W O R L D W I D E
wwPDB
P R O T E I N   D A T A   B A N K

wwPDB X-ray Structure Validation Summary Report (i)

Feb 18, 2024 – 05:38 PM EST

PDB ID	:	4FYM
Title	:	Crystal structure of Plasmodium falciparum orotate phosphoribosyltransferase
Authors	:	Rathod, P.K.; Kumar, S.
Deposited on	:	2012-07-05
Resolution	:	2.60 Å(reported)

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

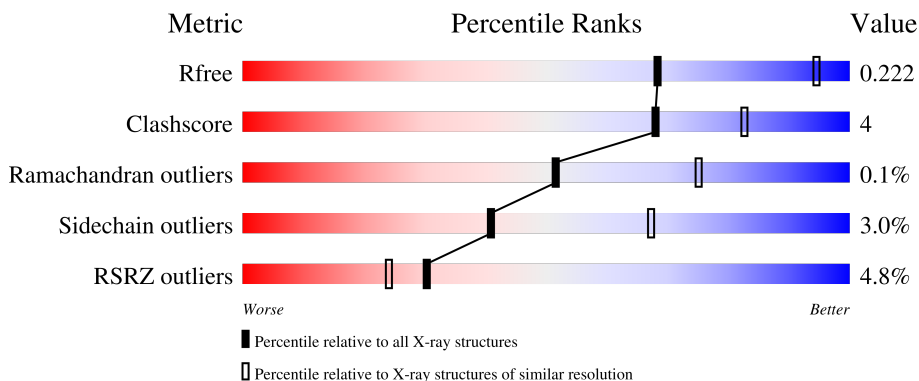|		
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	Percentile Ranks	Value
Rfree		0.222
Clashscore		4
Ramachandran outliers		0.1%
Sidechain outliers		3.0%
RSRZ outliers		4.8%

Worse   *Better*

▮ Percentile relative to all X-ray structures

▯ Percentile relative to X-ray structures of similar resolution

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

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 of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	266	4% — 76% 7% • 16%
1	B	266	2% — 73% 8% • 18%
1	C	266	4% — 75% 7% • 17%
1	D	266	3% — 76% 9% • 15%
1	E	266	5% — 71% 12% • 17%

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Mol	Chain	Length	Quality of chain
1	F	266	5% / 69% / 9% / 21%
1	G	266	2% / 74% / 7% / 19%
1	H	266	7% / 65% / 11% / • / 21%

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
2	SO4	B	305	-	-	X	-
2	SO4	E	301	-	-	X	-

2 Entry composition ⓘ

There are 3 unique types of molecules in this entry. The entry contains 14882 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 Orotate phosphoribosyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	223	Total	C	N	O	S	0	0	0
			1846	1214	288	337	7			
1	B	219	Total	C	N	O	S	0	0	0
			1811	1191	284	329	7			
1	C	222	Total	C	N	O	S	0	0	0
			1835	1210	290	328	7			
1	D	227	Total	C	N	O	S	0	0	0
			1882	1239	297	339	7			
1	E	222	Total	C	N	O	S	0	0	0
			1838	1213	290	328	7			
1	F	209	Total	C	N	O	S	0	0	0
			1725	1142	268	308	7			
1	G	216	Total	C	N	O	S	0	0	0
			1791	1180	280	324	7			
1	H	209	Total	C	N	O	S	0	0	0
			1725	1140	269	310	6			

There are 240 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	expression tag	UNP Q8N0R1
A	1	ALA	-	expression tag	UNP Q8N0R1
A	?	-	ASN	deletion	UNP Q8N0R1
A	?	-	ASN	deletion	UNP Q8N0R1
A	?	-	ASN	deletion	UNP Q8N0R1
A	?	-	ILE	deletion	UNP Q8N0R1
A	?	-	ASP	deletion	UNP Q8N0R1
A	?	-	ASN	deletion	UNP Q8N0R1
A	?	-	VAL	deletion	UNP Q8N0R1
A	?	-	ASN	deletion	UNP Q8N0R1
A	?	-	PHE	deletion	UNP Q8N0R1
A	?	-	ASN	deletion	UNP Q8N0R1
A	?	-	ASP	deletion	UNP Q8N0R1

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Chain	Residue	Modelled	Actual	Comment	Reference
A	?	-	ASP	deletion	UNP Q8N0R1
A	?	-	ASP	deletion	UNP Q8N0R1
A	?	-	ASP	deletion	UNP Q8N0R1
A	?	-	ASN	deletion	UNP Q8N0R1
A	?	-	ASN	deletion	UNP Q8N0R1
A	?	-	TYR	deletion	UNP Q8N0R1
A	?	-	ASP	deletion	UNP Q8N0R1
A	?	-	ASP	deletion	UNP Q8N0R1
A	?	-	ASP	deletion	UNP Q8N0R1
A	?	-	GLY	deletion	UNP Q8N0R1
A	?	-	ASN	deletion	UNP Q8N0R1
A	282	HIS	-	expression tag	UNP Q8N0R1
A	283	HIS	-	expression tag	UNP Q8N0R1
A	284	HIS	-	expression tag	UNP Q8N0R1
A	285	HIS	-	expression tag	UNP Q8N0R1
A	286	HIS	-	expression tag	UNP Q8N0R1
A	287	HIS	-	expression tag	UNP Q8N0R1
B	0	MET	-	expression tag	UNP Q8N0R1
B	1	ALA	-	expression tag	UNP Q8N0R1
B	?	-	ASN	deletion	UNP Q8N0R1
B	?	-	ASN	deletion	UNP Q8N0R1
B	?	-	ASN	deletion	UNP Q8N0R1
B	?	-	ILE	deletion	UNP Q8N0R1
B	?	-	ASP	deletion	UNP Q8N0R1
B	?	-	ASN	deletion	UNP Q8N0R1
B	?	-	VAL	deletion	UNP Q8N0R1
B	?	-	ASN	deletion	UNP Q8N0R1
B	?	-	PHE	deletion	UNP Q8N0R1
B	?	-	ASN	deletion	UNP Q8N0R1
B	?	-	ASP	deletion	UNP Q8N0R1
B	?	-	ASP	deletion	UNP Q8N0R1
B	?	-	ASP	deletion	UNP Q8N0R1
B	?	-	ASP	deletion	UNP Q8N0R1
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B	?	-	ASN	deletion	UNP Q8N0R1
B	?	-	TYR	deletion	UNP Q8N0R1
B	?	-	ASP	deletion	UNP Q8N0R1
B	?	-	ASP	deletion	UNP Q8N0R1
B	?	-	ASP	deletion	UNP Q8N0R1
B	?	-	GLY	deletion	UNP Q8N0R1
B	?	-	ASN	deletion	UNP Q8N0R1
B	282	HIS	-	expression tag	UNP Q8N0R1

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Chain	Residue	Modelled	Actual	Comment	Reference
B	283	HIS	-	expression tag	UNP Q8N0R1
B	284	HIS	-	expression tag	UNP Q8N0R1
B	285	HIS	-	expression tag	UNP Q8N0R1
B	286	HIS	-	expression tag	UNP Q8N0R1
B	287	HIS	-	expression tag	UNP Q8N0R1
C	0	MET	-	expression tag	UNP Q8N0R1
C	1	ALA	-	expression tag	UNP Q8N0R1
C	?	-	ASN	deletion	UNP Q8N0R1
C	?	-	ASN	deletion	UNP Q8N0R1
C	?	-	ASN	deletion	UNP Q8N0R1
C	?	-	ILE	deletion	UNP Q8N0R1
C	?	-	ASP	deletion	UNP Q8N0R1
C	?	-	ASN	deletion	UNP Q8N0R1
C	?	-	VAL	deletion	UNP Q8N0R1
C	?	-	ASN	deletion	UNP Q8N0R1
C	?	-	PHE	deletion	UNP Q8N0R1
C	?	-	ASN	deletion	UNP Q8N0R1
C	?	-	ASP	deletion	UNP Q8N0R1
C	?	-	ASP	deletion	UNP Q8N0R1
C	?	-	ASP	deletion	UNP Q8N0R1
C	?	-	ASP	deletion	UNP Q8N0R1
C	?	-	ASN	deletion	UNP Q8N0R1
C	?	-	ASN	deletion	UNP Q8N0R1
C	?	-	TYR	deletion	UNP Q8N0R1
C	?	-	ASP	deletion	UNP Q8N0R1
C	?	-	ASP	deletion	UNP Q8N0R1
C	?	-	ASP	deletion	UNP Q8N0R1
C	?	-	GLY	deletion	UNP Q8N0R1
C	?	-	ASN	deletion	UNP Q8N0R1
C	282	HIS	-	expression tag	UNP Q8N0R1
C	283	HIS	-	expression tag	UNP Q8N0R1
C	284	HIS	-	expression tag	UNP Q8N0R1
C	285	HIS	-	expression tag	UNP Q8N0R1
C	286	HIS	-	expression tag	UNP Q8N0R1
C	287	HIS	-	expression tag	UNP Q8N0R1
D	0	MET	-	expression tag	UNP Q8N0R1
D	1	ALA	-	expression tag	UNP Q8N0R1
D	?	-	ASN	deletion	UNP Q8N0R1
D	?	-	ASN	deletion	UNP Q8N0R1
D	?	-	ASN	deletion	UNP Q8N0R1
D	?	-	ILE	deletion	UNP Q8N0R1
D	?	-	ASP	deletion	UNP Q8N0R1

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Chain	Residue	Modelled	Actual	Comment	Reference
D	?	-	ASN	deletion	UNP Q8N0R1
D	?	-	VAL	deletion	UNP Q8N0R1
D	?	-	ASN	deletion	UNP Q8N0R1
D	?	-	PHE	deletion	UNP Q8N0R1
D	?	-	ASN	deletion	UNP Q8N0R1
D	?	-	ASP	deletion	UNP Q8N0R1
D	?	-	ASP	deletion	UNP Q8N0R1
D	?	-	ASP	deletion	UNP Q8N0R1
D	?	-	ASP	deletion	UNP Q8N0R1
D	?	-	ASN	deletion	UNP Q8N0R1
D	?	-	ASN	deletion	UNP Q8N0R1
D	?	-	TYR	deletion	UNP Q8N0R1
D	?	-	ASP	deletion	UNP Q8N0R1
D	?	-	ASP	deletion	UNP Q8N0R1
D	?	-	ASP	deletion	UNP Q8N0R1
D	?	-	GLY	deletion	UNP Q8N0R1
D	?	-	ASN	deletion	UNP Q8N0R1
D	282	HIS	-	expression tag	UNP Q8N0R1
D	283	HIS	-	expression tag	UNP Q8N0R1
D	284	HIS	-	expression tag	UNP Q8N0R1
D	285	HIS	-	expression tag	UNP Q8N0R1
D	286	HIS	-	expression tag	UNP Q8N0R1
D	287	HIS	-	expression tag	UNP Q8N0R1
E	0	MET	-	expression tag	UNP Q8N0R1
E	1	ALA	-	expression tag	UNP Q8N0R1
E	?	-	ASN	deletion	UNP Q8N0R1
E	?	-	ASN	deletion	UNP Q8N0R1
E	?	-	ASN	deletion	UNP Q8N0R1
E	?	-	ILE	deletion	UNP Q8N0R1
E	?	-	ASP	deletion	UNP Q8N0R1
E	?	-	ASN	deletion	UNP Q8N0R1
E	?	-	VAL	deletion	UNP Q8N0R1
E	?	-	ASN	deletion	UNP Q8N0R1
E	?	-	PHE	deletion	UNP Q8N0R1
E	?	-	ASN	deletion	UNP Q8N0R1
E	?	-	ASP	deletion	UNP Q8N0R1
E	?	-	ASP	deletion	UNP Q8N0R1
E	?	-	ASP	deletion	UNP Q8N0R1
E	?	-	ASP	deletion	UNP Q8N0R1
E	?	-	ASN	deletion	UNP Q8N0R1
E	?	-	ASN	deletion	UNP Q8N0R1
E	?	-	TYR	deletion	UNP Q8N0R1

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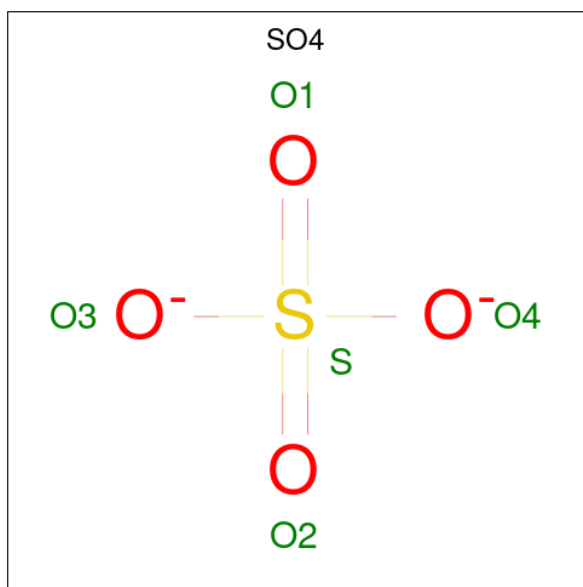| Chain	Residue	Modelled	Actual	Comment	Reference
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E	?	-	ASP	deletion	UNP Q8N0R1
E	?	-	ASP	deletion	UNP Q8N0R1
E	?	-	GLY	deletion	UNP Q8N0R1
E	?	-	ASN	deletion	UNP Q8N0R1
E	282	HIS	-	expression tag	UNP Q8N0R1
E	283	HIS	-	expression tag	UNP Q8N0R1
E	284	HIS	-	expression tag	UNP Q8N0R1
E	285	HIS	-	expression tag	UNP Q8N0R1
E	286	HIS	-	expression tag	UNP Q8N0R1
E	287	HIS	-	expression tag	UNP Q8N0R1
F	0	MET	-	expression tag	UNP Q8N0R1
F	1	ALA	-	expression tag	UNP Q8N0R1
F	?	-	ASN	deletion	UNP Q8N0R1
F	?	-	ASN	deletion	UNP Q8N0R1
F	?	-	ASN	deletion	UNP Q8N0R1
F	?	-	ILE	deletion	UNP Q8N0R1
F	?	-	ASP	deletion	UNP Q8N0R1
F	?	-	ASN	deletion	UNP Q8N0R1
F	?	-	VAL	deletion	UNP Q8N0R1
F	?	-	ASN	deletion	UNP Q8N0R1
F	?	-	PHE	deletion	UNP Q8N0R1
F	?	-	ASN	deletion	UNP Q8N0R1
F	?	-	ASP	deletion	UNP Q8N0R1
F	?	-	ASP	deletion	UNP Q8N0R1
F	?	-	ASP	deletion	UNP Q8N0R1
F	?	-	ASP	deletion	UNP Q8N0R1
F	?	-	ASN	deletion	UNP Q8N0R1
F	?	-	ASN	deletion	UNP Q8N0R1
F	?	-	TYR	deletion	UNP Q8N0R1
F	?	-	ASP	deletion	UNP Q8N0R1
F	?	-	ASP	deletion	UNP Q8N0R1
F	?	-	ASP	deletion	UNP Q8N0R1
F	?	-	GLY	deletion	UNP Q8N0R1
F	?	-	ASN	deletion	UNP Q8N0R1
F	282	HIS	-	expression tag	UNP Q8N0R1
F	283	HIS	-	expression tag	UNP Q8N0R1
F	284	HIS	-	expression tag	UNP Q8N0R1
F	285	HIS	-	expression tag	UNP Q8N0R1
F	286	HIS	-	expression tag	UNP Q8N0R1
F	287	HIS	-	expression tag	UNP Q8N0R1
G	0	MET	-	expression tag	UNP Q8N0R1

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Chain	Residue	Modelled	Actual	Comment	Reference
G	1	ALA	-	expression tag	UNP Q8N0R1
G	?	-	ASN	deletion	UNP Q8N0R1
G	?	-	ASN	deletion	UNP Q8N0R1
G	?	-	ASN	deletion	UNP Q8N0R1
G	?	-	ILE	deletion	UNP Q8N0R1
G	?	-	ASP	deletion	UNP Q8N0R1
G	?	-	ASN	deletion	UNP Q8N0R1
G	?	-	VAL	deletion	UNP Q8N0R1
G	?	-	ASN	deletion	UNP Q8N0R1
G	?	-	PHE	deletion	UNP Q8N0R1
G	?	-	ASN	deletion	UNP Q8N0R1
G	?	-	ASP	deletion	UNP Q8N0R1
G	?	-	ASP	deletion	UNP Q8N0R1
G	?	-	ASP	deletion	UNP Q8N0R1
G	?	-	ASP	deletion	UNP Q8N0R1
G	?	-	ASN	deletion	UNP Q8N0R1
G	?	-	ASN	deletion	UNP Q8N0R1
G	?	-	TYR	deletion	UNP Q8N0R1
G	?	-	ASP	deletion	UNP Q8N0R1
G	?	-	ASP	deletion	UNP Q8N0R1
G	?	-	ASP	deletion	UNP Q8N0R1
G	?	-	GLY	deletion	UNP Q8N0R1
G	?	-	ASN	deletion	UNP Q8N0R1
G	282	HIS	-	expression tag	UNP Q8N0R1
G	283	HIS	-	expression tag	UNP Q8N0R1
G	284	HIS	-	expression tag	UNP Q8N0R1
G	285	HIS	-	expression tag	UNP Q8N0R1
G	286	HIS	-	expression tag	UNP Q8N0R1
G	287	HIS	-	expression tag	UNP Q8N0R1
H	0	MET	-	expression tag	UNP Q8N0R1
H	1	ALA	-	expression tag	UNP Q8N0R1
H	?	-	ASN	deletion	UNP Q8N0R1
H	?	-	ASN	deletion	UNP Q8N0R1
H	?	-	ASN	deletion	UNP Q8N0R1
H	?	-	ILE	deletion	UNP Q8N0R1
H	?	-	ASP	deletion	UNP Q8N0R1
H	?	-	ASN	deletion	UNP Q8N0R1
H	?	-	VAL	deletion	UNP Q8N0R1
H	?	-	ASN	deletion	UNP Q8N0R1
H	?	-	PHE	deletion	UNP Q8N0R1
H	?	-	ASN	deletion	UNP Q8N0R1
H	?	-	ASP	deletion	UNP Q8N0R1

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Chain	Residue	Modelled	Actual	Comment	Reference
H	?	-	ASP	deletion	UNP Q8N0R1
H	?	-	ASP	deletion	UNP Q8N0R1
H	?	-	ASP	deletion	UNP Q8N0R1
H	?	-	ASN	deletion	UNP Q8N0R1
H	?	-	ASN	deletion	UNP Q8N0R1
H	?	-	TYR	deletion	UNP Q8N0R1
H	?	-	ASP	deletion	UNP Q8N0R1
H	?	-	ASP	deletion	UNP Q8N0R1
H	?	-	ASP	deletion	UNP Q8N0R1
H	?	-	GLY	deletion	UNP Q8N0R1
H	?	-	ASN	deletion	UNP Q8N0R1
H	282	HIS	-	expression tag	UNP Q8N0R1
H	283	HIS	-	expression tag	UNP Q8N0R1
H	284	HIS	-	expression tag	UNP Q8N0R1
H	285	HIS	-	expression tag	UNP Q8N0R1
H	286	HIS	-	expression tag	UNP Q8N0R1
H	287	HIS	-	expression tag	UNP Q8N0R1

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O_4S).

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

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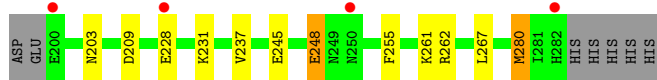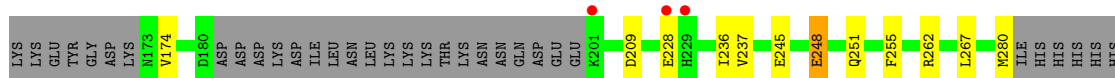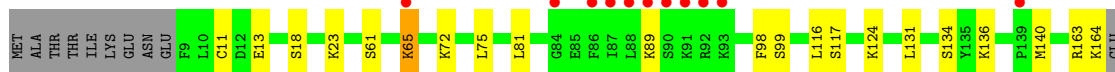| Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total 5	O 4	S 1	0	0
2	B	1	Total 5	O 4	S 1	0	0
2	B	1	Total 5	O 4	S 1	0	0
2	B	1	Total 5	O 4	S 1	0	0
2	B	1	Total 5	O 4	S 1	0	0
2	B	1	Total 5	O 4	S 1	0	0
2	C	1	Total 5	O 4	S 1	0	0
2	C	1	Total 5	O 4	S 1	0	0
2	C	1	Total 5	O 4	S 1	0	0
2	C	1	Total 5	O 4	S 1	0	0
2	C	1	Total 5	O 4	S 1	0	0
2	D	1	Total 5	O 4	S 1	0	0
2	D	1	Total 5	O 4	S 1	0	0
2	D	1	Total 5	O 4	S 1	0	0
2	D	1	Total 5	O 4	S 1	0	0
2	D	1	Total 5	O 4	S 1	0	0
2	E	1	Total 5	O 4	S 1	0	0
2	E	1	Total 5	O 4	S 1	0	0
2	E	1	Total 5	O 4	S 1	0	0
2	E	1	Total 5	O 4	S 1	0	0
2	F	1	Total 5	O 4	S 1	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	F	1	Total 5	O 4	S 1	0	0
2	F	1	Total 5	O 4	S 1	0	0
2	F	1	Total 5	O 4	S 1	0	0
2	G	1	Total 5	O 4	S 1	0	0
2	G	1	Total 5	O 4	S 1	0	0
2	G	1	Total 5	O 4	S 1	0	0
2	H	1	Total 5	O 4	S 1	0	0
2	H	1	Total 5	O 4	S 1	0	0
2	H	1	Total 5	O 4	S 1	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	56	Total 56	O 56	0	0
3	B	68	Total 68	O 68	0	0
3	C	39	Total 39	O 39	0	0
3	D	53	Total 53	O 53	0	0
3	E	13	Total 13	O 13	0	0
3	F	15	Total 15	O 15	0	0
3	G	19	Total 19	O 19	0	0
3	H	1	Total 1	O 1	0	0

3 Residue-property plots ⓘ

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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.
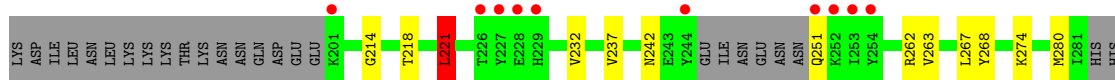
- Molecule 1: Orotate phosphoribosyltransferase

Chain A:

- Molecule 1: Orotate phosphoribosyltransferase

Chain B:

- Molecule 1: Orotate phosphoribosyltransferase

Chain C:

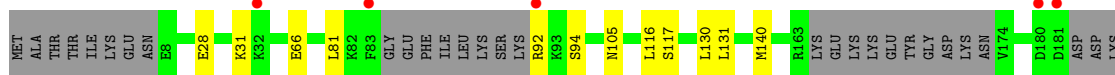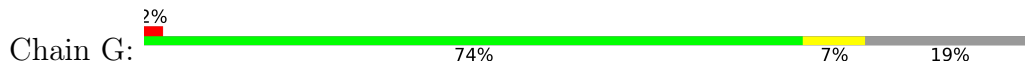- Molecule 1: Orotate phosphoribosyltransferase

Chain D:

- Molecule 1: Orotate phosphoribosyltransferase

Chain E:

- Molecule 1: Orotate phosphoribosyltransferase
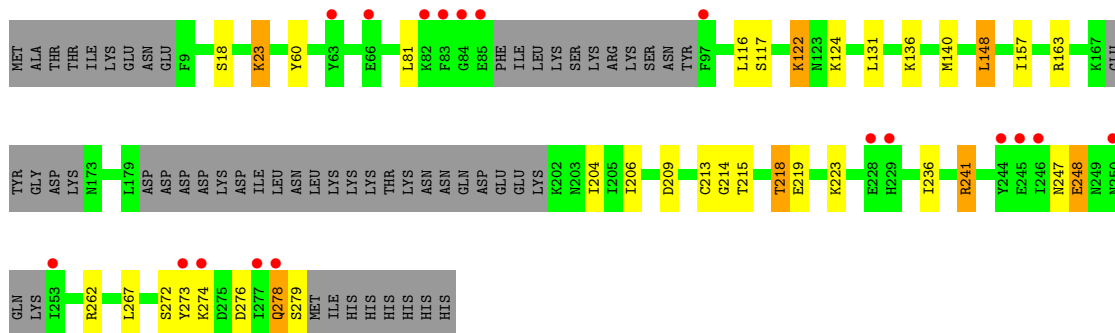
Chain F:

- Molecule 1: Orotate phosphoribosyltransferase

Chain G:

- Molecule 1: Orotate phosphoribosyltransferase

Chain H:

4 Data and refinement statistics ⓘ

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	114.77Å 152.49Å 167.75Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.47 – 2.60 40.44 – 2.59	Depositor EDS
% Data completeness (in resolution range)	98.5 (40.47-2.60) 98.6 (40.44-2.59)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.23	Depositor
$< I/\sigma(I) >$ [1]	1.65 (at 2.58Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.197 , 0.218 0.199 , 0.222	Depositor DCC
R_{free} test set	4588 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å^2)	51.0	Xtriage
Anisotropy	0.306	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.30 , 36.9	EDS
L-test for twinning[2]	$< \lvert L \rvert > = 0.50, < L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	14882	wwPDB-VP
Average B, all atoms (Å^2)	62.0	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 45.68 % 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 1.2757e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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[1]Intensities estimated from amplitudes.
[2]Theoretical values of $< \lvert L \rvert >$, $< L^2 >$ 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: 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	0.91	0/1878	0.87	4/2517 (0.2%)
1	B	0.99	4/1842 (0.2%)	0.92	4/2468 (0.2%)
1	C	0.82	0/1865	0.85	3/2495 (0.1%)
1	D	0.85	0/1915	0.90	2/2564 (0.1%)
1	E	0.76	0/1870	0.88	3/2503 (0.1%)
1	F	0.80	0/1754	0.87	4/2348 (0.2%)
1	G	0.82	2/1821 (0.1%)	0.89	3/2439 (0.1%)
1	H	0.71	0/1753	0.93	8/2346 (0.3%)
All	All	0.84	6/14698 (0.0%)	0.89	31/19680 (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	C	0	1

The worst 5 of 6 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	126	SER	CB-OG	-6.12	1.34	1.42
1	B	180	ASP	CB-CG	5.75	1.63	1.51
1	B	14	GLU	CD-OE2	5.67	1.31	1.25
1	G	66	GLU	CD-OE2	-5.58	1.19	1.25
1	B	14	GLU	CG-CD	5.48	1.60	1.51

The worst 5 of 31 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	262	ARG	NE-CZ-NH1	9.62	125.11	120.30
1	C	280	MET	CG-SD-CE	9.34	115.14	100.20
1	F	221	LEU	CA-CB-CG	7.39	132.29	115.30
1	H	276	ASP	CB-CG-OD2	7.32	124.89	118.30
1	A	277	ILE	CG1-CB-CG2	-7.09	95.81	111.40

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	83	PHE	Peptide

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1846	0	1901	12	0
1	B	1811	0	1873	16	0
1	C	1835	0	1927	13	0
1	D	1882	0	1960	15	0
1	E	1838	0	1926	17	0
1	F	1725	0	1801	19	0
1	G	1791	0	1862	9	0
1	H	1725	0	1801	17	0
2	A	20	0	0	0	0
2	B	25	0	0	2	0
2	C	25	0	0	1	0
2	D	25	0	0	0	0
2	E	20	0	0	4	0
2	F	20	0	0	1	0
2	G	15	0	0	0	0
2	H	15	0	0	1	0
3	A	56	0	0	2	0
3	B	68	0	0	4	0
3	C	39	0	0	2	0
3	D	53	0	0	1	0
3	E	13	0	0	0	0
3	F	15	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	G	19	0	0	0	0
3	H	1	0	0	0	0
All	All	14882	0	15051	113	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.

The worst 5 of 113 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:72:LYS:HD2	1:A:280:MET:HE2	1.56	0.88
1:C:88:LEU:HD22	1:C:97:PHE:HB2	1.58	0.84
1:F:134:SER:OG	1:F:162:ASP:O	1.96	0.82
1:A:72:LYS:HD2	1:A:280:MET:CE	2.14	0.77
1:C:88:LEU:CD2	1:C:97:PHE:HB2	2.16	0.74

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	215/266 (81%)	210 (98%)	5 (2%)	0	100	100
1	B	211/266 (79%)	209 (99%)	2 (1%)	0	100	100
1	C	214/266 (80%)	211 (99%)	3 (1%)	0	100	100
1	D	221/266 (83%)	215 (97%)	6 (3%)	0	100	100
1	E	216/266 (81%)	212 (98%)	4 (2%)	0	100	100
1	F	199/266 (75%)	197 (99%)	2 (1%)	0	100	100
1	G	208/266 (78%)	205 (99%)	3 (1%)	0	100	100
1	H	199/266 (75%)	194 (98%)	4 (2%)	1 (0%)	29	52

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	1683/2128 (79%)	1653 (98%)	29 (2%)	1 (0%)	51 75

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	H	278	GLN

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	211/252 (84%)	206 (98%)	5 (2%)	49 74
1	B	207/252 (82%)	201 (97%)	6 (3%)	42 68
1	C	210/252 (83%)	206 (98%)	4 (2%)	57 79
1	D	215/252 (85%)	208 (97%)	7 (3%)	38 64
1	E	210/252 (83%)	200 (95%)	10 (5%)	25 49
1	F	197/252 (78%)	192 (98%)	5 (2%)	47 73
1	G	205/252 (81%)	201 (98%)	4 (2%)	55 78
1	H	197/252 (78%)	188 (95%)	9 (5%)	27 51
All	All	1652/2016 (82%)	1602 (97%)	50 (3%)	41 67

5 of 50 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	236	ILE
1	F	237	VAL
1	H	272	SER
1	E	237	VAL
1	F	117	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	178	ASN
1	F	178	ASN
1	F	242	ASN
1	H	173	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

5.6 Ligand geometry ⓘ

33 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	SO4	B	305	-	4,4,4	0.74	0	6,6,6	0.96	0
2	SO4	E	303	-	4,4,4	0.63	0	6,6,6	0.40	0
2	SO4	E	304	-	4,4,4	0.52	0	6,6,6	0.66	0
2	SO4	C	304	-	4,4,4	0.55	0	6,6,6	0.47	0
2	SO4	C	305	-	4,4,4	0.39	0	6,6,6	0.76	0
2	SO4	G	302	-	4,4,4	0.34	0	6,6,6	0.27	0
2	SO4	D	302	-	4,4,4	0.30	0	6,6,6	0.63	0
2	SO4	F	303	-	4,4,4	0.53	0	6,6,6	0.80	0
2	SO4	A	302	-	4,4,4	0.34	0	6,6,6	0.19	0
2	SO4	D	303	-	4,4,4	0.26	0	6,6,6	0.58	0
2	SO4	G	301	-	4,4,4	0.57	0	6,6,6	0.73	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	#|Z| > 2	Counts	RMSZ	#|Z| > 2
2	SO4	H	301	-	4,4,4	0.57	0	6,6,6	0.51	0
2	SO4	D	304	-	4,4,4	0.14	0	6,6,6	0.65	0
2	SO4	F	302	-	4,4,4	0.48	0	6,6,6	0.77	0
2	SO4	E	301	-	4,4,4	0.42	0	6,6,6	1.11	1 (16%)
2	SO4	D	301	-	4,4,4	0.59	0	6,6,6	0.56	0
2	SO4	B	303	-	4,4,4	0.41	0	6,6,6	0.82	0
2	SO4	A	303	-	4,4,4	0.91	0	6,6,6	1.49	2 (33%)
2	SO4	B	302	-	4,4,4	0.29	0	6,6,6	0.40	0
2	SO4	H	303	-	4,4,4	0.55	0	6,6,6	0.61	0
2	SO4	F	301	-	4,4,4	0.32	0	6,6,6	1.14	0
2	SO4	C	303	-	4,4,4	0.42	0	6,6,6	0.31	0
2	SO4	A	304	-	4,4,4	0.88	0	6,6,6	1.03	1 (16%)
2	SO4	F	304	-	4,4,4	0.42	0	6,6,6	0.59	0
2	SO4	D	305	-	4,4,4	0.79	0	6,6,6	1.85	2 (33%)
2	SO4	H	302	-	4,4,4	0.55	0	6,6,6	0.64	0
2	SO4	B	301	-	4,4,4	0.56	0	6,6,6	0.87	0
2	SO4	A	301	-	4,4,4	0.50	0	6,6,6	1.14	0
2	SO4	B	304	-	4,4,4	0.78	0	6,6,6	1.42	2 (33%)
2	SO4	G	303	-	4,4,4	0.31	0	6,6,6	0.48	0
2	SO4	E	302	-	4,4,4	0.37	0	6,6,6	0.53	0
2	SO4	C	301	-	4,4,4	0.61	0	6,6,6	0.72	0
2	SO4	C	302	-	4,4,4	0.61	0	6,6,6	0.70	0

There are no bond length outliers.

The worst 5 of 8 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	D	305	SO4	O4-S-O3	-2.79	97.16	109.06
2	D	305	SO4	O4-S-O2	2.75	123.66	109.31
2	B	304	SO4	O4-S-O2	2.42	121.92	109.31
2	A	303	SO4	O3-S-O2	2.30	121.29	109.31
2	B	304	SO4	O4-S-O3	-2.28	99.35	109.06

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	305	SO4	2	0
2	E	304	SO4	1	0
2	C	305	SO4	1	0
2	E	301	SO4	3	0
2	F	304	SO4	1	0
2	H	302	SO4	1	0

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, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	A	223/266 (83%)	-0.19	10 (4%)	33	26	27, 44, 87, 130	0
1	B	219/266 (82%)	-0.27	6 (2%)	54	48	29, 43, 79, 108	0
1	C	222/266 (83%)	-0.17	10 (4%)	33	26	35, 52, 98, 144	0
1	D	227/266 (85%)	-0.15	7 (3%)	49	42	32, 47, 84, 120	0
1	E	222/266 (83%)	0.03	14 (6%)	20	15	45, 65, 105, 131	0
1	F	209/266 (78%)	-0.06	13 (6%)	20	15	39, 62, 105, 131	0
1	G	216/266 (81%)	-0.20	5 (2%)	60	54	41, 58, 100, 124	0
1	H	209/266 (78%)	0.46	18 (8%)	10	7	48, 82, 133, 169	0
All	All	1747/2128 (82%)	-0.07	83 (4%)	30	24	27, 56, 109, 169	0

The worst 5 of 83 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	86	PHE	5.7
1	H	244	TYR	5.6
1	E	87	ILE	5.3
1	H	246	ILE	5.3
1	H	250	ASN	5.2

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 monosaccharides in this entry.

WORLDWIDE
wwPDB
PROTEIN DATA BANK

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, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	SO4	E	304	5/5	0.71	0.27	103,103,132,141	0
2	SO4	C	304	5/5	0.80	0.34	121,122,126,127	0
2	SO4	E	302	5/5	0.85	0.20	112,116,123,124	0
2	SO4	D	305	5/5	0.85	0.24	74,79,102,106	0
2	SO4	H	303	5/5	0.91	0.12	84,93,100,104	0
2	SO4	B	305	5/5	0.94	0.11	77,82,86,94	0
2	SO4	H	302	5/5	0.94	0.14	86,88,89,91	0
2	SO4	A	303	5/5	0.94	0.16	53,57,66,70	0
2	SO4	G	303	5/5	0.95	0.10	83,83,86,94	0
2	SO4	C	305	5/5	0.95	0.11	72,75,87,87	0
2	SO4	G	301	5/5	0.95	0.11	72,86,89,89	0
2	SO4	E	301	5/5	0.96	0.11	69,69,73,82	0
2	SO4	C	301	5/5	0.96	0.11	69,70,76,82	0
2	SO4	B	304	5/5	0.96	0.12	54,62,62,65	0
2	SO4	F	303	5/5	0.96	0.08	76,79,81,81	0
2	SO4	H	301	5/5	0.97	0.07	84,87,89,98	0
2	SO4	F	302	5/5	0.97	0.17	78,79,89,97	0
2	SO4	F	304	5/5	0.97	0.13	67,67,73,87	0
2	SO4	D	304	5/5	0.98	0.12	64,72,77,77	0
2	SO4	B	302	5/5	0.98	0.11	55,59,62,64	0
2	SO4	B	303	5/5	0.98	0.08	47,54,57,60	0
2	SO4	A	301	5/5	0.98	0.12	54,54,58,59	0
2	SO4	A	304	5/5	0.98	0.10	53,54,60,60	0
2	SO4	F	301	5/5	0.98	0.12	65,68,74,75	0
2	SO4	D	302	5/5	0.98	0.10	66,71,79,79	0
2	SO4	B	301	5/5	0.99	0.09	42,43,48,50	0
2	SO4	A	302	5/5	0.99	0.07	46,46,48,52	0
2	SO4	D	301	5/5	0.99	0.17	53,55,58,61	0
2	SO4	G	302	5/5	0.99	0.09	48,48,51,52	0
2	SO4	E	303	5/5	0.99	0.09	48,51,52,54	0
2	SO4	C	302	5/5	0.99	0.08	46,47,48,48	0
2	SO4	D	303	5/5	0.99	0.14	44,45,48,48	0
2	SO4	C	303	5/5	0.99	0.08	63,66,69,70	0

6.5 Other polymers ⓘ

There are no such residues in this entry.