



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 12:02 PM UTC

PDB ID : 1Y97 / pdb_00001y97
Title : The human TREX2 3' exonuclease structure suggests a mechanism for efficient non-processive DNA catalysis
Authors : Perrino, F.W.; Harvey, S.; McMillin, S.; Hollis, T.
Deposited on : 2004-12-14
Resolution : 2.50 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

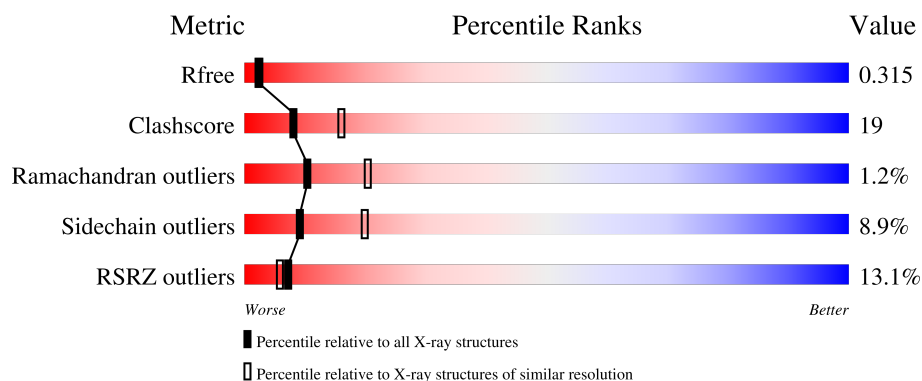
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	238	11% 62% 23% 5% 10%
1	B	238	12% 56% 27% • • 11%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3340 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 Three prime repair exonuclease 2.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	214	Total	C	N	O	S	Se	0	0	0
			1638	1044	288	297	6	3			
1	B	211	Total	C	N	O	S	Se	0	0	0
			1625	1036	282	299	6	2			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	ALA	-	cloning artifact	UNP Q9BQ50
A	0	GLY	-	cloning artifact	UNP Q9BQ50
A	1	MSE	MET	modified residue	UNP Q9BQ50
A	64	MSE	MET	modified residue	UNP Q9BQ50
A	225	MSE	MET	modified residue	UNP Q9BQ50
B	-1	ALA	-	cloning artifact	UNP Q9BQ50
B	0	GLY	-	cloning artifact	UNP Q9BQ50
B	1	MSE	MET	modified residue	UNP Q9BQ50
B	64	MSE	MET	modified residue	UNP Q9BQ50
B	225	MSE	MET	modified residue	UNP Q9BQ50

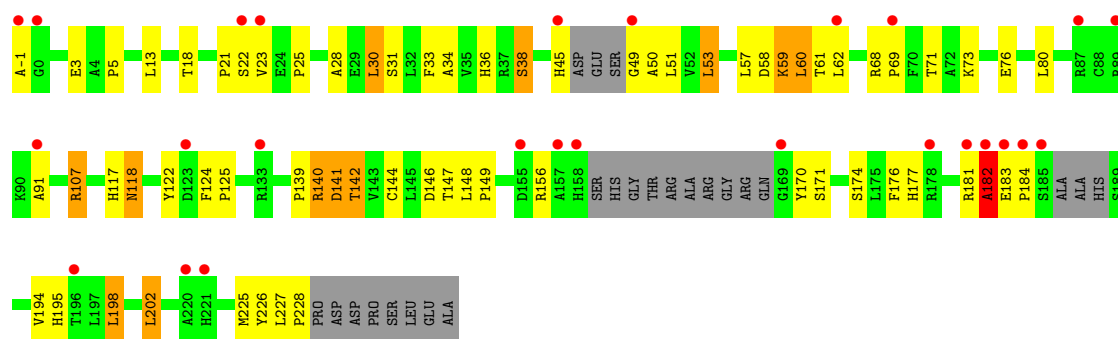
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	45	Total	O	0	0
			45	45		
2	B	32	Total	O	0	0
			32	32		

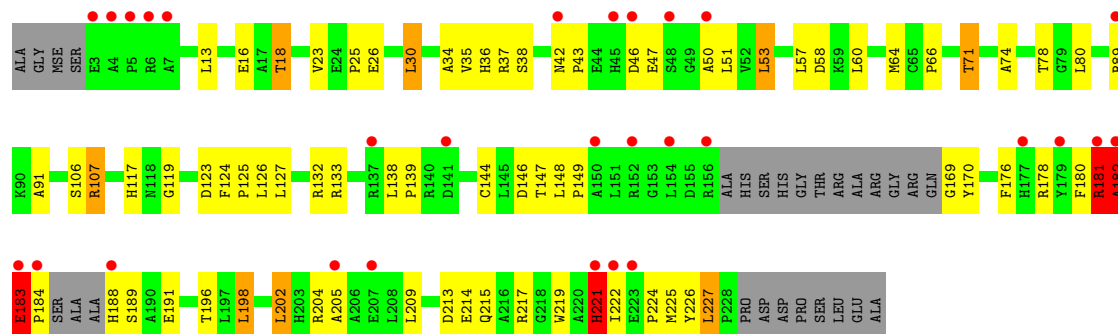
3 Residue-property plots

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: Three prime repair exonuclease 2



• Molecule 1: Three prime repair exonuclease 2



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	52.50Å 77.20Å 101.70Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.50 30.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	97.0 (30.00-2.50) 96.6 (30.00-2.50)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	53.90 (at 2.51Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.205 , 0.261 0.265 , 0.315	Depositor DCC
R_{free} test set	720 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	19.8	Xtriage
Anisotropy	0.763	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 45.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.86	EDS
Total number of atoms	3340	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.07% 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	0.84	1/1672 (0.1%)	1.17	11/2264 (0.5%)
1	B	0.92	2/1661 (0.1%)	1.38	19/2254 (0.8%)
All	All	0.88	3/3333 (0.1%)	1.28	30/4518 (0.7%)

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	B	0	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	182	ALA	C-N	-11.87	1.08	1.33
1	B	183	GLU	CA-C	-5.59	1.45	1.52
1	A	49	GLY	N-CA	5.03	1.53	1.45

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	182	ALA	O-C-N	-29.45	83.42	122.59
1	B	215	GLN	OE1-CD-NE2	-9.44	113.16	122.60
1	B	183	GLU	O-C-N	7.51	129.96	121.32
1	A	183	GLU	CA-C-N	7.16	128.79	119.84
1	A	183	GLU	C-N-CA	7.16	128.79	119.84
1	B	182	ALA	CA-C-N	7.05	139.01	121.80
1	B	182	ALA	C-N-CA	7.05	139.01	121.80
1	B	182	ALA	N-CA-C	6.72	125.11	110.80
1	B	205	ALA	N-CA-C	6.49	118.02	111.07
1	B	215	GLN	CG-CD-NE2	6.43	126.05	116.40
1	B	18	THR	N-CA-C	6.38	121.07	113.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	47	GLU	N-CA-C	-6.38	104.37	111.71
1	B	183	GLU	CA-C-O	5.69	127.96	120.16
1	B	13	LEU	N-CA-C	5.59	117.67	109.24
1	A	71	THR	N-CA-C	-5.52	102.31	110.48
1	A	171	SER	N-CA-C	-5.49	102.39	110.52
1	A	38	SER	CB-CA-C	-5.47	101.66	110.74
1	A	28	ALA	N-CA-C	-5.43	107.02	113.97
1	A	118	ASN	N-CA-C	-5.28	106.79	113.23
1	A	156	ARG	N-CA-C	-5.24	106.72	113.01
1	B	181	ARG	CA-C-N	-5.23	111.54	121.54
1	B	181	ARG	C-N-CA	-5.23	111.54	121.54
1	A	18	THR	N-CA-C	5.23	119.51	113.18
1	B	227	LEU	N-CA-C	-5.20	103.63	110.39
1	A	194	VAL	N-CA-C	-5.11	105.43	111.00
1	A	182	ALA	N-CA-C	5.10	121.67	110.80
1	B	42	ASN	N-CA-C	5.09	117.89	108.94
1	B	209	LEU	N-CA-C	-5.09	105.81	111.36
1	B	106	SER	N-CA-C	-5.04	106.23	112.38
1	B	221	HIS	CB-CA-C	5.00	119.02	110.56

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	182	ALA	Mainchain

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	1638	0	1617	56	2
1	B	1625	0	1590	74	2
2	A	45	0	0	1	0
2	B	32	0	0	1	0
All	All	3340	0	3207	126	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 19.

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:B:183:GLU:HB3	1:B:184:PRO:CD	1.34	1.40
1:B:183:GLU:CB	1:B:184:PRO:HD2	1.22	1.35
1:B:181:ARG:O	1:B:182:ALA:HB2	1.36	1.15
1:A:25:PRO:CD	1:A:225:MSE:HE1	1.84	1.08
1:A:25:PRO:HD2	1:A:225:MSE:HE1	1.14	1.07
1:B:181:ARG:O	1:B:182:ALA:CB	1.96	1.07
1:B:183:GLU:CG	1:B:184:PRO:HD2	1.84	1.06
1:B:183:GLU:CB	1:B:184:PRO:CD	2.06	1.05
1:A:25:PRO:HD2	1:A:225:MSE:CE	1.86	1.04
1:B:183:GLU:CD	1:B:184:PRO:CD	2.47	0.86
1:B:132:ARG:HH21	1:B:224:PRO:HD3	1.38	0.85
1:B:183:GLU:CD	1:B:184:PRO:HD2	2.04	0.82
1:B:180:PHE:O	1:B:181:ARG:CB	2.26	0.82
1:B:124:PHE:HB3	1:B:222:ILE:HD13	1.63	0.80
1:B:183:GLU:CD	1:B:184:PRO:HD3	2.06	0.80
1:B:126:LEU:HD13	1:B:225:MSE:CE	2.12	0.79
1:B:183:GLU:HB3	1:B:184:PRO:CG	2.14	0.77
1:B:71:THR:HG22	1:B:74:ALA:H	1.50	0.76
1:B:117:HIS:HD2	1:B:147:THR:OG1	1.71	0.73
1:A:139:PRO:O	1:A:142:THR:HG23	1.87	0.73
1:A:139:PRO:O	1:A:142:THR:CG2	2.39	0.71
1:A:36:HIS:CD2	1:A:38:SER:H	2.08	0.71
1:A:25:PRO:CG	1:A:225:MSE:HE1	2.19	0.71
1:A:34:ALA:HB3	1:A:58:ASP:HB2	1.74	0.70
1:A:140:ARG:O	1:A:141:ASP:OD1	2.09	0.70
1:B:71:THR:HG22	1:B:74:ALA:CB	2.22	0.69
1:B:225:MSE:HE3	1:B:226:TYR:HE2	1.57	0.69
1:B:169:GLY:O	1:B:178:ARG:NH2	2.24	0.69
1:A:181:ARG:O	1:A:182:ALA:HB2	1.93	0.68
1:B:217:ARG:HD3	1:B:221:HIS:HE1	1.60	0.67
1:A:117:HIS:HD2	1:A:147:THR:OG1	1.78	0.67
1:B:25:PRO:HG2	1:B:225:MSE:HE1	1.78	0.66
1:B:71:THR:CG2	1:B:74:ALA:H	2.08	0.66
1:B:30:LEU:C	1:B:30:LEU:HD23	2.23	0.64
1:A:148:LEU:HB3	1:A:149:PRO:HD3	1.79	0.64
1:B:126:LEU:HD13	1:B:225:MSE:HE2	1.79	0.64
1:B:126:LEU:HD13	1:B:225:MSE:HE1	1.79	0.64
1:B:183:GLU:HB3	1:B:184:PRO:HD2	0.64	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:148:LEU:HB3	1:B:149:PRO:HD3	1.83	0.61
1:B:217:ARG:HD3	1:B:221:HIS:CE1	2.35	0.61
1:B:35:VAL:HG22	1:B:57:LEU:CD1	2.31	0.60
1:A:73:LYS:HD2	1:A:76:GLU:OE1	2.02	0.59
1:A:5:PRO:HG3	1:A:141:ASP:HB2	1.85	0.59
1:B:35:VAL:HG22	1:B:57:LEU:HD11	1.83	0.59
1:B:225:MSE:HE3	1:B:226:TYR:CE2	2.38	0.59
1:A:30:LEU:HD23	1:A:30:LEU:O	2.03	0.59
1:A:140:ARG:CG	1:A:140:ARG:HH11	2.17	0.57
1:B:36:HIS:HD2	1:B:38:SER:OG	1.87	0.57
1:B:46:ASP:HB3	1:B:50:ALA:H	1.69	0.57
1:B:183:GLU:O	1:B:184:PRO:O	2.22	0.57
1:B:18:THR:O	1:B:71:THR:HB	2.04	0.56
1:A:45:HIS:HA	1:A:50:ALA:O	2.06	0.56
1:B:107:ARG:HD3	2:B:525:HOH:O	2.06	0.56
1:A:36:HIS:HD2	1:A:38:SER:H	1.51	0.55
1:B:182:ALA:O	1:B:183:GLU:O	2.25	0.55
1:A:140:ARG:HH11	1:A:140:ARG:HG3	1.73	0.54
1:A:51:LEU:HD21	1:A:202:LEU:HB3	1.89	0.54
1:B:36:HIS:CD2	1:B:38:SER:H	2.26	0.54
1:A:59:LYS:NZ	1:B:191:GLU:OE1	2.30	0.54
1:A:176:PHE:HD2	1:A:182:ALA:O	1.91	0.53
1:A:177:HIS:CE1	1:A:184:PRO:HD3	2.43	0.53
1:A:122:TYR:O	1:A:125:PRO:HD2	2.08	0.52
1:B:26:GLU:HB2	1:B:66:PRO:HB3	1.91	0.52
1:B:37:ARG:NH2	1:B:213:ASP:OD2	2.41	0.52
1:B:183:GLU:OE1	1:B:184:PRO:CD	2.57	0.52
1:A:198:LEU:HD22	1:A:202:LEU:HD22	1.92	0.52
1:A:33:PHE:CE1	1:A:59:LYS:HD3	2.46	0.51
1:A:228:PRO:HD3	2:A:539:HOH:O	2.10	0.51
1:B:71:THR:HG22	1:B:74:ALA:HB3	1.92	0.51
1:B:132:ARG:NH2	1:B:224:PRO:HD3	2.17	0.51
1:B:138:LEU:HD12	1:B:219:TRP:CE2	2.46	0.50
1:A:181:ARG:O	1:A:182:ALA:CB	2.59	0.49
1:A:107:ARG:NH2	1:B:91:ALA:HB3	2.28	0.49
1:B:138:LEU:O	1:B:139:PRO:C	2.55	0.49
1:B:53:LEU:HG	1:B:198:LEU:HD12	1.93	0.49
1:A:36:HIS:CD2	1:A:38:SER:OG	2.65	0.48
1:B:217:ARG:CD	1:B:221:HIS:HE1	2.25	0.48
1:A:140:ARG:CG	1:A:140:ARG:NH1	2.75	0.48
1:A:21:PRO:HA	1:A:225:MSE:CE	2.43	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:59:LYS:HE2	1:B:191:GLU:OE2	2.12	0.48
1:B:198:LEU:HD22	1:B:202:LEU:HD22	1.96	0.48
1:A:177:HIS:ND1	1:A:182:ALA:C	2.72	0.48
1:A:36:HIS:HD2	1:A:38:SER:CB	2.27	0.47
1:B:30:LEU:HD23	1:B:30:LEU:O	2.14	0.47
1:B:71:THR:HG22	1:B:74:ALA:HB2	1.97	0.47
1:B:183:GLU:OE1	1:B:184:PRO:HD3	2.12	0.47
1:A:68:ARG:HB3	1:A:69:PRO:HD2	1.96	0.47
1:A:36:HIS:HD2	1:A:38:SER:OG	1.98	0.47
1:A:225:MSE:HE2	1:A:226:TYR:CE2	2.50	0.47
1:B:144:CYS:N	1:B:217:ARG:O	2.48	0.46
1:A:30:LEU:HD22	1:A:62:LEU:HB2	1.96	0.46
1:B:117:HIS:HE1	1:B:170:TYR:O	1.99	0.46
1:A:3:GLU:HA	1:A:3:GLU:OE1	2.16	0.46
1:A:13:LEU:C	1:A:13:LEU:HD12	2.40	0.45
1:A:23:VAL:HG12	1:A:23:VAL:O	2.15	0.45
1:A:225:MSE:HE2	1:A:226:TYR:CD2	2.51	0.45
1:B:71:THR:HG22	1:B:74:ALA:N	2.26	0.45
1:B:117:HIS:CD2	1:B:147:THR:OG1	2.61	0.45
1:B:43:PRO:HB3	1:B:51:LEU:HD21	1.98	0.45
1:A:146:ASP:CG	1:A:149:PRO:HD2	2.42	0.45
1:B:181:ARG:O	1:B:182:ALA:HB3	2.07	0.45
1:A:91:ALA:HB3	1:B:107:ARG:NH2	2.32	0.45
1:A:124:PHE:HB2	1:A:125:PRO:HD3	1.99	0.44
1:B:46:ASP:HB2	1:B:50:ALA:O	2.18	0.44
1:A:31:SER:HA	1:A:60:LEU:O	2.18	0.44
1:B:124:PHE:CE1	1:B:144:CYS:HB3	2.53	0.44
1:A:30:LEU:O	1:A:61:THR:HA	2.17	0.44
1:A:21:PRO:HA	1:A:225:MSE:HE3	2.01	0.43
1:B:202:LEU:HD13	1:B:202:LEU:HA	1.89	0.42
1:B:37:ARG:HH22	1:B:213:ASP:CG	2.26	0.42
1:B:146:ASP:O	1:B:149:PRO:HD2	2.20	0.42
1:B:225:MSE:HG2	1:B:226:TYR:HD2	1.85	0.42
1:B:34:ALA:HB3	1:B:58:ASP:HB2	2.02	0.42
1:A:176:PHE:CD2	1:A:182:ALA:O	2.70	0.41
1:A:53:LEU:HD21	1:A:195:HIS:CD2	2.56	0.41
1:B:132:ARG:HH21	1:B:224:PRO:CD	2.22	0.41
1:A:30:LEU:HD23	1:A:30:LEU:C	2.45	0.41
1:A:117:HIS:HE1	1:A:170:TYR:O	2.03	0.41
1:A:124:PHE:CE1	1:A:144:CYS:HB3	2.55	0.41
1:A:117:HIS:CD2	1:A:147:THR:OG1	2.65	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:89:ARG:HG3	1:B:89:ARG:NH1	2.36	0.41
1:B:16:GLU:OE1	1:B:188:HIS:N	2.54	0.40
1:B:74:ALA:O	1:B:78:THR:HG23	2.22	0.40
1:B:119:GLY:O	1:B:125:PRO:HD2	2.21	0.40
1:A:117:HIS:O	1:A:118:ASN:HB3	2.22	0.40
1:B:176:PHE:CE2	1:B:183:GLU:HA	2.57	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:22:SER:OG	1:B:214:GLU:OE2[2_764]	2.04	0.16
1:A:-1:ALA:N	1:B:123:ASP:OD1[2_664]	2.05	0.15

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	206/238 (87%)	197 (96%)	8 (4%)	1 (0%)	24	43
1	B	205/238 (86%)	196 (96%)	5 (2%)	4 (2%)	6	10
All	All	411/476 (86%)	393 (96%)	13 (3%)	5 (1%)	10	20

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	181	ARG
1	B	182	ALA
1	A	182	ALA
1	B	189	SER
1	B	183	GLU

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	169/187 (90%)	155 (92%)	14 (8%)	10	22
1	B	169/187 (90%)	153 (90%)	16 (10%)	8	17
All	All	338/374 (90%)	308 (91%)	30 (9%)	9	20

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30	LEU
1	A	53	LEU
1	A	57	LEU
1	A	59	LYS
1	A	60	LEU
1	A	80	LEU
1	A	107	ARG
1	A	140	ARG
1	A	141	ASP
1	A	142	THR
1	A	174	SER
1	A	198	LEU
1	A	202	LEU
1	A	227	LEU
1	B	23	VAL
1	B	30	LEU
1	B	53	LEU
1	B	60	LEU
1	B	64	MSE
1	B	71	THR
1	B	80	LEU
1	B	107	ARG
1	B	127	LEU
1	B	133	ARG
1	B	196	THR
1	B	198	LEU
1	B	202	LEU

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Mol	Chain	Res	Type
1	B	204	ARG
1	B	221	HIS
1	B	227	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	36	HIS
1	A	117	HIS
1	A	158	HIS
1	B	36	HIS
1	B	42	ASN
1	B	45	HIS
1	B	108	GLN
1	B	117	HIS
1	B	221	HIS

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

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	B	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	B	182:ALA	C	183:GLU	N	1.09

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	211/238 (88%)	1.17	26 (12%) 8 7	8, 17, 44, 76	0
1	B	209/238 (87%)	1.18	29 (13%) 6 5	10, 18, 45, 82	0
All	All	420/476 (88%)	1.18	55 (13%) 7 6	8, 18, 47, 82	0

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	184	PRO	6.6
1	A	-1	ALA	6.1
1	A	185	SER	5.4
1	A	183	GLU	4.4
1	A	182	ALA	4.3
1	B	188	HIS	3.9
1	B	5	PRO	3.8
1	A	87	ARG	3.5
1	B	46	ASP	3.5
1	B	182	ALA	3.5
1	B	183	GLU	3.4
1	A	45	HIS	3.2
1	A	184	PRO	3.2
1	A	49	GLY	3.1
1	A	155	ASP	2.9
1	A	181	ARG	2.9
1	A	91	ALA	2.8
1	A	23	VAL	2.8
1	B	3	GLU	2.8
1	B	152	ARG	2.8
1	B	45	HIS	2.7
1	B	181	ARG	2.7
1	A	178	ARG	2.6
1	B	156	ARG	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	158	HIS	2.6
1	A	89	ARG	2.5
1	A	0	GLY	2.5
1	B	50	ALA	2.5
1	A	157	ALA	2.5
1	B	89	ARG	2.4
1	A	196	THR	2.4
1	B	4	ALA	2.4
1	B	6	ARG	2.4
1	B	207	GLU	2.4
1	B	48	SER	2.4
1	B	222	ILE	2.3
1	B	137	ARG	2.2
1	A	62	LEU	2.2
1	A	133	ARG	2.2
1	B	205	ALA	2.2
1	A	221	HIS	2.2
1	B	42	ASN	2.2
1	A	169	GLY	2.2
1	B	177	HIS	2.1
1	B	179	TYR	2.1
1	A	22	SER	2.1
1	A	123	ASP	2.1
1	B	7	ALA	2.1
1	B	150	ALA	2.1
1	B	154	LEU	2.1
1	B	223	GLU	2.1
1	A	69	PRO	2.1
1	A	220	ALA	2.0
1	B	141	ASP	2.0
1	B	221	HIS	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.