



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 30, 2017 – 11:36 PM EST

PDB ID : 1C03  
Title : CRYSTAL STRUCTURE OF YPD1P (TRICLINIC FORM)  
Authors : Song, H.K.; Lee, J.Y.; Lee, M.G.; Suh, S.W.  
Deposited on : 1999-07-14  
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  
Mogul : unknown  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20028442  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
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) : rb-20028442

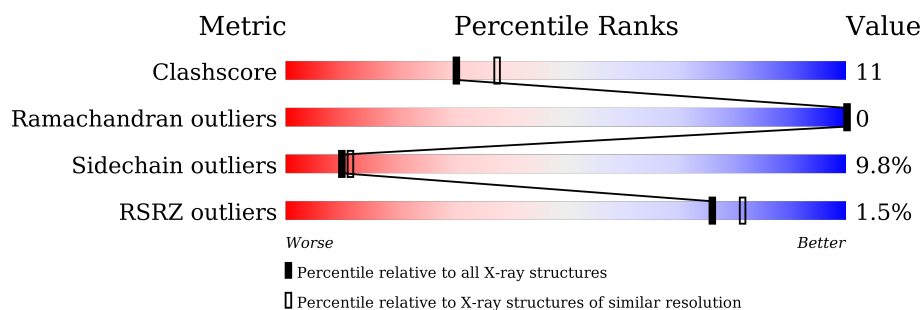
# 1 Overall quality at a glance

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))
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (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	168	<div> <div>2%</div> <div>73% 17% 7% .</div> </div>
1	B	168	<div> <div>2%</div> <div>72% 18% 7% .</div> </div>
1	C	168	<div> <div>2%</div> <div>74% 18% 5% .</div> </div>
1	D	168	<div> <div>%</div> <div>72% 20% 5% .</div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5396 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 HYPOTHETICAL PROTEIN YDL235C.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	163	Total	C	N	O	S	0	0	0
			1305	821	220	260	4			
1	B	163	Total	C	N	O	S	0	0	0
			1305	821	220	260	4			
1	C	163	Total	C	N	O	S	0	0	0
			1305	821	220	260	4			
1	D	163	Total	C	N	O	S	0	0	0
			1305	821	220	260	4			

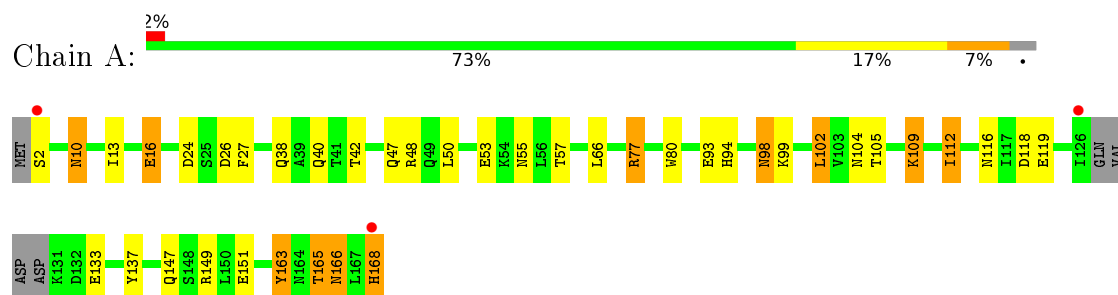
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	49	Total	O	0	0
			49	49		
2	B	44	Total	O	0	0
			44	44		
2	C	48	Total	O	0	0
			48	48		
2	D	35	Total	O	0	0
			35	35		

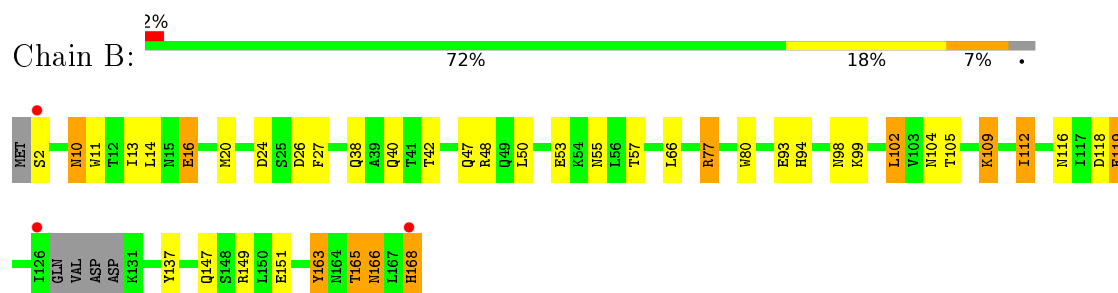
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

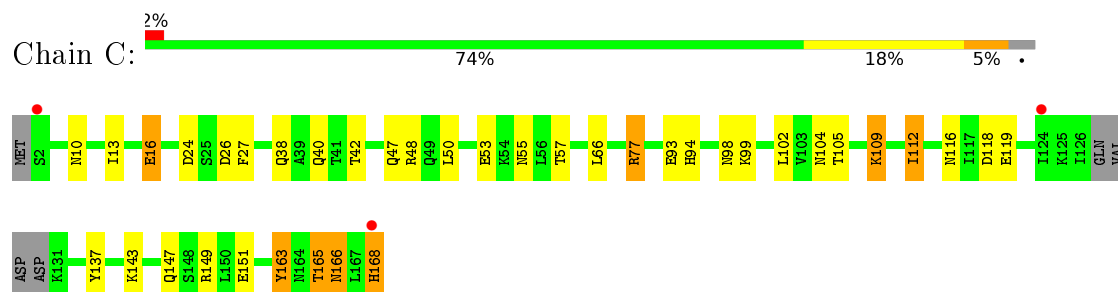
#### • Molecule 1: HYPOTHETICAL PROTEIN YDL235C



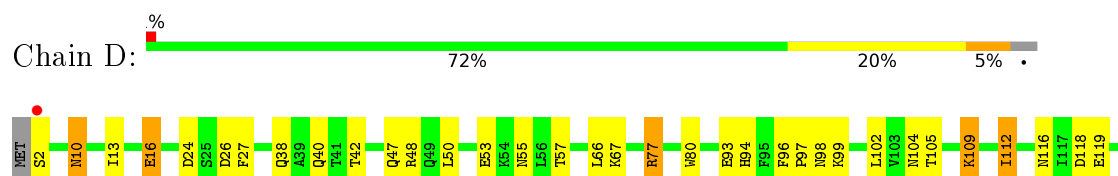
#### • Molecule 1: HYPOTHETICAL PROTEIN YDL235C



#### • Molecule 1: HYPOTHETICAL PROTEIN YDL235C



#### • Molecule 1: HYPOTHETICAL PROTEIN YDL235C



I126	GLN		Y137		Q147		Y163		M164		T165		M166		L167		H168
	VAL				S148												
	ASP				R149												
	ASP				L150												
	K131				E151												

## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	66.34Å 66.48Å 66.49Å 106.37° 106.66° 115.33°	Depositor
Resolution (Å)	8.00 – 2.30 19.92 – 2.30	Depositor EDS
% Data completeness (in resolution range)	0.9 (8.00-2.30) 71.6 (19.92-2.30)	Depositor EDS
$R_{merge}$	0.04	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.45 (at 2.30Å)	Xtriage
Refinement program	X-PLOR 3.851	Depositor
R, $R_{free}$	0.206 , 0.248 0.201 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	DCC
Wilson B-factor (Å <sup>2</sup> )	35.9	Xtriage
Anisotropy	0.146	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 70.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.52$ , $\langle L^2 \rangle = 0.36$	Xtriage
Estimated twinning fraction	0.014 for -l,h+k+l,-k 0.014 for h+k+l,-l,-h 0.470 for k,h,-h-k-l 0.014 for -k,-h,-l 0.477 for l,-h-k-l,h 0.477 for -h-k-l,l,k 0.015 for -h,-k,h+k+l	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	5396	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.99% 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	0.44	1/1322 (0.1%)	0.53	0/1783
1	B	0.44	1/1322 (0.1%)	0.53	0/1783
1	C	0.44	1/1322 (0.1%)	0.53	0/1783
1	D	0.44	1/1322 (0.1%)	0.53	0/1783
All	All	0.44	4/5288 (0.1%)	0.53	0/7132

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	A	0	1
1	B	0	1
1	C	0	1
1	D	0	1
All	All	0	4

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	93	GLU	CD-OE2	7.25	1.33	1.25
1	D	93	GLU	CD-OE2	7.08	1.33	1.25
1	C	93	GLU	CD-OE2	6.99	1.33	1.25
1	A	93	GLU	CD-OE2	6.59	1.32	1.25

There are no bond angle outliers.

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	163	TYR	Sidechain

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Mol	Chain	Res	Type	Group
1	B	163	TYR	Sidechain
1	C	163	TYR	Sidechain
1	D	163	TYR	Sidechain

## 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	1305	0	1280	33	0
1	B	1305	0	1280	33	0
1	C	1305	0	1280	30	0
1	D	1305	0	1280	31	0
2	A	49	0	0	3	0
2	B	44	0	0	1	0
2	C	48	0	0	2	0
2	D	35	0	0	1	0
All	All	5396	0	5120	114	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 11.

All (114) 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:47:GLN:HE21	1:D:149:ARG:HH12	1.28	0.80
1:C:116:ASN:O	1:C:119:GLU:HG3	1.82	0.80
1:B:47:GLN:HE21	1:B:149:ARG:HH12	1.29	0.80
1:A:116:ASN:O	1:A:119:GLU:HG3	1.82	0.79
1:D:116:ASN:O	1:D:119:GLU:HG3	1.81	0.79
1:B:47:GLN:NE2	1:B:149:ARG:HH12	1.82	0.78
1:A:47:GLN:HE21	1:A:149:ARG:HH12	1.31	0.78
1:B:116:ASN:O	1:B:119:GLU:HG3	1.83	0.78
1:D:47:GLN:NE2	1:D:149:ARG:HH12	1.81	0.78
1:A:47:GLN:NE2	1:A:149:ARG:HH12	1.83	0.77
1:C:47:GLN:NE2	1:C:149:ARG:HH12	1.83	0.77
1:C:47:GLN:HE21	1:C:149:ARG:HH12	1.34	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:94:HIS:HD2	1:C:137:TYR:OH	1.71	0.74
1:D:94:HIS:HD2	1:D:137:TYR:OH	1.70	0.73
1:B:94:HIS:HD2	1:B:137:TYR:OH	1.72	0.72
1:A:94:HIS:HD2	1:A:137:TYR:OH	1.71	0.72
1:A:99:LYS:HG3	1:A:147:GLN:HE22	1.58	0.69
1:A:119:GLU:HB3	2:A:641:HOH:O	1.94	0.66
1:B:48:ARG:HD3	1:C:40:GLN:HE21	1.60	0.66
1:A:48:ARG:HD3	1:D:40:GLN:HE21	1.62	0.64
1:B:16:GLU:HB3	1:D:13:ILE:HD11	1.80	0.64
1:C:99:LYS:HG3	1:C:147:GLN:HE22	1.62	0.63
1:D:99:LYS:HG3	1:D:147:GLN:HE22	1.63	0.62
1:A:109:LYS:HB3	1:A:112:ILE:HD11	1.82	0.62
1:A:16:GLU:HB3	1:C:13:ILE:HD11	1.83	0.61
1:B:99:LYS:HG3	1:B:147:GLN:HE22	1.63	0.61
1:B:40:GLN:HE21	1:C:48:ARG:HD3	1.66	0.61
1:B:13:ILE:HD11	1:D:16:GLU:HB3	1.83	0.60
1:B:166:ASN:HD21	1:B:168:HIS:CE1	2.19	0.60
1:D:166:ASN:HD21	1:D:168:HIS:CE1	2.19	0.60
1:A:163:TYR:O	1:A:165:THR:HG22	2.02	0.60
1:C:109:LYS:HB3	1:C:112:ILE:HD11	1.83	0.60
1:A:166:ASN:HD21	1:A:168:HIS:CE1	2.20	0.59
1:B:109:LYS:HB3	1:B:112:ILE:HD11	1.83	0.59
1:D:109:LYS:HB3	1:D:112:ILE:HD11	1.85	0.59
1:A:40:GLN:HE21	1:D:48:ARG:HD3	1.67	0.59
1:A:42:THR:HG22	2:A:740:HOH:O	2.02	0.59
1:D:163:TYR:O	1:D:165:THR:HG22	2.03	0.59
1:B:163:TYR:O	1:B:165:THR:HG22	2.04	0.57
1:C:166:ASN:HD21	1:C:168:HIS:CE1	2.21	0.57
1:B:42:THR:HG22	2:B:536:HOH:O	2.05	0.56
1:A:13:ILE:HD11	1:C:16:GLU:HB3	1.87	0.56
1:C:163:TYR:O	1:C:165:THR:HG22	2.05	0.56
1:A:133:GLU:HG2	2:A:633:HOH:O	2.05	0.56
1:C:42:THR:HG22	2:C:757:HOH:O	2.09	0.53
1:C:24:ASP:HB3	1:C:27:PHE:HB2	1.91	0.53
1:D:42:THR:OG1	1:D:66:LEU:HD22	2.09	0.52
1:A:42:THR:OG1	1:A:66:LEU:HD22	2.08	0.52
1:B:24:ASP:HB3	1:B:27:PHE:HB2	1.92	0.52
1:C:13:ILE:HG23	2:C:692:HOH:O	2.09	0.51
1:D:109:LYS:O	1:D:112:ILE:HG12	2.11	0.51
1:B:109:LYS:O	1:B:112:ILE:HG12	2.11	0.51
1:C:109:LYS:O	1:C:112:ILE:HG12	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:147:GLN:O	1:A:151:GLU:HG2	2.11	0.51
1:A:94:HIS:CD2	1:A:137:TYR:OH	2.60	0.51
1:D:24:ASP:HB3	1:D:27:PHE:HB2	1.93	0.50
1:A:24:ASP:HB3	1:A:27:PHE:HB2	1.93	0.50
1:B:42:THR:OG1	1:B:66:LEU:HD22	2.11	0.50
1:C:38:GLN:O	1:C:42:THR:HG23	2.11	0.50
1:A:38:GLN:O	1:A:42:THR:HG23	2.12	0.50
1:C:42:THR:OG1	1:C:66:LEU:HD22	2.12	0.50
1:D:38:GLN:O	1:D:42:THR:HG23	2.12	0.50
1:B:147:GLN:O	1:B:151:GLU:HG2	2.11	0.49
1:C:147:GLN:O	1:C:151:GLU:HG2	2.13	0.48
1:C:99:LYS:HG3	1:C:147:GLN:NE2	2.28	0.48
1:A:48:ARG:HD3	1:D:40:GLN:NE2	2.29	0.48
1:C:94:HIS:CD2	1:C:137:TYR:OH	2.61	0.48
1:D:147:GLN:O	1:D:151:GLU:HG2	2.13	0.48
1:B:94:HIS:CD2	1:B:137:TYR:OH	2.61	0.48
1:D:94:HIS:CD2	1:D:137:TYR:OH	2.59	0.47
1:D:42:THR:HG22	2:D:766:HOH:O	2.14	0.47
1:A:109:LYS:O	1:A:112:ILE:HG12	2.13	0.47
1:C:116:ASN:O	1:C:118:ASP:N	2.48	0.47
1:B:116:ASN:O	1:B:118:ASP:N	2.47	0.47
1:A:116:ASN:O	1:A:118:ASP:N	2.48	0.47
1:B:77:ARG:NH1	1:B:105:THR:OG1	2.49	0.46
1:B:38:GLN:O	1:B:42:THR:HG23	2.15	0.46
1:D:116:ASN:O	1:D:118:ASP:N	2.48	0.46
1:D:77:ARG:NH1	1:D:105:THR:OG1	2.49	0.46
1:A:99:LYS:HG3	1:A:147:GLN:NE2	2.26	0.45
1:B:48:ARG:HD3	1:C:40:GLN:NE2	2.30	0.45
1:C:77:ARG:NH1	1:C:105:THR:OG1	2.49	0.45
1:D:109:LYS:HB3	1:D:109:LYS:HZ3	1.83	0.44
1:A:98:ASN:HD22	1:A:98:ASN:HA	1.65	0.44
1:A:53:GLU:HG2	1:A:55:ASN:HB2	1.99	0.43
1:B:20:MET:HE2	1:D:67:LYS:HE2	2.01	0.43
1:A:40:GLN:NE2	1:D:48:ARG:HD3	2.32	0.43
1:C:116:ASN:O	1:C:119:GLU:CG	2.61	0.43
1:A:10:ASN:HD22	1:A:10:ASN:C	2.21	0.43
1:C:53:GLU:HG2	1:C:55:ASN:HB2	2.00	0.43
1:A:77:ARG:NH1	1:A:105:THR:OG1	2.52	0.42
1:B:40:GLN:NE2	1:C:48:ARG:HD3	2.33	0.42
1:D:53:GLU:HG2	1:D:55:ASN:HB2	2.01	0.42
1:B:53:GLU:HG2	1:B:55:ASN:HB2	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:143:LYS:HB3	1:C:143:LYS:HE2	1.80	0.42
1:B:20:MET:HB3	1:B:20:MET:HE2	1.92	0.42
1:C:109:LYS:HB3	1:C:109:LYS:HZ3	1.85	0.42
1:B:109:LYS:HZ3	1:B:109:LYS:HB3	1.85	0.42
1:B:10:ASN:C	1:B:10:ASN:HD22	2.23	0.42
1:A:116:ASN:O	1:A:119:GLU:CG	2.61	0.41
1:B:11:TRP:CZ3	1:B:14:LEU:HD23	2.54	0.41
1:D:109:LYS:NZ	1:D:109:LYS:HB3	2.35	0.41
1:D:10:ASN:HD22	1:D:10:ASN:C	2.24	0.41
1:B:2:SER:HB2	1:B:80:TRP:CZ2	2.55	0.41
1:D:96:PHE:HA	1:D:97:PRO:HD3	1.92	0.41
1:B:99:LYS:HG3	1:B:147:GLN:NE2	2.31	0.41
1:D:116:ASN:O	1:D:119:GLU:CG	2.62	0.41
1:D:2:SER:HB2	1:D:80:TRP:CZ2	2.55	0.41
1:B:102:LEU:HA	1:B:102:LEU:HD12	1.83	0.41
1:A:102:LEU:HD12	1:A:102:LEU:HA	1.82	0.41
1:A:109:LYS:HZ3	1:A:109:LYS:HB3	1.85	0.41
1:B:166:ASN:HA	1:B:166:ASN:HD22	1.68	0.41
1:C:166:ASN:HA	1:C:166:ASN:HD22	1.65	0.40
1:A:2:SER:HB2	1:A:80:TRP:CZ2	2.56	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	159/168 (95%)	152 (96%)	7 (4%)	0	100	100
1	B	159/168 (95%)	152 (96%)	7 (4%)	0	100	100
1	C	159/168 (95%)	152 (96%)	7 (4%)	0	100	100
1	D	159/168 (95%)	152 (96%)	7 (4%)	0	100	100
All	All	636/672 (95%)	608 (96%)	28 (4%)	0	100	100

There are no Ramachandran outliers to report.

### 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	145/153 (95%)	131 (90%)	14 (10%)	10	12
1	B	145/153 (95%)	130 (90%)	15 (10%)	9	10
1	C	145/153 (95%)	131 (90%)	14 (10%)	10	12
1	D	145/153 (95%)	131 (90%)	14 (10%)	10	12
All	All	580/612 (95%)	523 (90%)	57 (10%)	10	11

All (57) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	ASN
1	A	16	GLU
1	A	26	ASP
1	A	50	LEU
1	A	57	THR
1	A	77	ARG
1	A	98	ASN
1	A	102	LEU
1	A	104	ASN
1	A	109	LYS
1	A	112	ILE
1	A	165	THR
1	A	166	ASN
1	A	168	HIS
1	B	10	ASN
1	B	16	GLU
1	B	26	ASP
1	B	50	LEU
1	B	57	THR
1	B	77	ARG
1	B	98	ASN
1	B	102	LEU

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Mol	Chain	Res	Type
1	B	104	ASN
1	B	109	LYS
1	B	112	ILE
1	B	119	GLU
1	B	165	THR
1	B	166	ASN
1	B	168	HIS
1	C	10	ASN
1	C	16	GLU
1	C	26	ASP
1	C	50	LEU
1	C	57	THR
1	C	77	ARG
1	C	98	ASN
1	C	102	LEU
1	C	104	ASN
1	C	109	LYS
1	C	112	ILE
1	C	165	THR
1	C	166	ASN
1	C	168	HIS
1	D	10	ASN
1	D	16	GLU
1	D	26	ASP
1	D	50	LEU
1	D	57	THR
1	D	77	ARG
1	D	98	ASN
1	D	102	LEU
1	D	104	ASN
1	D	109	LYS
1	D	112	ILE
1	D	165	THR
1	D	166	ASN
1	D	168	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (44) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	10	ASN
1	A	15	ASN
1	A	38	GLN

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Mol	Chain	Res	Type
1	A	40	GLN
1	A	47	GLN
1	A	87	ASN
1	A	94	HIS
1	A	98	ASN
1	A	116	ASN
1	A	147	GLN
1	A	166	ASN
1	B	10	ASN
1	B	15	ASN
1	B	38	GLN
1	B	40	GLN
1	B	47	GLN
1	B	87	ASN
1	B	94	HIS
1	B	98	ASN
1	B	116	ASN
1	B	147	GLN
1	B	166	ASN
1	C	10	ASN
1	C	15	ASN
1	C	38	GLN
1	C	40	GLN
1	C	47	GLN
1	C	87	ASN
1	C	94	HIS
1	C	98	ASN
1	C	116	ASN
1	C	147	GLN
1	C	166	ASN
1	D	10	ASN
1	D	15	ASN
1	D	38	GLN
1	D	40	GLN
1	D	47	GLN
1	D	87	ASN
1	D	94	HIS
1	D	98	ASN
1	D	116	ASN
1	D	147	GLN
1	D	166	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 [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	163/168 (97%)	-0.05	3 (1%) 71 78	20, 37, 69, 96	0
1	B	163/168 (97%)	0.00	3 (1%) 71 78	20, 38, 70, 94	0
1	C	163/168 (97%)	0.01	3 (1%) 71 78	21, 37, 67, 98	0
1	D	163/168 (97%)	-0.04	1 (0%) 90 93	20, 38, 68, 94	0
All	All	652/672 (97%)	-0.02	10 (1%) 76 81	20, 38, 72, 98	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	2	SER	10.1
1	C	2	SER	9.4
1	D	2	SER	8.9
1	B	2	SER	8.7
1	B	126	ILE	3.1
1	B	168	HIS	3.0
1	A	126	ILE	2.6
1	C	168	HIS	2.4
1	A	168	HIS	2.3
1	C	124	ILE	2.1

### 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.