



# Full wwPDB X-ray Structure Validation Report ⓘ

May 29, 2020 – 04:59 pm BST

PDB ID : 2X1C  
Title : The crystal structure of precursor acyl coenzyme A:isopenicillin N acyltransferase from *Penicillium chrysogenum*  
Authors : Bokhove, M.; Yoshida, H.; Hensgens, C.M.H.; van der Laan, J.M.; Sutherland, J.D.; Dijkstra, B.W.  
Deposited on : 2009-12-23  
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

<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 : 1.8.5 (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

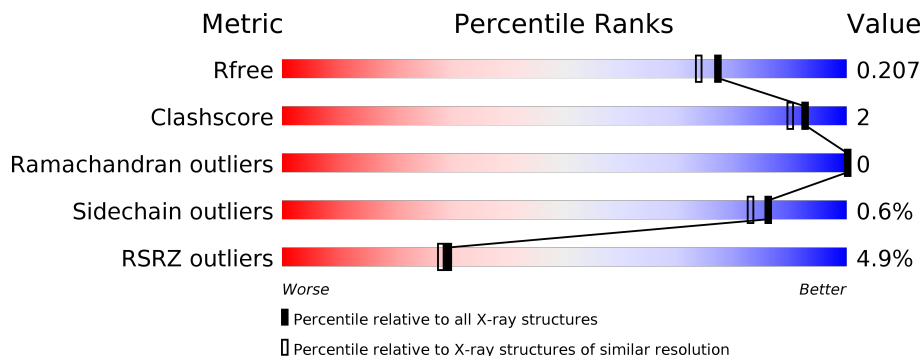
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 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))
$R_{free}$	130704	2469 (1.86-1.86)
Clashscore	141614	2625 (1.86-1.86)
Ramachandran outliers	138981	2592 (1.86-1.86)
Sidechain outliers	138945	2592 (1.86-1.86)
RSRZ outliers	127900	2436 (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 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	357	<p>3% 96% 5%</p>
1	B	357	<p>% 95% 5%</p>
1	C	357	<p>4% 96% 5%</p>
1	D	357	<p>11% 96% 5%</p>

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 12046 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 ACYL-COENZYME.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	355	2833	1777	512	531	13	0	5	0
1	B	356	2877	1803	518	542	14	0	9	0
1	C	355	2842	1782	509	538	13	0	7	0
1	D	355	2836	1778	513	533	12	0	5	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	103	ALA	CYS	engineered mutation	UNP P15802
B	103	ALA	CYS	engineered mutation	UNP P15802
C	103	ALA	CYS	engineered mutation	UNP P15802
D	103	ALA	CYS	engineered mutation	UNP P15802

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

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

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	2	Total Cl 2 2	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	161	Total O 161 161	0	0
5	B	221	Total O 221 221	0	0

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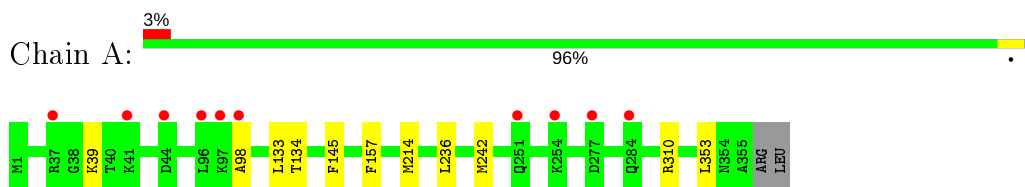
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<b>Mol</b>	<b>Chain</b>	<b>Residues</b>	<b>Atoms</b>		<b>ZeroOcc</b>	<b>AltConf</b>
5	C	138	Total 138	O 138	0	0
5	D	65	Total 65	O 65	0	0

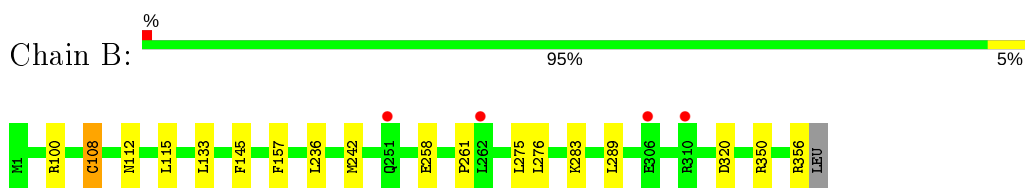
### 3 Residue-property plots [i](#)

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 ( $RSRZ > 2$ ). 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.

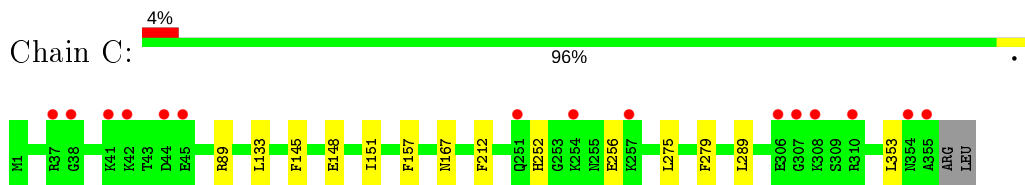
- Molecule 1: ACYL-COENZYME



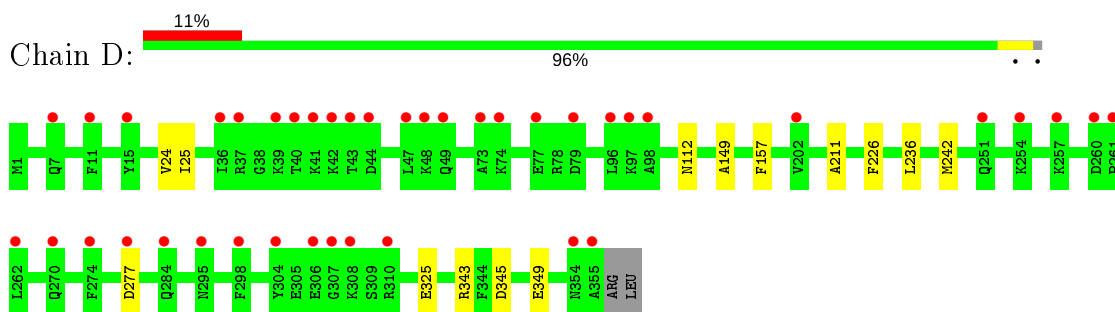
- Molecule 1: ACYL-COENZYME



- Molecule 1: ACYL-COENZYME



- Molecule 1: ACYL-COENZYME



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	231.22Å 68.23Å 151.28Å 90.00° 129.57° 90.00°	Depositor
Resolution (Å)	48.80 – 1.85 48.80 – 1.85	Depositor EDS
% Data completeness (in resolution range)	98.5 (48.80-1.85) 98.4 (48.80-1.85)	Depositor EDS
$R_{merge}$	0.04	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.30 (at 1.86Å)	Xtrriage
Refinement program	REFMAC 5.5.0102	Depositor
R, $R_{free}$	0.169 , 0.193 0.189 , 0.207	Depositor DCC
$R_{free}$ test set	7647 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	25.1	Xtrriage
Anisotropy	0.062	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.40 , 46.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.012 for -h-2*1,-k,l	Xtrriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	12046	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	29.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.52% 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: 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	0.52	0/2890	0.58	0/3900
1	B	0.56	1/2934 (0.0%)	0.60	0/3959
1	C	0.48	0/2899	0.55	0/3915
1	D	0.42	0/2893	0.53	0/3905
All	All	0.50	1/11616 (0.0%)	0.56	0/15679

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	108	CYS	CB-SG	-5.52	1.72	1.81

There are no bond angle outliers.

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	2833	0	2782	10	0
1	B	2877	0	2815	12	0
1	C	2842	0	2778	7	0
1	D	2836	0	2780	6	0
2	A	20	0	0	0	0
2	B	10	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	5	0	0	0	0
3	A	6	0	8	1	0
3	B	18	0	24	3	0
3	C	12	0	16	0	0
4	B	2	0	0	0	0
5	A	161	0	0	0	0
5	B	221	0	0	1	0
5	C	138	0	0	0	0
5	D	65	0	0	0	0
All	All	12046	0	11203	34	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 (34) 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:236:LEU:HD23	1:A:242[B]:MET:HG3	1.54	0.88
1:B:350:ARG:HD3	1:B:356:ARG:HB3	1.68	0.76
1:D:236:LEU:HD23	1:D:242:MET:HG3	1.75	0.68
1:B:236:LEU:HD23	1:B:242[A]:MET:HG3	1.76	0.66
1:B:283:LYS:HE2	1:B:320:ASP:OD1	1.98	0.62
1:B:283:LYS:HE2	1:B:320:ASP:CG	2.21	0.61
1:C:275:LEU:HB3	1:C:289:LEU:HD22	1.86	0.56
1:A:236:LEU:HD23	1:A:242[B]:MET:CG	2.29	0.56
1:B:100:ARG:HE	3:B:1360:GOL:C1	2.21	0.54
1:B:100:ARG:HE	3:B:1360:GOL:H12	1.71	0.54
1:B:133:LEU:HD12	1:B:145:PHE:CZ	2.44	0.52
1:A:133:LEU:HD12	1:A:145:PHE:CZ	2.45	0.52
3:B:1360:GOL:H11	5:B:2074:HOH:O	2.10	0.52
1:C:89:ARG:NH2	1:C:148:GLU:OE1	2.42	0.51
1:A:353:LEU:HG	1:D:112[B]:ASN:ND2	2.27	0.50
1:B:258:GLU:HG2	1:B:261:PRO:HG3	1.94	0.49
1:A:39:LYS:NZ	1:A:98:ALA:H	2.10	0.49
1:B:115:LEU:HD21	1:B:283:LYS:HE3	1.95	0.48
1:A:98:ALA:HB1	1:A:310:ARG:HH12	1.79	0.47
1:B:275:LEU:HB3	1:B:289:LEU:HD22	1.97	0.47
1:C:252:HIS:HB3	1:C:256:GLU:HG3	1.98	0.46
1:A:134:THR:HG21	3:A:1360:GOL:H32	1.98	0.46
1:B:112:ASN:ND2	1:C:353:LEU:HG	2.31	0.45
1:D:345:ASP:O	1:D:349:GLU:HG2	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:236:LEU:CD2	1:A:242[B]:MET:HG3	2.37	0.44
1:C:167:ASN:HB2	1:C:212:PHE:HB3	2.00	0.43
1:C:133:LEU:HD12	1:C:145:PHE:CZ	2.53	0.43
1:C:148:GLU:HB2	1:C:151:ILE:HD12	2.00	0.43
1:D:211:ALA:HB3	1:D:226:PHE:CZ	2.54	0.42
1:B:108:CYS:SG	1:B:276:LEU:HD22	2.60	0.42
1:D:25:ILE:HG13	1:D:149:ALA:HB1	2.02	0.41
1:A:242[B]:MET:HE3	1:A:242[B]:MET:HB3	1.81	0.41
1:A:214:MET:CE	1:A:242[B]:MET:HE2	2.52	0.40
1:D:325:GLU:HG2	1:D:343:ARG:HG2	2.03	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	358/357 (100%)	347 (97%)	11 (3%)	0	100	100
1	B	363/357 (102%)	355 (98%)	8 (2%)	0	100	100
1	C	360/357 (101%)	351 (98%)	9 (2%)	0	100	100
1	D	358/357 (100%)	351 (98%)	7 (2%)	0	100	100
All	All	1439/1428 (101%)	1404 (98%)	35 (2%)	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	294/291 (101%)	293 (100%)	1 (0%)	92	91
1	B	299/291 (103%)	298 (100%)	1 (0%)	92	91
1	C	296/291 (102%)	294 (99%)	2 (1%)	84	79
1	D	294/291 (101%)	291 (99%)	3 (1%)	76	69
All	All	1183/1164 (102%)	1176 (99%)	7 (1%)	86	83

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

Mol	Chain	Res	Type
1	A	157	PHE
1	B	157	PHE
1	C	157	PHE
1	C	279	PHE
1	D	24	VAL
1	D	157	PHE
1	D	277	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	B	251	GLN

### 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, 2 are monoatomic - leaving 13 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
3	GOL	B	1359	-	5,5,5	0.44	0	5,5,5	0.73	0
2	SO4	D	1356	-	4,4,4	0.16	0	6,6,6	0.29	0
3	GOL	C	1356	-	5,5,5	0.46	0	5,5,5	0.50	0
3	GOL	B	1360	-	5,5,5	0.41	0	5,5,5	0.42	0
2	SO4	A	1359	-	4,4,4	0.17	0	6,6,6	0.05	0
3	GOL	C	1357	-	5,5,5	0.36	0	5,5,5	0.35	0
2	SO4	B	1357	-	4,4,4	0.06	0	6,6,6	0.23	0
3	GOL	B	1361	-	5,5,5	0.44	0	5,5,5	0.42	0
3	GOL	A	1360	-	5,5,5	0.40	0	5,5,5	0.33	0
2	SO4	B	1358	-	4,4,4	0.16	0	6,6,6	0.08	0
2	SO4	A	1356	-	4,4,4	0.11	0	6,6,6	0.40	0
2	SO4	A	1358	-	4,4,4	0.11	0	6,6,6	0.13	0
2	SO4	A	1357	-	4,4,4	0.16	0	6,6,6	0.30	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	C	1356	-	-	4/4/4/4	-
3	GOL	B	1360	-	-	1/4/4/4	-
3	GOL	A	1360	-	-	0/4/4/4	-
3	GOL	C	1357	-	-	4/4/4/4	-
3	GOL	B	1361	-	-	4/4/4/4	-
3	GOL	B	1359	-	-	4/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	1356	GOL	O1-C1-C2-C3
3	B	1359	GOL	C1-C2-C3-O3
3	B	1359	GOL	O2-C2-C3-O3
3	C	1357	GOL	C1-C2-C3-O3
3	B	1361	GOL	O1-C1-C2-C3
3	B	1361	GOL	C1-C2-C3-O3
3	C	1357	GOL	O1-C1-C2-O2
3	C	1356	GOL	C1-C2-C3-O3
3	B	1359	GOL	O1-C1-C2-C3
3	C	1357	GOL	O1-C1-C2-C3
3	C	1356	GOL	O2-C2-C3-O3
3	B	1359	GOL	O1-C1-C2-O2
3	C	1357	GOL	O2-C2-C3-O3
3	B	1361	GOL	O1-C1-C2-O2
3	B	1361	GOL	O2-C2-C3-O3
3	C	1356	GOL	O1-C1-C2-O2
3	B	1360	GOL	C1-C2-C3-O3

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1360	GOL	3	0
3	A	1360	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	355/357 (99%)	-0.20	10 (2%) 53 52	15, 24, 42, 58	0
1	B	356/357 (99%)	-0.11	4 (1%) 80 81	14, 21, 36, 50	0
1	C	355/357 (99%)	0.09	15 (4%) 36 34	16, 28, 49, 66	0
1	D	355/357 (99%)	0.69	41 (11%) 4 4	22, 37, 63, 79	0
All	All	1421/1428 (99%)	0.12	70 (4%) 29 28	14, 27, 50, 79	0

All (70) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	98	ALA	8.6
1	C	355	ALA	6.2
1	D	355	ALA	5.4
1	D	298	PHE	5.2
1	C	310	ARG	4.7
1	D	98	ALA	4.6
1	D	44	ASP	4.6
1	D	41	LYS	4.5
1	C	254	LYS	4.3
1	D	36	ILE	4.3
1	D	15	TYR	4.2
1	D	307	GLY	4.2
1	C	307	GLY	3.8
1	D	304	TYR	3.8
1	C	354	ASN	3.8
1	C	41	LYS	3.8
1	D	274	PHE	3.6
1	D	47	LEU	3.6
1	D	43	THR	3.5
1	C	306	GLU	3.5
1	D	262	LEU	3.4

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
1	D	354	ASN	3.3
1	D	37	ARG	3.3
1	D	310	ARG	3.3
1	D	49	GLN	3.2
1	D	306	GLU	3.1
1	B	306	GLU	3.1
1	D	77	GLU	3.0
1	D	79	ASP	3.0
1	A	41	LYS	3.0
1	D	11	PHE	2.9
1	D	39	LYS	2.9
1	D	40	THR	2.9
1	C	37	ARG	2.9
1	D	260	ASP	2.8
1	D	73	ALA	2.8
1	D	308	LYS	2.7
1	A	254	LYS	2.6
1	A	284	GLN	2.6
1	D	270	GLN	2.5
1	D	254	LYS	2.5
1	D	261	PRO	2.4
1	A	44	ASP	2.4
1	D	277	ASP	2.4
1	D	74	LYS	2.4
1	D	251	GLN	2.4
1	A	251	GLN	2.4
1	C	308	LYS	2.3
1	D	295	ASN	2.3
1	D	202	VAL	2.3
1	B	310	ARG	2.3
1	A	97	LYS	2.3
1	C	257	LYS	2.3
1	B	262	LEU	2.3
1	B	251	GLN	2.3
1	C	38	GLY	2.2
1	D	257	LYS	2.2
1	D	284	GLN	2.2
1	A	37	ARG	2.2
1	A	96	LEU	2.2
1	D	48	LYS	2.2
1	D	97	LYS	2.1
1	D	7	GLN	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	251	GLN	2.1
1	C	44	ASP	2.1
1	C	45	GLU	2.1
1	D	42	LYS	2.1
1	D	96	LEU	2.1
1	C	42	LYS	2.1
1	A	277	ASP	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, 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( $\text{\AA}^2$ )	Q<0.9
3	GOL	C	1357	6/6	0.77	0.19	49,52,53,54	0
3	GOL	B	1360	6/6	0.81	0.19	43,44,46,47	0
3	GOL	B	1359	6/6	0.83	0.22	36,42,43,45	0
3	GOL	B	1361	6/6	0.86	0.17	31,36,38,40	0
2	SO4	B	1358	5/5	0.92	0.23	73,73,73,73	0
3	GOL	A	1360	6/6	0.93	0.17	34,36,36,36	0
3	GOL	C	1356	6/6	0.93	0.11	38,41,41,42	0
2	SO4	A	1359	5/5	0.94	0.30	72,72,72,72	0
2	SO4	A	1358	5/5	0.94	0.17	38,38,40,40	0
2	SO4	A	1357	5/5	0.96	0.09	32,32,34,35	0
2	SO4	B	1357	5/5	0.97	0.09	25,28,30,30	0
2	SO4	D	1356	5/5	0.97	0.06	35,35,35,37	0
4	CL	B	1363	1/1	0.99	0.31	50,50,50,50	1
2	SO4	A	1356	5/5	0.99	0.07	25,27,27,28	0
4	CL	B	1362	1/1	1.00	0.13	16,16,16,16	1

## 6.5 Other polymers [i](#)

There are no such residues in this entry.