



wwPDB X-ray Structure Validation Summary Report i

Sep 9, 2023 – 07:55 PM EDT

PDB ID : 4I3H
Title : A three-gate structure of topoisomerase IV from Streptococcus pneumoniae
Authors : Laponogov, I.; Veselkov, D.A.; Pan, X.-S.; Crevel, I.; Fisher, L.M.; Sanderson, M.R.
Deposited on : 2012-11-26
Resolution : 3.70 Å (reported)

This is a wwPDB X-ray Structure Validation Summary 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 : 2.35.1
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) : 2.35.1

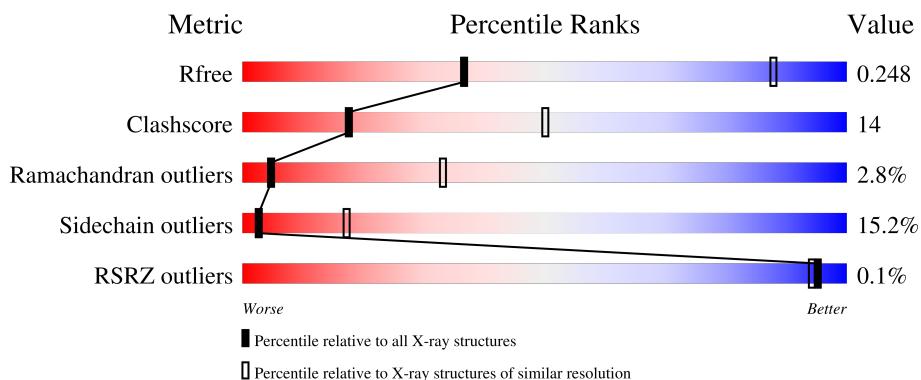
1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.70 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}	130704	1049 (3.88-3.52)
Clashscore	141614	1027 (3.86-3.54)
Ramachandran outliers	138981	1069 (3.88-3.52)
Sidechain outliers	138945	1065 (3.88-3.52)
RSRZ outliers	127900	1578 (3.90-3.50)

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 $\leq 5\%$. 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.



Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain		
3	B	1144		62%	24% • 10%

2 Entry composition i

There are 5 unique types of molecules in this entry. The entry contains 16228 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 DNA chain called DNA (5'-D(*CP*AP*AP*AP*GP*GP*CP*GP*GP*TP*AP*AP*TP*AP*CP*GP*GP*TP*TP*AP*TP*CP*CP*AP*CP*AP*GP*AP*AP*TP*CP*AP*GP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	E	14	Total	C	N	O	P	0	0	0
			282	133	52	83	14			

1	G	20	Total	C	N	O	P	0	0	0
			407	191	77	119	20			

- Molecule 2 is a DNA chain called DNA (5'-D(*CP*CP*TP*GP*AP*TP*TP*CP*TP*GP*TP*GP*AP*TP*AP*AP*CP*CP*GP*TP*AP*TP*TP*AP*CP*CP*GP*CP*CP*TP*TP*TP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	F	14	Total	C	N	O	P	0	0	0
			283	133	51	85	14			

2	H	20	Total	C	N	O	P	0	0	0
			402	189	72	121	20			

- Molecule 3 is a protein called Topoisomerase IV subunit B, DNA topoisomerase 4 subunit A chimera.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	A	1037	Total	C	N	O	S	0	0	0
			7477	4740	1286	1430	21			

3	B	1035	Total	C	N	O	S	0	0	0
			7372	4657	1280	1413	22			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1000	HIS	-	linker	UNP Q3HZ71
B	1000	HIS	-	linker	UNP Q3HZ71

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total Mg 1 1	0	0
4	B	1	Total Mg 1 1	0	0
4	H	1	Total Mg 1 1	0	0

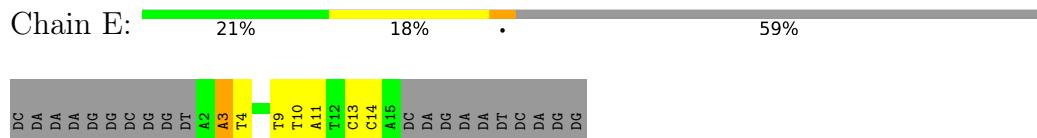
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	E	2	Total O 2 2	0	0

3 Residue-property plots

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.

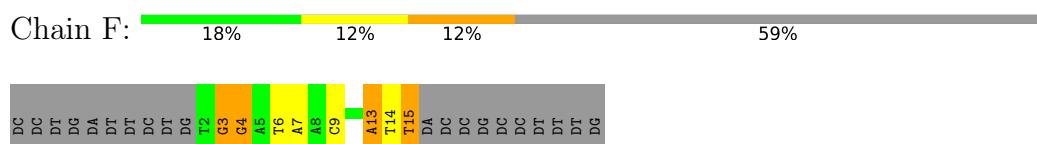
- Molecule 1: DNA (5'-D(*CP*AP*AP*AP*GP*GP*CP*GP*GP*TP*AP*AP*TP*AP*CP*GP*GP*TP*TP*AP*TP*CP*CP*AP*CP*AP*GP*AP*TP*CP*AP*GP*G)-3')



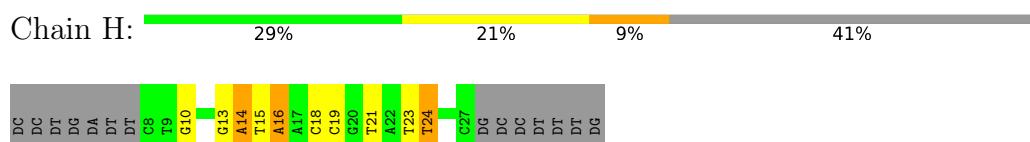
- Molecule 1: DNA (5'-D(*CP*AP*AP*AP*GP*GP*CP*GP*GP*TP*AP*AP*TP*AP*CP*GP*GP*TP*TP*AP*TP*CP*CP*AP*CP*AP*GP*AP*TP*CP*AP*GP*G)-3')



- Molecule 2: DNA (5'-D(*CP*CP*TP*GP*AP*TP*TP*CP*TP*GP*TP*GP*GP*AP*TP*AP*AP*CP*CP*GP*TP*AP*TP*TP*AP*CP*CP*GP*CP*CP*TP*TP*G)-3')

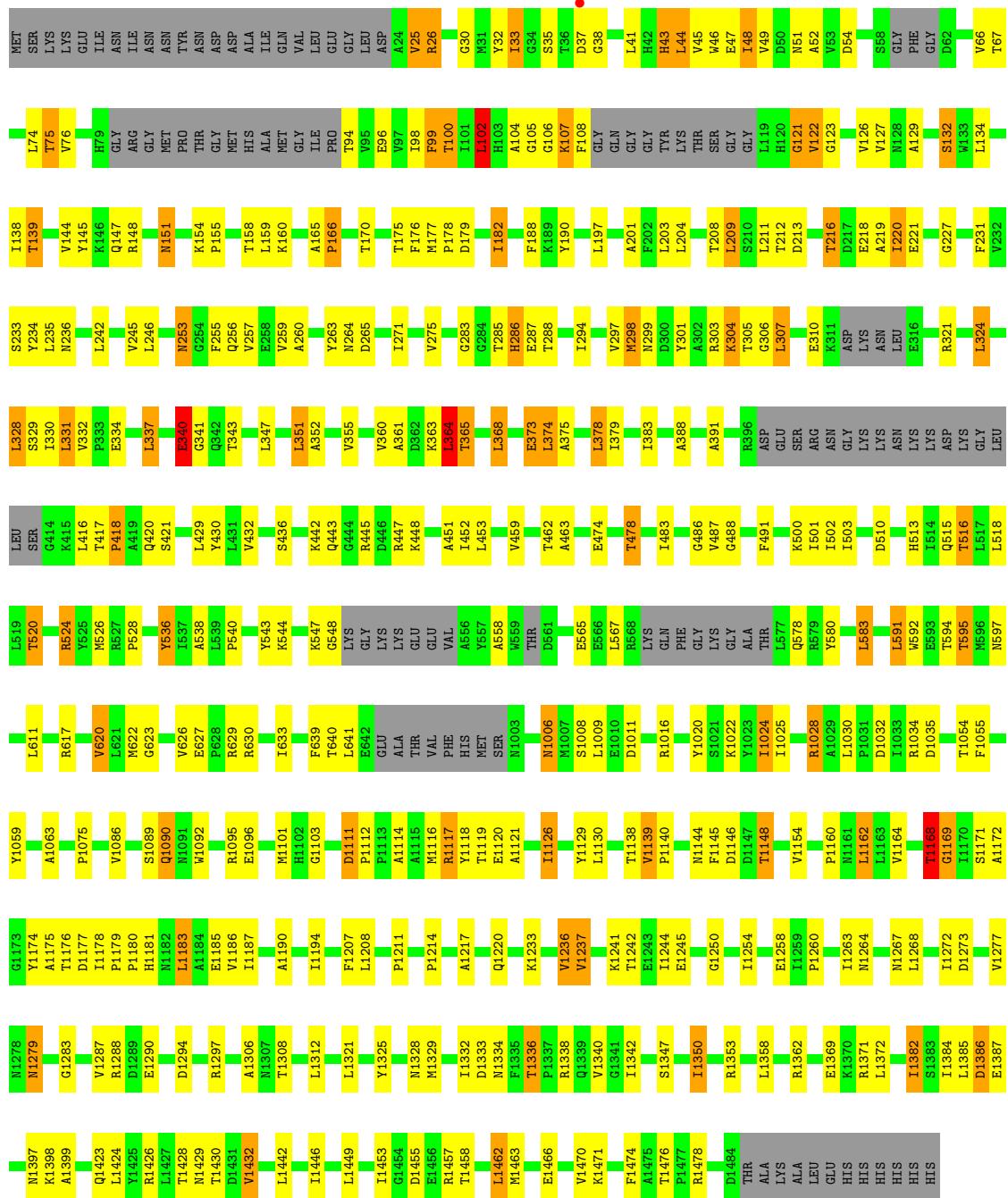


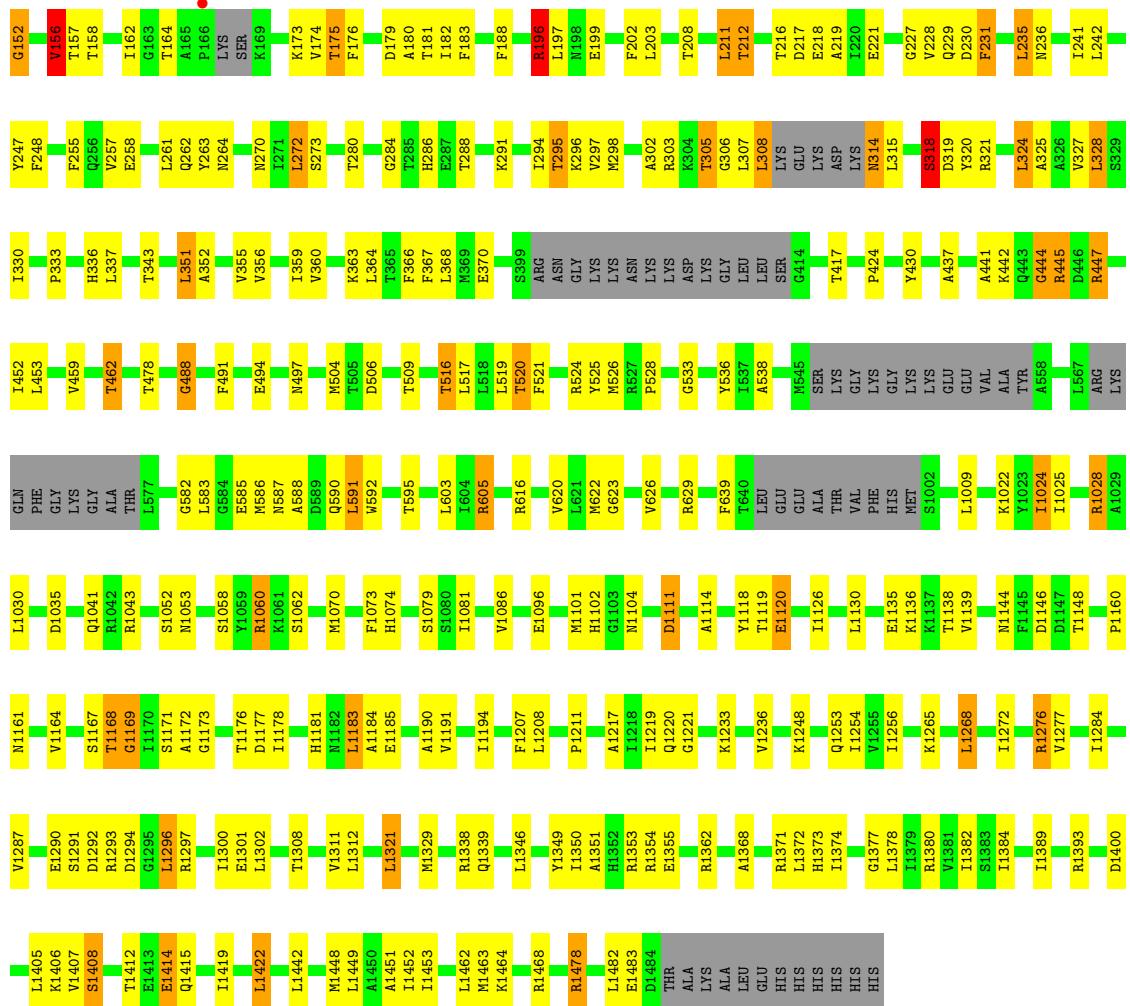
- Molecule 2: DNA (5'-D(*CP*CP*TP*GP*AP*TP*TP*CP*TP*GP*TP*GP*GP*AP*TP*AP*AP*CP*CP*GP*TP*AP*TP*TP*AP*CP*CP*GP*CP*CP*TP*TP*G)-3')



- Molecule 3: Topoisomerase IV subunit B, DNA topoisomerase 4 subunit A chimera







4 Data and refinement statistics (i)

Property	Value			Source
Space group	P 42 21 2			Depositor
Cell constants a, b, c, α , β , γ	160.60 Å 90.00°	160.60 Å 90.00°	280.56 Å 90.00°	Depositor
Resolution (Å)	71.82	–	3.70	Depositor
	71.82	–	3.70	EDS
% Data completeness (in resolution range)	99.7 (71.82-3.70) 99.7 (71.82-3.70)			Depositor EDS
R_{merge}	0.12			Depositor
R_{sym}	0.12			Depositor
$\langle I/\sigma(I) \rangle^1$	4.10 (at 3.67 Å)			Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.2_869)			Depositor
R , R_{free}	0.185 0.183	,	0.248 0.248	Depositor DCC
R_{free} test set	2006 reflections (5.04%)			wwPDB-VP
Wilson B-factor (Å ²)	95.5			Xtriage
Anisotropy	0.205			Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 82.0			EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$			Xtriage
Estimated twinning fraction	No twinning to report.			Xtriage
F_o, F_c correlation	0.91			EDS
Total number of atoms	16228			wwPDB-VP
Average B, all atoms (Å ²)	91.0			wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.52% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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

Bond lengths and bond angles in the following residue types are not validated in this section: MG

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	E	0.83	1/316 (0.3%)	1.85	12/483 (2.5%)
1	G	0.96	0/457	1.94	12/701 (1.7%)
2	F	0.80	0/317	1.91	10/485 (2.1%)
2	H	0.86	0/450	1.83	13/689 (1.9%)
3	A	0.45	0/7600	0.67	2/10359 (0.0%)
3	B	0.43	0/7491	0.65	2/10214 (0.0%)
All	All	0.50	1/16631 (0.0%)	0.87	51/22931 (0.2%)

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
3	A	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	4	DT	C1'-N1	5.18	1.55	1.49

The worst 5 of 51 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	20	DA	O4'-C1'-N9	15.76	119.03	108.00
2	H	16	DA	O4'-C4'-C3'	-10.97	99.42	106.00
1	G	24	DA	O4'-C1'-N9	-8.97	101.72	108.00
2	H	21	DT	O4'-C4'-C3'	-8.15	101.11	106.00
2	H	24	DT	O4'-C1'-N1	-8.15	102.30	108.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	A	102	LEU	Peptide

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	E	282	0	147	8	0
1	G	407	0	211	10	0
2	F	283	0	145	11	0
2	H	402	0	209	8	0
3	A	7477	0	6788	217	0
3	B	7372	0	6577	186	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	H	1	0	0	0	0
5	E	2	0	0	0	0
All	All	16228	0	14077	414	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.

The worst 5 of 414 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:E:3:DA:N6	2:F:13:DA:N6	2.20	0.89
1:E:3:DA:N1	2:F:13:DA:N1	2.21	0.89
3:A:520:THR:HG21	3:A:622:MET:HG3	1.56	0.88
3:B:48:ILE:HG12	3:B:127:VAL:HG21	1.60	0.82
1:E:3:DA:N6	2:F:13:DA:C6	2.47	0.82

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
3	A	1017/1144 (89%)	891 (88%)	101 (10%)	25 (2%)	5 35
3	B	1015/1144 (89%)	868 (86%)	115 (11%)	32 (3%)	4 31
All	All	2032/2288 (89%)	1759 (87%)	216 (11%)	57 (3%)	5 33

5 of 57 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	305	THR
3	A	373	GLU
3	A	1398	LYS
3	B	35	SER
3	B	99	PHE

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
3	A	672/969 (69%)	565 (84%)	107 (16%)	2 16
3	B	640/969 (66%)	548 (86%)	92 (14%)	3 19
All	All	1312/1938 (68%)	1113 (85%)	199 (15%)	3 17

5 of 199 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	B	100	THR

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
3	B	314	ASN
3	B	151	ASN
3	B	212	THR
3	B	445	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
3	A	51	ASN
3	A	264	ASN
3	B	264	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 [\(i\)](#)

Of 3 ligands modelled in this entry, 3 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	E	14/34 (41%)	-0.60	0 [100] [100]	144, 201, 233, 259	14 (100%)
1	G	20/34 (58%)	-0.50	0 [100] [100]	66, 82, 159, 192	0
2	F	14/34 (41%)	-0.62	0 [100] [100]	138, 196, 225, 254	14 (100%)
2	H	20/34 (58%)	-0.56	0 [100] [100]	60, 85, 168, 199	0
3	A	1037/1144 (90%)	-0.51	1 (0%) [95] [94]	35, 83, 129, 225	0
3	B	1035/1144 (90%)	-0.49	1 (0%) [95] [94]	44, 89, 136, 212	0
All	All	2140/2424 (88%)	-0.50	2 (0%) [95] [94]	35, 86, 142, 259	28 (1%)

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	B	166	PRO	3.3
3	A	37	ASP	2.9

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 [\(i\)](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	MG	B	1700	1/1	0.85	0.07	109,109,109,109	0
4	MG	H	1901	1/1	0.90	0.15	72,72,72,72	0
4	MG	A	1700	1/1	0.92	0.23	94,94,94,94	0

6.5 Other polymers [\(i\)](#)

There are no such residues in this entry.