Data File C:\Chem32\...ina-Xaris-23.10.23 2023-10-23 13-22-49\030-P2-C8-dimos 23\_10 2.72.D

Sample Name: dimos 23\_10 2.72

\_\_\_\_\_\_

Acq. Operator : SYSTEM Seq. Line : 30

Acq. Instrument : HPLC-0XTLAB Location : P2-C-08

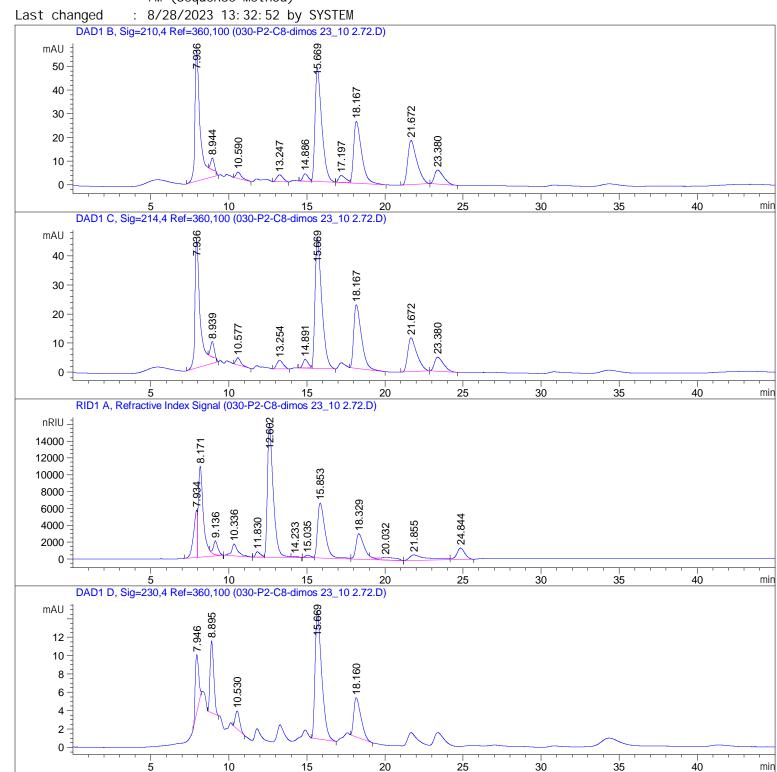
Inj Volume :  $20.000 \mu l$ 

 $Sequence\ File\ :\ C: \Chem32\1\Data\Dimos-Ntina-Xaris-23.\ 10.\ 23\ \ 2023-10-23\ \ 13-22-49\Dimos-Ntina-Xaris-23.\ \ 10.\ \ 2023-10-23\ \ 13-22-49\Dimos-Ntina-Xaris-23.\ \ 10.\ \ 2023-10-23\ \ \ 10-22-49\Dimos-Ntina-Xaris-23.\ \ 10.\ \ 10-22-49\Dimos-Ntina-Xaris-23.\ \ 10-23\Dimos-Ntina-Xaris-23.\ \ 10-23\Dimos-Ntina-Xaris-$ 

-Xari s-23. 10. 23. S

Method : C:\Chem32\1\Data\Dimos-Ntina-Xaris-23.10.23 2023-10-23 13-22-49\LACTIC\_TEMP

.M (Sequence Method)



Data File C:\Chem32\...ina-Xaris-23.10.23 2023-10-23 13-22-49\030-P2-C8-dimos 23\_10 2.72.D

Sample Name: dimos 23\_10 2.72

\_\_\_\_\_\_

## 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	Type	Area	Amt/Area	Amount	Grp Name
[min]		[mAU*s]		[g/L]	
15. 669	VB R	1429. 47583	6. 56750e-4	9. 38808e-1	Lactic acid
18. 167	VB R	926. 91150	9. 16360e-4	8. 49384e-1	Acetic acid
21. 672	BB	762. 31635	0.00000	0.00000	Propi oni c

Totals: 1. 78819

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

Signal 3: RID1 A, Refractive Index Signal

RetTime	Type	Area	Amt/Area	Amount	Grp	Name
[min]		[nRIU*s]		[g/L]		
					-	
9.863		-	_	_	S	uccrose
11. 408		-	-	-	G	Lucose
11.830	BV E	1. 39523e4	3. 22913e-7	4.50538e-3	F	ructose
24.844	VB	4. 92541e4	7. 25463e-6	3.57320e-1	Е	thanol

Total s : 3. 61825e-1

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

3 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), (Propionic)

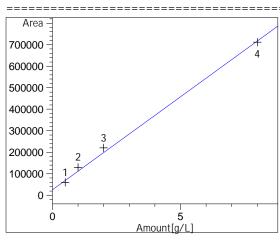
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Data File C:\Chem32\...ina-Xaris-23.10.23 2023-10-23 13-22-49\030-P2-C8-dimos 23\_10 2.72.D

Sample Name: dimos 23\_10 2.72

## -----

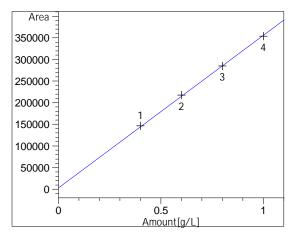
## Calibration Curves



Succrose at exp. RT: 9.863
RID1 A, Refractive Index Signal
Correlation: 0.99759
Residual Std. Dev.: 22755.41185

Formula: y = mx + b m: 86671.29673 b: 24839.01752 x: Amount[g/L]

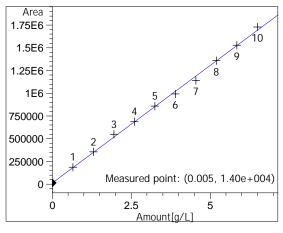
y: Area



Glucose at exp. RT: 11.408
RID1 A, Refractive Index Signal
Correlation: 0.99981
Residual Std. Dev.: 3023.36360

Formula: y = mx + b m: 353737.16216 b: 2639.18919 x: Amount[g/L]

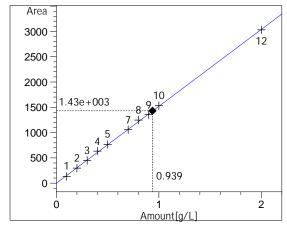
y: Area



Fructose at exp. RT: 12.100
RID1 A, Refractive Index Signal
Correlation: 0.99909
Residual Std. Dev.: 25081.95145

Formula: y = mx + b m: 258721.41259 b: 12786.68182 x: Amount[g/L]

y: Area



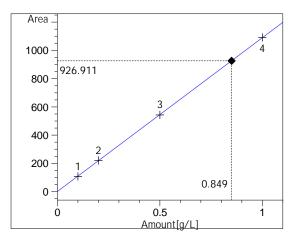
Lactic acid at exp. RT: 15.467 DAD1 B, Sig=210, 4 Ref=360, 100 Correlation: 0.99983 Residual Std. Dev.: 16.65547 Formula: v = mx + b

Formula: y = mx + bm: 1521.64235 b: 9.45650e-1

b: 9.45650e-7 x: Amount[g/L]

y: Area

Data File C:\Chem32\...ina-Xaris-23.10.23 2023-10-23 13-22-49\030-P2-C8-dimos 23\_10 2.72.D Sample Name: dimos 23\_10 2.72

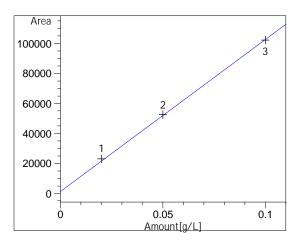


Acetic acid at exp. RT: 18.226 DAD1 B, Sig=210, 4 Ref=360, 100 Correlation: 0.99999 Residual Std. Dev.: 2.57101

Formula: y = mx + b m: 1092.07931 b: -6.83711e-1

x: Amount[g/L]
y: Area

y. Ale



Propionic at exp. RT: 21.787

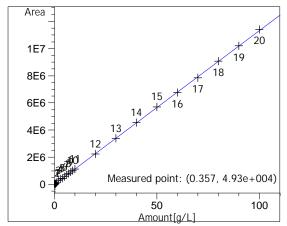
DAD1 B, Sig=210, 4 Ref=360, 100

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 + bm: 113284.07454 b: 8775.42396 x: Amount[g/L]

y: Area

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