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D C					
	ence Material Product Information Sheet				
Epichem's Quality System confe	orms to ISO9001:2015 as certified by ECAAS Pty Ltd - Certification number 616061.				
	$ \begin{array}{c}                                     $				
Name	(2-amino-1 <i>H</i> -benzimidazol-5-yl)(phenyl)methanone				
<b>BP/EP Name</b>	Mebendazole Impurity A				
Synonym(s)	2-amino-5-benzoylbenzimidazole				
Epichem Item #	EPL-AA32 Batch 7				
CAS #	52329-60-9				
Molecular Formula	C14H11N3O				
Molecular Weight	237.26 g/mol				
Appearance	Off-white powder				
Melting Point	193.6-196.2°C (decomposition)				
<b>Combustion Analysis</b>	Required (%): C:70.9; H:4.7; N:17.7. Found (%): C:71.1; H:4.8; N:17.9.				
Purity*	99.4%				
Date of Manufacture	27 November 2009				
Storage Requirements	Protect from heat, light and moisture.				
Special Precautions	This compound is for laboratory use only. Its toxicological properties may not				
-	have been fully established. It should be handled only by suitably qualified personnel.				
Intended Use	This compound is suitable for the identification of impurities and degradants in pharmaceutical materials. The purity assay is considered as relative contribution.				
Date of Shipment	TBA         This certificate is valid for one year from the date of shipment provided the substance is stored under the recommended conditions.				
Retest Date	TBA (Proper Storage and Handling Required)				

\* NATA accreditation does not cover the performance of this service

**Revision 4** 

 Epichem Pty Ltd, Suite 5, 3 Brodie-Hall Drive, Bentley WA 6102, Australia

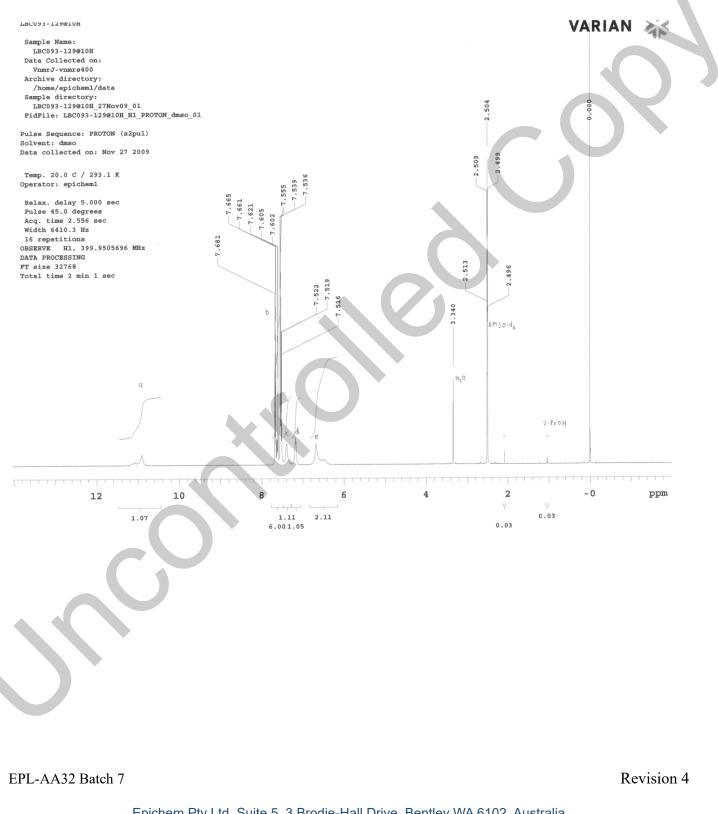
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 www.epichem.com.au
 ABN 80 106 769 902

# I. Identity

The identity of this product was established using the following analyses:

# Ia. <sup>1</sup>HNMR Spectrum

Conditions: 400 MHz, DMSO-d<sub>6</sub> <sup>1</sup>HNMR spectrum consistent with chemical structure.



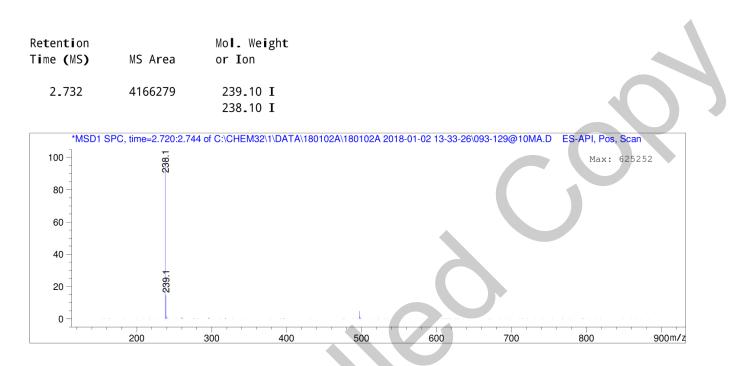
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#### **Ib. Mass Spectrum**

The mass spectrum of this material was analysed by Liquid Chromatography Mass Spectroscopy (LCMS) using in-house EM005.WI08.

Method: 5% to 100% ACN in water gradient (+0.1% formic acid) Zorbax Eclipse XDB-C8, 3.0 x 100 mm, 3.5 micron



Theoretical value: 238.1 [M+H]+.

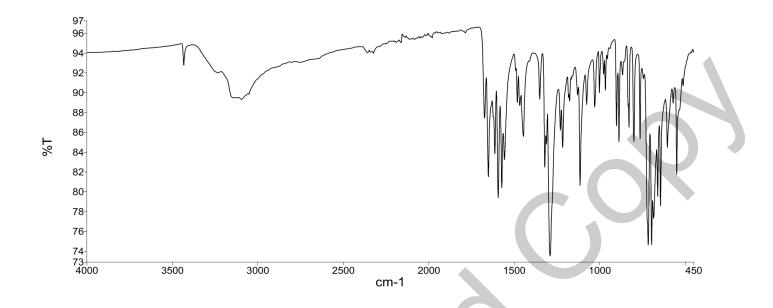
The signal of the Mass Spectrum is consistent with the theoretical value and its interpretation is consistent with the structural formula.

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## Ic. IR Spectrum

The infra-red spectrum of this material was analysed by Fourier-Transform Infrared Spectroscopy (FTIR) using in-house EM005.WI09.



The interpretation of the signals of the Fourier-Transform Infrared Spectrum is consistent with the structural formula.

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# **II.** Purity

The purity of this material was analysed by high performance liquid chromatography (HPLC) using inhouse EM005.WI07.

# **HPLC Conditions:**

C         C           e         % Line A (Water +           n)         0.1% (v/v) TFA)           0         90           0         60           0         5           0         5           0         90           0         5           0         90           0         90           0         90           0         90	% Line B (Acetonitrile + 0.1% (v/v) TFA) 10 40 95 95 10 10 10	Flow rate (mL/min) 1.0 1.0 1.0 1.0 1.0 1.0	DAD 293nm	Auto 1.0 μL 0.30 mg/mL in 50% acetonitrile 50% water (+0.1% TFA)
n)       0.1% (v/v) TFA)         0)       90         0)       60         0)       5         0)       5         0)       90         0)       5         0)       90         0)       90         0)       90         0)       90         0)       90         0)       90         4EM32/TA\221115A\221115A 2022-11-1	+ 0.1% (v/v) TFA) 10 40 95 95 10 10	(mL/min) 1.0 1.0 1.0 1.0 1.0	293nm	0.30 mg/mL in 50% acetonitrile 50% water
0         90           0         60           0         5           0         5           0         90           0         90           4EM32LTA\221115A\221115A 2022-11-1	10 40 95 95 10 10	1.0 1.0 1.0 1.0 1.0		50% acetonitrile 50% water
0         60           0         5           0         5           0         90           0         90	40 95 95 10 10	1.0 1.0 1.0 1.0		50% acetonitrile 50% water
0 5 0 5 0 90 0 90 HEM32\TA\221115A\221115A 2022-11-1	95 95 10 10	1.0 1.0 1.0		50% water
0 5 0 90 0 90 HEM32\TA\221115A\221115A 2022-11-1	95 10 10	1.0 1.0		
0 90 0 90	10 10	1.0		
0 90 HEM32\TA\221115A\221115A 2022-11-1	10			
ieM32\TA\221115A\221115A 2022-11-1		1.0		
5.301		)		
	10			
5				
4.748 5.769 6.769 6.056	8.585			
4 6	8 10	12	14	16 18 min
			- I Y C Y I C C I L C Y Y I C C Y Y C C Y	· [ x c x ] x c c [ x x x ] c c x 1 c x x [ x x x x

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#### Area Percent Report – Sorted by Signal

Peak Number	Retention Time (rounded)	Area	Area % (rounded)
1	3.27	0.25	0.03
2	4.75	0.08	0.01
3	5.30	809.82	99.66
4	5.77	0.16	0.02
5	6.06	1.94	0.24
6	7.67	0.15	0.02
7	8.58	0.13	0.02
8	10.23	0.03	0.00
Totals			100 (rounded)

For the calculation the system peaks were ignored. The content of the analyte was determined as a ratio of the peak area of the analyte and the cumulative areas of the purities, added up to 100%.

#### **Results:**

Average

99.7% (average of 10 duplicate runs)

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## **III. Water Content**

Method: Karl-Fischer titration using in-house EM005.WI04. **Results:** Average 0.1%

#### **IV. Ash Content**

Method: Combustion adjuvant added. **Result:** Contains 0.1% ash.

#### **V. Residual Solvents**

Method: <sup>1</sup>HNMR **Result:** Contains 0.1% Isopropanol by <sup>1</sup>H NMR analysis.

## VI. Final Result

Chromatographic purity (HPLC)	99.7%	
Water content	0.1%	
Ash content	0.1%	
Residual solvents	0.1%	
Purity*	99.4%	

This purity is assessed to be 99.4%.

Product Reviewed By:

Product Released By:

Jacob Heppell, PhD Chemist Carol Worth, PhD Quality Manager Release Date: 18 November 2022

\**NATA accreditation does not cover the performance of this service.* The calculation of the purity follows the formula:

 $Purity(\%) = \frac{((Chromatographicpurity[HPLC])x(100 - (watercontent + ashcontent + volatilecontents)))}{100}$ 

100

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