

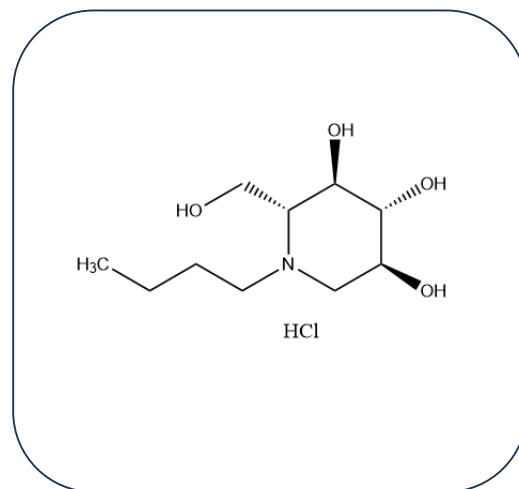
Certificate of Analysis

Catalogue Number

KL-01-00343

Lot Number

KLP0019AH01

Structure

Product Name

Miglustat hydrochloride

Synonyms

(2R,3R,4R,5S)-1-Butyl-2-(hydroxymethyl) piperidine-3,4,5-triol hydrochloride

CAS Number

210110-90-0

Molecular Formula
 $C_{10}H_{21}NO_4 \cdot HCl$
Molecular Weight

219.3 : 36.5 g/mol

Mfg. Date

30-08-2025

Re-Test Date

29-08-2028

Shipping Condition

All Product are stable to be shipped at room temperature, unless otherwise specified.

Long Term Storage Conditions

2°C ~ 8°C Protected from light and Sealed

Analytical Information

Test	Specifications	Results
Appearance	White to Off-White Solid (Hygroscopic)	Off-White Solid (Hygroscopic)
Solubility	Soluble in Methanol, Water	Methanol (Slightly), Water (Slightly)
Purity by HPLC	Not less than 90.00%	99.18%
Identification by 1H NMR	Conforms to structure	Complies
Identification by Mass	Conforms Molecular Weight	Complies



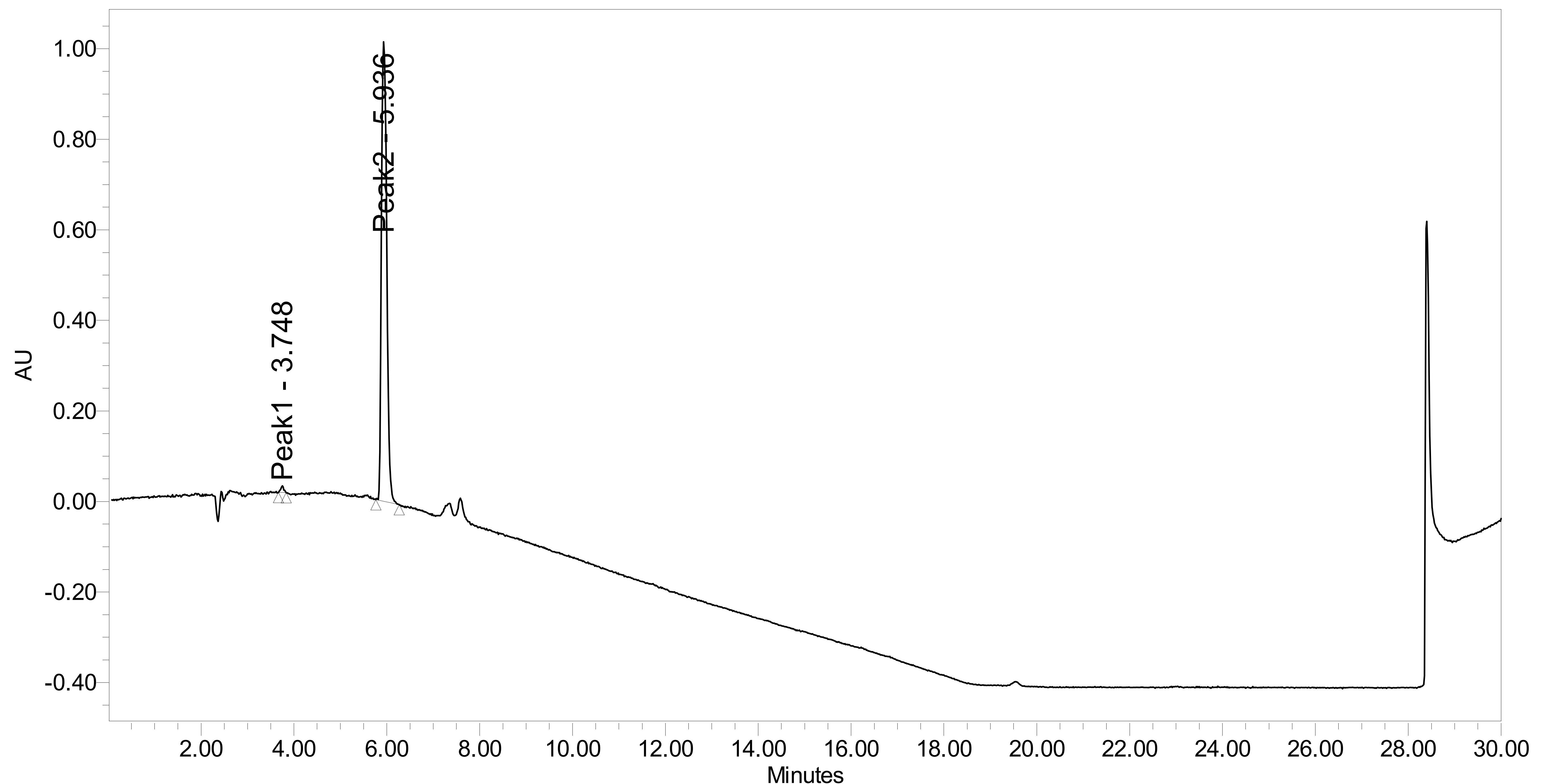
Note: Use for research & analytical purpose only. Not for human consumption

 Approved By B Kishore

SAMPLE INFORMATION

Sample Name: KLP0019AH01	Acquired By: System
Sample Type: Unknown	Sample Set Name: 30_08_2025
Vial: 43	Acq. Method Set: GEN_METHOD_06ml
Injection #: 1	Processing Method: MGS_
Injection Volume: 10.00 ul	Channel Name: 207.0nm
Run Time: 30.0 Minutes	Proc. Chnl. Descr.: PDA 207.0 nm
Date Acquired: 8/30/2025 7:33:09 PM IST	
Date Processed: 9/1/2025 2:22:03 PM IST	

MP-A:- 10MM AMMONIUM ACETATE in 1000ml of water
 MP-B:-100%ACN
 TIME:-%B:--0/10,15/90,25/90,25.1/10,30/10
 COLUM:---CRED CHROM C18 5um,,4.6X150mm
 FLOW RATE:-0.6ml/min
 RUN TIME:-30min



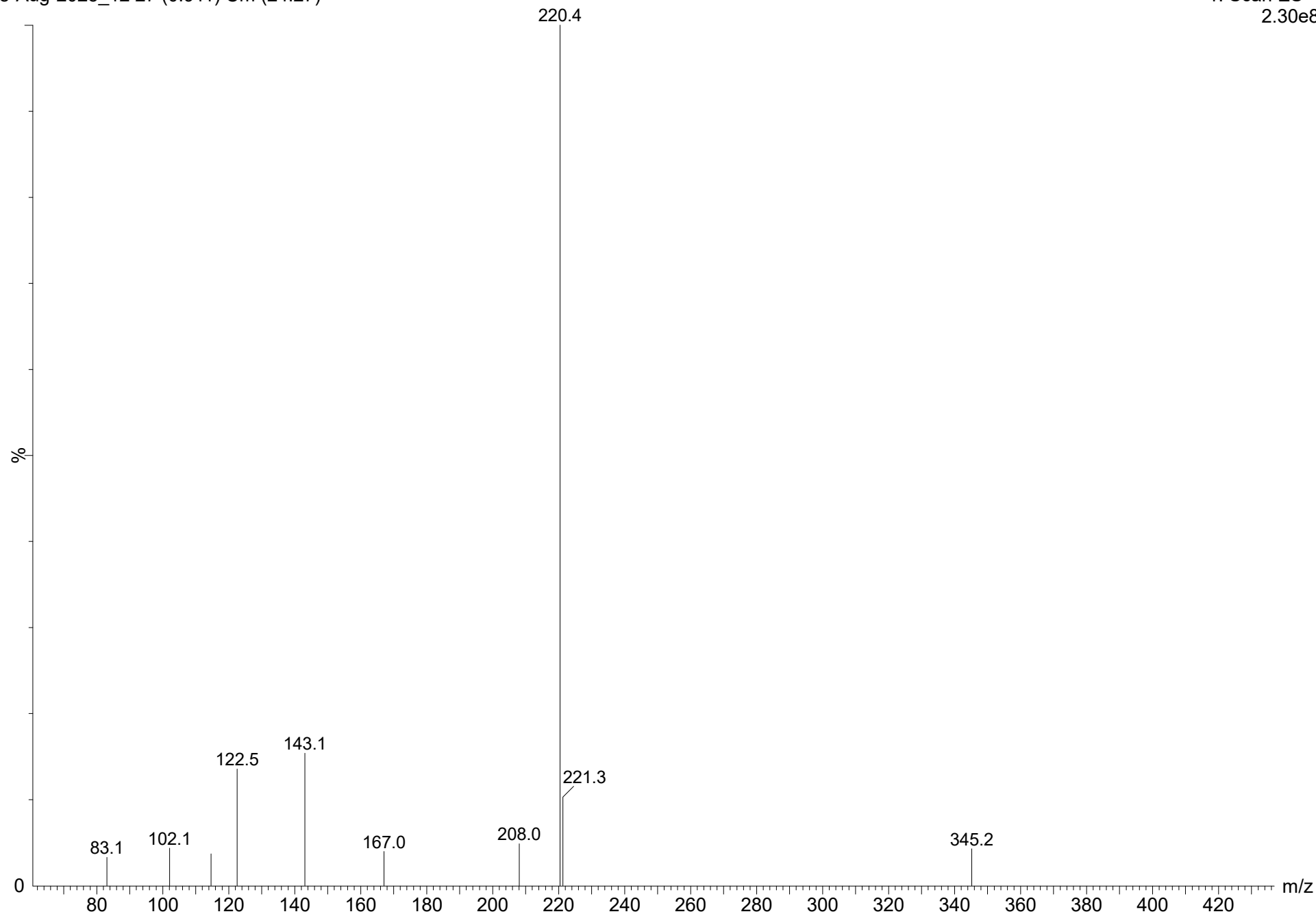
	Peak Name	RT	Area	% Area	Height
1	Peak1	3.748	66807	0.82	14262
2	Peak2	5.936	8068174	99.18	1011260

KLP0019AH01
037289

28-Aug-2025
15:18:26
ABNL10\1042

28-Aug-2025_12 27 (0.911) Cm (24:27)

1: Scan ES+
2.30e8



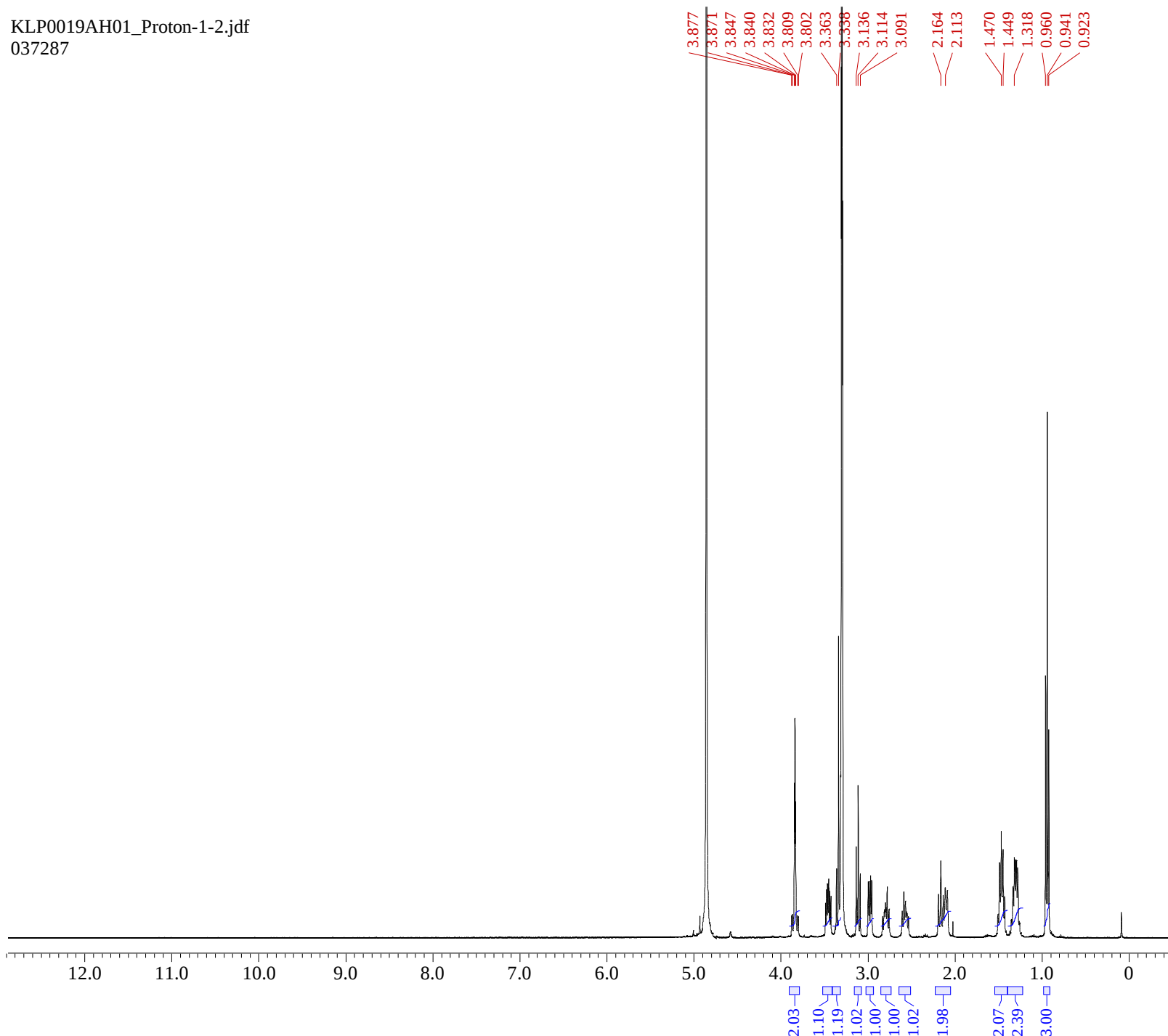
KLP0019AH01_Proton-1-2.jdf
037287



```
---- PROCESSING PARAMETERS ----  
sexp( 0.27483[Hz], 0.0[s] )  
trapezoid( 0[%], 0[%], 80[%], 100[%] )  
zerofill( 2, TRUE )  
fft( 1, TRUE, TRUE )  
machinephase  
ppm  
auto_reference( 5[%], TRUE )  
thresh( 0.15175[%], 1 )
```

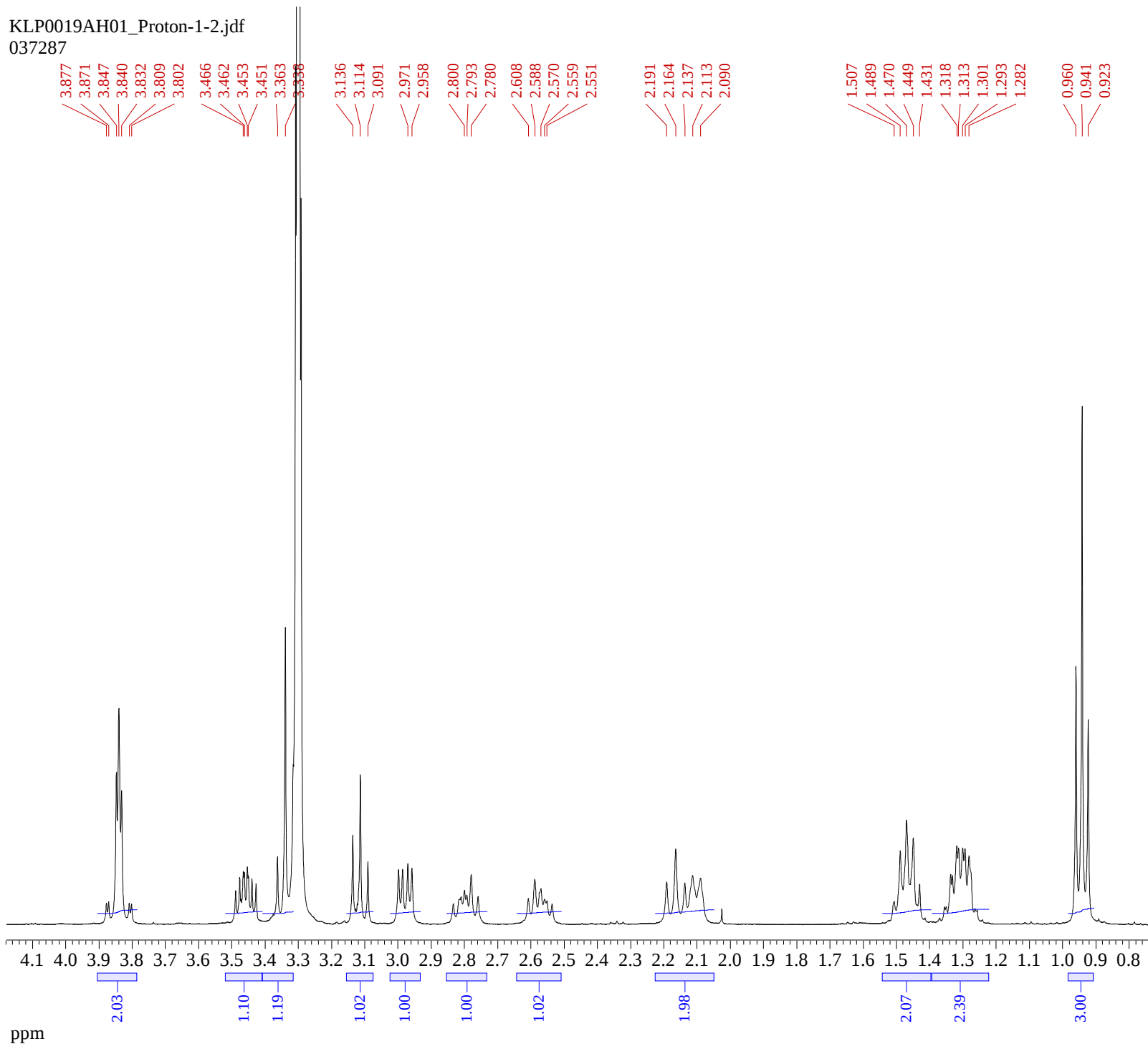
Derived from: MGS-IHRS_Proton-1-1.jdf

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Creation_Time   = 28-AUG-2025 18:00:31  
Instrument      = NMR-400MHz (JEOL)  
Spectrometer    = NM-70010S4L1  
Instrument id   = ABNL/QC/NMR-01  
Author         = 1033  
Reviewed by    = RMG  
Solvent        = METHANOL-D4  
Acquisition Parameters  
Experiment     = proton.jxp  
X_Offset       = 7[ppm]  
X_Sweep        = 9.00576369[kHz]  
Relaxation_Delay = 2[s]  
Scans          = 64
```



ppm

KLP0019AH01_Proton-1-2.jdf
037287



---- PROCESSING PARAMETERS ----
sexp(0.27483[Hz], 0.0[s])
trapezoid(0[%], 0[%], 80[%], 100[%])
zerofill(2, TRUE)
fft(1, TRUE, TRUE)
machinephase
ppm
auto_reference(5[%], TRUE)
thresh(0.15175[%], 1)

Derived from: MGS-IHRS_Proton-1-1.jdf

Creation_Time = 28-AUG-2025 18:00:31
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Solvent = METHANOL-D4
Acquisition Parameters
Experiment = proton.jxp
X_Offset = 7[ppm]
X_Sweep = 9.00576369[kHz]
Relaxation_Delay = 2[s]
Scans = 64

J-Coupling Analysis Report

Path = \\ABNL-01\Share folder\NMR Raw Data\Aug-2025\28-Aug-2025\KLP0019AH01_Proton-1-2.jdf

Position	Integral	Pattern	J
3.84 [ppm]	2	m	
3.46 [ppm]	1	m	
3.35 [ppm]	1	d	J1 = 9.619 [Hz]
3.11 [ppm]	1	t	J1 = 9.070 [Hz]
2.98 [ppm]	1	dd	J1 = 10.993 [Hz], J2 = 4.947 [Hz]
2.80 [ppm]	1	m	
2.57 [ppm]	1	m	
2.14 [ppm]	2	m	
1.47 [ppm]	2	m	
1.31 [ppm]	2	m	
0.94 [ppm]	3	t	J1 = 7.283 [Hz]