



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 09:14 AM GMT

PDB ID : 3HO5  
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](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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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 : 1.7 (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

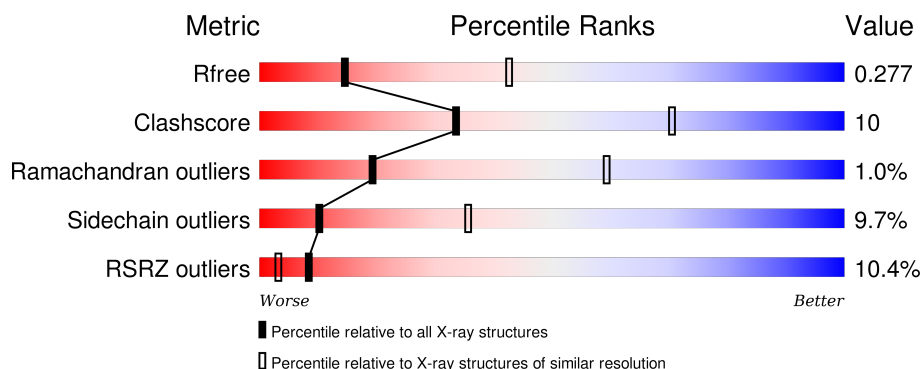
# 1 Overall quality at a glance ⓘ

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))
$R_{free}$	91344	1773 (3.04-3.00)
Clashscore	102246	2117 (3.04-3.00)
Ramachandran outliers	100387	2050 (3.04-3.00)
Sidechain outliers	100360	2053 (3.04-3.00)
RSRZ outliers	91569	1788 (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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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  $\leq 5\%$ . 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	<div> <div>8%</div> <div>64%</div> <div>24%</div> <div>•</div> <div>7%</div> </div>
1	B	481	<div> <div>5%</div> <div>64%</div> <div>25%</div> <div>5%</div> <div>6%</div> </div>
2	H	169	<div> <div>26%</div> <div>82%</div> <div>9%</div> <div>9%</div> </div>

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	401	-	-	-	X
4	CA	H	402	-	-	-	X

## 2 Entry composition

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 (three-letter code: ZN) (formula: Zn).

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

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

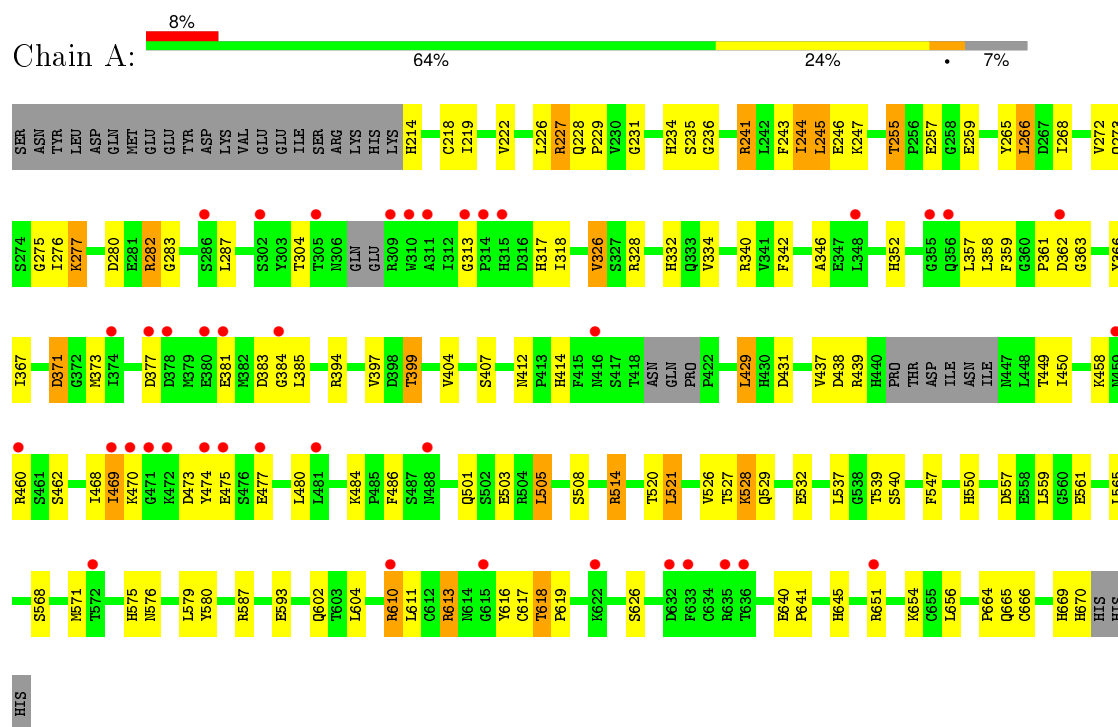
- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

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

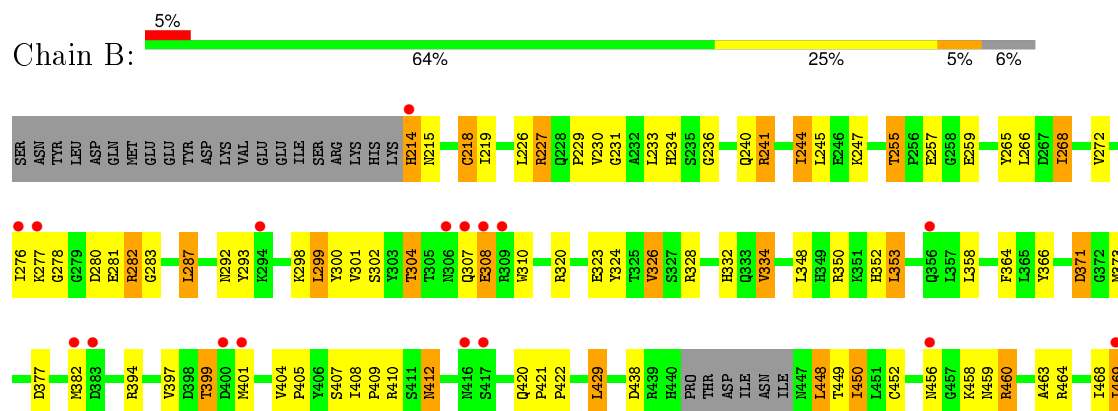
### 3 Residue-property plots

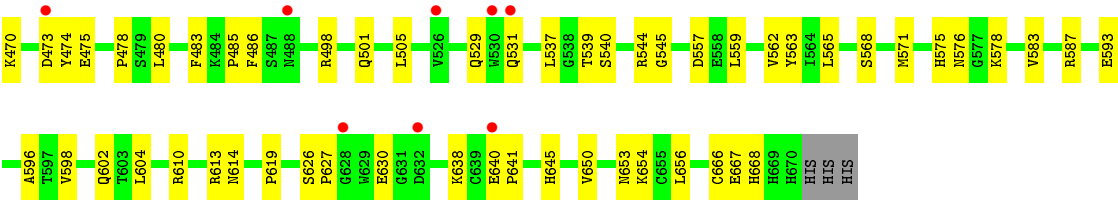
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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

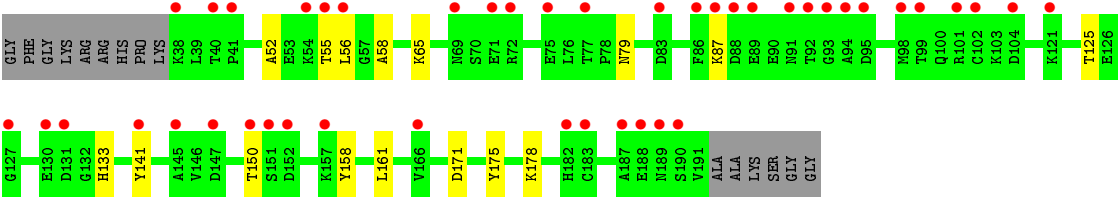
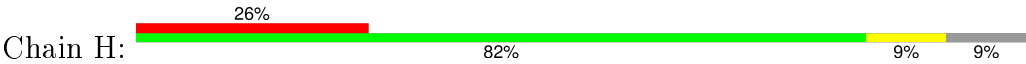


#### • Molecule 1: Hedgehog-interacting protein





● 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, $\alpha$ , $\beta$ , $\gamma$	101.62Å 101.62Å 302.83Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 – 3.01 29.33 – 3.01	Depositor EDS
% Data completeness (in resolution range)	99.8 (30.00-3.01) 99.8 (29.33-3.01)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.09	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.20 (at 3.00Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.231 , 0.287 0.224 , 0.277	Depositor DCC
$R_{free}$ test set	1864 reflections (5.32%)	DCC
Wilson B-factor (Å <sup>2</sup> )	79.7	Xtriage
Anisotropy	0.589	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 120.4	EDS
Estimated twinning fraction	0.017 for -h,-k,l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.35$	Xtriage
Outliers	1 of 37023 reflections (0.003%)	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	8255	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	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.*

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

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  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

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

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	0.45	3/3569 (0.1%)	0.63	3/4816 (0.1%)
1	B	0.47	4/3614 (0.1%)	0.64	3/4882 (0.1%)
2	H	0.34	0/1261	0.48	0/1701
All	All	0.44	7/8444 (0.1%)	0.62	6/11399 (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 (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	460	ARG	CA-CB	-10.50	1.30	1.53
1	A	460	ARG	CA-CB	-9.56	1.32	1.53
1	B	458	LYS	CA-CB	-6.67	1.39	1.53
1	B	410	ARG	CA-CB	-6.46	1.39	1.53
1	B	218	CYS	CB-SG	-6.38	1.71	1.82
1	A	528	LYS	CE-NZ	5.97	1.64	1.49
1	A	218	CYS	CB-SG	-5.73	1.72	1.81

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	460	ARG	N-CA-CB	-10.71	91.33	110.60
1	B	299	LEU	CA-CB-CG	7.91	133.50	115.30
1	A	458	LYS	CB-CA-C	-7.21	95.98	110.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	460	ARG	N-CA-CB	6.78	122.80	110.60
1	B	458	LYS	CB-CA-C	5.89	122.18	110.40
1	A	245	LEU	CA-CB-CG	5.34	127.58	115.30

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	78	0
1	B	3529	0	3424	85	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	165	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 10.

All (165) 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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:429:LEU:HD13	1:B:480:LEU:HD21	1.61	0.80
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:B:501:GLN:O	1:B:587:ARG:HD3	1.86	0.76
1:A:257:GLU:OE2	1:B:539:THR:HG21	1.85	0.76
1:B:568:SER:H	1:B:571:MET:HB2	1.54	0.72
1:B:397:VAL:HG12	1:B:619:PRO:HG2	1.72	0.71
1:A:399:THR:HG23	1:A:407:SER:HB2	1.72	0.71
1:B:463:ALA:HB2	1:B:486:PHE:HE2	1.55	0.69
1:A:414:HIS:HE1	1:A:477:GLU:HG3	1.57	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: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:A:257:GLU:CD	1:B:539:THR:HG21	2.18	0.64
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:B:421:PRO:HG3	2:H:178:LYS:NZ	2.17	0.59
1:A:257:GLU:OE2	1:B:539:THR:CG2	2.51	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: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: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
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:240:GLN:NE2	1:B:596:ALA:HB1	2.16	0.52
1:A:414:HIS:CE1	1:A:477:GLU:HG3	2.42	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:348:LEU:HD11	1:B:382:MET:HE1	1.92	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:421:PRO:HD3	2:H:178:LYS:CE	2.40	0.51
1:B:282:ARG:CG	1:B:304:THR:HG21	2.41	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
2:H:158:TYR:HA	2:H:161:LEU:HB3	1.94	0.49
1:A:664:PRO:HB2	1:A:665:GLN:NE2	2.27	0.49
1:A:383:ASP:C	1:A:385:LEU:H	2.16	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:282:ARG:HG2	1:A:304:THR:HG21	1.95	0.48
1:B:241:ARG:HH11	1:B:332:HIS:CD2	2.31	0.48
1:A:514:ARG:HG2	1:A:550:HIS:HB3	1.94	0.48
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:348:LEU:HD11	1:B:382:MET:CE	2.46	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:399:THR:HG22	1:B:401:MET:H	1.80	0.46
1:B:282:ARG:HD2	1:B:350:ARG:HG2	1.97	0.45
1:B:231:GLY:HA2	1:B:565:LEU:HD13	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:255:THR:HG23	1:B:257:GLU:H	1.81	0.45
1:B:218:CYS:HB3	1:B:583:VAL:HB	1.99	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: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
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:B:568:SER:N	1:B:571:MET:HB2	2.29	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:A:227:ARG:O	1:A:228:GLN:C	2.56	0.43
1:A:265:TYR:O	1:A:334:VAL:HG13	2.18	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: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: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:B:429:LEU:HD12	1:B:464:ARG:HD2	2.03	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: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:B:409:PRO:HG2	1:B:412:ASN:CG	2.41	0.41
1:A:361:PRO:C	1:A:363:GLY:H	2.24	0.41
1:A:222:VAL:HB	1:A:579:LEU:HD23	2.03	0.41
1:B:539:THR:HB	1:B:545:GLY:H	1.86	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: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:B:364:PHE:CZ	1:B:409:PRO:HB3	2.55	0.40
1:A:357:LEU:HD22	1:A:367:ILE:HG13	2.03	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
1:A:557:ASP:HB3	1:A:559:LEU:H	1.85	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	438/481 (91%)	412 (94%)	21 (5%)	5 (1%)	17	56
1	B	447/481 (93%)	412 (92%)	30 (7%)	5 (1%)	17	56
2	H	152/169 (90%)	147 (97%)	5 (3%)	0	100	100
All	All	1037/1131 (92%)	971 (94%)	56 (5%)	10 (1%)	19	59

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

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%)	351 (91%)	36 (9%)	11	38
1	B	392/422 (93%)	343 (88%)	49 (12%)	6	23

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	H	132/141 (94%)	129 (98%)	3 (2%)	58	87
All	All	911/985 (92%)	823 (90%)	88 (10%)	10	36

All (88) 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	326	VAL
1	A	328	ARG
1	A	362	ASP
1	A	371	ASP
1	A	399	THR
1	A	404	VAL
1	A	429	LEU
1	A	450	ILE
1	A	473	ASP
1	A	475	GLU
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	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	656	LEU

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Mol	Chain	Res	Type
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	352	HIS
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	473	ASP
1	B	475	GLU
1	B	498	ARG
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

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Mol	Chain	Res	Type
1	B	602	GLN
1	B	604	LEU
1	B	610	ARG
1	B	626	SER
1	B	650	VAL
1	B	654	LYS
1	B	656	LEU
2	H	56	LEU
2	H	79	ASN
2	H	175	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	234	HIS
1	A	332	HIS
1	A	412	ASN
1	A	414	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	645	HIS
1	A	665	GLN
1	B	214	HIS
1	B	234	HIS
1	B	240	GLN
1	B	315	HIS
1	B	332	HIS
1	B	412	ASN
1	B	414	HIS
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 carbohydrates 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, 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	446/481 (92%)	0.52	40 (8%)	12 4	67, 98, 136, 168	4 (0%)
1	B	451/481 (93%)	0.28	25 (5%)	29 11	69, 94, 128, 154	4 (0%)
2	H	154/169 (91%)	1.40	44 (28%)	1 0	62, 78, 92, 103	0
All	All	1051/1131 (92%)	0.54	109 (10%)	8 3	62, 94, 131, 168	8 (0%)

All (109) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	314	PRO	10.2
2	H	56	LEU	9.9
1	A	311	ALA	6.3
1	B	526	VAL	6.0
1	A	310	TRP	6.0
1	A	377	ASP	5.8
2	H	89	GLU	5.5
2	H	151	SER	5.0
1	A	378	ASP	4.9
2	H	55	THR	4.7
1	A	380	GLU	4.6
1	B	276	ILE	4.6
2	H	69	ASN	4.5
1	A	362	ASP	4.4
2	H	92	THR	4.4
1	A	632	ASP	4.3
2	H	152	ASP	4.3
2	H	121	LYS	4.3
1	A	459	ASN	4.1
1	B	530	TRP	4.1
2	H	147	ASP	4.0
1	B	307	GLN	4.0
2	H	101	ARG	3.9

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Mol	Chain	Res	Type	RSRZ
2	H	87	LYS	3.9
2	H	93	GLY	3.8
1	A	313	GLY	3.7
2	H	183	CYS	3.6
2	H	54	LYS	3.6
1	A	348	LEU	3.5
2	H	190	SER	3.5
1	A	471	GLY	3.5
2	H	102	CYS	3.5
1	B	416	ASN	3.5
1	A	460	ARG	3.5
1	A	633	PHE	3.5
1	A	286	SER	3.4
1	B	640	GLU	3.4
2	H	94	ALA	3.4
1	A	356	GLN	3.4
1	A	636	THR	3.4
1	A	381	GLU	3.3
2	H	145	ALA	3.3
1	B	473	ASP	3.2
2	H	72	ARG	3.1
1	B	456	ASN	3.0
1	A	488	ASN	3.0
1	A	309	ARG	3.0
1	B	214	HIS	3.0
2	H	157	LYS	2.9
2	H	189	ASN	2.8
1	B	356	GLN	2.8
2	H	104	ASP	2.8
1	B	308	GLU	2.8
2	H	40	THR	2.8
1	B	469	ILE	2.8
2	H	141	TYR	2.8
1	B	632	ASP	2.7
2	H	187	ALA	2.7
1	B	306	ASN	2.7
1	B	383	ASP	2.7
1	A	481	LEU	2.7
1	A	469	ILE	2.7
1	A	610	ARG	2.7
1	B	401	MET	2.6
1	A	635	ARG	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	277	LYS	2.6
2	H	130	GLU	2.5
1	A	355	GLY	2.5
1	B	309	ARG	2.5
1	A	475	GLU	2.5
1	A	472	LYS	2.5
1	A	384	GLY	2.5
2	H	75	GLU	2.5
1	A	615	GLY	2.5
1	B	628	GLY	2.5
2	H	38	LYS	2.5
2	H	83	ASP	2.5
1	A	416	ASN	2.5
2	H	91	ASN	2.4
1	B	417	SER	2.4
2	H	150	THR	2.4
2	H	41	PRO	2.4
1	A	651	ARG	2.3
2	H	98	MET	2.3
2	H	131	ASP	2.3
1	A	572	THR	2.3
1	B	400	ASP	2.3
2	H	86	PHE	2.3
1	A	477	GLU	2.2
1	B	382	MET	2.2
2	H	77	THR	2.2
2	H	99	THR	2.2
2	H	166	VAL	2.2
2	H	182	HIS	2.2
2	H	95	ASP	2.1
1	B	531	GLN	2.1
2	H	188	GLU	2.1
1	A	374	ILE	2.1
1	B	488	ASN	2.1
1	A	622	LYS	2.1
1	B	294	LYS	2.1
1	A	474	TYR	2.1
1	A	302	SER	2.1
1	A	305	THR	2.1
2	H	71	GLU	2.0
2	H	127	GLY	2.0
1	A	315	HIS	2.0

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Mol	Chain	Res	Type	RSRZ
2	H	88	ASP	2.0
1	A	470	LYS	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
4	CA	H	401	1/1	0.92	0.53	0.42	90,90,90,90	0
4	CA	H	402	1/1	0.55	0.42	0.03	91,91,91,91	0
3	ZN	H	400	1/1	0.89	0.28	-1.80	75,75,75,75	0
3	ZN	A	902	1/1	0.76	0.23	-	151,151,151,151	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.