

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082225\
 Data File : BF143582.D
 Acq On : 22 Aug 2025 19:50
 Operator : RC/JU
 Sample : Q2924-05
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-2B-082025

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: BF143582.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.252	3	10	17	rBV	1865232	3045920	9.64%	2.387%
2	5.440	544	552	557	rBV	5088336	8154404	25.82%	6.390%
3	5.557	563	572	579	rBV2	1323360	2507568	7.94%	1.965%
4	5.787	605	611	616	rBV	627813	802606	2.54%	0.629%
5	6.104	659	665	670	rBV	987835	1240375	3.93%	0.972%
6	6.387	709	713	720	rVB	304352	392468	1.24%	0.308%
7	6.528	733	737	739	rBV	314798	375601	1.19%	0.294%
8	6.563	739	743	748	rVV	851313	1217823	3.86%	0.954%
9	6.616	748	752	756	rVB	514687	642457	2.03%	0.503%
10	6.751	766	775	783	rVB	795168	1136128	3.60%	0.890%
11	6.928	800	805	809	rBV	434967	584600	1.85%	0.458%
12	6.987	809	815	820	rVV	1333733	2147064	6.80%	1.682%
13	7.034	820	823	829	rVB2	209386	396790	1.26%	0.311%
14	7.134	829	840	845	rBV	8880118	21634120	68.49%	16.952%
15	7.193	845	850	854	rVB2	674733	992119	3.14%	0.777%
16	7.493	888	901	910	rBV3	1877585	3385447	10.72%	2.653%
17	7.722	937	940	944	rVB	273379	332180	1.05%	0.260%
18	7.881	960	967	972	rVB	1892993	2529269	8.01%	1.982%
19	7.951	972	979	983	rBV3	2212379	3463267	10.96%	2.714%
20	7.998	983	987	992	rVV	1144970	1981735	6.27%	1.553%
21	8.051	992	996	1008	rVB	377995	531753	1.68%	0.417%
22	8.269	1017	1033	1043	rBV	9306882	31585166	100.00%	24.749%
23	8.469	1063	1067	1074	rVB2	311755	562692	1.78%	0.441%
24	8.675	1093	1102	1105	rBV4	165653	440651	1.40%	0.345%
25	8.798	1113	1123	1128	rVV2	396014	577477	1.83%	0.452%
26	8.875	1128	1136	1139	rVV	354686	509459	1.61%	0.399%
27	8.928	1139	1145	1150	rVV	3904521	5620762	17.80%	4.404%
28	9.034	1156	1163	1168	rVV	5178518	8424282	26.67%	6.601%
29	9.281	1199	1205	1210	rBV	2709711	3513710	11.12%	2.753%
30	9.381	1216	1222	1227	rBV	534116	704372	2.23%	0.552%
31	9.475	1235	1238	1244	rVV2	307957	525538	1.66%	0.412%
32	9.545	1244	1250	1255	rVV	599317	974129	3.08%	0.763%
33	9.622	1258	1263	1266	rVV	752509	1169606	3.70%	0.916%
34	9.645	1266	1267	1272	rVB2	413598	370153	1.17%	0.290%
35	9.739	1278	1283	1290	rVV	361556	555003	1.76%	0.435%
36	9.963	1311	1321	1324	rBV	795048	1302769	4.12%	1.021%

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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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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	9.998	1324	1327	1333	rVV	2302932	2974475	9.42%	2.331%
38	10.057	1333	1337	1342	rVB	328895	444539	1.41%	0.348%
39	10.510	1410	1414	1419	rVB	457175	592878	1.88%	0.465%
40	10.586	1419	1427	1431	rBV3	243257	553162	1.75%	0.433%
41	10.757	1447	1456	1467	rVB2	1757545	2582320	8.18%	2.023%
42	11.451	1569	1574	1576	rBV	689751	916636	2.90%	0.718%
43	11.475	1576	1578	1582	rVB	597211	631423	2.00%	0.495%
44	11.980	1661	1664	1668	rBV2	230371	339194	1.07%	0.266%
45	13.039	1838	1844	1849	rVB	2460817	3194323	10.11%	2.503%
46	14.092	2018	2023	2036	rVB	423195	564759	1.79%	0.443%
47	15.592	2272	2278	2292	rBV	316305	496686	1.57%	0.389%

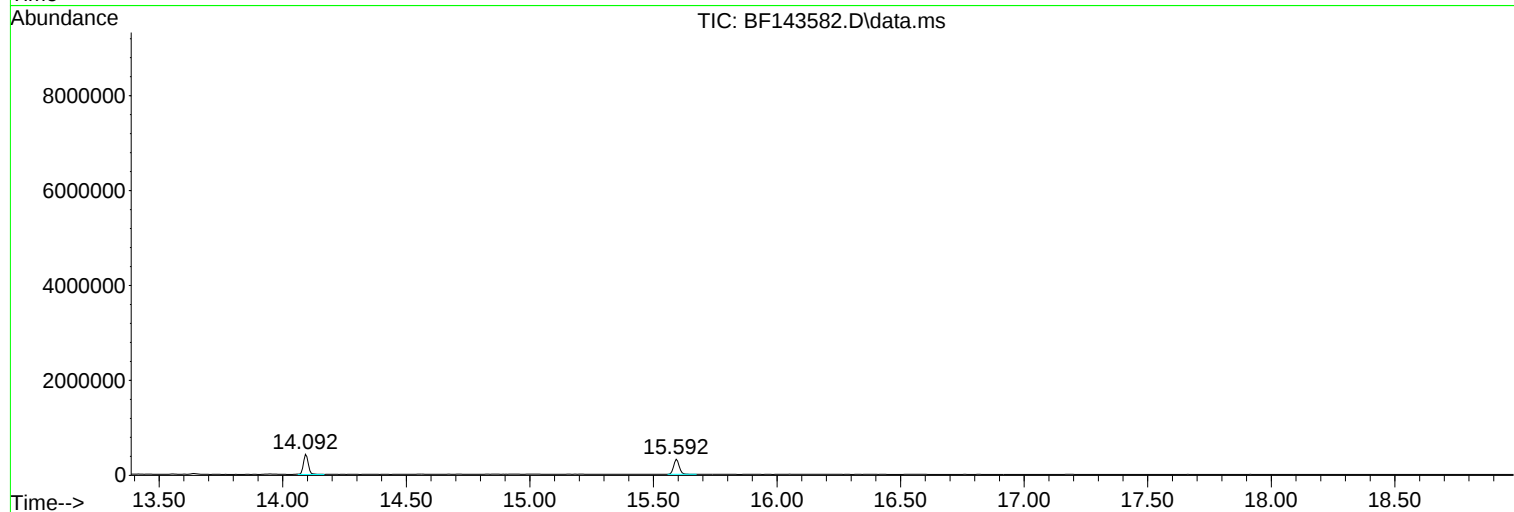
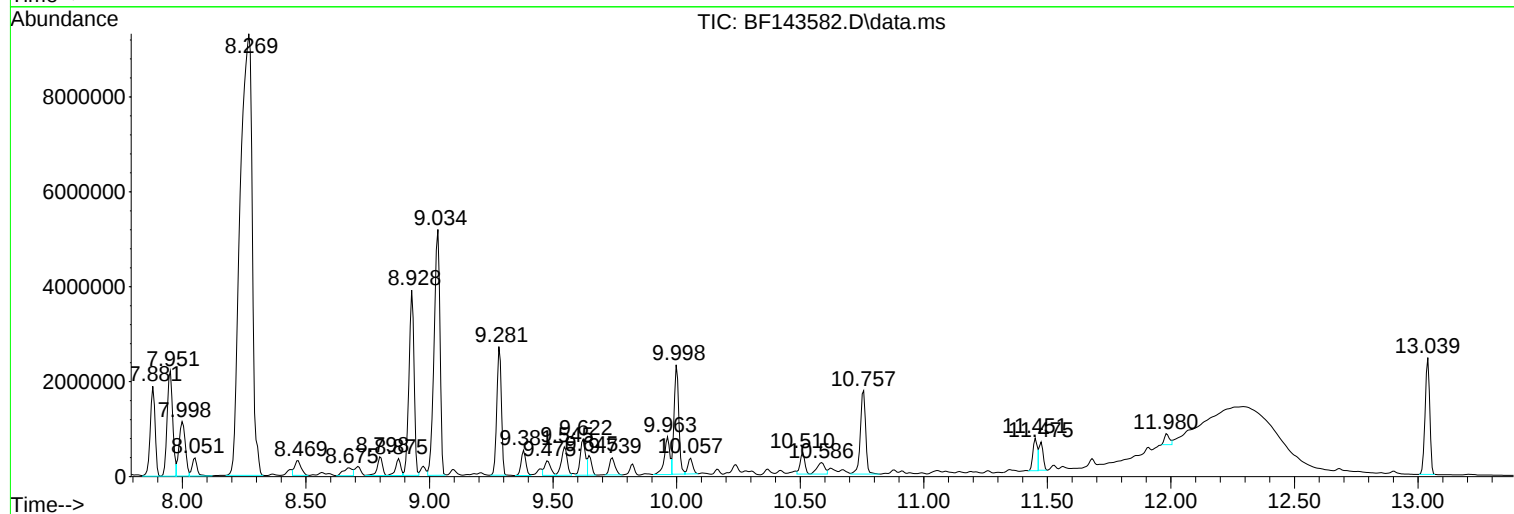
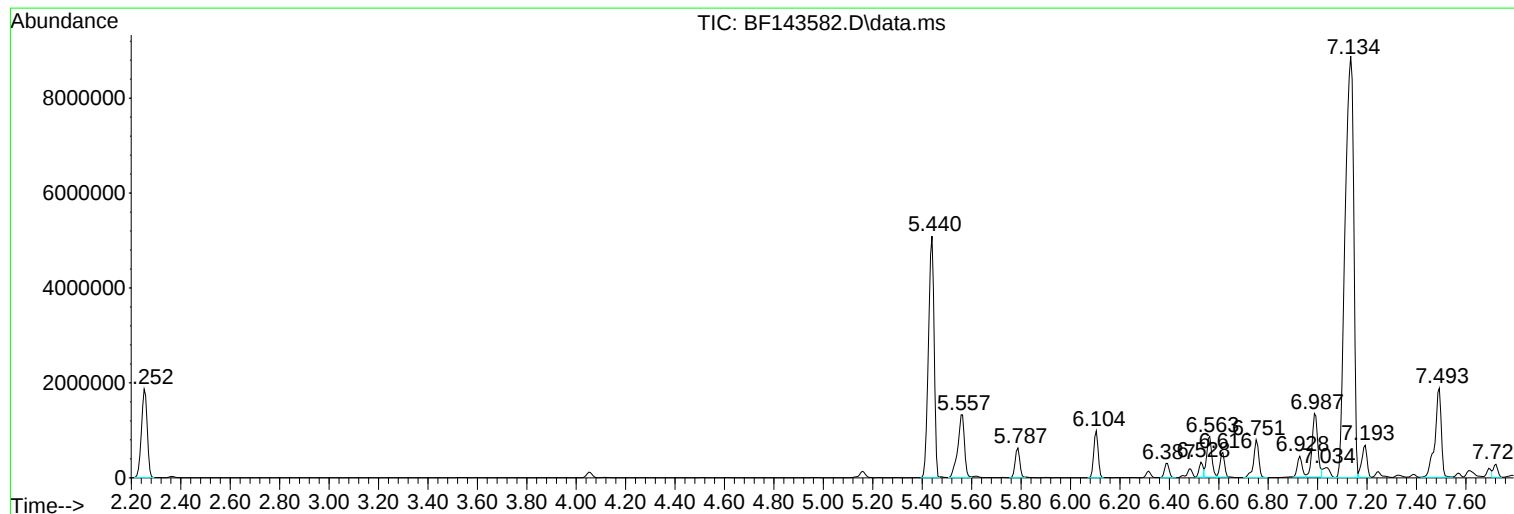
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF082025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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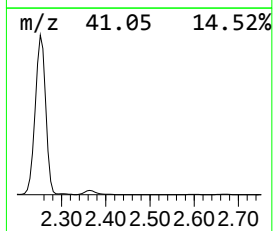
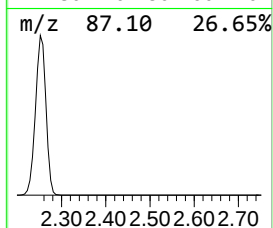
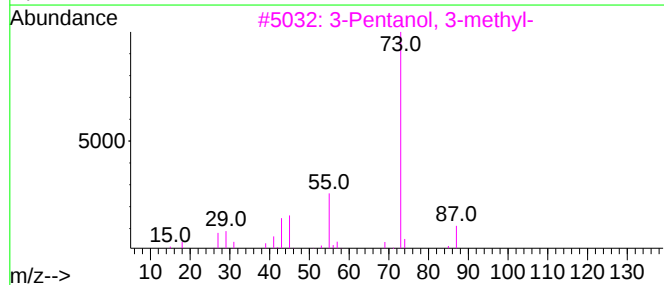
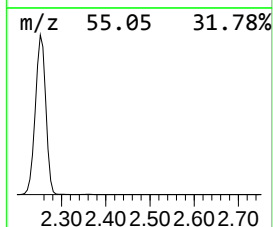
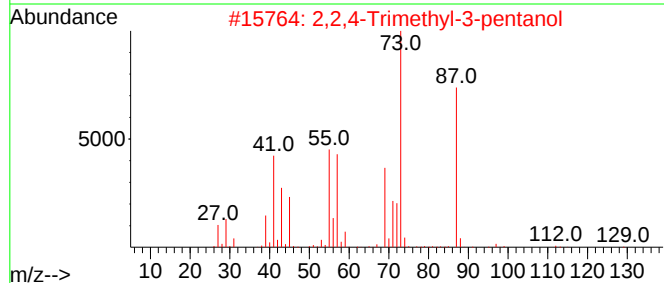
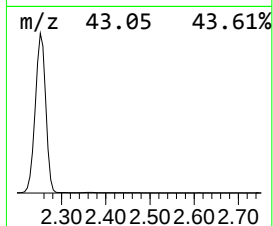
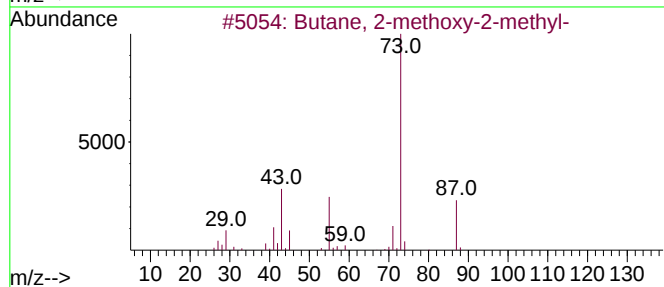
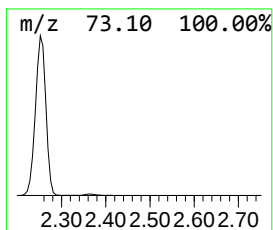
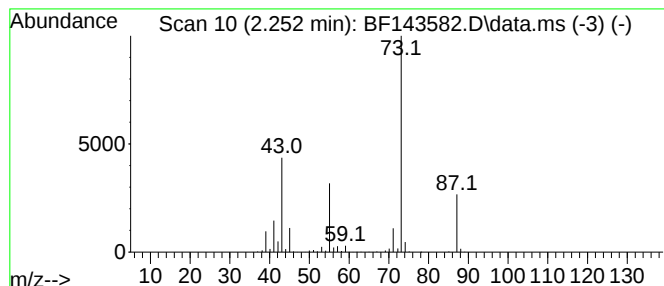
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF082025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.252	104.21 ng	3045920	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			3-Pentanol, 3-methyl-	102	C6H14O	000077-74-7	17
4			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12
5			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9



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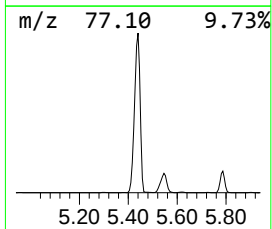
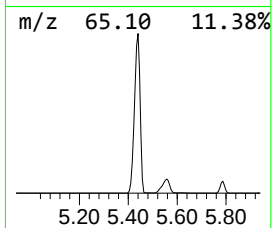
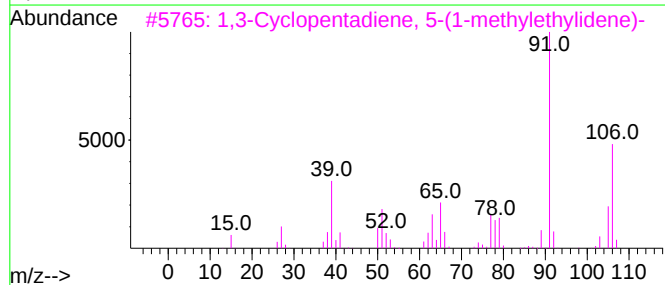
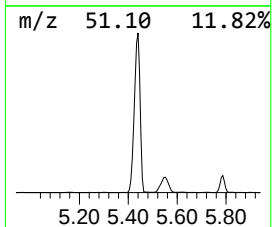
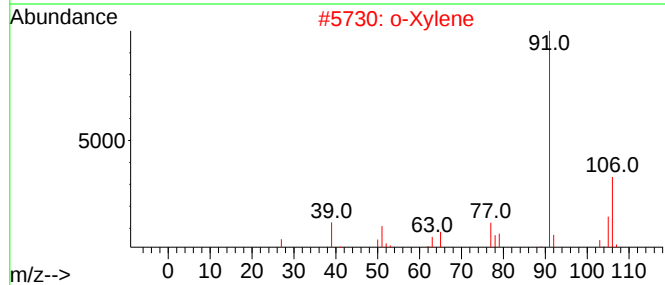
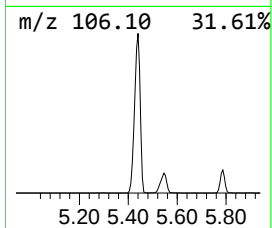
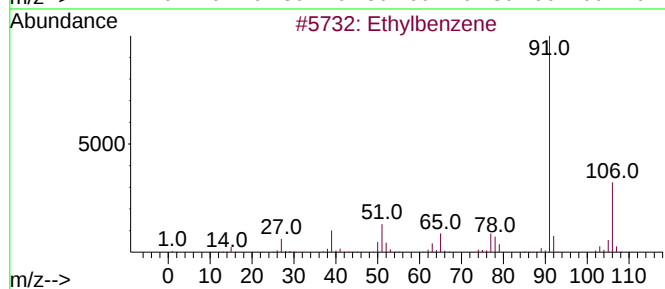
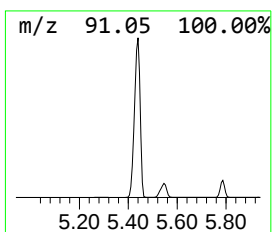
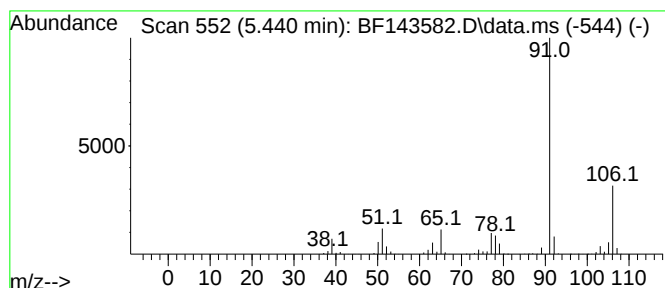
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF082025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 1,3-Cyclopentadiene, 5-(1-m... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.440	278.97 ng	8154400	1,4-Dichlorobenzene-d4	6.928

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Ethylbenzene	106	C8H10	000100-41-4	94
2		o-Xylene	106	C8H10	000095-47-6	81
3		1,3-Cyclopentadiene, 5-(1-methyl...	106	C8H10	002175-91-9	74
4		p-Xylene	106	C8H10	000106-42-3	72
5		Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	64



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ALS Vial : 21 Sample Multiplier: 1

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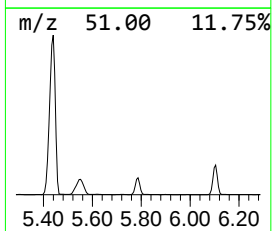
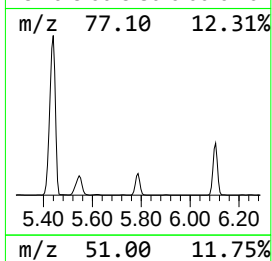
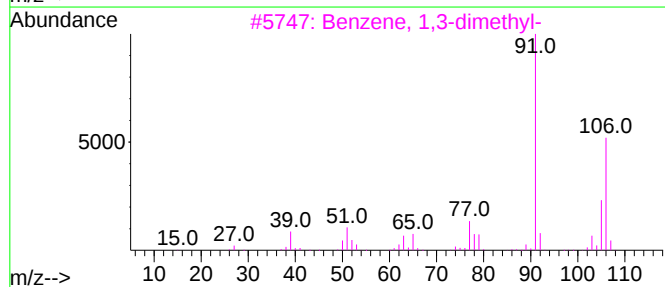
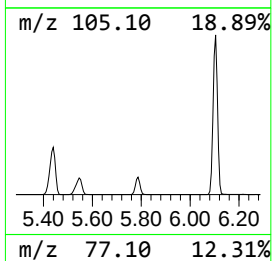
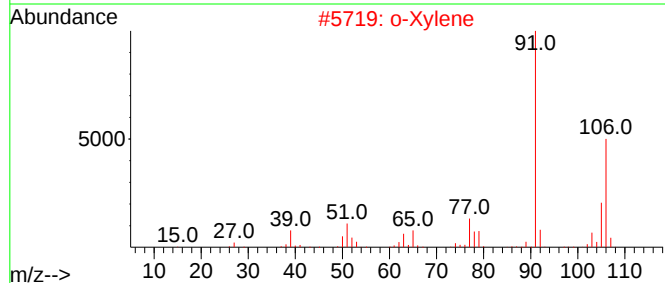
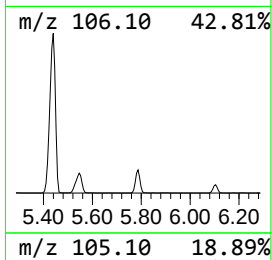
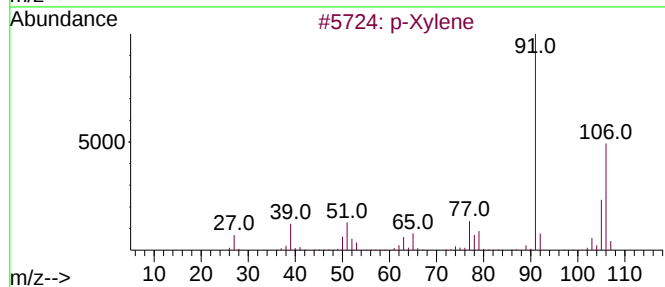
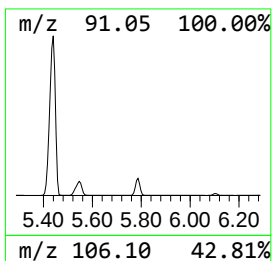
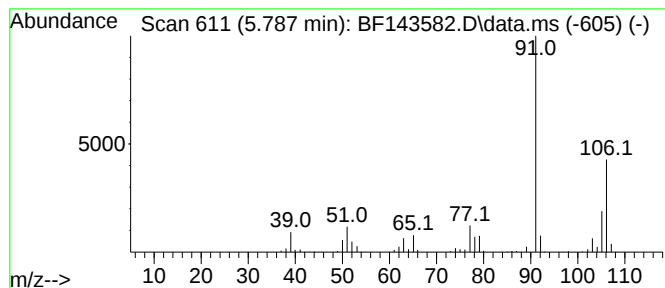
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, 1,3-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.787	27.46 ng	802606	1,4-Dichlorobenzene-d4	6.928

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			p-Xylene	106	C8H10	000106-42-3	97
2			o-Xylene	106	C8H10	000095-47-6	97
3			Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
4			1,3-Cyclopentadiene, 5-(1-methyl...	106	C8H10	002175-91-9	91
5			Ethylbenzene	106	C8H10	000100-41-4	91



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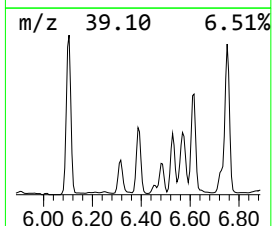
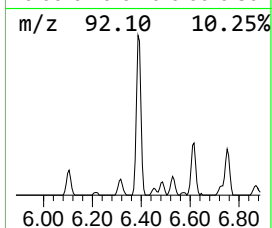
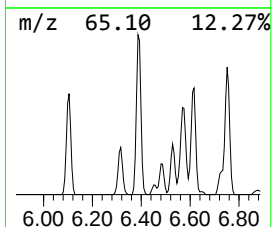
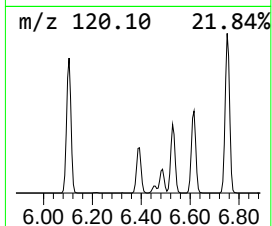
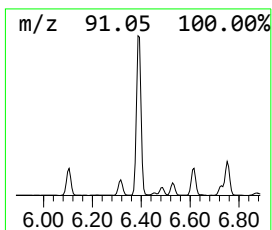
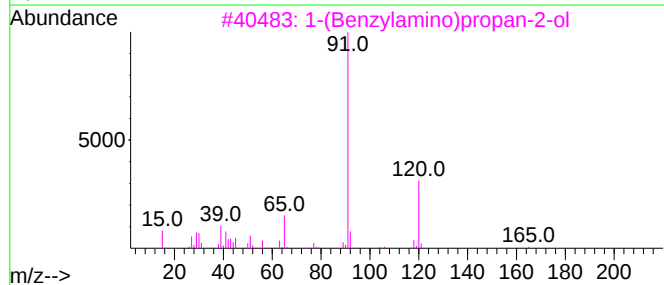
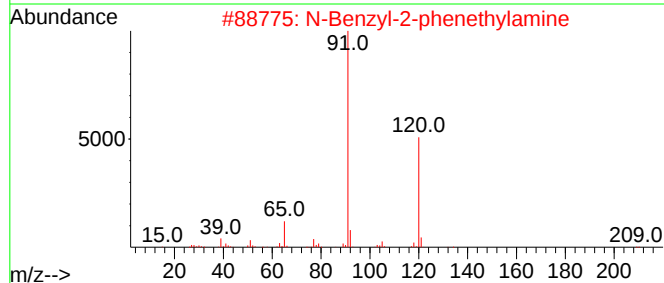
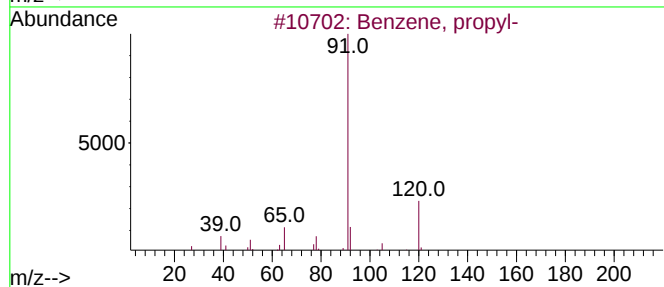
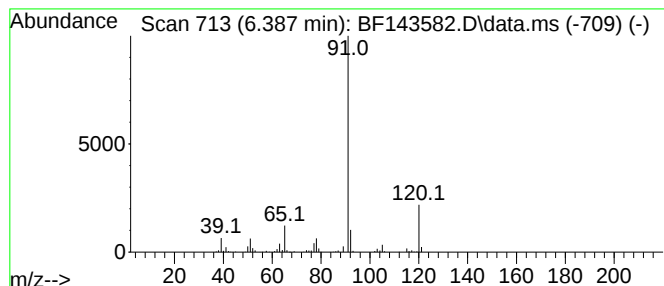
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 unknown6.387 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.387	13.43 ng	392468	1,4-Dichlorobenzene-d4	6.928

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			1-(Benzylamino)propan-2-ol	165	C10H15NO	1000486-84-7	72
4			1,2-Ethanediamine, N,N'-bis(phen...	240	C16H20N2	000140-28-3	72
5			Benzeneacetaldehyde	120	C8H8O	000122-78-1	62



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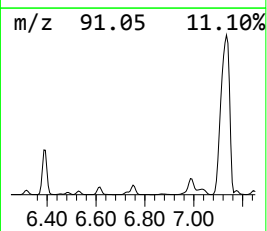
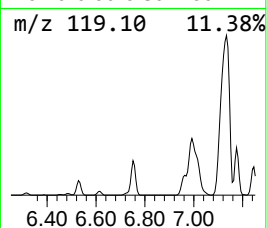
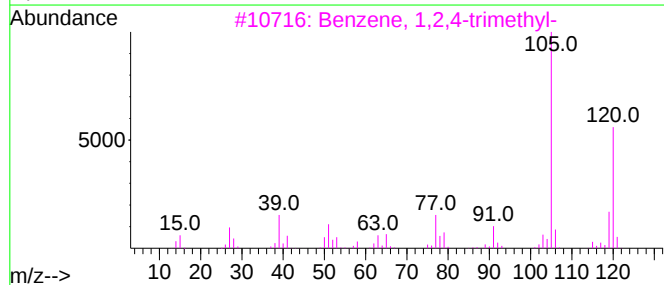
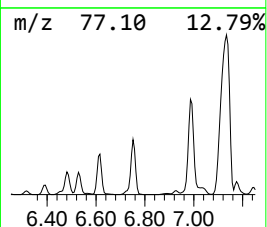
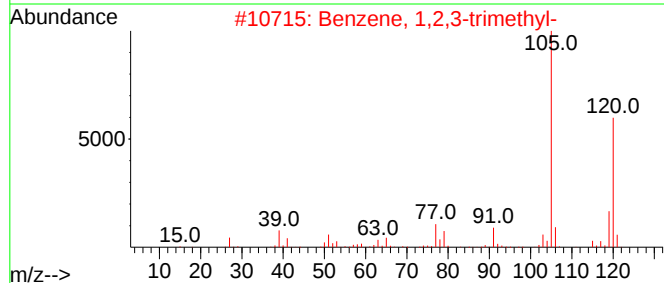
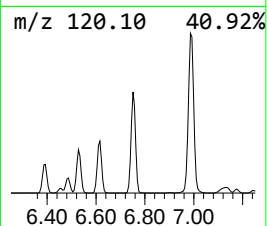
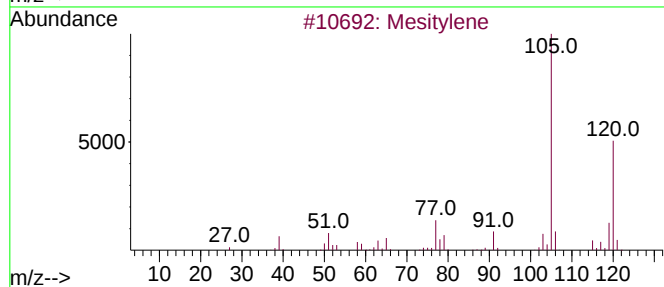
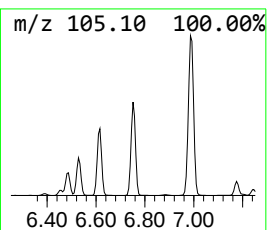
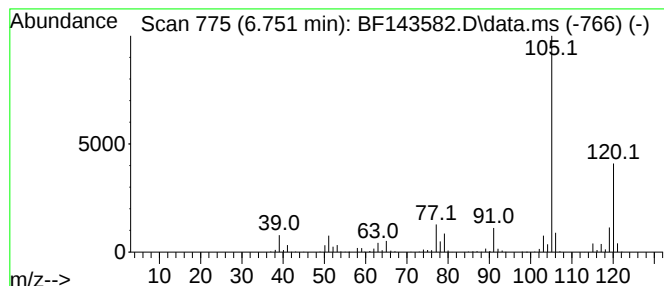
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 unknown6.751 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.751	38.87 ng	1136130	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	97
3			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	95
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082225\
Data File : BF143582.D
Acq On : 22 Aug 2025 19:50
Operator : RC/JU
Sample : Q2924-05
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-2B-082025

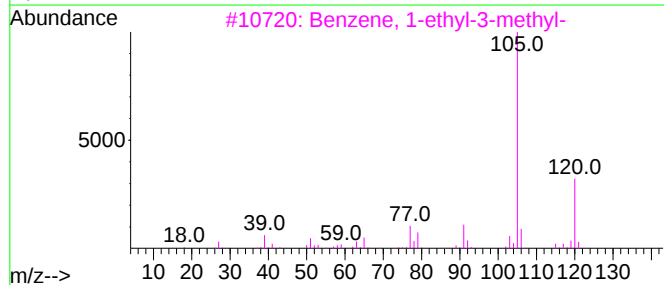
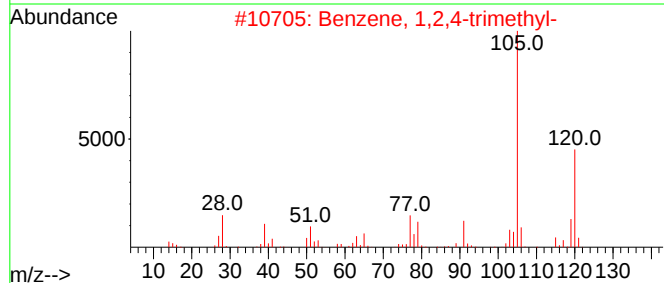
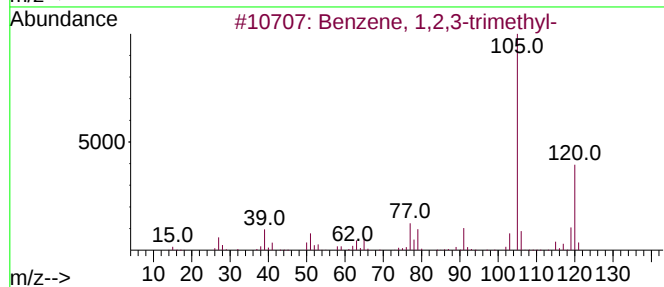
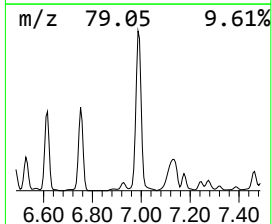
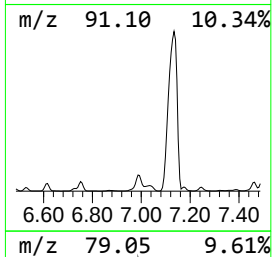
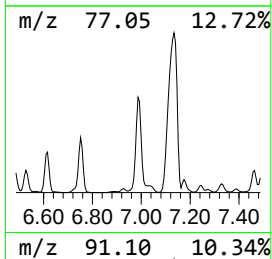
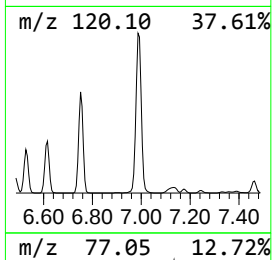
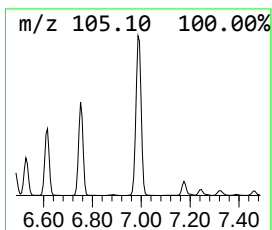
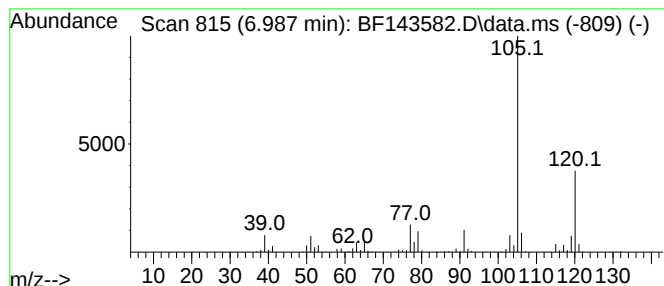
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown6.987 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.987	73.45 ng	2147060	1,4-Dichlorobenzene-d4	6.928

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082225\
Data File : BF143582.D
Acq On : 22 Aug 2025 19:50
Operator : RC/JU
Sample : Q2924-05
Misc :
ALS Vial : 21 Sample Multiplier: 1

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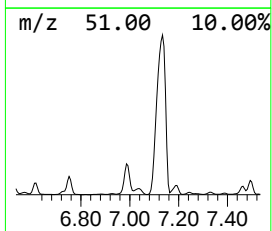
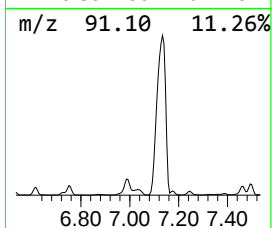
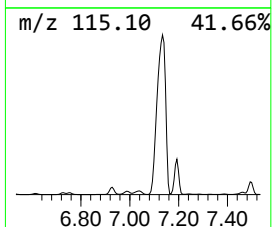
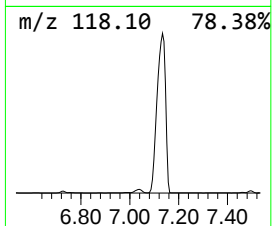
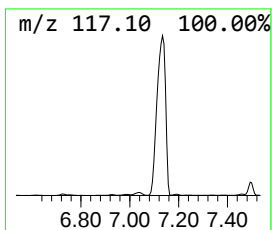
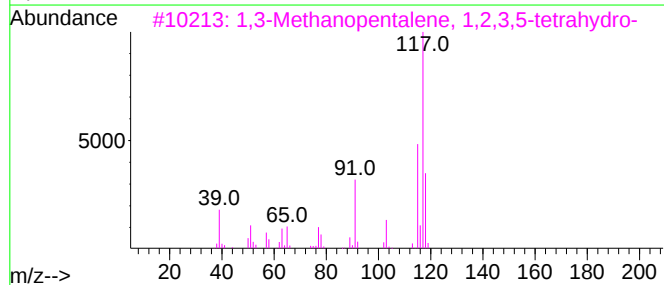
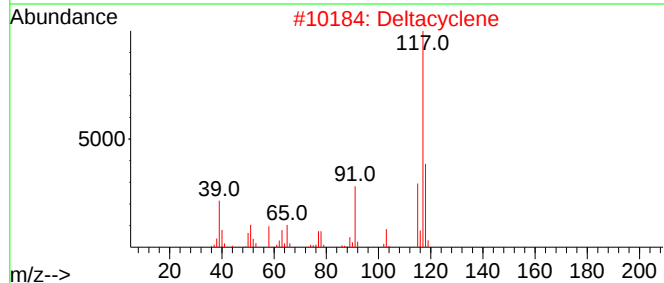
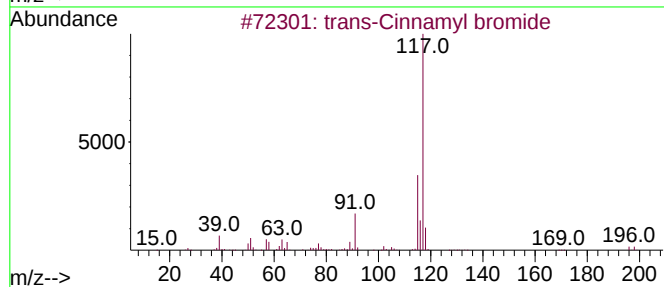
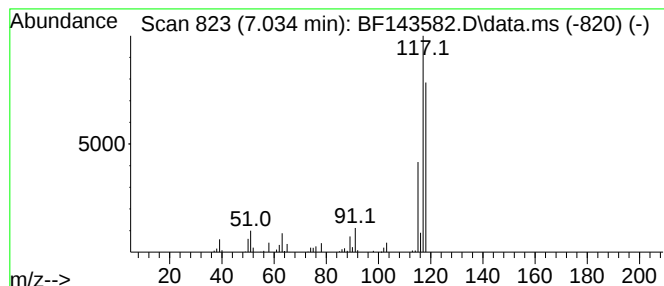
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown7.034 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.034	13.57 ng	396790	1,4-Dichlorobenzene-d4	6.928

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			trans-Cinnamyl bromide	196	C9H9Br	026146-77-0	47
2			Deltacyclene	118	C9H10	007785-10-6	43
3			1,3-Methanopentalene, 1,2,3,5-te...	118	C9H10	128600-88-4	37
4			Benzene, 1,1'-(1,5-hexadiene-1,6...	234	C18H18	004439-45-6	32
5			Benzene, (2-bromocyclopropyl)-	196	C9H9Br	036617-02-4	32



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082225\
Data File : BF143582.D
Acq On : 22 Aug 2025 19:50
Operator : RC/JU
Sample : Q2924-05
Misc :
ALS Vial : 21 Sample Multiplier: 1

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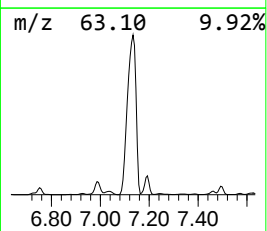
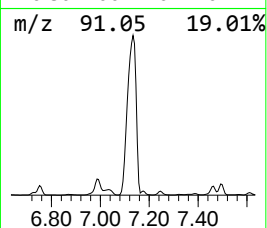
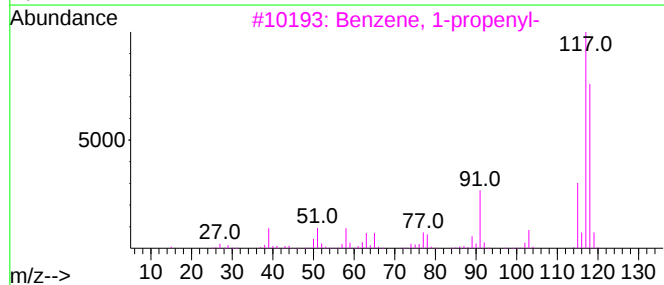
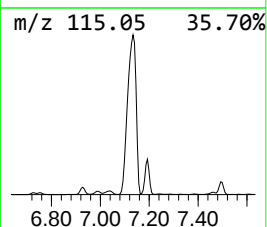
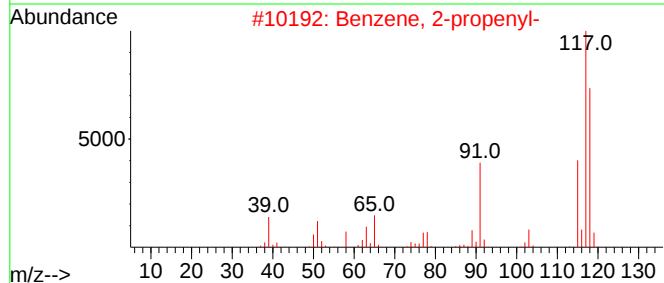
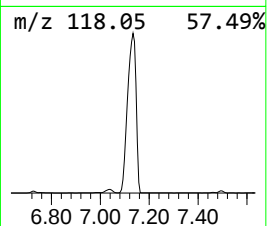
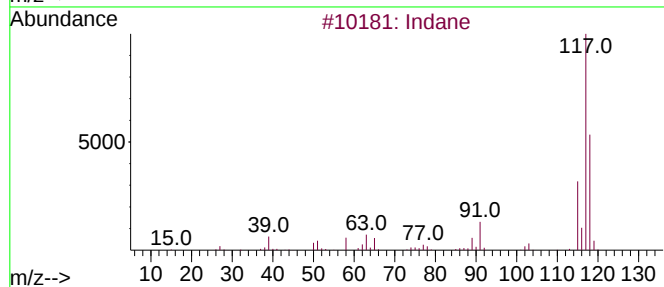
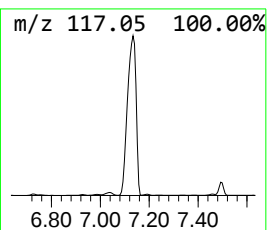
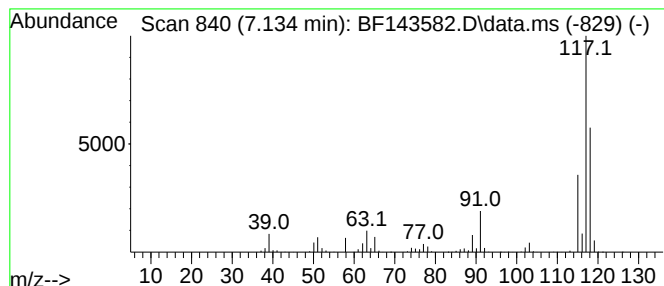
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 unknown7.134 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.134	740.13 ng	21634100	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-propenyl-	118	C9H10	000637-50-3	76
4			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	74
5			Benzene, cyclopropyl-	118	C9H10	000873-49-4	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082225\
Data File : BF143582.D
Acq On : 22 Aug 2025 19:50
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Sample : Q2924-05
Misc :
ALS Vial : 21 Sample Multiplier: 1

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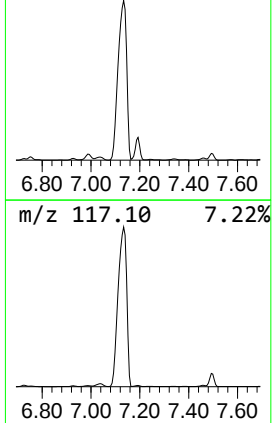
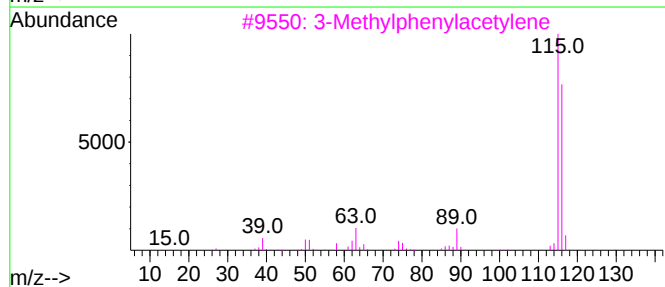
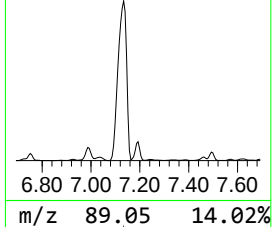
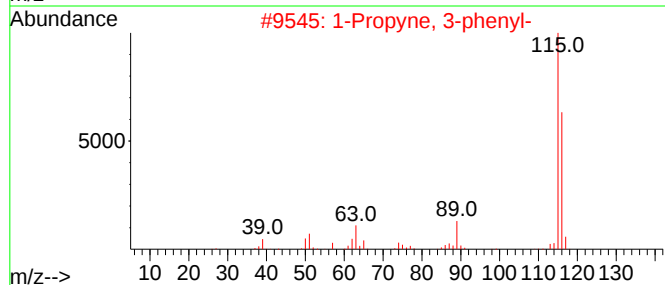
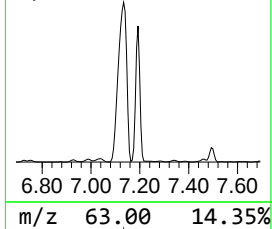
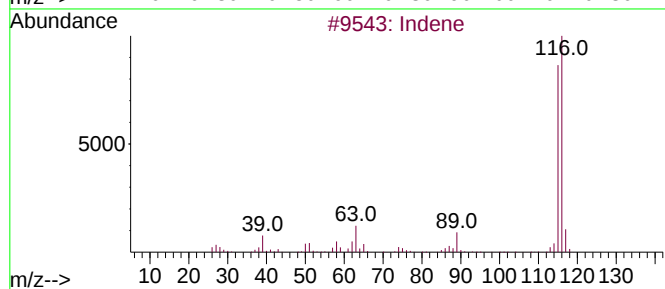
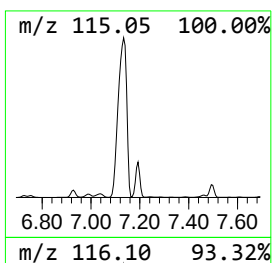
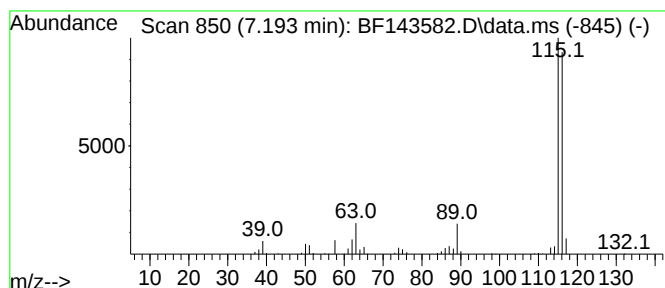
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown7.193 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.193	33.94 ng	992119	1,4-Dichlorobenzene-d4	6.928

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indene	116	C9H8	000095-13-6	94
2			1-Propyne, 3-phenyl-	116	C9H8	010147-11-2	94
3			3-Methylphenylacetylene	116	C9H8	000766-82-5	91
4			Benzene, 1,2-propadienyl-	116	C9H8	002327-99-3	91
5			Benzene, 1-propynyl-	116	C9H8	000673-32-5	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082225\
Data File : BF143582.D
Acq On : 22 Aug 2025 19:50
Operator : RC/JU
Sample : Q2924-05
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
MW-2B-082025

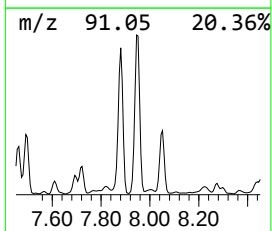
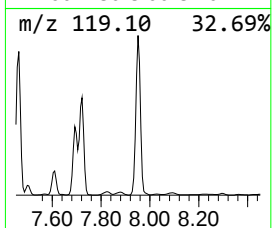
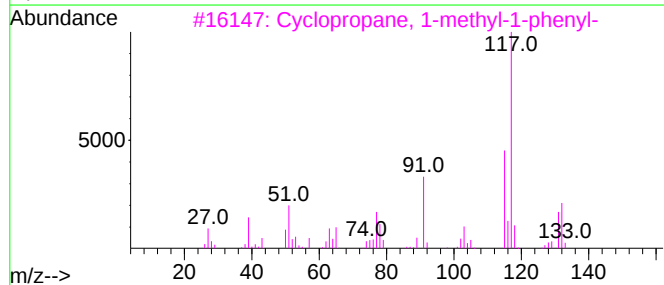
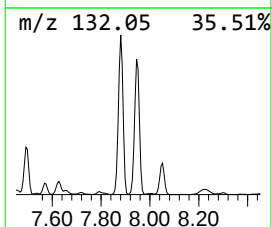
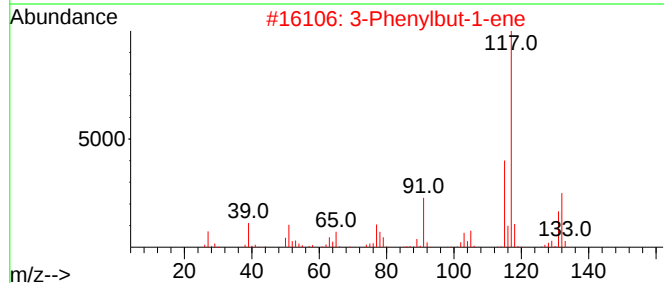
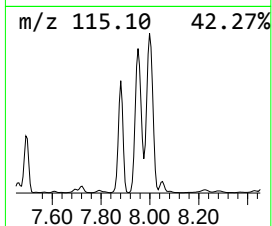
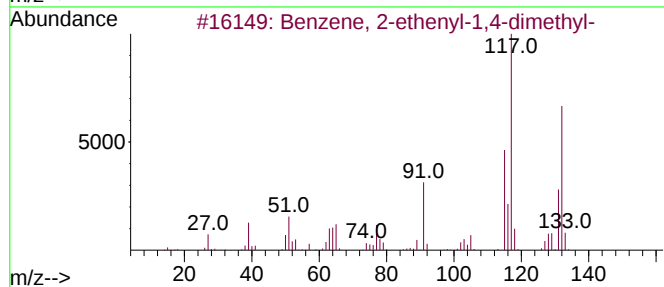
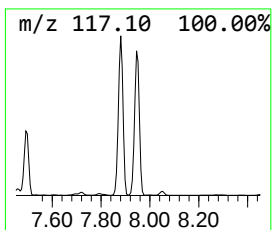
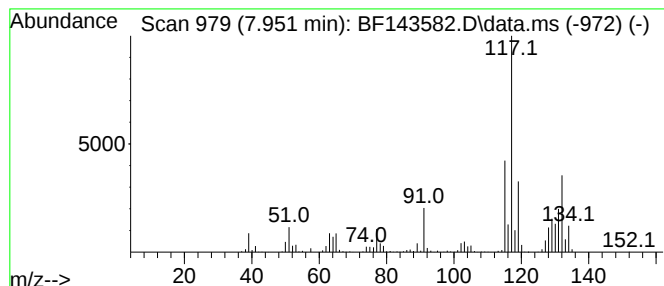
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown7.951 Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.951	2.19 ng	3463270	Naphthalene-d8	8.228

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	86
2		3-Phenylbut-1-ene	132	C10H12	000934-10-1	76
3		Cyclopropane, 1-methyl-1-phenyl-	132	C10H12	002214-14-4	70
4		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	64
5		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082225\
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Sample : Q2924-05
Misc :
ALS Vial : 21 Sample Multiplier: 1

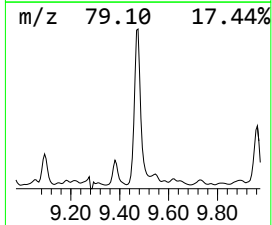
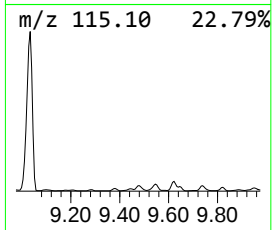
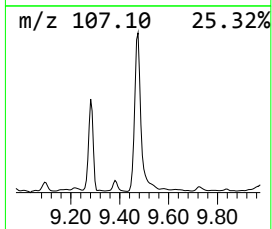
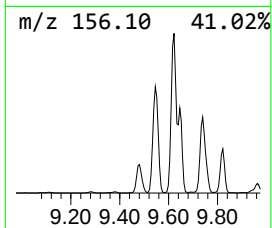
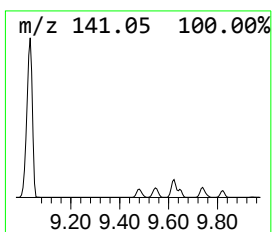
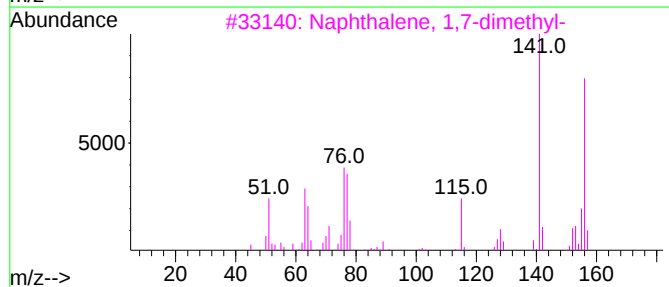
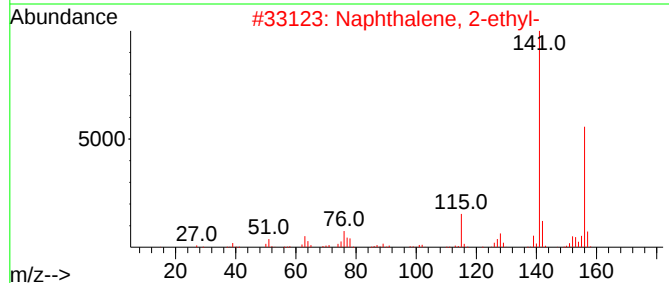
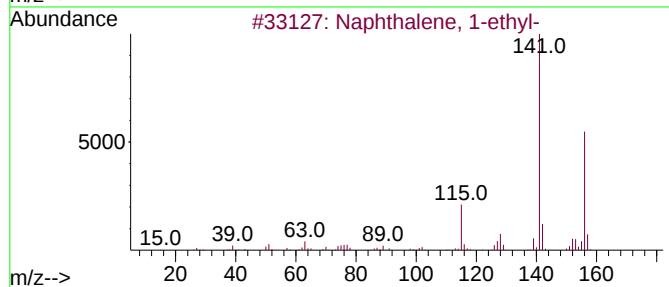
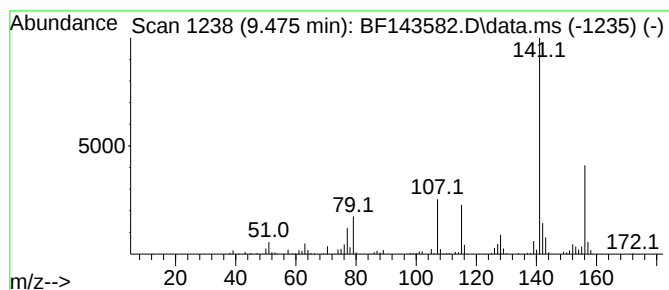
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MW-2B-082025

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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 12 Naphthalene, 1-ethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.	
9.475	8.07 ng	525538	Acenaphthene-d10	9.963	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	95
2	Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	95
3	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	72
4	2-Amino-4-tertbutylthiazole	156	C7H12N2S	074370-93-7	53
5	2-Thiophenecarboxylic acid, 5-et...	156	C7H8O2S	023229-72-3	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082225\
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Acq On : 22 Aug 2025 19:50
Operator : RC/JU
Sample : Q2924-05
Misc :
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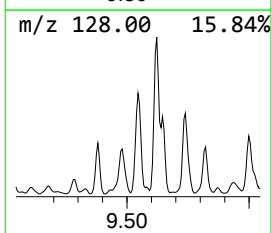
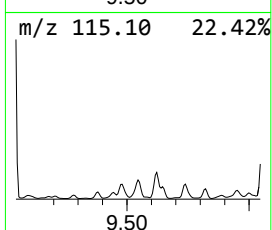
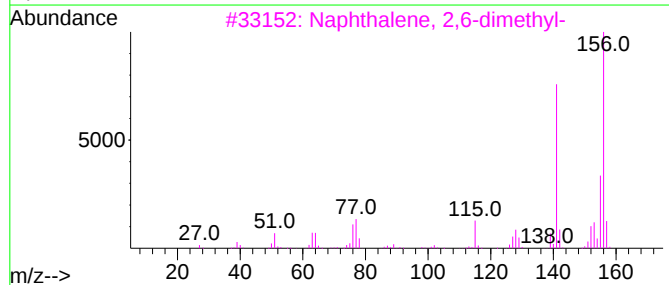
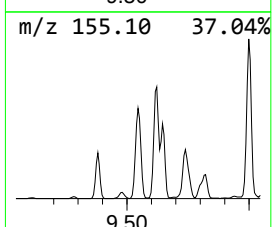
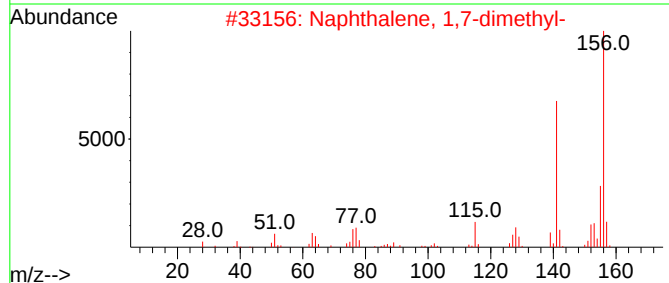
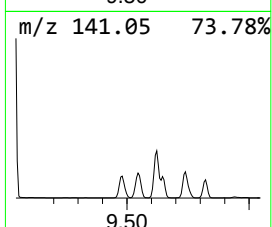
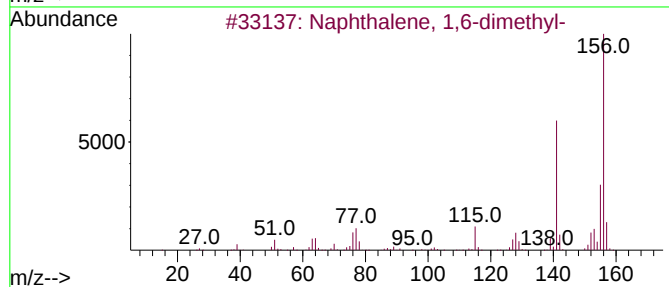
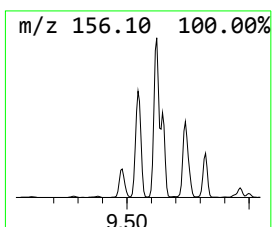
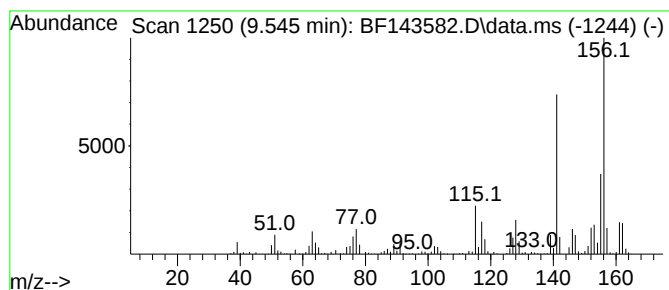
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ClientSampleId :
MW-2B-082025

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 13 Naphthalene, 1,6-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD			R.T.
9.545	14.95 ng	974129	Acenaphthene-d10			9.963
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Naphthalene, 1,6-dimethyl-		156	C12H12	000575-43-9	96
2	Naphthalene, 1,7-dimethyl-		156	C12H12	000575-37-1	96
3	Naphthalene, 2,6-dimethyl-		156	C12H12	000581-42-0	96
4	Naphthalene, 2,7-dimethyl-		156	C12H12	000582-16-1	96
5	Naphthalene, 2,3-dimethyl-		156	C12H12	000581-40-8	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082225\
Data File : BF143582.D
Acq On : 22 Aug 2025 19:50
Operator : RC/JU
Sample : Q2924-05
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-2B-082025

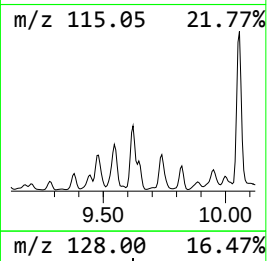
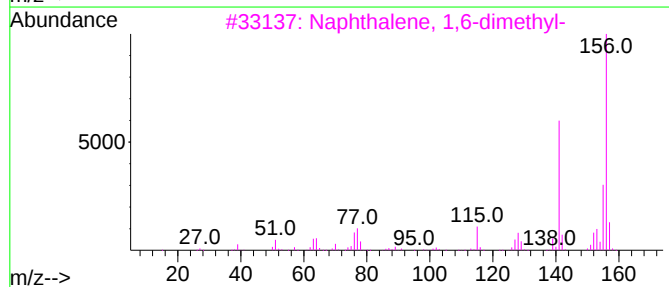
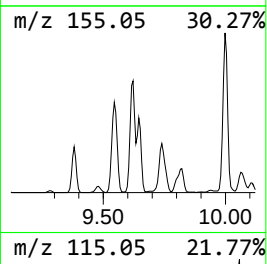
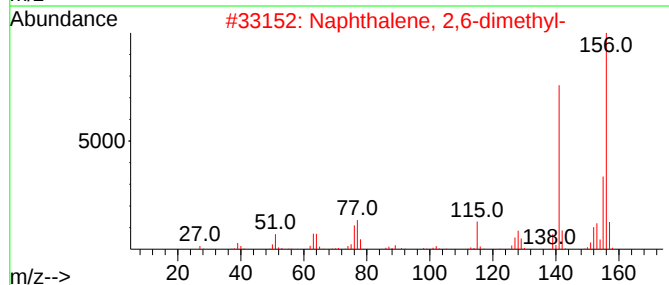
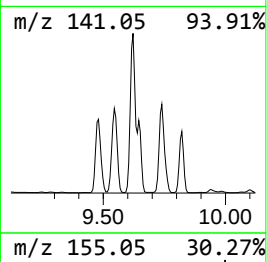
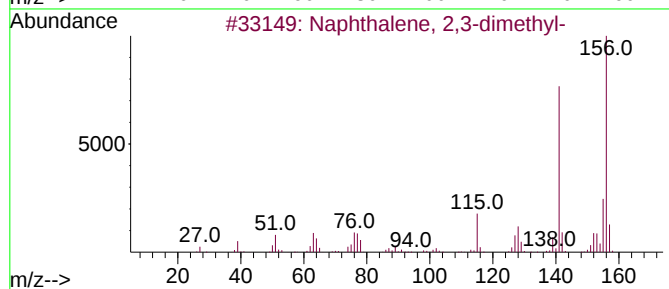
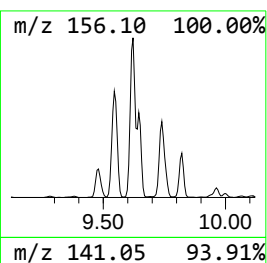
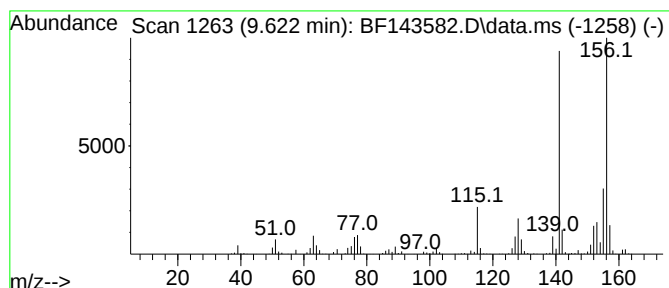
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Naphthalene, 2,3-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.622	17.96 ng	1169610	Acenaphthene-d10	9.963

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
2		Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
3		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
4		Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	96
5		Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082225\
Data File : BF143582.D
Acq On : 22 Aug 2025 19:50
Operator : RC/JU
Sample : Q2924-05
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-2B-082025

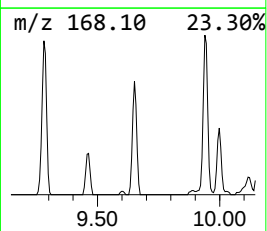
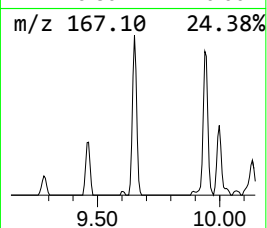
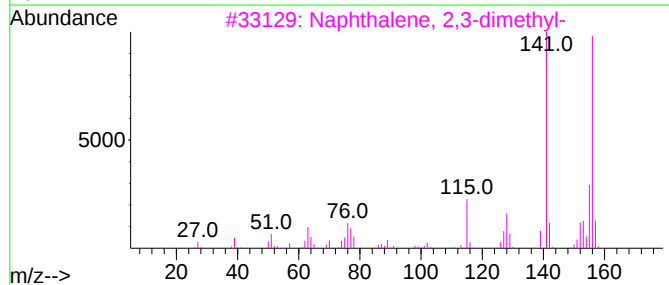
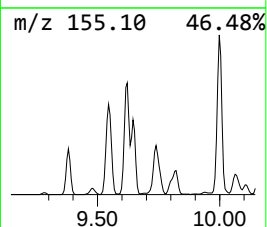
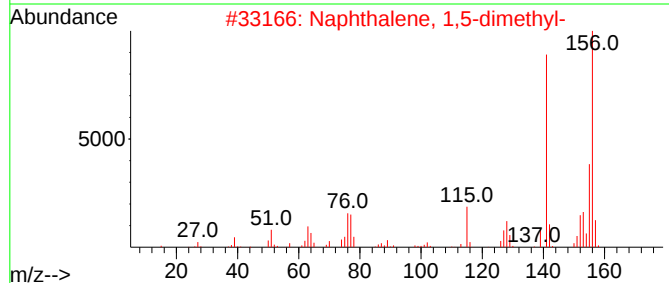
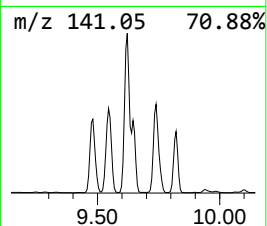
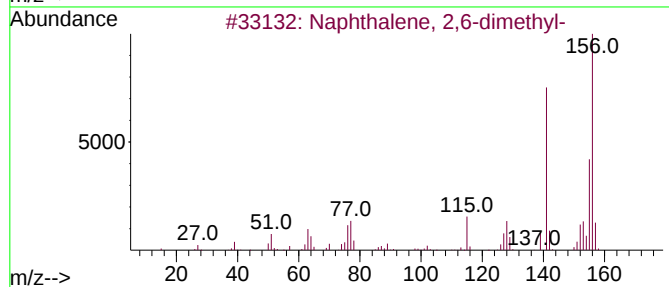
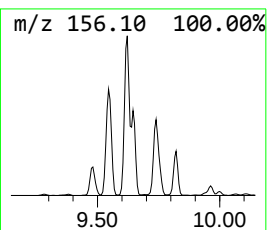
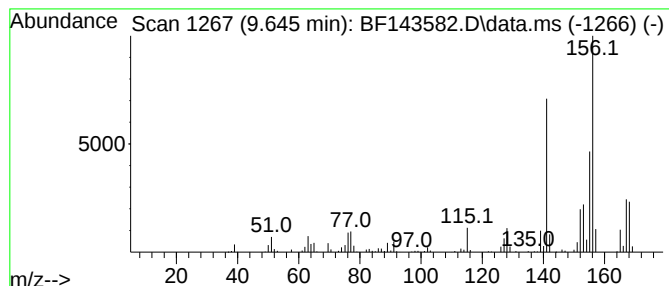
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Naphthalene, 2,6-dimethyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.645	5.68 ng	370153	Acenaphthene-d10	9.963

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	86
2			Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	83
3			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	78
4			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	78
5			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082225\
Data File : BF143582.D
Acq On : 22 Aug 2025 19:50
Operator : RC/JU
Sample : Q2924-05
Misc :
ALS Vial : 21 Sample Multiplier: 1

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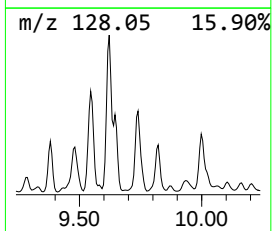
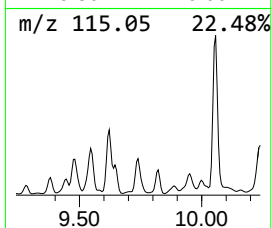
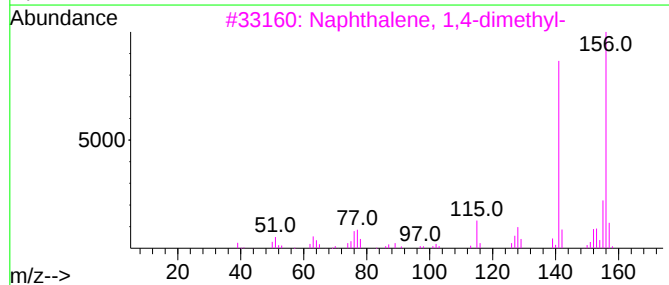
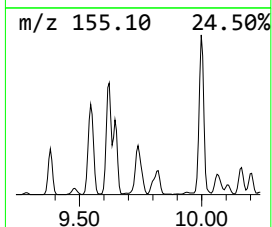
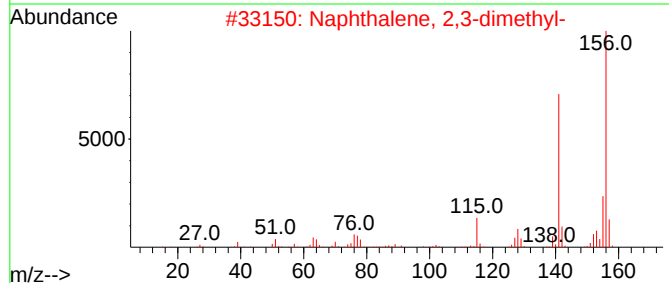
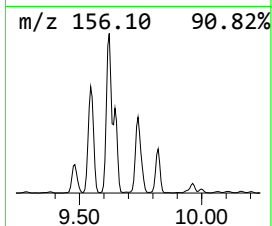
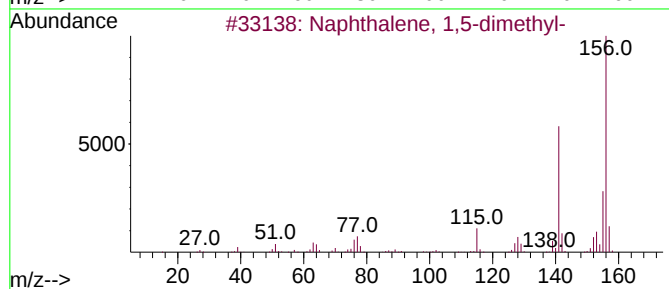
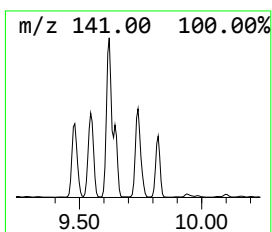
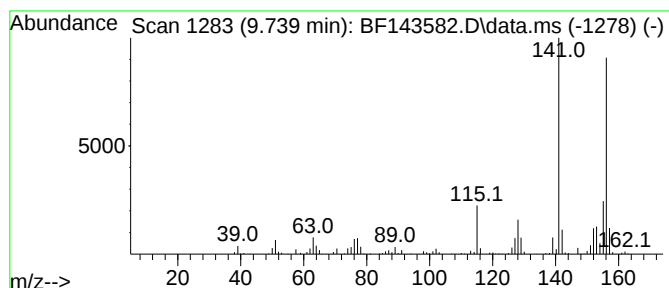
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Naphthalene, 1,5-dimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.739	8.52 ng	555003	Acenaphthene-d10	9.963

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	97
2			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
3			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	96
4			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96
5			Naphthalene, 1,8-dimethyl-	156	C12H12	000569-41-5	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082225\
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Acq On : 22 Aug 2025 19:50
Operator : RC/JU
Sample : Q2924-05
Misc :
ALS Vial : 21 Sample Multiplier: 1

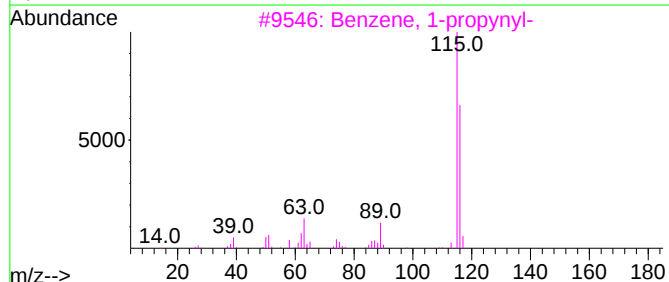
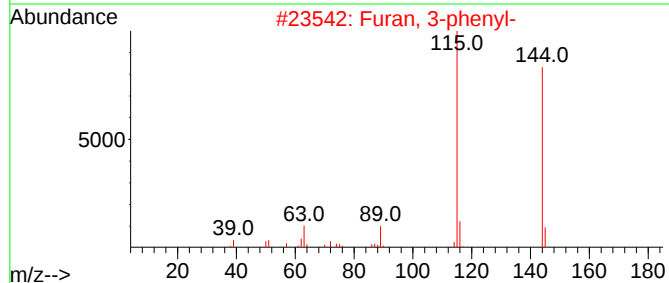
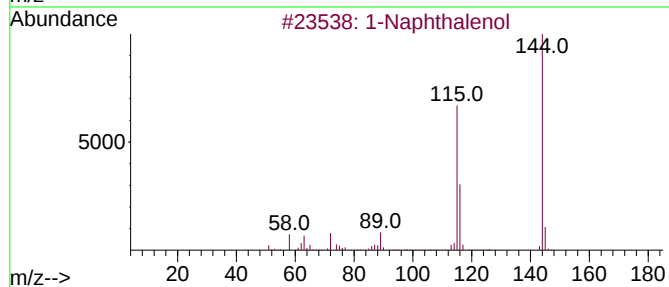
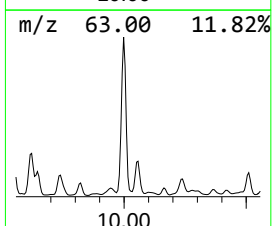
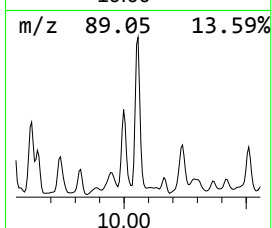
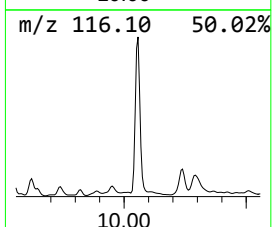
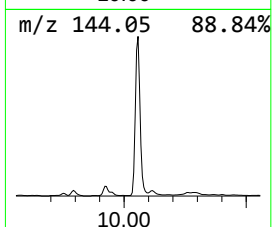
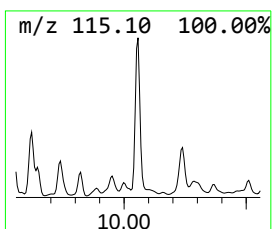
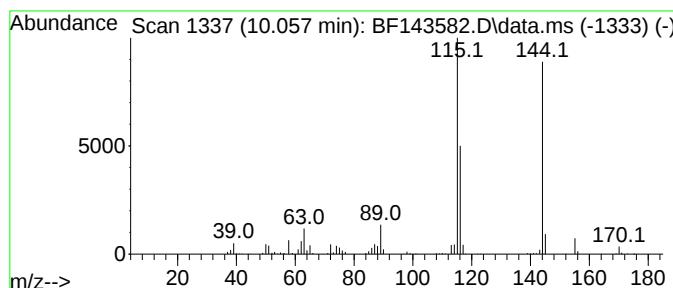
Instrument :
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 1-Naphthalenol Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD			R.T.
10.057	6.82 ng	444539	Acenaphthene-d10			9.963
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	1-Naphthalenol		144	C10H8O	000090-15-3	97
2	Furan, 3-phenyl-		144	C10H8O	013679-41-9	93
3	Benzene, 1-propynyl-		116	C9H8	000673-32-5	91
4	1-Propyne, 3-phenyl-		116	C9H8	010147-11-2	90
5	2-Naphthalenol		144	C10H8O	000135-19-3	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082225\
Data File : BF143582.D
Acq On : 22 Aug 2025 19:50
Operator : RC/JU
Sample : Q2924-05
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
MW-2B-082025

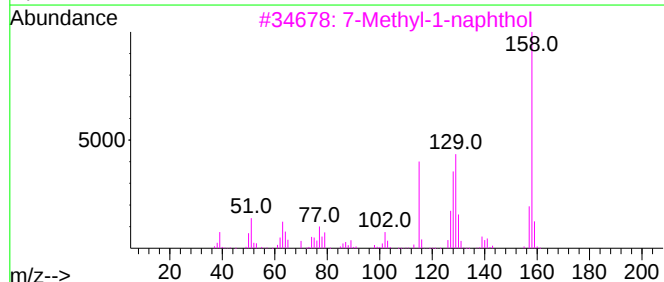
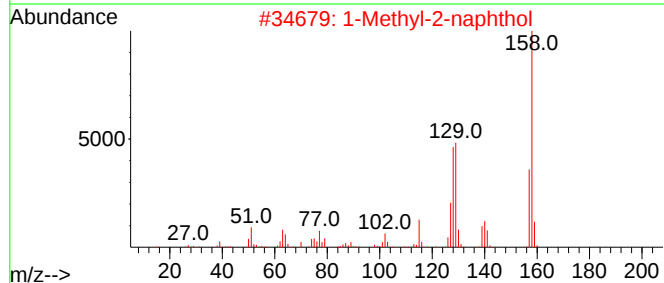
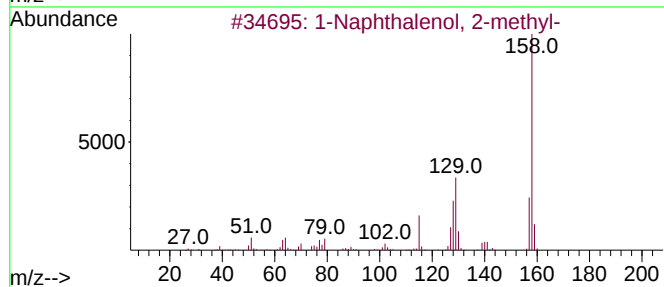
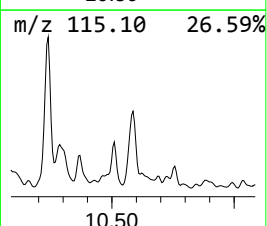
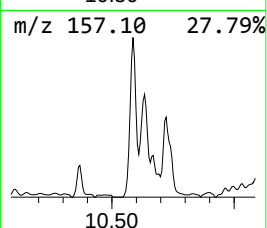
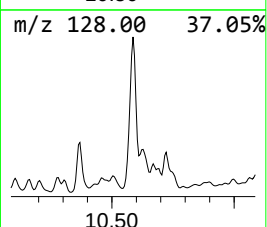
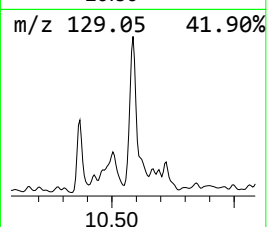
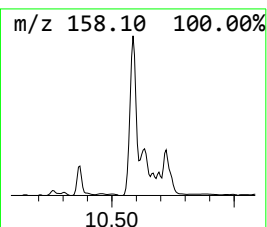
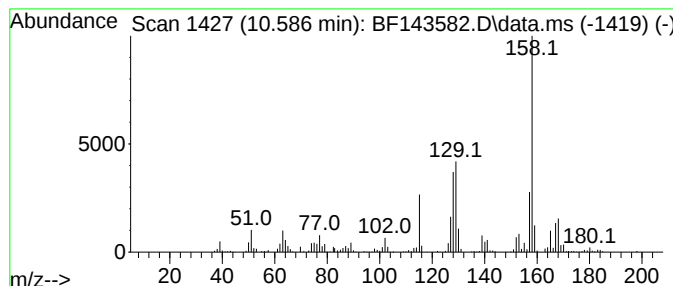
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 1-Naphthalenol, 2-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.586	8.49 ng	553162	Acenaphthene-d10	9.963

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Naphthalenol, 2-methyl-	158	C11H10O	007469-77-4	94
2		1-Methyl-2-naphthol	158	C11H10O	001076-26-2	90
3		7-Methyl-1-naphthol	158	C11H10O	006939-33-9	89
4		Cinnoline, 3,4-dimethyl-	158	C10H10N2	003929-83-7	68
5		1-Naphthalenol, 4-methyl-	158	C11H10O	010240-08-1	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082225\
Data File : BF143582.D
Acq On : 22 Aug 2025 19:50
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Sample : Q2924-05
Misc :
ALS Vial : 21 Sample Multiplier: 1

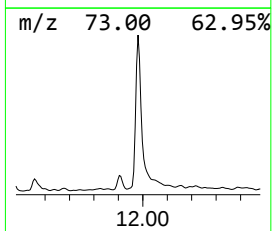
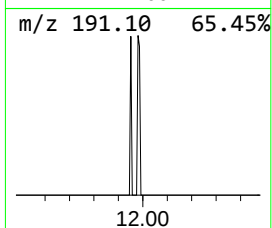
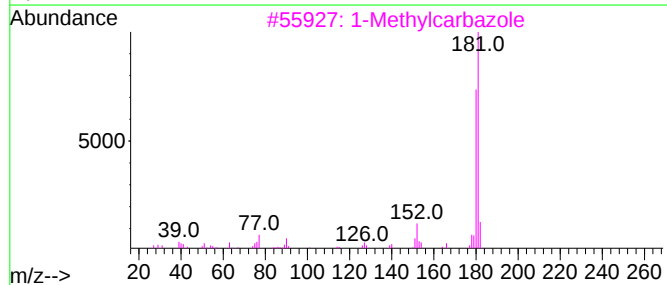
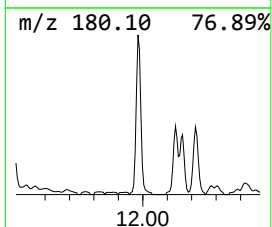
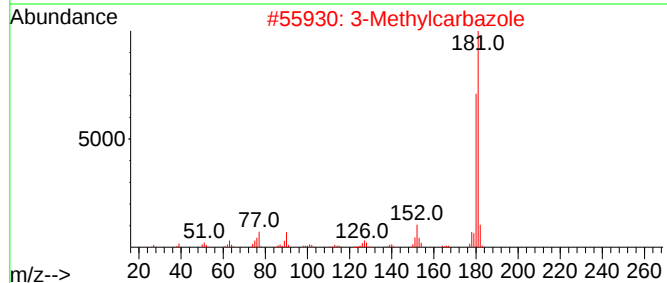
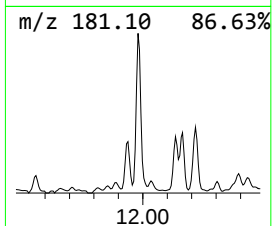
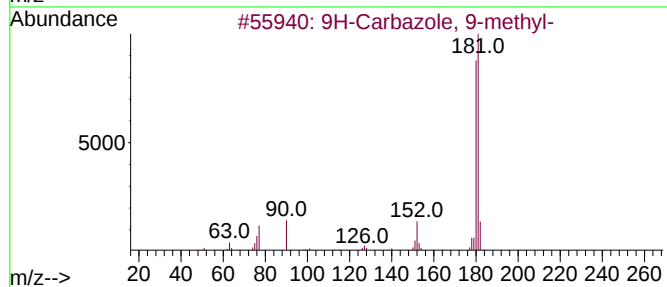
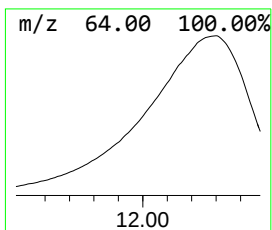
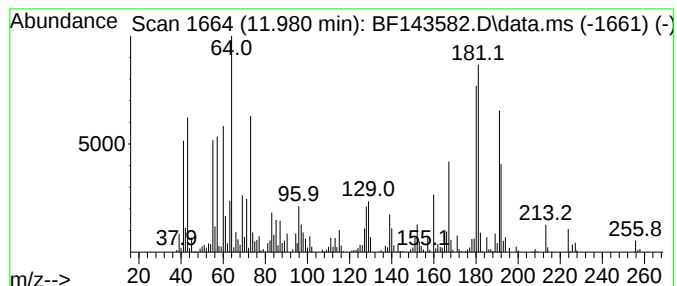
Instrument :
BNA_F
ClientSampleId :
MW-2B-082025

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 19 9H-Carbazole, 9-methyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.980	7.40 ng	339194	Phenanthrene-d10		11.451	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	9H-Carbazole, 9-methyl-		181	C13H11N	001484-12-4	56
2	3-Methylcarbazole		181	C13H11N	004630-20-0	56
3	1-Methylcarbazole		181	C13H11N	006510-65-2	44
4	Hexathiane		192	S6	013798-23-7	38
5	Cyclic octaatomic sulfur		256	S8	010544-50-0	32



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082225\
Data File : BF143582.D
Acq On : 22 Aug 2025 19:50
Operator : RC/JU
Sample : Q2924-05
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-2B-082025

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF082025.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
Butane, 2-metho...	2.252	104.2	ng	3045920	1	6.928	584600	20.0
1,3-Cyclopentad...	5.440	279.0	ng	8154400	1	6.928	584600	20.0
Benzene, 1,3-di...	5.787	27.5	ng	802606	1	6.928	584600	20.0
unknown6.387	6.387	13.4	ng	392468	1	6.928	584600	20.0
unknown6.751	6.751	38.9	ng	1136130	1	6.928	584600	20.0
unknown6.987	6.987	73.5	ng	2147060	1	6.928	584600	20.0
unknown7.034	7.034	13.6	ng	396790	1	6.928	584600	20.0
unknown7.134	7.134	740.1	ng	21634100	1	6.928	584600	20.0
unknown7.193	7.193	33.9	ng	992119	1	6.928	584600	20.0
unknown7.951	7.951	2.2	ng	3463270	2	8.228	31585200	20.0
Naphthalene, 1-...	9.475	8.1	ng	525538	3	9.963	1302770	20.0
Naphthalene, 1,...	9.545	14.9	ng	974129	3	9.963	1302770	20.0
Naphthalene, 2,...	9.622	18.0	ng	1169610	3	9.963	1302770	20.0
Naphthalene, 2,...	9.645	5.7	ng	370153	3	9.963	1302770	20.0
Naphthalene, 1,...	9.739	8.5	ng	555003	3	9.963	1302770	20.0
1-Naphthalenol	10.057	6.8	ng	444539	3	9.963	1302770	20.0
1-Naphthalenol,...	10.586	8.5	ng	553162	3	9.963	1302770	20.0
9H-Carbazole, 9...	11.980	7.4	ng	339194	4	11.451	916636	20.0