



Full wwPDB X-ray Structure Validation Report ⓘ

May 25, 2020 – 03:52 am BST

PDB ID : 4RNY
Title : Structure of Helicobacter pylori Csd3 from the orthorhombic crystal
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Deposited on : 2014-10-27
Resolution : 2.00 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 ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
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) : 2.11

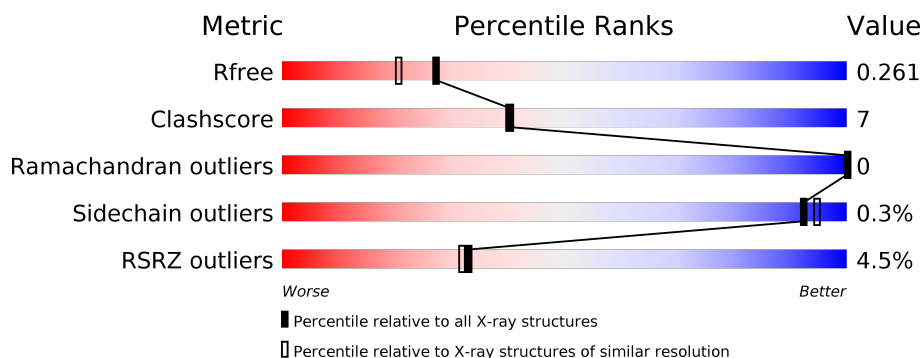
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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 ≥ 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	371	<div> <div>3%</div> <div>88%</div> <div>9%</div> <div>.</div> </div>
1	B	371	<div> <div>6%</div> <div>87%</div> <div>9%</div> <div>..</div> </div>

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	SO4	A	505	-	-	X	-
3	SO4	B	505	-	-	X	-
4	GOL	A	507	-	-	X	X
4	GOL	A	508	-	-	X	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 6124 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 Conserved hypothetical secreted protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	359	Total	C	N	O	S	0	0	0
			2911	1880	509	516	6			
1	B	359	Total	C	N	O	S	0	0	0
			2911	1880	509	516	6			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	41	MET	-	INITIATING METHIONINE	UNP O25247
A	404	LEU	-	EXPRESSION TAG	UNP O25247
A	405	GLU	-	EXPRESSION TAG	UNP O25247
A	406	HIS	-	EXPRESSION TAG	UNP O25247
A	407	HIS	-	EXPRESSION TAG	UNP O25247
A	408	HIS	-	EXPRESSION TAG	UNP O25247
A	409	HIS	-	EXPRESSION TAG	UNP O25247
A	410	HIS	-	EXPRESSION TAG	UNP O25247
A	411	HIS	-	EXPRESSION TAG	UNP O25247
B	41	MET	-	INITIATING METHIONINE	UNP O25247
B	404	LEU	-	EXPRESSION TAG	UNP O25247
B	405	GLU	-	EXPRESSION TAG	UNP O25247
B	406	HIS	-	EXPRESSION TAG	UNP O25247
B	407	HIS	-	EXPRESSION TAG	UNP O25247
B	408	HIS	-	EXPRESSION TAG	UNP O25247
B	409	HIS	-	EXPRESSION TAG	UNP O25247
B	410	HIS	-	EXPRESSION TAG	UNP O25247
B	411	HIS	-	EXPRESSION TAG	UNP O25247

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Zn	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



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

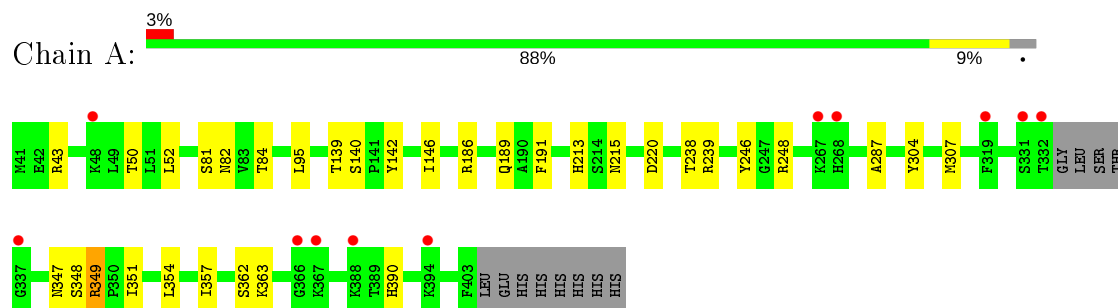
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	104	Total	O	0	0
			104	104		
5	B	121	Total	O	0	0
			121	121		

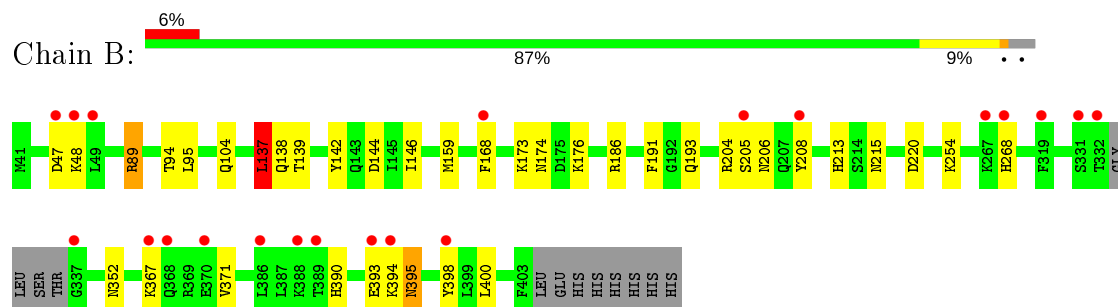
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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 ($\text{RSRZ} > 2$). 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: Conserved hypothetical secreted protein



- Molecule 1: Conserved hypothetical secreted protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	62.62Å 112.15Å 112.94Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.00 33.58 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.5 (50.00-2.00) 99.6 (33.58-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.88 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.203 , 0.256 0.210 , 0.261	Depositor DCC
R_{free} test set	2754 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	31.3	Xtriage
Anisotropy	0.017	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 41.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.039 for -h,l,k	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6124	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.36% 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

5.1 Standard geometry

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

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	0.69	0/2980	0.80	2/4013 (0.0%)
1	B	0.70	0/2980	0.81	3/4013 (0.1%)
All	All	0.70	0/5960	0.80	5/8026 (0.1%)

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	B	0	1

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	349	ARG	NE-CZ-NH2	-10.81	114.89	120.30
1	A	349	ARG	NE-CZ-NH1	9.69	125.14	120.30
1	B	89	ARG	NE-CZ-NH2	-8.91	115.84	120.30
1	B	137	LEU	CA-CB-CG	6.78	130.89	115.30
1	B	89	ARG	NE-CZ-NH1	6.34	123.47	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	393	GLU	Peptide

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	2911	0	2960	49	0
1	B	2911	0	2960	34	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	20	0	0	6	0
3	B	25	0	0	3	0
4	A	24	0	32	21	0
4	B	6	0	8	3	0
5	A	104	0	0	3	0
5	B	121	0	0	2	0
All	All	6124	0	5960	85	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 7.

All (85) 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:246:TYR:HE1	4:A:508:GOL:HO2	0.88	0.86
1:A:246:TYR:HE1	4:A:508:GOL:O2	1.58	0.85
1:A:390:HIS:CE1	3:A:505:SO4:O3	2.33	0.81
1:B:213:HIS:HD2	1:B:215:ASN:H	1.29	0.81
1:A:52:LEU:HB2	4:A:507:GOL:H31	1.68	0.74
1:A:390:HIS:NE2	3:A:505:SO4:O3	2.21	0.73
1:A:246:TYR:CE1	4:A:508:GOL:O2	2.37	0.73
1:A:52:LEU:HD12	4:A:507:GOL:H32	1.70	0.73
1:A:50:THR:HB	4:A:507:GOL:H2	1.72	0.70
1:B:89:ARG:HG3	5:B:652:HOH:O	1.91	0.70
1:A:81:SER:O	1:A:82:ASN:HB2	1.91	0.70
1:A:52:LEU:HD12	4:A:507:GOL:C3	2.23	0.68
1:A:213:HIS:HD2	1:A:215:ASN:H	1.46	0.64
1:A:50:THR:CB	4:A:507:GOL:H2	2.29	0.63
1:A:52:LEU:CB	4:A:507:GOL:H31	2.30	0.62
1:B:137:LEU:CD1	1:B:139:THR:H	2.14	0.61
1:B:142:TYR:O	1:B:146:ILE:HG12	2.02	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:186:ARG:HG3	1:B:191:PHE:CD1	2.37	0.60
1:A:287:ALA:HB3	4:A:507:GOL:H12	1.83	0.59
1:A:287:ALA:CB	4:A:507:GOL:H12	2.34	0.58
1:A:142:TYR:CZ	1:A:146:ILE:HD11	2.38	0.58
1:B:173:LYS:O	1:B:174:ASN:HB2	2.05	0.57
1:B:367:LYS:O	1:B:371:VAL:HG23	2.06	0.56
1:B:104:GLN:NE2	1:B:193:GLN:H	2.04	0.55
1:B:159:MET:HB2	1:B:168:PHE:CE2	2.41	0.55
1:B:213:HIS:CD2	1:B:215:ASN:H	2.18	0.55
1:B:204:ARG:O	1:B:205:SER:HB2	2.06	0.54
1:A:186:ARG:HG3	1:A:191:PHE:CD1	2.43	0.54
1:A:390:HIS:HE1	3:A:505:SO4:O4	1.91	0.54
1:A:349:ARG:HH11	4:A:508:GOL:H11	1.73	0.53
1:A:142:TYR:CZ	1:A:146:ILE:CD1	2.92	0.53
1:A:351:ILE:C	1:A:351:ILE:HD12	2.30	0.52
1:A:390:HIS:HE1	3:A:505:SO4:S	2.31	0.52
1:B:352:ASN:HD22	4:B:507:GOL:C1	2.22	0.52
1:A:351:ILE:O	1:A:351:ILE:HD12	2.10	0.51
1:A:362:SER:O	1:A:363:LYS:HG3	2.10	0.51
1:A:220:ASP:HB2	3:A:503:SO4:O3	2.11	0.51
1:A:287:ALA:H	4:A:507:GOL:H12	1.76	0.51
1:A:213:HIS:CD2	1:A:215:ASN:H	2.27	0.50
1:B:137:LEU:HD12	1:B:139:THR:N	2.26	0.50
1:A:390:HIS:CE1	3:A:505:SO4:S	3.05	0.50
4:A:508:GOL:C1	5:A:612:HOH:O	2.59	0.50
1:B:137:LEU:HD12	1:B:138:GLN:N	2.27	0.49
1:B:268:HIS:CD2	5:B:716:HOH:O	2.64	0.49
4:A:508:GOL:H12	5:A:612:HOH:O	2.13	0.49
1:B:206:ASN:HB2	1:B:208:TYR:OH	2.12	0.48
1:B:390:HIS:NE2	3:B:505:SO4:S	2.83	0.48
1:B:139:THR:HB	1:B:144:ASP:OD1	2.15	0.47
1:A:304:TYR:HB3	1:A:307:MET:HE2	1.96	0.47
1:A:349:ARG:HD2	4:A:508:GOL:O1	2.15	0.47
1:B:394:LYS:HE3	1:B:395:ASN:HA	1.96	0.47
1:B:137:LEU:CD1	1:B:139:THR:N	2.78	0.46
1:A:238:THR:OG1	1:A:239:ARG:N	2.49	0.46
1:A:347:ASN:O	1:A:348:SER:OG	2.29	0.46
1:B:352:ASN:HD22	4:B:507:GOL:H11	1.81	0.46
1:B:137:LEU:HD12	1:B:139:THR:H	1.79	0.45
1:A:139:THR:OG1	1:A:140:SER:N	2.50	0.45
1:A:248:ARG:HD3	5:A:659:HOH:O	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:137:LEU:HD11	1:B:139:THR:O	2.17	0.45
1:A:52:LEU:HD12	4:A:507:GOL:H31	1.98	0.45
1:B:220:ASP:HB2	3:B:506:SO4:O3	2.16	0.45
1:A:142:TYR:CE2	1:A:146:ILE:CD1	3.00	0.44
1:B:174:ASN:O	1:B:176:LYS:NZ	2.48	0.44
1:A:189:GLN:HE21	1:A:189:GLN:HB3	1.59	0.43
1:A:354:LEU:HD23	1:A:357:ILE:HD12	1.99	0.43
1:B:390:HIS:HD2	1:B:398:TYR:CD2	2.36	0.43
1:B:89:ARG:HA	1:B:94:THR:O	2.19	0.43
1:B:47:ASP:O	1:B:48:LYS:HB2	2.18	0.43
1:B:173:LYS:O	1:B:174:ASN:CB	2.67	0.43
1:A:52:LEU:CG	4:A:507:GOL:H31	2.48	0.43
1:B:352:ASN:HD22	4:B:507:GOL:H12	1.84	0.42
1:B:95:LEU:HD13	1:B:95:LEU:C	2.39	0.42
1:A:43:ARG:NE	1:A:189:GLN:OE1	2.53	0.42
4:A:509:GOL:H31	1:B:254:LYS:NZ	2.34	0.42
1:A:81:SER:O	1:A:82:ASN:CB	2.61	0.42
1:A:304:TYR:HB2	1:A:307:MET:HE3	2.00	0.42
1:A:304:TYR:HB3	1:A:307:MET:CE	2.50	0.42
1:A:43:ARG:HG3	1:A:84:THR:CG2	2.50	0.41
1:B:400:LEU:HD22	3:B:505:SO4:O4	2.20	0.41
1:A:142:TYR:CE1	1:A:146:ILE:HD11	2.55	0.41
1:A:349:ARG:CZ	4:A:508:GOL:H31	2.51	0.41
1:B:394:LYS:HD2	1:B:394:LYS:HA	1.87	0.41
1:A:95:LEU:C	1:A:95:LEU:HD13	2.41	0.41
1:A:347:ASN:C	1:A:348:SER:HG	2.23	0.40
1:A:287:ALA:H	4:A:507:GOL:C1	2.35	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	355/371 (96%)	344 (97%)	11 (3%)	0	100	100
1	B	355/371 (96%)	344 (97%)	11 (3%)	0	100	100
All	All	710/742 (96%)	688 (97%)	22 (3%)	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	315/326 (97%)	315 (100%)	0	100	100
1	B	315/326 (97%)	313 (99%)	2 (1%)	86	90
All	All	630/652 (97%)	628 (100%)	2 (0%)	92	95

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	137	LEU
1	B	395	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	62	GLN
1	A	67	ASN
1	A	82	ASN
1	A	97	GLN
1	A	213	HIS
1	A	390	HIS
1	B	67	ASN
1	B	97	GLN
1	B	104	GLN
1	B	206	ASN
1	B	213	HIS

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Mol	Chain	Res	Type
1	B	368	GLN
1	B	395	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 16 ligands modelled in this entry, 2 are monoatomic - leaving 14 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	$\# Z > 2$
4	GOL	A	508	-	5,5,5	0.42	0	5,5,5	1.55	1 (20%)
3	SO4	B	506	-	4,4,4	0.30	0	6,6,6	0.34	0
4	GOL	B	507	-	5,5,5	0.38	0	5,5,5	0.37	0
4	GOL	A	507	-	5,5,5	0.48	0	5,5,5	1.25	1 (20%)
3	SO4	B	504	-	4,4,4	0.28	0	6,6,6	0.83	0
3	SO4	A	502	-	4,4,4	0.37	0	6,6,6	0.36	0
3	SO4	A	505	-	4,4,4	0.45	0	6,6,6	0.45	0
4	GOL	A	509	-	5,5,5	0.47	0	5,5,5	0.92	0
3	SO4	B	503	-	4,4,4	0.38	0	6,6,6	0.50	0
3	SO4	A	503	-	4,4,4	0.32	0	6,6,6	0.37	0
3	SO4	B	502	-	4,4,4	0.21	0	6,6,6	0.57	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	SO4	A	504	-	4,4,4	1.09	0	6,6,6	0.45	0
4	GOL	A	506	-	5,5,5	0.37	0	5,5,5	0.21	0
3	SO4	B	505	-	4,4,4	0.37	0	6,6,6	0.18	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
4	GOL	A	508	-	-	2/4/4/4	-
4	GOL	A	506	-	-	2/4/4/4	-
4	GOL	A	509	-	-	2/4/4/4	-
4	GOL	B	507	-	-	2/4/4/4	-
4	GOL	A	507	-	-	4/4/4/4	-

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	508	GOL	C3-C2-C1	2.91	123.03	111.70
4	A	507	GOL	O3-C3-C2	2.66	122.93	110.20

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	509	GOL	O1-C1-C2-C3
4	A	508	GOL	O1-C1-C2-O2
4	A	508	GOL	O1-C1-C2-C3
4	A	507	GOL	C1-C2-C3-O3
4	A	509	GOL	O1-C1-C2-O2
4	A	507	GOL	O2-C2-C3-O3
4	A	506	GOL	O1-C1-C2-O2
4	B	507	GOL	O1-C1-C2-C3
4	A	507	GOL	O1-C1-C2-C3
4	A	506	GOL	O1-C1-C2-C3
4	B	507	GOL	O1-C1-C2-O2
4	A	507	GOL	O1-C1-C2-O2

There are no ring outliers.

8 monomers are involved in 33 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	508	GOL	8	0
3	B	506	SO4	1	0
4	B	507	GOL	3	0
4	A	507	GOL	12	0
3	A	505	SO4	5	0
4	A	509	GOL	1	0
3	A	503	SO4	1	0
3	B	505	SO4	2	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

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, 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	359/371 (96%)	0.07	11 (3%) 49 48	21, 34, 55, 73	0
1	B	359/371 (96%)	0.18	21 (5%) 23 22	20, 35, 58, 74	0
All	All	718/742 (96%)	0.13	32 (4%) 33 32	20, 34, 57, 74	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	48	LYS	5.4
1	B	337	GLY	4.8
1	A	337	GLY	4.2
1	A	331	SER	4.2
1	A	48	LYS	4.1
1	B	168	PHE	3.7
1	B	268	HIS	3.6
1	B	319	PHE	3.5
1	B	332	THR	3.4
1	A	388	LYS	3.4
1	B	331	SER	3.1
1	B	208	TYR	3.1
1	B	388	LYS	3.0
1	A	367	LYS	2.9
1	A	332	THR	2.7
1	B	389	THR	2.7
1	B	394	LYS	2.7
1	B	393	GLU	2.5
1	A	267	LYS	2.4
1	A	394	LYS	2.4
1	B	47	ASP	2.4
1	A	366	GLY	2.4
1	B	49	LEU	2.3
1	B	368	GLN	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	319	PHE	2.2
1	B	398	TYR	2.2
1	B	386	LEU	2.1
1	B	205	SER	2.1
1	B	367	LYS	2.1
1	A	268	HIS	2.1
1	B	370	GLU	2.1
1	B	267	LYS	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 carbohydrates 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	GOL	A	507	6/6	0.69	0.40	48,53,57,57	0
3	SO4	B	505	5/5	0.75	0.26	49,50,54,54	5
4	GOL	B	507	6/6	0.79	0.19	50,53,55,55	0
4	GOL	A	509	6/6	0.80	0.22	49,50,55,61	0
4	GOL	A	508	6/6	0.84	0.33	23,36,39,44	0
4	GOL	A	506	6/6	0.89	0.29	39,50,59,60	0
3	SO4	A	505	5/5	0.92	0.16	47,54,66,68	0
3	SO4	B	506	5/5	0.92	0.22	50,56,68,70	0
3	SO4	A	504	5/5	0.98	0.09	44,46,48,48	0
3	SO4	B	504	5/5	0.98	0.14	40,40,43,45	0
2	ZN	A	501	1/1	0.99	0.12	26,26,26,26	0
3	SO4	B	503	5/5	0.99	0.11	20,21,22,23	0
3	SO4	A	503	5/5	0.99	0.15	42,48,49,56	0
3	SO4	B	502	5/5	1.00	0.11	19,20,22,22	0
2	ZN	B	501	1/1	1.00	0.11	25,25,25,25	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	SO4	A	502	5/5	1.00	0.10	22,23,27,27	0

6.5 Other polymers [i](#)

There are no such residues in this entry.