



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 8, 2014 – 11:24 AM EDT

PDB ID : 4HQ6
Title : BC domain in the presence of citrate
Authors : Heo, Y.S.
Deposited on : 2012-10-25
Resolution : 2.70 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

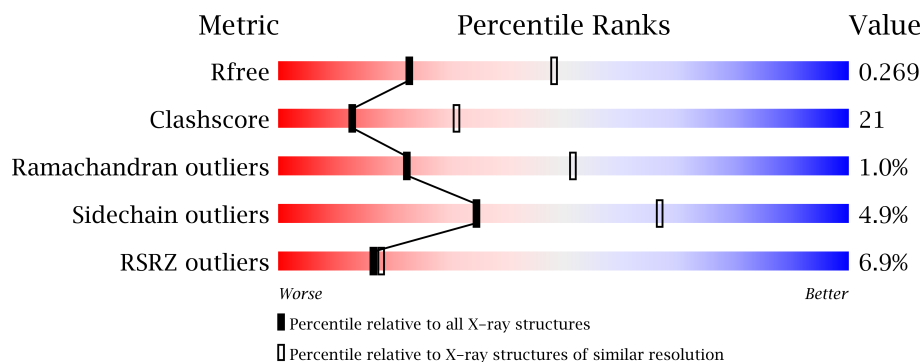
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.16 November 2013
Xtriage (Phenix) : dev-1439
EDS : stable23489
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) : stable23489

1 Overall quality at a glance

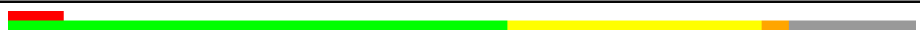
The reported resolution of this entry is 2.70 Å.

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	1557 (2.70-2.70)
Clashscore	79885	1939 (2.70-2.70)
Ramachandran outliers	78287	1905 (2.70-2.70)
Sidechain outliers	78261	1905 (2.70-2.70)
RSRZ outliers	66119	1559 (2.70-2.70)

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. 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	573	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3915 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 Acetyl-CoA carboxylase 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	494	Total	C	N	O	S	0	0	0
			3871	2468	671	714	18			

There are 15 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	211	MET	-	EXPRESSION TAG	UNP O00763
A	212	ARG	-	EXPRESSION TAG	UNP O00763
A	213	GLY	-	EXPRESSION TAG	UNP O00763
A	214	SER	-	EXPRESSION TAG	UNP O00763
A	215	GLY	-	EXPRESSION TAG	UNP O00763
A	216	SER	-	EXPRESSION TAG	UNP O00763
A	236	ILE	LEU	ENGINEERED MUTATION	UNP O00763
A	776	LEU	VAL	ENGINEERED MUTATION	UNP O00763
A	777	GLU	-	EXPRESSION TAG	UNP O00763
A	778	HIS	-	EXPRESSION TAG	UNP O00763
A	779	HIS	-	EXPRESSION TAG	UNP O00763
A	780	HIS	-	EXPRESSION TAG	UNP O00763
A	781	HIS	-	EXPRESSION TAG	UNP O00763
A	782	HIS	-	EXPRESSION TAG	UNP O00763
A	783	HIS	-	EXPRESSION TAG	UNP O00763

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	44	Total	O	0	0
			44	44		

4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	75.59Å 75.59Å 188.78Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	19.98 – 2.70 32.71 – 2.70	Depositor EDS
% Data completeness (in resolution range)	86.4 (19.98-2.70) 86.2 (32.71-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.53 (at 2.68Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.221 , 0.273 0.219 , 0.269	Depositor DCC
R_{free} test set	760 reflections (4.94%)	DCC
Wilson B-factor (Å ²)	50.5	Xtriage
Anisotropy	0.569	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 33.5	EDS
Estimated twinning fraction	0.039 for -h,-k,l	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 16508 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3915	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.45% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities 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	0.41	0/3961	0.60	0/5376

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	623	TYR	Sidechain

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	3871	0	3826	159	0
2	A	44	0	0	5	0
All	All	3915	0	3826	159	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 (159) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:464:ILE:O	1:A:465:ARG:HD3	1.78	0.84
1:A:739:ASN:O	1:A:743:THR:HG23	1.78	0.83
1:A:753:THR:O	1:A:756:LEU:HD13	1.82	0.79
1:A:478:ARG:HA	1:A:481:GLN:HE21	1.48	0.79
1:A:709:ASN:ND2	1:A:712:GLU:H	1.82	0.78
1:A:267:ASN:ND2	1:A:268:ASN:H	1.83	0.76
1:A:254:PHE:HA	1:A:615:ARG:HH12	1.50	0.76
1:A:267:ASN:HD22	1:A:268:ASN:H	1.29	0.76
1:A:525:ARG:HB3	2:A:827:HOH:O	1.88	0.72
1:A:391:VAL:HG13	1:A:422:ILE:HD13	1.69	0.71
1:A:436:ASP:OD1	1:A:438:ASP:HB2	1.91	0.70
1:A:538:ILE:HD11	1:A:751:ILE:HA	1.73	0.70
1:A:296:VAL:HG21	1:A:342:ILE:HD13	1.74	0.69
1:A:655:SER:O	1:A:733:THR:HG21	1.93	0.69
1:A:254:PHE:HA	1:A:615:ARG:NH1	2.10	0.66
1:A:708:GLU:CD	1:A:708:GLU:H	1.99	0.66
1:A:621:LEU:HD21	1:A:627:PRO:HG3	1.78	0.65
1:A:385:LYS:HG2	1:A:490:PHE:CZ	2.32	0.65
1:A:273:VAL:HG21	1:A:669:VAL:HG11	1.78	0.65
1:A:435:LYS:HE3	1:A:439:GLU:OE2	1.96	0.65
1:A:552:ILE:HG22	1:A:556:LYS:HD2	1.77	0.65
1:A:276:MET:HE2	1:A:314:ALA:HA	1.79	0.63
1:A:267:ASN:HD22	1:A:268:ASN:N	1.96	0.62
1:A:378:ALA:N	1:A:421:ARG:NH1	2.47	0.62
1:A:364:LEU:HD22	1:A:369:VAL:HG11	1.81	0.62
1:A:280:ARG:HB3	1:A:290:GLU:HG2	1.81	0.61
1:A:318:VAL:HG21	1:A:338:ILE:HD13	1.82	0.61
1:A:709:ASN:HD22	1:A:709:ASN:C	2.03	0.61
1:A:325:ASN:HA	1:A:328:ASN:OD1	2.01	0.60
1:A:435:LYS:HG2	1:A:439:GLU:OE2	2.01	0.59
1:A:276:MET:HE1	1:A:314:ALA:HB2	1.84	0.59
1:A:403:SER:OG	1:A:432:GLY:O	2.18	0.59
1:A:473:PHE:HB3	1:A:474:PRO:HD3	1.85	0.58
1:A:378:ALA:N	1:A:421:ARG:HH11	2.02	0.58
1:A:276:MET:HE3	1:A:295:PHE:HB3	1.86	0.57
1:A:594:MET:CE	1:A:681:TRP:HE1	2.18	0.57
1:A:289:ASN:ND2	1:A:291:ARG:H	2.03	0.57
1:A:291:ARG:HD2	2:A:808:HOH:O	2.05	0.57
1:A:351:TRP:HA	1:A:584:ARG:HD2	1.86	0.57

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:419:GLY:O	1:A:420:LYS:HB3	2.03	0.57
1:A:452:MET:SD	1:A:494:LEU:HD13	2.45	0.57
1:A:523:ILE:HD12	1:A:523:ILE:N	2.20	0.57
1:A:655:SER:HB2	1:A:684:PHE:CZ	2.40	0.57
1:A:584:ARG:HG3	1:A:584:ARG:HH11	1.70	0.56
1:A:276:MET:CE	1:A:314:ALA:HA	2.35	0.56
1:A:462:LYS:HE2	1:A:483:GLU:HG2	1.88	0.56
1:A:756:LEU:HD12	1:A:756:LEU:H	1.70	0.55
1:A:262:LYS:HG3	1:A:344:VAL:HG12	1.88	0.55
1:A:534:ALA:HB2	1:A:591:CYS:HB3	1.87	0.55
1:A:737:LEU:O	1:A:741:LEU:HG	2.05	0.55
1:A:756:LEU:HD12	1:A:756:LEU:N	2.21	0.55
1:A:530:ILE:HG22	1:A:531:VAL:HG23	1.88	0.55
1:A:727:ARG:HB3	1:A:729:ASP:OD1	2.07	0.54
1:A:364:LEU:HD22	1:A:369:VAL:CG1	2.38	0.53
1:A:305:LYS:HA	1:A:305:LYS:HE3	1.90	0.53
1:A:594:MET:HE1	1:A:681:TRP:HE1	1.74	0.53
1:A:561:VAL:O	1:A:562:SER:HB3	2.09	0.53
1:A:504:GLN:HG2	1:A:592:THR:HG21	1.92	0.53
1:A:709:ASN:HD21	1:A:712:GLU:H	1.54	0.53
1:A:724:LEU:HD22	1:A:730:PHE:CG	2.44	0.52
1:A:594:MET:HE1	1:A:704:PHE:HB3	1.91	0.52
1:A:436:ASP:HB2	2:A:803:HOH:O	2.10	0.51
1:A:276:MET:HE1	1:A:314:ALA:CB	2.40	0.51
1:A:296:VAL:HG11	1:A:342:ILE:HD11	1.91	0.51
1:A:385:LYS:HD2	1:A:389:THR:OG1	2.10	0.51
1:A:594:MET:HE1	1:A:704:PHE:CB	2.39	0.51
1:A:534:ALA:HB3	1:A:648:VAL:HB	1.93	0.51
1:A:325:ASN:N	1:A:325:ASN:HD22	2.07	0.50
1:A:377:GLU:C	1:A:421:ARG:HH11	2.14	0.50
1:A:594:MET:HE1	1:A:681:TRP:NE1	2.26	0.50
1:A:594:MET:CE	1:A:704:PHE:HB3	2.42	0.50
1:A:450:PRO:HB3	2:A:812:HOH:O	2.11	0.50
1:A:569:LEU:O	1:A:576:PHE:HA	2.12	0.49
1:A:289:ASN:HD22	1:A:290:GLU:N	2.09	0.49
1:A:340:LYS:O	1:A:343:PRO:HB3	2.12	0.49
1:A:572:GLN:CD	1:A:572:GLN:H	2.15	0.49
1:A:301:PRO:HD3	1:A:320:VAL:O	2.13	0.49
1:A:584:ARG:HH12	1:A:586:GLN:HG2	1.77	0.49
1:A:247:PRO:O	1:A:251:VAL:HG23	2.14	0.48
1:A:262:LYS:HG2	1:A:344:VAL:HA	1.95	0.48
1:A:708:GLU:OE1	1:A:712:GLU:HG3	2.14	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:296:VAL:HA	1:A:316:HIS:O	2.14	0.48
1:A:299:VAL:HG12	1:A:304:LEU:HB2	1.95	0.47
1:A:676:SER:HB3	1:A:719:VAL:HG12	1.95	0.47
1:A:397:VAL:HG22	1:A:550:CYS:HB3	1.96	0.47
1:A:669:VAL:HG13	1:A:669:VAL:O	2.14	0.47
1:A:465:ARG:HG3	1:A:476:LEU:HD22	1.96	0.47
1:A:589:HIS:ND1	1:A:590:PRO:HD3	2.30	0.47
1:A:710:ARG:O	1:A:714:ILE:HG13	2.15	0.47
1:A:728:GLY:O	1:A:731:ARG:HG2	2.14	0.47
1:A:286:MET:HE2	1:A:286:MET:HA	1.96	0.47
1:A:457:GLU:HB2	1:A:487:SER:HB2	1.97	0.47
1:A:287:PHE:C	1:A:289:ASN:N	2.67	0.46
1:A:325:ASN:N	1:A:325:ASN:ND2	2.64	0.46
1:A:495:ALA:HB1	1:A:569:LEU:HD21	1.98	0.46
1:A:647:HIS:CD2	1:A:710:ARG:HA	2.51	0.46
1:A:454:LYS:HG2	1:A:464:ILE:HG23	1.97	0.46
1:A:457:GLU:HB2	1:A:487:SER:CB	2.45	0.46
1:A:465:ARG:CG	1:A:476:LEU:HD22	2.45	0.46
1:A:457:GLU:HB2	1:A:487:SER:OG	2.16	0.45
1:A:656:GLU:OE1	1:A:656:GLU:HA	2.17	0.45
1:A:380:TRP:HA	1:A:380:TRP:CE3	2.52	0.45
1:A:300:THR:HG22	1:A:320:VAL:HG23	1.99	0.44
1:A:523:ILE:CD1	1:A:523:ILE:N	2.80	0.44
1:A:439:GLU:O	1:A:442:GLU:HB2	2.17	0.44
1:A:357:ASN:O	1:A:361:PRO:HD2	2.17	0.44
1:A:589:HIS:N	1:A:590:PRO:CD	2.80	0.44
1:A:614:HIS:HD2	2:A:830:HOH:O	2.00	0.44
1:A:267:ASN:ND2	1:A:268:ASN:N	2.59	0.44
1:A:286:MET:HA	1:A:286:MET:CE	2.48	0.44
1:A:523:ILE:H	1:A:523:ILE:CD1	2.31	0.44
1:A:268:ASN:HA	1:A:272:ALA:HB2	2.00	0.44
1:A:496:GLN:O	1:A:497:HIS:C	2.56	0.44
1:A:336:VAL:HG21	1:A:363:LEU:CB	2.48	0.43
1:A:545:GLU:O	1:A:548:GLU:HB2	2.18	0.43
1:A:668:THR:O	1:A:685:SER:HA	2.18	0.43
1:A:264:LEU:HB3	1:A:347:VAL:HG22	1.99	0.43
1:A:357:ASN:HD21	1:A:359:LYS:HE3	1.82	0.43
1:A:268:ASN:HB2	1:A:310:TYR:CE2	2.53	0.43
1:A:499:ARG:HD3	1:A:751:ILE:O	2.17	0.43
1:A:756:LEU:H	1:A:756:LEU:CD1	2.31	0.43
1:A:683:TYR:HE1	1:A:702:HIS:HB2	1.84	0.43
1:A:377:GLU:HB2	1:A:421:ARG:NH1	2.34	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:655:SER:N	1:A:684:PHE:HZ	2.16	0.43
1:A:751:ILE:HG23	1:A:752:ASP:N	2.34	0.43
1:A:520:ASP:HB3	1:A:533:GLU:HB2	2.00	0.42
1:A:410:GLU:H	1:A:410:GLU:HG3	1.62	0.42
1:A:262:LYS:CG	1:A:344:VAL:HA	2.49	0.42
1:A:318:VAL:HA	1:A:319:PRO:HD3	1.83	0.42
1:A:409:VAL:HG11	1:A:423:SER:O	2.19	0.42
1:A:584:ARG:NH1	1:A:584:ARG:HG3	2.34	0.42
1:A:318:VAL:O	1:A:318:VAL:HG23	2.20	0.42
1:A:361:PRO:HB3	1:A:371:PHE:CD2	2.55	0.42
1:A:303:ASP:OD1	1:A:668:THR:HG21	2.20	0.42
1:A:725:SER:HA	1:A:730:PHE:O	2.20	0.42
1:A:630:VAL:HG12	1:A:630:VAL:O	2.19	0.42
1:A:561:VAL:O	1:A:562:SER:CB	2.68	0.41
1:A:523:ILE:H	1:A:523:ILE:HD12	1.86	0.41
1:A:709:ASN:C	1:A:709:ASN:ND2	2.73	0.41
1:A:276:MET:HE3	1:A:295:PHE:CB	2.49	0.41
1:A:496:GLN:O	1:A:497:HIS:O	2.39	0.41
1:A:587:VAL:HG13	1:A:588:GLU:HG3	2.01	0.41
1:A:496:GLN:HE21	1:A:496:GLN:HB2	1.62	0.41
1:A:736:TYR:CE2	1:A:737:LEU:HG	2.56	0.41
1:A:336:VAL:HG21	1:A:363:LEU:HB3	2.03	0.41
1:A:332:VAL:O	1:A:336:VAL:HG23	2.21	0.41
1:A:446:ARG:HH11	1:A:446:ARG:HG2	1.85	0.41
1:A:299:VAL:O	1:A:319:PRO:HA	2.21	0.41
1:A:683:TYR:CE1	1:A:702:HIS:HB2	2.56	0.41
1:A:289:ASN:C	1:A:289:ASN:HD22	2.24	0.41
1:A:418:GLN:HG2	1:A:418:GLN:O	2.20	0.41
1:A:551:ALA:HB1	1:A:566:VAL:HG21	2.03	0.41
1:A:275:CYS:HB2	1:A:348:TRP:CH2	2.57	0.40
1:A:402:TRP:HB2	1:A:491:LEU:O	2.20	0.40
1:A:402:TRP:CH2	1:A:404:GLY:HA3	2.57	0.40
1:A:709:ASN:HD22	1:A:712:GLU:H	1.64	0.40
1:A:726:ILE:O	1:A:727:ARG:C	2.58	0.40
1:A:257:ASP:O	1:A:258:ARG:HB2	2.20	0.40
1:A:426:GLU:O	1:A:426:GLU:HG2	2.21	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	486/573 (85%)	448 (92%)	33 (7%)	5 (1%)	22	51

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	497	HIS
1	A	562	SER
1	A	573	ASP
1	A	574	GLY
1	A	420	LYS

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	412/477 (86%)	392 (95%)	20 (5%)	35	67

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	267	ASN
1	A	281	ARG
1	A	289	ASN
1	A	305	LYS
1	A	325	ASN
1	A	343	PRO
1	A	410	GLU
1	A	445	GLU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	465	ARG
1	A	497	HIS
1	A	528	GLN
1	A	572	GLN
1	A	573	ASP
1	A	615	ARG
1	A	708	GLU
1	A	709	ASN
1	A	711	GLU
1	A	724	LEU
1	A	729	ASP
1	A	749	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	267	ASN
1	A	289	ASN
1	A	325	ASN
1	A	353	HIS
1	A	367	ASN
1	A	481	GLN
1	A	496	GLN
1	A	524	GLN
1	A	528	GLN
1	A	699	GLN
1	A	709	ASN
1	A	716	ASN
1	A	739	ASN
1	A	747	GLN
1	A	748	ASN
1	A	749	ASN

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, 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	494/573 (86%)	0.37	34 (6%) 17 18	31, 53, 79, 103	0

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	572	GLN	5.2
1	A	686	VAL	4.7
1	A	564	GLY	4.4
1	A	563	ALA	4.4
1	A	667	GLY	4.0
1	A	420	LYS	3.7
1	A	565	THR	3.4
1	A	699	GLN	3.3
1	A	418	GLN	3.0
1	A	730	PHE	2.9
1	A	590	PRO	2.9
1	A	685	SER	2.7
1	A	700	PHE	2.7
1	A	504	GLN	2.7
1	A	419	GLY	2.6
1	A	666	SER	2.6
1	A	582	ASN	2.5
1	A	240	ARG	2.5
1	A	589	HIS	2.5
1	A	668	THR	2.5
1	A	562	SER	2.5
1	A	586	GLN	2.4
1	A	587	VAL	2.3
1	A	526	ARG	2.3
1	A	525	ARG	2.2
1	A	759	LEU	2.2
1	A	366	LYS	2.2

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	363	LEU	2.2
1	A	684	PHE	2.2
1	A	411	TRP	2.1
1	A	600	LEU	2.1
1	A	275	CYS	2.1
1	A	585	LEU	2.1
1	A	656	GLU	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.