

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082025\
 Data File : BP025518.D
 Acq On : 20 Aug 2025 22:32
 Operator : CG/JU
 Sample : Q2819-09
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 84SB-E

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_P\Methods\8270E-BP081425.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP025518.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.043	13	18	33	rVB	1354160	1723255	24.31%	3.938%
2	4.943	335	341	348	rBV	202236	294684	4.16%	0.673%
3	5.408	413	420	433	rBV	2251831	3325388	46.92%	7.600%
4	6.966	675	685	694	rBV	2012748	3387963	47.80%	7.743%
5	7.784	817	824	834	rBV	613529	971840	13.71%	2.221%
6	8.925	1009	1018	1030	rBV	1410989	2411017	34.02%	5.510%
7	10.555	1285	1295	1307	rBV	745669	1288898	18.19%	2.946%
8	13.013	1705	1713	1724	rBV	3155027	4642109	65.50%	10.609%
9	14.395	1941	1948	1957	rBV	1044236	1690993	23.86%	3.865%
10	15.778	2176	2183	2190	rBV	336598	489796	6.91%	1.119%
11	15.913	2198	2206	2225	rBV	2485643	4001499	56.46%	9.145%
12	16.354	2276	2281	2287	rVB	71056	98611	1.39%	0.225%
13	17.195	2417	2424	2434	rBV2	1384770	2038796	28.77%	4.660%
14	18.136	2578	2584	2593	rBV	425310	627821	8.86%	1.435%
15	19.366	2789	2793	2800	rVB	323621	392986	5.54%	0.898%
16	19.460	2804	2809	2818	rBV	163942	294323	4.15%	0.673%
17	19.619	2831	2836	2841	rBV	866729	1032860	14.57%	2.361%
18	19.736	2853	2856	2861	rVB	107488	137420	1.94%	0.314%
19	19.913	2880	2886	2894	rBV	6003462	7087305	100.00%	16.198%
20	20.495	2981	2985	2989	rVB	231951	265347	3.74%	0.606%
21	20.713	3018	3022	3028	rBV	762399	863099	12.18%	1.973%
22	20.783	3031	3034	3037	rVV	93238	94778	1.34%	0.217%
23	21.313	3119	3124	3135	rBV	490039	827965	11.68%	1.892%
24	21.636	3173	3179	3187	rVB2	1419429	2419088	34.13%	5.529%
25	21.901	3221	3224	3237	rVB2	148693	250898	3.54%	0.573%
26	22.554	3332	3335	3351	rVB4	160190	385776	5.44%	0.882%
27	25.013	3745	3753	3769	rVB	982410	2710118	38.24%	6.194%

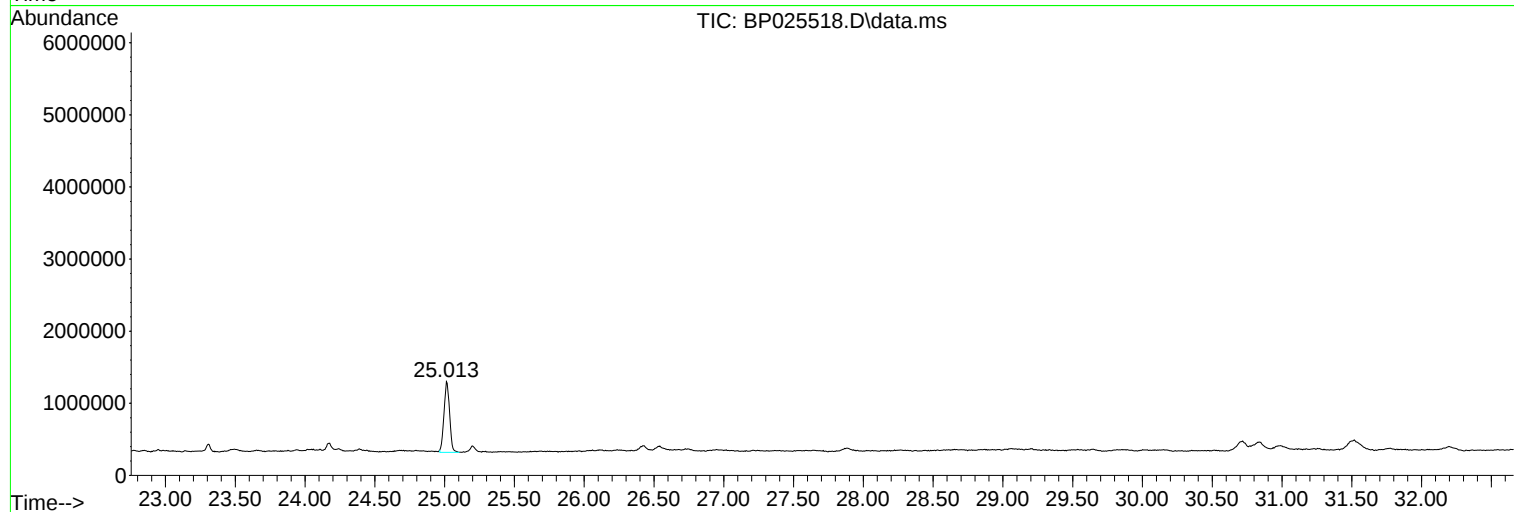
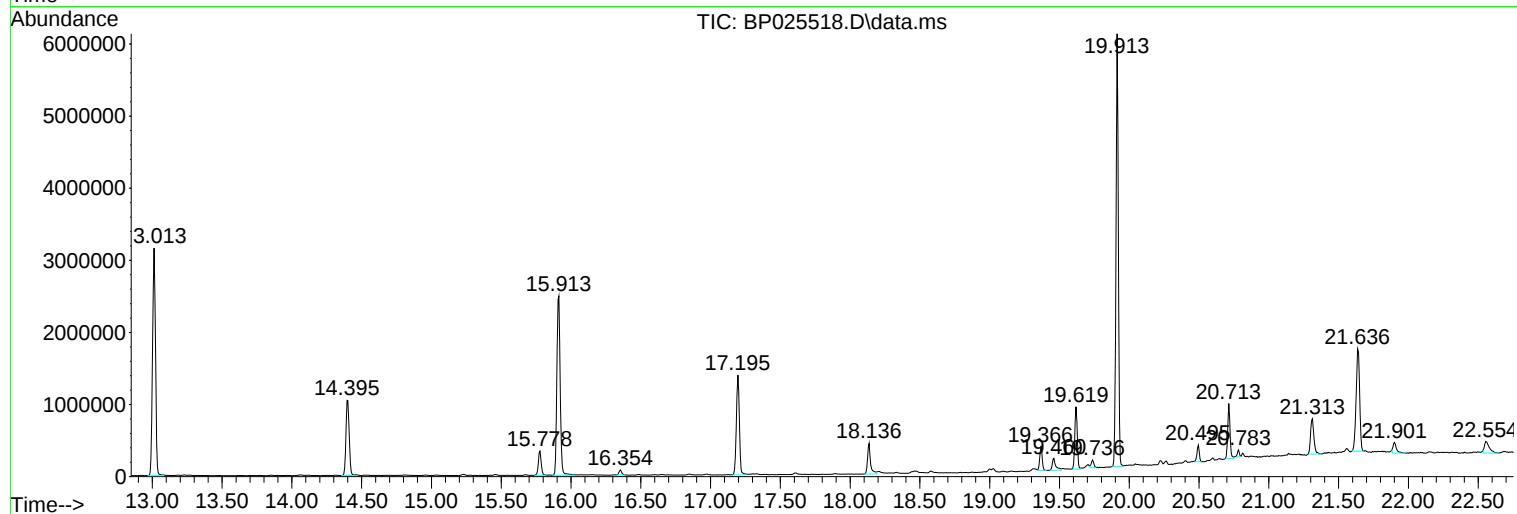
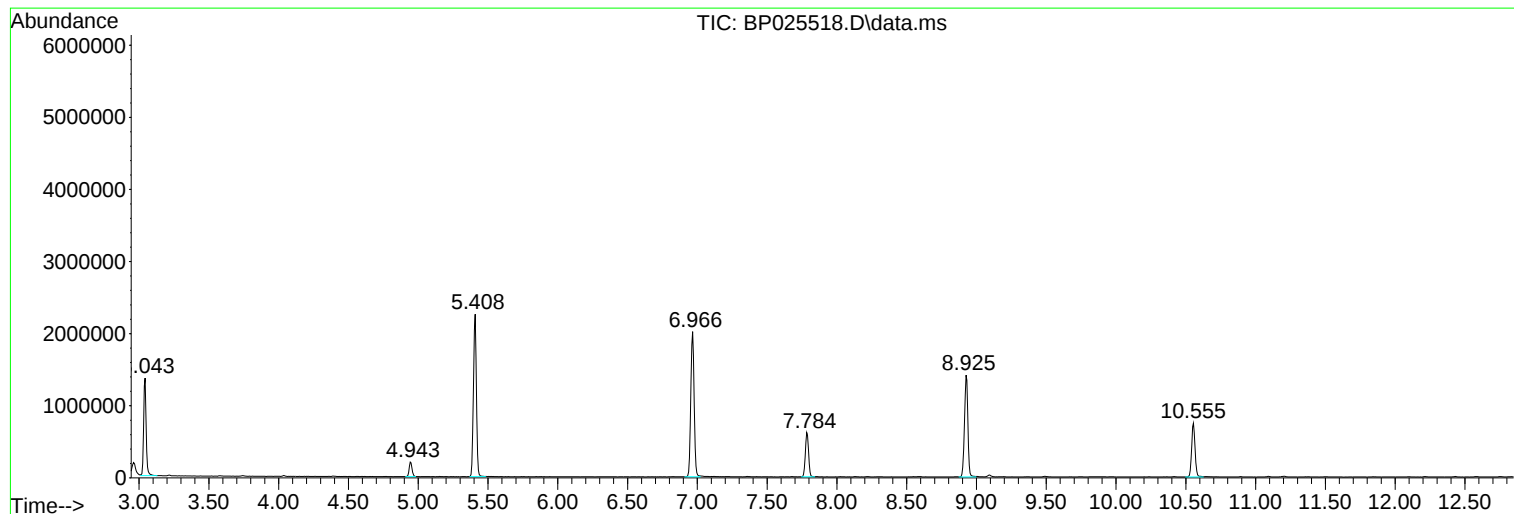
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Instrument :
BNA_P
ClientSampleId :
84SB-E

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP081425.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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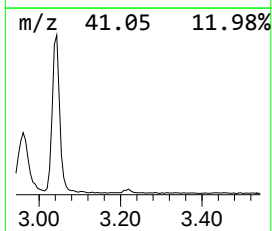
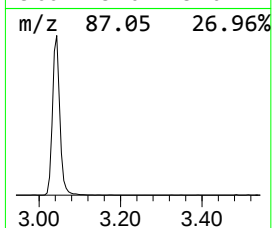
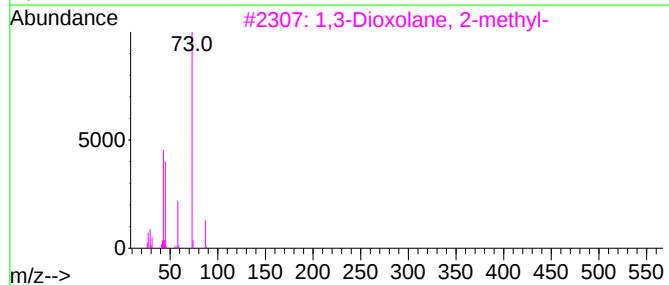
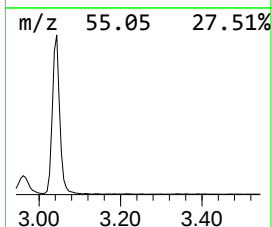
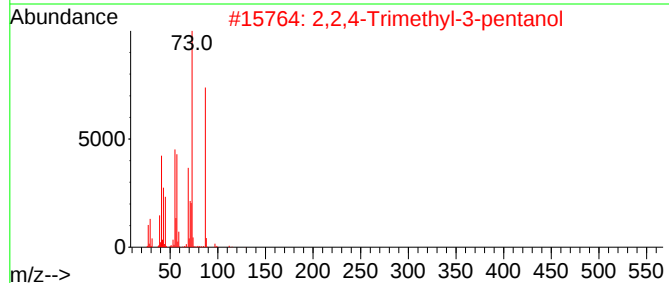
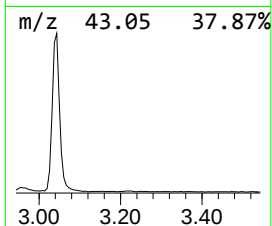
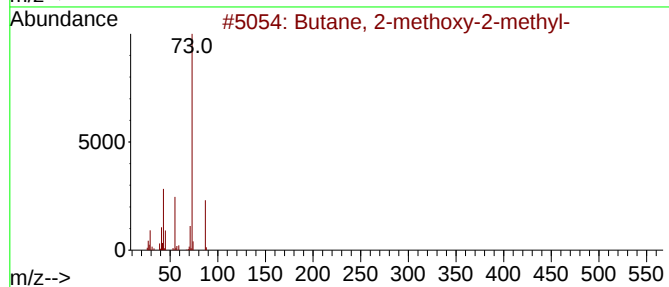
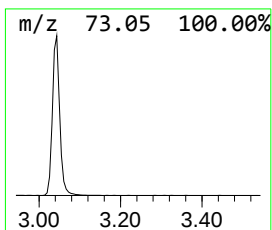
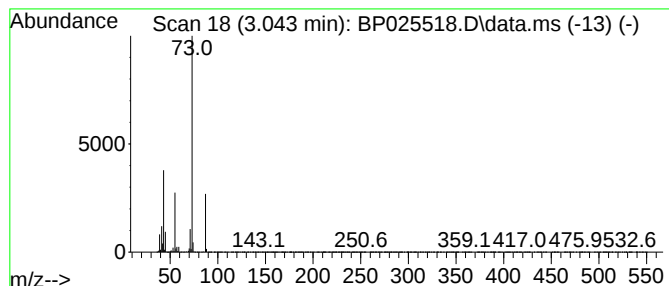
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.043	35.46 ng	1723260	1,4-Dichlorobenzene-d4	7.784

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			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	23
4			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	22
5			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12



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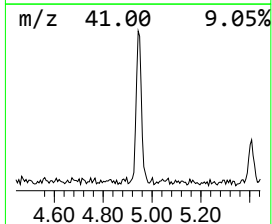
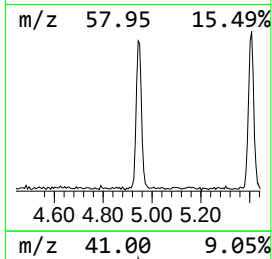
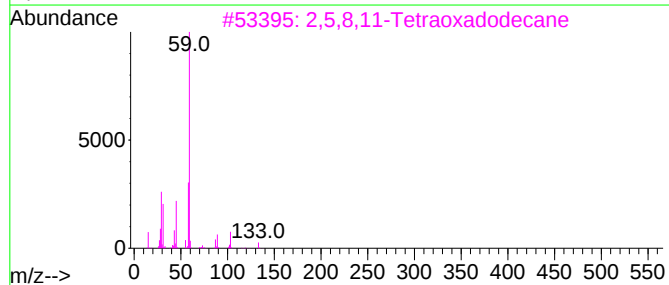
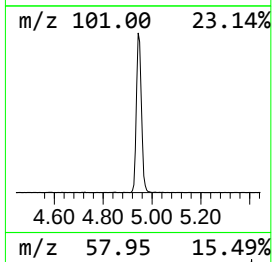
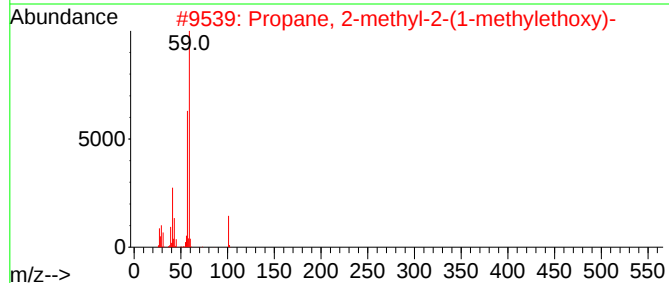
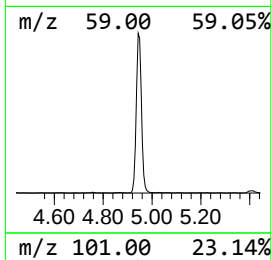
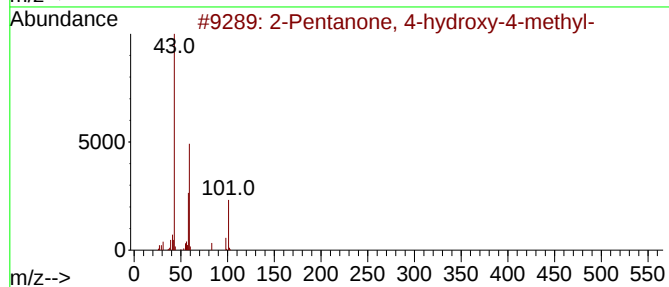
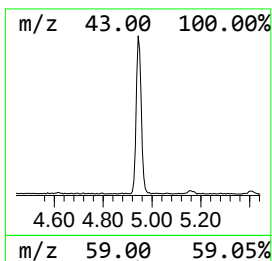
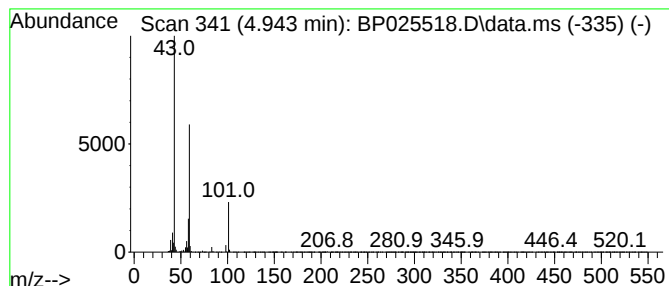
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.943	6.06 ng	294684	1,4-Dichlorobenzene-d4	7.784

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	53
3			2,5,8,11-Tetraoxadodecane	178	C8H18O4	000112-49-2	23
4			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
5			Formamide, N,N-diethyl-	101	C5H11NO	000617-84-5	9



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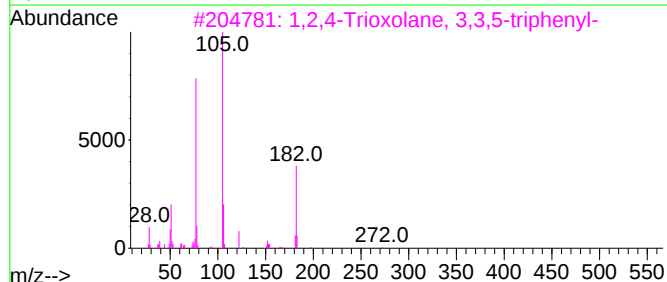
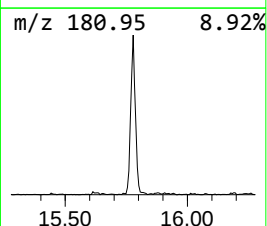
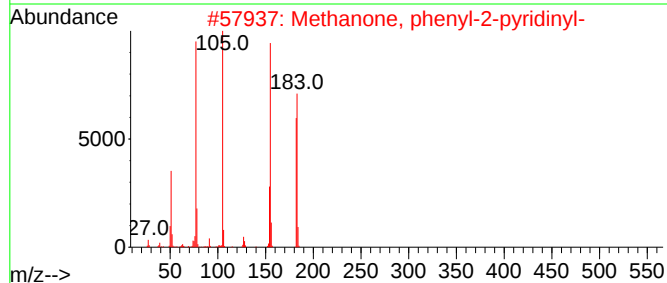
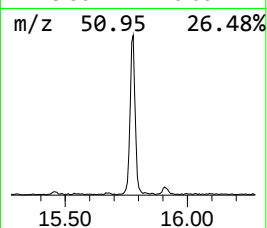
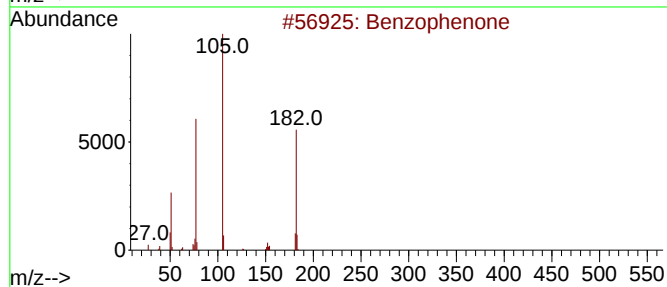
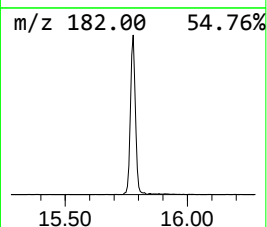
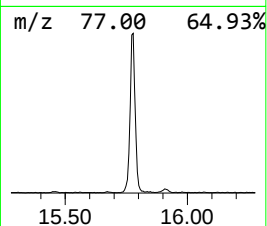
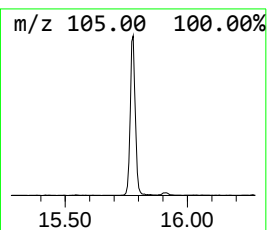
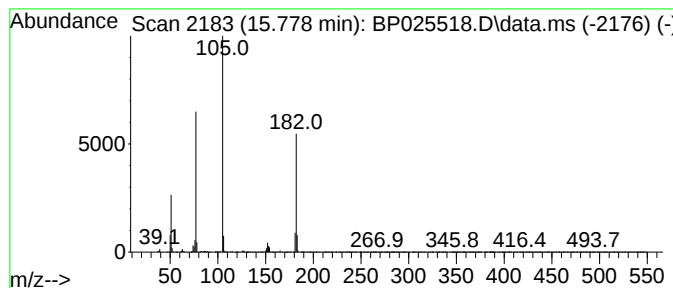
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzophenone Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.778	5.79 ng	489796	Acenaphthene-d10	14.401

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Methanone, phenyl-2-pyridinyl-	183	C12H9NO	000091-02-1	56
3			1,2,4-Trioxolane, 3,3,5-triphenyl-	304	C20H16O3	023246-12-0	45
4			Benzenamine, N-(3-pyridinylmethy...	182	C12H10N2	029722-97-2	45
5			Azobenzene	182	C12H10N2	000103-33-3	42



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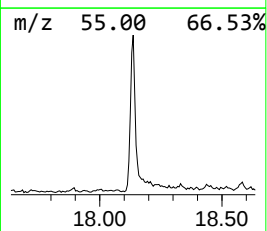
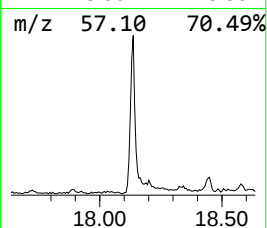
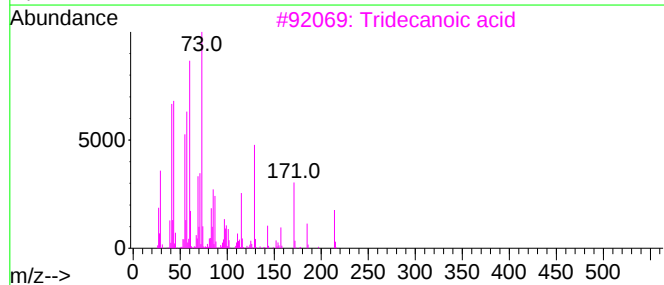
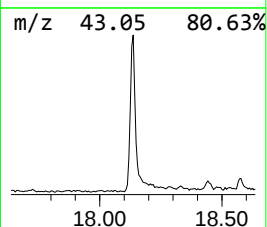
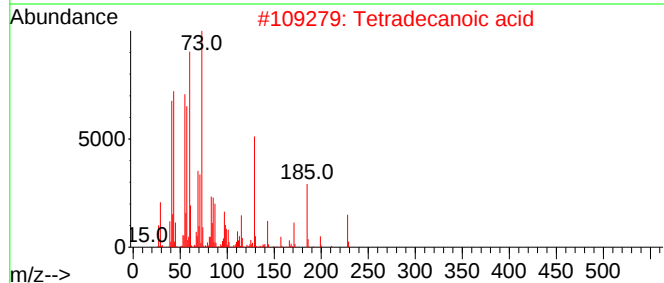
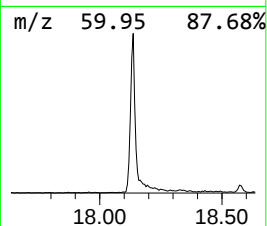
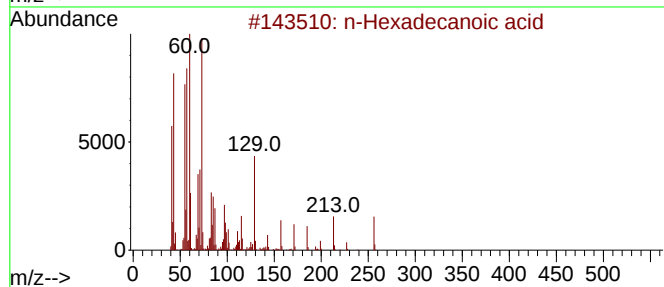
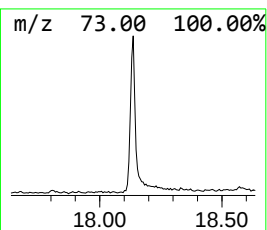
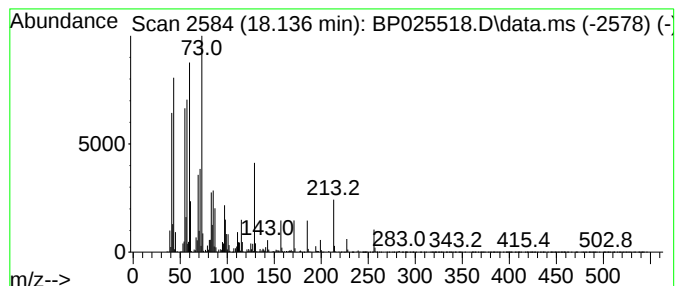
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Peak Number 4 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.137	6.16 ng	627821	Phenanthrene-d10	17.195

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	93
3		Tridecanoic acid	214	C13H26O2	000638-53-9	90
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	90
5		n-Decanoic acid	172	C10H20O2	000334-48-5	68



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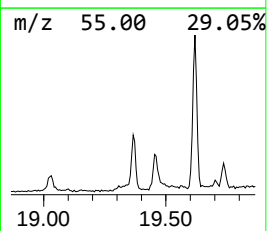
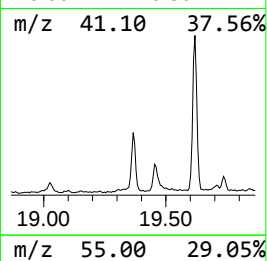
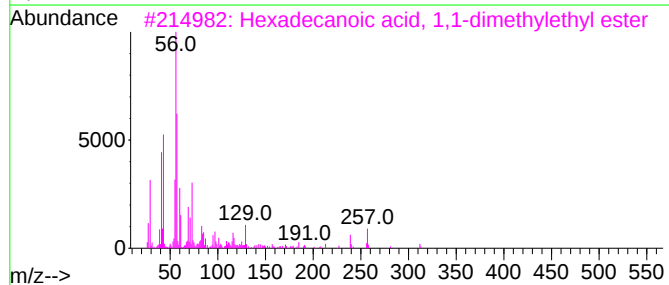
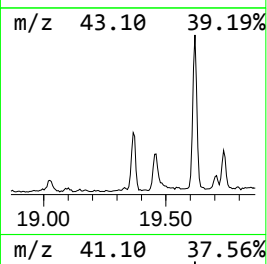
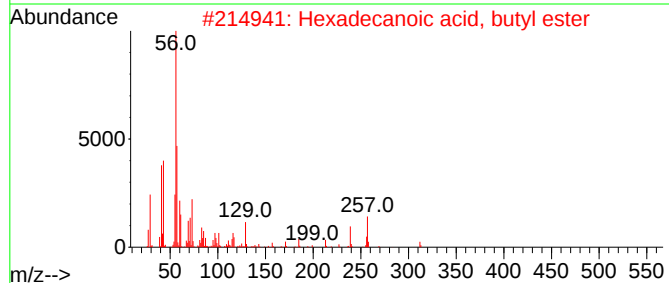
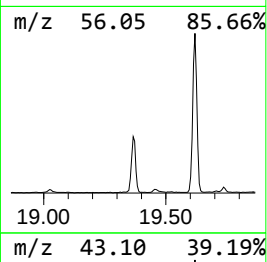
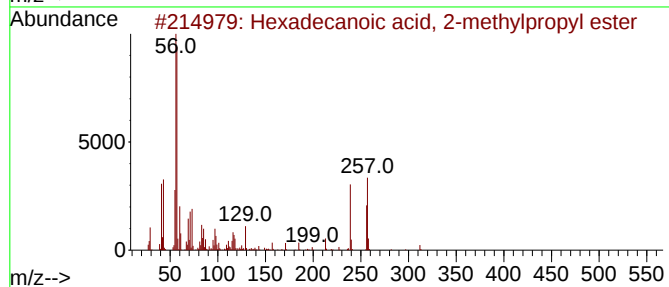
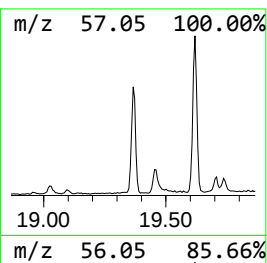
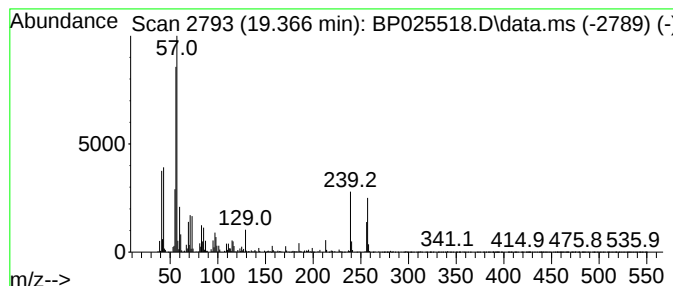
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Hexadecanoic acid, 2-methyl... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.366	3.86 ng	392986	Phenanthrene-d10	17.195

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecanoic acid, 2-methylpropy...	312	C20H40O2	000110-34-9	90
2			Hexadecanoic acid, butyl ester	312	C20H40O2	000111-06-8	64
3			Hexadecanoic acid, 1,1-dimethyle...	312	C20H40O2	031158-91-5	47
4			Urea, N,N'-di-2-propenyl-	140	C7H12N2O	001801-72-5	35
5			Decane, 2,2,3-trimethyl-	184	C13H28	062338-09-4	35



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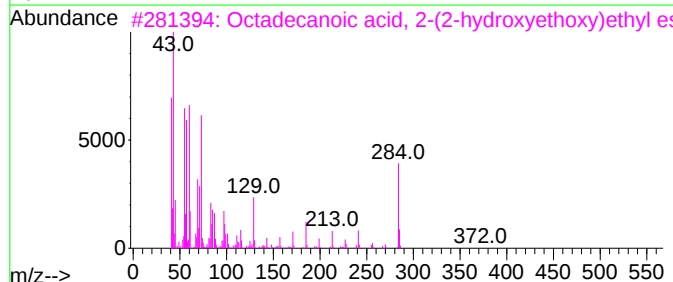
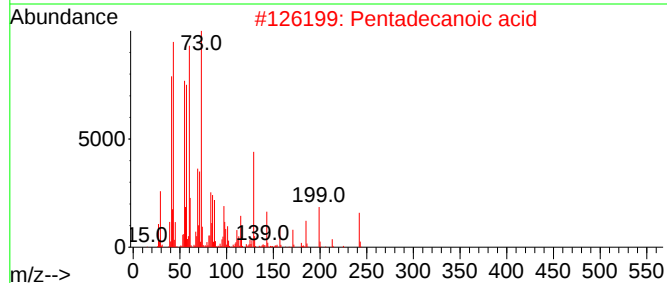
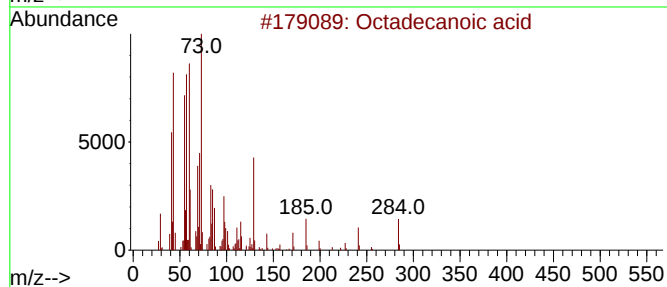
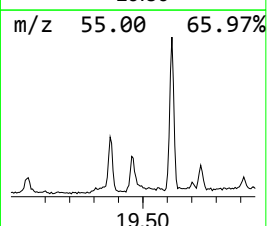
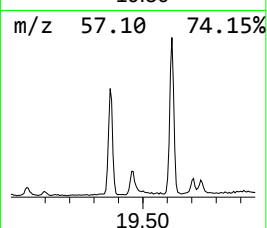
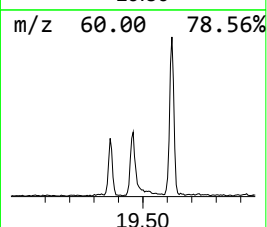
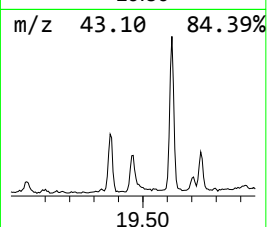
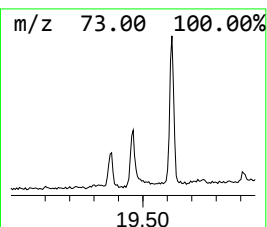
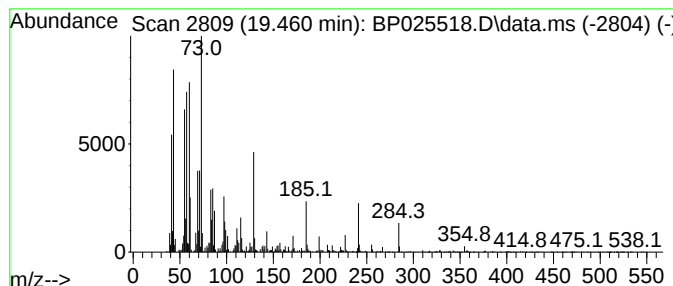
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Octadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.460	2.43 ng	294323	Chrysene-d12	21.636

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Pentadecanoic acid	242	C15H30O2	001002-84-2	89
3			Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	72
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	64
5			n-Decanoic acid	172	C10H20O2	000334-48-5	64



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Acq On : 20 Aug 2025 22:32
Operator : CG/JU
Sample : Q2819-09
Misc :
ALS Vial : 8 Sample Multiplier: 1

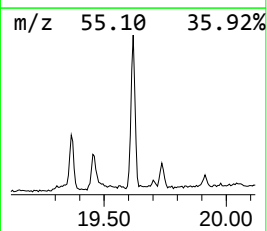
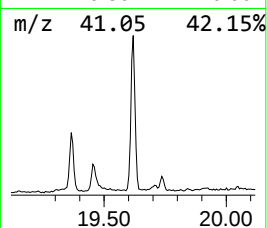
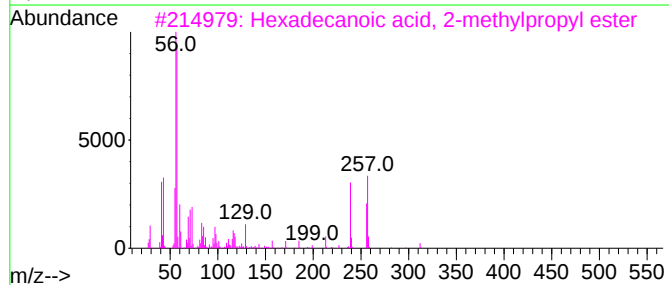
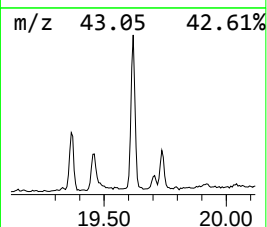
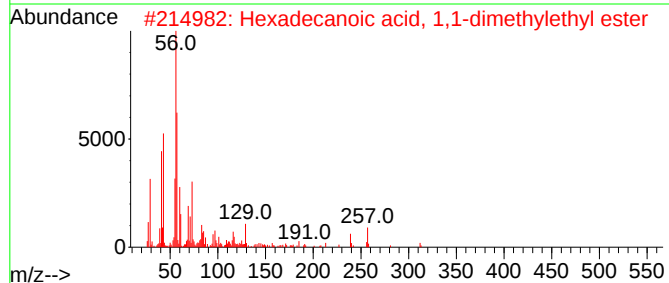
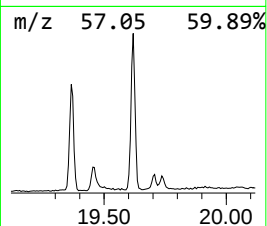
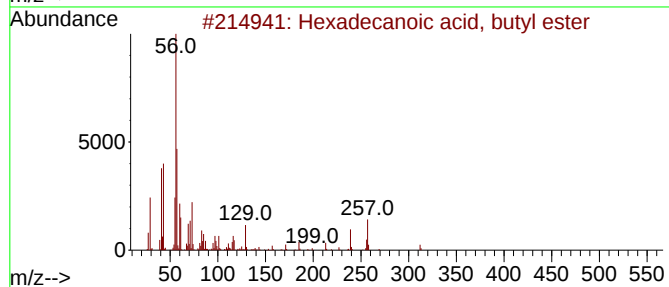
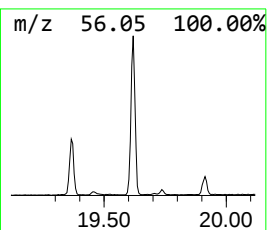
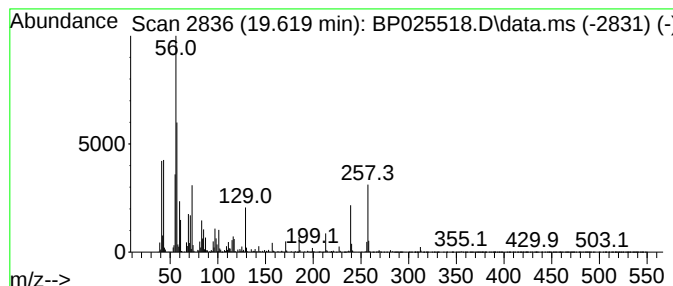
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Hexadecanoic acid, butyl ester Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.619	8.54 ng	1032860	Chrysene-d12	21.636	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecanoic acid, butyl ester	312	C20H40O2	000111-06-8	98
2	Hexadecanoic acid, 1,1-dimethyle...	312	C20H40O2	031158-91-5	96
3	Hexadecanoic acid, 2-methylpropy...	312	C20H40O2	000110-34-9	90
4	Cyclohexanol, 4-amino-, trans-	115	C6H13NO	027489-62-9	27
5	1-Heptene, 2-methyl-	112	C8H16	015870-10-7	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082025\
Data File : BP025518.D
Acq On : 20 Aug 2025 22:32
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Sample : Q2819-09
Misc :
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Instrument :
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ClientSampleId :
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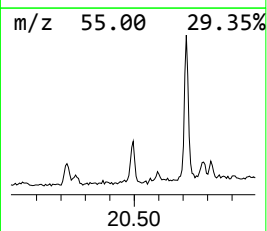
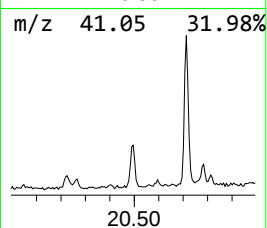
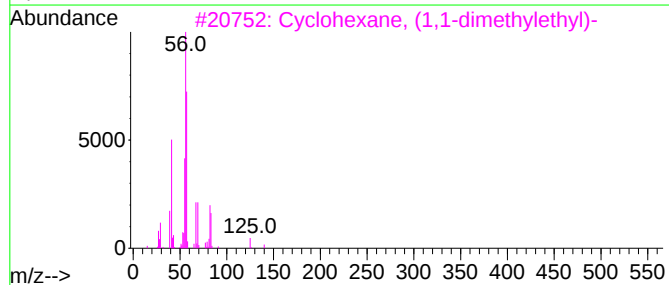
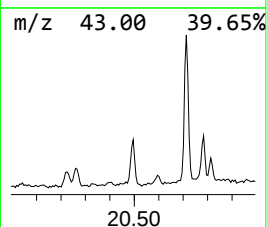
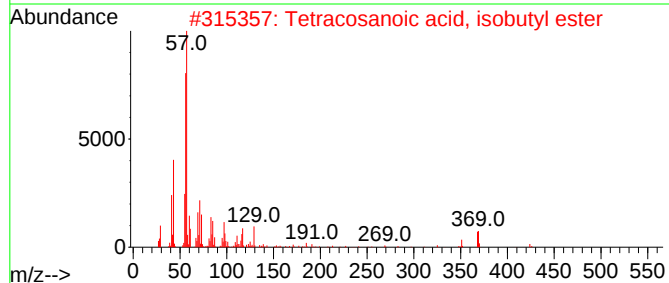
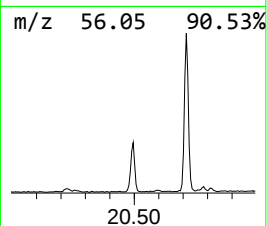
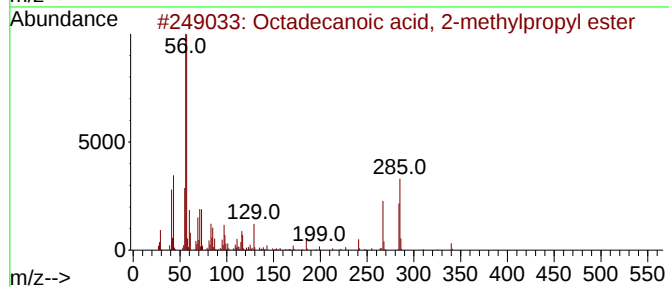
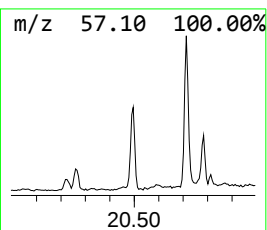
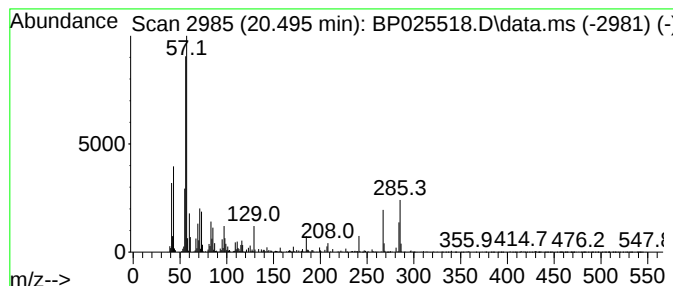
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Octadecanoic acid, 2-methyl... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.495	2.19 ng	265347	Chrysene-d12	21.636

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid, 2-methylpropy...	340	C22H44O2	000646-13-9	91
2			Tetracosanoic acid, isobutyl ester	424	C28H56O2	1010405-17-3	60
3			Cyclohexane, (1,1-dimethylethyl)-	140	C10H20	003178-22-1	38
4			Decane, 2,2,3-trimethyl-	184	C13H28	062338-09-4	38
5			Cyclopentanecarboxylic acid, 2-a...	129	C6H11NO2	037910-65-9	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082025\
Data File : BP025518.D
Acq On : 20 Aug 2025 22:32
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Misc :
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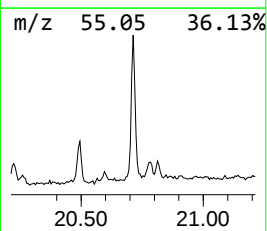
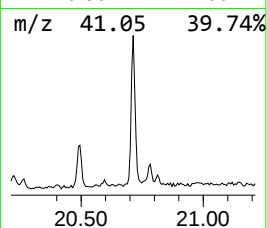
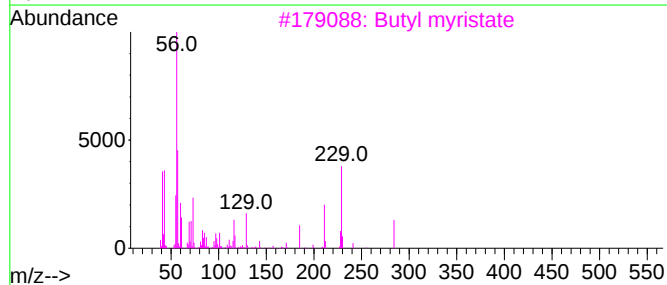
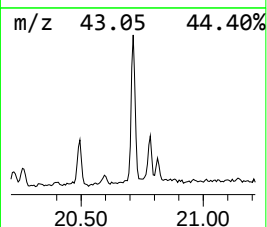
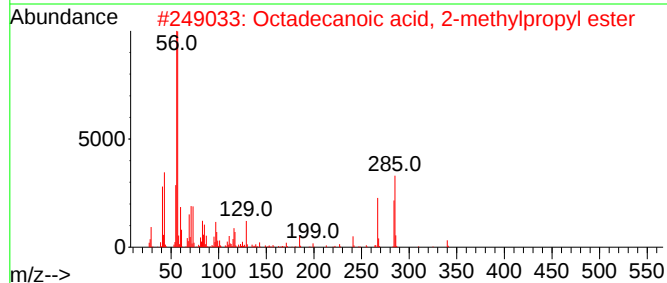
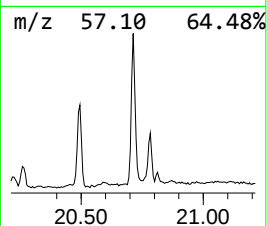
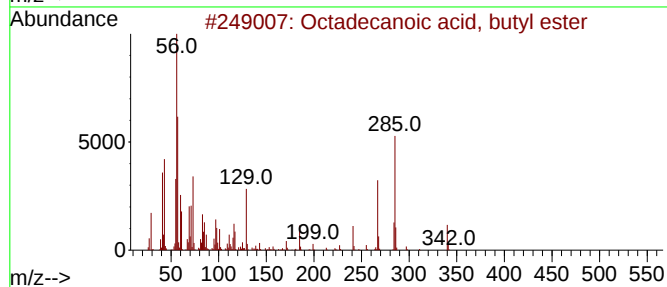
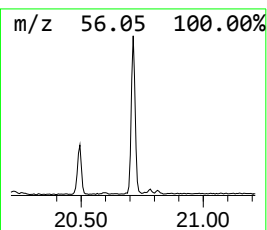
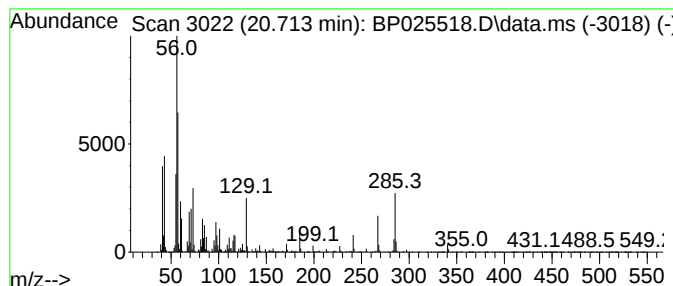
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Octadecanoic acid, butyl ester Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.713	7.14 ng	863099	Chrysene-d12	21.636

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid, butyl ester	340	C22H44O2	000123-95-5	99
2			Octadecanoic acid, 2-methylpropyl ester	340	C22H44O2	000646-13-9	89
3			Butyl myristate	284	C18H36O2	000110-36-1	50
4			Docosanoic acid, isobutyl ester	396	C26H52O2	1000405-17-1	43
5			Piperidin-4-carboxylic acid	129	C6H11NO2	000498-94-2	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082025\
Data File : BP025518.D
Acq On : 20 Aug 2025 22:32
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Sample : Q2819-09
Misc :
ALS Vial : 8 Sample Multiplier: 1

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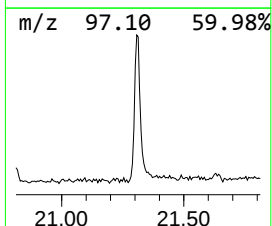
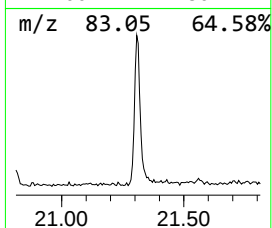
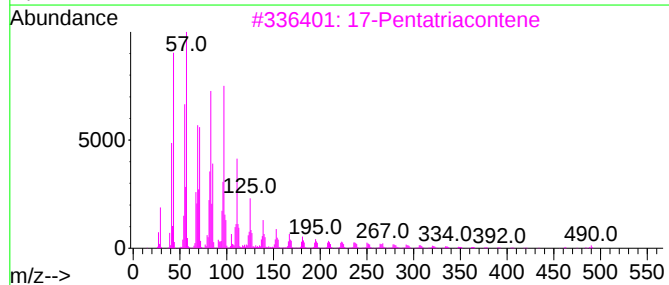
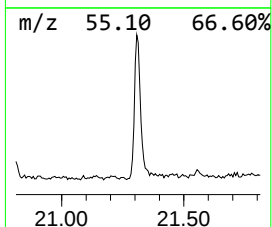
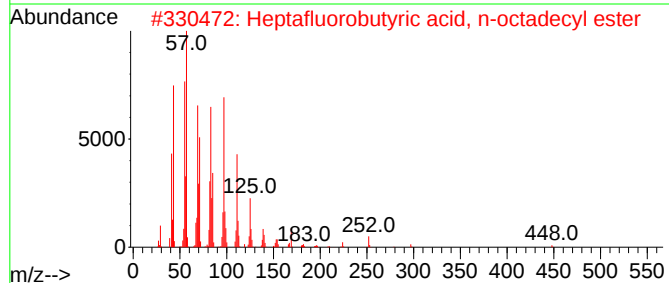
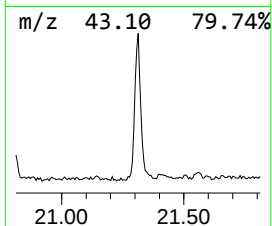
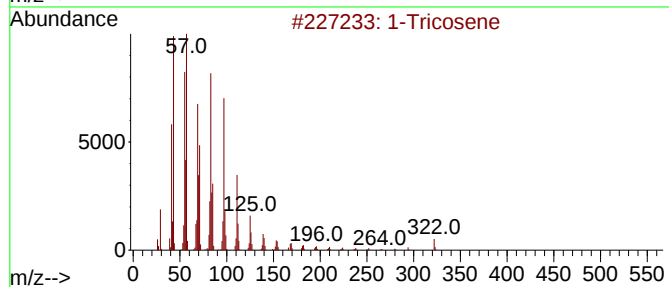
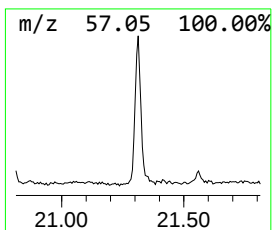
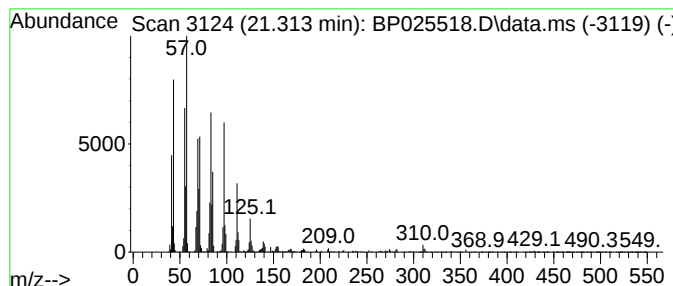
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 1-Tricosene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.313	6.85 ng	827965	Chrysene-d12	21.636

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Tricosene	322	C23H46	018835-32-0	95
2			Heptafluorobutyric acid, n-octad...	466	C22H37F7O2	000400-57-7	94
3			17-Pentatriacontene	491	C35H70	006971-40-0	94
4			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	94
5			1-Hexacosene	364	C26H52	018835-33-1	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082025\
Data File : BP025518.D
Acq On : 20 Aug 2025 22:32
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Sample : Q2819-09
Misc :
ALS Vial : 8 Sample Multiplier: 1

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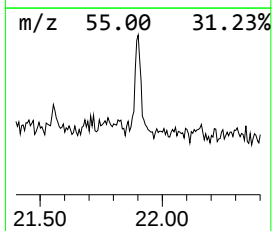
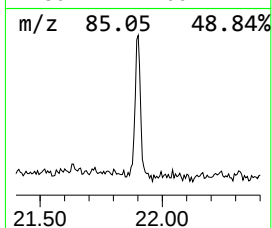
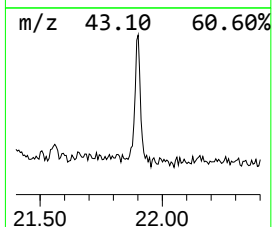
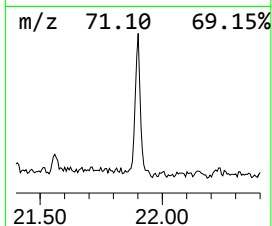
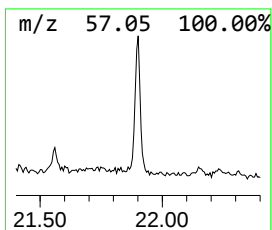
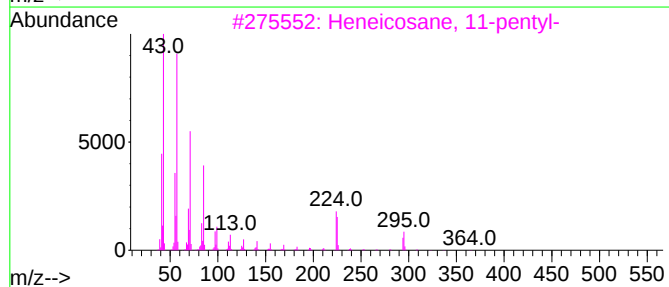
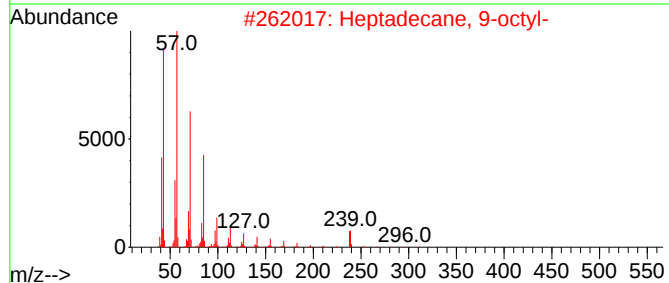
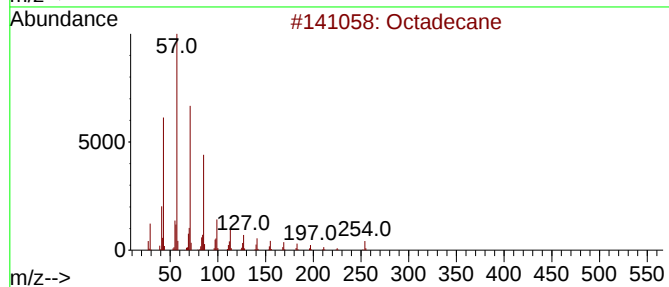
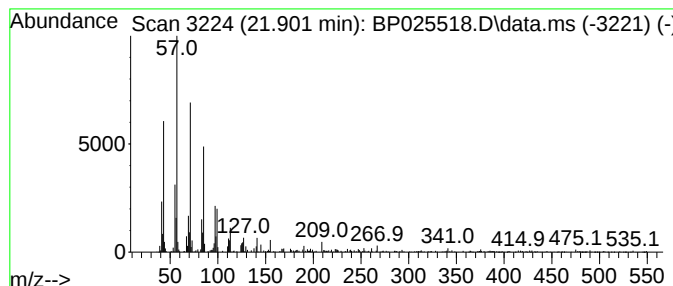
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Octadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.901	2.07 ng	250898	Chrysene-d12	21.636

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecane	254	C18H38	000593-45-3	89
2		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	89
3		Heneicosane, 11-pentyl-	366	C26H54	014739-72-1	86
4		Docosane, 11-butyl-	366	C26H54	013475-76-8	86
5		Tetracosane	338	C24H50	000646-31-1	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082025\
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Misc :
ALS Vial : 8 Sample Multiplier: 1

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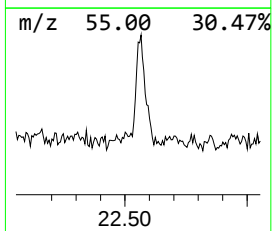
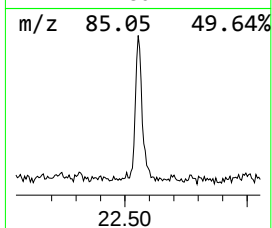
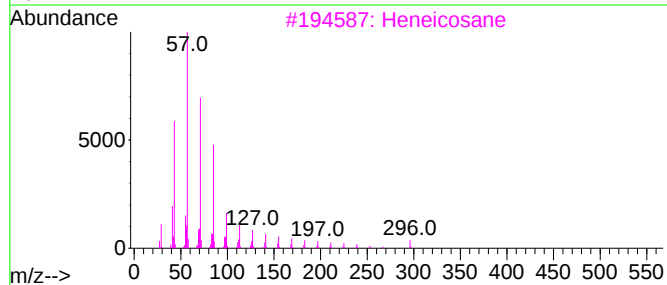
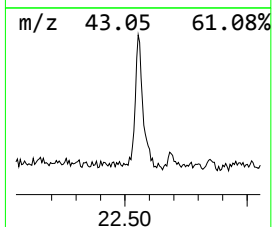
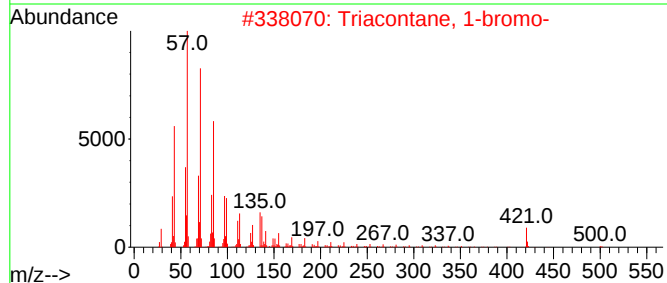
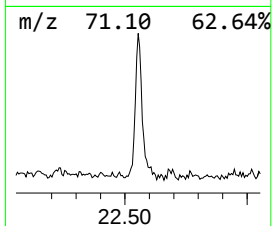
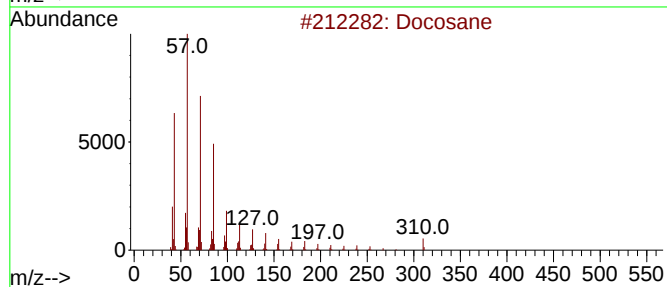
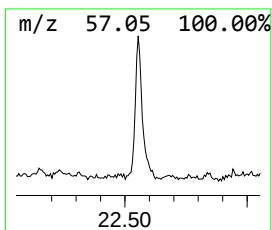
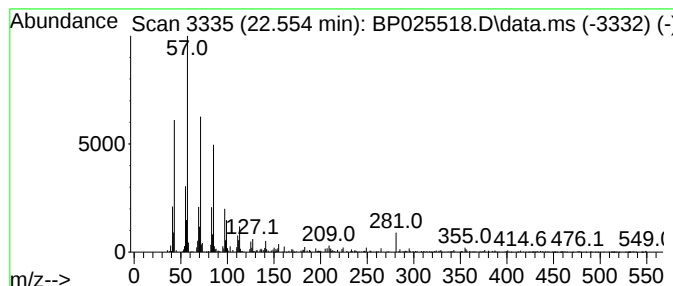
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Docosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.554	3.19 ng	385776	Chrysene-d12	21.636

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Docosane	310	C22H46	000629-97-0	93
2		Triacontane, 1-bromo-	500	C30H61Br	004209-22-7	91
3		Heneicosane	296	C21H44	000629-94-7	91
4		Tetratriacontane, 17-hexadecyl-	703	C50H102	055256-07-0	90
5		Hentriacontane	437	C31H64	000630-04-6	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082025\
Data File : BP025518.D
Acq On : 20 Aug 2025 22:32
Operator : CG/JU
Sample : Q2819-09
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
84SB-E

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP081425.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...	3.043	35.5	ng	1723260	1	7.784	971840	20.0
2-Pentanone, 4-...	4.943	6.1	ng	294684	1	7.784	971840	20.0
Benzophenone	15.778	5.8	ng	489796	3	14.401	1690990	20.0
n-Hexadecanoic ...	18.137	6.2	ng	627821	4	17.195	2038800	20.0
Hexadecanoic ac...	19.366	3.9	ng	392986	4	17.195	2038800	20.0
Octadecanoic acid	19.460	2.4	ng	294323	5	21.636	2419090	20.0
Hexadecanoic ac...	19.619	8.5	ng	1032860	5	21.636	2419090	20.0
Octadecanoic ac...	20.495	2.2	ng	265347	5	21.636	2419090	20.0
Octadecanoic ac...	20.713	7.1	ng	863099	5	21.636	2419090	20.0
1-Tricosene	21.313	6.8	ng	827965	5	21.636	2419090	20.0
Octadecane	21.901	2.1	ng	250898	5	21.636	2419090	20.0
Docosane	22.554	3.2	ng	385776	5	21.636	2419090	20.0