

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062723\
 Data File : BG057875.D
 Acq On : 27 Jun 2023 20:39
 Operator : CG/JU
 Sample : 03339-01
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MW-06

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-BG062723.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG057875.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.491	403	409	417	rBV	310309	477475	13.56%	3.078%
2	6.408	559	565	576	rBV	99766	159826	4.54%	1.030%
3	6.901	642	649	657	rBV	184820	284792	8.09%	1.836%
4	7.077	672	679	688	rBV	208223	345382	9.81%	2.227%
5	7.923	816	823	830	rBV	108306	175979	5.00%	1.134%
6	8.294	880	886	895	rBV	226350	377903	10.73%	2.436%
7	8.411	897	906	912	rBV	74025	118773	3.37%	0.766%
8	8.564	924	932	936	rBV2	46376	76668	2.18%	0.494%
9	8.611	936	940	946	rVB2	24245	37273	1.06%	0.240%
10	9.028	1003	1011	1015	rBV	129420	217107	6.17%	1.400%
11	9.081	1015	1020	1033	rVB2	420608	886351	25.18%	5.714%
12	9.551	1094	1100	1107	rVB	78430	130778	3.71%	0.843%
13	10.121	1185	1197	1205	rBV3	34886	70631	2.01%	0.455%
14	10.726	1288	1300	1304	rBV	156810	303820	8.63%	1.959%
15	10.773	1304	1308	1315	rVB	207456	364930	10.37%	2.353%
16	12.600	1610	1619	1625	rVB	70306	119637	3.40%	0.771%
17	13.182	1711	1718	1725	rVB	1126515	1708264	48.52%	11.012%
18	14.557	1945	1952	1957	rBV	278828	423548	12.03%	2.730%
19	14.622	1957	1963	1969	rVB	228060	344443	9.78%	2.220%
20	14.951	2013	2019	2026	rVB	105032	153088	4.35%	0.987%
21	15.597	2124	2129	2137	rBV	59139	94856	2.69%	0.611%
22	15.896	2175	2180	2186	rVB4	27215	49378	1.40%	0.318%
23	16.043	2196	2205	2213	rVB	1066808	1610197	45.74%	10.380%
24	16.273	2234	2244	2249	rBV3	22750	38343	1.09%	0.247%
25	17.301	2413	2419	2423	rBV	437549	651987	18.52%	4.203%
26	17.342	2423	2426	2434	rVV2	49684	76267	2.17%	0.492%
27	17.430	2434	2441	2446	rVV	46240	77047	2.19%	0.497%
28	17.483	2446	2450	2460	rVB2	26030	47013	1.34%	0.303%
29	17.700	2482	2487	2492	rVB	102450	150392	4.27%	0.970%
30	18.188	2566	2570	2572	rBV3	35535	45199	1.28%	0.291%
31	18.217	2572	2575	2584	rVV5	27612	55681	1.58%	0.359%
32	18.370	2595	2601	2606	rBV	43995	71874	2.04%	0.463%
33	19.351	2761	2768	2773	rBV	128059	184926	5.25%	1.192%
34	19.716	2820	2830	2837	rVB	95118	170448	4.84%	1.099%
35	19.915	2858	2864	2870	rBV	2644996	3520546	100.00%	22.696%
36	20.144	2898	2903	2909	rBV	36433	60164	1.71%	0.388%

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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	20.244	2916	2920	2926	rVB	47231	63368	1.80%	0.409%
38	21.202	3080	3083	3092	rVB5	30981	68618	1.95%	0.442%
39	21.284	3093	3097	3105	rVB	45177	66167	1.88%	0.427%
40	21.549	3135	3142	3152	rBV	473645	795500	22.60%	5.128%
41	24.698	3668	3678	3690	rVB	294390	837435	23.79%	5.399%

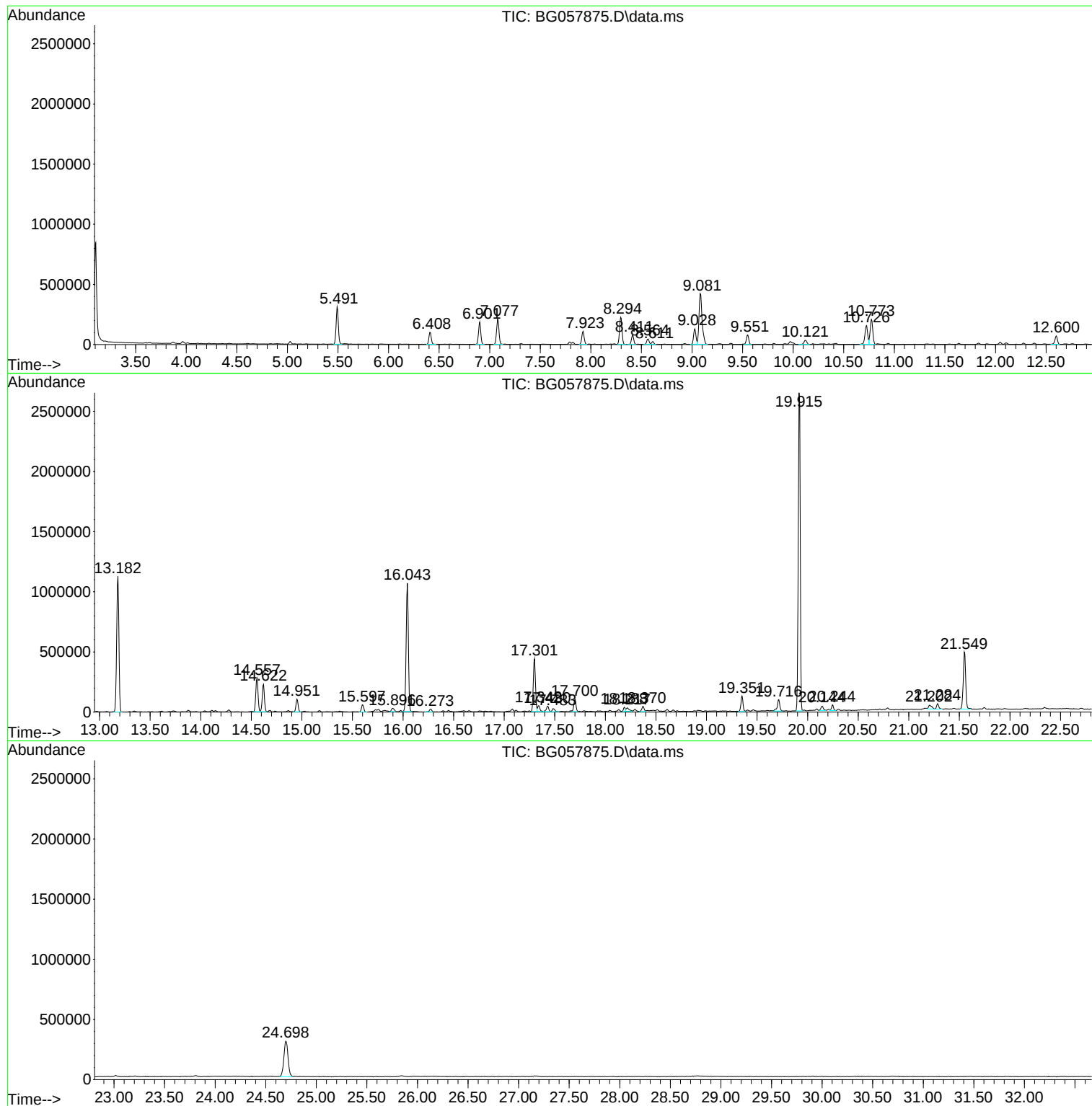
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG062723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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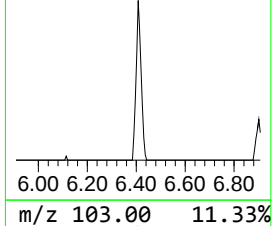
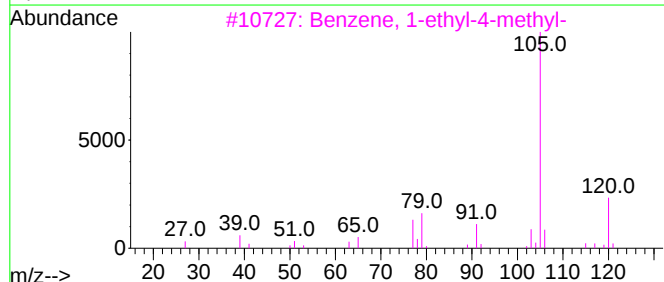
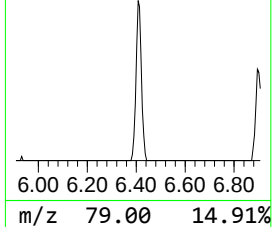
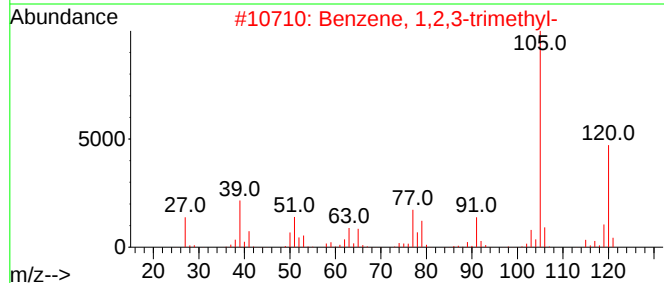
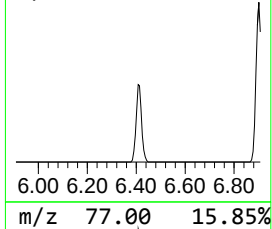
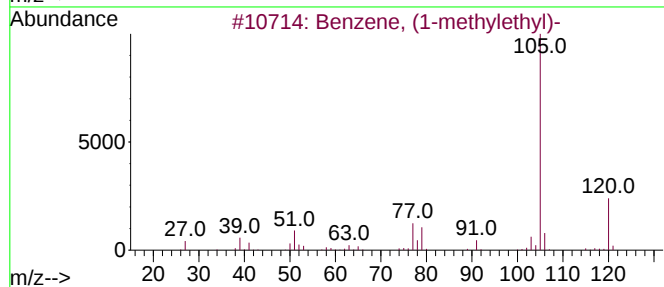
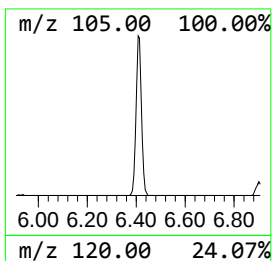
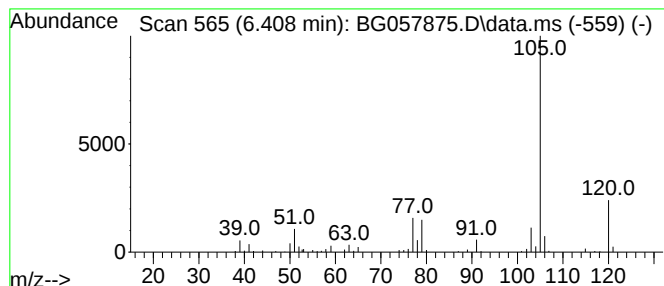
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TIC Library : C:\Database\NIST20.L
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Peak Number 1 Benzene, (1-methylethyl)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.408	18.16 ng	159826	1,4-Dichlorobenzene-d4	7.923

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	94
2			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
3			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	80
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	64
5			Mesitylene	120	C9H12	000108-67-8	64



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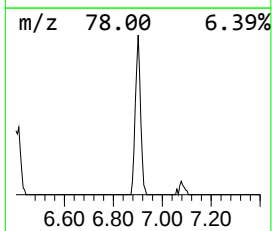
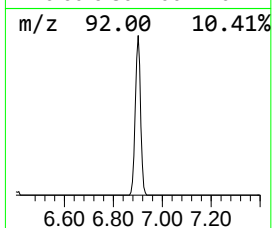
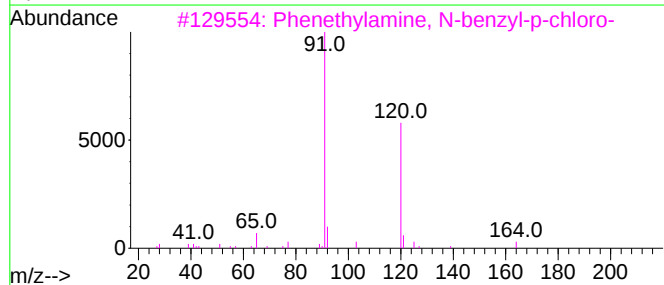
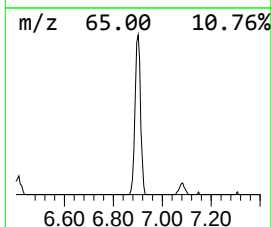
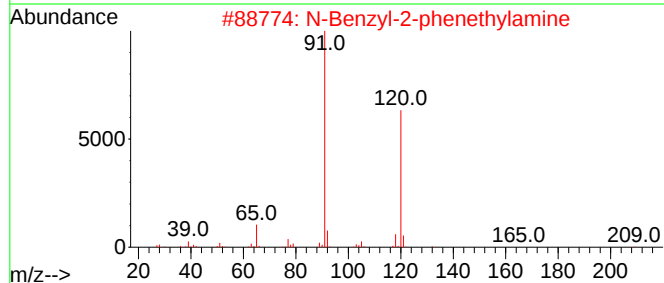
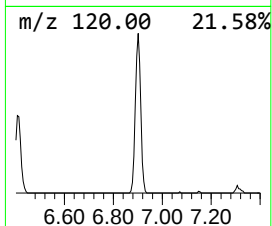
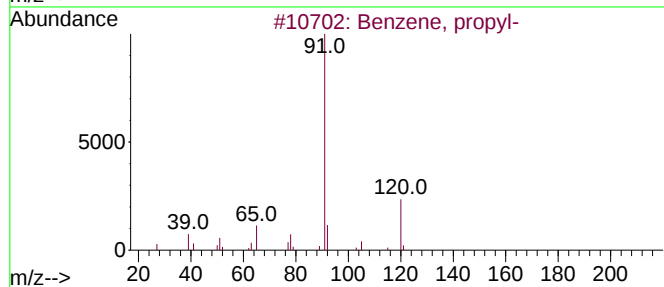
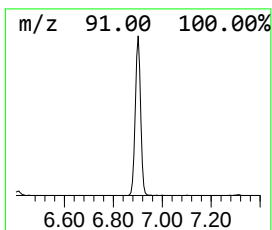
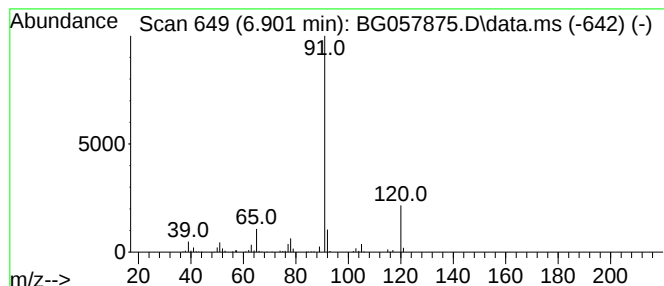
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG062723.M
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TIC Library : C:\Database\NIST20.L
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Peak Number 2 Benzene, propyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.901	32.37 ng	284792	1,4-Dichlorobenzene-d4	7.923

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, propyl-	120	C9H12	000103-65-1	94
2		N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	83
3		Phenethylamine, N-benzyl-p-chloro-	245	C15H16ClN	013622-43-0	72
4		N-Butylbenzylamine	163	C11H17N	002403-22-7	72
5		1,2-Ethanediamine, N,N'-bis(phen...	240	C16H20N2	000140-28-3	64



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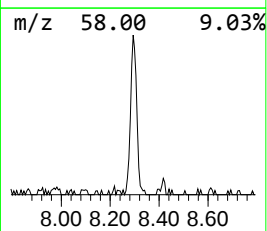
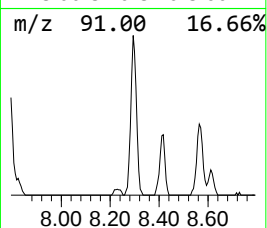
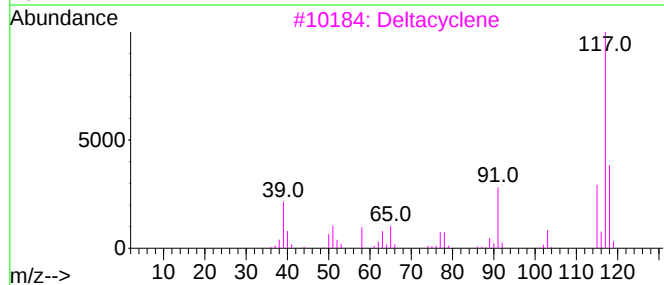
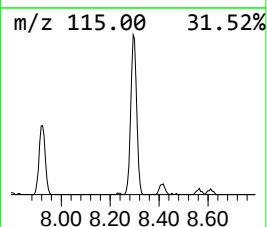
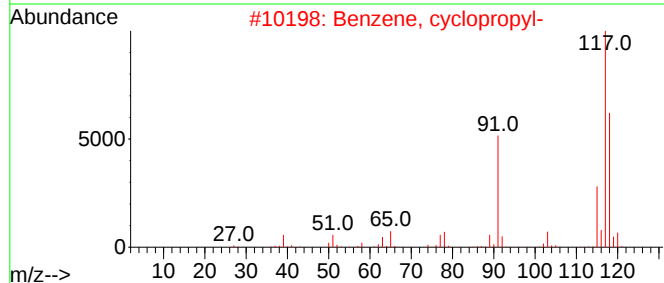
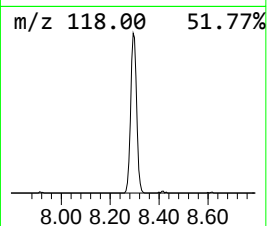
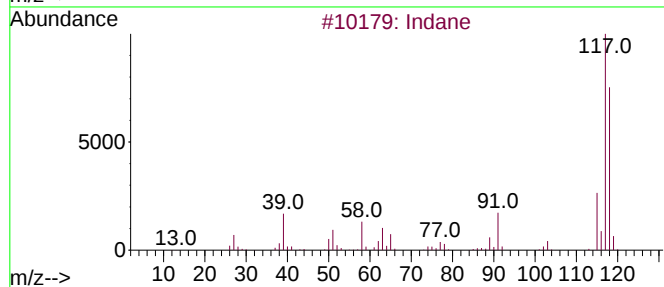
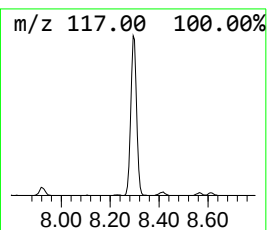
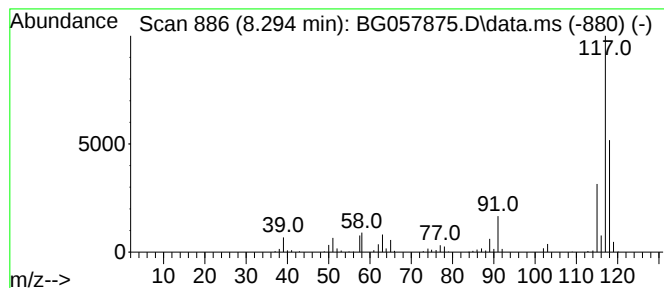
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TIC Library : C:\Database\NIST20.L
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Peak Number 3 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.294	42.95 ng	377903	1,4-Dichlorobenzene-d4	7.923

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	96
2			Benzene, cyclopropyl-	118	C9H10	000873-49-4	74
3			Deltacyclene	118	C9H10	007785-10-6	72
4			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	60
5			Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	58



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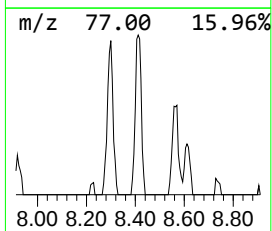
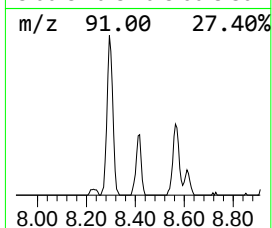
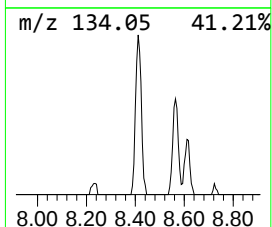
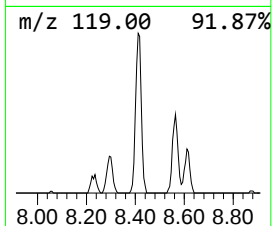
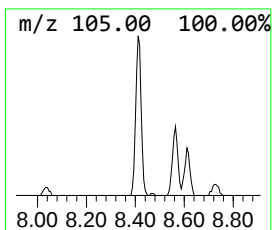
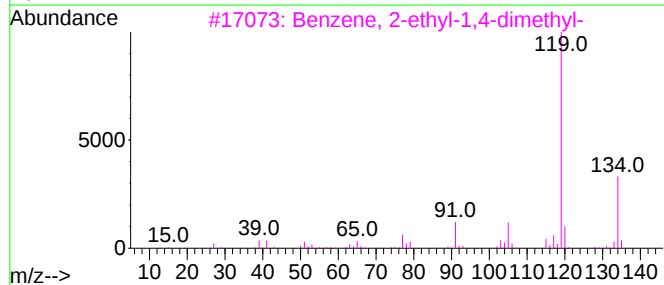
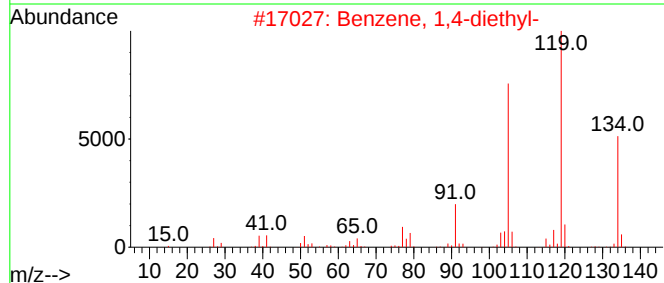
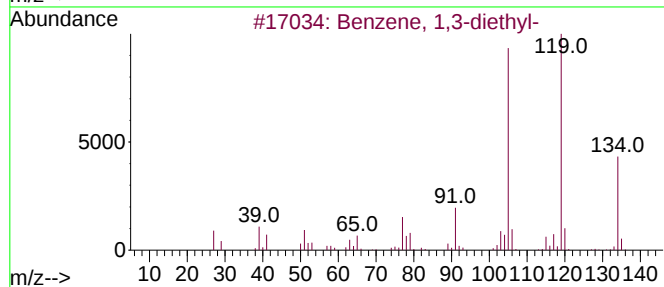
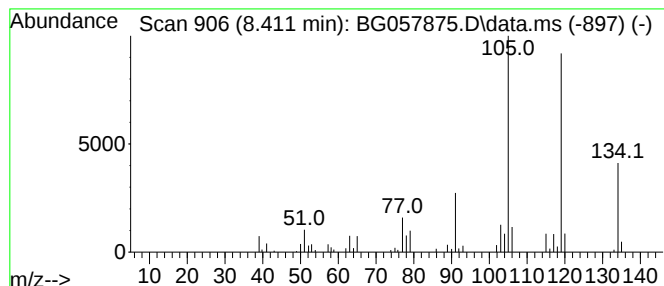
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TIC Library : C:\Database\NIST20.L
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Peak Number 4 Benzene, 1,3-diethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.411	13.50 ng	118773	1,4-Dichlorobenzene-d4	7.923

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	97
2		Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	91
3		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91
4		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	90
5		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	81



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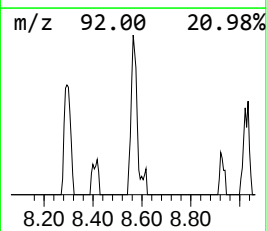
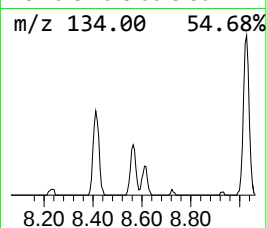
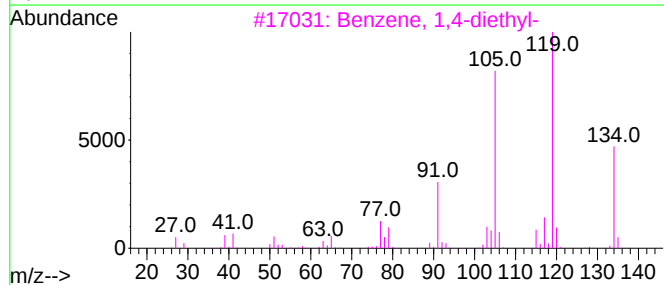
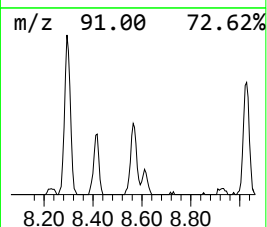
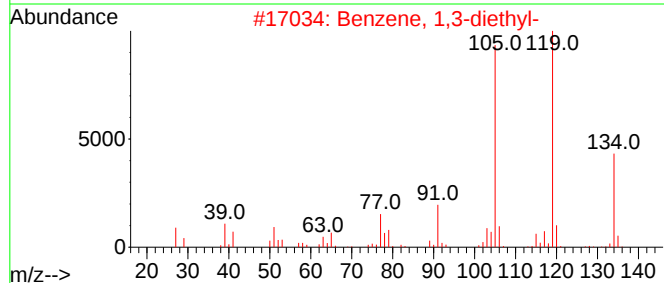
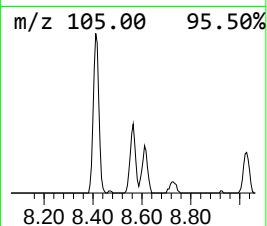
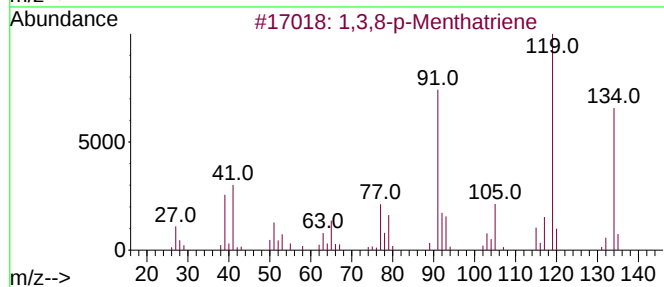
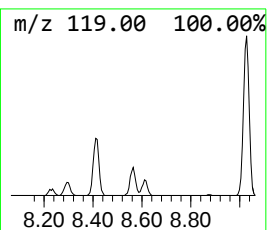
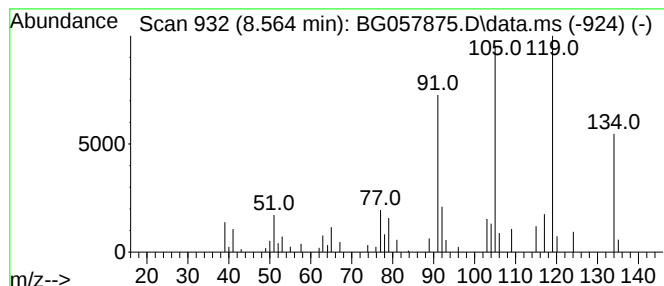
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Peak Number 5 1,3,8-p-Menthatriene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.564	8.71 ng	76668	1,4-Dichlorobenzene-d4	7.923

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3,8-p-Menthatriene	134	C10H14	018368-95-1	90
2			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	81
3			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	81
4			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	81
5			2,6-Dimethyl-1,3,5,7-octatetraen...	134	C10H14	000460-01-5	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062723\
Data File : BG057875.D
Acq On : 27 Jun 2023 20:39
Operator : CG/JU
Sample : 03339-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
MW-06

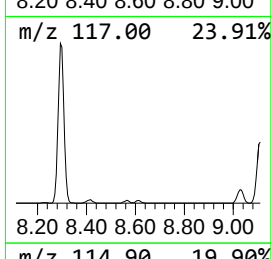
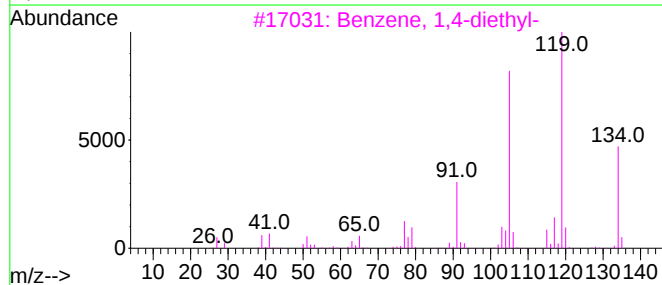
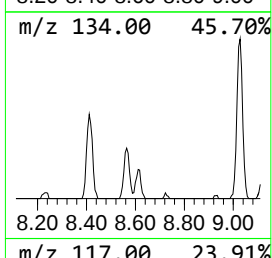
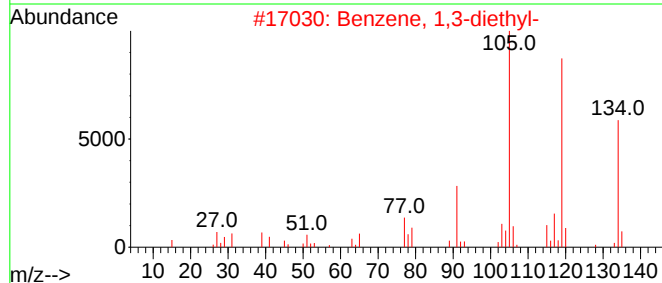
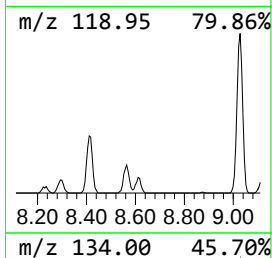
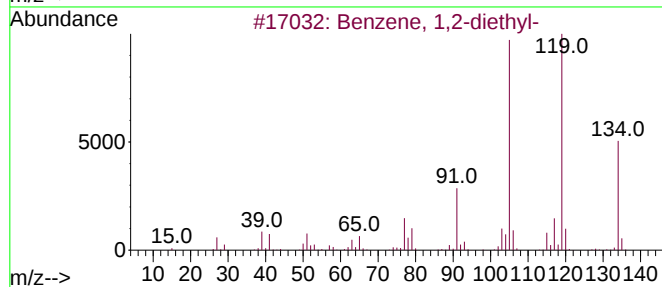
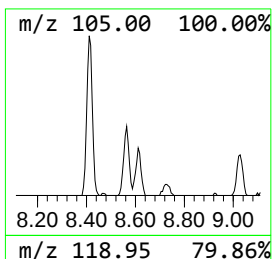
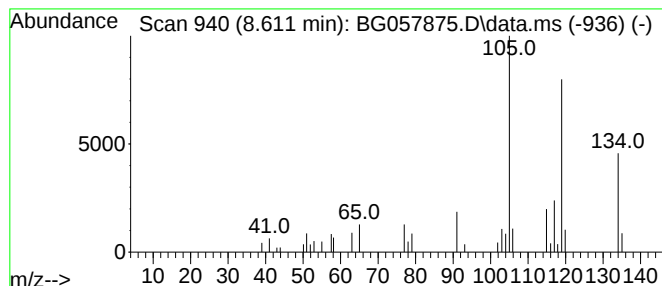
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Benzene, 1,2-diethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.611	4.24 ng	37273	1,4-Dichlorobenzene-d4	7.923

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	93
2			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	87
3			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	87
4			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	81
5			2,6-Dimethyl-1,3,5,7-octatetraen...	134	C10H14	000460-01-5	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062723\
Data File : BG057875.D
Acq On : 27 Jun 2023 20:39
Operator : CG/JU
Sample : 03339-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

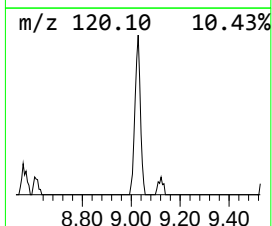
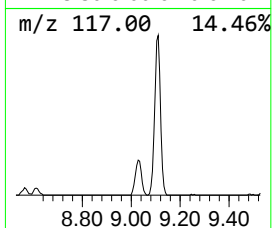
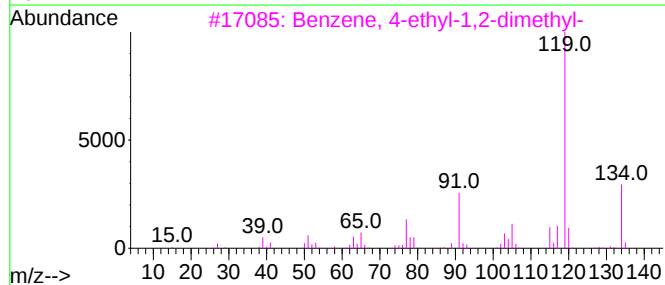
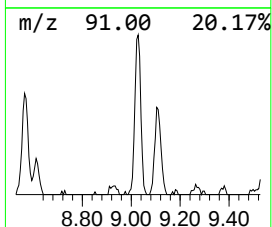
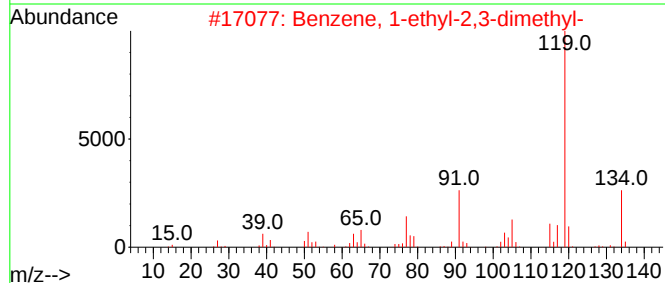
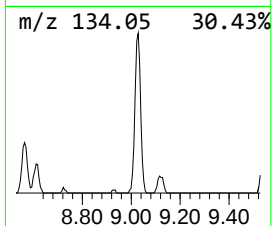
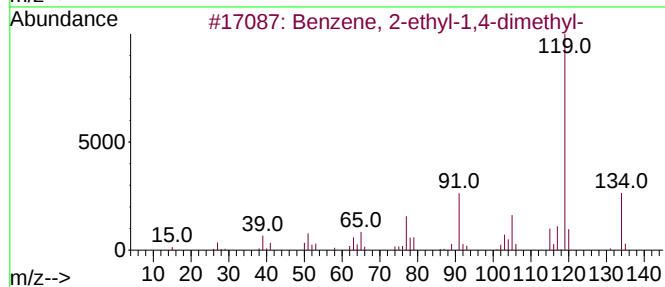
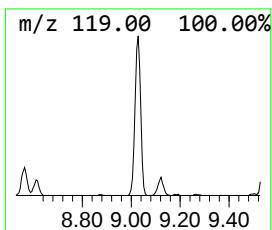
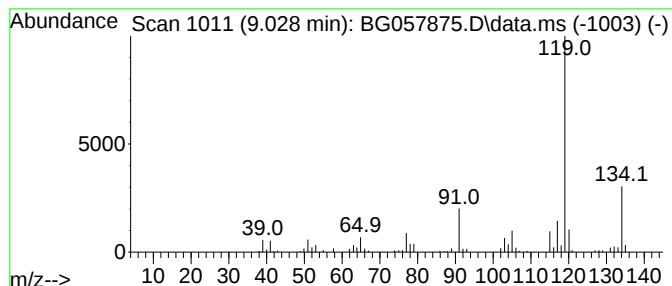
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BNA_G
ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
9.028	24.67 ng	217107	1,4-Dichlorobenzene-d4	7.923	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
3	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
4	o-Cymene	134	C10H14	000527-84-4	94
5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062723\
Data File : BG057875.D
Acq On : 27 Jun 2023 20:39
Operator : CG/JU
Sample : 03339-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
MW-06

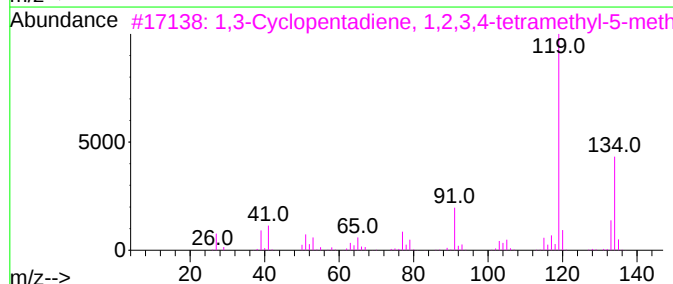
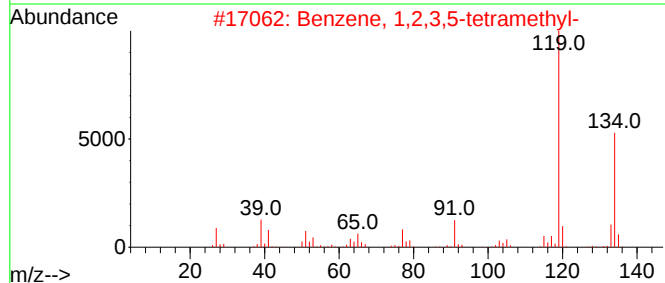
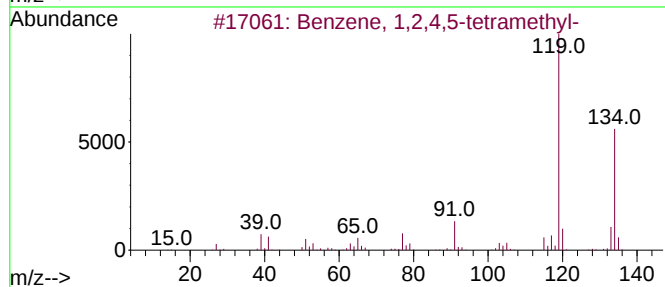
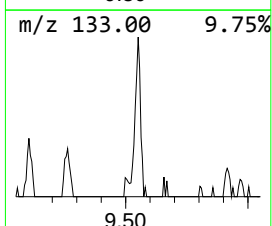
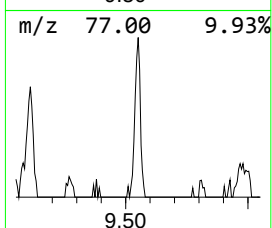
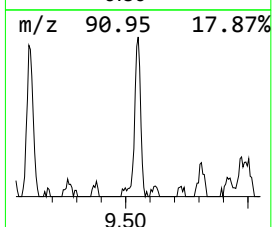
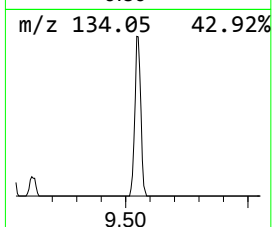
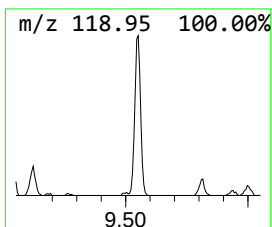
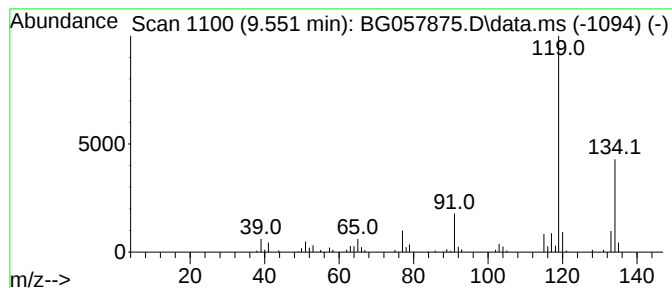
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.551	8.61 ng	130778	Naphthalene-d8	10.726

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
2			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
3			1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	94
4			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94
5			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062723\
Data File : BG057875.D
Acq On : 27 Jun 2023 20:39
Operator : CG/JU
Sample : 03339-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

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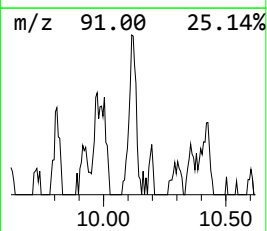
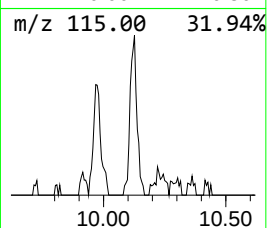
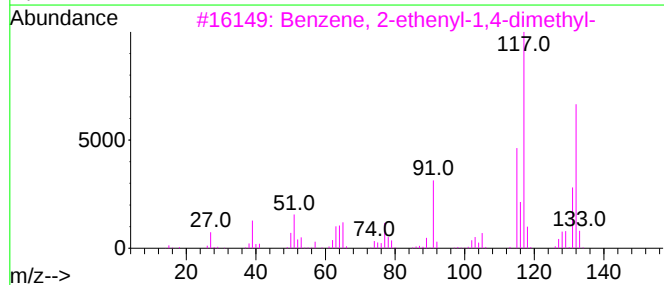
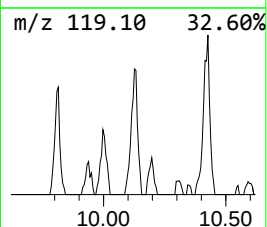
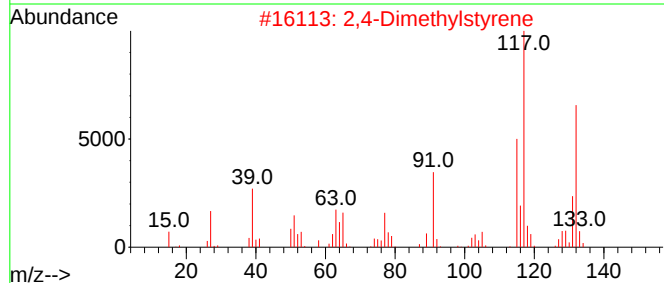
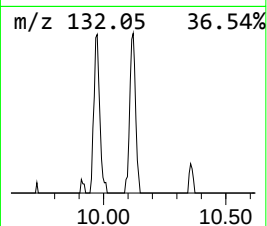
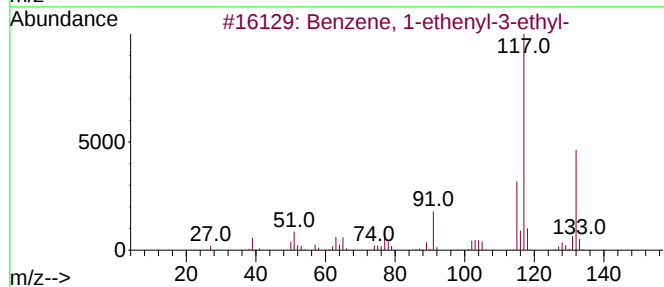
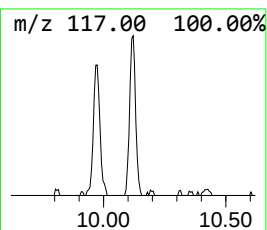
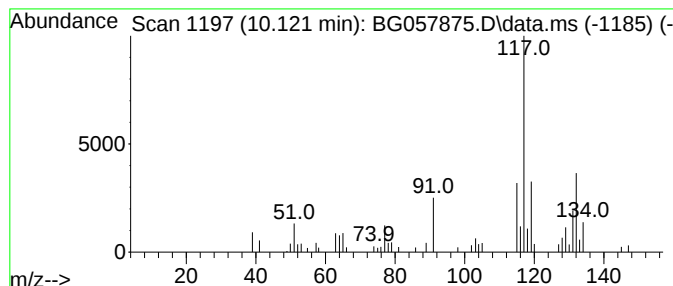
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.121	4.65 ng	70631	Naphthalene-d8	10.726

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	93
2			2,4-Dimethylstyrene	132	C10H12	002234-20-0	91
3			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	91
4			Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	89
5			3-Phenylbut-1-ene	132	C10H12	000934-10-1	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062723\
Data File : BG057875.D
Acq On : 27 Jun 2023 20:39
Operator : CG/JU
Sample : 03339-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

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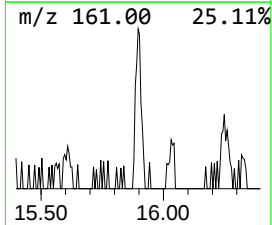
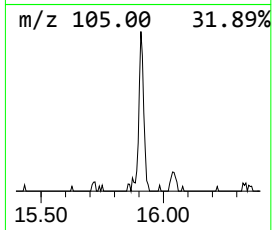
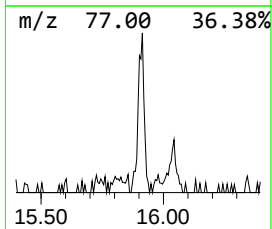
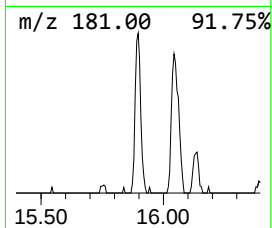
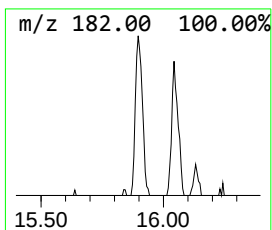
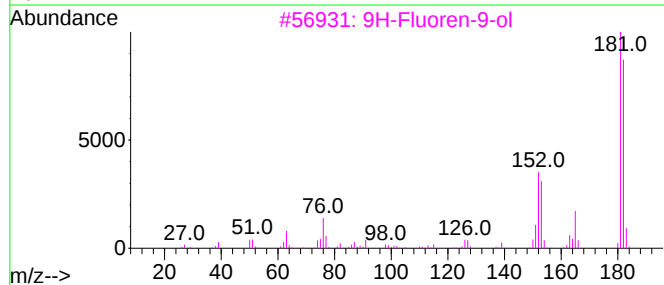
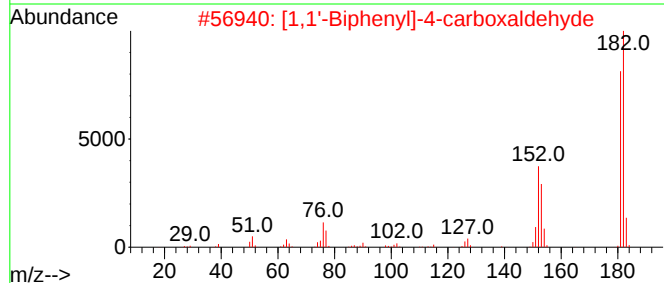
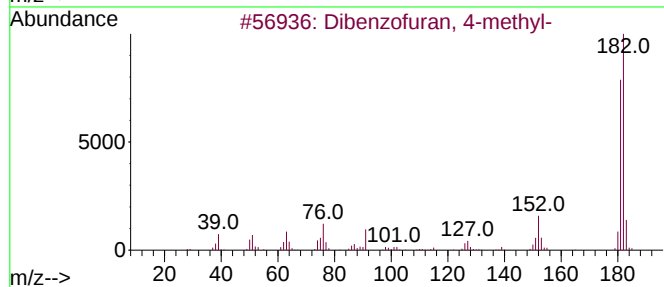
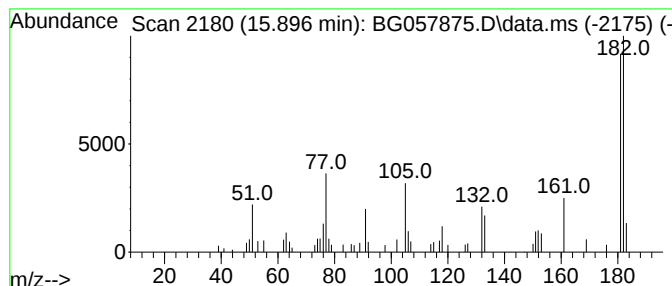
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Dibenzofuran, 4-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.897	2.33 ng	49378	Acenaphthene-d10	14.557

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dibenzofuran, 4-methyl-	182	C13H10O	007320-53-8	62
2			[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	58
3			9H-Fluoren-9-ol	182	C13H10O	001689-64-1	58
4			2,4,6-Cycloheptatrien-1-one, 2-p...	182	C13H10O	014562-09-5	53
5			Pyridine, 4,4'-(1,2-ethenediyl)bis-	182	C12H10N2	001135-32-6	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062723\
Data File : BG057875.D
Acq On : 27 Jun 2023 20:39
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Misc :
ALS Vial : 12 Sample Multiplier: 1

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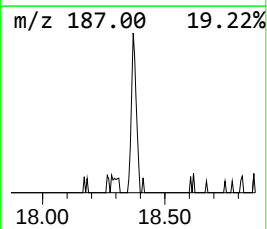
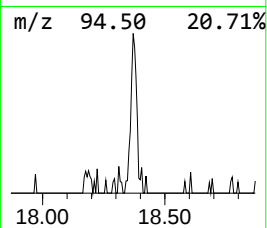
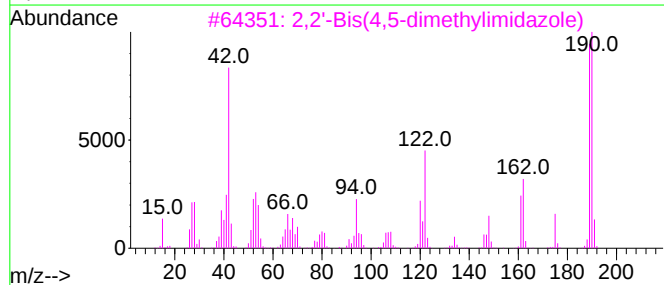
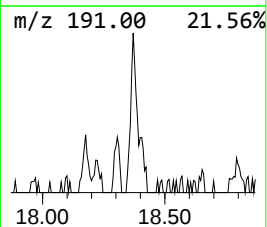
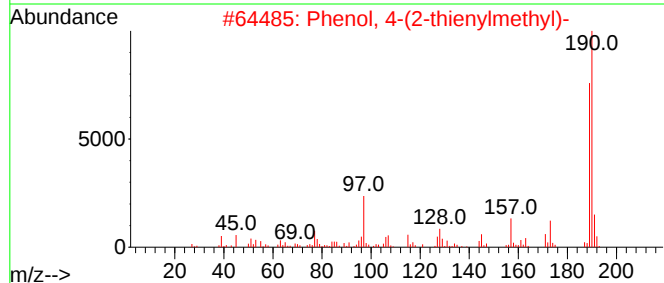
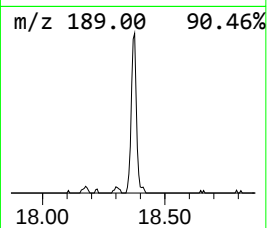
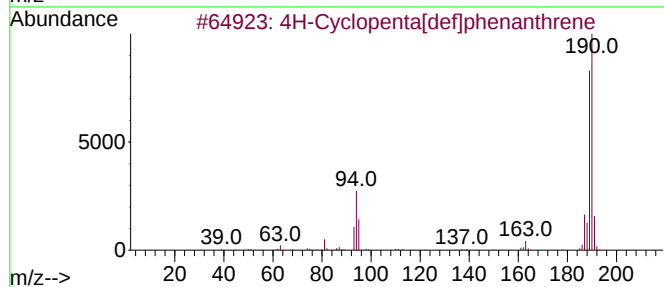
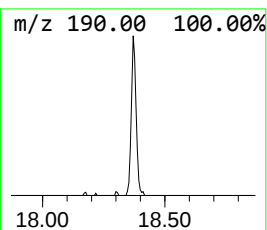
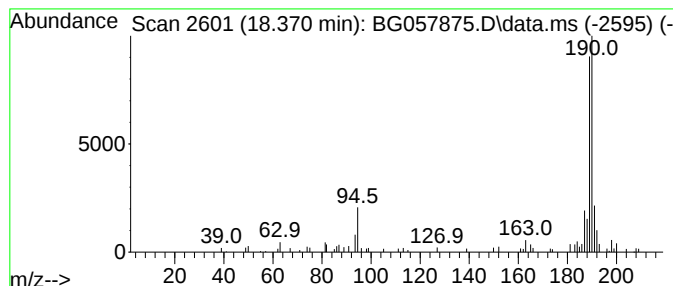
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 4H-Cyclopenta[def]phenanthrene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.370	2.20 ng	71874	Phenanthrene-d10	17.301

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	90
2			Phenol, 4-(2-thienylmethyl)-	190	C11H10OS	091680-55-6	59
3			2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	45
4			1,4-Naphthalenedione, 2,5-dihydr...	190	C10H6O4	004923-55-1	25
5			2,4(1H,3H)-Pyrimidinedione, 5-br...	190	C4H3BrN2O2	000051-20-7	16



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062723\
Data File : BG057875.D
Acq On : 27 Jun 2023 20:39
Operator : CG/JU
Sample : 03339-01
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-BG062723.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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Benzene, propyl-	6.901	32.4	ng	284792	1	7.923	175979	20.0
Indane	8.294	43.0	ng	377903	1	7.923	175979	20.0
Benzene, 1,3-di...	8.411	13.5	ng	118773	1	7.923	175979	20.0
1,3,8-p-Menthat...	8.564	8.7	ng	76668	1	7.923	175979	20.0
Benzene, 1,2-di...	8.611	4.2	ng	37273	1	7.923	175979	20.0
Benzene, 2-ethy...	9.028	24.7	ng	217107	1	7.923	175979	20.0
Benzene, 1,2,4,...	9.551	8.6	ng	130778	2	10.726	303820	20.0
Benzene, 1-ethe...	10.121	4.7	ng	70631	2	10.726	303820	20.0
Dibenzofuran, 4...	15.897	2.3	ng	49378	3	14.557	423548	20.0
4H-Cyclopenta[d...	18.370	2.2	ng	71874	4	17.301	651987	20.0