

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082125\
 Data File : BF143550.D
 Acq On : 22 Aug 2025 01:51
 Operator : RC/JU
 Sample : Q2903-07
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-12B-081925

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_F\Methods\8270-BF082025.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF143550.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.275	5	14	21	rBV	2116716	3543643	49.01%	10.808%
2	5.157	495	504	512	rBV2	101292	174391	2.41%	0.532%
3	5.563	564	573	596	rBV	1190681	1605602	22.20%	4.897%
4	6.104	660	665	671	rBV	200561	247717	3.43%	0.756%
5	6.557	734	742	748	rBV	822001	1101733	15.24%	3.360%
6	6.616	748	752	764	rVB	187818	247452	3.42%	0.755%
7	6.928	800	805	810	rBV	478780	623756	8.63%	1.902%
8	6.987	810	815	824	rVB	110182	158848	2.20%	0.484%
9	7.116	830	837	842	rBV	4926633	7231141	100.00%	22.055%
10	7.487	889	900	910	rBV2	1479780	2177270	30.11%	6.641%
11	7.875	960	966	972	rVB	76597	105280	1.46%	0.321%
12	7.945	972	978	982	rBV2	271674	377233	5.22%	1.151%
13	7.993	982	986	992	rVV	204125	349117	4.83%	1.065%
14	8.228	1015	1026	1031	rBV2	2223797	3814031	52.74%	11.633%
15	8.269	1031	1033	1045	rVB2	61935	90272	1.25%	0.275%
16	8.463	1062	1066	1079	rVB3	46725	106821	1.48%	0.326%
17	8.792	1115	1122	1127	rBV2	100372	143148	1.98%	0.437%
18	9.016	1155	1160	1165	rBV	264287	341077	4.72%	1.040%
19	9.098	1170	1174	1181	rVB	70171	108498	1.50%	0.331%
20	9.281	1199	1205	1210	rBV	2382226	3005482	41.56%	9.167%
21	9.963	1315	1321	1325	rBV	671702	863805	11.95%	2.635%
22	10.228	1361	1366	1373	rBV	89168	139114	1.92%	0.424%
23	10.410	1392	1397	1403	rVB	97593	140135	1.94%	0.427%
24	10.669	1437	1441	1448	rVB	93550	127847	1.77%	0.390%
25	10.751	1448	1455	1466	rBV2	1384436	1898202	26.25%	5.790%
26	11.451	1569	1574	1581	rVB	518162	664507	9.19%	2.027%
27	11.975	1658	1663	1675	rBV2	50553	107339	1.48%	0.327%
28	13.033	1837	1843	1848	rBV	1611376	2054876	28.42%	6.267%
29	14.092	2019	2023	2032	rVB	452273	608781	8.42%	1.857%
30	15.592	2272	2278	2290	rBV	402132	629265	8.70%	1.919%

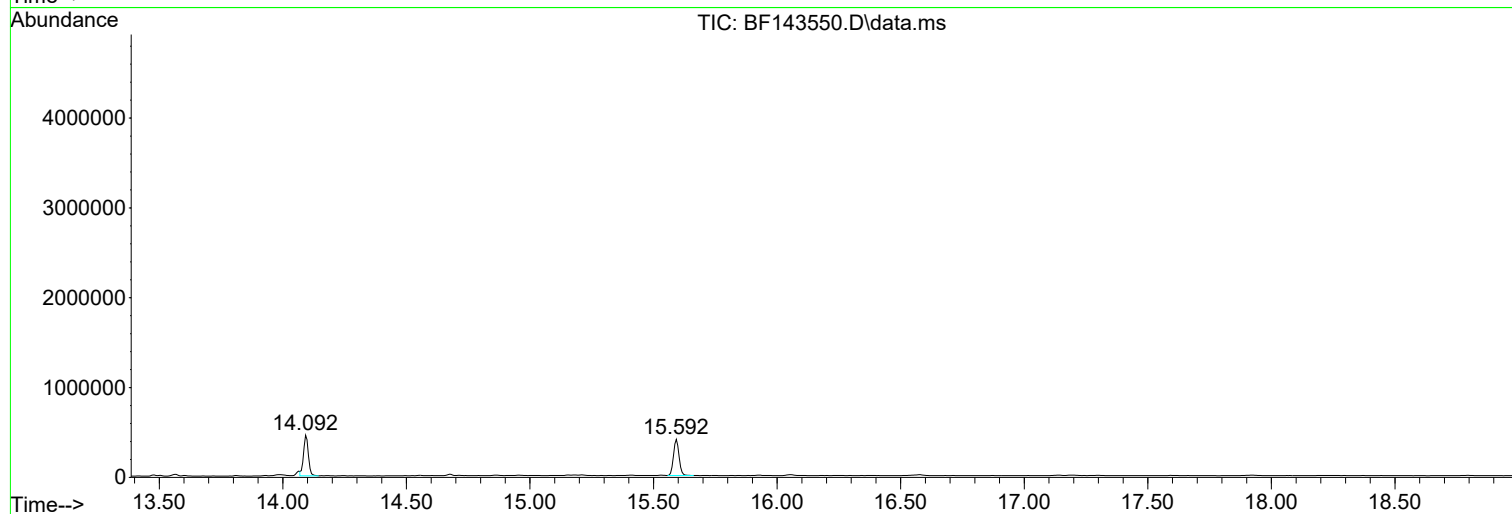
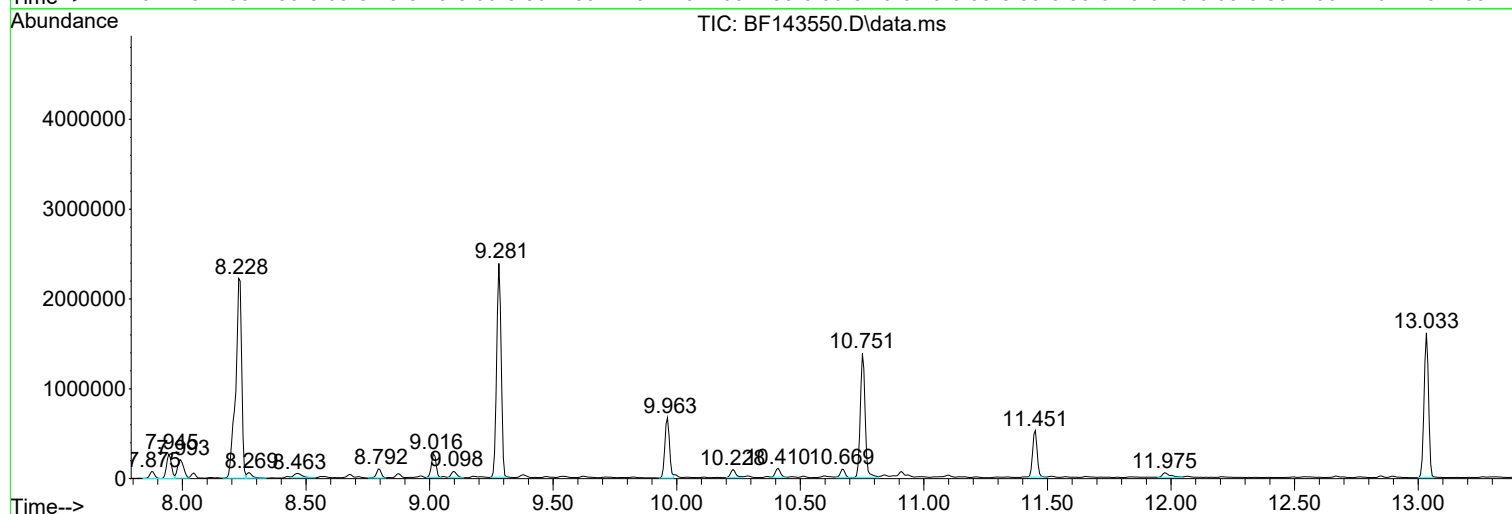
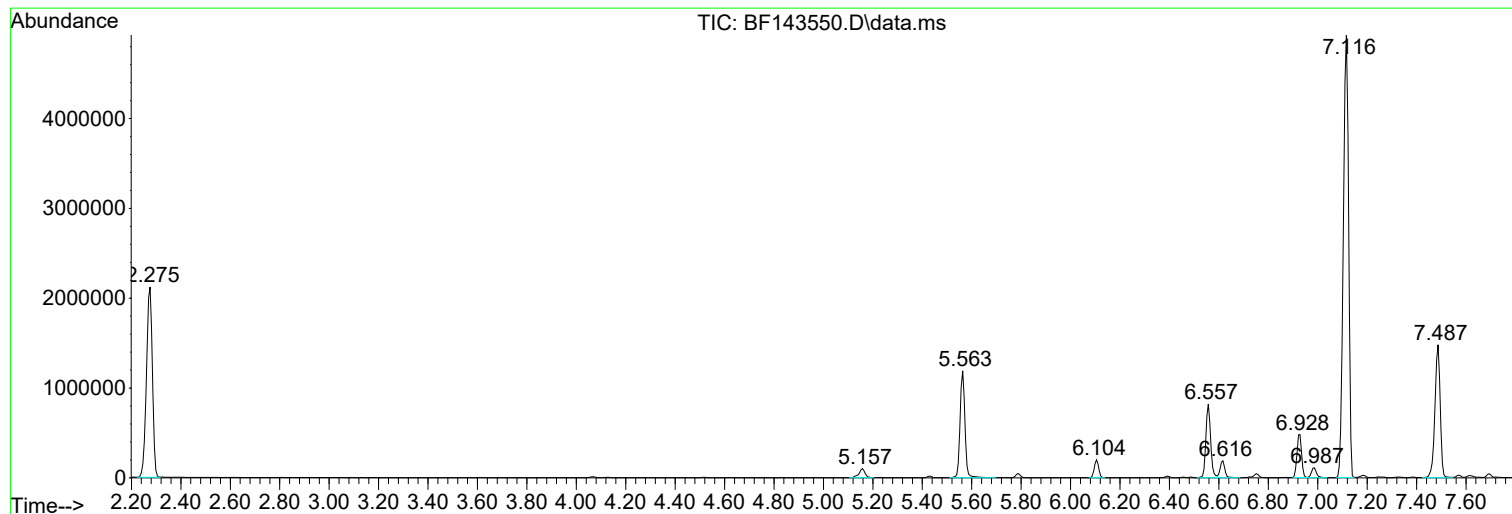
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF082025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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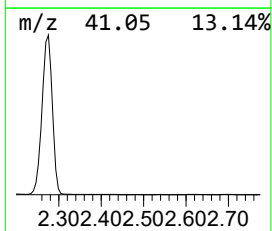
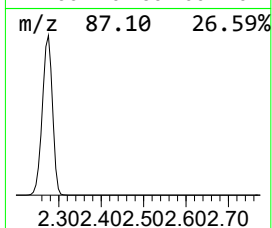
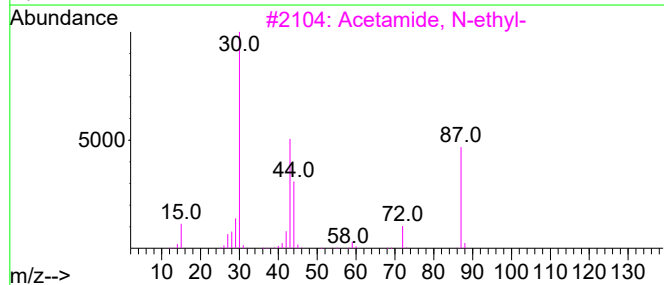
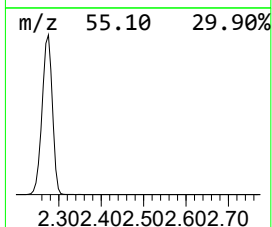
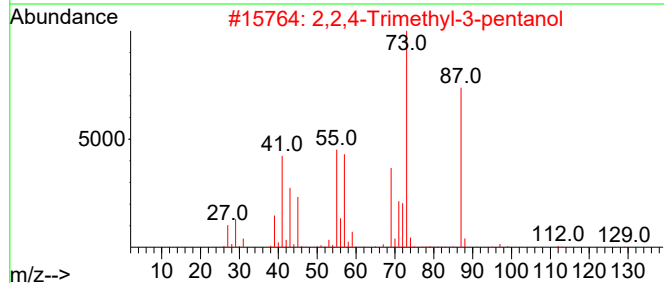
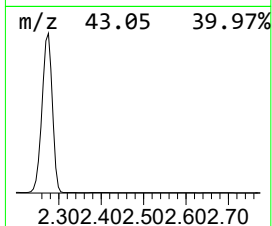
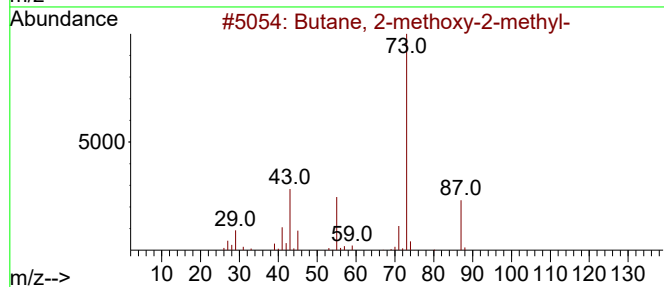
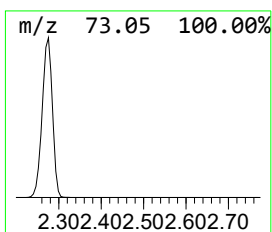
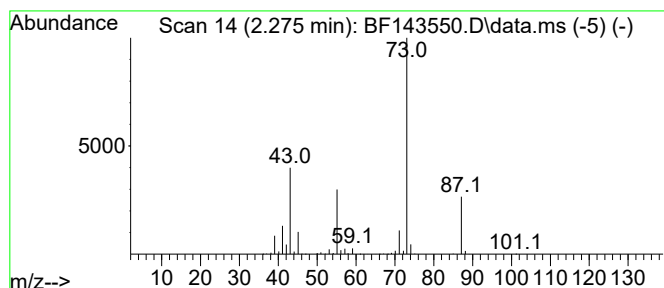
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.275	113.62 ng	3543640	1,4-Dichlorobenzene-d4	6.928

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	27
4			Octanal, 7-methoxy-3,7-dimethyl-	186	C11H22O2	003613-30-7	17
5			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12



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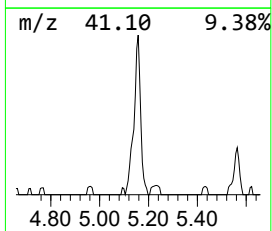
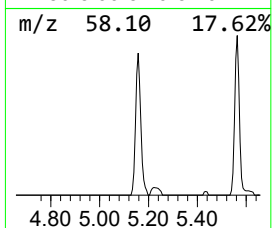
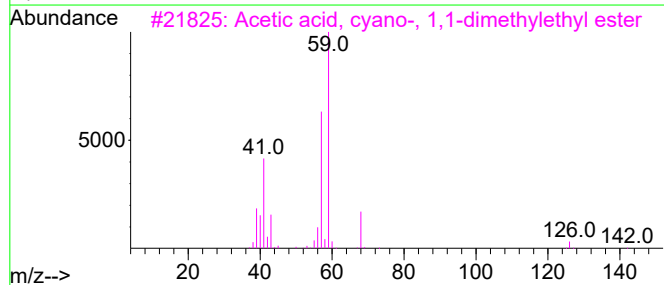
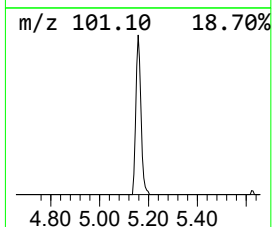
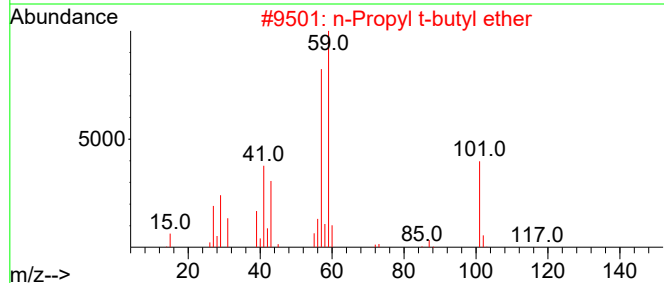
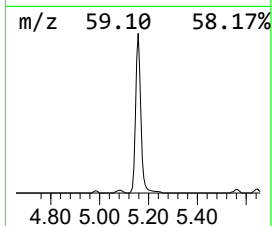
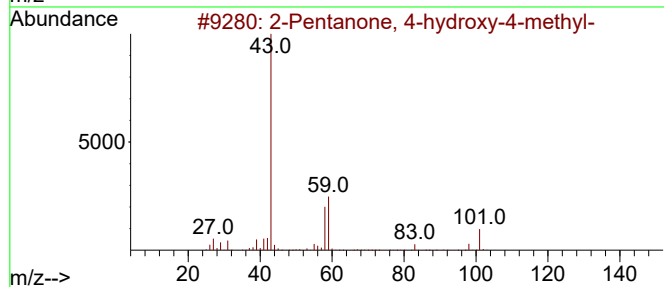
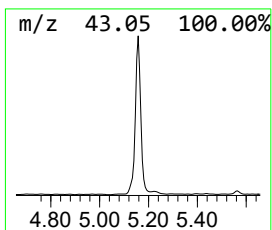
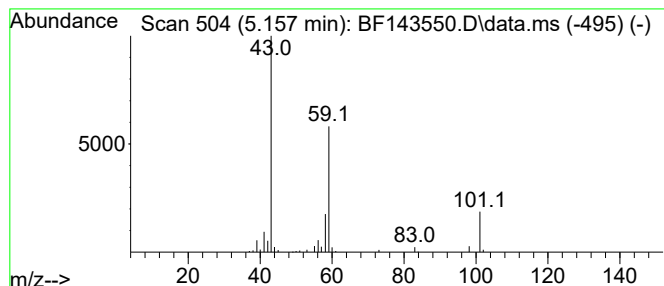
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.157	5.59 ng	174391	1,4-Dichlorobenzene-d4	6.928

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		n-Propyl t-butyl ether	116	C7H16O	1000334-81-9	33
3		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
4		1,2-Butanediol	90	C4H10O2	000584-03-2	25
5		Acetone	58	C3H6O	000067-64-1	12



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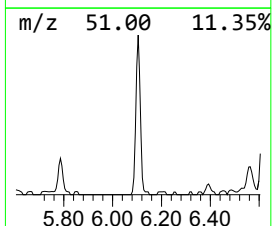
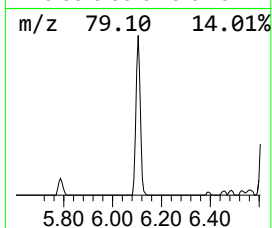
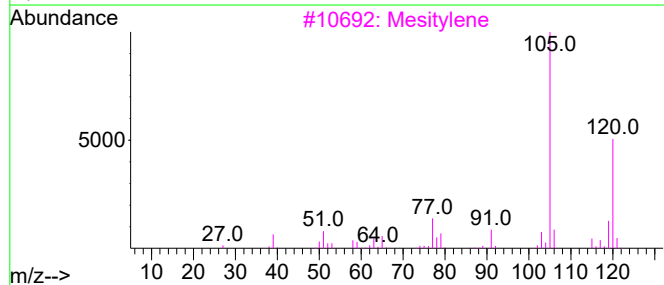
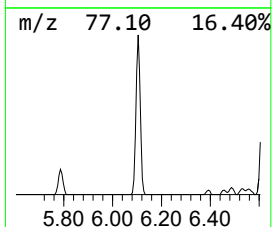
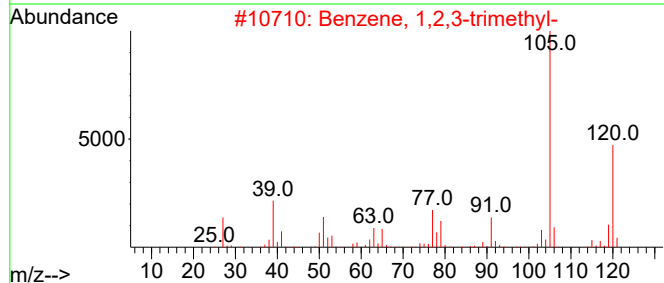
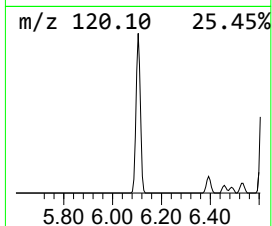
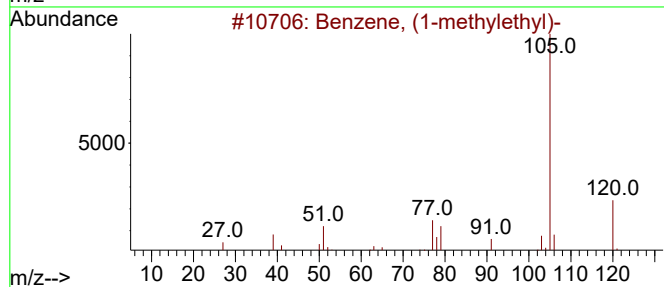
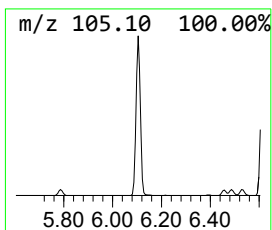
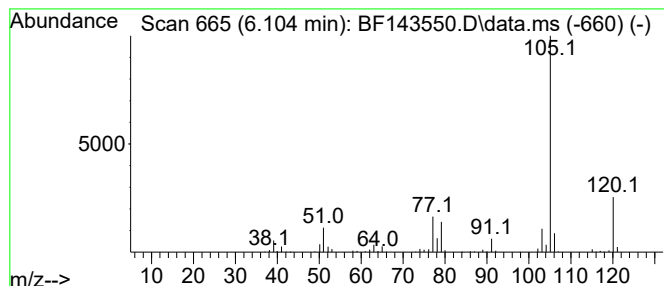
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TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, (1-methylethyl)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.104	7.94 ng	247717	1,4-Dichlorobenzene-d4	6.928

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	95
2			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
3			Mesitylene	120	C9H12	000108-67-8	87
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	83
5			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	83



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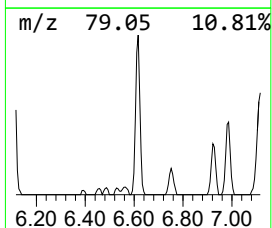
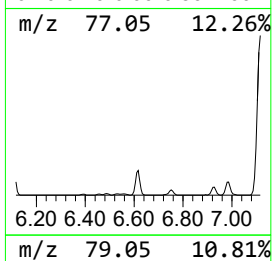
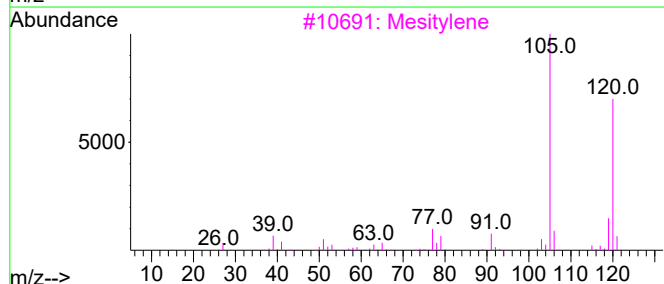
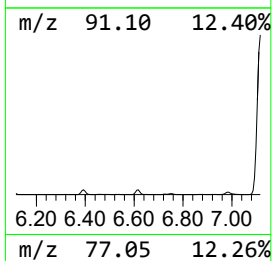
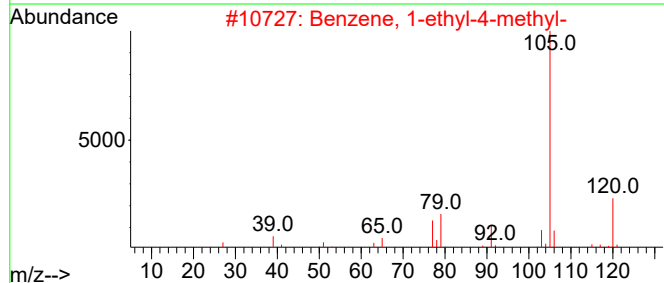
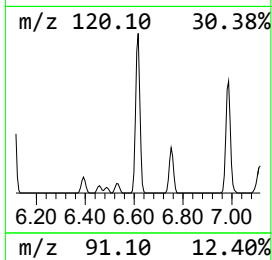
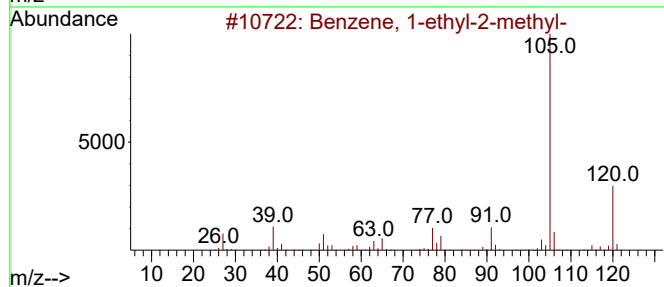
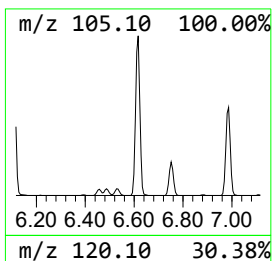
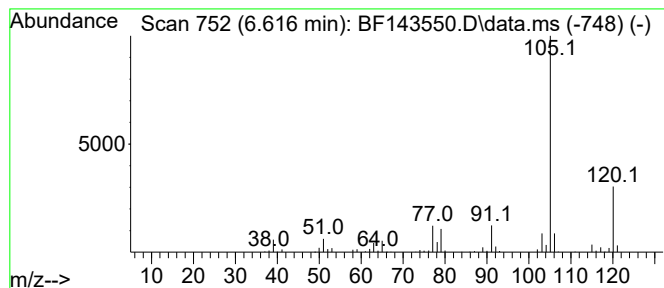
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Peak Number 4 Benzene, 1-ethyl-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.616	7.93 ng	247452	1,4-Dichlorobenzene-d4	6.928

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	94
3			Mesitylene	120	C9H12	000108-67-8	91
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



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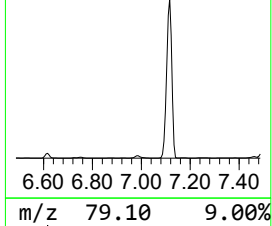
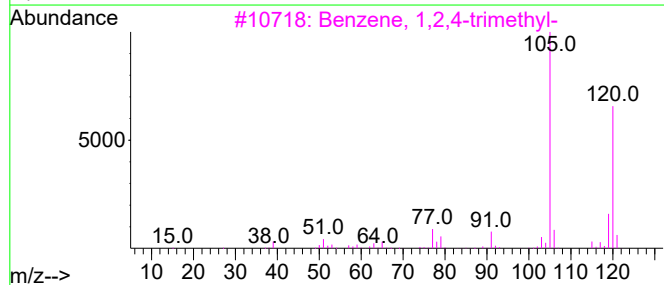
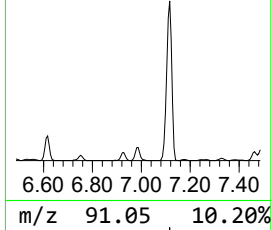
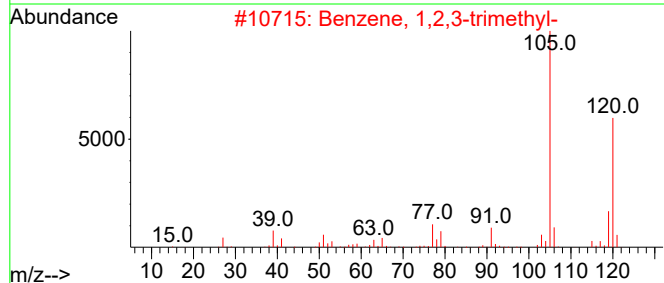
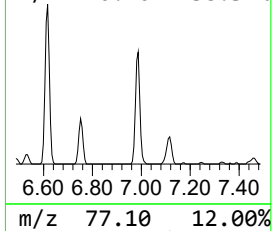
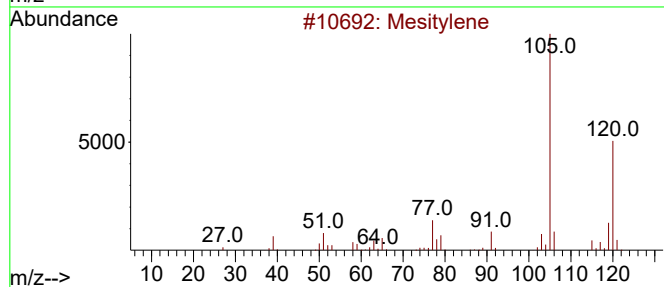
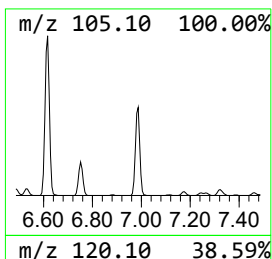
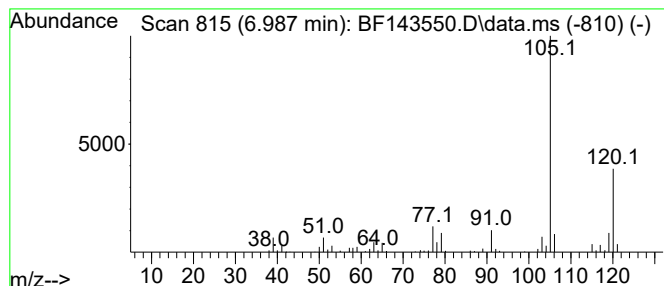
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, 1,2,3-trimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.987	5.09 ng	158848	1,4-Dichlorobenzene-d4	6.928

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Mesitylene	120	C9H12	000108-67-8	97
2			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93
5			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91



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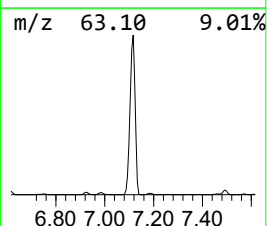
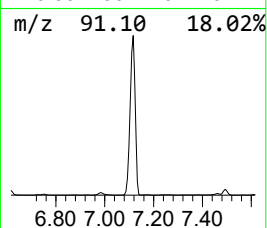
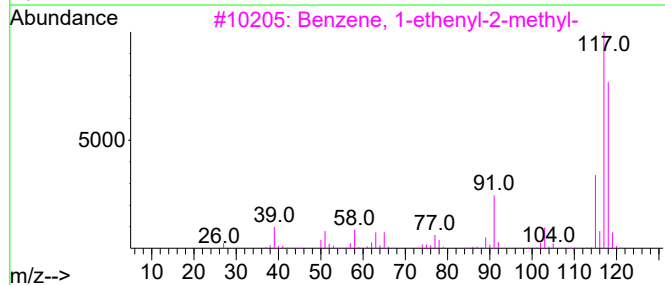
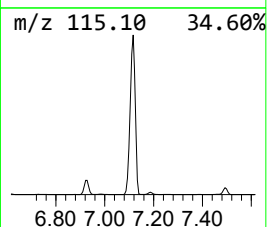
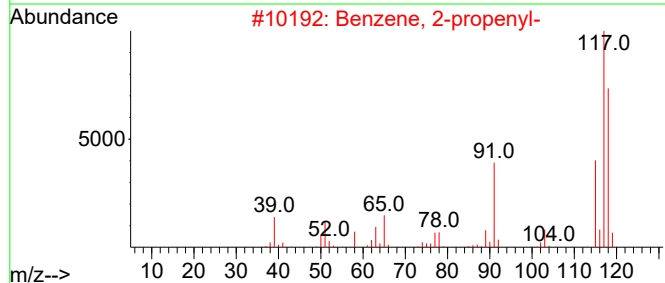
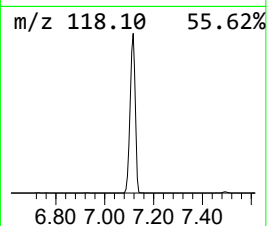
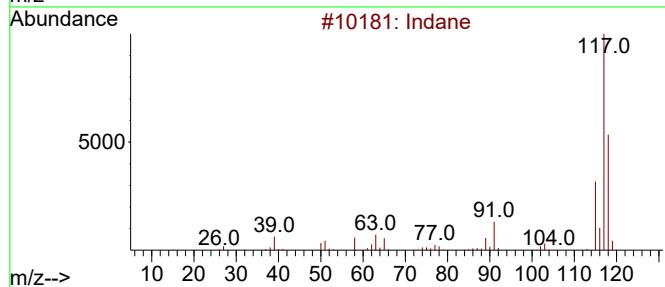
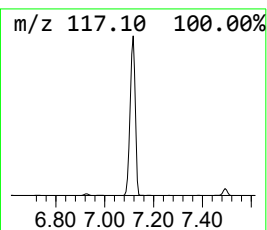
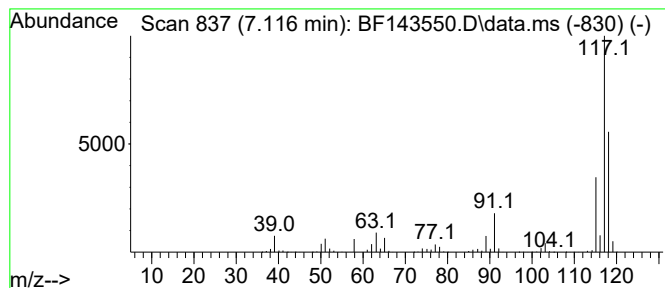
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.116	231.86 ng	7231140	1,4-Dichlorobenzene-d4	6.928

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	93
2			Benzene, 2-propenyl-	118	C9H10	000300-57-2	81
3			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	64
4			Benzene, cyclopropyl-	118	C9H10	000873-49-4	64
5			Benzeneethanol, .beta.-ethenyl-....	162	C11H14O	040596-14-3	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082125\
Data File : BF143550.D
Acq On : 22 Aug 2025 01:51
Operator : RC/JU
Sample : Q2903-07
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-12B-081925

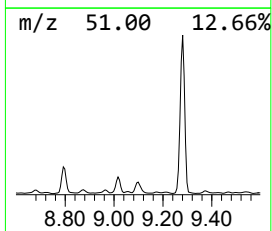
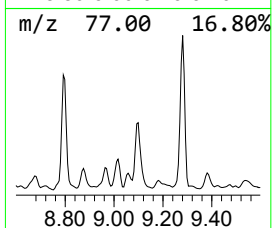
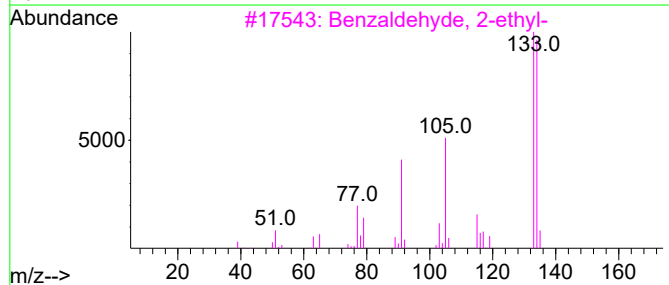
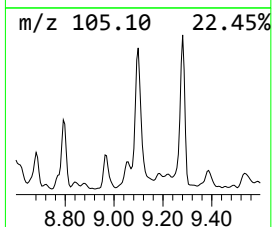
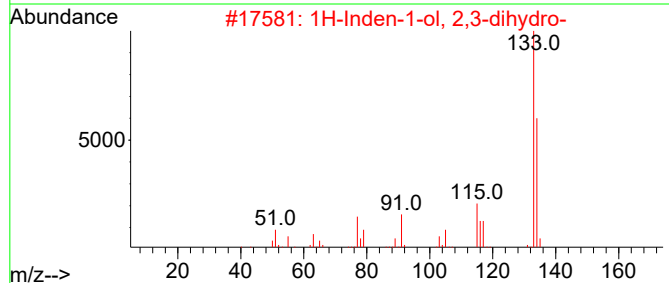
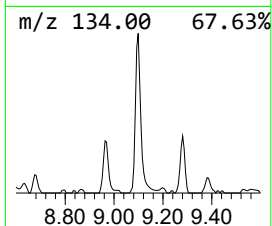
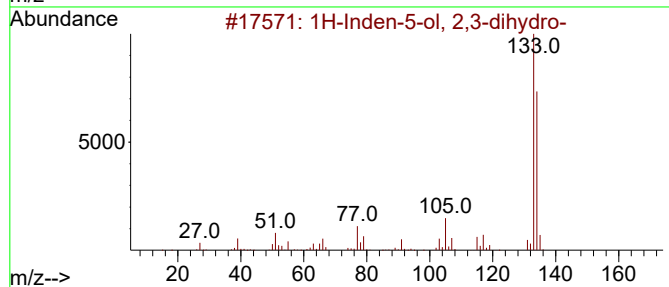
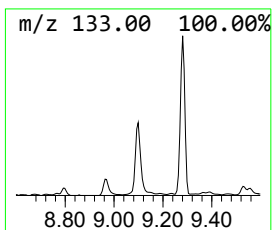
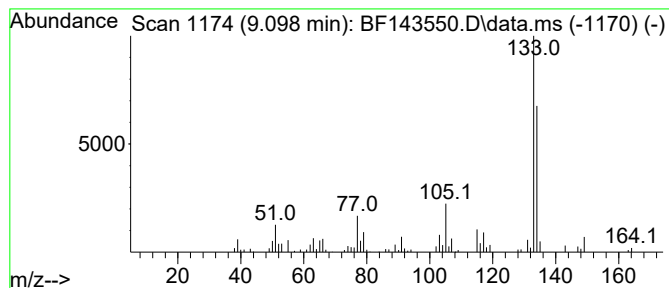
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 1H-Inden-5-ol, 2,3-dihydro- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.098	2.51 ng	108498	Acenaphthene-d10	9.963

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Inden-5-ol, 2,3-dihydro-	134	C9H10O	001470-94-6	94
2		1H-Inden-1-ol, 2,3-dihydro-	134	C9H10O	006351-10-6	72
3		Benzaldehyde, 2-ethyl-	134	C9H10O	022927-13-5	64
4		Benzaldehyde, 3,4-dimethyl-	134	C9H10O	005973-71-7	64
5		Phenol, 4-(2-propenyl)-	134	C9H10O	000501-92-8	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082125\
Data File : BF143550.D
Acq On : 22 Aug 2025 01:51
Operator : RC/JU
Sample : Q2903-07
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
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ClientSampleId :
MW-12B-081925

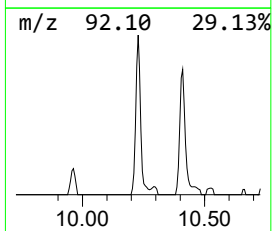
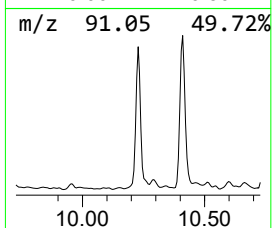
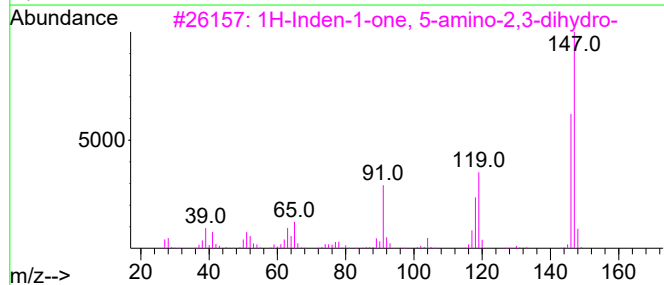
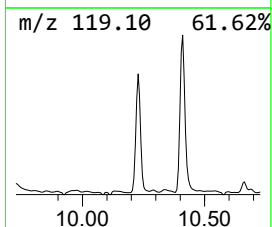
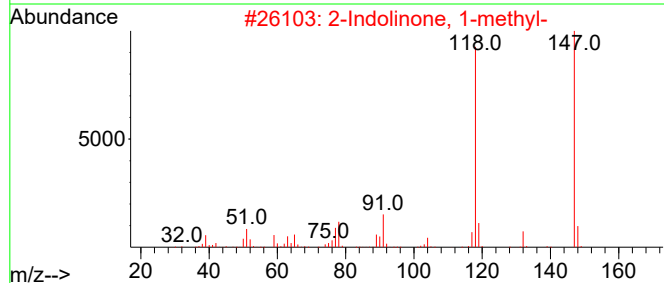
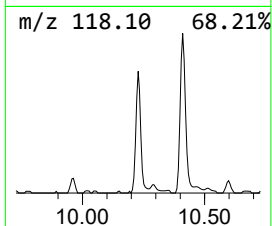
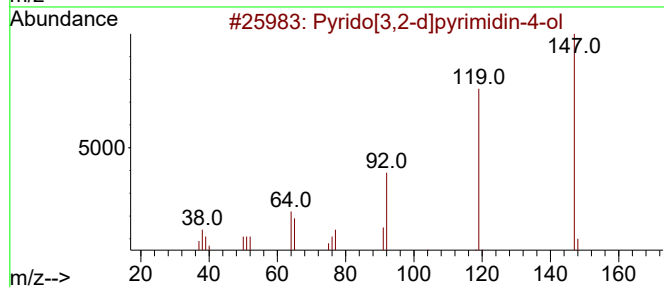
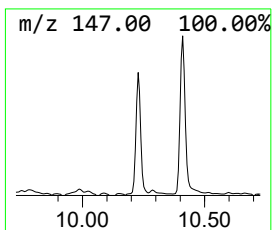
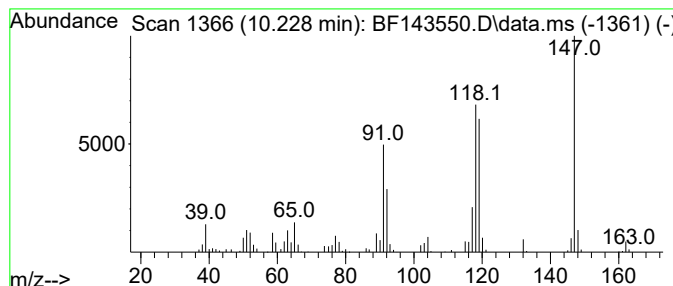
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Pyrido[3,2-d]pyrimidin-4-ol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.228	3.22 ng	139114	Acenaphthene-d10	9.963

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pyrido[3,2-d]pyrimidin-4-ol	147	C7H5N3O	037538-67-3	62
2			2-Indolinone, 1-methyl-	147	C9H9NO	000061-70-1	58
3			1H-Inden-1-one, 5-amino-2,3-dihy...	147	C9H9NO	003470-54-0	49
4			2-Indolizinemethanol	147	C9H9NO	022320-21-4	49
5			Pyrido[3,4-d]pyrimidin-4(3H)-one	147	C7H5N3O	019178-25-7	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082125\
Data File : BF143550.D
Acq On : 22 Aug 2025 01:51
Operator : RC/JU
Sample : Q2903-07
Misc :
ALS Vial : 31 Sample Multiplier: 1

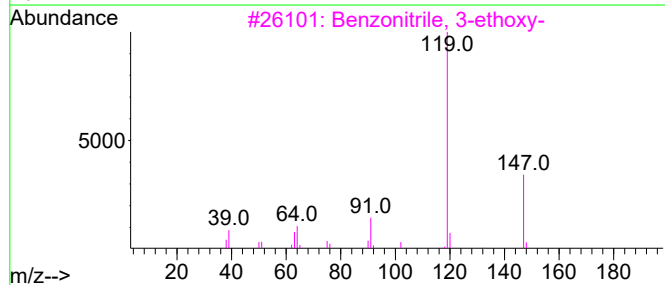
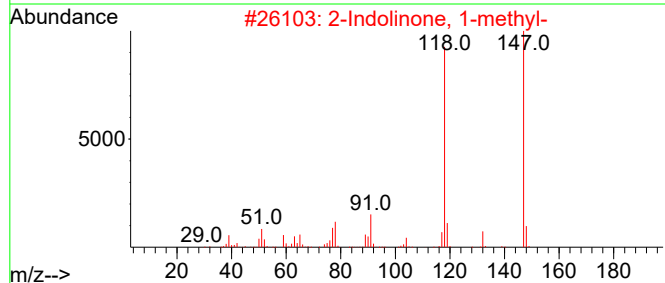
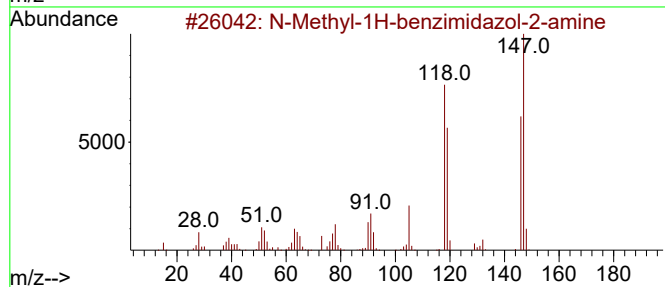
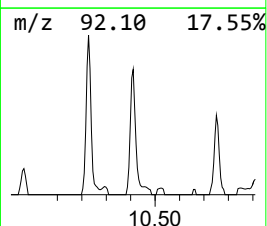
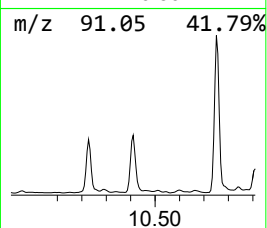
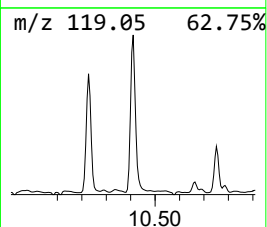
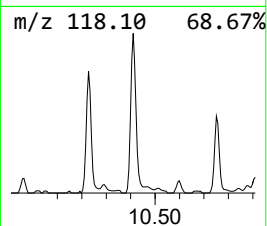
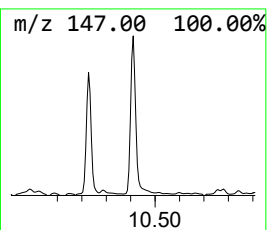
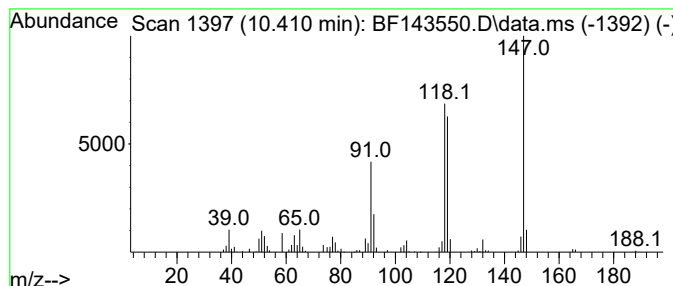
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ClientSampleId :
MW-12B-081925

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF082025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 N-Methyl-1H-benzimidazol-2-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.	
10.410	3.24 ng	140135	Acenaphthene-d10		9.963	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	N-Methyl-1H-benzimidazol-2-amine		147	C8H9N3	017228-38-5	86
2	2-Indolinone, 1-methyl-		147	C9H9NO	000061-70-1	68
3	Benzonitrile, 3-ethoxy-		147	C9H9NO	025117-75-3	52
4	5-Cyanotropolone		147	C8H5NO2	003266-92-0	52
5	3-(4,5-Dihydro-1H-imidazol-2-yl)...		147	C8H9N3	006302-53-0	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082125\
Data File : BF143550.D
Acq On : 22 Aug 2025 01:51
Operator : RC/JU
Sample : Q2903-07
Misc :
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Instrument :
BNA_F
ClientSampleId :
MW-12B-081925

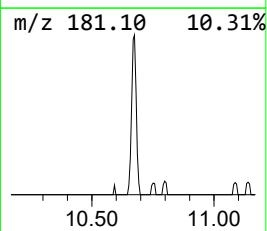
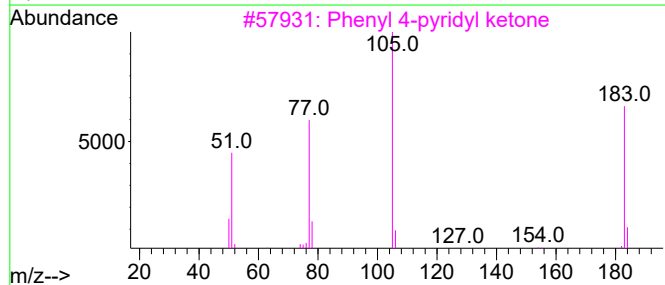
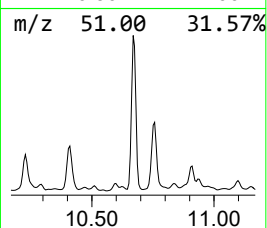
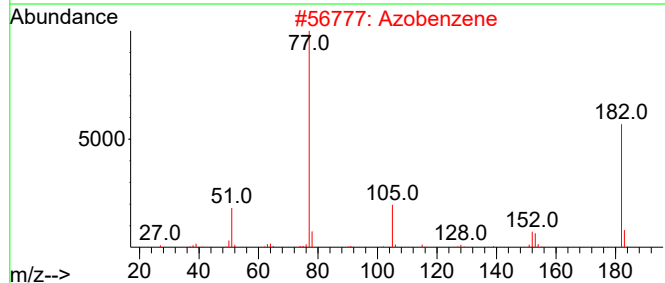
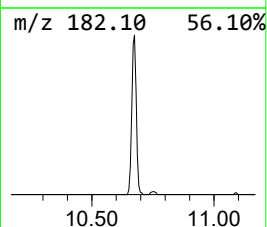
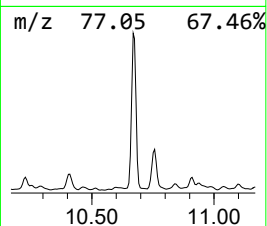
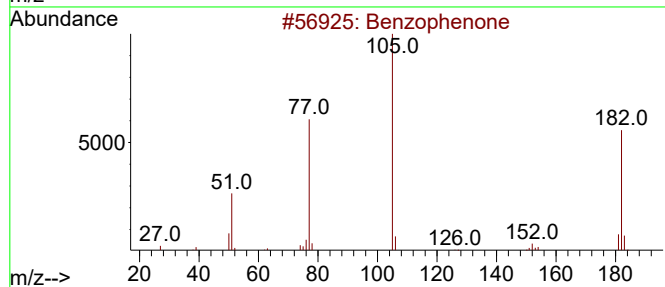
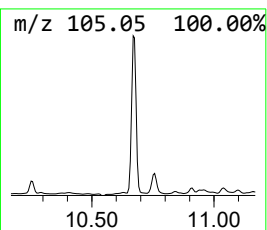
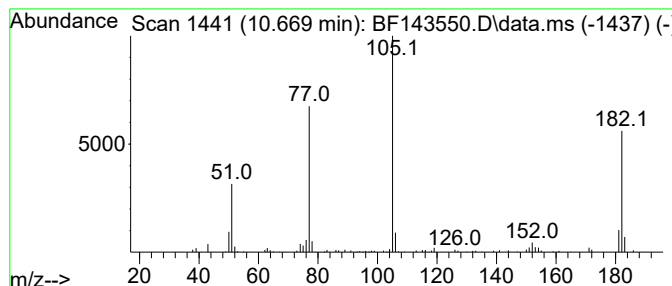
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Benzophenone Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.669	2.96 ng	127847	Acenaphthene-d10	9.963

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Azobenzene	182	C12H10N2	000103-33-3	58
3			Phenyl 4-pyridyl ketone	183	C12H9NO	014548-46-0	53
4			Benzoylformic acid	150	C8H6O3	000611-73-4	49
5			Acetophenone, 2-chloro-	154	C8H7ClO	000532-27-4	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082125\
Data File : BF143550.D
Acq On : 22 Aug 2025 01:51
Operator : RC/JU
Sample : Q2903-07
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
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ClientSampleId :
MW-12B-081925

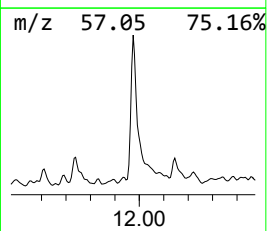
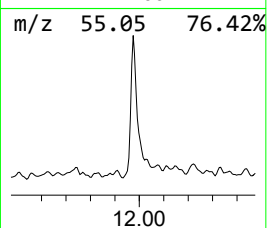
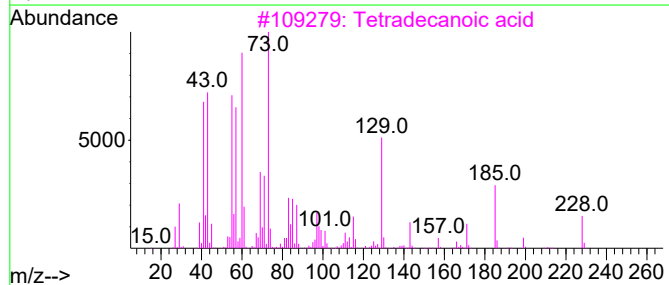
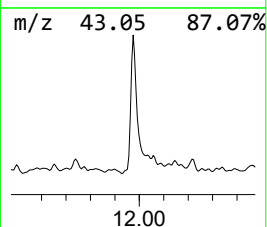
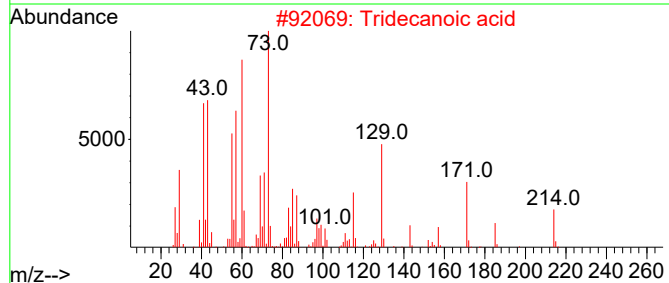
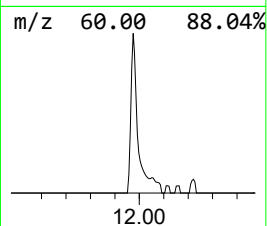
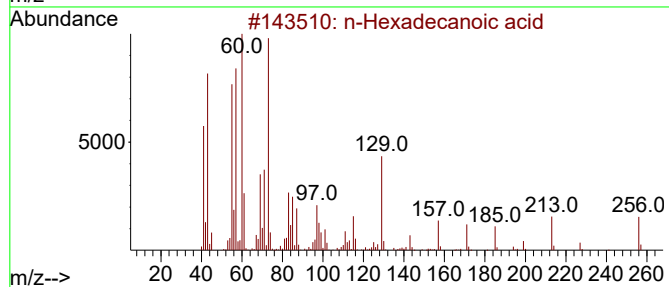
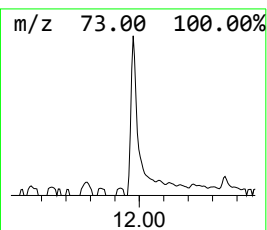
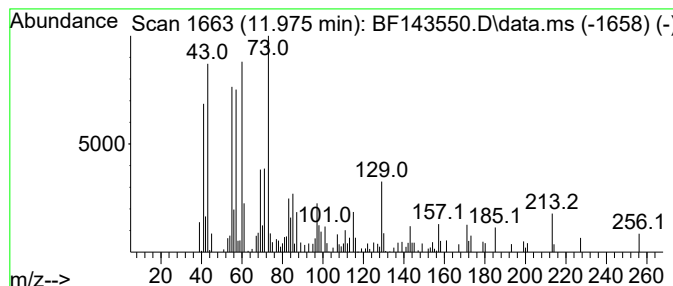
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 n-Hexadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.975	3.23 ng	107339	Phenanthrene-d10	11.451

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			Tridecanoic acid	214	C13H26O2	000638-53-9	95
3			Tetradecanoic acid	228	C14H28O2	000544-63-8	80
4			Pentadecanoic acid	242	C15H30O2	001002-84-2	76
5			Dodecanoic acid	200	C12H24O2	000143-07-7	62



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082125\
Data File : BF143550.D
Acq On : 22 Aug 2025 01:51
Operator : RC/JU
Sample : Q2903-07
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
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ClientSampleId :
MW-12B-081925

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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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2-Pentanone, 4-...	5.157	5.6	ng	174391	1	6.928	623756	20.0
Benzene, (1-met...	6.104	7.9	ng	247717	1	6.928	623756	20.0
Benzene, 1-ethy...	6.616	7.9	ng	247452	1	6.928	623756	20.0
Benzene, 1,2,3-...	6.987	5.1	ng	158848	1	6.928	623756	20.0
Indane	7.116	231.9	ng	7231140	1	6.928	623756	20.0
1H-Inden-5-ol, ...	9.098	2.5	ng	108498	3	9.963	863805	20.0
Pyrido[3,2-d]py...	10.228	3.2	ng	139114	3	9.963	863805	20.0
N-Methyl-1H-ben...	10.410	3.2	ng	140135	3	9.963	863805	20.0
Benzophenone	10.669	3.0	ng	127847	3	9.963	863805	20.0
n-Hexadecanoic ...	11.975	3.2	ng	107339	4	11.451	664507	20.0