



# Full wwPDB X-ray Structure Validation Report ⓘ

May 23, 2020 – 02:29 pm BST

PDB ID : 2HJB  
Title : Crystal structure of *Alcaligenes faecalis* AADH in complex with p-methoxybenzylamine  
Authors : Roujeinikova, A.; Leys, D.  
Deposited on : 2006-06-30  
Resolution : 1.85 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.11

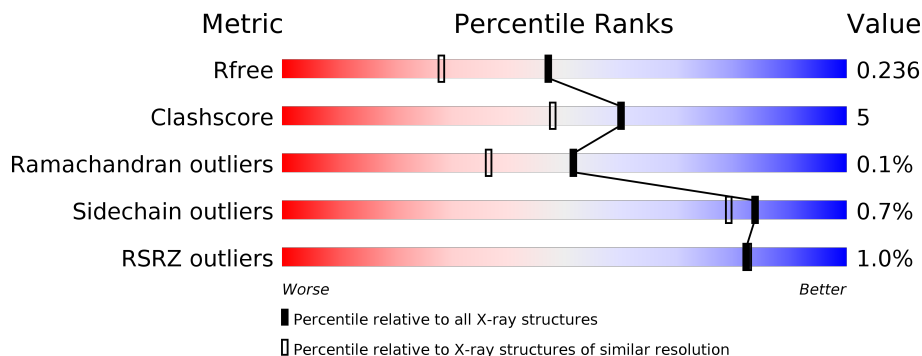
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.85 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	2469 (1.86-1.86)
Clashscore	141614	2625 (1.86-1.86)
Ramachandran outliers	138981	2592 (1.86-1.86)
Sidechain outliers	138945	2592 (1.86-1.86)
RSRZ outliers	127900	2436 (1.86-1.86)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	D	135	 2% (red), 70% (green), 10% (yellow), 19% (grey)
1	H	135	 2% (red), 80% (green), 10% (yellow), 10% (grey)
2	A	361	 % (red), 91% (green), 8% (yellow), % (grey)
2	B	361	 % (red), 92% (green), 6% (yellow), % (grey)

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 8466 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

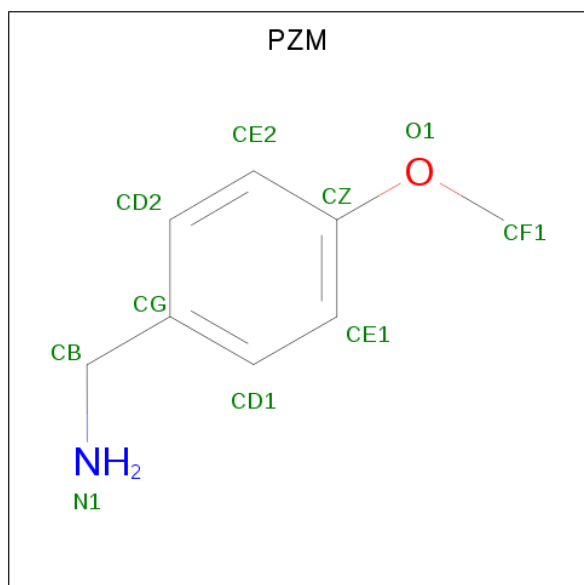
- Molecule 1 is a protein called Aromatic amine dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	D	110	Total 841	C 513	N 148	O 165	S 15	0	0	0
1	H	122	Total 925	C 564	N 162	O 184	S 15	0	0	0

- Molecule 2 is a protein called Aromatic amine dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	A	358	Total 2795	C 1764	N 485	O 532	S 14	0	1	0
2	B	358	Total 2794	C 1763	N 487	O 530	S 14	0	1	0

- Molecule 3 is 1-(4-METHOXYPHENYL)METHANAMINE (three-letter code: PZM) (formula: C<sub>8</sub>H<sub>11</sub>NO).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	D	1	Total	C	N	O	0	0
			10	8	1	1		
3	A	1	Total	C	N	O	0	0
			10	8	1	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	85	Total	O	0	0
			85	85		
4	H	116	Total	O	0	0
			116	116		
4	A	423	Total	O	0	0
			423	423		
4	B	467	Total	O	0	0
			467	467		

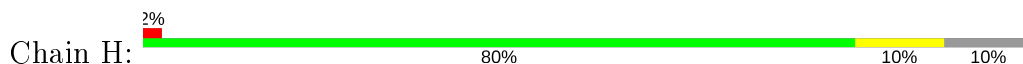
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

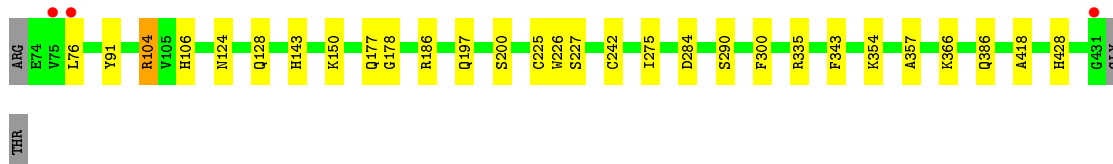
- Molecule 1: Aromatic amine dehydrogenase



- Molecule 1: Aromatic amine dehydrogenase



- Molecule 2: Aromatic amine dehydrogenase



- Molecule 2: Aromatic amine dehydrogenase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	91.33Å 96.60Å 119.67Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 1.85 23.48 – 1.85	Depositor EDS
% Data completeness (in resolution range)	88.2 (15.00-1.85) 88.2 (23.48-1.85)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.16 (at 1.85Å)	Xtrriage
Refinement program	REFMAC 5.2.0005	Depositor
R, $R_{free}$	0.170 , 0.232 0.179 , 0.236	Depositor DCC
$R_{free}$ test set	4003 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	23.1	Xtrriage
Anisotropy	0.357	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 58.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	8466	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	29.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.87% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: TRQ, PZM

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
1	D	0.51	0/847	0.66	0/1153
1	H	0.58	0/932	0.68	0/1271
2	A	0.57	0/2863	0.70	0/3880
2	B	0.59	0/2862	0.71	2/3880 (0.1%)
All	All	0.57	0/7504	0.69	2/10184 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
2	A	0	1
2	B	0	1
All	All	0	2

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
2	B	182	ASP	CB-CG-OD1	5.21	122.99	118.30
2	B	190	ASP	CB-CG-OD1	5.18	122.96	118.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	A	200	SER	Peptide
2	B	200	SER	Peptide

## 5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	841	0	737	16	0
1	H	925	0	811	11	0
2	A	2795	0	2736	30	0
2	B	2794	0	2731	21	0
3	A	10	0	11	5	0
3	D	10	0	11	3	0
4	A	423	0	0	3	0
4	B	467	0	0	3	0
4	D	85	0	0	2	0
4	H	116	0	0	1	0
All	All	8466	0	7037	66	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 5.

All (66) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:141:ASN:C	1:D:141:ASN:HD22	1.91	0.73
1:H:105:SER:HA	1:H:164:ASN:HD21	1.56	0.70
1:D:105:SER:HA	1:D:164:ASN:HD21	1.56	0.69
2:B:91:TYR:OH	2:B:428:HIS:HD2	1.76	0.69
2:A:91:TYR:OH	2:A:428:HIS:HD2	1.76	0.68
2:A:104:ARG:HH11	2:A:106:HIS:HE1	1.41	0.66
2:A:335:ARG:H	2:A:386:GLN:HE22	1.43	0.65
1:D:173:THR:HG21	2:A:76:LEU:HD13	1.78	0.65
1:D:109:TRQ:CH2	3:D:1102:PZM:HN12	2.10	0.65
2:A:197:GLN:NE2	2:A:227:SER:H	1.95	0.64
1:D:173:THR:HG21	2:A:76:LEU:CD1	2.30	0.62
3:D:1102:PZM:HF11	2:B:124:ASN:CG	2.21	0.60
2:B:257:LYS:NZ	4:B:877:HOH:O	2.36	0.58
2:B:197:GLN:NE2	2:B:227:SER:H	2.02	0.58
2:A:275:ILE:HA	2:A:290:SER:HA	1.86	0.58
1:D:71:GLU:N	4:D:1121:HOH:O	2.37	0.57
2:A:335:ARG:H	2:A:386:GLN:NE2	2.02	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:167:SER:HB2	2:B:177:GLN:HE22	1.69	0.57
2:A:124:ASN:CG	3:A:1101:PZM:HF11	2.24	0.57
2:A:197:GLN:HE21	2:A:227:SER:H	1.52	0.57
1:D:141:ASN:ND2	1:D:141:ASN:C	2.59	0.56
1:H:173:THR:HG21	2:B:76:LEU:HD13	1.89	0.55
1:H:159:ASN:HD22	2:A:177:GLN:NE2	2.04	0.55
1:H:109:TRQ:HB2	1:H:160:TRP:NE1	2.22	0.55
2:A:343:PHE:HB3	2:A:357:ALA:HB2	1.89	0.55
1:D:128:ASP:OD2	4:D:1186:HOH:O	2.18	0.54
1:D:109:TRQ:HB2	1:D:160:TRP:NE1	2.23	0.54
2:A:366:LYS:HE3	4:A:1515:HOH:O	2.08	0.53
1:D:159:ASN:HD22	2:B:177:GLN:NE2	2.06	0.53
1:H:149:GLY:HA3	2:A:226:TRP:CD1	2.45	0.53
2:A:143:HIS:HA	2:A:150:LYS:O	2.09	0.52
2:A:177:GLN:HB3	3:A:1101:PZM:HF13	1.92	0.52
2:A:284:ASP:HA	2:A:300:PHE:CE1	2.44	0.52
2:A:124:ASN:CB	3:A:1101:PZM:HF11	2.40	0.52
2:B:73:ARG:N	4:B:647:HOH:O	2.44	0.51
2:B:335:ARG:H	2:B:386:GLN:HE22	1.56	0.51
1:D:114:HIS:ND1	1:D:121:ASP:OD2	2.44	0.49
1:D:109:TRQ:O6	3:D:1102:PZM:N1	2.35	0.49
1:H:155:HIS:CE1	1:H:157:ASP:HB2	2.47	0.49
2:A:91:TYR:OH	2:A:428:HIS:CD2	2.62	0.48
2:B:104:ARG:HH11	2:B:106:HIS:HE1	1.61	0.48
2:A:178:GLY:HA2	3:A:1101:PZM:HE2	1.95	0.47
2:B:197:GLN:HE21	2:B:227:SER:H	1.61	0.47
2:B:231:GLN:NE2	2:B:234:ARG:HH11	2.13	0.47
2:B:257:LYS:NZ	4:B:865:HOH:O	2.48	0.47
2:B:91:TYR:OH	2:B:428:HIS:CD2	2.63	0.46
1:D:109:TRQ:HB2	1:D:160:TRP:HE1	1.79	0.46
2:A:124:ASN:OD1	3:A:1101:PZM:HF11	2.16	0.46
1:H:167:SER:HB2	2:A:177:GLN:HE22	1.81	0.45
2:A:225:CYS:HA	2:A:242:CYS:HA	1.99	0.45
2:B:275:ILE:HA	2:B:290:SER:HA	2.00	0.44
2:A:128:GLN:HE22	2:A:186:ARG:C	2.22	0.44
2:B:335:ARG:H	2:B:386:GLN:NE2	2.16	0.43
1:H:145:ARG:HD2	4:H:243:HOH:O	2.18	0.43
1:D:143:GLN:NE2	2:B:354:LYS:NZ	2.67	0.43
2:A:104:ARG:HH11	2:A:106:HIS:CE1	2.27	0.42
2:A:366:LYS:CE	4:A:1515:HOH:O	2.67	0.42
2:A:106:HIS:HD2	2:A:418:ALA:O	2.03	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:143:GLN:CD	2:A:354:LYS:HD2	2.40	0.42
1:D:127:HIS:HB2	2:A:76:LEU:HD22	2.02	0.42
2:B:106:HIS:HD2	2:B:418:ALA:O	2.03	0.41
1:H:63:ASN:HA	1:H:64:PRO:HD2	1.93	0.40
2:A:143:HIS:HD2	4:A:1203:HOH:O	2.04	0.40
2:B:200:SER:N	2:B:201:PRO:HA	2.36	0.40
1:H:173:THR:HG21	2:B:76:LEU:CD1	2.51	0.40
2:B:128:GLN:HE22	2:B:186:ARG:C	2.24	0.40

There are no symmetry-related clashes.

### 5.3 Torsion angles [i](#)

#### 5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	107/135 (79%)	102 (95%)	5 (5%)	0	100	100
1	H	119/135 (88%)	117 (98%)	2 (2%)	0	100	100
2	A	357/361 (99%)	344 (96%)	13 (4%)	0	100	100
2	B	357/361 (99%)	344 (96%)	12 (3%)	1 (0%)	41	26
All	All	940/992 (95%)	907 (96%)	32 (3%)	1 (0%)	51	36

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	201	PRO

#### 5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	95/112 (85%)	92 (97%)	3 (3%)	39	22
1	H	104/112 (93%)	104 (100%)	0	100	100
2	A	304/305 (100%)	303 (100%)	1 (0%)	92	91
2	B	303/305 (99%)	301 (99%)	2 (1%)	84	79
All	All	806/834 (97%)	800 (99%)	6 (1%)	84	79

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	121	ASP
1	D	141	ASN
1	D	176	LEU
2	A	104	ARG
2	B	104	ARG
2	B	218	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (34) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	139	GLN
1	D	141	ASN
1	D	143	GLN
1	D	164	ASN
1	H	143	GLN
1	H	164	ASN
2	A	86	GLN
2	A	106	HIS
2	A	128	GLN
2	A	143	HIS
2	A	177	GLN
2	A	180	ASN
2	A	187	GLN
2	A	197	GLN
2	A	231	GLN
2	A	261	GLN
2	A	266	GLN
2	A	386	GLN
2	A	424	GLN

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Mol	Chain	Res	Type
2	A	426	GLN
2	A	428	HIS
2	B	86	GLN
2	B	106	HIS
2	B	128	GLN
2	B	143	HIS
2	B	177	GLN
2	B	187	GLN
2	B	197	GLN
2	B	231	GLN
2	B	322	ASN
2	B	386	GLN
2	B	424	GLN
2	B	426	GLN
2	B	428	HIS

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

2 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
1	TRQ	D	109	1	13,17,18	5.43	5 (38%)	14,24,26	2.19	3 (21%)
1	TRQ	H	109	1	13,17,18	5.36	6 (46%)	14,24,26	2.30	4 (28%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	TRQ	D	109	1	-	0/4/19/21	0/2/2/2
1	TRQ	H	109	1	-	0/4/19/21	0/2/2/2

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	109	TRQ	CH2-CZ2	-15.63	1.36	1.54
1	H	109	TRQ	CH2-CZ2	-14.63	1.37	1.54
1	D	109	TRQ	CE2-CZ2	-7.56	1.40	1.50
1	H	109	TRQ	CE2-CZ2	-7.14	1.40	1.50
1	H	109	TRQ	CZ3-CE3	6.33	1.45	1.34
1	D	109	TRQ	CZ3-CE3	5.98	1.44	1.34
1	H	109	TRQ	O7-CZ2	5.77	1.35	1.23
1	D	109	TRQ	O7-CZ2	5.08	1.34	1.23
1	H	109	TRQ	O6-CH2	4.50	1.36	1.24
1	D	109	TRQ	O6-CH2	3.67	1.34	1.24
1	H	109	TRQ	CD2-CE3	-2.05	1.40	1.44

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	109	TRQ	CZ2-CE2-NE1	5.99	129.50	119.94
1	D	109	TRQ	CZ2-CE2-NE1	5.84	129.27	119.94
1	D	109	TRQ	O6-CH2-CZ2	-3.49	116.13	118.51
1	H	109	TRQ	O7-CZ2-CH2	3.17	122.72	119.00
1	D	109	TRQ	CD2-CE3-CZ3	-3.05	117.40	121.09
1	H	109	TRQ	CD2-CE3-CZ3	-2.81	117.69	121.09
1	H	109	TRQ	O7-CZ2-CE2	-2.70	118.98	121.84

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	D	109	TRQ	4	0
1	H	109	TRQ	1	0

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry

2 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	PZM	D	1102	-	10,10,10	4.75	3 (30%)	12,12,12	1.72	2 (16%)
3	PZM	A	1101	-	10,10,10	4.00	3 (30%)	12,12,12	0.85	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	PZM	D	1102	-	-	2/4/4/4	0/1/1/1
3	PZM	A	1101	-	-	2/4/4/4	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	1102	PZM	CD2-CE2	14.36	1.64	1.38
3	A	1101	PZM	CD2-CE2	11.98	1.60	1.38
3	A	1101	PZM	CE1-CD1	2.79	1.43	1.38
3	D	1102	PZM	O1-CF1	-2.44	1.35	1.42
3	A	1101	PZM	O1-CZ	2.39	1.42	1.37
3	D	1102	PZM	CE1-CD1	2.37	1.43	1.38

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	1102	PZM	CF1-O1-CZ	-4.57	107.58	117.51
3	D	1102	PZM	CD2-CE2-CZ	-2.28	116.94	119.73

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	1102	PZM	CE2-CZ-O1-CF1
3	D	1102	PZM	CE1-CZ-O1-CF1
3	A	1101	PZM	CE2-CZ-O1-CF1
3	A	1101	PZM	CE1-CZ-O1-CF1

There are no ring outliers.

2 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	1102	PZM	3	0
3	A	1101	PZM	5	0

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	D	109/135 (80%)	-0.08	3 (2%) 53 52	20, 30, 53, 64	0
1	H	121/135 (89%)	-0.13	3 (2%) 57 56	18, 27, 45, 64	0
2	A	358/361 (99%)	-0.31	3 (0%) 86 86	17, 27, 43, 71	0
2	B	358/361 (99%)	-0.54	0 100 100	17, 24, 37, 55	0
All	All	946/992 (95%)	-0.35	9 (0%) 82 82	17, 26, 43, 71	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	A	75	VAL	6.0
2	A	431	GLY	5.2
2	A	76	LEU	3.1
1	D	179	LEU	2.7
1	H	59	HIS	2.6
1	H	180	ALA	2.4
1	H	60	ILE	2.4
1	D	117	HIS	2.2
1	D	119	GLY	2.1

### 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
1	TRQ	D	109	16/17	0.95	0.10	27,30,35,43	0
1	TRQ	H	109	16/17	0.96	0.12	21,27,31,38	0



### 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

### 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	PZM	A	1101	10/10	0.94	0.16	23,30,40,44	0
3	PZM	D	1102	10/10	0.97	0.09	11,32,35,40	0

### 6.5 Other polymers [i](#)

There are no such residues in this entry.