



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 10:00 AM UTC

PDB ID : 7TQQ / pdb_00007tqq
Title : Structure of human TREX1-DNA complex
Authors : Zhou, W.; Richmond-Buccola, D.; Kranzusch, P.J.
Deposited on : 2022-01-26
Resolution : 2.20 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

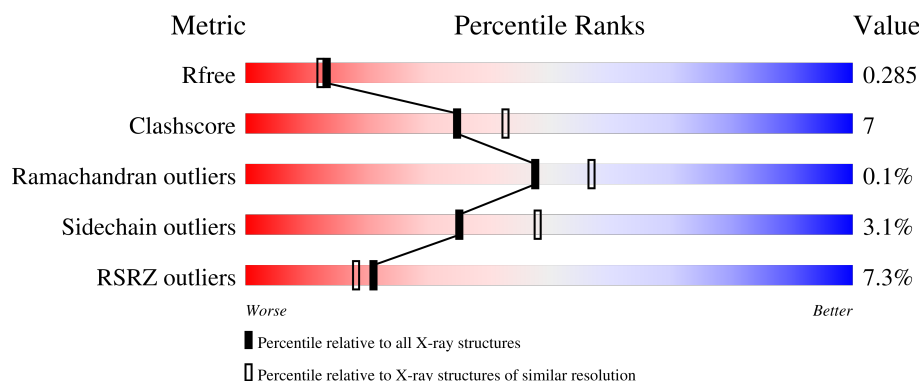
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	243	10% 75% 11% • 13%
1	B	243	% 76% 11% • 12%
1	E	243	4% 75% 13% 12%
1	F	243	8% 78% 10% 12%
2	C	22	14% 41% 32% 27%

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Mol	Chain	Length	Quality of chain
2	D	22	9% 45% 23% 32%
2	G	22	14% 36% 23% 41%
2	H	22	14% 36% 36% 27%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 8139 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 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	212	Total	C	N	O	S	0	0	0
			1608	1027	277	292	12			
1	B	213	Total	C	N	O	S	0	1	0
			1621	1038	279	291	13			
1	E	214	Total	C	N	O	S	0	0	0
			1631	1043	283	293	12			
1	F	215	Total	C	N	O	S	0	0	0
			1631	1043	281	295	12			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	SER	-	expression tag	UNP Q9NSU2
B	0	SER	-	expression tag	UNP Q9NSU2
E	0	SER	-	expression tag	UNP Q9NSU2
F	0	SER	-	expression tag	UNP Q9NSU2

- Molecule 2 is a DNA chain called DNA (5'-D(P*TP*TP*GP*GP*CP*CP*GP*GP*CP*CP*AP*TP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	16	Total	C	N	O	P	0	0	0
			327	154	59	98	16			
2	D	15	Total	C	N	O	P	0	0	0
			305	144	54	92	15			
2	G	13	Total	C	N	O	P	0	0	0
			261	124	44	80	13			
2	H	16	Total	C	N	O	P	0	0	0
			327	154	59	98	16			

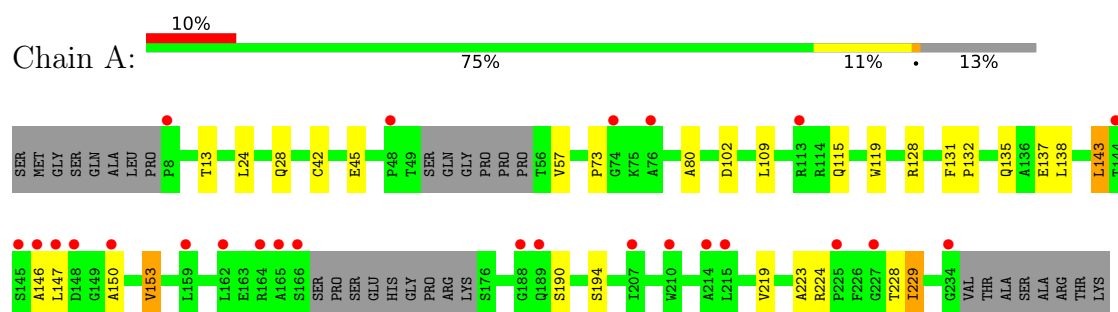
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	65	Total O 65 65	0	0
3	B	126	Total O 126 126	0	0
3	C	13	Total O 13 13	0	0
3	D	21	Total O 21 21	0	0
3	E	91	Total O 91 91	0	0
3	F	83	Total O 83 83	0	0
3	G	13	Total O 13 13	0	0
3	H	16	Total O 16 16	0	0

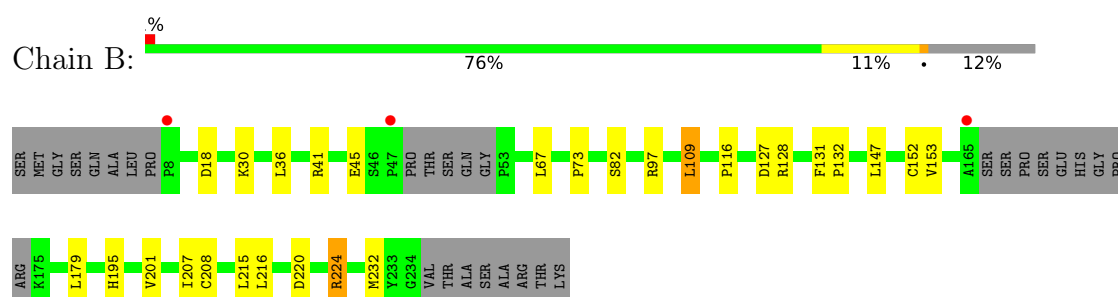
3 Residue-property plots

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.

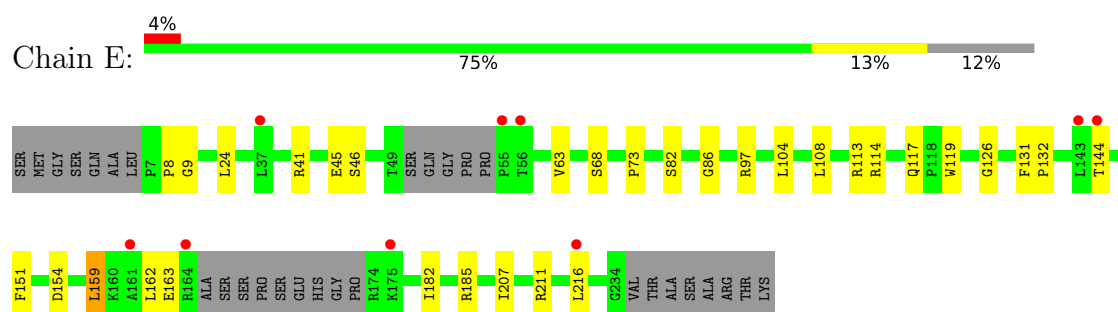
- Molecule 1: Three-prime repair exonuclease 1



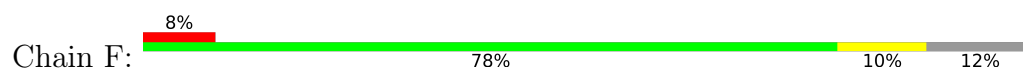
- Molecule 1: Three-prime repair exonuclease 1

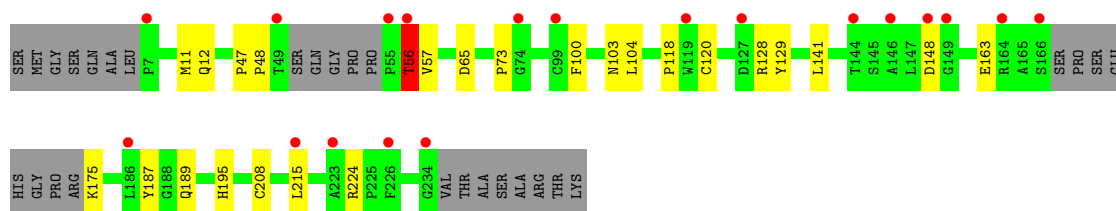


- Molecule 1: Three-prime repair exonuclease 1

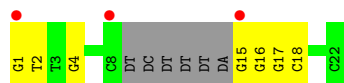
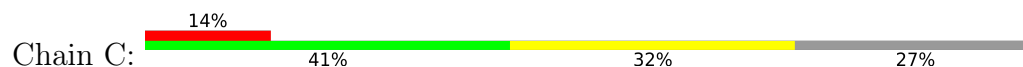


- Molecule 1: Three-prime repair exonuclease 1

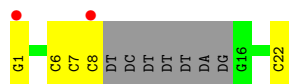




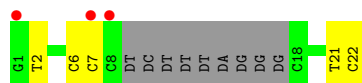
- Molecule 2: DNA (5'-D(P*TP*TP*GP*GP*CP*CP*GP*GP*CP*CP*AP*TP*C)-3')



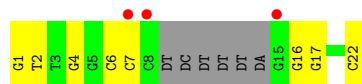
- Molecule 2: DNA (5'-D(P*TP*TP*GP*GP*CP*CP*GP*GP*CP*CP*AP*TP*C)-3')



- Molecule 2: DNA (5'-D(P*TP*TP*GP*GP*CP*CP*GP*GP*CP*CP*AP*TP*C)-3')



- Molecule 2: DNA (5'-D(P*TP*TP*GP*GP*CP*CP*GP*GP*CP*CP*AP*TP*C)-3')



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	43.68Å 91.36Å 157.39Å 90.00° 90.71° 90.00°	Depositor
Resolution (Å)	45.68 – 2.20 45.68 – 2.20	Depositor EDS
% Data completeness (in resolution range)	98.1 (45.68-2.20) 98.1 (45.68-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.26 (at 2.20Å)	Xtriage
Refinement program	PHENIX 1.19.1_4122	Depositor
R, R_{free}	0.226 , 0.261 (Not available) , 0.285	Depositor DCC
R_{free} test set	2003 reflections (3.19%)	wwPDB-VP
Wilson B-factor (Å ²)	37.9	Xtriage
Anisotropy	0.192	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 52.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	0.165 for h,-k,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	8139	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.59% 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	0.21	0/1649	0.38	0/2251
1	B	0.22	0/1667	0.42	0/2275
1	E	0.22	0/1674	0.43	0/2284
1	F	0.22	0/1674	0.42	0/2285
2	C	0.22	0/364	0.37	0/557
2	D	0.33	0/339	0.50	0/518
2	G	0.35	0/289	0.57	0/440
2	H	0.38	0/364	0.51	0/557
All	All	0.24	0/8020	0.43	0/11167

There are no bond length outliers.

There are no bond angle outliers.

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	1608	0	1609	18	0
1	B	1621	0	1630	20	0
1	E	1631	0	1640	24	0
1	F	1631	0	1637	15	0
2	C	327	0	181	9	0
2	D	305	0	170	10	0
2	G	261	0	148	9	0
2	H	327	0	181	10	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	65	0	0	8	1
3	B	126	0	0	5	0
3	C	13	0	0	0	0
3	D	21	0	0	0	1
3	E	91	0	0	7	0
3	F	83	0	0	5	0
3	G	13	0	0	1	0
3	H	16	0	0	0	0
All	All	8139	0	7196	107	1

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 7.

All (107) 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:E:113:ARG:NH2	3:E:301:HOH:O	1.92	1.00
2:H:1:DG:H2''	2:H:2:DT:H5'	1.59	0.85
1:A:57:VAL:O	3:A:301:HOH:O	1.94	0.85
1:B:128:ARG:HH22	2:D:6:DC:H5'	1.44	0.82
1:E:9:GLY:O	3:E:302:HOH:O	2.00	0.78
1:E:126:GLY:N	1:E:154:ASP:OD1	2.18	0.76
1:B:195:HIS:ND1	3:B:303:HOH:O	2.19	0.76
1:E:73:PRO:O	3:E:303:HOH:O	2.03	0.75
1:A:73:PRO:O	3:A:302:HOH:O	2.04	0.75
1:E:86:GLY:O	3:E:304:HOH:O	2.07	0.72
1:B:41:ARG:NH2	1:B:220:ASP:OD2	2.23	0.71
1:B:73:PRO:O	3:B:302:HOH:O	2.10	0.69
1:E:159:LEU:HD23	1:E:182:ILE:HD13	1.75	0.68
1:E:63:VAL:O	3:E:305:HOH:O	2.11	0.67
1:E:46:SER:OG	3:E:306:HOH:O	2.13	0.67
2:C:2:DT:H5'	2:C:2:DT:H6	1.58	0.66
1:F:73:PRO:O	3:F:303:HOH:O	2.13	0.65
2:G:7:DC:N3	3:G:101:HOH:O	2.28	0.65
1:B:128:ARG:NH2	2:D:6:DC:H5'	2.11	0.64
2:G:22:DC:H5''	2:G:22:DC:H6	1.63	0.63
1:F:195:HIS:ND1	3:F:301:HOH:O	1.81	0.63
2:C:1:DG:H8	2:C:1:DG:H5''	1.64	0.63
2:H:22:DC:H6	2:H:22:DC:H5''	1.64	0.62
1:A:147:LEU:N	3:A:306:HOH:O	2.27	0.61
2:H:6:DC:H42	2:H:17:DG:H1	1.49	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:17:DG:C8	2:H:17:DG:H5'	2.36	0.60
2:D:7:DC:H2''	2:D:8:DC:H5'	1.84	0.60
1:E:8:PRO:HG3	1:E:151:PHE:CZ	2.36	0.60
1:F:224:ARG:HD3	3:F:333:HOH:O	2.02	0.59
2:D:22:DC:H6	2:D:22:DC:H5''	1.67	0.59
1:E:45:GLU:HG3	1:E:216:LEU:HD12	1.85	0.58
1:A:224:ARG:NH2	3:A:308:HOH:O	2.36	0.58
2:G:6:DC:H1'	2:G:7:DC:H4'	1.84	0.58
2:C:15:DG:H1'	2:C:16:DG:H5'	1.86	0.57
1:B:127:ASP:OD2	1:B:224:ARG:NH2	2.37	0.56
2:C:1:DG:H2'	2:C:2:DT:H5''	1.88	0.55
1:E:41:ARG:NH2	1:E:45:GLU:OE2	2.38	0.55
1:E:162:LEU:O	1:E:211:ARG:NH1	2.40	0.55
2:G:6:DC:C2	2:G:7:DC:H1'	2.42	0.54
1:A:24:LEU:HD13	2:C:2:DT:C2	2.44	0.53
2:D:22:DC:H5''	2:D:22:DC:C6	2.44	0.52
2:H:16:DG:H2''	2:H:17:DG:C8	2.44	0.52
1:E:24:LEU:HD13	2:G:2:DT:C2	2.45	0.51
1:F:129:TYR:HB3	3:F:321:HOH:O	2.10	0.51
2:H:22:DC:H5''	2:H:22:DC:C6	2.45	0.51
1:B:195:HIS:NE2	3:B:301:HOH:O	1.96	0.50
1:F:208:CYS:HB3	1:F:215:LEU:HD22	1.93	0.50
1:F:100:PHE:HA	1:F:104:LEU:HD23	1.94	0.50
1:A:146:ALA:N	3:A:306:HOH:O	2.45	0.49
1:B:179:LEU:HD12	1:B:207:ILE:HD12	1.95	0.48
1:E:68:SER:OG	3:E:307:HOH:O	2.20	0.48
2:D:7:DC:OP2	2:D:7:DC:H2'	2.14	0.48
1:B:131:PHE:CZ	1:B:152:CYS:HB2	2.49	0.47
1:E:114:ARG:HD2	3:F:311:HOH:O	2.14	0.47
1:A:153:VAL:HG11	1:A:219:VAL:HG22	1.96	0.47
1:B:208:CYS:HB3	1:B:215:LEU:HD22	1.96	0.47
2:G:22:DC:H6	2:G:22:DC:C5'	2.27	0.47
1:E:163:GLU:CD	1:E:185:ARG:HH12	2.22	0.46
2:H:7:DC:H42	2:H:16:DG:H1	1.64	0.46
1:B:36:LEU:HD12	1:B:67:LEU:HD23	1.98	0.46
2:C:16:DG:H2''	2:C:17:DG:C8	2.50	0.46
2:D:6:DC:H2''	2:D:7:DC:C6	2.50	0.46
1:E:45:GLU:HG3	1:E:216:LEU:CD1	2.45	0.46
2:C:4:DG:O6	2:C:18:DC:N4	2.49	0.45
1:E:131:PHE:HB2	1:E:132:PRO:HD3	1.98	0.45
1:F:12:GLN:HG2	1:F:118:PRO:HD2	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:135:GLN:NE2	3:A:313:HOH:O	2.49	0.45
2:G:6:DC:C1'	2:G:7:DC:H4'	2.46	0.45
1:E:216:LEU:HD23	1:E:216:LEU:HA	1.78	0.44
1:B:132:PRO:HB2	1:B:232:MET:HB2	1.99	0.44
1:E:117:GLN:HG2	1:E:119:TRP:CE2	2.52	0.44
1:A:131:PHE:HB2	1:A:132:PRO:HD3	1.99	0.44
2:G:6:DC:H2''	2:G:7:DC:OP2	2.17	0.44
1:A:13:THR:HG21	1:A:115:GLN:HB3	1.99	0.44
1:B:109:LEU:HD13	1:B:109:LEU:HA	1.86	0.44
1:E:104:LEU:O	1:E:108:LEU:HG	2.18	0.44
2:H:6:DC:H2''	2:H:7:DC:OP1	2.16	0.44
1:A:137:GLU:OE1	3:A:304:HOH:O	2.21	0.43
1:A:119:TRP:HE3	1:A:150:ALA:HB2	1.83	0.43
1:B:97:ARG:NH2	3:B:314:HOH:O	2.51	0.43
1:B:116:PRO:HD2	3:B:308:HOH:O	2.18	0.43
2:D:7:DC:H2''	2:D:8:DC:H3'	2.01	0.43
1:F:187:TYR:O	1:F:189:GLN:HG2	2.18	0.43
1:B:30:LYS:HE3	1:B:30:LYS:HB3	1.90	0.43
1:B:41:ARG:NH1	1:B:45:GLU:OE2	2.52	0.43
2:C:2:DT:H5'	2:C:2:DT:C6	2.46	0.43
1:A:224:ARG:NH1	1:A:228:THR:OG1	2.46	0.43
1:F:128:ARG:NH2	2:H:4:DG:H4'	2.34	0.43
1:F:47:PRO:HA	1:F:48:PRO:HD3	1.93	0.43
1:A:137:GLU:HA	3:A:304:HOH:O	2.19	0.42
1:B:18:ASP:HB3	1:B:201:VAL:HG23	2.01	0.42
1:B:216:LEU:HD23	1:B:216:LEU:HA	1.77	0.42
1:F:11:MET:SD	1:F:120:CYS:HB2	2.59	0.42
1:E:113:ARG:NH1	1:F:103:ASN:OD1	2.53	0.42
1:A:80:ALA:HB3	2:C:1:DG:C2	2.55	0.42
1:A:138:LEU:HB3	1:A:143:LEU:O	2.20	0.42
2:D:7:DC:C2'	2:D:8:DC:H5'	2.47	0.41
1:E:159:LEU:HD21	1:E:207:ILE:HG21	2.01	0.41
1:F:56:THR:HG23	1:F:57:VAL:H	1.85	0.41
2:G:21:DT:H2'	2:G:22:DC:H5''	2.01	0.41
1:A:131:PHE:HB3	1:A:229:ILE:HD12	2.01	0.41
1:B:147:LEU:HD23	1:B:147:LEU:HA	1.75	0.41
2:D:1:DG:H8	2:D:1:DG:H5''	1.86	0.41
1:A:219:VAL:O	1:A:223:ALA:N	2.53	0.41
2:H:17:DG:H5'	2:H:17:DG:H8	1.85	0.41
1:E:97:ARG:NH1	1:F:65:ASP:OD2	2.55	0.40
1:F:100:PHE:HD1	1:F:141:LEU:HD11	1.87	0.40

All (1) 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)
3:A:333:HOH:O	3:D:110:HOH:O[2_545]	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	206/243 (85%)	203 (98%)	3 (2%)	0	100	100
1	B	208/243 (86%)	203 (98%)	5 (2%)	0	100	100
1	E	208/243 (86%)	203 (98%)	5 (2%)	0	100	100
1	F	209/243 (86%)	205 (98%)	3 (1%)	1 (0%)	24	27
All	All	831/972 (86%)	814 (98%)	16 (2%)	1 (0%)	48	57

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	F	56	THR

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	174/199 (87%)	163 (94%)	11 (6%)	16	19
1	B	176/199 (88%)	172 (98%)	4 (2%)	44	59

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	E	177/199 (89%)	174 (98%)	3 (2%)	53	69
1	F	177/199 (89%)	173 (98%)	4 (2%)	44	59
All	All	704/796 (88%)	682 (97%)	22 (3%)	35	48

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	GLN
1	A	42	CYS
1	A	45	GLU
1	A	102	ASP
1	A	109	LEU
1	A	128	ARG
1	A	143	LEU
1	A	153	VAL
1	A	190	SER
1	A	194	SER
1	A	229	ILE
1	B	82	SER
1	B	109	LEU
1	B	153	VAL
1	B	224	ARG
1	E	82	SER
1	E	144	THR
1	E	159	LEU
1	F	56	THR
1	F	148	ASP
1	F	163	GLU
1	F	175	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	103	ASN
1	A	124	HIS
1	B	12	GLN
1	E	12	GLN
1	E	103	ASN
1	F	12	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	212/243 (87%)	0.86	25 (11%) 9 7	30, 46, 73, 95	0
1	B	213/243 (87%)	0.40	3 (1%) 73 71	26, 36, 55, 69	1 (0%)
1	E	214/243 (88%)	0.57	9 (4%) 40 37	28, 39, 61, 78	0
1	F	215/243 (88%)	0.66	19 (8%) 15 13	29, 39, 66, 80	0
2	C	16/22 (72%)	1.08	3 (18%) 3 2	41, 96, 143, 145	0
2	D	15/22 (68%)	0.83	2 (13%) 7 5	34, 54, 119, 120	0
2	G	13/22 (59%)	1.14	3 (23%) 2 1	38, 89, 133, 142	0
2	H	16/22 (72%)	0.85	3 (18%) 3 2	37, 73, 143, 149	0
All	All	914/1060 (86%)	0.64	67 (7%) 21 18	26, 40, 75, 149	1 (0%)

All (67) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	C	1	DG	4.2
1	E	164	ARG	4.0
1	E	175	LYS	3.6
1	A	234	GLY	3.6
1	F	144	THR	3.6
1	A	144	THR	3.5
1	B	165	ALA	3.3
1	E	161	ALA	3.2
1	F	146	ALA	3.2
2	D	8	DC	3.2
2	G	8	DC	3.0
1	A	165	ALA	3.0
1	F	223	ALA	2.9
1	F	119	TRP	2.9
1	A	150	ALA	2.9
1	A	188	GLY	2.8

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Mol	Chain	Res	Type	RSRZ
1	F	74	GLY	2.8
1	A	189	GLN	2.7
1	A	166	SER	2.7
1	B	8	PRO	2.7
1	E	56	THR	2.6
1	A	147	LEU	2.6
1	E	143	LEU	2.6
1	F	49	THR	2.6
1	A	8	PRO	2.6
1	F	234	GLY	2.6
1	A	207	ILE	2.6
1	A	162	LEU	2.5
1	F	56	THR	2.5
2	H	7	DC	2.5
1	A	148	ASP	2.5
1	A	74	GLY	2.5
1	F	7	PRO	2.5
1	A	145	SER	2.5
1	E	144	THR	2.5
2	C	8	DC	2.5
1	F	149	GLY	2.3
1	F	164	ARG	2.3
1	A	225	PRO	2.3
2	D	1	DG	2.3
1	A	146	ALA	2.3
1	A	164	ARG	2.3
1	F	55	PRO	2.2
1	F	215	LEU	2.2
1	A	210	TRP	2.2
2	H	8	DC	2.2
2	G	1	DG	2.2
1	E	37	LEU	2.2
1	F	226	PHE	2.2
1	F	99	CYS	2.2
1	A	113	ARG	2.1
1	E	216	LEU	2.1
1	A	48	PRO	2.1
2	H	15	DG	2.1
1	A	215	LEU	2.1
1	F	186	LEU	2.1
1	E	55	PRO	2.1
1	A	227	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	76	ALA	2.1
1	A	214	ALA	2.1
2	C	15	DG	2.1
1	F	148	ASP	2.1
1	B	47	PRO	2.1
1	F	166	SER	2.0
2	G	7	DC	2.0
1	A	159	LEU	2.0
1	F	127	ASP	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.