



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2014 – 03:19 AM GMT

PDB ID : 4IQF  
Title : Crystal Structure of Methionyl-tRNA Formyltransferase from *Bacillus anthracis*  
Authors : Kim, Y.; Makowska-Grzyska, M.; Kwon, K.; Anderson, W.F.; Joachimiak, A.;  
Center for Structural Genomics of Infectious Diseases (CSGID)  
Deposited on : 2013-01-11  
Resolution : 2.38 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

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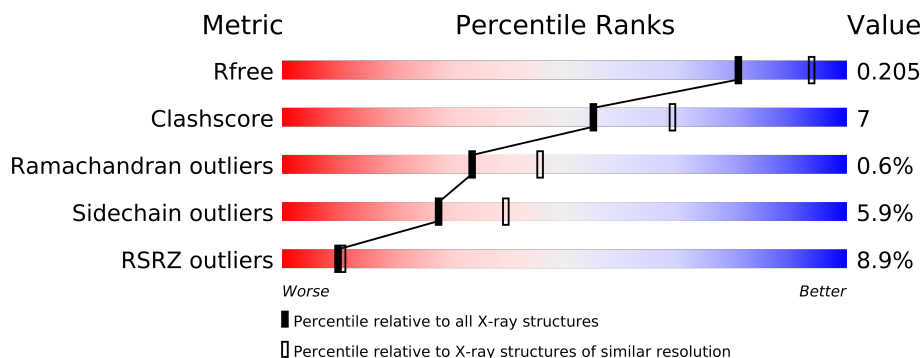
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 2.38 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))
$R_{free}$	66092	2963 (2.40-2.36)
Clashscore	79885	3668 (2.40-2.36)
Ramachandran outliers	78287	3600 (2.40-2.36)
Sidechain outliers	78261	3602 (2.40-2.36)
RSRZ outliers	66119	2966 (2.40-2.36)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	317	
1	B	317	
1	C	317	
1	D	317	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
3	GOL	A	402	-	X
4	SO4	B	402	-	X

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 10413 atoms, of which 0 are hydrogen and 0 are deuterium.

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.

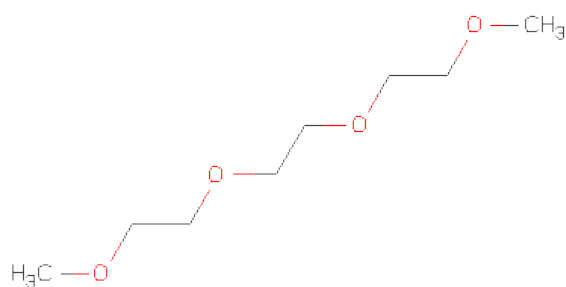
- Molecule 1 is a protein called Methionyl-tRNA formyltransferase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	314	Total	C	N	O	S	Se	0	1	0
			2450	1567	409	466	2	6			
1	B	314	Total	C	N	O	S	Se	0	1	0
			2458	1575	410	465	2	6			
1	C	315	Total	C	N	O	S	Se	0	1	0
			2464	1578	411	467	2	6			
1	D	315	Total	C	N	O	S	Se	10	3	0
			2483	1589	415	471	2	6			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	EXPRESSION TAG	UNP Q81WH2
A	-1	ASN	-	EXPRESSION TAG	UNP Q81WH2
A	0	ALA	-	EXPRESSION TAG	UNP Q81WH2
B	-2	SER	-	EXPRESSION TAG	UNP Q81WH2
B	-1	ASN	-	EXPRESSION TAG	UNP Q81WH2
B	0	ALA	-	EXPRESSION TAG	UNP Q81WH2
C	-2	SER	-	EXPRESSION TAG	UNP Q81WH2
C	-1	ASN	-	EXPRESSION TAG	UNP Q81WH2
C	0	ALA	-	EXPRESSION TAG	UNP Q81WH2
D	-2	SER	-	EXPRESSION TAG	UNP Q81WH2
D	-1	ASN	-	EXPRESSION TAG	UNP Q81WH2
D	0	ALA	-	EXPRESSION TAG	UNP Q81WH2

- Molecule 2 is 1-METHOXY-2-[2-(2-METHOXY-ETHOXY)-ETHANE (three-letter code: PG5) (formula: C<sub>8</sub>H<sub>18</sub>O<sub>4</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			10	6	4		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula:  $C_3H_8O_3$ ).



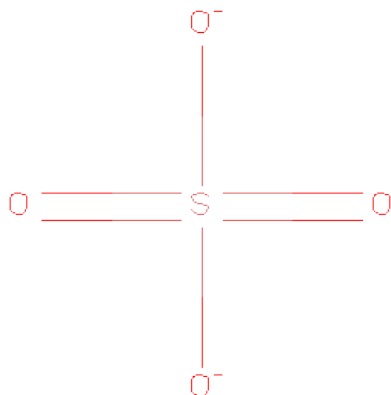
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	C	1	Total	C	O	0	0
			6	3	3		

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

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



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

- Molecule 5 is water.

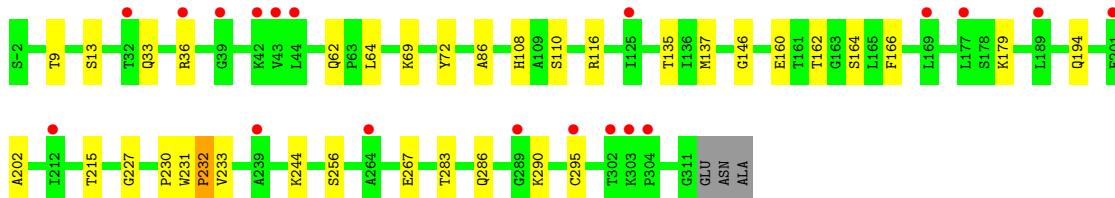
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	150	Total	O	0	0
			150	150		
5	B	122	Total	O	0	0
			122	122		
5	C	97	Total	O	0	0
			97	97		
5	D	135	Total	O	0	0
			135	135		

### 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 errors 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.

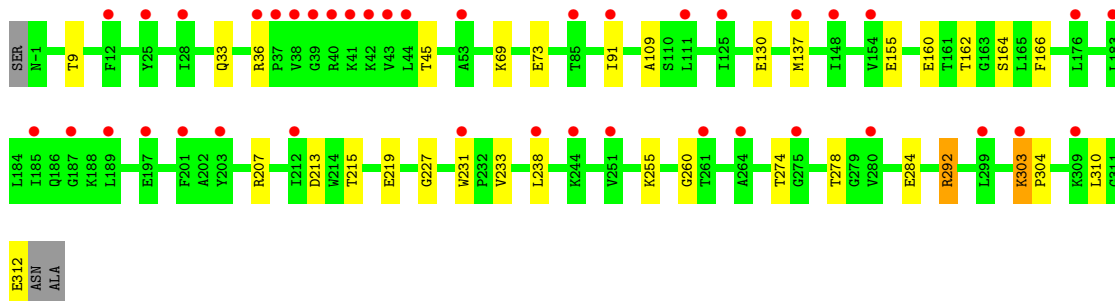
- Molecule 1: Methionyl-tRNA formyltransferase

Chain A: 



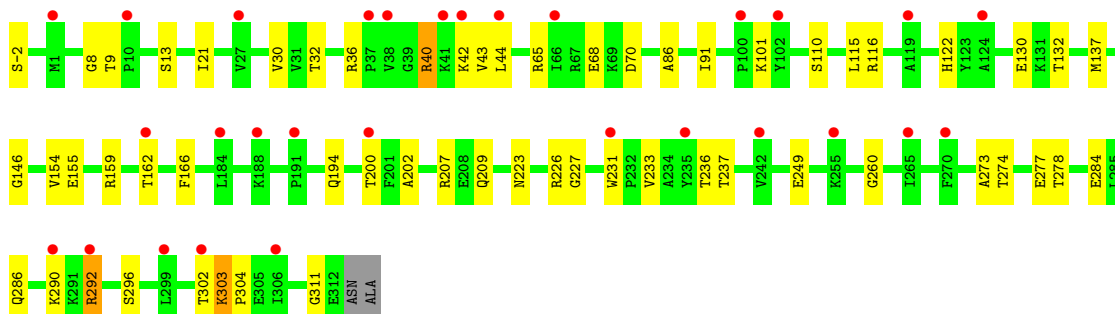
- Molecule 1: Methionyl-tRNA formyltransferase

Chain B: 



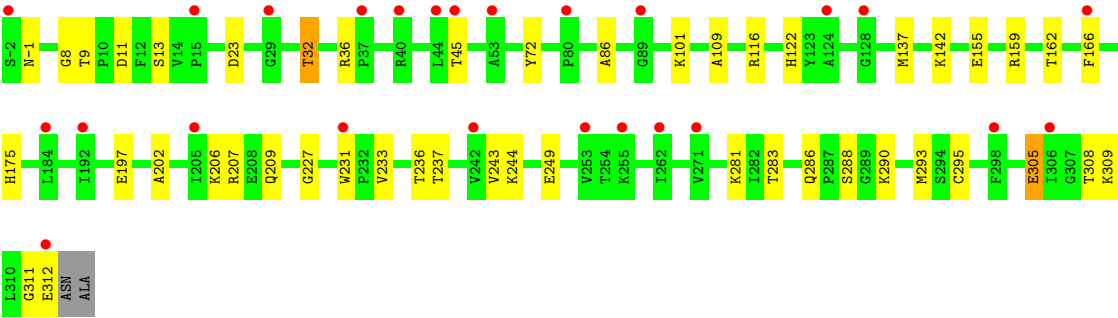
- Molecule 1: Methionyl-tRNA formyltransferase

Chain C: 



- Molecule 1: Methionyl-tRNA formyltransferase

Chain D: 



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	62.86Å 191.50Å 89.95Å 90.00° 89.95° 90.00°	Depositor
Resolution (Å)	36.59 – 2.38 36.59 – 2.38	Depositor EDS
% Data completeness (in resolution range)	97.5 (36.59-2.38) 97.5 (36.59-2.38)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.11	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.72 (at 2.37Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.1_1161)	Depositor
R, $R_{free}$	0.181 , 0.210 0.180 , 0.205	Depositor DCC
$R_{free}$ test set	4126 reflections (5.24%)	DCC
Wilson B-factor (Å <sup>2</sup> )	44.4	Xtriage
Anisotropy	0.677	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 41.4	EDS
Estimated twinning fraction	0.438 for h,-k,-l	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	1 of 82905 reflections (0.001%)	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	10413	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 35.26 % 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 5.9941e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<sup>1</sup>Intensities estimated from amplitudes.



## 5 Model quality ⓘ

### 5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, PG5, 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	0.30	2/2491 (0.1%)	0.46	0/3369
1	B	0.27	1/2501 (0.0%)	0.45	1/3385 (0.0%)
1	C	0.28	1/2507 (0.0%)	0.45	1/3393 (0.0%)
1	D	0.23	0/2527	0.44	0/3420
All	All	0.27	4/10026 (0.0%)	0.45	2/13567 (0.0%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	232	PRO	N-CD	5.41	1.55	1.47
1	C	304	PRO	N-CD	5.20	1.55	1.47
1	B	304	PRO	N-CD	5.13	1.55	1.47
1	A	230	PRO	N-CD	5.02	1.54	1.47

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	303	LYS	C-N-CD	5.59	140.14	128.40
1	C	303	LYS	C-N-CD	5.55	140.06	128.40

There are no chirality outliers.

There are no planarity outliers.

### 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the

chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2450	0	0	14	0
1	B	2458	0	0	10	0
1	C	2464	0	0	22	0
1	D	2483	0	0	20	0
2	A	10	0	12	0	0
3	A	6	0	8	1	0
3	B	6	0	8	1	0
3	C	6	0	8	0	0
3	D	6	0	8	1	0
4	A	5	0	0	0	0
4	B	5	0	0	0	0
4	D	10	0	0	0	0
5	A	150	0	0	4	0
5	B	122	0	0	2	0
5	C	97	0	0	4	0
5	D	135	0	0	4	0
All	All	10413	0	44	67	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 7.

All (67) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:237:THR:O	1:C:311:GLY:N	2.26	0.69
1:B:109:ALA:N	3:B:401:GOL:HO1	1.92	0.68
1:A:69:LYS:NZ	1:A:72:TYR:OH	2.30	0.64
1:D:122:HIS:NE2	1:D:231[B]:TRP:O	2.33	0.61
1:D:122:HIS:NE2	1:D:231[A]:TRP:O	2.34	0.60
1:C:146:GLY:O	1:C:194:GLN:NE2	2.37	0.58
1:D:116:ARG:O	1:D:202:ALA:N	2.37	0.58
1:C:13:SER:OG	1:C:86:ALA:O	2.23	0.57
1:C:159:ARG:NH1	1:C:249:GLU:OE2	2.38	0.56
1:B:213:ASP:OD1	1:B:215:THR:OG1	2.25	0.55
1:A:13:SER:OG	1:A:86:ALA:O	2.24	0.55
1:C:284:GLU:OE2	1:C:292:ARG:NH2	2.40	0.53
1:C:122:HIS:NE2	1:C:231[A]:TRP:O	2.42	0.53
1:B:260:GLY:O	1:B:310:LEU:N	2.42	0.53
1:C:162:THR:OG1	1:C:227:GLY:O	2.27	0.53
1:A:62:GLN:NE2	5:A:501:HOH:O	2.41	0.52
1:C:122:HIS:NE2	1:C:231[B]:TRP:O	2.44	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:284:GLU:OE2	1:B:292:ARG:NH2	2.44	0.51
1:B:162:THR:OG1	1:B:227:GLY:O	2.27	0.51
1:A:179:LYS:NZ	5:A:574:HOH:O	2.43	0.51
1:C:40:ARG:NE	5:C:570:HOH:O	2.43	0.51
1:D:305:GLU:O	1:D:308:THR:OG1	2.28	0.51
1:D:159:ARG:NH1	1:D:249:GLU:OE2	2.44	0.50
1:C:116:ARG:O	1:C:202:ALA:N	2.44	0.50
1:D:237:THR:OG1	1:D:311:GLY:CA	2.59	0.50
1:D:206:LYS:N	1:D:209:GLN:OE1	2.45	0.50
1:B:160:GLU:OE1	1:B:164:SER:OG	2.30	0.49
1:D:13:SER:OG	1:D:86:ALA:O	2.30	0.49
1:D:283:THR:O	1:D:295:CYS:N	2.46	0.49
1:C:237:THR:OG1	1:C:311:GLY:CA	2.61	0.48
1:A:146:GLY:O	1:A:194:GLN:NE2	2.47	0.48
1:C:65:ARG:NH2	1:C:68:GLU:OE2	2.46	0.47
1:A:116:ARG:O	1:A:202:ALA:N	2.47	0.47
1:B:219:GLU:OE2	1:B:278:THR:OG1	2.32	0.47
1:D:309:LYS:O	5:D:535:HOH:O	2.21	0.47
1:A:110:SER:O	1:A:135:THR:OG1	2.33	0.46
1:A:283:THR:O	1:A:295:CYS:N	2.49	0.46
1:B:213:ASP:N	5:B:607:HOH:O	2.49	0.46
1:D:290:LYS:NZ	5:D:630:HOH:O	2.48	0.46
1:C:-2:SER:N	5:C:553:HOH:O	2.48	0.46
1:D:162:THR:OG1	1:D:227:GLY:O	2.33	0.46
1:D:142:LYS:NZ	5:D:528:HOH:O	2.48	0.46
1:C:260:GLY:N	1:C:273:ALA:O	2.50	0.45
1:A:162:THR:OG1	1:A:227:GLY:O	2.35	0.45
1:C:132:THR:O	1:C:154:VAL:N	2.48	0.45
1:C:274:THR:OG1	1:C:278:THR:O	2.33	0.45
1:C:8:GLY:O	1:C:32:THR:OG1	2.35	0.45
1:D:109:ALA:N	3:D:401:GOL:HO3	2.15	0.45
1:C:223:ASN:OD1	1:C:226:ARG:NH2	2.50	0.45
1:A:231:TRP:CD1	1:A:231:TRP:O	2.70	0.44
1:B:274:THR:N	5:B:541:HOH:O	2.51	0.44
1:D:8:GLY:O	1:D:32:THR:OG1	2.36	0.43
1:D:244:LYS:N	1:D:286:GLN:O	2.51	0.43
1:A:160:GLU:OE1	1:A:164:SER:OG	2.37	0.43
1:A:244:LYS:O	1:A:286:GLN:N	2.52	0.43
1:D:249:GLU:N	1:D:281:LYS:O	2.52	0.42
1:C:209:GLN:NE2	5:C:521:HOH:O	2.52	0.42
1:B:69:LYS:NZ	1:B:73:GLU:OE1	2.53	0.42
1:C:21:ILE:O	5:C:509:HOH:O	2.21	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:236:THR:O	1:D:243:VAL:N	2.53	0.41
1:D:281:LYS:NZ	5:D:602:HOH:O	2.52	0.41
1:A:290:LYS:NZ	5:A:636:HOH:O	2.54	0.41
1:C:115:LEU:N	1:C:200:THR:OG1	2.52	0.41
1:C:286:GLN:NE2	1:C:290:LYS:O	2.54	0.41
1:A:108:HIS:N	1:A:135:THR:O	2.53	0.40
3:A:402:GOL:H31	5:A:624:HOH:O	2.21	0.40
1:D:244:LYS:O	1:D:286:GLN:N	2.55	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	313/317 (99%)	302 (96%)	9 (3%)	2 (1%)	33	46
1	B	313/317 (99%)	300 (96%)	11 (4%)	2 (1%)	33	46
1	C	314/317 (99%)	299 (95%)	13 (4%)	2 (1%)	33	46
1	D	316/317 (100%)	300 (95%)	14 (4%)	2 (1%)	33	46
All	All	1256/1268 (99%)	1201 (96%)	47 (4%)	8 (1%)	33	46

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	233	VAL
1	A	233	VAL
1	D	9	THR
1	D	233	VAL
1	A	9	THR
1	B	9	THR
1	B	233	VAL
1	C	9	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	269/264 (102%)	260 (97%)	9 (3%)	50	70
1	B	269/264 (102%)	253 (94%)	16 (6%)	28	39
1	C	270/264 (102%)	249 (92%)	21 (8%)	18	25
1	D	272/264 (103%)	254 (93%)	18 (7%)	24	33
All	All	1080/1056 (102%)	1016 (94%)	64 (6%)	28	39

All (64) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	33	GLN
1	A	36	ARG
1	A	64	LEU
1	A	137	MSE
1	A	166	PHE
1	A	215	THR
1	A	232	PRO
1	A	256	SER
1	A	267	GLU
1	B	33	GLN
1	B	36	ARG
1	B	45	THR
1	B	91	ILE
1	B	130	GLU
1	B	137	MSE
1	B	155	GLU
1	B	166	PHE
1	B	207	ARG
1	B	231[A]	TRP
1	B	231[B]	TRP
1	B	238	LEU
1	B	255	LYS
1	B	292	ARG
1	B	303	LYS
1	B	312	GLU
1	C	30	VAL

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Mol	Chain	Res	Type
1	C	36	ARG
1	C	40	ARG
1	C	42	LYS
1	C	43	VAL
1	C	44	LEU
1	C	70	ASP
1	C	91	ILE
1	C	101	LYS
1	C	110	SER
1	C	130	GLU
1	C	137	MSE
1	C	155	GLU
1	C	166	PHE
1	C	207	ARG
1	C	236	THR
1	C	277	GLU
1	C	292	ARG
1	C	296	SER
1	C	302	THR
1	C	303	LYS
1	D	-1	ASN
1	D	11	ASP
1	D	23	ASP
1	D	32	THR
1	D	36	ARG
1	D	45	THR
1	D	72	TYR
1	D	101	LYS
1	D	137	MSE
1	D	155	GLU
1	D	166	PHE
1	D	175	HIS
1	D	197	GLU
1	D	207	ARG
1	D	288	SER
1	D	293	MSE
1	D	305	GLU
1	D	312	GLU

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

### 5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

9 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	PG5	A	401	-	9,9,11	0.60	0	8,8,10	1.52	1 (12%)
3	GOL	A	402	-	5,5,5	0.35	0	5,5,5	0.40	0
4	SO4	A	403	-	4,4,4	0.21	0	6,6,6	0.07	0
3	GOL	B	401	-	5,5,5	0.35	0	5,5,5	0.23	0
4	SO4	B	402	-	4,4,4	0.21	0	6,6,6	0.08	0
3	GOL	C	401	-	5,5,5	0.37	0	5,5,5	0.28	0
3	GOL	D	401	-	5,5,5	0.38	0	5,5,5	0.20	0
4	SO4	D	402	-	4,4,4	0.22	0	6,6,6	0.06	0
4	SO4	D	403	-	4,4,4	0.20	0	6,6,6	0.07	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	PG5	A	401	-	-	0/7/7/9	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GOL	A	402	-	-	0/4/4/4	0/0/0/0
4	SO4	A	403	-	-	0/0/0/0	0/0/0/0
3	GOL	B	401	-	-	0/4/4/4	0/0/0/0
4	SO4	B	402	-	-	0/0/0/0	0/0/0/0
3	GOL	C	401	-	-	0/4/4/4	0/0/0/0
3	GOL	D	401	-	-	0/4/4/4	0/0/0/0
4	SO4	D	402	-	-	0/0/0/0	0/0/0/0
4	SO4	D	403	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	PG5	O3-C6-C7	2.02	118.56	109.44

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 5.7 Other polymers ⓘ

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	314/317 (99%)	0.80	19 (6%) 21 22	29, 48, 84, 118	0
1	B	314/317 (99%)	1.03	40 (12%) 4 5	36, 59, 106, 145	0
1	C	315/317 (99%)	0.87	29 (9%) 9 10	35, 56, 97, 125	1 (0%)
1	D	315/317 (99%)	0.91	25 (7%) 13 14	39, 54, 78, 103	1 (0%)
All	All	1258/1268 (99%)	0.90	113 (8%) 10 10	29, 54, 92, 145	2 (0%)

All (113) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	303	LYS	7.2
1	B	40	ARG	6.8
1	B	36	ARG	6.4
1	A	36	ARG	6.4
1	C	37	PRO	5.6
1	A	43	VAL	5.4
1	B	39	GLY	5.4
1	B	42	LYS	4.8
1	B	43	VAL	4.6
1	B	38	VAL	4.4
1	B	275	GLY	4.0
1	B	238	LEU	4.0
1	C	41	LYS	3.9
1	B	148	ILE	3.8
1	C	242	VAL	3.8
1	D	231[A]	TRP	3.7
1	C	255	LYS	3.7
1	B	264	ALA	3.6
1	D	-2	SER	3.6
1	C	302	THR	3.6
1	D	45	THR	3.6

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Mol	Chain	Res	Type	RSRZ
1	D	255	LYS	3.5
1	B	189	LEU	3.5
1	A	302	THR	3.4
1	C	102	TYR	3.4
1	C	184	LEU	3.4
1	C	119	ALA	3.4
1	C	292	ARG	3.4
1	B	251	VAL	3.4
1	A	189	LEU	3.4
1	A	239	ALA	3.2
1	B	53	ALA	3.2
1	C	188	LYS	3.0
1	B	37	PRO	3.0
1	D	37	PRO	3.0
1	D	80	PRO	3.0
1	A	304	PRO	2.9
1	D	89	GLY	2.9
1	B	44	LEU	2.9
1	D	312	GLU	2.9
1	B	231[A]	TRP	2.9
1	B	187	GLY	2.9
1	C	231[A]	TRP	2.9
1	D	40	ARG	2.9
1	B	203	TYR	2.9
1	A	39	GLY	2.8
1	D	44	LEU	2.8
1	B	41	LYS	2.8
1	C	38	VAL	2.7
1	D	15	PRO	2.7
1	D	184	LEU	2.7
1	B	299	LEU	2.7
1	D	166	PHE	2.7
1	B	261	THR	2.7
1	B	12	PHE	2.7
1	D	242	VAL	2.6
1	C	191	PRO	2.6
1	D	192	ILE	2.6
1	B	185	ILE	2.5
1	A	201	PHE	2.5
1	B	91	ILE	2.5
1	C	124	ALA	2.5
1	D	124	ALA	2.5

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Mol	Chain	Res	Type	RSRZ
1	C	265	ILE	2.4
1	B	85	THR	2.4
1	C	290	LYS	2.4
1	C	235	TYR	2.4
1	B	176	LEU	2.4
1	D	253	VAL	2.4
1	A	125	ILE	2.4
1	A	169	LEU	2.4
1	C	270	PHE	2.4
1	A	44	LEU	2.4
1	A	264	ALA	2.4
1	B	28	ILE	2.4
1	C	27	VAL	2.4
1	A	212	ILE	2.4
1	D	298	PHE	2.3
1	C	200	THR	2.3
1	B	183	LEU	2.3
1	C	100	PRO	2.3
1	C	306	ILE	2.3
1	A	295	CYS	2.2
1	C	10	PRO	2.2
1	D	306	ILE	2.2
1	B	197	GLU	2.2
1	D	205	ILE	2.2
1	D	29	GLY	2.2
1	A	42	LYS	2.2
1	B	309	LYS	2.2
1	B	154	VAL	2.2
1	B	125	ILE	2.2
1	B	137	MSE	2.2
1	B	201	PHE	2.1
1	C	162	THR	2.1
1	C	299	LEU	2.1
1	C	42	LYS	2.1
1	C	1	MSE	2.1
1	B	280	VAL	2.1
1	A	303	LYS	2.1
1	B	111	LEU	2.1
1	D	128	GLY	2.1
1	B	212	ILE	2.1
1	A	289	GLY	2.1
1	C	66	ILE	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	44	LEU	2.1
1	B	25	TYR	2.1
1	D	53	ALA	2.1
1	B	244	LYS	2.0
1	D	262	ILE	2.0
1	A	177	LEU	2.0
1	D	271	VAL	2.0
1	A	32	THR	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

## 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands ⓘ

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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
4	SO4	B	402	5/5	0.21	3.78	70,72,77,84	0
3	GOL	A	402	6/6	0.22	3.08	36,46,52,58	0
3	GOL	D	401	6/6	0.20	0.42	35,37,49,58	0
3	GOL	C	401	6/6	0.20	0.35	39,45,47,49	0
3	GOL	B	401	6/6	0.13	-1.38	39,44,44,47	0
4	SO4	D	402	5/5	0.15	-1.38	94,100,102,103	0
2	PG5	A	401	10/12	0.11	-2.22	46,50,55,57	0
4	SO4	D	403	5/5	0.13	-2.40	85,89,92,96	0
4	SO4	A	403	5/5	0.12	-3.26	81,81,92,96	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.