

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062023\
 Data File : BG057777.D
 Acq On : 20 Jun 2023 15:30
 Operator : CG/JU
 Sample : 03289-02
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SB-01-4.0-6.0

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-BG061923.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG057777.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.039	325	332	340	rBV	109378	171636	14.42%	1.272%
2	5.503	404	411	422	rVB	391771	630942	53.02%	4.675%
3	7.089	673	681	691	rBV	445351	745447	62.65%	5.523%
4	7.929	816	824	832	rVB	166543	274931	23.11%	2.037%
5	9.087	1014	1021	1029	rBV	262719	453654	38.12%	3.361%
6	10.732	1292	1301	1308	rBV	222640	401695	33.76%	2.976%
7	13.188	1712	1719	1729	rVB	614312	936766	78.73%	6.941%
8	14.287	1900	1906	1915	rBV2	10263	21353	1.79%	0.158%
9	14.563	1945	1953	1959	rBV	334977	506126	42.53%	3.750%
10	14.627	1959	1964	1969	rVB2	18931	28951	2.43%	0.215%
11	14.962	2015	2021	2027	rBV2	17755	27630	2.32%	0.205%
12	15.609	2126	2131	2137	rBV	20526	30620	2.57%	0.227%
13	15.914	2177	2183	2188	rVB2	88035	135622	11.40%	1.005%
14	16.043	2197	2205	2214	rBV	385148	592724	49.81%	4.392%
15	16.272	2240	2244	2251	rVB9	8885	17750	1.49%	0.132%
16	16.831	2334	2339	2344	rBV7	6851	13656	1.15%	0.101%
17	16.954	2356	2360	2366	rVB2	24635	39412	3.31%	0.292%
18	17.124	2384	2389	2395	rVB	25913	36199	3.04%	0.268%
19	17.307	2414	2420	2423	rBV	381990	580392	48.78%	4.300%
20	17.348	2423	2427	2433	rVV	407139	584444	49.12%	4.330%
21	17.436	2437	2442	2446	rVV	74734	113730	9.56%	0.843%
22	17.489	2448	2451	2457	rVB5	10814	19596	1.65%	0.145%
23	17.706	2484	2488	2492	rVB	34138	45628	3.83%	0.338%
24	17.882	2516	2518	2523	rVB3	15271	19573	1.64%	0.145%
25	18.194	2564	2571	2574	rBV2	135119	252270	21.20%	1.869%
26	18.229	2574	2577	2582	rVV2	89135	131255	11.03%	0.973%
27	18.305	2586	2590	2595	rVV	26390	46010	3.87%	0.341%
28	18.370	2596	2601	2606	rVV3	86978	186338	15.66%	1.381%
29	18.411	2606	2608	2614	rVB	47598	59341	4.99%	0.440%
30	18.681	2649	2654	2657	rBV	47887	72823	6.12%	0.540%
31	18.717	2657	2660	2666	rVB3	45937	68248	5.74%	0.506%
32	18.922	2693	2695	2699	rVB2	18938	20043	1.68%	0.149%
33	19.099	2720	2725	2729	rBV2	56549	103726	8.72%	0.769%
34	19.234	2744	2748	2757	rBV	58363	114647	9.63%	0.849%
35	19.357	2764	2769	2776	rVB	798438	1109369	93.23%	8.220%
36	19.692	2822	2826	2827	rBV2	104071	138242	11.62%	1.024%

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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

Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG061923.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	19.721	2827	2831	2838	rVB	721142	1031347	86.67%	7.642%
38	19.921	2860	2865	2869	rBV	821821	1043424	87.69%	7.731%
39	21.549	3136	3142	3148	rBV4	458992	1189918	100.00%	8.817%
40	21.601	3148	3151	3158	rVB	330089	530197	44.56%	3.928%
41	23.717	3505	3511	3518	rBV	185658	550357	46.25%	4.078%
42	24.569	3651	3656	3666	rVB	164370	420195	35.31%	3.113%

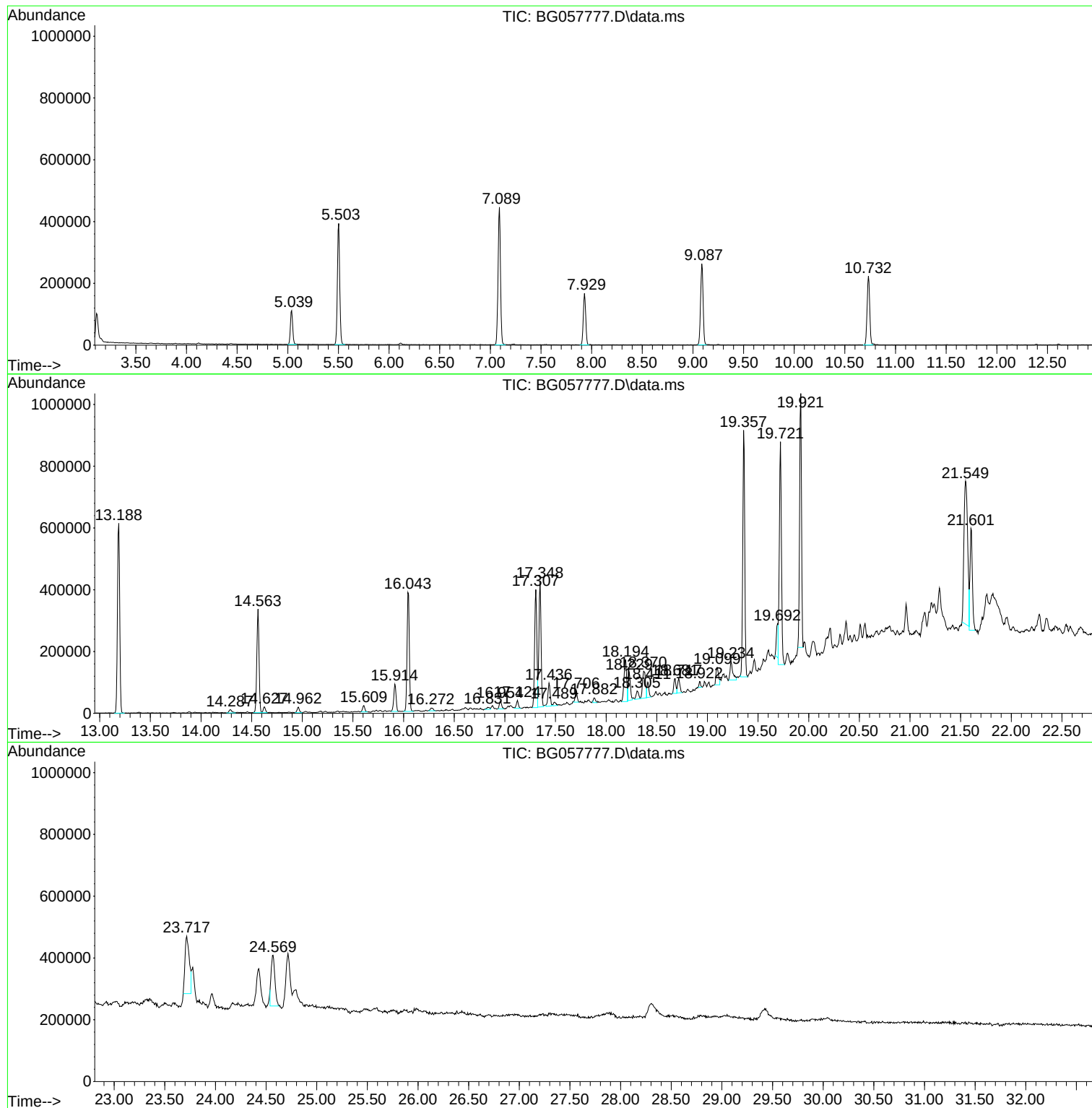
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG061923.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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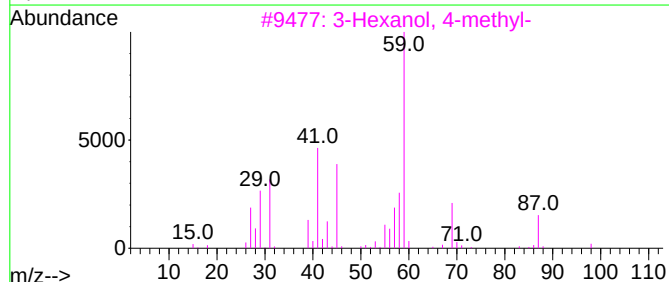
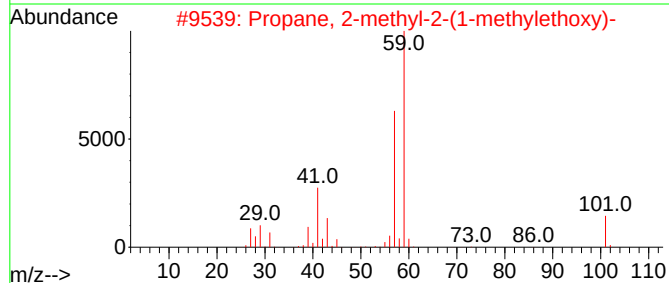
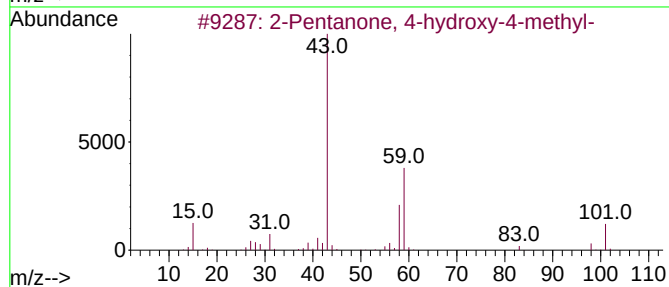
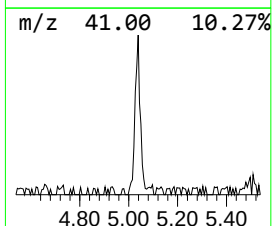
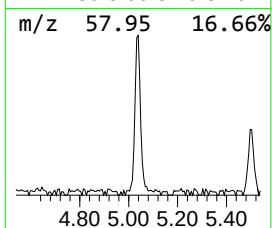
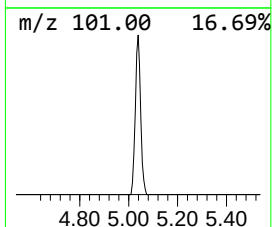
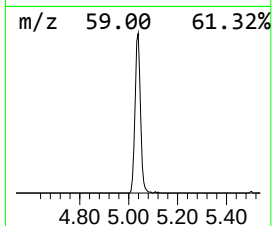
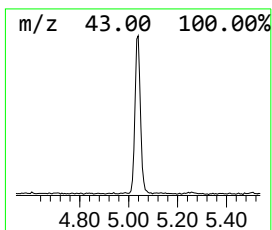
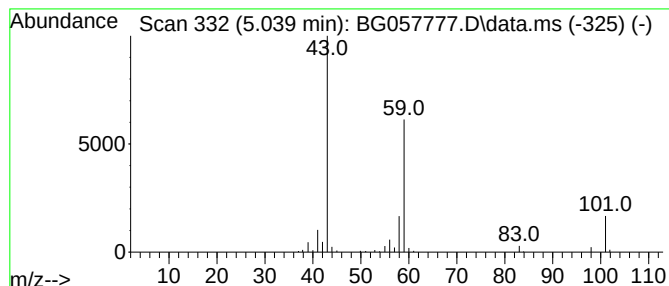
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG061923.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.039	12.49 ng	171636	1,4-Dichlorobenzene-d4	7.929

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53	
2	Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	36	
3	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33	
4	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	32	
5	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28	



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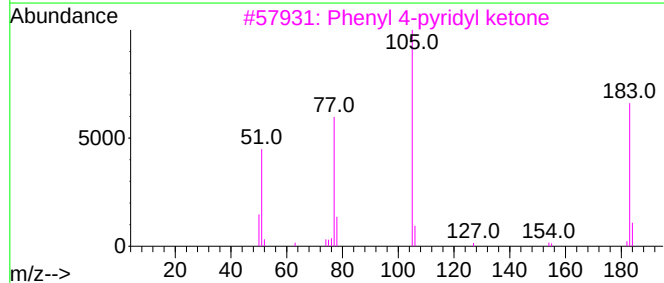
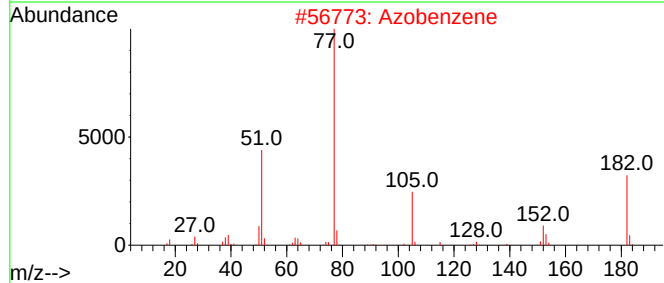
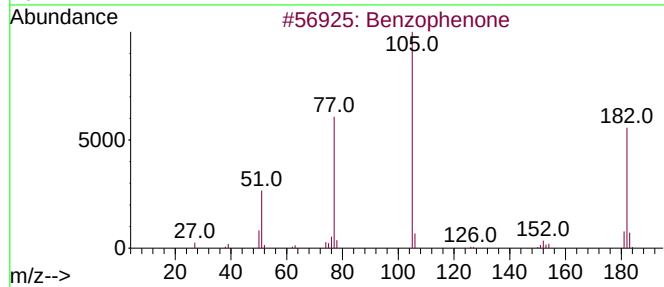
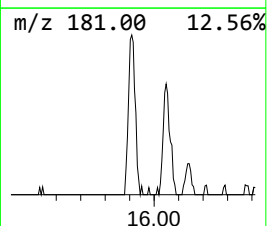
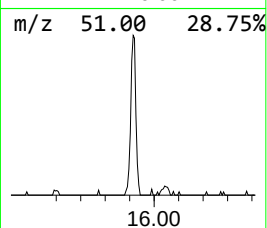
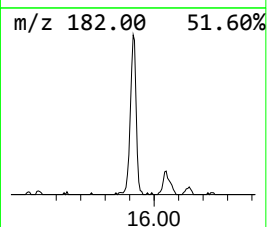
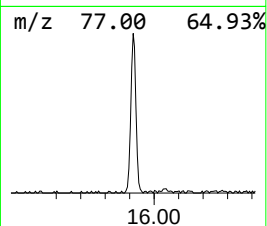
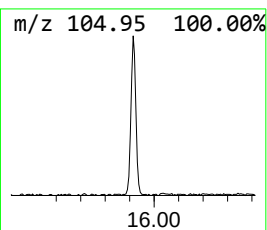
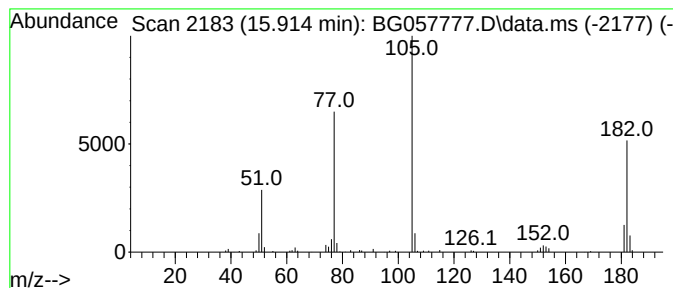
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG061923.M
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TIC Library : C:\Database\NIST20.L
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Peak Number 2 Benzophenone Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.914	5.36 ng	135622	Acenaphthene-d10	14.563

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Azobenzene	182	C12H10N2	000103-33-3	74
3			Phenyl 4-pyridyl ketone	183	C12H9NO	014548-46-0	58
4			Benzenepropanenitrile, .beta.-oxo-	145	C9H7NO	000614-16-4	49
5			N-Methoxy-N-methylbenzamide	165	C9H11NO2	006919-61-5	47



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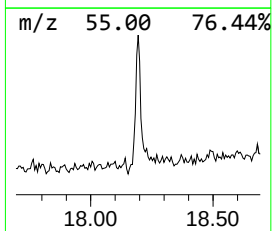
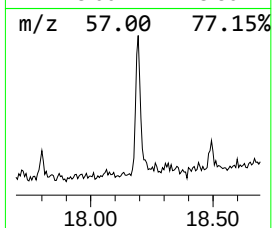
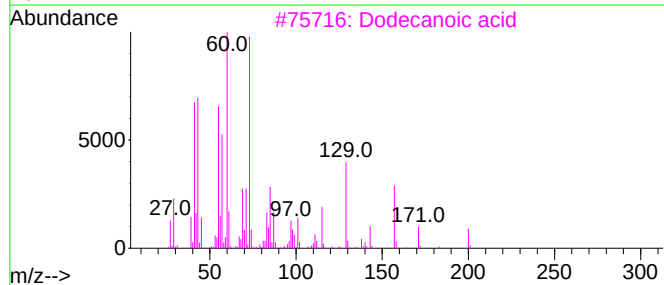
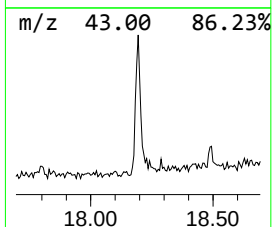
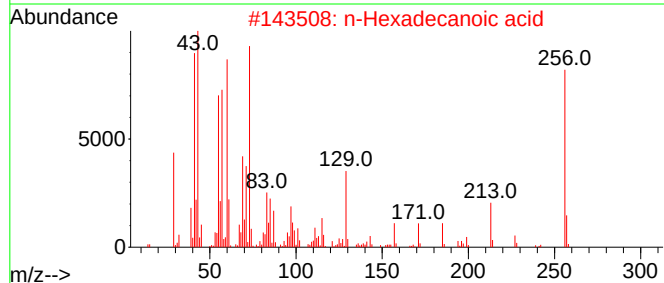
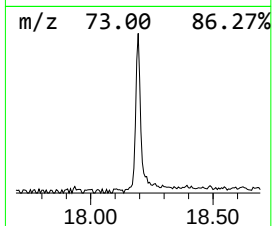
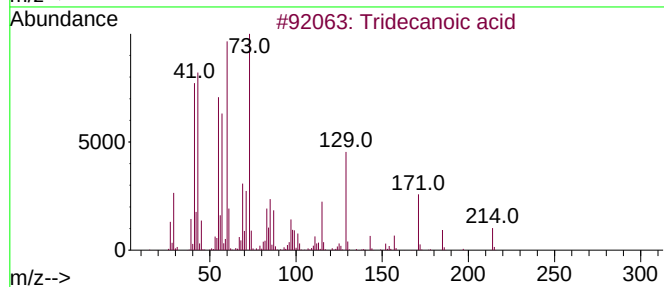
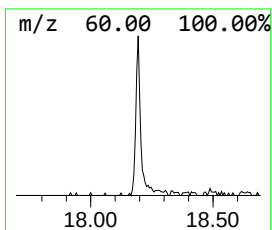
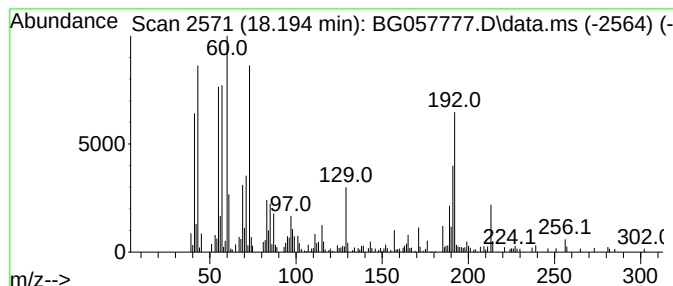
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG061923.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Tridecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.194	8.69 ng	252270	Phenanthrene-d10	17.307

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecanoic acid	214	C13H26O2	000638-53-9	89
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	86
3			Dodecanoic acid	200	C12H24O2	000143-07-7	45
4			Octadecanoic acid	284	C18H36O2	000057-11-4	42
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	38



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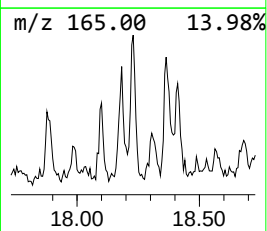
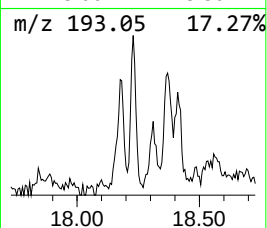
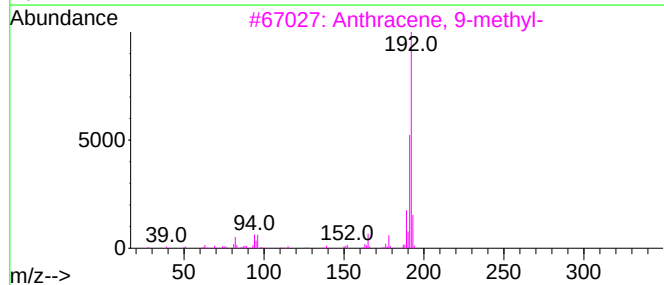
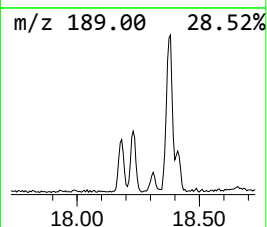
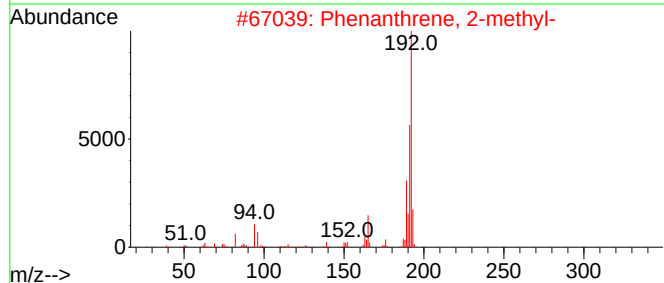
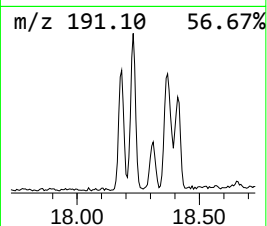
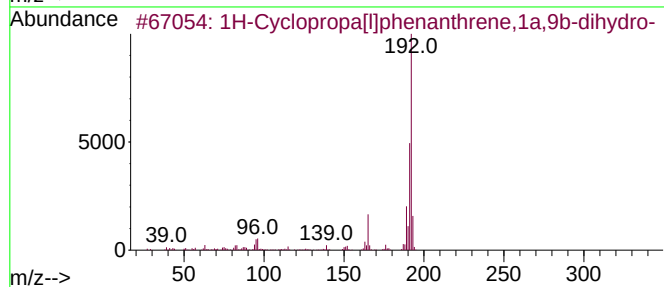
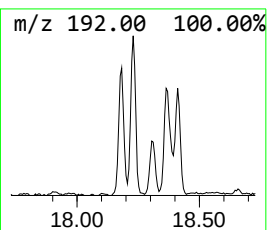
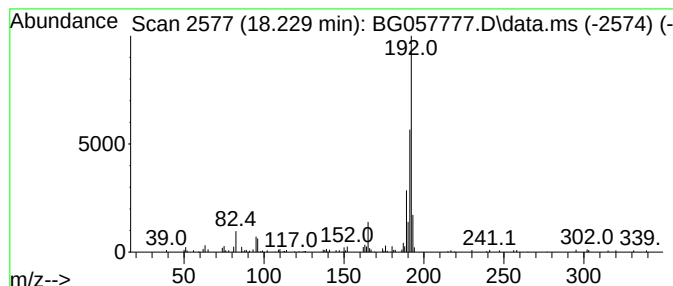
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 1H-Cyclopropa[l]phenanthren... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.229	4.52 ng	131255	Phenanthrene-d10	17.307

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	96
2		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
3		Anthracene, 9-methyl-	192	C15H12	000779-02-2	95
4		Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	94
5		Phenanthrene, 3-methyl-	192	C15H12	000832-71-3	94



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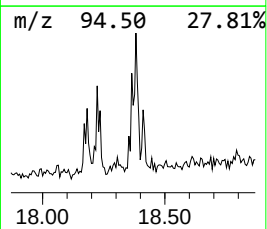
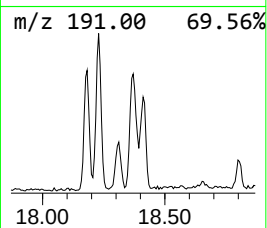
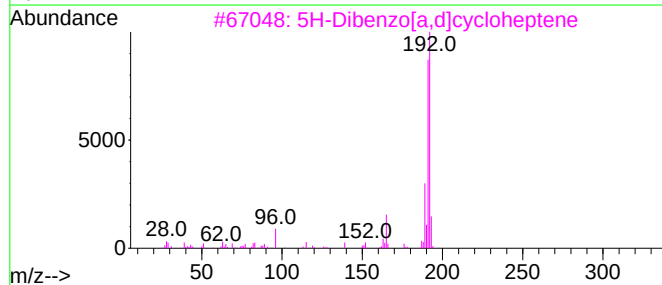
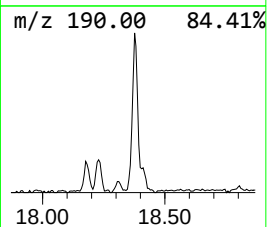
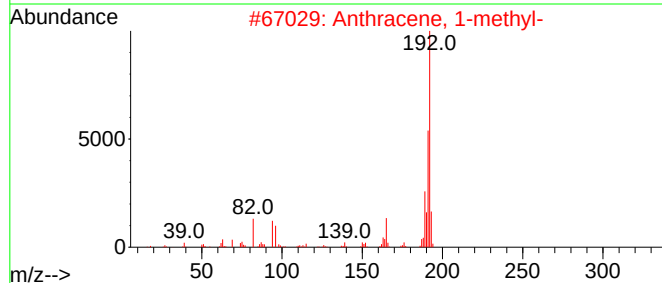
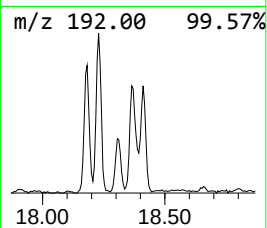
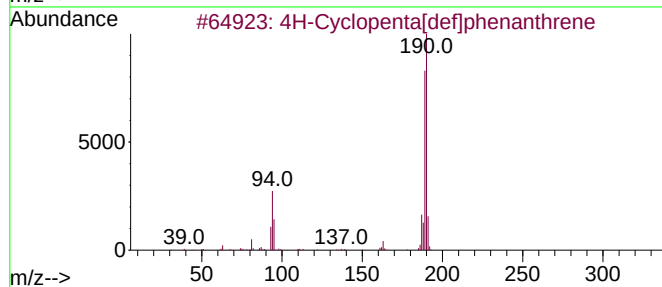
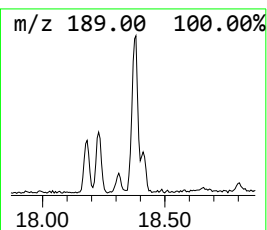
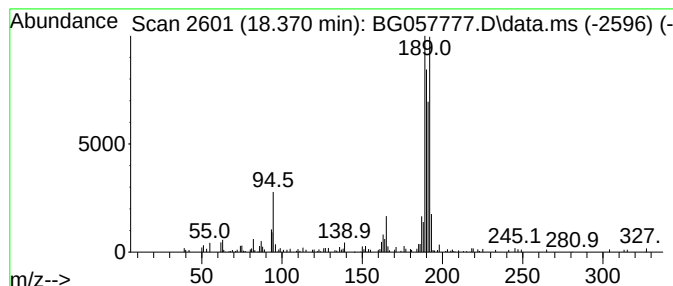
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG061923.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 unknown18.370 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.370	6.42 ng	186338	Phenanthrene-d10	17.307

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	49
2			Anthracene, 1-methyl-	192	C15H12	000610-48-0	46
3			5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	42
4			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	42
5			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	42



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062023\
Data File : BG057777.D
Acq On : 20 Jun 2023 15:30
Operator : CG/JU
Sample : 03289-02
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SB-01-4.0-6.0

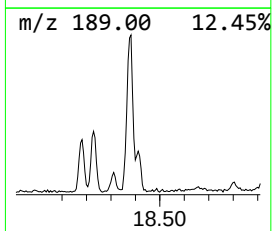
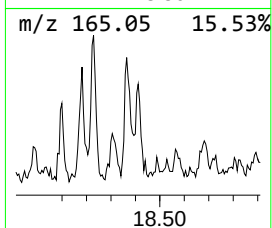
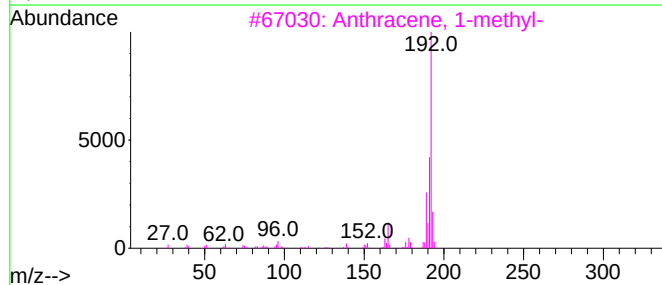
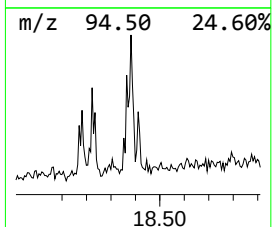
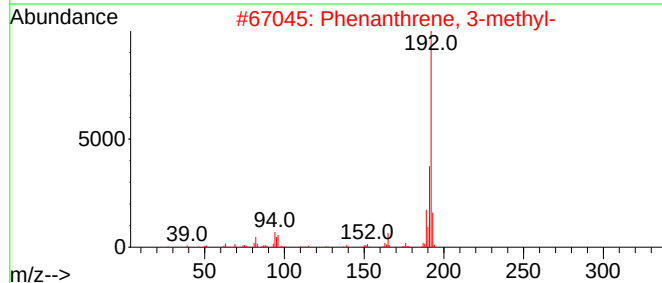
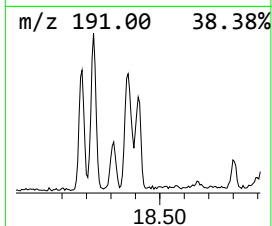
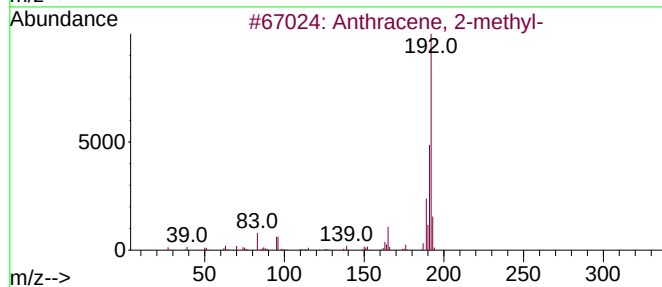
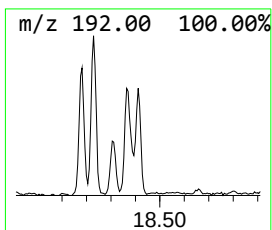
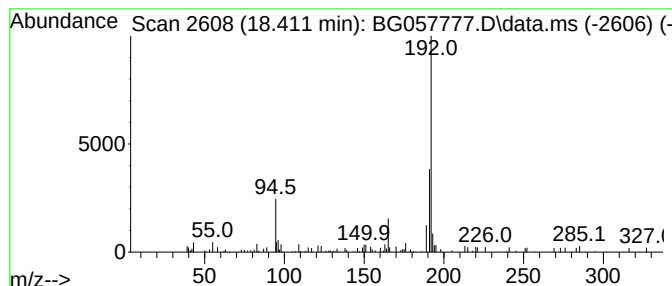
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Anthracene, 2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.411	2.04 ng	59341	Phenanthrene-d10	17.307

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene, 2-methyl-	192	C15H12	000613-12-7	64
2			Phenanthrene, 3-methyl-	192	C15H12	000832-71-3	64
3			Anthracene, 1-methyl-	192	C15H12	000610-48-0	64
4			1H-Indene, 1-phenyl-	192	C15H12	001961-96-2	64
5			Anthracene, 9-methyl-	192	C15H12	000779-02-2	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062023\
Data File : BG057777.D
Acq On : 20 Jun 2023 15:30
Operator : CG/JU
Sample : 03289-02
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ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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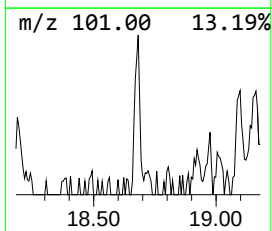
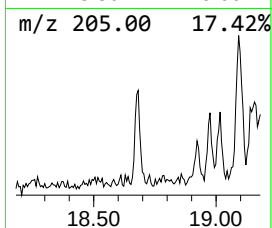
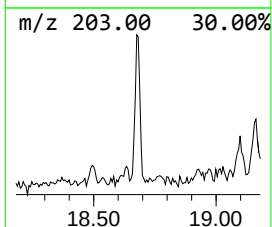
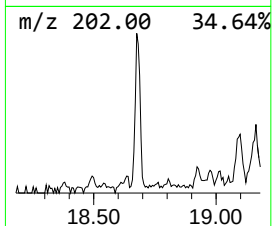
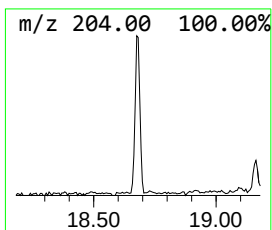
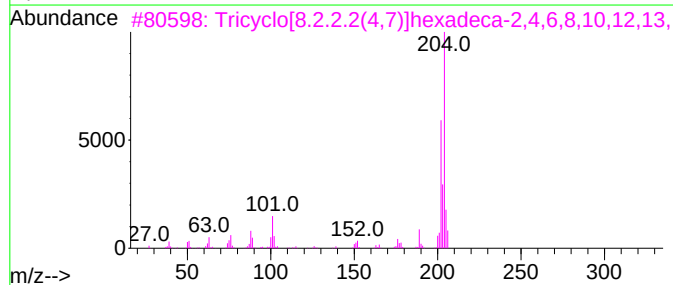
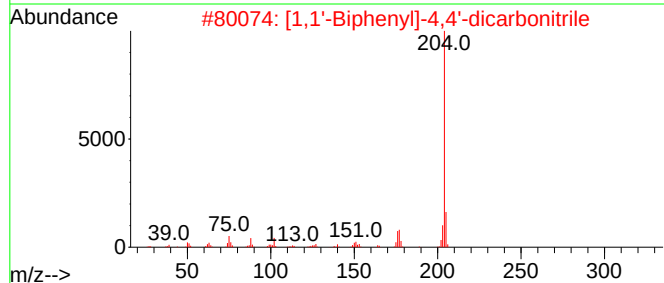
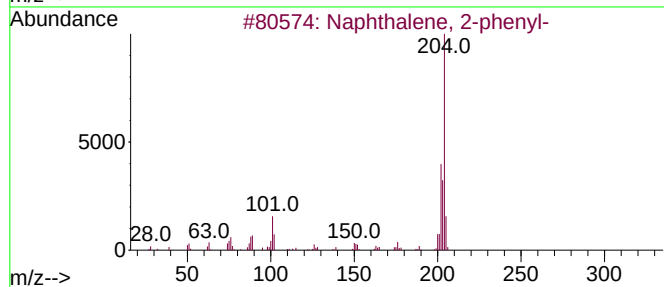
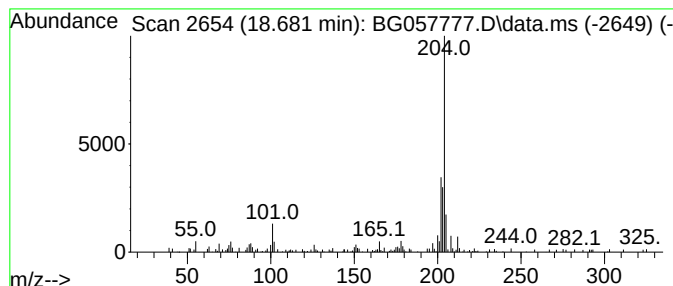
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Naphthalene, 2-phenyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.681	2.51 ng	72823	Phenanthrene-d10	17.307

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	76
2			[1,1'-Biphenyl]-4,4'-dicarbonitrile	204	C14H8N2	001591-30-6	60
3			Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	59
4			6-Cyanophenanthridine	204	C14H8N2	002176-28-5	53
5			Naphthalene, 1-phenyl-	204	C16H12	000605-02-7	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062023\
Data File : BG057777.D
Acq On : 20 Jun 2023 15:30
Operator : CG/JU
Sample : 03289-02
Misc :
ALS Vial : 12 Sample Multiplier: 1

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ClientSampleId :
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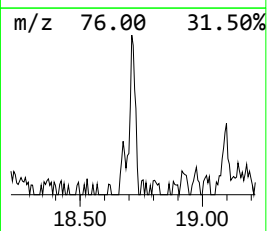
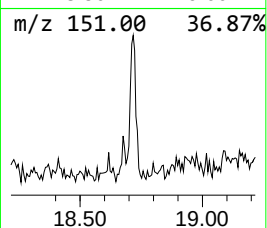
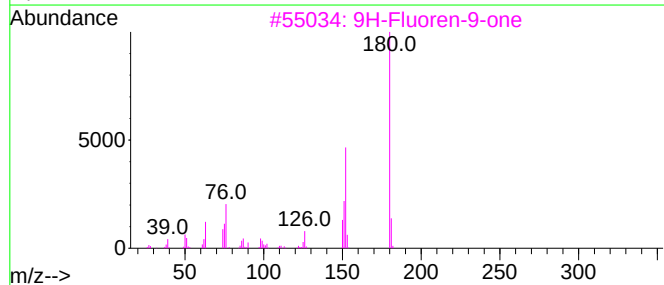
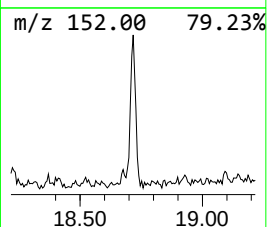
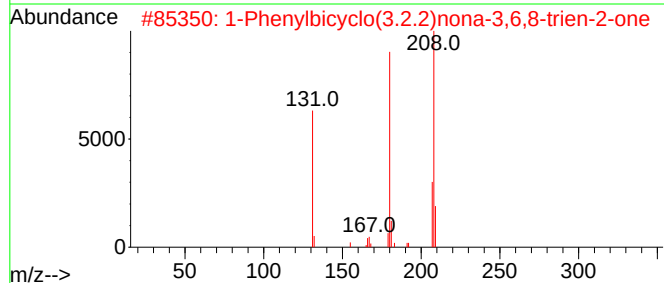
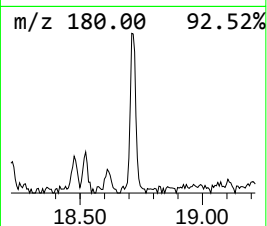
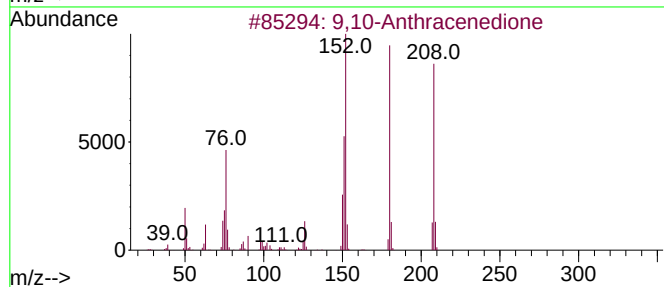
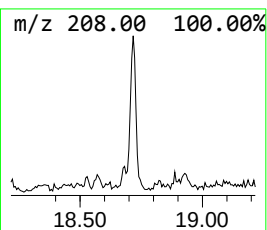
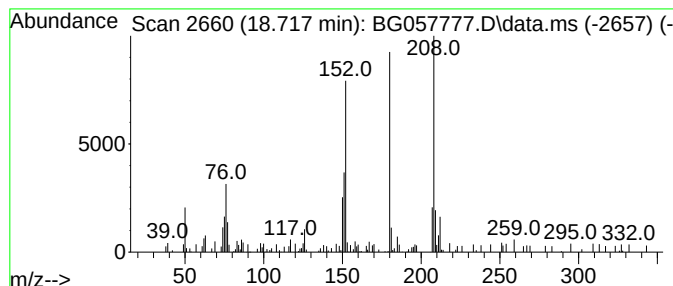
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 9,10-Anthracenedione Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.717	2.35 ng	68248	Phenanthrene-d10	17.307

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9,10-Anthracenedione	208	C14H8O2	000084-65-1	93
2			1-Phenylbicyclo(3.2.2)nona-3,6,8...	208	C15H12O	1010427-19-9	53
3			9H-Fluoren-9-one	180	C13H8O	000486-25-9	53
4			Benzo[c]cinnoline	180	C12H8N2	000230-17-1	49
5			5,6-Dimethyl-1,10-phenanthroline	208	C14H12N2	003002-81-1	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062023\
Data File : BG057777.D
Acq On : 20 Jun 2023 15:30
Operator : CG/JU
Sample : 03289-02
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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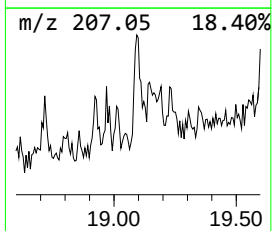
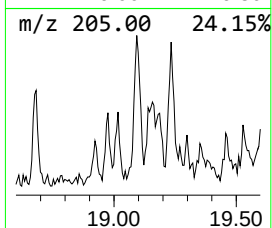
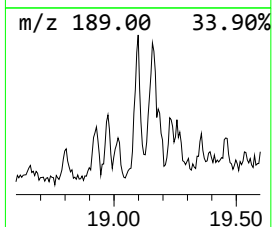
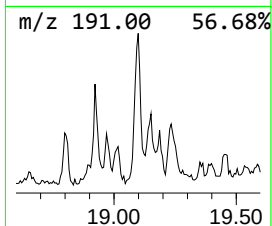
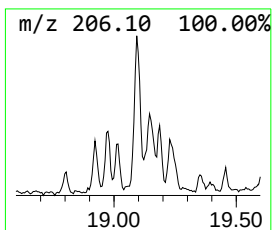
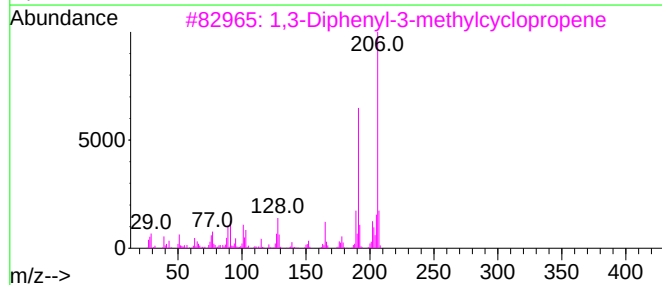
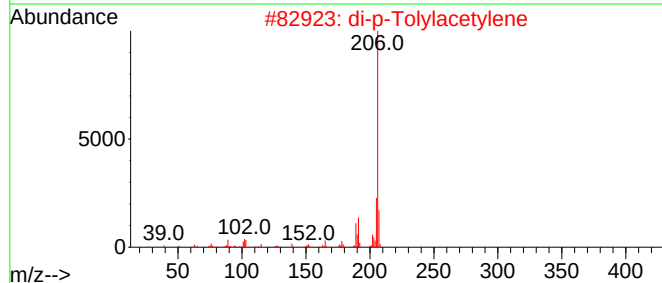
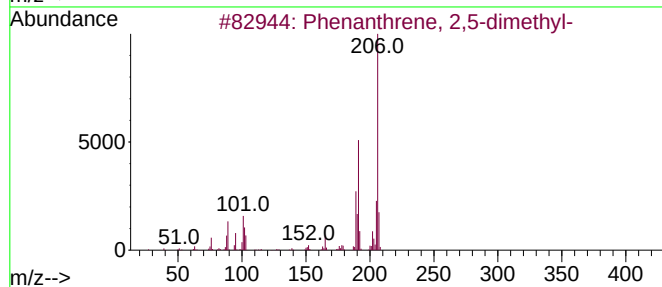
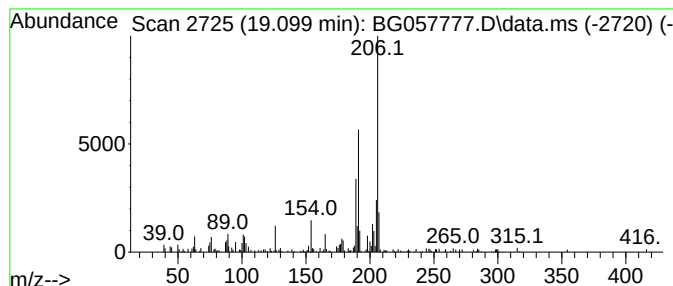
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.099	3.57 ng	103726	Phenanthrene-d10	17.307

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	93
2			di-p-Tolylacetylene	206	C16H14	002789-88-0	91
3			1,3-Diphenyl-3-methylcyclopropene	206	C16H14	086544-79-8	90
4			9,10-Dimethylantracene	206	C16H14	000781-43-1	90
5			3-Methyl-1-phenyl-1H-indene	206	C16H14	022360-63-0	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062023\
Data File : BG057777.D
Acq On : 20 Jun 2023 15:30
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ALS Vial : 12 Sample Multiplier: 1

Instrument :
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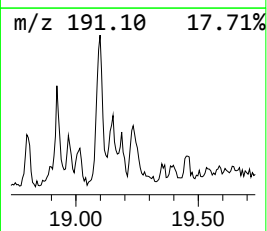
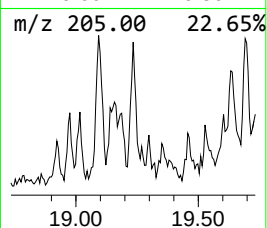
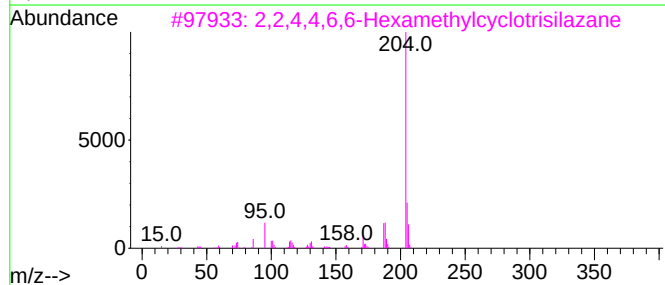
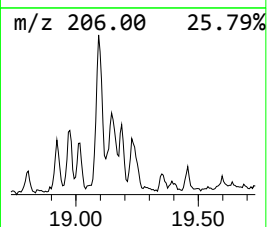
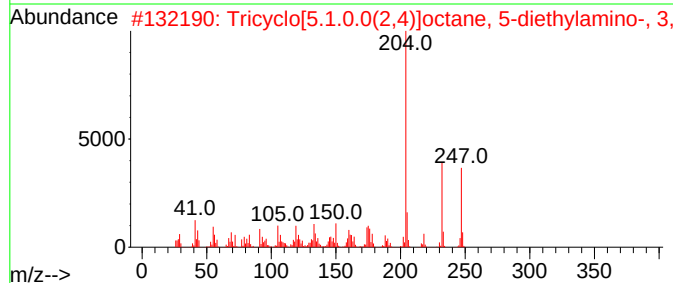
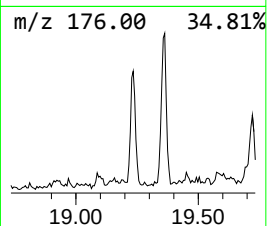
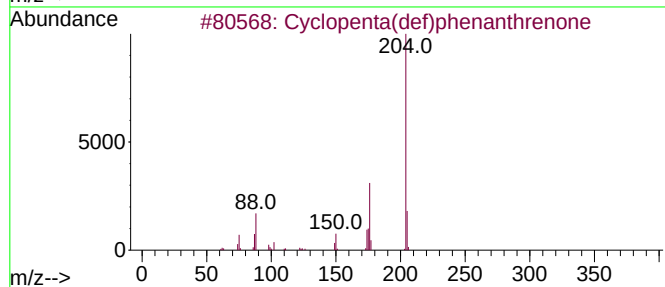
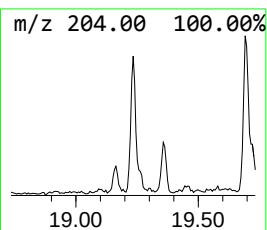
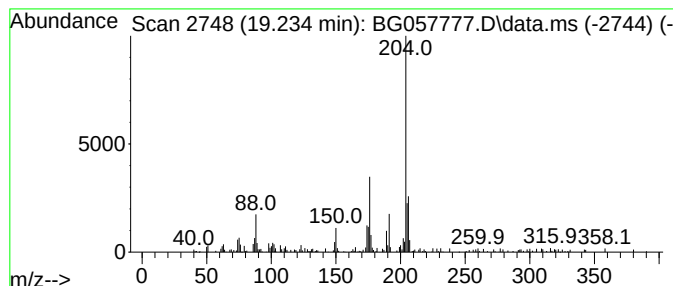
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Cyclopenta(def)phenanthrenone Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.234	3.95 ng	114647	Phenanthrene-d10	17.307

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	96
2		Tricyclo[5.1.0.0(2,4)]octane, 5-...	247	C17H29N	1000063-59-3	47
3		2,2,4,4,6,6-Hexamethylcyclotrisi...	219	C6H21N3Si3	001009-93-4	47
4		Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	46
5		2-Chloro-4-phenylphenol	204	C12H9ClO	000092-04-6	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062023\
Data File : BG057777.D
Acq On : 20 Jun 2023 15:30
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Sample : 03289-02
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG061923.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
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Benzophenone	15.914	5.4	ng	135622	3	14.563	506126	20.0
Tridecanoic acid	18.194	8.7	ng	252270	4	17.307	580392	20.0
1H-Cyclopropa[1...	18.229	4.5	ng	131255	4	17.307	580392	20.0
unknown18.370	18.370	6.4	ng	186338	4	17.307	580392	20.0
Anthracene, 2-m...	18.411	2.0	ng	59341	4	17.307	580392	20.0
Naphthalene, 2-...	18.681	2.5	ng	72823	4	17.307	580392	20.0
9,10-Anthracene...	18.717	2.4	ng	68248	4	17.307	580392	20.0
Phenanthrene, 2...	19.099	3.6	ng	103726	4	17.307	580392	20.0
Cyclopenta(def)...	19.234	4.0	ng	114647	4	17.307	580392	20.0