



# Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 04:19 am BST

PDB ID : 3OOW  
Title : Octameric structure of the phosphoribosylaminoimidazole carboxylase catalytic subunit from *Francisella tularensis* subsp. *tularensis* SCHU S4.  
Authors : Filippova, E.V.; Wawrzak, Z.; Kudritska, M.; Edwards, A.; Savchenko, A.; Anderson, F.W.; Center for Structural Genomics of Infectious Diseases (CS-GID)  
Deposited on : 2010-08-31  
Resolution : 1.75 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

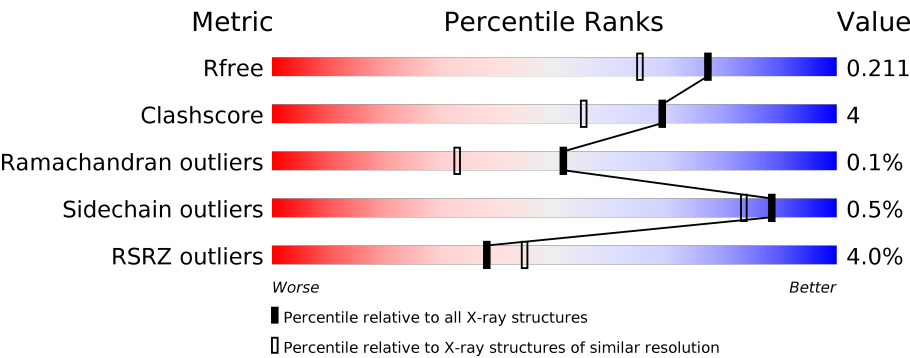
# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.75 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	2340 (1.76-1.76)
Clashscore	141614	2466 (1.76-1.76)
Ramachandran outliers	138981	2437 (1.76-1.76)
Sidechain outliers	138945	2437 (1.76-1.76)
RSRZ outliers	127900	2298 (1.76-1.76)

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	166	<div><div>4%</div><div><div></div><div>88%</div><div>8%</div><div></div></div><div></div></div>
1	B	166	<div><div>4%</div><div><div></div><div>89%</div><div>9%</div><div></div></div><div></div></div>
1	C	166	<div><div>4%</div><div><div></div><div>89%</div><div>8%</div><div></div></div><div></div></div>
1	D	166	<div><div>2%</div><div><div></div><div>89%</div><div>9%</div><div></div></div><div></div></div>
1	E	166	<div><div>4%</div><div><div></div><div>89%</div><div>8%</div><div></div></div><div></div></div>
1	F	166	<div><div>5%</div><div><div></div><div>84%</div><div>13%</div><div></div></div><div></div></div>

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Mol	Chain	Length	Quality of chain
1	G	166	 2% 90% 5% . .
1	H	166	 5% 89% 7% . .

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
5	MPD	C	166	X	-	-	-
5	MPD	D	166	X	-	-	-
5	MPD	F	166	X	-	-	-
5	MPD	G	165	X	-	-	-

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 10270 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 Phosphoribosylaminoimidazole carboxylase,catalytic subunit.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	160	Total	C	N	O	S	Se	0	3	0
			1196	753	204	228	4	7			
1	B	162	Total	C	N	O	S	Se	0	1	0
			1199	756	204	228	3	8			
1	C	161	Total	C	N	O	S	Se	0	2	0
			1197	754	204	228	4	7			
1	D	162	Total	C	N	O	S	Se	0	1	0
			1201	757	206	228	3	7			
1	E	160	Total	C	N	O	S	Se	0	1	0
			1182	746	202	224	3	7			
1	F	162	Total	C	N	O	S	Se	0	2	0
			1212	763	210	229	3	7			
1	G	159	Total	C	N	O	S	Se	0	0	0
			1163	735	197	222	3	6			
1	H	161	Total	C	N	O	S	Se	0	2	0
			1198	755	204	229	3	7			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	EXPRESSION TAG	UNP Q5NGE9
A	-1	ASN	-	EXPRESSION TAG	UNP Q5NGE9
A	0	ALA	-	EXPRESSION TAG	UNP Q5NGE9
B	-2	SER	-	EXPRESSION TAG	UNP Q5NGE9
B	-1	ASN	-	EXPRESSION TAG	UNP Q5NGE9
B	0	ALA	-	EXPRESSION TAG	UNP Q5NGE9
C	-2	SER	-	EXPRESSION TAG	UNP Q5NGE9
C	-1	ASN	-	EXPRESSION TAG	UNP Q5NGE9
C	0	ALA	-	EXPRESSION TAG	UNP Q5NGE9
D	-2	SER	-	EXPRESSION TAG	UNP Q5NGE9
D	-1	ASN	-	EXPRESSION TAG	UNP Q5NGE9
D	0	ALA	-	EXPRESSION TAG	UNP Q5NGE9
E	-2	SER	-	EXPRESSION TAG	UNP Q5NGE9

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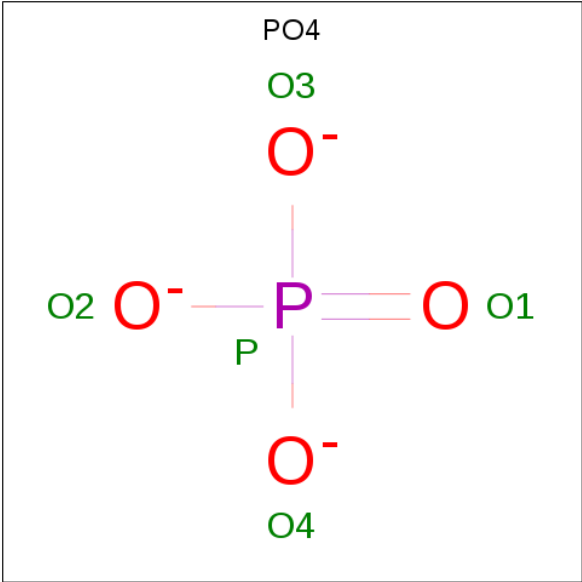
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Chain	Residue	Modelled	Actual	Comment	Reference
E	-1	ASN	-	EXPRESSION TAG	UNP Q5NGE9
E	0	ALA	-	EXPRESSION TAG	UNP Q5NGE9
F	-2	SER	-	EXPRESSION TAG	UNP Q5NGE9
F	-1	ASN	-	EXPRESSION TAG	UNP Q5NGE9
F	0	ALA	-	EXPRESSION TAG	UNP Q5NGE9
G	-2	SER	-	EXPRESSION TAG	UNP Q5NGE9
G	-1	ASN	-	EXPRESSION TAG	UNP Q5NGE9
G	0	ALA	-	EXPRESSION TAG	UNP Q5NGE9
H	-2	SER	-	EXPRESSION TAG	UNP Q5NGE9
H	-1	ASN	-	EXPRESSION TAG	UNP Q5NGE9
H	0	ALA	-	EXPRESSION TAG	UNP Q5NGE9

- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	G	1	Total Cl 1 1	0	0
2	D	2	Total Cl 2 2	0	0
2	E	2	Total Cl 2 2	0	0
2	H	2	Total Cl 2 2	0	0
2	B	1	Total Cl 1 1	0	0
2	C	2	Total Cl 2 2	0	0
2	A	2	Total Cl 2 2	0	0
2	F	2	Total Cl 2 2	0	0

- Molecule 3 is PHOSPHATE ION (three-letter code: PO4) (formula: O<sub>4</sub>P).



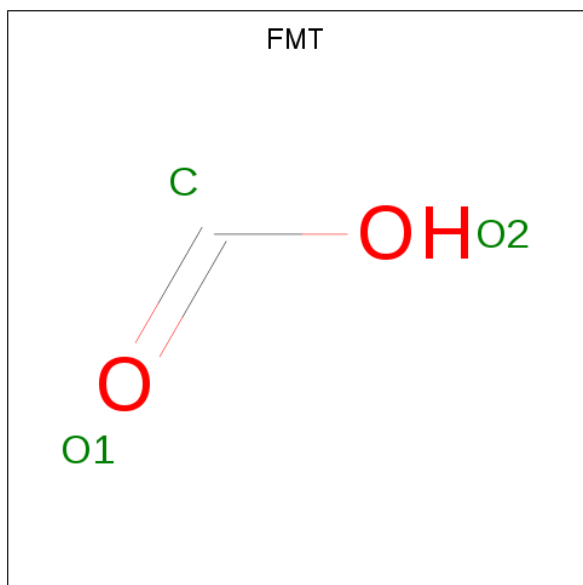
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	P	0	0
			5	4	1		
3	A	1	Total	O	P	0	0
			5	4	1		
3	A	1	Total	O	P	0	0
			5	4	1		
3	B	1	Total	O	P	0	0
			5	4	1		
3	B	1	Total	O	P	0	0
			5	4	1		
3	C	1	Total	O	P	0	0
			5	4	1		
3	C	1	Total	O	P	0	0
			5	4	1		
3	D	1	Total	O	P	0	0
			5	4	1		
3	D	1	Total	O	P	0	0
			5	4	1		
3	E	1	Total	O	P	0	0
			5	4	1		
3	E	1	Total	O	P	0	0
			5	4	1		
3	F	1	Total	O	P	0	0
			5	4	1		
3	F	1	Total	O	P	0	0
			5	4	1		
3	G	1	Total	O	P	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	G	1	Total	O	P	0	0
			5	4	1		
3	H	1	Total	O	P	0	0
			5	4	1		
3	H	1	Total	O	P	0	0
			5	4	1		

- Molecule 4 is FORMIC ACID (three-letter code: FMT) (formula: CH<sub>2</sub>O<sub>2</sub>).



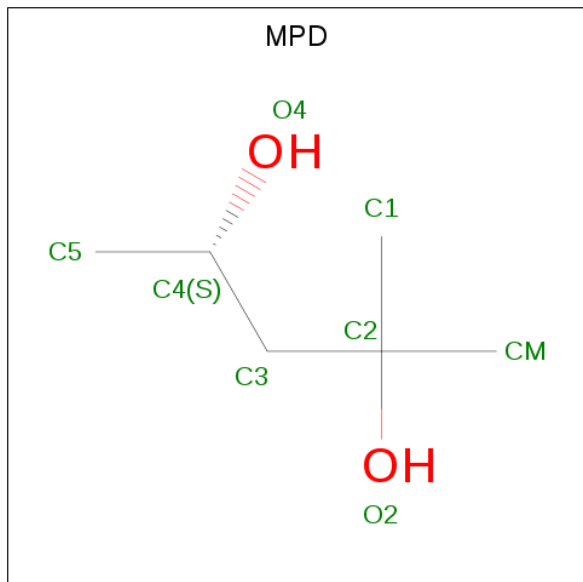
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			3	1	2		
4	A	1	Total	C	O	0	0
			3	1	2		
4	B	1	Total	C	O	0	0
			3	1	2		
4	B	1	Total	C	O	0	0
			3	1	2		
4	B	1	Total	C	O	0	0
			3	1	2		
4	C	1	Total	C	O	0	0
			3	1	2		
4	D	1	Total	C	O	0	0
			3	1	2		
4	E	1	Total	C	O	0	0
			3	1	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	E	1	Total	C	O	0	0
			3	1	2		
4	F	1	Total	C	O	0	0
			3	1	2		
4	G	1	Total	C	O	0	0
			3	1	2		
4	H	1	Total	C	O	0	0
			3	1	2		
4	H	1	Total	C	O	0	0
			3	1	2		

- Molecule 5 is (4S)-2-METHYL-2,4-PENTANEDIOL (three-letter code: MPD) (formula:  $C_6H_{14}O_2$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	C	1	Total	C	O	0	0
			8	6	2		
5	D	1	Total	C	O	0	0
			8	6	2		
5	F	1	Total	C	O	0	0
			8	6	2		
5	G	1	Total	C	O	0	0
			8	6	2		

- Molecule 6 is water.

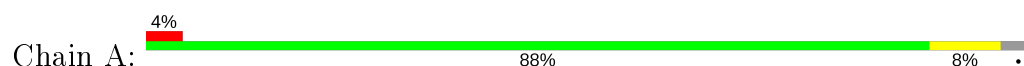


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	60	Total O 61 61	0	1
6	B	64	Total O 65 65	0	1
6	C	66	Total O 69 69	0	3
6	D	83	Total O 83 83	0	0
6	E	60	Total O 60 60	0	0
6	F	65	Total O 65 65	0	0
6	G	79	Total O 80 80	0	1
6	H	68	Total O 69 69	0	1

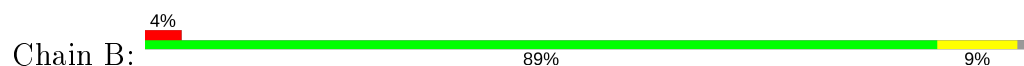
### 3 Residue-property plots [i](#)

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.

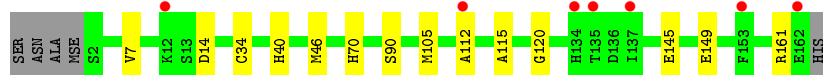
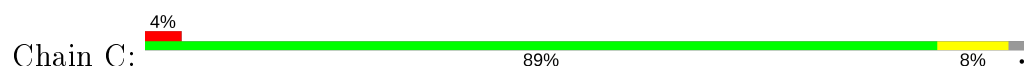
- Molecule 1: Phosphoribosylaminoimidazole carboxylase,catalytic subunit



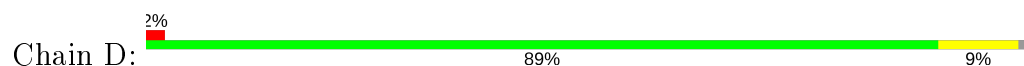
- Molecule 1: Phosphoribosylaminoimidazole carboxylase,catalytic subunit



- Molecule 1: Phosphoribosylaminoimidazole carboxylase,catalytic subunit



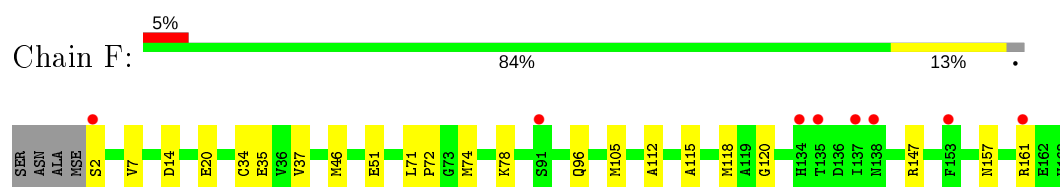
- Molecule 1: Phosphoribosylaminoimidazole carboxylase,catalytic subunit



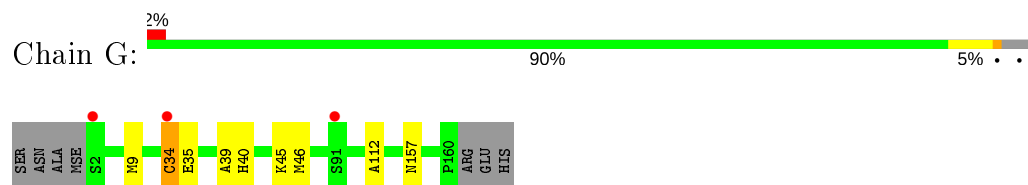
- Molecule 1: Phosphoribosylaminoimidazole carboxylase,catalytic subunit



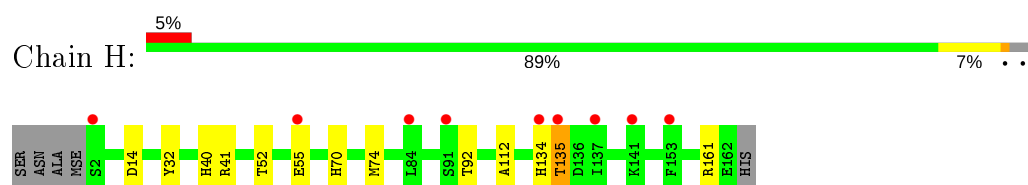
- Molecule 1: Phosphoribosylaminoimidazole carboxylase,catalytic subunit



- Molecule 1: Phosphoribosylaminoimidazole carboxylase,catalytic subunit



- Molecule 1: Phosphoribosylaminoimidazole carboxylase,catalytic subunit



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	88.56 Å 96.30 Å 128.62 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	77.09 – 1.75 29.38 – 1.75	Depositor EDS
% Data completeness (in resolution range)	29.4 (77.09-1.75) 98.7 (29.38-1.75)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.27 (at 1.75 Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, $R_{free}$	0.168 , 0.205 0.176 , 0.211	Depositor DCC
$R_{free}$ test set	5493 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	22.5	Xtriage
Anisotropy	0.026	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.39 , 45.3	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	0.97	EDS
Total number of atoms	10270	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	26.0	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 62.27 % 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 1.1548e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<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: PO4, MPD, FMT, 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	0.75	0/1209	0.74	1/1628 (0.1%)
1	B	0.78	0/1212	0.74	0/1631
1	C	0.74	0/1210	0.72	0/1629
1	D	0.79	1/1215 (0.1%)	0.76	1/1636 (0.1%)
1	E	0.77	0/1195	0.74	2/1609 (0.1%)
1	F	0.71	0/1226	0.71	1/1650 (0.1%)
1	G	0.77	1/1176 (0.1%)	0.74	0/1585
1	H	0.75	0/1211	0.79	1/1631 (0.1%)
All	All	0.76	2/9654 (0.0%)	0.74	6/12999 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	51	GLU	CB-CG	-5.82	1.41	1.52
1	G	34	CYS	CB-SG	-5.24	1.73	1.81

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	97	ASP	CB-CG-OD1	5.81	123.53	118.30
1	H	41	ARG	NE-CZ-NH2	-5.63	117.48	120.30
1	E	152	ARG	NE-CZ-NH1	5.51	123.06	120.30
1	F	147	ARG	NE-CZ-NH1	5.43	123.01	120.30
1	E	147	ARG	NE-CZ-NH1	5.40	123.00	120.30
1	D	41	ARG	NE-CZ-NH2	-5.08	117.76	120.30

There are no chirality outliers.

There are no planarity outliers.

## 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	1196	0	1211	12	0
1	B	1199	0	1222	11	0
1	C	1197	0	1214	11	0
1	D	1201	0	1217	10	0
1	E	1182	0	1204	8	0
1	F	1212	0	1229	16	0
1	G	1163	0	1183	10	0
1	H	1198	0	1216	14	0
2	A	2	0	0	0	0
2	B	1	0	0	0	0
2	C	2	0	0	1	0
2	D	2	0	0	0	0
2	E	2	0	0	0	0
2	F	2	0	0	0	0
2	G	1	0	0	0	0
2	H	2	0	0	0	0
3	A	15	0	0	0	0
3	B	10	0	0	0	0
3	C	10	0	0	2	0
3	D	10	0	0	0	0
3	E	10	0	0	1	0
3	F	10	0	0	1	0
3	G	10	0	0	1	0
3	H	10	0	0	1	0
4	A	6	0	2	0	0
4	B	9	0	3	0	0
4	C	3	0	1	0	0
4	D	3	0	1	0	0
4	E	6	0	2	0	0
4	F	3	0	1	0	0
4	G	3	0	1	0	0
4	H	6	0	2	0	0
5	C	8	0	14	1	0
5	D	8	0	14	0	0
5	F	8	0	14	1	0
5	G	8	0	14	0	0
6	A	61	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	B	65	0	0	1	0
6	C	69	0	0	2	0
6	D	83	0	0	3	0
6	E	60	0	0	0	0
6	F	65	0	0	1	0
6	G	80	0	0	0	0
6	H	69	0	0	2	0
All	All	10270	0	9765	72	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 4.

All (72) 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:H:92[A]:THR:HG22	6:H:412:HOH:O	1.76	0.86
1:F:46[A]:MSE:HE1	1:F:74:MSE:HB3	1.55	0.85
6:D:189:HOH:O	1:H:92[A]:THR:HG21	1.77	0.84
1:G:34:CYS:SG	1:H:161:ARG:CZ	2.68	0.81
1:A:34[B]:CYS:SG	1:F:161:ARG:NE	2.66	0.68
1:A:135:THR:HG21	1:H:134:HIS:CE1	2.28	0.68
1:D:34:CYS:SG	1:E:161:ARG:NH1	2.68	0.66
1:D:90:SER:HB2	6:D:183:HOH:O	1.98	0.64
1:A:112:ALA:HB1	1:H:112:ALA:HB1	1.79	0.63
2:C:165:CL:CL	6:C:301:HOH:O	2.52	0.63
1:B:34:CYS:SG	1:C:161:ARG:NH2	2.72	0.62
1:G:34:CYS:SG	1:H:161:ARG:NE	2.73	0.61
1:F:112:ALA:HB1	1:G:112:ALA:HB1	1.86	0.57
1:D:24:ILE:HD12	1:D:122:LYS:HA	1.87	0.57
1:F:20:GLU:HG3	1:F:118:MSE:HG3	1.87	0.56
1:A:135:THR:HG21	1:H:134:HIS:HE1	1.72	0.55
1:F:2:SER:N	6:F:439:HOH:O	2.41	0.54
1:B:112:ALA:HB1	1:E:112:ALA:HB1	1.90	0.54
5:C:166:MPD:H52	5:C:166:MPD:HM1	1.89	0.54
1:F:14:ASP:OD2	3:F:168:PO4:O4	2.26	0.54
1:D:71:LEU:HB3	1:D:72:PRO:HD3	1.92	0.51
1:G:39:ALA:HA	1:G:46:MSE:SE	2.59	0.51
1:G:35:GLU:OE2	1:G:45:LYS:NZ	2.42	0.49
1:D:7:VAL:O	1:D:34:CYS:HA	2.14	0.47
1:E:40:HIS:CD2	1:F:105:MSE:HE1	2.49	0.47
1:C:14:ASP:OD2	3:C:168:PO4:O2	2.33	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:14:ASP:OD2	3:E:166:PO4:O1	2.33	0.47
1:B:24:ILE:HD11	1:B:122:LYS:HB3	1.97	0.46
1:H:14:ASP:OD2	3:H:166:PO4:O4	2.32	0.46
1:G:40:HIS:ND1	3:G:167:PO4:O4	2.48	0.46
1:F:7:VAL:O	1:F:34:CYS:HA	2.14	0.46
1:A:52:THR:O	1:A:55:GLU:HG2	2.15	0.46
1:A:74:MSE:CE	1:G:46:MSE:HE1	2.46	0.46
1:C:145:GLU:O	1:C:149:GLU:HG3	2.16	0.46
1:A:2:SER:N	6:A:538:HOH:O	2.50	0.45
1:B:161:ARG:NH2	6:B:293:HOH:O	2.49	0.45
1:F:115:ALA:O	1:F:120:GLY:HA3	2.16	0.45
1:H:161:ARG:NH1	6:H:487:HOH:O	2.42	0.45
1:C:112:ALA:HB1	1:D:112:ALA:HB1	1.98	0.45
1:B:110:PRO:HB3	1:E:123:ASN:HB3	1.99	0.44
1:C:115:ALA:O	1:C:120:GLY:HA3	2.18	0.44
1:D:56:ARG:HD2	6:D:195:HOH:O	2.16	0.44
1:E:7:VAL:O	1:E:34:CYS:HA	2.17	0.44
1:B:115:ALA:O	1:B:120:GLY:HA3	2.18	0.43
1:E:75:VAL:O	1:E:79:THR:HG22	2.19	0.43
1:A:74:MSE:HE2	1:G:46:MSE:HE1	2.01	0.43
1:B:24:ILE:CD1	1:B:122:LYS:HB3	2.49	0.43
1:F:35:GLU:HG2	1:F:37:VAL:HG23	1.99	0.43
1:F:71:LEU:HB3	1:F:72:PRO:HD3	2.01	0.42
1:F:96:GLN:NE2	5:F:166:MPD:H53	2.34	0.42
1:F:51:GLU:HG3	1:F:78:LYS:HE3	2.00	0.42
1:D:161:ARG:NH2	1:H:32:TYR:OH	2.52	0.42
1:D:118:MSE:HB2	1:D:118:MSE:HE3	1.81	0.42
1:A:12:LYS:HE3	1:F:157:ASN:HB3	2.02	0.42
1:B:52:THR:O	1:B:55:GLU:HG2	2.20	0.42
1:C:105:MSE:HE1	1:H:40:HIS:CD2	2.54	0.42
1:A:12:LYS:HE3	1:F:157:ASN:CB	2.49	0.42
1:A:71:LEU:HB3	1:A:72:PRO:HD3	2.03	0.41
1:B:71:LEU:HB3	1:B:72:PRO:HD3	2.01	0.41
1:F:46[A]:MSE:HE1	1:F:74:MSE:CB	2.38	0.41
1:C:70:HIS:CD2	1:H:70:HIS:CD2	3.08	0.41
1:H:52:THR:O	1:H:55:GLU:HG2	2.21	0.41
1:A:46[A]:MSE:SE	1:A:71:LEU:HD12	2.70	0.41
1:D:115:ALA:O	1:D:120:GLY:HA3	2.21	0.41
1:B:7:VAL:O	1:B:34:CYS:HA	2.21	0.41
1:E:12:LYS:HE3	1:G:157:ASN:HB3	2.01	0.41
1:C:40:HIS:ND1	3:C:167:PO4:O3	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:118:MSE:HE3	1:B:118:MSE:HB2	1.90	0.41
1:C:46[B]:MSE:HE1	1:H:74:MSE:CE	2.51	0.40
1:G:9:MSE:HE2	1:G:9:MSE:HB3	2.02	0.40
1:C:7:VAL:O	1:C:34[A]:CYS:HA	2.21	0.40
1:C:90:SER:HB2	6:C:191[B]:HOH:O	2.20	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	161/166 (97%)	159 (99%)	2 (1%)	0	100	100
1	B	161/166 (97%)	158 (98%)	3 (2%)	0	100	100
1	C	161/166 (97%)	160 (99%)	1 (1%)	0	100	100
1	D	161/166 (97%)	160 (99%)	1 (1%)	0	100	100
1	E	159/166 (96%)	157 (99%)	2 (1%)	0	100	100
1	F	162/166 (98%)	159 (98%)	3 (2%)	0	100	100
1	G	157/166 (95%)	156 (99%)	1 (1%)	0	100	100
1	H	161/166 (97%)	158 (98%)	2 (1%)	1 (1%)	25	10
All	All	1283/1328 (97%)	1267 (99%)	15 (1%)	1 (0%)	51	33

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	H	135	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	127/122 (104%)	124 (98%)	3 (2%)	49	26
1	B	127/122 (104%)	126 (99%)	1 (1%)	81	72
1	C	127/122 (104%)	127 (100%)	0	100	100
1	D	127/122 (104%)	127 (100%)	0	100	100
1	E	125/122 (102%)	124 (99%)	1 (1%)	81	72
1	F	128/122 (105%)	128 (100%)	0	100	100
1	G	123/122 (101%)	123 (100%)	0	100	100
1	H	127/122 (104%)	126 (99%)	1 (1%)	81	72
All	All	1011/976 (104%)	1005 (99%)	6 (1%)	88	79

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	48[A]	ASP
1	A	48[B]	ASP
1	A	161	ARG
1	B	1	MSE
1	E	137	ILE
1	H	135	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	27	ASN
1	B	27	ASN
1	B	133	GLN
1	C	27	ASN
1	D	27	ASN
1	F	27	ASN
1	G	27	ASN
1	G	96	GLN

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Mol	Chain	Res	Type
1	H	27	ASN
1	H	134	HIS

### 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 48 ligands modelled in this entry, 14 are monoatomic - leaving 34 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	PO4	F	167	-	4,4,4	0.93	0	6,6,6	0.44	0
4	FMT	B	168	-	0,2,2	0.00	-	0,1,1	0.00	-
4	FMT	E	168	-	0,2,2	0.00	-	0,1,1	0.00	-
3	PO4	D	167	-	4,4,4	0.83	0	6,6,6	0.97	0
3	PO4	C	167	-	4,4,4	0.97	0	6,6,6	0.50	0
4	FMT	C	169	-	0,2,2	0.00	-	0,1,1	0.00	-
3	PO4	B	165	-	4,4,4	0.72	0	6,6,6	0.48	0
4	FMT	H	168	-	0,2,2	0.00	-	0,1,1	0.00	-
3	PO4	A	166	-	4,4,4	0.65	0	6,6,6	1.06	0
3	PO4	G	166	-	4,4,4	1.13	0	6,6,6	0.50	0
4	FMT	B	169	-	0,2,2	0.00	-	0,1,1	0.00	-
4	FMT	A	170	-	0,2,2	0.00	-	0,1,1	0.00	-

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	PO4	F	168	-	4,4,4	1.03	0	6,6,6	1.06	0
3	PO4	C	168	-	4,4,4	1.25	0	6,6,6	0.36	0
4	FMT	E	169	-	0,2,2	0.00	-	0,1,1	0.00	-
5	MPD	D	166	-	7,7,7	0.38	0	9,10,10	0.50	0
5	MPD	F	166	-	7,7,7	0.33	0	9,10,10	0.71	0
4	FMT	H	169	-	0,2,2	0.00	-	0,1,1	0.00	-
3	PO4	E	167	-	4,4,4	0.87	0	6,6,6	0.89	0
3	PO4	G	167	-	4,4,4	0.90	0	6,6,6	1.08	0
4	FMT	B	167	-	0,2,2	0.00	-	0,1,1	0.00	-
3	PO4	A	168	-	4,4,4	0.72	0	6,6,6	0.38	0
5	MPD	G	165	-	7,7,7	0.32	0	9,10,10	0.38	0
4	FMT	A	169	-	0,2,2	0.00	-	0,1,1	0.00	-
3	PO4	H	167	-	4,4,4	0.90	0	6,6,6	0.80	0
5	MPD	C	166	-	7,7,7	0.28	0	9,10,10	0.76	0
4	FMT	G	168	-	0,2,2	0.00	-	0,1,1	0.00	-
4	FMT	D	169	-	0,2,2	0.00	-	0,1,1	0.00	-
3	PO4	E	166	-	4,4,4	1.05	0	6,6,6	0.90	0
3	PO4	A	167	-	4,4,4	1.05	0	6,6,6	0.47	0
3	PO4	H	166	-	4,4,4	1.05	0	6,6,6	1.00	0
3	PO4	D	168	-	4,4,4	1.30	0	6,6,6	0.96	0
4	FMT	F	169	-	0,2,2	0.00	-	0,1,1	0.00	-
3	PO4	B	166	-	4,4,4	0.77	0	6,6,6	0.85	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
5	MPD	F	166	-	1/1/2/2	1/5/5/5	-
5	MPD	C	166	-	1/1/2/2	2/5/5/5	-
5	MPD	D	166	-	1/1/2/2	0/5/5/5	-
5	MPD	G	165	-	1/1/2/2	0/5/5/5	-

There are no bond length outliers.

There are no bond angle outliers.

All (4) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
5	F	166	MPD	C4
5	D	166	MPD	C4

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Mol	Chain	Res	Type	Atom
5	G	165	MPD	C4
5	C	166	MPD	C4

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	C	166	MPD	C2-C3-C4-C5
5	F	166	MPD	C2-C3-C4-C5
5	C	166	MPD	C2-C3-C4-O4

There are no ring outliers.

8 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	167	PO4	1	0
3	F	168	PO4	1	0
3	C	168	PO4	1	0
5	F	166	MPD	1	0
3	G	167	PO4	1	0
5	C	166	MPD	1	0
3	E	166	PO4	1	0
3	H	166	PO4	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	154/166 (92%)	0.27	6 (3%)	39	45	15, 25, 39, 54	0
1	B	155/166 (93%)	0.21	6 (3%)	39	45	16, 24, 36, 45	0
1	C	155/166 (93%)	0.21	7 (4%)	33	39	15, 24, 38, 55	0
1	D	156/166 (93%)	0.08	4 (2%)	56	62	15, 23, 34, 41	0
1	E	154/166 (92%)	0.16	6 (3%)	39	45	16, 24, 36, 45	0
1	F	156/166 (93%)	0.26	8 (5%)	28	34	16, 25, 43, 50	0
1	G	153/166 (92%)	0.06	3 (1%)	65	72	16, 24, 35, 42	0
1	H	155/166 (93%)	0.19	9 (5%)	23	28	15, 23, 38, 54	0
All	All	1238/1328 (93%)	0.18	49 (3%)	38	45	15, 24, 38, 55	0

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	137	ILE	5.9
1	A	2	SER	5.5
1	B	91	SER	5.4
1	C	162	GLU	4.7
1	C	137	ILE	4.6
1	H	2	SER	4.6
1	F	2	SER	4.5
1	G	91	SER	3.8
1	A	153	PHE	3.8
1	F	153	PHE	3.8
1	D	91	SER	3.6
1	A	91	SER	3.5
1	A	34[A]	CYS	3.5
1	H	137	ILE	3.4
1	H	134	HIS	3.3
1	F	91	SER	3.3

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Mol	Chain	Res	Type	RSRZ
1	B	2	SER	3.2
1	B	90	SER	3.1
1	F	135	THR	3.0
1	C	134	HIS	2.9
1	E	2	SER	2.9
1	B	162	GLU	2.7
1	G	34	CYS	2.7
1	H	84	LEU	2.6
1	F	134	HIS	2.6
1	H	55	GLU	2.4
1	E	34	CYS	2.3
1	F	161	ARG	2.3
1	C	135	THR	2.3
1	H	135	THR	2.3
1	G	2	SER	2.3
1	E	127	PHE	2.3
1	B	141	LYS	2.2
1	C	12	LYS	2.2
1	H	91	SER	2.2
1	F	138	ASN	2.2
1	E	55	GLU	2.2
1	E	141	LYS	2.1
1	C	112	ALA	2.1
1	D	90	SER	2.1
1	A	94	ASN	2.1
1	C	153	PHE	2.1
1	H	141	LYS	2.1
1	E	91	SER	2.1
1	D	137	ILE	2.1
1	D	92	THR	2.0
1	A	149	GLU	2.0
1	B	34	CYS	2.0
1	H	153	PHE	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 ⓘ

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
4	FMT	A	169	3/3	0.79	0.21	52,52,52,52	0
4	FMT	B	168	3/3	0.80	0.18	38,38,39,42	0
5	MPD	G	165	8/8	0.89	0.18	33,41,45,45	0
5	MPD	D	166	8/8	0.89	0.14	33,39,40,41	0
4	FMT	H	168	3/3	0.90	0.13	38,38,39,39	0
5	MPD	C	166	8/8	0.90	0.16	38,40,43,46	0
4	FMT	H	169	3/3	0.91	0.07	46,46,50,51	0
3	PO4	A	168	5/5	0.91	0.34	52,54,55,55	0
4	FMT	E	169	3/3	0.91	0.17	58,58,58,59	0
5	MPD	F	166	8/8	0.91	0.12	35,39,41,46	0
3	PO4	G	167	5/5	0.91	0.14	45,46,47,50	0
4	FMT	B	167	3/3	0.92	0.12	29,29,33,36	0
4	FMT	C	169	3/3	0.93	0.11	37,37,38,39	0
4	FMT	E	168	3/3	0.93	0.10	37,37,37,38	0
3	PO4	B	166	5/5	0.93	0.12	47,50,52,53	0
3	PO4	H	167	5/5	0.94	0.13	44,47,51,51	0
4	FMT	A	170	3/3	0.94	0.08	39,39,39,39	0
3	PO4	D	167	5/5	0.94	0.13	56,56,58,58	0
3	PO4	D	168	5/5	0.94	0.15	43,48,50,52	0
3	PO4	E	167	5/5	0.94	0.16	48,49,50,51	0
3	PO4	F	167	5/5	0.95	0.14	46,49,49,51	0
3	PO4	A	166	5/5	0.95	0.11	42,44,47,47	0
4	FMT	D	169	3/3	0.96	0.17	37,37,37,38	0
3	PO4	C	167	5/5	0.96	0.12	62,62,63,64	0
4	FMT	B	169	3/3	0.96	0.06	34,34,34,35	0
4	FMT	G	168	3/3	0.97	0.07	34,34,35,35	0
3	PO4	F	168	5/5	0.97	0.13	39,44,45,47	0
3	PO4	H	166	5/5	0.97	0.16	40,43,44,46	0
3	PO4	G	166	5/5	0.97	0.10	38,40,42,42	0
3	PO4	C	168	5/5	0.97	0.19	46,51,51,53	0
2	CL	F	165	1/1	0.98	0.17	35,35,35,35	0
3	PO4	E	166	5/5	0.98	0.12	40,43,46,46	0
3	PO4	B	165	5/5	0.98	0.22	34,35,37,37	0
4	FMT	F	169	3/3	0.98	0.07	39,39,40,41	0
3	PO4	A	167	5/5	0.98	0.14	37,38,42,43	0
2	CL	H	165	1/1	0.99	0.04	34,34,34,34	0
2	CL	C	165	1/1	0.99	0.07	34,34,34,34	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	CL	D	165	1/1	0.99	0.06	30,30,30,30	0
2	CL	E	165	1/1	0.99	0.05	40,40,40,40	0
2	CL	C	164	1/1	0.99	0.03	29,29,29,29	0
2	CL	H	164	1/1	0.99	0.05	28,28,28,28	0
2	CL	A	165	1/1	0.99	0.07	37,37,37,37	0
2	CL	F	164	1/1	0.99	0.03	29,29,29,29	0
2	CL	B	164	1/1	0.99	0.12	33,33,33,33	0
2	CL	A	164	1/1	0.99	0.03	29,29,29,29	0
2	CL	G	164	1/1	0.99	0.04	30,30,30,30	0
2	CL	E	164	1/1	0.99	0.02	29,29,29,29	0
2	CL	D	164	1/1	0.99	0.02	31,31,31,31	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.