

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082225\
 Data File : BF143572.D
 Acq On : 22 Aug 2025 14:56
 Operator : RC/JU
 Sample : Q2903-07
 Misc :
 ALS Vial : 11 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: BF143572.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.257	3	11	24	rVB	2021337	3337920	44.12%	8.606%
2	5.157	495	504	512	rBV2	106263	181546	2.40%	0.468%
3	5.563	563	573	597	rBV	1201318	1648218	21.79%	4.249%
4	6.104	660	665	670	rBV	192036	246942	3.26%	0.637%
5	6.557	733	742	748	rBV	862167	1159667	15.33%	2.990%
6	6.610	748	751	764	rVB	183489	244377	3.23%	0.630%
7	6.922	799	804	810	rBV	504220	634651	8.39%	1.636%
8	6.981	810	814	824	rVB	114126	163943	2.17%	0.423%
9	7.116	829	837	842	rBV	5045509	7565055	100.00%	19.504%
10	7.487	888	900	910	rBV2	1557989	2333853	30.85%	6.017%
11	7.875	960	966	972	rVB	83435	110860	1.47%	0.286%
12	7.945	972	978	982	rBV2	282665	400923	5.30%	1.034%
13	7.992	982	986	991	rVV	210310	362425	4.79%	0.934%
14	8.228	1014	1026	1031	rBV2	2396219	4126826	54.55%	10.639%
15	8.269	1031	1033	1045	rVB2	66732	92000	1.22%	0.237%
16	8.463	1061	1066	1078	rVB2	82016	150322	1.99%	0.388%
17	8.792	1114	1122	1127	rBV2	131216	179344	2.37%	0.462%
18	9.016	1154	1160	1164	rBV	300762	391198	5.17%	1.009%
19	9.092	1169	1173	1180	rVB	95815	147341	1.95%	0.380%
20	9.281	1199	1205	1210	rBV	2736639	3545090	46.86%	9.140%
21	9.375	1216	1221	1228	rBV4	44605	82003	1.08%	0.211%
22	9.963	1315	1321	1325	rBV	846772	1100338	14.55%	2.837%
23	10.233	1361	1367	1374	rBV	142188	239384	3.16%	0.617%
24	10.416	1393	1398	1404	rVB	168830	233695	3.09%	0.602%
25	10.675	1437	1442	1447	rVB	151187	199794	2.64%	0.515%
26	10.757	1449	1456	1466	rBV2	2085264	3044597	40.25%	7.849%
27	10.910	1479	1482	1486	rBV	81288	98547	1.30%	0.254%
28	11.451	1569	1574	1581	rVB	874584	1119499	14.80%	2.886%
29	11.974	1658	1663	1675	rBV3	173979	310526	4.10%	0.801%
30	12.504	1748	1753	1767	rVB5	37758	107744	1.42%	0.278%
31	12.851	1807	1812	1816	rVV	75329	105936	1.40%	0.273%
32	13.033	1837	1843	1849	rBV	2570163	3588838	47.44%	9.252%
33	13.257	1876	1881	1897	rVB	48665	106024	1.40%	0.273%
34	13.939	1992	1997	2013	rBV3	55883	136859	1.81%	0.353%
35	14.092	2019	2023	2031	rVB	518990	688070	9.10%	1.774%
36	15.592	2272	2278	2292	rVB	378940	603725	7.98%	1.556%

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Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

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

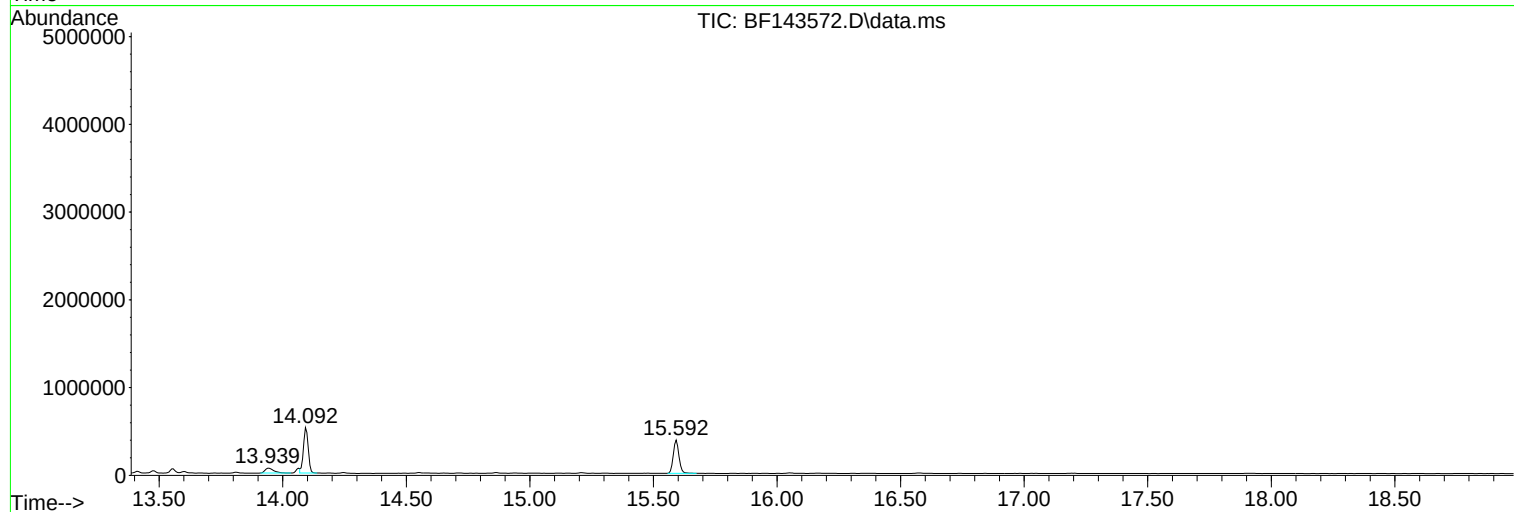
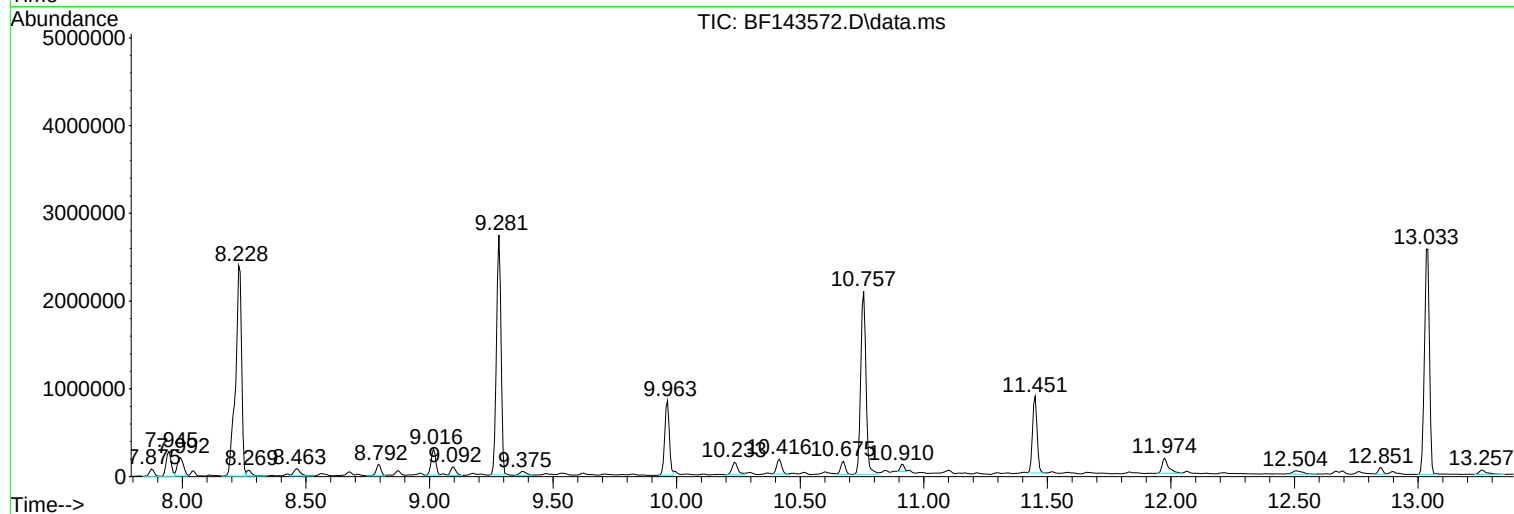
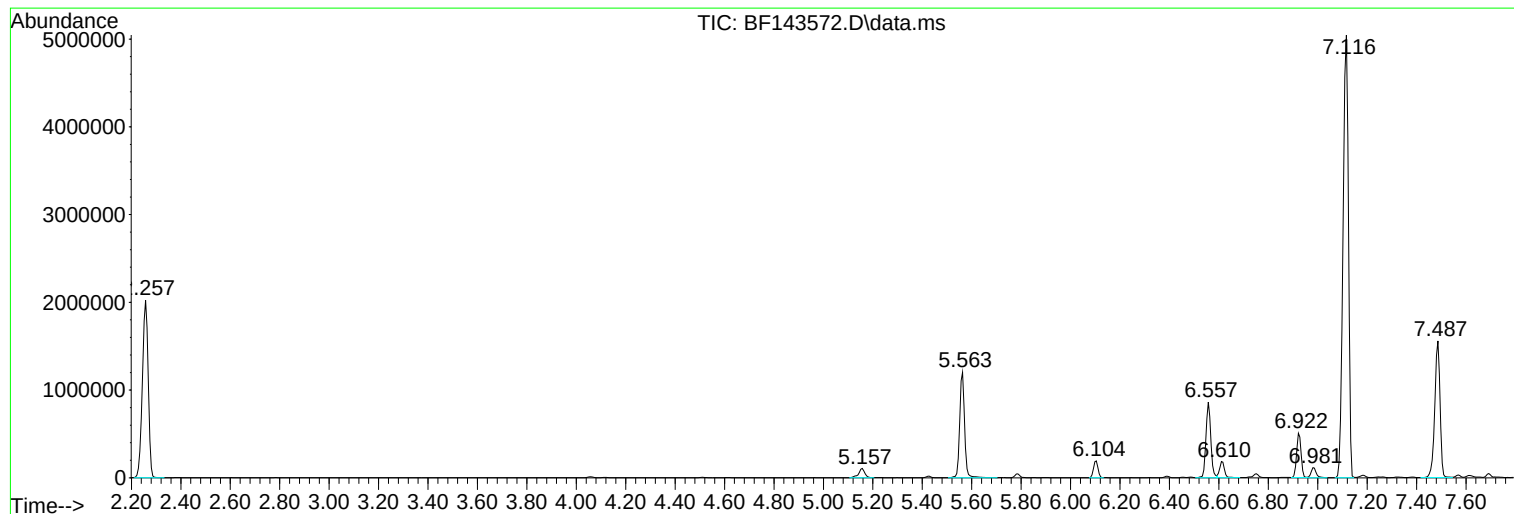
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF082025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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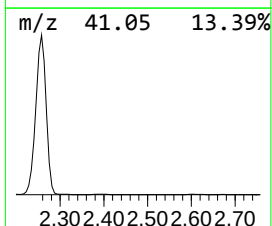
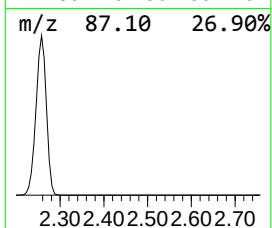
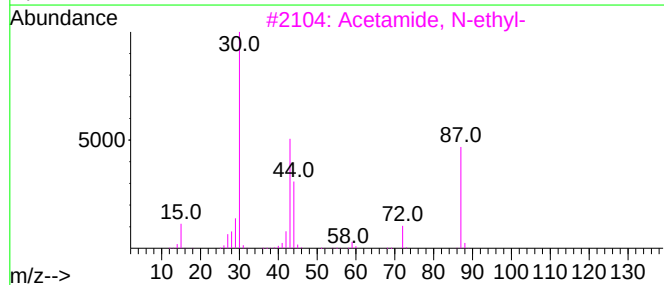
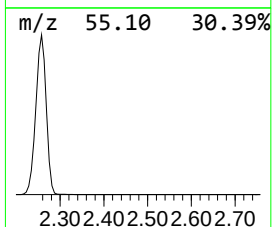
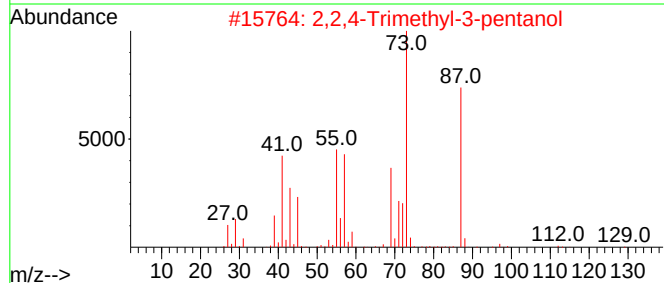
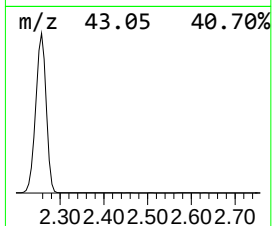
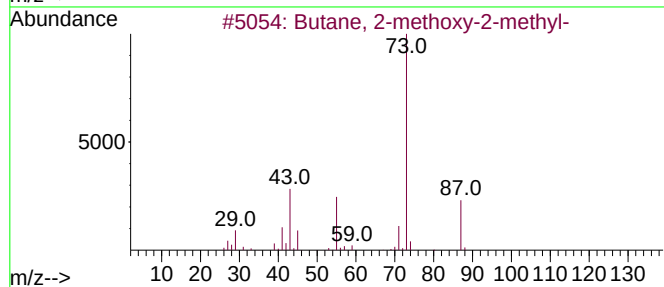
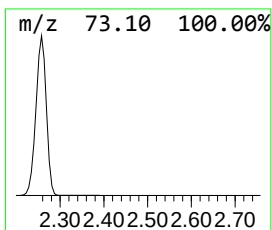
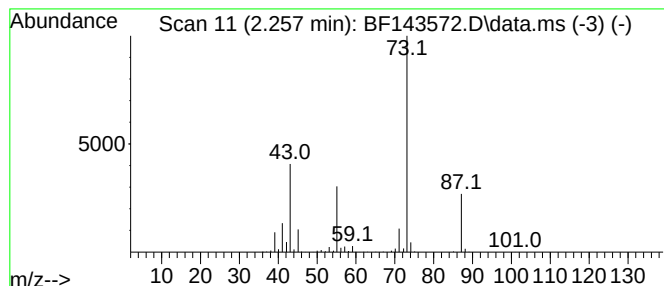
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.
2.257	105.19 ng	3337920	1,4-Dichlorobenzene-d4	6.922

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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Instrument :
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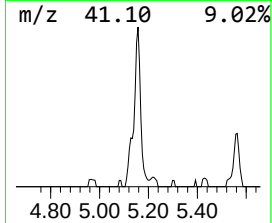
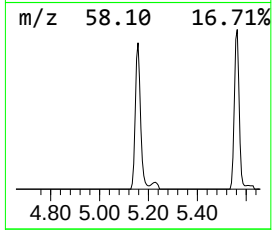
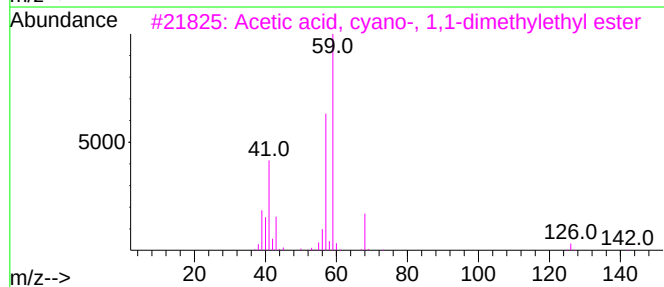
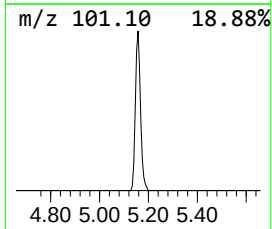
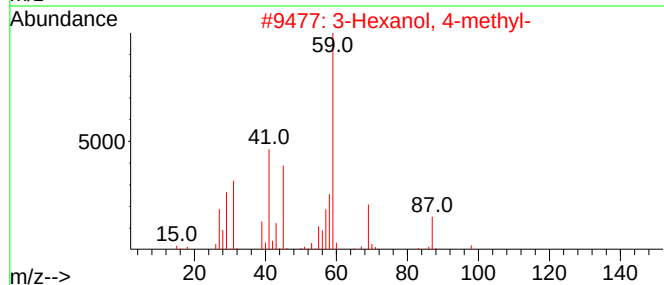
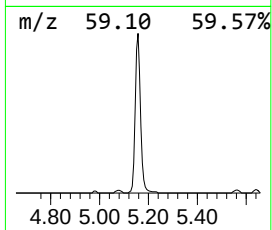
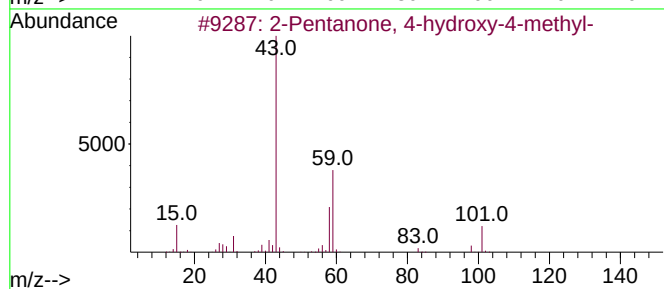
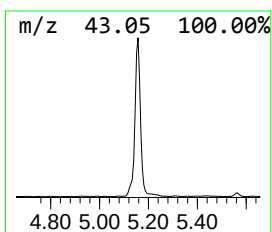
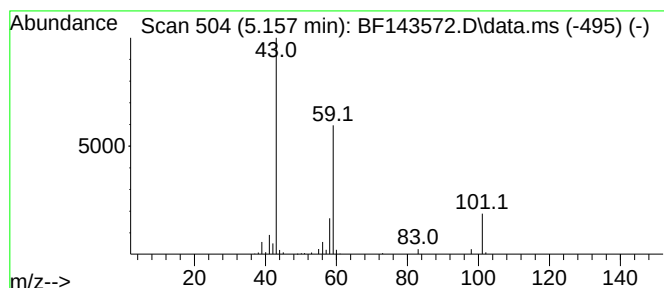
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF082025.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.157	5.72 ng	181546	1,4-Dichlorobenzene-d4	6.922

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
3		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
4		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
5		Acetone	58	C3H6O	000067-64-1	12



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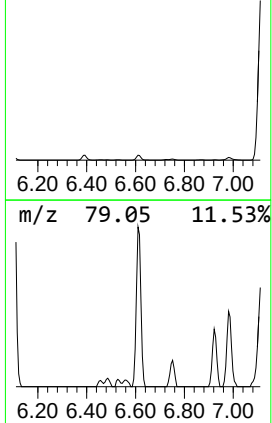
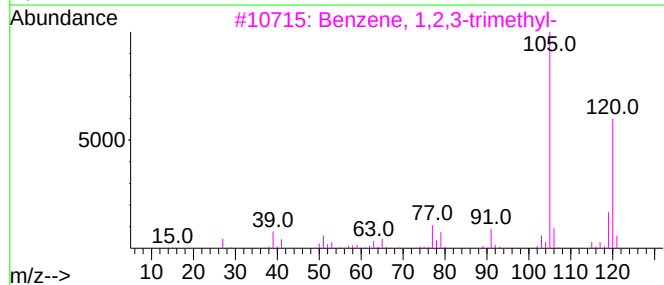
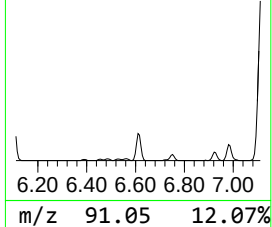
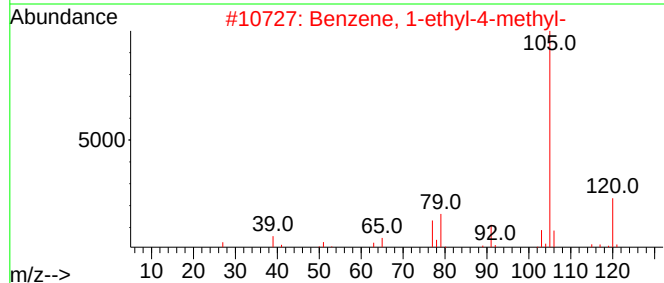
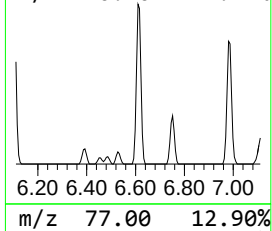
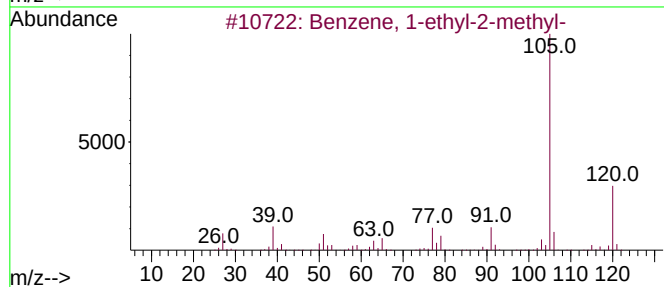
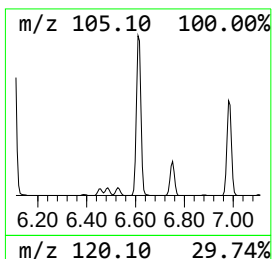
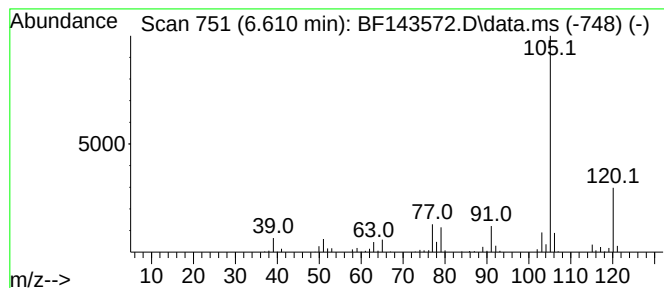
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF082025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.610	7.70 ng	244377	1,4-Dichlorobenzene-d4	6.922

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			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
4			Mesitylene	120	C9H12	000108-67-8	91
5			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91



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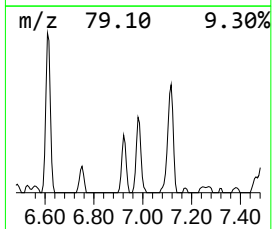
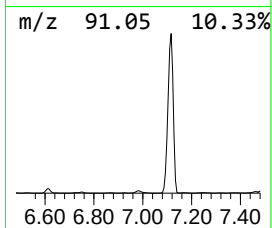
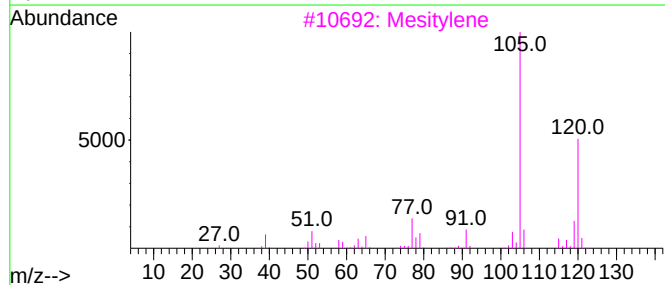
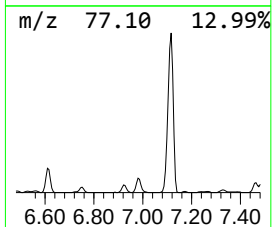
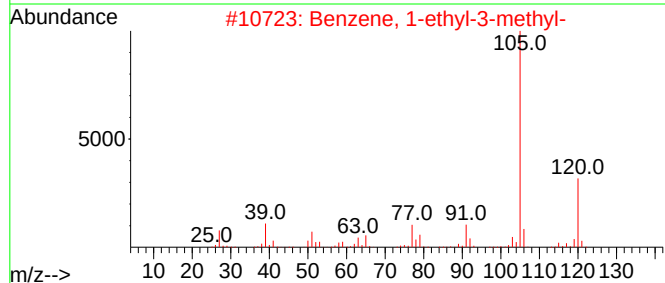
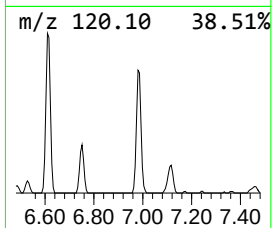
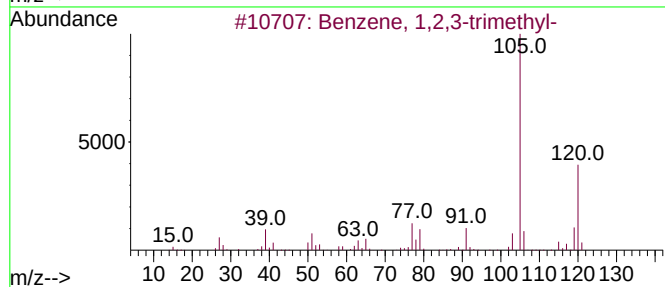
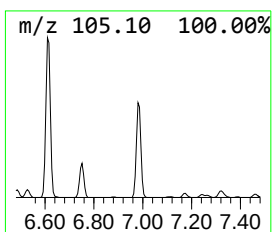
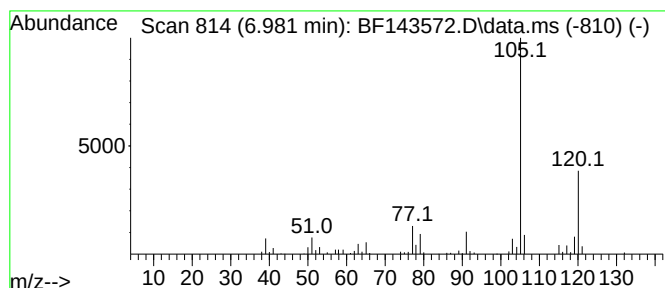
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 unknown6.981 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.981	5.17 ng	163943	1,4-Dichlorobenzene-d4	6.922

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
3			Mesitylene	120	C9H12	000108-67-8	93
4			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91



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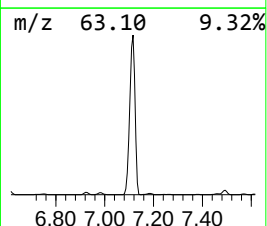
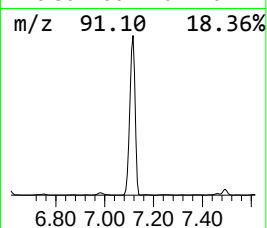
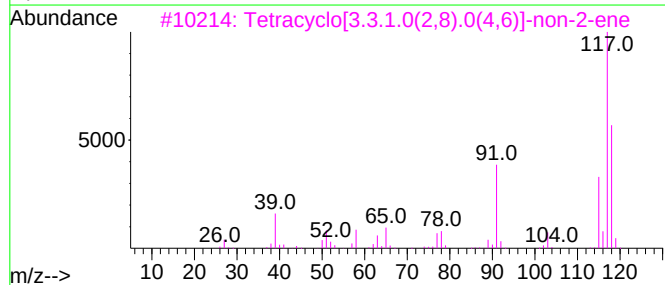
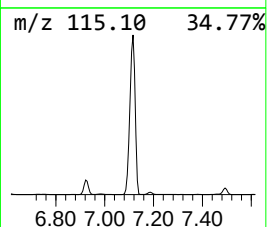
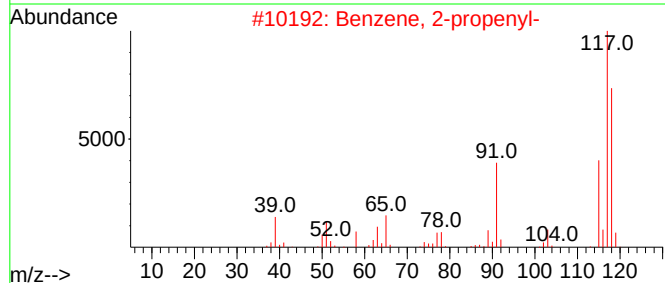
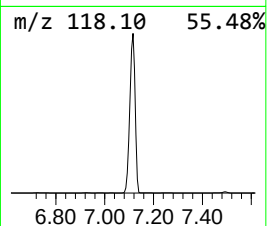
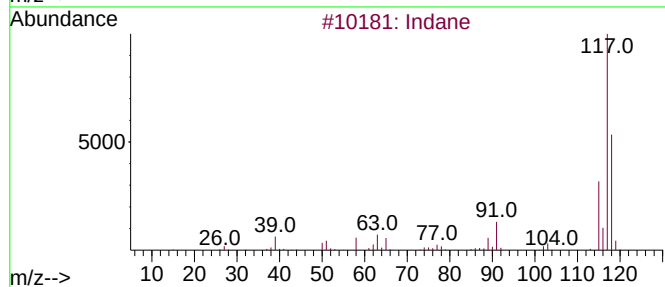
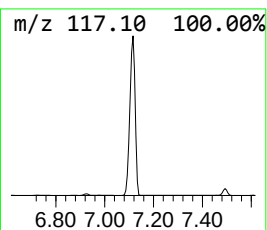
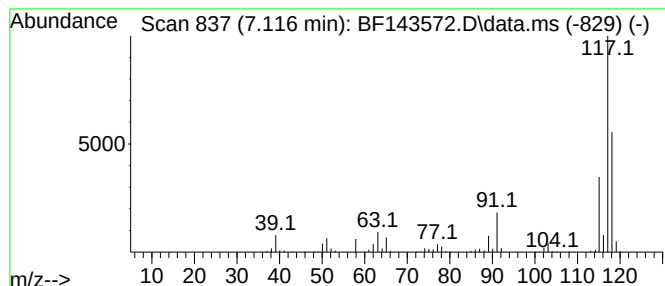
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzene, 2-propenyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.116	238.40 ng	7565060	1,4-Dichlorobenzene-d4	6.922

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	94
2			Benzene, 2-propenyl-	118	C9H10	000300-57-2	81
3			Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	80
4			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	64
5			Benzene, (isocyanomethyl)-	117	C8H7N	010340-91-7	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082225\
Data File : BF143572.D
Acq On : 22 Aug 2025 14:56
Operator : RC/JU
Sample : Q2903-07
Misc :
ALS Vial : 11 Sample Multiplier: 1

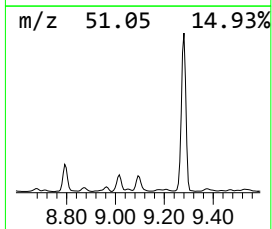
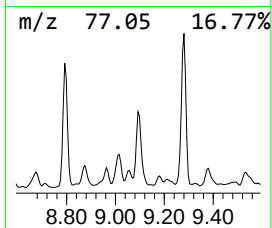
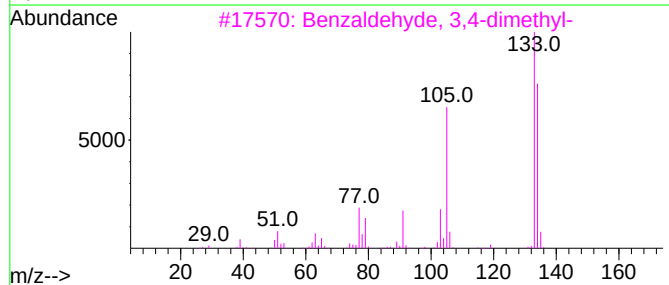
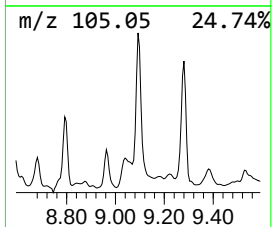
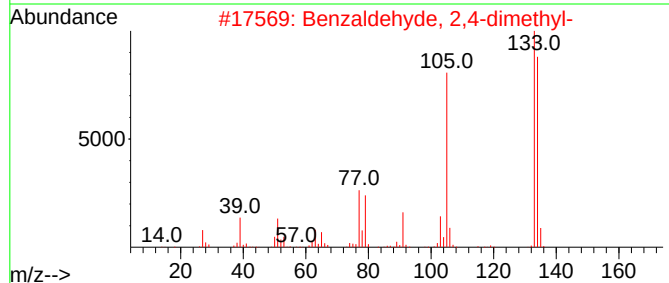
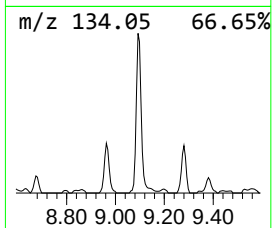
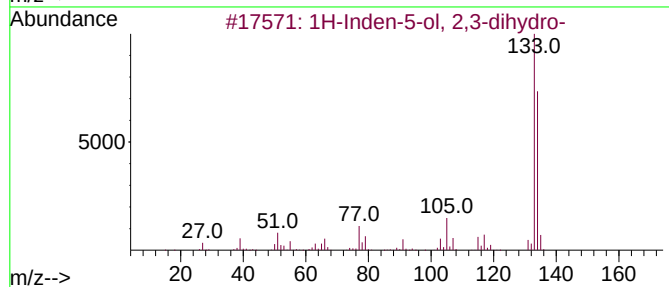
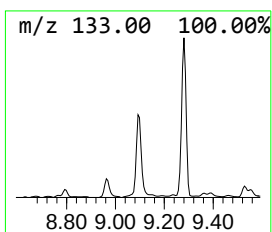
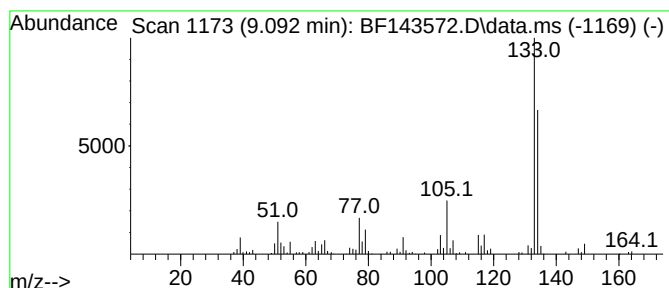
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ClientSampleId :
MW-12B-081925

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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 7 1H-Inden-5-ol, 2,3-dihydro- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD		R.T.	
9.092	2.68 ng	147341	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	87
2	Benzaldehyde, 2,4-dimethyl-		134	C9H10O	015764-16-6	74
3	Benzaldehyde, 3,4-dimethyl-		134	C9H10O	005973-71-7	72
4	Benzaldehyde, 4-ethyl-		134	C9H10O	004748-78-1	72
5	1H-Inden-1-ol, 2,3-dihydro-		134	C9H10O	006351-10-6	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082225\
Data File : BF143572.D
Acq On : 22 Aug 2025 14:56
Operator : RC/JU
Sample : Q2903-07
Misc :
ALS Vial : 11 Sample Multiplier: 1

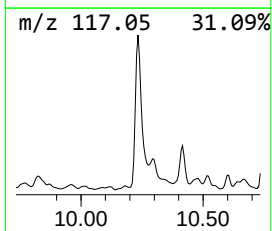
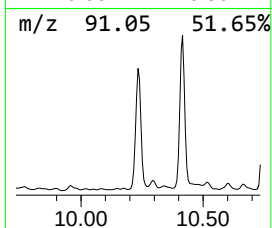
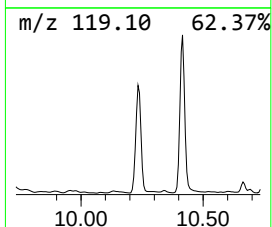
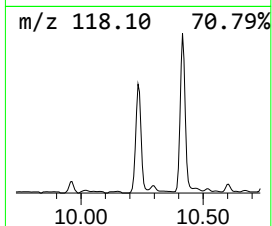
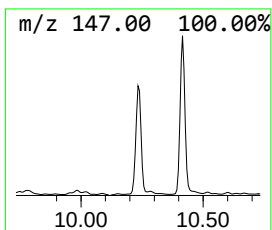
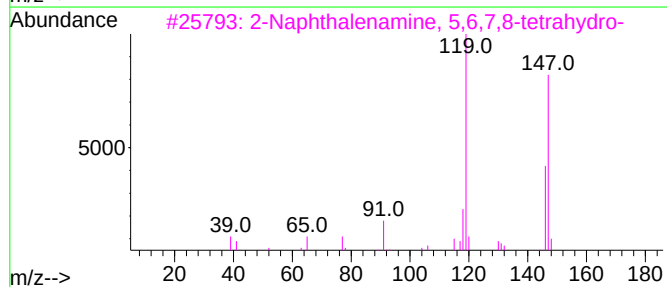
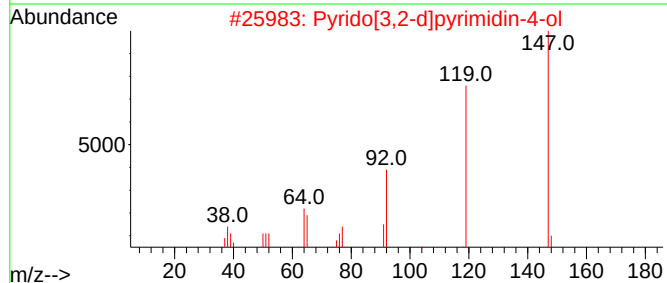
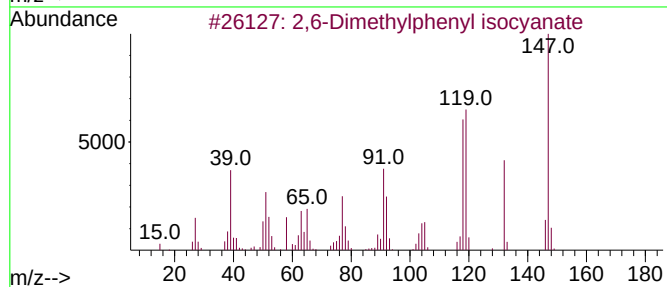
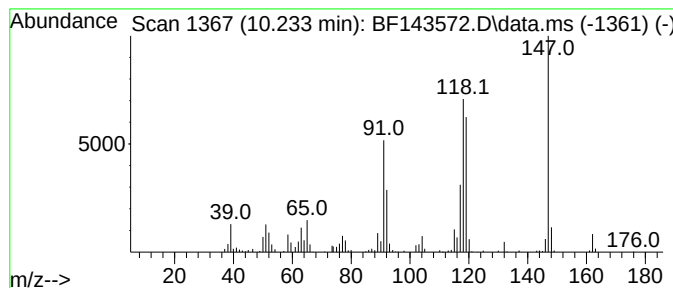
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF082025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 2,6-Dimethylphenyl isocyanate Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD		R.T.	
10.233	4.35 ng	239384	Acenaphthene-d10		9.963	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	2,6-Dimethylphenyl isocyanate		147	C9H9NO	028556-81-2	83
2	Pyrido[3,2-d]pyrimidin-4-ol		147	C7H5N3O	037538-67-3	70
3	2-Naphthalenamine, 5,6,7,8-tetra...		147	C10H13N	002217-43-8	64
4	1H-Inden-1-one, 5-amino-2,3-dihy...		147	C9H9NO	003470-54-0	53
5	1H-Indole-1-carboxaldehyde, 2,3-...		147	C9H9NO	002861-59-8	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082225\
Data File : BF143572.D
Acq On : 22 Aug 2025 14:56
Operator : RC/JU
Sample : Q2903-07
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
MW-12B-081925

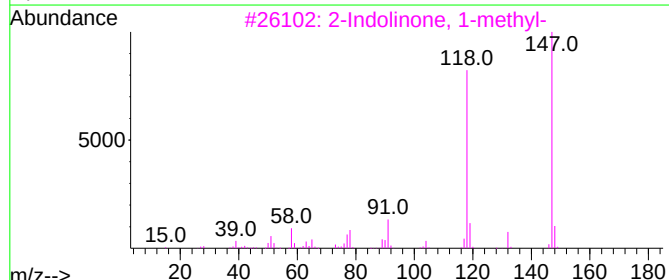
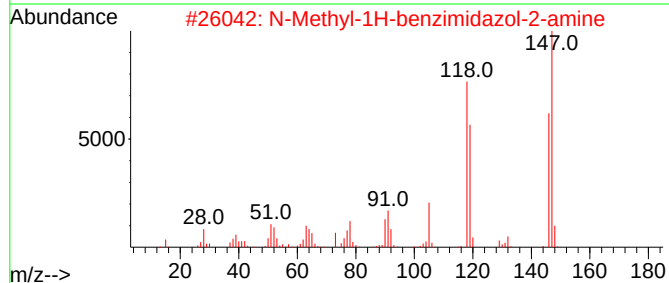
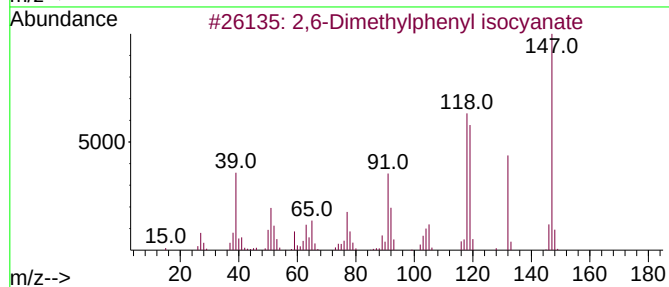
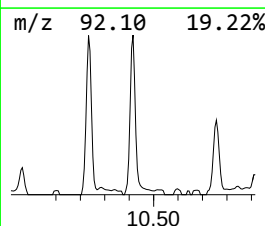
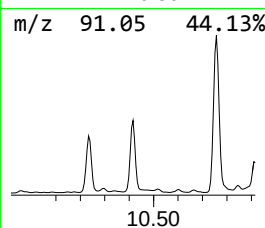
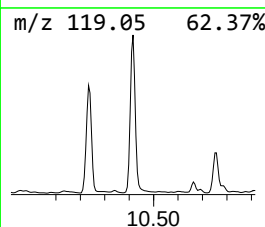
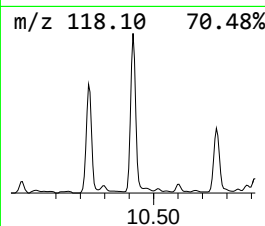
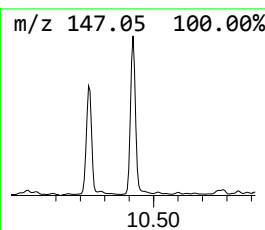
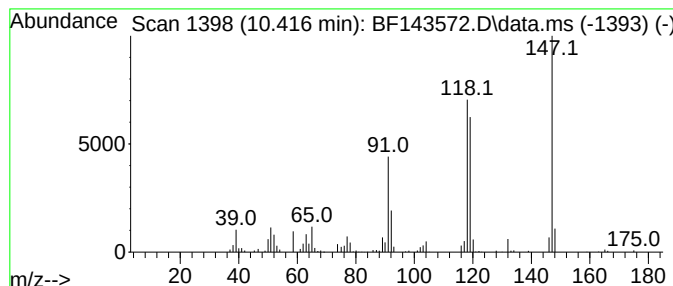
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.416	4.25 ng	233695	Acenaphthene-d10	9.963

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,6-Dimethylphenyl isocyanate	147	C9H9NO	028556-81-2	86
2			N-Methyl-1H-benzimidazol-2-amine	147	C8H9N3	017228-38-5	80
3			2-Indolinone, 1-methyl-	147	C9H9NO	000061-70-1	76
4			1H-Indole, 2,3-dihydro-	119	C8H9N	000496-15-1	38
5			Pyridine, 5-ethenyl-2-methyl-	119	C8H9N	000140-76-1	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082225\
Data File : BF143572.D
Acq On : 22 Aug 2025 14:56
Operator : RC/JU
Sample : Q2903-07
Misc :
ALS Vial : 11 Sample Multiplier: 1

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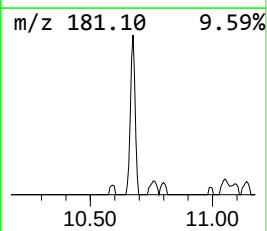
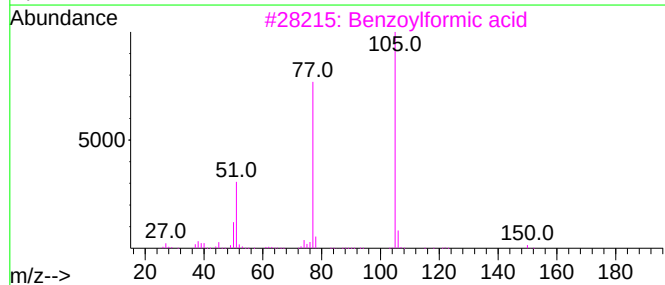
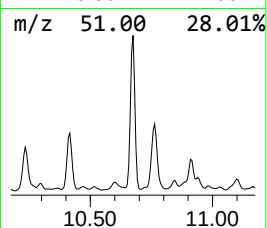
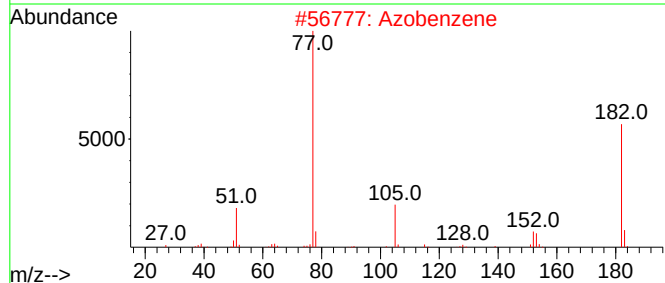
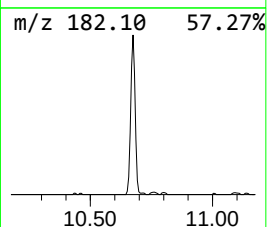
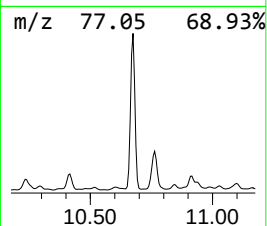
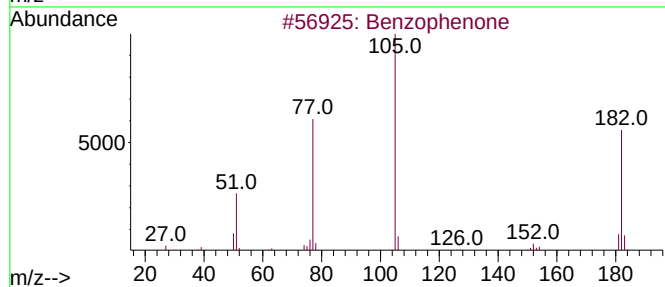
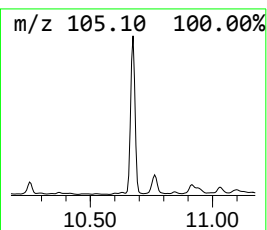
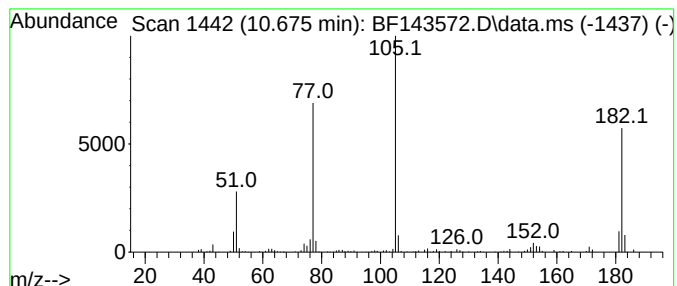
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Benzophenone Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.675	3.63 ng	199794	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	59
3			Benzoylformic acid	150	C8H6O3	000611-73-4	49
4			Acetophenone, 2-chloro-	154	C8H7ClO	000532-27-4	46
5			2-Chloro-1-phenyl-1-propanone	168	C9H9ClO	006084-17-9	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082225\
Data File : BF143572.D
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Operator : RC/JU
Sample : Q2903-07
Misc :
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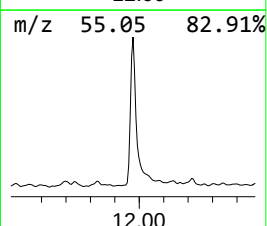
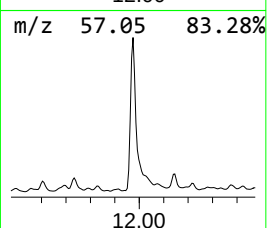
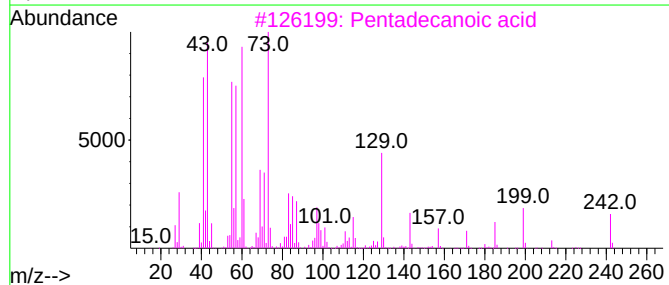
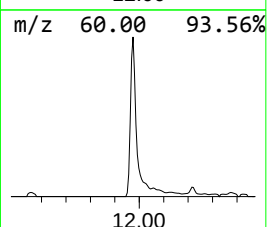
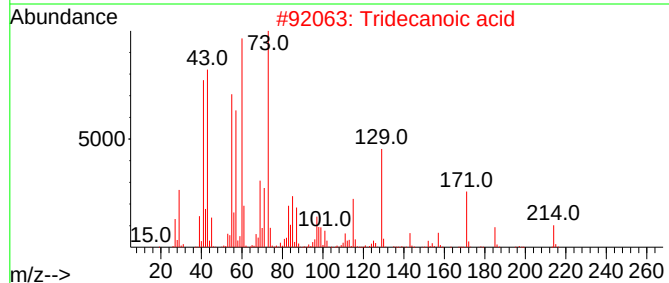
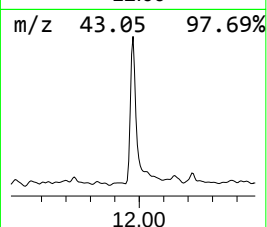
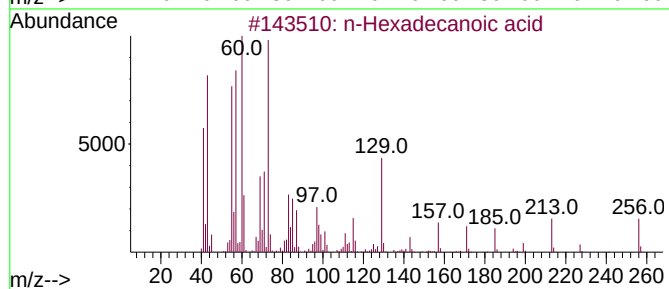
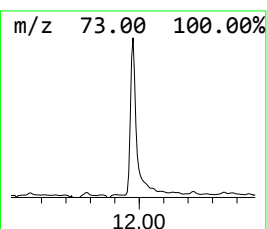
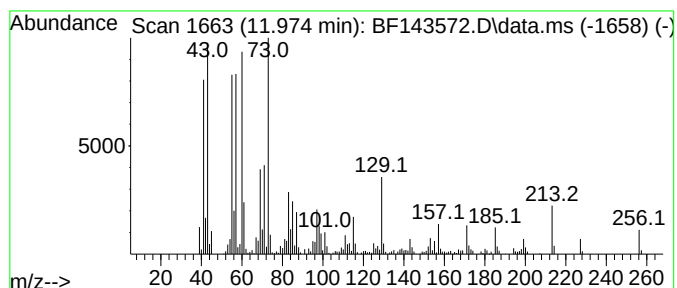
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.975	5.55 ng	310526	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	93
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	87
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	83
5	Undecanoic acid	186	C11H22O2	000112-37-8	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082225\
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Acq On : 22 Aug 2025 14:56
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Sample : Q2903-07
Misc :
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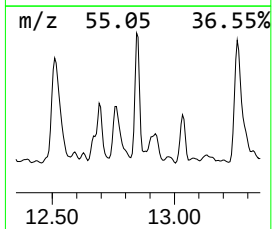
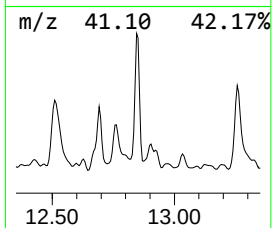
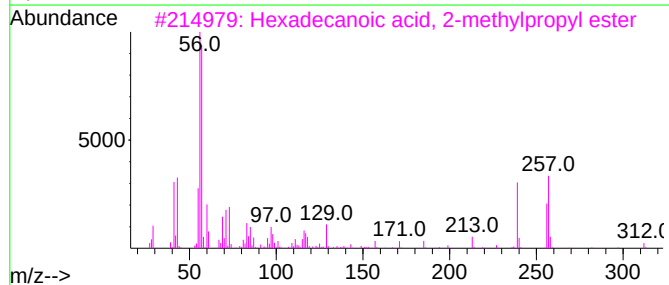
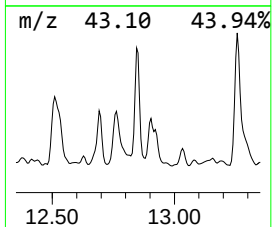
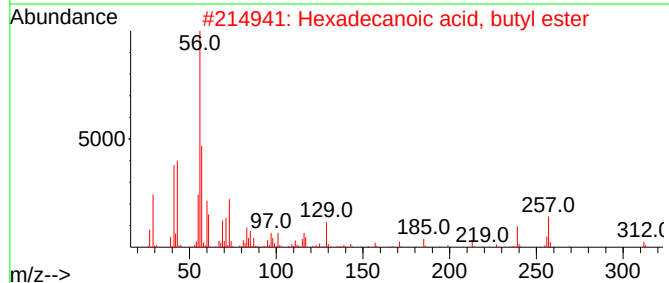
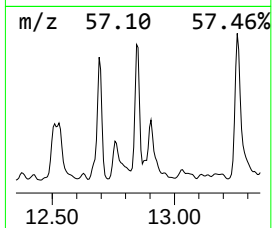
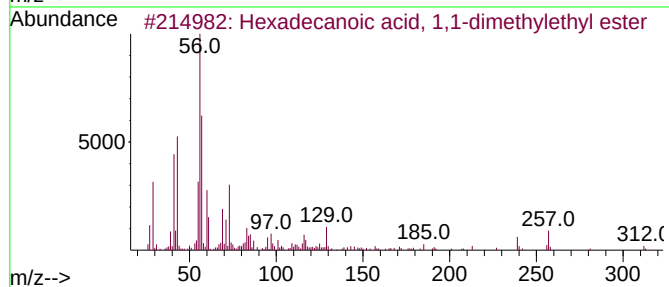
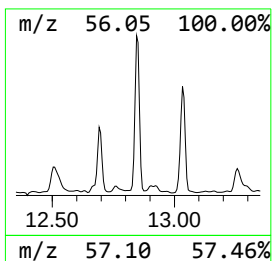
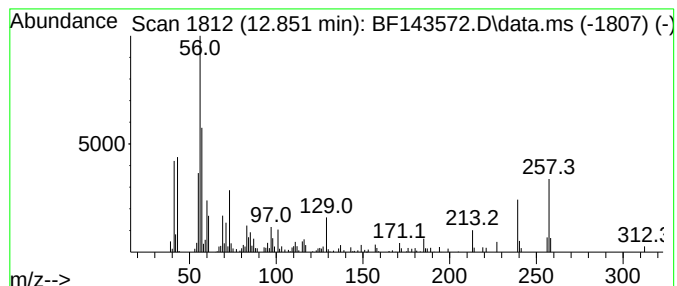
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Hexadecanoic acid, 1,1-dime... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.851	3.08 ng	105936	Chrysene-d12	14.092	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecanoic acid, 1,1-dimethyle...	312	C20H40O2	031158-91-5	93
2	Hexadecanoic acid, butyl ester	312	C20H40O2	000111-06-8	93
3	Hexadecanoic acid, 2-methylpropy...	312	C20H40O2	000110-34-9	70
4	1-Heptene, 2-methyl-	112	C8H16	015870-10-7	30
5	1-Hexene, 2-methyl-	98	C7H14	006094-02-6	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082225\
Data File : BF143572.D
Acq On : 22 Aug 2025 14:56
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ClientSampleId :
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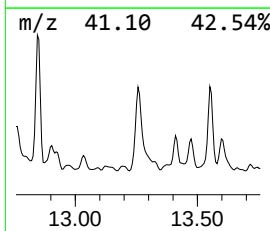
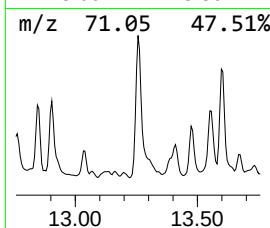
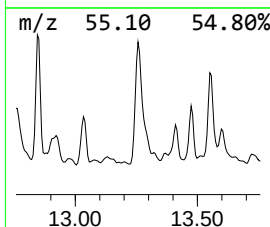
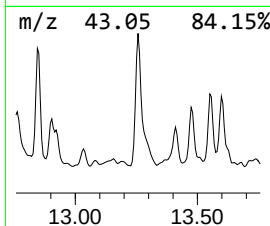
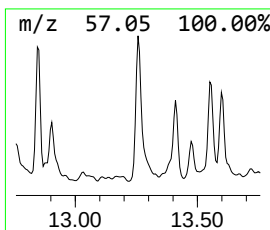
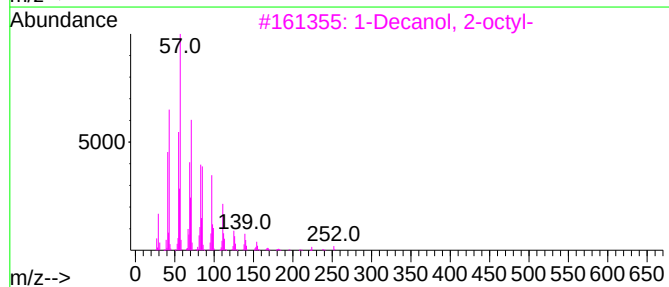
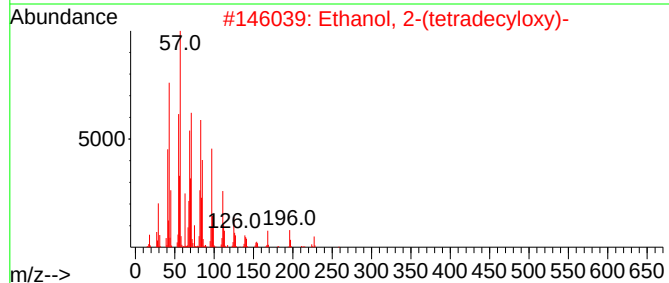
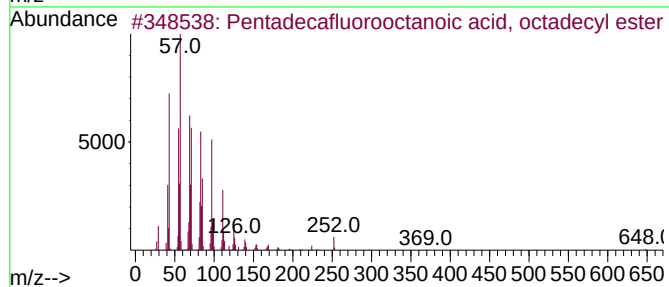
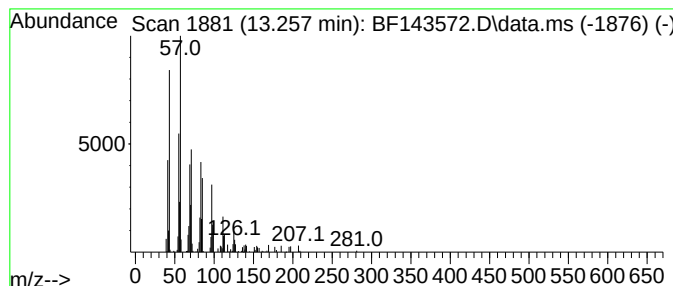
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Pentadecafluorooctanoic aci... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.257	3.08 ng	106024	Chrysene-d12	14.092

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	91
2			Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	91
3			1-Decanol, 2-octyl-	270	C18H38O	045235-48-1	91
4			Oxalic acid, isobutyl hexadecyl ...	370	C22H42O4	1000309-38-1	91
5			Oxalic acid, isobutyl tetradecyl...	342	C20H38O4	1000309-37-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082225\
Data File : BF143572.D
Acq On : 22 Aug 2025 14:56
Operator : RC/JU
Sample : Q2903-07
Misc :
ALS Vial : 11 Sample Multiplier: 1

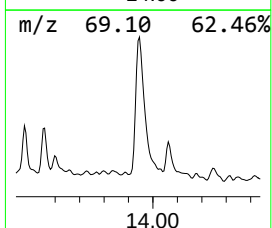
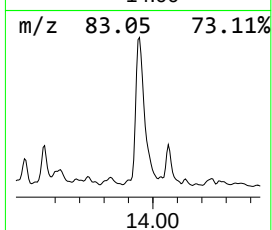
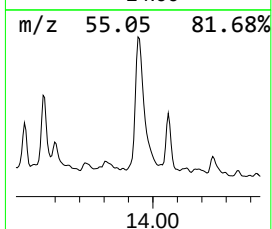
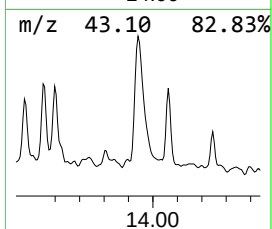
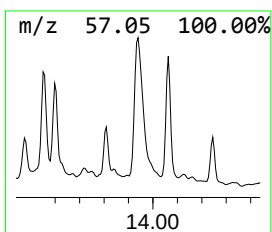
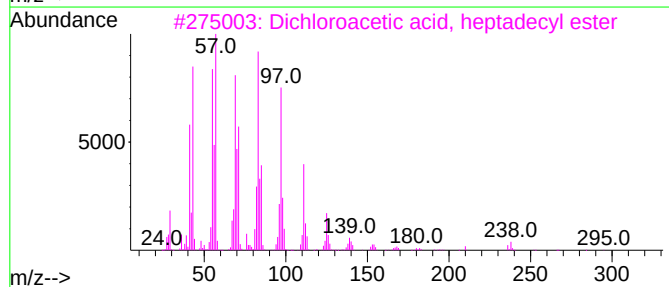
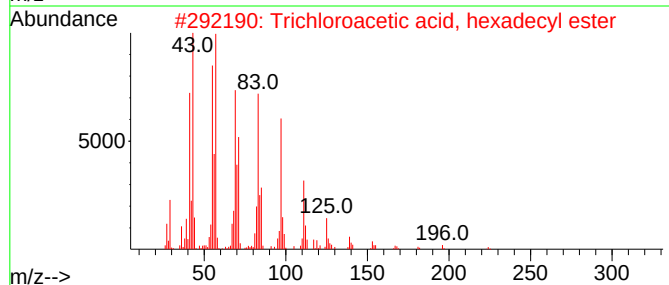
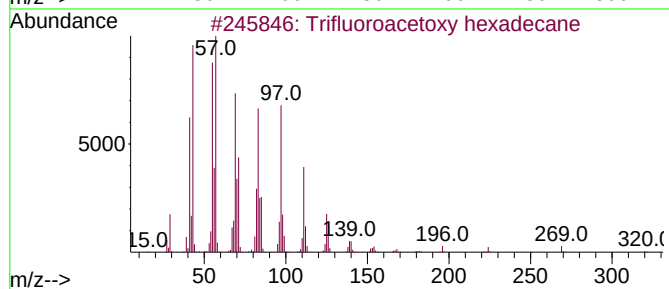
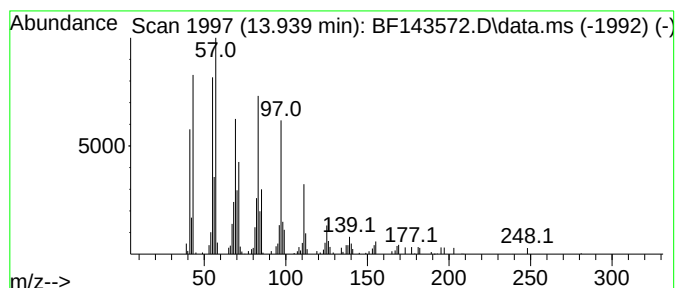
Instrument :
BNA_F
ClientSampleId :
MW-12B-081925

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 14 Trifluoroacetoxy hexadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.939	3.98 ng	136859	Chrysene-d12	14.092	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	95
2	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	95
3	Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	94
4	Heptafluorobutyric acid, hexadec...	438	C20H33F7O2	006385-15-5	94
5	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082225\
Data File : BF143572.D
Acq On : 22 Aug 2025 14:56
Operator : RC/JU
Sample : Q2903-07
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
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\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	2.257	105.2	ng	3337920	1	6.922	634651	20.0
2-Pentanone, 4-...	5.157	5.7	ng	181546	1	6.922	634651	20.0
Benzene, 1-ethy...	6.610	7.7	ng	244377	1	6.922	634651	20.0
unknown6.981	6.981	5.2	ng	163943	1	6.922	634651	20.0
Benzene, 2-prop...	7.116	238.4	ng	7565060	1	6.922	634651	20.0
1H-Inden-5-ol, ...	9.092	2.7	ng	147341	3	9.963	1100340	20.0
2,6-Dimethylphe...	10.233	4.3	ng	239384	3	9.963	1100340	20.0
N-Methyl-1H-ben...	10.416	4.3	ng	233695	3	9.963	1100340	20.0
Benzophenone	10.675	3.6	ng	199794	3	9.963	1100340	20.0
n-Hexadecanoic ...	11.975	5.5	ng	310526	4	11.451	1119500	20.0
Hexadecanoic ac...	12.851	3.1	ng	105936	5	14.092	688070	20.0
Pentadecafluoro...	13.257	3.1	ng	106024	5	14.092	688070	20.0
Trifluoroacetox...	13.939	4.0	ng	136859	5	14.092	688070	20.0