



# Full wwPDB X-ray Structure Validation Report ⓘ

Nov 12, 2024 – 02:34 PM EST

PDB ID : 3M5W  
Title : Crystal Structure of Tryptophanyl-tRNA Synthetase from *Campylobacter jejuni*  
Authors : Kim, Y.; Zhou, M.; Hasseman, J.; Anderson, W.F.; Joachimiak, A.; Center for Structural Genomics of Infectious Diseases (CSGID)  
Deposited on : 2010-03-14  
Resolution : 2.32 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.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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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	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.39

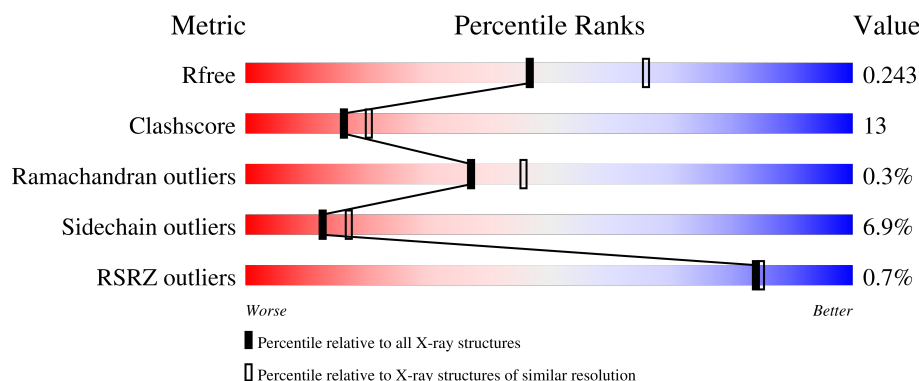
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.32 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))
$R_{free}$	164625	7250 (2.34-2.30)
Clashscore	180529	8063 (2.34-2.30)
Ramachandran outliers	177936	7993 (2.34-2.30)
Sidechain outliers	177891	7993 (2.34-2.30)
RSRZ outliers	164620	7250 (2.34-2.30)

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 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	322	
1	B	322	

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
3	GOL	A	323	-	-	X	-

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 5331 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 Tryptophanyl-tRNA synthetase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	309	Total	C	N	O	S	Se	0	3	0
			2489	1593	415	470	1	10			
1	B	319	Total	C	N	O	S	Se	0	5	0
			2584	1651	437	485	1	10			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	expression tag	UNP Q9PIB4
A	-1	ASN	-	expression tag	UNP Q9PIB4
A	0	ALA	-	expression tag	UNP Q9PIB4
B	-2	SER	-	expression tag	UNP Q9PIB4
B	-1	ASN	-	expression tag	UNP Q9PIB4
B	0	ALA	-	expression tag	UNP Q9PIB4

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total O S 5 4 1	0	0
2	A	1	Total O S 5 4 1	0	0
2	A	1	Total O S 5 4 1	0	0
2	B	1	Total O S 5 4 1	0	0
2	B	1	Total O S 5 4 1	0	0

- Molecule 3 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
3	A	1	Total C O 6 3 3	0	0
3	B	1	Total C O 6 3 3	0	0
3	B	1	Total C O 6 3 3	0	0

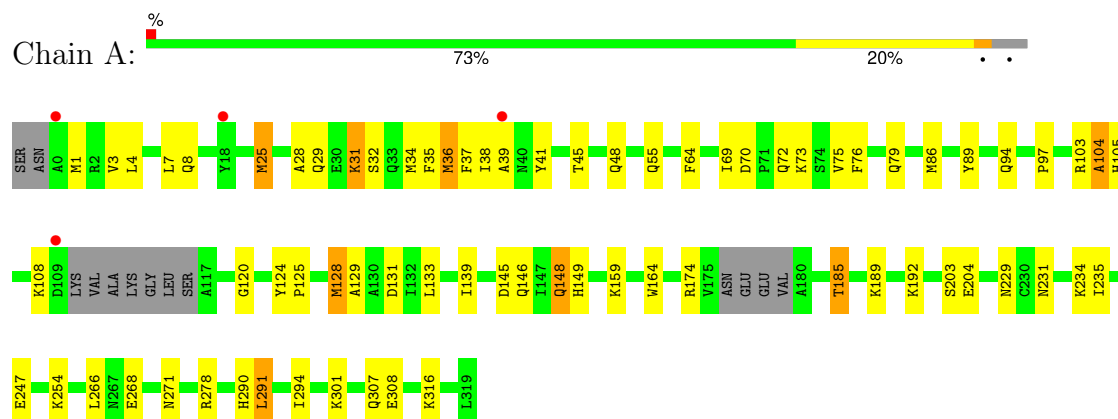
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	115	Total O 115 115	0	0
4	B	100	Total O 100 100	0	0

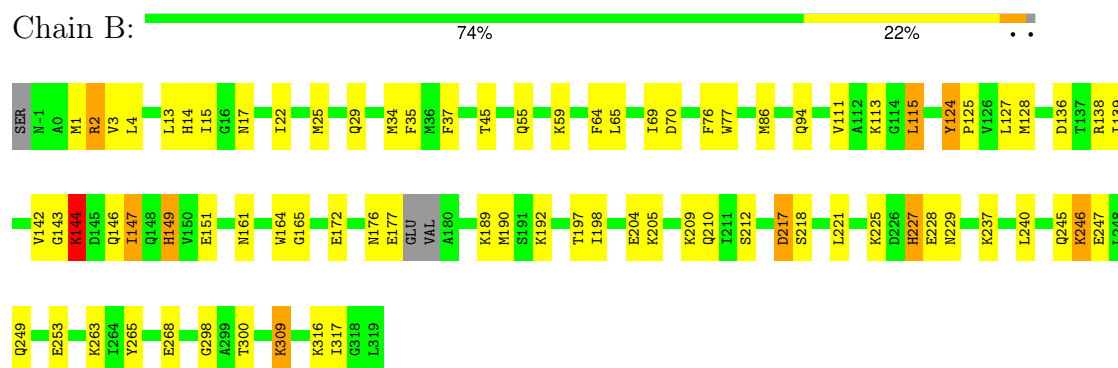
### 3 Residue-property plots [i](#)

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: Tryptophanyl-tRNA synthetase



#### • Molecule 1: Tryptophanyl-tRNA synthetase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	66.15Å 50.60Å 108.64Å 90.00° 107.61° 90.00°	Depositor
Resolution (Å)	47.81 – 2.32 47.81 – 2.32	Depositor EDS
% Data completeness (in resolution range)	96.7 (47.81-2.32) 96.6 (47.81-2.32)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.13	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.57 (at 2.32Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6 _289)	Depositor
R, $R_{free}$	0.183 , 0.243 0.180 , 0.243	Depositor DCC
$R_{free}$ test set	1997 reflections (6.88%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	40.5	Xtriage
Anisotropy	0.218	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 54.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.029 for h,-k,-h-l	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	5331	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.07% 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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, GOL

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	0.68	0/2525	0.70	0/3377
1	B	0.64	0/2622	0.72	1/3506 (0.0%)
All	All	0.66	0/5147	0.71	1/6883 (0.0%)

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

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	B	4	LEU	CA-CB-CG	5.35	127.59	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	104	ALA	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	A	2489	0	2494	74	0
1	B	2584	0	2599	61	0
2	A	15	0	0	0	0
2	B	10	0	0	0	0
3	A	6	0	8	7	0
3	B	12	0	16	1	0
4	A	115	0	0	4	0
4	B	100	0	0	1	0
All	All	5331	0	5117	132	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 13.

All (132) 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:A:25:MSE:HE2	1:A:36:MSE:CE	1.77	1.15
1:A:25:MSE:CE	1:A:36:MSE:HG2	1.80	1.11
1:B:1:MSE:HE3	1:B:139:ILE:HD11	1.19	1.11
1:B:2:ARG:HD2	1:B:35:PHE:CE1	1.85	1.09
1:A:25:MSE:HE2	1:A:36:MSE:HE3	1.14	1.06
1:B:29:GLN:HE22	1:B:70:ASP:H	1.02	0.96
1:A:25:MSE:CE	1:A:36:MSE:HE3	1.96	0.94
1:A:1:MSE:HE3	1:A:139:ILE:HD11	1.52	0.89
1:A:36:MSE:HE2	1:A:69:ILE:HD11	1.52	0.88
1:A:104:ALA:O	1:A:108:LYS:HB2	1.75	0.87
1:A:204:GLU:HB2	3:A:323:GOL:H31	1.55	0.87
1:A:25:MSE:HE1	1:A:36:MSE:HG2	1.55	0.86
1:A:36:MSE:HE2	1:A:69:ILE:CD1	2.05	0.86
1:A:204:GLU:H	3:A:323:GOL:H32	1.43	0.83
1:A:1:MSE:CE	1:A:139:ILE:HD11	2.08	0.82
1:A:204:GLU:N	3:A:323:GOL:H32	1.95	0.82
1:A:25:MSE:HE2	1:A:36:MSE:HG2	1.60	0.81
1:A:32:SER:OG	1:A:34:MSE:HE3	1.83	0.78
1:B:64:PHE:HB3	1:B:69:ILE:HD12	1.65	0.78
1:A:36:MSE:HE1	1:A:69:ILE:HG12	1.64	0.78
1:A:192:LYS:HD2	4:A:373:HOH:O	1.84	0.76
1:A:86:MSE:HE3	1:A:86:MSE:HA	1.66	0.75
1:A:29:GLN:HE22	1:A:70:ASP:H	1.35	0.72
1:A:231:ASN:O	1:A:235:ILE:HD12	1.90	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1:MSE:HE3	1:A:139:ILE:CD1	2.23	0.69
1:A:203:SER:HB2	3:A:323:GOL:H11	1.74	0.68
1:B:1:MSE:CE	1:B:139:ILE:HD11	2.10	0.68
1:A:189:LYS:HE3	4:A:431:HOH:O	1.93	0.68
1:A:64:PHE:HB3	1:A:69:ILE:HD12	1.76	0.68
1:A:79:GLN:NE2	1:A:131:ASP:OD2	2.27	0.67
1:A:145:ASP:OD2	1:A:146:GLN:HG2	1.94	0.67
1:B:128:MSE:HG3	3:B:325:GOL:O1	1.96	0.66
1:A:271[A]:ASN:HA	1:A:278:ARG:HH12	1.62	0.65
1:B:1:MSE:HE3	1:B:139:ILE:CD1	2.12	0.65
1:B:29:GLN:HE22	1:B:70:ASP:N	1.85	0.65
1:A:25:MSE:HE2	1:A:36:MSE:CG	2.27	0.65
1:B:86:MSE:HA	1:B:86:MSE:HE2	1.79	0.64
1:A:28:ALA:HA	1:A:31:LYS:HE3	1.79	0.63
1:A:125:PRO:HA	1:A:128:MSE:CE	2.29	0.63
1:B:2:ARG:HD2	1:B:35:PHE:CD1	2.32	0.63
1:A:125:PRO:HA	1:A:128:MSE:HE2	1.81	0.62
1:A:146:GLN:OE1	1:A:149:HIS:HD2	1.81	0.62
1:A:25:MSE:HE2	1:A:36:MSE:SE	2.50	0.61
1:A:36:MSE:CE	1:A:69:ILE:CD1	2.77	0.61
1:B:29:GLN:NE2	1:B:70:ASP:H	1.86	0.60
1:A:48:GLN:HB2	4:A:433:HOH:O	2.02	0.60
1:B:147:ILE:HD11	1:B:151:GLU:OE2	2.03	0.59
1:A:185:THR:HG21	4:A:360:HOH:O	2.03	0.59
1:B:198:ILE:HA	1:B:210:GLN:OE1	2.02	0.58
1:A:37:PHE:CE1	1:A:39:ALA:HB2	2.38	0.58
1:A:37:PHE:HE1	1:A:39:ALA:HB2	1.69	0.58
1:A:38:ILE:HD12	1:A:38:ILE:H	1.69	0.58
1:B:164:TRP:O	1:B:316:LYS:HE3	2.04	0.57
1:A:271[B]:ASN:HA	1:A:278:ARG:HH12	1.69	0.57
1:A:204:GLU:H	3:A:323:GOL:C3	2.17	0.56
1:B:265:TYR:O	1:B:268:GLU:HG2	2.05	0.56
1:A:70:ASP:OD2	1:A:73:LYS:HG2	2.06	0.56
1:A:3:VAL:HB	1:A:34:MSE:CE	2.36	0.56
1:B:17:ASN:O	1:B:22:ILE:HG13	2.06	0.56
1:B:17:ASN:HD21	1:B:192:LYS:NZ	2.05	0.55
1:A:129:ALA:O	1:A:133:LEU:HG	2.07	0.55
1:B:249:GLN:O	1:B:253:GLU:HG2	2.07	0.54
1:B:25:MSE:HG3	1:B:34:MSE:CE	2.37	0.54
1:B:146:GLN:O	1:B:149[B]:HIS:ND1	2.30	0.54
1:B:217:ASP:OD1	1:B:217:ASP:N	2.34	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:204:GLU:HB2	3:A:323:GOL:C3	2.33	0.54
1:B:136:ASP:HA	4:B:341:HOH:O	2.08	0.54
1:A:120:GLY:HA3	1:B:94:GLN:OE1	2.09	0.53
1:A:290:HIS:O	1:A:294:ILE:HG13	2.09	0.53
1:A:36:MSE:CE	1:A:69:ILE:HD11	2.34	0.52
1:B:237:LYS:HG2	1:B:245:GLN:OE1	2.09	0.52
1:B:227:HIS:ND1	1:B:227:HIS:N	2.56	0.52
1:A:203:SER:CB	3:A:323:GOL:H11	2.39	0.51
1:A:94:GLN:OE1	1:B:45:THR:HG22	2.11	0.51
1:B:176:ASN:O	1:B:177:GLU:C	2.49	0.51
1:B:128:MSE:SE	1:B:149[A]:HIS:HD2	2.44	0.50
1:A:104:ALA:HA	1:A:148:GLN:OE1	2.11	0.50
1:A:3:VAL:HB	1:A:34:MSE:HE2	1.93	0.50
1:A:164:TRP:O	1:A:316:LYS:HE3	2.12	0.49
1:B:113:LYS:HB2	1:B:115:LEU:HD12	1.94	0.49
1:B:124:TYR:N	1:B:125:PRO:CD	2.75	0.49
1:B:13:LEU:HA	1:B:17:ASN:HD22	1.76	0.49
1:B:161:ASN:HA	1:B:165:GLY:O	2.13	0.49
1:B:237:LYS:HG2	1:B:245:GLN:HG2	1.95	0.49
1:A:229:ASN:HA	1:A:234:LYS:NZ	2.27	0.48
1:A:41:TYR:HB2	1:A:79:GLN:OE1	2.13	0.48
1:B:55[B]:GLN:HG3	1:B:59:LYS:HE3	1.95	0.48
1:A:97:PRO:HG2	1:A:159:LYS:HE2	1.95	0.48
1:B:142:VAL:HG13	1:B:143:GLY:N	2.28	0.48
1:A:1:MSE:SE	1:A:139:ILE:HD11	2.63	0.47
1:B:86:MSE:CE	1:B:127:LEU:HD13	2.44	0.47
1:B:240:LEU:O	1:B:245:GLN:NE2	2.47	0.47
1:B:142:VAL:HG22	1:B:149[B]:HIS:HE1	1.80	0.47
1:B:76:PHE:O	1:B:298:GLY:HA3	2.15	0.47
1:A:124:TYR:N	1:A:125:PRO:CD	2.78	0.47
1:B:14:HIS:H	1:B:17:ASN:HD22	1.62	0.46
1:B:128:MSE:SE	1:B:149[B]:HIS:CD2	3.18	0.46
1:A:25:MSE:SE	1:A:36:MSE:HE3	2.65	0.46
1:A:45:THR:HA	1:B:94:GLN:OE1	2.16	0.46
1:B:144:LYS:HB3	1:B:144:LYS:NZ	2.30	0.46
1:B:205:LYS:HE3	1:B:209:LYS:NZ	2.31	0.45
1:A:124:TYR:CD2	1:A:125:PRO:HD3	2.52	0.45
1:B:15:ILE:HG12	1:B:198:ILE:O	2.16	0.45
1:B:2:ARG:CD	1:B:35:PHE:CE1	2.78	0.44
1:B:147:ILE:O	1:B:151:GLU:HG3	2.18	0.44
1:B:212:SER:HA	1:B:263:LYS:HD3	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:3:VAL:HG22	1:B:139:ILE:HB	1.99	0.44
1:B:1:MSE:HG3	1:B:139:ILE:CD1	2.48	0.43
1:B:309:LYS:HB2	1:B:309:LYS:HE2	1.42	0.43
1:A:36:MSE:HE1	1:A:69:ILE:CG1	2.43	0.43
1:A:36:MSE:CE	1:A:69:ILE:HG12	2.42	0.43
1:A:86:MSE:O	1:A:89:TYR:HB3	2.18	0.43
1:A:291:LEU:HD23	1:A:291:LEU:HA	1.84	0.43
1:B:246:LYS:O	1:B:247:GLU:C	2.57	0.42
1:A:86:MSE:HA	1:A:86:MSE:CE	2.43	0.42
1:A:35:PHE:C	1:A:36:MSE:HG3	2.39	0.42
1:B:317:ILE:HG13	1:B:317:ILE:O	2.19	0.42
1:A:3:VAL:HB	1:A:34:MSE:HE1	2.02	0.42
1:B:37:PHE:HA	1:B:77:TRP:O	2.20	0.42
1:B:149[A]:HIS:N	1:B:149[A]:HIS:ND1	2.67	0.41
1:B:138:ARG:HB3	1:B:139:ILE:HD12	2.02	0.41
1:A:75:VAL:HG22	1:A:301:LYS:HE3	2.01	0.41
1:B:128:MSE:SE	1:B:149[A]:HIS:CD2	3.24	0.41
1:B:190:MSE:HG3	1:B:197:THR:HA	2.03	0.41
1:A:38:ILE:HD11	1:A:76:PHE:CZ	2.56	0.41
1:A:235:ILE:HG22	1:A:266:LEU:HD11	2.02	0.41
1:A:35:PHE:O	1:A:36:MSE:HG3	2.21	0.40
1:B:65:LEU:HD23	1:B:65:LEU:HA	1.94	0.40
1:B:70:ASP:C	1:B:70:ASP:OD1	2.60	0.40
1:A:25:MSE:CE	1:A:36:MSE:CG	2.70	0.40
1:A:254:LYS:HB3	1:A:254:LYS:HE2	1.72	0.40
1:B:221:LEU:HD23	1:B:221:LEU:C	2.42	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone

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	306/322 (95%)	295 (96%)	11 (4%)	0	100	100
1	B	320/322 (99%)	313 (98%)	5 (2%)	2 (1%)	22	26
All	All	626/644 (97%)	608 (97%)	16 (3%)	2 (0%)	37	46

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	144	LYS
1	B	228	GLU

### 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	268/266 (101%)	249 (93%)	19 (7%)	12	16
1	B	278/266 (104%)	259 (93%)	19 (7%)	13	17
All	All	546/532 (103%)	508 (93%)	38 (7%)	13	16

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	LEU
1	A	7	LEU
1	A	8	GLN
1	A	25	MSE
1	A	31	LYS
1	A	36	MSE
1	A	55	GLN
1	A	72	GLN
1	A	103	ARG
1	A	105	HIS
1	A	128	MSE
1	A	148	GLN
1	A	174	ARG
1	A	185	THR

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Mol	Chain	Res	Type
1	A	247	GLU
1	A	268	GLU
1	A	291	LEU
1	A	307	GLN
1	A	308	GLU
1	B	2	ARG
1	B	111	VAL
1	B	115	LEU
1	B	124	TYR
1	B	144	LYS
1	B	147	ILE
1	B	149[A]	HIS
1	B	149[B]	HIS
1	B	172	GLU
1	B	189	LYS
1	B	204	GLU
1	B	217	ASP
1	B	218	SER
1	B	225	LYS
1	B	227	HIS
1	B	229	ASN
1	B	246	LYS
1	B	300	THR
1	B	309	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	29	GLN
1	A	56	ASN
1	A	149	HIS
1	B	8	GLN
1	B	17	ASN
1	B	29	GLN
1	B	229	ASN

### 5.3.3 RNA ⓘ

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 oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

8 ligands are modelled in this entry.

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	SO4	B	321	-	4,4,4	0.27	0	6,6,6	0.41	0
2	SO4	A	322	-	4,4,4	0.24	0	6,6,6	0.18	0
2	SO4	B	323	-	4,4,4	0.25	0	6,6,6	0.17	0
2	SO4	A	324	-	4,4,4	0.23	0	6,6,6	0.20	0
3	GOL	B	324	-	5,5,5	0.40	0	5,5,5	0.42	0
2	SO4	A	321	-	4,4,4	0.22	0	6,6,6	0.59	0
3	GOL	A	323	-	5,5,5	0.46	0	5,5,5	0.21	0
3	GOL	B	325	-	5,5,5	0.43	0	5,5,5	0.36	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. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GOL	B	324	-	-	4/4/4/4	-
3	GOL	A	323	-	-	4/4/4/4	-
3	GOL	B	325	-	-	0/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	324	GOL	O1-C1-C2-C3
3	B	324	GOL	C1-C2-C3-O3
3	A	323	GOL	O1-C1-C2-C3
3	A	323	GOL	C1-C2-C3-O3
3	B	324	GOL	O1-C1-C2-O2
3	B	324	GOL	O2-C2-C3-O3
3	A	323	GOL	O1-C1-C2-O2
3	A	323	GOL	O2-C2-C3-O3

There are no ring outliers.

2 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	323	GOL	7	0
3	B	325	GOL	1	0

## 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	299/322 (92%)	-0.23	4 (1%) 74 75	22, 47, 80, 104	3 (1%)
1	B	309/322 (95%)	-0.21	0 100 100	20, 48, 85, 120	5 (1%)
All	All	608/644 (94%)	-0.22	4 (0%) 84 85	20, 47, 84, 120	8 (1%)

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	18	TYR	4.8
1	A	0	ALA	2.7
1	A	109	ASP	2.5
1	A	39	ALA	2.2

### 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
3	GOL	A	323	6/6	0.65	0.16	71,78,81,85	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	SO4	B	323	5/5	0.73	0.10	117,121,122,125	0
2	SO4	A	324	5/5	0.75	0.11	125,126,129,131	0
3	GOL	B	324	6/6	0.75	0.17	93,98,103,107	0
2	SO4	A	322	5/5	0.81	0.11	118,121,121,123	0
3	GOL	B	325	6/6	0.89	0.17	63,69,74,84	0
2	SO4	B	321	5/5	0.95	0.08	64,70,76,84	0
2	SO4	A	321	5/5	0.98	0.06	47,48,53,65	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.