

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050225\
 Data File : BP024516.D
 Acq On : 02 May 2025 19:44
 Operator : RC/JU
 Sample : Q1932-03
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 COMP-5

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-BP041425.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP024516.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.875	299	304	317	rVB	121391	173103	4.03%	0.545%
2	5.328	375	381	401	rBV	1352330	2000644	46.55%	6.293%
3	6.893	638	647	667	rBV	1225933	2091391	48.66%	6.579%
4	7.704	779	785	794	rBV	549444	865985	20.15%	2.724%
5	8.851	973	980	998	rBV	748126	1317509	30.66%	4.144%
6	10.475	1248	1256	1272	rBV	687887	1244951	28.97%	3.916%
7	12.763	1640	1645	1651	rVB5	32690	51791	1.21%	0.163%
8	12.951	1668	1677	1686	rBV	1707432	2743871	63.85%	8.631%
9	13.110	1700	1704	1714	rVB2	61549	104897	2.44%	0.330%
10	13.757	1809	1814	1822	rBV2	245694	374616	8.72%	1.178%
11	13.875	1830	1834	1842	rVB3	65658	124994	2.91%	0.393%
12	14.080	1864	1869	1874	rBV4	127574	252091	5.87%	0.793%
13	14.133	1875	1878	1880	rVV2	75260	90749	2.11%	0.285%
14	14.169	1880	1884	1891	rVB	408787	549194	12.78%	1.728%
15	14.245	1893	1897	1902	rBV4	73350	140653	3.27%	0.442%
16	14.310	1903	1908	1909	rBV	204068	278510	6.48%	0.876%
17	14.345	1909	1914	1922	rVB	1127989	1768289	41.15%	5.562%
18	14.433	1925	1929	1934	rBV3	86932	159203	3.70%	0.501%
19	14.669	1965	1969	1973	rBV2	101810	197280	4.59%	0.621%
20	14.880	2001	2005	2008	rBV2	161506	250735	5.83%	0.789%
21	15.010	2024	2027	2032	rVB2	187770	262267	6.10%	0.825%
22	15.074	2033	2038	2044	rVV4	180508	313605	7.30%	0.986%
23	15.145	2045	2050	2058	rBV2	452457	832315	19.37%	2.618%
24	15.863	2167	2172	2179	rBV	1853130	3152775	73.36%	9.918%
25	17.151	2387	2391	2397	rBV	1406395	2096720	48.79%	6.596%
26	18.092	2548	2551	2555	rVB3	490547	562435	13.09%	1.769%
27	19.868	2849	2853	2860	rVB	3271765	4297593	100.00%	13.519%
28	21.592	3141	3146	3152	rVB	1565681	2341775	54.49%	7.366%
29	21.762	3172	3175	3185	rVB	266948	449945	10.47%	1.415%
30	24.950	3707	3717	3730	rVB	907531	2699784	62.82%	8.493%

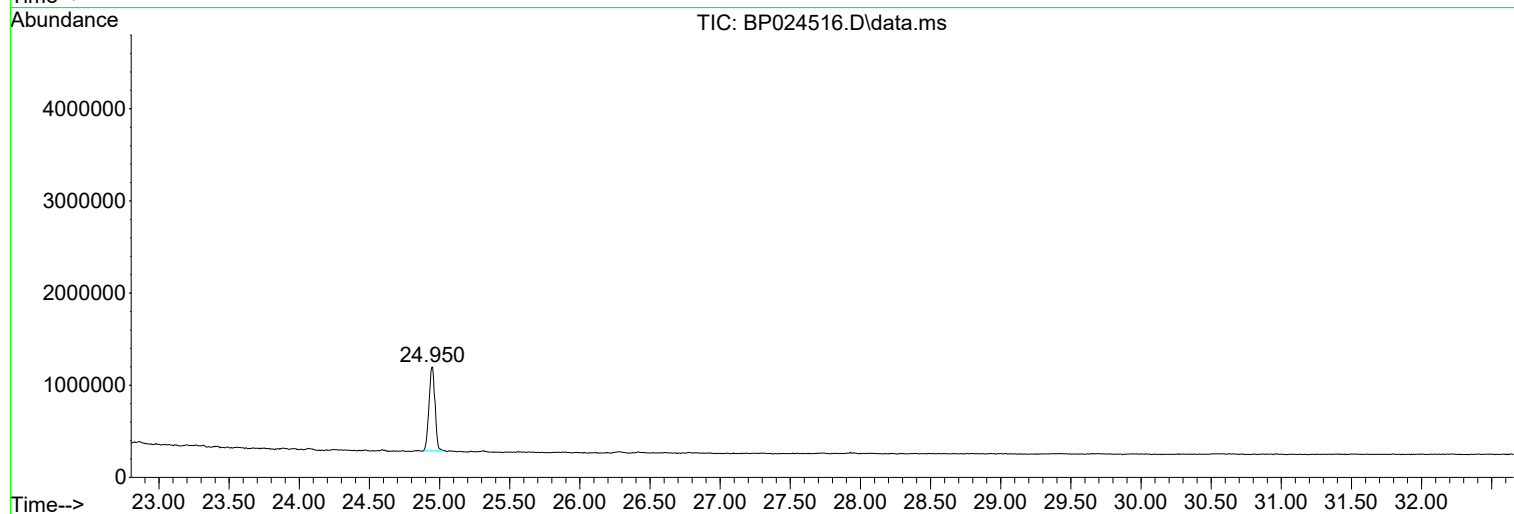
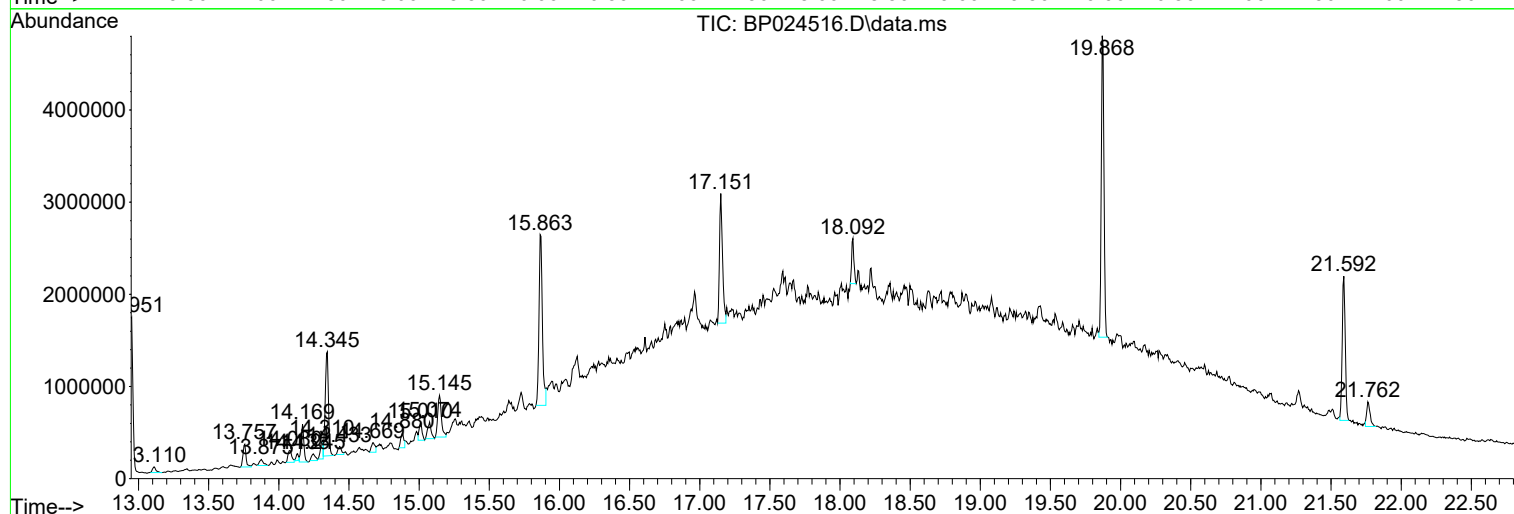
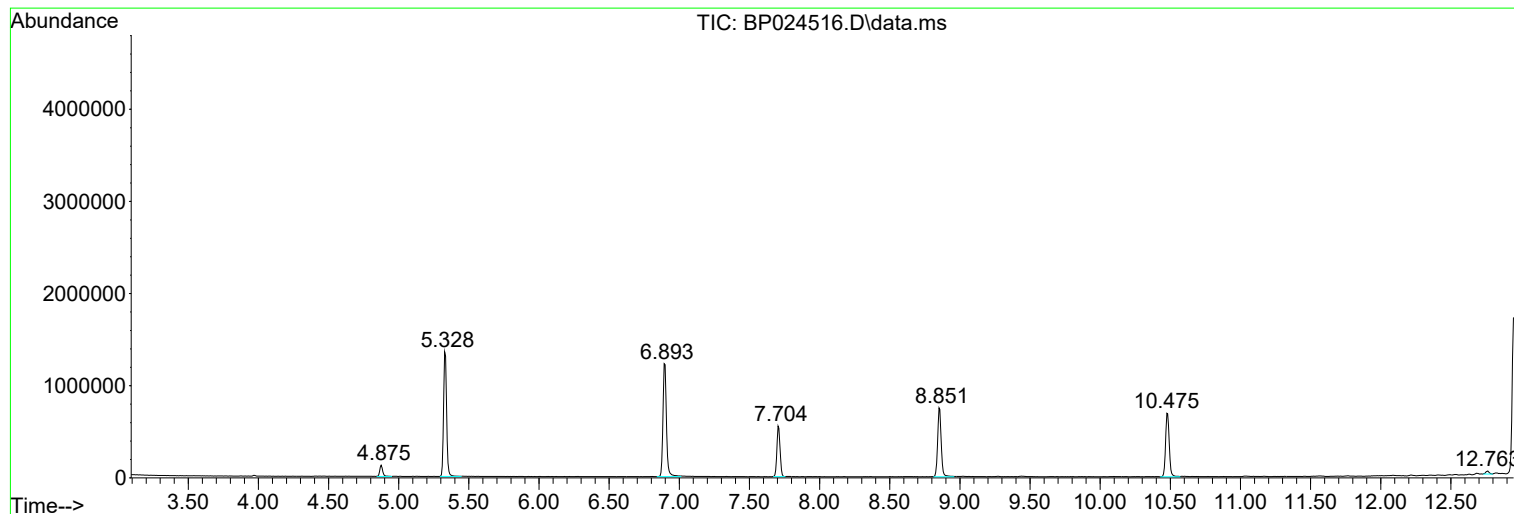
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP041425.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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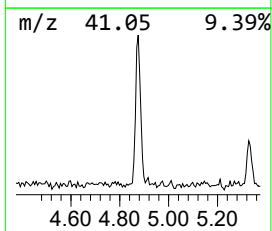
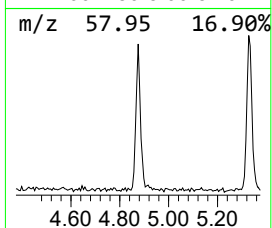
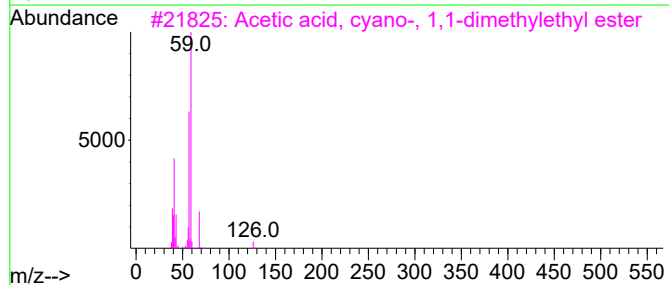
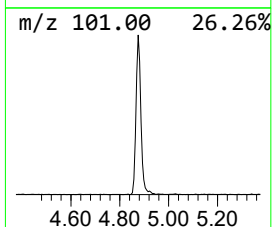
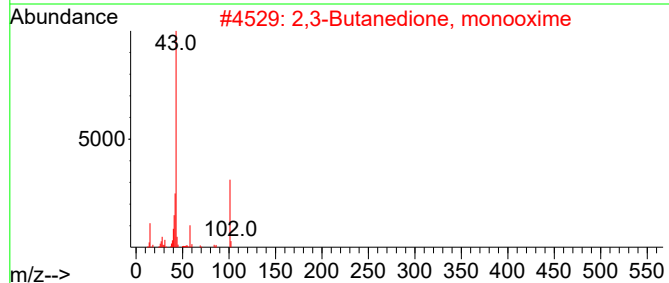
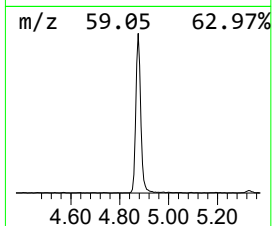
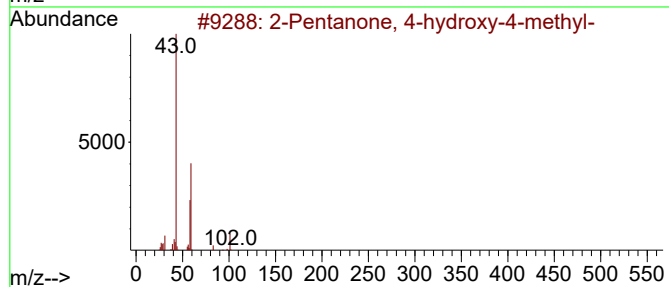
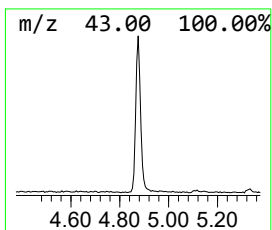
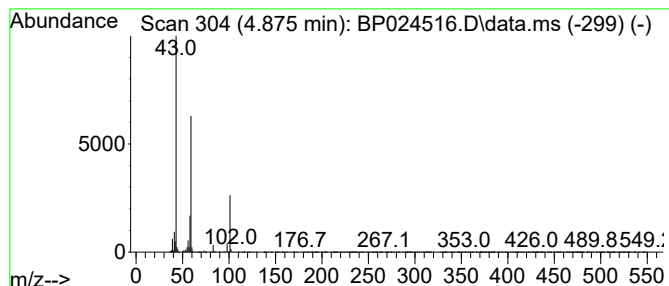
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TIC Library : C:\Database\NIST20.L
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.875	4.00 ng	173103	1,4-Dichlorobenzene-d4	7.704

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56		
2	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25		
3	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17		
4	Acetone	58	C3H6O	000067-64-1	10		
5	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10		



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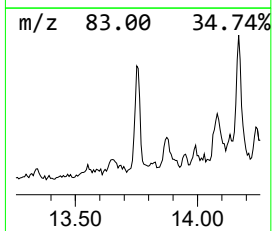
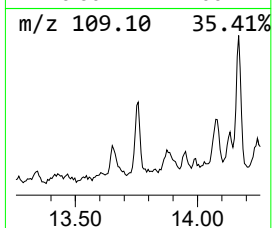
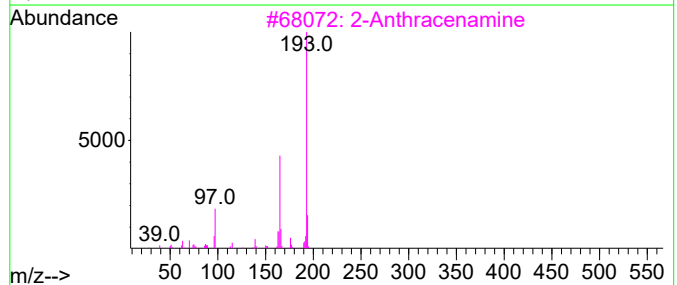
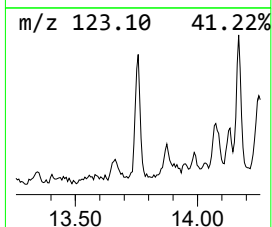
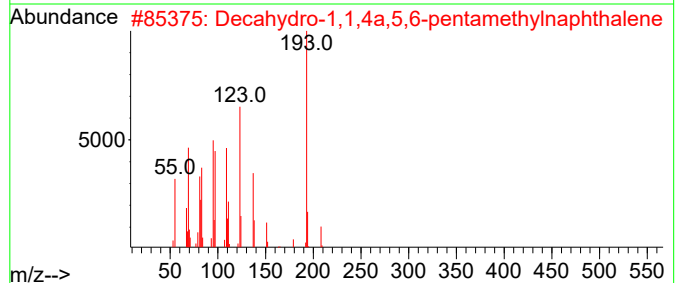
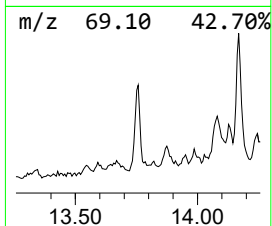
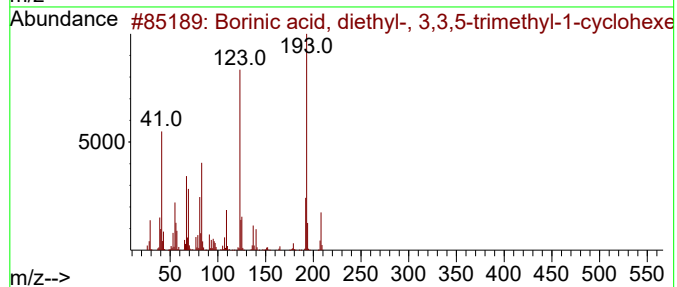
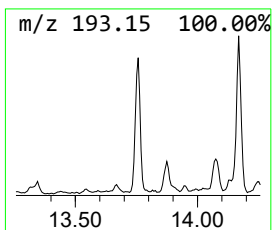
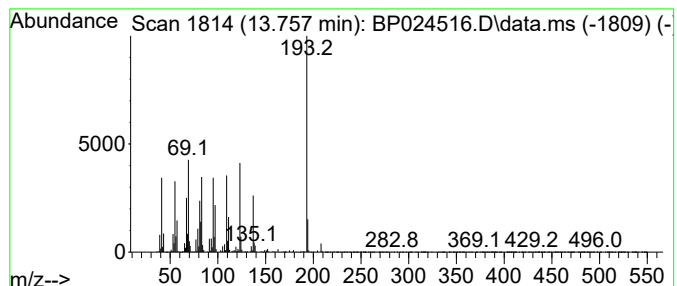
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Peak Number 2 Borinic acid, diethyl-, 3,3... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.757	4.24 ng	374616	Acenaphthene-d10	14.345

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Borinic acid, diethyl-, 3,3,5-tr...	208	C13H25BO	057387-76-5	50
2			Decahydro-1,1,4a,5,6-pentamethyl...	208	C15H28	080655-44-3	38
3			2-Anthracenamine	193	C14H11N	000613-13-8	38
4			2-Pentanone, 4-cyclohexylidene-3...	222	C15H26O	313253-65-5	37
5			1H-Pyrazole, 3,5-bis(1,1-dimethy...	208	C13H24N2	125281-21-2	35



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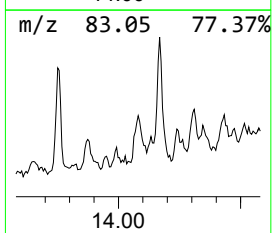
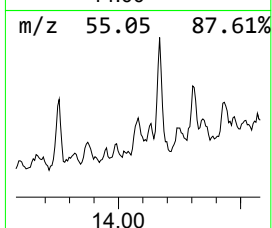
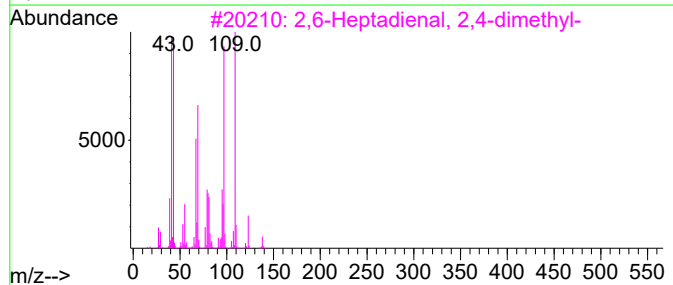
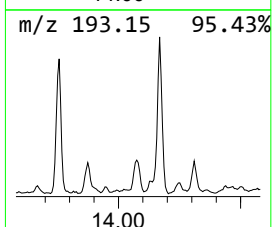
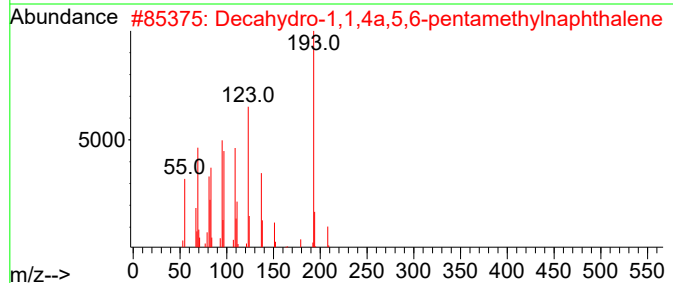
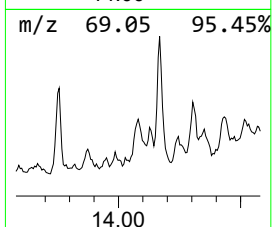
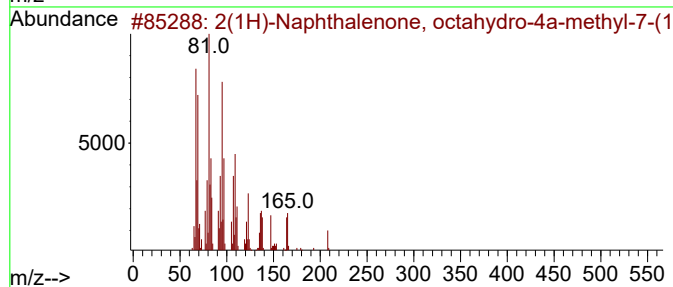
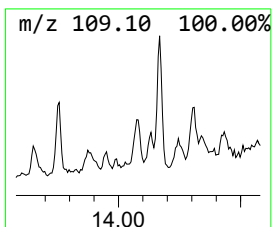
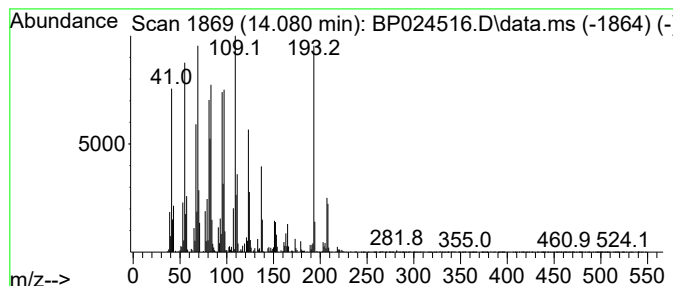
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Peak Number 3 unknown14.080 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.080	2.85 ng	252091	Acenaphthene-d10	14.345

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2(1H)-Naphthalenone, octahydro-4...	208	C14H24O	054594-42-2	46		
2	Decahydro-1,1,4a,5,6-pentamethyl...	208	C15H28	080655-44-3	38		
3	2,6-Heptadienal, 2,4-dimethyl-	138	C9H14O	085136-08-9	38		
4	1H-Indene, 5-butyl-6-hexyloctahy...	264	C19H36	055044-36-5	27		
5	1-Cyclohexanol, 2-(3-methyl-1,3-...	208	C14H24O	1000196-01-6	25		



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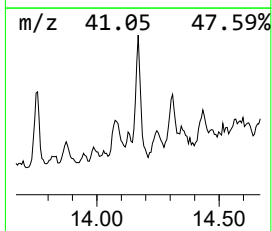
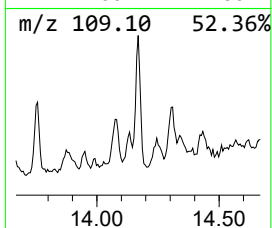
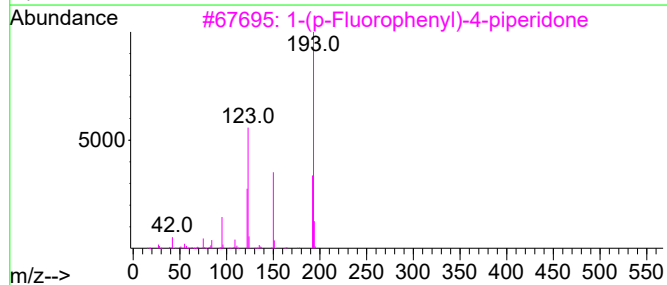
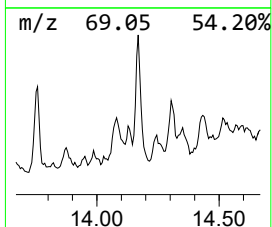
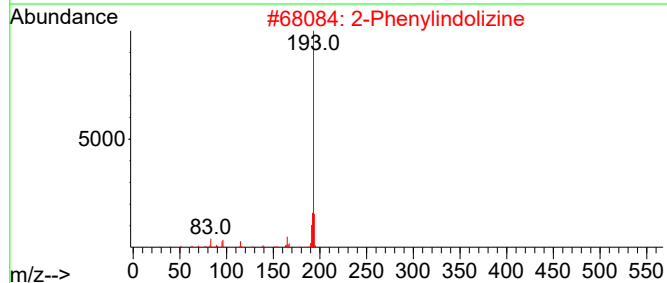
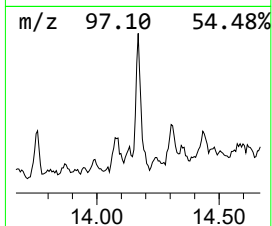
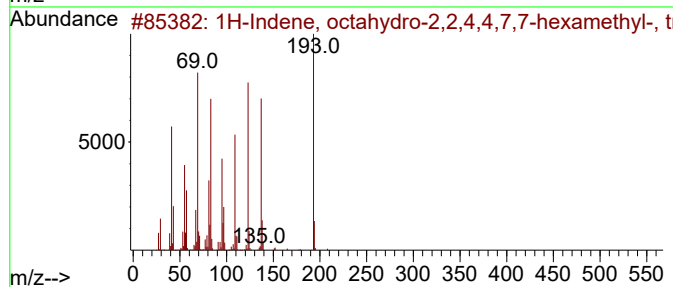
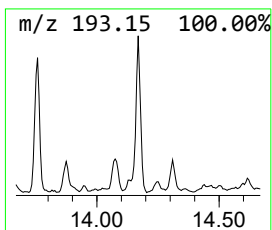
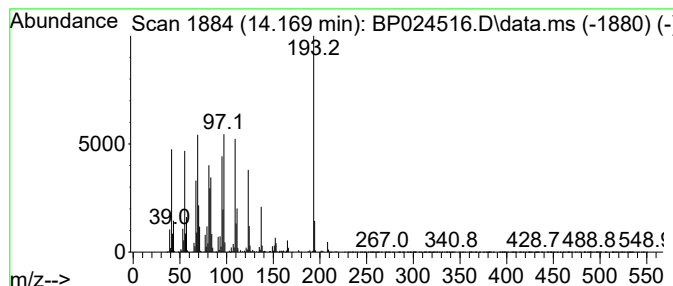
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Peak Number 4 unknown14.169 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.169	6.21 ng	549194	Acenaphthene-d10	14.345

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, octahydro-2,2,4,7,7...	208	C15H28	054832-83-6	37
2			2-Phenylindolizine	193	C14H11N	025379-20-8	27
3			1-(p-Fluorophenyl)-4-piperidone	193	C11H12FNO	1000238-56-7	27
4			6-Methylphenanthridine	193	C14H11N	003955-65-5	27
5			Decahydro-1,1,4a,5,6-pentamethyl...	208	C15H28	080655-44-3	25



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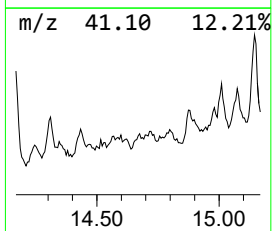
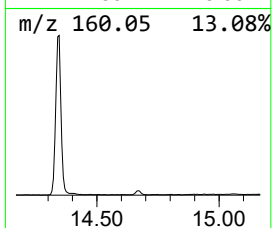
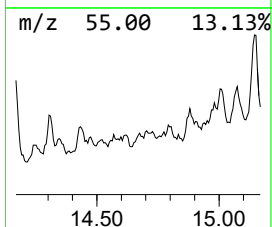
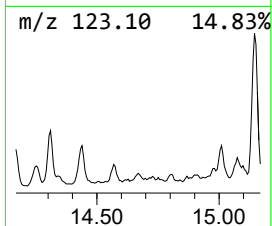
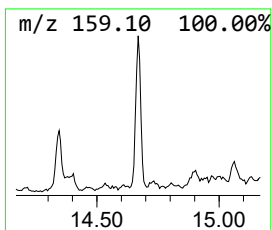
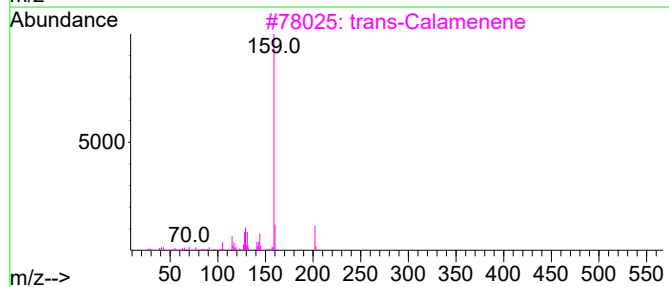
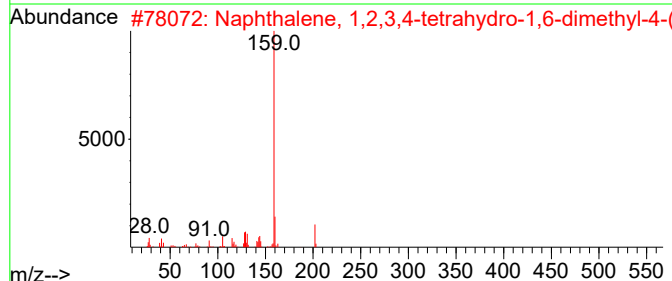
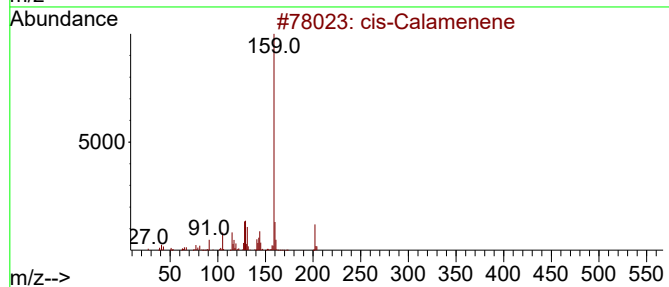
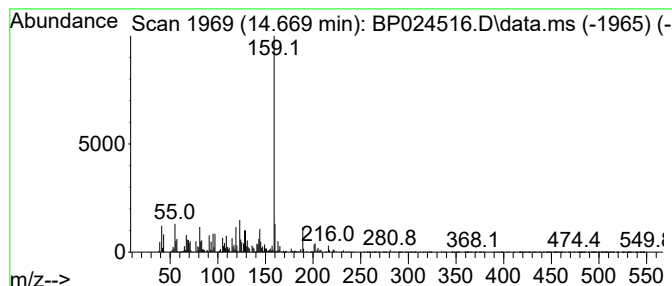
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Peak Number 5 cis-Calamenene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.669	2.23 ng	197280	Acenaphthene-d10	14.345

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			cis-Calamenene	202	C15H22	072937-55-4	76
2			Naphthalene, 1,2,3,4-tetrahydro-...	202	C15H22	000483-77-2	76
3			trans-Calamenene	202	C15H22	073209-42-4	76
4			Naphthalene, 1,2,3,4-tetrahydro-...	174	C13H18	000475-03-6	59
5			Naphthalene, 1,2,3,4-tetrahydro-...	174	C13H18	021693-51-6	59



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ALS Vial : 14 Sample Multiplier: 1

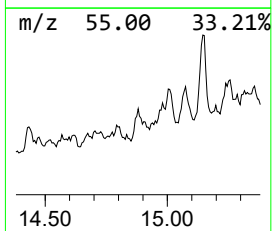
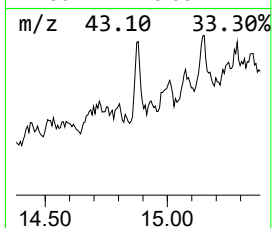
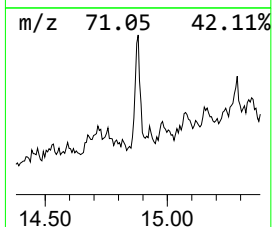
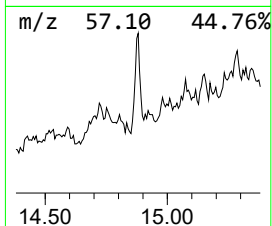
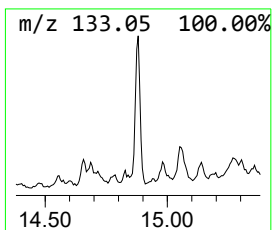
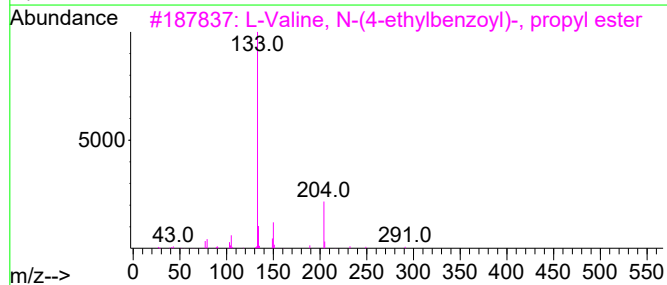
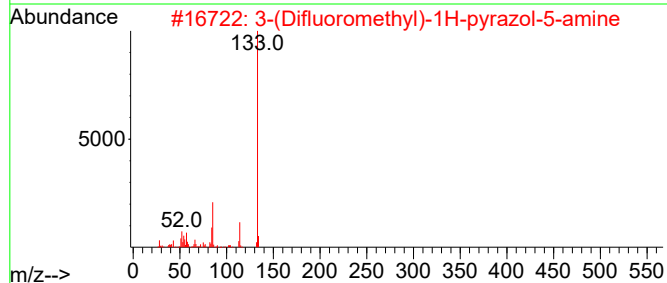
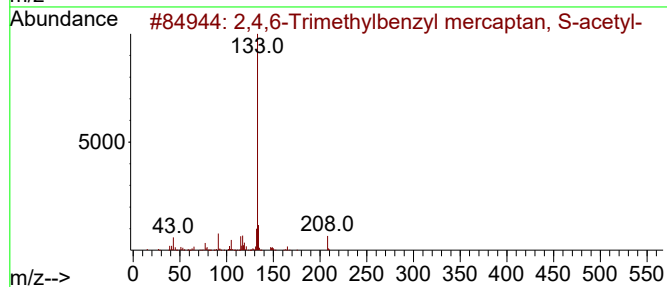
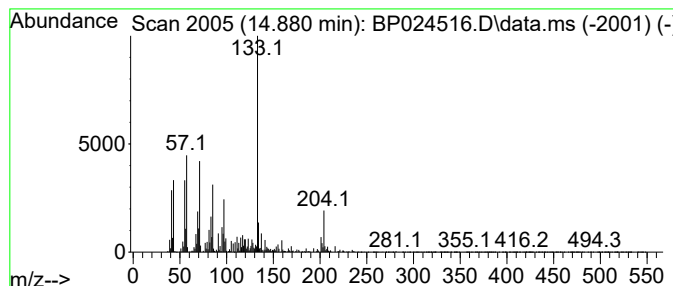
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ClientSampleId :
COMP-5

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP041425.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 unknown14.880 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD			R.T.
14.880	2.84 ng	250735	Acenaphthene-d10			14.345
Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,6-Trimethylbenzyl mercaptan,...	208	C12H16OS		1439439-12-9	41
2	3-(Difluoromethyl)-1H-pyrazol-5-...	133	C4H5F2N3		1284220-49-0	38
3	L-Valine, N-(4-ethylbenzoyl)-, p...	291	C17H25NO3		1000346-62-9	38
4	Benzamide, 4,N-diethyl-N-(2-pent...	247	C16H25NO		1000489-43-6	38
5	Glycine, N-(4-ethylbenzoyl)-, pe...	277	C16H23NO3		1000314-51-6	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050225\
Data File : BP024516.D
Acq On : 02 May 2025 19:44
Operator : RC/JU
Sample : Q1932-03
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
COMP-5

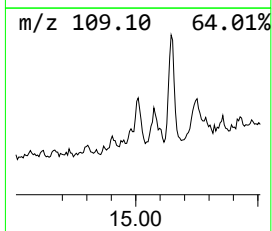
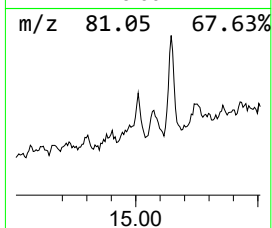
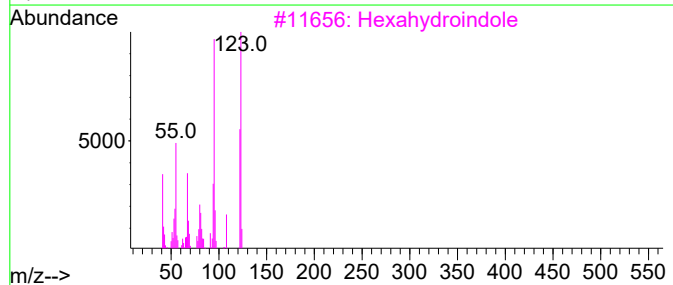
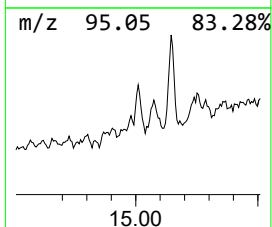
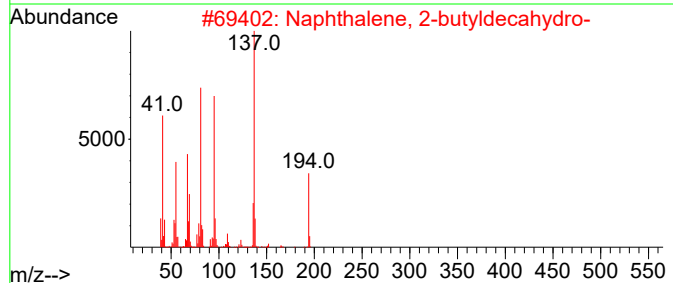
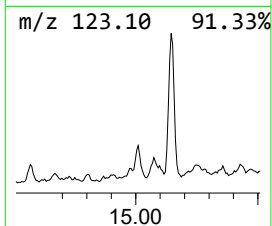
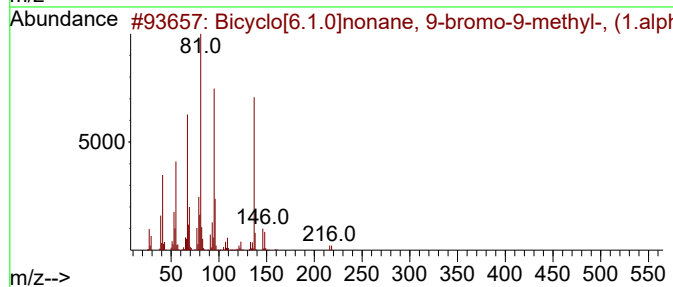
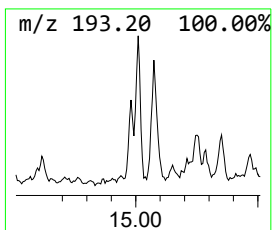
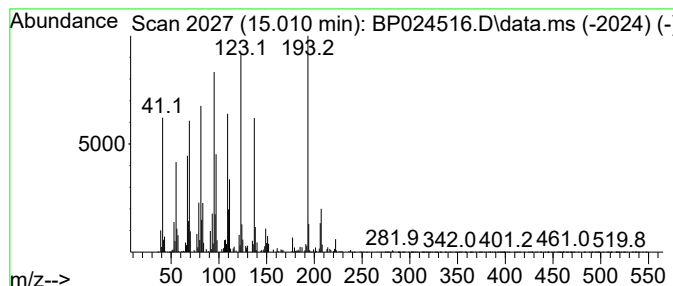
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown15.010 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.010	2.97 ng	262267	Acenaphthene-d10	14.345

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bicyclo[6.1.0]nonane, 9-bromo-9-...	216	C10H17Br	059474-03-2	44
2			Naphthalene, 2-butyldecahydro-	194	C14H26	006305-52-8	35
3			Hexahydroindole	123	C8H13N	018159-32-5	30
4			Silane, dimethyldi(2,4-cyclopent...	188	C12H16Si	018053-74-2	27
5			Naphthalene, 1,1'-ethylidenebis[...	302	C22H38	054934-70-2	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050225\
Data File : BP024516.D
Acq On : 02 May 2025 19:44
Operator : RC/JU
Sample : Q1932-03
Misc :
ALS Vial : 14 Sample Multiplier: 1

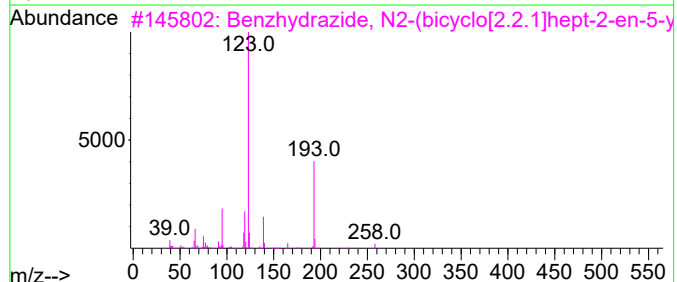
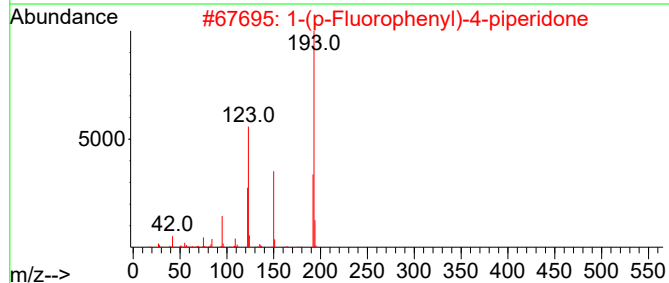
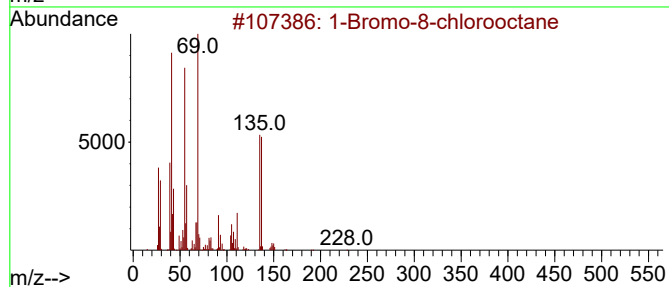
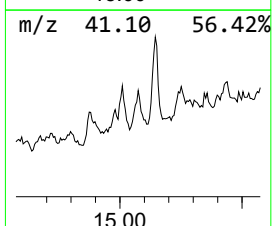
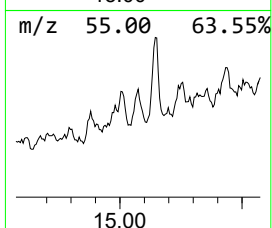
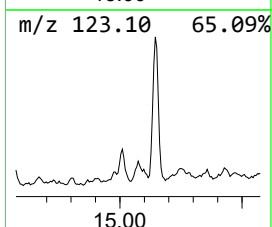
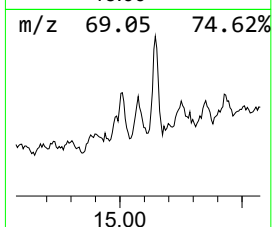
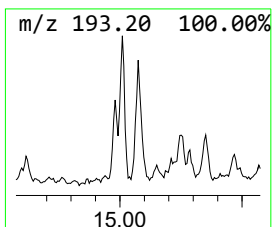
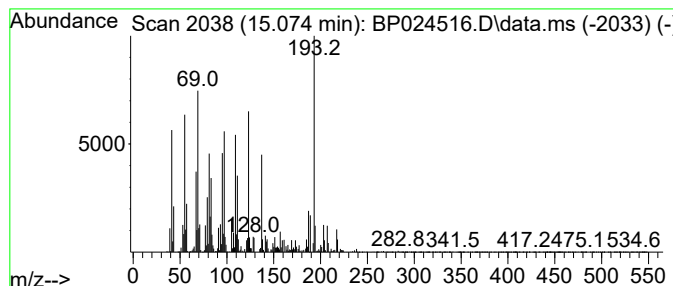
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP041425.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown15.075 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.	
15.075	3.55 ng	313605	Acenaphthene-d10		14.345	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	1-Bromo-8-chlorooctane		226	C8H16BrCl	1000438-83-6	35
2	1-(p-Fluorophenyl)-4-piperidone		193	C11H12FNO	1000238-56-7	25
3	Benzhydrazide, N2-(bicyclo[2.2.1...]		258	C15H15FN2O	1000274-01-8	22
4	Decahydro-1,1,4a,5,6-pentamethyl...		208	C15H28	080655-44-3	18
5	Geranyl bromide		216	C10H17Br	006138-90-5	14



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050225\
Data File : BP024516.D
Acq On : 02 May 2025 19:44
Operator : RC/JU
Sample : Q1932-03
Misc :
ALS Vial : 14 Sample Multiplier: 1

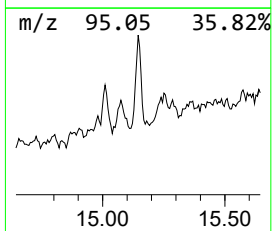
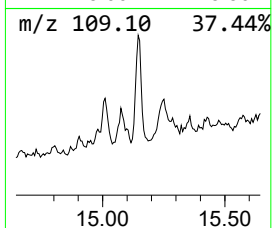
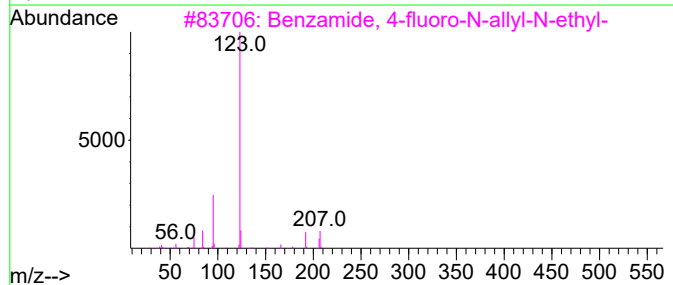
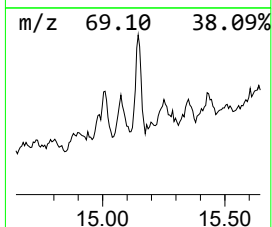
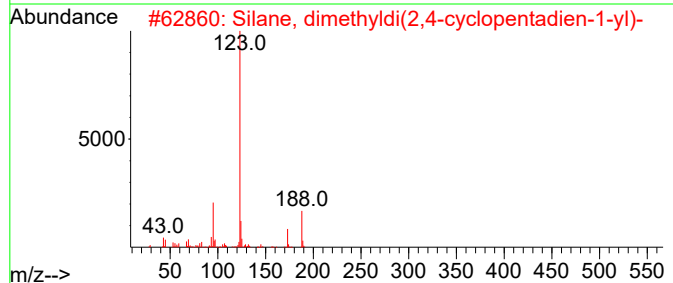
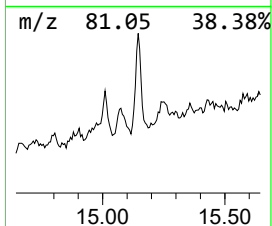
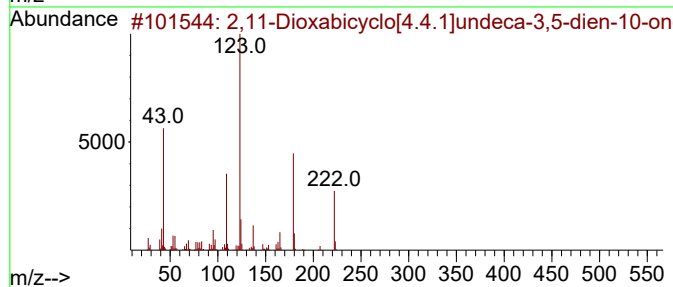
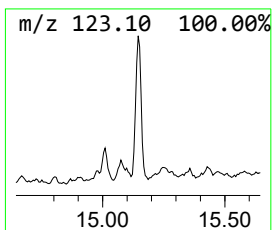
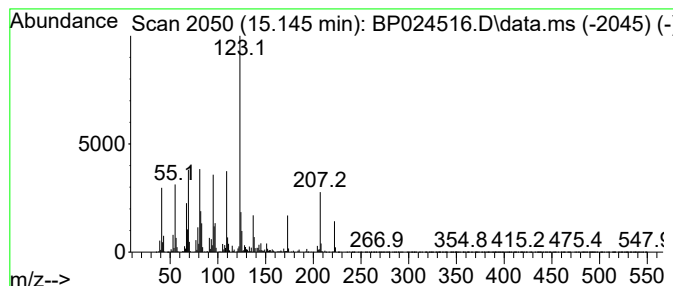
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 2,11-Dioxabicyclo[4.4.1]und... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD		R.T.	
15.145	9.41 ng	832315	Acenaphthene-d10		14.345	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	2,11-Dioxabicyclo[4.4.1]undeca-3...		222	C13H18O3	070412-52-1	53
2	Silane, dimethyldi(2,4-cyclopent...		188	C12H16Si	018053-74-2	50
3	Benzamide, 4-fluoro-N-allyl-N-et...		207	C12H14FNO	1000446-42-7	47
4	1-Butanone, 1-bicyclo[4.1.0]hept...		166	C11H18O	054764-62-4	47
5	3-Fluorobenzoic acid, 3-methylbu...		208	C12H13FO2	154559-38-3	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050225\
Data File : BP024516.D
Acq On : 02 May 2025 19:44
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Sample : Q1932-03
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Instrument :
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ClientSampleId :
COMP-5

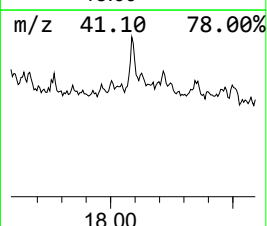
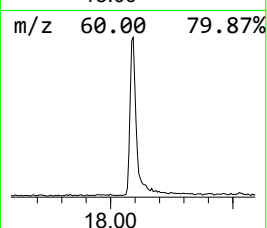
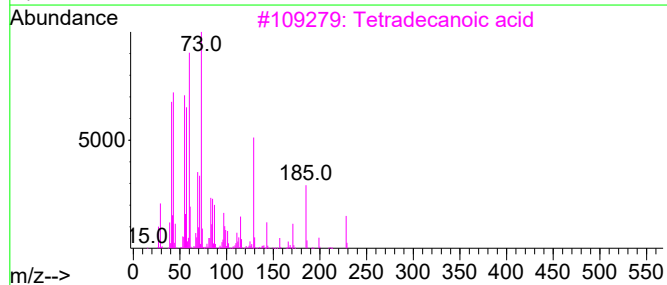
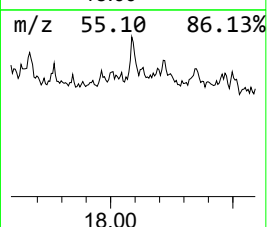
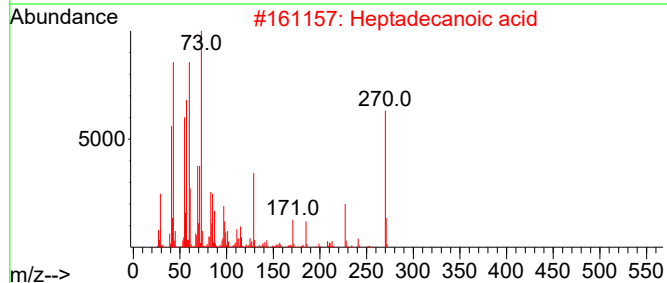
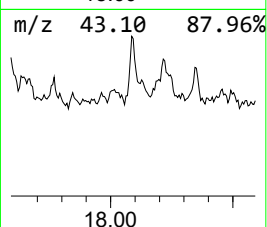
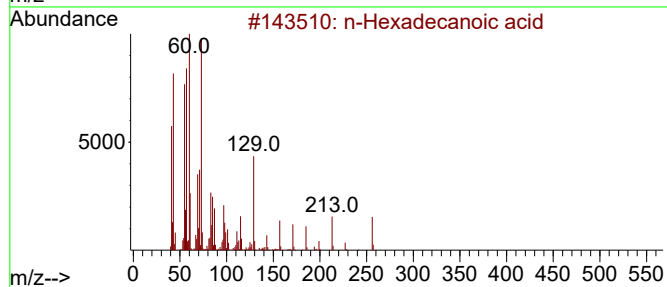
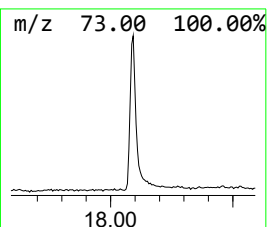
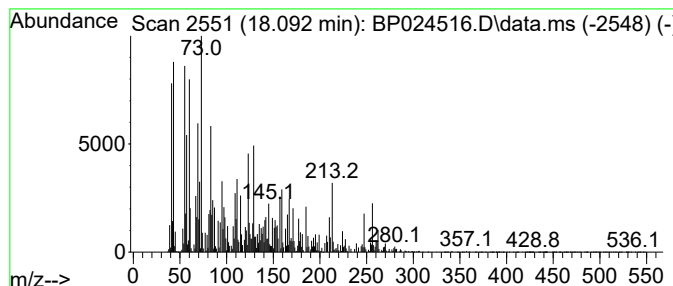
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.092	5.36 ng	562435	Phenanthrene-d10	17.151

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
2		Heptadecanoic acid	270	C17H34O2	000506-12-7	64
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	50
4		n-Decanoic acid	172	C10H20O2	000334-48-5	46
5		Tridecanoic acid	214	C13H26O2	000638-53-9	41



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050225\
Data File : BP024516.D
Acq On : 02 May 2025 19:44
Operator : RC/JU
Sample : Q1932-03
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
COMP-5

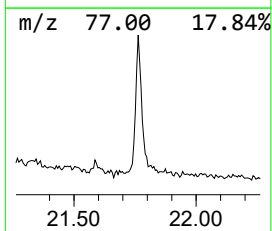
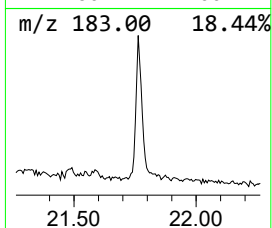
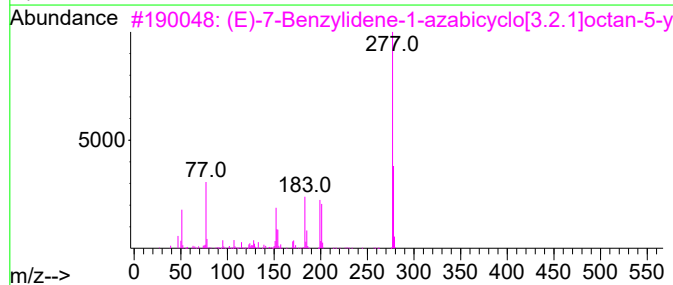
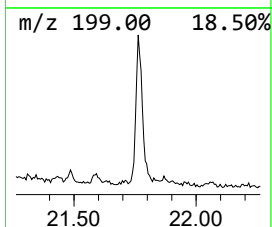
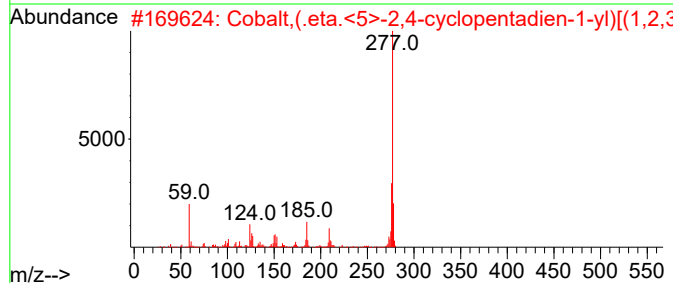
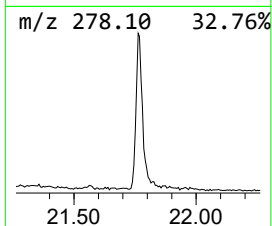
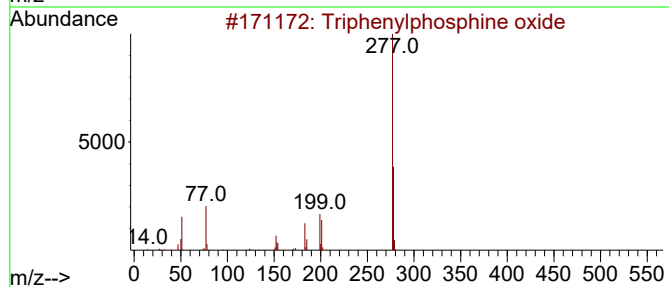
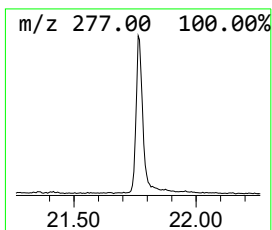
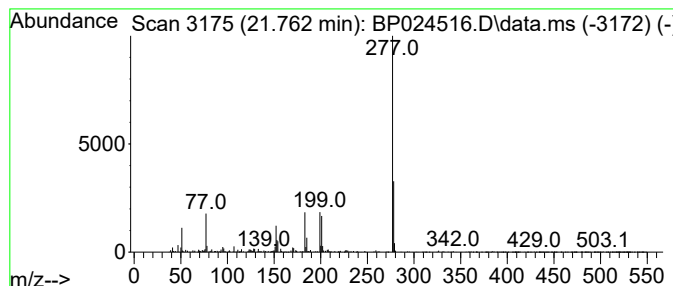
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Triphenylphosphine oxide Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.762	3.84 ng	449945	Chrysene-d12	21.592

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Triphenylphosphine oxide	278	C18H15OP	000791-28-6	91
2			Cobalt,(.eta.<5>-2,4-cyclopentad...	277	C16H15BCo	036682-12-9	59
3			(E)-7-Benzylidene-1-azabicyclo[3...	293	C15H19N03S	1000483-83-4	52
4			Formaldehyde, (triphenylphosphor...	304	C19H17N2P	015990-54-2	45
5			(Carbethoxymethyl)-triphenylphos...	428	C22H22Br02P	001530-45-6	42



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050225\
Data File : BP024516.D
Acq On : 02 May 2025 19:44
Operator : RC/JU
Sample : Q1932-03
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP041425.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-...	4.875	4.0	ng	173103	1	7.704	865985	20.0
Borinic acid, d...	13.757	4.2	ng	374616	3	14.345	1768290	20.0
unknown14.080	14.080	2.9	ng	252091	3	14.345	1768290	20.0
unknown14.169	14.169	6.2	ng	549194	3	14.345	1768290	20.0
cis-Calamenene	14.669	2.2	ng	197280	3	14.345	1768290	20.0
unknown14.880	14.880	2.8	ng	250735	3	14.345	1768290	20.0
unknown15.010	15.010	3.0	ng	262267	3	14.345	1768290	20.0
unknown15.075	15.075	3.5	ng	313605	3	14.345	1768290	20.0
2,11-Dioxabicyc...	15.145	9.4	ng	832315	3	14.345	1768290	20.0
n-Hexadecanoic ...	18.092	5.4	ng	562435	4	17.151	2096720	20.0
Triphenylphosph...	21.762	3.8	ng	449945	5	21.592	2341780	20.0