



# Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 01:57 PM GMT

PDB ID : 3CDI  
Title : Crystal structure of E. coli PNPase  
Authors : Shi, Z.; Yang, W.Z.; Lin-Chao, S.; Chak, K.F.; Yuan, H.S.  
Deposited on : 2008-02-27  
Resolution : 2.60 Å(reported)

This is a full wwPDB 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 <http://wwpdb.org/ValidationPDFNotes.html>

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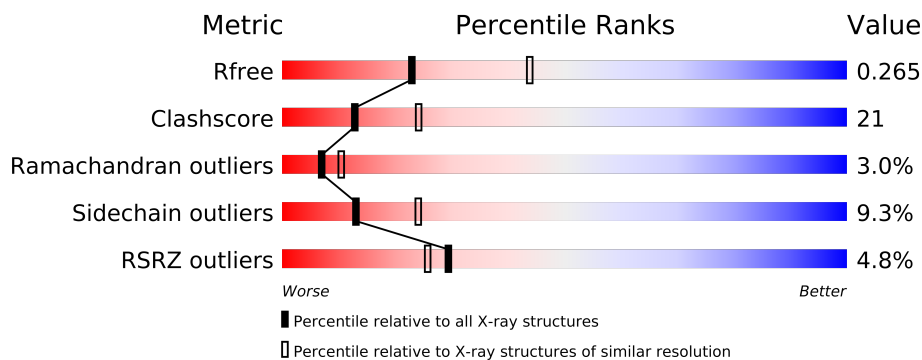
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 2.60 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))
$R_{free}$	66092	1718 (2.60-2.60)
Clashscore	79885	2154 (2.60-2.60)
Ramachandran outliers	78287	2113 (2.60-2.60)
Sidechain outliers	78261	2113 (2.60-2.60)
RSRZ outliers	66119	1718 (2.60-2.60)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	723	

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4025 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Polynucleotide phosphorylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	516	Total	C	N	O	S	0	0	0
			3924	2462	684	759	19			

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	101	Total	O	0	0
			101	101		



## 4 Data and refinement statistics

Property	Value	Source
Space group	F 41 3 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	270.12Å 270.12Å 270.12Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.60 35.17 – 2.60	Depositor EDS
% Data completeness (in resolution range)	93.0 (50.00-2.60) 99.3 (35.17-2.60)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	5.58 (at 2.61Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.232 , 0.277 0.253 , 0.265	Depositor DCC
$R_{free}$ test set	2644 reflections (10.06%)	DCC
Wilson B-factor (Å <sup>2</sup> )	60.6	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 34.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	1 of 26284 reflections (0.004%)	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	4025	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.75% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

## 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	0/3980	0.66	1/5393 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	397	ASP	N-CA-C	-5.60	95.87	111.00

There are no chirality outliers.

There are no planarity outliers.

### 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3924	0	3955	164	0
2	A	101	0	0	5	0
All	All	4025	0	3955	164	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 21.

All (164) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:106:ARG:HB3	1:A:107:PRO:HA	1.25	1.09

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:146:GLN:HG3	1:A:189:THR:HB	1.42	1.00
1:A:171:PRO:HG2	1:A:239:ILE:HG23	1.45	0.99
1:A:396:THR:O	1:A:397:ASP:HB3	1.62	0.97
1:A:106:ARG:HB3	1:A:107:PRO:CA	1.98	0.92
1:A:437:VAL:HG12	1:A:480:ILE:HB	1.52	0.90
1:A:77:GLN:HB3	1:A:131:VAL:HG11	1.57	0.86
1:A:109:GLU:HB2	1:A:424:ILE:HD11	1.58	0.85
1:A:309:GLU:HB2	2:A:817:HOH:O	1.76	0.83
1:A:40:THR:HG23	1:A:57:SER:HB3	1.60	0.81
1:A:557:LEU:O	1:A:561:GLU:HG3	1.80	0.79
1:A:202:VAL:HG12	1:A:212:VAL:HG22	1.64	0.79
1:A:283:ILE:HD11	1:A:289:ARG:HA	1.67	0.77
1:A:66:THR:HG23	2:A:763:HOH:O	1.86	0.75
1:A:372:GLN:HB3	1:A:456:GLU:HB2	1.67	0.74
1:A:294:ASP:HB3	2:A:835:HOH:O	1.87	0.74
1:A:148:ASN:ND2	1:A:150:ASP:H	1.85	0.73
1:A:348:ARG:HD3	1:A:503:ASP:OD1	1.89	0.72
1:A:320:LEU:O	1:A:323:ILE:HD13	1.88	0.72
1:A:148:ASN:HD22	1:A:149:PRO:N	1.88	0.72
1:A:437:VAL:CG1	1:A:480:ILE:HB	2.19	0.71
1:A:74:LYS:HB3	1:A:75:PRO:HD2	1.75	0.69
1:A:258:PRO:O	1:A:259:GLU:HB3	1.91	0.68
1:A:308:ASP:CG	1:A:309:GLU:H	1.95	0.67
1:A:403:TYR:OH	1:A:461:SER:HB2	1.95	0.67
1:A:437:VAL:HG13	1:A:482:ALA:H	1.60	0.66
1:A:513:ASP:HA	1:A:533:LYS:HE2	1.77	0.66
1:A:351:ASP:HB3	1:A:367:THR:HG23	1.77	0.66
1:A:285:ASP:HB2	1:A:288:GLU:HB3	1.79	0.65
1:A:410:VAL:HG12	1:A:410:VAL:O	1.97	0.65
1:A:198:LEU:HD22	1:A:199:ASP:N	2.10	0.64
1:A:30:ARG:HH21	1:A:244:GLU:HG3	1.62	0.64
1:A:200:LEU:HD21	1:A:202:VAL:HG13	1.80	0.63
1:A:59:ASP:O	1:A:60:ASP:HB2	1.97	0.63
1:A:40:THR:CG2	1:A:57:SER:HB3	2.27	0.63
1:A:522:ARG:NH1	1:A:561:GLU:OE1	2.32	0.62
1:A:81:PRO:HB2	1:A:133:GLU:HG3	1.83	0.61
1:A:537:ILE:O	1:A:537:ILE:HD13	2.01	0.61
1:A:489:MET:HB2	1:A:516:PHE:CE2	2.36	0.61
1:A:131:VAL:HG12	1:A:131:VAL:O	2.01	0.60
1:A:280:ALA:O	1:A:283:ILE:HG12	2.00	0.60
1:A:438:MET:HE1	1:A:473:LEU:HD13	1.82	0.60
1:A:268:ARG:O	1:A:272:LEU:HD21	2.02	0.60

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:37:HIS:HD2	1:A:59:ASP:OD2	1.86	0.59
1:A:148:ASN:HD22	1:A:148:ASN:C	2.05	0.59
1:A:242:ILE:O	1:A:246:VAL:HG23	2.02	0.59
1:A:92:ALA:HB1	1:A:143:VAL:HG23	1.85	0.59
1:A:74:LYS:HB2	1:A:77:GLN:HG2	1.83	0.59
1:A:35:GLY:O	1:A:36:GLN:HB3	2.03	0.59
1:A:537:ILE:C	1:A:537:ILE:HD13	2.23	0.58
1:A:291:ALA:O	1:A:295:VAL:HG23	2.02	0.58
1:A:395:ARG:HG2	1:A:398:THR:CG2	2.33	0.58
1:A:55:MET:HG2	1:A:64:PHE:CD1	2.38	0.58
1:A:256:TRP:CZ2	1:A:258:PRO:HA	2.39	0.58
1:A:395:ARG:HD3	1:A:399:PHE:HB2	1.86	0.57
1:A:463:MET:HA	1:A:463:MET:CE	2.34	0.57
1:A:150:ASP:OD1	1:A:151:ILE:N	2.37	0.57
1:A:340:ASP:OD1	1:A:342:ARG:HD3	2.04	0.57
1:A:148:ASN:HD22	1:A:150:ASP:H	1.51	0.55
1:A:537:ILE:O	1:A:537:ILE:HG23	2.07	0.55
1:A:491:LEU:HB3	1:A:514:MET:HB3	1.87	0.55
1:A:463:MET:HA	1:A:463:MET:HE2	1.88	0.55
1:A:395:ARG:NH1	1:A:439:PRO:O	2.41	0.54
1:A:308:ASP:CG	1:A:309:GLU:N	2.60	0.54
1:A:422:ARG:HA	1:A:422:ARG:NE	2.22	0.54
1:A:30:ARG:NH2	1:A:244:GLU:HG3	2.23	0.54
1:A:77:GLN:O	1:A:78:ASP:HB2	2.07	0.54
1:A:277:LEU:HD12	1:A:323:ILE:HD11	1.91	0.53
1:A:55:MET:HG2	1:A:64:PHE:HD1	1.74	0.53
1:A:395:ARG:O	1:A:396:THR:O	2.27	0.53
1:A:79:PHE:O	1:A:131:VAL:HG13	2.09	0.52
1:A:256:TRP:CH2	1:A:258:PRO:HA	2.45	0.52
1:A:321:HIS:NE2	1:A:325:LYS:HE2	2.25	0.52
1:A:308:ASP:O	1:A:309:GLU:HB3	2.09	0.52
1:A:441:MET:HE3	1:A:445:PRO:HA	1.92	0.52
1:A:292:GLN:O	1:A:296:ILE:HG13	2.10	0.51
1:A:259:GLU:HG3	1:A:259:GLU:O	2.11	0.51
1:A:225:LEU:HD21	1:A:542:MET:HB3	1.91	0.51
1:A:202:VAL:HG11	1:A:228:VAL:HA	1.92	0.51
1:A:229:VAL:O	1:A:233:GLU:HG3	2.11	0.51
1:A:79:PHE:C	1:A:131:VAL:HG13	2.30	0.51
1:A:354:THR:CG2	1:A:355:GLY:N	2.73	0.51
1:A:500:VAL:HB	1:A:548:GLN:NE2	2.26	0.51
1:A:276:ARG:HH21	1:A:296:ILE:HA	1.76	0.50
1:A:396:THR:C	1:A:398:THR:H	2.11	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:120:ARG:NH2	1:A:213:GLU:HB2	2.26	0.50
1:A:329:ARG:NH2	1:A:513:ASP:OD1	2.44	0.50
1:A:27:PRO:HB3	1:A:45:MET:HB2	1.92	0.50
1:A:352:VAL:HG13	1:A:471:LEU:HD13	1.93	0.50
1:A:376:THR:HG22	1:A:377:ALA:N	2.27	0.50
1:A:272:LEU:HD23	1:A:272:LEU:N	2.26	0.50
1:A:474:MET:SD	1:A:564:ILE:HD12	2.51	0.50
1:A:190:GLN:HG3	1:A:191:ASP:N	2.26	0.50
1:A:90:THR:OG1	1:A:97:PRO:HD3	2.12	0.49
1:A:395:ARG:HG2	1:A:398:THR:HG22	1.94	0.49
1:A:283:ILE:CD1	1:A:289:ARG:HA	2.40	0.49
1:A:30:ARG:HH21	1:A:244:GLU:CG	2.25	0.49
1:A:213:GLU:OE1	1:A:529:GLN:HG3	2.13	0.49
1:A:130:PHE:CZ	1:A:132:ASN:HB2	2.48	0.49
1:A:351:ASP:OD2	1:A:353:ARG:NH1	2.47	0.48
1:A:258:PRO:O	1:A:259:GLU:CB	2.59	0.47
1:A:277:LEU:CD1	1:A:323:ILE:HD11	2.45	0.47
1:A:261:VAL:O	1:A:263:GLU:N	2.46	0.47
1:A:285:ASP:CB	1:A:288:GLU:HB3	2.45	0.47
1:A:56:VAL:HG21	1:A:156:GLY:HA2	1.97	0.47
1:A:484:VAL:HG23	1:A:520:GLY:C	2.35	0.47
1:A:85:ASN:ND2	1:A:135:GLN:HE21	2.13	0.47
1:A:536:GLY:O	1:A:537:ILE:C	2.51	0.47
1:A:397:ASP:OD2	1:A:397:ASP:C	2.54	0.46
1:A:308:ASP:O	1:A:309:GLU:CB	2.63	0.46
1:A:381:THR:O	1:A:381:THR:HG22	2.15	0.46
1:A:79:PHE:O	1:A:81:PRO:HD3	2.16	0.46
1:A:432:ARG:HA	1:A:435:LEU:HD12	1.97	0.46
1:A:342:ARG:NH2	1:A:503:ASP:HB3	2.30	0.46
1:A:438:MET:CE	1:A:473:LEU:HD13	2.46	0.46
1:A:66:THR:HG22	1:A:137:ILE:HB	1.97	0.46
1:A:55:MET:HE1	1:A:62:ALA:HB1	1.98	0.45
1:A:422:ARG:O	1:A:424:ILE:N	2.49	0.45
1:A:422:ARG:NE	1:A:422:ARG:CA	2.79	0.45
1:A:266:ASN:HB3	1:A:319:ILE:HD13	1.99	0.45
1:A:395:ARG:HG2	1:A:398:THR:HG21	1.98	0.45
1:A:92:ALA:CB	1:A:143:VAL:HG23	2.47	0.45
1:A:120:ARG:HH22	1:A:213:GLU:HB2	1.82	0.45
1:A:297:LYS:HG2	1:A:320:LEU:HD13	1.99	0.45
1:A:315:GLU:O	1:A:318:GLU:HB3	2.18	0.44
1:A:422:ARG:C	1:A:424:ILE:H	2.21	0.44
1:A:441:MET:CE	1:A:445:PRO:HA	2.47	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:30:ARG:HB2	1:A:245:LEU:HD13	1.99	0.44
1:A:25:LEU:HG	1:A:45:MET:HG3	1.98	0.44
1:A:144:ASN:O	1:A:146:GLN:N	2.51	0.44
1:A:454:ILE:HG13	1:A:461:SER:HB3	1.99	0.44
1:A:564:ILE:O	1:A:565:ASN:C	2.55	0.44
1:A:180:ILE:HG12	1:A:195:GLU:O	2.18	0.44
1:A:342:ARG:HB3	1:A:346:MET:HB2	2.01	0.43
1:A:71:LYS:HG2	2:A:833:HOH:O	2.17	0.43
1:A:282:ARG:HH11	1:A:282:ARG:HG3	1.84	0.43
1:A:316:LEU:HD23	1:A:316:LEU:HA	1.85	0.43
1:A:30:ARG:HG2	2:A:830:HOH:O	2.17	0.43
1:A:52:ALA:HB2	1:A:164:SER:HA	2.01	0.43
1:A:265:LEU:O	1:A:269:VAL:HG23	2.19	0.42
1:A:313:GLU:O	1:A:316:LEU:HB2	2.19	0.42
1:A:443:LYS:HE2	1:A:479:PRO:HB2	2.02	0.42
1:A:357:LEU:HA	1:A:358:PRO:HD3	1.75	0.42
1:A:67:VAL:HG22	1:A:136:VAL:HG22	2.01	0.42
1:A:189:THR:HG23	1:A:192:GLU:OE1	2.20	0.42
1:A:148:ASN:HA	1:A:149:PRO:HD3	1.90	0.42
1:A:340:ASP:OD1	1:A:342:ARG:CD	2.67	0.42
1:A:428:ARG:O	1:A:432:ARG:HG3	2.20	0.42
1:A:289:ARG:O	1:A:293:VAL:HG23	2.20	0.41
1:A:505:LEU:HB2	1:A:508:GLU:HG3	2.01	0.41
1:A:395:ARG:O	1:A:396:THR:C	2.59	0.41
1:A:265:LEU:C	1:A:267:ALA:H	2.23	0.41
1:A:120:ARG:NH1	1:A:531:ASP:OD2	2.51	0.41
1:A:184:TYR:OH	1:A:226:GLY:HA3	2.21	0.41
1:A:512:GLY:C	1:A:514:MET:N	2.73	0.41
1:A:437:VAL:O	1:A:437:VAL:HG12	2.20	0.41
1:A:62:ALA:HB3	1:A:142:SER:OG	2.20	0.41
1:A:127:PRO:HD2	1:A:167:PRO:O	2.21	0.40
1:A:532:ILE:N	1:A:532:ILE:HD12	2.35	0.40
1:A:106:ARG:CB	1:A:107:PRO:CA	2.83	0.40
1:A:277:LEU:HD23	1:A:296:ILE:HD12	2.03	0.40
1:A:148:ASN:HD22	1:A:149:PRO:CD	2.34	0.40
1:A:78:ASP:O	1:A:79:PHE:HB2	2.22	0.40
1:A:262:ASN:OD1	1:A:265:LEU:HB2	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	508/723 (70%)	454 (89%)	39 (8%)	15 (3%)	7 10

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	396	THR
1	A	423	GLU
1	A	537	ILE
1	A	35	GLY
1	A	260	PRO
1	A	309	GLU
1	A	458	ASN
1	A	397	ASP
1	A	257	GLN
1	A	259	GLU
1	A	262	ASN
1	A	440	ASP
1	A	145	PRO
1	A	79	PHE
1	A	459	GLY

### 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	420/587 (72%)	381 (91%)	39 (9%)	13 24

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	26	ASN
1	A	30	ARG
1	A	38	THR
1	A	40	THR
1	A	55	MET
1	A	66	THR
1	A	71	LYS
1	A	83	THR
1	A	109	GLU
1	A	112	THR
1	A	116	ARG
1	A	139	THR
1	A	148	ASN
1	A	151	ILE
1	A	190	GLN
1	A	198	LEU
1	A	211	MET
1	A	221	GLU
1	A	222	ASP
1	A	259	GLU
1	A	260	PRO
1	A	265	LEU
1	A	272	LEU
1	A	287	GLN
1	A	323	ILE
1	A	365	LEU
1	A	367	THR
1	A	371	THR
1	A	374	LEU
1	A	397	ASP
1	A	398	THR
1	A	400	LEU
1	A	491	LEU
1	A	504	ILE
1	A	528	LEU
1	A	537	ILE
1	A	539	LYS
1	A	542	MET
1	A	564	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	33	GLN
1	A	37	HIS
1	A	85	ASN
1	A	146	GLN
1	A	148	ASN
1	A	257	GLN
1	A	266	ASN
1	A	287	GLN
1	A	426	HIS
1	A	510	HIS
1	A	543	GLN
1	A	548	GLN

### 5.3.3 RNA ⓘ

There are no RNA chains 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<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	516/723 (71%)	0.28	25 (4%)	29 26	27, 49, 88, 97	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	287	GLN	4.2
1	A	75	PRO	3.5
1	A	314	ASN	3.5
1	A	98	GLY	3.1
1	A	395	ARG	2.9
1	A	415	MET	2.8
1	A	306	ALA	2.8
1	A	78	ASP	2.8
1	A	397	ASP	2.7
1	A	310	THR	2.6
1	A	305	LEU	2.6
1	A	290	TYR	2.5
1	A	153	ALA	2.5
1	A	396	THR	2.4
1	A	258	PRO	2.3
1	A	301	ILE	2.3
1	A	265	LEU	2.3
1	A	304	LEU	2.3
1	A	25	LEU	2.3
1	A	381	THR	2.3
1	A	157	ALA	2.2
1	A	424	ILE	2.2
1	A	317	GLY	2.1
1	A	313	GLU	2.0
1	A	312	ASP	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.