

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF012325\
 Data File : BF141251.D
 Acq On : 23 Jan 2025 15:49
 Operator : RC/JU
 Sample : Q1154-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 VNJ-220

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF011025.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF141251.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.234	17	24	29	rBV	334919	564028	8.57%	1.696%
2	5.004	481	495	501	rBV	186683	293236	4.45%	0.882%
3	5.416	557	565	571	rBV	1791700	2509831	38.13%	7.548%
4	6.087	670	679	684	rBV	2815179	3614854	54.91%	10.871%
5	6.251	699	707	712	rBV	277673	437642	6.65%	1.316%
6	6.345	713	723	728	rBV	268265	577556	8.77%	1.737%
7	6.440	732	739	743	rBV	1602413	2488100	37.80%	7.482%
8	6.698	777	783	786	rBV	543960	1116047	16.95%	3.356%
9	6.804	796	801	808	rVB	598740	762532	11.58%	2.293%
10	6.881	808	814	816	rBV	270405	408931	6.21%	1.230%
11	6.922	816	821	832	rVB5	569790	1278428	19.42%	3.845%
12	7.210	862	870	877	rBV2	371251	695748	10.57%	2.092%
13	7.363	887	896	901	rBV2	1237512	2112053	32.08%	6.351%
14	7.581	907	933	938	rBV	1362023	6582720	100.00%	19.796%
15	7.622	938	940	945	rVV	359789	452103	6.87%	1.360%
16	7.675	945	949	954	rVB2	189023	268319	4.08%	0.807%
17	7.845	972	978	982	rVB4	230129	428110	6.50%	1.287%
18	7.998	999	1004	1008	rBV	410211	559606	8.50%	1.683%
19	8.087	1013	1019	1021	rBV2	738255	1123357	17.07%	3.378%
20	8.116	1021	1024	1028	rVB	1504143	1850584	28.11%	5.565%
21	8.234	1039	1044	1051	rBV	278596	601734	9.14%	1.810%
22	8.386	1066	1070	1073	rBV6	161972	259470	3.94%	0.780%
23	8.475	1081	1085	1088	rBV3	202915	316421	4.81%	0.952%
24	8.592	1102	1105	1108	rBV3	168521	234704	3.57%	0.706%
25	9.169	1199	1203	1207	rBV	1809015	2404983	36.53%	7.232%
26	9.251	1214	1217	1221	rBV4	829322	1311895	19.93%	3.945%

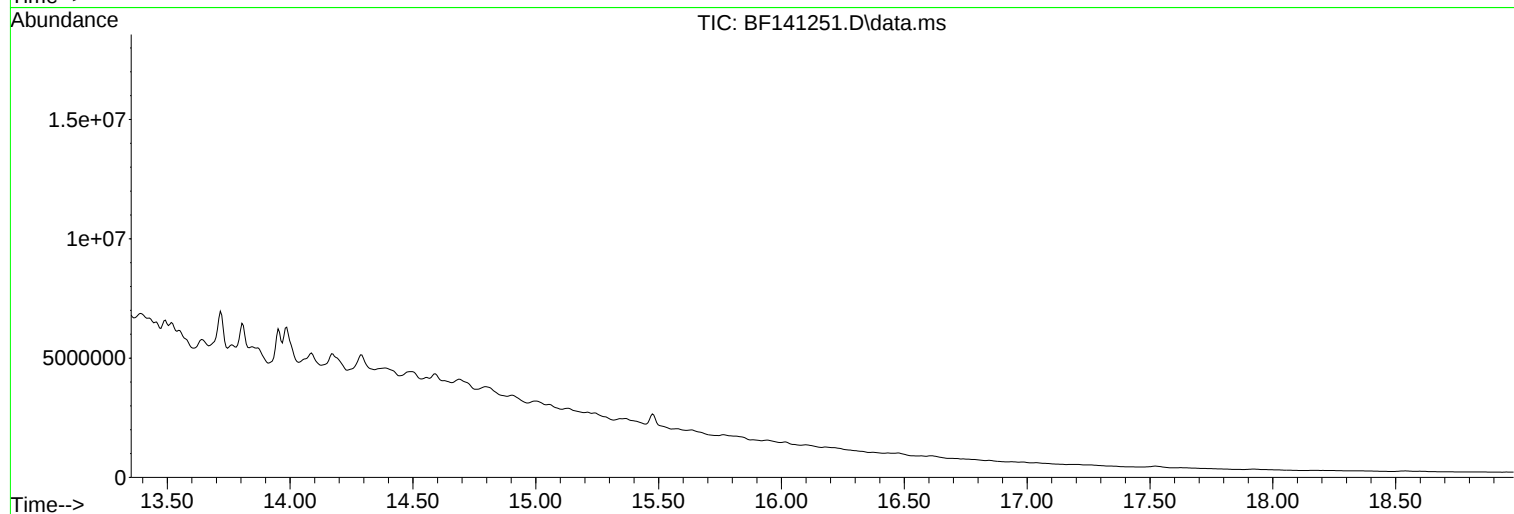
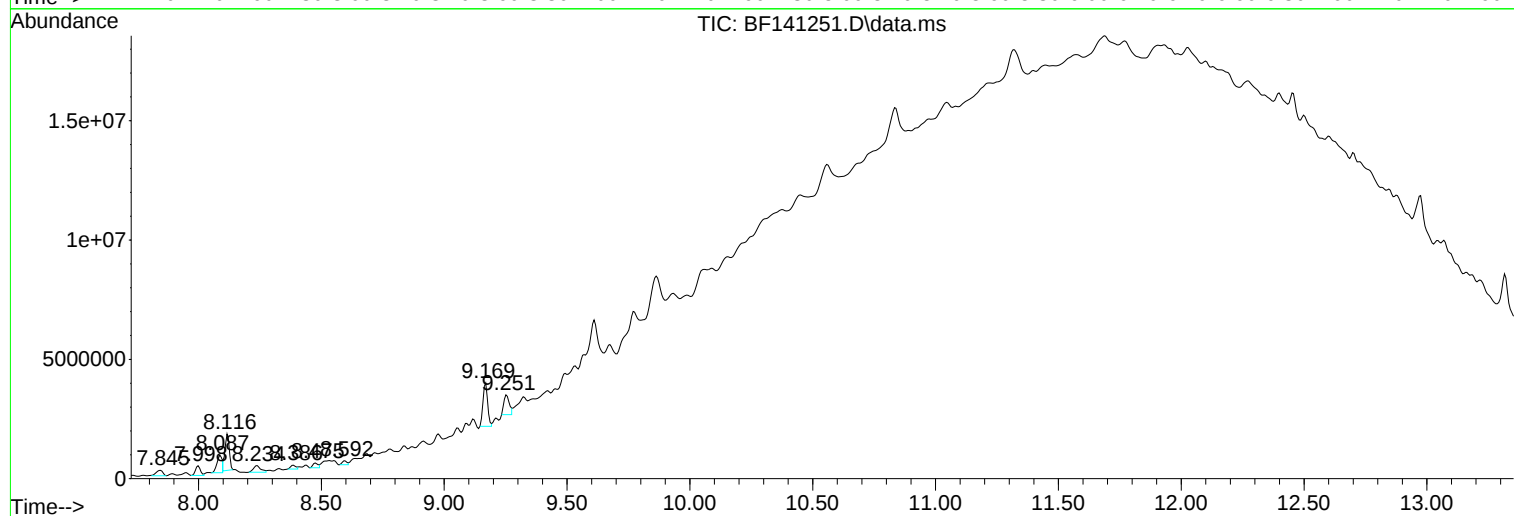
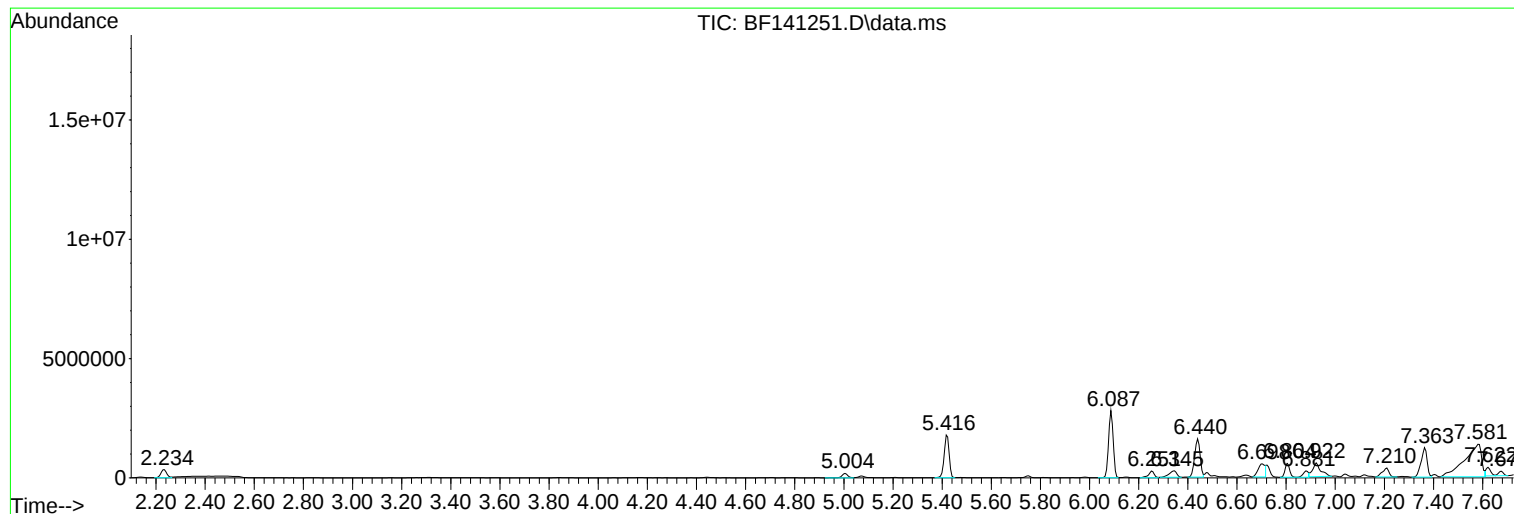
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BNA_F
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF011025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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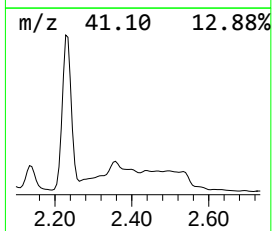
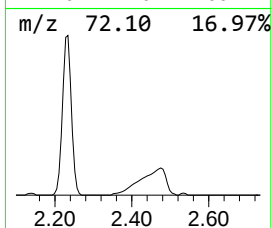
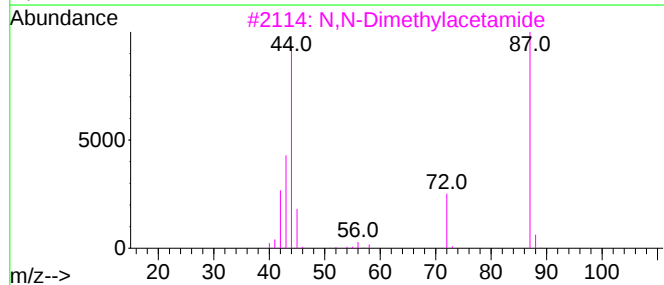
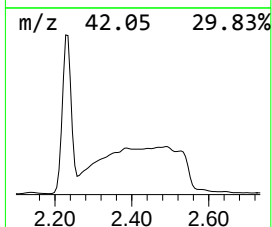
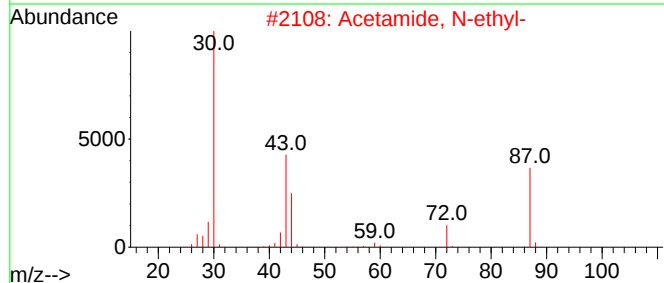
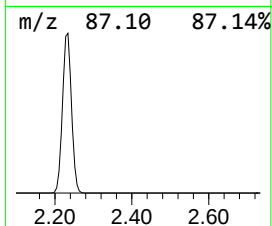
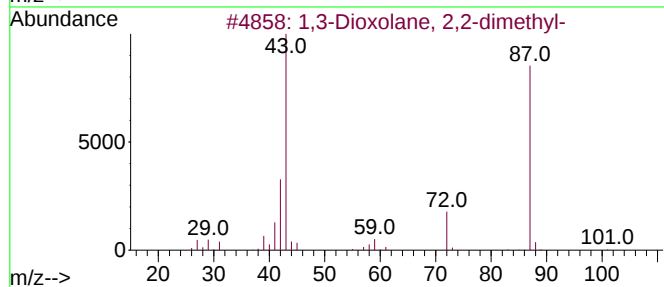
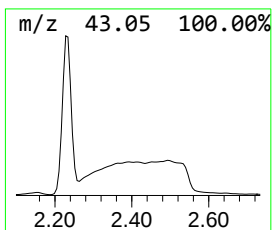
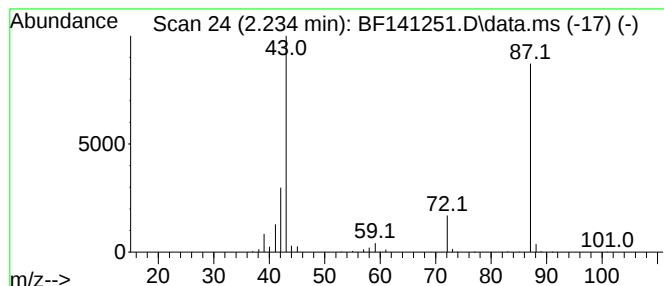
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF011025.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 1,3-Dioxolane, 2,2-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.234	14.79 ng	564028	1,4-Dichlorobenzene-d4	6.804

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3-Dioxolane, 2,2-dimethyl-	102	C5H10O2	002916-31-6	90
2			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	72
3			N,N-Dimethylacetamide	87	C4H9NO	000127-19-5	59
4			2-OXOPROPIONAMIDE	87	C3H5NO2	000631-66-3	42
5			Ethanol, 2,2'-oxybis-, diacetate	190	C8H14O5	000628-68-2	36



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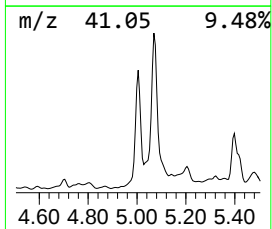
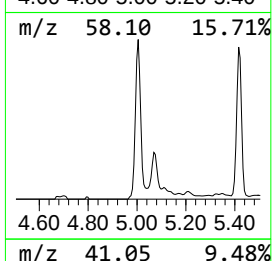
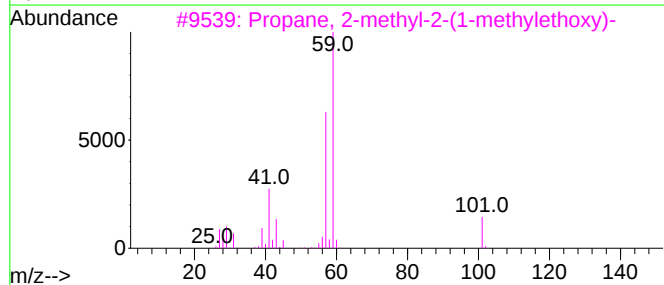
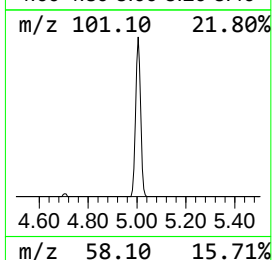
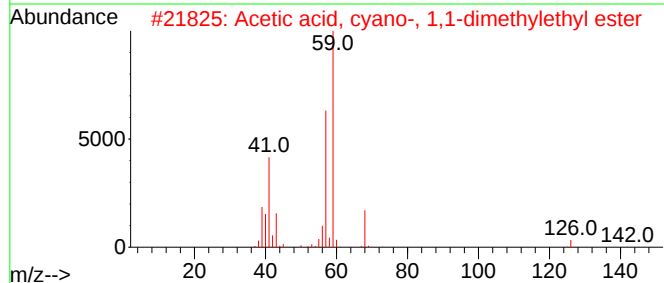
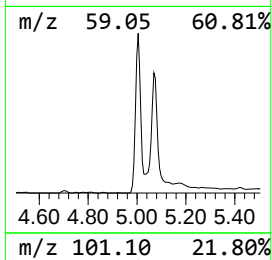
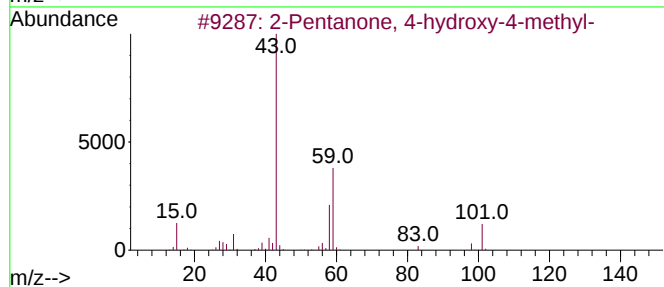
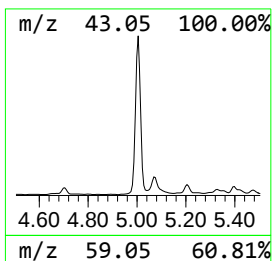
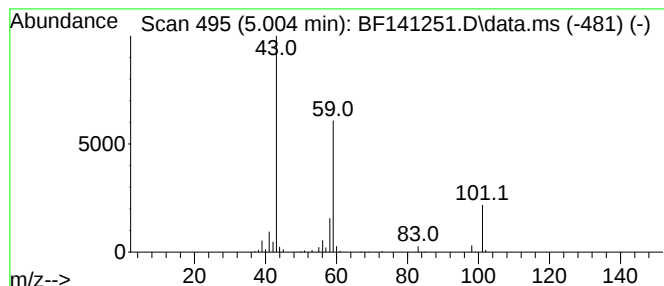
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.004	7.69 ng	293236	1,4-Dichlorobenzene-d4	6.804

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	32
3			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	23
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	12
5			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10



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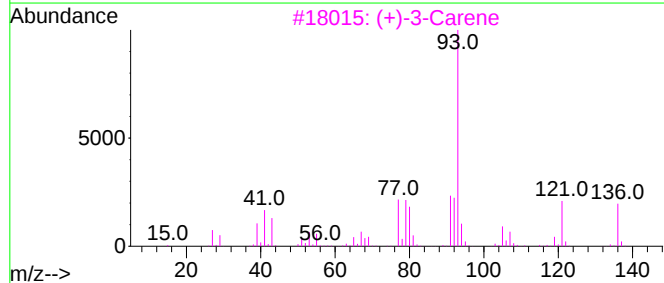
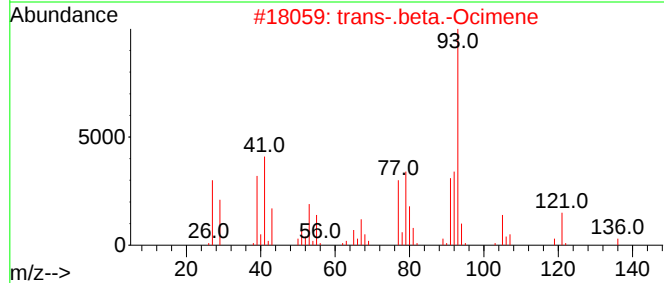
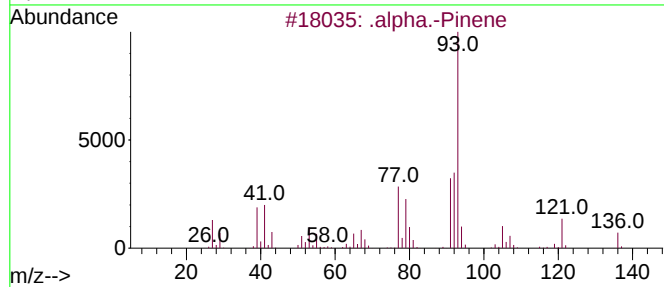
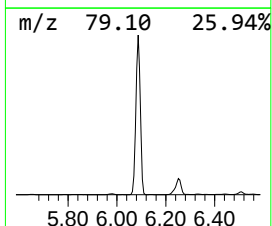
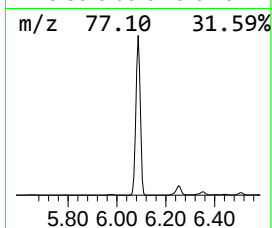
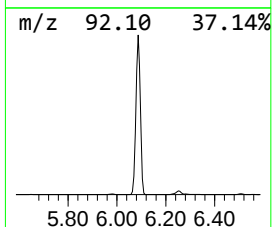
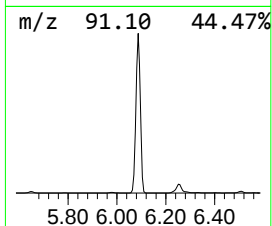
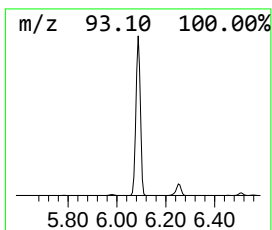
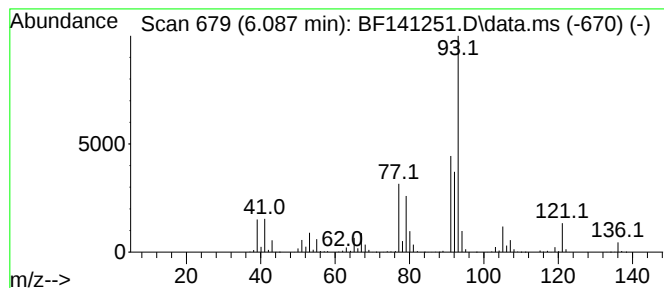
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 .alpha.-Pinene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.087	94.81 ng	3614850	1,4-Dichlorobenzene-d4	6.804

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	.		.alpha.-Pinene	136	C10H16	000080-56-8	95
2	trans-.		beta.-Ocimene	136	C10H16	003779-61-1	94
3	(+)-3-		Carene	136	C10H16	000498-15-7	91
4	Tricyclo[2.2.1.0(2,6)]heptane, 1...			136	C10H16	000508-32-7	91
5	.gamma.-		Terpinene	136	C10H16	000099-85-4	91



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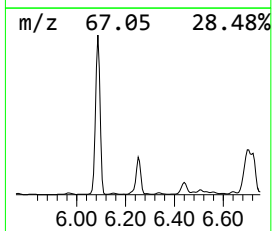
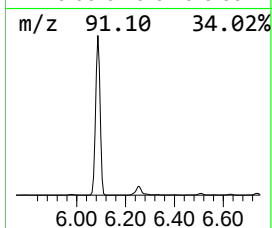
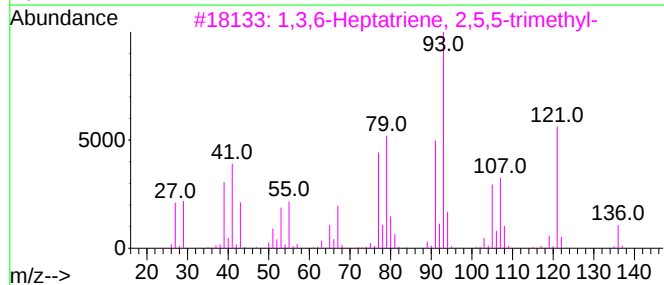
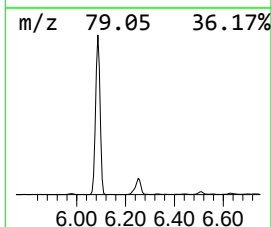
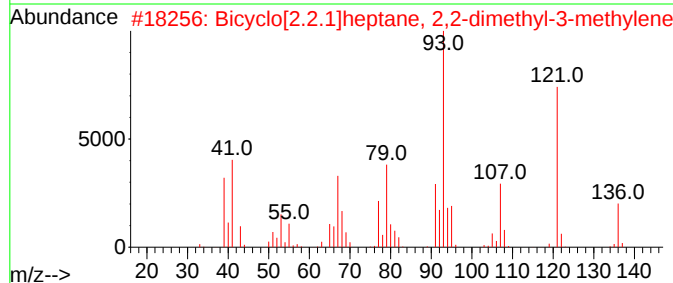
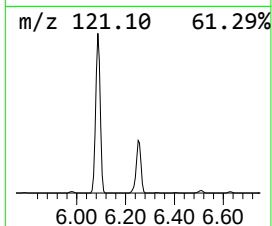
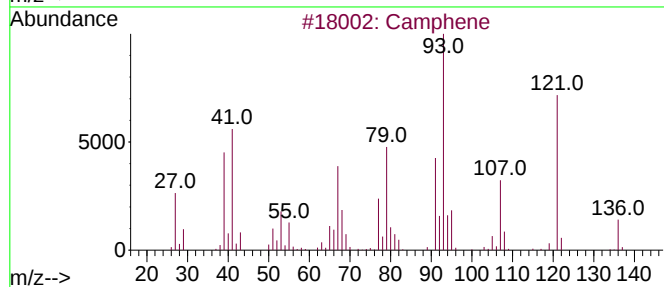
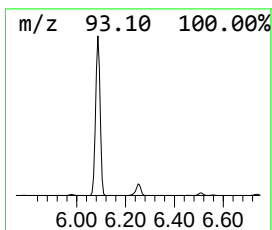
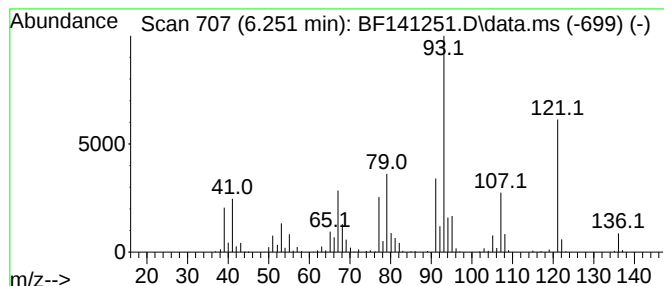
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Peak Number 4 Camphene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.251	11.48 ng	437642	1,4-Dichlorobenzene-d4	6.804

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Camphene	136	C10H16	000079-92-5	94
2			Bicyclo[2.2.1]heptane, 2,2-dimet...	136	C10H16	005794-04-7	94
3			1,3,6-Heptatriene, 2,5,5-trimethyl-	136	C10H16	029548-02-5	90
4			Cyclohexene, 3-methyl-6-(1-methy...	136	C10H16	000586-63-0	80
5			(+)-4-Carene	136	C10H16	029050-33-7	74



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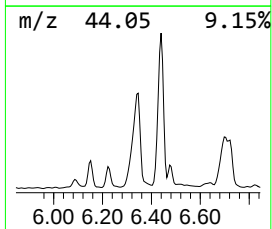
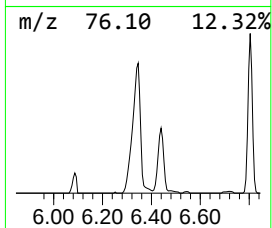
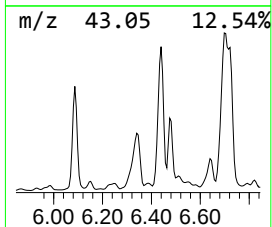
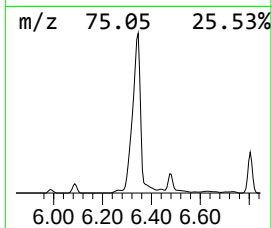
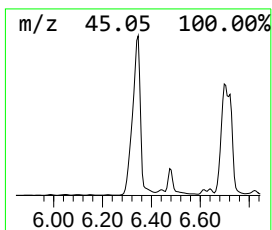
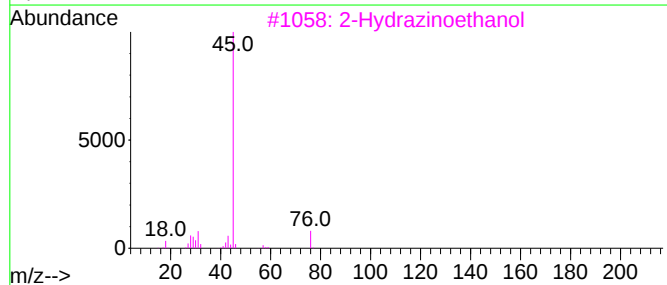
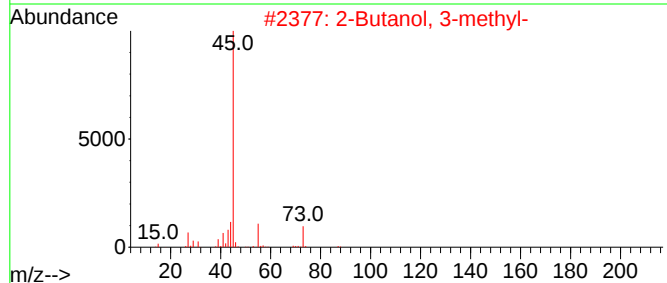
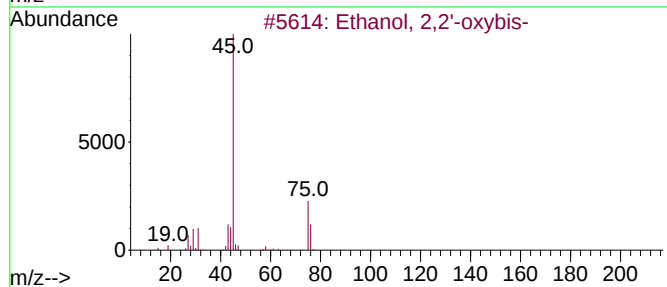
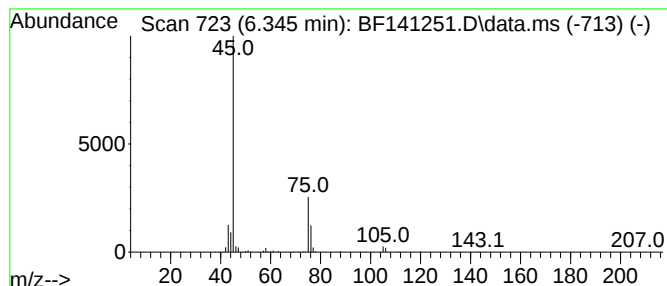
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Peak Number 5 Ethanol, 2,2'-oxybis- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.345	15.15 ng	577556	1,4-Dichlorobenzene-d4	6.804

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2,2'-oxybis-	106	C4H10O3	000111-46-6	83
2			2-Butanol, 3-methyl-	88	C5H12O	000598-75-4	42
3			2-Hydrazinoethanol	76	C2H8N2O	000109-84-2	9
4			Isopropyl Alcohol	60	C3H8O	000067-63-0	9
5			Ethanol, 2-methoxy-	76	C3H8O2	000109-86-4	9



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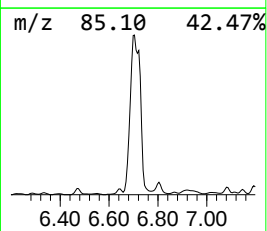
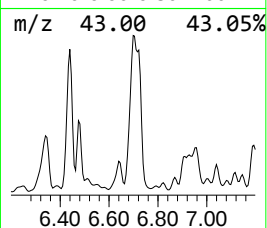
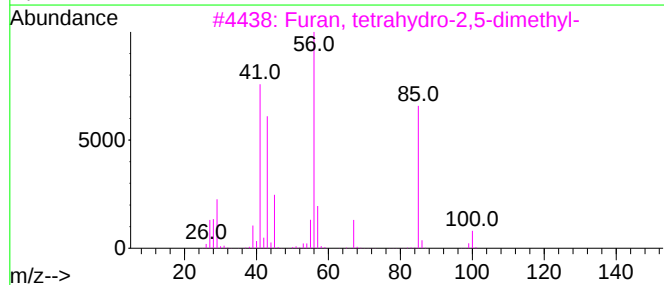
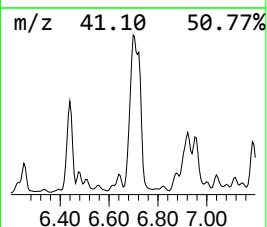
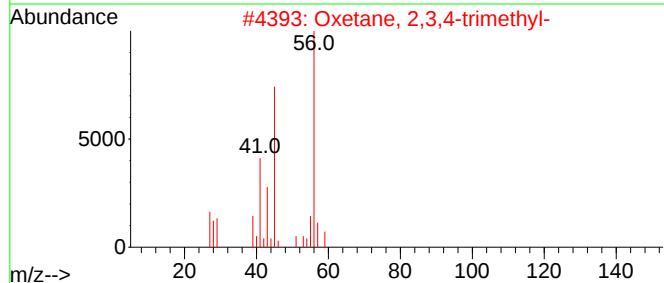
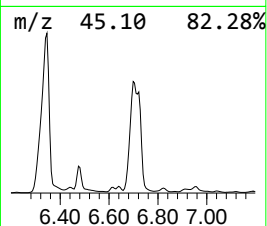
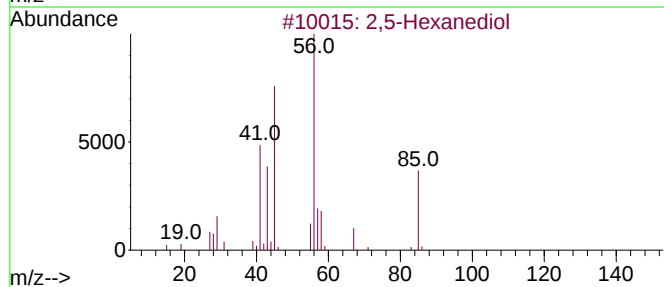
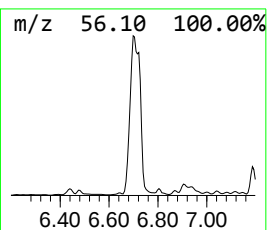
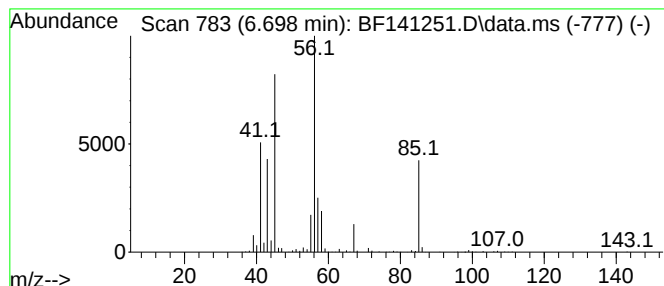
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 2,5-Hexanediol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.698	29.27 ng	1116050	1,4-Dichlorobenzene-d4	6.804

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,5-Hexanediol	118	C6H14O2	002935-44-6	83
2			Oxetane, 2,3,4-trimethyl-	100	C6H12O	053778-61-3	53
3			Furan, tetrahydro-2,5-dimethyl-	100	C6H12O	001003-38-9	47
4			4-Penten-2-ol, 4-methyl-	100	C6H12O	002004-67-3	42
5			Hexane, 3,4-dimethyl-	114	C8H18	000583-48-2	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF012325\
Data File : BF141251.D
Acq On : 23 Jan 2025 15:49
Operator : RC/JU
Sample : Q1154-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
VNJ-220

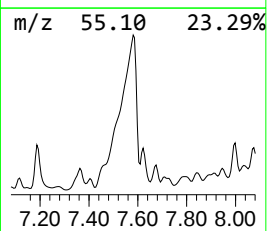
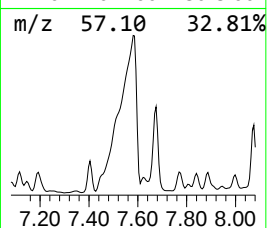
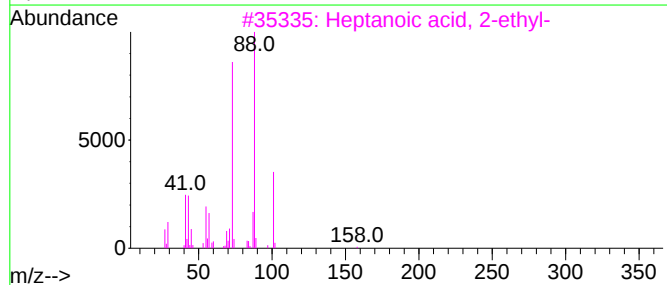
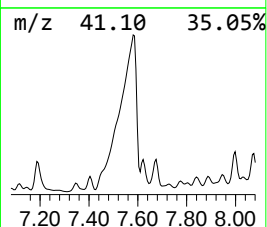
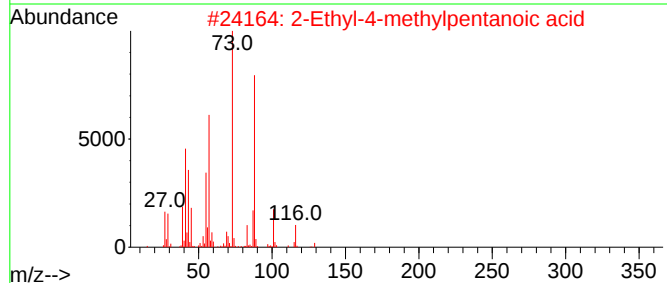
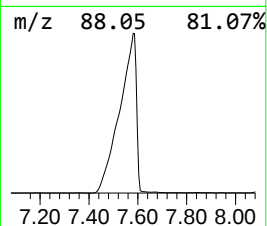
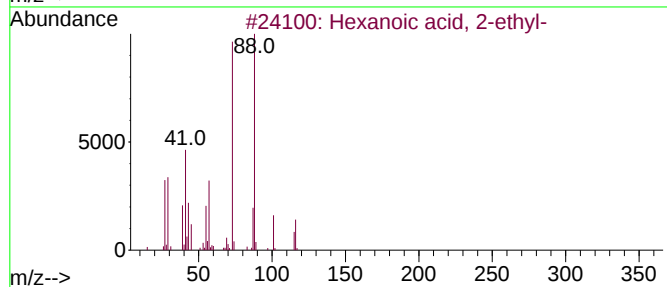
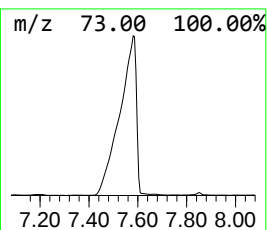
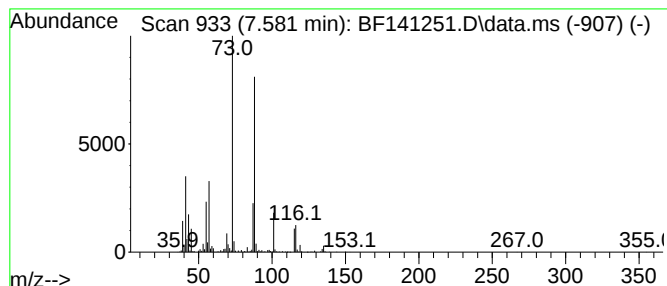
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Hexanoic acid, 2-ethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.581	117.20 ng	6582720	Naphthalene-d8	8.087

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	90
2			2-Ethyl-4-methylpentanoic acid	144	C8H16O2	000108-81-6	72
3			Heptanoic acid, 2-ethyl-	158	C9H18O2	003274-29-1	53
4			Methyl-2,3,4--tri-O-methyl.alpha...	220	C10H20O5	003253-23-4	50
5			Diisopropanolamine	133	C6H15NO2	000110-97-4	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF012325\
Data File : BF141251.D
Acq On : 23 Jan 2025 15:49
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Sample : Q1154-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

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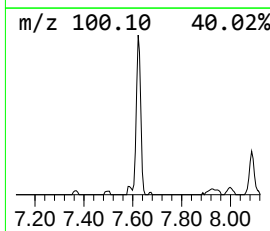
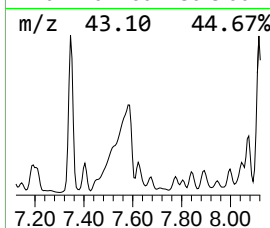
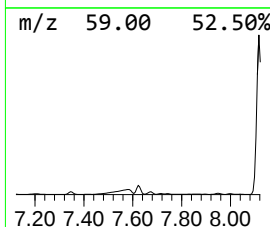
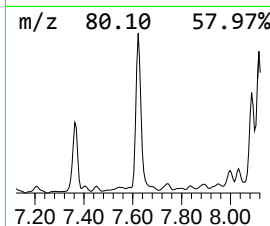
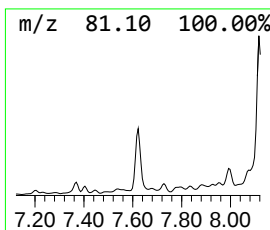
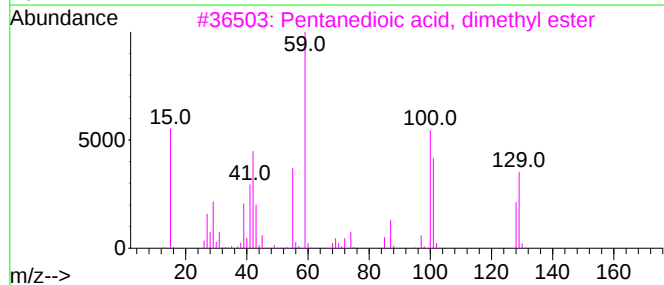
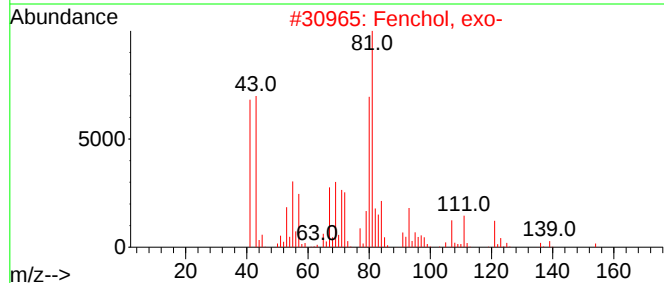
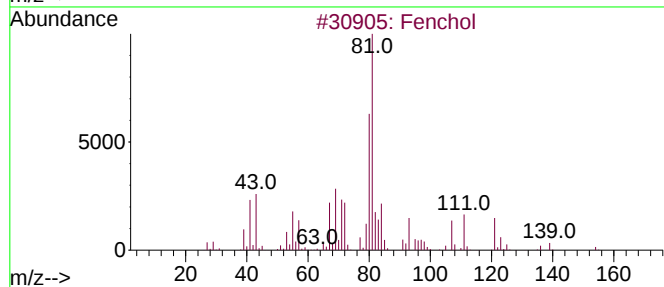
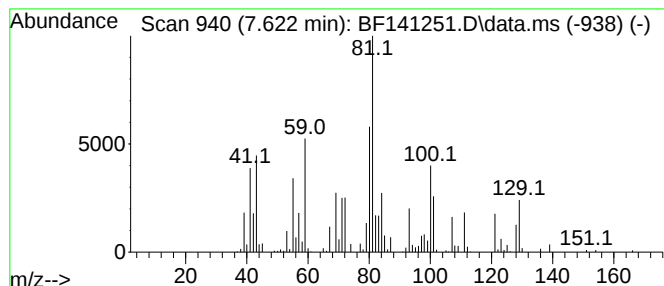
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Fenchol Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.622	8.05 ng	452103	Naphthalene-d8	8.087

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Fenchol	154	C10H18O	001632-73-1	50
2			Fenchol, exo-	154	C10H18O	022627-95-8	37
3			Pentanedioic acid, dimethyl ester	160	C7H12O4	001119-40-0	25
4			Bicyclo[2.2.1]heptan-2-ol, 1,3,3...	196	C12H20O2	076109-40-5	17
5			Butanedioic acid, methyl-, dimet...	160	C7H12O4	001604-11-1	14



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF012325\
Data File : BF141251.D
Acq On : 23 Jan 2025 15:49
Operator : RC/JU
Sample : Q1154-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

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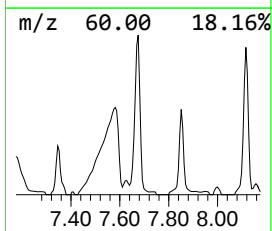
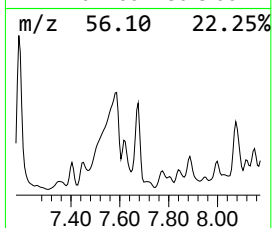
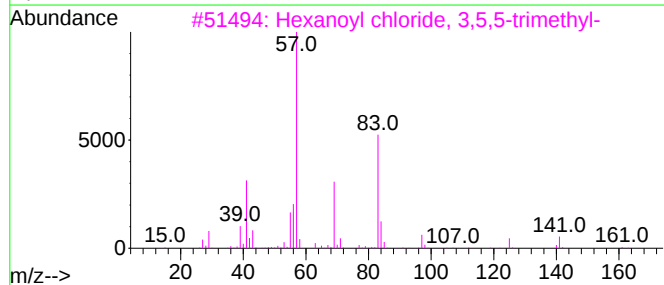
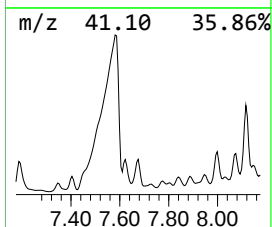
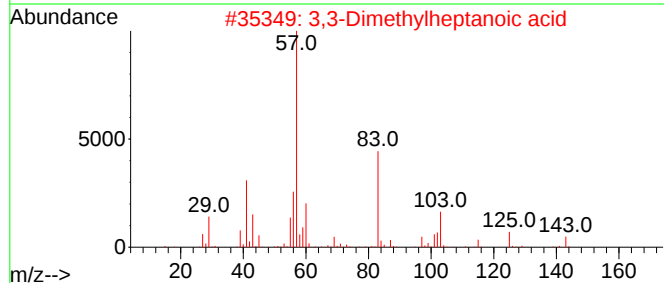
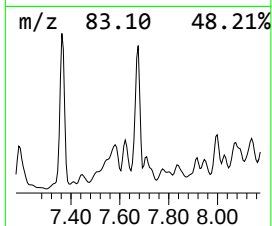
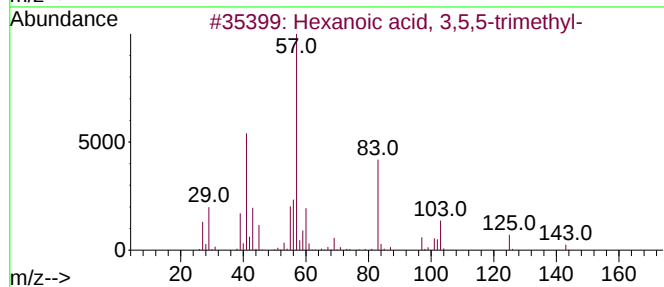
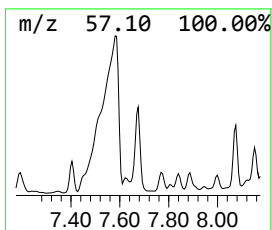
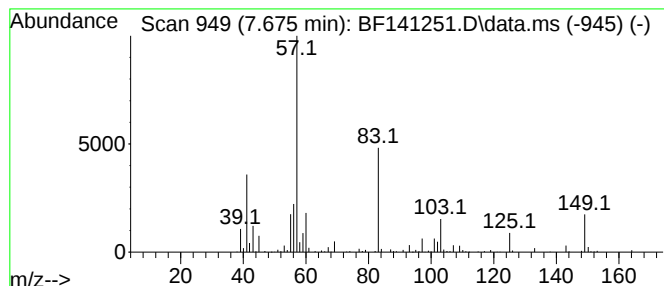
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Hexanoic acid, 3,5,5-trimet... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.675	4.78 ng	268319	Naphthalene-d8	8.087

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexanoic acid, 3,5,5-trimethyl-	158	C9H18O2	003302-10-1	86
2			3,3-Dimethylheptanoic acid	158	C9H18O2	067061-30-7	78
3			Hexanoyl chloride, 3,5,5-trimethyl-	176	C9H17ClO	036727-29-4	42
4			6,6-Dimethylheptanoic acid	158	C9H18O2	015898-92-7	37
5			4,4-Dimethyl-2-pentanol, chlorod...	228	C9H15ClF2O2	1000376-27-0	32



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF012325\
Data File : BF141251.D
Acq On : 23 Jan 2025 15:49
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Sample : Q1154-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

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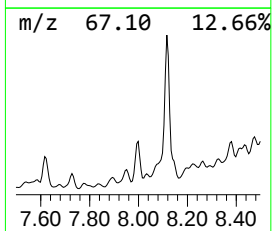
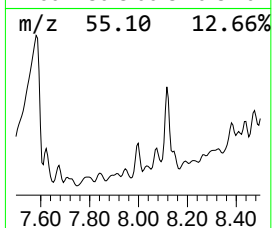
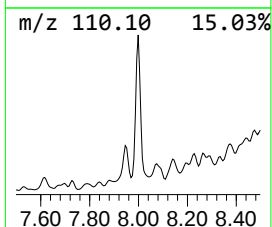
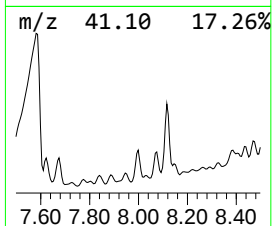
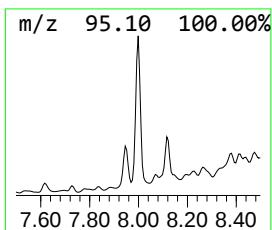
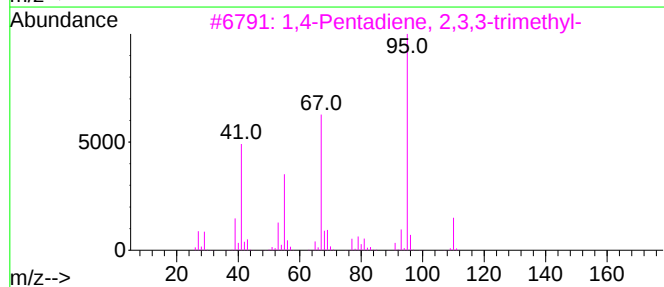
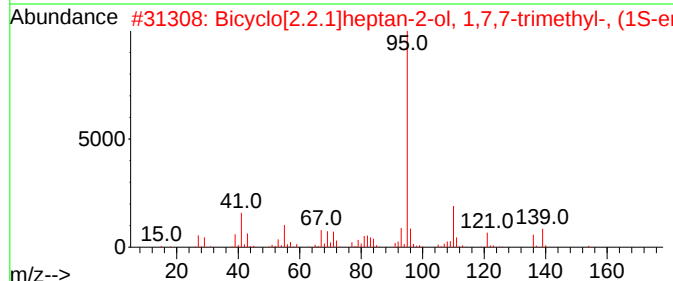
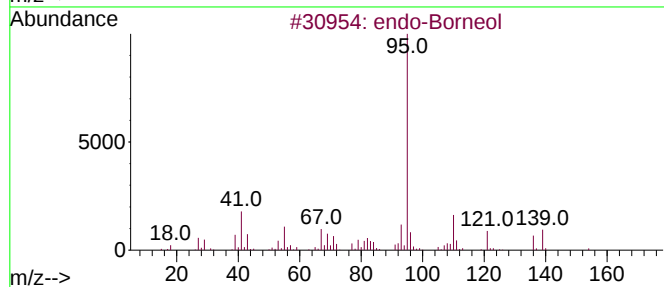
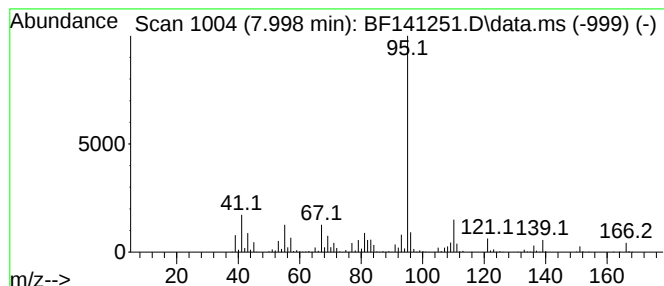
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 endo-Borneol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.998	9.96 ng	559606	Naphthalene-d8	8.087

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		endo-Borneol	154	C10H18O	000507-70-0	90
2		Bicyclo[2.2.1]heptan-2-ol, 1,7,7...	154	C10H18O	000464-45-9	87
3		1,4-Pentadiene, 2,3,3-trimethyl-	110	C8H14	000756-02-5	64
4		1,4-Pentadiene, 2,3,4-trimethyl-	110	C8H14	072014-90-5	64
5		Perilla ketone	166	C10H14O2	000553-84-4	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF012325\
Data File : BF141251.D
Acq On : 23 Jan 2025 15:49
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Misc :
ALS Vial : 9 Sample Multiplier: 1

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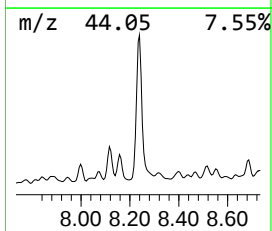
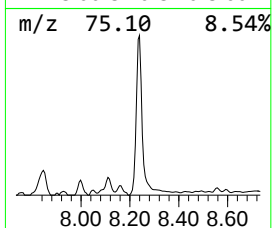
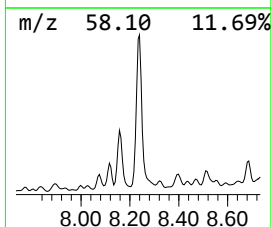
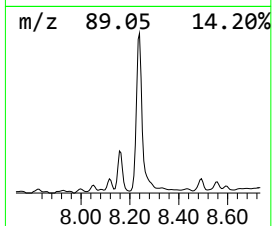
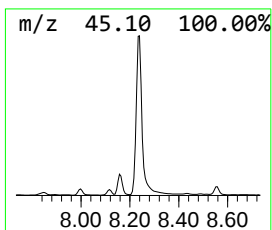
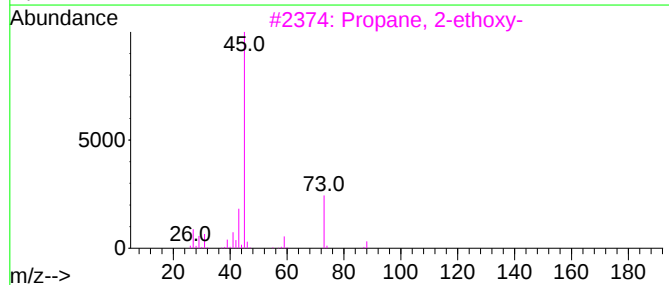
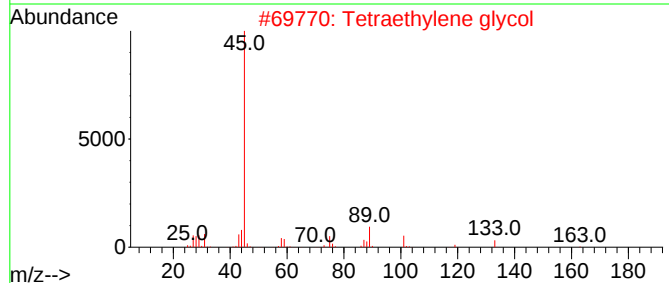
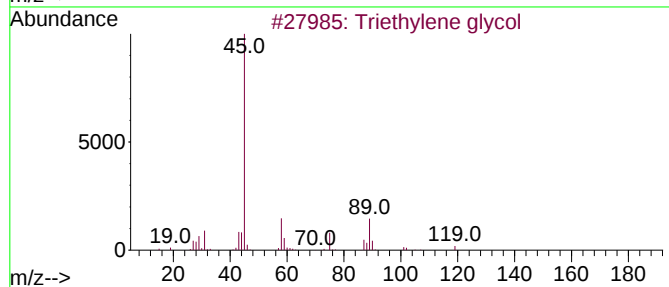
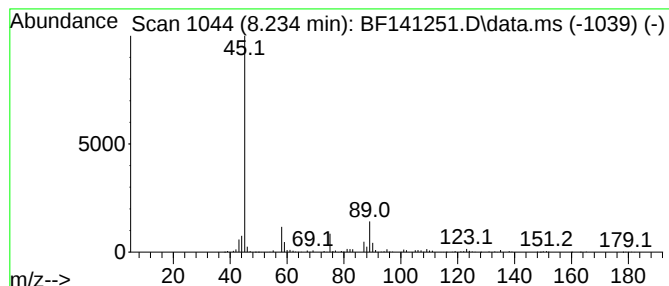
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Triethylene glycol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.234	10.71 ng	601734	Naphthalene-d8	8.087

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Triethylene glycol	150	C6H14O4	000112-27-6	90
2			Tetraethylene glycol	194	C8H18O5	000112-60-7	72
3			Propane, 2-ethoxy-	88	C5H12O	000625-54-7	42
4			Formic acid, 1-methylpropyl ester	102	C5H10O2	000589-40-2	42
5			Isopropyl Alcohol	60	C3H8O	000067-63-0	42



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF012325\
Data File : BF141251.D
Acq On : 23 Jan 2025 15:49
Operator : RC/JU
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Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
VNJ-220

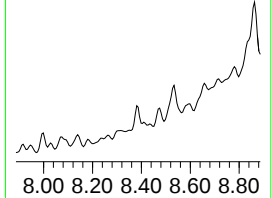
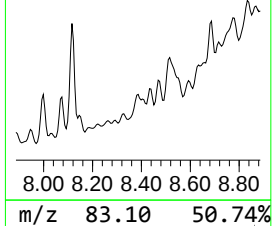
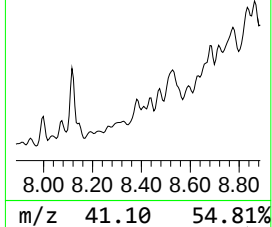
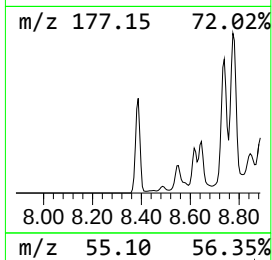
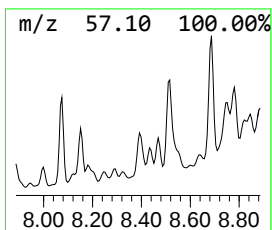
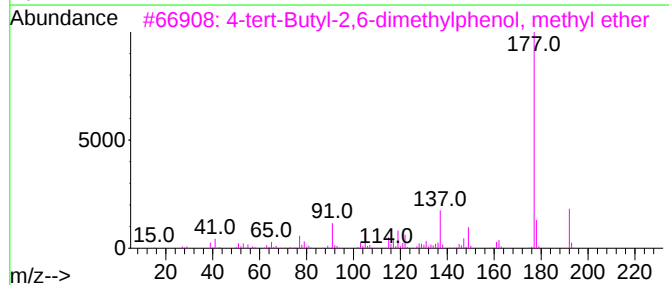
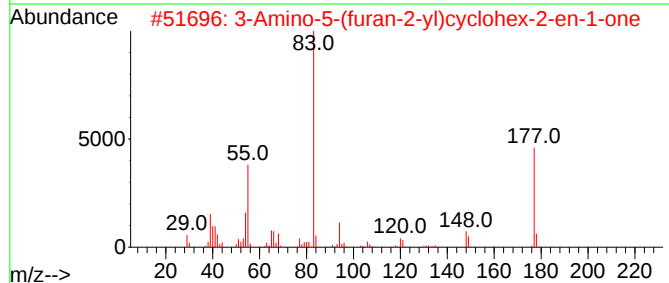
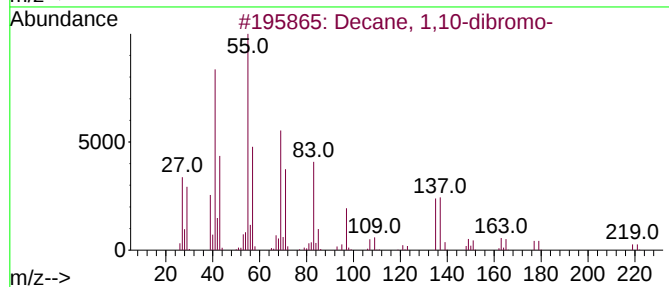
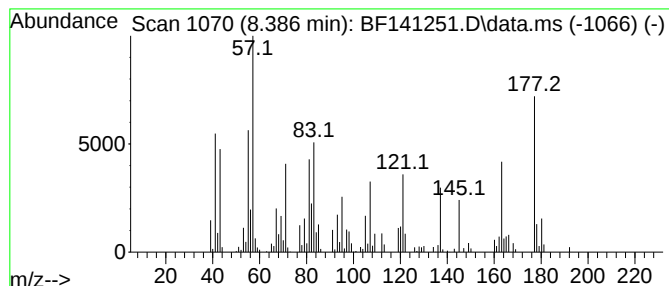
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown8.386 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.386	4.62 ng	259470	Naphthalene-d8	8.087

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 1,10-dibromo-	298	C10H20Br2	004101-68-2	25
2			3-Amino-5-(furan-2-yl)cyclohex-2-en-1-one	177	C10H11NO2	090921-87-2	22
3			4-tert-Butyl-2,6-dimethylphenol,...	192	C13H20O	092100-74-8	15
4			Benzaldehyde diisopentyl acetal	264	C17H28O2	1010431-61-7	12
5			trans-.beta.-Ionone	192	C13H20O	000079-77-6	10



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF012325\
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Acq On : 23 Jan 2025 15:49
Operator : RC/JU
Sample : Q1154-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
VNJ-220

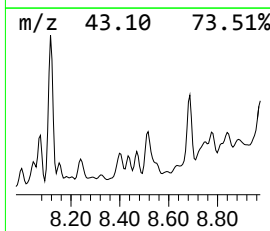
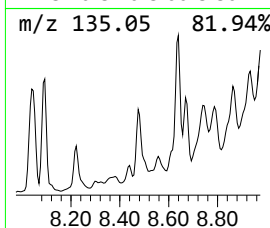
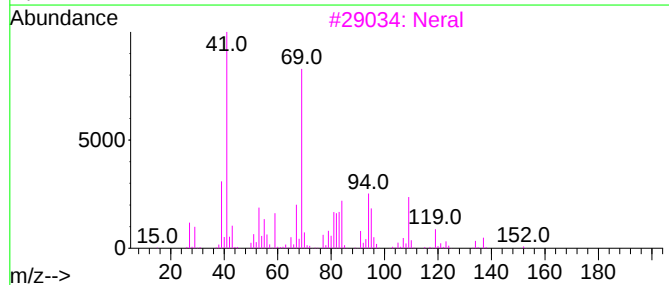
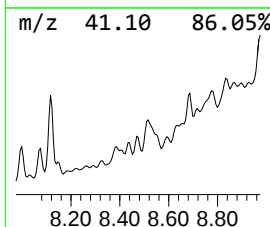
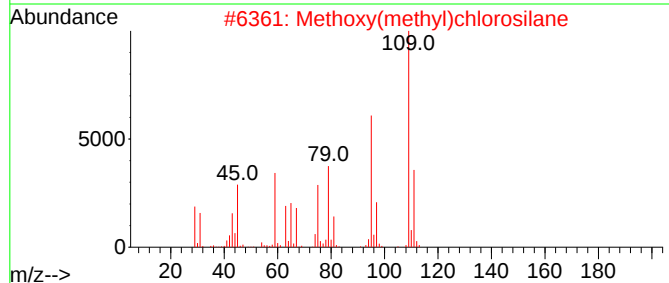
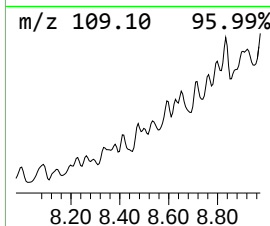
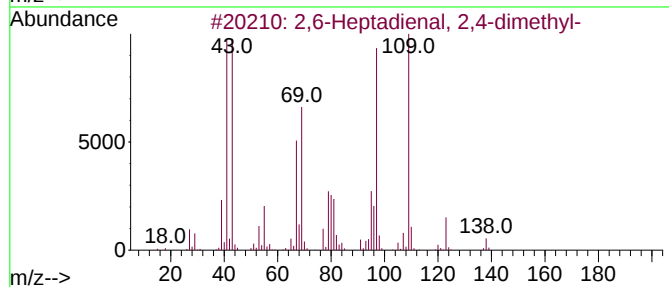
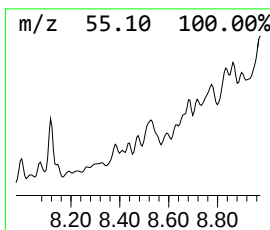
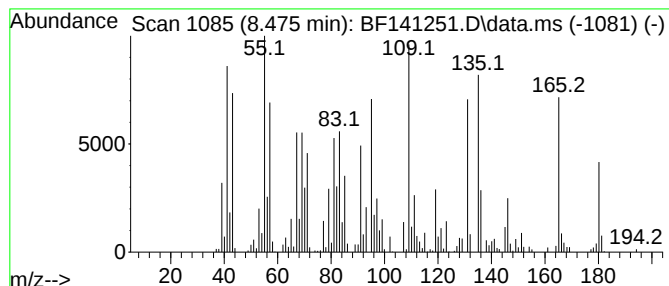
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown8.475 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.475	5.63 ng	316421	Naphthalene-d8	8.087

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,6-Heptadienal, 2,4-dimethyl-	138	C9H14O	085136-08-9	44
2			Methoxy(methyl)chlorosilane	110	C2H7ClOSi	018151-27-4	25
3			Neral	152	C10H16O	000106-26-3	25
4			Propionic acid, 2-(m-tolyloxy)-	180	C10H12O3	025140-95-8	18
5			Benzenepropanol, 4-methoxy-.gamm...	180	C11H16O2	018229-94-2	18



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF012325\
Data File : BF141251.D
Acq On : 23 Jan 2025 15:49
Operator : RC/JU
Sample : Q1154-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
VNJ-220

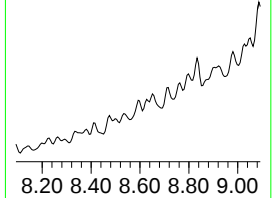
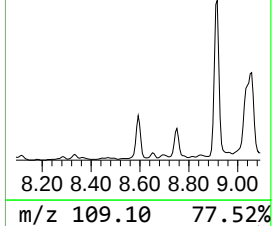
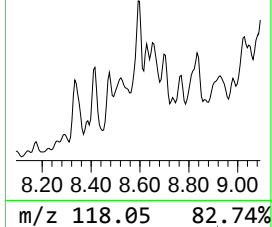
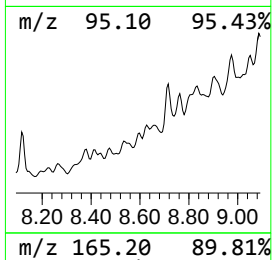
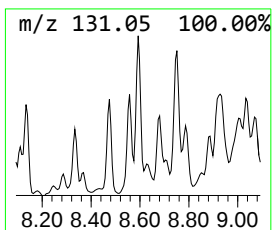
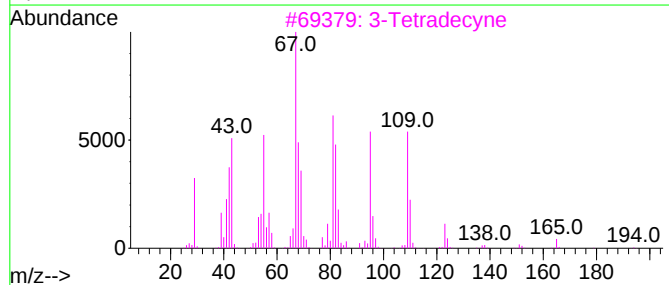
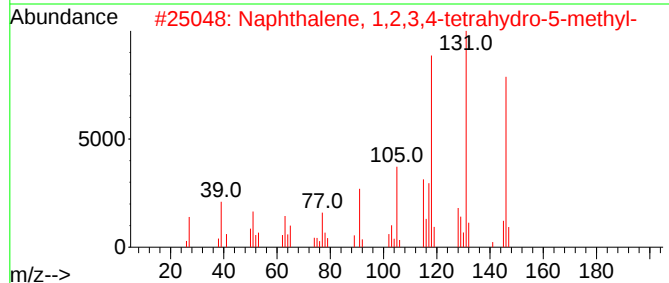
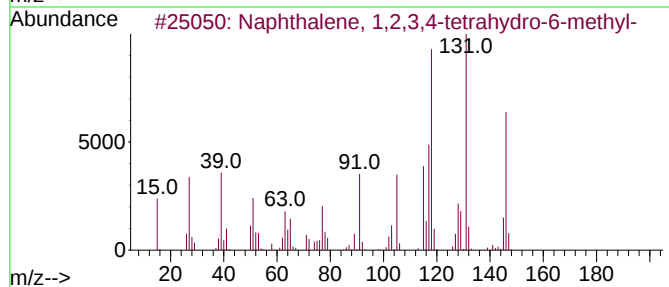
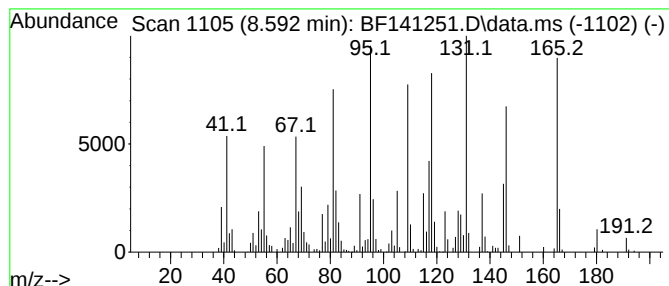
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.592	4.18 ng	234704	Naphthalene-d8	8.087

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	86
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	55
3			3-Tetradecyne	194	C14H26	060212-32-0	30
4			2,3,4,5,6,7-Hexahydro-1H-cyclope...	146	C11H14	1010189-31-0	25
5			cis,trans-1,6-Dimethylspiro[4.5]...	166	C12H22	1000111-72-3	18



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF012325\
Data File : BF141251.D
Acq On : 23 Jan 2025 15:49
Operator : RC/JU
Sample : Q1154-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
VNJ-220

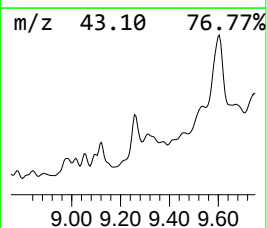
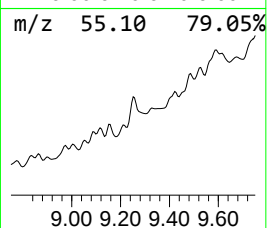
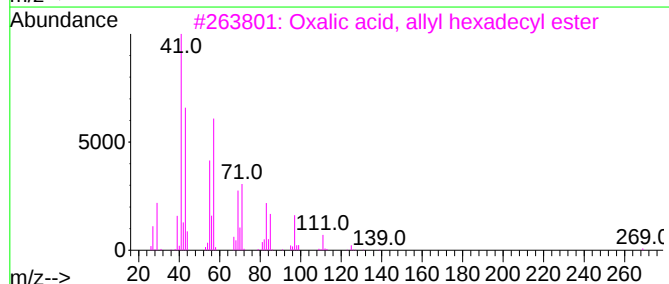
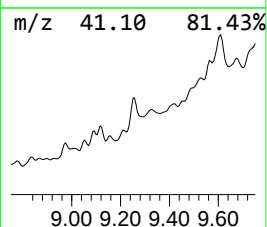
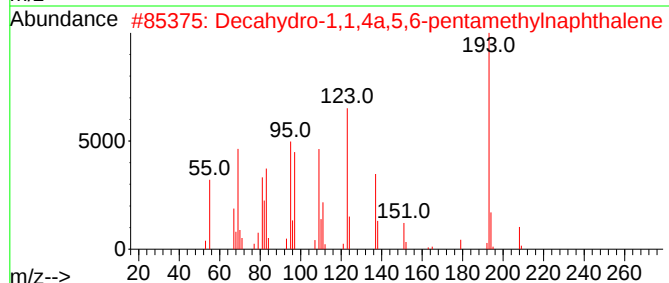
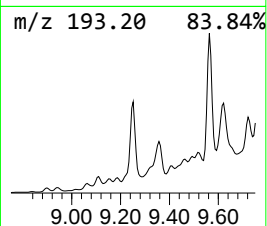
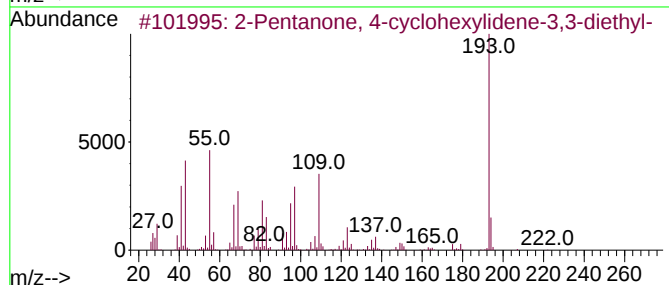
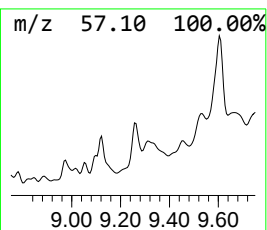
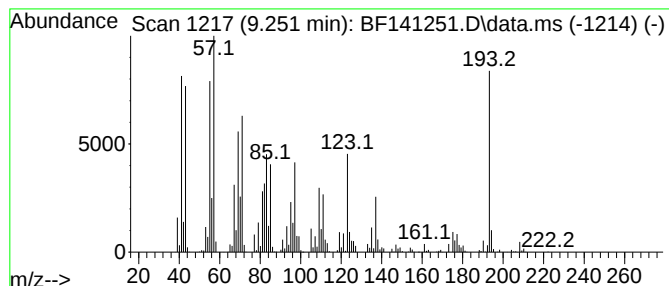
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 unknown9.251 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.251	20.00 ng	1311900	Acenaphthene-d10	9.857

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-cyclohexylidene-3...	222	C15H26O	313253-65-5	38
2			Decahydro-1,1,4a,5,6-pentamethyl...	208	C15H28	080655-44-3	38
3			Oxalic acid, allyl hexadecyl ester	354	C21H38O4	1000309-24-4	38
4			Oxalic acid, allyl pentadecyl ester	340	C20H36O4	1000309-24-3	30
5			Oxalic acid, cyclobutyl hexadecy...	368	C22H40O4	1000309-70-6	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF012325\
Data File : BF141251.D
Acq On : 23 Jan 2025 15:49
Operator : RC/JU
Sample : Q1154-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
VNJ-220

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF011025.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
1,3-Dioxolane, ...	2.234	14.8	ng	564028	1	6.804	762532	20.0
2-Pentanone, 4-...	5.004	7.7	ng	293236	1	6.804	762532	20.0
.alpha.-Pinene	6.087	94.8	ng	3614850	1	6.804	762532	20.0
Camphene	6.251	11.5	ng	437642	1	6.804	762532	20.0
Ethanol, 2,2'-o...	6.345	15.2	ng	577556	1	6.804	762532	20.0
2,5-Hexanediol	6.698	29.3	ng	1116050	1	6.804	762532	20.0
Hexanoic acid, ...	7.581	117.2	ng	6582720	2	8.087	1123360	20.0
Fenchol	7.622	8.1	ng	452103	2	8.087	1123360	20.0
Hexanoic acid, ...	7.675	4.8	ng	268319	2	8.087	1123360	20.0
endo-Borneol	7.998	10.0	ng	559606	2	8.087	1123360	20.0
Triethylene glycol	8.234	10.7	ng	601734	2	8.087	1123360	20.0
unknown8.386	8.386	4.6	ng	259470	2	8.087	1123360	20.0
unknown8.475	8.475	5.6	ng	316421	2	8.087	1123360	20.0
Naphthalene, 1,...	8.592	4.2	ng	234704	2	8.087	1123360	20.0
unknown9.251	9.251	20.0	ng	1311900	3	9.857	1311900	20.0