



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 09:00 am GMT

PDB ID : 2ZPX  
Title : TNF Receptor Subtype One-selective TNF Mutant with Antagonistic Activity;  
R1antTNF-T8  
Authors : Mukai, Y.; Nakamura, T.; Yamagata, Y.; Tsutsumi, Y.  
Deposited on : 2008-07-29  
Resolution : 2.83 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](mailto:validation@mail.wwpdb.org)

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Xtriage (Phenix) : 1.9-1692  
EDS : trunk28620  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : recalc28949

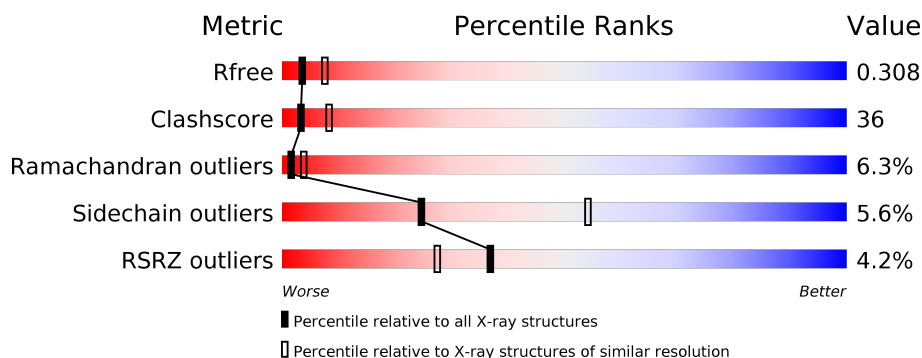
# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.83 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))
$R_{free}$	100719	3466 (2.88-2.80)
Clashscore	112137	3975 (2.88-2.80)
Ramachandran outliers	110173	3902 (2.88-2.80)
Sidechain outliers	110143	3905 (2.88-2.80)
RSRZ outliers	101464	3501 (2.88-2.80)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	157	<div> <div>3%</div> <div> <div></div> <div>39%</div> <div>46%</div> <div>•</div> <div>10%</div> </div> </div>
1	B	157	<div> <div>3%</div> <div> <div></div> <div>40%</div> <div>41%</div> <div>8%</div> <div>•</div> <div>10%</div> </div> </div>
1	C	157	<div> <div>6%</div> <div> <div></div> <div>39%</div> <div>42%</div> <div>7%</div> <div>•</div> <div>10%</div> </div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3319 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 Tumor necrosis factor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	142	Total	C	N	O	S	0	0	0
			1104	704	192	205	3			
1	B	142	Total	C	N	O	S	0	0	0
			1104	704	192	205	3			
1	C	142	Total	C	N	O	S	0	0	0
			1104	704	192	205	3			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	11	MET	LYS	ENGINEERED	UNP P01375
A	65	SER	LYS	ENGINEERED	UNP P01375
A	84	THR	ALA	ENGINEERED	UNP P01375
A	85	PRO	VAL	ENGINEERED	UNP P01375
A	86	ALA	SER	ENGINEERED	UNP P01375
A	87	ILE	TYR	ENGINEERED	UNP P01375
A	88	ASN	GLN	ENGINEERED	UNP P01375
A	89	ARG	THR	ENGINEERED	UNP P01375
A	90	PRO	LYS	ENGINEERED	UNP P01375
A	98	ARG	LYS	ENGINEERED	UNP P01375
A	112	ASN	LYS	ENGINEERED	UNP P01375
A	128	PRO	LYS	ENGINEERED	UNP P01375
B	11	MET	LYS	ENGINEERED	UNP P01375
B	65	SER	LYS	ENGINEERED	UNP P01375
B	84	THR	ALA	ENGINEERED	UNP P01375
B	85	PRO	VAL	ENGINEERED	UNP P01375
B	86	ALA	SER	ENGINEERED	UNP P01375
B	87	ILE	TYR	ENGINEERED	UNP P01375
B	88	ASN	GLN	ENGINEERED	UNP P01375
B	89	ARG	THR	ENGINEERED	UNP P01375
B	90	PRO	LYS	ENGINEERED	UNP P01375
B	98	ARG	LYS	ENGINEERED	UNP P01375
B	112	ASN	LYS	ENGINEERED	UNP P01375

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Chain	Residue	Modelled	Actual	Comment	Reference
B	128	PRO	LYS	ENGINEERED	UNP P01375
C	11	MET	LYS	ENGINEERED	UNP P01375
C	65	SER	LYS	ENGINEERED	UNP P01375
C	84	THR	ALA	ENGINEERED	UNP P01375
C	85	PRO	VAL	ENGINEERED	UNP P01375
C	86	ALA	SER	ENGINEERED	UNP P01375
C	87	ILE	TYR	ENGINEERED	UNP P01375
C	88	ASN	GLN	ENGINEERED	UNP P01375
C	89	ARG	THR	ENGINEERED	UNP P01375
C	90	PRO	LYS	ENGINEERED	UNP P01375
C	98	ARG	LYS	ENGINEERED	UNP P01375
C	112	ASN	LYS	ENGINEERED	UNP P01375
C	128	PRO	LYS	ENGINEERED	UNP P01375

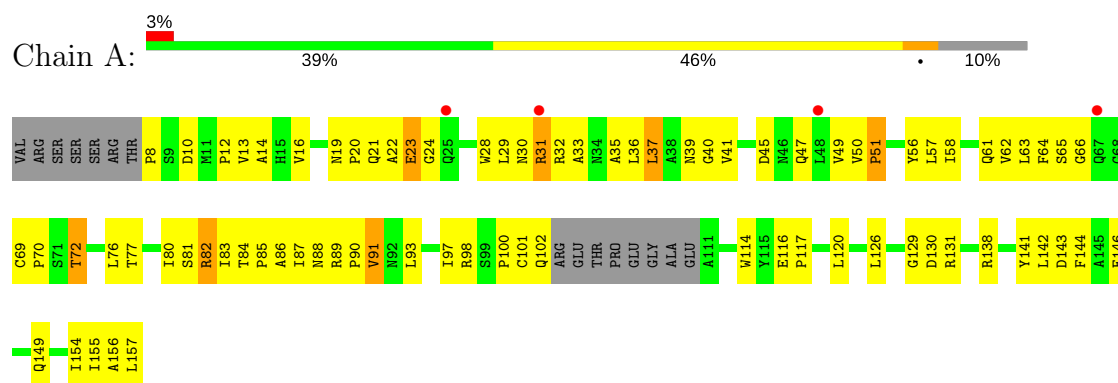
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	5	Total O 5 5	0	0
2	C	2	Total O 2 2	0	0

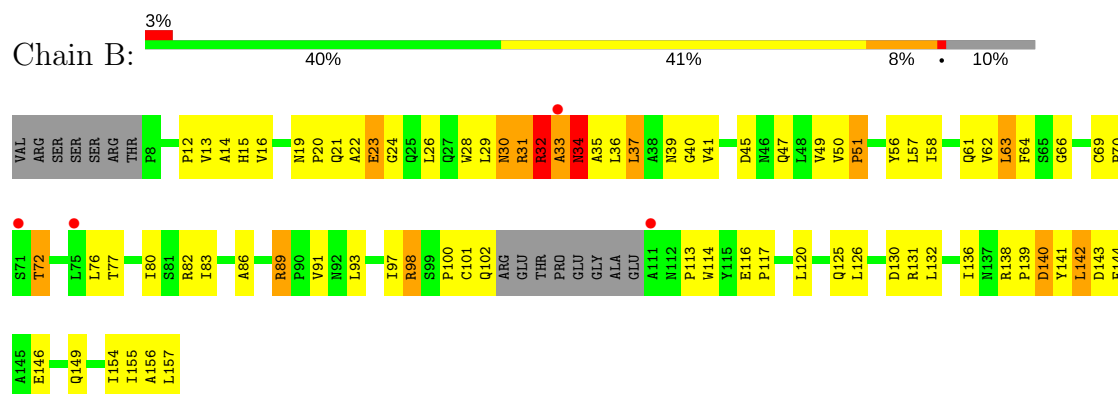
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

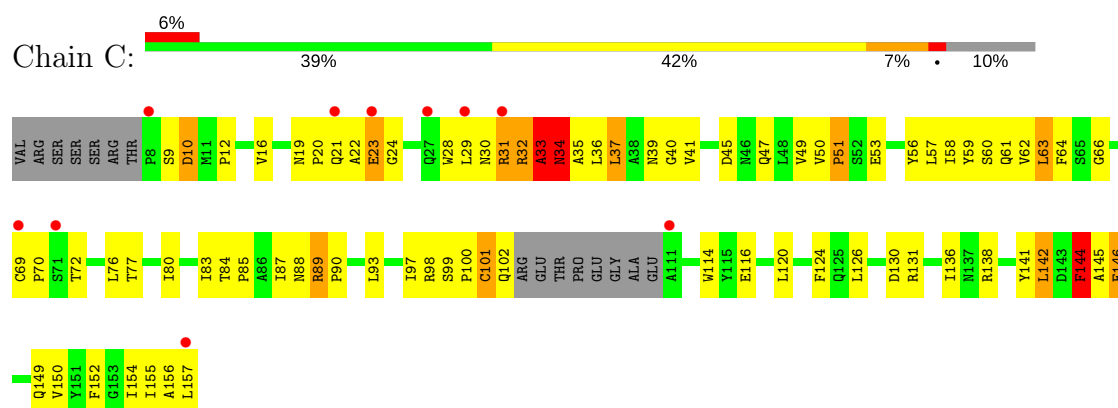
#### • Molecule 1: Tumor necrosis factor



#### • Molecule 1: Tumor necrosis factor



#### • Molecule 1: Tumor necrosis factor



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	49.59Å 111.33Å 75.38Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.83 55.67 – 2.83	Depositor EDS
% Data completeness (in resolution range)	93.2 (50.00-2.83) 93.2 (55.67-2.83)	Depositor EDS
$R_{merge}$	0.24	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.22 (at 2.81Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, $R_{free}$	0.268 , 0.309 0.268 , 0.308	Depositor DCC
$R_{free}$ test set	1009 reflections (10.34%)	DCC
Wilson B-factor (Å <sup>2</sup> )	74.0	Xtriage
Anisotropy	0.250	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 67.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	3319	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	76.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.79% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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	0.43	0/1131	0.62	0/1545
1	B	0.67	2/1131 (0.2%)	1.02	9/1545 (0.6%)
1	C	0.66	3/1131 (0.3%)	1.01	7/1545 (0.5%)
All	All	0.60	5/3393 (0.1%)	0.90	16/4635 (0.3%)

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	2	2
1	C	0	2
All	All	2	4

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	34	ASN	CB-CG	-10.53	1.26	1.51
1	B	32	ARG	N-CA	-9.40	1.27	1.46
1	C	33	ALA	C-O	5.73	1.34	1.23
1	C	33	ALA	N-CA	5.50	1.57	1.46
1	B	31	ARG	CA-CB	-5.23	1.42	1.53

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	34	ASN	N-CA-CB	17.41	141.94	110.60
1	B	33	ALA	N-CA-C	15.40	152.57	111.00
1	C	35	ALA	N-CA-C	-11.84	79.02	111.00
1	B	33	ALA	CA-C-N	-10.62	93.83	117.20
1	B	32	ARG	C-N-CA	-9.44	98.10	121.70
1	B	34	ASN	N-CA-CB	8.67	126.20	110.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	31	ARG	C-N-CA	-8.39	100.73	121.70
1	C	34	ASN	N-CA-C	-8.11	89.09	111.00
1	C	33	ALA	C-N-CA	-8.00	101.69	121.70
1	B	34	ASN	CB-CA-C	7.57	125.53	110.40
1	B	33	ALA	N-CA-CB	-7.18	100.05	110.10
1	C	32	ARG	C-N-CA	-6.75	104.83	121.70
1	B	33	ALA	O-C-N	6.68	133.38	122.70
1	B	31	ARG	CA-C-N	-6.66	102.56	117.20
1	C	32	ARG	N-CA-CB	-6.21	99.43	110.60
1	B	32	ARG	N-CA-C	5.85	126.79	111.00

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	B	33	ALA	CA
1	B	34	ASN	CA

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	32	ARG	Peptide
1	B	33	ALA	Peptide
1	C	33	ALA	Mainchain,Peptide

## 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	1104	0	1087	75	1
1	B	1104	0	1087	93	1
1	C	1104	0	1087	94	2
2	A	5	0	0	2	0
2	C	2	0	0	1	0
All	All	3319	0	3261	236	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 36.



All (236) 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:125:GLN:HB3	1:C:34:ASN:ND2	1.50	1.27
1:C:34:ASN:HB2	1:C:36:LEU:HD21	1.36	1.07
2:A:162:HOH:O	1:B:34:ASN:HB3	1.58	1.02
1:B:89:ARG:HB3	1:B:89:ARG:HH11	1.30	0.95
1:C:34:ASN:HB2	1:C:36:LEU:CD2	1.96	0.94
1:C:62:VAL:HA	1:C:150:VAL:HG22	1.50	0.91
1:B:30:ASN:OD1	1:B:37:LEU:HD23	1.69	0.91
1:C:34:ASN:CB	1:C:36:LEU:HD21	2.03	0.88
1:B:15:HIS:O	1:B:35:ALA:HB1	1.78	0.83
1:B:125:GLN:CB	1:C:34:ASN:ND2	2.38	0.83
1:C:93:LEU:HD22	1:C:124:PHE:CE2	2.20	0.77
1:B:125:GLN:HB3	1:C:34:ASN:HD21	1.47	0.77
1:B:66:GLY:HA3	1:B:114:TRP:CE2	2.21	0.76
2:A:162:HOH:O	1:B:34:ASN:CB	2.26	0.72
1:B:30:ASN:OD1	1:B:37:LEU:CD2	2.37	0.72
1:A:142:LEU:HD23	1:A:143:ASP:N	2.04	0.71
1:C:100:PRO:HG3	1:C:116:GLU:HG3	1.73	0.71
1:B:16:VAL:HG21	1:B:28:TRP:HB3	1.72	0.71
1:C:101:CYS:O	1:C:102:GLN:HG3	1.91	0.70
1:A:16:VAL:HG21	1:A:28:TRP:HB3	1.75	0.68
1:B:26:LEU:HG	1:B:142:LEU:HD21	1.76	0.68
1:A:98:ARG:CZ	1:B:98:ARG:HH22	2.07	0.67
1:C:16:VAL:HG21	1:C:28:TRP:HB3	1.76	0.67
1:B:26:LEU:HG	1:B:142:LEU:CD2	2.24	0.67
1:A:154:ILE:HD12	1:A:154:ILE:C	2.17	0.66
1:A:98:ARG:HH22	1:C:98:ARG:CZ	2.08	0.66
1:C:154:ILE:C	1:C:154:ILE:HD12	2.16	0.66
1:A:142:LEU:HD23	1:A:143:ASP:H	1.60	0.65
1:A:66:GLY:HA3	1:A:114:TRP:CE2	2.32	0.65
1:C:80:ILE:O	1:C:93:LEU:HB2	1.97	0.64
1:A:84:THR:HG23	1:A:85:PRO:HD2	1.81	0.63
1:B:154:ILE:HD12	1:B:154:ILE:C	2.19	0.62
1:A:146:GLU:H	1:A:149:GLN:NE2	1.97	0.62
1:A:98:ARG:NH1	1:B:98:ARG:HH12	1.98	0.62
1:B:125:GLN:H	1:C:34:ASN:CG	2.04	0.61
1:A:35:ALA:O	1:A:36:LEU:HD23	2.02	0.60
1:B:125:GLN:CB	1:C:34:ASN:HD21	2.10	0.60
1:A:87:ILE:HD12	1:A:87:ILE:H	1.67	0.59
1:B:82:ARG:HG2	1:B:83:ILE:N	2.17	0.59
1:C:58:ILE:HA	1:C:154:ILE:HG22	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:125:GLN:HB3	1:C:34:ASN:HD22	1.58	0.59
1:B:157:LEU:HD13	1:C:155:ILE:HG12	1.84	0.59
1:A:12:PRO:HD2	1:A:156:ALA:HB3	1.84	0.59
1:B:58:ILE:HA	1:B:154:ILE:HG22	1.84	0.59
1:B:98:ARG:CZ	1:C:98:ARG:HH22	2.16	0.58
1:B:89:ARG:CB	1:B:89:ARG:HH11	2.12	0.58
1:C:32:ARG:O	1:C:32:ARG:HG3	2.01	0.58
1:A:58:ILE:HA	1:A:154:ILE:HG22	1.85	0.58
1:C:87:ILE:HD13	1:C:87:ILE:N	2.19	0.58
1:B:102:GLN:HG3	1:C:102:GLN:OE1	2.04	0.58
1:B:125:GLN:CA	1:C:34:ASN:HD21	2.17	0.58
1:C:30:ASN:O	1:C:30:ASN:ND2	2.38	0.57
1:B:72:THR:HA	1:B:101:CYS:CB	2.35	0.56
1:A:36:LEU:HD21	2:C:158:HOH:O	2.05	0.56
1:A:64:PHE:CE1	1:A:76:LEU:HD12	2.41	0.56
1:C:30:ASN:OD1	1:C:37:LEU:HB3	2.05	0.56
1:A:16:VAL:CG2	1:A:28:TRP:HB3	2.36	0.56
1:C:34:ASN:CB	1:C:36:LEU:CD2	2.76	0.55
1:C:64:PHE:CE1	1:C:76:LEU:HD12	2.42	0.55
1:A:82:ARG:O	1:A:91:VAL:HG23	2.07	0.55
1:C:12:PRO:HD2	1:C:156:ALA:HB3	1.87	0.55
1:C:32:ARG:O	1:C:33:ALA:C	2.44	0.55
1:C:154:ILE:O	1:C:154:ILE:HD12	2.07	0.55
1:A:30:ASN:O	1:A:32:ARG:N	2.40	0.54
1:B:16:VAL:CG2	1:B:28:TRP:HB3	2.36	0.54
1:B:57:LEU:HB3	1:B:155:ILE:CG2	2.36	0.54
1:B:76:LEU:HD23	1:B:100:PRO:CG	2.37	0.54
1:A:57:LEU:HB3	1:A:155:ILE:CG2	2.37	0.54
1:B:57:LEU:HB3	1:B:155:ILE:HG22	1.88	0.54
1:A:97:ILE:O	1:B:117:PRO:HD3	2.08	0.54
1:B:19:ASN:ND2	1:B:29:LEU:HD11	2.23	0.54
1:B:64:PHE:CE1	1:B:76:LEU:HD12	2.43	0.54
1:C:19:ASN:ND2	1:C:29:LEU:HD11	2.23	0.54
1:A:98:ARG:HH12	1:C:98:ARG:NH1	2.05	0.53
1:B:76:LEU:HD23	1:B:100:PRO:HG2	1.90	0.53
1:B:82:ARG:HB2	1:B:93:LEU:HD11	1.89	0.53
1:C:62:VAL:CA	1:C:150:VAL:HG22	2.33	0.53
1:A:12:PRO:HD2	1:A:156:ALA:CB	2.38	0.53
1:A:88:ASN:O	1:A:89:ARG:HG3	2.09	0.53
1:B:12:PRO:HD2	1:B:156:ALA:HB3	1.89	0.53
1:B:66:GLY:O	1:B:113:PRO:HA	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:8:PRO:HB3	1:C:53:GLU:OE1	2.09	0.53
1:C:76:LEU:HD23	1:C:100:PRO:CD	2.38	0.53
1:A:57:LEU:HB3	1:A:155:ILE:HG22	1.91	0.53
1:C:16:VAL:CG2	1:C:28:TRP:HB3	2.38	0.52
1:C:93:LEU:HD22	1:C:124:PHE:CD2	2.45	0.52
1:C:20:PRO:HB2	1:C:21:GLN:NE2	2.25	0.52
1:B:20:PRO:HB2	1:B:21:GLN:NE2	2.25	0.51
1:A:154:ILE:HD12	1:A:154:ILE:O	2.10	0.51
1:A:49:VAL:HG22	1:A:131:ARG:HG2	1.93	0.51
1:B:45:ASP:O	1:B:47:GLN:HG3	2.11	0.51
1:C:76:LEU:HD23	1:C:100:PRO:HG2	1.93	0.51
1:B:98:ARG:NH1	1:C:98:ARG:HH12	2.08	0.51
1:C:12:PRO:HD2	1:C:156:ALA:CB	2.41	0.51
1:C:57:LEU:HB3	1:C:155:ILE:CG2	2.40	0.51
1:B:66:GLY:HA3	1:B:114:TRP:NE1	2.25	0.51
1:C:76:LEU:HD23	1:C:100:PRO:CG	2.40	0.51
1:B:30:ASN:H	1:B:30:ASN:HD22	1.59	0.51
1:B:76:LEU:HD11	1:B:141:TYR:HD1	1.76	0.51
1:C:66:GLY:HA3	1:C:114:TRP:NE1	2.26	0.50
1:B:76:LEU:HD23	1:B:100:PRO:CD	2.42	0.50
1:A:117:PRO:HD2	1:C:98:ARG:HD3	1.93	0.50
1:C:19:ASN:CB	1:C:29:LEU:HD11	2.42	0.50
1:B:72:THR:HA	1:B:101:CYS:HB2	1.92	0.50
1:C:87:ILE:O	1:C:88:ASN:OD1	2.30	0.50
1:B:154:ILE:HD12	1:B:154:ILE:O	2.12	0.50
1:A:57:LEU:O	1:A:154:ILE:HA	2.12	0.49
1:A:19:ASN:CB	1:A:29:LEU:HD11	2.42	0.49
1:A:65:SER:HB3	1:A:143:ASP:HB2	1.94	0.49
1:B:138:ARG:NH2	1:B:141:TYR:OH	2.45	0.49
1:C:63:LEU:HB2	1:C:149:GLN:O	2.12	0.49
1:A:98:ARG:HD3	1:B:117:PRO:HD2	1.94	0.49
1:B:12:PRO:HD2	1:B:156:ALA:CB	2.41	0.49
1:A:82:ARG:HB2	1:A:93:LEU:HD11	1.95	0.49
1:C:57:LEU:HB3	1:C:155:ILE:HG22	1.93	0.49
1:B:61:GLN:HG2	1:B:62:VAL:N	2.27	0.49
1:A:61:GLN:HG2	1:A:62:VAL:N	2.28	0.49
1:A:64:PHE:CD1	1:A:76:LEU:HD12	2.47	0.49
1:B:76:LEU:CD2	1:B:100:PRO:HG2	2.44	0.48
1:C:45:ASP:O	1:C:47:GLN:HG3	2.13	0.48
1:B:19:ASN:CB	1:B:29:LEU:HD11	2.43	0.48
1:C:144:PHE:CD2	1:C:144:PHE:N	2.79	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:100:PRO:HG3	1:A:116:GLU:HG3	1.94	0.48
1:C:61:GLN:HG2	1:C:62:VAL:N	2.28	0.48
1:C:64:PHE:CD1	1:C:76:LEU:HD12	2.49	0.48
1:B:40:GLY:C	1:B:51:PRO:HG3	2.34	0.48
1:B:64:PHE:CD1	1:B:76:LEU:HD12	2.49	0.48
1:B:140:ASP:OD1	1:B:140:ASP:N	2.47	0.48
1:C:66:GLY:HA3	1:C:114:TRP:CE2	2.49	0.48
1:C:34:ASN:HB2	1:C:36:LEU:HD23	1.93	0.48
1:A:50:VAL:HG11	1:A:126:LEU:HD13	1.95	0.47
1:A:40:GLY:C	1:A:51:PRO:HG3	2.35	0.47
1:A:12:PRO:HA	1:A:39:ASN:HB2	1.96	0.47
1:A:80:ILE:O	1:A:93:LEU:HB2	2.15	0.47
1:B:72:THR:HA	1:B:101:CYS:HB3	1.95	0.47
1:B:82:ARG:O	1:B:83:ILE:HG12	2.14	0.47
1:C:49:VAL:HG22	1:C:131:ARG:HG2	1.96	0.47
1:B:63:LEU:HD12	1:B:149:GLN:HG2	1.96	0.47
1:C:12:PRO:HA	1:C:39:ASN:HB2	1.97	0.47
1:B:100:PRO:HG3	1:B:116:GLU:HG3	1.96	0.46
1:B:50:VAL:HG11	1:B:126:LEU:HD13	1.96	0.46
1:A:19:ASN:ND2	1:A:29:LEU:HD11	2.30	0.46
1:C:50:VAL:HG11	1:C:126:LEU:HD13	1.97	0.46
1:B:138:ARG:C	1:B:140:ASP:H	2.19	0.46
1:B:26:LEU:HG	1:B:142:LEU:HD22	1.97	0.46
1:B:57:LEU:O	1:B:154:ILE:HA	2.15	0.46
1:C:57:LEU:O	1:C:154:ILE:HA	2.16	0.46
1:C:87:ILE:HG22	1:C:88:ASN:H	1.80	0.46
1:A:20:PRO:HB2	1:A:21:GLN:NE2	2.31	0.46
1:B:22:ALA:O	1:B:24:GLY:N	2.48	0.46
1:B:35:ALA:O	1:B:36:LEU:HD23	2.16	0.46
1:B:12:PRO:HA	1:B:39:ASN:HB2	1.97	0.46
1:C:12:PRO:HB3	1:C:41:VAL:HG23	1.98	0.46
1:C:60:SER:HB2	1:C:152:PHE:HD1	1.80	0.46
1:C:22:ALA:O	1:C:24:GLY:N	2.49	0.45
1:A:45:ASP:O	1:A:47:GLN:HG3	2.15	0.45
1:C:89:ARG:HA	1:C:90:PRO:HD3	1.65	0.45
1:A:117:PRO:HD3	1:C:97:ILE:O	2.17	0.45
1:A:72:THR:HA	1:A:101:CYS:HB3	1.99	0.45
1:C:21:GLN:C	1:C:23:GLU:H	2.19	0.45
1:C:83:ILE:O	1:C:83:ILE:HG22	2.16	0.45
1:A:138:ARG:NH2	1:A:141:TYR:HE1	2.15	0.45
1:B:56:TYR:CG	1:B:126:LEU:HD12	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:49:VAL:HG22	1:B:131:ARG:HG2	1.98	0.45
1:B:125:GLN:C	1:C:34:ASN:HD21	2.20	0.45
1:A:30:ASN:ND2	1:A:37:LEU:HD23	2.32	0.45
1:C:19:ASN:CG	1:C:29:LEU:HD11	2.38	0.44
1:C:40:GLY:C	1:C:51:PRO:HG3	2.37	0.44
1:B:125:GLN:H	1:C:34:ASN:ND2	2.15	0.44
1:B:21:GLN:C	1:B:23:GLU:H	2.20	0.44
1:C:144:PHE:C	1:C:146:GLU:H	2.20	0.44
1:A:56:TYR:CG	1:A:126:LEU:HD12	2.53	0.44
1:A:83:ILE:HD12	1:A:83:ILE:N	2.32	0.44
1:A:19:ASN:HB2	1:A:29:LEU:HD11	2.00	0.44
1:C:19:ASN:HD22	1:C:29:LEU:HD21	1.82	0.44
1:C:84:THR:OG1	1:C:85:PRO:HD2	2.18	0.44
1:B:30:ASN:O	1:B:31:ARG:C	2.51	0.44
1:A:72:THR:HB	1:A:102:GLN:C	2.38	0.44
1:A:13:VAL:HG22	1:A:14:ALA:N	2.32	0.43
1:B:13:VAL:HG22	1:B:14:ALA:N	2.33	0.43
1:A:89:ARG:HA	1:A:90:PRO:HD3	1.89	0.43
1:A:72:THR:HA	1:A:101:CYS:CB	2.48	0.43
1:A:50:VAL:HA	1:A:51:PRO:HD3	1.81	0.43
1:B:82:ARG:CG	1:B:83:ILE:N	2.82	0.43
1:C:100:PRO:CG	1:C:116:GLU:HG3	2.44	0.43
1:A:81:SER:HA	1:A:91:VAL:O	2.18	0.43
1:C:56:TYR:CG	1:C:126:LEU:HD12	2.54	0.43
1:A:50:VAL:CG2	1:A:130:ASP:HB2	2.49	0.43
1:B:19:ASN:CG	1:B:29:LEU:HD11	2.39	0.43
1:B:80:ILE:O	1:B:93:LEU:HB2	2.19	0.43
1:B:143:ASP:OD1	1:B:149:GLN:NE2	2.51	0.43
1:A:66:GLY:HA3	1:A:114:TRP:CZ2	2.54	0.42
1:A:69:CYS:HA	1:A:70:PRO:HD3	1.86	0.42
1:B:19:ASN:HB2	1:B:29:LEU:HD11	2.00	0.42
1:B:126:LEU:CD1	1:B:132:LEU:HD11	2.50	0.42
1:B:136:ILE:O	1:B:136:ILE:HG13	2.19	0.42
1:B:77:THR:HG22	1:B:97:ILE:HG23	2.01	0.42
1:A:114:TRP:HA	1:C:99:SER:OG	2.18	0.42
1:A:22:ALA:O	1:A:24:GLY:N	2.52	0.42
1:B:66:GLY:HA3	1:B:114:TRP:CZ2	2.53	0.42
1:C:63:LEU:H	1:C:150:VAL:CG2	2.33	0.42
1:C:77:THR:HG22	1:C:97:ILE:HG23	2.00	0.42
1:A:154:ILE:C	1:A:154:ILE:CD1	2.86	0.42
1:A:85:PRO:HG3	1:A:129:GLY:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:69:CYS:HA	1:B:70:PRO:HD3	1.84	0.42
1:A:58:ILE:HG23	1:A:154:ILE:HG22	2.02	0.42
1:B:58:ILE:HG23	1:B:154:ILE:HG22	2.02	0.42
1:C:154:ILE:C	1:C:154:ILE:CD1	2.86	0.42
1:A:155:ILE:HG12	1:C:157:LEU:HD13	2.01	0.42
1:B:30:ASN:N	1:B:30:ASN:HD22	2.17	0.42
1:B:146:GLU:H	1:B:149:GLN:NE2	2.18	0.41
1:C:142:LEU:HD13	1:C:144:PHE:CD1	2.56	0.41
1:C:58:ILE:HG23	1:C:154:ILE:HG22	2.02	0.41
1:A:88:ASN:C	1:A:89:ARG:HG3	2.41	0.41
1:A:21:GLN:C	1:A:23:GLU:H	2.23	0.41
1:C:141:TYR:O	1:C:142:LEU:C	2.59	0.41
1:A:12:PRO:HB3	1:A:41:VAL:HG23	2.01	0.41
1:A:19:ASN:CG	1:A:29:LEU:HD11	2.41	0.41
1:C:50:VAL:CG2	1:C:130:ASP:HB2	2.51	0.41
1:A:157:LEU:HD13	1:B:155:ILE:HG12	2.02	0.41
1:C:138:ARG:NH2	1:C:141:TYR:HE1	2.19	0.41
1:C:76:LEU:CD2	1:C:100:PRO:HG2	2.51	0.41
1:B:50:VAL:HA	1:B:51:PRO:HD3	1.83	0.40
1:A:77:THR:HG22	1:A:97:ILE:HG23	2.02	0.40
1:B:50:VAL:CG2	1:B:130:ASP:HB2	2.52	0.40
1:B:13:VAL:CG2	1:B:14:ALA:N	2.85	0.40
1:C:136:ILE:O	1:C:136:ILE:HG13	2.21	0.40
1:C:9:SER:HB3	1:C:10:ASP:H	1.67	0.40
1:C:34:ASN:CG	1:C:36:LEU:HD21	2.37	0.40
1:A:142:LEU:CD2	1:A:143:ASP:N	2.80	0.40
1:A:76:LEU:HD23	1:A:100:PRO:CD	2.51	0.40
1:C:19:ASN:HB2	1:C:29:LEU:HD11	2.02	0.40
1:C:59:TYR:O	1:C:60:SER:HB2	2.22	0.40
1:C:69:CYS:HA	1:C:70:PRO:HD3	1.84	0.40
1:A:98:ARG:CZ	1:B:98:ARG:NH2	2.81	0.40
1:B:12:PRO:HB3	1:B:41:VAL:HG23	2.03	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:21:GLN:NE2	1:C:31:ARG:NH2[2_554]	1.83	0.37
1:B:32:ARG:NH2	1:C:10:ASP:OD1[4_445]	2.00	0.20

## 5.3 Torsion angles ⓘ

### 5.3.1 Protein backbone ⓘ

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	138/157 (88%)	111 (80%)	19 (14%)	8 (6%)	2	5
1	B	138/157 (88%)	118 (86%)	10 (7%)	10 (7%)	1	2
1	C	138/157 (88%)	111 (80%)	19 (14%)	8 (6%)	2	5
All	All	414/471 (88%)	340 (82%)	48 (12%)	26 (6%)	1	4

All (26) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	23	GLU
1	A	37	LEU
1	A	51	PRO
1	A	72	THR
1	B	23	GLU
1	B	32	ARG
1	B	37	LEU
1	B	51	PRO
1	B	72	THR
1	B	86	ALA
1	C	10	ASP
1	C	23	GLU
1	C	37	LEU
1	C	51	PRO
1	C	72	THR
1	C	101	CYS
1	C	144	PHE
1	A	10	ASP
1	A	31	ARG
1	A	33	ALA
1	B	34	ASN
1	A	86	ALA
1	B	142	LEU
1	C	145	ALA

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Mol	Chain	Res	Type
1	B	91	VAL
1	B	139	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	120/133 (90%)	114 (95%)	6 (5%)	28	59
1	B	120/133 (90%)	113 (94%)	7 (6%)	23	53
1	C	120/133 (90%)	113 (94%)	7 (6%)	23	53
All	All	360/399 (90%)	340 (94%)	20 (6%)	25	54

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	31	ARG
1	A	63	LEU
1	A	82	ARG
1	A	91	VAL
1	A	120	LEU
1	A	144	PHE
1	B	30	ASN
1	B	63	LEU
1	B	89	ARG
1	B	98	ARG
1	B	120	LEU
1	B	140	ASP
1	B	144	PHE
1	C	34	ASN
1	C	63	LEU
1	C	89	ARG
1	C	120	LEU
1	C	142	LEU
1	C	144	PHE
1	C	146	GLU



Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	21	GLN
1	A	112	ASN
1	A	125	GLN
1	A	149	GLN
1	B	21	GLN
1	B	125	GLN
1	B	149	GLN
1	C	21	GLN
1	C	88	ASN
1	C	125	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 carbohydrates 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 [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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	142/157 (90%)	0.21	4 (2%)	53	44	46, 74, 111, 128	0
1	B	142/157 (90%)	0.28	4 (2%)	53	44	47, 73, 112, 129	0
1	C	142/157 (90%)	0.43	10 (7%)	17	10	50, 75, 110, 129	0
All	All	426/471 (90%)	0.31	18 (4%)	37	27	46, 74, 112, 129	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	33	ALA	4.1
1	C	21	GLN	3.9
1	C	157	LEU	3.6
1	C	23	GLU	3.2
1	B	71	SER	3.2
1	C	69	CYS	3.1
1	C	71	SER	2.8
1	C	27	GLN	2.7
1	B	75	LEU	2.5
1	A	25	GLN	2.3
1	C	8	PRO	2.3
1	C	31	ARG	2.3
1	A	67	GLN	2.3
1	C	111	ALA	2.2
1	A	48	LEU	2.1
1	B	111	ALA	2.0
1	A	31	ARG	2.0
1	C	29	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 carbohydrates 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.