

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052223\
 Data File : BG057504.D
 Acq On : 22 May 2023 22:06
 Operator : CG/JU
 Sample : 02864-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 ETGI-289

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-BG042723.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG057504.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.390	309	315	320	rBV	11583	18311	2.07%	0.274%
2	5.871	390	397	405	rBV	207480	348260	39.37%	5.212%
3	6.318	467	473	480	rBV3	5922	10998	1.24%	0.165%
4	7.499	665	674	688	rBV	224738	402930	45.55%	6.030%
5	8.356	814	820	829	rVB	72882	125394	14.18%	1.877%
6	9.531	1012	1020	1030	rBV	140224	263069	29.74%	3.937%
7	11.117	1284	1290	1295	rBV2	8590	14089	1.59%	0.211%
8	11.194	1295	1303	1308	rVV	97802	181035	20.47%	2.709%
9	11.241	1308	1311	1317	rVV	14617	23892	2.70%	0.358%
10	11.305	1317	1322	1328	rVB	6035	9943	1.12%	0.149%
11	12.157	1459	1467	1471	rBV	10853	16876	1.91%	0.253%
12	12.539	1524	1532	1539	rVB2	12052	18392	2.08%	0.275%
13	12.821	1573	1580	1586	rVB2	21245	33446	3.78%	0.501%
14	13.038	1612	1617	1623	rBV	14268	22242	2.51%	0.333%
15	13.473	1685	1691	1697	rBV2	6548	9409	1.06%	0.141%
16	13.590	1701	1711	1718	rBV	408885	621139	70.22%	9.296%
17	13.755	1733	1739	1744	rBV2	14057	19826	2.24%	0.297%
18	14.131	1798	1803	1805	rBV3	6151	8858	1.00%	0.133%
19	14.295	1824	1831	1834	rBV2	11120	21412	2.42%	0.320%
20	14.342	1834	1839	1843	rVV3	8670	12850	1.45%	0.192%
21	14.407	1843	1850	1854	rVV5	8278	16333	1.85%	0.244%
22	14.524	1864	1870	1874	rBV4	9214	13664	1.54%	0.204%
23	14.812	1912	1919	1924	rVB2	14628	20732	2.34%	0.310%
24	14.971	1940	1946	1952	rVB	154830	248490	28.09%	3.719%
25	15.036	1952	1957	1961	rVB3	5780	10414	1.18%	0.156%
26	15.364	2006	2013	2018	rVV3	13761	24631	2.78%	0.369%
27	15.423	2018	2023	2031	rVB4	8247	17165	1.94%	0.257%
28	15.570	2043	2048	2052	rBV2	6738	10861	1.23%	0.163%
29	15.623	2052	2057	2061	rVV4	6385	9019	1.02%	0.135%
30	15.758	2067	2080	2085	rBV4	17556	39481	4.46%	0.591%
31	15.993	2115	2120	2128	rBV6	8640	20442	2.31%	0.306%
32	16.163	2142	2149	2152	rBV2	5518	9284	1.05%	0.139%
33	16.304	2167	2173	2179	rVB2	39758	62285	7.04%	0.932%
34	16.451	2192	2198	2205	rBV	308008	477589	53.99%	7.148%
35	16.616	2221	2226	2228	rBV	17266	25445	2.88%	0.381%
36	16.639	2228	2230	2233	rVB2	14992	15545	1.76%	0.233%

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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

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37	17.227	2325	2330	2334	rBV5	8359	16090	1.82%	0.241%
38	17.362	2348	2353	2357	rBV3	11546	18613	2.10%	0.279%
39	17.403	2357	2360	2367	rVV2	14859	27409	3.10%	0.410%
40	17.456	2367	2369	2377	rVB3	6671	8920	1.01%	0.133%
41	17.708	2405	2412	2416	rBV	215609	326012	36.86%	4.879%
42	17.749	2416	2419	2426	rVB	49502	69878	7.90%	1.046%
43	18.020	2462	2465	2469	rVB2	8536	13130	1.48%	0.197%
44	18.143	2481	2486	2491	rVB4	16468	24426	2.76%	0.366%
45	18.290	2507	2511	2520	rVB6	11197	20463	2.31%	0.306%
46	18.425	2532	2534	2543	rVB8	4984	9714	1.10%	0.145%
47	18.548	2551	2555	2557	rBV2	28424	40258	4.55%	0.603%
48	18.572	2557	2559	2563	rVV	24724	36472	4.12%	0.546%
49	18.619	2563	2567	2574	rVB2	36287	61493	6.95%	0.920%
50	18.760	2585	2591	2596	rBV4	25110	43981	4.97%	0.658%
51	18.807	2596	2599	2601	rBV2	10782	14789	1.67%	0.221%
52	18.954	2621	2624	2630	rVB7	8067	11793	1.33%	0.176%
53	19.059	2637	2642	2647	rBV2	25263	39827	4.50%	0.596%
54	19.112	2647	2651	2659	rVB8	10254	20205	2.28%	0.302%
55	19.300	2675	2683	2687	rBV4	11722	22441	2.54%	0.336%
56	19.353	2688	2692	2696	rVB2	12467	18018	2.04%	0.270%
57	19.394	2696	2699	2702	rVB	10796	12940	1.46%	0.194%
58	19.447	2704	2708	2709	rBV2	15760	18274	2.07%	0.273%
59	19.471	2709	2712	2716	rVV2	29355	38727	4.38%	0.580%
60	19.518	2718	2720	2725	rVV3	9537	14364	1.62%	0.215%
61	19.565	2725	2728	2731	rVB3	10059	12351	1.40%	0.185%
62	19.623	2732	2738	2742	rBV3	10482	22460	2.54%	0.336%
63	19.741	2750	2758	2762	rBV	82008	132526	14.98%	1.983%
64	19.788	2762	2766	2770	rBV3	17171	27670	3.13%	0.414%
65	20.011	2798	2804	2808	rVB8	15796	25097	2.84%	0.376%
66	20.064	2808	2813	2815	rBV2	18410	26071	2.95%	0.390%
67	20.099	2815	2819	2825	rVV	80054	123380	13.95%	1.847%
68	20.158	2825	2829	2835	rVB2	16786	27481	3.11%	0.411%
69	20.275	2843	2849	2854	rBV	719770	884563	100.00%	13.239%
70	20.534	2890	2893	2896	rBV3	20299	27752	3.14%	0.415%
71	20.581	2899	2901	2906	rVB3	18661	21831	2.47%	0.327%
72	20.681	2915	2918	2922	rBV3	7837	12492	1.41%	0.187%
73	20.740	2926	2928	2931	rVB2	11499	11648	1.32%	0.174%
74	20.886	2949	2953	2957	rBV	14295	23234	2.63%	0.348%
75	21.039	2976	2979	2982	rVB3	13710	15678	1.77%	0.235%
76	21.374	3032	3036	3041	rVB	26151	40602	4.59%	0.608%

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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

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77	21.568	3065	3069	3074	rBV4	31969	57044	6.45%	0.854%
78	22.014	3136	3145	3150	rBV3	208579	432385	48.88%	6.471%
79	22.061	3150	3153	3163	rVB	41744	78983	8.93%	1.182%
80	22.220	3178	3180	3185	rBV6	6415	12714	1.44%	0.190%
81	22.790	3275	3277	3283	rVB3	21938	33627	3.80%	0.503%
82	23.530	3401	3403	3408	rVB6	9287	11714	1.32%	0.175%
83	24.393	3543	3550	3557	rBV3	33795	99265	11.22%	1.486%
84	25.345	3705	3712	3721	rVB3	20058	59288	6.70%	0.887%
85	25.510	3731	3740	3750	rVB2	110032	329890	37.29%	4.937%

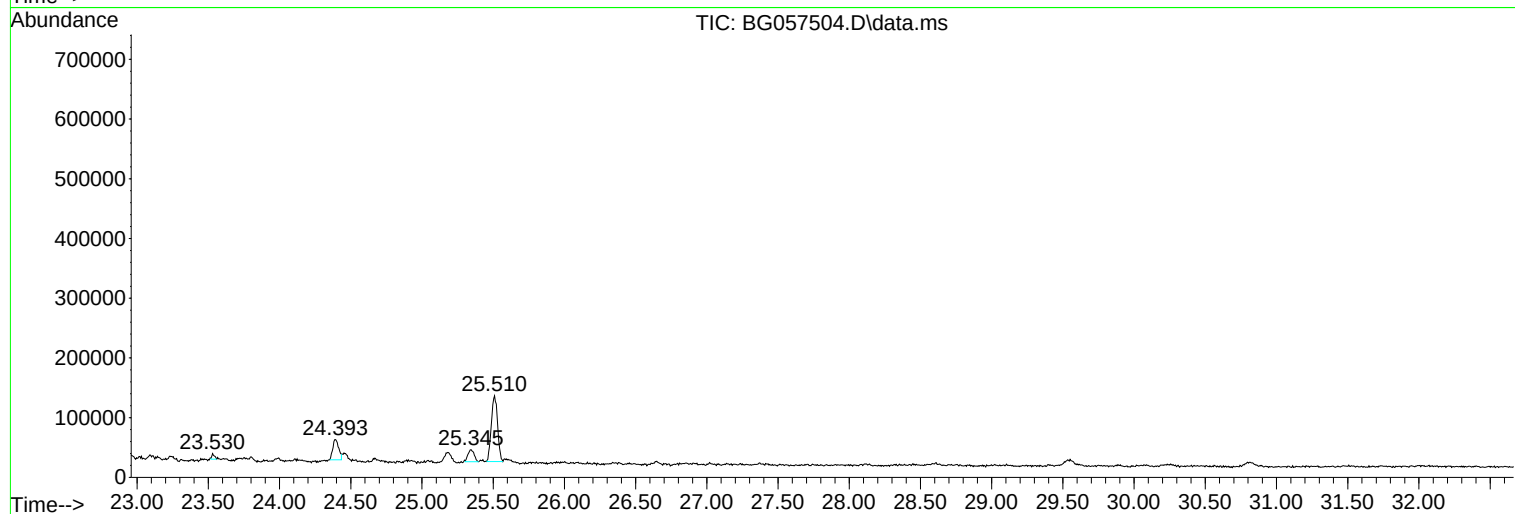
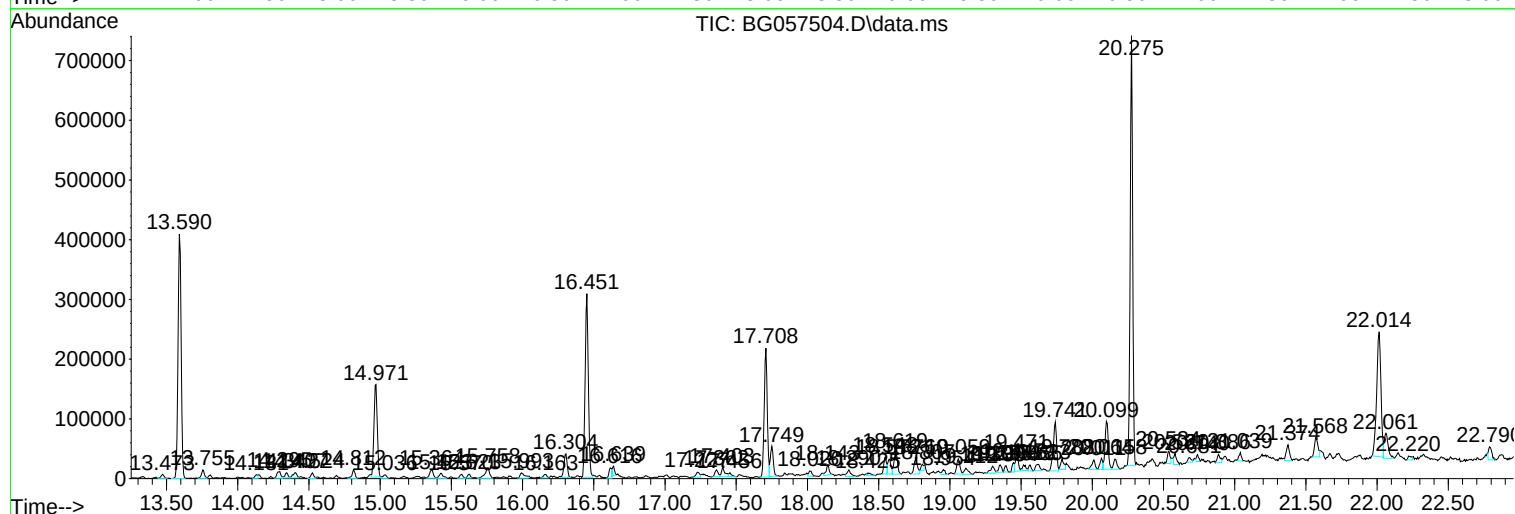
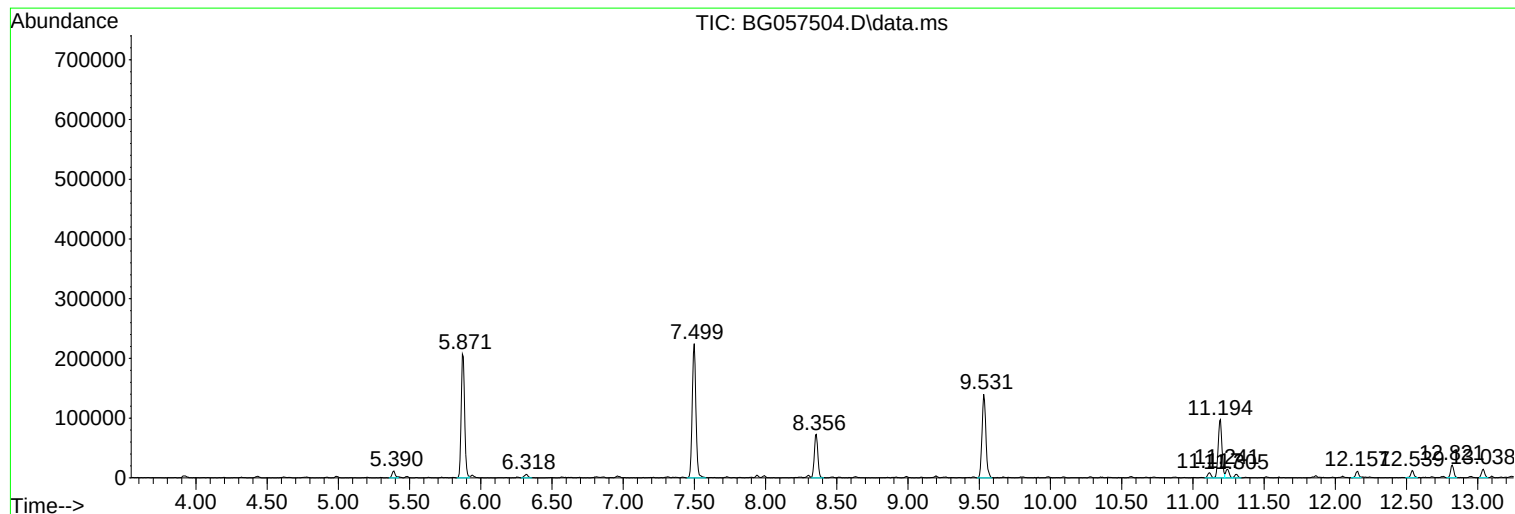
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG042723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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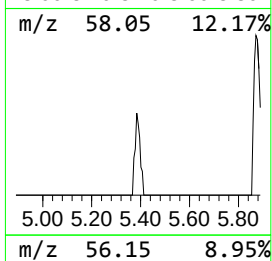
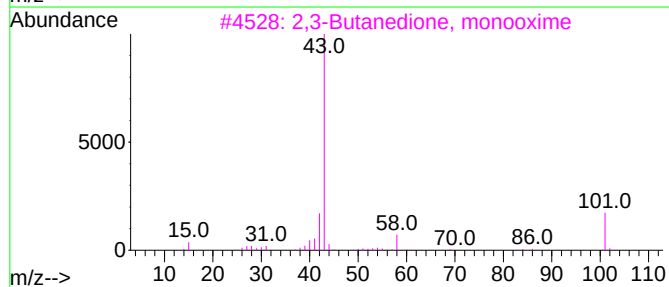
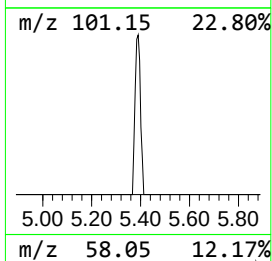
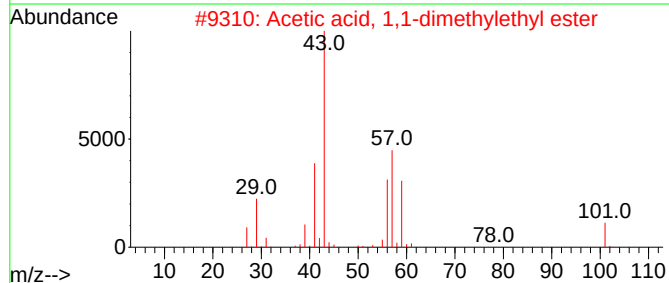
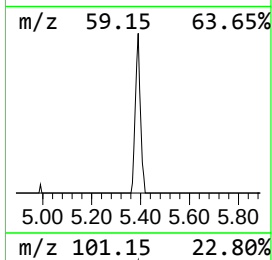
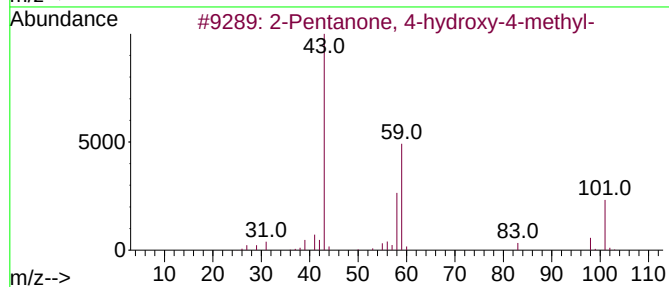
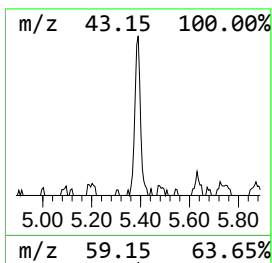
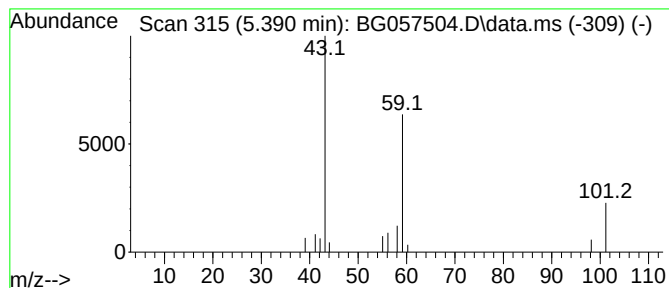
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG042723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.390	2.92 ng	18311	1,4-Dichlorobenzene-d4	8.350

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64		
2	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	33		
3	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9		
4	2,3-Dimethyloxirane-2-carboxylic...	130	C6H10O3	1000306-64-4	9		
5	2-Nonanone, 9-hydroxy-	158	C9H18O2	025368-56-3	9		



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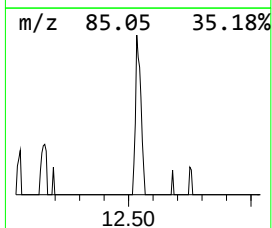
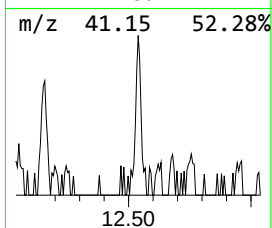
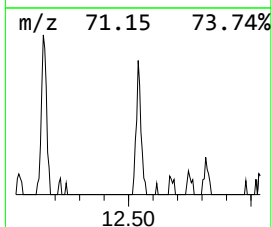
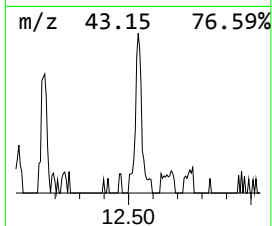
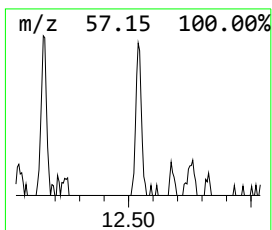
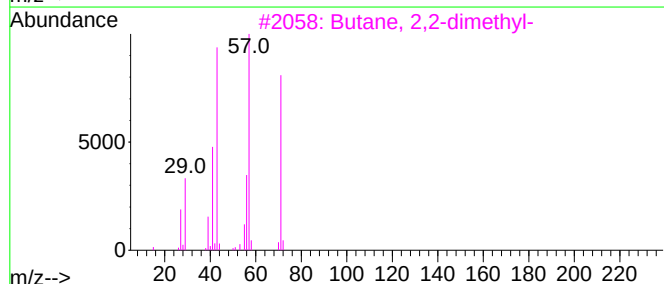
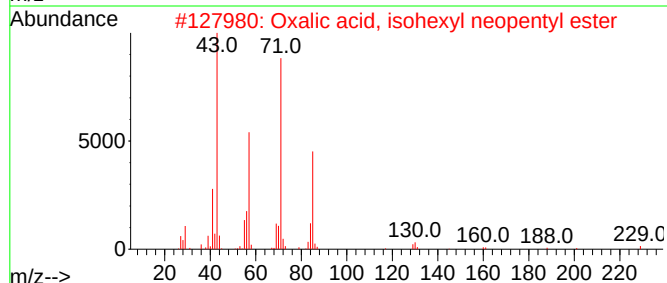
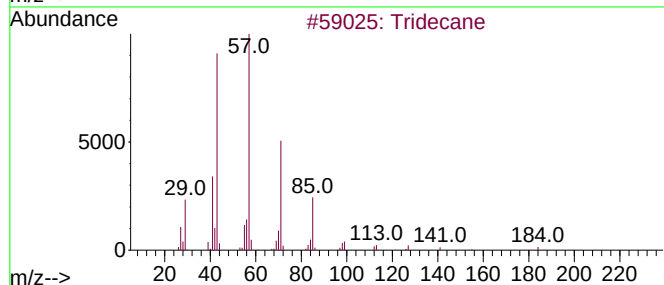
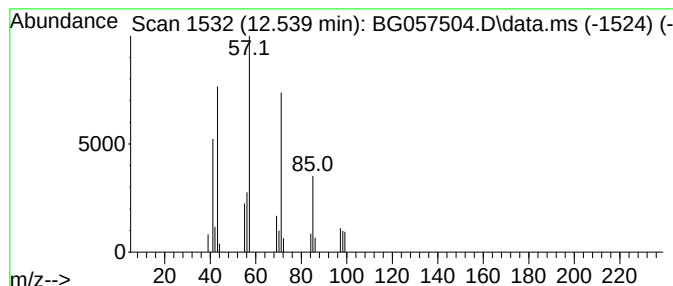
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG042723.M
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TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Tridecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.539	2.03 ng	18392	Naphthalene-d8	11.194

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane	184	C13H28	000629-50-5	72
2			Oxalic acid, isohexyl neopentyl ...	244	C13H24O4	1000309-73-0	50
3			Butane, 2,2-dimethyl-	86	C6H14	000075-83-2	50
4			Carbonic acid, nonyl prop-1-en-2...	228	C13H24O3	1000382-53-9	50
5			Hexadecane	226	C16H34	000544-76-3	45



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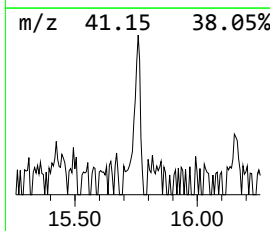
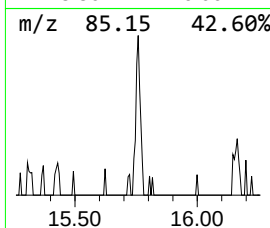
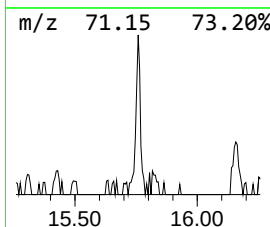
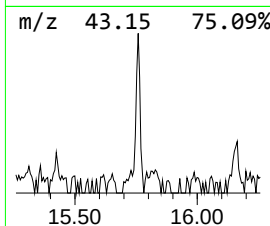
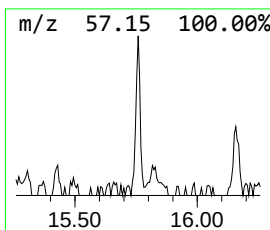
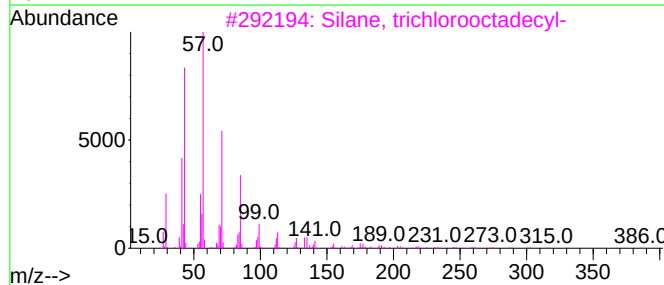
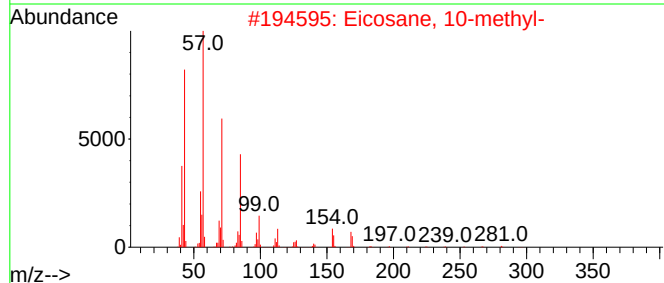
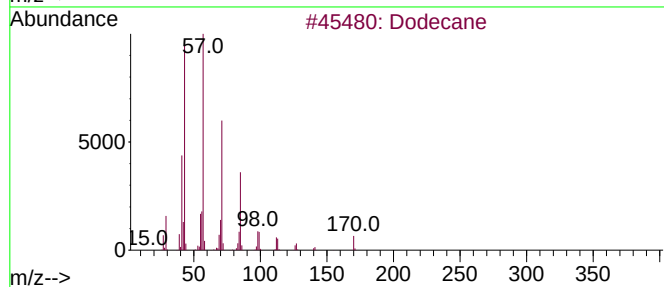
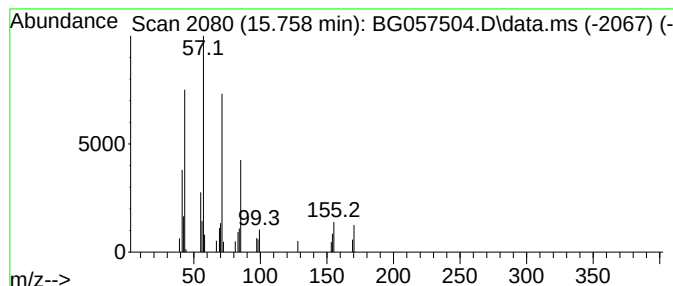
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG042723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Dodecane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.758	3.18 ng	39481	Acenaphthene-d10	14.965

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane	170	C12H26	000112-40-3	83
2			Eicosane, 10-methyl-	296	C21H44	054833-23-7	80
3			Silane, trichlorooctadecyl-	386	C18H37Cl3Si	000112-04-9	72
4			Pentacosane	352	C25H52	000629-99-2	72
5			Hexadecane	226	C16H34	000544-76-3	72



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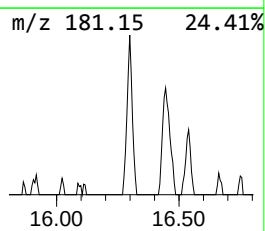
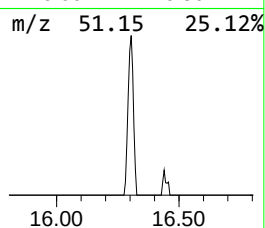
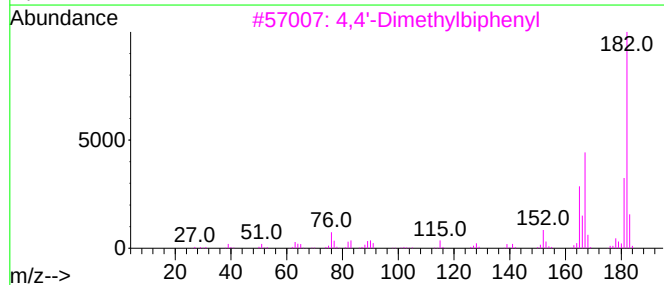
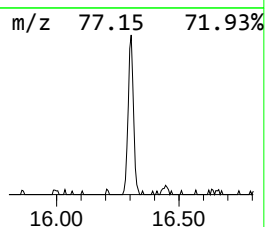
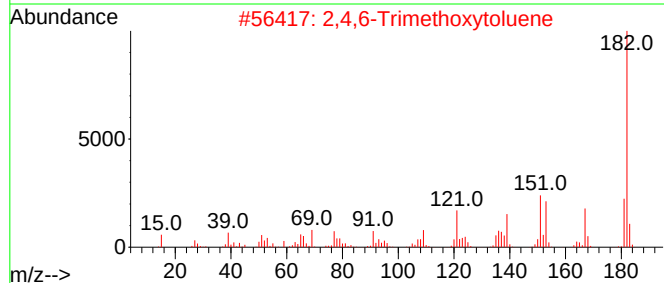
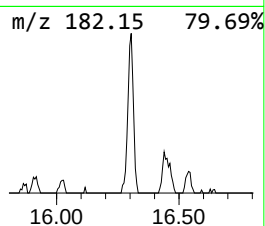
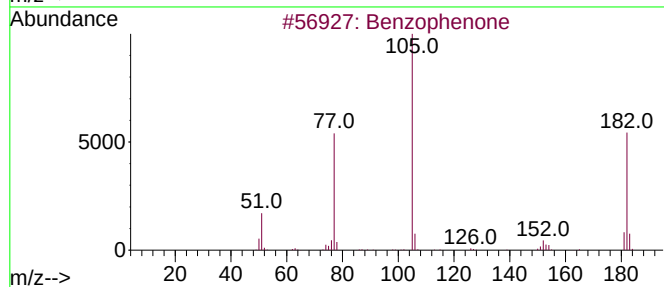
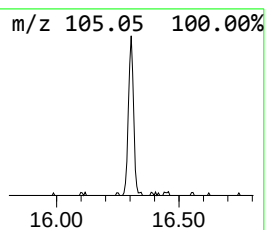
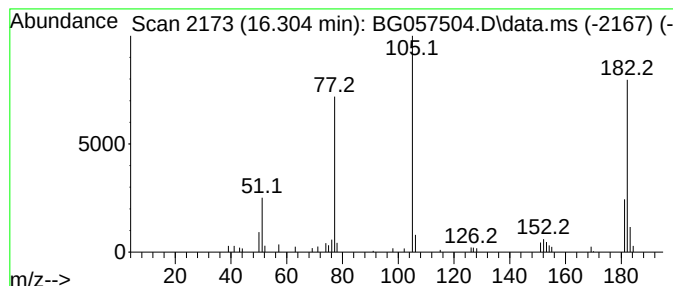
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Benzophenone Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.304	5.01 ng	62285	Acenaphthene-d10	14.965

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	94
2			2,4,6-Trimethoxytoluene	182	C10H14O3	014107-97-2	49
3			4,4'-Dimethylbiphenyl	182	C14H14	000613-33-2	43
4			Azobenzene	182	C12H10N2	000103-33-3	38
5			2-Mercapto-7-methyl-1H-imidazolo...	182	C6H6N4OS	128920-50-3	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052223\
Data File : BG057504.D
Acq On : 22 May 2023 22:06
Operator : CG/JU
Sample : 02864-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
ETGI-289

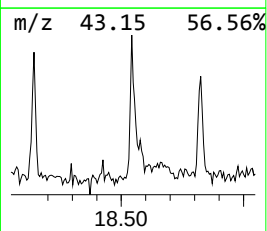
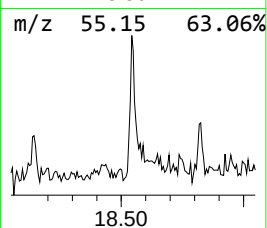
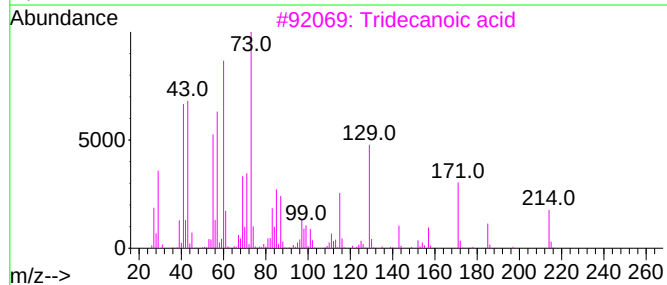
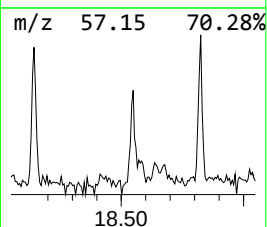
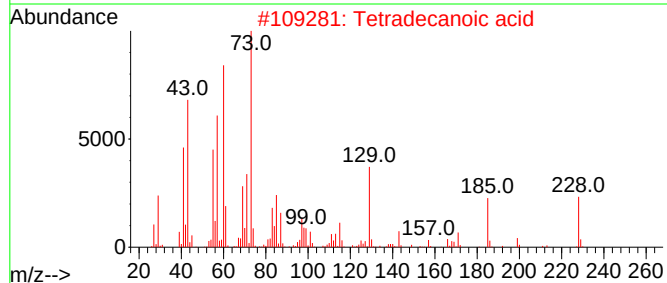
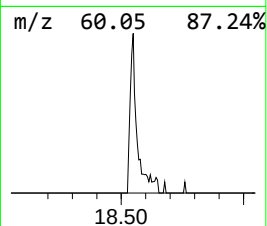
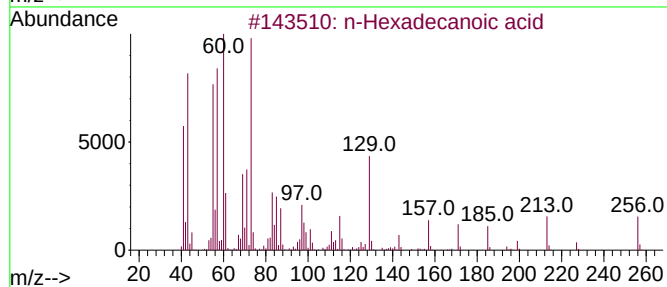
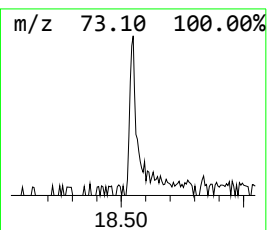
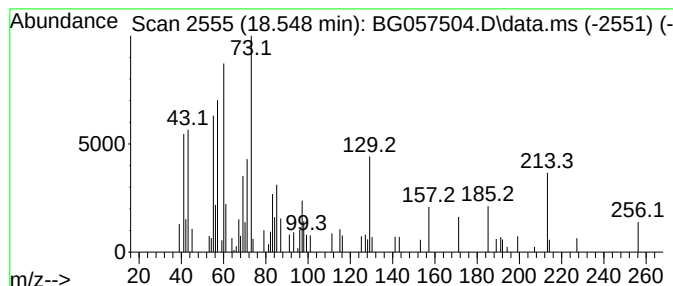
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.548	2.47 ng	40258	Phenanthrene-d10	17.708

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2			Tetradecanoic acid	228	C14H28O2	000544-63-8	74
3			Tridecanoic acid	214	C13H26O2	000638-53-9	62
4			Pentadecanoic acid	242	C15H30O2	001002-84-2	59
5			Decanoic acid, silver(1+) salt	278	C10H19AgO2	013126-67-5	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052223\
Data File : BG057504.D
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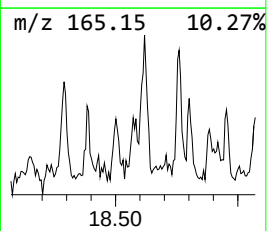
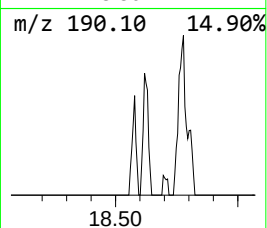
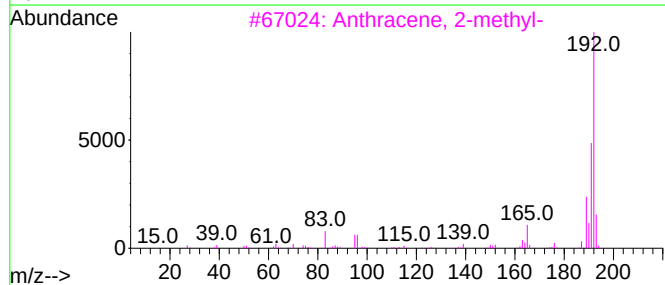
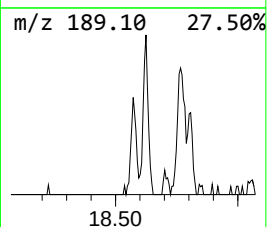
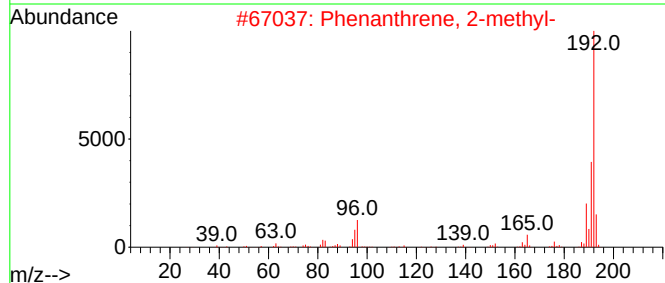
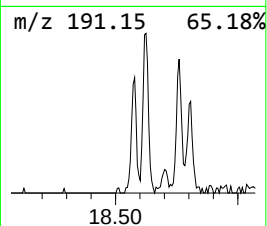
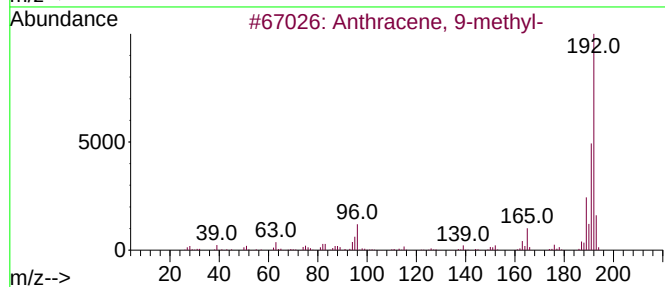
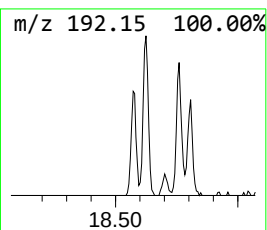
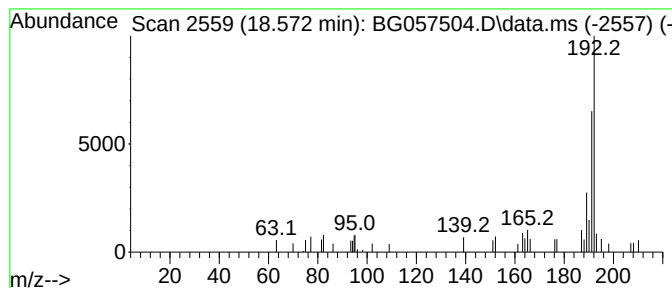
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Anthracene, 9-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.572	2.24 ng	36472	Phenanthrene-d10	17.708

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Anthracene, 9-methyl-	192	C15H12	000779-02-2	64
2		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	59
3		Anthracene, 2-methyl-	192	C15H12	000613-12-7	59
4		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	58
5		Naphtho[2,3-b]norbornadiene	192	C15H12	107426-38-0	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052223\
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Sample : 02864-01
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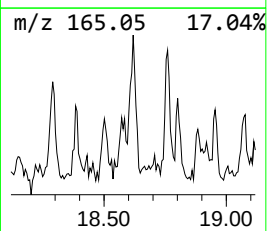
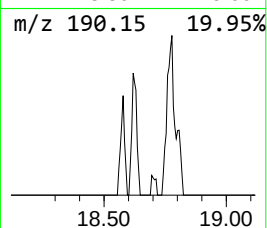
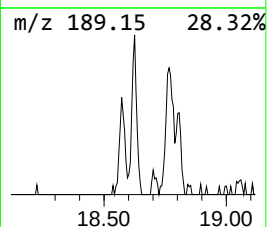
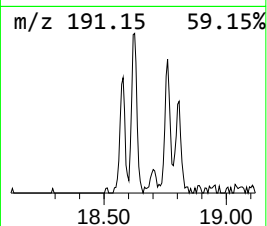
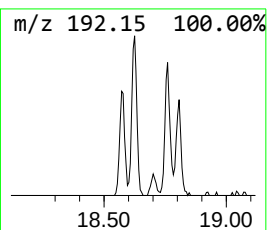
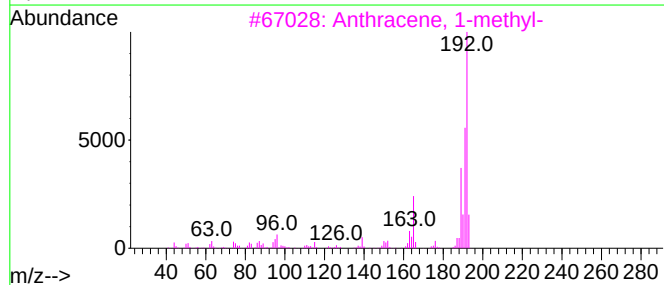
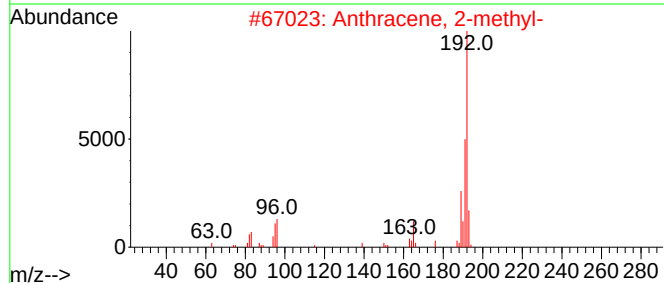
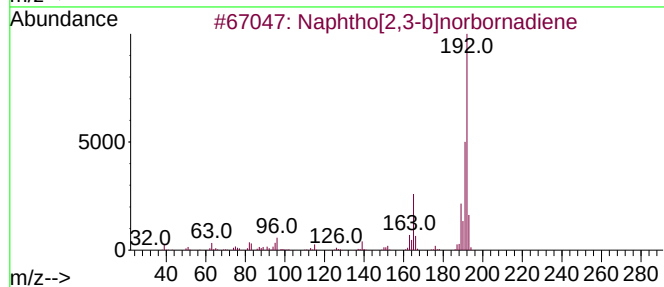
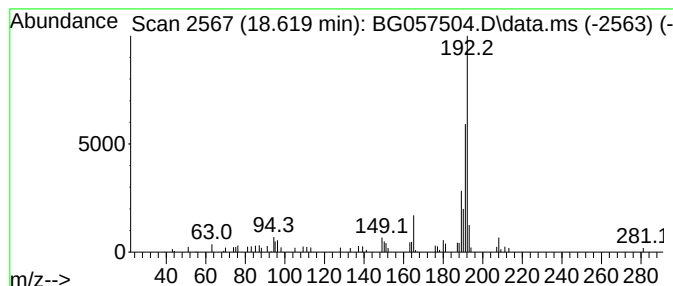
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Naphtho[2,3-b]norbornadiene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.619	3.77 ng	61493	Phenanthrene-d10	17.708	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphtho[2,3-b]norbornadiene	192	C15H12	107426-38-0	76
2	Anthracene, 2-methyl-	192	C15H12	000613-12-7	76
3	Anthracene, 1-methyl-	192	C15H12	000610-48-0	76
4	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	76
5	Anthracene, 9-methyl-	192	C15H12	000779-02-2	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052223\
Data File : BG057504.D
Acq On : 22 May 2023 22:06
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Sample : 02864-01
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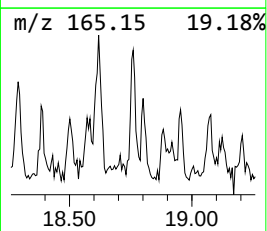
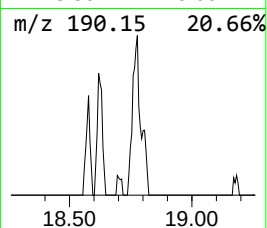
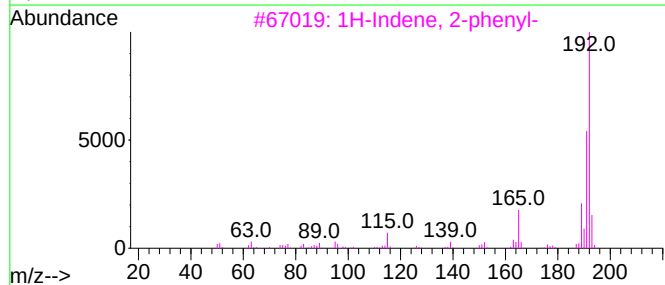
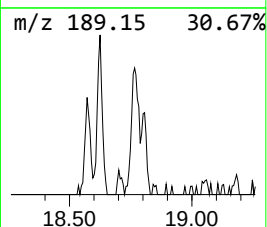
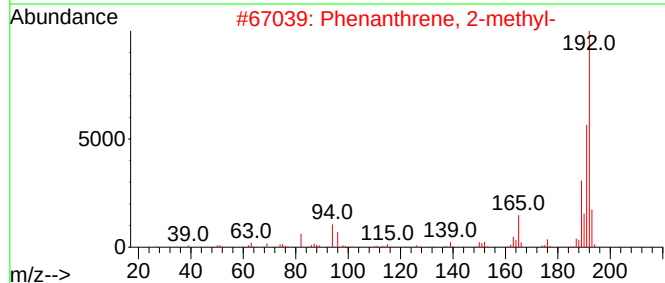
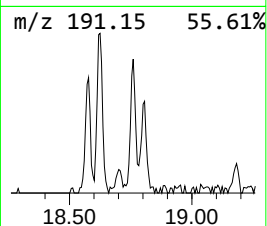
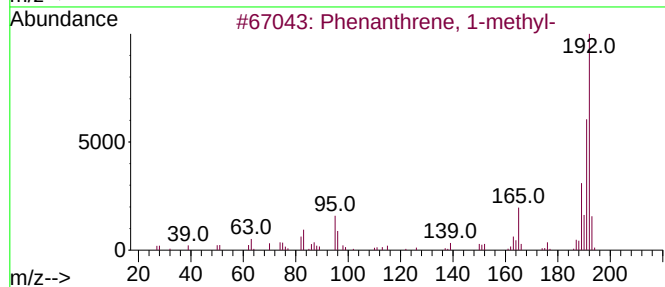
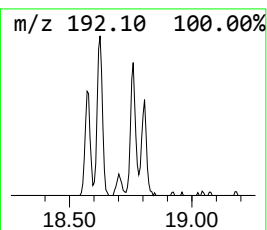
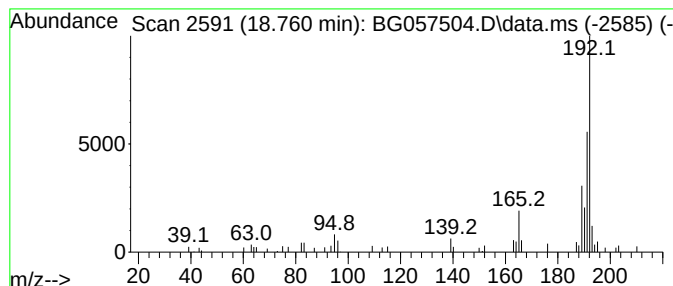
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Phenanthrene, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.760	2.70 ng	43981	Phenanthrene-d10	17.708

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	86
2		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	76
3		1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	76
4		Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	70
5		Anthracene, 1-methyl-	192	C15H12	000610-48-0	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052223\
Data File : BG057504.D
Acq On : 22 May 2023 22:06
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Sample : 02864-01
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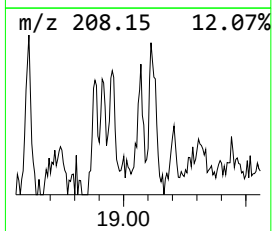
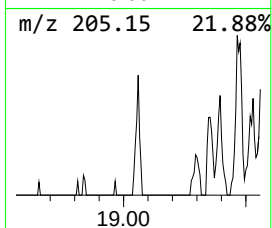
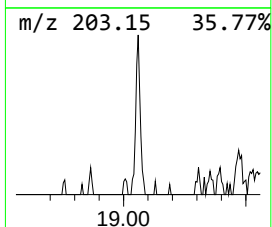
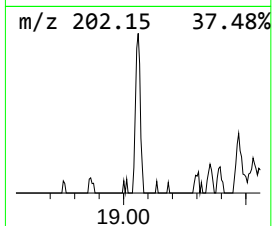
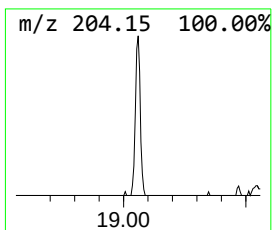
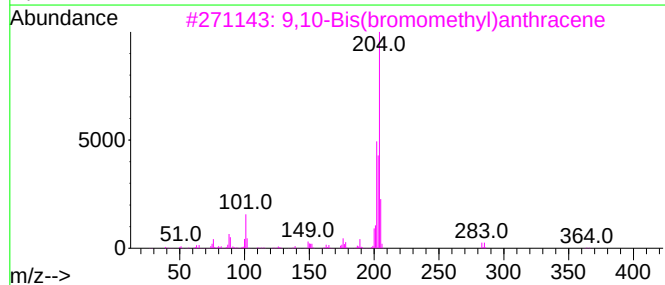
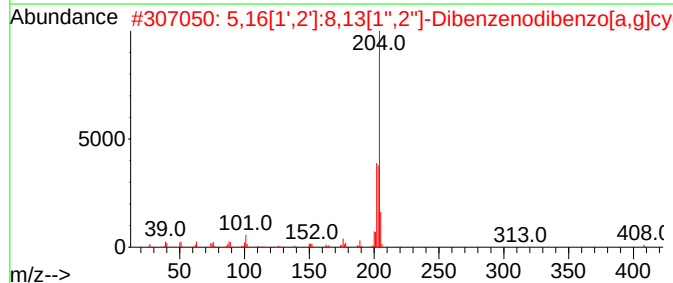
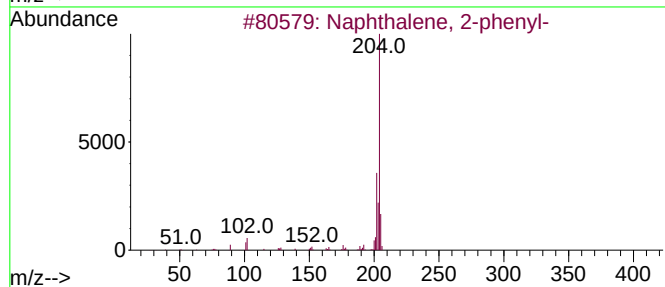
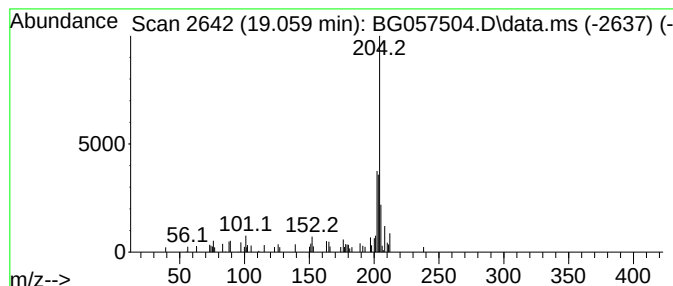
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Naphthalene, 2-phenyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.060	2.44 ng	39827	Phenanthrene-d10	17.708

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	70
2			5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	64
3			9,10-Bis(bromomethyl)anthracene	362	C16H12Br2	034373-96-1	64
4			Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	62
5			[1,1'-Biphenyl]-4,4'-dicarbonitrile	204	C14H8N2	001591-30-6	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052223\
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ALS Vial : 11 Sample Multiplier: 1

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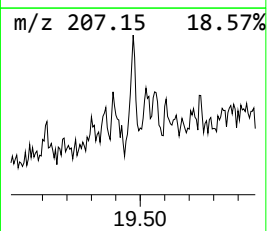
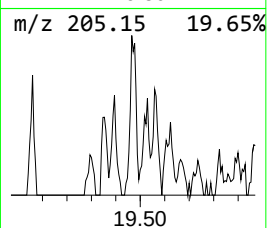
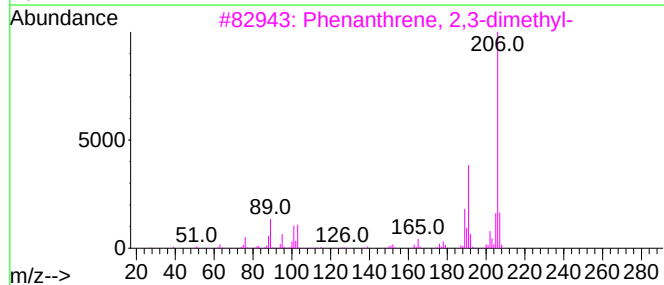
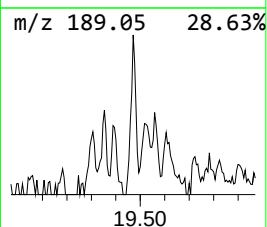
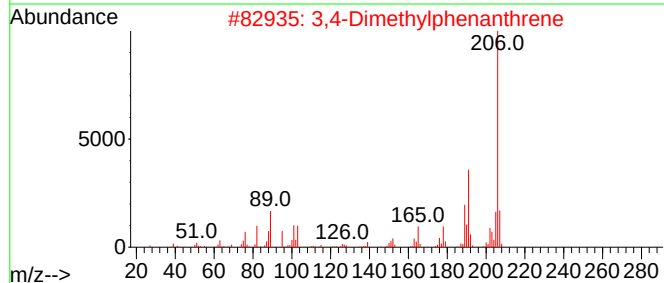
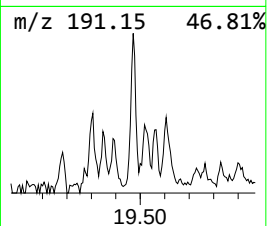
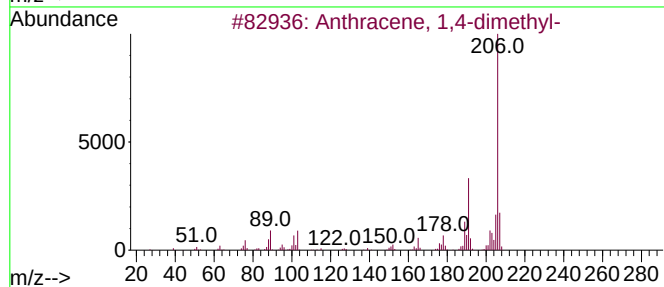
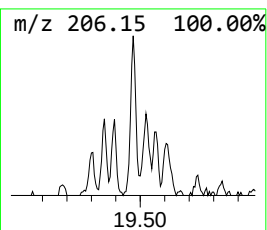
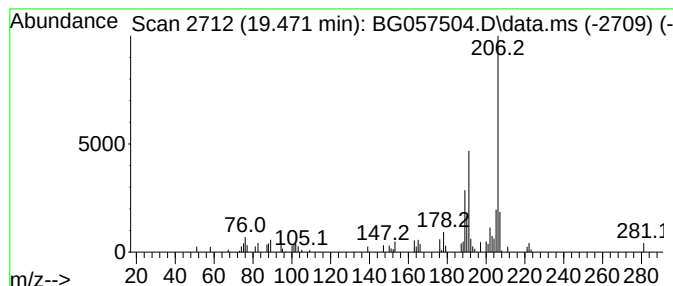
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Anthracene, 1,4-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.471	2.38 ng	38727	Phenanthrene-d10	17.708

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	91
2		3,4-Dimethylphenanthrene	206	C16H14	1000452-24-5	87
3		Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	87
4		Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	83
5		di-p-Tolylacetylene	206	C16H14	002789-88-0	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052223\
Data File : BG057504.D
Acq On : 22 May 2023 22:06
Operator : CG/JU
Sample : 02864-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

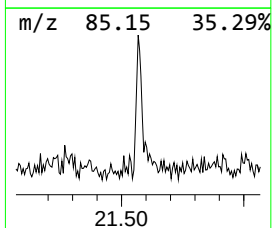
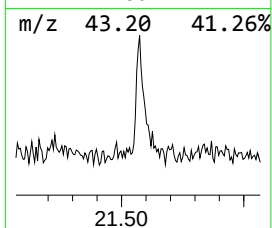
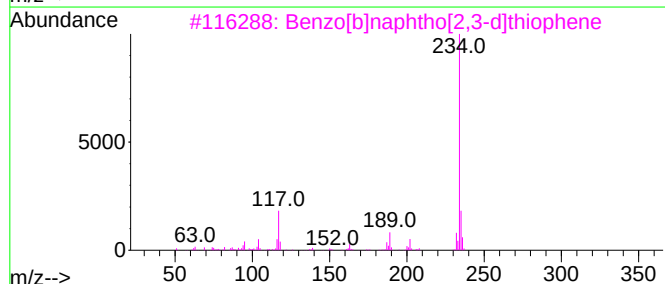
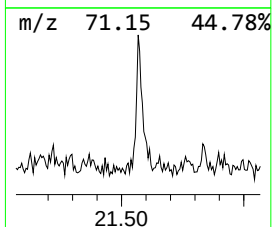
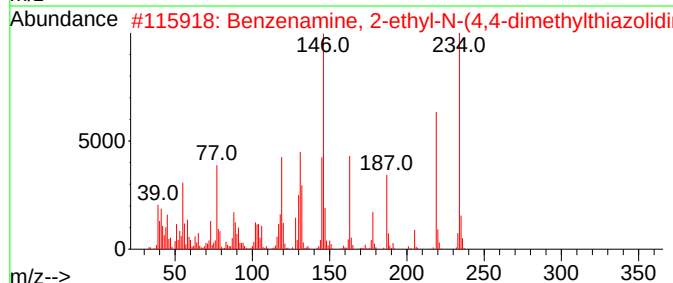
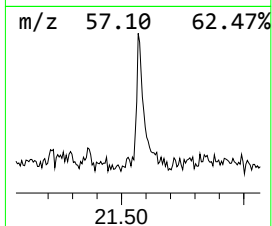
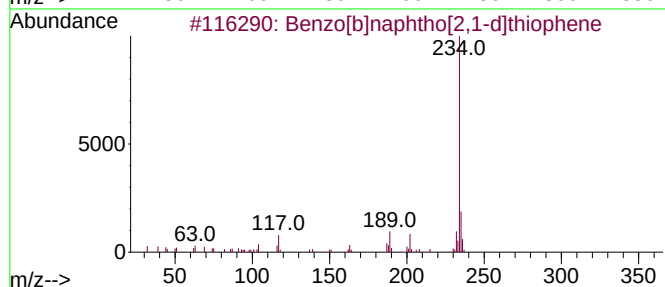
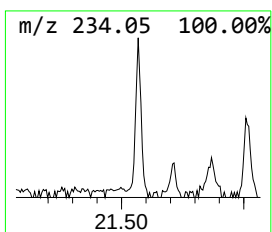
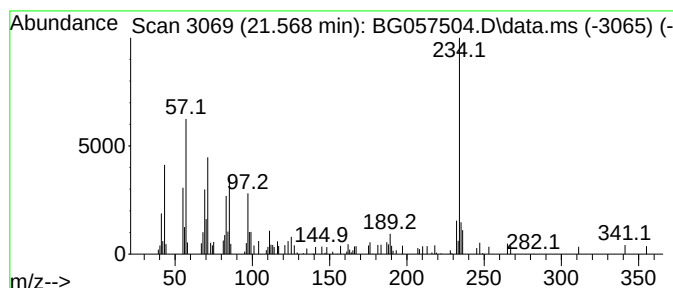
Instrument :
BNA_G
ClientSampleId :
ETGI-289

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG042723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Benzo[b]naphtho[2,1-d]thiop... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.568	2.64 ng	57044	Chrysene-d12	22.014	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[b]naphtho[2,1-d]thiophene	234	C16H10S	000239-35-0	50
2	Benzenamine, 2-ethyl-N-(4,4-dime...	234	C13H18N2S	1000272-26-2	42
3	Benzo[b]naphtho[2,3-d]thiophene	234	C16H10S	000243-46-9	41
4	Benzo[b]naphtho[1,2-d]thiophene	234	C16H10S	000205-43-6	41
5	Thiazol-2-amine, 4-(1-adamantyl)-	234	C13H18N2S	019735-74-1	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052223\
Data File : BG057504.D
Acq On : 22 May 2023 22:06
Operator : CG/JU
Sample : 02864-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
ETGI-289

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG042723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	5.390	2.9	ng	18311	1	8.350	125394	20.0
Tridecane	12.539	2.0	ng	18392	2	11.194	181035	20.0
Dodecane	15.758	3.2	ng	39481	3	14.965	248490	20.0
Benzophenone	16.304	5.0	ng	62285	3	14.965	248490	20.0
n-Hexadecanoic ...	18.548	2.5	ng	40258	4	17.708	326012	20.0
Anthracene, 9-m...	18.572	2.2	ng	36472	4	17.708	326012	20.0
Naphtho[2,3-b]n...	18.619	3.8	ng	61493	4	17.708	326012	20.0
Phenanthrene, 1...	18.760	2.7	ng	43981	4	17.708	326012	20.0
Naphthalene, 2-...	19.060	2.4	ng	39827	4	17.708	326012	20.0
Anthracene, 1,4...	19.471	2.4	ng	38727	4	17.708	326012	20.0
Benzo[b]naphtho...	21.568	2.6	ng	57044	5	22.014	432385	20.0