

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF063025\
 Data File : BF142912.D
 Acq On : 30 Jun 2025 12:17
 Operator : RC/JU
 Sample : Q2430-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MH-E/F

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-BF061925.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF142912.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.263	3	12	24	rVB	263880	423059	13.97%	2.184%
2	5.122	485	498	514	rBV	129299	222465	7.35%	1.148%
3	5.516	558	565	570	rBV	1528661	2110716	69.69%	10.895%
4	6.510	727	734	740	rBV	1431527	2132586	70.42%	11.007%
5	6.875	791	796	802	rBV	481728	611569	20.19%	3.157%
6	7.439	885	892	897	rBV	1010890	1407734	46.48%	7.266%
7	8.157	1009	1014	1024	rBV	628295	810245	26.75%	4.182%
8	9.239	1191	1198	1203	rBV	2066922	2789568	92.11%	14.399%
9	9.916	1308	1313	1318	rBV	776178	956600	31.59%	4.938%
10	10.633	1430	1435	1440	rBV	191060	239948	7.92%	1.239%
11	10.710	1442	1448	1453	rBV	1310360	1699766	56.12%	8.773%
12	11.404	1561	1566	1572	rBV	796902	1031763	34.07%	5.326%
13	12.998	1830	1837	1842	rBV	2216332	3028537	100.00%	15.632%
14	13.886	1982	1988	2001	rBV	82463	146385	4.83%	0.756%
15	14.051	2011	2016	2024	rVB	585019	744647	24.59%	3.844%
16	14.804	2139	2144	2154	rBV	185853	290148	9.58%	1.498%
17	15.545	2263	2270	2281	rBV	468689	728264	24.05%	3.759%

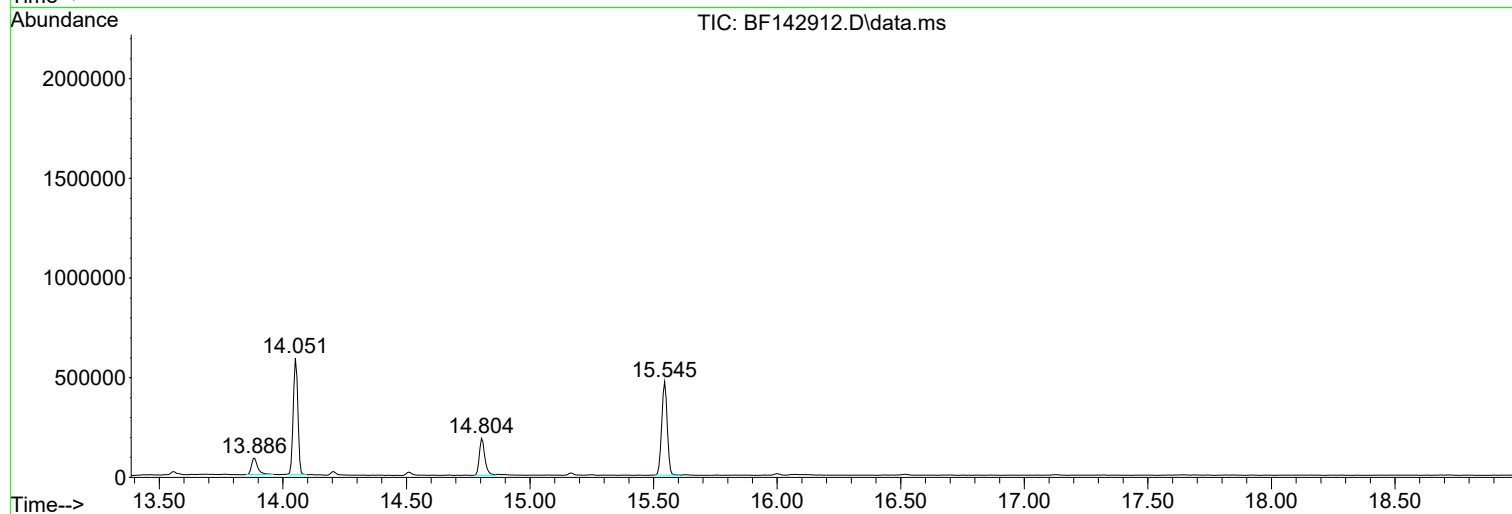
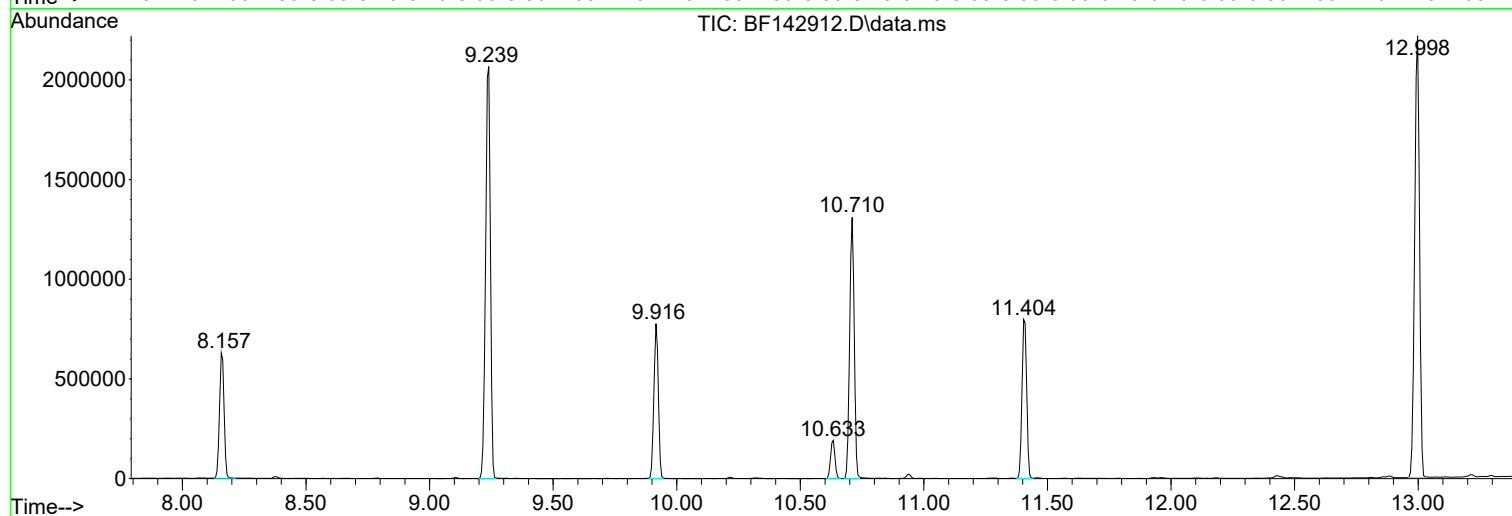
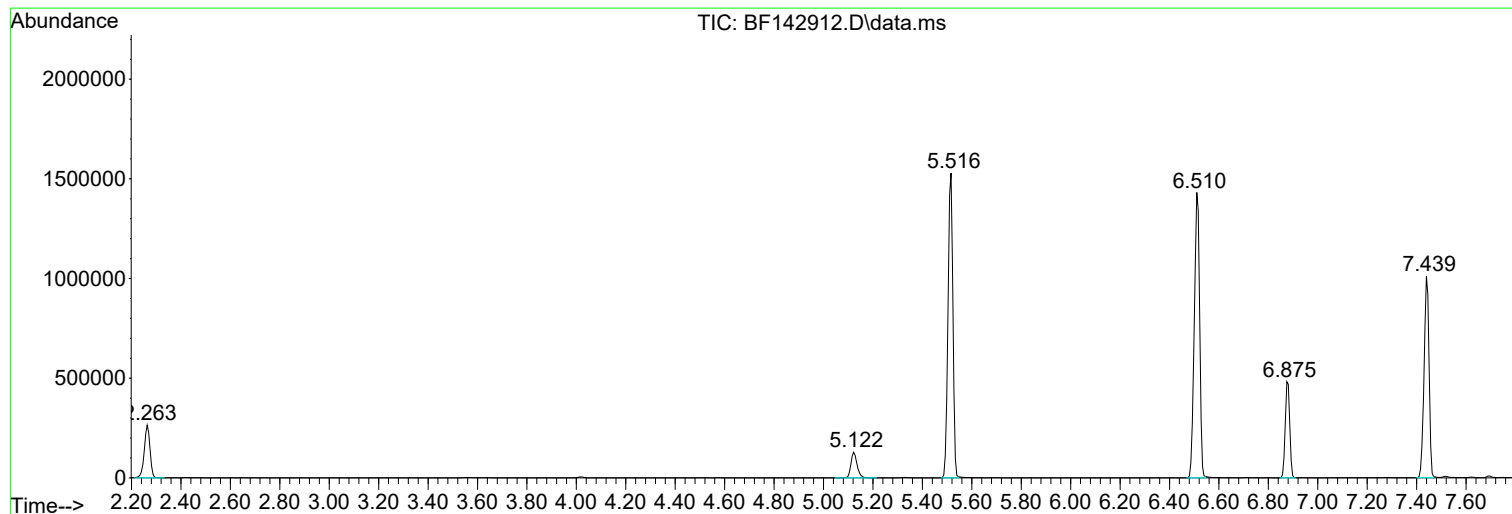
Sum of corrected areas: 19374000

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF063025\
Data File : BF142912.D
Acq On : 30 Jun 2025 12:17
Operator : RC/JU
Sample : Q2430-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MH-E/F

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF061925.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF063025\
Data File : BF142912.D
Acq On : 30 Jun 2025 12:17
Operator : RC/JU
Sample : Q2430-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MH-E/F

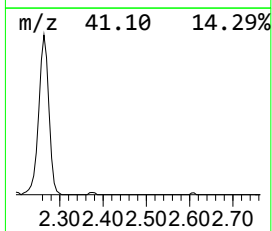
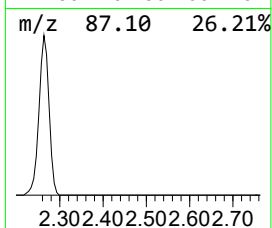
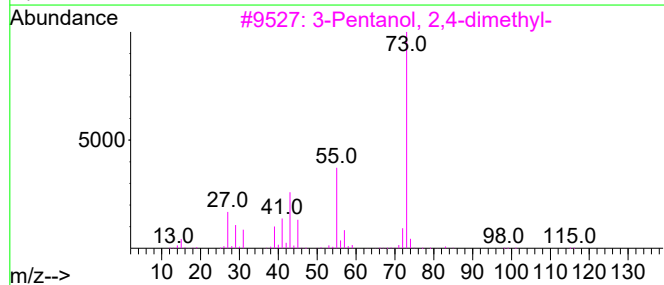
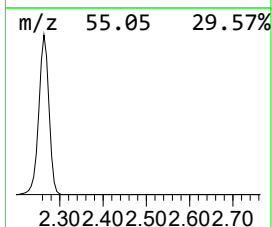
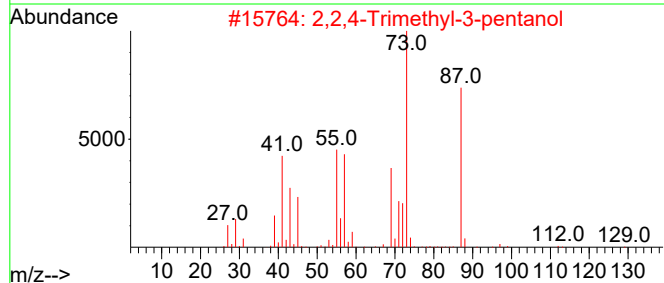
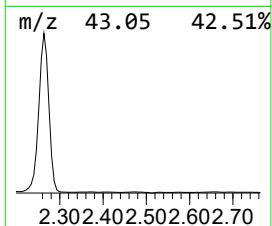
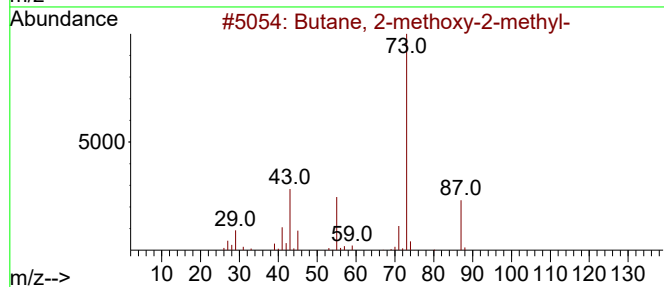
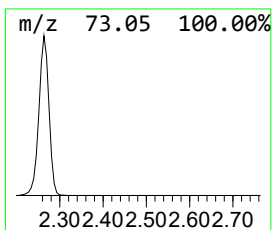
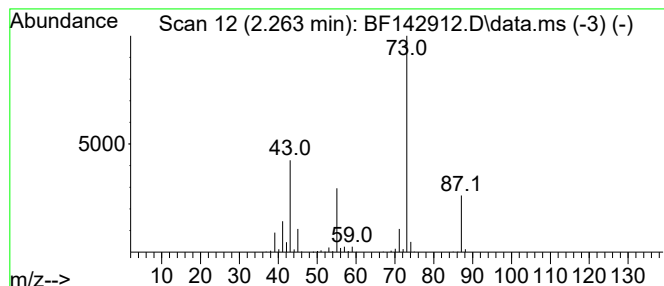
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF061925.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.263	13.84 ng	423059	1,4-Dichlorobenzene-d4	6.881

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	40
3			3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	37
4			1,3-Dioxolane, 2-propyl-	116	C6H12O2	003390-13-4	23
5			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	23



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF063025\
Data File : BF142912.D
Acq On : 30 Jun 2025 12:17
Operator : RC/JU
Sample : Q2430-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MH-E/F

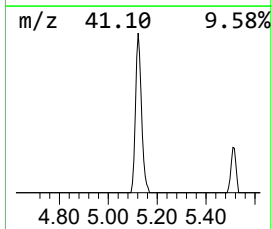
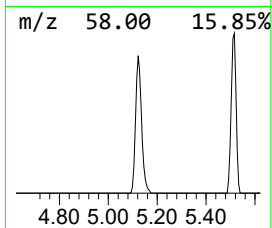
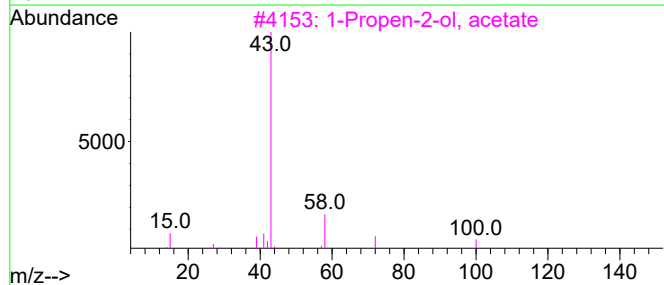
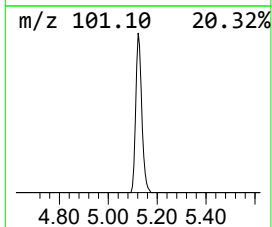
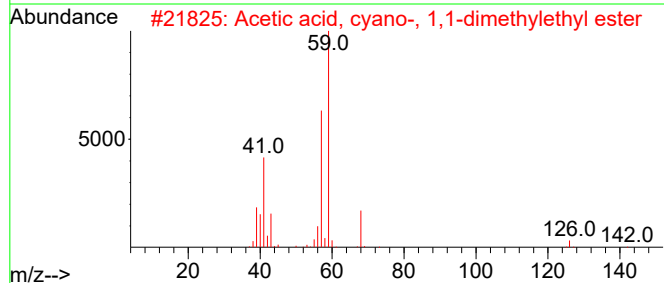
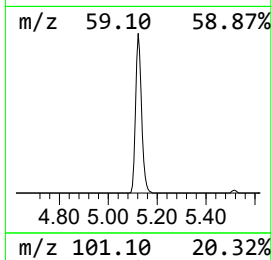
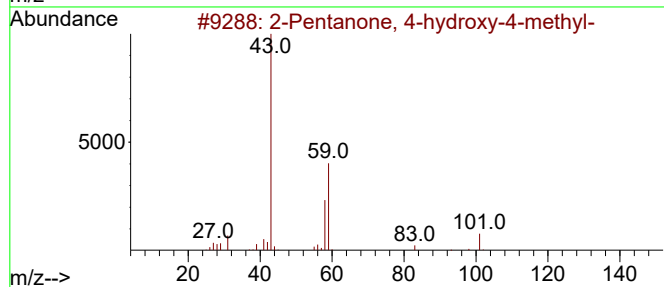
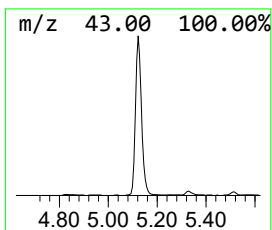
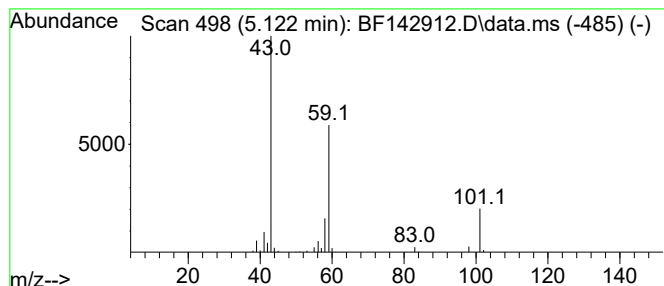
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF061925.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.122	7.28 ng	222465	1,4-Dichlorobenzene-d4	6.881

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
3			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
4			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF063025\
Data File : BF142912.D
Acq On : 30 Jun 2025 12:17
Operator : RC/JU
Sample : Q2430-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MH-E/F

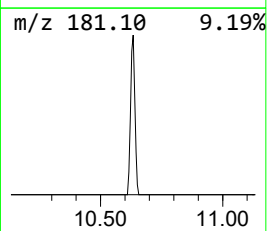
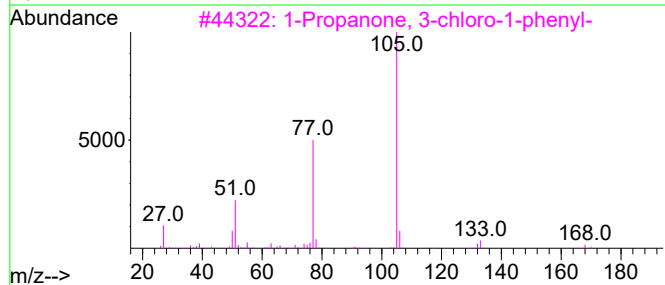
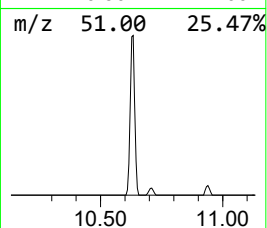
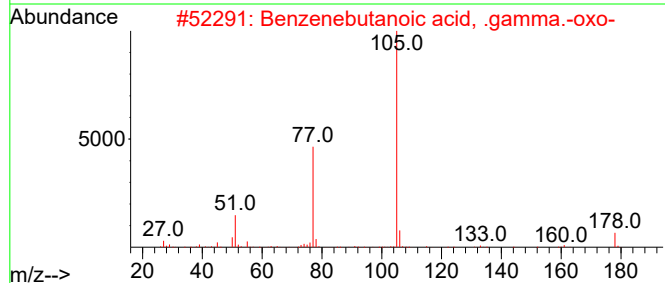
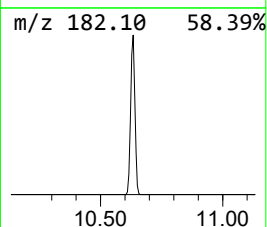
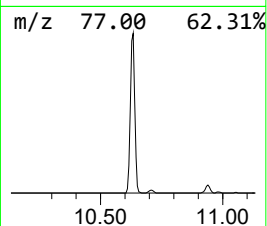
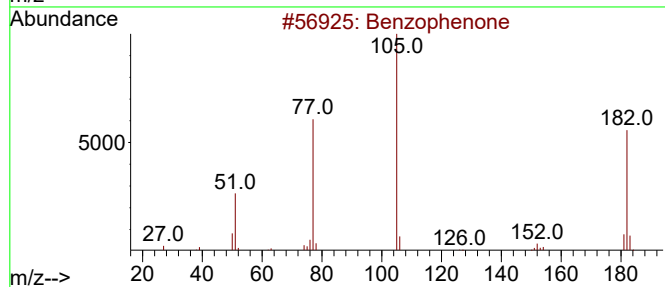
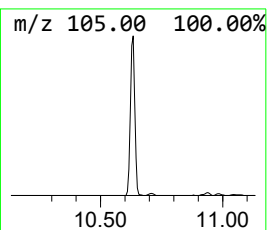
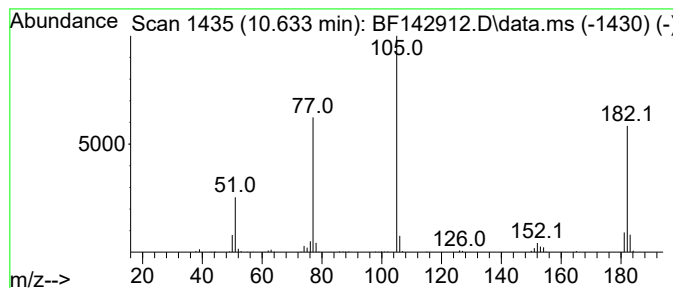
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF061925.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Benzophenone Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.633	5.02 ng	239948	Acenaphthene-d10	9.916

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzenebutanoic acid, .gamma.-oxo-	178	C10H10O3	002051-95-8	47
3			1-Propanone, 3-chloro-1-phenyl-	168	C9H9ClO	000936-59-4	47
4			Vinyl benzoate	148	C9H8O2	000769-78-8	47
5			2-(2-Oxo-2-phenyl-ethyl)-malonon...	184	C11H8N2O	1000296-76-9	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF063025\
Data File : BF142912.D
Acq On : 30 Jun 2025 12:17
Operator : RC/JU
Sample : Q2430-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MH-E/F

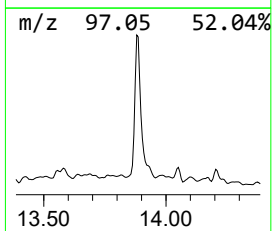
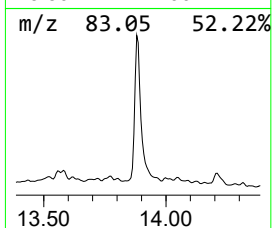
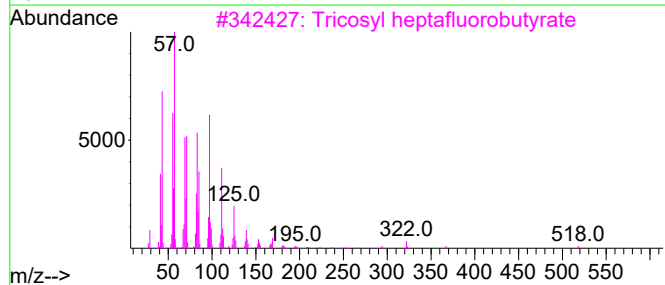
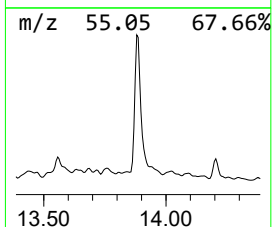
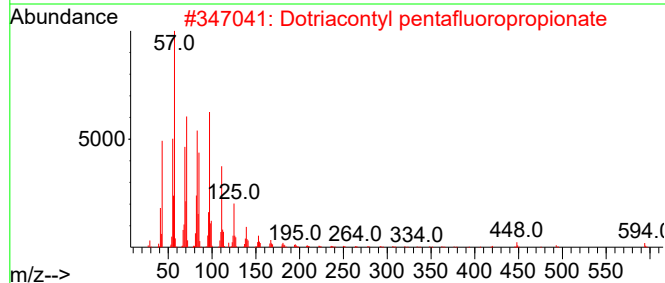
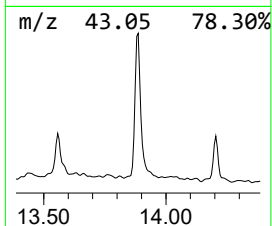
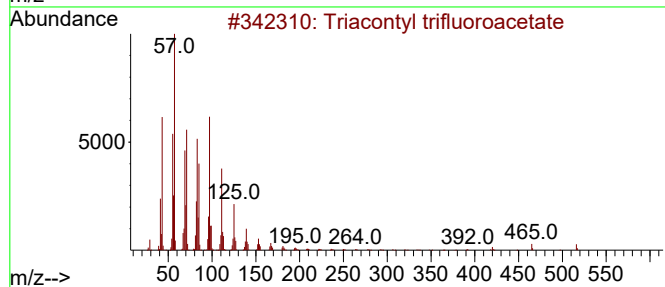
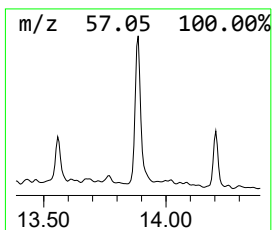
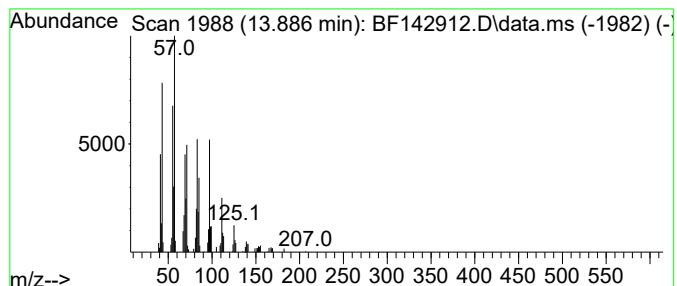
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF061925.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Triacontyl trifluoroacetate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.886	3.93 ng	146385	Chrysene-d12	14.051

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Triacontyl trifluoroacetate	534	C32H61F3O2	1000351-75-7	91
2			Dotriacontyl pentafluoropropionate	612	C35H65F5O2	1010351-81-4	91
3			Tricosyl heptafluorobutyrate	536	C27H47F7O2	1000351-83-4	91
4			Octadecyl trifluoroacetate	366	C20H37F3O2	079392-43-1	91
5			Tricosyl pentafluoropropionate	486	C26H47F5O2	1000351-81-0	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF063025\
Data File : BF142912.D
Acq On : 30 Jun 2025 12:17
Operator : RC/JU
Sample : Q2430-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MH-E/F

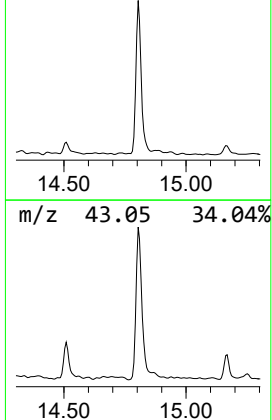
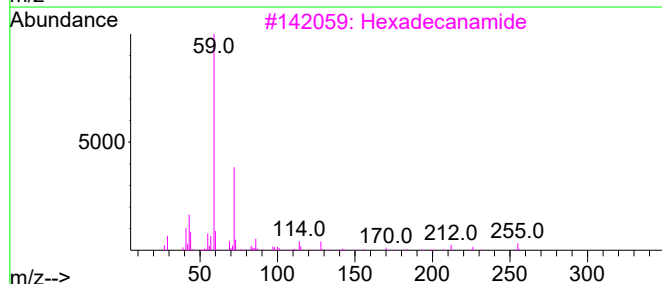
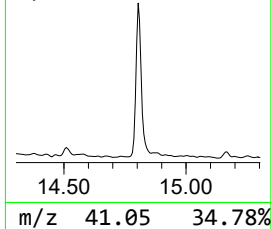
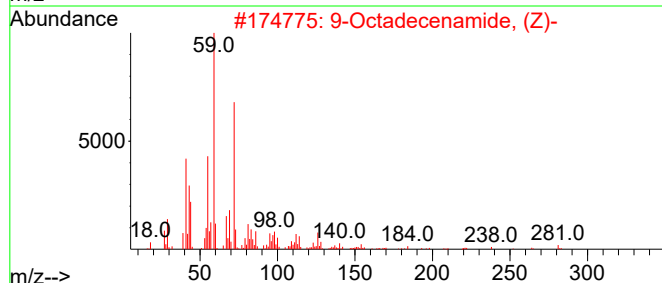
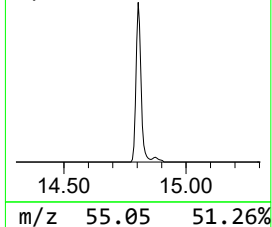
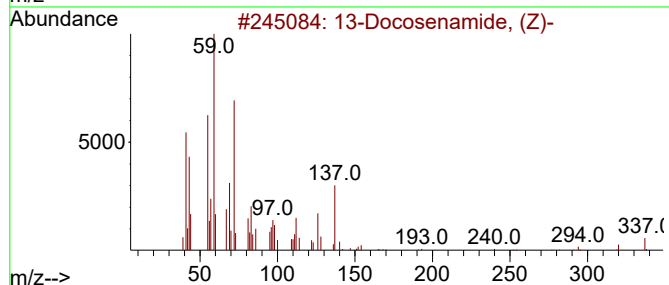
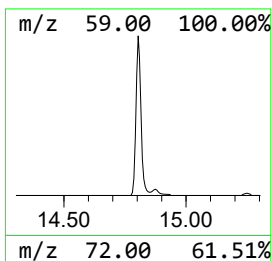
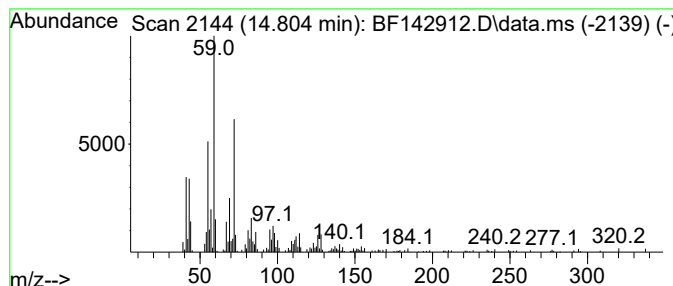
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF061925.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 13-Docosenamide, (Z)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.804	7.97 ng	290148	Perylene-d12	15.545

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	93
2			9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	91
3			Hexadecanamide	255	C16H33NO	000629-54-9	72
4			Benzeneethanamine, 2-fluoro-.bet...	229	C11H16FNO3	061338-98-5	59
5			2-Propanone, 1-cyclohexyl-	140	C9H16O	000103-78-6	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF063025\
Data File : BF142912.D
Acq On : 30 Jun 2025 12:17
Operator : RC/JU
Sample : Q2430-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MH-E/F

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF061925.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	2.263	13.8	ng	423059	1	6.881	611569	20.0
2-Pentanone, 4-...	5.122	7.3	ng	222465	1	6.881	611569	20.0
Benzophenone	10.633	5.0	ng	239948	3	9.916	956600	20.0
Triacetyl trif...	13.886	3.9	ng	146385	5	14.051	744647	20.0
13-Docosenamide...	14.804	8.0	ng	290148	6	15.545	728264	20.0