

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071425\
 Data File : BP025135.D
 Acq On : 15 Jul 2025 01:22
 Operator : CG/JU
 Sample : Q2564-05
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 ARS20-0039

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-BP070325.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP025135.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.661	287	293	302	rVB	654013	871687	3.47%	0.444%
2	5.102	362	368	375	rBV	7062160	9998164	39.78%	5.092%
3	6.631	618	628	635	rBV	7014321	10322678	41.07%	5.257%
4	7.437	758	765	772	rBV	2566878	3968185	15.79%	2.021%
5	8.566	946	957	964	rBV	4456034	6822938	27.15%	3.475%
6	10.184	1224	1232	1238	rBV	2893924	5725727	22.78%	2.916%
7	10.231	1238	1240	1246	rVB	473557	617399	2.46%	0.314%
8	11.784	1500	1504	1510	rBV6	391468	822967	3.27%	0.419%
9	11.960	1528	1534	1537	rBV3	685932	1304777	5.19%	0.665%
10	12.272	1584	1587	1591	rBV3	363685	572775	2.28%	0.292%
11	12.372	1600	1604	1609	rBV5	564219	1231616	4.90%	0.627%
12	12.431	1610	1614	1620	rVV7	689850	1538896	6.12%	0.784%
13	12.501	1621	1626	1632	rVV	1700936	2817133	11.21%	1.435%
14	12.690	1652	1658	1663	rBV	11089354	16516664	65.72%	8.412%
15	12.860	1684	1687	1691	rVB2	1663518	2232695	8.88%	1.137%
16	13.366	1771	1773	1778	rVB4	2070748	2723031	10.83%	1.387%
17	13.419	1779	1782	1786	rVV2	1588933	2736082	10.89%	1.393%
18	13.460	1786	1789	1793	rVB2	2768147	3775903	15.02%	1.923%
19	13.513	1793	1798	1803	rBV	10544720	15018694	59.76%	7.649%
20	13.837	1849	1853	1858	rBV5	4481670	8140764	32.39%	4.146%
21	13.937	1866	1870	1876	rVB	12330750	19366112	77.05%	9.863%
22	14.084	1890	1895	1900	rBV2	7780785	17507052	69.66%	8.916%
23	14.454	1954	1958	1961	rBV	3730048	6958944	27.69%	3.544%
24	14.925	2033	2038	2044	rVB2	15968104	25133112	100.00%	12.800%
25	21.319	3121	3125	3129	rBV	13458886	19683401	78.32%	10.025%
26	22.266	3283	3286	3290	rBV	2446736	3529192	14.04%	1.797%
27	29.771	4558	4562	4573	rVB3	2251954	6409974	25.50%	3.265%

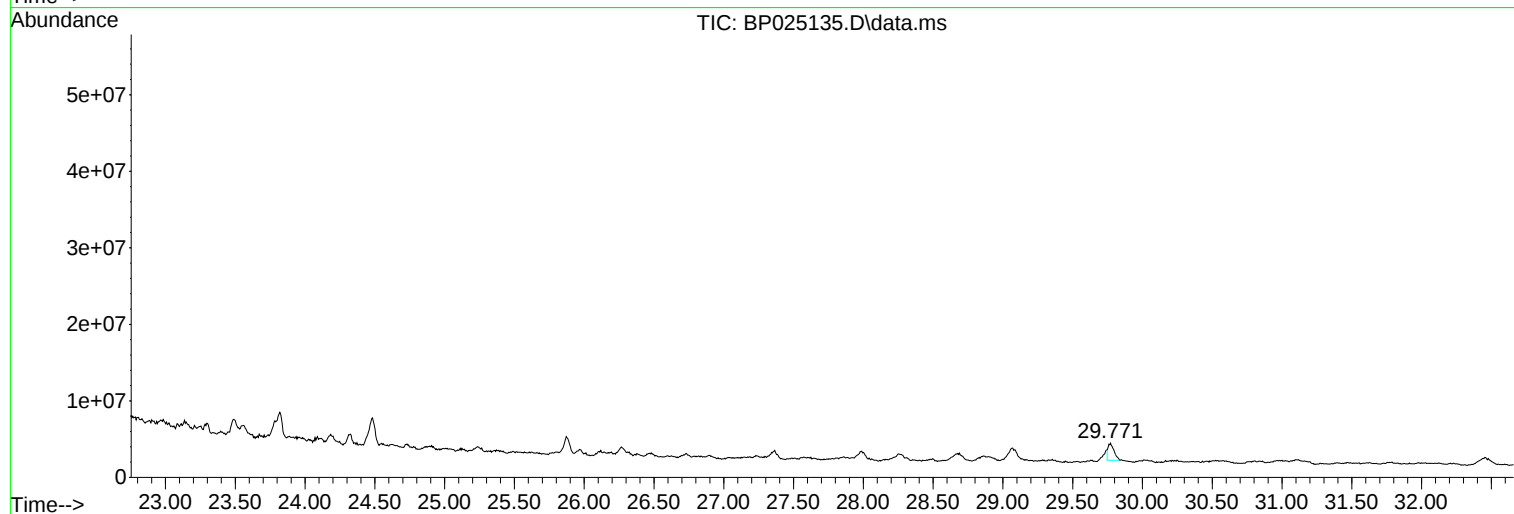
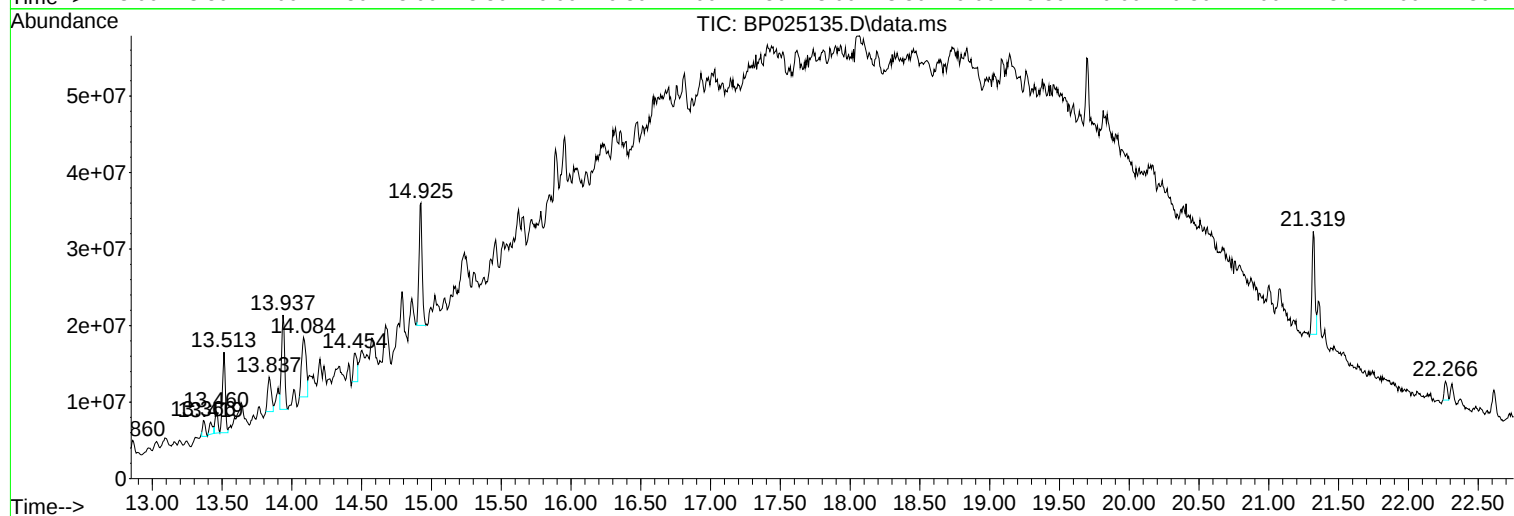
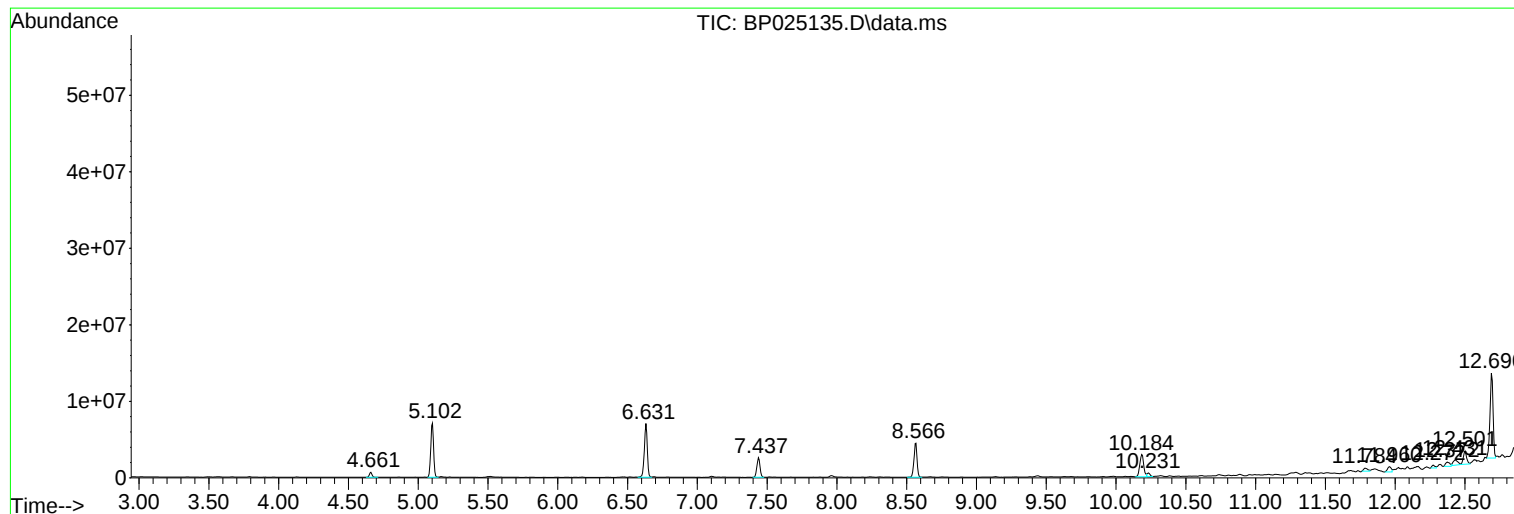
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP070325.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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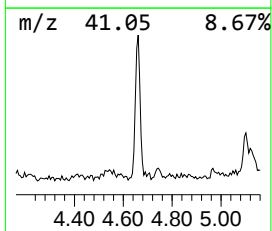
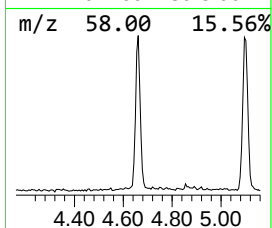
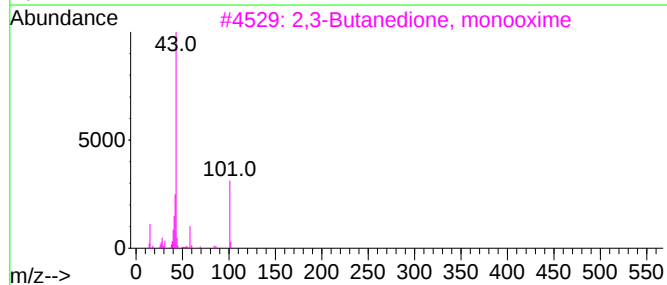
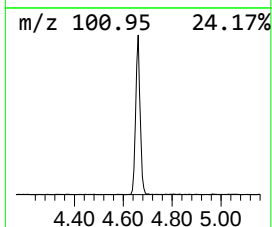
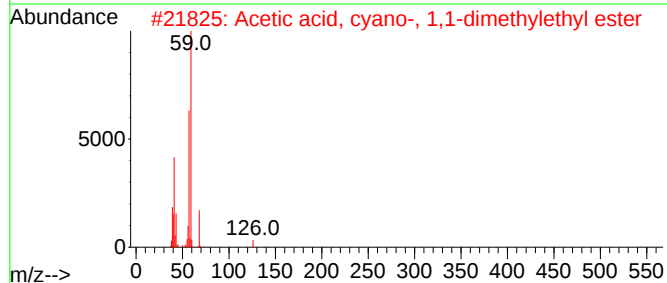
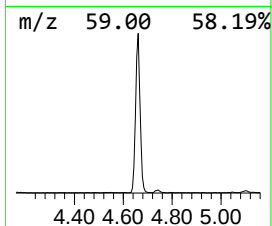
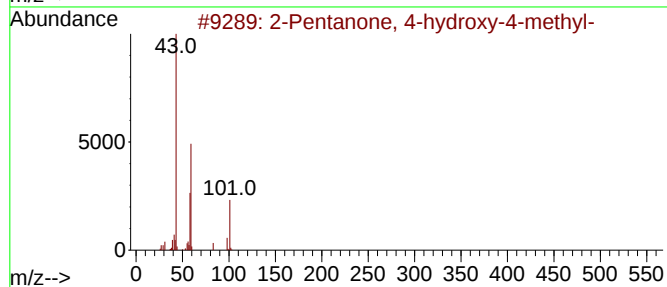
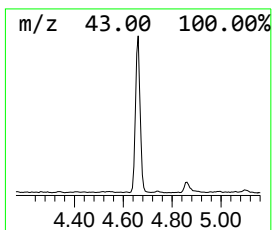
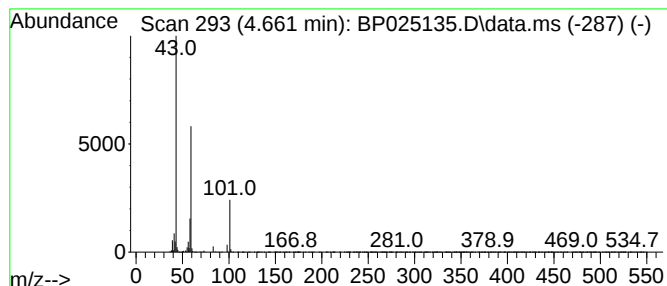
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.661	4.39 ng	871687	1,4-Dichlorobenzene-d4	7.437

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	47
2			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	17
4			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
5			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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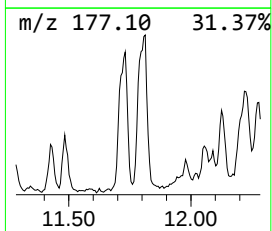
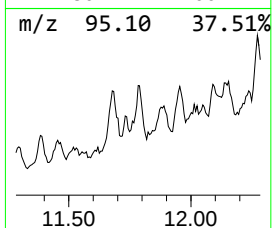
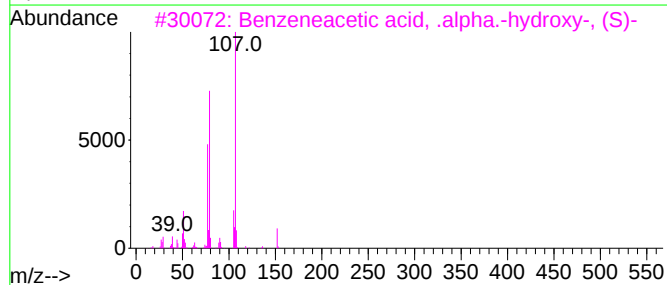
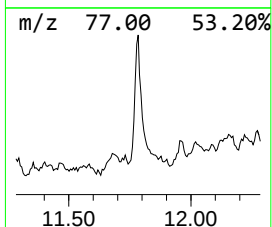
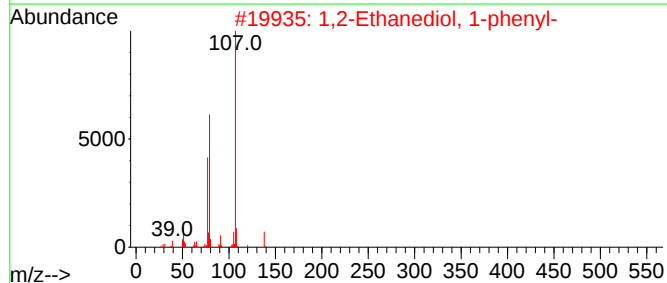
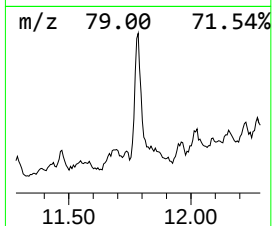
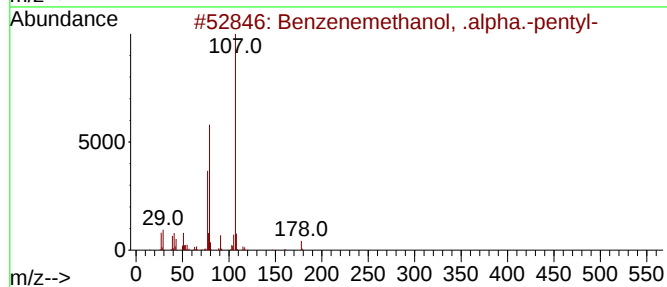
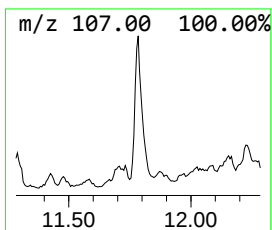
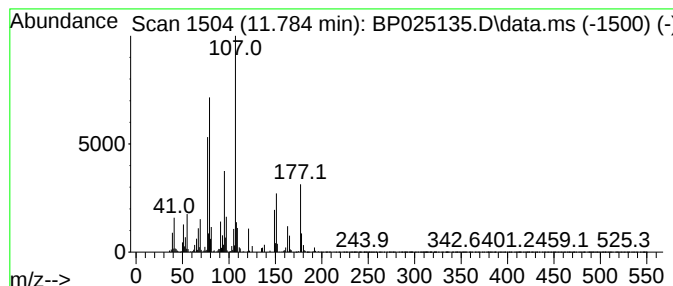
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TIC Library : C:\Database\NIST20.L
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Peak Number 2 Benzenemethanol, .alpha.-pe... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.784	2.87 ng	822967	Naphthalene-d8	10.184

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenemethanol, .alpha.-pentyl-	178	C12H18O	004471-05-0	60
2			1,2-Ethanediol, 1-phenyl-	138	C8H10O2	000093-56-1	60
3			Benzeneacetic acid, .alpha.-hydr...	152	C8H8O3	017199-29-0	53
4			Benzeneacetic acid, .alpha.-hydr...	152	C8H8O3	000611-71-2	49
5			Mandelic acid	152	C8H8O3	000090-64-2	49



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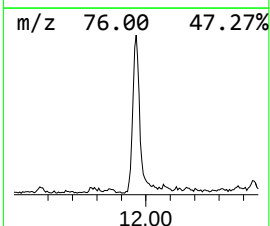
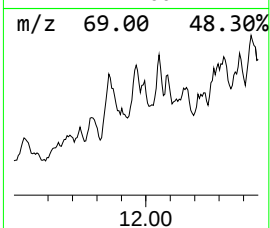
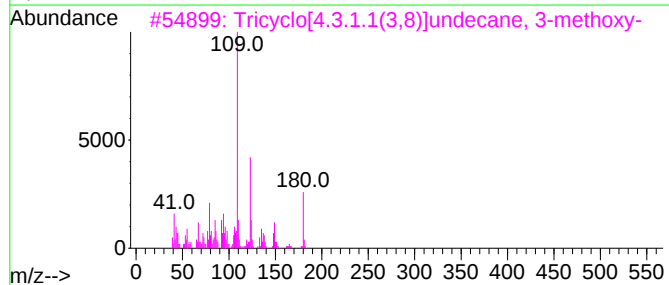
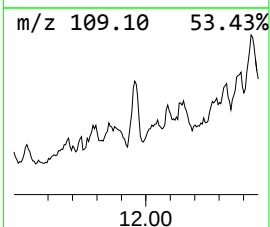
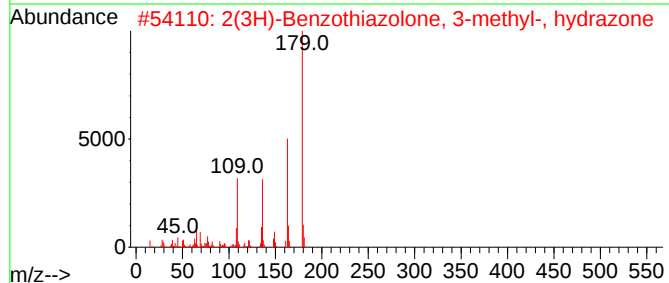
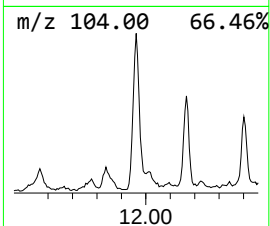
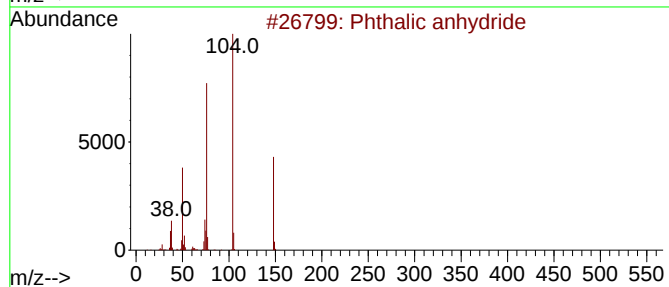
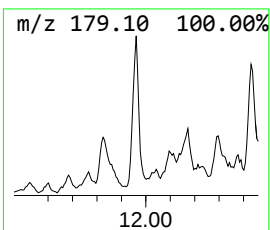
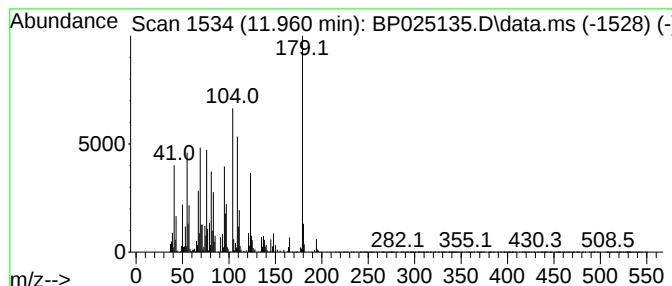
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TIC Library : C:\Database\NIST20.L
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Peak Number 3 unknown11.960 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.960	4.56 ng	1304780	Naphthalene-d8	10.184

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phthalic anhydride	148	C8H4O3	000085-44-9	40
2			2(3H)-Benzothiazolone, 3-methyl-...	179	C8H9N3S	001128-67-2	27
3			Tricyclo[4.3.1.1(3,8)]undecane, ...	180	C12H20O	021898-92-0	25
4			4,2-Cresotic acid, 6-methoxy-, b...	538	C29H30O10	019314-74-0	22
5			Carbamic acid, methylphenyl-, et...	179	C10H13NO2	002621-79-6	14



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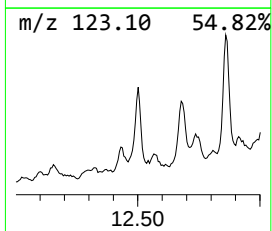
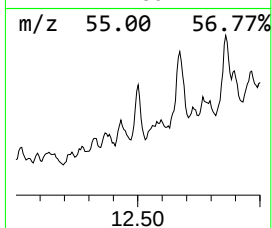
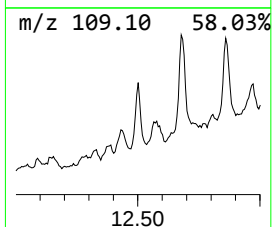
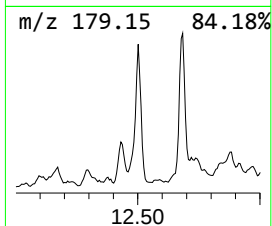
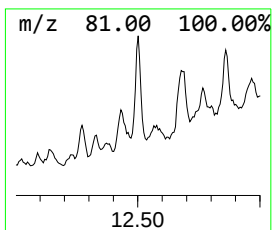
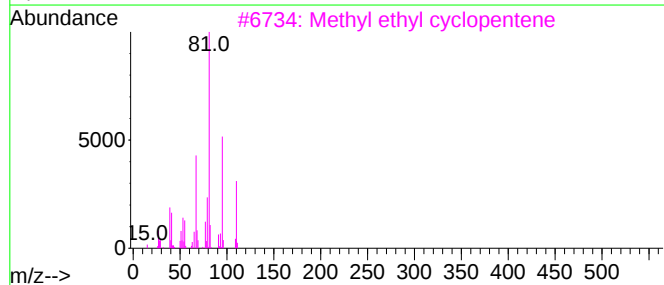
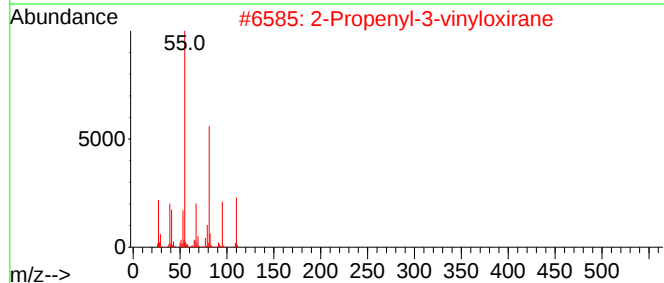
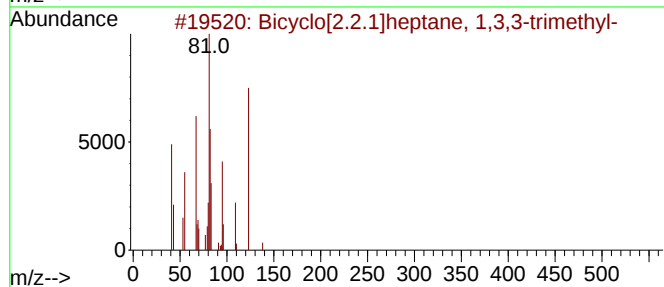
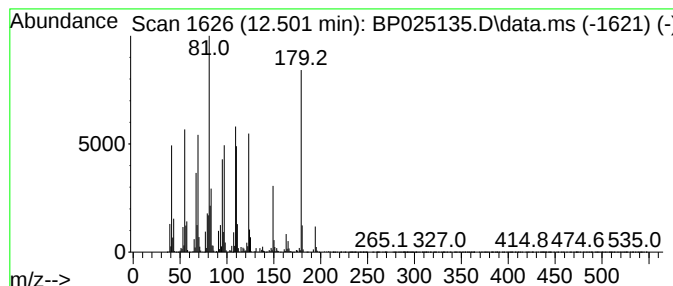
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Peak Number 4 unknown12.501 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.501	3.22 ng	2817130	Acenaphthene-d10	14.089

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bicyclo[2.2.1]heptane, 1,3,3-tri...	138	C10H18	006248-88-0	42
2			2-Propenyl-3-vinyloxirane	110	C7H10O	070597-49-8	27
3			Methyl ethyl cyclopentene	110	C8H14	019780-56-4	22
4			Cyclohexene, 3-ethyl-	110	C8H14	002808-71-1	18
5			Cyclohexene, 1-ethyl-	110	C8H14	001453-24-3	18



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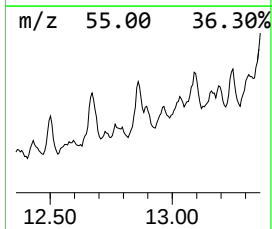
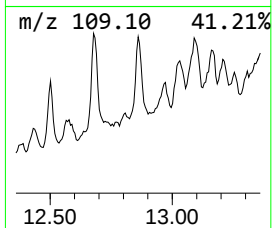
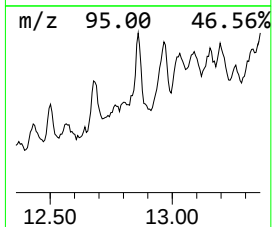
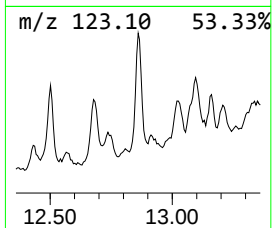
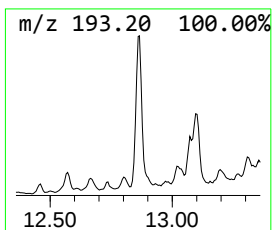
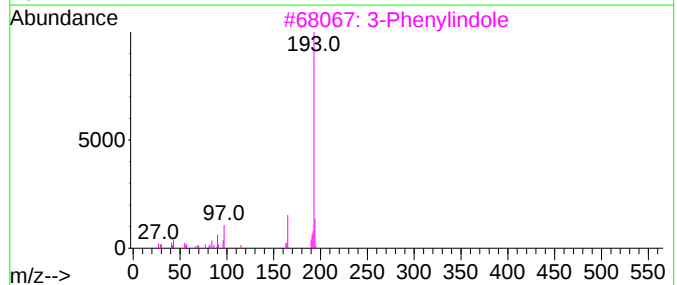
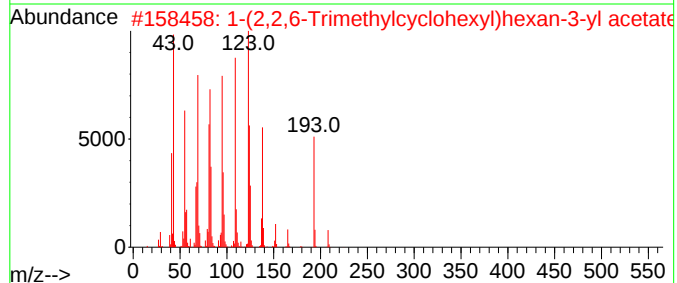
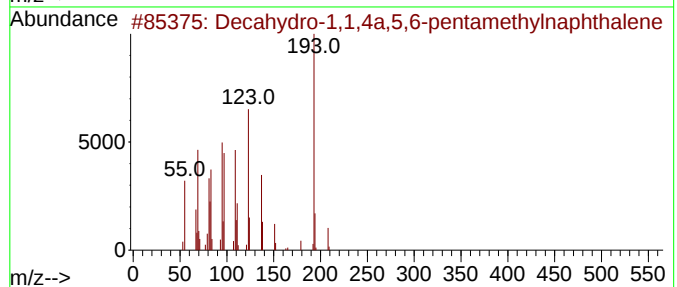
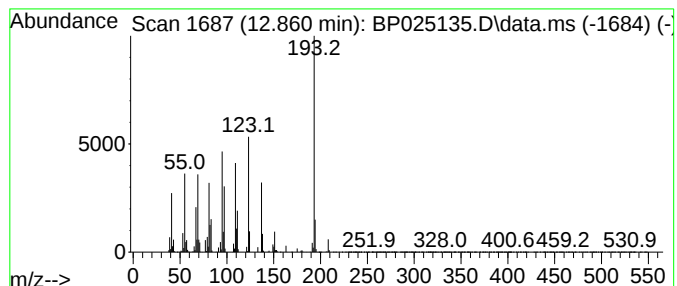
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Peak Number 5 Decahydro-1,1,4a,5,6-pentam... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.860	2.55 ng	2232700	Acenaphthene-d10	14.089

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decahydro-1,1,4a,5,6-pentamethyl...	208	C15H28	080655-44-3	64
2			1-(2,2,6-Trimethylcyclohexyl)hex...	268	C17H32O2	1000498-96-1	37
3			3-Phenylindole	193	C14H11N	001504-16-1	22
4			6-Methylphenanthridine	193	C14H11N	003955-65-5	22
5			2-Pentanone, 4-cyclohexylidene-3...	222	C15H26O	313253-65-5	22



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ARS20-0039

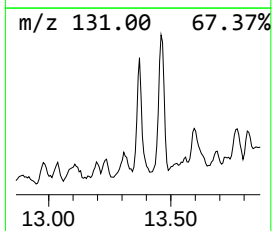
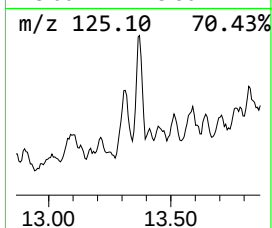
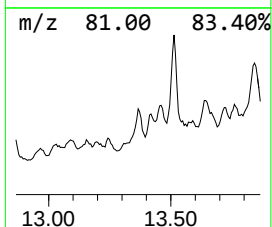
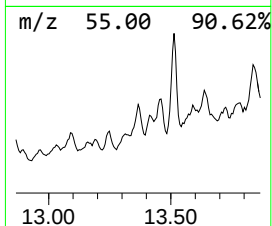
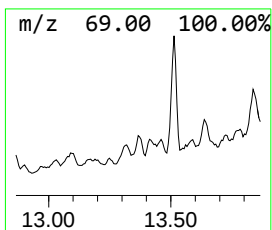
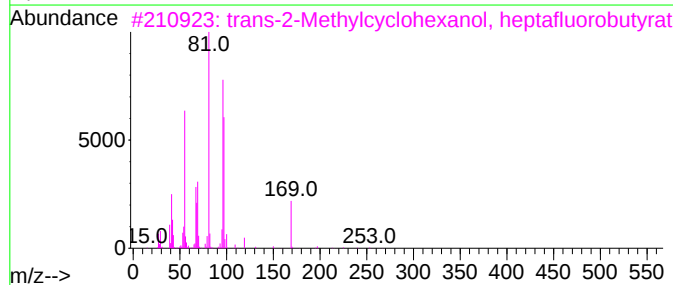
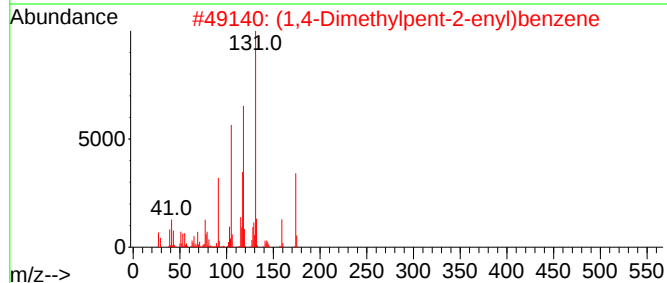
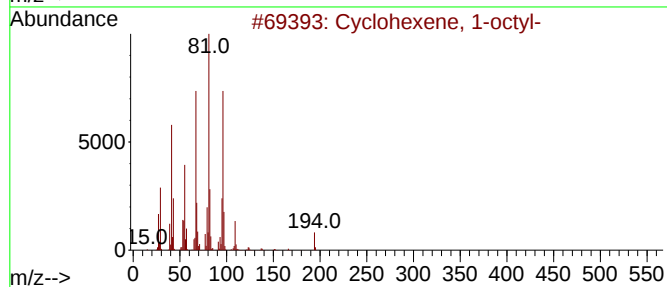
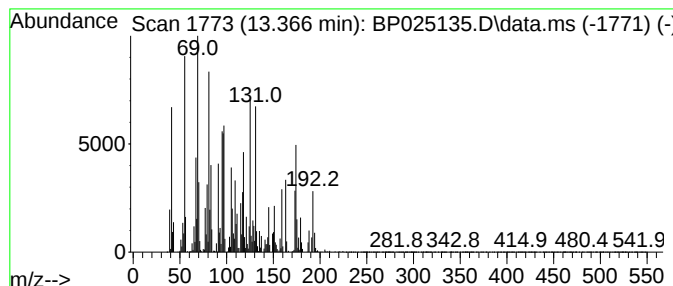
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Cyclohexene, 1-octyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.366	3.11 ng	2723030	Acenaphthene-d10	14.089

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexene, 1-octyl-	194	C14H26	015232-87-8	60
2			(1,4-Dimethylpent-2-enyl)benzene	174	C13H18	1000184-97-9	56
3			trans-2-Methylcyclohexanol, hept...	310	C11H13F7O2	1010364-13-5	22
4			Glycine, furfuryl ester	155	C7H9NO3	1000306-78-7	18
5			Cyclohexene, 3-(1-methylethyl)-	124	C9H16	003983-08-2	11



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071425\
Data File : BP025135.D
Acq On : 15 Jul 2025 01:22
Operator : CG/JU
Sample : Q2564-05
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ARS20-0039

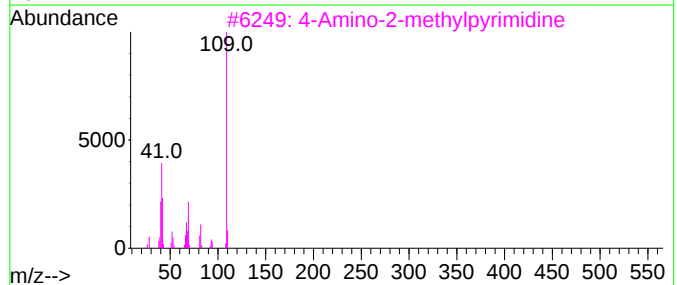
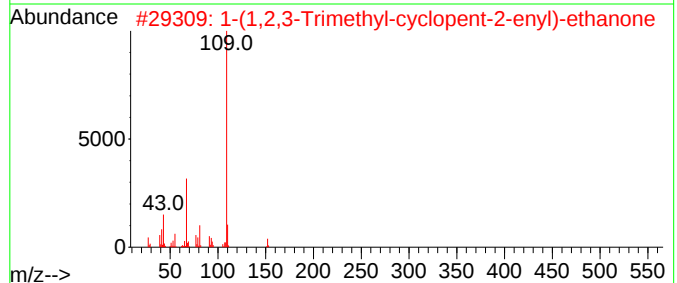
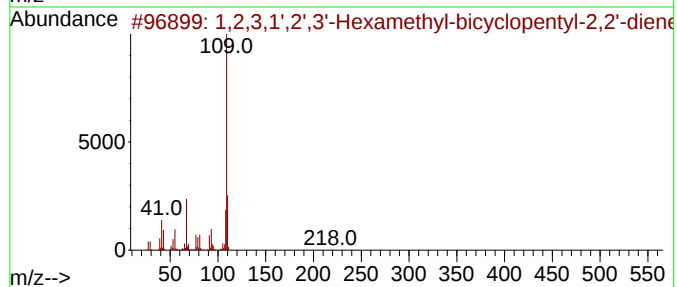
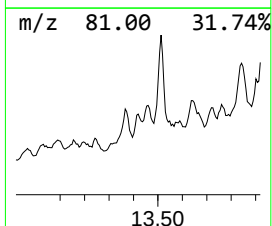
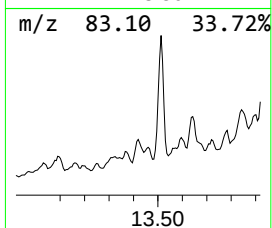
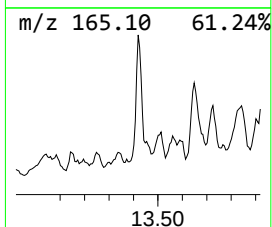
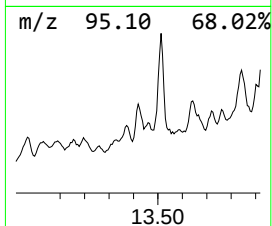
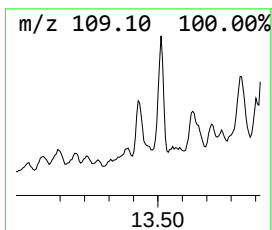
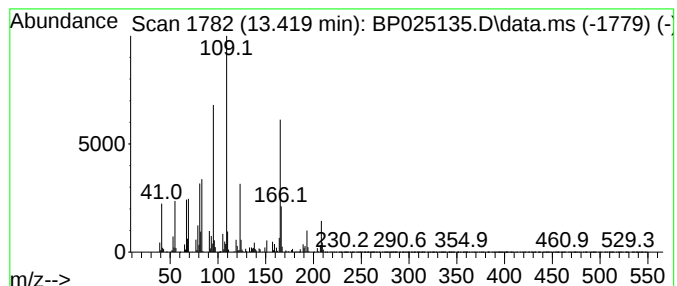
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown13.419 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.419	3.13 ng	2736080	Acenaphthene-d10	14.089

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2,3,1',2',3'-Hexamethyl-bicycl...	218	C16H26	051999-35-0	30		
2	1-(1,2,3-Trimethyl-cyclopent-2-e...	152	C10H16O	070987-81-4	27		
3	4-Amino-2-methylpyrimidine	109	C5H7N3	000074-69-1	25		
4	Cyclopentane, 2-ethylidene-1,1-d...	124	C9H16	056324-66-4	25		
5	Benzenamine, 4-(hexyloxy)-	193	C12H19NO	039905-57-2	25		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071425\
Data File : BP025135.D
Acq On : 15 Jul 2025 01:22
Operator : CG/JU
Sample : Q2564-05
Misc :
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Instrument :
BNA_P
ClientSampleId :
ARS20-0039

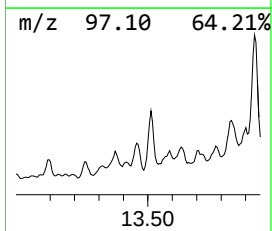
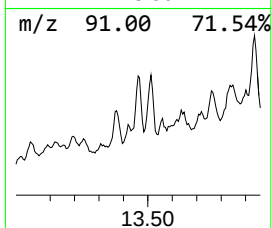
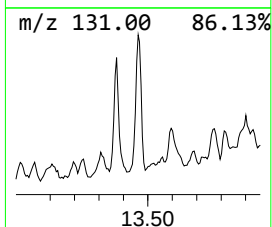
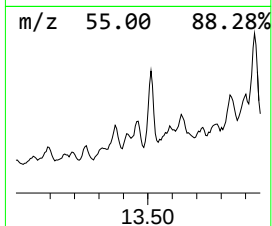
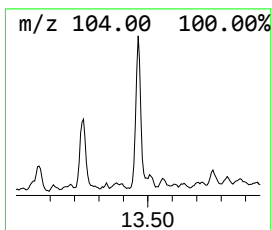
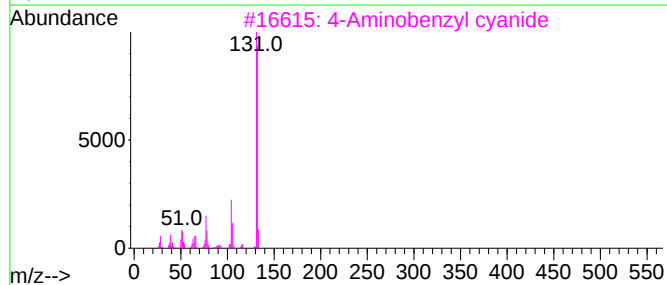
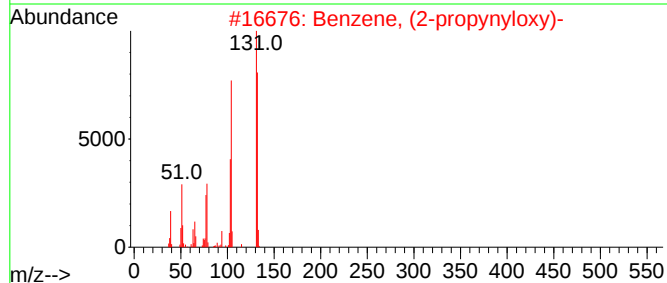
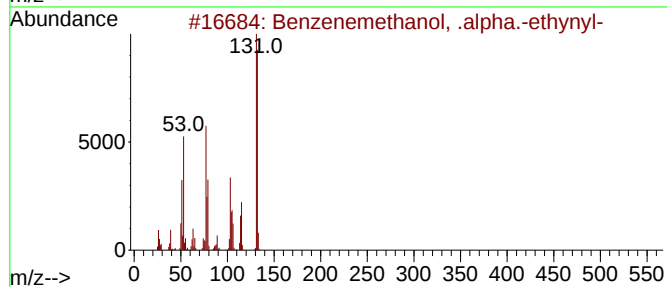
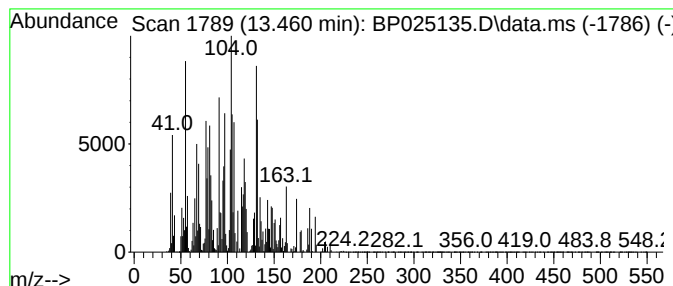
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Benzenemethanol, .alpha.-et... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.460	4.31 ng	3775900	Acenaphthene-d10	14.089

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenemethanol, .alpha.-ethynyl-	132	C9H8O	004187-87-5	81
2			Benzene, (2-propynyloxy)-	132	C9H8O	013610-02-1	55
3			4-Aminobenzyl cyanide	132	C8H8N2	003544-25-0	46
4			Naphthalene, 2-(1,1-dimethylethy...	188	C14H20	042044-22-4	46
5			1H-Indene-1,2-diol, 2,3-dihydro-	150	C9H10O2	004370-02-9	45



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071425\
Data File : BP025135.D
Acq On : 15 Jul 2025 01:22
Operator : CG/JU
Sample : Q2564-05
Misc :
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Instrument :
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ClientSampleId :
ARS20-0039

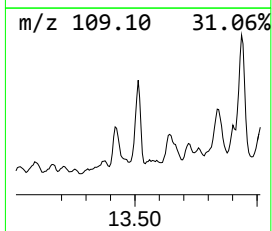
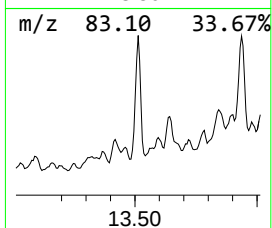
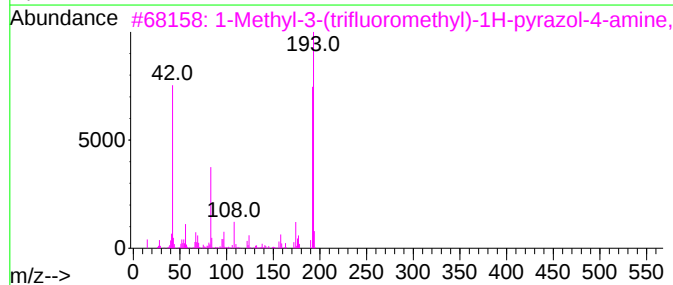
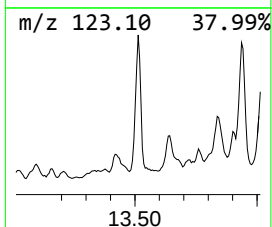
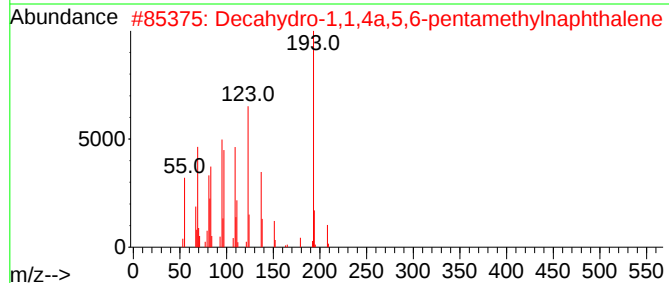
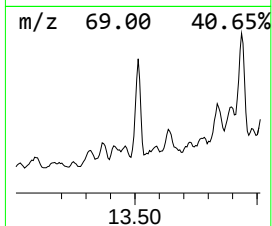
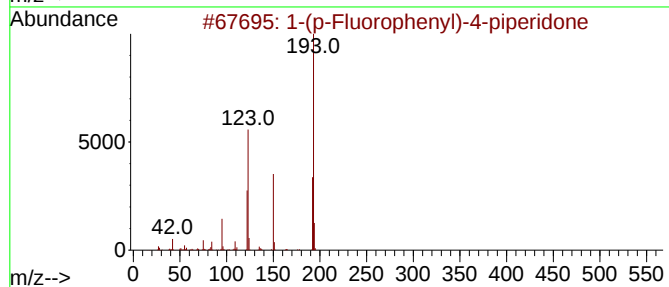
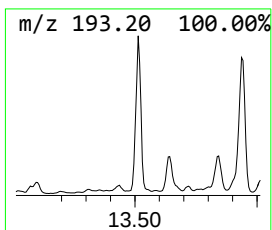
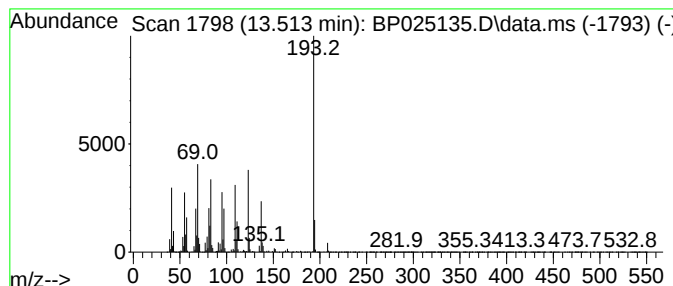
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 unknown13.513 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.513	17.16 ng	15018700	Acenaphthene-d10	14.089

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-(p-Fluorophenyl)-4-piperidone	193	C11H12FNO	1000238-56-7	47
2			Decahydro-1,1,4a,5,6-pentamethyl...	208	C15H28	080655-44-3	43
3			1-Methyl-3-(trifluoromethyl)-1H...	193	C7H10F3N3	1000511-73-1	38
4			1H-Indene, octahydro-2,2,4,4,7,7...	208	C15H28	054832-83-6	38
5			Borinic acid, diethyl-, 3,3,5-tr...	208	C13H25BO	057387-76-5	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071425\
Data File : BP025135.D
Acq On : 15 Jul 2025 01:22
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Sample : Q2564-05
Misc :
ALS Vial : 16 Sample Multiplier: 1

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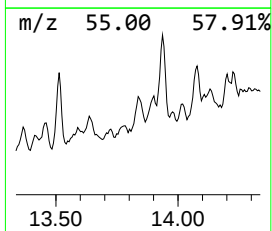
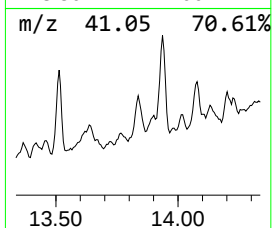
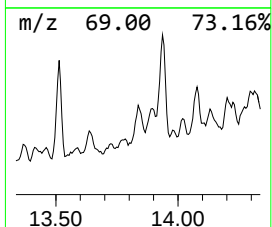
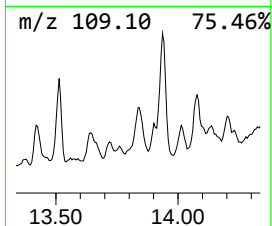
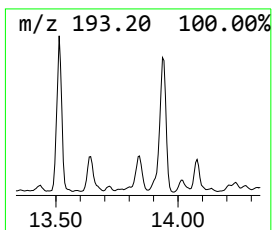
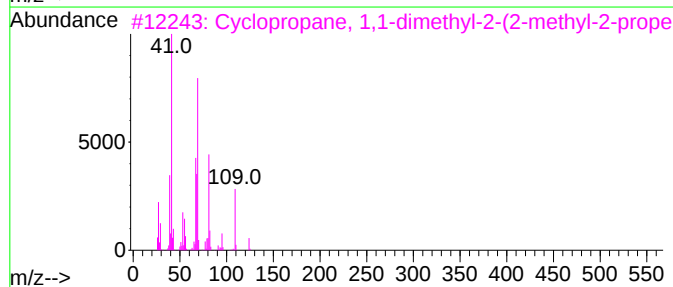
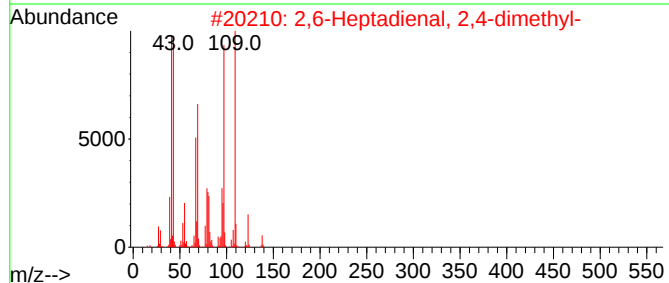
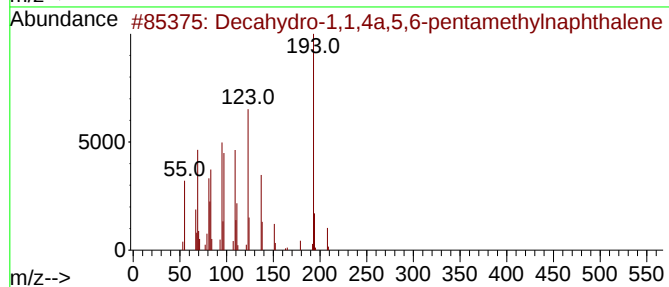
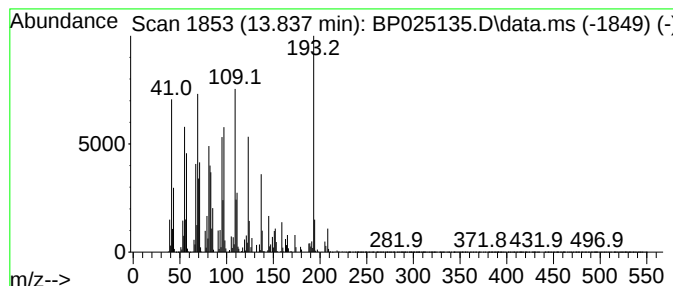
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown13.837 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.837	9.30 ng	8140760	Acenaphthene-d10	14.089

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decahydro-1,1,4a,5,6-pentamethyl...	208	C15H28	080655-44-3	60
2			2,6-Heptadienal, 2,4-dimethyl-	138	C9H14O	085136-08-9	38
3			Cyclopropane, 1,1-dimethyl-2-(2-...	124	C9H16	069147-03-1	30
4			4-Amino-2-methylpyrimidine	109	C5H7N3	000074-69-1	25
5			1a,2,5,5-Tetramethyl-cis-1a,4a,5...	194	C13H22O	1000215-77-7	15



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071425\
Data File : BP025135.D
Acq On : 15 Jul 2025 01:22
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Misc :
ALS Vial : 16 Sample Multiplier: 1

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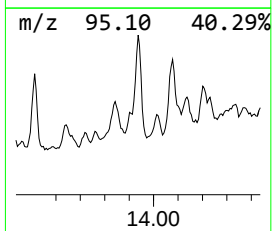
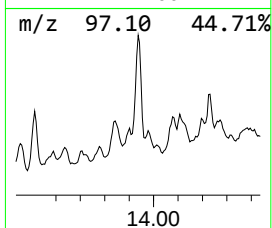
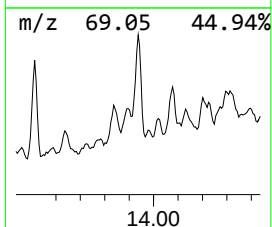
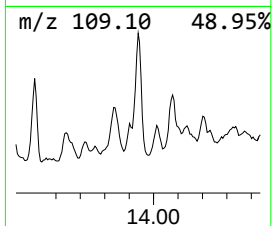
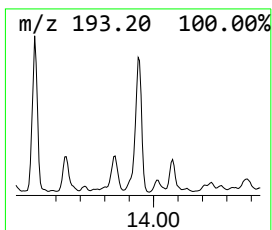
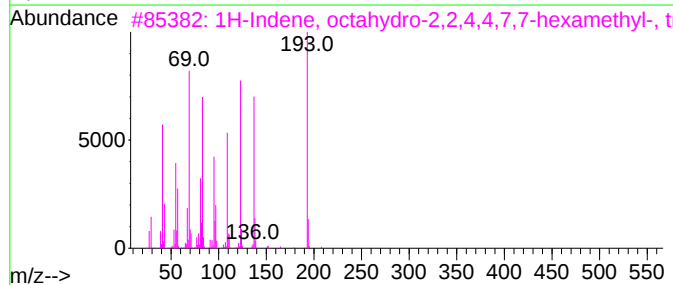
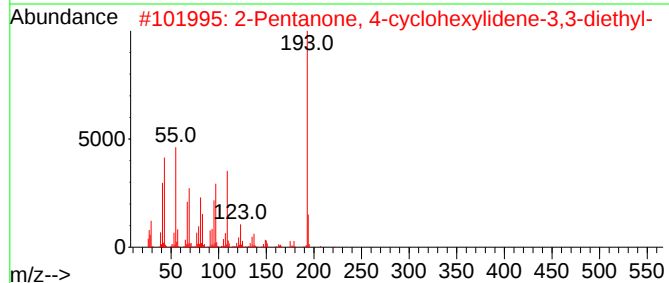
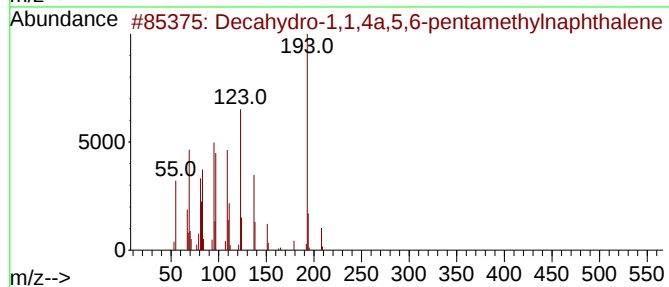
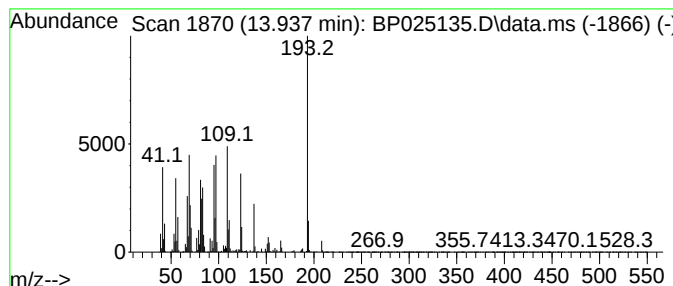
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown13.937 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.937	22.12 ng	19366100	Acenaphthene-d10	14.089

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decahydro-1,1,4a,5,6-pentamethyl...	208	C15H28	080655-44-3	49
2			2-Pentanone, 4-cyclohexylidene-3...	222	C15H26O	313253-65-5	45
3			1H-Indene, octahydro-2,2,4,4,7,7...	208	C15H28	054832-83-6	42
4			Butan-2-one, 1-[3-(3,3-dimethyl-...	278	C16H26N2O2	1000303-34-2	27
5			Borinic acid, diethyl-, 3,3,5-tr...	208	C13H25BO	057387-76-5	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071425\
Data File : BP025135.D
Acq On : 15 Jul 2025 01:22
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Misc :
ALS Vial : 16 Sample Multiplier: 1

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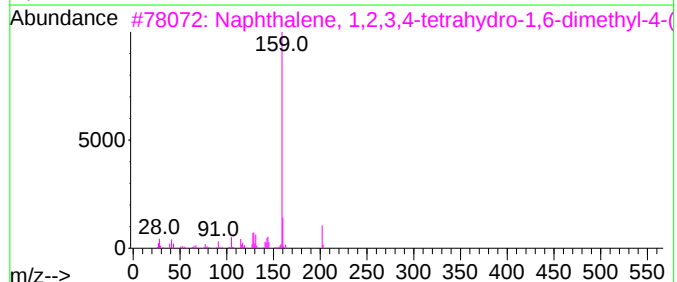
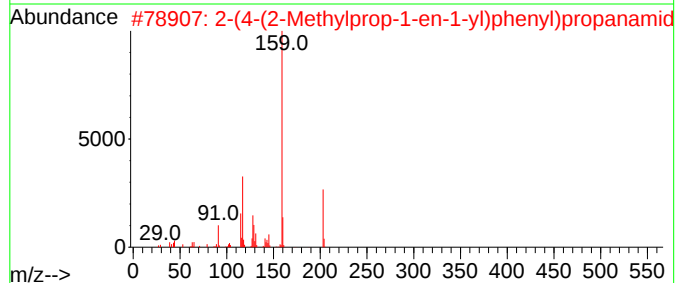
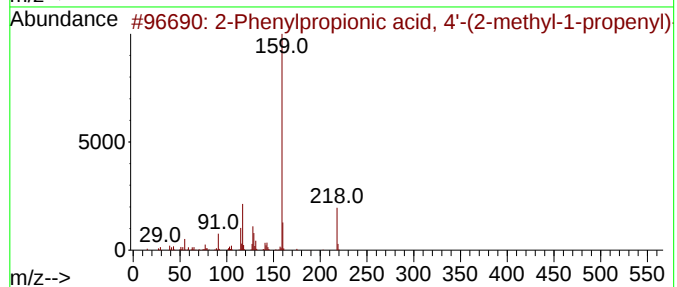
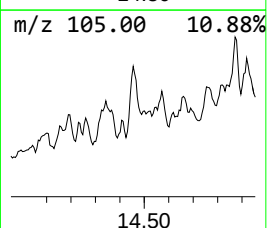
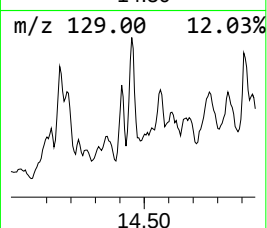
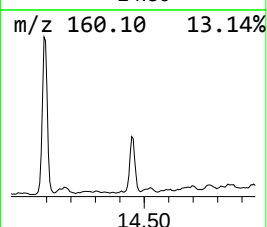
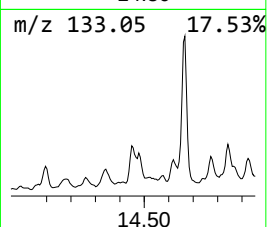
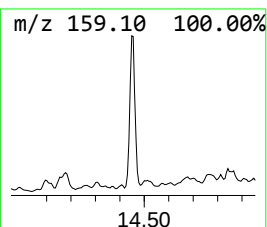
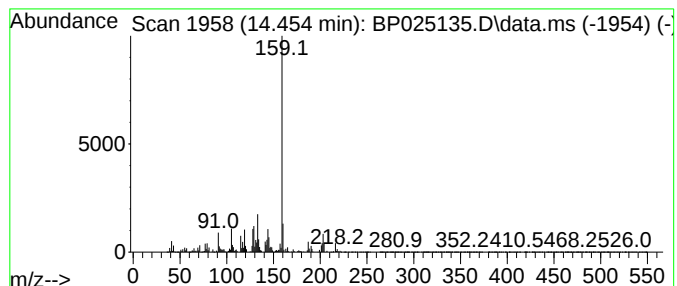
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 2-Phenylpropionic acid, 4'-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.454	7.95 ng	6958940	Acenaphthene-d10	14.089

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Phenylpropionic acid, 4'-(2-me...	218	C14H18O2	1010454-94-3	86
2			2-(4-(2-Methylprop-1-en-1-yl)phe...	203	C13H17NO	1000466-09-8	76
3			Naphthalene, 1,2,3,4-tetrahydro-...	202	C15H22	000483-77-2	76
4			Naphthalene, 1,2,3,4-tetrahydro-...	174	C13H18	030316-36-0	64
5			Naphthalene, 1,2,3,4-tetrahydro-...	174	C13H18	021693-51-6	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071425\
Data File : BP025135.D
Acq On : 15 Jul 2025 01:22
Operator : CG/JU
Sample : Q2564-05
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ARS20-0039

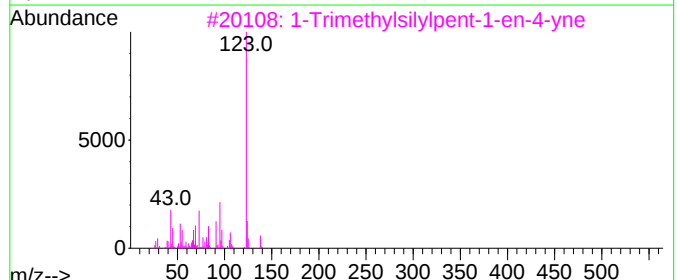
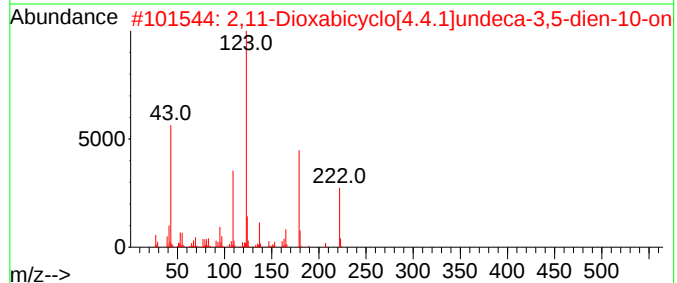
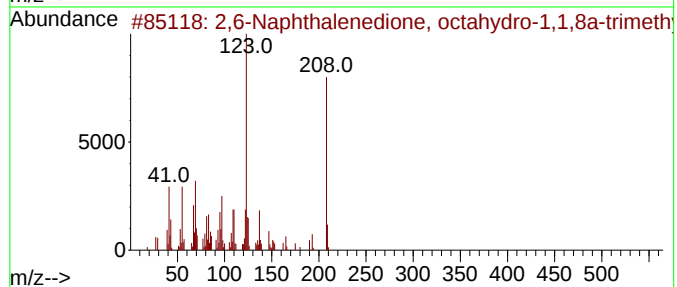
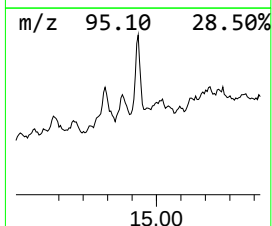
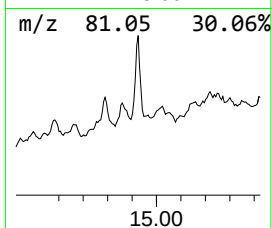
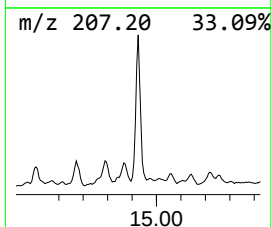
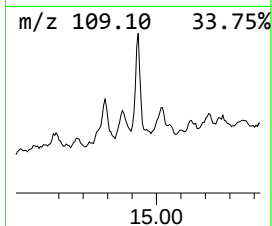
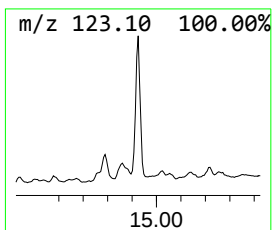
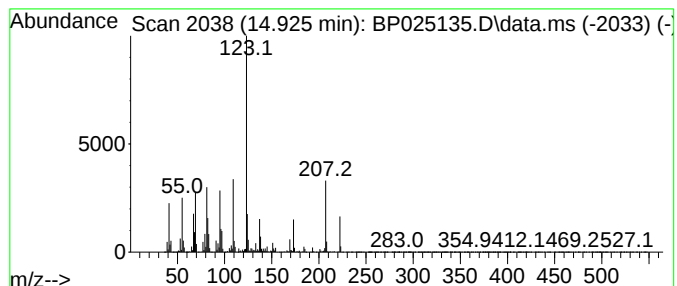
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 2,6-Naphthalenedione, octah... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.925	28.71 ng	25133100	Acenaphthene-d10	14.089

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,6-Naphthalenedione, octahydro-...	208	C13H20O2	057289-16-4	50		
2	2,11-Dioxabicyclo[4.4.1]undeca-3...	222	C13H18O3	070412-52-1	50		
3	1-Trimethylsilylpent-1-en-4-yne	138	C8H14Si	079516-25-9	47		
4	Benzamide, 3-fluoro-N-(2-phenyle...	313	C20H24FNO	1000451-30-5	47		
5	(Phenylthio)acetic acid, cyclobu...	222	C12H14O2S	1010299-95-5	43		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071425\
Data File : BP025135.D
Acq On : 15 Jul 2025 01:22
Operator : CG/JU
Sample : Q2564-05
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ARS20-0039

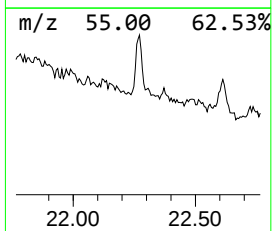
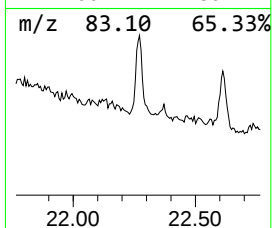
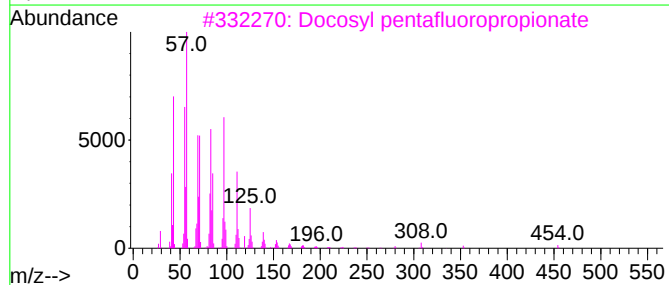
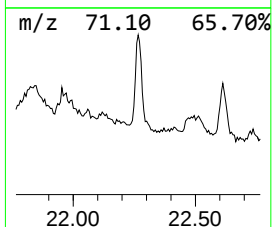
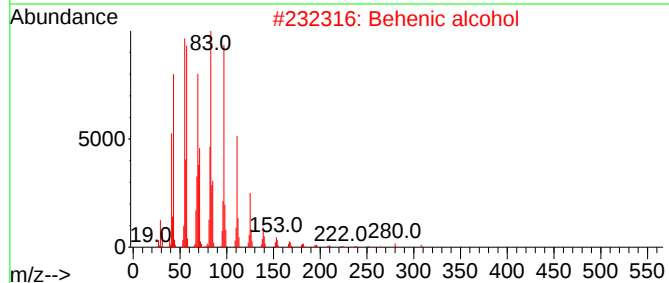
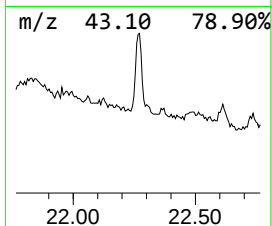
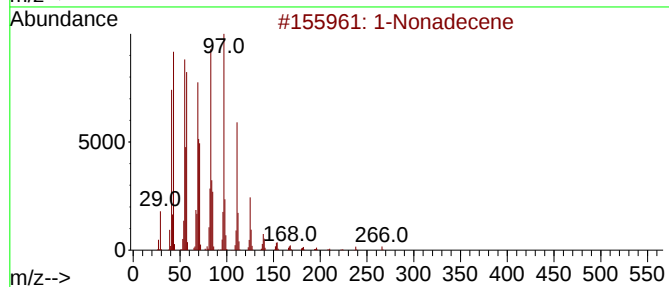
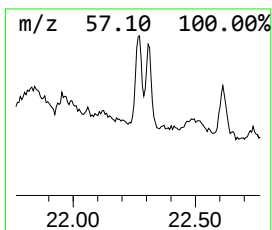
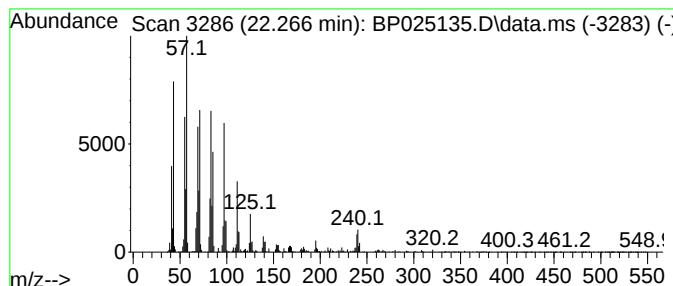
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 1-Nonadecene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.266	3.59 ng	3529190	Chrysene-d12	21.360

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Nonadecene	266	C19H38	018435-45-5	97
2		Behenic alcohol	326	C22H46O	000661-19-8	93
3		Docosyl pentafluoropropionate	472	C25H45F5O2	1000351-80-9	91
4		Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	90
5		Oxalic acid, hexadecyl propyl ester	356	C21H40O4	1000309-26-9	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071425\
Data File : BP025135.D
Acq On : 15 Jul 2025 01:22
Operator : CG/JU
Sample : Q2564-05
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ARS20-0039

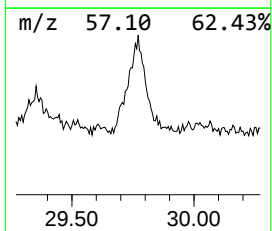
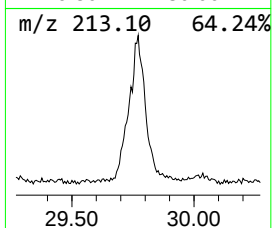
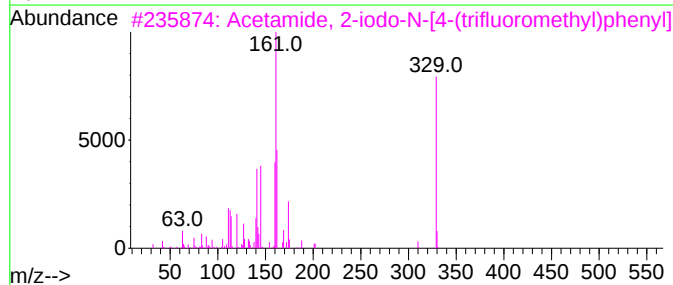
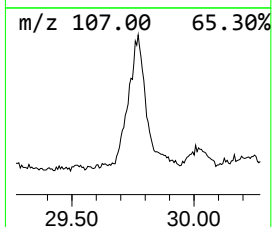
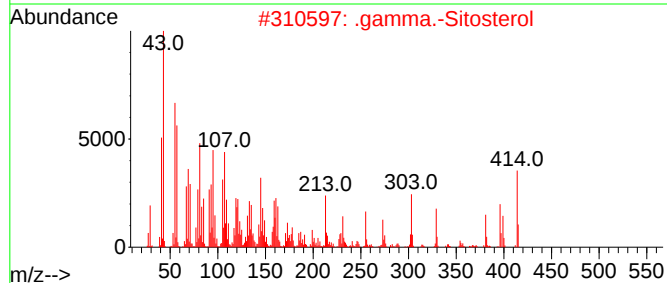
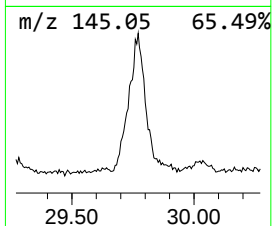
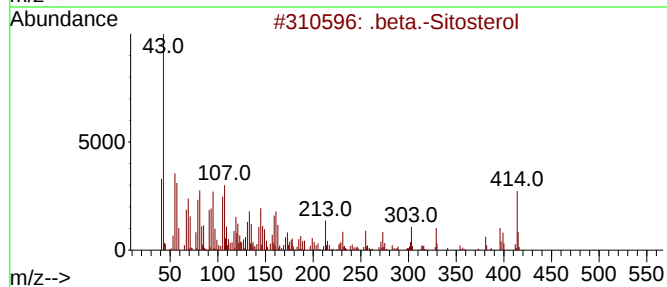
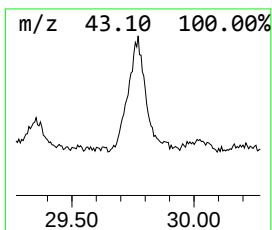
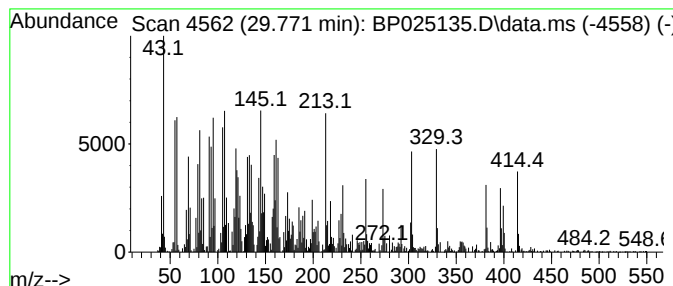
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 .beta.-Sitosterol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
29.771	36.33 ng	6409970	Perylene-d12	24.483

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			.beta.-Sitosterol	414	C29H50O	000083-46-5	78
2			.gamma.-Sitosterol	414	C29H50O	000083-47-6	64
3			Acetamide, 2-iodo-N-[4-(trifluor...	329	C9H7F3INO	1000362-86-0	11
4			(3S,8S,9S,10R,13R,14S,17R)-17-(...	414	C29H50O	029365-29-5	10
5			(3S,8S,9S,10R,13R,14S,17R)-17-(...	428	C30H52O	020194-50-7	10



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071425\
Data File : BP025135.D
Acq On : 15 Jul 2025 01:22
Operator : CG/JU
Sample : Q2564-05
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ARS20-0039

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP070325.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
2-Pentanone, 4-...	4.661	4.4	ng	871687	1	7.437	3968190	20.0
Benzenemethanol...	11.784	2.9	ng	822967	2	10.184	5725730	20.0
unknown11.960	11.960	4.6	ng	1304780	2	10.184	5725730	20.0
unknown12.501	12.501	3.2	ng	2817130	3	14.089	17507100	20.0
Decahydro-1,1,4...	12.860	2.5	ng	2232700	3	14.089	17507100	20.0
Cyclohexene, 1-...	13.366	3.1	ng	2723030	3	14.089	17507100	20.0
unknown13.419	13.419	3.1	ng	2736080	3	14.089	17507100	20.0
Benzenemethanol...	13.460	4.3	ng	3775900	3	14.089	17507100	20.0
unknown13.513	13.513	17.2	ng	15018700	3	14.089	17507100	20.0
unknown13.837	13.837	9.3	ng	8140760	3	14.089	17507100	20.0
unknown13.937	13.937	22.1	ng	19366100	3	14.089	17507100	20.0
2-Phenylpropion...	14.454	8.0	ng	6958940	3	14.089	17507100	20.0
2,6-Naphthalene...	14.925	28.7	ng	25133100	3	14.089	17507100	20.0
1-Nonadecene	22.266	3.6	ng	3529190	5	21.360	19683400	20.0
.beta.-Sitosterol	29.771	36.3	ng	6409970	6	24.483	3529190	20.0