



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 24, 2022 – 12:16 pm GMT

PDB ID : 5OWO  
Title : Human cytoplasmic Dynein N-Terminus dimerization domain at 1.8 Angstrom resolution  
Authors : Urnavicius, L.; Lau, C.K.; Elshenawy, M.M.; Morales-Rios, E.; Motz, C.; Yildiz, A.; Carter, A.P.  
Deposited on : 2017-09-01  
Resolution : 1.79 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

<https://www.wwpdb.org/validation/2017/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 : 1.8.4, CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.27  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0267  
CCP4 : 7.1.010 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.27

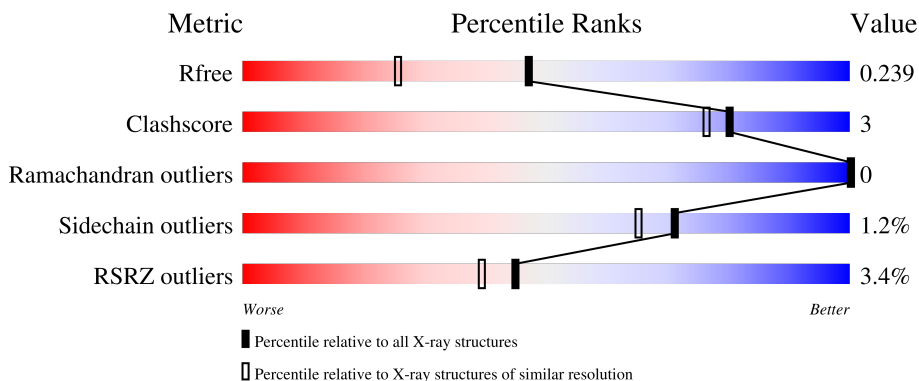
# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.79 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	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (1.80-1.80)

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  $\geq 3$ , 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.

Mol	Chain	Length	Quality of chain
1	A	201	<div style="display: flex; align-items: center;"> <div style="width: 2%; height: 10px; background-color: red; margin-right: 2px;"></div> <div style="width: 76%; height: 10px; background-color: green; margin-right: 2px;"></div> <div style="width: 5%; height: 10px; background-color: yellow; margin-right: 2px;"></div> <div style="width: 19%; height: 10px; background-color: grey;"></div> </div> <p style="font-size: small; margin-top: 5px;">2% 76% 5% 19%</p>
1	B	201	<div style="display: flex; align-items: center;"> <div style="width: 3%; height: 10px; background-color: red; margin-right: 2px;"></div> <div style="width: 77%; height: 10px; background-color: green; margin-right: 2px;"></div> <div style="width: 1%; height: 10px; background-color: yellow; margin-right: 2px;"></div> <div style="width: 19%; height: 10px; background-color: grey;"></div> </div> <p style="font-size: small; margin-top: 5px;">3% 77% 1% 19%</p>
1	C	201	<div style="display: flex; align-items: center;"> <div style="width: 2%; height: 10px; background-color: red; margin-right: 2px;"></div> <div style="width: 74%; height: 10px; background-color: green; margin-right: 2px;"></div> <div style="width: 5%; height: 10px; background-color: yellow; margin-right: 2px;"></div> <div style="width: 20%; height: 10px; background-color: grey;"></div> </div> <p style="font-size: small; margin-top: 5px;">2% 74% 5% 20%</p>
1	D	201	<div style="display: flex; align-items: center;"> <div style="width: 3%; height: 10px; background-color: red; margin-right: 2px;"></div> <div style="width: 68%; height: 10px; background-color: green; margin-right: 2px;"></div> <div style="width: 10%; height: 10px; background-color: yellow; margin-right: 2px;"></div> <div style="width: 22%; height: 10px; background-color: grey;"></div> </div> <p style="font-size: small; margin-top: 5px;">3% 68% 10% 22%</p>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	GOL	A	301	-	X	-	-
2	GOL	B	303	-	-	-	X
6	K	D	306	-	-	-	X

## 2 Entry composition [i](#)

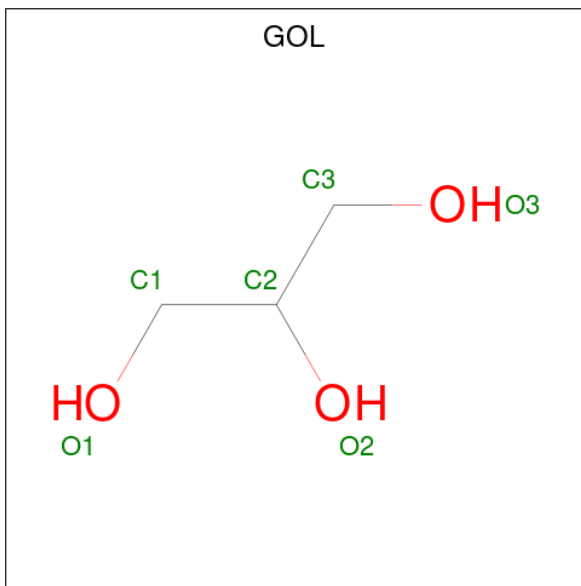
There are 7 unique types of molecules in this entry. The entry contains 5380 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Cytoplasmic dynein 1 heavy chain 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	Se			
1	A	162	Total 1283	C 819	N 216	O 245	Se 3	0	0	0
1	B	158	Total 1247	C 797	N 209	O 238	Se 3	0	0	0
1	C	160	Total 1261	C 806	N 213	O 239	Se 3	0	0	0
1	D	157	Total 1231	C 788	N 207	O 233	Se 3	0	0	0

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
2	A	1	Total 6	C 3	O 3	0	0
2	A	1	Total 6	C 3	O 3	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 6 3 3	0	0
2	B	1	Total C O 6 3 3	0	0
2	B	1	Total C O 6 3 3	0	0
2	B	1	Total C O 6 3 3	0	0
2	C	1	Total C O 6 3 3	0	0
2	C	1	Total C O 6 3 3	0	0
2	D	1	Total C O 6 3 3	0	0

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	8	Total Mg 8 8	0	0
3	B	8	Total Mg 8 8	0	0
3	C	7	Total Mg 7 7	0	0
3	D	3	Total Mg 3 3	0	0

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total Ca 1 1	0	0
4	D	1	Total Ca 1 1	0	0

- Molecule 5 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	B	1	Total Na 1 1	0	0

- Molecule 6 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	D	1	Total K 1 1	0	0

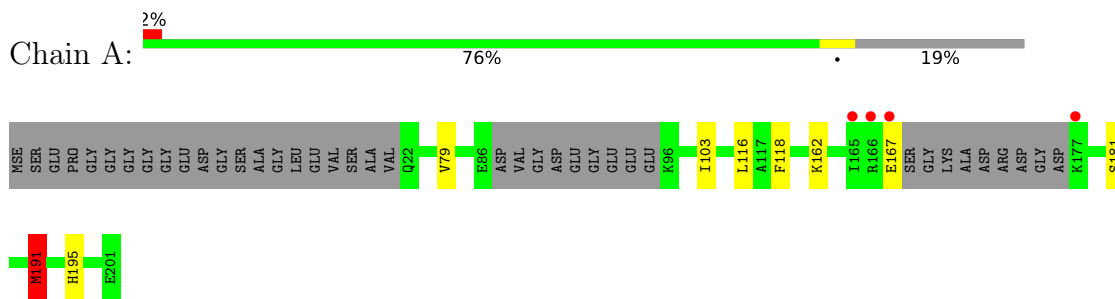
- Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	74	Total O 74 74	0	0
7	B	59	Total O 59 59	0	0
7	C	72	Total O 72 72	0	0
7	D	69	Total O 69 69	0	0

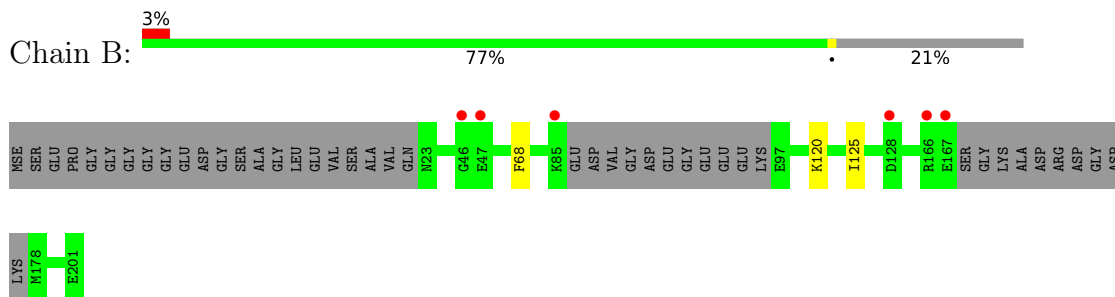
### 3 Residue-property plots [i](#)

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.

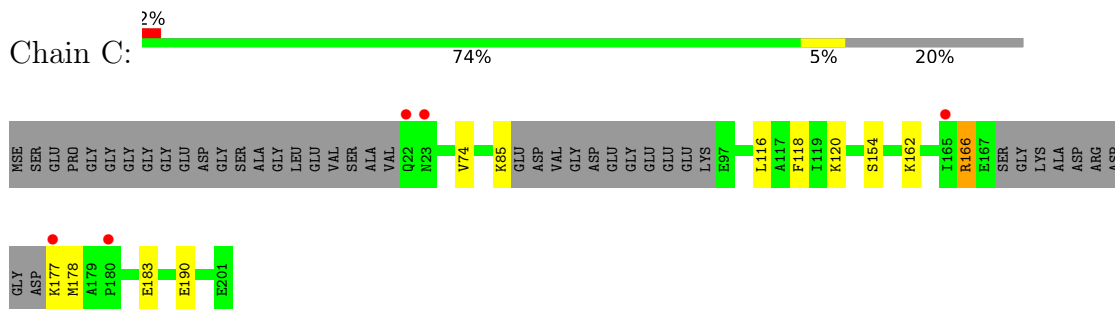
- Molecule 1: Cytoplasmic dynein 1 heavy chain 1



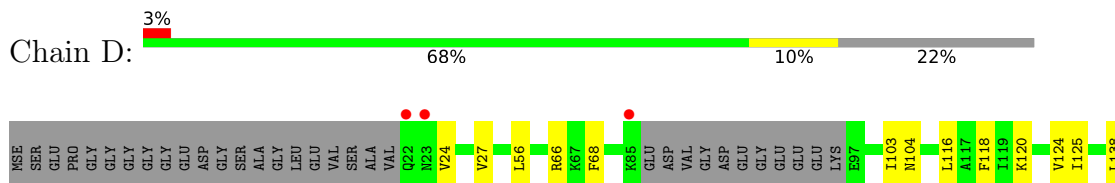
- Molecule 1: Cytoplasmic dynein 1 heavy chain 1

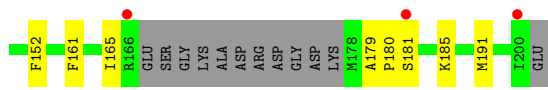


- Molecule 1: Cytoplasmic dynein 1 heavy chain 1



- Molecule 1: Cytoplasmic dynein 1 heavy chain 1







## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	101.24Å 175.56Å 50.44Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	87.78 – 1.79 48.64 – 1.79	Depositor EDS
% Data completeness (in resolution range)	99.7 (87.78-1.79) 99.7 (48.64-1.79)	Depositor EDS
$R_{merge}$	0.20	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.66 (at 1.79Å)	Xtrriage
Refinement program	REFMAC 5.8.0158	Depositor
R, $R_{free}$	0.196 , 0.229 0.210 , 0.239	Depositor DCC
$R_{free}$ test set	4245 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	26.7	Xtrriage
Anisotropy	0.066	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	(Not available) , (Not available)	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	5380	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	31.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 45.46 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.3176e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NA, MG, K, GOL, CA

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	1.00	0/1300	0.94	1/1749 (0.1%)
1	B	0.92	0/1265	0.90	0/1706
1	C	0.96	0/1278	0.95	0/1721
1	D	0.97	0/1249	0.98	0/1686
All	All	0.96	0/5092	0.94	1/6862 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	191	MSE	CG-SE-CE	-6.28	85.09	98.90

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	1283	0	1313	4	0
1	B	1247	0	1272	2	0
1	C	1261	0	1290	9	0
1	D	1231	0	1254	13	0
2	A	18	0	23	0	0
2	B	18	0	24	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	12	0	16	0	0
2	D	6	0	8	0	0
3	A	8	0	0	0	0
3	B	8	0	0	0	0
3	C	7	0	0	0	0
3	D	3	0	0	0	0
4	A	1	0	0	0	0
4	D	1	0	0	0	0
5	B	1	0	0	0	0
6	D	1	0	0	0	0
7	A	74	0	0	1	0
7	B	59	0	0	0	0
7	C	72	0	0	2	0
7	D	69	0	0	1	0
All	All	5380	0	5200	28	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 3.

All (28) 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:C:166:ARG:HH11	1:C:166:ARG:HG3	1.37	0.87
1:D:104:ASN:HB3	7:D:423:HOH:O	1.82	0.78
1:A:79:VAL:HG22	1:A:103:ILE:HD12	1.71	0.72
1:C:166:ARG:HH11	1:C:166:ARG:CG	2.05	0.69
1:D:138:LEU:HD11	1:D:152:PHE:CZ	2.32	0.64
1:A:162:LYS:HG3	7:A:456:HOH:O	2.04	0.57
1:C:116:LEU:HG	1:C:118:PHE:CE1	2.39	0.57
1:D:138:LEU:HD11	1:D:152:PHE:CE2	2.40	0.56
1:D:116:LEU:HG	1:D:118:PHE:CE1	2.44	0.53
1:D:161:PHE:CE2	1:D:165:ILE:HD13	2.45	0.52
1:D:27:VAL:CG2	1:D:66:ARG:HG2	2.40	0.52
1:D:120:LYS:HG2	1:D:125:ILE:HD11	1.91	0.52
1:D:138:LEU:CD1	1:D:152:PHE:CZ	2.92	0.52
1:C:162:LYS:HE2	1:C:183:GLU:OE2	2.11	0.51
1:B:120:LYS:HG2	1:B:125:ILE:HD11	1.92	0.50
1:D:68:PHE:CE1	1:D:120:LYS:HG3	2.47	0.50
1:C:154:SER:HB2	1:C:190:GLU:CD	2.32	0.49
1:C:74:VAL:O	1:C:120:LYS:HE2	2.13	0.48
1:C:177:LYS:NZ	7:C:402:HOH:O	2.47	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:162:LYS:HG3	7:C:450:HOH:O	2.13	0.47
1:D:181:SER:O	1:D:185:LYS:HG2	2.14	0.47
1:D:24:VAL:HG22	1:D:124:VAL:CG1	2.45	0.46
1:A:116:LEU:HG	1:A:118:PHE:CE1	2.50	0.46
1:A:191:MSE:HE2	1:A:195:HIS:CE1	2.50	0.46
1:D:56:LEU:HD23	1:D:103:ILE:CD1	2.46	0.45
1:D:179:ALA:HB3	1:D:180:PRO:HD3	1.98	0.44
1:B:68:PHE:CE1	1:B:120:LYS:HG3	2.56	0.41
1:C:166:ARG:CG	1:C:166:ARG:NH1	2.73	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	156/201 (78%)	155 (99%)	1 (1%)	0	100	100
1	B	152/201 (76%)	151 (99%)	1 (1%)	0	100	100
1	C	154/201 (77%)	152 (99%)	2 (1%)	0	100	100
1	D	151/201 (75%)	150 (99%)	1 (1%)	0	100	100
All	All	613/804 (76%)	608 (99%)	5 (1%)	0	100	100

There are no Ramachandran outliers to report.

#### 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	145/166 (87%)	142 (98%)	3 (2%)	53	42
1	B	141/166 (85%)	141 (100%)	0	100	100
1	C	142/166 (86%)	139 (98%)	3 (2%)	53	42
1	D	138/166 (83%)	137 (99%)	1 (1%)	84	81
All	All	566/664 (85%)	559 (99%)	7 (1%)	71	65

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	167	GLU
1	A	181	SER
1	A	191	MSE
1	C	85	LYS
1	C	166	ARG
1	C	178	MSE
1	D	191	MSE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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 39 ligands modelled in this entry, 30 are monoatomic - leaving 9 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The

Link column lists molecule types, if any, to which the group is linked. 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  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	GOL	C	301	-	5,5,5	1.27	1 (20%)	5,5,5	2.11	2 (40%)
2	GOL	A	303	3	5,5,5	0.32	0	5,5,5	0.28	0
2	GOL	A	302	-	5,5,5	1.11	0	5,5,5	1.23	1 (20%)
2	GOL	B	302	3	5,5,5	0.60	0	5,5,5	1.13	0
2	GOL	B	303	-	5,5,5	0.50	0	5,5,5	0.35	0
2	GOL	C	302	-	5,5,5	1.40	0	5,5,5	1.51	1 (20%)
2	GOL	A	301	-	5,5,5	0.74	0	5,5,5	2.55	2 (40%)
2	GOL	D	301	-	5,5,5	0.71	0	5,5,5	1.15	1 (20%)
2	GOL	B	301	-	5,5,5	0.82	0	5,5,5	0.79	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '2' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GOL	C	301	-	-	0/4/4/4	-
2	GOL	A	303	3	-	0/4/4/4	-
2	GOL	A	302	-	-	4/4/4/4	-
2	GOL	B	302	3	-	1/4/4/4	-
2	GOL	B	303	-	-	2/4/4/4	-
2	GOL	C	302	-	-	4/4/4/4	-
2	GOL	A	301	-	-	4/4/4/4	-
2	GOL	D	301	-	-	1/4/4/4	-
2	GOL	B	301	-	-	4/4/4/4	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	301	GOL	O2-C2	2.37	1.50	1.43

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	301	GOL	O1-C1-C2	-4.43	88.95	110.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	301	GOL	O1-C1-C2	-3.72	92.37	110.20
2	A	301	GOL	O2-C2-C3	2.93	122.04	109.12
2	C	302	GOL	O3-C3-C2	2.70	123.13	110.20
2	C	301	GOL	O2-C2-C3	2.20	118.81	109.12
2	D	301	GOL	C3-C2-C1	-2.15	103.36	111.70
2	A	302	GOL	O3-C3-C2	2.14	120.47	110.20

There are no chirality outliers.

All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	301	GOL	C1-C2-C3-O3
2	A	301	GOL	O2-C2-C3-O3
2	A	302	GOL	O1-C1-C2-C3
2	A	302	GOL	C1-C2-C3-O3
2	A	302	GOL	O2-C2-C3-O3
2	B	301	GOL	O1-C1-C2-C3
2	B	301	GOL	C1-C2-C3-O3
2	C	302	GOL	O1-C1-C2-C3
2	C	302	GOL	C1-C2-C3-O3
2	B	301	GOL	O1-C1-C2-O2
2	B	302	GOL	C1-C2-C3-O3
2	B	303	GOL	C1-C2-C3-O3
2	A	302	GOL	O1-C1-C2-O2
2	B	301	GOL	O2-C2-C3-O3
2	C	302	GOL	O1-C1-C2-O2
2	C	302	GOL	O2-C2-C3-O3
2	A	301	GOL	O1-C1-C2-O2
2	B	303	GOL	O2-C2-C3-O3
2	D	301	GOL	O2-C2-C3-O3
2	A	301	GOL	O1-C1-C2-C3

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

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	159/201 (79%)	0.44	4 (2%) 57 52	18, 26, 53, 79	0
1	B	155/201 (77%)	0.42	6 (3%) 39 33	16, 29, 49, 85	0
1	C	157/201 (78%)	0.39	5 (3%) 47 41	17, 25, 50, 90	0
1	D	154/201 (76%)	0.49	6 (3%) 39 33	17, 27, 48, 88	0
All	All	625/804 (77%)	0.43	21 (3%) 45 39	16, 27, 51, 90	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	166	ARG	5.6
1	C	177	LYS	4.7
1	A	166	ARG	4.2
1	C	22	GLN	4.2
1	B	166	ARG	4.1
1	D	22	GLN	3.6
1	B	128	ASP	3.4
1	A	177	LYS	3.3
1	A	167	GLU	3.2
1	D	200	ILE	3.2
1	D	23	ASN	3.1
1	B	85	LYS	2.9
1	D	85	LYS	2.6
1	C	165	ILE	2.5
1	C	23	ASN	2.3
1	B	47	GLU	2.2
1	B	167	GLU	2.2
1	A	165	ILE	2.2
1	D	181	SER	2.1
1	B	46	GLY	2.1
1	C	180	PRO	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 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
2	GOL	B	303	6/6	0.01	0.54	92,101,107,110	0
6	K	D	306	1/1	0.30	1.69	180,180,180,180	1
2	GOL	B	302	6/6	0.57	0.24	71,77,81,83	0
3	MG	C	306	1/1	0.62	0.33	71,71,71,71	0
2	GOL	A	302	6/6	0.77	0.23	43,48,55,60	0
2	GOL	A	303	6/6	0.82	0.31	77,79,83,84	0
2	GOL	C	302	6/6	0.83	0.25	43,48,56,57	0
3	MG	B	310	1/1	0.84	0.09	56,56,56,56	0
3	MG	A	309	1/1	0.86	0.13	48,48,48,48	0
2	GOL	C	301	6/6	0.86	0.17	27,32,42,44	0
3	MG	B	304	1/1	0.87	0.23	51,51,51,51	0
3	MG	A	308	1/1	0.87	0.16	62,62,62,62	0
3	MG	B	309	1/1	0.88	0.25	57,57,57,57	0
3	MG	C	307	1/1	0.88	0.11	44,44,44,44	0
3	MG	D	303	1/1	0.88	0.09	52,52,52,52	0
2	GOL	B	301	6/6	0.88	0.17	34,36,47,47	0
3	MG	A	311	1/1	0.89	0.23	52,52,52,52	0
3	MG	C	308	1/1	0.90	0.12	46,46,46,46	0
3	MG	C	309	1/1	0.90	0.14	45,45,45,45	0
2	GOL	D	301	6/6	0.90	0.12	32,36,43,45	0
3	MG	B	311	1/1	0.90	0.25	54,54,54,54	0
3	MG	A	307	1/1	0.92	0.20	65,65,65,65	0
3	MG	C	305	1/1	0.92	0.17	47,47,47,47	0
3	MG	C	304	1/1	0.93	0.13	48,48,48,48	0
3	MG	D	302	1/1	0.93	0.14	48,48,48,48	0
3	MG	B	308	1/1	0.94	0.24	50,50,50,50	0
2	GOL	A	301	6/6	0.94	0.12	26,34,43,52	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	MG	C	303	1/1	0.94	0.14	44,44,44,44	0
3	MG	A	310	1/1	0.95	0.21	72,72,72,72	0
3	MG	A	306	1/1	0.96	0.14	48,48,48,48	0
3	MG	B	305	1/1	0.96	0.17	42,42,42,42	0
3	MG	D	304	1/1	0.97	0.12	47,47,47,47	0
3	MG	B	307	1/1	0.97	0.06	51,51,51,51	0
3	MG	A	304	1/1	0.98	0.08	49,49,49,49	0
3	MG	A	305	1/1	0.98	0.10	41,41,41,41	0
4	CA	A	312	1/1	0.99	0.12	21,21,21,21	0
4	CA	D	305	1/1	0.99	0.31	45,45,45,45	1
5	NA	B	312	1/1	0.99	0.38	28,28,28,28	1
3	MG	B	306	1/1	0.99	0.17	37,37,37,37	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.