Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 10:38 AM GMT

PDB ID	:	3MIZ
Title	:	Crystal structure of a putative transcriptional regulator protein, LacI family from Rhizobium etli
Authors	:	Palani, K.; Burley, S.K.; Swaminathan, S.; New York SGX Research Center for Structural Genomics (NYSGXRC)
Deposited on	:	2010-04-12
Resolution	:	1.91 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:

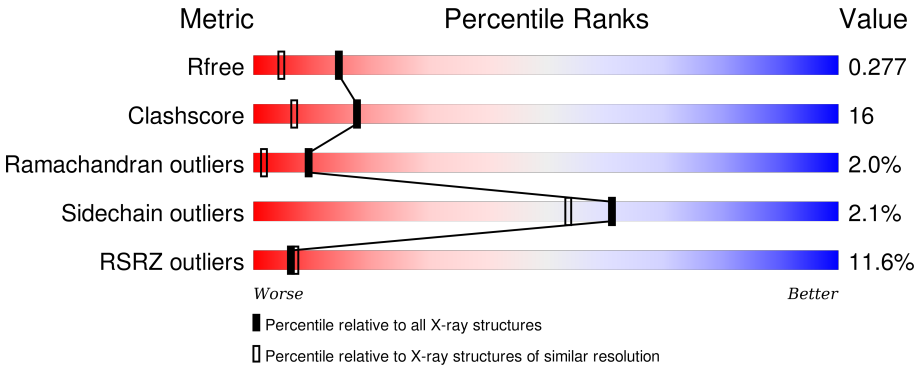
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	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 1.91 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	5832 (1.94-1.90)
Clashscore	102246	6540 (1.94-1.90)
Ramachandran outliers	100387	6464 (1.94-1.90)
Sidechain outliers	100360	6465 (1.94-1.90)
RSRZ outliers	91569	5846 (1.94-1.90)

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	301	7%, 67%, 24%, 7%
1	B	301	14%, 59%, 31%, 7%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4578 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 Putative transcriptional regulator protein, LacI family.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	281	Total	C	N	O	S	Se	0	0	0
			2199	1384	382	422	2	9			
1	B	281	Total	C	N	O	S	Se	0	0	0
			2192	1380	379	422	2	9			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	47	MSE	-	EXPRESSION TAG	UNP Q2K0Z9
A	48	SER	-	EXPRESSION TAG	UNP Q2K0Z9
A	49	LEU	-	EXPRESSION TAG	UNP Q2K0Z9
A	340	GLU	-	EXPRESSION TAG	UNP Q2K0Z9
A	341	GLY	-	EXPRESSION TAG	UNP Q2K0Z9
A	342	HIS	-	EXPRESSION TAG	UNP Q2K0Z9
A	343	HIS	-	EXPRESSION TAG	UNP Q2K0Z9
A	344	HIS	-	EXPRESSION TAG	UNP Q2K0Z9
A	345	HIS	-	EXPRESSION TAG	UNP Q2K0Z9
A	346	HIS	-	EXPRESSION TAG	UNP Q2K0Z9
A	347	HIS	-	EXPRESSION TAG	UNP Q2K0Z9
B	47	MSE	-	EXPRESSION TAG	UNP Q2K0Z9
B	48	SER	-	EXPRESSION TAG	UNP Q2K0Z9
B	49	LEU	-	EXPRESSION TAG	UNP Q2K0Z9
B	340	GLU	-	EXPRESSION TAG	UNP Q2K0Z9
B	341	GLY	-	EXPRESSION TAG	UNP Q2K0Z9
B	342	HIS	-	EXPRESSION TAG	UNP Q2K0Z9
B	343	HIS	-	EXPRESSION TAG	UNP Q2K0Z9
B	344	HIS	-	EXPRESSION TAG	UNP Q2K0Z9
B	345	HIS	-	EXPRESSION TAG	UNP Q2K0Z9
B	346	HIS	-	EXPRESSION TAG	UNP Q2K0Z9
B	347	HIS	-	EXPRESSION TAG	UNP Q2K0Z9

- Molecule 2 is water.

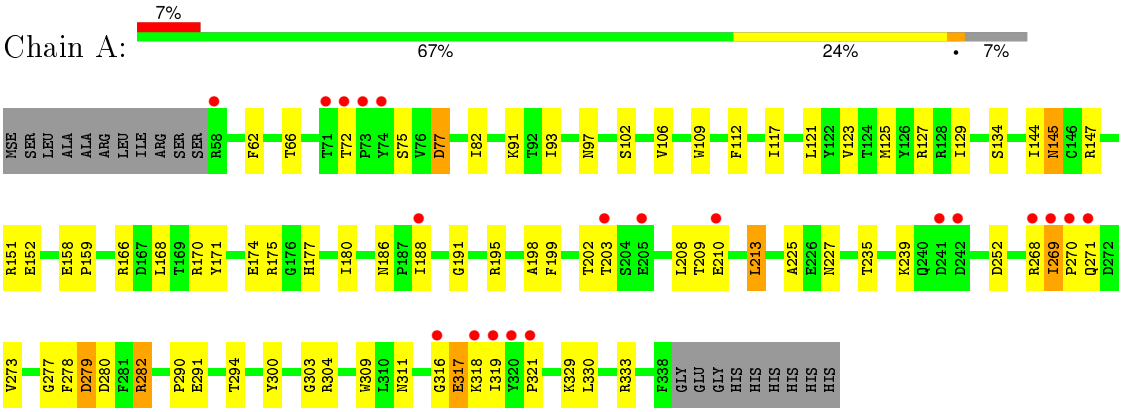
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	123	Total O 123 123	0	0
2	B	64	Total O 64 64	0	0

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.

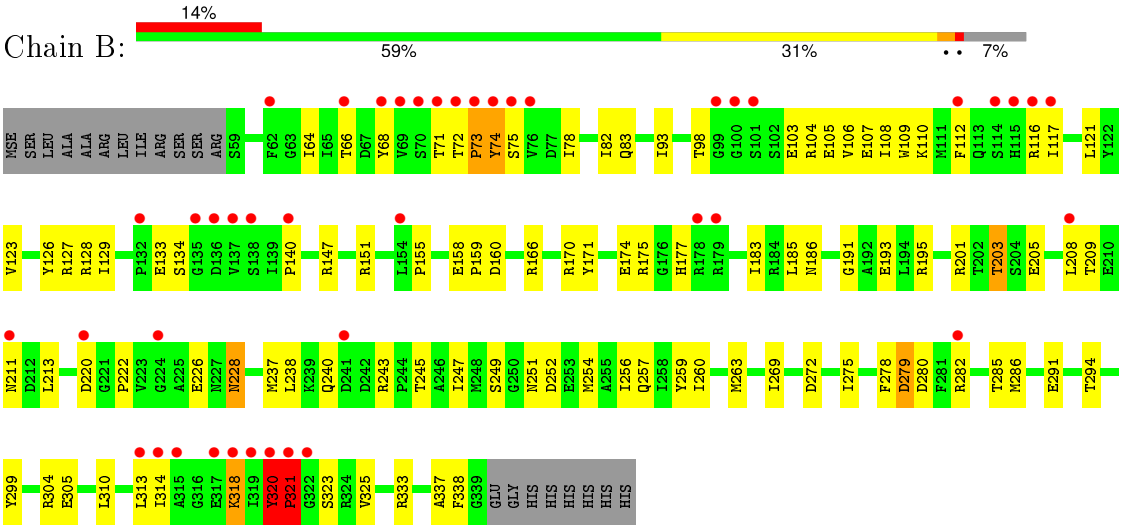
- Molecule 1: Putative transcriptional regulator protein, LacI family

Chain A:



- Molecule 1: Putative transcriptional regulator protein, LacI family

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α, β, γ	69.51Å 71.91Å 212.38Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.22 – 1.91 45.22 – 1.91	Depositor EDS
% Data completeness (in resolution range)	98.9 (45.22-1.91) 89.8 (45.22-1.91)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	6.02 (at 1.91Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.238 , 0.276 0.238 , 0.277	Depositor DCC
R_{free} test set	2067 reflections (5.01%)	DCC
Wilson B-factor (Å^2)	21.6	Xtriage
Anisotropy	0.622	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 30.9	EDS
Estimated twinning fraction	0.146 for -k,-h,-l	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Outliers	0 of 41835 reflections	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	4578	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 6.82% 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	0.32	0/2230	0.64	1/3010 (0.0%)
1	B	0.29	0/2223	0.60	1/3001 (0.0%)
All	All	0.30	0/4453	0.62	2/6011 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	227	ASN	N-CA-C	-6.17	94.34	111.00
1	B	320	TYR	C-N-CD	-5.54	108.41	120.60

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	2199	0	2190	70	0
1	B	2192	0	2180	76	0
2	A	123	0	0	2	0
2	B	64	0	0	5	0
All	All	4578	0	4370	139	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 16.

All (139) 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:82:ILE:HD12	1:A:121:LEU:HD13	1.59	0.84
1:B:203:THR:HG21	1:B:213:LEU:HD11	1.62	0.81
1:A:309:TRP:HE1	1:A:319:ILE:HG21	1.48	0.78
1:B:291:GLU:HB3	1:B:333:ARG:HD3	1.67	0.76
1:A:82:ILE:HD13	1:A:303:GLY:O	1.85	0.76
1:B:185:LEU:HD21	1:B:251:ASN:ND2	2.02	0.75
1:B:203:THR:HG23	1:B:208:LEU:HB2	1.69	0.74
1:B:237:MSE:SE	1:B:247:ILE:HD12	2.39	0.72
1:A:177:HIS:HB3	1:A:180:ILE:HD11	1.72	0.72
1:B:313:LEU:CB	1:B:318:LYS:HD3	2.21	0.71
1:A:199:PHE:O	1:A:203:THR:HG22	1.91	0.71
1:B:313:LEU:HB2	1:B:318:LYS:HD3	1.72	0.71
1:B:310:LEU:O	1:B:314:ILE:HG12	1.93	0.69
1:A:318:LYS:NZ	1:A:318:LYS:HB3	2.09	0.68
1:A:72:THR:HG21	1:A:125:MSE:HB3	1.75	0.68
1:A:291:GLU:HB3	1:A:333:ARG:HD3	1.77	0.67
1:A:147:ARG:HD3	2:B:382:HOH:O	1.95	0.66
1:B:140:PRO:HB3	1:B:313:LEU:HD22	1.76	0.66
1:A:171:TYR:O	1:A:175:ARG:HG2	1.96	0.66
1:B:237:MSE:HA	1:B:240:GLN:HE21	1.62	0.65
1:B:282:ARG:O	1:B:286:MSE:HG2	1.97	0.64
1:A:309:TRP:HE1	1:A:319:ILE:CG2	2.10	0.63
1:B:186:ASN:ND2	1:B:226:GLU:HG2	2.14	0.63
1:A:235:THR:HG22	1:A:239:LYS:HE2	1.79	0.63
1:B:320:TYR:HA	1:B:321:PRO:O	1.98	0.63
1:A:269:ILE:HB	1:A:270:PRO:CD	2.29	0.63
1:A:66:THR:HG23	1:A:97:ASN:OD1	1.98	0.62
1:A:270:PRO:HD3	1:A:290:PRO:HG3	1.80	0.62
1:B:186:ASN:HD22	1:B:226:GLU:HG2	1.65	0.61
1:B:201:ARG:O	1:B:205:GLU:HG3	2.01	0.61
1:A:210:GLU:HA	1:A:213:LEU:HD23	1.83	0.61
1:A:309:TRP:NE1	1:A:319:ILE:HG21	2.16	0.60
1:B:78:ILE:HD11	1:B:299:TYR:CD1	2.36	0.60
1:A:318:LYS:HZ2	1:A:318:LYS:HB3	1.66	0.60
1:B:323:SER:HA	2:B:393:HOH:O	2.02	0.60
1:A:166:ARG:HA	1:A:202:THR:HG21	1.83	0.60
1:A:170:ARG:O	1:A:174:GLU:HG3	2.02	0.59
1:A:177:HIS:CB	1:A:180:ILE:HD11	2.33	0.59
1:B:170:ARG:O	1:B:174:GLU:HG3	2.02	0.59
1:B:222:PRO:O	1:B:226:GLU:HG3	2.03	0.59
1:A:177:HIS:HB3	1:A:180:ILE:CD1	2.32	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:116:ARG:O	1:B:116:ARG:HG3	2.03	0.58
1:A:91:LYS:HE3	1:A:311:ASN:OD1	2.03	0.58
1:B:78:ILE:HD13	1:B:299:TYR:O	2.02	0.58
1:B:64:ILE:HG22	1:B:66:THR:HG23	1.85	0.58
1:A:269:ILE:O	1:A:273:VAL:O	2.22	0.57
1:B:171:TYR:O	1:B:175:ARG:HG2	2.03	0.57
1:B:78:ILE:HD11	1:B:299:TYR:HD1	1.68	0.57
1:B:98:THR:HG21	1:B:105:GLU:HG3	1.86	0.56
1:B:107:GLU:HA	1:B:110:LYS:HD3	1.89	0.55
1:A:168:LEU:HD13	1:A:330:LEU:HD13	1.88	0.55
1:B:71:THR:HG22	1:B:71:THR:O	2.07	0.54
1:B:177:HIS:HE1	1:B:337:ALA:O	1.91	0.54
1:B:278:PHE:O	1:B:279:ASP:HB2	2.08	0.54
1:A:282:ARG:HD2	1:A:282:ARG:H	1.72	0.54
1:B:127:ARG:HD2	1:B:158:GLU:OE1	2.07	0.53
1:A:186:ASN:OD1	1:A:188:ILE:HG22	2.09	0.53
1:B:278:PHE:O	1:B:279:ASP:CB	2.57	0.53
1:B:72:THR:HA	1:B:73:PRO:O	2.09	0.53
1:B:155:PRO:HG2	1:B:318:LYS:HE2	1.90	0.52
1:A:209:THR:O	1:A:213:LEU:HD22	2.09	0.52
1:A:278:PHE:O	1:A:279:ASP:CB	2.57	0.52
1:A:203:THR:OG1	1:A:208:LEU:HB2	2.10	0.52
1:A:166:ARG:HH12	1:B:147:ARG:NH1	2.08	0.51
1:A:329:LYS:HG2	1:B:151:ARG:CZ	2.41	0.51
1:B:104:ARG:O	1:B:108:ILE:HG12	2.11	0.51
1:B:160:ASP:HB2	1:B:325:VAL:HG12	1.91	0.51
1:B:83:GLN:NE2	1:B:93:ILE:HB	2.27	0.50
1:A:166:ARG:CZ	1:B:129:ILE:HD13	2.42	0.50
1:A:62:PHE:HB2	1:A:93:ILE:HD13	1.94	0.50
1:B:269:ILE:HD13	1:B:275:ILE:HD12	1.93	0.50
1:A:77:ASP:HB2	2:A:19:HOH:O	2.12	0.49
1:B:72:THR:OG1	1:B:73:PRO:HA	2.11	0.49
1:B:304:ARG:HG2	1:B:304:ARG:HH11	1.76	0.49
1:A:198:ALA:O	1:A:202:THR:HG23	2.13	0.49
1:A:72:THR:HB	1:A:125:MSE:HG3	1.95	0.48
1:A:268:ARG:HD2	1:A:271:GLN:CD	2.33	0.48
1:A:121:LEU:HD22	1:A:144:ILE:HD11	1.95	0.48
1:B:209:THR:OG1	1:B:211:ASN:OD1	2.31	0.48
1:B:203:THR:HG23	1:B:208:LEU:CB	2.43	0.47
1:A:127:ARG:HA	1:A:145:ASN:O	2.14	0.47
1:A:112:PHE:CD1	1:A:117:ILE:HG13	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:278:PHE:O	1:A:279:ASP:HB2	2.14	0.47
1:B:238:LEU:O	1:B:243:ARG:HG3	2.14	0.47
1:A:269:ILE:O	1:A:270:PRO:C	2.51	0.47
1:A:151:ARG:HG2	1:A:152:GLU:HG3	1.95	0.47
1:A:269:ILE:HB	1:A:270:PRO:HD2	1.97	0.47
1:A:316:GLY:O	1:A:317:GLU:O	2.32	0.47
1:A:129:ILE:HD13	1:B:166:ARG:CZ	2.45	0.47
1:A:75:SER:HB2	1:A:125:MSE:SE	2.65	0.46
1:B:109:TRP:CD1	1:B:134:SER:HA	2.51	0.46
1:A:166:ARG:NH2	1:B:147:ARG:HD2	2.31	0.46
1:A:102:SER:O	1:A:106:VAL:HG23	2.16	0.46
1:A:62:PHE:HB2	1:A:93:ILE:CD1	2.47	0.45
1:A:282:ARG:H	1:A:282:ARG:CD	2.30	0.45
1:A:166:ARG:HH22	1:B:147:ARG:HH11	1.65	0.45
1:A:252:ASP:OD2	1:A:280:ASP:N	2.49	0.45
1:A:300:TYR:CE2	1:A:304:ARG:HD2	2.51	0.45
1:B:252:ASP:OD2	1:B:280:ASP:N	2.49	0.45
1:A:66:THR:OG1	1:A:125:MSE:HE3	2.17	0.44
1:B:75:SER:HB2	1:B:78:ILE:HG12	2.00	0.44
1:A:82:ILE:HD12	1:A:121:LEU:CD1	2.39	0.44
1:A:186:ASN:CG	1:A:188:ILE:HG22	2.38	0.44
1:A:147:ARG:HE	1:A:147:ARG:HB3	1.56	0.44
1:A:147:ARG:NH1	1:B:166:ARG:HH21	2.16	0.44
1:B:243:ARG:NH2	1:B:272:ASP:O	2.51	0.44
1:A:269:ILE:CB	1:A:270:PRO:CD	2.96	0.43
1:B:82:ILE:HD13	1:B:121:LEU:HD13	2.00	0.43
1:A:168:LEU:HD21	1:A:277:GLY:O	2.18	0.43
1:B:106:VAL:HG22	1:B:133:GLU:HG3	1.99	0.43
1:B:280:ASP:HA	1:B:285:THR:HG21	2.00	0.43
1:B:112:PHE:CD1	1:B:117:ILE:HG13	2.54	0.43
1:B:256:ILE:O	1:B:260:ILE:HG12	2.19	0.43
1:A:268:ARG:HD2	1:A:271:GLN:NE2	2.34	0.43
1:B:73:PRO:O	1:B:74:TYR:HB2	2.19	0.42
1:B:175:ARG:HG2	1:B:175:ARG:HH11	1.84	0.42
1:B:191:GLY:O	1:B:195:ARG:HG3	2.19	0.42
1:B:193:GLU:HB3	2:B:387:HOH:O	2.19	0.42
1:B:228:ASN:HD21	1:B:257:GLN:HE22	1.66	0.42
1:B:249:SER:OG	1:B:254:MSE:HB2	2.20	0.42
1:B:205:GLU:HB2	2:B:392:HOH:O	2.19	0.42
1:A:191:GLY:O	1:A:195:ARG:HG3	2.19	0.42
1:A:294:THR:O	1:A:330:LEU:HD12	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:158:GLU:HB3	1:B:159:PRO:HD2	2.01	0.41
1:A:158:GLU:HB3	1:A:159:PRO:HD2	2.03	0.41
1:B:126:TYR:O	1:B:128:ARG:HG3	2.20	0.41
1:B:175:ARG:HD3	1:B:338:PHE:CE1	2.55	0.41
1:B:259:TYR:HB3	1:B:263:MSE:HE3	2.03	0.41
1:B:177:HIS:HD2	1:B:245:THR:OG1	2.04	0.41
1:A:268:ARG:HB2	1:A:271:GLN:HB3	2.02	0.41
1:A:109:TRP:CD1	1:A:134:SER:HA	2.56	0.41
1:A:202:THR:HG22	2:A:376:HOH:O	2.20	0.41
1:B:280:ASP:HA	1:B:294:THR:HG21	2.02	0.41
1:B:66:THR:HG22	1:B:123:VAL:HB	2.03	0.40
1:B:323:SER:HB2	2:B:380:HOH:O	2.20	0.40
1:A:66:THR:HA	1:A:123:VAL:O	2.20	0.40
1:B:103:GLU:HG3	1:B:104:ARG:N	2.36	0.40
1:B:177:HIS:CE1	1:B:337:ALA:O	2.73	0.40
1:B:183:ILE:HG22	1:B:254:MSE:CE	2.51	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	A	279/301 (93%)	264 (95%)	10 (4%)	5 (2%)	11 2
1	B	279/301 (93%)	258 (92%)	15 (5%)	6 (2%)	8 1
All	All	558/602 (93%)	522 (94%)	25 (4%)	11 (2%)	9 2

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	269	ILE
1	A	317	GLU

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Mol	Chain	Res	Type
1	A	321	PRO
1	B	68	TYR
1	B	73	PRO
1	B	321	PRO
1	A	279	ASP
1	B	279	ASP
1	B	318	LYS
1	A	225	ALA
1	B	74	TYR

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	237/243 (98%)	233 (98%)	4 (2%)	68 63
1	B	236/243 (97%)	230 (98%)	6 (2%)	55 46
All	All	473/486 (97%)	463 (98%)	10 (2%)	61 54

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	77	ASP
1	A	145	ASN
1	A	213	LEU
1	A	282	ARG
1	B	203	THR
1	B	220	ASP
1	B	228	ASN
1	B	305	GLU
1	B	320	TYR
1	B	321	PRO

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

Mol	Chain	Res	Type
1	A	115	HIS
1	A	145	ASN
1	B	89	ASN
1	B	177	HIS
1	B	228	ASN
1	B	240	GLN

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

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	272/301 (90%)	0.57	20 (7%) 17 20	11, 20, 39, 59	0
1	B	272/301 (90%)	1.08	43 (15%) 3 3	14, 28, 54, 64	0
All	All	544/602 (90%)	0.83	63 (11%) 6 7	11, 24, 50, 64	0

All (63) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	320	TYR	13.7
1	B	318	LYS	12.7
1	B	319	ILE	10.2
1	B	71	THR	10.1
1	B	73	PRO	8.2
1	A	321	PRO	7.0
1	B	69	VAL	6.8
1	A	319	ILE	6.2
1	A	269	ILE	6.0
1	A	74	TYR	6.0
1	B	68	TYR	5.9
1	A	73	PRO	5.8
1	B	70	SER	5.3
1	B	99	GLY	5.2
1	A	71	THR	5.1
1	B	321	PRO	4.9
1	B	320	TYR	4.7
1	A	72	THR	4.3
1	B	315	ALA	3.9
1	B	112	PHE	3.8
1	B	74	TYR	3.7
1	B	322	GLY	3.7
1	B	317	GLU	3.6
1	B	135	GLY	3.6

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Mol	Chain	Res	Type	RSRZ
1	A	241	ASP	3.5
1	B	75	SER	3.5
1	B	138	SER	3.2
1	B	76	VAL	3.0
1	B	314	ILE	3.0
1	B	116	ARG	2.9
1	B	137	VAL	2.9
1	B	224	GLY	2.9
1	B	72	THR	2.8
1	A	268	ARG	2.8
1	A	210	GLU	2.8
1	B	117	ILE	2.8
1	B	100	GLY	2.7
1	B	62	PHE	2.7
1	A	270	PRO	2.6
1	B	154	LEU	2.6
1	A	318	LYS	2.6
1	B	282	ARG	2.5
1	A	203	THR	2.5
1	B	136	ASP	2.5
1	A	205	GLU	2.4
1	B	313	LEU	2.4
1	B	241	ASP	2.4
1	B	115	HIS	2.4
1	A	316	GLY	2.3
1	B	208	LEU	2.3
1	B	220	ASP	2.3
1	A	271	GLN	2.3
1	A	58	ARG	2.2
1	B	140	PRO	2.2
1	B	211	ASN	2.2
1	B	114	SER	2.2
1	A	242	ASP	2.1
1	B	179	ARG	2.1
1	B	132	PRO	2.1
1	B	101	SER	2.1
1	B	66	THR	2.1
1	B	178	ARG	2.1
1	A	188	ILE	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

There are no carbohydrates in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.