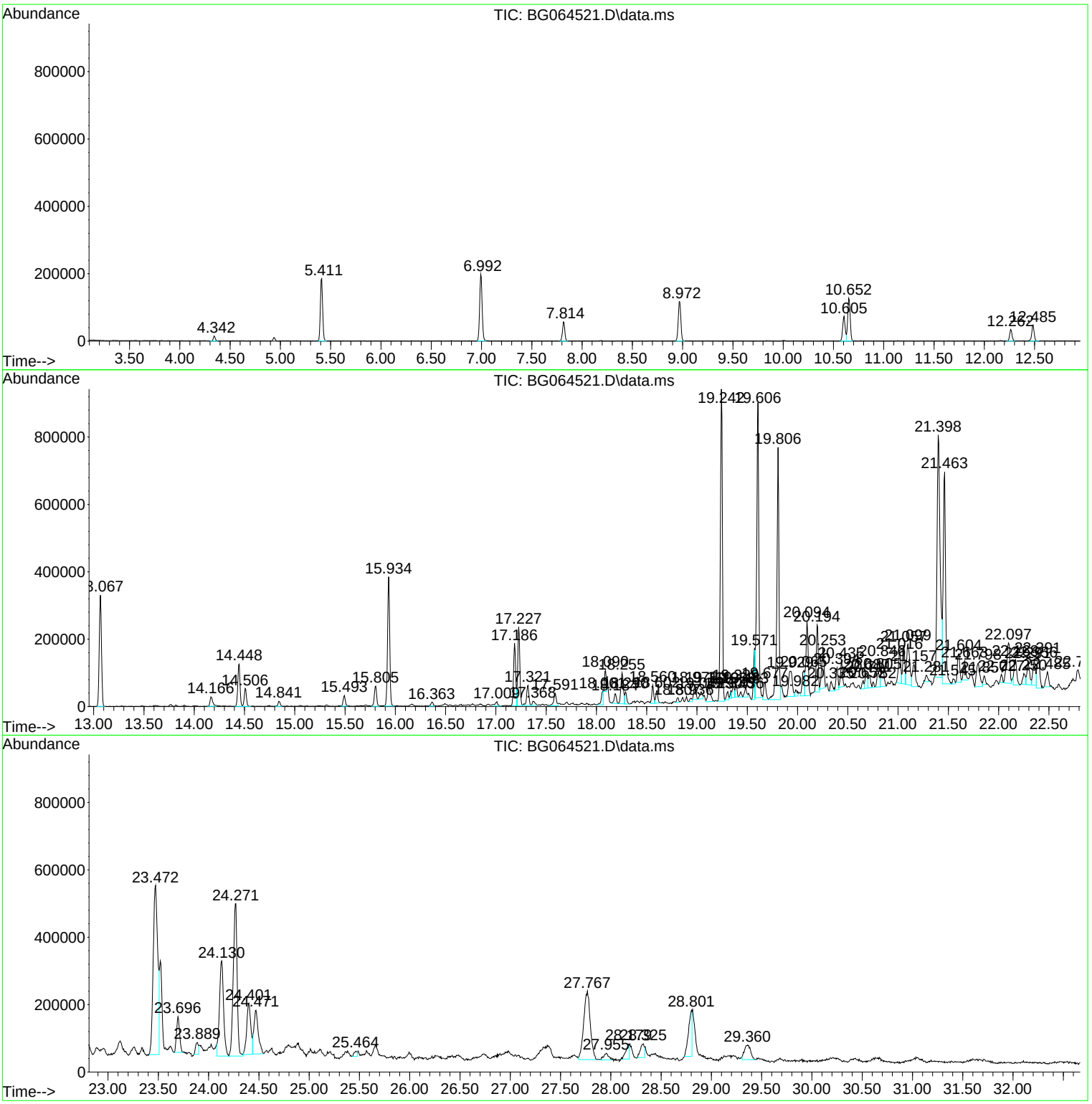
Sum of corrected areas: 21778111

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG101625\
Data File : BG064521.D
Acq On : 16 Oct 2025 18:20
Operator : RC/JU
Sample : Q3348-05
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
WCS-TP2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100225.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\BG101625\
Data File : BG064521.D
Acq On : 16 Oct 2025 18:20
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100225.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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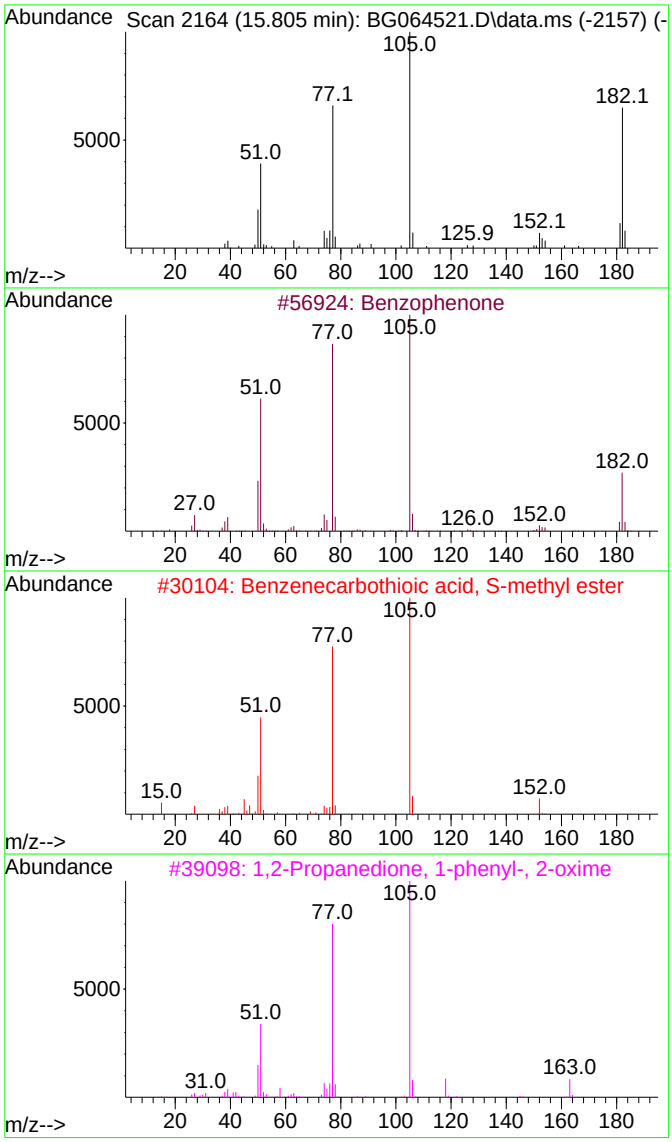
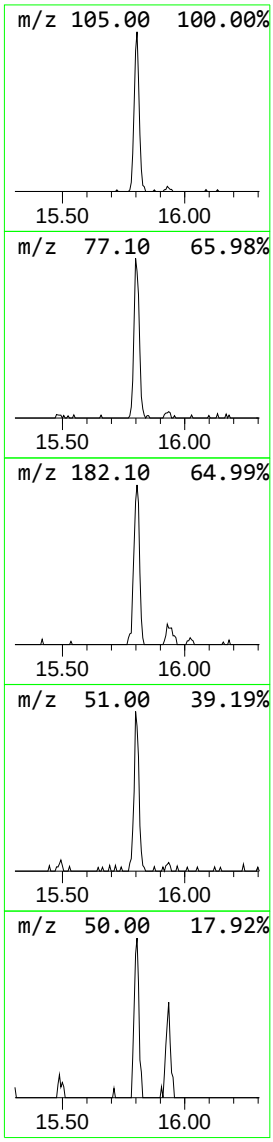
Peak Number 2 Benzophenone Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.805	9.97 ng	97014	Acenaphthene-d10	14.448

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzophenone	182	C13H10O	000119-61-9	95
2	Benzenecarbothioic acid, S-methy...	152	C8H8OS	005925-68-8	49
3	1,2-Propanedione, 1-phenyl-, 2-o...	163	C9H9NO2	000119-51-7	47
4	Ethanone, 2-bromo-1-phenyl-	198	C8H7BrO	000070-11-1	47
5	Benzenepropanenitrile, .beta.-oxo-	145	C9H7NO	000614-16-4	47





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG101625\
Data File : BG064521.D
Acq On : 16 Oct 2025 18:20
Operator : RC/JU
Sample : Q3348-05
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
WCS-TP2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100225.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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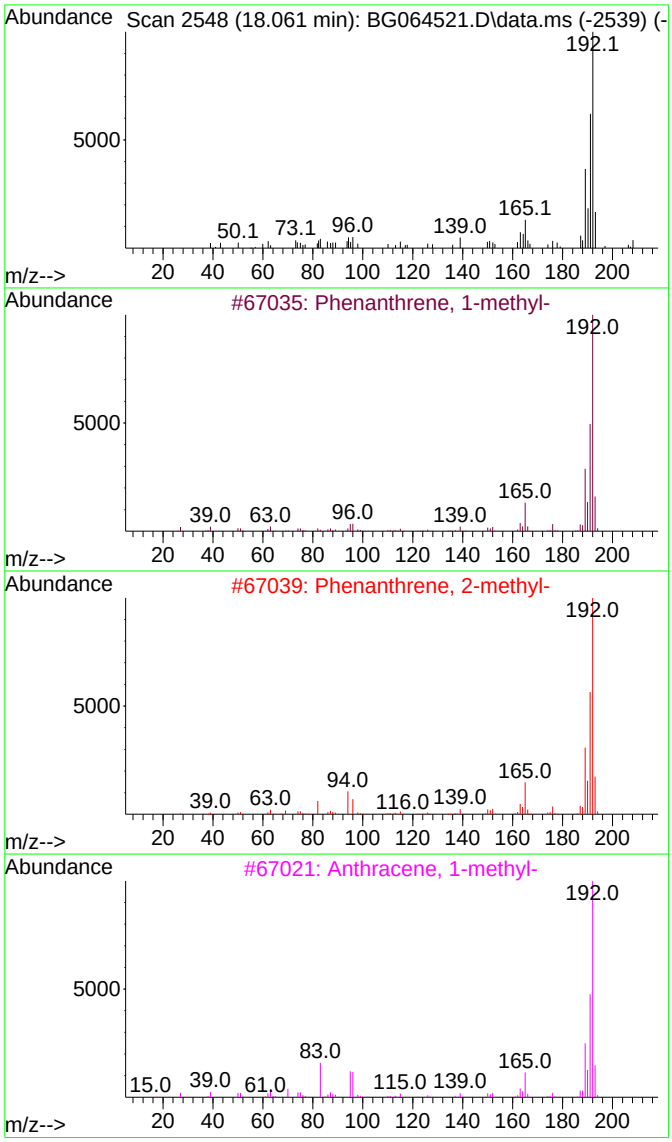
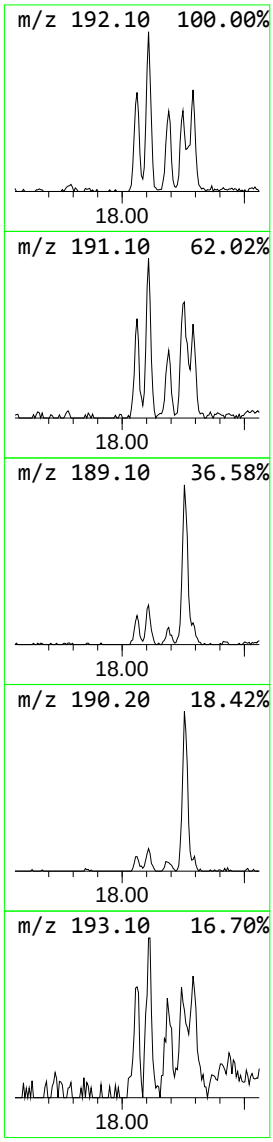
Peak Number 3 Phenanthrene, 1-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.061	4.05 ng	58147	Phenanthrene-d10	17.186

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	94
2	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
3	Anthracene, 1-methyl-	192	C15H12	000610-48-0	87
4	Anthracene, 9-methyl-	192	C15H12	000779-02-2	87
5	Anthracene, 2-methyl-	192	C15H12	000613-12-7	87





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG101625\
Data File : BG064521.D
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Operator : RC/JU
Sample : Q3348-05
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
WCS-TP2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100225.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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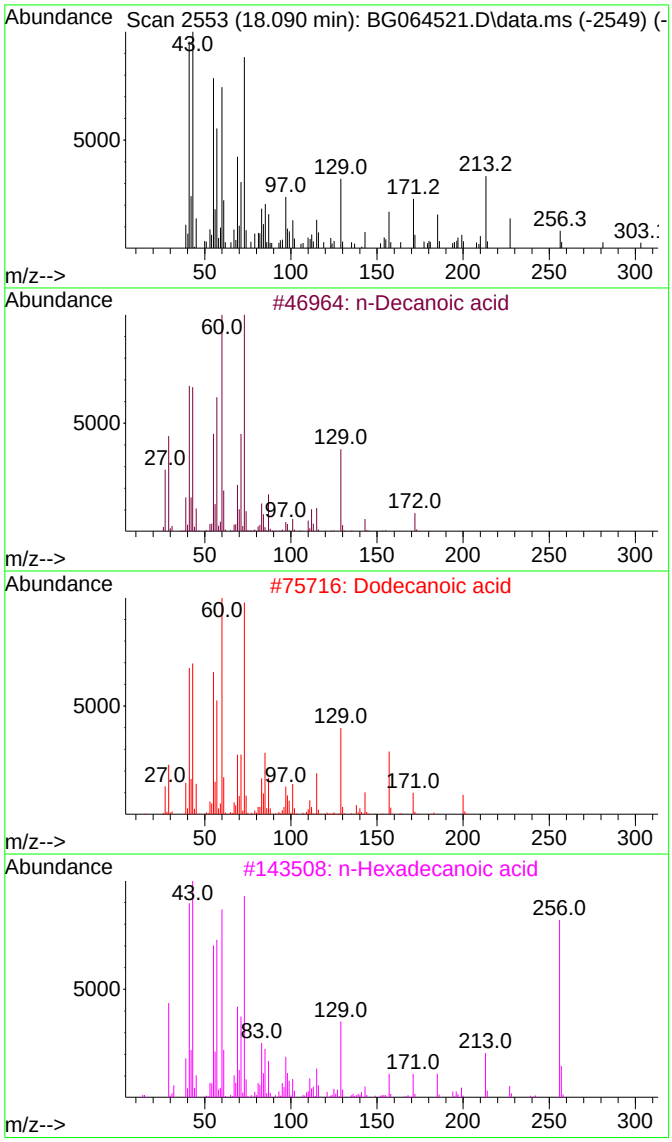
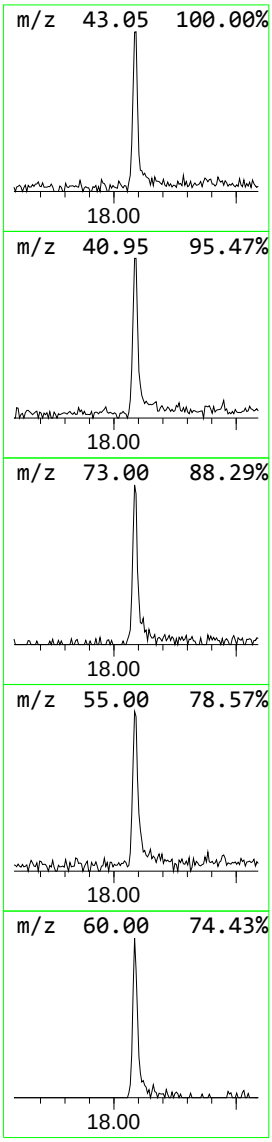
Peak Number 4 n-Decanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.090	5.94 ng	85283	Phenanthrene-d10	17.186

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Decanoic acid	172	C10H20O2	000334-48-5	87
2	Dodecanoic acid	200	C12H24O2	000143-07-7	83
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	76
4	Octadecanoic acid	284	C18H36O2	000057-11-4	68
5	Tetradecanoic acid	228	C14H28O2	000544-63-8	68





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Acq On : 16 Oct 2025 18:20
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ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
WCS-TP2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100225.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

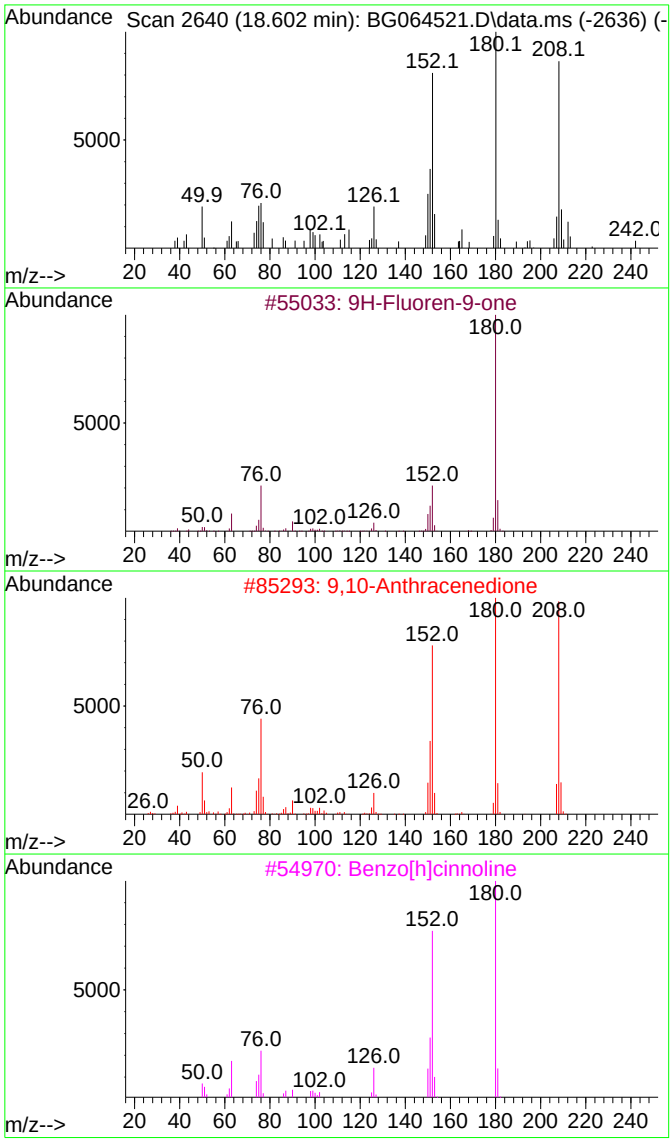
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

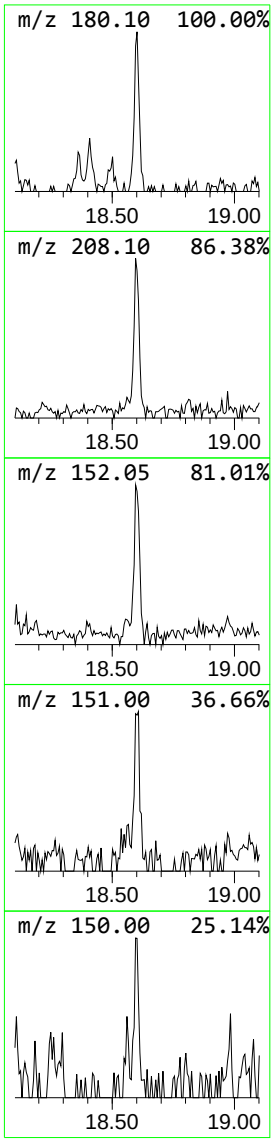
****

Peak Number 7 9H-Fluoren-9-one Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.602	3.66 ng	52490	Phenanthrene-d10	17.186

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9H-Fluoren-9-one	180	C13H8O	000486-25-9	95
2	9,10-Anthracenedione	208	C14H8O2	000084-65-1	93
3	Benzo[h]cinnoline	180	C12H8N2	000230-31-9	64
4	1H-Phenalen-1-one	180	C13H8O	000548-39-0	55
5	9,10-Phenanthrenedione	208	C14H8O2	000084-11-7	52





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG101625\
Data File : BG064521.D
Acq On : 16 Oct 2025 18:20
Operator : RC/JU
Sample : Q3348-05
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
WCS-TP2

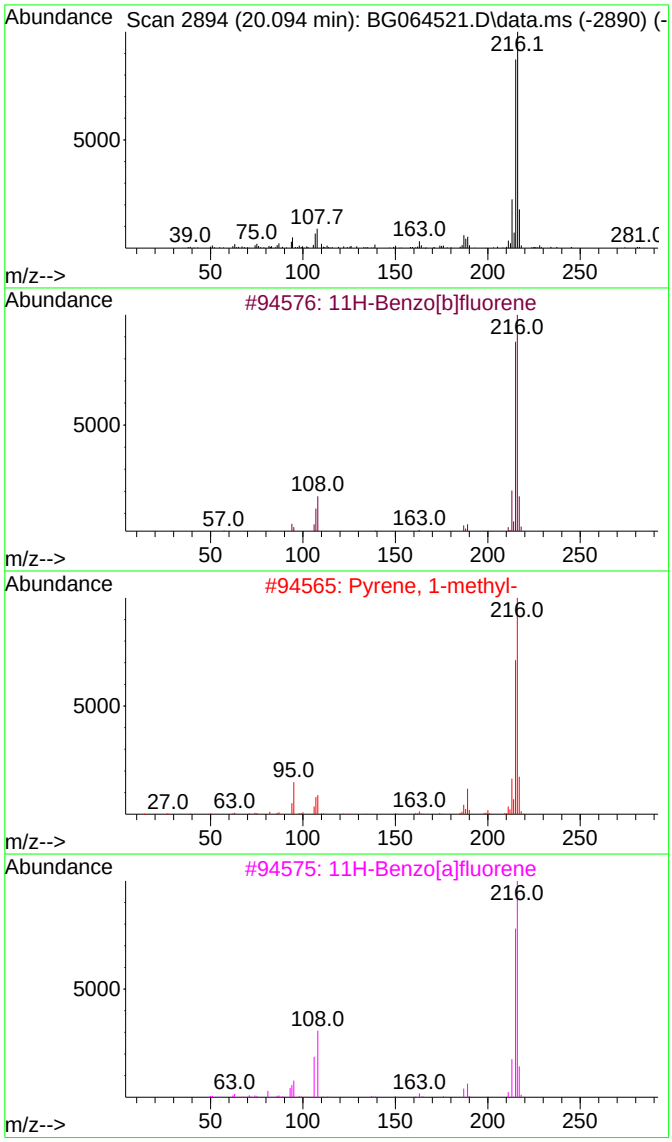
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100225.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

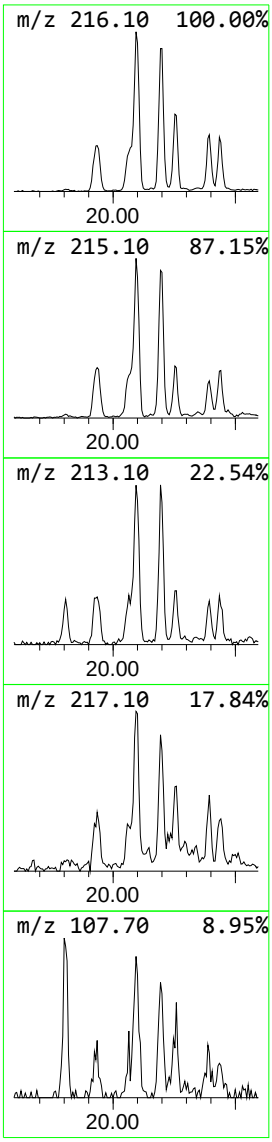
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 11H-Benzo[b]fluorene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.094	4.44 ng	343030	Chrysene-d12	21.416

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	93
2	Pyrene, 1-methyl-	216	C17H12	002381-21-7	93
3	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	90
4	7H-Benzo[c]fluorene	216	C17H12	000205-12-9	83
5	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	83





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG101625\
Data File : BG064521.D
Acq On : 16 Oct 2025 18:20
Operator : RC/JU
Sample : Q3348-05
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
WCS-TP2

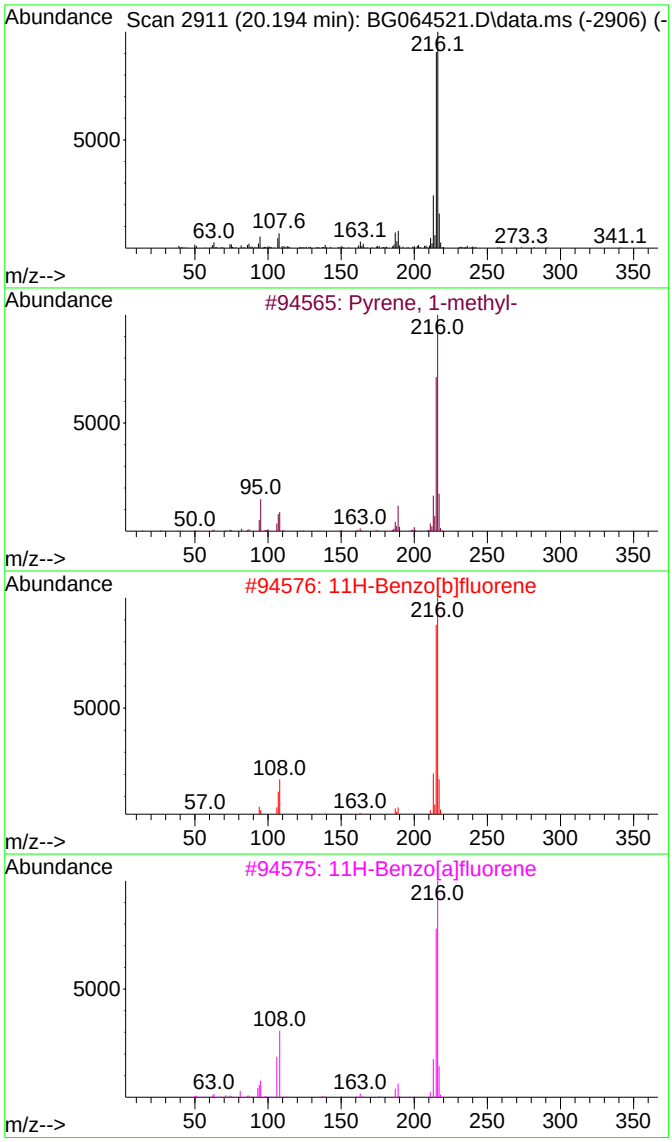
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100225.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

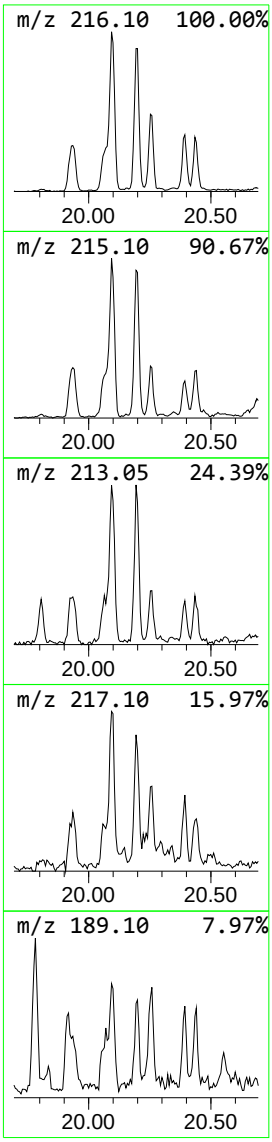
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 Pyrene, 1-methyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.194	3.53 ng	272842	Chrysene-d12	21.416

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-	216	C17H12	002381-21-7	93
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	72
5	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	70





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Data File : BG064521.D
Acq On : 16 Oct 2025 18:20
Operator : RC/JU
Sample : Q3348-05
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
WCS-TP2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100225.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

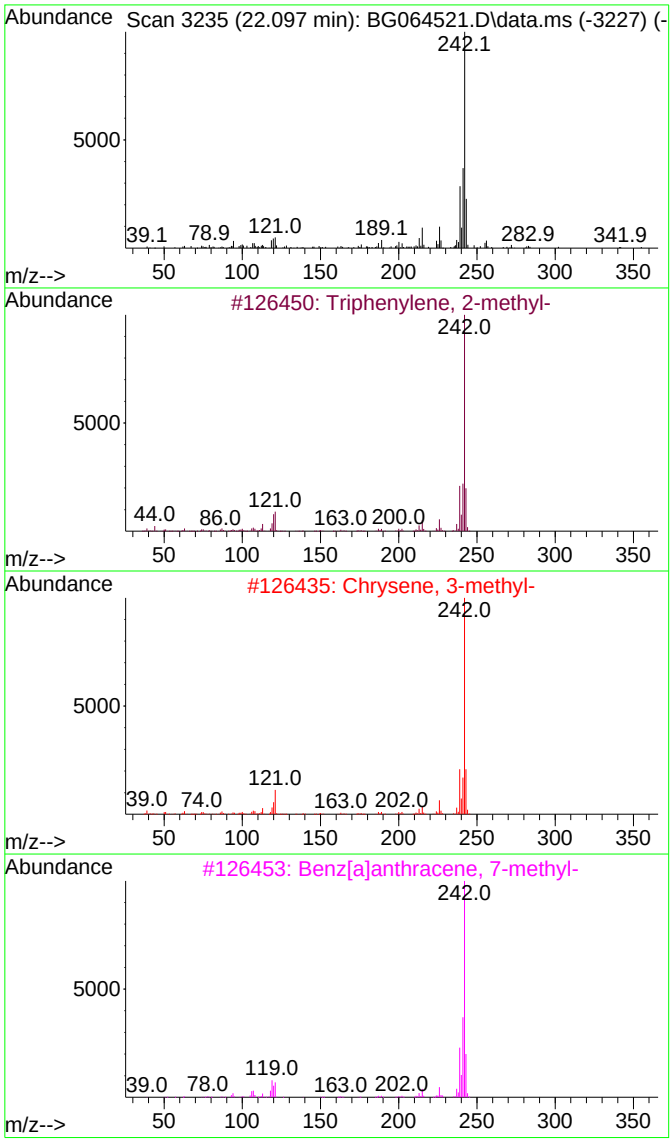
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

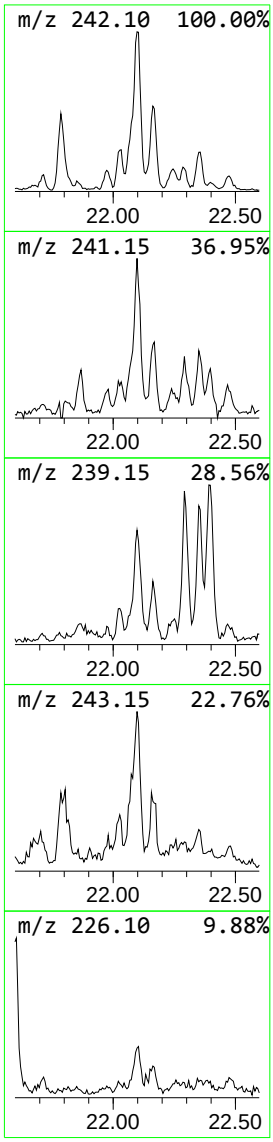
****

Peak Number 13 Triphenylene, 2-methyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.097	3.35 ng	258634	Chrysene-d12	21.416

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Triphenylene, 2-methyl-	242	C19H14	001705-84-6	94
2	Chrysene, 3-methyl-	242	C19H14	003351-31-3	93
3	Benz[a]anthracene, 7-methyl-	242	C19H14	002541-69-7	92
4	Benz[a]anthracene, 8-methyl-	242	C19H14	002381-31-9	92
5	Benz[a]anthracene, 1-methyl-	242	C19H14	002498-77-3	91





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Data File : BG064521.D
Acq On : 16 Oct 2025 18:20
Operator : RC/JU
Sample : Q3348-05
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
WCS-TP2

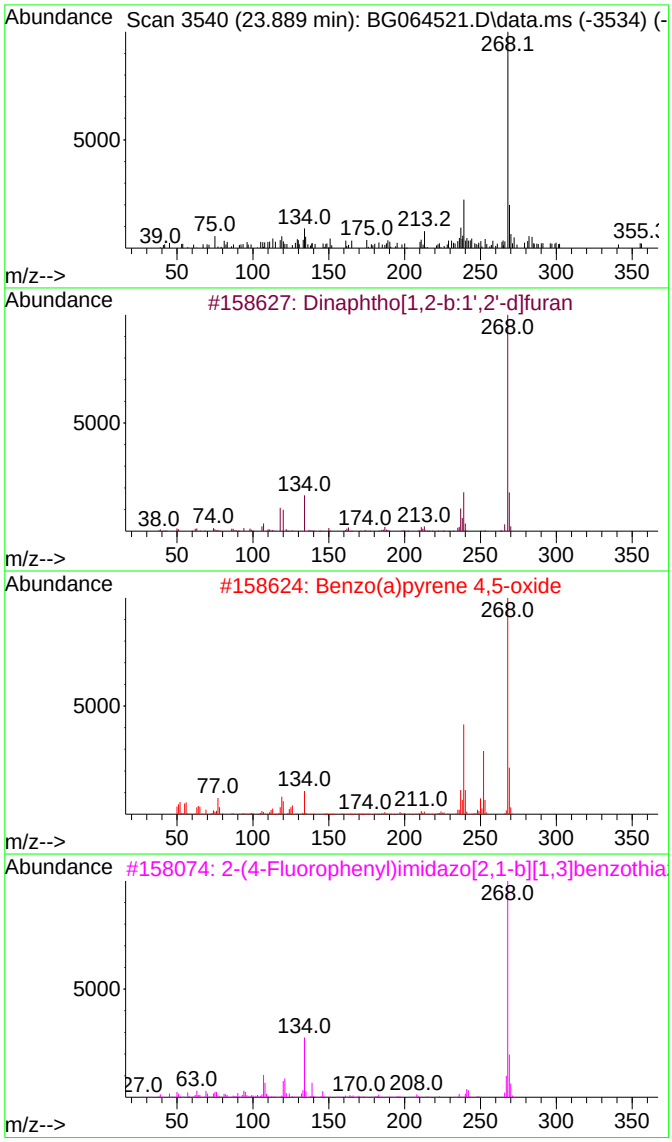
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100225.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

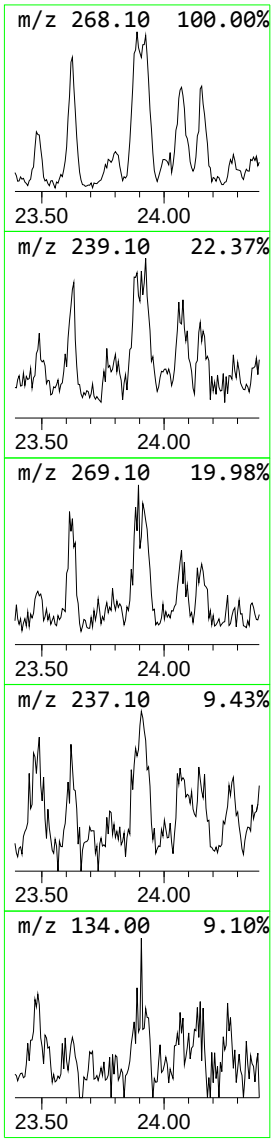
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 Dinaphtho[1,2-b:1',2'-d]furan Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.889	3.32 ng	71950	Perylene-d12	24.395

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dinaphtho[1,2-b:1',2'-d]furan	268	C20H12O	000207-93-2	87
2	Benzo(a)pyrene 4,5-oxide	268	C20H12O	037574-47-3	58
3	2-(4-Fluorophenyl)imidazo[2,1-b]...	268	C15H9FN2S	007067-87-0	53
4	2.6-Dihydroxyanthraquinone, dime...	268	C16H12O4	000963-96-2	52
5	Benzo(a)pyren-7-ol	268	C20H12O	037994-82-4	52





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Sample : Q3348-05
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
WCS-TP2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100225.M
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TIC Library : C:\Database\NIST20.L
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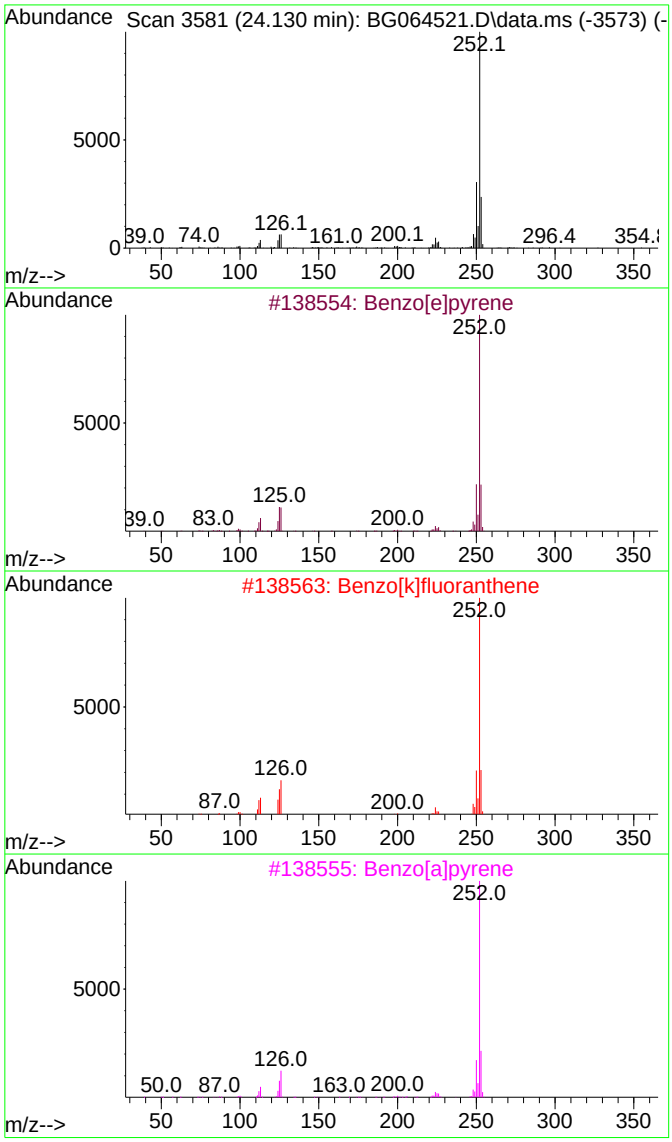
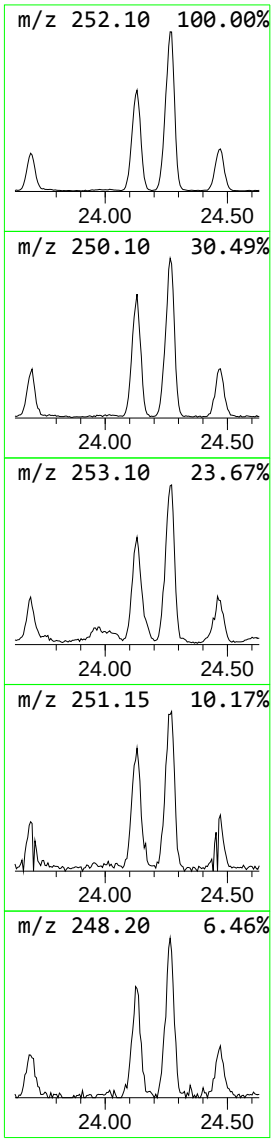
Peak Number 16 Benzo[e]pyrene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.130	36.87 ng	798847	Perylene-d12	24.395

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[e]pyrene	252	C20H12	000192-97-2	95
2	Benzo[k]fluoranthene	252	C20H12	000207-08-9	94
3	Benzo[a]pyrene	252	C20H12	000050-32-8	90
4	Benzo[b]fluoranthene	252	C20H12	000205-99-2	87
5	Benzo[j]fluoranthene	252	C20H12	000205-82-3	81





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Data File : BG064521.D
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Sample : Q3348-05
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
WCS-TP2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100225.M
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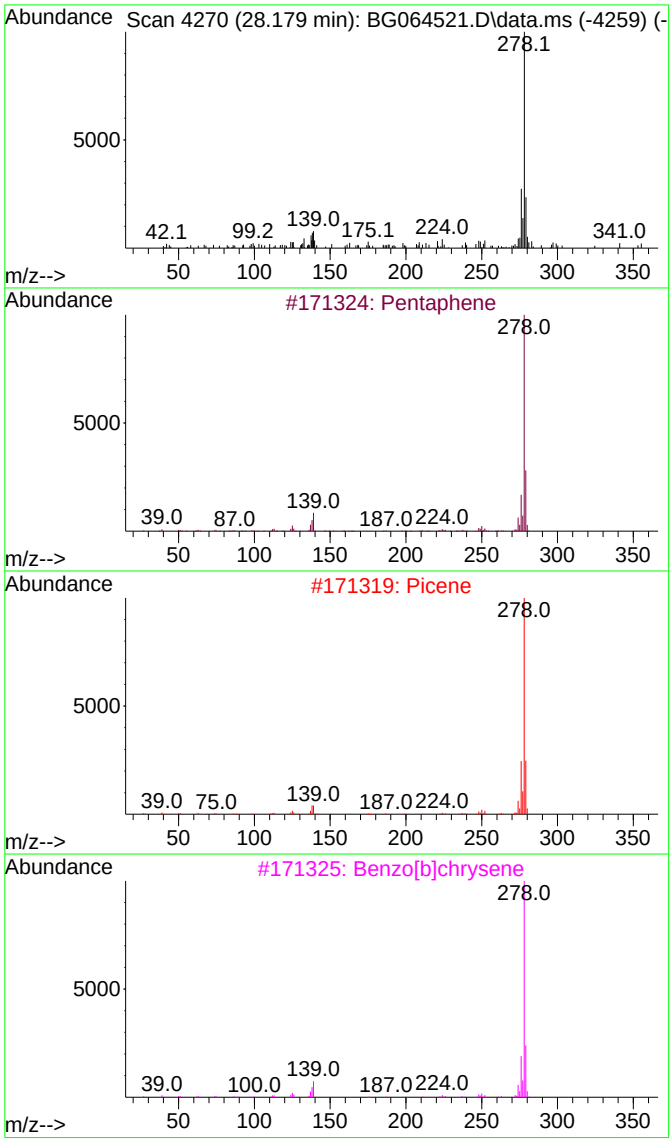
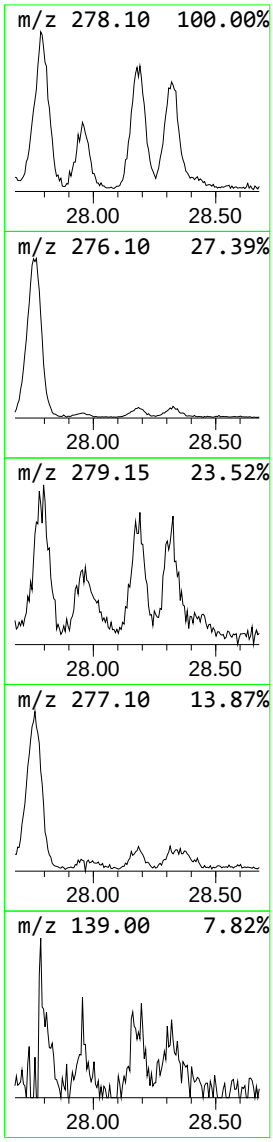
Peak Number 18 Pentaphene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
28.179	5.08 ng	109960	Perylene-d12	24.395

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentaphene	278	C22H14	000222-93-5	90
2	Picene	278	C22H14	000213-46-7	90
3	Benzo[b]chrysene	278	C22H14	000214-17-5	90
4	Naphtho[1,2-a]anthracene	278	C22H14	000195-06-2	76
5	Pyrene, 1-phenyl-	278	C22H14	005101-27-9	76





Tentatively Identified Compound (LSC) summary

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ClientSampleId :
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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
Benzophenone	15.805	10.0	ng	97014	3	14.448	194705	20.0
Phenanthrene, 1...	18.061	4.0	ng	58147	4	17.186	287209	20.0
n-Decanoic acid	18.090	5.9	ng	85283	4	17.186	287209	20.0
9H-Fluoren-9-one	18.602	3.7	ng	52490	4	17.186	287209	20.0
11H-Benzo[b]flu...	20.094	4.4	ng	343030	5	21.416	1545190	20.0
Pyrene, 1-methyl-	20.194	3.5	ng	272842	5	21.416	1545190	20.0
Triphenylene, 2...	22.097	3.4	ng	258634	5	21.416	1545190	20.0
Dinaphtho[1,2-b...	23.889	3.3	ng	71950	6	24.395	433334	20.0
Benzo[e]pyrene	24.130	36.9	ng	798847	6	24.395	433334	20.0
Pentaphene	28.179	5.1	ng	109960	6	24.395	433334	20.0