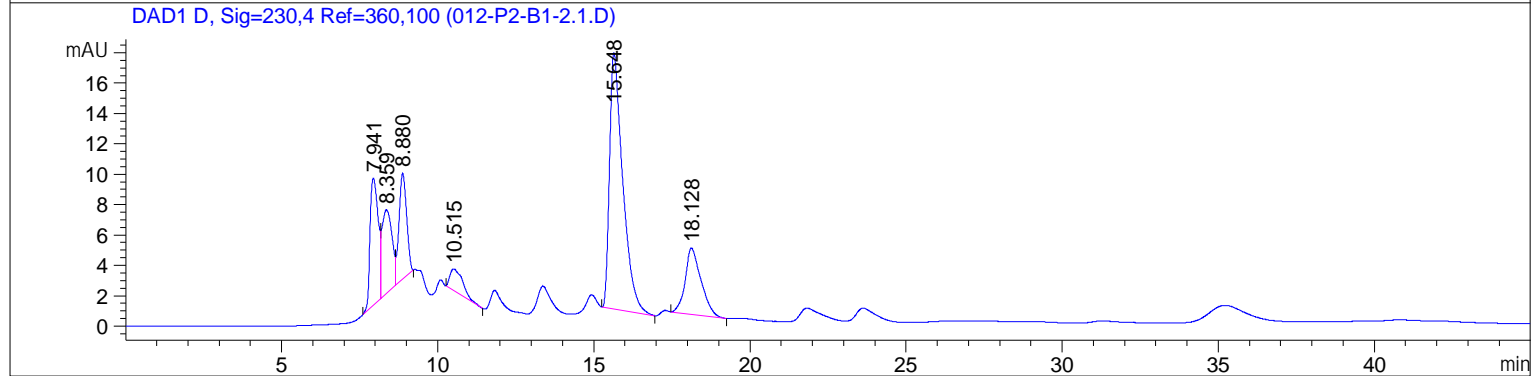
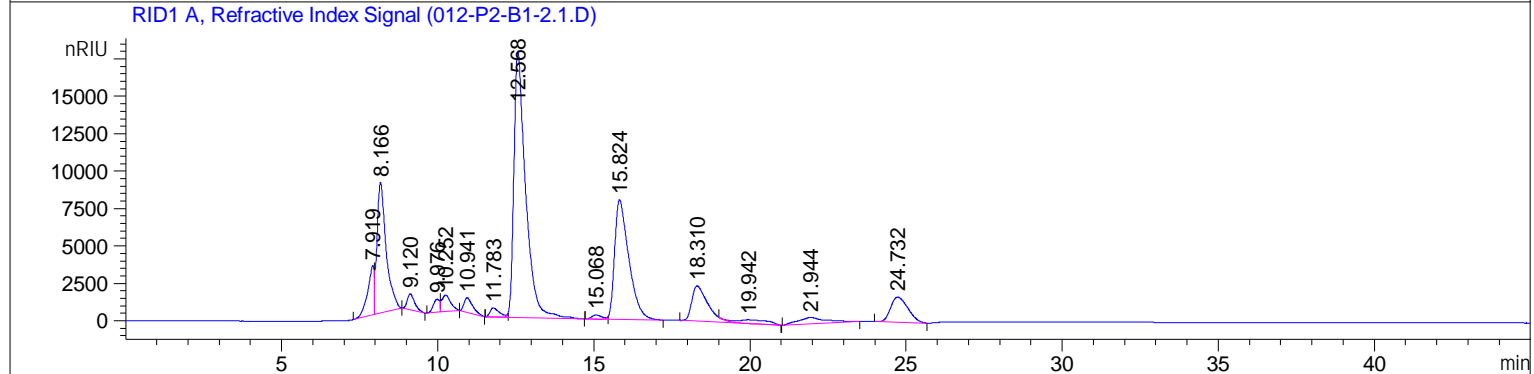
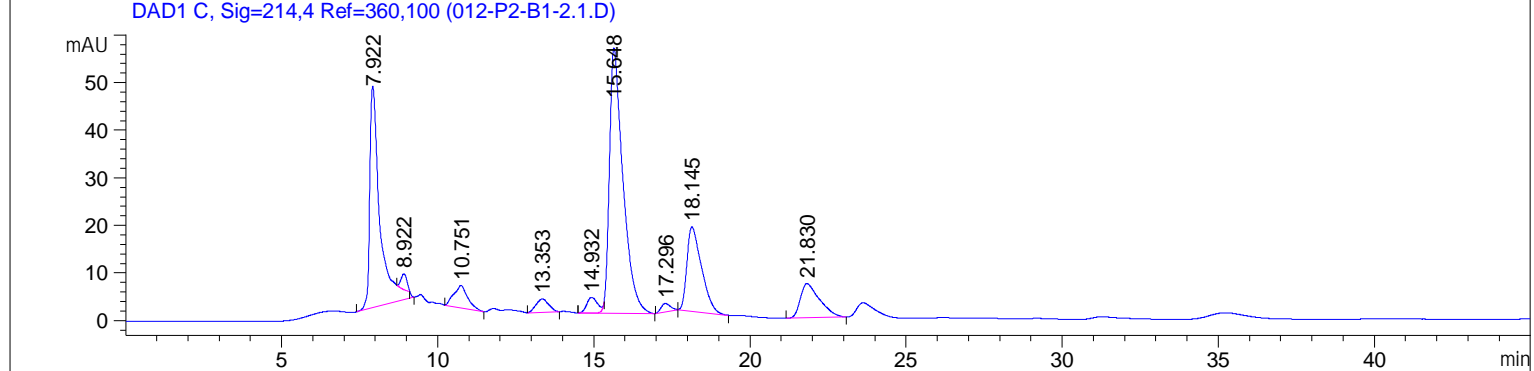


DAD1 B, Sig=210,4 Ref=360,100 (012-P2-B1-2.1.D)

The chromatogram displays absorbance (mAU) on the y-axis (0 to 50) against time (min) on the x-axis (0 to 45). The baseline is relatively flat with minor noise. Several distinct peaks are observed and labeled with their retention times. The peak at 15.648 min is the most intense, reaching approximately 50 mAU. Other significant peaks include those at 7.922 min (approx. 55 mAU) and 18.146 min (approx. 25 mAU). The peak at 21.830 min is also prominent, reaching about 15 mAU. The peak at 23.628 min is smaller, around 5 mAU. The peaks at 8.927, 10.752, 13.350, 14.926, and 17.295 min are all relatively small, generally below 10 mAU.

Retention Time (min)	Approximate Absorbance (mAU)
7.922	55
8.927	5
10.752	8
13.350	4
14.926	3
15.648	50
17.295	2
18.146	25
21.830	15
23.628	5



External Standard Report

Sorted By : Signal  
Calib. Data Modified : 7/1/2021 14:39:58  
Multiplier : 1.0000  
Dilution : 1.0000  
Do not use Multiplier & Dilution Factor with ISTDs

Signal 1: DAD1 B, Sig=210, 4 Ref=360, 100

RetTime [min]	Type	Area [mAU*s]	Amt/Area	Amount [g/L]	Grp	Name
15.648	VB R	1778.46887	6.56835e-4	1.16816		Lactic acid
18.146	BB	662.85474	9.16629e-4	6.07592e-1		Acetic acid
21.830	BB	490.69052	0.00000	0.00000		Propionic

Totals : 1.77575

Signal 2: DAD1 C, Sig=214, 4 Ref=360, 100

Signal 3: RID1 A, Refractive Index Signal

RetTime [min]	Type	Area [nRIU*s]	Amt/Area	Amount [g/L]	Grp	Name
9.976	BV	1.21446e4	0.00000	0.00000		Succrose
11.408		-	-	-		Glucose
12.568	VB R	4.65522e5	3.75900e-6	1.74989		Fructose
24.732	BB	6.44077e4	7.62465e-6	4.91087e-1		Ethanol

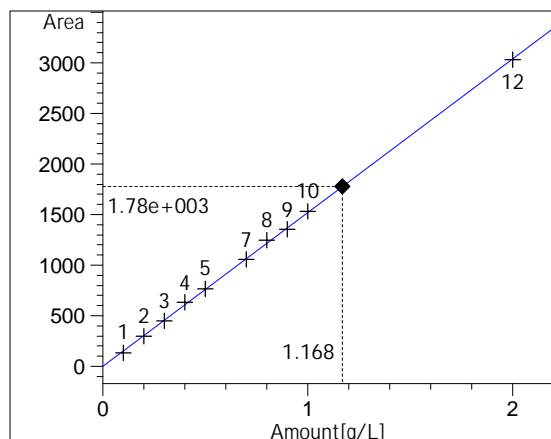
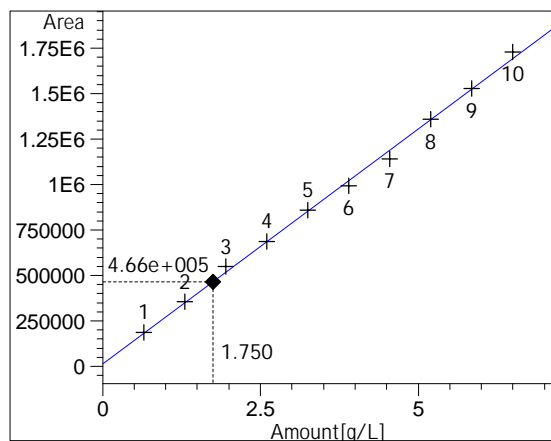
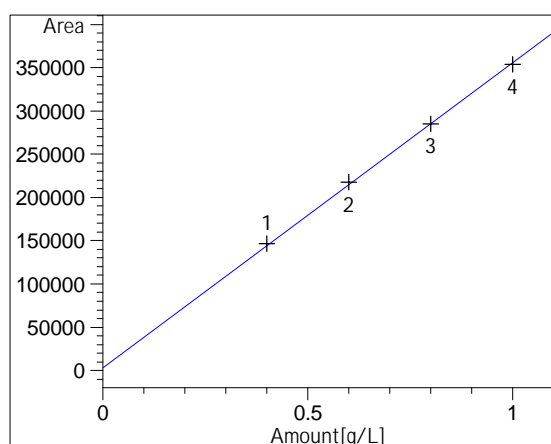
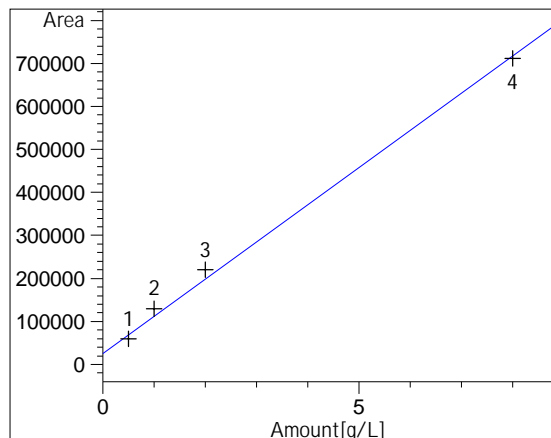
Totals : 2.24098

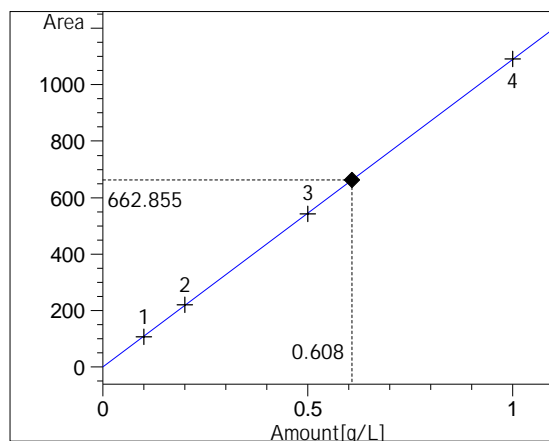
Signal 4: DAD1 D, Sig=230, 4 Ref=360, 100

4 Warnings or Errors :

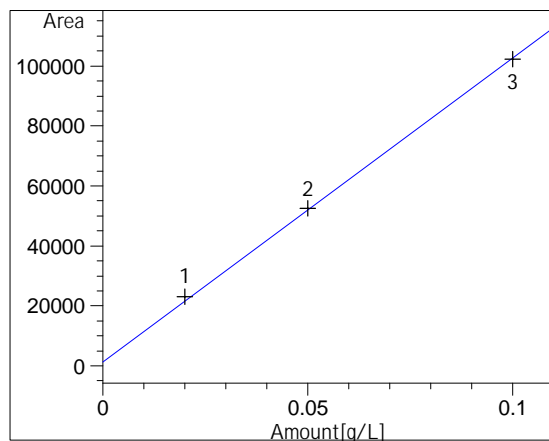
Warning : Calibration warnings (see calibration table listing)  
Warning : Calibrated compound(s) not found  
Warning : Negative results set to zero (cal. curve intercept), (Succrose)  
Warning : Negative results set to zero (cal. curve intercept), (Propionic)

# Calibration Curves

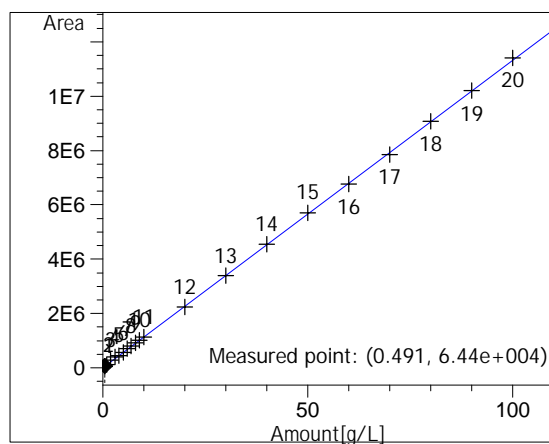




Acetic acid at exp. RT: 18.226  
DAD1 B, Sig=210,4 Ref=360,100  
Correlation: 0.99999  
Residual Std. Dev.: 2.57101  
Formula:  $y = mx + b$   
m: 1092.07931  
b: -6.83711e-1  
x: Amount[g/L]  
y: Area



Propionic at exp. RT: 21.787  
DAD1 B, Sig=210,4 Ref=360,100  
Correlation: 0.99961  
Residual Std. Dev.: 1507.31144  
Formula:  $y = mx + b$   
m: 1.01431e6  
b: 1395.13216  
x: Amount[g/L]  
y: Area



Ethanol at exp. RT: 24.978  
RID1 A, Refractive Index Signal  
Correlation: 0.99996  
Residual Std. Dev.: 35437.06772  
Formula:  $y = mx + b$   
m: 113284.07454  
b: 8775.42396  
x: Amount[g/L]  
y: Area

\*\*\* End of Report \*\*\*