



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 03:42 PM UTC

PDB ID : 1UF4 / pdb_00001uf4
Title : Crystal structure of C171A/V236A Mutant of N-carbamyl-D-amino acid amidohydrolase
Authors : Hashimoto, H.; Aoki, M.; Shimizu, T.; Nakai, T.; Morikawa, H.; Ikenaka, Y.; Takahashi, S.; Sato, M.
Deposited on : 2003-05-23
Resolution : 2.15 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

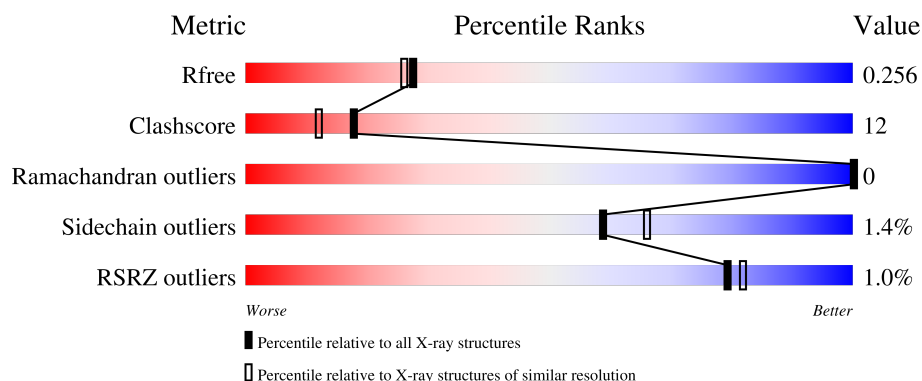
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.15 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	2057 (2.16-2.16)
Clashscore	190562	2159 (2.16-2.16)
Ramachandran outliers	187476	2134 (2.16-2.16)
Sidechain outliers	187428	2133 (2.16-2.16)
RSRZ outliers	180081	2059 (2.16-2.16)

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	303	% 71% 27% .
1	B	303	% 73% 25% .

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5232 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 N-carbamyl-D-amino acid amidohydrolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	303	Total	C	N	O	S	0	0	0
			2404	1533	425	433	13			
1	B	303	Total	C	N	O	S	0	0	0
			2404	1533	425	433	13			

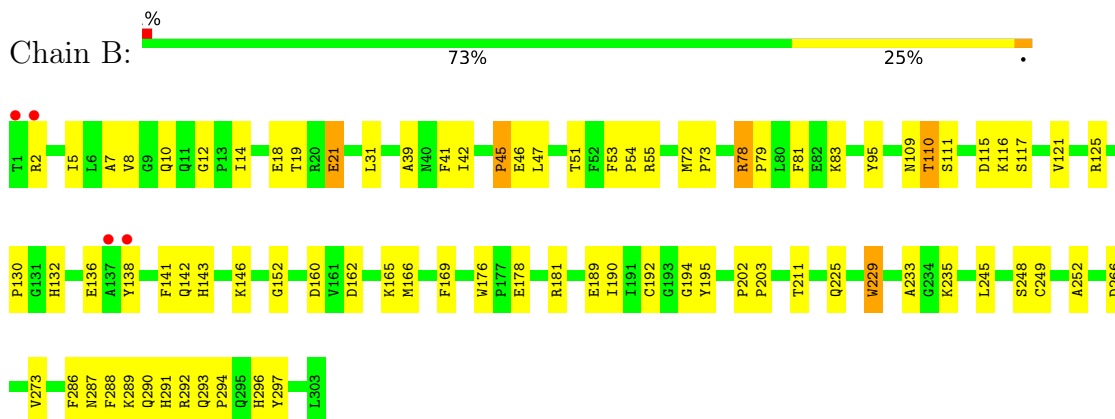
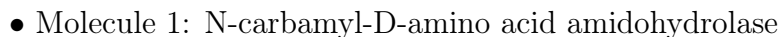
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	171	ALA	CYS	engineered mutation	UNP P60327
A	236	ALA	VAL	engineered mutation	UNP P60327
B	171	ALA	CYS	engineered mutation	UNP P60327
B	236	ALA	VAL	engineered mutation	UNP P60327

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	214	Total	O	0	0
			214	214		
2	B	210	Total	O	0	0
			210	210		

- Molecule 1: N-carbamyl-D-amino acid amidohydrolase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	67.26Å 136.85Å 67.76Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	68.42 – 2.15 68.42 – 2.15	Depositor EDS
% Data completeness (in resolution range)	98.2 (68.42-2.15) 98.3 (68.42-2.15)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.57 (at 2.14Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.201 , 0.256 0.202 , 0.256	Depositor DCC
R_{free} test set	3413 reflections (9.98%)	wwPDB-VP
Wilson B-factor (Å ²)	26.9	Xtriage
Anisotropy	0.561	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 55.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.000 for l,-k,h	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5232	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.21% 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

5.1 Standard geometry

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	0.37	0/2467	0.92	11/3341 (0.3%)
1	B	0.39	0/2467	0.92	11/3341 (0.3%)
All	All	0.38	0/4934	0.92	22/6682 (0.3%)

There are no bond length outliers.

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	235	LYS	N-CA-C	-7.15	95.33	107.99
1	B	229	TRP	N-CA-C	-6.65	101.76	110.53
1	B	249	CYS	N-CA-C	6.45	118.56	108.96
1	A	110	THR	N-CA-C	6.43	120.11	110.14
1	A	109	ASN	N-CA-C	-6.32	97.40	108.20
1	A	249	CYS	N-CA-C	6.19	117.63	108.60
1	A	248	SER	N-CA-C	-6.03	102.00	110.50
1	B	109	ASN	N-CA-C	-5.97	96.64	107.75
1	B	51	THR	N-CA-C	-5.96	103.08	110.41
1	B	45	PRO	N-CA-C	5.88	119.45	111.22
1	B	110	THR	N-CA-C	5.79	119.11	110.14
1	B	192	CYS	N-CA-C	5.74	117.69	109.14
1	B	12	GLY	N-CA-C	-5.50	101.11	112.34
1	B	235	LYS	N-CA-C	-5.36	98.50	107.99
1	A	51	THR	N-CA-C	-5.35	103.46	110.53
1	A	12	GLY	N-CA-C	-5.25	101.63	112.34
1	A	229	TRP	N-CA-C	-5.24	102.72	110.48
1	A	195	TYR	N-CA-C	5.21	116.76	108.79
1	B	195	TYR	N-CA-C	5.10	116.28	108.42
1	A	46	GLU	N-CA-C	-5.05	103.39	110.50
1	A	189	GLU	N-CA-C	-5.03	107.28	113.41
1	B	248	SER	N-CA-C	-5.01	102.86	110.28

There are no chirality outliers.

There are no planarity outliers.

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	A	2404	0	2360	69	0
1	B	2404	0	2360	54	0
2	A	214	0	0	5	0
2	B	210	0	0	2	0
All	All	5232	0	4720	117	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 12.

All (117) 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:78:ARG:HH21	1:B:78:ARG:HB2	1.47	0.79
1:B:78:ARG:HB2	1:B:78:ARG:NH2	2.03	0.73
1:B:73:PRO:HB3	1:B:81:PHE:CE1	2.24	0.72
1:A:9:GLY:N	1:A:42:ILE:HD11	2.04	0.72
1:A:21:GLU:HG3	1:A:76:VAL:HG13	1.72	0.71
1:A:1:THR:HG22	1:A:3:GLN:H	1.56	0.71
1:A:21:GLU:O	1:A:25:VAL:HG23	1.96	0.64
1:A:73:PRO:HB3	1:A:81:PHE:CE1	2.33	0.64
1:B:31:LEU:C	1:B:31:LEU:HD23	2.24	0.63
1:B:211:THR:HG22	1:B:245:LEU:HD22	1.80	0.62
1:A:42:ILE:HD13	1:A:43:VAL:N	2.15	0.61
1:B:225:GLN:HA	1:B:286:PHE:CE1	2.36	0.61
1:A:42:ILE:HD13	1:A:42:ILE:C	2.26	0.60
1:A:166:MET:HG2	1:A:190:ILE:HB	1.82	0.59
1:B:202:PRO:HB2	1:B:203:PRO:HD3	1.84	0.59
1:A:200:HIS:HD2	1:A:208:ASP:OD2	1.86	0.58
1:A:1:THR:HG22	1:A:2:ARG:H	1.68	0.58
1:A:60:ASP:O	1:A:63:GLU:HB2	2.03	0.58
1:B:132:HIS:ND1	1:B:136:GLU:HG3	2.20	0.57
1:A:1:THR:HG22	1:A:2:ARG:N	2.19	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:110:THR:HG22	1:A:111:SER:N	2.19	0.56
1:A:10:GLN:O	1:A:266:ASP:HA	2.06	0.55
1:A:78:ARG:HD2	2:A:391:HOH:O	2.07	0.55
1:B:110:THR:HG22	1:B:111:SER:N	2.23	0.54
1:A:9:GLY:CA	1:A:42:ILE:HD11	2.37	0.54
1:B:289:LYS:HZ2	1:B:290:GLN:HG2	1.73	0.53
1:B:19:THR:OG1	1:B:21:GLU:HG2	2.08	0.53
1:B:292:ARG:C	1:B:294:PRO:HD3	2.33	0.53
1:B:116:LYS:HG2	1:B:162:ASP:CG	2.34	0.53
1:B:138:TYR:OH	1:B:203:PRO:HG3	2.08	0.53
1:B:14:ILE:HA	1:B:18:GLU:OE1	2.08	0.53
1:A:2:ARG:HB2	2:A:496:HOH:O	2.08	0.52
1:B:2:ARG:HH11	1:B:2:ARG:HG2	1.75	0.52
1:B:294:PRO:HA	1:B:297:TYR:CD2	2.45	0.52
1:B:55:ARG:HA	1:B:142:GLN:O	2.09	0.51
1:B:45:PRO:HA	1:B:194:GLY:HA3	1.92	0.51
1:A:146:LYS:NZ	1:B:291:HIS:HD2	2.08	0.51
1:A:240:GLU:O	1:A:241:ASN:HB2	2.09	0.51
1:A:9:GLY:HA3	1:A:42:ILE:HD11	1.93	0.51
1:A:55:ARG:HA	1:A:142:GLN:O	2.11	0.51
1:A:136:GLU:HB2	1:A:139:ARG:CD	2.41	0.51
1:A:291:HIS:HD2	1:B:146:LYS:NZ	2.07	0.51
1:B:5:ILE:HB	1:B:39:ALA:HA	1.93	0.50
1:B:229:TRP:CE3	1:B:252:ALA:HA	2.47	0.50
1:A:110:THR:CG2	1:A:111:SER:N	2.75	0.50
1:A:40:ASN:C	1:A:89:ILE:HG13	2.37	0.49
1:A:130:PRO:HG2	1:A:143:HIS:NE2	2.26	0.49
1:A:41:PHE:CZ	1:A:166:MET:HE1	2.46	0.49
1:A:294:PRO:HA	1:A:297:TYR:CD2	2.47	0.49
1:A:136:GLU:HB2	1:A:139:ARG:HD2	1.95	0.48
1:A:303:LEU:HD22	1:A:303:LEU:N	2.28	0.48
1:B:292:ARG:HD3	2:B:440:HOH:O	2.14	0.48
1:A:125:ARG:O	1:A:152:GLY:HA3	2.13	0.48
1:A:113:LEU:HG	1:A:168:MET:SD	2.54	0.48
1:A:220:GLN:HG2	1:A:255:GLY:O	2.14	0.48
1:A:176:TRP:CZ3	1:B:292:ARG:HG2	2.49	0.48
1:B:53:PHE:N	1:B:54:PRO:CD	2.77	0.47
1:B:194:GLY:HA2	1:B:233:ALA:O	2.14	0.47
1:B:287:ASN:OD1	1:B:289:LYS:HG2	2.14	0.47
1:A:116:LYS:HG2	1:A:162:ASP:CG	2.39	0.47
1:A:297:TYR:CD2	1:B:178:GLU:HG2	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:101:GLU:OE1	1:A:106:ARG:HD2	2.15	0.46
1:B:160:ASP:OD2	1:B:165:LYS:NZ	2.48	0.46
1:B:293:GLN:HB3	1:B:296:HIS:HD2	1.81	0.46
1:A:89:ILE:HG23	1:A:89:ILE:O	2.15	0.46
1:B:125:ARG:O	1:B:152:GLY:HA3	2.16	0.45
1:A:42:ILE:HD12	1:A:44:PHE:CE2	2.51	0.45
1:B:169:PHE:C	1:B:169:PHE:CD1	2.93	0.45
1:A:200:HIS:CD2	1:A:208:ASP:OD2	2.69	0.45
1:A:100:VAL:HG12	2:A:454:HOH:O	2.17	0.45
1:A:133:LYS:HE3	1:A:146:LYS:O	2.17	0.45
1:B:10:GLN:O	1:B:266:ASP:HA	2.17	0.45
1:B:121:VAL:HG11	1:B:160:ASP:O	2.17	0.45
1:B:41:PHE:CE2	1:B:166:MET:HE1	2.51	0.45
1:A:225:GLN:HA	1:A:286:PHE:CE1	2.52	0.44
1:B:72:MET:HE3	1:B:95:TYR:CD2	2.52	0.44
1:A:46:GLU:O	1:A:47:LEU:C	2.61	0.44
1:B:141:PHE:HA	2:B:444:HOH:O	2.16	0.44
1:A:201:ASN:ND2	1:A:203:PRO:HG2	2.32	0.44
1:A:292:ARG:HD3	2:A:339:HOH:O	2.17	0.44
1:A:297:TYR:HB3	1:A:300:ILE:HD12	2.00	0.43
1:A:53:PHE:N	1:A:54:PRO:CD	2.81	0.43
1:A:20:ARG:O	1:A:24:VAL:HG23	2.17	0.43
1:B:46:GLU:O	1:B:47:LEU:C	2.61	0.43
1:A:201:ASN:HD21	1:A:203:PRO:HG2	1.83	0.43
1:A:4:MET:HB2	1:A:40:ASN:HD21	1.82	0.43
1:A:57:HIS:CD2	1:A:135:TYR:HB2	2.54	0.43
1:B:166:MET:HG2	1:B:190:ILE:HB	1.99	0.43
1:A:21:GLU:HG3	1:A:76:VAL:CG1	2.47	0.42
1:A:45:PRO:HA	1:A:194:GLY:HA3	2.01	0.42
1:B:115:ASP:C	1:B:117:SER:H	2.27	0.42
1:A:229:TRP:CE3	1:A:252:ALA:HA	2.54	0.42
1:A:169:PHE:C	1:A:169:PHE:CD1	2.97	0.42
1:B:7:ALA:O	1:B:42:ILE:HA	2.19	0.42
1:A:132:HIS:H	1:A:132:HIS:CD2	2.38	0.42
1:B:132:HIS:CE1	1:B:136:GLU:HG3	2.54	0.42
1:B:2:ARG:NH2	1:B:189:GLU:OE2	2.50	0.42
1:A:14:ILE:HG13	1:A:235:LYS:HD2	2.02	0.41
1:B:8:VAL:HG13	1:B:233:ALA:HB2	2.03	0.41
1:B:83:LYS:HD2	1:B:83:LYS:HA	1.87	0.41
1:A:7:ALA:O	1:A:42:ILE:HA	2.20	0.41
1:A:176:TRP:HZ3	1:B:292:ARG:HG2	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:288:PHE:CG	1:B:292:ARG:HD2	2.55	0.41
1:A:133:LYS:HE3	1:A:146:LYS:C	2.45	0.41
1:B:78:ARG:HB3	1:B:79:PRO:HD3	2.01	0.41
1:B:72:MET:HA	1:B:73:PRO:HA	1.82	0.41
1:A:176:TRP:HA	1:A:177:PRO:HD2	1.97	0.41
1:A:41:PHE:CE2	1:A:166:MET:HE1	2.56	0.41
1:A:78:ARG:HB3	1:A:79:PRO:HD3	2.03	0.41
1:A:78:ARG:N	1:A:79:PRO:CD	2.84	0.41
1:A:292:ARG:HG2	1:B:176:TRP:CZ3	2.56	0.41
1:A:75:PRO:HD2	2:A:415:HOH:O	2.21	0.40
1:B:130:PRO:HG2	1:B:143:HIS:NE2	2.37	0.40
1:A:78:ARG:NH1	1:A:78:ARG:HG3	2.36	0.40
1:A:229:TRP:CE2	1:A:273:VAL:HG13	2.56	0.40
1:B:110:THR:CG2	1:B:111:SER:N	2.84	0.40
1:B:229:TRP:CE2	1:B:273:VAL:HG13	2.56	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	301/303 (99%)	288 (96%)	13 (4%)	0	100	100
1	B	301/303 (99%)	290 (96%)	11 (4%)	0	100	100
All	All	602/606 (99%)	578 (96%)	24 (4%)	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	249/249 (100%)	245 (98%)	4 (2%)	55	61
1	B	249/249 (100%)	246 (99%)	3 (1%)	63	70
All	All	498/498 (100%)	491 (99%)	7 (1%)	59	66

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	42	ILE
1	A	63	GLU
1	A	133	LYS
1	A	181	ARG
1	B	21	GLU
1	B	78	ARG
1	B	181	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	57	HIS
1	A	200	HIS
1	A	201	ASN
1	A	290	GLN
1	A	291	HIS
1	A	296	HIS
1	B	57	HIS
1	B	290	GLN
1	B	291	HIS
1	B	296	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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	303/303 (100%)	0.08	2 (0%) 84 85	16, 26, 40, 53	0
1	B	303/303 (100%)	0.04	4 (1%) 75 78	15, 25, 41, 57	0
All	All	606/606 (100%)	0.06	6 (0%) 79 82	15, 26, 41, 57	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1	THR	6.3
1	B	1	THR	4.2
1	B	138	TYR	3.2
1	A	163	ALA	2.1
1	B	2	ARG	2.1
1	B	137	ALA	2.1

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.