



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 04:26 am GMT

PDB ID : 4LL4  
Title : The structure of the TRX and TXNIP complex  
Authors : Hwang, J.; Kim, M.H.  
Deposited on : 2013-07-09  
Resolution : 2.70 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

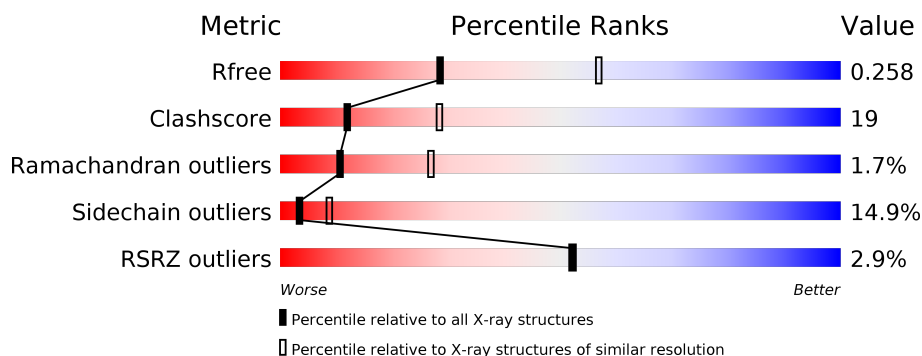
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.70 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	2259 (2.70-2.70)
Clashscore	112137	2590 (2.70-2.70)
Ramachandran outliers	110173	2550 (2.70-2.70)
Sidechain outliers	110143	2550 (2.70-2.70)
RSRZ outliers	101464	2275 (2.70-2.70)

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	315	<div> <div>2%</div> <div> <div>54%</div> <div>27%</div> <div>7%</div> <div>12%</div> </div> </div>
1	C	315	<div> <div>3%</div> <div> <div>54%</div> <div>28%</div> <div>7%</div> <div>11%</div> </div> </div>
2	B	105	<div> <div>5%</div> <div> <div>54%</div> <div>40%</div> <div>6%</div> </div> </div>
2	D	105	<div> <div>67%</div> <div>30%</div> </div>

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 6070 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 Thioredoxin-interacting protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	277	Total	C	N	O	S	0	0	0
			2174	1384	378	401	11			
1	C	281	Total	C	N	O	S	0	0	0
			2197	1397	381	408	11			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	170	SER	CYS	ENGINEERED MUTATION	UNP Q9H3M7
A	205	SER	CYS	ENGINEERED MUTATION	UNP Q9H3M7
A	267	SER	CYS	ENGINEERED MUTATION	UNP Q9H3M7
C	170	SER	CYS	ENGINEERED MUTATION	UNP Q9H3M7
C	205	SER	CYS	ENGINEERED MUTATION	UNP Q9H3M7
C	267	SER	CYS	ENGINEERED MUTATION	UNP Q9H3M7

- Molecule 2 is a protein called Thioredoxin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	105	Total	C	N	O	S	0	0	0
			820	525	128	160	7			
2	D	105	Total	C	N	O	S	0	0	0
			820	525	128	160	7			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	35	ALA	CYS	ENGINEERED MUTATION	UNP P10599
D	35	ALA	CYS	ENGINEERED MUTATION	UNP P10599

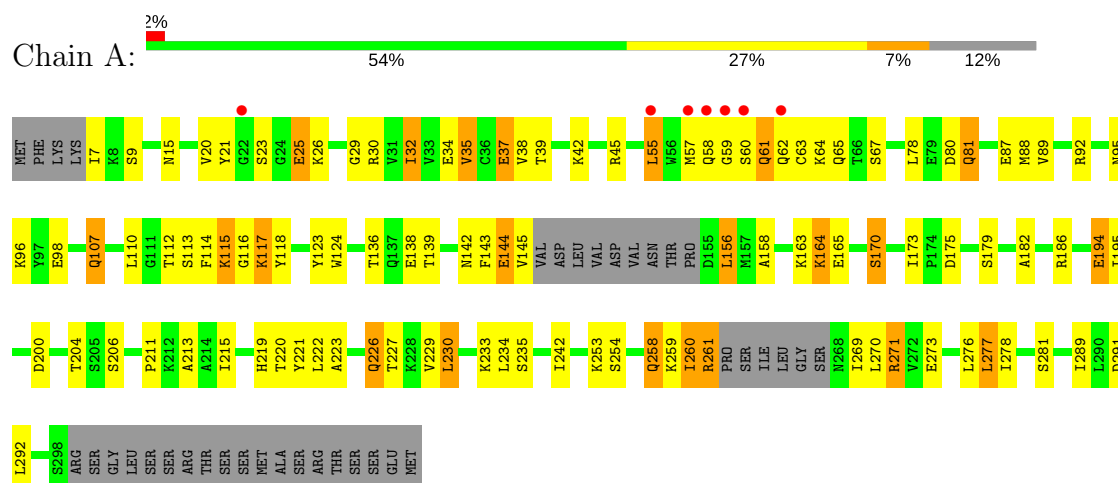
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	22	Total 22	O 22	0	0
3	B	4	Total 4	O 4	0	0
3	C	26	Total 26	O 26	0	0
3	D	7	Total 7	O 7	0	0

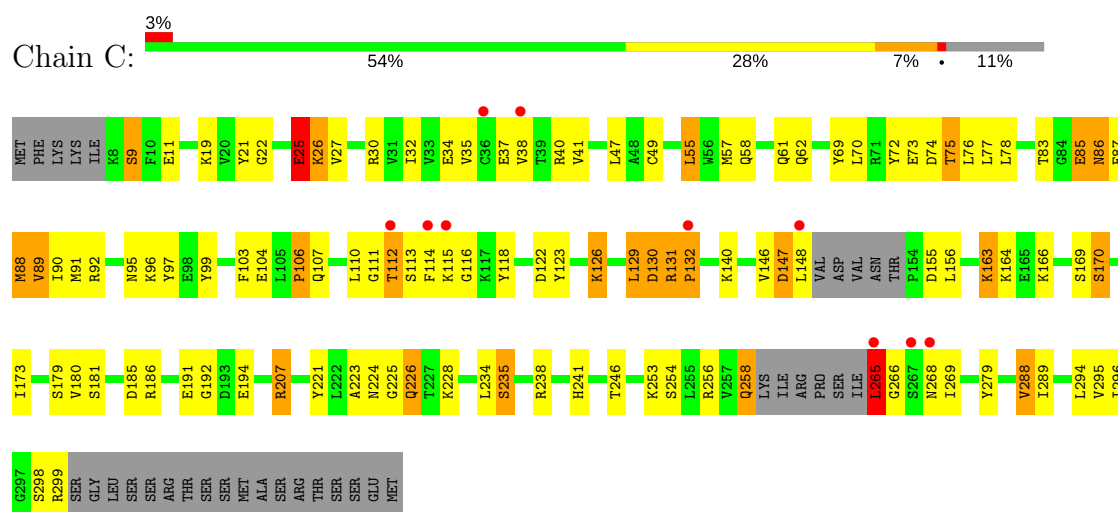
### 3 Residue-property plots [i](#)

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.

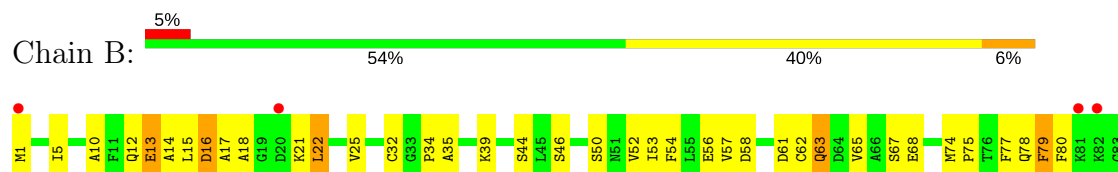
#### • Molecule 1: Thioredoxin-interacting protein

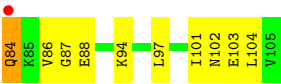


#### • Molecule 1: Thioredoxin-interacting protein



#### • Molecule 2: Thioredoxin





● Molecule 2: Thioredoxin



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	79.83Å 64.99Å 88.42Å 90.00° 90.88° 90.00°	Depositor
Resolution (Å)	40.00 – 2.70 31.88 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.5 (40.00-2.70) 99.5 (31.88-2.70)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.25 (at 2.72Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.196 , 0.264 0.199 , 0.258	Depositor DCC
$R_{free}$ test set	1942 reflections (8.46%)	DCC
Wilson B-factor (Å <sup>2</sup> )	57.4	Xtriage
Anisotropy	0.017	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 51.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.031 for h,-k,-l	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	6070	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.90% 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	1.02	1/2212 (0.0%)	1.05	4/2979 (0.1%)
1	C	1.05	0/2236	1.07	3/3013 (0.1%)
2	B	1.03	2/835 (0.2%)	0.93	1/1121 (0.1%)
2	D	1.04	0/835	0.98	1/1121 (0.1%)
All	All	1.03	3/6118 (0.0%)	1.03	9/8234 (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
1	C	0	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	21	LYS	CD-CE	8.92	1.73	1.51
2	B	21	LYS	CE-NZ	5.53	1.62	1.49
1	A	273	GLU	CD-OE1	5.06	1.31	1.25

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	57	MET	CG-SD-CE	7.38	112.01	100.20
1	C	148	LEU	CA-CB-CG	6.33	129.85	115.30
1	A	276	LEU	CB-CG-CD2	-5.87	101.03	111.00
1	C	265	LEU	CA-CB-CG	5.53	128.02	115.30
1	A	92	ARG	NE-CZ-NH1	-5.31	117.64	120.30
1	A	200	ASP	CB-CG-OD1	5.28	123.05	118.30
1	A	32	ILE	CG1-CB-CG2	-5.20	99.96	111.40
2	B	74	MET	CG-SD-CE	5.16	108.45	100.20
2	D	25	VAL	CB-CA-C	-5.12	101.67	111.40



There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	107	GLN	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	2174	0	2230	83	0
1	C	2197	0	2245	96	0
2	B	820	0	811	34	0
2	D	820	0	810	22	0
3	A	22	0	0	2	0
3	B	4	0	0	1	0
3	C	26	0	0	4	0
3	D	7	0	0	0	0
All	All	6070	0	6096	231	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 19.

All (231) 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:A:226:GLN:CB	1:A:227:THR:HG22	1.60	1.29
1:A:226:GLN:CG	1:A:227:THR:HG22	1.86	1.06
2:B:78:GLN:HG2	2:B:88:GLU:HG3	1.34	1.05
1:A:20:VAL:HG22	1:A:144:GLU:HB2	1.36	1.04
1:C:234:LEU:O	1:C:235:SER:HB2	1.60	1.02
1:C:40:ARG:HG2	1:C:88:MET:CE	1.92	1.00
1:A:226:GLN:HG3	1:A:227:THR:CG2	1.95	0.95
1:A:226:GLN:CA	1:A:227:THR:HG22	1.97	0.94
2:D:12:GLN:NE2	2:D:12:GLN:HA	1.80	0.93
1:C:221:TYR:CD2	1:C:265:LEU:HD11	2.04	0.93
1:A:226:GLN:HB3	1:A:227:THR:HG22	1.50	0.91
1:A:226:GLN:CG	1:A:227:THR:CG2	2.52	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:170:SER:HB3	1:C:173:ILE:H	1.39	0.86
1:A:25:GLU:HG3	1:A:26:LYS:H	1.42	0.85
1:A:226:GLN:CB	1:A:227:THR:CG2	2.51	0.85
1:A:165:GLU:HG3	1:A:179:SER:HB3	1.59	0.84
1:C:163:LYS:HB3	1:C:181:SER:HB3	1.59	0.84
1:C:40:ARG:HG2	1:C:88:MET:HE1	1.59	0.83
1:C:89:VAL:HG23	1:C:90:ILE:H	1.43	0.82
1:A:81:GLN:HG3	1:A:81:GLN:O	1.80	0.82
2:B:12:GLN:HA	2:B:15:LEU:HD12	1.61	0.81
1:C:40:ARG:HG2	1:C:88:MET:HE2	1.61	0.80
1:C:69:TYR:O	1:C:106:PRO:HG2	1.83	0.78
1:A:57:MET:HA	1:A:61:GLN:O	1.83	0.77
2:D:12:GLN:HE21	2:D:12:GLN:HA	1.50	0.76
2:D:78:GLN:HG2	2:D:88:GLU:HG3	1.66	0.76
1:A:226:GLN:HB3	1:A:227:THR:CG2	2.17	0.74
1:C:111:GLY:O	1:C:112:THR:HG23	1.87	0.73
1:C:163:LYS:HB3	1:C:181:SER:CB	2.18	0.73
1:C:192:GLY:O	1:C:256:ARG:NH2	2.23	0.72
1:C:258:GLN:CA	1:C:258:GLN:HE21	2.03	0.72
1:C:89:VAL:CG2	1:C:90:ILE:N	2.54	0.71
2:D:100:THR:HG22	2:D:104:LEU:HD12	1.73	0.71
1:A:260:ILE:HG22	1:A:261:ARG:N	2.04	0.71
1:C:114:PHE:HE1	1:C:116:GLY:HA3	1.56	0.70
1:C:266:GLY:O	1:C:269:ILE:HG22	1.92	0.70
1:C:131:ARG:HB3	1:C:132:PRO:HD2	1.71	0.70
1:C:26:LYS:HB2	1:C:104:GLU:HG2	1.73	0.70
1:A:226:GLN:HA	1:A:227:THR:HG22	1.73	0.69
1:A:42:LYS:HA	1:A:88:MET:HG2	1.72	0.69
1:A:112:THR:O	1:A:115:LYS:N	2.15	0.69
1:A:233:LYS:HD2	1:C:87:GLU:OE1	1.93	0.69
1:C:129:LEU:C	1:C:130:ASP:OD2	2.32	0.68
1:C:146:VAL:O	1:C:146:VAL:HG13	1.94	0.68
1:C:41:VAL:HG21	1:C:91:MET:CE	2.24	0.68
1:C:89:VAL:HG23	1:C:90:ILE:N	2.09	0.68
1:C:90:ILE:HG22	1:C:90:ILE:O	1.93	0.68
1:C:191:GLU:O	1:C:299:ARG:NH1	2.28	0.67
1:C:221:TYR:CG	1:C:265:LEU:HD11	2.29	0.67
2:B:84:GLN:HE21	2:B:84:GLN:HA	1.58	0.67
1:C:223:ALA:HB3	1:C:228:LYS:HD2	1.76	0.67
1:C:114:PHE:CE1	1:C:116:GLY:HA3	2.30	0.66
2:D:48:LYS:HD3	2:D:49:TYR:CE2	2.29	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:25:GLU:HG3	1:A:26:LYS:N	2.08	0.66
1:A:21:TYR:O	1:A:145:VAL:HA	1.96	0.66
1:A:221:TYR:HB3	1:A:269:ILE:HG22	1.78	0.66
1:C:73:GLU:O	1:C:74:ASP:HB2	1.95	0.66
1:C:265:LEU:HB2	3:C:415:HOH:O	1.97	0.65
1:C:111:GLY:O	1:C:112:THR:CG2	2.44	0.65
1:C:74:ASP:OD1	1:C:75:THR:N	2.30	0.64
1:A:34:GLU:OE2	1:A:96:LYS:HG3	1.98	0.64
2:D:63:GLN:H	2:D:63:GLN:HE21	1.45	0.63
1:C:131:ARG:CB	1:C:132:PRO:HD2	2.28	0.63
1:C:234:LEU:O	1:C:235:SER:CB	2.36	0.63
1:A:115:LYS:O	1:A:115:LYS:HG3	1.98	0.63
1:C:112:THR:HB	1:C:147:ASP:H	1.64	0.63
1:A:81:GLN:OE1	1:A:89:VAL:HG22	1.99	0.62
1:A:222:LEU:HD22	1:C:225:GLY:HA3	1.81	0.62
2:D:12:GLN:NE2	2:D:12:GLN:CA	2.61	0.62
1:C:258:GLN:HE21	1:C:258:GLN:C	2.02	0.62
2:B:102:ASN:O	2:B:104:LEU:N	2.32	0.62
1:A:23:SER:OG	1:A:107:GLN:HA	2.00	0.62
1:A:260:ILE:HG22	1:A:261:ARG:H	1.64	0.62
1:A:219:HIS:O	1:A:229:VAL:HA	2.01	0.61
1:C:129:LEU:O	1:C:130:ASP:OD2	2.19	0.60
1:A:117:LYS:HD3	1:A:118:TYR:CE2	2.36	0.60
1:A:223:ALA:O	1:A:226:GLN:HB2	2.01	0.60
1:C:110:LEU:HB3	1:C:123:TYR:OH	2.01	0.60
1:A:32:ILE:HG12	1:A:98:GLU:HG2	1.83	0.60
1:C:258:GLN:NE2	1:C:258:GLN:H	2.00	0.59
1:A:124:TRP:HB3	1:A:142:ASN:HA	1.84	0.59
1:C:258:GLN:CA	1:C:258:GLN:NE2	2.65	0.59
2:D:100:THR:CG2	2:D:104:LEU:HD12	2.33	0.59
1:C:72:TYR:CE1	1:C:103:PHE:HB3	2.38	0.59
2:B:46:SER:HA	2:B:54:PHE:CE1	2.38	0.58
1:A:23:SER:HA	1:A:110:LEU:HD11	1.86	0.57
1:C:170:SER:HB3	1:C:173:ILE:N	2.17	0.57
1:A:110:LEU:HD13	1:A:145:VAL:HG21	1.88	0.56
1:A:123:TYR:HD1	1:A:143:PHE:CZ	2.23	0.56
1:C:169:SER:O	1:C:207:ARG:NH2	2.37	0.56
2:B:35:ALA:HA	2:B:75:PRO:HG3	1.88	0.55
1:C:288:VAL:HA	3:C:414:HOH:O	2.06	0.55
2:B:84:GLN:NE2	2:B:84:GLN:HA	2.21	0.55
1:A:234:LEU:HB3	1:A:258:GLN:HE21	1.72	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:26:LYS:HA	1:C:104:GLU:HA	1.89	0.55
1:A:260:ILE:CG2	1:A:261:ARG:N	2.70	0.54
2:B:16:ASP:OD1	2:B:16:ASP:N	2.41	0.54
1:A:58:GLN:O	1:A:60:SER:N	2.41	0.54
2:B:87:GLY:O	2:B:88:GLU:HB2	2.07	0.54
1:A:118:TYR:CD2	1:A:156:LEU:HD23	2.42	0.54
2:B:22:LEU:HD23	2:B:52:VAL:HG13	1.90	0.54
1:A:170:SER:HB3	1:A:173:ILE:H	1.73	0.53
2:D:12:GLN:HE21	2:D:12:GLN:CA	2.18	0.53
2:B:46:SER:HA	2:B:54:PHE:CZ	2.44	0.53
1:A:114:PHE:CE1	1:A:156:LEU:HB2	2.44	0.53
1:A:195:ILE:HD12	1:A:215:ILE:HD11	1.91	0.53
2:D:26:ASP:HB2	2:D:77:PHE:CE1	2.44	0.53
1:A:158:ALA:O	1:A:186:ARG:HA	2.09	0.52
1:C:126:LYS:HD2	1:C:140:LYS:HB2	1.91	0.52
1:C:114:PHE:HE1	1:C:116:GLY:CA	2.22	0.52
1:C:41:VAL:HG21	1:C:91:MET:HE2	1.92	0.52
1:C:22:GLY:O	1:C:25:GLU:HG3	2.10	0.52
2:D:59:VAL:O	2:D:63:GLN:HG3	2.09	0.52
1:C:126:LYS:CD	1:C:140:LYS:HB2	2.40	0.52
1:A:226:GLN:HG2	3:A:417:HOH:O	2.10	0.51
1:A:7:ILE:HD13	1:A:35:VAL:HG13	1.92	0.51
1:A:219:HIS:CD2	1:A:230:LEU:HD23	2.45	0.51
1:A:226:GLN:HE21	1:A:227:THR:HG21	1.75	0.51
1:A:114:PHE:HE1	1:A:156:LEU:HB2	1.73	0.51
1:C:258:GLN:NE2	1:C:258:GLN:N	2.58	0.51
1:C:163:LYS:CB	1:C:181:SER:HB3	2.36	0.51
1:C:91:MET:SD	1:C:97:TYR:CD2	3.04	0.51
2:B:84:GLN:CA	2:B:84:GLN:HE21	2.24	0.51
1:A:260:ILE:CG2	1:A:261:ARG:H	2.24	0.51
1:C:111:GLY:C	1:C:112:THR:CG2	2.79	0.50
2:B:101:ILE:HG22	2:B:102:ASN:N	2.26	0.50
2:B:94:LYS:HG3	3:B:202:HOH:O	2.10	0.50
1:A:182:ALA:HB1	1:A:278:ILE:HD13	1.93	0.50
1:A:55:LEU:HB3	1:A:64:LYS:HG3	1.93	0.50
1:C:164:LYS:HE2	1:C:289:ILE:O	2.10	0.50
2:B:18:ALA:HB2	2:B:53:ILE:HG13	1.92	0.50
1:C:27:VAL:HB	1:C:103:PHE:CE1	2.47	0.49
1:C:223:ALA:O	1:C:226:GLN:O	2.30	0.49
1:A:80:ASP:O	1:A:80:ASP:CG	2.50	0.49
1:C:72:TYR:CZ	1:C:103:PHE:HB3	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:57:VAL:HG11	2:B:65:VAL:CG1	2.43	0.49
1:C:269:ILE:O	1:C:269:ILE:HG23	2.12	0.49
1:A:277:LEU:HD12	1:A:291:ASP:HA	1.95	0.48
1:C:112:THR:O	1:C:112:THR:OG1	2.29	0.48
1:C:266:GLY:O	1:C:269:ILE:CG2	2.60	0.48
1:C:131:ARG:HB3	1:C:132:PRO:CD	2.41	0.48
1:A:226:GLN:HA	1:A:227:THR:CG2	2.43	0.48
1:A:78:LEU:HB2	1:A:81:GLN:HB3	1.95	0.48
1:C:90:ILE:CG2	1:C:90:ILE:O	2.61	0.48
2:B:22:LEU:HA	2:B:80:PHE:O	2.13	0.48
1:C:19:LYS:HD3	1:C:21:TYR:CE1	2.49	0.48
1:A:227:THR:O	1:A:227:THR:HG23	2.13	0.48
2:B:62:CYS:O	2:B:63:GLN:C	2.51	0.48
1:C:221:TYR:CD2	1:C:265:LEU:CD1	2.87	0.48
1:A:81:GLN:O	1:A:81:GLN:CG	2.58	0.48
2:B:32:CYS:SG	2:B:34:PRO:HD2	2.55	0.47
2:D:27:PHE:CG	2:D:71:VAL:HG21	2.49	0.47
2:D:39:LYS:HE3	2:D:39:LYS:HB3	1.47	0.47
1:C:112:THR:HG23	3:C:422:HOH:O	2.14	0.47
1:C:246:THR:HG22	2:D:91:GLY:HA2	1.95	0.47
1:C:295:VAL:O	1:C:296:ILE:HG13	2.14	0.47
1:A:30:ARG:HH11	1:A:30:ARG:HG2	1.80	0.47
2:B:102:ASN:C	2:B:104:LEU:H	2.17	0.47
1:A:35:VAL:HG11	1:A:39:THR:OG1	2.14	0.47
1:C:238:ARG:HH22	1:C:241:HIS:CD2	2.33	0.46
1:A:211:PRO:HG3	1:A:242:ILE:HD11	1.96	0.46
1:A:30:ARG:NH1	1:A:30:ARG:HG2	2.30	0.46
1:A:114:PHE:CE1	1:A:156:LEU:HD13	2.50	0.46
2:B:14:ALA:HA	2:B:17:ALA:HB3	1.97	0.46
2:D:32:CYS:HB3	2:D:35:ALA:HB3	1.98	0.46
1:C:191:GLU:O	1:C:191:GLU:HG2	2.16	0.46
1:A:110:LEU:HD13	1:A:145:VAL:CG2	2.45	0.46
1:C:146:VAL:O	1:C:146:VAL:CG1	2.63	0.46
1:C:166:LYS:NZ	3:C:405:HOH:O	2.49	0.46
1:C:76:LEU:HD23	1:C:99:TYR:HB3	1.96	0.46
2:B:54:PHE:CD1	2:B:54:PHE:N	2.84	0.46
1:C:40:ARG:O	1:C:132:PRO:HD3	2.16	0.46
1:C:55:LEU:HD11	1:C:62:GLN:NE2	2.31	0.46
1:C:70:LEU:O	1:C:70:LEU:HG	2.16	0.46
1:C:279:TYR:CD1	1:C:279:TYR:N	2.84	0.45
1:C:30:ARG:HE	1:C:32:ILE:HD11	1.80	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:79:PHE:N	2:B:79:PHE:CD1	2.84	0.45
1:A:123:TYR:HB3	1:A:143:PHE:CE2	2.52	0.45
2:D:32:CYS:SG	2:D:34:PRO:HD2	2.56	0.45
1:A:226:GLN:HG3	1:A:227:THR:HG21	1.92	0.45
1:A:139:THR:O	1:A:139:THR:HG23	2.17	0.45
1:C:49:CYS:O	1:C:123:TYR:HA	2.15	0.44
1:C:130:ASP:N	1:C:130:ASP:OD2	2.50	0.44
1:A:67:SER:HB3	1:A:271:ARG:HB3	1.99	0.44
2:D:37:MET:O	2:D:40:PRO:HD2	2.17	0.44
1:C:111:GLY:C	1:C:112:THR:HG22	2.36	0.44
2:D:46:SER:HA	2:D:54:PHE:CE1	2.53	0.44
1:A:173:ILE:HD12	1:A:206:SER:HB3	1.98	0.44
2:B:58:ASP:HB3	2:B:61:ASP:HB2	1.99	0.44
1:C:49:CYS:HA	1:C:70:LEU:O	2.18	0.44
1:C:9:SER:OG	1:C:34:GLU:HB2	2.16	0.44
2:D:95:GLU:OE1	2:D:95:GLU:N	2.47	0.44
1:A:81:GLN:HE21	1:A:81:GLN:HA	1.82	0.44
1:A:114:PHE:O	1:A:116:GLY:N	2.51	0.44
1:C:224:ASN:HA	1:C:225:GLY:HA2	1.57	0.44
1:C:86:ASN:HA	1:C:86:ASN:HD22	1.57	0.44
1:A:118:TYR:CG	1:A:156:LEU:HD23	2.53	0.44
1:A:15:ASN:OD1	1:A:29:GLY:HA2	2.17	0.44
2:B:5:ILE:HB	2:B:56:GLU:O	2.17	0.44
2:D:41:PHE:CE2	2:D:94:LYS:HD3	2.53	0.43
1:C:83:THR:O	1:C:87:GLU:HG3	2.18	0.43
1:C:89:VAL:CG2	1:C:90:ILE:H	2.13	0.43
1:A:204:THR:HG22	2:B:34:PRO:HG3	2.01	0.43
1:A:213:ALA:HA	1:A:277:LEU:O	2.19	0.43
2:B:22:LEU:HD23	2:B:52:VAL:CG1	2.48	0.43
2:B:18:ALA:HB2	2:B:53:ILE:CG1	2.49	0.43
1:A:226:GLN:CA	1:A:227:THR:CG2	2.85	0.42
2:D:89:PHE:CD1	2:D:89:PHE:C	2.92	0.42
1:C:55:LEU:HD21	1:C:62:GLN:NE2	2.35	0.42
1:A:23:SER:CA	1:A:110:LEU:HD11	2.49	0.42
2:B:25:VAL:O	2:B:77:PHE:HA	2.19	0.42
1:A:37:GLU:HG3	1:A:37:GLU:H	1.48	0.42
1:C:194:GLU:OE1	1:C:256:ARG:NH1	2.52	0.42
1:A:25:GLU:CG	1:A:26:LYS:N	2.81	0.42
1:A:112:THR:HB	1:A:115:LYS:HG2	2.01	0.42
1:A:164:LYS:NZ	1:A:289:ILE:O	2.51	0.41
1:C:118:TYR:HE2	1:C:155:ASP:OD1	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:10:ALA:HA	2:B:13:GLU:OE1	2.20	0.41
1:A:34:GLU:HA	3:A:411:HOH:O	2.21	0.41
1:C:58:GLN:OE1	1:C:186:ARG:NH1	2.53	0.41
1:C:85:GLU:C	1:C:87:GLU:H	2.23	0.41
2:B:12:GLN:CA	2:B:15:LEU:HD12	2.42	0.41
2:B:77:PHE:O	2:B:88:GLU:HA	2.21	0.41
2:B:94:LYS:O	2:B:97:LEU:HB3	2.20	0.41
1:A:156:LEU:O	1:A:156:LEU:HG	2.10	0.40
1:C:47:LEU:HD11	1:C:49:CYS:SG	2.61	0.40
2:B:80:PHE:N	2:B:80:PHE:CD1	2.90	0.40
2:D:19:GLY:H	2:D:21:LYS:HG3	1.86	0.40
1:A:194:GLU:HG2	1:A:254:SER:OG	2.20	0.40
1:C:114:PHE:CE1	1:C:116:GLY:CA	3.01	0.40
1:C:22:GLY:H	1:C:25:GLU:HG3	1.87	0.40

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	A	271/315 (86%)	249 (92%)	18 (7%)	4 (2%)	12	30
1	C	275/315 (87%)	249 (90%)	18 (6%)	8 (3%)	5	13
2	B	103/105 (98%)	86 (84%)	16 (16%)	1 (1%)	18	43
2	D	103/105 (98%)	98 (95%)	5 (5%)	0	100	100
All	All	752/840 (90%)	682 (91%)	57 (8%)	13 (2%)	11	27

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	156	LEU

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Mol	Chain	Res	Type
2	B	103	GLU
1	C	38	VAL
1	C	113	SER
1	C	132	PRO
1	C	235	SER
1	A	59	GLY
1	C	25	GLU
1	C	156	LEU
1	A	113	SER
1	A	115	LYS
1	C	131	ARG
1	C	106	PRO

### 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	242/277 (87%)	204 (84%)	38 (16%)	3	8
1	C	245/277 (88%)	205 (84%)	40 (16%)	3	7
2	B	91/91 (100%)	78 (86%)	13 (14%)	4	9
2	D	91/91 (100%)	82 (90%)	9 (10%)	9	21
All	All	669/736 (91%)	569 (85%)	100 (15%)	3	9

All (100) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	SER
1	A	25	GLU
1	A	35	VAL
1	A	37	GLU
1	A	38	VAL
1	A	45	ARG
1	A	55	LEU
1	A	61	GLN
1	A	62	GLN

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Mol	Chain	Res	Type
1	A	63	CYS
1	A	65	GLN
1	A	81	GLN
1	A	87	GLU
1	A	95	ASN
1	A	107	GLN
1	A	117	LYS
1	A	136	THR
1	A	138	GLU
1	A	144	GLU
1	A	163	LYS
1	A	164	LYS
1	A	170	SER
1	A	175	ASP
1	A	194	GLU
1	A	220	THR
1	A	226	GLN
1	A	230	LEU
1	A	235	SER
1	A	253	LYS
1	A	258	GLN
1	A	259	LYS
1	A	260	ILE
1	A	261	ARG
1	A	270	LEU
1	A	271	ARG
1	A	277	LEU
1	A	281	SER
1	A	292	LEU
2	B	1	MET
2	B	13	GLU
2	B	16	ASP
2	B	22	LEU
2	B	39	LYS
2	B	44	SER
2	B	50	SER
2	B	63	GLN
2	B	67	SER
2	B	68	GLU
2	B	79	PHE
2	B	84	GLN
2	B	86	VAL

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Mol	Chain	Res	Type
1	C	9	SER
1	C	11	GLU
1	C	25	GLU
1	C	26	LYS
1	C	35	VAL
1	C	37	GLU
1	C	55	LEU
1	C	61	GLN
1	C	75	THR
1	C	77	LEU
1	C	78	LEU
1	C	85	GLU
1	C	86	ASN
1	C	88	MET
1	C	89	VAL
1	C	92	ARG
1	C	95	ASN
1	C	96	LYS
1	C	112	THR
1	C	115	LYS
1	C	122	ASP
1	C	126	LYS
1	C	129	LEU
1	C	130	ASP
1	C	147	ASP
1	C	163	LYS
1	C	170	SER
1	C	179	SER
1	C	180	VAL
1	C	185	ASP
1	C	207	ARG
1	C	226	GLN
1	C	253	LYS
1	C	254	SER
1	C	258	GLN
1	C	265	LEU
1	C	268	ASN
1	C	288	VAL
1	C	294	LEU
1	C	298	SER
2	D	8	LYS
2	D	12	GLN

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Mol	Chain	Res	Type
2	D	15	LEU
2	D	21	LYS
2	D	45	LEU
2	D	51	ASN
2	D	63	GLN
2	D	82	LYS
2	D	105	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	81	GLN
1	A	219	HIS
1	A	226	GLN
1	A	258	GLN
2	B	78	GLN
2	B	84	GLN
2	B	102	ASN
1	C	15	ASN
1	C	62	GLN
1	C	81	GLN
1	C	86	ASN
1	C	95	ASN
1	C	134	GLN
1	C	219	HIS
1	C	224	ASN
1	C	232	GLN
1	C	258	GLN
2	D	12	GLN
2	D	43	HIS
2	D	63	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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 ⓘ

### 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, 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	277/315 (87%)	0.05	7 (2%) 58 58	33, 59, 89, 110	0
1	C	281/315 (89%)	-0.08	10 (3%) 43 42	27, 51, 92, 104	0
2	B	105/105 (100%)	0.13	5 (4%) 31 29	28, 70, 125, 143	0
2	D	105/105 (100%)	-0.49	0 100 100	23, 45, 78, 91	0
All	All	768/840 (91%)	-0.06	22 (2%) 52 52	23, 55, 97, 143	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	57	MET	4.6
1	A	60	SER	4.5
1	A	58	GLN	4.5
2	B	20	ASP	4.4
1	C	267	SER	4.4
1	A	59	GLY	3.7
1	A	55	LEU	3.4
1	C	148	LEU	3.2
1	C	265	LEU	3.1
2	B	81	LYS	3.0
1	C	36	CYS	2.9
1	A	22	GLY	2.9
1	C	114	PHE	2.7
1	C	38	VAL	2.6
1	C	112	THR	2.5
2	B	82	LYS	2.4
2	B	1	MET	2.4
1	C	132	PRO	2.4
1	C	268	ASN	2.4
1	A	62	GLN	2.4
1	C	115	LYS	2.2
2	B	84	GLN	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.