

Certificate of Analysis

SR-9011

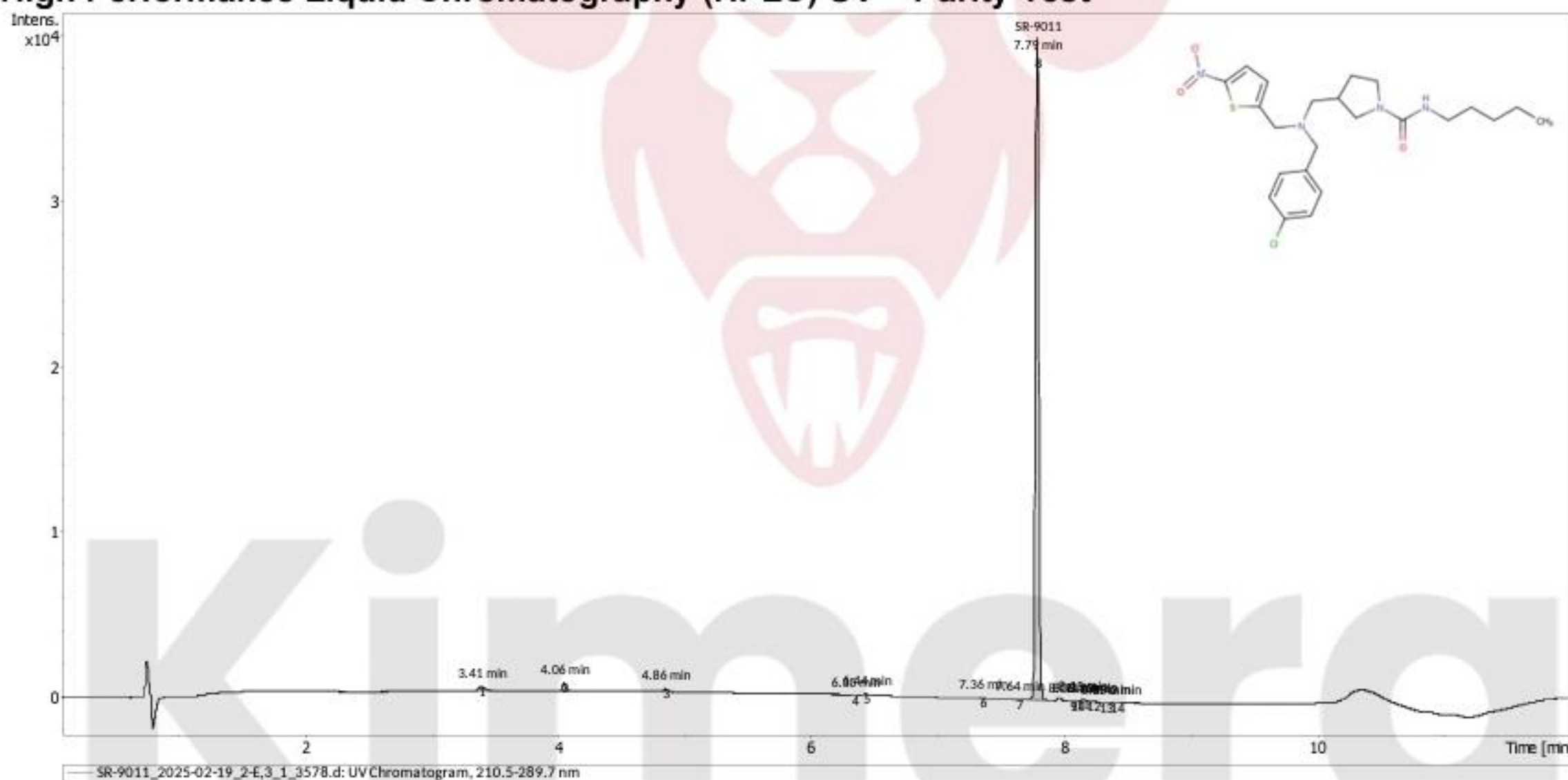
3-[[[(4-chlorophenyl)methyl-[(5-nitrothiophen-2-yl)methyl]amino)methyl]-N-pentylpyrrolidine-1-carboxamide

Compound : SR-9011
Lot number : 2025-02-19
Analysis date : 2025-02-27
Purity % : 96.82%
Method : HPLC-UV-MS

Client : Kimera Chems
Kimerachems.co

PubChem CID: 57394021
<https://pubchem.ncbi.nlm.nih.gov/compound/57394021>

High Performance Liquid Chromatography (HPLC) UV – Purity Test



SR-9011 DETECTED ON PEAK 8
SEE NEXT PAGE FOR PEAK LIST

Analysis Performed by
Ken Pendarvis, ChE
Analytical Chemist
MZ Biolabs
contact@mzbiolabs.com



2025-03-07

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PEAK LIST		Number of detected peaks: 14		
	Time (min)	Area	%Area	
1	3.41	8.09E+02	1.05	
2	4.06	7.68E+02	1.00	
3	4.86	2.81E+02	0.37	
4	6.35	1.54E+01	0.02	
5	6.44	1.26E+02	0.16	
6	7.36	4.56E+01	0.06	
7	7.64	8.39E+01	0.11	
8	7.79	7.44E+04	96.82	SR-9011
9	8.06	7.04E+01	0.09	
10	8.11	9.91E+00	0.01	
11	8.15	1.57E+02	0.20	
12	8.23	2.76E+01	0.04	
13	8.33	2.47E+01	0.03	
14	8.42	2.79E+01	0.04	

Kimera
CHEMS

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Mass Spectrometry (MS) – Identity Test

Identity confirmed using HPLC-MS

Molecular weight calculated using monoisotopic m/z values from mass spectrum

Expected monoisotopic mass : 478.18 Da

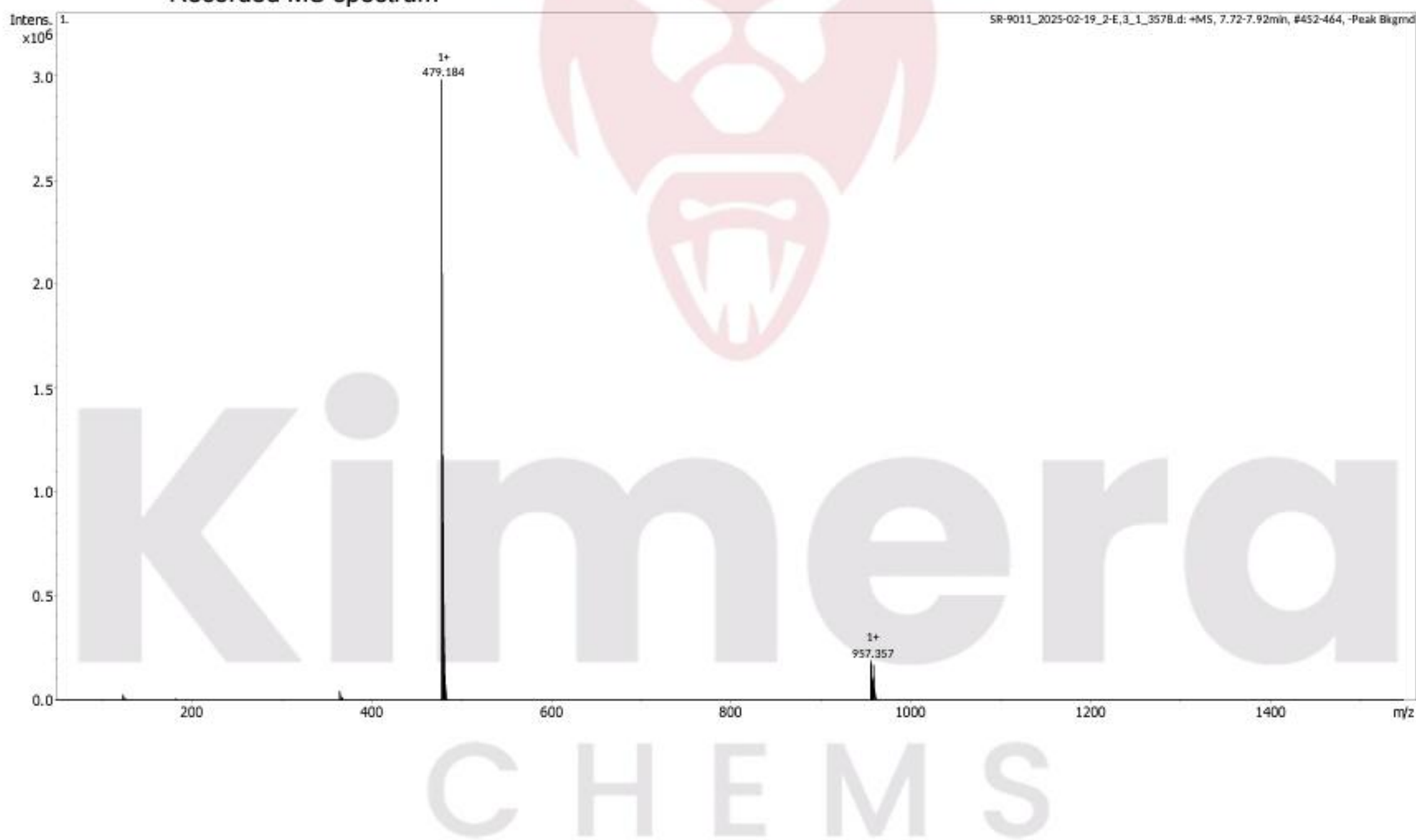
Measured monoisotopic mass : 478.18 Da

Molecular weight confirmed

Note : Monoisotopic m/z values are not easily seen in full spectrum view for larger molecules and peptides.

The dominant isotopic peak (base peak) shown in the spectrum below can be used to approximate the average molecular weight frequently reported by vendors and databases as a secondary means of confirmation.

Recorded MS spectrum



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