

Certificate of Analysis

Y-134

[6-hydroxy-2-(4-hydroxyphenyl)-1-benzothiophen-3-yl]-[4-(4-propan-2-ylpiperazin-1-yl)phenyl]methanone

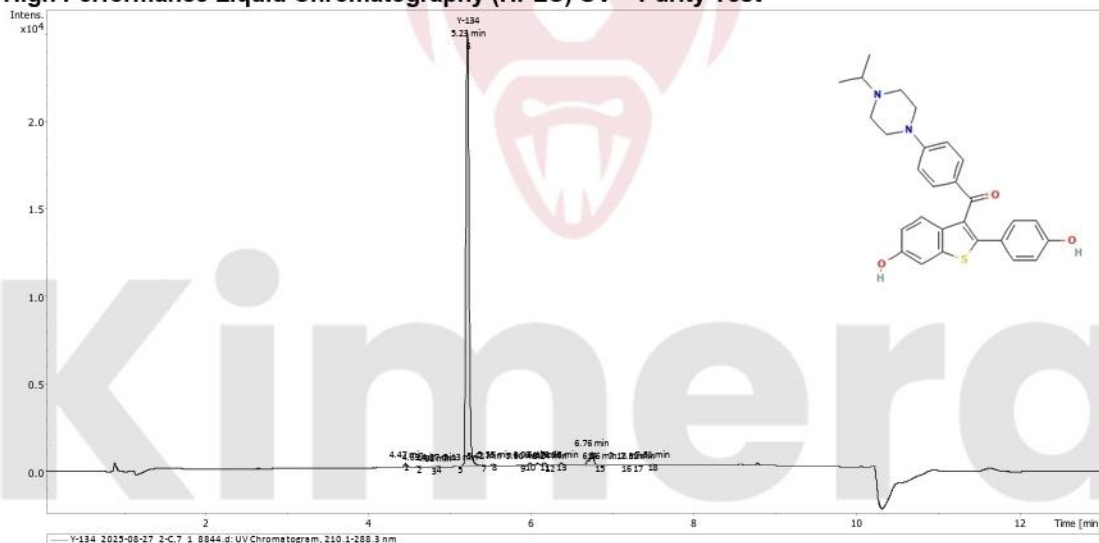
Compound : Y-134
Lot number : KC-Y34-02
Analysis date : 2025-09-30 retest
Purity % : 93.49%
Method : HPLC-UV-MS

Client : Kimera Chems
Kimerachems.co

PubChem CID: 11784736

<https://pubchem.ncbi.nlm.nih.gov/compound/11784736>

High Performance Liquid Chromatography (HPLC) UV – Purity Test



PEAK LIST
Time (min) Number of detected peaks: 18
Area %Area

Detailed peak list on next page

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



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PEAK LIST		Number of detected peaks: 18		
	Time (min)	Area	%Area	
1	4.47	4.64E+02	0.63	
2	4.63	1.23E+02	0.17	
3	4.81	2.69E+01	0.04	
4	4.87	2.42E+01	0.03	
5	5.13	3.55E+01	0.05	
6	5.23	6.89E+04	93.49	Y-134
7	5.42	3.56E+01	0.05	
8	5.55	2.58E+02	0.35	
9	5.90	1.06E+02	0.14	
10	6.00	3.03E+02	0.41	
11	6.17	2.78E+02	0.38	
12	6.24	2.46E+01	0.03	
13	6.38	3.56E+02	0.48	
14	6.76	2.58E+03	3.50	
15	6.86	2.15E+01	0.03	
16	7.18	4.50E+01	0.06	
17	7.32	5.38E+01	0.07	
18	7.51	6.23E+01	0.08	

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 : 472.18 Da

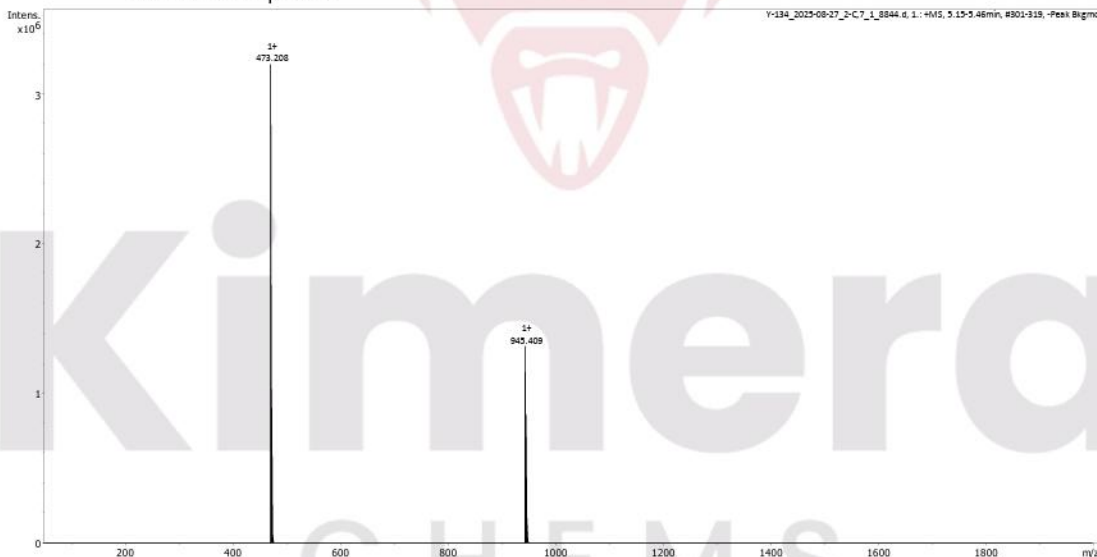
Measured monoisotopic mass : 472.21 Da

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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