Full wwPDB X-ray Structure Validation Report (i)

Jun 13, 2024 – 08:24 AM EDT

PDB ID	:	4DSR
Title	:	Crystal structure of peroxiredoxin Ahp1 from Saccharomyces cerevisiae in reduced form
Authors	:	Lian, F.M.; Yu, J.; Ma, X.X.; Yu, X.J.; Chen, Y.; Zhou, C.Z.
Deposited on	:	2012-02-19
Resolution	:	2.91 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) 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 (1)) were used in the production of this report:

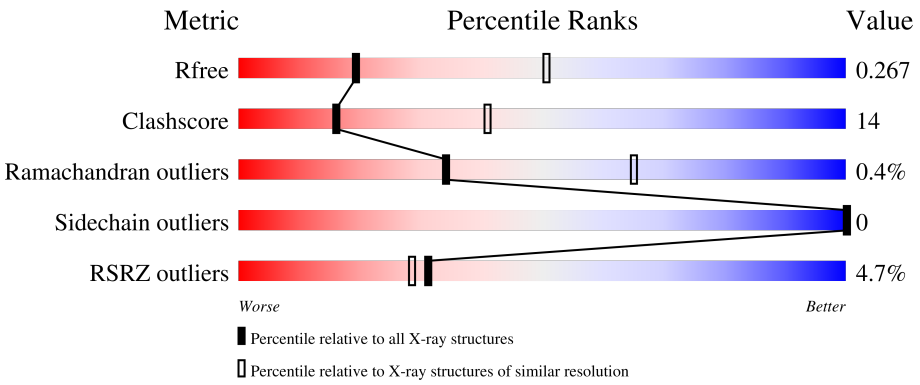
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.20.1
EDS	:	2.36.2
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

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 2.91 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}	130704	2307 (2.94-2.90)
Clashscore	141614	2531 (2.94-2.90)
Ramachandran outliers	138981	2462 (2.94-2.90)
Sidechain outliers	138945	2464 (2.94-2.90)
RSRZ outliers	127900	2248 (2.94-2.90)

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 <=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	184	4% 74% 20% 5%
1	B	184	2% 79% 15% 5%
1	C	184	7% 70% 25% 5%
1	D	184	5% 58% 37% 5%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 5339 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 Peroxiredoxin type-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	174	1332	852	213	262	5	0	0	0
1	B	174	1332	852	213	262	5	0	0	0
1	C	174	1332	852	213	262	5	0	0	0
1	D	174	1332	852	213	262	5	0	0	0

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	MET	-	expression tag	UNP P38013
A	-6	GLY	-	expression tag	UNP P38013
A	-5	HIS	-	expression tag	UNP P38013
A	-4	HIS	-	expression tag	UNP P38013
A	-3	HIS	-	expression tag	UNP P38013
A	-2	HIS	-	expression tag	UNP P38013
A	-1	HIS	-	expression tag	UNP P38013
A	0	HIS	-	expression tag	UNP P38013
B	-7	MET	-	expression tag	UNP P38013
B	-6	GLY	-	expression tag	UNP P38013
B	-5	HIS	-	expression tag	UNP P38013
B	-4	HIS	-	expression tag	UNP P38013
B	-3	HIS	-	expression tag	UNP P38013
B	-2	HIS	-	expression tag	UNP P38013
B	-1	HIS	-	expression tag	UNP P38013
B	0	HIS	-	expression tag	UNP P38013
C	-7	MET	-	expression tag	UNP P38013
C	-6	GLY	-	expression tag	UNP P38013
C	-5	HIS	-	expression tag	UNP P38013
C	-4	HIS	-	expression tag	UNP P38013
C	-3	HIS	-	expression tag	UNP P38013

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-2	HIS	-	expression tag	UNP P38013
C	-1	HIS	-	expression tag	UNP P38013
C	0	HIS	-	expression tag	UNP P38013
D	-7	MET	-	expression tag	UNP P38013
D	-6	GLY	-	expression tag	UNP P38013
D	-5	HIS	-	expression tag	UNP P38013
D	-4	HIS	-	expression tag	UNP P38013
D	-3	HIS	-	expression tag	UNP P38013
D	-2	HIS	-	expression tag	UNP P38013
D	-1	HIS	-	expression tag	UNP P38013
D	0	HIS	-	expression tag	UNP P38013

- Molecule 2 is water.

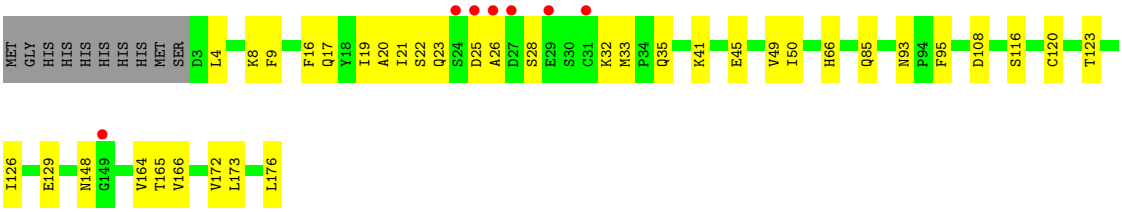
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	5	5	5	0	0
			Total	O		
2	B	1	1	1	0	0
			Total	O		
2	C	3	3	3	0	0
			Total	O		
2	D	2	2	2	0	0

3 Residue-property plots (i)

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.

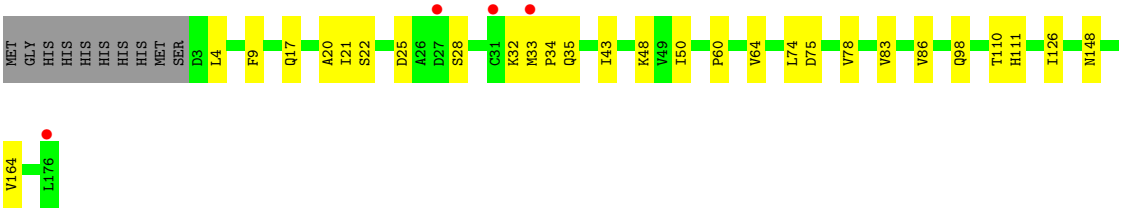
• Molecule 1: Peroxiredoxin type-2

Chain A:



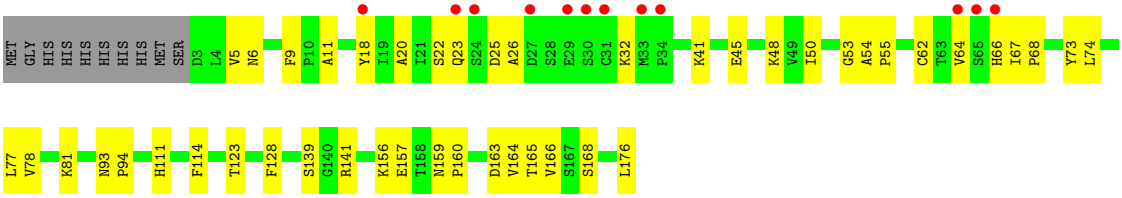
• Molecule 1: Peroxiredoxin type-2

Chain B:



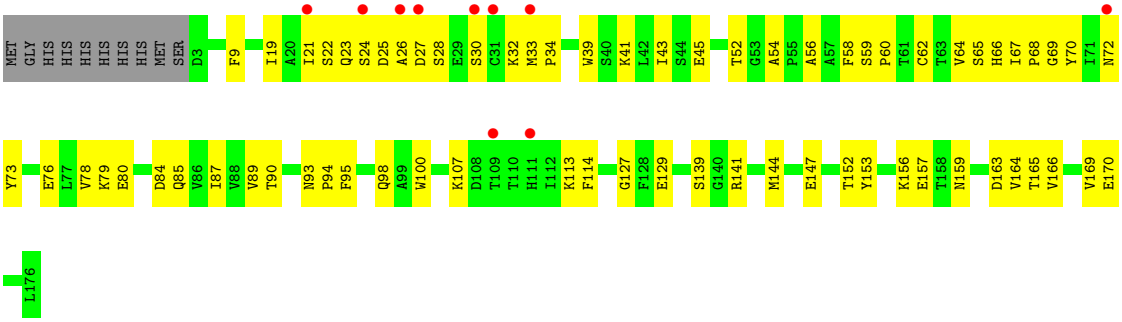
• Molecule 1: Peroxiredoxin type-2

Chain C:



• Molecule 1: Peroxiredoxin type-2

Chain D:


MET
GLY
HIS
HIS
HIS
HIS
HIS
MET
SER
D3
F9
I19
A20
I21
S22
Q23
S24
D25
A26
D27
S28
E29
S30
C31
K32
M33
P34
W39
S40
K41
L42
I43
S44
E45
T52
G53
A54
P55
A56
A57
F58
S59
P60
T61
C62
T63
V64
S65
H66
I67
P68
G69
Y70
I71
N72
Y73
E76
L77
V78
K79
E80
D84
Q85
V86
I87
V88
V89
T90
N93
P94
F95
Q98
A99
W100
K107
D108
T109
T110
H111
I112
K113
F114
G127
F128
E129
S139
G140
R141
M144
E147
T152
Y153
K156
E157
T158
N159
D163
V164
T165
V166
V169
E170
L176

4 Data and refinement statistics (i)

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α, β, γ	154.46Å 154.46Å 90.97Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	44.20 – 2.91 44.19 – 2.91	Depositor EDS
% Data completeness (in resolution range)	91.3 (44.20-2.91) 91.3 (44.19-2.91)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	0.13	Depositor
$< I/\sigma(I) >$ [1]	2.85 (at 2.90Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.239 , 0.274 0.234 , 0.267	Depositor DCC
R_{free} test set	835 reflections (5.16%)	wwPDB-VP
Wilson B-factor (Å^2)	62.3	Xtriage
Anisotropy	0.082	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.29 , 27.5	EDS
L-test for twinning[2]	$< \|L\| > = 0.46$, $< L^2 > = 0.29$	Xtriage
Estimated twinning fraction	0.028 for -2/3*h-1/3*k-4/3*l,-1/3*h-2/3*k+4/3*l,-1/3*h+1/3*k+1/3*l 0.026 for -h,1/3*h-1/3*k-4/3*l,-1/3*h-2/3*k+1/3*l 0.028 for -1/3*h+1/3*k+4/3*l,-k,2/3*h+1/3*k+1/3*l 0.047 for -h,2/3*h+1/3*k+4/3*l,1/3*h+2/3*k-1/3*l 0.078 for -1/3*h-2/3*k+4/3*l,-2/3*h-1/3*k-4/3*l,1/3*h-1/3*k-1/3*l 0.047 for 1/3*h+2/3*k-4/3*l,-k,-2/3*h-1/3*k-1/3*l 0.043 for h,-h-k,-l	Xtriage
F_o,F_c correlation	0.90	EDS
Total number of atoms	5339	wwPDB-VP
Average B, all atoms (Å^2)	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.23% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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	0.57	0/1363	0.66	0/1856
1	B	0.60	0/1363	0.70	0/1856
1	C	0.65	0/1363	0.67	0/1856
1	D	0.70	0/1363	0.68	0/1856
All	All	0.63	0/5452	0.68	0/7424

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1332	0	1308	34	0
1	B	1332	0	1308	27	0
1	C	1332	0	1308	37	0
1	D	1332	0	1308	56	0
2	A	5	0	0	0	0
2	B	1	0	0	0	0
2	C	3	0	0	0	0
2	D	2	0	0	0	0
All	All	5339	0	5232	153	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 14.

All (153) 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:20:ALA:HB3	1:B:32:LYS:CE	1.72	1.18
1:D:22:SER:HB2	1:D:25:ASP:HB3	1.25	1.15
1:B:20:ALA:CB	1:B:32:LYS:HE3	1.79	1.12
1:D:85:GLN:HE22	1:D:113:LYS:NZ	1.48	1.10
1:D:22:SER:CB	1:D:25:ASP:HB3	1.87	1.04
1:A:22:SER:OG	1:A:25:ASP:HB2	1.57	1.03
1:A:20:ALA:CB	1:A:32:LYS:HE2	2.01	0.90
1:A:20:ALA:HB3	1:A:32:LYS:HD3	1.53	0.90
1:A:20:ALA:HB3	1:A:32:LYS:CD	2.06	0.86
1:D:85:GLN:NE2	1:D:113:LYS:NZ	2.23	0.85
1:A:20:ALA:HB3	1:A:32:LYS:CE	2.05	0.85
1:A:22:SER:HG	1:A:25:ASP:HB2	1.40	0.85
1:A:20:ALA:O	1:A:32:LYS:HD2	1.77	0.84
1:A:20:ALA:HB3	1:A:32:LYS:HE2	1.61	0.83
1:B:20:ALA:HB3	1:B:32:LYS:HE3	0.87	0.83
1:D:85:GLN:HE22	1:D:113:LYS:HZ3	1.28	0.80
1:D:85:GLN:HE22	1:D:113:LYS:HZ2	1.31	0.75
1:B:164:VAL:O	1:B:164:VAL:HG13	1.86	0.74
1:C:5:VAL:HG12	1:C:6:ASN:ND2	2.03	0.73
1:A:22:SER:OG	1:A:25:ASP:CB	2.34	0.73
1:A:25:ASP:O	1:A:28:SER:N	2.21	0.73
1:D:9:PHE:CD2	1:D:43:ILE:HG13	2.24	0.72
1:D:22:SER:HB2	1:D:25:ASP:CB	2.12	0.72
1:C:48:LYS:HE2	1:C:176:LEU:CD2	2.19	0.71
1:A:4:LEU:HD11	1:A:126:ILE:HB	1.72	0.71
1:B:148:ASN:O	1:B:148:ASN:OD1	2.09	0.71
1:D:85:GLN:NE2	1:D:113:LYS:HZ3	1.86	0.70
1:B:25:ASP:C	1:B:28:SER:O	2.31	0.69
1:C:48:LYS:HE2	1:C:176:LEU:HD22	1.75	0.69
1:A:20:ALA:HB1	1:A:32:LYS:HE2	1.73	0.68
1:C:64:VAL:O	1:C:68:PRO:HG2	1.93	0.68
1:D:56:ALA:HB3	1:D:59:SER:HB3	1.75	0.68
1:C:22:SER:OG	1:C:25:ASP:HB2	1.94	0.67
1:D:59:SER:OG	1:D:62:CYS:HB2	1.95	0.66
1:C:128:PHE:HD1	1:C:156:LYS:HD3	1.63	0.64
1:B:25:ASP:OD1	1:B:28:SER:CB	2.46	0.63
1:B:22:SER:OG	1:B:25:ASP:HB2	1.99	0.63
1:B:25:ASP:OD1	1:B:28:SER:HB2	1.99	0.62
1:D:159:ASN:O	1:D:163:ASP:HB2	2.00	0.62
1:C:74:LEU:O	1:C:78:VAL:HG23	1.99	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:84:ASP:O	1:D:85:GLN:HG2	2.01	0.60
1:C:73:TYR:O	1:C:74:LEU:C	2.41	0.59
1:D:85:GLN:HE22	1:D:113:LYS:CE	2.16	0.58
1:D:23:GLN:HB3	1:D:107:LYS:HB3	1.85	0.57
1:D:85:GLN:NE2	1:D:113:LYS:HZ2	1.94	0.57
1:A:25:ASP:O	1:A:28:SER:C	2.44	0.56
1:D:23:GLN:HG3	1:D:24:SER:H	1.70	0.56
1:C:164:VAL:HG13	1:C:164:VAL:O	2.05	0.56
1:B:9:PHE:CD2	1:B:43:ILE:HG13	2.41	0.56
1:C:128:PHE:CD1	1:C:156:LYS:HD3	2.41	0.56
1:A:49:VAL:HG22	1:A:85:GLN:HB2	1.89	0.55
1:B:25:ASP:CA	1:B:28:SER:O	2.55	0.54
1:A:23:GLN:O	1:A:26:ALA:HB3	2.06	0.54
1:D:52:THR:OG1	1:D:66:HIS:HE1	1.89	0.54
1:D:21:ILE:HG12	1:D:98:GLN:HE21	1.73	0.54
1:C:77:LEU:HA	1:C:81:LYS:HB2	1.90	0.54
1:C:165:THR:OG1	1:C:166:VAL:N	2.40	0.53
1:B:9:PHE:CE2	1:B:43:ILE:HG13	2.44	0.53
1:D:60:PRO:O	1:D:64:VAL:HG23	2.08	0.52
1:B:33:MET:HB3	1:B:34:PRO:HD2	1.91	0.52
1:B:164:VAL:O	1:B:164:VAL:CG1	2.55	0.52
1:D:67:ILE:O	1:D:70:TYR:N	2.41	0.52
1:A:41:LYS:O	1:A:45:GLU:HB2	2.10	0.52
1:D:39:TRP:O	1:D:43:ILE:HG12	2.09	0.51
1:D:67:ILE:O	1:D:68:PRO:C	2.48	0.51
1:D:69:GLY:O	1:D:73:TYR:HD1	1.94	0.51
1:D:147:GLU:OE1	1:D:152:THR:HG21	2.10	0.51
1:C:165:THR:O	1:C:168:SER:CB	2.59	0.51
1:D:93:ASN:HB2	1:D:94:PRO:HD2	1.93	0.51
1:A:108:ASP:OD1	1:A:108:ASP:C	2.49	0.50
1:B:110:THR:HG22	1:B:111:HIS:CD2	2.46	0.50
1:C:165:THR:O	1:C:168:SER:HB3	2.12	0.50
1:B:25:ASP:OD1	1:B:28:SER:HB3	2.11	0.50
1:C:5:VAL:HG12	1:C:6:ASN:CG	2.32	0.49
1:D:159:ASN:HB2	1:D:163:ASP:OD2	2.11	0.49
1:C:93:ASN:HB2	1:C:94:PRO:HD2	1.94	0.48
1:A:25:ASP:O	1:A:28:SER:O	2.31	0.48
1:C:22:SER:CB	1:C:25:ASP:HB2	2.43	0.48
1:D:67:ILE:HG21	1:D:114:PHE:HZ	1.79	0.47
1:A:173:LEU:O	1:A:176:LEU:HB2	2.15	0.47
1:A:21:ILE:O	1:A:108:ASP:HA	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:74:LEU:HD11	1:C:111:HIS:HB2	1.96	0.47
1:A:25:ASP:O	1:A:28:SER:CA	2.62	0.47
1:C:66:HIS:ND1	1:C:67:ILE:HD13	2.30	0.47
1:D:58:PHE:CZ	1:D:100:TRP:HA	2.49	0.47
1:A:8:LYS:HG2	1:A:9:PHE:N	2.30	0.47
1:C:163:ASP:N	1:C:163:ASP:OD1	2.47	0.47
1:B:50:ILE:HB	1:B:86:VAL:HG13	1.97	0.46
1:D:26:ALA:C	1:D:28:SER:N	2.69	0.46
1:A:173:LEU:HA	1:A:176:LEU:HD12	1.97	0.46
1:D:59:SER:OG	1:D:62:CYS:SG	2.74	0.46
1:D:87:ILE:HG22	1:D:89:VAL:HG23	1.97	0.46
1:B:25:ASP:HA	1:B:28:SER:HB2	1.98	0.46
1:C:141:ARG:O	1:C:157:GLU:HG2	2.15	0.45
1:D:52:THR:OG1	1:D:66:HIS:CE1	2.69	0.45
1:B:17:GLN:HA	1:B:35:GLN:O	2.17	0.45
1:B:50:ILE:HD11	1:B:83:VAL:HG21	1.99	0.45
1:C:9:PHE:HD2	1:C:11:ALA:HB2	1.81	0.45
1:C:164:VAL:O	1:C:164:VAL:HG22	2.17	0.45
1:D:54:ALA:O	1:D:90:THR:HG22	2.17	0.45
1:C:23:GLN:O	1:C:26:ALA:HB3	2.17	0.45
1:D:144:MET:HA	1:D:153:TYR:O	2.17	0.45
1:C:67:ILE:N	1:C:68:PRO:CD	2.80	0.44
1:A:50:ILE:HD13	1:A:172:VAL:HG11	1.99	0.44
1:D:9:PHE:CE2	1:D:43:ILE:HG13	2.52	0.44
1:D:41:LYS:O	1:D:45:GLU:HB2	2.18	0.44
1:A:19:ILE:HG23	1:A:32:LYS:HB3	1.99	0.44
1:B:48:LYS:HB3	1:B:83:VAL:HA	2.00	0.44
1:D:165:THR:OG1	1:D:166:VAL:N	2.50	0.44
1:A:120:CYS:SG	1:A:123:THR:HB	2.58	0.44
1:D:22:SER:OG	1:D:25:ASP:HB3	2.17	0.44
1:A:66:HIS:HA	1:A:164:VAL:HG21	2.00	0.43
1:B:60:PRO:O	1:B:64:VAL:HB	2.18	0.43
1:D:56:ALA:O	1:D:59:SER:OG	2.33	0.43
1:A:129:GLU:OE2	1:A:129:GLU:N	2.42	0.43
1:A:165:THR:OG1	1:A:166:VAL:N	2.48	0.43
1:C:53:GLY:O	1:C:141:ARG:HA	2.18	0.43
1:A:33:MET:HE2	1:A:95:PHE:HE1	1.83	0.43
1:D:156:LYS:HG2	1:D:157:GLU:O	2.18	0.43
1:A:16:PHE:HB2	1:A:116:SER:O	2.19	0.43
1:A:93:ASN:OD1	1:A:95:PHE:HB2	2.19	0.43
1:D:69:GLY:HA2	1:D:72:ASN:HD22	1.84	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:128:PHE:CE1	1:C:156:LYS:HB2	2.54	0.42
1:C:18:TYR:HB2	1:C:114:PHE:O	2.18	0.42
1:D:127:GLY:HA2	1:D:129:GLU:OE2	2.19	0.42
1:B:74:LEU:O	1:B:78:VAL:HG23	2.18	0.42
1:C:54:ALA:HA	1:C:55:PRO:HD3	1.94	0.42
1:B:21:ILE:HG12	1:B:98:GLN:HE21	1.84	0.42
1:B:33:MET:HB3	1:B:34:PRO:CD	2.48	0.42
1:D:19:ILE:CD1	1:D:95:PHE:HA	2.50	0.42
1:D:23:GLN:HG3	1:D:24:SER:N	2.34	0.42
1:D:78:VAL:HG12	1:D:79:LYS:HG3	2.02	0.42
1:D:54:ALA:O	1:D:139:SER:HB3	2.19	0.42
1:D:65:SER:O	1:D:164:VAL:HG21	2.20	0.42
1:D:59:SER:OG	1:D:62:CYS:CB	2.65	0.41
1:A:17:GLN:HA	1:A:35:GLN:O	2.19	0.41
1:B:75:ASP:CB	1:C:159:ASN:HD21	2.34	0.41
1:C:20:ALA:O	1:C:32:LYS:HG2	2.21	0.41
1:A:33:MET:CE	1:A:95:PHE:HE1	2.33	0.41
1:C:50:ILE:HD11	1:C:176:LEU:HD11	2.02	0.41
1:C:55:PRO:HD2	1:C:62:CYS:SG	2.61	0.41
1:C:123:THR:HG21	1:C:139:SER:HA	2.03	0.41
1:D:30:SER:O	1:D:32:LYS:HD2	2.21	0.41
1:D:70:TYR:HA	1:D:169:VAL:HG21	2.02	0.40
1:D:76:GLU:O	1:D:80:GLU:HB2	2.21	0.40
1:B:4:LEU:HD11	1:B:126:ILE:O	2.21	0.40
1:C:141:ARG:NE	1:C:160:PRO:O	2.48	0.40
1:C:41:LYS:O	1:C:45:GLU:HG3	2.20	0.40
1:C:81:LYS:HA	1:C:81:LYS:HD2	1.74	0.40
1:D:19:ILE:HD12	1:D:95:PHE:HA	2.02	0.40
1:D:26:ALA:O	1:D:28:SER:O	2.40	0.40
1:D:33:MET:HB3	1:D:34:PRO:CD	2.51	0.40
1:D:141:ARG:HB3	1:D:157:GLU:CG	2.52	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	172/184 (94%)	154 (90%)	17 (10%)	1 (1%)	25	57
1	B	172/184 (94%)	160 (93%)	12 (7%)	0	100	100
1	C	172/184 (94%)	152 (88%)	20 (12%)	0	100	100
1	D	172/184 (94%)	150 (87%)	20 (12%)	2 (1%)	13	38
All	All	688/736 (94%)	616 (90%)	69 (10%)	3 (0%)	34	65

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	148	ASN
1	D	170	GLU
1	D	27	ASP

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	147/156 (94%)	147 (100%)	0	100	100
1	B	147/156 (94%)	147 (100%)	0	100	100
1	C	147/156 (94%)	147 (100%)	0	100	100
1	D	147/156 (94%)	147 (100%)	0	100	100
All	All	588/624 (94%)	588 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	6	ASN
1	A	97	ASN
1	B	17	GLN

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Mol	Chain	Res	Type
1	B	23	GLN
1	B	97	ASN
1	B	175	HIS
1	C	6	ASN
1	C	97	ASN
1	C	98	GLN
1	C	159	ASN
1	D	66	HIS
1	D	72	ASN
1	D	85	GLN
1	D	97	ASN
1	D	98	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 monosaccharides 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, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	A	174/184 (94%)	0.05	7 (4%)	38	35	52, 61, 68, 73	8 (4%)
1	B	174/184 (94%)	0.14	4 (2%)	60	59	51, 60, 68, 75	8 (4%)
1	C	174/184 (94%)	0.28	12 (6%)	16	13	54, 64, 72, 76	8 (4%)
1	D	174/184 (94%)	0.19	10 (5%)	23	20	54, 65, 72, 75	8 (4%)
All	All	696/736 (94%)	0.17	33 (4%)	31	28	51, 62, 72, 76	32 (4%)

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	31	CYS	8.5
1	B	31	CYS	6.5
1	C	31	CYS	6.5
1	D	33	MET	5.6
1	C	23	GLN	5.3
1	B	27	ASP	4.9
1	D	21	ILE	4.2
1	C	24	SER	3.9
1	A	149	GLY	3.3
1	D	27	ASP	3.3
1	C	29	GLU	3.2
1	C	34	PRO	3.0
1	D	24	SER	2.9
1	A	29	GLU	2.9
1	D	26	ALA	2.8
1	A	24	SER	2.8
1	C	65	SER	2.8
1	C	18	TYR	2.7
1	C	66	HIS	2.7
1	C	33	MET	2.7
1	A	25	ASP	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	27	ASP	2.6
1	A	31	CYS	2.4
1	C	64	VAL	2.3
1	A	26	ALA	2.3
1	C	27	ASP	2.2
1	D	30	SER	2.1
1	B	33	MET	2.1
1	C	30	SER	2.1
1	D	111	HIS	2.1
1	D	109	THR	2.0
1	D	72	ASN	2.0
1	B	176	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 monosaccharides 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.