



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 17, 2026 – 06:13 PM UTC

PDB ID : 3HO5 / pdb_00003ho5
Title : Crystal structure of Hedgehog-interacting protein (HHIP) and Sonic hedgehog (SHH) complex
Authors : Hymowitz, S.G.; Bosanac, I.
Deposited on : 2009-06-01
Resolution : 3.01 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

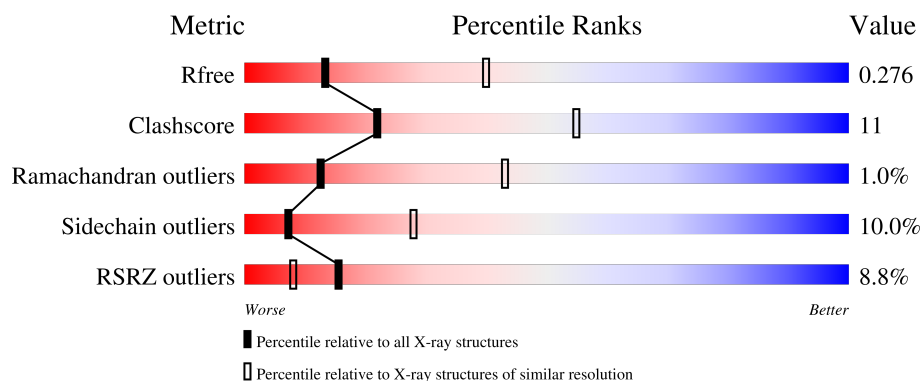
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.01 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}	180053	3131 (3.04-3.00)
Clashscore	190562	3444 (3.04-3.00)
Ramachandran outliers	187476	3319 (3.04-3.00)
Sidechain outliers	187428	3322 (3.04-3.00)
RSRZ outliers	180081	3130 (3.04-3.00)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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 $\leq 5\%$. 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	481	11% 63% 26% 7%
1	B	481	5% 64% 25% 6%
2	H	169	10% 83% 8% 9%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	CA	H	402	-	-	-	X

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 8255 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 Hedgehog-interacting protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	446	Total	C	N	O	S	19	0	0
			3487	2189	623	647	28			
1	B	451	Total	C	N	O	S	19	0	0
			3529	2213	631	657	28			

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

- Molecule 2 is a protein called Sonic hedgehog protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	154	Total	C	N	O	S	0	0	0
			1235	771	220	239	5			

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Zn	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	H	1	Total	Zn	0	0
			1	1		

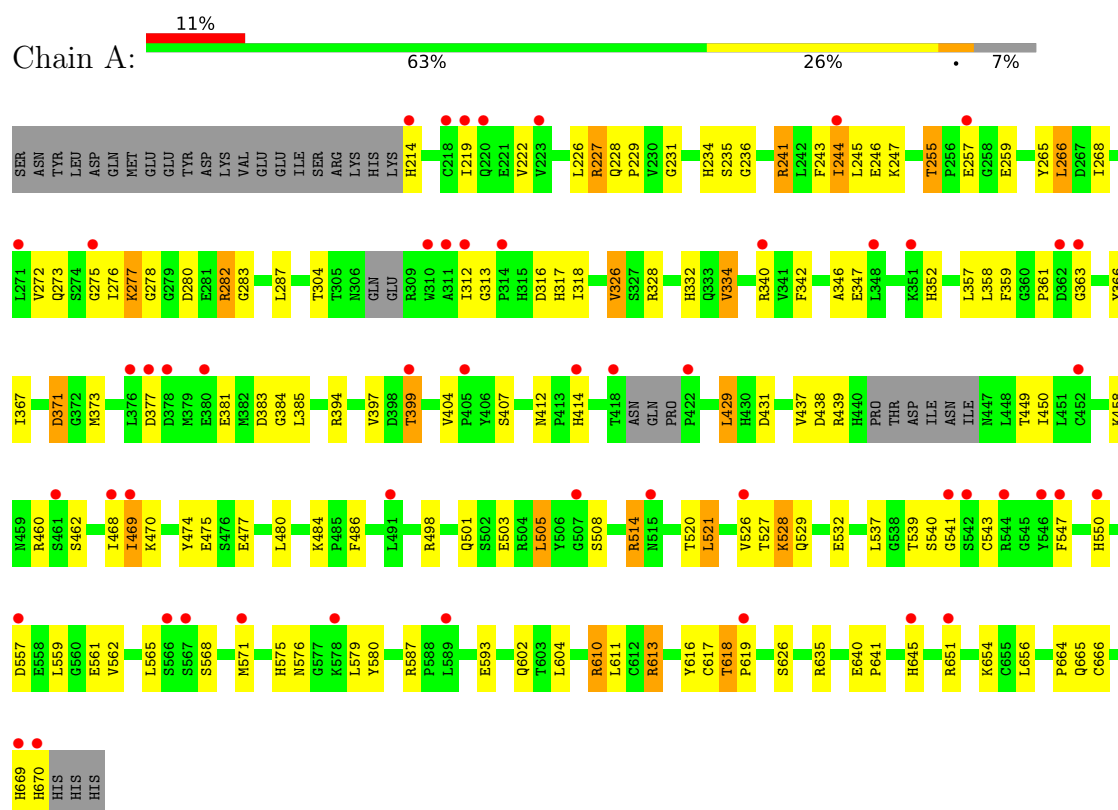
- Molecule 4 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	H	2	Total	Ca	0	0
			2	2		

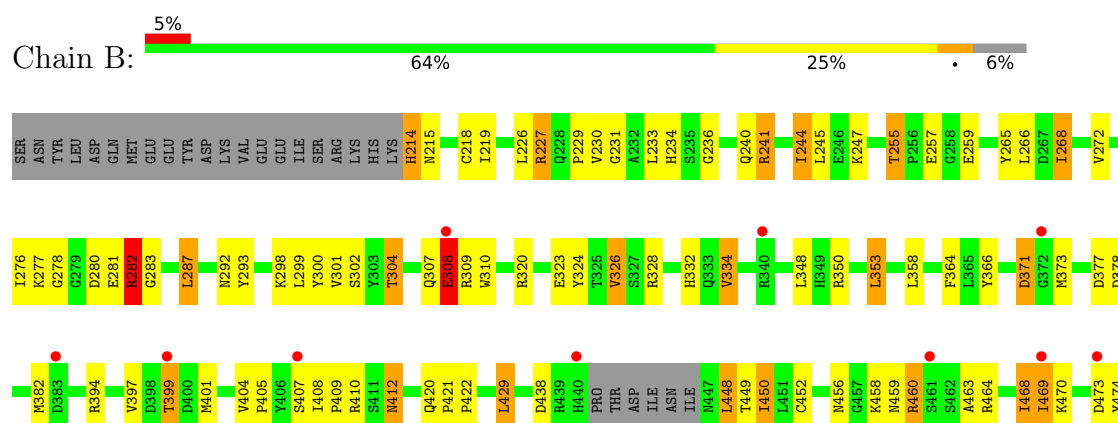
3 Residue-property plots

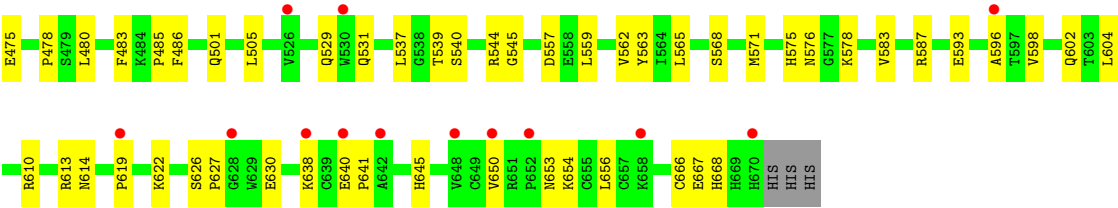
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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

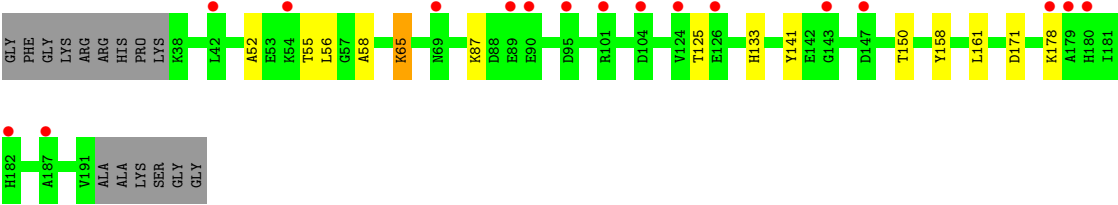
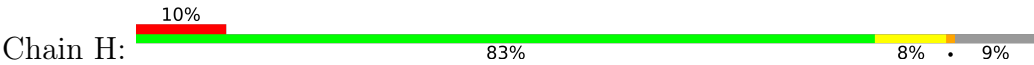


• Molecule 1: Hedgehog-interacting protein





● Molecule 2: Sonic hedgehog protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	101.62Å 101.62Å 302.83Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 – 3.01 30.00 – 3.01	Depositor EDS
% Data completeness (in resolution range)	99.8 (30.00-3.01) 99.7 (30.00-3.01)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.20 (at 3.00Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.231 , 0.287 0.224 , 0.276	Depositor DCC
R_{free} test set	1874 reflections (4.53%)	wwPDB-VP
Wilson B-factor (Å ²)	79.7	Xtriage
Anisotropy	0.589	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 124.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.017 for -h,-k,l	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	8255	wwPDB-VP
Average B, all atoms (Å ²)	96.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.00% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CA, ZN

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	0.55	2/3569 (0.1%)	0.85	4/4816 (0.1%)
1	B	0.58	3/3614 (0.1%)	0.86	3/4882 (0.1%)
2	H	0.42	0/1261	0.75	0/1701
All	All	0.55	5/8444 (0.1%)	0.84	7/11399 (0.1%)

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

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	460	ARG	CA-CB	-13.38	1.30	1.53
1	A	460	ARG	CA-CB	-12.16	1.32	1.53
1	B	458	LYS	CA-CB	-8.40	1.39	1.53
1	B	410	ARG	CA-CB	-7.67	1.39	1.53
1	A	458	LYS	CA-CB	6.61	1.64	1.53

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	460	ARG	N-CA-CB	-11.34	91.33	110.49
1	A	460	ARG	N-CA-CB	7.28	122.80	110.49
1	A	458	LYS	CB-CA-C	-7.25	95.98	110.42
1	B	458	LYS	CB-CA-C	5.91	122.18	110.42
1	A	316	ASP	N-CA-C	-5.71	105.99	113.12

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	528	LYS	N-CA-C	5.56	122.64	110.80
1	B	282	ARG	N-CA-C	5.02	119.03	113.01

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	275	GLY	Peptide

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	3487	0	3388	81	0
1	B	3529	0	3424	90	0
2	H	1235	0	1197	11	0
3	A	1	0	0	0	0
3	H	1	0	0	0	0
4	H	2	0	0	0	0
All	All	8255	0	8009	173	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 11.

All (173) 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:277:LYS:HB2	1:A:280:ASP:HB2	1.51	0.91
1:B:240:GLN:HE22	1:B:596:ALA:HB1	1.38	0.89
1:A:429:LEU:HD13	1:A:480:LEU:HD21	1.55	0.86
1:B:469:ILE:HG22	1:B:470:LYS:H	1.41	0.85
1:A:501:GLN:O	1:A:587:ARG:HD3	1.78	0.83
1:A:645:HIS:HD2	1:A:666:CYS:O	1.63	0.81
1:A:318:ILE:HG22	1:A:346:ALA:HA	1.64	0.80
1:B:429:LEU:HD13	1:B:480:LEU:HD21	1.61	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:527:THR:O	1:A:529:GLN:N	2.15	0.78
1:B:382:MET:HG2	2:H:133:HIS:HB3	1.66	0.78
1:A:469:ILE:HG22	1:A:470:LYS:H	1.49	0.78
1:B:421:PRO:HD3	2:H:178:LYS:HE2	1.66	0.76
1:A:257:GLU:OE2	1:B:539:THR:HG21	1.85	0.76
1:B:501:GLN:O	1:B:587:ARG:HD3	1.86	0.76
1:B:348:LEU:HD11	1:B:382:MET:HE1	1.70	0.74
1:B:568:SER:H	1:B:571:MET:HB2	1.54	0.72
1:A:399:THR:HG23	1:A:407:SER:HB2	1.72	0.71
1:B:397:VAL:HG12	1:B:619:PRO:HG2	1.72	0.71
1:A:414:HIS:HE1	1:A:477:GLU:HG3	1.57	0.69
1:B:463:ALA:HB2	1:B:486:PHE:HE2	1.55	0.69
1:A:257:GLU:CD	1:B:539:THR:HG21	2.18	0.69
1:B:282:ARG:O	1:B:304:THR:HB	1.93	0.69
1:B:667:GLU:HG3	1:B:668:HIS:CD2	2.30	0.67
1:B:378:ASP:HB2	1:B:382:MET:HE2	1.78	0.66
1:A:394:ARG:HE	1:A:412:ASN:HD21	1.45	0.65
1:B:255:THR:HG22	1:B:259:GLU:H	1.60	0.65
1:B:421:PRO:HG3	2:H:178:LYS:HZ1	1.60	0.65
1:B:226:LEU:HD13	1:B:244:ILE:CD1	2.28	0.64
1:B:300:TYR:OH	1:B:397:VAL:HG11	1.97	0.64
1:A:255:THR:HG22	1:A:259:GLU:H	1.60	0.64
1:B:300:TYR:CE2	1:B:323:GLU:HG3	2.34	0.63
1:A:640:GLU:HB2	1:A:641:PRO:HD3	1.79	0.63
1:A:313:GLY:HA3	1:A:384:GLY:HA3	1.81	0.62
1:A:359:PHE:O	1:A:439:ARG:NH1	2.29	0.62
1:A:241:ARG:HH11	1:A:332:HIS:HD2	1.46	0.61
2:H:65:LYS:HG3	2:H:141:TYR:HB3	1.84	0.60
1:A:257:GLU:OE2	1:B:539:THR:CG2	2.51	0.59
1:B:421:PRO:HG3	2:H:178:LYS:NZ	2.17	0.59
1:B:645:HIS:HD2	1:B:666:CYS:O	1.86	0.58
1:B:640:GLU:HB2	1:B:641:PRO:HD3	1.87	0.57
1:A:234:HIS:HD2	1:A:236:GLY:H	1.52	0.56
1:B:408:ILE:HD13	1:B:422:PRO:HB2	1.86	0.56
1:B:348:LEU:CD1	1:B:382:MET:HE1	2.34	0.56
1:A:234:HIS:HD2	1:A:236:GLY:N	2.04	0.56
1:B:226:LEU:HD13	1:B:244:ILE:HD13	1.87	0.55
1:B:244:ILE:HD12	1:B:565:LEU:CD2	2.37	0.54
1:B:568:SER:OG	1:B:571:MET:HE3	2.07	0.54
1:A:468:ILE:O	1:A:474:TYR:OH	2.24	0.54
1:A:340:ARG:HD3	1:A:616:TYR:CG	2.42	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:383:ASP:C	1:A:385:LEU:H	2.16	0.53
1:B:293:TYR:OH	1:B:326:VAL:HG11	2.08	0.53
1:B:469:ILE:HG22	1:B:470:LYS:N	2.20	0.53
1:B:557:ASP:HB3	1:B:559:LEU:H	1.73	0.53
1:A:414:HIS:CE1	1:A:477:GLU:HG3	2.42	0.52
1:B:240:GLN:NE2	1:B:596:ALA:HB1	2.16	0.52
1:A:651:ARG:HD2	1:A:654:LYS:HD3	1.91	0.52
1:A:326:VAL:HA	1:A:334:VAL:HA	1.91	0.52
1:A:571:MET:HB2	1:A:576:ASN:HB3	1.91	0.52
1:B:463:ALA:HB2	1:B:486:PHE:CE2	2.40	0.51
1:B:613:ARG:O	1:B:614:ASN:HB2	2.11	0.51
1:B:630:GLU:HB3	1:B:638:LYS:HE3	1.93	0.51
1:A:244:ILE:HD12	1:A:565:LEU:CD2	2.41	0.51
1:B:247:LYS:HG2	1:B:283:GLY:HA3	1.92	0.51
1:B:282:ARG:CG	1:B:304:THR:HG21	2.41	0.51
1:B:421:PRO:HD3	2:H:178:LYS:CE	2.40	0.51
1:B:571:MET:HG2	1:B:576:ASN:O	2.11	0.51
1:A:539:THR:HB	1:B:257:GLU:OE2	2.11	0.50
1:A:227:ARG:HG2	1:A:575:HIS:HD2	1.77	0.50
1:B:429:LEU:CD1	1:B:480:LEU:HD21	2.39	0.50
1:A:394:ARG:HE	1:A:412:ASN:ND2	2.07	0.50
1:A:429:LEU:CD1	1:A:480:LEU:HD21	2.35	0.50
1:B:282:ARG:HG2	1:B:304:THR:HG21	1.92	0.50
1:A:547:PHE:HB2	1:A:580:TYR:CE1	2.46	0.49
1:B:241:ARG:HH11	1:B:332:HIS:HD2	1.60	0.49
1:A:371:ASP:HB3	1:A:373:MET:H	1.76	0.49
1:B:366:TYR:CZ	1:B:394:ARG:HD2	2.48	0.49
1:A:664:PRO:HB2	1:A:665:GLN:NE2	2.27	0.49
2:H:158:TYR:HA	2:H:161:LEU:HB3	1.94	0.49
1:B:234:HIS:HD2	1:B:236:GLY:H	1.60	0.49
1:B:645:HIS:CD2	1:B:666:CYS:O	2.65	0.48
1:A:231:GLY:HA2	1:A:565:LEU:HD13	1.94	0.48
1:B:302:SER:HA	1:B:320:ARG:O	2.13	0.48
1:B:268:ILE:HD11	1:B:324:TYR:OH	2.14	0.48
1:B:227:ARG:HB3	1:B:575:HIS:CD2	2.48	0.48
1:A:342:PHE:HE1	1:A:397:VAL:HG12	1.78	0.48
1:B:474:TYR:CE2	1:B:478:PRO:HG3	2.48	0.48
1:A:227:ARG:O	1:A:228:GLN:C	2.56	0.48
1:A:282:ARG:HG2	1:A:304:THR:HG21	1.95	0.48
1:A:514:ARG:HG2	1:A:550:HIS:HB3	1.94	0.48
1:B:241:ARG:HH11	1:B:332:HIS:CD2	2.31	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:371:ASP:HB3	1:B:373:MET:H	1.78	0.47
1:A:247:LYS:HG2	1:A:283:GLY:HA3	1.96	0.47
1:A:486:PHE:CE1	1:A:521:LEU:HD21	2.49	0.47
1:A:618:THR:HB	1:A:619:PRO:HD2	1.96	0.47
1:A:282:ARG:CG	1:A:304:THR:HG21	2.44	0.47
1:B:529:GLN:HG2	1:B:531:GLN:HG2	1.96	0.47
1:A:438:ASP:HB3	1:A:449:THR:HB	1.96	0.47
1:A:568:SER:OG	1:A:571:MET:HG2	2.15	0.47
1:A:469:ILE:HG22	1:A:470:LYS:N	2.24	0.47
1:B:241:ARG:NH1	1:B:332:HIS:CD2	2.83	0.46
1:A:282:ARG:O	1:A:304:THR:HB	2.14	0.46
1:A:645:HIS:CD2	1:A:666:CYS:O	2.54	0.46
1:B:307:GLN:O	1:B:310:TRP:N	2.48	0.46
1:A:352:HIS:CD2	1:A:431:ASP:HB2	2.51	0.46
1:A:241:ARG:NH1	1:A:332:HIS:HD2	2.13	0.46
1:B:399:THR:HG22	1:B:401:MET:H	1.80	0.46
1:B:420:GLN:HB3	1:B:421:PRO:HD2	1.98	0.46
1:B:448:LEU:HB3	1:B:468:ILE:HG12	1.98	0.46
1:B:409:PRO:HG2	1:B:412:ASN:CG	2.41	0.45
1:A:361:PRO:C	1:A:363:GLY:H	2.24	0.45
1:B:231:GLY:HA2	1:B:565:LEU:HD13	1.97	0.45
1:B:282:ARG:HD2	1:B:350:ARG:HG2	1.97	0.45
1:B:450:ILE:HD11	1:B:452:CYS:SG	2.56	0.45
2:H:52:ALA:HB3	2:H:55:THR:HG23	1.97	0.45
1:B:276:ILE:HG13	1:B:277:LYS:H	1.82	0.45
1:A:610:ARG:HG3	1:A:611:LEU:HD13	1.99	0.45
1:B:218:CYS:HB3	1:B:583:VAL:HB	1.99	0.45
1:B:255:THR:HG23	1:B:257:GLU:H	1.81	0.45
1:B:214:HIS:HB2	1:B:215:ASN:H	1.55	0.45
1:B:640:GLU:HB2	1:B:653:ASN:HD21	1.82	0.45
1:B:280:ASP:OD2	1:B:282:ARG:NH1	2.50	0.45
1:A:227:ARG:HG3	1:A:246:GLU:OE1	2.17	0.44
1:B:307:GLN:O	1:B:308:GLU:C	2.60	0.44
1:B:300:TYR:CZ	1:B:397:VAL:HG11	2.51	0.44
1:A:255:THR:HG23	1:A:257:GLU:H	1.82	0.44
1:B:244:ILE:HD12	1:B:565:LEU:HD23	1.99	0.44
1:A:394:ARG:HH21	1:A:412:ASN:HD22	1.65	0.44
2:H:87:LYS:HD2	2:H:125:THR:HG22	1.98	0.44
1:B:281:GLU:HB2	1:B:353:LEU:HD22	1.99	0.44
1:B:229:PRO:HB2	1:B:565:LEU:HB3	1.99	0.44
1:B:438:ASP:HB3	1:B:449:THR:HB	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:456:ASN:HB2	1:B:459:ASN:HB3	2.00	0.44
1:B:298:LYS:HD2	1:B:619:PRO:HG3	1.99	0.43
1:A:226:LEU:HD13	1:A:244:ILE:CD1	2.48	0.43
1:A:304:THR:HG23	1:A:317:HIS:ND1	2.34	0.43
1:A:669:HIS:O	1:A:670:HIS:HB3	2.18	0.43
1:A:273:GLN:O	1:A:282:ARG:O	2.36	0.43
1:B:562:VAL:HG12	1:B:563:TYR:N	2.32	0.43
1:A:358:LEU:HG	1:A:366:TYR:HB2	2.01	0.43
1:B:326:VAL:HA	1:B:334:VAL:HA	2.00	0.43
1:B:568:SER:N	1:B:571:MET:HB2	2.29	0.43
1:A:265:TYR:O	1:A:334:VAL:HG13	2.18	0.43
1:A:527:THR:C	1:A:529:GLN:H	2.23	0.43
1:A:278:GLY:C	1:A:280:ASP:H	2.27	0.43
1:A:503:GLU:H	1:A:587:ARG:HD2	1.84	0.43
1:A:520:THR:O	1:A:532:GLU:HA	2.19	0.42
2:H:58:ALA:O	2:H:171:ASP:HB3	2.19	0.42
1:B:348:LEU:HD11	1:B:382:MET:CE	2.46	0.42
1:B:233:LEU:HD21	1:B:287:LEU:HB3	2.00	0.42
1:B:404:VAL:CG1	1:B:405:PRO:HD2	2.49	0.42
1:B:483:PHE:CE2	1:B:485:PRO:HG3	2.55	0.42
1:A:277:LYS:HB2	1:A:280:ASP:CB	2.36	0.42
1:A:541:GLY:C	1:A:543:CYS:H	2.27	0.42
1:B:227:ARG:HB3	1:B:575:HIS:HD2	1.85	0.42
2:H:150:THR:HG23	2:H:161:LEU:HD22	2.02	0.41
1:A:235:SER:HB2	1:A:243:PHE:HE1	1.85	0.41
1:A:377:ASP:O	1:A:381:GLU:HB2	2.20	0.41
1:B:429:LEU:HD12	1:B:464:ARG:HD2	2.03	0.41
1:A:229:PRO:HB2	1:A:565:LEU:HB3	2.02	0.41
1:A:358:LEU:HD22	1:A:437:VAL:HG21	2.03	0.41
1:A:222:VAL:HB	1:A:579:LEU:HD23	2.03	0.41
1:B:307:GLN:CD	1:B:309:ARG:H	2.29	0.41
1:A:366:TYR:CE2	1:A:394:ARG:HD3	2.56	0.41
1:A:462:SER:HB3	1:A:484:LYS:HG2	2.03	0.41
1:B:539:THR:HB	1:B:545:GLY:H	1.86	0.41
1:A:640:GLU:HB2	1:A:641:PRO:CD	2.50	0.41
1:B:265:TYR:O	1:B:334:VAL:HG13	2.21	0.41
1:A:357:LEU:HD22	1:A:367:ILE:HG13	2.03	0.40
1:B:364:PHE:CZ	1:B:409:PRO:HB3	2.55	0.40
1:A:505:LEU:HD22	1:A:508:SER:HB2	2.02	0.40
1:A:266:LEU:HD13	1:A:268:ILE:HB	2.03	0.40
1:A:557:ASP:HB2	1:A:561:GLU:H	1.87	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:557:ASP:HB3	1:A:559:LEU:H	1.85	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	438/481 (91%)	412 (94%)	21 (5%)	5 (1%)	11	41
1	B	447/481 (93%)	412 (92%)	30 (7%)	5 (1%)	11	41
2	H	152/169 (90%)	147 (97%)	5 (3%)	0	100	100
All	All	1037/1131 (92%)	971 (94%)	56 (5%)	10 (1%)	12	43

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	276	ILE
1	A	277	LYS
1	A	469	ILE
1	A	528	LYS
1	A	613	ARG
1	B	460	ARG
1	B	469	ILE
1	B	308	GLU
1	B	627	PRO
1	B	278	GLY

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	387/422 (92%)	347 (90%)	40 (10%)	7	27
1	B	392/422 (93%)	343 (88%)	49 (12%)	4	19
2	H	132/141 (94%)	130 (98%)	2 (2%)	57	79
All	All	911/985 (92%)	820 (90%)	91 (10%)	7	28

All (91) 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	241	ARG
1	A	244	ILE
1	A	245	LEU
1	A	255	THR
1	A	266	LEU
1	A	272	VAL
1	A	282	ARG
1	A	287	LEU
1	A	312	ILE
1	A	326	VAL
1	A	328	ARG
1	A	334	VAL
1	A	347	GLU
1	A	371	ASP
1	A	399	THR
1	A	404	VAL
1	A	429	LEU
1	A	450	ILE
1	A	475	GLU
1	A	498	ARG
1	A	505	LEU
1	A	514	ARG
1	A	521	LEU
1	A	526	VAL
1	A	537	LEU
1	A	540	SER
1	A	562	VAL

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Mol	Chain	Res	Type
1	A	593	GLU
1	A	602	GLN
1	A	604	LEU
1	A	610	ARG
1	A	613	ARG
1	A	617	CYS
1	A	618	THR
1	A	626	SER
1	A	635	ARG
1	A	656	LEU
1	B	214	HIS
1	B	219	ILE
1	B	227	ARG
1	B	230	VAL
1	B	241	ARG
1	B	244	ILE
1	B	245	LEU
1	B	255	THR
1	B	266	LEU
1	B	268	ILE
1	B	272	VAL
1	B	282	ARG
1	B	287	LEU
1	B	292	ASN
1	B	299	LEU
1	B	301	VAL
1	B	304	THR
1	B	308	GLU
1	B	326	VAL
1	B	328	ARG
1	B	334	VAL
1	B	353	LEU
1	B	358	LEU
1	B	371	ASP
1	B	377	ASP
1	B	399	THR
1	B	407	SER
1	B	412	ASN
1	B	429	LEU
1	B	448	LEU
1	B	450	ILE
1	B	468	ILE

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Mol	Chain	Res	Type
1	B	473	ASP
1	B	475	GLU
1	B	505	LEU
1	B	537	LEU
1	B	540	SER
1	B	544	ARG
1	B	578	LYS
1	B	593	GLU
1	B	598	VAL
1	B	602	GLN
1	B	604	LEU
1	B	610	ARG
1	B	622	LYS
1	B	626	SER
1	B	650	VAL
1	B	654	LYS
1	B	656	LEU
2	H	56	LEU
2	H	65	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (32) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	234	HIS
1	A	292	ASN
1	A	332	HIS
1	A	333	GLN
1	A	412	ASN
1	A	414	HIS
1	A	430	HIS
1	A	440	HIS
1	A	447	ASN
1	A	459	ASN
1	A	515	ASN
1	A	550	HIS
1	A	575	HIS
1	A	599	GLN
1	A	645	HIS
1	A	665	GLN
1	B	214	HIS
1	B	234	HIS
1	B	240	GLN

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Mol	Chain	Res	Type
1	B	332	HIS
1	B	412	ASN
1	B	414	HIS
1	B	440	HIS
1	B	467	GLN
1	B	575	HIS
1	B	614	ASN
1	B	645	HIS
1	B	665	GLN
1	B	668	HIS
2	H	69	ASN
2	H	79	ASN
2	H	91	ASN

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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

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	446/481 (92%)	0.95	52 (11%) 9 5	37, 98, 135, 168	4 (0%)
1	B	451/481 (93%)	0.58	23 (5%) 33 17	37, 94, 128, 154	4 (0%)
2	H	154/169 (91%)	0.78	17 (11%) 10 6	62, 78, 92, 103	0
All	All	1051/1131 (92%)	0.77	92 (8%) 15 8	37, 94, 131, 168	8 (0%)

All (92) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	380	GLU	4.1
1	A	311	ALA	4.0
1	A	348	LEU	4.0
1	A	310	TRP	3.8
2	H	95	ASP	3.5
1	B	383	ASP	3.4
2	H	89	GLU	3.4
1	A	314	PRO	3.4
1	A	362	ASP	3.3
2	H	179	ALA	3.2
2	H	69	ASN	3.2
2	H	147	ASP	3.0
1	B	526	VAL	3.0
2	H	178	LYS	3.0
1	A	376	LEU	3.0
1	A	461	SER	3.0
1	B	658	LYS	3.0
1	B	638	LYS	2.9
2	H	101	ARG	2.9
1	A	422	PRO	2.9
1	B	619	PRO	2.9
1	B	340	ARG	2.9
1	A	468	ILE	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	469	ILE	2.8
1	B	640	GLU	2.8
1	B	399	THR	2.8
1	A	271	LEU	2.8
2	H	182	HIS	2.8
1	A	340	ARG	2.8
1	A	541	GLY	2.7
1	A	619	PRO	2.7
1	A	363	GLY	2.7
1	B	461	SER	2.7
1	B	642	ALA	2.7
1	A	312	ILE	2.6
1	A	257	GLU	2.6
1	B	652	PRO	2.6
1	A	275	GLY	2.6
1	B	628	GLY	2.6
1	A	491	LEU	2.6
1	B	648	VAL	2.6
1	A	571	MET	2.6
1	A	507	GLY	2.6
1	A	547	PHE	2.6
1	A	542	SER	2.5
1	A	351	LYS	2.5
1	A	399	THR	2.5
1	A	515	ASN	2.5
1	A	378	ASP	2.4
1	A	546	TYR	2.4
1	A	219	ILE	2.4
1	B	440	HIS	2.4
1	B	308	GLU	2.4
1	A	377	ASP	2.4
1	A	223	VAL	2.4
1	B	670	HIS	2.3
1	A	651	ARG	2.3
2	H	104	ASP	2.3
1	A	452	CYS	2.3
2	H	90	GLU	2.3
1	B	469	ILE	2.3
1	A	670	HIS	2.3
1	A	220	GLN	2.3
1	B	650	VAL	2.3
1	A	405	PRO	2.3

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Mol	Chain	Res	Type	RSRZ
2	H	42	LEU	2.3
1	B	596	ALA	2.2
1	A	414	HIS	2.2
1	A	244	ILE	2.2
1	A	544	ARG	2.2
1	B	530	TRP	2.2
2	H	180	HIS	2.2
1	B	372	GLY	2.2
1	A	589	LEU	2.1
1	A	218	CYS	2.1
2	H	124	VAL	2.1
1	A	550	HIS	2.1
2	H	54	LYS	2.1
1	A	645	HIS	2.1
2	H	143	GLY	2.1
2	H	126	GLU	2.1
1	B	407	SER	2.1
1	A	669	HIS	2.1
1	A	566	SER	2.0
1	A	567	SER	2.0
1	A	214	HIS	2.0
1	A	578	LYS	2.0
1	A	418	THR	2.0
2	H	187	ALA	2.0
1	A	557	ASP	2.0
1	B	473	ASP	2.0
1	A	526	VAL	2.0

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 oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,

median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	CA	H	402	1/1	0.79	0.44	91,91,91,91	0
3	ZN	A	902	1/1	0.85	0.17	151,151,151,151	0
4	CA	H	401	1/1	0.97	0.46	90,90,90,90	0
3	ZN	H	400	1/1	0.98	0.34	75,75,75,75	0

6.5 Other polymers [i](#)

There are no such residues in this entry.