



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2014 – 07:02 AM GMT

PDB ID : 1HQE  
Title : HUMAN IMMUNODEFICIENCY VIRUS TYPE 1  
Authors : Ding, J.; Hsiou, Y.; Arnold, E.  
Deposited on : 2000-12-15  
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](mailto:validation@mail.wwpdb.org)  
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

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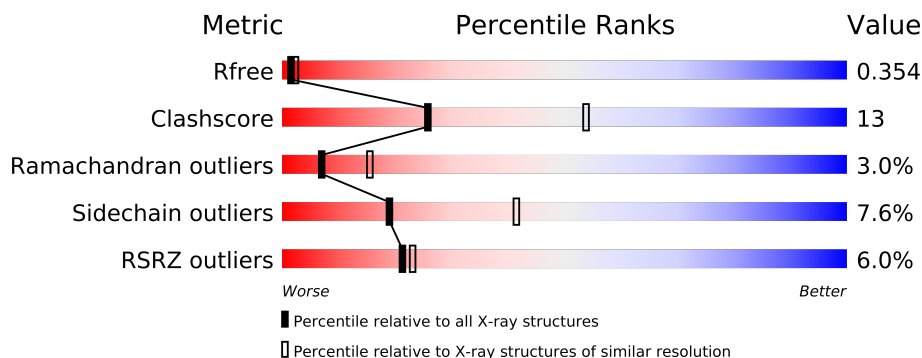
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.70 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  $\geq 3$ , 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	560	
2	B	430	

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7832 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 POL POLYPROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	554	Total	C	N	O	S	0	0	0
			4349	2820	721	802	6			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	103	ASN	LYS	ENGINEERED	UNP P03366
A	280	SER	CYS	ENGINEERED	UNP P03366

- Molecule 2 is a protein called POL POLYPROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	415	Total	C	N	O	S	34	0	0
			3301	2142	547	607	5			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	103	ASN	LYS	ENGINEERED	UNP P03366
B	280	SER	CYS	ENGINEERED	UNP P03366

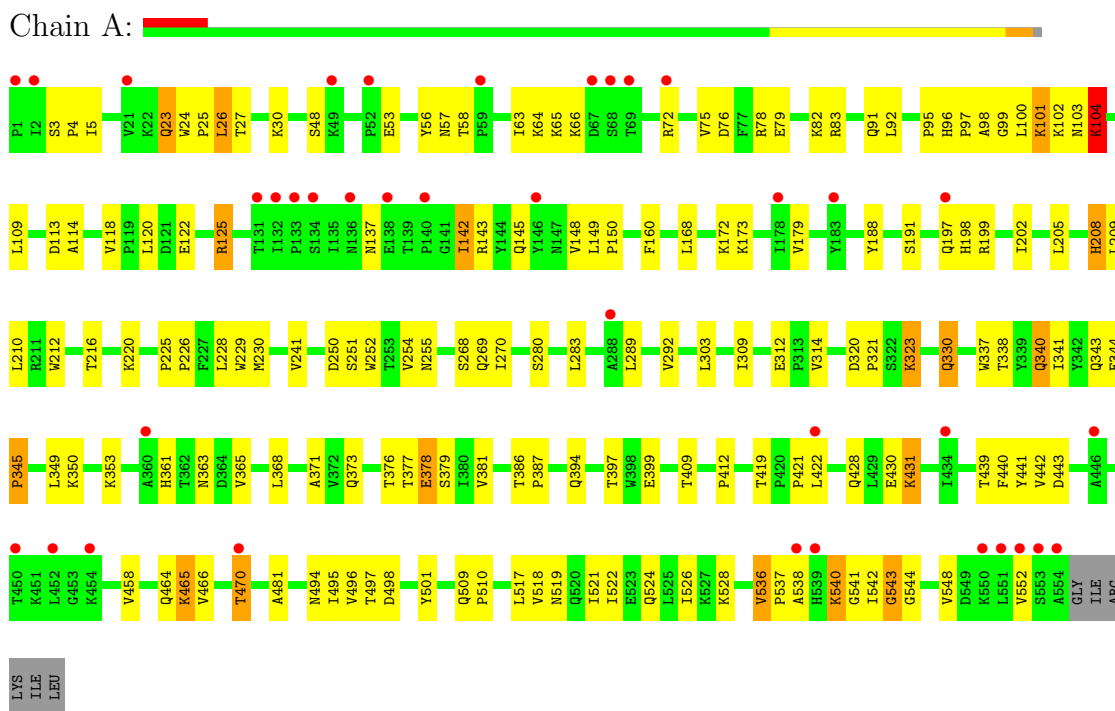
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	78	Total	O	0	0
			78	78		
3	B	104	Total	O	0	0
			104	104		

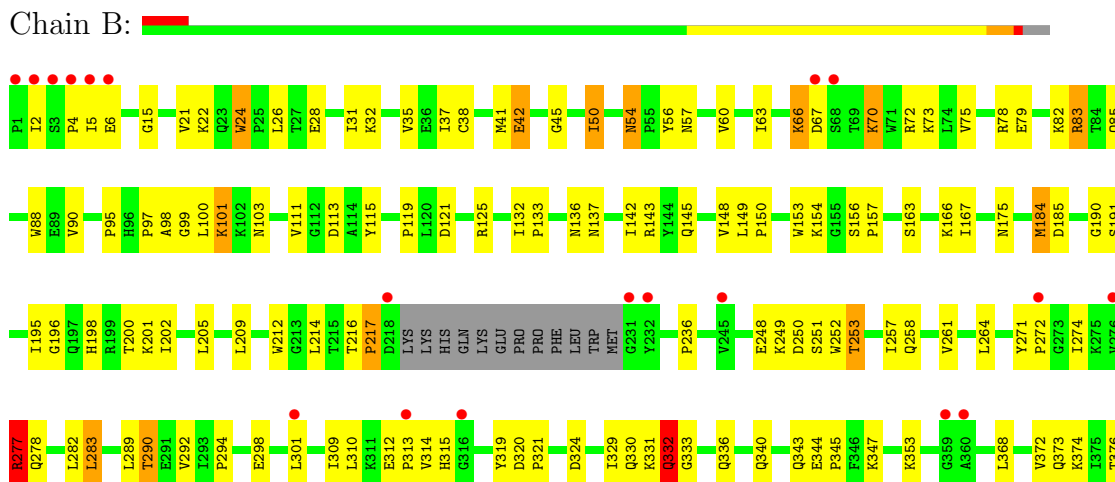
### 3 Residue-property plots

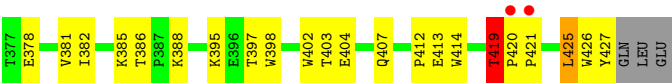
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according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



#### • Molecule 2: POL POLYPROTEIN





## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	236.10Å 70.40Å 93.40Å 90.00° 106.20° 90.00°	Depositor
Resolution (Å)	25.00 – 2.70 28.34 – 2.70	Depositor EDS
% Data completeness (in resolution range)	88.4 (25.00-2.70) 88.4 (28.34-2.70)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.98 (at 2.68Å)	Xtriage
Refinement program	X-PLOR 3.843	Depositor
R, $R_{free}$	0.250 , 0.331 0.304 , 0.354	Depositor DCC
$R_{free}$ test set	1440 reflections (4.16%)	DCC
Wilson B-factor (Å <sup>2</sup> )	49.5	Xtriage
Anisotropy	0.099	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.25 , 34.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.29$	Xtriage
Outliers	1 of 37612 reflections (0.003%)	Xtriage
$F_o, F_c$ correlation	0.85	EDS
Total number of atoms	7832	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	64.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.83% 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	0.49	0/4467	0.76	1/6098 (0.0%)
2	B	0.52	0/3394	0.82	5/4626 (0.1%)
All	All	0.50	0/7861	0.79	6/10724 (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	2

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	63	ILE	N-CA-C	5.85	126.80	111.00
2	B	196	GLY	N-CA-C	-5.64	99.00	113.10
1	A	538	ALA	N-CA-C	5.58	126.07	111.00
2	B	419	THR	N-CA-C	5.49	125.82	111.00
2	B	425	LEU	N-CA-C	5.23	125.12	111.00
2	B	277	ARG	N-CA-C	5.19	125.02	111.00

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	188	TYR	Sidechain
1	A	501	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	4349	0	4251	103	2
2	B	3301	0	3228	104	1
3	A	78	0	0	7	0
3	B	104	0	0	7	1
All	All	7832	0	7479	203	2

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 13.

All (203) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:465:LYS:HG2	3:A:620:HOH:O	0.93	1.10
2:B:57:ASN:HD22	2:B:143:ARG:HH21	1.05	0.97
1:A:142:ILE:H	1:A:142:ILE:HD13	1.27	0.96
2:B:332:GLN:HE22	2:B:427:TYR:H	1.12	0.96
2:B:60:VAL:HG12	2:B:75:VAL:HG22	1.54	0.87
2:B:314:VAL:HG12	2:B:315:HIS:H	1.39	0.87
2:B:57:ASN:HD22	2:B:143:ARG:NH2	1.80	0.79
2:B:253:THR:O	2:B:257:ILE:HD12	1.87	0.75
1:A:114:ALA:HB1	1:A:160:PHE:CZ	2.23	0.74
1:A:465:LYS:O	1:A:466:VAL:HG23	1.88	0.73
2:B:101:LYS:NZ	2:B:101:LYS:H	1.85	0.73
1:A:101:LYS:HD3	1:A:102:LYS:H	1.53	0.73
2:B:373:GLN:HE22	2:B:407:GLN:H	1.35	0.72
2:B:332:GLN:NE2	2:B:427:TYR:H	1.87	0.69
2:B:50:ILE:HD12	2:B:143:ARG:HB3	1.75	0.68
1:A:442:VAL:CG1	1:A:481:ALA:HB1	2.25	0.67
1:A:330:GLN:HE22	1:A:340:GLN:NE2	1.93	0.67
1:A:114:ALA:HB1	1:A:160:PHE:CE1	2.30	0.66
2:B:314:VAL:HG12	2:B:315:HIS:N	2.09	0.66
2:B:163:SER:O	2:B:167:ILE:HG13	1.95	0.66
2:B:70:LYS:HD3	2:B:70:LYS:H	1.60	0.65
1:A:544:GLY:O	1:A:548:VAL:HG23	1.96	0.65
2:B:101:LYS:H	2:B:101:LYS:HZ2	1.43	0.65

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:442:VAL:HG12	1:A:443:ASP:N	2.12	0.64
2:B:70:LYS:N	2:B:70:LYS:HD3	2.12	0.63
2:B:24:TRP:CZ3	2:B:403:THR:HG21	2.33	0.63
2:B:57:ASN:ND2	2:B:143:ARG:HH21	1.88	0.63
1:A:101:LYS:HD3	1:A:102:LYS:N	2.13	0.62
1:A:3:SER:OG	1:A:4:PRO:HD2	2.00	0.62
2:B:31:ILE:O	2:B:35:VAL:HG23	2.00	0.62
1:A:536:VAL:HG12	1:A:537:PRO:HD2	1.82	0.61
1:A:79:GLU:HG3	1:A:83:ARG:HE	1.65	0.60
2:B:100:LEU:HG	2:B:381:VAL:HG13	1.84	0.59
2:B:332:GLN:HE22	2:B:427:TYR:N	1.93	0.58
1:A:199:ARG:HH22	1:A:220:LYS:HD2	1.68	0.58
1:A:464:GLN:O	1:A:465:LYS:HB2	2.02	0.58
2:B:344:GLU:O	2:B:347:LYS:HB2	2.03	0.58
2:B:115:TYR:OH	2:B:157:PRO:HB3	2.04	0.58
1:A:78:ARG:O	1:A:82:LYS:HG3	2.03	0.57
1:A:26:LEU:HD12	1:A:26:LEU:H	1.69	0.57
2:B:157:PRO:HG3	2:B:184:MET:HA	1.87	0.57
1:A:343:GLN:HG2	1:A:349:LEU:HD11	1.86	0.57
2:B:277:ARG:HB3	2:B:278:GLN:HE21	1.70	0.57
1:A:96:HIS:ND1	1:A:97:PRO:HD2	2.19	0.56
2:B:373:GLN:NE2	2:B:407:GLN:H	2.03	0.56
1:A:378:GLU:O	1:A:378:GLU:HG2	2.04	0.56
1:A:540:LYS:O	1:A:542:ILE:N	2.35	0.56
1:A:254:VAL:HB	1:A:289:LEU:HA	1.86	0.56
2:B:38:CYS:SG	2:B:132:ILE:HD11	2.46	0.56
2:B:24:TRP:N	2:B:24:TRP:CD1	2.73	0.56
2:B:42:GLU:HB3	3:B:472:HOH:O	2.06	0.56
1:A:517:LEU:O	1:A:521:ILE:HG13	2.06	0.56
1:A:518:VAL:O	1:A:522:ILE:HG12	2.07	0.55
2:B:309:ILE:HG22	2:B:309:ILE:O	2.07	0.55
1:A:98:ALA:HB1	1:A:349:LEU:HB3	1.88	0.54
2:B:137:ASN:HD22	2:B:137:ASN:N	2.05	0.54
2:B:312:GLU:OE2	2:B:313:PRO:HD2	2.08	0.54
1:A:341:ILE:HD12	1:A:350:LYS:HB3	1.90	0.54
2:B:38:CYS:O	2:B:42:GLU:HB2	2.08	0.53
1:A:95:PRO:HB3	2:B:136:ASN:O	2.08	0.53
1:A:376:THR:HG23	1:A:386:THR:HG22	1.91	0.53
1:A:142:ILE:N	1:A:142:ILE:HD13	2.11	0.53
2:B:50:ILE:HG12	2:B:145:GLN:HE21	1.73	0.53
2:B:101:LYS:O	2:B:236:PRO:HB2	2.09	0.52
1:A:225:PRO:HG2	1:A:226:PRO:HD3	1.91	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:57:ASN:HD22	1:A:143:ARG:NH2	2.08	0.52
1:A:465:LYS:O	1:A:466:VAL:CG2	2.57	0.51
2:B:374:LYS:O	2:B:378:GLU:HB2	2.10	0.51
2:B:264:LEU:HD13	2:B:274:ILE:HD11	1.92	0.51
1:A:97:PRO:O	1:A:100:LEU:HB2	2.11	0.50
2:B:313:PRO:HG2	3:B:506:HOH:O	2.10	0.50
2:B:21:VAL:HG12	2:B:22:LYS:N	2.26	0.50
1:A:118:VAL:HG21	1:A:160:PHE:HD1	1.76	0.50
2:B:103:ASN:H	2:B:103:ASN:HD22	1.60	0.50
2:B:72:ARG:NH2	3:B:449:HOH:O	2.43	0.50
1:A:439:THR:HG22	1:A:494:ASN:HB2	1.93	0.50
2:B:149:LEU:HB3	2:B:156:SER:HB3	1.94	0.50
1:A:27:THR:OG1	1:A:30:LYS:HB2	2.13	0.49
1:A:268:SER:C	1:A:270:ILE:H	2.16	0.49
2:B:314:VAL:CG1	2:B:315:HIS:H	2.19	0.49
1:A:97:PRO:C	1:A:99:GLY:H	2.14	0.49
1:A:419:THR:HG22	1:A:421:PRO:HD2	1.94	0.48
1:A:142:ILE:H	1:A:142:ILE:CD1	2.08	0.48
1:A:56:TYR:O	1:A:143:ARG:NH2	2.43	0.48
1:A:309:ILE:O	1:A:312:GLU:HB3	2.12	0.48
1:A:498:ASP:HA	1:A:536:VAL:O	2.13	0.48
2:B:376:THR:CG2	2:B:386:THR:HG22	2.44	0.48
2:B:198:HIS:O	2:B:202:ILE:HG12	2.13	0.48
1:A:522:ILE:O	1:A:526:ILE:HG13	2.13	0.48
2:B:331:LYS:O	2:B:333:GLY:N	2.47	0.48
2:B:163:SER:O	2:B:166:LYS:HB2	2.14	0.48
2:B:54:ASN:HD21	2:B:56:TYR:HB2	1.79	0.48
1:A:430:GLU:HA	1:A:431:LYS:HZ1	1.78	0.48
1:A:543:GLY:H	2:B:283:LEU:HD12	1.77	0.48
2:B:37:ILE:O	2:B:41:MET:HG3	2.14	0.48
2:B:251:SER:HB3	2:B:294:PRO:HA	1.95	0.48
2:B:78:ARG:O	2:B:82:LYS:HG3	2.14	0.47
1:A:208:HIS:HE1	1:A:212:TRP:HE1	1.61	0.47
2:B:153:TRP:O	2:B:184:MET:SD	2.72	0.47
2:B:205:LEU:O	2:B:209:LEU:HG	2.15	0.47
1:A:363:ASN:OD1	1:A:365:VAL:HG22	2.14	0.47
1:A:442:VAL:CG1	1:A:443:ASP:N	2.76	0.47
1:A:109:LEU:HB3	1:A:216:THR:HG21	1.97	0.47
1:A:495:ILE:HG22	1:A:496:VAL:N	2.30	0.46
1:A:199:ARG:HH12	1:A:220:LYS:HZ3	1.64	0.46
2:B:330:GLN:OE1	2:B:340:GLN:NE2	2.33	0.46
2:B:38:CYS:O	2:B:42:GLU:N	2.48	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:338:THR:HG22	1:A:353:LYS:HB3	1.97	0.46
1:A:350:LYS:HA	3:A:616:HOH:O	2.16	0.45
1:A:118:VAL:O	1:A:148:VAL:HG22	2.17	0.45
2:B:175:ASN:HD21	2:B:201:LYS:NZ	2.14	0.45
1:A:198:HIS:O	1:A:202:ILE:HG12	2.16	0.45
1:A:439:THR:HG21	2:B:289:LEU:H	1.82	0.45
1:A:509:GLN:N	1:A:510:PRO:HD3	2.31	0.45
2:B:388:LYS:NZ	2:B:413:GLU:HG3	2.31	0.45
1:A:377:THR:O	1:A:381:VAL:HG23	2.17	0.45
2:B:26:LEU:HD12	2:B:133:PRO:HG3	1.98	0.45
2:B:66:LYS:CG	2:B:67:ASP:H	2.30	0.45
2:B:119:PRO:HA	2:B:148:VAL:HA	1.97	0.45
1:A:24:TRP:HA	1:A:25:PRO:HD2	1.77	0.45
1:A:191:SER:OG	1:A:198:HIS:CD2	2.70	0.45
2:B:184:MET:HB3	2:B:185:ASP:H	1.62	0.44
2:B:419:THR:HG22	2:B:421:PRO:HD2	1.98	0.44
2:B:149:LEU:HA	2:B:150:PRO:HD3	1.76	0.44
2:B:419:THR:HA	2:B:420:PRO:HD2	1.90	0.44
1:A:241:VAL:HG23	1:A:314:VAL:O	2.17	0.44
2:B:216:THR:HA	2:B:217:PRO:HD2	1.68	0.44
1:A:373:GLN:NE2	2:B:397:THR:HG23	2.32	0.44
1:A:58:THR:HG23	1:A:76:ASP:O	2.18	0.44
2:B:258:GLN:HG2	2:B:283:LEU:HD21	2.00	0.44
1:A:495:ILE:N	1:A:495:ILE:HD12	2.33	0.44
2:B:60:VAL:CG1	2:B:75:VAL:HG22	2.36	0.44
1:A:168:LEU:O	1:A:172:LYS:HG3	2.18	0.44
2:B:385:LYS:HG2	2:B:386:THR:N	2.32	0.43
1:A:53:GLU:CD	1:A:53:GLU:N	2.71	0.43
1:A:344:GLU:HB3	1:A:345:PRO:HD2	2.00	0.43
1:A:337:TRP:O	1:A:353:LYS:HA	2.18	0.43
1:A:120:LEU:O	1:A:125:ARG:NE	2.51	0.43
1:A:524:GLN:O	1:A:528:LYS:HG2	2.18	0.43
2:B:154:LYS:O	2:B:157:PRO:HD2	2.19	0.43
1:A:519:ASN:H	1:A:519:ASN:HD22	1.65	0.43
1:A:101:LYS:CD	1:A:102:LYS:H	2.26	0.43
2:B:78:ARG:NH1	2:B:412:PRO:O	2.51	0.43
2:B:209:LEU:HB3	2:B:214:LEU:HB2	2.01	0.43
1:A:441:TYR:CD1	1:A:441:TYR:N	2.86	0.43
1:A:441:TYR:HD1	1:A:441:TYR:N	2.15	0.43
1:A:63:ILE:HG22	1:A:64:LYS:N	2.33	0.43
2:B:319:TYR:O	2:B:321:PRO:HD3	2.19	0.43
1:A:118:VAL:O	1:A:148:VAL:CG2	2.66	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:298:GLU:HG2	2:B:298:GLU:H	1.60	0.43
1:A:97:PRO:C	1:A:99:GLY:N	2.72	0.42
2:B:5:ILE:HG22	2:B:5:ILE:O	2.19	0.42
1:A:103:ASN:O	1:A:104:LYS:C	2.57	0.42
1:A:205:LEU:O	1:A:209:LEU:HG	2.19	0.42
2:B:257:ILE:O	2:B:261:VAL:HG23	2.19	0.42
2:B:345:PRO:C	2:B:347:LYS:H	2.23	0.42
2:B:101:LYS:HE3	2:B:382:ILE:HA	2.02	0.42
1:A:280:SER:HA	1:A:283:LEU:HD12	2.02	0.42
1:A:199:ARG:HH12	1:A:220:LYS:NZ	2.18	0.42
2:B:190:GLY:HA2	3:B:450:HOH:O	2.18	0.42
1:A:229:TRP:O	1:A:230:MET:HB2	2.19	0.42
2:B:191:SER:OG	2:B:198:HIS:HD2	2.02	0.42
2:B:330:GLN:NE2	2:B:340:GLN:OE1	2.46	0.42
2:B:310:LEU:HD23	2:B:310:LEU:HA	1.85	0.42
2:B:32:LYS:HD2	3:B:480:HOH:O	2.19	0.42
2:B:329:ILE:HG22	2:B:330:GLN:N	2.34	0.42
1:A:228:LEU:HD23	3:A:570:HOH:O	2.20	0.42
1:A:96:HIS:HA	1:A:97:PRO:HD3	1.82	0.41
1:A:122:GLU:HB2	3:A:604:HOH:O	2.19	0.41
2:B:79:GLU:HG3	2:B:83:ARG:HH11	1.84	0.41
1:A:252:TRP:O	1:A:292:VAL:HG13	2.19	0.41
2:B:45:GLY:HA2	3:B:484:HOH:O	2.20	0.41
2:B:66:LYS:HG2	2:B:67:ASP:N	2.36	0.41
2:B:121:ASP:O	2:B:125:ARG:HG3	2.20	0.41
1:A:149:LEU:HA	1:A:150:PRO:HD3	1.77	0.41
2:B:320:ASP:H	2:B:343:GLN:HE22	1.67	0.41
2:B:137:ASN:ND2	2:B:137:ASN:N	2.67	0.41
2:B:272:PRO:C	2:B:274:ILE:H	2.24	0.41
1:A:440:PHE:CD1	1:A:440:PHE:N	2.87	0.41
1:A:320:ASP:HA	1:A:321:PRO:HD2	1.84	0.41
1:A:394:GLN:HB2	1:A:397:THR:OG1	2.20	0.41
2:B:97:PRO:C	2:B:99:GLY:H	2.24	0.41
2:B:336:GLN:HG2	2:B:353:LYS:HD3	2.03	0.41
2:B:332:GLN:NE2	2:B:426:TRP:HB2	2.36	0.41
1:A:320:ASP:OD2	1:A:323:LYS:HG3	2.21	0.41
1:A:368:LEU:O	1:A:371:ALA:HB3	2.21	0.41
2:B:54:ASN:C	2:B:54:ASN:HD22	2.24	0.41
1:A:495:ILE:CG2	1:A:496:VAL:N	2.84	0.41
1:A:5:ILE:H	1:A:5:ILE:HG13	1.58	0.41
2:B:250:ASP:N	3:B:491:HOH:O	2.42	0.41
1:A:23:GLN:H	1:A:23:GLN:CD	2.24	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:28:GLU:O	2:B:32:LYS:HB2	2.21	0.41
2:B:368:LEU:O	2:B:372:VAL:HG23	2.21	0.41
1:A:330:GLN:NE2	1:A:340:GLN:HE22	2.20	0.40
2:B:150:PRO:HB2	2:B:153:TRP:HB2	2.02	0.40
1:A:91:GLN:O	1:A:92:LEU:HB3	2.21	0.40
1:A:379:SER:OG	1:A:387:PRO:HD3	2.22	0.40
2:B:252:TRP:O	2:B:253:THR:O	2.39	0.40
2:B:398:TRP:O	2:B:402:TRP:HD1	2.04	0.40
1:A:548:VAL:HG11	3:A:566:HOH:O	2.20	0.40
2:B:66:LYS:CG	2:B:67:ASP:N	2.84	0.40
1:A:466:VAL:N	3:A:620:HOH:O	2.53	0.40
1:A:517:LEU:HG	3:A:617:HOH:O	2.22	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:250:ASP:CB	3:B:498:HOH:O[4_647]	1.80	0.40
1:A:251:SER:OG	2:B:15:GLY:O[4_647]	1.83	0.37

## 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	552/560 (99%)	488 (88%)	51 (9%)	13 (2%)	9	22
2	B	411/430 (96%)	345 (84%)	50 (12%)	16 (4%)	5	10
All	All	963/990 (97%)	833 (86%)	101 (10%)	29 (3%)	7	15

All (29) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	458	VAL
1	A	541	GLY
2	B	66	LYS

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Mol	Chain	Res	Type
2	B	88	TRP
2	B	90	VAL
2	B	95	PRO
2	B	253	THR
2	B	277	ARG
2	B	332	GLN
1	A	345	PRO
2	B	2	ILE
2	B	425	LEU
1	A	104	LYS
1	A	125	ARG
1	A	137	ASN
1	A	412	PRO
1	A	470	THR
2	B	98	ALA
2	B	217	PRO
1	A	66	LYS
1	A	465	LYS
2	B	4	PRO
2	B	290	THR
2	B	292	VAL
2	B	395	LYS
2	B	419	THR
1	A	269	GLN
1	A	543	GLY
1	A	552	VAL

### 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	452/500 (90%)	420 (93%)	32 (7%)	21	46
2	B	347/392 (88%)	318 (92%)	29 (8%)	16	34
All	All	799/892 (90%)	738 (92%)	61 (8%)	19	41

All (61) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	23	GLN
1	A	26	LEU
1	A	48	SER
1	A	65	LYS
1	A	72	ARG
1	A	75	VAL
1	A	101	LYS
1	A	104	LYS
1	A	113	ASP
1	A	142	ILE
1	A	145	GLN
1	A	173	LYS
1	A	179	VAL
1	A	197	GLN
1	A	208	HIS
1	A	210	LEU
1	A	255	ASN
1	A	303	LEU
1	A	323	LYS
1	A	330	GLN
1	A	340	GLN
1	A	361	HIS
1	A	378	GLU
1	A	399	GLU
1	A	409	THR
1	A	422	LEU
1	A	428	GLN
1	A	431	LYS
1	A	470	THR
1	A	497	THR
1	A	536	VAL
1	A	540	LYS
2	B	6	GLU
2	B	24	TRP
2	B	42	GLU
2	B	50	ILE
2	B	54	ASN
2	B	70	LYS
2	B	73	LYS
2	B	83	ARG
2	B	85	GLN
2	B	101	LYS
2	B	111	VAL

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Mol	Chain	Res	Type
2	B	113	ASP
2	B	142	ILE
2	B	184	MET
2	B	195	ILE
2	B	200	THR
2	B	212	TRP
2	B	248	GLU
2	B	249	LYS
2	B	271	TYR
2	B	282	LEU
2	B	283	LEU
2	B	290	THR
2	B	301	LEU
2	B	324	ASP
2	B	332	GLN
2	B	404	GLU
2	B	414	TRP
2	B	419	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	23	GLN
1	A	57	ASN
1	A	145	GLN
1	A	198	HIS
1	A	222	GLN
1	A	306	ASN
1	A	330	GLN
1	A	340	GLN
1	A	373	GLN
1	A	428	GLN
1	A	471	ASN
1	A	475	GLN
1	A	519	ASN
2	B	54	ASN
2	B	57	ASN
2	B	85	GLN
2	B	96	HIS
2	B	103	ASN
2	B	137	ASN
2	B	145	GLN

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Mol	Chain	Res	Type
2	B	147	ASN
2	B	198	HIS
2	B	332	GLN
2	B	334	GLN
2	B	373	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	554/560 (98%)	0.52	37 (6%) 17 19	22, 69, 100, 100	0
2	B	408/430 (94%)	0.34	21 (5%) 27 30	8, 58, 100, 100	0
All	All	962/990 (97%)	0.44	58 (6%) 21 23	8, 65, 100, 100	0

All (58) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	231	GLY	9.9
1	A	552	VAL	6.2
2	B	1	PRO	4.9
2	B	2	ILE	4.7
1	A	551	LEU	4.5
2	B	4	PRO	4.5
1	A	72	ARG	4.3
1	A	131	THR	4.3
1	A	2	ILE	4.2
2	B	67	ASP	4.1
1	A	197	GLN	4.1
1	A	1	PRO	4.0
1	A	134	SER	3.7
2	B	5	ILE	3.6
2	B	232	TYR	3.5
1	A	539	HIS	3.4
2	B	218	ASP	3.4
1	A	452	LEU	3.4
1	A	554	ALA	3.3
1	A	21	VAL	3.3
2	B	360	ALA	3.3
2	B	276	VAL	3.2
1	A	470	THR	3.2
1	A	68	SER	3.1

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Mol	Chain	Res	Type	RSRZ
2	B	3	SER	3.1
2	B	245	VAL	2.9
1	A	138	GLU	2.8
1	A	183	TYR	2.7
1	A	178	ILE	2.7
1	A	553	SER	2.7
1	A	288	ALA	2.6
1	A	446	ALA	2.6
1	A	454	LYS	2.6
1	A	69	THR	2.6
1	A	434	ILE	2.5
1	A	550	LYS	2.5
2	B	359	GLY	2.5
1	A	132	ILE	2.5
1	A	52	PRO	2.4
1	A	67	ASP	2.4
1	A	538	ALA	2.3
1	A	450	THR	2.3
1	A	133	PRO	2.2
2	B	272	PRO	2.2
1	A	140	PRO	2.2
1	A	422	LEU	2.2
2	B	420	PRO	2.2
1	A	136	ASN	2.1
2	B	316	GLY	2.1
2	B	313	PRO	2.1
2	B	6	GLU	2.1
2	B	301	LEU	2.1
1	A	360	ALA	2.1
1	A	146	TYR	2.1
1	A	59	PRO	2.1
2	B	421	PRO	2.1
1	A	49	LYS	2.0
2	B	68	SER	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.