



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 02:30 PM UTC

PDB ID : 2GKI / pdb_00002gki
Title : Heavy and light chain variable single domains of an anti-DNA binding antibody hydrolyze both double- and single-stranded DNAs without sequence specificity
Authors : Kim, Y.R.; Kim, J.S.; Lee, S.H.; Lee, W.R.; Sohn, J.N.; Chung, Y.C.; Shim, H.K.; Lee, S.C.; Kwon, M.H.; Kim, Y.S.
Deposited on : 2006-04-02
Resolution : 2.88 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

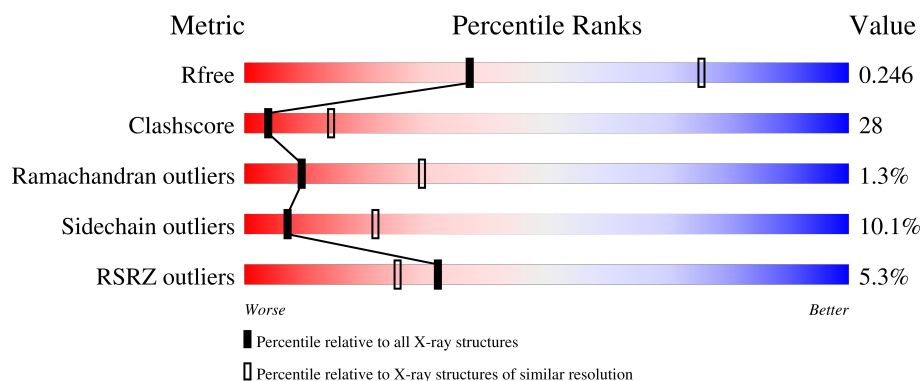
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.88 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	3557 (2.90-2.86)
Clashscore	190562	3801 (2.90-2.86)
Ramachandran outliers	187476	3699 (2.90-2.86)
Sidechain outliers	187428	3702 (2.90-2.86)
RSRZ outliers	180081	3558 (2.90-2.86)

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	291	4% 43% 31% 5% • 19%
1	B	291	4% 41% 34% 6% • 18%

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 3754 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 nuclease.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	235	Total	C	N	O	S	0	0	0
			1832	1157	303	361	11			
1	B	238	Total	C	N	O	S	0	0	0
			1852	1168	308	365	11			

There are 120 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	cloning artifact	GB 8777863
A	2	LYS	-	cloning artifact	GB 8777863
A	3	GLN	-	cloning artifact	GB 8777863
A	4	SER	-	cloning artifact	GB 8777863
A	5	THR	-	cloning artifact	GB 8777863
A	6	ILE	-	cloning artifact	GB 8777863
A	7	ALA	-	cloning artifact	GB 8777863
A	8	LEU	-	cloning artifact	GB 8777863
A	9	ALA	-	cloning artifact	GB 8777863
A	10	LEU	-	cloning artifact	GB 8777863
A	11	LEU	-	cloning artifact	GB 8777863
A	12	PRO	-	cloning artifact	GB 8777863
A	13	LEU	-	cloning artifact	GB 8777863
A	14	LEU	-	cloning artifact	GB 8777863
A	15	PHE	-	cloning artifact	GB 8777863
A	16	THR	-	cloning artifact	GB 8777863
A	17	PRO	-	cloning artifact	GB 8777863
A	18	VAL	-	cloning artifact	GB 8777863
A	19	THR	-	cloning artifact	GB 8777863
A	20	LYS	-	cloning artifact	GB 8777863
A	21	ALA	-	cloning artifact	GB 8777863
A	22	ARG	-	cloning artifact	GB 8777863
A	143	ARG	-	linker	GB 8777863
A	144	GLY	-	linker	GB 8777863
A	145	GLY	-	linker	GB 8777863

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Chain	Residue	Modelled	Actual	Comment	Reference
A	146	GLY	-	linker	GB 8777863
A	147	GLY	-	linker	GB 8777863
A	148	SER	-	linker	GB 8777863
A	149	GLY	-	linker	GB 8777863
A	150	GLY	-	linker	GB 8777863
A	151	GLY	-	linker	GB 8777863
A	152	GLY	-	linker	GB 8777863
A	153	SER	-	linker	GB 8777863
A	154	GLY	-	linker	GB 8777863
A	155	GLY	-	linker	GB 8777863
A	156	GLY	-	linker	GB 8777863
A	157	GLY	-	linker	GB 8777863
A	158	SER	-	linker	GB 8777863
A	160	LEU	ILE	SEE REMARK 999	GB 8777865
A	231	SER	PRO	SEE REMARK 999	GB 8777865
A	272	HIS	-	cloning artifact	GB 8777865
A	273	HIS	-	cloning artifact	GB 8777865
A	274	HIS	-	cloning artifact	GB 8777865
A	275	HIS	-	cloning artifact	GB 8777865
A	276	HIS	-	cloning artifact	GB 8777865
A	277	GLY	-	cloning artifact	GB 8777865
A	278	LEU	-	cloning artifact	GB 8777865
A	279	VAL	-	cloning artifact	GB 8777865
A	280	PRO	-	cloning artifact	GB 8777865
A	281	ARG	-	cloning artifact	GB 8777865
A	282	GLY	-	cloning artifact	GB 8777865
A	283	SER	-	cloning artifact	GB 8777865
A	284	GLY	-	cloning artifact	GB 8777865
A	285	ASP	-	cloning artifact	GB 8777865
A	286	PRO	-	cloning artifact	GB 8777865
A	287	LYS	-	cloning artifact	GB 8777865
A	288	ALA	-	cloning artifact	GB 8777865
A	289	ASP	-	cloning artifact	GB 8777865
A	290	ASN	-	cloning artifact	GB 8777865
A	291	LYS	-	cloning artifact	GB 8777865
B	1	MET	-	cloning artifact	GB 8777863
B	2	LYS	-	cloning artifact	GB 8777863
B	3	GLN	-	cloning artifact	GB 8777863
B	4	SER	-	cloning artifact	GB 8777863
B	5	THR	-	cloning artifact	GB 8777863
B	6	ILE	-	cloning artifact	GB 8777863
B	7	ALA	-	cloning artifact	GB 8777863

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Chain	Residue	Modelled	Actual	Comment	Reference
B	8	LEU	-	cloning artifact	GB 8777863
B	9	ALA	-	cloning artifact	GB 8777863
B	10	LEU	-	cloning artifact	GB 8777863
B	11	LEU	-	cloning artifact	GB 8777863
B	12	PRO	-	cloning artifact	GB 8777863
B	13	LEU	-	cloning artifact	GB 8777863
B	14	LEU	-	cloning artifact	GB 8777863
B	15	PHE	-	cloning artifact	GB 8777863
B	16	THR	-	cloning artifact	GB 8777863
B	17	PRO	-	cloning artifact	GB 8777863
B	18	VAL	-	cloning artifact	GB 8777863
B	19	THR	-	cloning artifact	GB 8777863
B	20	LYS	-	cloning artifact	GB 8777863
B	21	ALA	-	cloning artifact	GB 8777863
B	22	ARG	-	cloning artifact	GB 8777863
B	143	ARG	-	linker	GB 8777863
B	144	GLY	-	linker	GB 8777863
B	145	GLY	-	linker	GB 8777863
B	146	GLY	-	linker	GB 8777863
B	147	GLY	-	linker	GB 8777863
B	148	SER	-	linker	GB 8777863
B	149	GLY	-	linker	GB 8777863
B	150	GLY	-	linker	GB 8777863
B	151	GLY	-	linker	GB 8777863
B	152	GLY	-	linker	GB 8777863
B	153	SER	-	linker	GB 8777863
B	154	GLY	-	linker	GB 8777863
B	155	GLY	-	linker	GB 8777863
B	156	GLY	-	linker	GB 8777863
B	157	GLY	-	linker	GB 8777863
B	158	SER	-	linker	GB 8777863
B	160	LEU	ILE	SEE REMARK 999	GB 8777865
B	231	SER	PRO	SEE REMARK 999	GB 8777865
B	272	HIS	-	cloning artifact	GB 8777865
B	273	HIS	-	cloning artifact	GB 8777865
B	274	HIS	-	cloning artifact	GB 8777865
B	275	HIS	-	cloning artifact	GB 8777865
B	276	HIS	-	cloning artifact	GB 8777865
B	277	GLY	-	cloning artifact	GB 8777865
B	278	LEU	-	cloning artifact	GB 8777865
B	279	VAL	-	cloning artifact	GB 8777865
B	280	PRO	-	cloning artifact	GB 8777865

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Chain	Residue	Modelled	Actual	Comment	Reference
B	281	ARG	-	cloning artifact	GB 8777865
B	282	GLY	-	cloning artifact	GB 8777865
B	283	SER	-	cloning artifact	GB 8777865
B	284	GLY	-	cloning artifact	GB 8777865
B	285	ASP	-	cloning artifact	GB 8777865
B	286	PRO	-	cloning artifact	GB 8777865
B	287	LYS	-	cloning artifact	GB 8777865
B	288	ALA	-	cloning artifact	GB 8777865
B	289	ASP	-	cloning artifact	GB 8777865
B	290	ASN	-	cloning artifact	GB 8777865
B	291	LYS	-	cloning artifact	GB 8777865

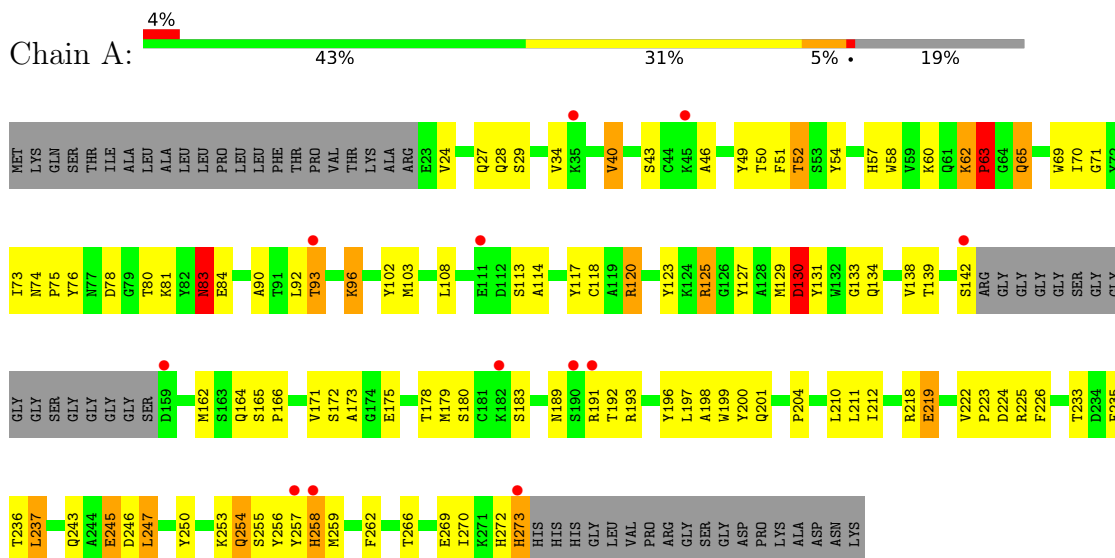
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	36	Total O 36 36	0	0
2	B	34	Total O 34 34	0	0

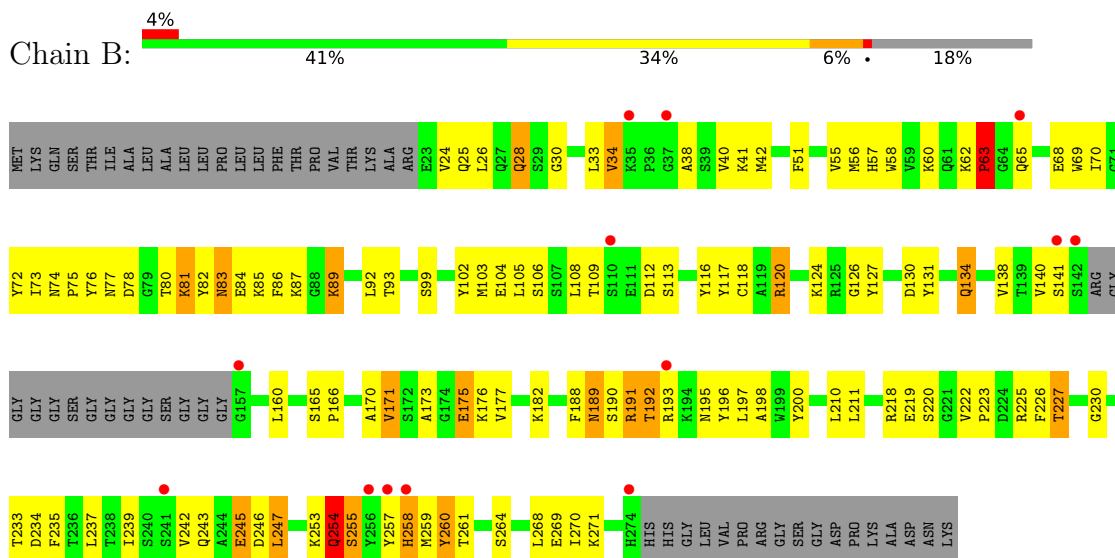
3 Residue-property plots

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: nuclease



• Molecule 1: nuclease



4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	179.74Å 179.74Å 184.23Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	49.58 – 2.88 49.58 – 2.88	Depositor EDS
% Data completeness (in resolution range)	99.4 (49.58-2.88) 99.4 (49.58-2.88)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.16 (at 2.86Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.212 , 0.242 0.216 , 0.246	Depositor DCC
R_{free} test set	4017 reflections (10.04%)	wwPDB-VP
Wilson B-factor (Å ²)	53.4	Xtriage
Anisotropy	0.066	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 94.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	3754	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.98% 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 [i](#)

5.1 Standard geometry [i](#)

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.39	0/1877	0.92	6/2535 (0.2%)
1	B	0.38	0/1898	0.95	8/2563 (0.3%)
All	All	0.39	0/3775	0.94	14/5098 (0.3%)

There are no bond length outliers.

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	83	ASN	N-CA-C	-6.85	98.78	109.25
1	B	255	SER	N-CA-C	6.64	119.44	108.49
1	A	65	GLN	N-CA-C	6.54	118.29	110.19
1	A	118	CYS	N-CA-C	-6.53	98.76	109.40
1	B	89	LYS	N-CA-C	5.61	118.59	111.69
1	B	254	GLN	N-CA-C	-5.57	101.60	109.96
1	A	62	LYS	CA-C-N	5.50	126.71	119.84
1	A	62	LYS	C-N-CA	5.50	126.71	119.84
1	B	83	ASN	N-CA-C	-5.47	101.07	109.76
1	B	30	GLY	CA-C-N	5.36	125.30	119.78
1	B	30	GLY	C-N-CA	5.36	125.30	119.78
1	B	34	VAL	N-CA-C	5.17	114.84	108.53
1	B	76	TYR	N-CA-C	5.07	116.50	111.07
1	A	133	GLY	N-CA-C	-5.02	104.77	112.85

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	1832	0	1764	105	1
1	B	1852	0	1779	103	0
2	A	36	0	0	0	0
2	B	34	0	0	2	1
All	All	3754	0	3543	204	2

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 28.

All (204) 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:B:34:VAL:HG21	1:B:108:LEU:HD13	1.38	0.99
1:A:70:ILE:HD13	1:A:103:MET:HE1	1.44	0.97
1:A:40:VAL:HG22	1:A:108:LEU:HD11	1.51	0.92
1:A:34:VAL:HG21	1:A:108:LEU:HD13	1.59	0.84
1:B:62:LYS:HB2	1:B:65:GLN:HG2	1.61	0.82
1:B:103:MET:HE3	1:B:105:LEU:HD21	1.61	0.81
1:A:57:HIS:HD2	1:A:69:TRP:HE1	1.28	0.81
1:B:253:LYS:NZ	1:B:255:SER:HB3	1.96	0.81
1:B:253:LYS:HZ2	1:B:255:SER:HB3	1.48	0.79
1:B:189:ASN:HD22	1:B:189:ASN:C	1.92	0.77
1:B:55:VAL:HG21	1:B:72:TYR:CD1	2.19	0.76
1:A:189:ASN:CG	1:A:192:THR:HG22	2.11	0.75
1:B:192:THR:HG21	1:B:196:TYR:OH	1.87	0.74
1:A:253:LYS:NZ	1:A:255:SER:HB3	2.05	0.71
1:B:171:VAL:HG11	1:B:177:VAL:CG1	2.21	0.70
1:A:200:TYR:CE2	1:A:210:LEU:HD23	2.26	0.70
1:A:70:ILE:CD1	1:A:103:MET:HE1	2.21	0.70
1:A:225:ARG:HD3	1:A:243:GLN:HE21	1.57	0.69
1:A:200:TYR:CE1	1:A:253:LYS:HE3	2.28	0.69
1:B:171:VAL:HG11	1:B:177:VAL:HG12	1.75	0.68
1:A:257:TYR:O	1:A:258:HIS:ND1	2.26	0.68
1:A:183:SER:O	1:A:233:THR:HG23	1.93	0.68
1:A:189:ASN:OD1	1:A:191:ARG:HB3	1.95	0.67
1:B:218:ARG:HH12	1:B:227:THR:HG23	1.59	0.67
1:B:34:VAL:HG21	1:B:108:LEU:CD1	2.22	0.66
1:A:70:ILE:HG21	1:A:103:MET:CE	2.25	0.66
1:B:200:TYR:CE1	1:B:253:LYS:HE3	2.30	0.66
1:A:253:LYS:HD3	1:A:262:PHE:CE1	2.32	0.65
1:A:165:SER:HB2	1:A:166:PRO:HA	1.79	0.65
1:A:90:ALA:HB1	1:A:103:MET:HE2	1.78	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:225:ARG:NH2	1:B:246:ASP:OD1	2.31	0.63
1:A:52:THR:HG22	1:A:76:TYR:N	2.13	0.63
1:B:257:TYR:O	1:B:258:HIS:ND1	2.28	0.63
1:B:197:LEU:HD22	1:B:235:PHE:CG	2.33	0.63
1:B:270:ILE:HD12	1:B:270:ILE:N	2.14	0.62
1:B:225:ARG:HH21	1:B:246:ASP:CG	2.08	0.62
1:A:211:LEU:HB3	1:A:212:ILE:HD12	1.81	0.62
1:B:120:ARG:NH1	1:B:130:ASP:OD1	2.34	0.61
1:A:218:ARG:NH1	1:A:224:ASP:HA	2.15	0.61
1:A:270:ILE:HD12	1:A:270:ILE:N	2.16	0.61
1:A:96:LYS:HD3	1:A:96:LYS:H	1.64	0.61
1:A:164:GLN:HB2	1:A:180:SER:O	2.01	0.60
1:B:55:VAL:HG23	1:B:73:ILE:C	2.27	0.60
1:B:89:LYS:HD3	1:B:112:ASP:OD1	2.01	0.60
1:A:69:TRP:CZ2	1:A:71:GLY:HA2	2.37	0.60
1:A:73:ILE:HA	1:A:80:THR:HG22	1.83	0.59
1:B:233:THR:HG23	1:B:234:ASP:OD1	2.03	0.59
1:A:225:ARG:HD3	1:A:243:GLN:NE2	2.17	0.59
1:B:189:ASN:HD22	1:B:190:SER:N	2.00	0.59
1:B:189:ASN:C	1:B:189:ASN:ND2	2.60	0.58
1:A:257:TYR:CD1	1:A:258:HIS:HB3	2.39	0.58
1:A:63:PRO:HD3	1:A:114:ALA:HA	1.86	0.58
1:A:70:ILE:HG21	1:A:103:MET:HE1	1.84	0.58
1:A:34:VAL:HG11	1:A:40:VAL:HG13	1.87	0.57
1:A:223:PRO:HG2	1:A:226:PHE:CD2	2.39	0.57
1:B:160:LEU:C	1:B:160:LEU:HD23	2.30	0.57
1:A:57:HIS:HD2	1:A:69:TRP:NE1	2.00	0.57
1:B:200:TYR:HE1	1:B:253:LYS:HE3	1.69	0.57
1:A:210:LEU:HD12	1:A:219:GLU:CG	2.34	0.57
1:A:57:HIS:CD2	1:A:69:TRP:HE1	2.17	0.57
1:A:210:LEU:HD12	1:A:219:GLU:HG3	1.87	0.57
1:B:197:LEU:HG	1:B:198:ALA:N	2.20	0.56
1:A:196:TYR:HB3	1:A:255:SER:O	2.05	0.56
1:B:171:VAL:HG21	1:B:242:VAL:HG21	1.88	0.56
1:A:73:ILE:O	1:A:75:PRO:HD3	2.07	0.55
1:B:109:THR:O	1:B:140:VAL:HG21	2.07	0.55
1:A:52:THR:HA	1:A:75:PRO:HB2	1.89	0.55
1:A:201:GLN:HG3	1:A:250:TYR:CE2	2.42	0.55
1:A:34:VAL:HG21	1:A:108:LEU:CD1	2.33	0.54
1:B:70:ILE:HG12	1:B:86:PHE:CE1	2.41	0.54
1:B:173:ALA:O	1:B:242:VAL:O	2.24	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:90:ALA:CB	1:A:103:MET:HE2	2.36	0.54
1:B:257:TYR:CD1	1:B:258:HIS:HB3	2.42	0.54
1:A:123:TYR:C	1:A:125:ARG:H	2.16	0.54
1:A:192:THR:HG21	1:A:196:TYR:OH	2.08	0.54
1:A:269:GLU:C	1:A:270:ILE:HD12	2.32	0.53
1:B:223:PRO:HG2	1:B:226:PHE:CD2	2.43	0.53
1:A:142:SER:HB2	1:A:257:TYR:OH	2.08	0.53
1:B:188:PHE:HB2	1:B:195:ASN:OD1	2.09	0.53
1:B:93:THR:OG1	1:B:102:TYR:HB2	2.09	0.53
1:B:124:LYS:HE2	2:B:303:HOH:O	2.07	0.52
1:B:55:VAL:HG23	1:B:73:ILE:O	2.08	0.52
1:B:189:ASN:HD21	1:B:191:ARG:HB2	1.74	0.52
1:B:28:GLN:OE1	1:B:117:TYR:HA	2.09	0.52
1:B:243:GLN:HB3	1:B:245:GLU:OE2	2.10	0.52
1:A:253:LYS:HZ2	1:A:255:SER:HB3	1.73	0.52
1:B:25:GLN:O	1:B:26:LEU:HD23	2.09	0.52
1:A:162:MET:SD	1:A:183:SER:HB3	2.50	0.52
1:B:57:HIS:HD2	1:B:69:TRP:HE1	1.59	0.51
1:A:58:TRP:CD1	1:A:92:LEU:HD22	2.45	0.51
1:A:130:ASP:HB3	1:A:131:TYR:CD1	2.45	0.51
1:A:189:ASN:O	1:A:193:ARG:N	2.44	0.51
1:B:62:LYS:HB3	1:B:63:PRO:HD2	1.92	0.51
1:B:189:ASN:O	1:B:193:ARG:N	2.41	0.51
1:A:253:LYS:HD3	1:A:262:PHE:CZ	2.46	0.50
1:A:120:ARG:NH1	1:A:130:ASP:OD1	2.44	0.50
1:A:130:ASP:HB3	1:A:131:TYR:HD1	1.75	0.50
1:A:34:VAL:HG11	1:A:40:VAL:CG1	2.41	0.50
1:A:201:GLN:HB2	1:A:211:LEU:HD11	1.93	0.50
1:A:200:TYR:HE1	1:A:253:LYS:HE3	1.74	0.50
1:B:84:GLU:O	1:B:87:LYS:HB2	2.12	0.50
1:B:40:VAL:O	1:B:104:GLU:HA	2.12	0.50
1:A:245:GLU:H	1:A:245:GLU:CD	2.19	0.49
1:B:57:HIS:O	1:B:118:CYS:HA	2.12	0.49
1:A:225:ARG:NH2	1:A:246:ASP:OD1	2.39	0.49
1:A:179:MET:HG2	1:A:266:THR:HG21	1.94	0.49
1:B:230:GLY:HA3	1:B:235:PHE:HA	1.94	0.49
1:A:200:TYR:CZ	1:A:210:LEU:HD23	2.47	0.48
1:B:260:TYR:N	1:B:260:TYR:CD1	2.81	0.48
1:B:130:ASP:HB3	1:B:131:TYR:HD1	1.77	0.48
1:A:171:VAL:HG13	1:A:175:GLU:HB2	1.95	0.48
1:B:62:LYS:HB2	1:B:65:GLN:CG	2.39	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:51:PHE:CE2	1:B:75:PRO:HB3	2.48	0.48
1:B:70:ILE:HG12	1:B:86:PHE:CD1	2.48	0.48
1:B:259:MET:C	1:B:260:TYR:CD1	2.91	0.48
1:B:257:TYR:CD1	1:B:257:TYR:C	2.92	0.47
1:B:83:ASN:O	1:B:84:GLU:C	2.57	0.47
1:A:171:VAL:CG1	1:A:175:GLU:HB2	2.44	0.47
1:B:72:TYR:CD1	1:B:72:TYR:C	2.92	0.47
1:B:254:GLN:HG3	1:B:261:THR:HB	1.96	0.47
1:A:123:TYR:C	1:A:125:ARG:N	2.72	0.47
1:B:86:PHE:CD2	1:B:86:PHE:N	2.82	0.47
1:B:89:LYS:HE2	1:B:106:SER:O	2.14	0.47
1:A:197:LEU:HD22	1:A:235:PHE:CG	2.50	0.47
1:B:223:PRO:HG2	1:B:226:PHE:HD2	1.80	0.47
1:B:62:LYS:CB	1:B:65:GLN:HG2	2.39	0.47
1:A:96:LYS:HD3	1:A:96:LYS:N	2.28	0.46
1:B:103:MET:HE3	1:B:105:LEU:CD2	2.36	0.46
1:A:83:ASN:O	1:A:84:GLU:C	2.58	0.46
1:B:170:ALA:HB1	1:B:271:LYS:HD2	1.96	0.46
1:A:73:ILE:HD12	1:A:80:THR:CG2	2.46	0.46
1:A:211:LEU:CB	1:A:212:ILE:HD12	2.46	0.46
1:B:182:LYS:HE3	1:B:234:ASP:OD2	2.15	0.46
1:A:127:TYR:CZ	1:B:257:TYR:HB3	2.51	0.46
1:A:92:LEU:C	1:A:93:THR:HG22	2.39	0.46
1:B:247:LEU:HD23	1:B:268:LEU:O	2.16	0.46
1:A:129:MET:HE3	1:B:253:LYS:HE2	1.98	0.46
1:B:177:VAL:HG22	1:B:239:ILE:HB	1.97	0.46
1:A:223:PRO:HG2	1:A:226:PHE:HD2	1.79	0.45
1:A:51:PHE:CE2	1:A:75:PRO:HB3	2.52	0.45
1:B:219:GLU:HG3	1:B:220:SER:N	2.31	0.45
1:A:272:HIS:ND1	1:A:273:HIS:N	2.64	0.45
1:B:211:LEU:HA	1:B:222:VAL:HG21	1.98	0.45
1:A:254:GLN:OE1	1:A:259:MET:HG3	2.17	0.45
1:A:83:ASN:OD1	1:A:83:ASN:C	2.58	0.45
1:B:113:SER:HA	1:B:138:VAL:O	2.17	0.45
1:B:42:MET:HE3	1:B:103:MET:CE	2.46	0.45
1:B:58:TRP:CD1	1:B:92:LEU:HD22	2.51	0.45
1:A:73:ILE:HD12	1:A:80:THR:HG22	1.99	0.44
1:B:134:GLN:HE21	1:B:134:GLN:HB3	1.62	0.44
1:A:28:GLN:OE1	1:A:117:TYR:HA	2.18	0.44
1:A:29:SER:HB3	1:A:43:SER:OG	2.17	0.44
1:A:172:SER:O	1:A:173:ALA:C	2.60	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:218:ARG:HG2	1:A:222:VAL:HB	1.99	0.44
1:B:226:PHE:CD1	1:B:239:ILE:HG12	2.53	0.44
1:A:191:ARG:HH11	1:A:191:ARG:HG2	1.83	0.44
1:B:170:ALA:HA	1:B:269:GLU:O	2.18	0.44
1:A:257:TYR:CD1	1:A:257:TYR:C	2.96	0.44
1:B:51:PHE:CD2	1:B:99:SER:HA	2.52	0.44
1:B:65:GLN:CD	1:B:65:GLN:N	2.71	0.44
1:B:176:LYS:HG2	1:B:177:VAL:N	2.33	0.44
1:A:172:SER:O	1:A:175:GLU:HG2	2.17	0.43
1:B:73:ILE:HG13	1:B:80:THR:HG22	2.01	0.43
1:A:27:GLN:NE2	1:A:134:GLN:HE22	2.17	0.43
1:A:62:LYS:HE2	1:A:65:GLN:HE22	1.83	0.43
1:A:189:ASN:OD1	1:A:192:THR:HG22	2.17	0.43
1:A:223:PRO:HG2	1:A:226:PHE:CE2	2.54	0.43
1:B:34:VAL:HG11	1:B:40:VAL:CG1	2.48	0.43
1:A:93:THR:HG23	1:A:102:TYR:HB2	2.00	0.43
1:B:83:ASN:ND2	1:B:85:LYS:HB2	2.34	0.43
1:A:270:ILE:N	1:A:270:ILE:CD1	2.82	0.43
1:B:86:PHE:O	1:B:87:LYS:C	2.61	0.42
1:B:126:GLY:O	1:B:127:TYR:HB2	2.18	0.42
1:B:165:SER:HA	1:B:166:PRO:C	2.43	0.42
1:A:113:SER:HA	1:A:138:VAL:O	2.19	0.42
1:B:74:ASN:HB3	1:B:77:ASN:OD1	2.19	0.42
1:A:210:LEU:HD12	1:A:219:GLU:HG2	2.01	0.42
1:B:103:MET:HE1	1:B:116:TYR:CD2	2.55	0.42
1:B:175:GLU:O	1:B:242:VAL:HG23	2.19	0.42
1:A:54:TYR:CD1	1:A:54:TYR:N	2.87	0.42
1:A:60:LYS:HB2	1:A:70:ILE:HD11	2.01	0.42
1:B:245:GLU:H	1:B:245:GLU:CD	2.28	0.42
1:B:124:LYS:CE	2:B:303:HOH:O	2.66	0.42
1:A:127:TYR:CE2	1:B:257:TYR:HB3	2.54	0.42
1:B:26:LEU:CD2	1:B:56:MET:HE1	2.49	0.41
1:A:52:THR:HG22	1:A:75:PRO:C	2.45	0.41
1:A:127:TYR:HB2	1:B:255:SER:O	2.20	0.41
1:B:60:LYS:HB2	1:B:70:ILE:HD11	2.02	0.41
1:A:74:ASN:O	1:A:78:ASP:N	2.53	0.41
1:A:24:VAL:HG11	1:A:131:TYR:CD1	2.56	0.41
1:A:199:TRP:CE2	1:A:237:LEU:HB2	2.56	0.41
1:A:204:PRO:HG2	1:A:247:LEU:HD13	2.03	0.41
1:B:81:LYS:HD3	1:B:82:TYR:N	2.36	0.41
1:B:57:HIS:HD2	1:B:69:TRP:NE1	2.17	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:257:TYR:O	1:B:258:HIS:CB	2.69	0.41
1:B:24:VAL:HB	1:B:131:TYR:CE2	2.57	0.40
1:B:28:GLN:N	1:B:134:GLN:OE1	2.53	0.40
1:B:41:LYS:HA	1:B:103:MET:O	2.20	0.40
1:A:46:ALA:HB1	1:A:49:TYR:CE2	2.56	0.40
1:A:256:TYR:O	1:A:257:TYR:C	2.63	0.40
1:B:68:GLU:OE1	1:B:85:LYS:HE2	2.21	0.40
1:B:84:GLU:HA	1:B:87:LYS:HD3	2.03	0.40
1:A:52:THR:HG22	1:A:76:TYR:CA	2.51	0.40
1:A:197:LEU:HG	1:A:198:ALA:N	2.36	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	Interatomic distance (Å)	Clash overlap (Å)
1:A:123:TYR:OH	1:A:123:TYR:OH[12_557]	1.75	0.45
2:B:308:HOH:O	2:B:308:HOH:O[9_766]	2.11	0.09

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	231/291 (79%)	211 (91%)	18 (8%)	2 (1%)	14	38
1	B	234/291 (80%)	206 (88%)	24 (10%)	4 (2%)	7	23
All	All	465/582 (80%)	417 (90%)	42 (9%)	6 (1%)	9	29

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	63	PRO
1	B	175	GLU

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Mol	Chain	Res	Type
1	B	38	ALA
1	B	141	SER
1	A	130	ASP
1	B	63	PRO

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	201/238 (84%)	180 (90%)	21 (10%)	7	20
1	B	203/238 (85%)	183 (90%)	20 (10%)	7	22
All	All	404/476 (85%)	363 (90%)	41 (10%)	7	22

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	40	VAL
1	A	50	THR
1	A	52	THR
1	A	63	PRO
1	A	81	LYS
1	A	83	ASN
1	A	93	THR
1	A	96	LYS
1	A	120	ARG
1	A	125	ARG
1	A	130	ASP
1	A	139	THR
1	A	178	THR
1	A	219	GLU
1	A	236	THR
1	A	237	LEU
1	A	245	GLU
1	A	247	LEU
1	A	254	GLN

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Mol	Chain	Res	Type
1	A	258	HIS
1	A	273	HIS
1	B	28	GLN
1	B	33	LEU
1	B	63	PRO
1	B	78	ASP
1	B	81	LYS
1	B	120	ARG
1	B	134	GLN
1	B	171	VAL
1	B	189	ASN
1	B	191	ARG
1	B	192	THR
1	B	210	LEU
1	B	227	THR
1	B	237	LEU
1	B	245	GLU
1	B	247	LEU
1	B	254	GLN
1	B	258	HIS
1	B	260	TYR
1	B	264	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	27	GLN
1	A	61	GLN
1	A	65	GLN
1	A	185	GLN
1	A	201	GLN
1	A	243	GLN
1	B	57	HIS
1	B	83	ASN
1	B	185	GLN
1	B	189	ASN
1	B	202	GLN
1	B	254	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	235/291 (80%)	-0.07	12 (5%) 33 27	23, 44, 79, 140	0
1	B	238/291 (81%)	0.23	13 (5%) 30 24	21, 49, 97, 140	0
All	All	473/582 (81%)	0.08	25 (5%) 32 25	21, 46, 89, 140	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	274	HIS	7.7
1	B	142	SER	7.1
1	A	273	HIS	5.8
1	B	257	TYR	5.1
1	A	142	SER	4.2
1	B	258	HIS	4.2
1	A	191	ARG	3.8
1	B	256	TYR	3.6
1	A	190	SER	3.5
1	B	157	GLY	3.2
1	A	35	LYS	3.0
1	B	141	SER	3.0
1	B	110	SER	2.8
1	B	35	LYS	2.8
1	A	93	THR	2.8
1	B	241	SER	2.6
1	A	258	HIS	2.6
1	A	257	TYR	2.6
1	A	45	LYS	2.3
1	A	159	ASP	2.2
1	B	193	ARG	2.2
1	B	37	GLY	2.2
1	A	111	GLU	2.1
1	B	65	GLN	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	182	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 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.