



Full wwPDB X-ray Structure Validation Report

Feb 26, 2014 – 03:25 PM GMT

PDB ID : 3HO4
Title : Crystal structure of Hedgehog-interacting protein (HHIP)
Authors : Hymowitz, S.G.; Bosanac, I.
Deposited on : 2009-06-01
Resolution : 3.10 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>

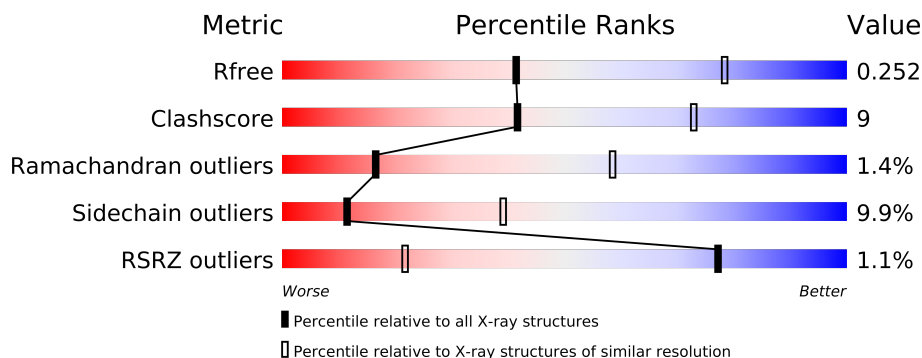
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 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 3.10 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	1007 (3.18-3.02)
Clashscore	79885	1078 (3.16-3.04)
Ramachandran outliers	78287	1044 (3.16-3.04)
Sidechain outliers	78261	1044 (3.16-3.04)
RSRZ outliers	66119	1008 (3.18-3.02)

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	481	
1	B	481	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 6868 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 Hedgehog-interacting protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	437	Total	C	N	O	S	Se	5	0	0
			3414	2143	608	636	22	5			
1	B	442	Total	C	N	O	S	Se	9	0	0
			3454	2166	616	644	22	6			

There are 12 discrepancies between the modelled and reference sequences:

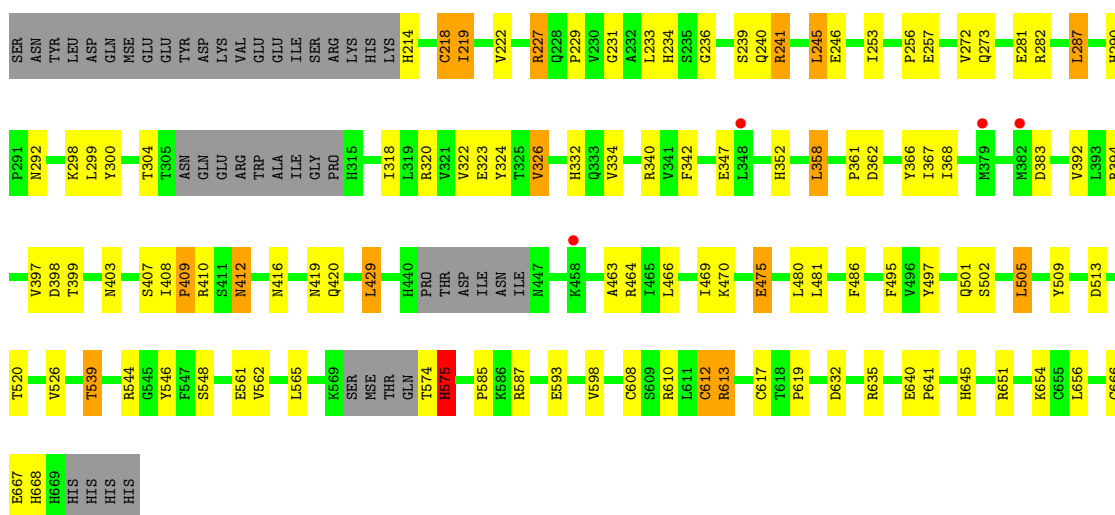
Chain	Residue	Modelled	Actual	Comment	Reference
A	668	HIS	-	EXPRESSION TAG	UNP Q96QV1
A	669	HIS	-	EXPRESSION TAG	UNP Q96QV1
A	670	HIS	-	EXPRESSION TAG	UNP Q96QV1
A	671	HIS	-	EXPRESSION TAG	UNP Q96QV1
A	672	HIS	-	EXPRESSION TAG	UNP Q96QV1
A	673	HIS	-	EXPRESSION TAG	UNP Q96QV1
B	668	HIS	-	EXPRESSION TAG	UNP Q96QV1
B	669	HIS	-	EXPRESSION TAG	UNP Q96QV1
B	670	HIS	-	EXPRESSION TAG	UNP Q96QV1
B	671	HIS	-	EXPRESSION TAG	UNP Q96QV1
B	672	HIS	-	EXPRESSION TAG	UNP Q96QV1
B	673	HIS	-	EXPRESSION TAG	UNP Q96QV1

3 Residue-property plots

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.

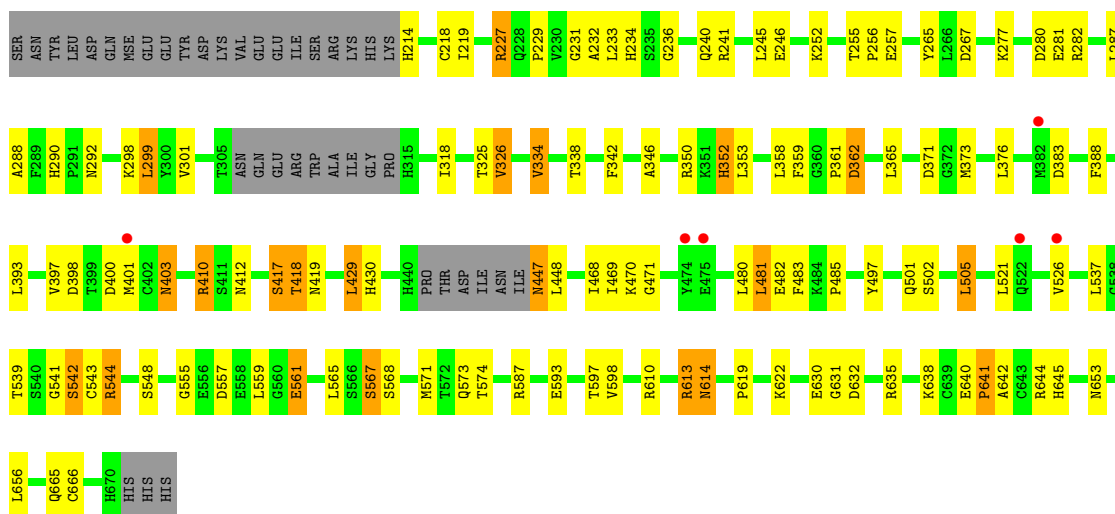
• Molecule 1: Hedgehog-interacting protein

Chain A: 



• Molecule 1: Hedgehog-interacting protein

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	101.04Å 101.04Å 304.92Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 – 3.10 29.17 – 3.10	Depositor EDS
% Data completeness (in resolution range)	99.9 (30.00-3.10) 99.9 (29.17-3.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.36 (at 3.11Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.217 , 0.252 0.226 , 0.252	Depositor DCC
R_{free} test set	1709 reflections (5.35%)	DCC
Wilson B-factor (Å ²)	76.8	Xtriage
Anisotropy	0.302	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 32.1	EDS
Estimated twinning fraction	0.027 for -h,-k,l	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 33679 reflections	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	6868	wwPDB-VP
Average B, all atoms (Å ²)	73.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.46% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality i

5.1 Standard geometry i

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	0.42	1/3488 (0.0%)	0.61	1/4699 (0.0%)
1	B	0.41	1/3529 (0.0%)	0.60	1/4754 (0.0%)
All	All	0.41	2/7017 (0.0%)	0.61	2/9453 (0.0%)

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	2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	218	CYS	CB-SG	-5.94	1.72	1.81
1	A	218	CYS	CB-SG	-5.68	1.72	1.81

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	277	LYS	CB-CA-C	-6.03	98.33	110.40
1	A	299	LEU	CA-CB-CG	5.52	128.00	115.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	281	GLU	Peptide
1	A	409	PRO	Peptide

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	3414	0	3317	61	0
1	B	3454	0	3354	64	0
All	All	6868	0	6671	122	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 9.

All (122) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:574:THR:O	1:A:575:HIS:HB2	1.80	0.82
1:B:544:ARG:HG2	1:B:544:ARG:HH11	1.46	0.80
1:A:241:ARG:HH11	1:A:332:HIS:HD2	1.31	0.77
1:B:281:GLU:HB3	1:B:353:LEU:HD22	1.69	0.75
1:B:544:ARG:CG	1:B:544:ARG:HH11	1.99	0.75
1:B:231:GLY:HA2	1:B:565:LEU:HD13	1.70	0.72
1:B:290:HIS:HD2	1:B:292:ASN:H	1.37	0.72
1:B:567:SER:HA	1:B:571:MSE:HE2	1.70	0.72
1:B:501:GLN:O	1:B:587:ARG:HD3	1.94	0.67
1:A:290:HIS:HD2	1:A:292:ASN:H	1.42	0.67
1:A:241:ARG:HH11	1:A:332:HIS:CD2	2.11	0.67
1:B:410:ARG:H	1:B:410:ARG:CD	2.08	0.67
1:A:290:HIS:CD2	1:A:292:ASN:H	2.13	0.66
1:A:539:THR:HB	1:A:544:ARG:HA	1.76	0.66
1:A:394:ARG:HE	1:A:412:ASN:HD21	1.44	0.65
1:A:469:ILE:HG12	1:A:470:LYS:H	1.62	0.64
1:B:290:HIS:CD2	1:B:292:ASN:H	2.15	0.63
1:B:429:LEU:HD13	1:B:480:LEU:HD21	1.79	0.63
1:A:234:HIS:CD2	1:A:236:GLY:H	2.16	0.62
1:A:234:HIS:HD2	1:A:236:GLY:H	1.47	0.62
1:B:630:GLU:HB3	1:B:638:LYS:HE3	1.81	0.62
1:A:358:LEU:HG	1:A:366:TYR:HB2	1.82	0.61
1:A:429:LEU:HD13	1:A:480:LEU:HD21	1.82	0.61
1:A:368:ILE:HG12	1:A:392:VAL:HG22	1.83	0.60
1:A:282:ARG:O	1:A:304:THR:HB	2.02	0.60

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:342:PHE:HE1	1:B:397:VAL:HG12	1.69	0.58
1:A:501:GLN:O	1:A:587:ARG:HD3	2.04	0.57
1:A:394:ARG:HE	1:A:412:ASN:ND2	2.03	0.57
1:B:613:ARG:O	1:B:614:ASN:HB2	2.05	0.56
1:A:231:GLY:HA2	1:A:565:LEU:HD13	1.87	0.56
1:B:401:MSE:HE3	1:B:403:ASN:HD21	1.71	0.56
1:A:495:PHE:CZ	1:A:562:VAL:HG11	2.41	0.55
1:B:568:SER:H	1:B:571:MSE:HB2	1.70	0.55
1:A:475:GLU:HA	1:A:475:GLU:OE1	2.06	0.55
1:A:495:PHE:CE2	1:A:562:VAL:HG11	2.42	0.55
1:B:645:HIS:HD2	1:B:666:CYS:O	1.90	0.54
1:A:466:LEU:HB2	1:A:480:LEU:HD12	1.89	0.54
1:B:544:ARG:NH1	1:B:544:ARG:CG	2.64	0.54
1:A:651:ARG:HH11	1:A:654:LYS:HG2	1.72	0.54
1:A:640:GLU:HB2	1:A:641:PRO:CD	2.38	0.53
1:A:651:ARG:NH1	1:A:654:LYS:HG2	2.23	0.53
1:B:326:VAL:HA	1:B:334:VAL:HA	1.90	0.52
1:B:448:LEU:HD23	1:B:468:ILE:HD12	1.92	0.51
1:A:245:LEU:HD11	1:A:287:LEU:HB2	1.92	0.51
1:A:463:ALA:HB2	1:A:486:PHE:CE2	2.46	0.51
1:A:241:ARG:NH1	1:A:332:HIS:CD2	2.79	0.50
1:B:290:HIS:HD2	1:B:292:ASN:N	2.08	0.50
1:A:300:TYR:OH	1:A:397:VAL:HG11	2.12	0.50
1:A:361:PRO:O	1:A:362:ASP:HB2	2.10	0.50
1:A:229:PRO:HB2	1:A:565:LEU:HB3	1.93	0.50
1:B:219:ILE:HD13	1:B:537:LEU:HD13	1.94	0.50
1:B:229:PRO:HB2	1:B:565:LEU:HB3	1.94	0.49
1:A:326:VAL:HA	1:A:334:VAL:HA	1.95	0.49
1:B:227:ARG:HG3	1:B:246:GLU:CD	2.32	0.49
1:B:557:ASP:HB3	1:B:559:LEU:H	1.78	0.49
1:B:298:LYS:NZ	1:B:325:THR:OG1	2.46	0.49
1:B:265:TYR:O	1:B:334:VAL:HG13	2.13	0.48
1:B:287:LEU:HD23	1:B:288:ALA:N	2.29	0.48
1:B:641:PRO:HG2	1:B:665:GLN:HG2	1.95	0.48
1:B:280:ASP:OD1	1:B:282:ARG:NH1	2.46	0.48
1:B:410:ARG:H	1:B:410:ARG:HD3	1.79	0.48
1:B:388:PHE:HB3	1:B:393:LEU:HD11	1.95	0.47
1:A:227:ARG:HG3	1:A:246:GLU:CD	2.35	0.47
1:B:447:ASN:HB3	1:B:469:ILE:HA	1.95	0.47
1:B:267:ASP:OD1	1:B:267:ASP:C	2.52	0.47
1:B:255:THR:HB	1:B:256:PRO:HD2	1.96	0.47
1:A:509:TYR:O	1:A:520:THR:HA	2.15	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:290:HIS:HE1	1:B:398:ASP:OD2	1.99	0.46
1:A:273:GLN:O	1:A:282:ARG:O	2.32	0.46
1:A:397:VAL:HG12	1:A:619:PRO:HG3	1.97	0.46
1:B:469:ILE:HG12	1:B:470:LYS:H	1.80	0.46
1:A:320:ARG:HG2	1:A:322:VAL:HG22	1.97	0.46
1:A:340:ARG:HD2	1:A:342:PHE:HE2	1.80	0.46
1:A:342:PHE:HE1	1:A:397:VAL:HG13	1.80	0.46
1:B:282:ARG:HD2	1:B:350:ARG:HG2	1.98	0.46
1:B:246:GLU:OE2	1:B:252:LYS:HE3	2.16	0.46
1:B:318:ILE:HG22	1:B:346:ALA:HA	1.97	0.46
1:A:219:ILE:HD11	1:A:546:TYR:C	2.37	0.45
1:A:613:ARG:HH11	1:A:613:ARG:HB3	1.81	0.45
1:A:240:GLN:HB2	1:A:256:PRO:HB3	1.97	0.45
1:A:645:HIS:HD2	1:A:666:CYS:O	2.00	0.45
1:A:300:TYR:CZ	1:A:397:VAL:HG11	2.52	0.45
1:A:222:VAL:HG22	1:B:542:SER:HB3	1.98	0.45
1:A:218:CYS:HB2	1:A:585:PRO:HG3	1.98	0.45
1:A:241:ARG:HG2	1:A:253:ILE:HG23	1.98	0.45
1:B:231:GLY:CA	1:B:565:LEU:HD13	2.43	0.44
1:A:608:CYS:O	1:A:612:CYS:HB3	2.17	0.44
1:B:234:HIS:HD2	1:B:236:GLY:H	1.64	0.44
1:B:400:ASP:O	1:B:622:LYS:HD3	2.17	0.44
1:B:418:THR:HG23	1:B:419:ASN:H	1.82	0.44
1:B:298:LYS:HD2	1:B:619:PRO:HG3	1.99	0.43
1:A:257:GLU:OE2	1:B:539:THR:HG21	2.19	0.43
1:B:497:TYR:OH	1:B:561:GLU:OE2	2.25	0.43
1:A:539:THR:HG21	1:B:257:GLU:CD	2.39	0.43
1:B:359:PHE:CE2	1:B:365:LEU:HB2	2.54	0.43
1:B:541:GLY:O	1:B:543:CYS:N	2.52	0.43
1:A:640:GLU:CB	1:A:641:PRO:CD	2.96	0.43
1:B:502:SER:OG	1:B:505:LEU:HB2	2.19	0.43
1:A:667:GLU:HG2	1:A:668:HIS:CD2	2.53	0.43
1:B:361:PRO:O	1:B:362:ASP:HB2	2.19	0.42
1:B:371:ASP:OD1	1:B:430:HIS:HA	2.20	0.42
1:A:290:HIS:HE1	1:A:398:ASP:OD2	2.03	0.42
1:A:342:PHE:CE1	1:A:397:VAL:HG13	2.54	0.42
1:A:239:SER:C	1:A:241:ARG:H	2.22	0.41
1:A:245:LEU:HD23	1:A:245:LEU:N	2.35	0.41
1:A:497:TYR:CD1	1:A:505:LEU:HB3	2.54	0.41
1:B:410:ARG:H	1:B:410:ARG:NE	2.18	0.41
1:B:417:SER:O	1:B:419:ASN:N	2.53	0.41
1:A:502:SER:OG	1:A:505:LEU:HB2	2.20	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:232:ALA:O	1:B:555:GLY:HA2	2.21	0.41
1:A:613:ARG:HB3	1:A:613:ARG:NH1	2.36	0.41
1:B:352:HIS:HB3	1:B:371:ASP:HB2	2.02	0.41
1:A:408:ILE:HA	1:A:409:PRO:HD3	1.82	0.41
1:B:573:GLN:HG3	1:B:574:THR:HG23	2.03	0.41
1:A:324:TYR:HB3	1:A:334:VAL:HG11	2.02	0.40
1:B:640:GLU:HG2	1:B:653:ASN:OD1	2.21	0.40
1:B:371:ASP:HB3	1:B:373:MSE:H	1.86	0.40
1:B:631:GLY:O	1:B:632:ASP:C	2.60	0.40
1:A:298:LYS:HE3	1:A:617:CYS:O	2.22	0.40
1:B:481:LEU:HD23	1:B:482:GLU:H	1.86	0.40
1:B:483:PHE:CE2	1:B:485:PRO:HG3	2.57	0.40
1:B:299:LEU:N	1:B:299:LEU:HD12	2.36	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	429/481 (89%)	399 (93%)	25 (6%)	5 (1%)	19	62
1	B	436/481 (91%)	395 (91%)	34 (8%)	7 (2%)	14	55
All	All	865/962 (90%)	794 (92%)	59 (7%)	12 (1%)	16	58

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	410	ARG
1	A	575	HIS
1	B	418	THR
1	B	471	GLY
1	A	612	CYS
1	B	542	SER
1	B	614	ASN

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Mol	Chain	Res	Type
1	B	642	ALA
1	A	613	ARG
1	B	417	SER
1	B	641	PRO
1	A	420	GLN

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	380/415 (92%)	341 (90%)	39 (10%)	10	36
1	B	385/415 (93%)	348 (90%)	37 (10%)	12	42
All	All	765/830 (92%)	689 (90%)	76 (10%)	11	39

All (76) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	214	HIS
1	A	219	ILE
1	A	227	ARG
1	A	233	LEU
1	A	241	ARG
1	A	245	LEU
1	A	272	VAL
1	A	287	LEU
1	A	318	ILE
1	A	323	GLU
1	A	326	VAL
1	A	347	GLU
1	A	352	HIS
1	A	358	LEU
1	A	367	ILE
1	A	383	ASP
1	A	399	THR
1	A	403	ASN
1	A	407	SER
1	A	412	ASN

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Mol	Chain	Res	Type
1	A	416	ASN
1	A	419	ASN
1	A	429	LEU
1	A	464	ARG
1	A	475	GLU
1	A	481	LEU
1	A	505	LEU
1	A	513	ASP
1	A	526	VAL
1	A	539	THR
1	A	548	SER
1	A	561	GLU
1	A	575	HIS
1	A	593	GLU
1	A	598	VAL
1	A	610	ARG
1	A	632	ASP
1	A	635	ARG
1	A	656	LEU
1	B	214	HIS
1	B	227	ARG
1	B	233	LEU
1	B	240	GLN
1	B	241	ARG
1	B	245	LEU
1	B	299	LEU
1	B	301	VAL
1	B	326	VAL
1	B	334	VAL
1	B	338	THR
1	B	352	HIS
1	B	358	LEU
1	B	362	ASP
1	B	376	LEU
1	B	383	ASP
1	B	403	ASN
1	B	410	ARG
1	B	412	ASN
1	B	429	LEU
1	B	447	ASN
1	B	481	LEU
1	B	505	LEU

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Mol	Chain	Res	Type
1	B	521	LEU
1	B	526	VAL
1	B	544	ARG
1	B	548	SER
1	B	561	GLU
1	B	567	SER
1	B	593	GLU
1	B	597	THR
1	B	598	VAL
1	B	610	ARG
1	B	613	ARG
1	B	635	ARG
1	B	644	ARG
1	B	656	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	214	HIS
1	A	234	HIS
1	A	290	HIS
1	A	332	HIS
1	A	412	ASN
1	A	416	ASN
1	A	419	ASN
1	A	427	HIS
1	A	602	GLN
1	A	645	HIS
1	A	668	HIS
1	B	234	HIS
1	B	269	HIS
1	B	290	HIS
1	B	332	HIS
1	B	403	ASN
1	B	412	ASN
1	B	419	ASN
1	B	522	GLN
1	B	576	ASN
1	B	602	GLN
1	B	645	HIS

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	437/481 (90%)	-0.01	4 (0%)	81 25	48, 67, 102, 139	8 (1%)
1	B	442/481 (91%)	0.09	6 (1%)	72 17	50, 70, 113, 136	6 (1%)
All	All	879/962 (91%)	0.04	10 (1%)	77 22	48, 68, 109, 139	14 (1%)

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	382	MSE	3.2
1	B	401	MSE	2.8
1	B	382	MSE	2.5
1	B	526	VAL	2.3
1	A	458	LYS	2.3
1	B	522	GLN	2.2
1	B	475	GLU	2.1
1	B	474	TYR	2.1
1	A	348	LEU	2.1
1	A	379	MSE	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.