



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 03:33 pm GMT

PDB ID : 1JZD  
Title : DsbC-DsbDalphi complex  
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Deposited on : 2001-09-15  
Resolution : 2.30 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

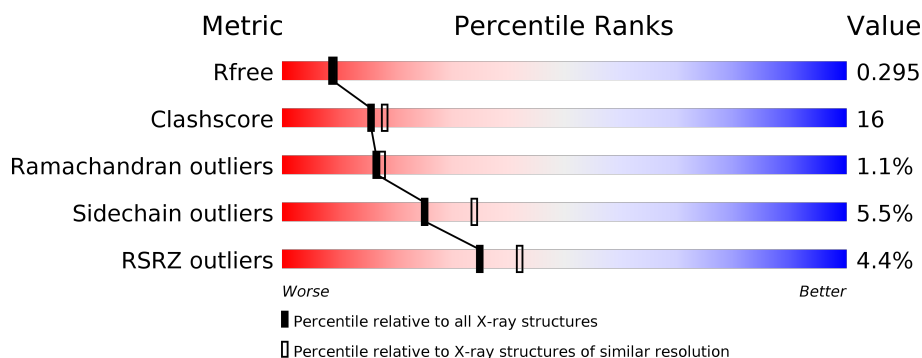
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.30 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	4130 (2.30-2.30)
Clashscore	112137	4751 (2.30-2.30)
Ramachandran outliers	110173	4705 (2.30-2.30)
Sidechain outliers	110143	4704 (2.30-2.30)
RSRZ outliers	101464	4156 (2.30-2.30)

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	220	<div> <div>5%</div> <div> <div></div> <div>67%</div> <div>27%</div> <div>5%</div> </div> </div>
1	B	220	<div> <div>3%</div> <div> <div></div> <div>68%</div> <div>29%</div> <div>•</div> </div> </div>
2	C	132	<div> <div>5%</div> <div> <div></div> <div>61%</div> <div>23%</div> <div>5%</div> <div>•</div> <div>11%</div> </div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4347 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 thiol:disulfide interchange protein dsbc.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	219	Total	C	N	O	S	0	0	0
			1652	1042	275	321	14			
1	B	214	Total	C	N	O	S	0	0	0
			1626	1027	270	316	13			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	GLY	-	EXPRESSION TAG	UNP P21892
A	-2	ALA	-	EXPRESSION TAG	UNP P21892
A	-1	MET	-	EXPRESSION TAG	UNP P21892
A	0	ALA	-	EXPRESSION TAG	UNP P21892
A	101	SER	CYS	ENGINEERED	UNP P21892
B	-3	GLY	-	EXPRESSION TAG	UNP P21892
B	-2	ALA	-	EXPRESSION TAG	UNP P21892
B	-1	MET	-	EXPRESSION TAG	UNP P21892
B	0	ALA	-	EXPRESSION TAG	UNP P21892
B	101	SER	CYS	ENGINEERED	UNP P21892

- Molecule 2 is a protein called thiol:disulfide interchange protein dsbd.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	118	Total	C	N	O	S	4	0	0
			945	605	158	181	1			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	103	ALA	CYS	ENGINEERED	UNP P36655

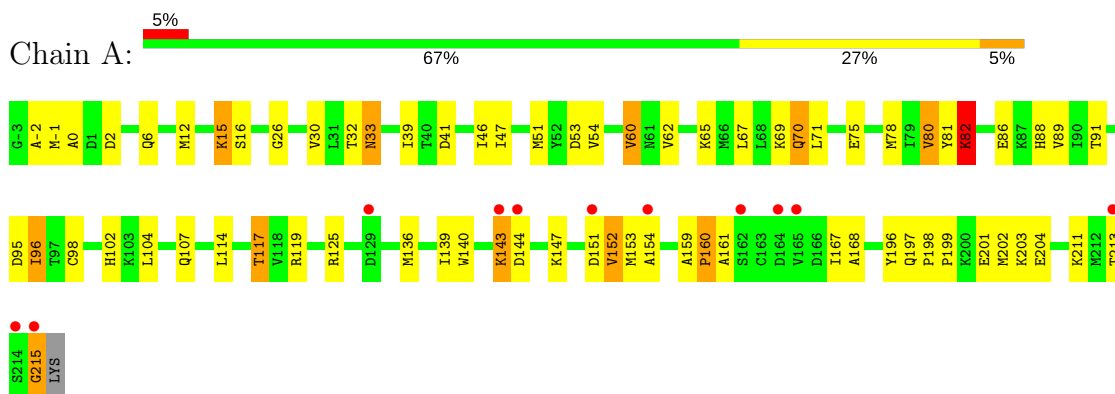
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	37	Total 37	O 37	0	0
3	B	46	Total 46	O 46	0	0
3	C	41	Total 41	O 41	0	0

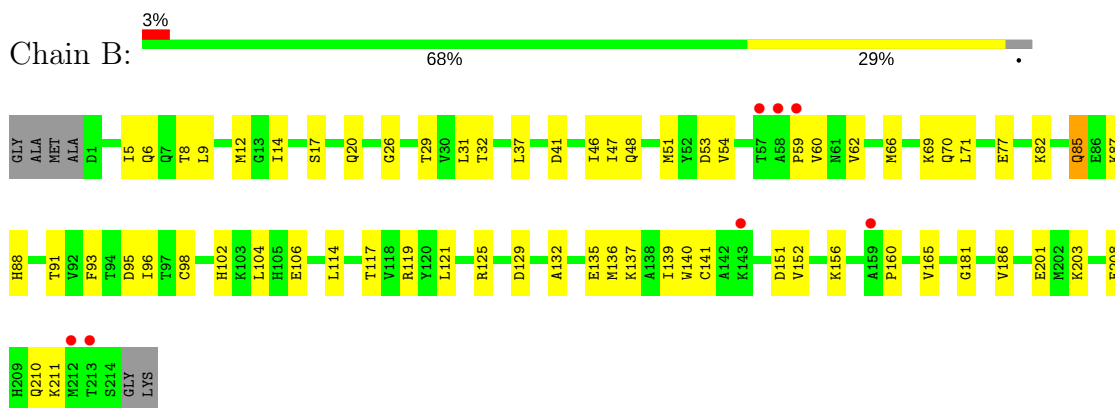
### 3 Residue-property plots [i](#)

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.

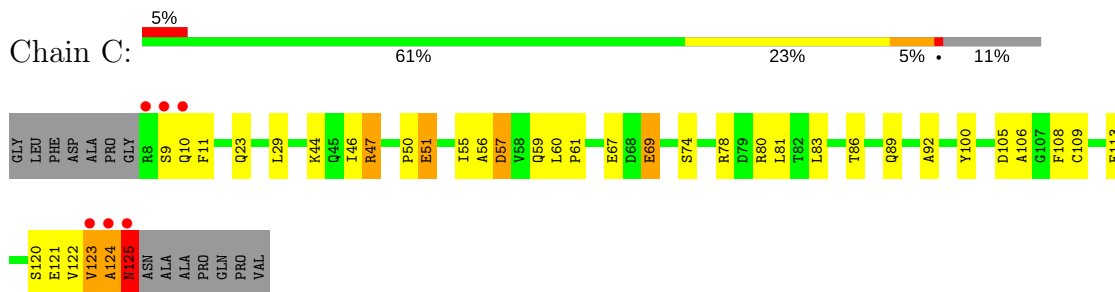
- Molecule 1: thiol:disulfide interchange protein dsbc



- Molecule 1: thiol:disulfide interchange protein dsbc



- Molecule 2: thiol:disulfide interchange protein dsbd



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	68.91Å 68.91Å 230.34Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.77 – 2.30 33.01 – 1.90	Depositor EDS
% Data completeness (in resolution range)	98.3 (29.77-2.30) 88.8 (33.01-1.90)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	0.07	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.13 (at 1.89Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, $R_{free}$	0.234 , 0.293 0.235 , 0.295	Depositor DCC
$R_{free}$ test set	1268 reflections (5.03%)	DCC
Wilson B-factor (Å <sup>2</sup> )	39.0	Xtriage
Anisotropy	0.136	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 47.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	4347	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.95% 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 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	0.48	3/1682 (0.2%)	0.68	3/2277 (0.1%)
1	B	0.36	0/1656	0.60	0/2243
2	C	0.59	2/971 (0.2%)	0.91	7/1326 (0.5%)
All	All	0.47	5/4309 (0.1%)	0.71	10/5846 (0.2%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	82	LYS	CB-CG	6.61	1.70	1.52
2	C	125	ASN	N-CA	5.86	1.58	1.46
1	A	80	VAL	CB-CG1	-5.55	1.41	1.52
2	C	123	VAL	CA-C	5.52	1.67	1.52
1	A	82	LYS	CD-CE	5.22	1.64	1.51

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	124	ALA	N-CA-C	12.36	144.38	111.00
2	C	123	VAL	N-CA-C	9.79	137.44	111.00
2	C	125	ASN	N-CA-C	9.05	135.44	111.00
2	C	124	ALA	CB-CA-C	-7.21	99.29	110.10
2	C	125	ASN	CB-CA-C	-6.68	97.05	110.40
1	A	215	GLY	N-CA-C	6.36	128.99	113.10
1	A	80	VAL	CG1-CB-CG2	-5.96	101.36	110.90
2	C	124	ALA	CA-C-O	-5.75	108.03	120.10
1	A	82	LYS	CB-CG-CD	5.30	125.37	111.60
2	C	124	ALA	CA-C-N	5.04	128.28	117.20

There are no chirality outliers.

There are no planarity outliers.

## 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	1652	0	1658	55	6
1	B	1626	0	1633	52	0
2	C	945	0	896	39	6
3	A	37	0	0	0	0
3	B	46	0	0	0	0
3	C	41	0	0	2	0
All	All	4347	0	4187	137	6

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 16.

All (137) 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)
2:C:124:ALA:C	2:C:125:ASN:ND2	1.72	1.42
2:C:124:ALA:O	2:C:125:ASN:ND2	1.85	1.10
1:A:53:ASP:HB3	1:A:60:VAL:HG13	1.50	0.92
2:C:124:ALA:CA	2:C:125:ASN:ND2	2.36	0.88
1:A:15:LYS:HD2	1:A:16:SER:H	1.40	0.83
2:C:124:ALA:HA	2:C:125:ASN:ND2	1.92	0.82
1:A:15:LYS:HD2	1:A:16:SER:N	1.95	0.81
1:B:14:ILE:HD11	1:B:32:THR:HG21	1.64	0.80
2:C:92:ALA:HB2	2:C:121:GLU:OE2	1.82	0.79
1:A:198:PRO:HB2	1:A:201:GLU:OE1	1.83	0.78
2:C:50:PRO:HG3	2:C:55:ILE:HD11	1.64	0.77
1:A:102:HIS:HA	1:A:153:MET:HE3	1.67	0.77
1:A:211:LYS:O	1:A:215:GLY:HA2	1.86	0.76
1:A:71:LEU:HD11	1:A:91:THR:HG21	1.69	0.75
1:A:102:HIS:HA	1:A:153:MET:CE	2.19	0.72
1:B:87:LYS:HE2	1:B:87:LYS:HA	1.71	0.71
2:C:123:VAL:HG12	2:C:124:ALA:N	2.07	0.70
2:C:125:ASN:N	2:C:125:ASN:ND2	2.40	0.69
1:B:14:ILE:HG12	1:B:17:SER:OG	1.93	0.68
1:B:14:ILE:HG12	1:B:14:ILE:O	1.94	0.67
2:C:78:ARG:HD3	3:C:158:HOH:O	1.95	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:51:MET:HE2	1:B:47:ILE:HD11	1.78	0.66
1:B:125:ARG:HH11	1:B:125:ARG:HG2	1.61	0.65
2:C:50:PRO:HG3	2:C:55:ILE:CD1	2.26	0.64
1:A:95:ASP:OD2	1:A:125:ARG:HG2	1.99	0.63
1:A:69:LYS:HB3	1:A:70:GLN:NE2	2.14	0.62
1:B:14:ILE:HD11	1:B:32:THR:CG2	2.29	0.62
1:B:151:ASP:HA	1:B:156:LYS:HE2	1.82	0.61
2:C:69:GLU:CD	2:C:69:GLU:H	2.05	0.60
1:A:2:ASP:O	1:A:6:GLN:HG3	2.00	0.60
2:C:46:ILE:O	2:C:47:ARG:HD2	2.02	0.60
2:C:78:ARG:HG3	2:C:78:ARG:HH11	1.66	0.60
1:A:96:ILE:HD12	1:A:136:MET:CG	2.31	0.59
2:C:120:SER:O	2:C:122:VAL:HG23	2.02	0.59
1:B:54:VAL:HG12	1:B:59:PRO:HB3	1.82	0.59
1:A:96:ILE:HD12	1:A:136:MET:HG3	1.83	0.58
1:A:211:LYS:O	1:A:215:GLY:CA	2.51	0.58
1:B:211:LYS:NZ	1:B:211:LYS:HB3	2.18	0.58
1:B:53:ASP:HB3	1:B:60:VAL:HG12	1.84	0.58
1:A:82:LYS:NZ	1:A:86:GLU:OE1	2.35	0.58
2:C:56:ALA:HB3	2:C:86:THR:HB	1.85	0.57
2:C:69:GLU:CD	2:C:69:GLU:N	2.58	0.57
2:C:124:ALA:O	2:C:125:ASN:CG	2.42	0.57
1:A:211:LYS:O	1:A:215:GLY:N	2.38	0.56
2:C:123:VAL:CG1	2:C:124:ALA:N	2.69	0.55
1:B:71:LEU:HD11	1:B:91:THR:HG21	1.89	0.55
1:A:54:VAL:HG11	1:B:12:MET:SD	2.47	0.54
1:A:39:ILE:HG12	1:A:46:ILE:HG12	1.89	0.53
1:A:67:LEU:O	1:A:70:GLN:HG2	2.08	0.52
1:B:77:GLU:HG2	1:B:165:VAL:HG22	1.92	0.52
1:B:82:LYS:HA	1:B:117:THR:HG23	1.92	0.51
1:B:37:LEU:CD2	1:B:48:GLN:HG2	2.40	0.51
1:A:213:THR:HG22	1:A:213:THR:O	2.10	0.50
1:B:12:MET:HE1	1:B:46:ILE:HD11	1.92	0.50
1:B:53:ASP:HB3	1:B:60:VAL:CG1	2.40	0.50
1:B:137:LYS:HE2	1:B:141:CYS:SG	2.51	0.50
1:A:51:MET:O	1:A:62:VAL:HG22	2.11	0.49
1:B:201:GLU:OE1	1:B:201:GLU:N	2.45	0.49
1:A:104:LEU:C	1:A:104:LEU:HD23	2.32	0.49
2:C:44:LYS:NZ	2:C:74:SER:HB3	2.28	0.49
1:B:125:ARG:NH1	1:B:125:ARG:HG2	2.27	0.49
2:C:89:GLN:HA	2:C:122:VAL:O	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:44:LYS:HG3	3:C:142:HOH:O	2.11	0.49
1:A:143:LYS:N	1:A:143:LYS:HD2	2.28	0.49
1:B:88:HIS:HE1	1:B:210:GLN:CG	2.26	0.48
1:A:89:VAL:HG13	1:A:117:THR:HG22	1.95	0.48
1:B:98:CYS:HG	2:C:109:CYS:HG	0.60	0.48
1:A:197:GLN:HB3	1:A:202:MET:HE2	1.96	0.48
1:A:-2:ALA:HB1	1:A:2:ASP:OD2	2.13	0.48
1:B:14:ILE:CD1	1:B:32:THR:HG21	2.37	0.48
2:C:23:GLN:O	2:C:29:LEU:HD12	2.14	0.48
1:A:30:VAL:HG21	1:A:39:ILE:HD12	1.96	0.48
1:A:15:LYS:CD	1:A:16:SER:N	2.74	0.47
1:B:14:ILE:HG12	1:B:17:SER:HG	1.79	0.47
1:B:93:PHE:HE1	1:B:186:VAL:HG21	1.78	0.47
1:B:20:GLN:HE22	1:B:69:LYS:HD3	1.79	0.47
1:A:47:ILE:HD11	1:B:51:MET:HE2	1.96	0.47
1:A:51:MET:CE	1:B:47:ILE:HD11	2.44	0.47
2:C:69:GLU:OE1	2:C:69:GLU:N	2.49	0.46
1:A:139:ILE:HD11	1:A:152:VAL:HG21	1.98	0.46
2:C:78:ARG:NH2	2:C:105:ASP:OD2	2.45	0.46
1:A:167:ILE:N	1:A:167:ILE:HD12	2.31	0.46
2:C:56:ALA:O	2:C:57:ASP:HB3	2.16	0.46
1:A:98:CYS:O	1:A:102:HIS:HD2	1.99	0.45
1:A:53:ASP:HB3	1:A:60:VAL:CG1	2.34	0.45
1:A:144:ASP:HB3	1:A:147:LYS:HB3	1.98	0.45
2:C:11:PHE:CE2	2:C:106:ALA:HB3	2.51	0.45
1:A:26:GLY:HA2	1:A:41:ASP:OD2	2.16	0.45
1:A:96:ILE:HG23	1:A:136:MET:HG2	1.98	0.45
2:C:60:LEU:HD23	2:C:83:LEU:HD11	1.99	0.45
2:C:92:ALA:N	2:C:121:GLU:HG2	2.31	0.45
1:A:12:MET:HE1	1:B:59:PRO:HB3	1.98	0.45
2:C:51:GLU:HA	2:C:51:GLU:OE1	2.17	0.45
1:B:132:ALA:O	1:B:136:MET:HG3	2.17	0.44
1:B:211:LYS:HZ2	1:B:211:LYS:HB3	1.81	0.44
1:B:51:MET:O	1:B:62:VAL:HG22	2.17	0.44
1:B:121:LEU:HD23	1:B:121:LEU:HA	1.80	0.44
1:B:114:LEU:HB3	1:B:203:LYS:HD3	2.00	0.43
1:A:151:ASP:O	1:A:154:ALA:HB3	2.18	0.43
1:B:104:LEU:HD23	1:B:104:LEU:C	2.38	0.43
1:A:144:ASP:O	1:A:147:LYS:HB3	2.17	0.43
1:A:114:LEU:O	1:A:203:LYS:HE3	2.18	0.43
1:A:88:HIS:O	1:A:117:THR:HB	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:5:ILE:O	1:B:9:LEU:HG	2.19	0.43
1:B:181:GLY:HA3	2:C:108:PHE:CZ	2.54	0.43
1:A:75:GLU:HA	1:A:78:MET:HG3	2.01	0.43
1:B:8:THR:HG22	1:B:12:MET:HE1	1.99	0.43
1:B:95:ASP:OD2	1:B:125:ARG:HG2	2.19	0.43
1:B:91:THR:OG1	1:B:119:ARG:NH2	2.49	0.42
1:A:104:LEU:O	1:A:104:LEU:HD23	2.19	0.42
1:A:96:ILE:O	1:A:102:HIS:NE2	2.48	0.42
1:A:107:GLN:NE2	1:A:196:TYR:OH	2.48	0.42
2:C:124:ALA:O	2:C:125:ASN:CB	2.68	0.42
2:C:78:ARG:HG3	2:C:78:ARG:NH1	2.32	0.42
1:B:102:HIS:HB2	2:C:69:GLU:HB2	2.02	0.42
1:B:96:ILE:HD11	1:B:135:GLU:HB2	2.01	0.41
1:B:69:LYS:HG3	1:B:70:GLN:N	2.34	0.41
2:C:57:ASP:OD2	2:C:59:GLN:NE2	2.52	0.41
2:C:80:ARG:HG2	2:C:81:LEU:N	2.35	0.41
1:B:152:VAL:HA	1:B:156:LYS:O	2.20	0.41
1:A:-2:ALA:O	1:A:0:ALA:N	2.54	0.41
1:A:32:THR:HG22	1:A:33:ASN:N	2.35	0.41
2:C:124:ALA:HA	2:C:125:ASN:HD21	1.76	0.41
1:B:8:THR:HG22	1:B:12:MET:CE	2.51	0.41
1:A:159:ALA:O	1:A:160:PRO:O	2.39	0.41
1:B:31:LEU:HD12	1:B:31:LEU:N	2.36	0.41
1:A:96:ILE:HD12	1:A:136:MET:HG2	2.03	0.41
2:C:60:LEU:HA	2:C:61:PRO:HD3	1.96	0.41
1:A:198:PRO:HA	1:A:199:PRO:HD3	1.95	0.40
1:B:29:THR:HG21	1:B:66:MET:CE	2.51	0.40
1:A:197:GLN:HA	1:A:198:PRO:HD3	1.95	0.40
1:A:80:VAL:HG13	1:A:119:ARG:HG2	2.04	0.40
1:A:54:VAL:HG21	1:B:46:ILE:HD11	2.03	0.40
1:B:26:GLY:HA2	1:B:41:ASP:OD2	2.22	0.40
1:B:85:GLN:HE21	1:B:85:GLN:HB3	1.58	0.40
1:B:139:ILE:HD11	1:B:152:VAL:HG21	2.03	0.40
2:C:100:TYR:CE1	2:C:113:GLU:HB2	2.56	0.40

All (6) 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:81:TYR:O	2:C:125:ASN:O[5_655]	1.65	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:81:TYR:C	2:C:125:ASN:O[5_655]	1.65	0.55
1:A:82:LYS:CB	2:C:124:ALA:O[5_655]	1.82	0.38
1:A:82:LYS:NZ	2:C:125:ASN:OD1[5_655]	2.00	0.20
1:A:82:LYS:N	2:C:125:ASN:O[5_655]	2.04	0.16
1:A:80:VAL:CG1	2:C:125:ASN:N[5_655]	2.05	0.15

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	217/220 (99%)	193 (89%)	19 (9%)	5 (2%)	7	5
1	B	212/220 (96%)	202 (95%)	9 (4%)	1 (0%)	32	39
2	C	116/132 (88%)	105 (90%)	11 (10%)	0	100	100
All	All	545/572 (95%)	500 (92%)	39 (7%)	6 (1%)	17	18

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	-1	MET
1	A	160	PRO
1	A	161	ALA
1	A	168	ALA
1	A	152	VAL
1	B	160	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	179/180 (99%)	168 (94%)	11 (6%)	22	29
1	B	178/180 (99%)	172 (97%)	6 (3%)	42	57
2	C	99/109 (91%)	91 (92%)	8 (8%)	14	17
All	All	456/469 (97%)	431 (94%)	25 (6%)	25	34

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	LYS
1	A	33	ASN
1	A	60	VAL
1	A	65	LYS
1	A	70	GLN
1	A	82	LYS
1	A	96	ILE
1	A	117	THR
1	A	140	TRP
1	A	143	LYS
1	A	204	GLU
1	B	6	GLN
1	B	85	GLN
1	B	106	GLU
1	B	129	ASP
1	B	140	TRP
1	B	208	GLU
2	C	9	SER
2	C	10	GLN
2	C	47	ARG
2	C	51	GLU
2	C	57	ASP
2	C	67	GLU
2	C	69	GLU
2	C	125	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	70	GLN
1	A	107	GLN

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Mol	Chain	Res	Type
1	A	112	ASN
1	A	197	GLN
1	B	20	GLN
1	B	85	GLN
1	B	88	HIS
1	B	112	ASN
1	B	126	GLN
1	B	176	GLN
1	B	209	HIS
2	C	23	GLN
2	C	25	ASN

### 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

### 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	219/220 (99%)	0.35	11 (5%) 30 36	26, 44, 79, 86	0
1	B	214/220 (97%)	0.09	7 (3%) 47 54	24, 41, 64, 90	0
2	C	118/132 (89%)	0.44	6 (5%) 29 36	19, 34, 63, 101	1 (0%)
All	All	551/572 (96%)	0.27	24 (4%) 35 42	19, 41, 71, 101	1 (0%)

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	C	124	ALA	23.6
2	C	125	ASN	20.2
1	A	215	GLY	7.5
1	A	129	ASP	6.3
1	B	213	THR	4.7
1	B	57	THR	4.2
1	A	162	SER	4.2
1	A	164	ASP	4.2
2	C	10	GLN	4.1
2	C	8	ARG	3.8
1	B	212	MET	3.7
2	C	123	VAL	3.3
1	B	59	PRO	3.2
1	B	143	LYS	3.2
1	B	58	ALA	2.7
1	B	159	ALA	2.6
1	A	213	THR	2.6
1	A	154	ALA	2.6
1	A	143	LYS	2.6
1	A	165	VAL	2.4
1	A	151	ASP	2.3
2	C	9	SER	2.3
1	A	214	SER	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	144	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 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.