



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 9, 2026 – 01:48 PM UTC

PDB ID : 3QRF / pdb_00003qrf
Title : Structure of a domain-swapped FOXP3 dimer
Authors : Bandukwala, H.S.; Wu, Y.; Feurer, M.; Chen, Y.; Barbosa, B.; Ghosh, S.;
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Deposited on : 2011-02-17
Resolution : 2.80 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

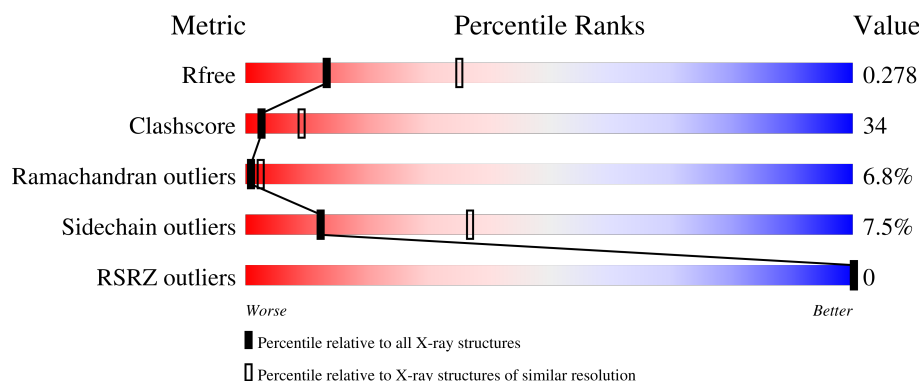
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

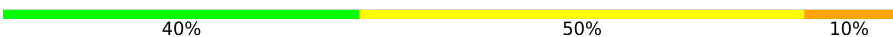
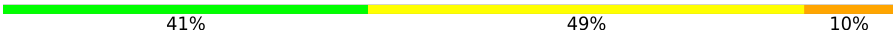
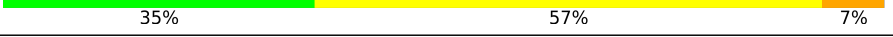
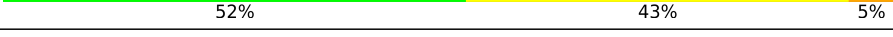
The reported resolution of this entry is 2.80 Å.

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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	M	286	
1	N	286	
2	F	82	
2	G	82	
2	H	82	

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Mol	Chain	Length	Quality of chain
2	I	82	52%43%5%
3	A	21	24%71%5%
3	C	21	19%76%5%
4	B	21	5%86%10%
4	D	21	5%86%10%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 9105 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 Nuclear factor of activated T-cells, cytoplasmic 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	N	286	Total	C	N	O	S	0	0	0
			2264	1421	413	421	9			
1	M	286	Total	C	N	O	S	0	0	0
			2264	1421	413	421	9			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
N	393	SER	-	expression tag	UNP Q13469
N	394	SER	-	expression tag	UNP Q13469
N	395	VAL	-	expression tag	UNP Q13469
N	629	ALA	ASP	conflict	UNP Q13469
N	630	ALA	LYS	conflict	UNP Q13469
N	631	ALA	ASP	conflict	UNP Q13469
N	632	ALA	LYS	conflict	UNP Q13469
N	633	ALA	SER	conflict	UNP Q13469
N	634	ALA	GLN	conflict	UNP Q13469
M	393	SER	-	expression tag	UNP Q13469
M	394	SER	-	expression tag	UNP Q13469
M	395	VAL	-	expression tag	UNP Q13469
M	629	ALA	ASP	conflict	UNP Q13469
M	630	ALA	LYS	conflict	UNP Q13469
M	631	ALA	ASP	conflict	UNP Q13469
M	632	ALA	LYS	conflict	UNP Q13469
M	633	ALA	SER	conflict	UNP Q13469
M	634	ALA	GLN	conflict	UNP Q13469

- Molecule 2 is a protein called Forkhead box protein P3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	F	82	Total	C	N	O	S	0	0	0
			714	465	132	114	3			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	G	82	Total	C	N	O	S	0	0	0
			714	465	132	114	3			
2	H	82	Total	C	N	O	S	0	0	0
			714	465	132	114	3			
2	I	82	Total	C	N	O	S	0	0	0
			714	465	132	114	3			

- Molecule 3 is a DNA chain called human hARRE2 DNA (Plus Strand).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	21	Total	C	N	O	P	0	0	0
			431	209	76	126	20			
3	A	21	Total	C	N	O	P	0	0	0
			431	209	76	126	20			

- Molecule 4 is a DNA chain called human hARRE2 DNA (Minus Strand).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	21	Total	C	N	O	P	0	0	0
			424	206	76	122	20			
4	B	21	Total	C	N	O	P	0	0	0
			424	206	76	122	20			

- Molecule 5 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	F	1	Total	Mg	0	0
			1	1		
5	G	1	Total	Mg	0	0
			1	1		
5	H	1	Total	Mg	0	0
			1	1		
5	I	1	Total	Mg	0	0
			1	1		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	N	1	Total	O	0	0
			1	1		
6	C	2	Total	O	0	0
			2	2		

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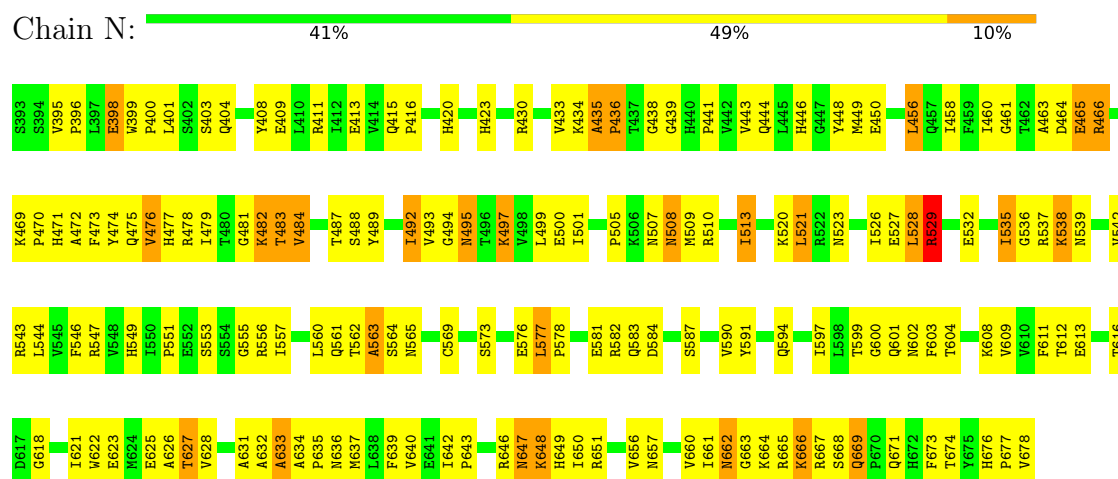
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	D	3	Total	O	0	0
			3	3		
6	B	1	Total	O	0	0
			1	1		

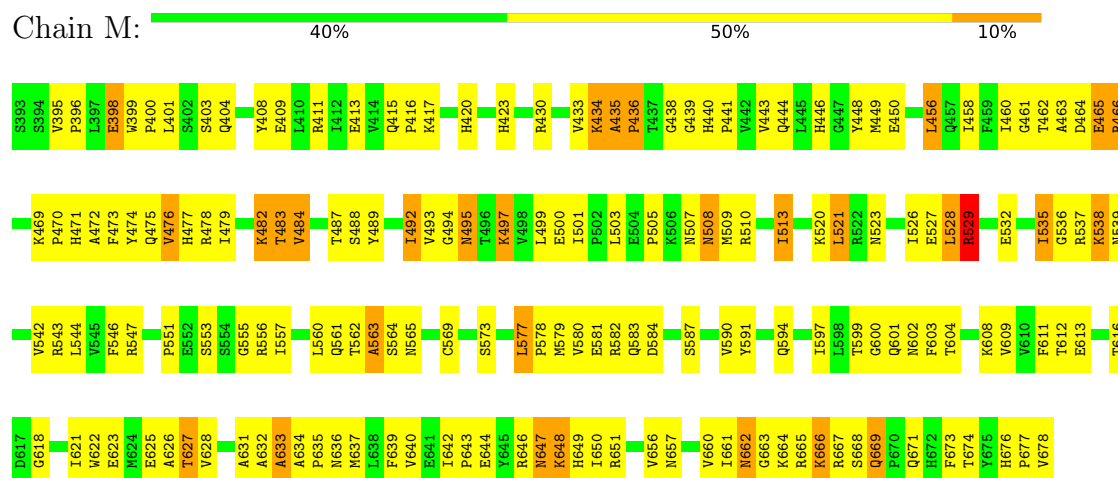
3 Residue-property plots

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: Nuclear factor of activated T-cells, cytoplasmic 2

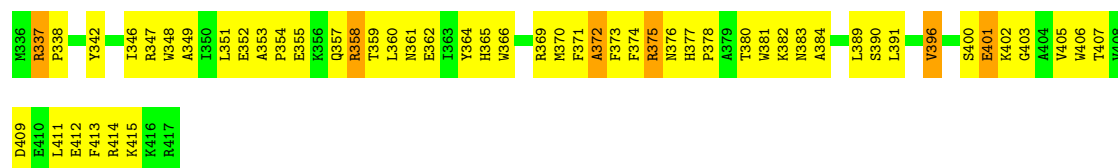


- Molecule 1: Nuclear factor of activated T-cells, cytoplasmic 2



- Molecule 2: Forkhead box protein P3

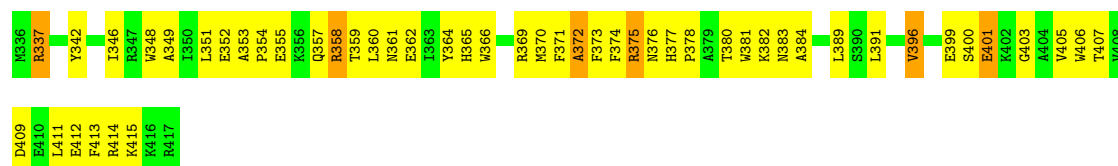




- Molecule 2: Forkhead box protein P3



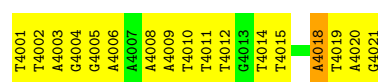
- Molecule 2: Forkhead box protein P3



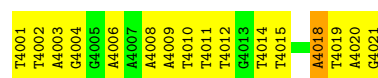
- Molecule 2: Forkhead box protein P3



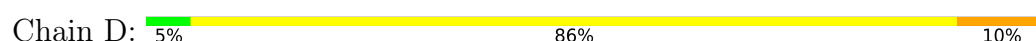
- Molecule 3: human hARRE2 DNA (Plus Strand)



- Molecule 3: human hARRE2 DNA (Plus Strand)



- Molecule 4: human hARRE2 DNA (Minus Strand)



- | |
|-------|
| A5001 |
| A5002 |
| C5003 |
| T5004 |
| A5005 |
| T5006 |
| G5007 |
| A5008 |
| A5009 |
| A5010 |
| C5011 |
| A5012 |
| A5013 |
| A5014 |
| T5015 |
| T5016 |
| T5017 |
| T5018 |
| C5019 |
| C5020 |
| T5021 |

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	83.67Å 131.23Å 68.67Å 90.00° 89.95° 90.00°	Depositor
Resolution (Å)	50.00 – 2.80 50.00 – 2.80	Depositor EDS
% Data completeness (in resolution range)	(Not available) (50.00-2.80) 96.0 (50.00-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.68 (at 2.79Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.248 , 0.283 0.244 , 0.278	Depositor DCC
R_{free} test set	3538 reflections (9.74%)	wwPDB-VP
Wilson B-factor (Å ²)	63.7	Xtriage
Anisotropy	0.645	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.23 , 30.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.480 for -h,-k,l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	9105	wwPDB-VP
Average B, all atoms (Å ²)	65.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.89% 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

Bond lengths and bond angles in the following residue types are not validated in this section: MG

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	M	0.45	0/2314	0.99	5/3134 (0.2%)
1	N	0.46	0/2314	0.99	6/3134 (0.2%)
2	F	0.56	0/738	0.94	1/996 (0.1%)
2	G	0.55	0/738	0.89	2/996 (0.2%)
2	H	0.57	0/738	0.94	1/996 (0.1%)
2	I	0.53	0/738	0.90	2/996 (0.2%)
3	A	0.37	0/483	0.95	1/745 (0.1%)
3	C	0.35	0/483	0.95	1/745 (0.1%)
4	B	0.37	0/475	0.87	0/730
4	D	0.37	0/475	0.86	0/730
All	All	0.47	0/9496	0.95	19/13202 (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
3	A	0	1
3	C	0	1
4	B	0	2
4	D	0	2
All	All	0	6

There are no bond length outliers.

The worst 5 of 19 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	M	435	ALA	CA-C-N	6.69	128.20	119.84
1	M	435	ALA	C-N-CA	6.69	128.20	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	N	435	ALA	CA-C-N	6.62	128.12	119.84
1	N	435	ALA	C-N-CA	6.62	128.12	119.84
3	A	4006	DA	N9-C1'-C2'	6.19	122.78	113.50

There are no chirality outliers.

5 of 6 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	A	4018	DA	Sidechain
4	B	5004	DT	Sidechain
3	C	4018	DA	Sidechain
4	D	5004	DT	Sidechain
4	D	5005	DA	Sidechain

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	M	2264	0	2271	176	0
1	N	2264	0	2271	170	0
2	F	714	0	708	56	0
2	G	714	0	708	42	0
2	H	714	0	708	54	0
2	I	714	0	708	42	0
3	A	431	0	242	24	0
3	C	431	0	242	25	0
4	B	424	0	240	24	0
4	D	424	0	240	25	0
5	F	1	0	0	0	0
5	G	1	0	0	0	0
5	H	1	0	0	0	0
5	I	1	0	0	0	0
6	B	1	0	0	0	0
6	C	2	0	0	0	0
6	D	3	0	0	0	0
6	N	1	0	0	0	0
All	All	9105	0	8338	593	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 34.

The worst 5 of 593 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:N:669:GLN:HE21	1:N:671:GLN:HE22	1.07	0.98
1:M:634:ALA:HB3	1:M:637:MET:HB2	1.44	0.98
3:A:4002:DT:H2''	3:A:4003:DA:H5''	1.44	0.98
1:N:475:GLN:HE22	1:N:497:LYS:H	1.04	0.98
3:C:4002:DT:H2''	3:C:4003:DA:H5''	1.45	0.97

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	M	284/286 (99%)	216 (76%)	42 (15%)	26 (9%)	0	1
1	N	284/286 (99%)	217 (76%)	41 (14%)	26 (9%)	0	1
2	F	80/82 (98%)	66 (82%)	10 (12%)	4 (5%)	1	5
2	G	80/82 (98%)	69 (86%)	11 (14%)	0	100	100
2	H	80/82 (98%)	66 (82%)	10 (12%)	4 (5%)	1	5
2	I	80/82 (98%)	69 (86%)	11 (14%)	0	100	100
All	All	888/900 (99%)	703 (79%)	125 (14%)	60 (7%)	1	2

5 of 60 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	N	465	GLU
1	N	563	ALA
1	N	583	GLN

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Mol	Chain	Res	Type
1	N	604	THR
1	N	627	THR

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	M	250/250 (100%)	231 (92%)	19 (8%)	12	36
1	N	250/250 (100%)	231 (92%)	19 (8%)	12	36
2	F	74/74 (100%)	71 (96%)	3 (4%)	27	62
2	G	74/74 (100%)	66 (89%)	8 (11%)	6	21
2	H	74/74 (100%)	71 (96%)	3 (4%)	27	62
2	I	74/74 (100%)	66 (89%)	8 (11%)	6	21
All	All	796/796 (100%)	736 (92%)	60 (8%)	12	37

5 of 60 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	G	396	VAL
2	I	380	THR
1	M	476	VAL
2	I	360	LEU
2	I	401	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 42 such sidechains are listed below:

Mol	Chain	Res	Type
1	M	539	ASN
1	M	669	GLN
1	M	583	GLN
1	M	619	GLN
2	H	377	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](#)

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	M	286/286 (100%)	-1.31	0 100 100	36, 78, 110, 117	0
1	N	286/286 (100%)	-1.30	0 100 100	36, 78, 110, 117	0
2	F	82/82 (100%)	-1.41	0 100 100	31, 50, 82, 98	0
2	G	82/82 (100%)	-1.40	0 100 100	28, 52, 95, 100	0
2	H	82/82 (100%)	-1.41	0 100 100	31, 50, 81, 98	0
2	I	82/82 (100%)	-1.37	0 100 100	29, 52, 95, 100	0
3	A	21/21 (100%)	-1.83	0 100 100	32, 42, 65, 71	0
3	C	21/21 (100%)	-1.80	0 100 100	31, 42, 65, 71	0
4	B	21/21 (100%)	-1.85	0 100 100	32, 43, 68, 81	0
4	D	21/21 (100%)	-1.76	0 100 100	32, 43, 68, 81	0
All	All	984/984 (100%)	-1.38	0 100 100	28, 66, 105, 117	0

There are no RSRZ outliers to report.

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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,

median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	MG	G	1	1/1	0.96	0.05	62,62,62,62	0
5	MG	F	1	1/1	0.97	0.03	43,43,43,43	0
5	MG	I	1	1/1	0.97	0.05	61,61,61,61	0
5	MG	H	1	1/1	0.98	0.03	43,43,43,43	0

6.5 Other polymers [i](#)

There are no such residues in this entry.