



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 01:41 AM UTC

PDB ID : 6Z90 / pdb_00006z90
Title : Crystal structure of MINDY1 mutant-P138A
Authors : Abdul Rehman, S.A.; Kulathu, Y.
Deposited on : 2020-06-03
Resolution : 3.59 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

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>.

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

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

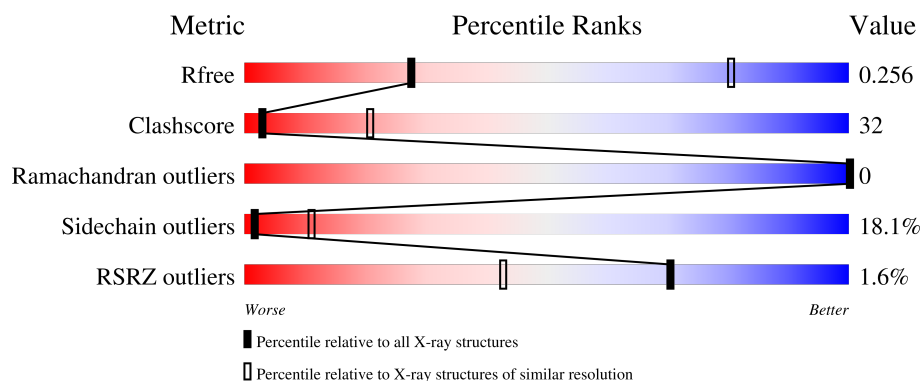
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.59 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}	180053	1747 (3.70-3.50)
Clashscore	190562	1827 (3.70-3.50)
Ramachandran outliers	187476	1773 (3.70-3.50)
Sidechain outliers	187428	1772 (3.70-3.50)
RSRZ outliers	180081	1745 (3.70-3.50)

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 ≥ 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	A	289	47% 35% 7% 11%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 1946 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 Ubiquitin carboxyl-terminal hydrolase MINDY-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	256	1946	1240	315	379	12	0	0	0

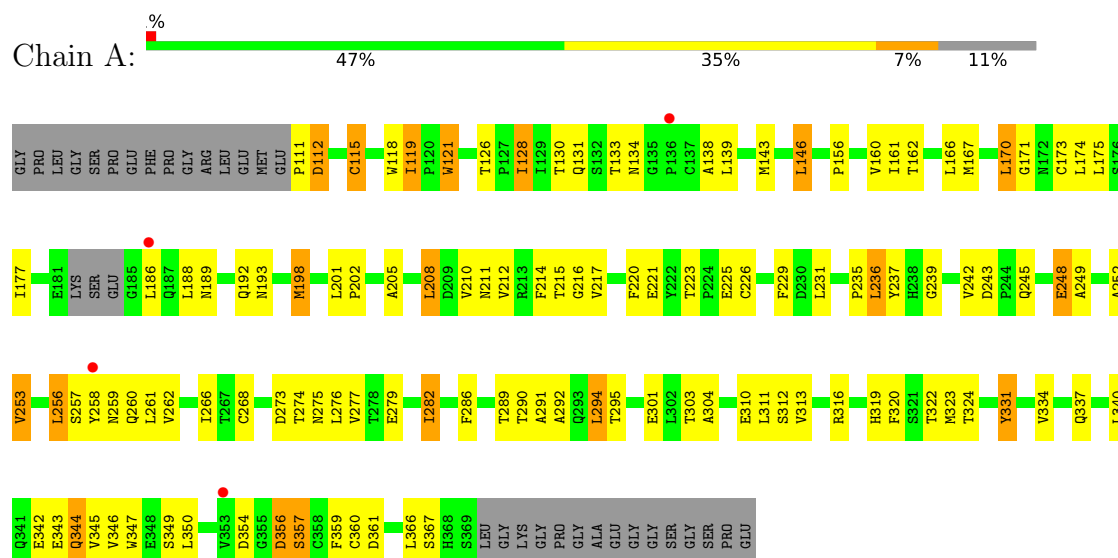
There are 15 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	96	GLY	-	expression tag	UNP Q8N5J2
A	97	PRO	-	expression tag	UNP Q8N5J2
A	98	LEU	-	expression tag	UNP Q8N5J2
A	99	GLY	-	expression tag	UNP Q8N5J2
A	100	SER	-	expression tag	UNP Q8N5J2
A	101	PRO	-	expression tag	UNP Q8N5J2
A	102	GLU	-	expression tag	UNP Q8N5J2
A	103	PHE	-	expression tag	UNP Q8N5J2
A	104	PRO	-	expression tag	UNP Q8N5J2
A	105	GLY	-	expression tag	UNP Q8N5J2
A	106	ARG	-	expression tag	UNP Q8N5J2
A	107	LEU	-	expression tag	UNP Q8N5J2
A	108	GLU	-	expression tag	UNP Q8N5J2
A	109	MET	-	expression tag	UNP Q8N5J2
A	138	ALA	PRO	engineered mutation	UNP Q8N5J2

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 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: Ubiquitin carboxyl-terminal hydrolase MINDY-1



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 2 2	Depositor
Cell constants a, b, c, α , β , γ	98.65Å 98.65Å 166.07Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.00 – 3.59 48.00 – 3.59	Depositor EDS
% Data completeness (in resolution range)	99.8 (48.00-3.59) 99.8 (48.00-3.59)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.08 (at 3.57Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.212 , 0.258 (Not available) , 0.256	Depositor DCC
R_{free} test set	525 reflections (5.18%)	wwPDB-VP
Wilson B-factor (Å ²)	149.3	Xtriage
Anisotropy	0.676	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 199.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	1946	wwPDB-VP
Average B, all atoms (Å ²)	191.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.00% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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	1.10	0/1993	1.57	7/2731 (0.3%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	331	TYR	CB-CA-C	5.69	119.57	109.72
1	A	361	ASP	CA-CB-CG	5.54	118.14	112.60
1	A	279	GLU	CA-C-N	5.27	126.73	120.14
1	A	279	GLU	C-N-CA	5.27	126.73	120.14
1	A	128	ILE	CA-C-O	-5.09	116.91	121.09
1	A	189	ASN	CA-C-N	5.07	127.38	120.54
1	A	189	ASN	C-N-CA	5.07	127.38	120.54

There are no chirality outliers.

There are no planarity outliers.

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	1946	0	1798	120	0
All	All	1946	0	1798	120	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 32.

All (120) 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:A:177:ILE:HD12	1:A:231:LEU:CD2	1.55	1.37
1:A:256:LEU:HD11	1:A:260:GLN:C	1.69	1.18
1:A:256:LEU:HD21	1:A:260:GLN:HB3	1.32	1.06
1:A:177:ILE:CD1	1:A:231:LEU:HD23	1.87	1.05
1:A:177:ILE:CD1	1:A:231:LEU:CD2	2.36	1.03
1:A:256:LEU:HD11	1:A:260:GLN:O	1.63	0.96
1:A:177:ILE:HD12	1:A:231:LEU:HD23	1.42	0.95
1:A:256:LEU:CD1	1:A:261:LEU:N	2.33	0.91
1:A:256:LEU:HD12	1:A:261:LEU:HA	1.53	0.91
1:A:177:ILE:HD12	1:A:231:LEU:HD21	1.51	0.90
1:A:256:LEU:CD1	1:A:260:GLN:C	2.43	0.90
1:A:256:LEU:CD2	1:A:260:GLN:HB3	2.04	0.86
1:A:170:LEU:HD12	1:A:170:LEU:O	1.76	0.86
1:A:256:LEU:CD1	1:A:261:LEU:CA	2.54	0.85
1:A:231:LEU:HD23	1:A:231:LEU:O	1.80	0.82
1:A:256:LEU:CD1	1:A:261:LEU:HA	2.13	0.78
1:A:256:LEU:HD12	1:A:261:LEU:CA	2.16	0.75
1:A:208:LEU:CD1	1:A:225:GLU:OE1	2.34	0.75
1:A:131:GLN:OE1	1:A:319:HIS:HE1	1.71	0.72
1:A:177:ILE:HD13	1:A:231:LEU:HD23	1.70	0.72
1:A:177:ILE:HD12	1:A:231:LEU:HD22	1.66	0.71
1:A:242:VAL:CG2	1:A:249:ALA:HB1	2.20	0.70
1:A:256:LEU:HD13	1:A:261:LEU:N	2.05	0.69
1:A:290:THR:HG22	1:A:290:THR:O	1.93	0.69
1:A:167:MET:HE2	1:A:205:ALA:HB2	1.76	0.68
1:A:346:VAL:HG12	1:A:347:TRP:CD1	2.29	0.67
1:A:211:ASN:ND2	1:A:259:ASN:HB2	2.10	0.67
1:A:256:LEU:CD2	1:A:260:GLN:CB	2.72	0.67
1:A:344:GLN:HG3	1:A:367:SER:HB2	1.77	0.66
1:A:208:LEU:HD11	1:A:225:GLU:OE1	1.98	0.64
1:A:170:LEU:HD12	1:A:170:LEU:C	2.21	0.64
1:A:337:GLN:HG3	1:A:340:LEU:HD22	1.81	0.61
1:A:356:ASP:N	1:A:356:ASP:OD1	2.34	0.60
1:A:256:LEU:HD22	1:A:257:SER:H	1.66	0.60
1:A:146:LEU:HD23	1:A:146:LEU:N	2.17	0.59
1:A:175:LEU:HG	1:A:198:MET:HE3	1.82	0.59
1:A:344:GLN:O	1:A:360:CYS:O	2.21	0.59
1:A:208:LEU:HD13	1:A:225:GLU:CG	2.33	0.58
1:A:131:GLN:OE1	1:A:319:HIS:CE1	2.54	0.58
1:A:217:VAL:O	1:A:237:TYR:HA	2.04	0.57
1:A:331:TYR:CD1	1:A:349:SER:HA	2.39	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:349:SER:OG	1:A:357:SER:CB	2.53	0.57
1:A:256:LEU:HD11	1:A:261:LEU:N	2.05	0.57
1:A:334:VAL:O	1:A:334:VAL:HG12	2.04	0.56
1:A:342:GLU:HA	1:A:342:GLU:OE1	2.05	0.56
1:A:290:THR:HG21	1:A:295:THR:HG23	1.87	0.55
1:A:256:LEU:HD21	1:A:260:GLN:CB	2.19	0.55
1:A:292:ALA:O	1:A:294:LEU:N	2.31	0.54
1:A:273:ASP:OD2	1:A:275:ASN:HB3	2.07	0.54
1:A:223:THR:HB	1:A:225:GLU:OE1	2.07	0.54
1:A:256:LEU:HD12	1:A:261:LEU:HD12	1.90	0.54
1:A:208:LEU:HD13	1:A:225:GLU:CD	2.33	0.53
1:A:290:THR:HG21	1:A:295:THR:CG2	2.39	0.53
1:A:331:TYR:HD1	1:A:349:SER:HA	1.74	0.53
1:A:118:TRP:CZ2	1:A:156:PRO:HA	2.43	0.53
1:A:139:LEU:HD21	1:A:166:LEU:HB3	1.91	0.52
1:A:211:ASN:CG	1:A:259:ASN:HB2	2.34	0.52
1:A:301:GLU:O	1:A:304:ALA:HB3	2.09	0.52
1:A:343:GLU:O	1:A:346:VAL:CG2	2.58	0.52
1:A:177:ILE:HD13	1:A:231:LEU:O	2.09	0.52
1:A:242:VAL:HG23	1:A:249:ALA:HB1	1.92	0.51
1:A:344:GLN:HG3	1:A:367:SER:CB	2.41	0.50
1:A:211:ASN:HD21	1:A:259:ASN:HB2	1.74	0.50
1:A:346:VAL:HG12	1:A:347:TRP:NE1	2.27	0.50
1:A:171:GLY:O	1:A:175:LEU:HG	2.12	0.49
1:A:258:TYR:CG	1:A:258:TYR:O	2.65	0.49
1:A:262:VAL:O	1:A:266:ILE:HG12	2.13	0.49
1:A:143:MET:HE3	1:A:170:LEU:HB2	1.94	0.49
1:A:208:LEU:CD1	1:A:225:GLU:CD	2.87	0.48
1:A:220:PHE:CZ	1:A:236:LEU:HD23	2.49	0.48
1:A:345:VAL:HG13	1:A:359:PHE:CD2	2.49	0.48
1:A:242:VAL:CG2	1:A:249:ALA:CB	2.91	0.47
1:A:214:PHE:HZ	1:A:258:TYR:HB2	1.79	0.47
1:A:323:MET:HG2	1:A:324:THR:N	2.29	0.47
1:A:115:CYS:O	1:A:130:THR:OG1	2.29	0.47
1:A:167:MET:CE	1:A:205:ALA:HB2	2.43	0.47
1:A:214:PHE:CZ	1:A:258:TYR:HB2	2.50	0.47
1:A:282:ILE:HG23	1:A:282:ILE:O	2.15	0.47
1:A:143:MET:HA	1:A:143:MET:HE2	1.96	0.46
1:A:231:LEU:HD23	1:A:231:LEU:C	2.40	0.46
1:A:345:VAL:HG13	1:A:359:PHE:HD2	1.80	0.46
1:A:239:GLY:HA3	1:A:294:LEU:HD12	1.95	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:256:LEU:HD22	1:A:260:GLN:CB	2.46	0.46
1:A:121:TRP:CD1	1:A:121:TRP:C	2.90	0.46
1:A:256:LEU:HD22	1:A:260:GLN:HB2	1.99	0.45
1:A:138:ALA:HB2	1:A:320:PHE:CD2	2.52	0.45
1:A:208:LEU:O	1:A:208:LEU:HG	2.16	0.45
1:A:249:ALA:O	1:A:253:VAL:HB	2.16	0.45
1:A:292:ALA:HB1	1:A:316:ARG:HH22	1.82	0.44
1:A:248:GLU:O	1:A:252:ALA:N	2.44	0.44
1:A:121:TRP:CD1	1:A:121:TRP:O	2.70	0.44
1:A:138:ALA:HB2	1:A:320:PHE:CG	2.52	0.44
1:A:289:THR:C	1:A:291:ALA:H	2.25	0.44
1:A:235:PRO:HD2	1:A:312:SER:HA	2.00	0.44
1:A:257:SER:O	1:A:261:LEU:N	2.50	0.43
1:A:175:LEU:CD2	1:A:198:MET:HE3	2.48	0.43
1:A:170:LEU:CD1	1:A:174:LEU:CD1	2.96	0.43
1:A:119:ILE:HD11	1:A:121:TRP:HB2	2.00	0.43
1:A:170:LEU:HD11	1:A:174:LEU:CD1	2.47	0.43
1:A:337:GLN:O	1:A:340:LEU:HB2	2.18	0.43
1:A:243:ASP:C	1:A:245:GLN:N	2.77	0.43
1:A:349:SER:OG	1:A:357:SER:HB3	2.18	0.42
1:A:167:MET:HE3	1:A:202:PRO:HA	2.01	0.42
1:A:290:THR:O	1:A:290:THR:CG2	2.63	0.42
1:A:111:PRO:HG3	1:A:133:THR:HG21	2.00	0.42
1:A:112:ASP:OD1	1:A:112:ASP:N	2.53	0.42
1:A:170:LEU:HD11	1:A:174:LEU:HD11	2.01	0.42
1:A:211:ASN:HB3	1:A:221:GLU:HB2	2.01	0.42
1:A:248:GLU:H	1:A:248:GLU:HG2	1.41	0.42
1:A:146:LEU:N	1:A:146:LEU:CD2	2.81	0.42
1:A:268:CYS:O	1:A:277:VAL:HG22	2.19	0.42
1:A:286:PHE:O	1:A:286:PHE:CG	2.73	0.42
1:A:131:GLN:NE2	1:A:134:ASN:HA	2.34	0.41
1:A:171:GLY:HA2	1:A:198:MET:HE1	2.02	0.41
1:A:229:PHE:O	1:A:231:LEU:N	2.53	0.41
1:A:128:ILE:HD12	1:A:128:ILE:HA	1.94	0.41
1:A:215:THR:OG1	1:A:216:GLY:N	2.53	0.41
1:A:201:LEU:N	1:A:202:PRO:CD	2.84	0.40
1:A:350:LEU:HD12	1:A:350:LEU:HA	1.86	0.40
1:A:343:GLU:O	1:A:346:VAL:HG23	2.22	0.40

There are no symmetry-related clashes.

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	217/289 (75%)	182 (84%)	35 (16%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains

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	210/252 (83%)	172 (82%)	38 (18%)	2	11

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	112	ASP
1	A	115	CYS
1	A	119	ILE
1	A	121	TRP
1	A	126	THR
1	A	146	LEU
1	A	160	VAL
1	A	161	ILE
1	A	162	THR
1	A	170	LEU
1	A	173	CYS
1	A	186	LEU
1	A	188	LEU

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Mol	Chain	Res	Type
1	A	192	GLN
1	A	193	ASN
1	A	198	MET
1	A	208	LEU
1	A	210	VAL
1	A	212	VAL
1	A	226	CYS
1	A	236	LEU
1	A	248	GLU
1	A	253	VAL
1	A	256	LEU
1	A	274	THR
1	A	276	LEU
1	A	282	ILE
1	A	294	LEU
1	A	303	THR
1	A	310	GLU
1	A	311	LEU
1	A	313	VAL
1	A	322	THR
1	A	344	GLN
1	A	354	ASP
1	A	356	ASP
1	A	357	SER
1	A	366	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	189	ASN
1	A	192	GLN
1	A	193	ASN
1	A	317	ASN
1	A	318	ASN
1	A	319	HIS

5.3.3 RNA ⓘ

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 [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, 95th 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 ²)	Q<0.9
1	A	256/289 (88%)	-0.07	4 (1%) 70 43	132, 189, 246, 293	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	258	TYR	5.2
1	A	136	PRO	3.4
1	A	353	VAL	2.5
1	A	186	LEU	2.4

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides 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.