Full wwPDB X-ray Structure Validation Report (i)

Jun 17, 2024 – 08:54 PM EDT

PDB ID	:	5Y82
Title	:	Crystal structure of the periplasmic domain of the Thermotoga maritima YidC
Authors	:	Huang, Y.; Xin, Y.
Deposited on	:	2017-08-18
Resolution	:	2.52 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:

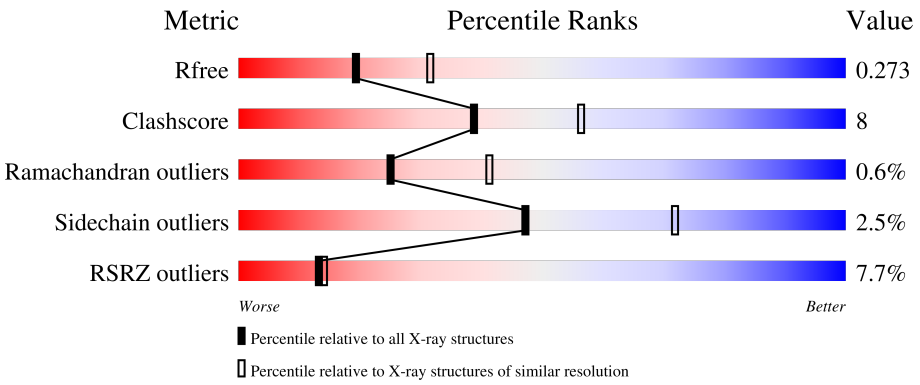
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

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 2.52 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	5743 (2.54-2.50)
Clashscore	141614	6463 (2.54-2.50)
Ramachandran outliers	138981	6335 (2.54-2.50)
Sidechain outliers	138945	6337 (2.54-2.50)
RSRZ outliers	127900	5630 (2.54-2.50)

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	205	11% 76% 20% • •
1	B	205	6% 74% 22% • •
1	C	205	7% 81% 15% •
1	D	205	5% 74% 20% • •

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 6581 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 Membrane protein insertase YidC.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	D	199	1611	1045	258	304	4	0	0	0
1	A	199	1611	1045	258	304	4	0	0	0
1	B	199	1611	1045	258	304	4	0	0	0
1	C	199	1611	1045	258	304	4	0	0	0

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	57	MET	LEU	engineered mutation	UNP Q9X1H2
D	105	MET	ILE	engineered mutation	UNP Q9X1H2
D	153	MET	LEU	engineered mutation	UNP Q9X1H2
D	223	HIS	-	expression tag	UNP Q9X1H2
D	224	HIS	-	expression tag	UNP Q9X1H2
D	225	HIS	-	expression tag	UNP Q9X1H2
D	226	HIS	-	expression tag	UNP Q9X1H2
D	227	HIS	-	expression tag	UNP Q9X1H2
D	228	HIS	-	expression tag	UNP Q9X1H2
A	57	MET	LEU	engineered mutation	UNP Q9X1H2
A	105	MET	ILE	engineered mutation	UNP Q9X1H2
A	153	MET	LEU	engineered mutation	UNP Q9X1H2
A	223	HIS	-	expression tag	UNP Q9X1H2
A	224	HIS	-	expression tag	UNP Q9X1H2
A	225	HIS	-	expression tag	UNP Q9X1H2
A	226	HIS	-	expression tag	UNP Q9X1H2
A	227	HIS	-	expression tag	UNP Q9X1H2
A	228	HIS	-	expression tag	UNP Q9X1H2
B	57	MET	LEU	engineered mutation	UNP Q9X1H2
B	105	MET	ILE	engineered mutation	UNP Q9X1H2
B	153	MET	LEU	engineered mutation	UNP Q9X1H2

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Chain	Residue	Modelled	Actual	Comment	Reference
B	223	HIS	-	expression tag	UNP Q9X1H2
B	224	HIS	-	expression tag	UNP Q9X1H2
B	225	HIS	-	expression tag	UNP Q9X1H2
B	226	HIS	-	expression tag	UNP Q9X1H2
B	227	HIS	-	expression tag	UNP Q9X1H2
B	228	HIS	-	expression tag	UNP Q9X1H2
C	57	MET	LEU	engineered mutation	UNP Q9X1H2
C	105	MET	ILE	engineered mutation	UNP Q9X1H2
C	153	MET	LEU	engineered mutation	UNP Q9X1H2
C	223	HIS	-	expression tag	UNP Q9X1H2
C	224	HIS	-	expression tag	UNP Q9X1H2
C	225	HIS	-	expression tag	UNP Q9X1H2
C	226	HIS	-	expression tag	UNP Q9X1H2
C	227	HIS	-	expression tag	UNP Q9X1H2
C	228	HIS	-	expression tag	UNP Q9X1H2

- Molecule 2 is water.

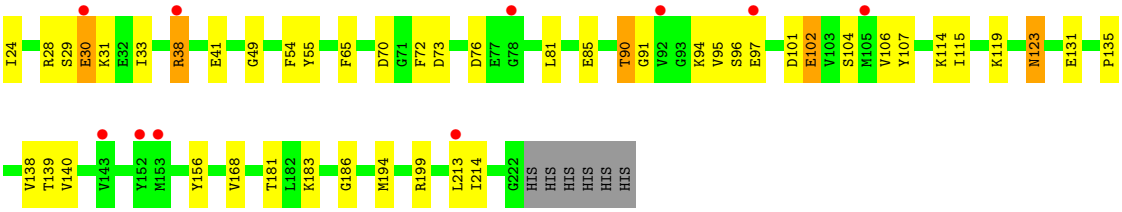
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	D	36	36	36	0	0
			Total	O		
2	A	28	28	28	0	0
			Total	O		
2	B	31	31	31	0	0
			Total	O		
2	C	42	42	42	0	0

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.

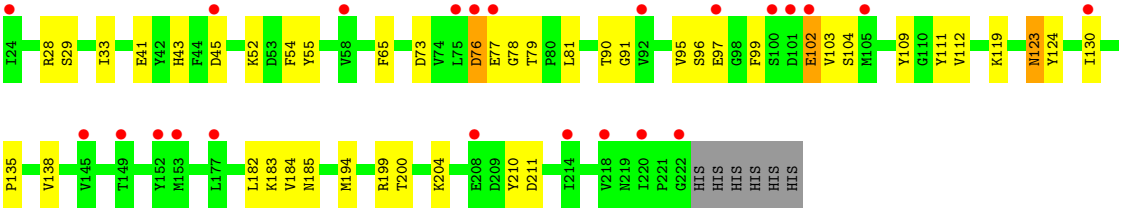
• Molecule 1: Membrane protein insertase YidC

Chain D:



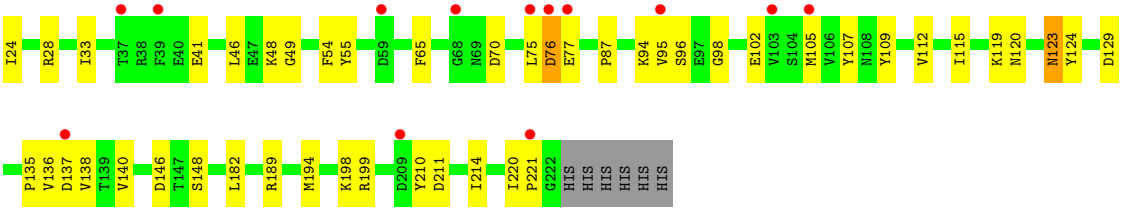
• Molecule 1: Membrane protein insertase YidC

Chain A:



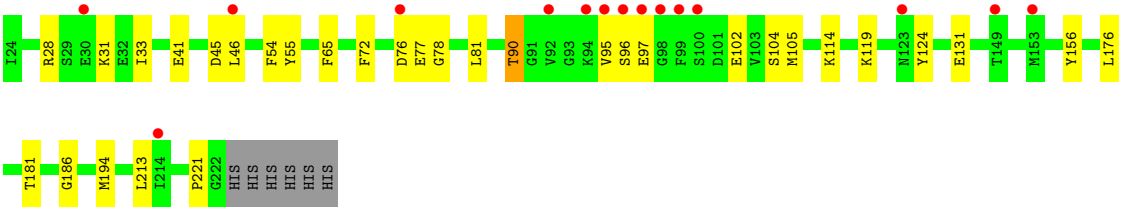
• Molecule 1: Membrane protein insertase YidC

Chain B:



• Molecule 1: Membrane protein insertase YidC

Chain C:


I24
R28
S29
E30
K31
E32
I33
E41
D45
L46
F54
Y55
F65
F72
D76
E77
G78
L81
T90
G91
V92
G93
K94
V95
S96
E97
G98
F99
S100
D101
E102
V103
S104
M105
K114
K119
N123
Y124
E131
T149
M153
Y156
L176
T181
G186
M194
L213
I214
P221
G222
HIS
HIS
HIS
HIS
HIS
HIS

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, α, β, γ	79.95Å 79.95Å 135.82Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.97 – 2.52 29.97 – 2.52	Depositor EDS
% Data completeness (in resolution range)	88.2 (29.97-2.52) 88.0 (29.97-2.52)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.94 (at 2.51Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.221 , 0.261 0.230 , 0.273	Depositor DCC
R_{free} test set	1349 reflections (5.11%)	wwPDB-VP
Wilson B-factor (Å^2)	33.7	Xtriage
Anisotropy	0.060	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 18.5	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	0.449 for h,-k,-l	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	6581	wwPDB-VP
Average B, all atoms (Å^2)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.09% 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	0.42	0/1646	0.55	0/2225
1	B	0.41	0/1646	0.56	0/2225
1	C	0.41	0/1646	0.57	0/2225
1	D	0.42	0/1646	0.56	0/2225
All	All	0.41	0/6584	0.56	0/8900

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	1611	0	1613	25	0
1	B	1611	0	1613	27	0
1	C	1611	0	1613	18	0
1	D	1611	0	1613	28	0
2	A	28	0	0	1	1
2	B	31	0	0	2	1
2	C	42	0	0	2	0
2	D	36	0	0	4	0
All	All	6581	0	6452	97	1

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 8.

All (97) 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:D:38:ARG:NH1	2:D:301:HOH:O	1.94	0.98
1:D:76:ASP:HB2	1:D:81:LEU:HD21	1.57	0.84
1:C:78:GLY:O	2:C:302:HOH:O	2.00	0.79
1:C:181:THR:OG1	2:C:301:HOH:O	1.94	0.78
1:D:94:LYS:NZ	2:D:303:HOH:O	2.13	0.74
1:D:135:PRO:O	2:D:302:HOH:O	2.05	0.73
1:C:76:ASP:HB2	1:C:81:LEU:HD21	1.76	0.66
1:A:211:ASP:N	1:A:211:ASP:OD1	2.28	0.65
1:C:131:GLU:HA	1:C:186:GLY:O	1.99	0.63
1:B:77:GLU:N	2:B:303:HOH:O	2.32	0.62
1:B:210:TYR:CE1	1:B:214:ILE:HD12	2.35	0.62
1:D:106:VAL:HG22	1:D:115:ILE:HG12	1.83	0.61
1:A:54:PHE:HB3	1:A:65:PHE:HB3	1.83	0.61
1:B:109:TYR:HB2	1:B:112:VAL:HB	1.81	0.61
1:A:76:ASP:OD2	1:A:111:TYR:OH	2.14	0.60
1:C:102:GLU:HG2	1:C:119:LYS:HG2	1.83	0.60
1:D:131:GLU:HA	1:D:186:GLY:O	2.02	0.60
1:B:41:GLU:HB3	1:B:55:TYR:HB2	1.85	0.59
1:A:77:GLU:N	2:A:301:HOH:O	2.37	0.58
1:D:54:PHE:HB3	1:D:65:PHE:HB3	1.84	0.58
1:C:95:VAL:HG13	1:C:97:GLU:H	1.69	0.57
1:C:28:ARG:HG3	1:C:33:ILE:HG12	1.87	0.56
1:C:54:PHE:HB3	1:C:65:PHE:HB3	1.87	0.56
1:B:54:PHE:HB3	1:B:65:PHE:HB3	1.87	0.55
1:C:124:TYR:HB3	1:C:194:MET:O	2.06	0.55
1:C:41:GLU:HB3	1:C:55:TYR:HB2	1.89	0.55
1:B:124:TYR:HB3	1:B:194:MET:O	2.06	0.55
1:A:109:TYR:HB2	1:A:112:VAL:HB	1.88	0.54
1:D:41:GLU:HB3	1:D:55:TYR:HB2	1.89	0.54
1:C:72:PHE:O	1:C:114:LYS:NZ	2.41	0.54
1:A:135:PRO:HB3	1:A:185:ASN:HB2	1.89	0.54
1:D:95:VAL:HG13	1:D:97:GLU:H	1.73	0.53
1:D:102:GLU:HG2	1:D:119:LYS:HG2	1.90	0.53
1:B:48:LYS:HD3	2:B:310:HOH:O	2.06	0.53
1:B:49:GLY:HA2	1:B:107:TYR:CE1	2.44	0.53
1:D:95:VAL:HG22	1:D:96:SER:H	1.73	0.53
1:A:130:ILE:HG23	1:A:184:VAL:HG11	1.90	0.53
1:C:95:VAL:HG22	1:C:96:SER:H	1.72	0.52
1:B:24:ILE:HD12	1:B:120:ASN:HD21	1.75	0.52
1:A:123:ASN:O	1:A:123:ASN:ND2	2.42	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:85:GLU:OE1	2:D:304:HOH:O	2.18	0.52
1:D:90:THR:HG23	1:D:104:SER:HB2	1.91	0.51
1:C:90:THR:HG23	1:C:104:SER:HB2	1.92	0.51
1:A:138:VAL:HB	1:A:182:LEU:HB2	1.93	0.51
1:A:135:PRO:HA	1:A:184:VAL:O	2.11	0.51
1:D:29:SER:C	1:D:31:LYS:H	2.14	0.51
1:A:124:TYR:HB3	1:A:194:MET:O	2.11	0.50
1:B:123:ASN:O	1:B:123:ASN:ND2	2.38	0.50
1:B:28:ARG:HG3	1:B:33:ILE:HG12	1.94	0.49
1:A:41:GLU:HB3	1:A:55:TYR:HB2	1.95	0.49
1:D:181:THR:OG1	1:D:183:LYS:NZ	2.45	0.49
1:D:123:ASN:O	1:D:123:ASN:ND2	2.41	0.48
1:D:70:ASP:HA	1:D:139:THR:OG1	2.14	0.48
1:B:138:VAL:HB	1:B:182:LEU:HB2	1.96	0.48
1:A:73:ASP:O	1:A:138:VAL:HA	2.14	0.47
1:A:102:GLU:HG2	1:A:119:LYS:HG2	1.97	0.47
1:A:199:ARG:HG2	1:A:210:TYR:OH	2.15	0.47
1:A:90:THR:HG23	1:A:104:SER:HB2	1.97	0.47
1:B:129:ASP:OD1	1:B:189:ARG:NE	2.28	0.47
1:D:73:ASP:O	1:D:138:VAL:HA	2.15	0.46
1:A:95:VAL:HG13	1:A:97:GLU:H	1.79	0.46
1:D:101:ASP:HB3	1:B:135:PRO:HG3	1.96	0.46
1:B:95:VAL:HG22	1:B:96:SER:H	1.80	0.46
1:C:31:LYS:HB3	1:C:31:LYS:HE2	1.70	0.46
1:A:28:ARG:HG3	1:A:33:ILE:HG12	1.98	0.46
1:A:200:THR:HG22	1:A:204:LYS:HE3	1.96	0.46
1:D:49:GLY:HA2	1:D:107:TYR:CE1	2.51	0.46
1:B:76:ASP:HA	1:B:136:VAL:HG12	1.97	0.46
1:D:90:THR:OG1	1:D:91:GLY:N	2.49	0.45
1:B:146:ASP:OD2	1:B:148:SER:HB3	2.17	0.45
1:C:176:LEU:HD23	1:C:176:LEU:HA	1.84	0.44
1:A:99:PHE:CD1	1:A:103:VAL:HG22	2.51	0.44
1:B:70:ASP:HB2	1:B:140:VAL:O	2.17	0.44
1:B:28:ARG:NH2	1:B:87:PRO:O	2.44	0.44
1:B:220:ILE:HG22	1:B:221:PRO:HD2	1.99	0.44
1:B:33:ILE:HG13	1:B:46:LEU:HD21	2.00	0.44
1:A:95:VAL:HG22	1:A:96:SER:H	1.83	0.43
1:A:183:LYS:HD2	1:A:183:LYS:H	1.82	0.43
1:C:33:ILE:HG13	1:C:46:LEU:HD21	2.01	0.43
1:D:156:TYR:CE2	1:D:213:LEU:HD11	2.54	0.42
1:C:156:TYR:CE2	1:C:213:LEU:HD11	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:43:HIS:O	1:A:52:LYS:N	2.50	0.42
1:D:28:ARG:HG3	1:D:33:ILE:HG12	2.00	0.42
1:B:105:MET:O	1:B:115:ILE:HA	2.19	0.42
1:C:33:ILE:HB	1:C:105:MET:HE1	2.02	0.42
1:D:199:ARG:HG3	1:D:214:ILE:HD13	2.02	0.41
1:B:75:LEU:HB2	1:B:137:ASP:HB2	2.02	0.41
1:B:146:ASP:N	1:B:146:ASP:OD1	2.53	0.41
1:B:95:VAL:N	1:B:98:GLY:O	2.51	0.41
1:D:72:PHE:O	1:D:114:LYS:NZ	2.53	0.41
1:B:198:LYS:HD2	1:B:198:LYS:HA	1.89	0.41
1:B:102:GLU:HG3	1:B:119:LYS:HG2	2.03	0.41
1:D:49:GLY:HA2	1:D:107:TYR:CZ	2.56	0.41
1:D:70:ASP:HB2	1:D:140:VAL:O	2.21	0.40
1:D:168:VAL:HG22	1:D:194:MET:HG3	2.03	0.40
1:A:76:ASP:HB2	1:A:81:LEU:HD21	2.03	0.40
1:A:90:THR:OG1	1:A:91:GLY:N	2.54	0.40

All (1) 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)
2:A:324:HOH:O	2:B:328:HOH:O[3_745]	1.85	0.35

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	197/205 (96%)	181 (92%)	14 (7%)	2 (1%)	15 27
1	B	197/205 (96%)	182 (92%)	13 (7%)	2 (1%)	15 27
1	C	197/205 (96%)	176 (89%)	21 (11%)	0	100 100
1	D	197/205 (96%)	180 (91%)	16 (8%)	1 (0%)	29 47

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	788/820 (96%)	719 (91%)	64 (8%)	5 (1%)	25 41

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	76	ASP
1	A	78	GLY
1	D	30	GLU
1	B	76	ASP
1	B	94	LYS

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	182/188 (97%)	177 (97%)	5 (3%)	44 69
1	B	182/188 (97%)	179 (98%)	3 (2%)	62 82
1	C	182/188 (97%)	178 (98%)	4 (2%)	52 75
1	D	182/188 (97%)	176 (97%)	6 (3%)	38 62
All	All	728/752 (97%)	710 (98%)	18 (2%)	47 72

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	24	ILE
1	D	30	GLU
1	D	38	ARG
1	D	90	THR
1	D	102	GLU
1	D	123	ASN
1	A	29	SER
1	A	45	ASP
1	A	79	THR
1	A	102	GLU

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Mol	Chain	Res	Type
1	A	123	ASN
1	B	123	ASN
1	B	199	ARG
1	B	211	ASP
1	C	45	ASP
1	C	77	GLU
1	C	90	THR
1	C	221	PRO

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	120	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 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	199/205 (97%)	0.79	23 (11%)	4	4	18, 33, 58, 69	0
1	B	199/205 (97%)	0.68	13 (6%)	18	20	18, 33, 59, 73	0
1	C	199/205 (97%)	0.59	15 (7%)	14	15	16, 30, 56, 86	0
1	D	199/205 (97%)	0.53	10 (5%)	28	31	18, 30, 59, 81	0
All	All	796/820 (97%)	0.65	61 (7%)	13	14	16, 32, 59, 86	0

All (61) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	97	GLU	5.9
1	B	77	GLU	5.6
1	A	97	GLU	4.9
1	A	145	VAL	4.8
1	C	76	ASP	4.3
1	D	92	VAL	4.1
1	A	76	ASP	4.1
1	A	218	VAL	3.8
1	C	96	SER	3.8
1	D	38	ARG	3.6
1	D	78	GLY	3.6
1	D	30	GLU	3.6
1	A	101	ASP	3.5
1	C	97	GLU	3.5
1	A	220	ILE	3.5
1	D	143	VAL	3.4
1	A	58	VAL	3.3
1	A	77	GLU	3.3
1	A	222	GLY	3.3
1	C	92	VAL	3.3
1	C	98	GLY	3.3

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Mol	Chain	Res	Type	RSRZ
1	B	59	ASP	3.3
1	C	99	PHE	3.2
1	A	24	ILE	3.1
1	C	95	VAL	3.0
1	B	95	VAL	3.0
1	A	102	GLU	3.0
1	A	177	LEU	2.9
1	C	100	SER	2.8
1	D	153	MET	2.8
1	A	153	MET	2.7
1	C	46	LEU	2.6
1	C	94	LYS	2.6
1	A	92	VAL	2.6
1	D	152	TYR	2.6
1	B	103	VAL	2.5
1	D	105	MET	2.5
1	B	39	PHE	2.5
1	B	75	LEU	2.4
1	C	123	ASN	2.4
1	A	214	ILE	2.4
1	B	76	ASP	2.4
1	C	149	THR	2.4
1	A	208	GLU	2.3
1	B	137	ASP	2.3
1	B	221	PRO	2.3
1	B	37	THR	2.3
1	B	68	GLY	2.3
1	C	30	GLU	2.2
1	A	105	MET	2.2
1	D	213	LEU	2.2
1	B	209	ASP	2.2
1	A	100	SER	2.2
1	A	152	TYR	2.1
1	A	149	THR	2.1
1	A	130	ILE	2.1
1	A	45	ASP	2.1
1	C	214	ILE	2.1
1	A	75	LEU	2.1
1	C	153	MET	2.0
1	B	105	MET	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.