



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 10:42 AM UTC

PDB ID : 1NXZ / pdb_00001nxz
Title : X-Ray Crystal Structure of Protein yggj_haein of Haemophilus influenzae.
Northeast Structural Genomics Consortium Target IR73.
Authors : Forouhar, F.; Shen, J.; Xiao, R.; Acton, T.B.; Rost, B.; Montelione, G.T.;
Tong, L.; Northeast Structural Genomics Consortium (NESG)
Deposited on : 2003-02-11
Resolution : 2.00 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
Mogul	:	2022.3.0, CSD as543be (2022)
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

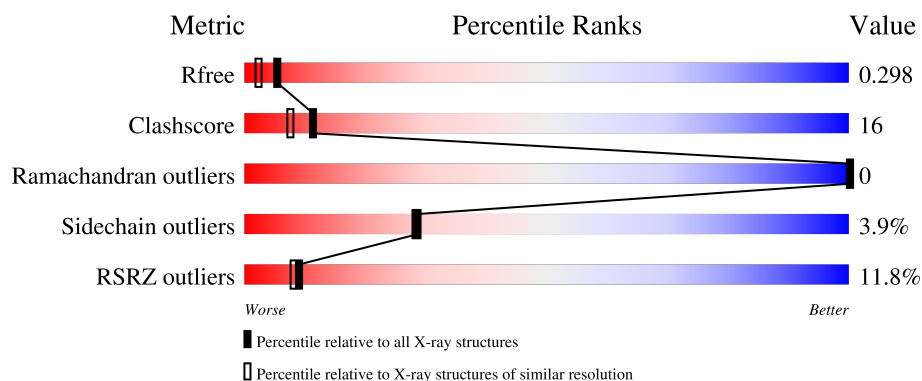
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	246	6% 70% 28% .
1	B	246	17% 65% 32% .

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4155 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 Hypothetical protein HI0303.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	246	Total	C	N	O	S	Se	0	0	0
			1920	1205	346	359	6	4			
1	B	245	Total	C	N	O	S	Se	0	0	0
			1915	1202	345	358	6	4			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	34	MSE	MET	modified residue	UNP P44627
A	92	MSE	MET	modified residue	UNP P44627
A	124	MSE	MET	modified residue	UNP P44627
A	154	MSE	MET	modified residue	UNP P44627
A	246	ALA	-	cloning artifact	UNP P44627
A	247	ALA	-	cloning artifact	UNP P44627
B	34	MSE	MET	modified residue	UNP P44627
B	92	MSE	MET	modified residue	UNP P44627
B	124	MSE	MET	modified residue	UNP P44627
B	154	MSE	MET	modified residue	UNP P44627
B	246	ALA	-	cloning artifact	UNP P44627
B	247	ALA	-	cloning artifact	UNP P44627

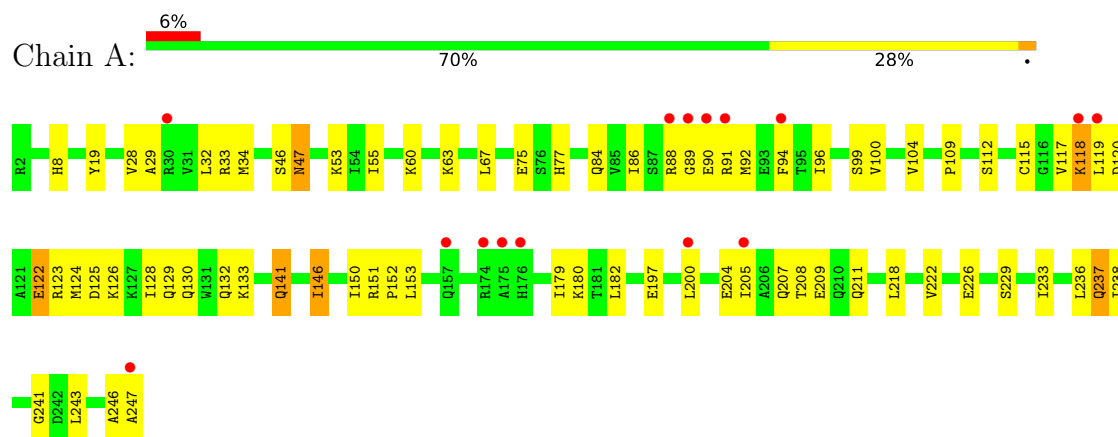
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	176	Total	O	0	0
			176	176		
2	B	144	Total	O	0	0
			144	144		

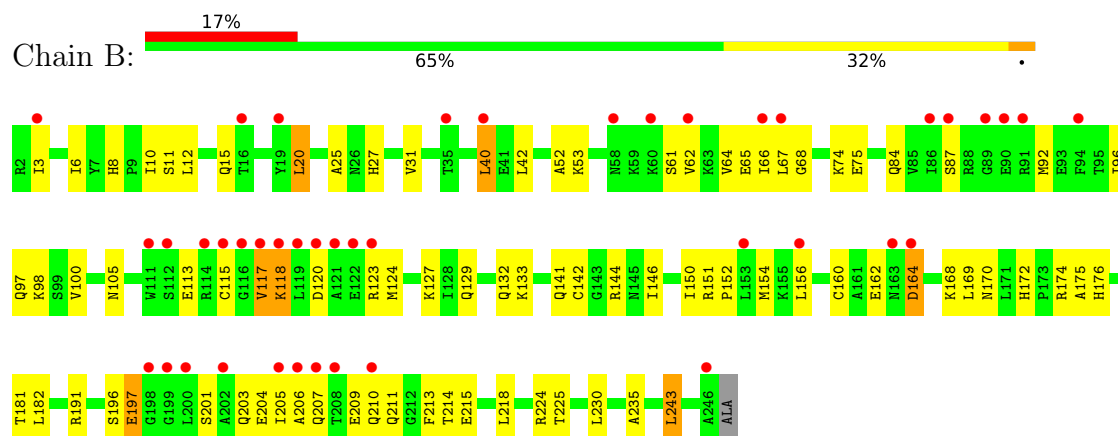
3 Residue-property plots

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: Hypothetical protein HI0303



• Molecule 1: Hypothetical protein HI0303



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	61.30Å 73.52Å 109.47Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.52 – 2.00 29.52 – 2.00	Depositor EDS
% Data completeness (in resolution range)	85.0 (29.52-2.00) 94.3 (29.52-2.00)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.87 (at 2.00Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.214 , 0.264 (Not available) , 0.298	Depositor DCC
R_{free} test set	6083 reflections (9.75%)	wwPDB-VP
Wilson B-factor (Å ²)	28.1	Xtriage
Anisotropy	0.054	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 47.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4155	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.94% 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

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	0.39	0/1945	0.89	6/2617 (0.2%)
1	B	0.37	0/1940	0.84	4/2610 (0.2%)
All	All	0.38	0/3885	0.86	10/5227 (0.2%)

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	8	HIS	CA-C-N	9.60	129.23	119.82
1	A	8	HIS	C-N-CA	9.60	129.23	119.82
1	B	8	HIS	CA-C-N	6.99	127.58	119.47
1	B	8	HIS	C-N-CA	6.99	127.58	119.47
1	A	146	ILE	N-CA-C	-6.94	98.39	108.11
1	A	141	GLN	N-CA-C	5.92	117.53	111.14
1	B	20	LEU	N-CA-C	-5.66	103.06	110.53
1	B	141	GLN	N-CA-C	5.24	116.79	111.14
1	A	182	LEU	CA-C-N	5.12	124.74	119.05
1	A	182	LEU	C-N-CA	5.12	124.74	119.05

There are no chirality outliers.

There are no planarity outliers.

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	1920	0	1964	64	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1915	0	1959	69	0
2	A	176	0	0	11	0
2	B	144	0	0	5	0
All	All	4155	0	3923	126	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 16.

All (126) 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:28:VAL:HG13	1:A:34:MSE:HE3	1.40	1.02
1:B:120:ASP:HB3	1:B:123:ARG:HB2	1.57	0.86
1:B:124:MSE:HB3	2:B:311:HOH:O	1.77	0.84
1:A:88:ARG:HD3	1:A:88:ARG:H	1.46	0.80
1:B:52:ALA:HA	1:B:67:LEU:HD23	1.63	0.79
1:A:120:ASP:HB3	1:A:123:ARG:HB3	1.67	0.76
1:B:12:LEU:HD22	1:B:66:ILE:HD11	1.68	0.75
1:A:88:ARG:HH11	1:A:89:GLY:H	1.35	0.73
1:A:222:VAL:HG13	1:B:142:CYS:HA	1.70	0.72
1:A:237:GLN:HE21	1:A:237:GLN:HA	1.54	0.72
1:B:117:VAL:HG12	1:B:118:LYS:H	1.54	0.71
1:B:66:ILE:HG22	1:B:68:GLY:H	1.55	0.70
1:A:197:GLU:HG2	2:A:367:HOH:O	1.95	0.66
1:A:60:LYS:HA	2:A:302:HOH:O	1.96	0.65
1:A:118:LYS:HD2	1:A:118:LYS:H	1.60	0.64
1:B:12:LEU:HA	1:B:15:GLN:HG3	1.79	0.64
1:B:132:GLN:HE22	1:B:150:ILE:H	1.46	0.63
1:A:32:LEU:HD12	1:A:34:MSE:HE2	1.82	0.62
1:B:25:ALA:CB	1:B:62:VAL:HG12	2.31	0.61
1:A:91:ARG:HA	2:A:413:HOH:O	2.01	0.60
1:A:207:GLN:HE22	1:A:211:GLN:HE21	1.49	0.60
1:A:125:ASP:O	1:A:129:GLN:HG3	2.02	0.60
1:B:64:VAL:HG12	1:B:65:GLU:N	2.17	0.59
1:B:87:SER:HB3	1:B:196:SER:HB2	1.84	0.59
1:A:237:GLN:HE21	1:A:241:GLY:HA3	1.67	0.58
1:A:53:LYS:HG3	1:A:67:LEU:HD11	1.85	0.58
1:B:87:SER:HA	1:B:117:VAL:HG22	1.86	0.58
1:B:97:GLN:HE21	1:B:98:LYS:NZ	2.02	0.58
1:B:61:SER:HB2	2:B:287:HOH:O	2.03	0.58
1:B:123:ARG:HG3	2:B:372:HOH:O	2.03	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:53:LYS:O	1:B:64:VAL:HG13	2.04	0.57
1:B:92:MSE:O	1:B:96:ILE:HG12	2.04	0.57
1:A:229:SER:O	1:A:233:ILE:HG12	2.05	0.57
1:A:179:ILE:HG23	1:A:236:LEU:HD23	1.87	0.57
1:A:55:ILE:HD11	1:A:63:LYS:HD2	1.87	0.56
1:A:90:GLU:HG2	1:A:91:ARG:N	2.21	0.56
1:A:75:GLU:HG3	1:A:146:ILE:HD12	1.86	0.56
1:A:96:ILE:O	1:A:100:VAL:HG23	2.07	0.53
1:A:247:ALA:HB3	1:B:176:HIS:HA	1.89	0.53
1:A:132:GLN:HE22	1:A:150:ILE:H	1.56	0.53
1:A:246:ALA:HB2	1:B:181:THR:HG21	1.90	0.53
1:B:205:ILE:O	1:B:209:GLU:HG3	2.09	0.53
1:A:207:GLN:NE2	1:A:211:GLN:HE21	2.07	0.52
1:B:96:ILE:O	1:B:100:VAL:HG23	2.09	0.52
1:A:46:SER:O	1:A:47:ASN:CB	2.58	0.52
1:B:172:HIS:ND1	1:B:174:ARG:HG2	2.24	0.52
1:B:230:LEU:C	1:B:230:LEU:HD23	2.35	0.52
1:B:151:ARG:HB3	1:B:152:PRO:HD2	1.92	0.51
1:A:19:TYR:OH	1:A:63:LYS:HE2	2.10	0.51
1:B:25:ALA:HB1	1:B:62:VAL:HG12	1.92	0.51
1:B:132:GLN:NE2	1:B:150:ILE:H	2.08	0.51
1:A:86:ILE:HD12	1:A:117:VAL:HG11	1.93	0.50
1:A:237:GLN:NE2	1:A:241:GLY:HA3	2.26	0.50
1:A:112:SER:HB3	1:A:115:CYS:HB3	1.95	0.49
1:B:169:LEU:HD23	1:B:214:THR:HB	1.93	0.49
1:A:122:GLU:O	1:A:126:LYS:HG2	2.12	0.49
1:A:119:LEU:O	1:A:119:LEU:HG	2.13	0.48
1:B:164:ASP:HB2	1:B:191:ARG:HE	1.77	0.48
1:A:132:GLN:NE2	1:A:150:ILE:H	2.10	0.48
1:A:130:GLN:HG2	2:A:382:HOH:O	2.14	0.48
1:B:3:ILE:HD11	1:B:40:LEU:C	2.39	0.47
1:B:115:CYS:SG	1:B:117:VAL:HG23	2.54	0.47
1:A:119:LEU:HB2	2:A:380:HOH:O	2.15	0.47
1:A:222:VAL:CG1	1:B:142:CYS:HA	2.42	0.47
1:B:62:VAL:HG13	1:B:62:VAL:O	2.15	0.47
1:B:225:THR:HG22	2:B:313:HOH:O	2.15	0.47
1:B:98:LYS:HD2	1:B:98:LYS:N	2.30	0.46
1:B:120:ASP:HB3	1:B:123:ARG:CB	2.38	0.46
1:B:160:CYS:HB3	1:B:211:GLN:HG3	1.97	0.46
1:B:170:ASN:HD22	1:B:213:PHE:HB3	1.80	0.46
1:A:88:ARG:HD3	1:A:88:ARG:N	2.22	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:204:GLU:O	1:A:208:THR:HG23	2.16	0.46
1:B:129:GLN:O	1:B:133:LYS:HD3	2.16	0.46
1:A:200:LEU:N	1:A:200:LEU:HD12	2.31	0.45
1:B:201:SER:O	1:B:205:ILE:HG13	2.16	0.45
2:A:312:HOH:O	1:B:224:ARG:HD3	2.16	0.45
1:A:118:LYS:HG2	1:A:123:ARG:HH21	1.82	0.44
1:A:151:ARG:HB3	1:A:152:PRO:HD2	2.00	0.44
1:B:156:LEU:HD13	1:B:204:GLU:HG2	1.99	0.44
1:B:203:GLN:O	1:B:206:ALA:HB3	2.18	0.44
1:A:118:LYS:HD2	1:A:118:LYS:N	2.31	0.43
1:A:207:GLN:HE21	1:A:211:GLN:HG2	1.83	0.43
1:A:237:GLN:HA	1:A:237:GLN:NE2	2.29	0.43
1:B:27:HIS:O	1:B:31:VAL:HB	2.18	0.43
1:A:47:ASN:HD22	1:A:47:ASN:HA	1.59	0.43
1:B:10:ILE:HG12	1:B:11:SER:H	1.84	0.43
1:B:124:MSE:HE3	1:B:124:MSE:HA	2.00	0.43
1:A:124:MSE:HG2	1:A:153:LEU:HD22	2.00	0.43
1:B:6:ILE:HB	1:B:42:LEU:HD23	2.00	0.43
1:A:120:ASP:OD1	1:A:122:GLU:HG2	2.18	0.43
1:A:128:ILE:HD13	1:A:152:PRO:HA	1.99	0.43
1:B:15:GLN:HB2	1:B:66:ILE:HD12	2.00	0.43
1:A:29:ALA:O	1:A:33:ARG:HA	2.19	0.43
1:B:168:LYS:HB3	1:B:213:PHE:CE1	2.53	0.43
1:B:170:ASN:O	1:B:215:GLU:HA	2.19	0.43
1:A:92:MSE:O	1:A:96:ILE:HG12	2.19	0.42
1:B:84:GLN:HG2	1:B:92:MSE:SE	2.70	0.42
1:A:205:ILE:O	1:A:209:GLU:HG3	2.19	0.42
1:B:64:VAL:HG12	1:B:65:GLU:H	1.81	0.42
1:B:75:GLU:HG3	1:B:146:ILE:HD12	2.01	0.42
1:A:118:LYS:NZ	1:A:118:LYS:HB3	2.34	0.42
1:A:243:LEU:HG	1:B:218:LEU:HD12	2.01	0.42
1:B:151:ARG:HH22	1:B:162:GLU:CD	2.28	0.42
1:B:243:LEU:HD23	1:B:243:LEU:HA	1.82	0.42
1:A:109:PRO:O	1:A:153:LEU:HD12	2.20	0.41
1:B:207:GLN:O	1:B:210:GLN:HG2	2.20	0.41
1:A:77:HIS:HE1	2:A:320:HOH:O	2.03	0.41
1:A:222:VAL:HG12	1:B:142:CYS:O	2.20	0.41
1:B:113:GLU:HB2	1:B:203:GLN:NE2	2.35	0.41
1:B:196:SER:O	1:B:197:GLU:C	2.64	0.41
1:A:133:LYS:HD2	2:A:382:HOH:O	2.20	0.41
1:B:64:VAL:CG1	1:B:65:GLU:N	2.83	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:84:GLN:HG3	1:A:92:MSE:HE1	2.01	0.41
1:B:6:ILE:HB	1:B:42:LEU:CD2	2.51	0.41
1:B:142:CYS:SG	1:B:144:ARG:HG3	2.60	0.41
1:A:19:TYR:CZ	1:A:63:LYS:HE2	2.56	0.41
1:A:141:GLN:NE2	2:A:260:HOH:O	2.51	0.41
1:A:77:HIS:HD2	2:A:264:HOH:O	2.03	0.41
1:B:105:ASN:HB3	2:B:355:HOH:O	2.20	0.41
1:B:124:MSE:CE	1:B:127:LYS:HG3	2.51	0.40
1:B:210:GLN:HG3	1:B:211:GLN:N	2.37	0.40
1:A:99:SER:HB3	1:A:104:VAL:HG21	2.03	0.40
1:A:180:LYS:O	1:A:180:LYS:HG2	2.22	0.40
1:A:238:ILE:HG13	1:B:235:ALA:CB	2.52	0.40
2:A:301:HOH:O	1:B:74:LYS:HE3	2.22	0.40
1:B:175:ALA:HB2	1:B:215:GLU:HB3	2.03	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	244/246 (99%)	232 (95%)	12 (5%)	0	100	100
1	B	243/246 (99%)	223 (92%)	20 (8%)	0	100	100
All	All	487/492 (99%)	455 (93%)	32 (7%)	0	100	100

There are no Ramachandran outliers to report.

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	207/203 (102%)	200 (97%)	7 (3%)	32	33
1	B	207/203 (102%)	198 (96%)	9 (4%)	26	25
All	All	414/406 (102%)	398 (96%)	16 (4%)	28	28

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	47	ASN
1	A	94	PHE
1	A	118	LYS
1	A	122	GLU
1	A	218	LEU
1	A	226	GLU
1	A	237	GLN
1	B	20	LEU
1	B	40	LEU
1	B	117	VAL
1	B	118	LYS
1	B	154	MSE
1	B	164	ASP
1	B	182	LEU
1	B	197	GLU
1	B	243	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	14	ASN
1	A	47	ASN
1	A	77	HIS
1	A	84	GLN
1	A	97	GLN
1	A	129	GLN
1	A	132	GLN
1	A	157	GLN
1	A	172	HIS
1	A	203	GLN
1	A	207	GLN

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Mol	Chain	Res	Type
1	A	211	GLN
1	A	237	GLN
1	B	15	GLN
1	B	39	GLN
1	B	97	GLN
1	B	130	GLN
1	B	132	GLN
1	B	141	GLN
1	B	207	GLN
1	B	210	GLN

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](#)

There are no ligands in this entry.

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	242/246 (98%)	0.41	15 (6%) 26 25	16, 28, 52, 69	0
1	B	241/246 (97%)	0.91	42 (17%) 4 3	17, 32, 67, 80	0
All	All	483/492 (98%)	0.66	57 (11%) 9 8	16, 31, 64, 80	0

All (57) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	120	ASP	5.7
1	B	246	ALA	5.5
1	B	111	TRP	4.5
1	B	121	ALA	4.5
1	A	119	LEU	4.4
1	B	117	VAL	4.3
1	A	176	HIS	4.3
1	B	202	ALA	4.2
1	B	116	GLY	4.1
1	A	175	ALA	4.0
1	B	119	LEU	3.9
1	B	198	GLY	3.9
1	A	89	GLY	3.9
1	A	88	ARG	3.8
1	A	91	ARG	3.7
1	B	118	LYS	3.6
1	A	90	GLU	3.5
1	B	199	GLY	3.5
1	B	205	ILE	3.4
1	B	115	CYS	3.4
1	B	89	GLY	3.4
1	B	206	ALA	3.3
1	B	86	ILE	3.2
1	B	164	ASP	3.1

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Mol	Chain	Res	Type	RSRZ
1	B	87	SER	3.1
1	B	200	LEU	3.1
1	A	118	LYS	2.9
1	A	94	PHE	2.7
1	B	114	ARG	2.7
1	B	58	ASN	2.7
1	B	3	ILE	2.7
1	B	153	LEU	2.6
1	B	66	ILE	2.6
1	B	210	GLN	2.6
1	B	208	THR	2.5
1	B	94	PHE	2.5
1	B	123	ARG	2.5
1	B	67	LEU	2.5
1	B	122	GLU	2.4
1	B	207	GLN	2.4
1	B	62	VAL	2.4
1	A	247	ALA	2.4
1	A	174	ARG	2.3
1	A	205	ILE	2.3
1	B	90	GLU	2.3
1	B	163	ASN	2.3
1	B	40	LEU	2.3
1	B	16	THR	2.2
1	B	91	ARG	2.2
1	B	156	LEU	2.2
1	A	30	ARG	2.2
1	B	60	LYS	2.2
1	B	112	SER	2.1
1	A	157	GLN	2.1
1	B	35	THR	2.1
1	B	19	TYR	2.0
1	A	200	LEU	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.