



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 06:21 am GMT

PDB ID : 4P3M  
Title : Crystal structure of serine hydroxymethyltransferase from *Psychromonas ingrahamii*  
Authors : Dworkowski, F.; Angelaccio, S.; Pascarella, S.; Capitani, G.  
Deposited on : 2014-03-09  
Resolution : 1.85 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 : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : 1.9-1692  
EDS : trunk28620  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
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) : recalc28949

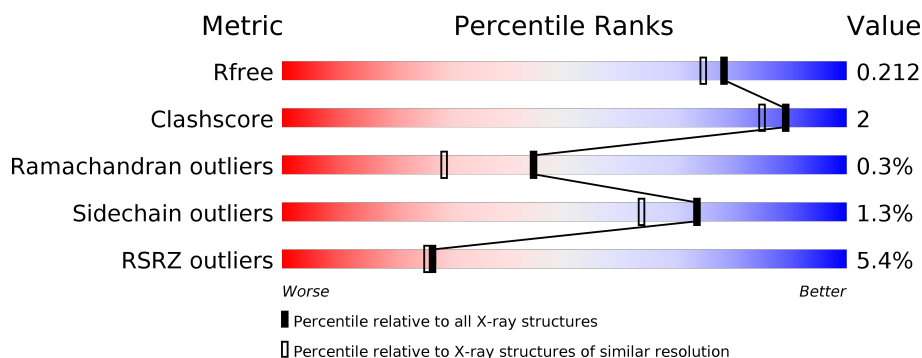
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.85 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))
$R_{free}$	100719	1923 (1.86-1.86)
Clashscore	112137	2083 (1.86-1.86)
Ramachandran outliers	110173	2060 (1.86-1.86)
Sidechain outliers	110143	2060 (1.86-1.86)
RSRZ outliers	101464	1932 (1.86-1.86)

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	429	
1	B	429	

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
4	GOL	A	509	-	-	-	X
4	GOL	A	510	-	-	-	X
4	GOL	A	511	-	-	-	X
4	GOL	B	502	-	-	-	X
4	GOL	B	503	-	-	-	X

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 12402 atoms, of which 5921 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 Serine hydroxymethyltransferase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	376	Total	C	H	N	O	S	0	10	0
			5806	1840	2907	494	550	15			
1	B	389	Total	C	H	N	O	S	0	2	0
			5922	1884	2958	503	561	16			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	422	LEU	-	expression tag	UNP A1SUU0
A	423	GLU	-	expression tag	UNP A1SUU0
A	424	HIS	-	expression tag	UNP A1SUU0
A	425	HIS	-	expression tag	UNP A1SUU0
A	426	HIS	-	expression tag	UNP A1SUU0
A	427	HIS	-	expression tag	UNP A1SUU0
A	428	HIS	-	expression tag	UNP A1SUU0
A	429	HIS	-	expression tag	UNP A1SUU0
B	422	LEU	-	expression tag	UNP A1SUU0
B	423	GLU	-	expression tag	UNP A1SUU0
B	424	HIS	-	expression tag	UNP A1SUU0
B	425	HIS	-	expression tag	UNP A1SUU0
B	426	HIS	-	expression tag	UNP A1SUU0
B	427	HIS	-	expression tag	UNP A1SUU0
B	428	HIS	-	expression tag	UNP A1SUU0
B	429	HIS	-	expression tag	UNP A1SUU0

- 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	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	3	Total	Cl	0	0
			3	3		

- Molecule 4 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
4	A	1	Total	C	H	O	0	0
			14	3	8	3		
4	A	1	Total	C	H	O	0	0
			14	3	8	3		
4	A	1	Total	C	H	O	0	0
			14	3	8	3		
4	A	1	Total	C	H	O	0	0
			14	3	8	3		
4	B	1	Total	C	H	O	0	0
			14	3	8	3		
4	B	1	Total	C	H	O	0	0
			14	3	8	3		
4	B	1	Total	C	H	O	0	0
			14	3	8	3		

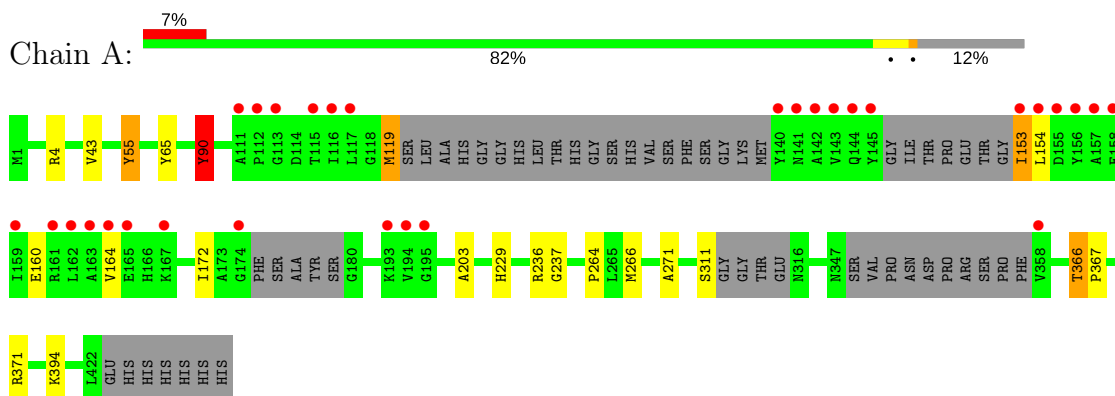
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	258	Total	O	0	0
			258	258		
5	B	290	Total	O	0	0
			290	290		

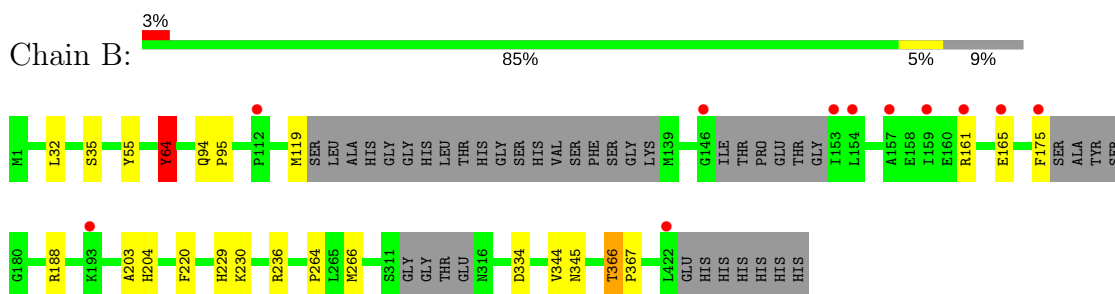
### 3 Residue-property plots

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



#### • Molecule 1: Serine hydroxymethyltransferase



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	146.82Å 64.86Å 116.94Å 90.00° 125.49° 90.00°	Depositor
Resolution (Å)	47.61 – 1.85 47.61 – 1.85	Depositor EDS
% Data completeness (in resolution range)	100.0 (47.61-1.85) 100.0 (47.61-1.85)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.61 (at 1.86Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, $R_{free}$	0.169 , 0.212 0.169 , 0.212	Depositor DCC
$R_{free}$ test set	1531 reflections (2.00%)	DCC
Wilson B-factor (Å <sup>2</sup> )	17.5	Xtriage
Anisotropy	0.468	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.45 , 55.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	12402	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.43% 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

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

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	0.83	5/2987 (0.2%)	0.90	9/4045 (0.2%)
1	B	0.90	9/3026 (0.3%)	0.77	10/4105 (0.2%)
All	All	0.86	14/6013 (0.2%)	0.84	19/8150 (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
1	A	0	1
1	B	0	1
All	All	0	2

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	64	TYR	CD1-CE1	29.54	1.83	1.39
1	A	90	TYR	CD2-CE2	29.08	1.82	1.39
1	A	90	TYR	CD1-CE1	-14.64	1.17	1.39
1	B	64	TYR	CG-CD1	11.85	1.54	1.39
1	A	90	TYR	CG-CD2	11.22	1.53	1.39
1	B	64	TYR	CD2-CE2	-11.00	1.22	1.39
1	B	175	PHE	CE1-CZ	10.91	1.58	1.37
1	B	64	TYR	CZ-OH	10.80	1.56	1.37
1	B	175	PHE	CG-CD1	9.70	1.53	1.38
1	B	175	PHE	CG-CD2	8.87	1.52	1.38
1	A	90	TYR	CE1-CZ	-7.80	1.28	1.38
1	A	90	TYR	CZ-OH	7.64	1.50	1.37
1	B	64	TYR	CE2-CZ	-7.46	1.28	1.38
1	B	64	TYR	CE1-CZ	-6.33	1.30	1.38

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	90	TYR	CE1-CZ-CE2	-23.88	81.58	119.80
1	A	90	TYR	CG-CD1-CE1	-16.91	107.77	121.30
1	A	90	TYR	CD1-CE1-CZ	16.01	134.21	119.80
1	B	64	TYR	CE1-CZ-CE2	-14.62	96.40	119.80
1	B	64	TYR	CD1-CG-CD2	-14.25	102.23	117.90
1	A	90	TYR	CD1-CG-CD2	-12.96	103.64	117.90
1	A	90	TYR	OH-CZ-CE2	-12.74	85.69	120.10
1	B	64	TYR	CZ-CE2-CD2	12.31	130.88	119.80
1	A	90	TYR	CB-CG-CD1	11.83	128.10	121.00
1	A	90	TYR	CG-CD2-CE2	-11.19	112.35	121.30
1	B	64	TYR	CG-CD1-CE1	-10.53	112.88	121.30
1	A	90	TYR	CE1-CZ-OH	10.44	148.28	120.10
1	A	90	TYR	CB-CG-CD2	9.93	126.96	121.00
1	B	64	TYR	CB-CG-CD1	9.25	126.55	121.00
1	B	64	TYR	OH-CZ-CE2	8.75	143.72	120.10
1	B	64	TYR	CG-CD2-CE2	-8.50	114.50	121.30
1	B	64	TYR	CE1-CZ-OH	-6.92	101.42	120.10
1	B	64	TYR	CD1-CE1-CZ	-5.83	114.56	119.80
1	B	64	TYR	CB-CG-CD2	5.38	124.23	121.00

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	90	TYR	Sidechain
1	B	64	TYR	Sidechain

## 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	2899	2907	2859	15	0
1	B	2964	2958	2950	11	0
2	A	20	0	0	1	0
2	B	5	0	0	0	0
3	A	3	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	24	32	32	0	0
4	B	18	24	24	1	0
5	A	258	0	0	3	0
5	B	290	0	0	1	0
All	All	6481	5921	5865	26	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 2.

All (26) 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:90:TYR:CE2	1:A:90:TYR:CD2	1.83	1.66
1:B:64:TYR:CD1	1:B:64:TYR:CE1	1.83	1.59
1:A:4:ARG:NH1	5:A:858:HOH:O	2.20	0.73
1:A:153:ILE:HD12	1:A:154:LEU:N	2.04	0.72
1:A:371[B]:ARG:NH2	5:A:703:HOH:O	2.22	0.71
1:A:153:ILE:HD12	1:A:154:LEU:H	1.63	0.63
1:A:55:TYR:OH	2:A:503:SO4:O4	2.12	0.58
1:A:43:VAL:HG13	1:A:271:ALA:HB1	1.89	0.54
1:A:229:HIS:CD2	1:A:236:ARG:HA	2.44	0.53
1:B:35:SER:HB2	1:B:230:LYS:HE3	1.92	0.52
1:B:229:HIS:CD2	1:B:236:ARG:HA	2.47	0.49
1:B:366:THR:N	1:B:367:PRO:CD	2.76	0.49
1:B:188:ARG:HD2	1:B:220:PHE:O	2.14	0.48
1:A:119:MET:HE1	1:A:172:ILE:HG22	1.97	0.46
1:B:161:ARG:NH1	1:B:165:GLU:OE1	2.49	0.45
1:A:203:ALA:O	5:A:653:HOH:O	2.21	0.45
1:A:65:TYR:CD2	1:B:345:ASN:ND2	2.85	0.45
1:A:160:GLU:O	1:A:164:VAL:HG23	2.17	0.44
1:B:264:PRO:HG2	1:B:266:MET:SD	2.58	0.44
1:A:229:HIS:HD2	1:A:237:GLY:H	1.64	0.44
1:B:203:ALA:O	1:B:204:HIS:HB2	2.17	0.44
1:B:32:LEU:HB2	1:B:344:VAL:HG12	2.01	0.43
1:A:366:THR:N	1:A:367:PRO:CD	2.82	0.43
4:B:502:GOL:H31	5:B:688:HOH:O	2.18	0.42
1:B:94:GLN:N	1:B:95:PRO:CD	2.83	0.41
1:A:264:PRO:HG2	1:A:266:MET:SD	2.61	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	374/429 (87%)	368 (98%)	5 (1%)	1 (0%)	44	29
1	B	381/429 (89%)	373 (98%)	7 (2%)	1 (0%)	44	29
All	All	755/858 (88%)	741 (98%)	12 (2%)	2 (0%)	44	29

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	366	THR
1	B	366	THR

### 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	300/335 (90%)	295 (98%)	5 (2%)	66	52
1	B	306/335 (91%)	303 (99%)	3 (1%)	80	74
All	All	606/670 (90%)	598 (99%)	8 (1%)	73	63

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	55	TYR
1	A	119	MET
1	A	153	ILE
1	A	311	SER

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Mol	Chain	Res	Type
1	A	394	LYS
1	B	55	TYR
1	B	119	MET
1	B	334	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	229	HIS

### 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](#)

Of 15 ligands modelled in this entry, 3 are monoatomic - leaving 12 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	SO4	A	501	-	4,4,4	0.19	0	6,6,6	0.17	0
2	SO4	A	502	-	4,4,4	0.22	0	6,6,6	0.62	0
2	SO4	A	503	-	4,4,4	0.38	0	6,6,6	0.53	0
2	SO4	A	504	-	4,4,4	0.22	0	6,6,6	0.18	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
4	GOL	A	508	-	5,5,5	0.38	0	5,5,5	0.68	0
4	GOL	A	509	-	5,5,5	0.49	0	5,5,5	0.77	0
4	GOL	A	510	-	5,5,5	0.30	0	5,5,5	0.50	0
4	GOL	A	511	-	5,5,5	0.70	0	5,5,5	1.65	2 (40%)
2	SO4	B	501	-	4,4,4	0.17	0	6,6,6	0.51	0
4	GOL	B	502	-	5,5,5	0.53	0	5,5,5	0.85	0
4	GOL	B	503	-	5,5,5	0.40	0	5,5,5	0.32	0
4	GOL	B	504	-	5,5,5	0.39	0	5,5,5	0.28	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
2	SO4	A	501	-	-	0/0/0/0	0/0/0/0
2	SO4	A	502	-	-	0/0/0/0	0/0/0/0
2	SO4	A	503	-	-	0/0/0/0	0/0/0/0
2	SO4	A	504	-	-	0/0/0/0	0/0/0/0
4	GOL	A	508	-	-	0/4/4/4	0/0/0/0
4	GOL	A	509	-	-	0/4/4/4	0/0/0/0
4	GOL	A	510	-	-	0/4/4/4	0/0/0/0
4	GOL	A	511	-	-	0/4/4/4	0/0/0/0
2	SO4	B	501	-	-	0/0/0/0	0/0/0/0
4	GOL	B	502	-	-	0/4/4/4	0/0/0/0
4	GOL	B	503	-	-	0/4/4/4	0/0/0/0
4	GOL	B	504	-	-	0/4/4/4	0/0/0/0

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	511	GOL	O1-C1-C2	-2.40	97.98	110.07
4	A	511	GOL	O2-C2-C1	-2.24	98.28	108.84

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	503	SO4	1	0
4	B	502	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

### 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	376/429 (87%)	-0.06	30 (7%)	13 13	10, 19, 50, 71	4 (1%)
1	B	389/429 (90%)	-0.31	11 (2%)	53 52	11, 19, 44, 63	3 (0%)
All	All	765/858 (89%)	-0.19	41 (5%)	26 26	10, 19, 48, 71	7 (0%)

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	162	LEU	5.2
1	A	165[A]	GLU	5.1
1	A	145	TYR	4.9
1	A	144	GLN	4.8
1	A	113	GLY	4.6
1	A	153	ILE	4.3
1	A	112	PRO	4.3
1	A	164	VAL	4.2
1	B	153	ILE	4.0
1	A	143	VAL	4.0
1	A	161	ARG	3.6
1	A	154	LEU	3.6
1	A	194	VAL	3.6
1	A	142	ALA	3.5
1	B	175	PHE	3.5
1	A	140	TYR	3.5
1	A	159	ILE	3.4
1	A	155	ASP	3.3
1	B	157	ALA	3.3
1	A	167	LYS	3.1
1	A	163	ALA	3.0
1	A	156	TYR	2.9
1	A	141	ASN	2.9
1	A	193	LYS	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	422	LEU	2.7
1	A	117	LEU	2.7
1	B	154	LEU	2.7
1	A	115	THR	2.6
1	B	193	LYS	2.6
1	A	116	ILE	2.5
1	A	111	ALA	2.5
1	A	157	ALA	2.4
1	A	358	VAL	2.4
1	B	146	GLY	2.4
1	A	158	GLU	2.4
1	B	161	ARG	2.3
1	A	174	GLY	2.3
1	A	195	GLY	2.2
1	B	165	GLU	2.1
1	B	112	PRO	2.1
1	B	159	ILE	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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
4	GOL	A	511	6/6	0.78	0.30	19.07	25,36,55,64	0
4	GOL	B	502	6/6	0.89	0.16	6.75	24,32,44,47	0
4	GOL	A	509	6/6	0.74	0.20	5.13	35,47,56,56	0
4	GOL	A	510	6/6	0.94	0.10	3.43	21,29,38,43	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
4	GOL	B	503	6/6	0.87	0.18	3.11	30,47,59,67	0
4	GOL	A	508	6/6	0.94	0.12	1.72	18,30,44,53	0
2	SO4	A	503	5/5	0.97	0.13	0.99	28,29,33,48	0
2	SO4	A	502	5/5	0.97	0.12	0.35	20,23,26,28	0
3	CL	A	507	1/1	1.00	0.09	-0.20	28,28,28,28	0
2	SO4	A	501	5/5	0.99	0.09	-0.35	23,27,29,36	0
2	SO4	B	501	5/5	0.95	0.10	-0.40	44,46,54,58	0
4	GOL	B	504	6/6	0.97	0.07	-0.88	19,26,46,56	0
3	CL	A	505	1/1	1.00	0.06	-2.47	23,23,23,23	0
3	CL	A	506	1/1	1.00	0.04	-3.54	21,21,21,21	0
2	SO4	A	504	5/5	0.97	0.08	-	37,45,54,65	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.