Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 04:36 PM GMT

PDB ID	:	4FJS
Title	:	Crystal structure of ureidoglycolate dehydrogenase enzyme in apo form
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Deposited on	:	2012-06-12
Resolution	:	2.13 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
http://wwpdb.org/validation/2016/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

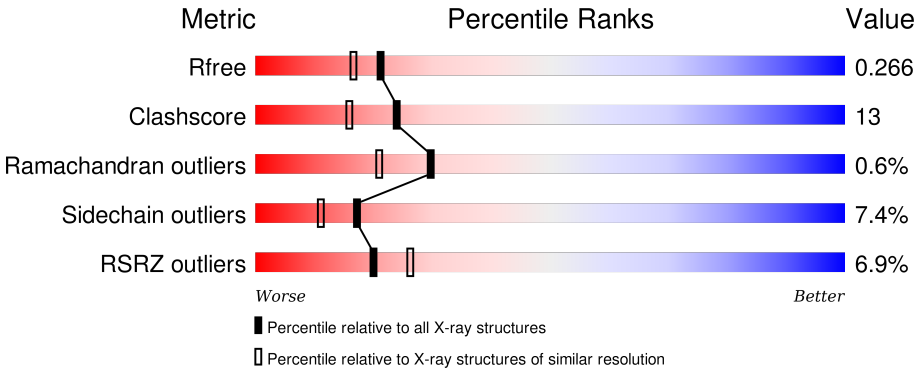
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.13 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}	91344	1693 (2.16-2.12)
Clashscore	102246	1824 (2.16-2.12)
Ramachandran outliers	100387	1798 (2.16-2.12)
Sidechain outliers	100360	1798 (2.16-2.12)
RSRZ outliers	91569	1699 (2.16-2.12)

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 >=3, 2, 1 and 0 types of geometric quality criteria. 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 <=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	351	7% 76% 18%
1	B	351	6% 72% 16% 10%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5459 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 Ureidoglycolate dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	337	Total	C	N	O	S	0	0	0
			2570	1621	450	483	16			
1	B	315	Total	C	N	O	S	0	0	0
			2389	1502	426	446	15			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	EXPRESSION TAG	UNP B1XGB5
A	0	HIS	-	EXPRESSION TAG	UNP B1XGB5
B	-1	GLY	-	EXPRESSION TAG	UNP B1XGB5
B	0	HIS	-	EXPRESSION TAG	UNP B1XGB5

- Molecule 2 is water.

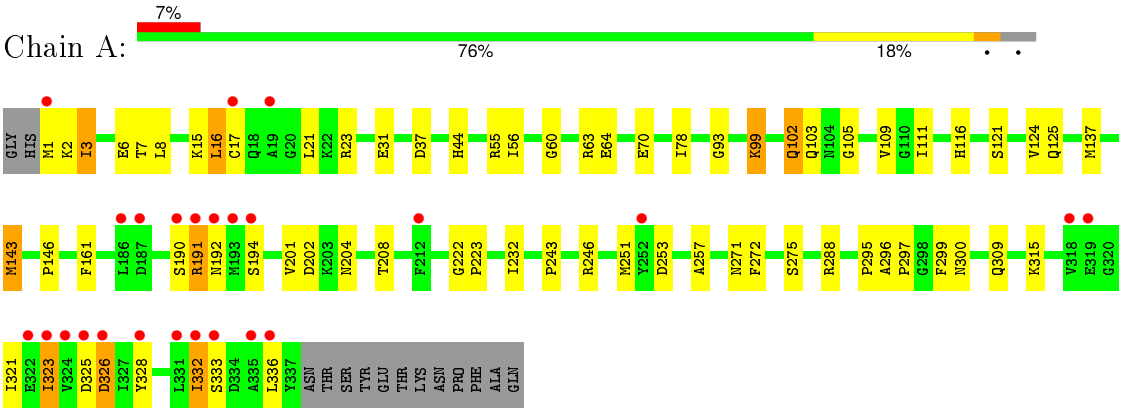
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	232	Total	O	0	0
			232	232		
2	B	268	Total	O	0	0
			268	268		

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 errors 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.

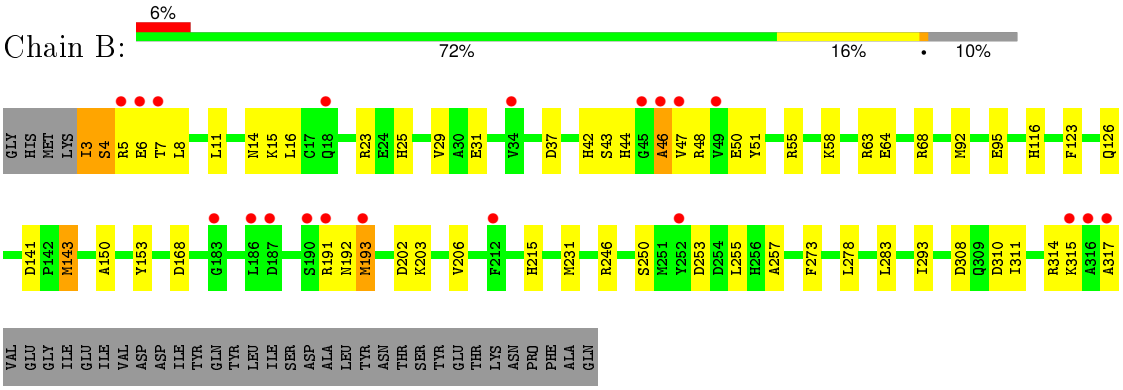
- Molecule 1: Ureidoglycolate dehydrogenase

Chain A:



- Molecule 1: Ureidoglycolate dehydrogenase

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 42 21 2	Depositor
Cell constants a, b, c, α, β, γ	162.76Å 162.76Å 61.45Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.18 – 2.13 49.04 – 2.12	Depositor EDS
% Data completeness (in resolution range)	99.0 (49.18-2.13) 91.6 (49.04-2.12)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	5.11 (at 2.12Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE: 1.7.2_869)	Depositor
R, R_{free}	0.222 , 0.240 0.245 , 0.266	Depositor DCC
R_{free} test set	1883 reflections (4.35%)	DCC
Wilson B-factor (Å^2)	21.8	Xtriage
Anisotropy	0.010	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 41.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.47$, $< L^2 > = 0.30$	Xtriage
Outliers	0 of 46309 reflections	Xtriage
F_o,F_c correlation	0.91	EDS
Total number of atoms	5459	wwPDB-VP
Average B, all atoms (Å^2)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.80% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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	0.44	0/2625	0.57	0/3558
1	B	0.47	1/2441 (0.0%)	0.56	0/3308
All	All	0.45	1/5066 (0.0%)	0.57	0/6866

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	193	MET	CG-SD	6.23	1.97	1.81

There are no bond angle outliers.

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	2570	0	2528	69	2
1	B	2389	0	2347	63	1
2	A	232	0	0	44	1
2	B	268	0	0	43	4
All	All	5459	0	4875	129	8

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 13.

All (129) 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:42:HIS:HA	2:B:639:HOH:O	1.27	1.24
1:B:231:MET:SD	2:B:565:HOH:O	1.99	1.20
1:A:3:ILE:HG13	2:A:615:HOH:O	1.40	1.17
1:A:271:ASN:ND2	2:A:441:HOH:O	1.81	1.12
1:A:288:ARG:NH1	2:A:596:HOH:O	1.88	1.05
1:B:15:LYS:NZ	2:B:587:HOH:O	1.86	0.99
1:A:299:PHE:HA	2:A:602:HOH:O	1.59	0.99
1:B:317:ALA:O	2:B:653:HOH:O	1.80	0.99
1:A:31:GLU:OE1	2:A:606:HOH:O	1.81	0.98
1:A:295:PRO:C	2:A:604:HOH:O	2.00	0.98
1:B:15:LYS:CE	2:B:587:HOH:O	2.09	0.98
1:A:6:GLU:O	2:A:560:HOH:O	1.80	0.98
1:B:68:ARG:NH1	2:B:603:HOH:O	1.94	0.97
1:A:299:PHE:CA	2:A:602:HOH:O	2.13	0.96
1:B:253:ASP:OD1	2:B:576:HOH:O	1.87	0.92
1:A:202:ASP:OD2	2:A:611:HOH:O	1.88	0.92
1:B:6:GLU:OE1	2:B:579:HOH:O	1.85	0.91
1:B:37:ASP:OD1	2:B:639:HOH:O	1.87	0.91
1:B:68:ARG:NH2	2:B:603:HOH:O	2.00	0.90
1:B:168:ASP:OD1	2:B:656:HOH:O	1.89	0.89
1:A:103:GLN:NE2	2:A:580:HOH:O	2.05	0.88
1:B:141:ASP:OD1	2:B:647:HOH:O	1.91	0.87
1:A:300:ASN:N	2:A:602:HOH:O	1.94	0.86
1:B:37:ASP:OD1	2:B:568:HOH:O	1.94	0.86
2:A:604:HOH:O	1:B:150:ALA:HB2	1.76	0.85
1:B:95:GLU:OE1	2:B:507:HOH:O	1.94	0.83
1:A:1:MET:HG3	1:A:2:LYS:N	1.94	0.81
1:A:295:PRO:O	2:A:604:HOH:O	1.97	0.80
1:A:64:GLU:OE1	2:A:572:HOH:O	1.99	0.80
2:A:604:HOH:O	1:B:150:ALA:CB	2.30	0.79
1:B:68:ARG:CZ	2:B:603:HOH:O	2.23	0.79
1:B:310:ASP:OD2	2:B:424:HOH:O	2.00	0.78
1:A:70:GLU:OE2	2:A:545:HOH:O	2.01	0.78
1:A:3:ILE:CG1	2:A:615:HOH:O	2.13	0.77
1:B:50:GLU:O	2:B:462:HOH:O	2.01	0.77
1:A:103:GLN:OE1	2:A:552:HOH:O	2.03	0.76
1:B:37:ASP:OD2	2:B:568:HOH:O	2.03	0.76
1:B:44:HIS:HD2	1:B:116:HIS:HE1	1.35	0.75
1:B:37:ASP:CG	2:B:568:HOH:O	2.26	0.73
1:A:309:GLN:OE1	2:A:550:HOH:O	2.07	0.73
1:B:314:ARG:O	2:B:578:HOH:O	2.06	0.72
1:B:5:ARG:HA	2:B:426:HOH:O	1.88	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:42:HIS:ND1	2:B:639:HOH:O	1.94	0.72
1:B:3:ILE:N	2:B:628:HOH:O	2.23	0.71
1:A:99:LYS:CE	2:A:577:HOH:O	2.35	0.70
1:A:326:ASP:N	1:A:326:ASP:OD2	2.22	0.70
1:B:4:SER:O	2:B:426:HOH:O	2.09	0.70
1:B:44:HIS:CD2	1:B:116:HIS:HE1	2.09	0.69
1:A:204:ASN:HB2	2:A:573:HOH:O	1.93	0.69
1:B:250:SER:HB2	2:B:660:HOH:O	1.94	0.68
1:B:50:GLU:C	2:B:462:HOH:O	2.31	0.67
1:A:253:ASP:OD1	2:A:490:HOH:O	2.13	0.66
1:A:3:ILE:CB	2:A:615:HOH:O	2.41	0.66
1:B:3:ILE:N	2:B:556:HOH:O	2.29	0.65
1:B:15:LYS:CD	2:B:587:HOH:O	2.39	0.65
1:A:99:LYS:NZ	2:A:577:HOH:O	1.91	0.65
1:B:23:ARG:NH1	2:B:573:HOH:O	1.95	0.64
1:A:17:CYS:O	2:A:526:HOH:O	2.15	0.64
1:A:102:GLN:NE2	2:A:556:HOH:O	1.95	0.62
1:A:99:LYS:HE3	2:A:577:HOH:O	1.98	0.62
1:A:78:ILE:HD11	1:B:273:PHE:HA	1.81	0.62
1:A:60:GLY:C	2:A:553:HOH:O	2.37	0.62
1:A:55:ARG:HD3	2:A:553:HOH:O	2.00	0.61
1:A:44:HIS:HD2	1:A:116:HIS:HE1	1.49	0.60
1:A:288:ARG:NE	2:A:551:HOH:O	2.13	0.58
1:A:336:LEU:HD11	2:A:513:HOH:O	2.05	0.56
1:A:6:GLU:HA	2:A:605:HOH:O	2.06	0.55
1:A:3:ILE:HB	2:A:615:HOH:O	2.06	0.55
1:A:6:GLU:CB	2:A:605:HOH:O	2.56	0.53
1:A:1:MET:HG3	1:A:2:LYS:H	1.74	0.52
1:A:44:HIS:HD2	1:A:116:HIS:CE1	2.28	0.52
1:B:15:LYS:HD3	2:B:587:HOH:O	2.03	0.52
1:A:246:ARG:NH2	2:A:425:HOH:O	2.16	0.52
1:B:14:ASN:HA	2:B:573:HOH:O	2.10	0.51
1:A:297:PRO:HA	2:A:563:HOH:O	2.09	0.51
1:B:14:ASN:HB2	2:B:609:HOH:O	2.10	0.51
1:A:37:ASP:HB3	1:A:321:ILE:HG21	1.93	0.51
1:A:6:GLU:HB3	2:A:605:HOH:O	2.11	0.50
1:A:93:GLY:HA3	1:A:111:ILE:HD11	1.93	0.50
1:B:153:TYR:CD1	1:B:246:ARG:HG3	2.47	0.50
1:B:51:TYR:HA	2:B:462:HOH:O	2.10	0.50
1:A:146:PRO:HG3	1:B:293:ILE:HD12	1.92	0.49
1:B:46:ALA:N	2:B:568:HOH:O	2.45	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:8:LEU:HB2	2:B:426:HOH:O	2.11	0.49
1:A:55:ARG:NH1	1:A:257:ALA:O	2.45	0.48
1:A:296:ALA:N	2:A:604:HOH:O	2.33	0.47
1:B:3:ILE:HD11	1:B:8:LEU:HD13	1.96	0.47
1:A:1:MET:HB3	1:A:323:ILE:O	2.15	0.47
1:A:102:GLN:NE2	2:A:580:HOH:O	2.46	0.46
1:A:191:ARG:HH11	1:A:191:ARG:HB3	1.80	0.46
1:A:6:GLU:CA	2:A:605:HOH:O	2.63	0.46
1:A:161:PHE:CD1	1:B:231:MET:HG3	2.50	0.46
1:B:50:GLU:N	2:B:641:HOH:O	2.47	0.46
1:B:8:LEU:HD12	1:B:8:LEU:HA	1.67	0.46
1:B:308:ASP:N	1:B:308:ASP:OD1	2.44	0.46
1:A:328:TYR:CE1	2:A:554:HOH:O	2.68	0.45
1:A:137:MET:HB2	1:A:232:ILE:HG23	1.98	0.45
1:B:3:ILE:CA	2:B:556:HOH:O	2.66	0.44
1:B:68:ARG:NH2	2:B:517:HOH:O	2.49	0.44
1:A:332:ILE:HG12	1:A:332:ILE:H	1.33	0.44
1:A:3:ILE:O	1:A:321:ILE:N	2.46	0.44
1:A:2:LYS:HE3	1:A:2:LYS:HB2	1.64	0.44
1:B:47:VAL:HB	2:B:608:HOH:O	2.18	0.43
1:B:4:SER:O	1:B:7:THR:HB	2.18	0.43
1:A:6:GLU:CD	1:A:6:GLU:H	2.22	0.43
1:A:222:GLY:HA3	1:A:223:PRO:HD3	1.90	0.43
1:B:44:HIS:CD2	1:B:116:HIS:CE1	2.99	0.43
1:B:203:LYS:HD3	1:B:215:HIS:HB2	2.01	0.43
1:A:16:LEU:HD12	1:A:16:LEU:HA	1.88	0.42
1:B:42:HIS:CB	2:B:639:HOH:O	2.56	0.42
1:B:202:ASP:OD1	1:B:206:VAL:HB	2.19	0.42
1:B:311:ILE:HA	1:B:314:ARG:HG3	2.02	0.42
1:B:278:LEU:HD23	1:B:278:LEU:HA	1.84	0.42
1:A:125:GLN:NE2	2:A:567:HOH:O	2.50	0.42
1:A:7:THR:OG1	2:A:615:HOH:O	1.97	0.41
1:A:121:SER:HA	1:A:124:VAL:HG22	2.01	0.41
1:B:123:PHE:O	1:B:126:GLN:HB2	2.20	0.41
1:B:253:ASP:CB	2:B:576:HOH:O	2.68	0.41
1:A:1:MET:CG	1:A:2:LYS:N	2.77	0.41
1:A:143:MET:HB2	1:A:143:MET:HE2	1.88	0.41
1:B:55:ARG:NH1	1:B:257:ALA:O	2.53	0.41
1:A:44:HIS:CD2	1:A:116:HIS:CE1	3.08	0.41
1:A:299:PHE:CB	2:A:602:HOH:O	2.61	0.41
1:A:21:LEU:HG	1:A:56:ILE:HD13	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:14:ASN:CB	2:B:609:HOH:O	2.68	0.41
1:B:25:HIS:O	1:B:29:VAL:HG23	2.21	0.41
1:A:105:GLY:HA2	1:A:272:PHE:CD1	2.57	0.40
1:B:58:LYS:NZ	2:B:410:HOH:O	2.31	0.40
1:A:3:ILE:HD11	1:A:8:LEU:HB2	2.03	0.40

All (8) 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)
2:B:594:HOH:O	2:B:610:HOH:O[7_555]	1.57	0.63
2:B:492:HOH:O	2:B:517:HOH:O[8_554]	1.82	0.38
1:A:1:MET:CE	1:A:204:ASN:ND2[3_544]	1.90	0.30
2:A:552:HOH:O	2:A:600:HOH:O[8_554]	1.96	0.24
2:B:520:HOH:O	2:B:638:HOH:O[7_555]	2.00	0.20
1:B:7:THR:OG1	1:B:7:THR:OG1[7_555]	2.06	0.14
2:B:508:HOH:O	2:B:508:HOH:O[8_555]	2.14	0.06
1:A:271:ASN:O	1:A:275:SER:OG[8_554]	2.18	0.02

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	335/351 (95%)	319 (95%)	15 (4%)	1 (0%)	46	41
1	B	313/351 (89%)	295 (94%)	15 (5%)	3 (1%)	19	10
All	All	648/702 (92%)	614 (95%)	30 (5%)	4 (1%)	30	21

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	333	SER

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Mol	Chain	Res	Type
1	B	192	ASN
1	B	46	ALA
1	B	143	MET

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	266/278 (96%)	244 (92%)	22 (8%)	14 8
1	B	246/278 (88%)	230 (94%)	16 (6%)	21 15
All	All	512/556 (92%)	474 (93%)	38 (7%)	17 11

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

Mol	Chain	Res	Type
1	A	3	ILE
1	A	15	LYS
1	A	16	LEU
1	A	23	ARG
1	A	63	ARG
1	A	99	LYS
1	A	102	GLN
1	A	109	VAL
1	A	143	MET
1	A	190	SER
1	A	191	ARG
1	A	192	ASN
1	A	194	SER
1	A	201	VAL
1	A	208	THR
1	A	243	PRO
1	A	251	MET
1	A	315	LYS
1	A	323	ILE
1	A	325	ASP

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Mol	Chain	Res	Type
1	A	326	ASP
1	A	332	ILE
1	B	3	ILE
1	B	4	SER
1	B	11	LEU
1	B	16	LEU
1	B	31	GLU
1	B	43	SER
1	B	48	ARG
1	B	63	ARG
1	B	64	GLU
1	B	92	MET
1	B	143	MET
1	B	191	ARG
1	B	193	MET
1	B	255	LEU
1	B	283	LEU
1	B	315	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	10	GLN
1	A	14	ASN
1	A	44	HIS
1	A	103	GLN
1	A	271	ASN
1	A	300	ASN
1	B	44	HIS
1	B	116	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)

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

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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^{th} 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	337/351 (96%)	0.41	25 (7%)	17	23	12, 30, 65, 79	0
1	B	315/351 (89%)	0.52	20 (6%)	23	31	11, 30, 61, 92	0
All	All	652/702 (92%)	0.47	45 (6%)	20	26	11, 30, 64, 92	0

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	193	MET	6.4
1	B	191	ARG	5.1
1	B	46	ALA	5.0
1	A	1	MET	4.8
1	A	326	ASP	4.1
1	A	193	MET	3.9
1	B	7	THR	3.9
1	A	318	VAL	3.9
1	B	252	TYR	3.8
1	A	19	ALA	3.6
1	B	190	SER	3.4
1	B	212	PHE	3.4
1	A	191	ARG	3.4
1	A	332	ILE	3.2
1	B	6	GLU	3.2
1	A	335	ALA	3.2
1	A	324	VAL	3.1
1	A	190	SER	3.1
1	B	186	LEU	3.1
1	B	47	VAL	3.0
1	A	331	LEU	3.0
1	A	17	CYS	2.9
1	B	45	GLY	2.8
1	A	212	PHE	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	336	LEU	2.8
1	B	183	GLY	2.7
1	A	322	GLU	2.7
1	A	323	ILE	2.6
1	A	192	ASN	2.6
1	B	5	ARG	2.4
1	A	328	TYR	2.4
1	A	187	ASP	2.4
1	B	34	VAL	2.4
1	A	319	GLU	2.4
1	A	325	ASP	2.4
1	A	194	SER	2.3
1	A	252	TYR	2.3
1	B	315	LYS	2.3
1	A	333	SER	2.3
1	B	316	ALA	2.2
1	B	49	VAL	2.2
1	B	18	GLN	2.1
1	A	186	LEU	2.1
1	B	317	ALA	2.1
1	B	187	ASP	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 carbohydrates 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.