



# Full wwPDB X-ray Structure Validation Report ⓘ

Nov 25, 2024 – 04:51 PM EST

PDB ID : 2OZV  
Title : Crystal structure of a predicted O-methyltransferase, protein Atu636 from *Agrobacterium tumefaciens*.  
Authors : Cuff, M.E.; Xu, X.; Zheng, X.; Edwards, A.; Savchenko, A.; Joachimiak, A.; Midwest Center for Structural Genomics (MCSG)  
Deposited on : 2007-02-27  
Resolution : 1.70 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.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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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	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.21
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.004 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.40

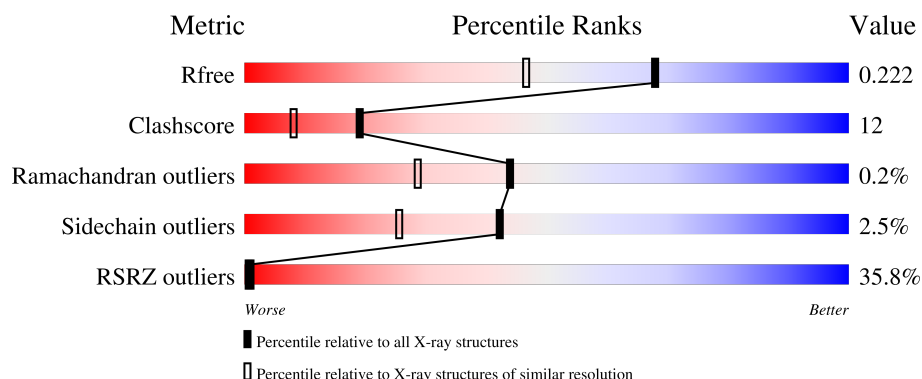
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.70 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))
$R_{free}$	164625	5161 (1.70-1.70)
Clashscore	180529	5671 (1.70-1.70)
Ramachandran outliers	177936	5594 (1.70-1.70)
Sidechain outliers	177891	5594 (1.70-1.70)
RSRZ outliers	164620	5159 (1.70-1.70)

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 of 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  $\leq 5\%$ .

Mol	Chain	Length	Quality of chain
1	A	260	<div> <div>28%</div> <div>67%</div> <div>13%</div> <div>20%</div> </div>
1	B	260	<div> <div>27%</div> <div>64%</div> <div>15%</div> <div>•</div> <div>19%</div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4033 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 Hypothetical protein Atu0636.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	208	Total	C	N	O	S	Se	0	23	0
			1759	1089	326	332	2	10			
1	B	210	Total	C	N	O	S	Se	0	23	0
			1763	1097	325	328	2	11			

There are 66 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	9	MSE	-	cloning artifact	UNP Q8UHP4
A	10	GLY	-	cloning artifact	UNP Q8UHP4
A	11	SER	-	cloning artifact	UNP Q8UHP4
A	12	SER	-	cloning artifact	UNP Q8UHP4
A	13	HIS	-	cloning artifact	UNP Q8UHP4
A	14	HIS	-	cloning artifact	UNP Q8UHP4
A	15	HIS	-	cloning artifact	UNP Q8UHP4
A	16	HIS	-	cloning artifact	UNP Q8UHP4
A	17	HIS	-	cloning artifact	UNP Q8UHP4
A	18	HIS	-	cloning artifact	UNP Q8UHP4
A	19	SER	-	cloning artifact	UNP Q8UHP4
A	20	SER	-	cloning artifact	UNP Q8UHP4
A	21	GLY	-	cloning artifact	UNP Q8UHP4
A	22	ARG	-	cloning artifact	UNP Q8UHP4
A	23	GLU	-	cloning artifact	UNP Q8UHP4
A	24	ASN	-	cloning artifact	UNP Q8UHP4
A	25	LEU	-	cloning artifact	UNP Q8UHP4
A	26	TYR	-	cloning artifact	UNP Q8UHP4
A	27	PHE	-	cloning artifact	UNP Q8UHP4
A	28	GLN	-	cloning artifact	UNP Q8UHP4
A	29	GLY	-	cloning artifact	UNP Q8UHP4
A	30	HIS	-	cloning artifact	UNP Q8UHP4
A	31	MSE	MET	modified residue	UNP Q8UHP4
A	34	MSE	MET	modified residue	UNP Q8UHP4
A	60	MSE	MET	modified residue	UNP Q8UHP4

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Chain	Residue	Modelled	Actual	Comment	Reference
A	80	MSE	MET	modified residue	UNP Q8UHP4
A	129	MSE	MET	modified residue	UNP Q8UHP4
A	152	MSE	MET	modified residue	UNP Q8UHP4
A	168	MSE	MET	modified residue	UNP Q8UHP4
A	213	MSE	MET	modified residue	UNP Q8UHP4
A	233	MSE	MET	modified residue	UNP Q8UHP4
A	267	GLY	-	cloning artifact	UNP Q8UHP4
A	268	SER	-	cloning artifact	UNP Q8UHP4
B	9	MSE	-	cloning artifact	UNP Q8UHP4
B	10	GLY	-	cloning artifact	UNP Q8UHP4
B	11	SER	-	cloning artifact	UNP Q8UHP4
B	12	SER	-	cloning artifact	UNP Q8UHP4
B	13	HIS	-	cloning artifact	UNP Q8UHP4
B	14	HIS	-	cloning artifact	UNP Q8UHP4
B	15	HIS	-	cloning artifact	UNP Q8UHP4
B	16	HIS	-	cloning artifact	UNP Q8UHP4
B	17	HIS	-	cloning artifact	UNP Q8UHP4
B	18	HIS	-	cloning artifact	UNP Q8UHP4
B	19	SER	-	cloning artifact	UNP Q8UHP4
B	20	SER	-	cloning artifact	UNP Q8UHP4
B	21	GLY	-	cloning artifact	UNP Q8UHP4
B	22	ARG	-	cloning artifact	UNP Q8UHP4
B	23	GLU	-	cloning artifact	UNP Q8UHP4
B	24	ASN	-	cloning artifact	UNP Q8UHP4
B	25	LEU	-	cloning artifact	UNP Q8UHP4
B	26	TYR	-	cloning artifact	UNP Q8UHP4
B	27	PHE	-	cloning artifact	UNP Q8UHP4
B	28	GLN	-	cloning artifact	UNP Q8UHP4
B	29	GLY	-	cloning artifact	UNP Q8UHP4
B	30	HIS	-	cloning artifact	UNP Q8UHP4
B	31	MSE	MET	modified residue	UNP Q8UHP4
B	34	MSE	MET	modified residue	UNP Q8UHP4
B	60	MSE	MET	modified residue	UNP Q8UHP4
B	80	MSE	MET	modified residue	UNP Q8UHP4
B	129	MSE	MET	modified residue	UNP Q8UHP4
B	152	MSE	MET	modified residue	UNP Q8UHP4
B	168	MSE	MET	modified residue	UNP Q8UHP4
B	213	MSE	MET	modified residue	UNP Q8UHP4
B	233	MSE	MET	modified residue	UNP Q8UHP4
B	267	GLY	-	cloning artifact	UNP Q8UHP4
B	268	SER	-	cloning artifact	UNP Q8UHP4

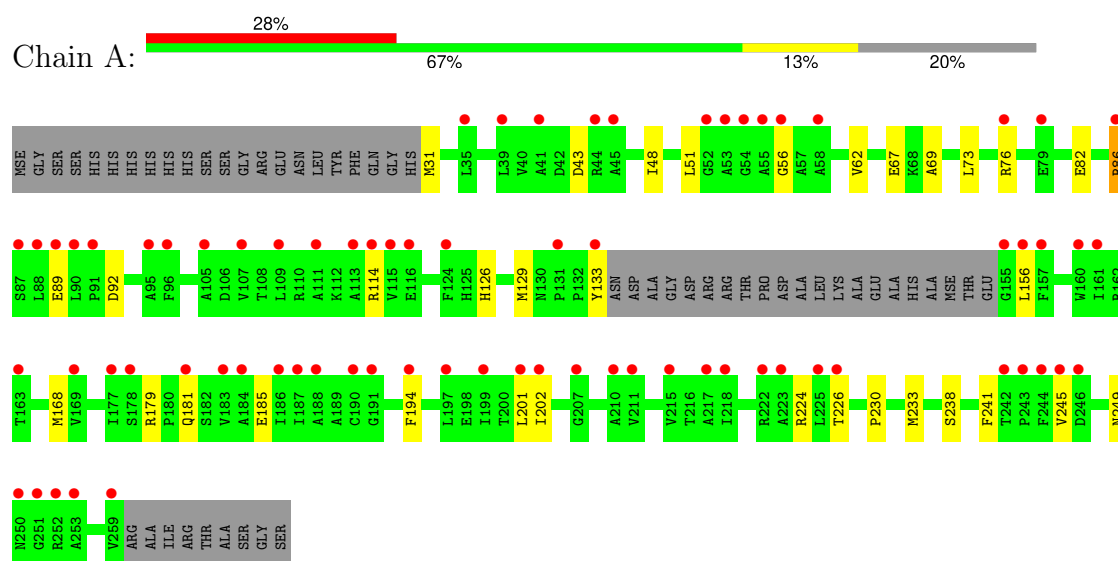
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	266	Total 266	O 266	0	0
2	B	245	Total 245	O 245	0	0

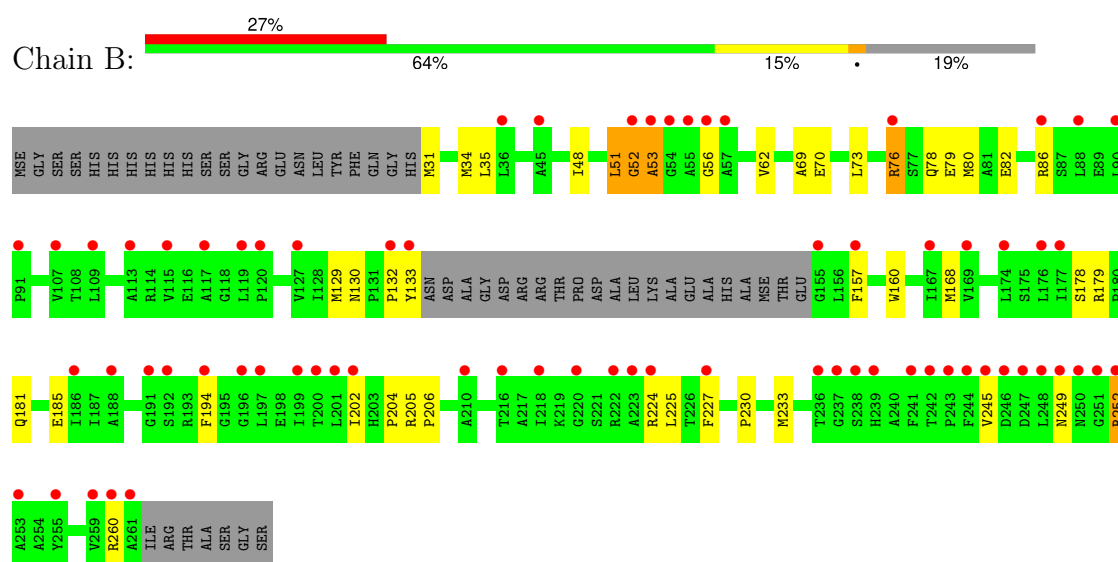
### 3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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. 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. 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: Hypothetical protein Atu0636



#### • Molecule 1: Hypothetical protein Atu0636



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	48.54Å 78.05Å 62.26Å 90.00° 94.55° 90.00°	Depositor
Resolution (Å)	31.03 – 1.70 31.03 – 1.70	Depositor EDS
% Data completeness (in resolution range)	95.5 (31.03-1.70) 95.5 (31.03-1.70)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.08	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.24 (at 1.70Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.178 , 0.214 0.189 , 0.222	Depositor DCC
$R_{free}$ test set	2490 reflections (5.12%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	26.5	Xtriage
Anisotropy	0.385	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 37.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	4033	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.66% 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](#)

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	A	0.65	0/1786	0.72	0/2395
1	B	0.64	0/1796	0.73	0/2408
All	All	0.64	0/3582	0.72	0/4803

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
1	B	0	2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	51	LEU	Peptide
1	B	52[B]	GLY	Peptide

### 5.2 Too-close contacts [i](#)

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	A	1759	0	1735	34	1
1	B	1763	0	1756	55	2

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	266	0	0	10	1
2	B	245	0	0	9	0
All	All	4033	0	3491	82	2

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 12.

All (82) 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:B:76:ARG:HG3	2:B:499:HOH:O	1.45	1.17
1:B:133[B]:TYR:CD2	1:B:179[B]:ARG:NH2	2.16	1.12
1:B:179[A]:ARG:NH1	1:B:181[A]:GLN:OE1	1.92	1.03
1:B:252:ARG:HG2	1:B:252:ARG:HH11	1.31	0.93
1:B:133[B]:TYR:OH	1:B:178[B]:SER:OG	1.91	0.89
1:B:224[B]:ARG:HH21	1:B:224[B]:ARG:HG2	1.39	0.87
1:B:34[B]:MSE:HG3	1:B:249:ASN:HA	1.57	0.87
1:A:185[B]:GLU:HG3	2:A:292:HOH:O	1.75	0.85
1:B:53[A]:ALA:HB3	1:B:130:ASN:O	1.76	0.85
1:A:156:LEU:HD23	2:A:526:HOH:O	1.76	0.84
1:A:224[A]:ARG:CD	1:B:181[A]:GLN:NE2	2.45	0.80
1:B:56:GLY:HA2	1:B:73[B]:LEU:HD11	1.65	0.78
1:A:133:TYR:HA	2:A:525:HOH:O	1.85	0.76
1:B:224[B]:ARG:HG2	1:B:224[B]:ARG:NH2	1.98	0.75
1:A:224[A]:ARG:HD2	1:B:181[A]:GLN:NE2	2.01	0.74
1:A:230:PRO:HG3	1:B:230:PRO:HG3	1.69	0.74
1:B:179[A]:ARG:CZ	1:B:181[A]:GLN:OE1	2.35	0.73
1:A:249[A]:ASN:HA	2:A:524:HOH:O	1.88	0.72
1:B:34[B]:MSE:CG	1:B:249:ASN:HA	2.19	0.71
1:B:252:ARG:HH11	1:B:252:ARG:CG	2.02	0.70
1:A:43[B]:ASP:OD1	1:A:43[B]:ASP:N	2.24	0.70
1:A:168[B]:MSE:SE	1:A:194:PHE:HE1	2.27	0.68
1:A:249[B]:ASN:HA	2:A:524:HOH:O	1.94	0.67
1:B:233[A]:MSE:HE2	1:B:245:VAL:HG21	1.75	0.67
1:A:226[A]:THR:HG22	2:A:521:HOH:O	1.95	0.67
1:B:48:ILE:HD13	1:B:62:VAL:HG11	1.77	0.66
1:A:181[B]:GLN:OE1	1:B:224[B]:ARG:NH2	2.30	0.65
1:A:86[A]:ARG:O	1:A:89[A]:GLU:HG2	1.96	0.64
1:A:202:ILE:HG23	1:A:233[A]:MSE:HE3	1.78	0.64
1:B:168[B]:MSE:SE	1:B:194:PHE:HE1	2.31	0.64
1:B:133[B]:TYR:CD2	1:B:179[B]:ARG:CZ	2.81	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:82:GLU:OE1	1:A:86[A]:ARG:NH2	2.31	0.63
1:B:133[B]:TYR:CE2	1:B:179[B]:ARG:NH2	2.66	0.63
1:B:73[B]:LEU:CD1	2:B:436:HOH:O	2.48	0.62
1:B:224[B]:ARG:HH21	1:B:224[B]:ARG:CG	2.12	0.61
1:B:252:ARG:HG2	1:B:252:ARG:NH1	2.11	0.61
1:B:51:LEU:HD12	1:B:129:MSE:HE2	1.83	0.60
1:B:252:ARG:HD2	2:B:501:HOH:O	2.01	0.60
1:B:35:LEU:HD13	1:B:202[B]:ILE:HD13	1.83	0.60
1:A:48:ILE:HD13	1:A:62:VAL:HG11	1.85	0.58
1:B:224[B]:ARG:HD3	1:B:225:LEU:O	2.03	0.58
1:A:224[A]:ARG:CD	1:B:181[A]:GLN:HE22	2.15	0.58
1:B:56:GLY:HA2	1:B:73[B]:LEU:CD1	2.34	0.58
1:A:31:MSE:HE1	1:A:241:PHE:CZ	2.39	0.57
1:A:233[A]:MSE:HE2	1:A:245:VAL:HG21	1.86	0.57
1:A:114[A]:ARG:NH2	2:A:518:HOH:O	2.33	0.57
1:B:53[A]:ALA:HB2	1:B:132:PRO:HD3	1.88	0.56
1:B:133[B]:TYR:HD2	1:B:179[B]:ARG:NH2	1.96	0.56
1:A:67[A]:GLU:HG2	2:A:389:HOH:O	2.05	0.55
1:B:224[B]:ARG:HG3	2:B:364:HOH:O	2.06	0.55
1:A:156:LEU:CD2	2:A:526:HOH:O	2.46	0.55
1:B:56:GLY:CA	1:B:73[B]:LEU:HD11	2.36	0.54
1:B:78:GLN:O	1:B:82[A]:GLU:HG3	2.09	0.52
1:A:51:LEU:HD12	1:A:129:MSE:HE2	1.92	0.51
1:B:34[B]:MSE:SE	1:B:249:ASN:HA	2.60	0.50
1:B:157:PHE:HB2	1:B:185[A]:GLU:HG2	1.95	0.49
1:A:168[B]:MSE:SE	1:A:194:PHE:CE1	3.12	0.49
1:B:168[B]:MSE:SE	1:B:194:PHE:CE1	3.14	0.48
1:A:89[B]:GLU:HG3	2:A:481:HOH:O	2.14	0.48
1:B:56:GLY:HA3	1:B:73[A]:LEU:HD13	1.96	0.48
1:B:48:ILE:HD12	1:B:69:ALA:HB1	1.96	0.48
1:B:76:ARG:O	1:B:76:ARG:HD2	2.14	0.47
1:A:224[A]:ARG:HD3	1:B:181[A]:GLN:NE2	2.26	0.47
1:B:179[A]:ARG:CZ	1:B:181[A]:GLN:CD	2.83	0.47
1:B:224[B]:ARG:NH2	2:B:495:HOH:O	2.47	0.46
1:A:56:GLY:HA3	1:A:73:LEU:HD13	1.96	0.46
1:B:252:ARG:CG	1:B:252:ARG:NH1	2.71	0.45
1:B:73[B]:LEU:HD11	2:B:436:HOH:O	2.13	0.45
1:A:48:ILE:HD12	1:A:69:ALA:HB1	1.98	0.44
1:A:201:LEU:HD11	1:B:227:PHE:CD2	2.52	0.44
1:A:133:TYR:CZ	1:A:179:ARG:HG3	2.52	0.43
1:B:52[A]:GLY:HA3	2:B:436:HOH:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:31:MSE:SE	1:A:233[A]:MSE:SE	3.36	0.43
1:A:31:MSE:HB2	1:A:31:MSE:HE3	1.69	0.43
1:B:205[A]:ARG:HB2	1:B:206:PRO:CD	2.48	0.43
1:B:70[A]:GLU:HG2	2:B:412:HOH:O	2.19	0.43
1:B:260:ARG:HG3	2:B:500:HOH:O	2.20	0.42
1:A:48:ILE:HG12	1:A:126:HIS:HB2	2.03	0.41
1:B:133[B]:TYR:CG	1:B:179[B]:ARG:CZ	3.04	0.41
1:A:233[A]:MSE:HE2	1:A:245:VAL:CG2	2.51	0.41
1:B:53[B]:ALA:H	1:B:160:TRP:HH2	1.69	0.40
1:B:204:PRO:HA	1:B:233[B]:MSE:SE	2.72	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:76:ARG:NH1	1:B:79[A]:GLU:OE2[1_556]	2.03	0.17
1:B:79[A]:GLU:OE1	2:A:471:HOH:O[1_554]	2.03	0.17

## 5.3 Torsion angles ⓘ

### 5.3.1 Protein backbone ⓘ

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	A	227/260 (87%)	222 (98%)	5 (2%)	0	100	100
1	B	228/260 (88%)	222 (97%)	4 (2%)	2 (1%)	14	4
All	All	455/520 (88%)	444 (98%)	9 (2%)	2 (0%)	44	17

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	53[A]	ALA
1	B	53[B]	ALA

### 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	A	179/188 (95%)	175 (98%)	4 (2%)	47	30
1	B	179/188 (95%)	174 (97%)	5 (3%)	38	21
All	All	358/376 (95%)	349 (98%)	9 (2%)	42	25

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	86[A]	ARG
1	A	86[B]	ARG
1	A	92	ASP
1	A	238	SER
1	B	31	MSE
1	B	76	ARG
1	B	80	MSE
1	B	86	ARG
1	B	252	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	250	ASN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

There are no ligands in this entry.

## 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

### 6.1 Protein, DNA and RNA chains

**Warning:** The R factor obtained from EDS is 0.2939, which does not match the depositor's R factor of 0.17767. Please interpret the results in this section carefully.

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	A	200/260 (76%)	1.86	74 (37%)  	22, 45, 57, 69	21 (10%)
1	B	202/260 (77%)	1.81	70 (34%)  	22, 47, 61, 76	20 (9%)
All	All	402/520 (77%)	1.84	144 (35%)  	22, 46, 60, 76	41 (10%)

All (144) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	133[A]	TYR	7.8
1	A	55[A]	ALA	7.2
1	A	259	VAL	6.2
1	B	191[A]	GLY	5.4
1	A	54[A]	GLY	5.1
1	A	133	TYR	5.1
1	B	53[A]	ALA	4.7
1	B	243	PRO	4.5
1	B	241	PHE	4.4
1	B	236	THR	4.1
1	B	52[A]	GLY	4.0
1	A	88[A]	LEU	4.0
1	A	243	PRO	4.0
1	B	55	ALA	4.0
1	B	252	ARG	3.9
1	B	259	VAL	3.9
1	A	76	ARG	3.9
1	B	261	ALA	3.9
1	A	155	GLY	3.8
1	B	250	ASN	3.8
1	B	239	HIS	3.8
1	B	260	ARG	3.6

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Mol	Chain	Res	Type	RSRZ
1	B	237	GLY	3.6
1	A	44[A]	ARG	3.6
1	A	91	PRO	3.6
1	A	53	ALA	3.5
1	A	157	PHE	3.4
1	B	54	GLY	3.4
1	A	183	VAL	3.3
1	A	52	GLY	3.3
1	A	90	LEU	3.3
1	A	244	PHE	3.2
1	A	181[A]	GLN	3.1
1	A	111	ALA	3.1
1	B	248	LEU	3.1
1	A	114[A]	ARG	3.0
1	B	247	ASP	3.0
1	A	186	ILE	3.0
1	A	191	GLY	2.8
1	A	96	PHE	2.8
1	A	199	ILE	2.8
1	A	156	LEU	2.7
1	A	194	PHE	2.7
1	A	202	ILE	2.7
1	A	107	VAL	2.7
1	A	184	ALA	2.6
1	A	217	ALA	2.6
1	B	238	SER	2.6
1	B	109	LEU	2.6
1	B	249	ASN	2.6
1	A	210	ALA	2.6
1	B	117	ALA	2.6
1	B	169	VAL	2.6
1	B	245	VAL	2.6
1	B	76	ARG	2.6
1	A	131	PRO	2.6
1	A	178[A]	SER	2.6
1	B	90	LEU	2.5
1	A	161	ILE	2.5
1	A	160	TRP	2.5
1	B	224[A]	ARG	2.5
1	A	218	ILE	2.5
1	B	57	ALA	2.5
1	B	157	PHE	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	194	PHE	2.5
1	A	89[A]	GLU	2.5
1	A	252	ARG	2.5
1	A	197	LEU	2.4
1	A	201	LEU	2.4
1	B	176	LEU	2.4
1	B	186	ILE	2.4
1	B	218	ILE	2.4
1	A	250	ASN	2.4
1	B	107	VAL	2.4
1	B	255	TYR	2.4
1	A	95	ALA	2.4
1	A	223	ALA	2.4
1	A	109	LEU	2.4
1	B	56	GLY	2.4
1	B	91	PRO	2.4
1	A	115	VAL	2.4
1	A	222[A]	ARG	2.4
1	A	177	ILE	2.4
1	B	188	ALA	2.4
1	B	246	ASP	2.4
1	B	216	THR	2.3
1	A	245	VAL	2.3
1	B	88	LEU	2.3
1	B	174	LEU	2.3
1	B	222[A]	ARG	2.3
1	B	227	PHE	2.3
1	A	188	ALA	2.3
1	A	226[A]	THR	2.3
1	A	242	THR	2.3
1	A	116[A]	GLU	2.3
1	A	58	ALA	2.3
1	B	192[A]	SER	2.3
1	B	167	ILE	2.2
1	B	199	ILE	2.2
1	A	45	ALA	2.2
1	A	113	ALA	2.2
1	A	79[A]	GLU	2.2
1	A	163	THR	2.2
1	B	36	LEU	2.2
1	A	86[A]	ARG	2.2
1	A	190	CYS	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	187	ILE	2.2
1	B	202[A]	ILE	2.2
1	B	86	ARG	2.2
1	A	105	ALA	2.2
1	B	210	ALA	2.2
1	B	223	ALA	2.2
1	A	87	SER	2.2
1	A	169	VAL	2.2
1	B	115	VAL	2.2
1	A	225	LEU	2.2
1	A	56	GLY	2.2
1	B	251	GLY	2.2
1	B	253	ALA	2.1
1	B	244	PHE	2.1
1	A	251	GLY	2.1
1	B	220	GLY	2.1
1	A	39	LEU	2.1
1	B	119	LEU	2.1
1	B	201	LEU	2.1
1	A	41	ALA	2.1
1	B	45	ALA	2.1
1	B	113	ALA	2.1
1	A	215	VAL	2.1
1	B	200	THR	2.1
1	A	246	ASP	2.1
1	A	35	LEU	2.1
1	A	211	VAL	2.1
1	B	127	VAL	2.1
1	A	253	ALA	2.0
1	B	242	THR	2.0
1	B	177	ILE	2.0
1	A	207	GLY	2.0
1	B	155	GLY	2.0
1	A	124	PHE	2.0
1	B	197	LEU	2.0
1	B	196	GLY	2.0
1	B	120	PRO	2.0
1	B	132	PRO	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

### 6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

### 6.4 Ligands [i](#)

There are no ligands in this entry.

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

There are no such residues in this entry.