



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:40 PM GMT

PDB ID : 1Q1G
Title : Crystal structure of Plasmodium falciparum PNP with 5'-methylthio-immucillin-H
Authors : Shi, W.; Ting, L.M.; Kicska, G.A.; Lewandowicz, A.; Tyler, P.C.; Evans, G.B.; Furneaux, R.H.; Kim, K.; Almo, S.C.; Schramm, V.L.
Deposited on : 2003-07-19
Resolution : 2.02 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
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

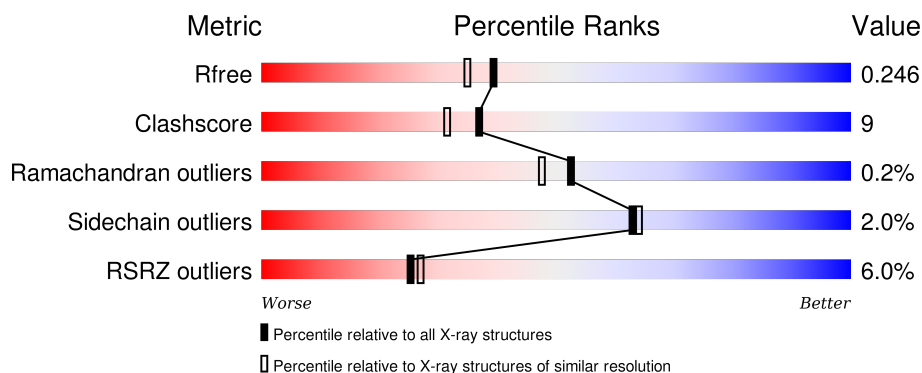
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.02 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}	91344	7858 (2.04-2.00)
Clashscore	102246	9060 (2.04-2.00)
Ramachandran outliers	100387	8952 (2.04-2.00)
Sidechain outliers	100360	8951 (2.04-2.00)
RSRZ outliers	91569	7873 (2.04-2.00)

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 ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	276	<div> <div>5%</div> <div>72% 15% 12%</div> </div>
1	B	276	<div> <div>9%</div> <div>71% 16% 12%</div> </div>
1	C	276	<div> <div>6%</div> <div>69% 18% 12%</div> </div>
1	D	276	<div> <div>4%</div> <div>67% 21% 12%</div> </div>
1	E	276	<div> <div>6%</div> <div>69% 18% 12%</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	276	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	IPA	A	503	-	-	-	X
4	IPA	B	502	-	-	-	X
4	IPA	B	508	-	-	-	X
4	IPA	D	507	-	-	-	X
4	IPA	E	506	-	-	-	X
4	IPA	F	505	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 11789 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 Uridine phosphorylase putative.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	243	Total	C	N	O	S	0	0	0
			1861	1179	319	347	16			
1	B	243	Total	C	N	O	S	0	0	0
			1861	1179	319	347	16			
1	C	243	Total	C	N	O	S	0	0	0
			1861	1179	319	347	16			
1	D	243	Total	C	N	O	S	0	0	0
			1861	1179	319	347	16			
1	E	243	Total	C	N	O	S	0	0	0
			1861	1179	319	347	16			
1	F	243	Total	C	N	O	S	0	0	0
			1861	1179	319	347	16			

There are 186 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	ALA	-	CLONING ARTIFACT	UNP Q8I3X4
A	1	LEU	-	CLONING ARTIFACT	UNP Q8I3X4
A	246	LYS	-	CLONING ARTIFACT	UNP Q8I3X4
A	247	GLY	-	CLONING ARTIFACT	UNP Q8I3X4
A	248	GLU	-	CLONING ARTIFACT	UNP Q8I3X4
A	249	PHE	-	CLONING ARTIFACT	UNP Q8I3X4
A	250	GLU	-	CLONING ARTIFACT	UNP Q8I3X4
A	251	ALA	-	CLONING ARTIFACT	UNP Q8I3X4
A	252	TYR	-	CLONING ARTIFACT	UNP Q8I3X4
A	253	VAL	-	CLONING ARTIFACT	UNP Q8I3X4
A	254	GLU	-	CLONING ARTIFACT	UNP Q8I3X4
A	255	GLN	-	CLONING ARTIFACT	UNP Q8I3X4
A	256	LYS	-	CLONING ARTIFACT	UNP Q8I3X4
A	257	LEU	-	CLONING ARTIFACT	UNP Q8I3X4
A	258	ILE	-	CLONING ARTIFACT	UNP Q8I3X4
A	259	SER	-	CLONING ARTIFACT	UNP Q8I3X4
A	260	GLU	-	CLONING ARTIFACT	UNP Q8I3X4

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Chain	Residue	Modelled	Actual	Comment	Reference
A	261	GLU	-	CLONING ARTIFACT	UNP Q8I3X4
A	262	ASP	-	CLONING ARTIFACT	UNP Q8I3X4
A	263	LEU	-	CLONING ARTIFACT	UNP Q8I3X4
A	264	ASN	-	CLONING ARTIFACT	UNP Q8I3X4
A	265	SER	-	CLONING ARTIFACT	UNP Q8I3X4
A	266	ALA	-	CLONING ARTIFACT	UNP Q8I3X4
A	267	VAL	-	CLONING ARTIFACT	UNP Q8I3X4
A	268	ASP	-	CLONING ARTIFACT	UNP Q8I3X4
A	269	HIS	-	EXPRESSION TAG	UNP Q8I3X4
A	270	HIS	-	EXPRESSION TAG	UNP Q8I3X4
A	271	HIS	-	EXPRESSION TAG	UNP Q8I3X4
A	272	HIS	-	EXPRESSION TAG	UNP Q8I3X4
A	273	HIS	-	EXPRESSION TAG	UNP Q8I3X4
A	274	HIS	-	EXPRESSION TAG	UNP Q8I3X4
B	0	ALA	-	CLONING ARTIFACT	UNP Q8I3X4
B	1	LEU	-	CLONING ARTIFACT	UNP Q8I3X4
B	246	LYS	-	CLONING ARTIFACT	UNP Q8I3X4
B	247	GLY	-	CLONING ARTIFACT	UNP Q8I3X4
B	248	GLU	-	CLONING ARTIFACT	UNP Q8I3X4
B	249	PHE	-	CLONING ARTIFACT	UNP Q8I3X4
B	250	GLU	-	CLONING ARTIFACT	UNP Q8I3X4
B	251	ALA	-	CLONING ARTIFACT	UNP Q8I3X4
B	252	TYR	-	CLONING ARTIFACT	UNP Q8I3X4
B	253	VAL	-	CLONING ARTIFACT	UNP Q8I3X4
B	254	GLU	-	CLONING ARTIFACT	UNP Q8I3X4
B	255	GLN	-	CLONING ARTIFACT	UNP Q8I3X4
B	256	LYS	-	CLONING ARTIFACT	UNP Q8I3X4
B	257	LEU	-	CLONING ARTIFACT	UNP Q8I3X4
B	258	ILE	-	CLONING ARTIFACT	UNP Q8I3X4
B	259	SER	-	CLONING ARTIFACT	UNP Q8I3X4
B	260	GLU	-	CLONING ARTIFACT	UNP Q8I3X4
B	261	GLU	-	CLONING ARTIFACT	UNP Q8I3X4
B	262	ASP	-	CLONING ARTIFACT	UNP Q8I3X4
B	263	LEU	-	CLONING ARTIFACT	UNP Q8I3X4
B	264	ASN	-	CLONING ARTIFACT	UNP Q8I3X4
B	265	SER	-	CLONING ARTIFACT	UNP Q8I3X4
B	266	ALA	-	CLONING ARTIFACT	UNP Q8I3X4
B	267	VAL	-	CLONING ARTIFACT	UNP Q8I3X4
B	268	ASP	-	CLONING ARTIFACT	UNP Q8I3X4
B	269	HIS	-	EXPRESSION TAG	UNP Q8I3X4
B	270	HIS	-	EXPRESSION TAG	UNP Q8I3X4
B	271	HIS	-	EXPRESSION TAG	UNP Q8I3X4

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Chain	Residue	Modelled	Actual	Comment	Reference
B	272	HIS	-	EXPRESSION TAG	UNP Q8I3X4
B	273	HIS	-	EXPRESSION TAG	UNP Q8I3X4
B	274	HIS	-	EXPRESSION TAG	UNP Q8I3X4
C	0	ALA	-	CLONING ARTIFACT	UNP Q8I3X4
C	1	LEU	-	CLONING ARTIFACT	UNP Q8I3X4
C	246	LYS	-	CLONING ARTIFACT	UNP Q8I3X4
C	247	GLY	-	CLONING ARTIFACT	UNP Q8I3X4
C	248	GLU	-	CLONING ARTIFACT	UNP Q8I3X4
C	249	PHE	-	CLONING ARTIFACT	UNP Q8I3X4
C	250	GLU	-	CLONING ARTIFACT	UNP Q8I3X4
C	251	ALA	-	CLONING ARTIFACT	UNP Q8I3X4
C	252	TYR	-	CLONING ARTIFACT	UNP Q8I3X4
C	253	VAL	-	CLONING ARTIFACT	UNP Q8I3X4
C	254	GLU	-	CLONING ARTIFACT	UNP Q8I3X4
C	255	GLN	-	CLONING ARTIFACT	UNP Q8I3X4
C	256	LYS	-	CLONING ARTIFACT	UNP Q8I3X4
C	257	LEU	-	CLONING ARTIFACT	UNP Q8I3X4
C	258	ILE	-	CLONING ARTIFACT	UNP Q8I3X4
C	259	SER	-	CLONING ARTIFACT	UNP Q8I3X4
C	260	GLU	-	CLONING ARTIFACT	UNP Q8I3X4
C	261	GLU	-	CLONING ARTIFACT	UNP Q8I3X4
C	262	ASP	-	CLONING ARTIFACT	UNP Q8I3X4
C	263	LEU	-	CLONING ARTIFACT	UNP Q8I3X4
C	264	ASN	-	CLONING ARTIFACT	UNP Q8I3X4
C	265	SER	-	CLONING ARTIFACT	UNP Q8I3X4
C	266	ALA	-	CLONING ARTIFACT	UNP Q8I3X4
C	267	VAL	-	CLONING ARTIFACT	UNP Q8I3X4
C	268	ASP	-	CLONING ARTIFACT	UNP Q8I3X4
C	269	HIS	-	EXPRESSION TAG	UNP Q8I3X4
C	270	HIS	-	EXPRESSION TAG	UNP Q8I3X4
C	271	HIS	-	EXPRESSION TAG	UNP Q8I3X4
C	272	HIS	-	EXPRESSION TAG	UNP Q8I3X4
C	273	HIS	-	EXPRESSION TAG	UNP Q8I3X4
C	274	HIS	-	EXPRESSION TAG	UNP Q8I3X4
D	0	ALA	-	CLONING ARTIFACT	UNP Q8I3X4
D	1	LEU	-	CLONING ARTIFACT	UNP Q8I3X4
D	246	LYS	-	CLONING ARTIFACT	UNP Q8I3X4
D	247	GLY	-	CLONING ARTIFACT	UNP Q8I3X4
D	248	GLU	-	CLONING ARTIFACT	UNP Q8I3X4
D	249	PHE	-	CLONING ARTIFACT	UNP Q8I3X4
D	250	GLU	-	CLONING ARTIFACT	UNP Q8I3X4
D	251	ALA	-	CLONING ARTIFACT	UNP Q8I3X4

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Chain	Residue	Modelled	Actual	Comment	Reference
D	252	TYR	-	CLONING ARTIFACT	UNP Q8I3X4
D	253	VAL	-	CLONING ARTIFACT	UNP Q8I3X4
D	254	GLU	-	CLONING ARTIFACT	UNP Q8I3X4
D	255	GLN	-	CLONING ARTIFACT	UNP Q8I3X4
D	256	LYS	-	CLONING ARTIFACT	UNP Q8I3X4
D	257	LEU	-	CLONING ARTIFACT	UNP Q8I3X4
D	258	ILE	-	CLONING ARTIFACT	UNP Q8I3X4
D	259	SER	-	CLONING ARTIFACT	UNP Q8I3X4
D	260	GLU	-	CLONING ARTIFACT	UNP Q8I3X4
D	261	GLU	-	CLONING ARTIFACT	UNP Q8I3X4
D	262	ASP	-	CLONING ARTIFACT	UNP Q8I3X4
D	263	LEU	-	CLONING ARTIFACT	UNP Q8I3X4
D	264	ASN	-	CLONING ARTIFACT	UNP Q8I3X4
D	265	SER	-	CLONING ARTIFACT	UNP Q8I3X4
D	266	ALA	-	CLONING ARTIFACT	UNP Q8I3X4
D	267	VAL	-	CLONING ARTIFACT	UNP Q8I3X4
D	268	ASP	-	CLONING ARTIFACT	UNP Q8I3X4
D	269	HIS	-	EXPRESSION TAG	UNP Q8I3X4
D	270	HIS	-	EXPRESSION TAG	UNP Q8I3X4
D	271	HIS	-	EXPRESSION TAG	UNP Q8I3X4
D	272	HIS	-	EXPRESSION TAG	UNP Q8I3X4
D	273	HIS	-	EXPRESSION TAG	UNP Q8I3X4
D	274	HIS	-	EXPRESSION TAG	UNP Q8I3X4
E	0	ALA	-	CLONING ARTIFACT	UNP Q8I3X4
E	1	LEU	-	CLONING ARTIFACT	UNP Q8I3X4
E	246	LYS	-	CLONING ARTIFACT	UNP Q8I3X4
E	247	GLY	-	CLONING ARTIFACT	UNP Q8I3X4
E	248	GLU	-	CLONING ARTIFACT	UNP Q8I3X4
E	249	PHE	-	CLONING ARTIFACT	UNP Q8I3X4
E	250	GLU	-	CLONING ARTIFACT	UNP Q8I3X4
E	251	ALA	-	CLONING ARTIFACT	UNP Q8I3X4
E	252	TYR	-	CLONING ARTIFACT	UNP Q8I3X4
E	253	VAL	-	CLONING ARTIFACT	UNP Q8I3X4
E	254	GLU	-	CLONING ARTIFACT	UNP Q8I3X4
E	255	GLN	-	CLONING ARTIFACT	UNP Q8I3X4
E	256	LYS	-	CLONING ARTIFACT	UNP Q8I3X4
E	257	LEU	-	CLONING ARTIFACT	UNP Q8I3X4
E	258	ILE	-	CLONING ARTIFACT	UNP Q8I3X4
E	259	SER	-	CLONING ARTIFACT	UNP Q8I3X4
E	260	GLU	-	CLONING ARTIFACT	UNP Q8I3X4
E	261	GLU	-	CLONING ARTIFACT	UNP Q8I3X4
E	262	ASP	-	CLONING ARTIFACT	UNP Q8I3X4

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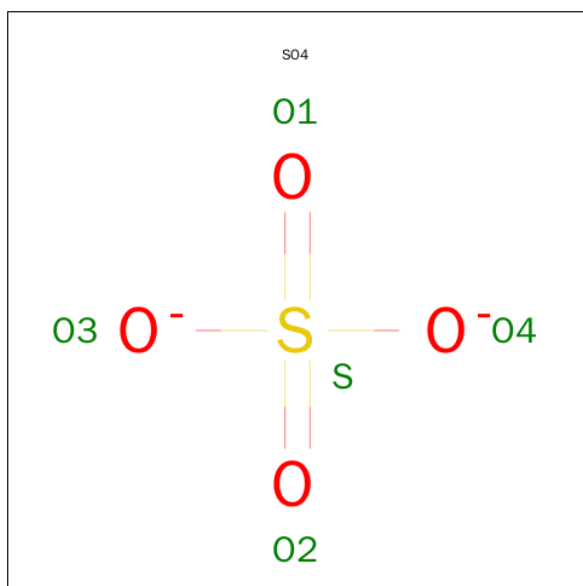
Chain	Residue	Modelled	Actual	Comment	Reference
E	263	LEU	-	CLONING ARTIFACT	UNP Q8I3X4
E	264	ASN	-	CLONING ARTIFACT	UNP Q8I3X4
E	265	SER	-	CLONING ARTIFACT	UNP Q8I3X4
E	266	ALA	-	CLONING ARTIFACT	UNP Q8I3X4
E	267	VAL	-	CLONING ARTIFACT	UNP Q8I3X4
E	268	ASP	-	CLONING ARTIFACT	UNP Q8I3X4
E	269	HIS	-	EXPRESSION TAG	UNP Q8I3X4
E	270	HIS	-	EXPRESSION TAG	UNP Q8I3X4
E	271	HIS	-	EXPRESSION TAG	UNP Q8I3X4
E	272	HIS	-	EXPRESSION TAG	UNP Q8I3X4
E	273	HIS	-	EXPRESSION TAG	UNP Q8I3X4
E	274	HIS	-	EXPRESSION TAG	UNP Q8I3X4
F	0	ALA	-	CLONING ARTIFACT	UNP Q8I3X4
F	1	LEU	-	CLONING ARTIFACT	UNP Q8I3X4
F	246	LYS	-	CLONING ARTIFACT	UNP Q8I3X4
F	247	GLY	-	CLONING ARTIFACT	UNP Q8I3X4
F	248	GLU	-	CLONING ARTIFACT	UNP Q8I3X4
F	249	PHE	-	CLONING ARTIFACT	UNP Q8I3X4
F	250	GLU	-	CLONING ARTIFACT	UNP Q8I3X4
F	251	ALA	-	CLONING ARTIFACT	UNP Q8I3X4
F	252	TYR	-	CLONING ARTIFACT	UNP Q8I3X4
F	253	VAL	-	CLONING ARTIFACT	UNP Q8I3X4
F	254	GLU	-	CLONING ARTIFACT	UNP Q8I3X4
F	255	GLN	-	CLONING ARTIFACT	UNP Q8I3X4
F	256	LYS	-	CLONING ARTIFACT	UNP Q8I3X4
F	257	LEU	-	CLONING ARTIFACT	UNP Q8I3X4
F	258	ILE	-	CLONING ARTIFACT	UNP Q8I3X4
F	259	SER	-	CLONING ARTIFACT	UNP Q8I3X4
F	260	GLU	-	CLONING ARTIFACT	UNP Q8I3X4
F	261	GLU	-	CLONING ARTIFACT	UNP Q8I3X4
F	262	ASP	-	CLONING ARTIFACT	UNP Q8I3X4
F	263	LEU	-	CLONING ARTIFACT	UNP Q8I3X4
F	264	ASN	-	CLONING ARTIFACT	UNP Q8I3X4
F	265	SER	-	CLONING ARTIFACT	UNP Q8I3X4
F	266	ALA	-	CLONING ARTIFACT	UNP Q8I3X4
F	267	VAL	-	CLONING ARTIFACT	UNP Q8I3X4
F	268	ASP	-	CLONING ARTIFACT	UNP Q8I3X4
F	269	HIS	-	EXPRESSION TAG	UNP Q8I3X4
F	270	HIS	-	EXPRESSION TAG	UNP Q8I3X4
F	271	HIS	-	EXPRESSION TAG	UNP Q8I3X4
F	272	HIS	-	EXPRESSION TAG	UNP Q8I3X4
F	273	HIS	-	EXPRESSION TAG	UNP Q8I3X4

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Chain	Residue	Modelled	Actual	Comment	Reference
F	274	HIS	-	EXPRESSION TAG	UNP Q8I3X4

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



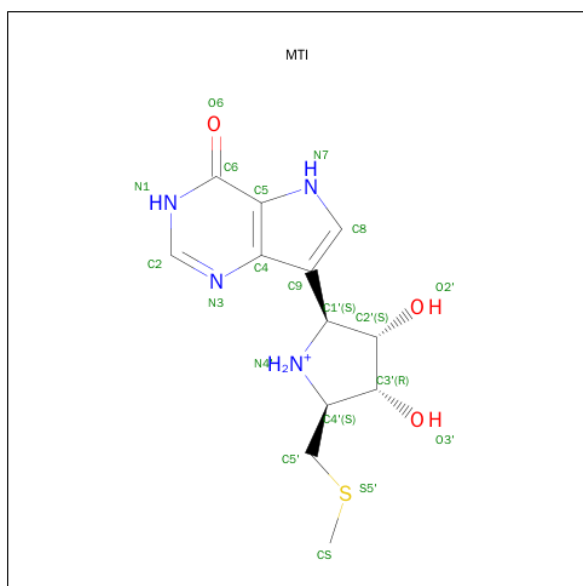
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	E	1	Total	O	S	0	0
			5	4	1		
2	F	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	F	1	Total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

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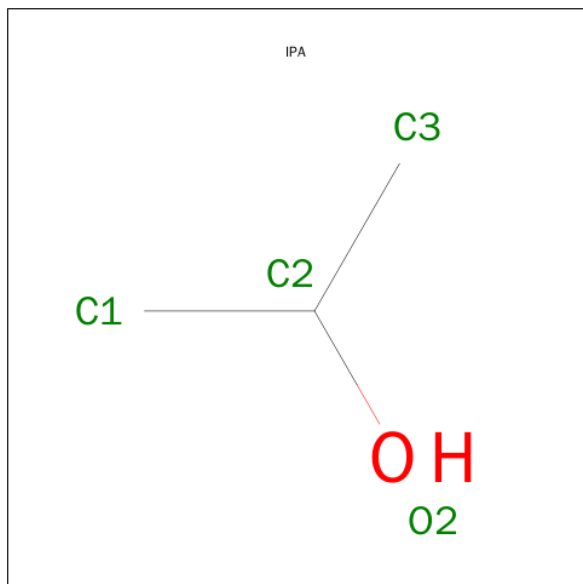
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	D	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	E	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	F	1	Total	O	S	0	0
			5	4	1		
2	E	1	Total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

- Molecule 3 is 3,4-DIHYDROXY-2-[(METHYLSULFANYL)METHYL]-5-(4-OXO-4,5-DIHYDRO-3H-PYRROLO[3,2-D]PYRIMIDIN-7-YL)PYRROLIDINIUM (three-letter code: MTI) (formula: C₁₂H₁₇N₄O₃S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			20	12	4	3	1		
3	B	1	Total	C	N	O	S	0	0
			20	12	4	3	1		
3	C	1	Total	C	N	O	S	0	0
			20	12	4	3	1		
3	D	1	Total	C	N	O	S	0	0
			20	12	4	3	1		
3	E	1	Total	C	N	O	S	0	0
			20	12	4	3	1		
3	F	1	Total	C	N	O	S	0	0
			20	12	4	3	1		

- Molecule 4 is ISOPROPYL ALCOHOL (three-letter code: IPA) (formula: C₃H₈O).



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

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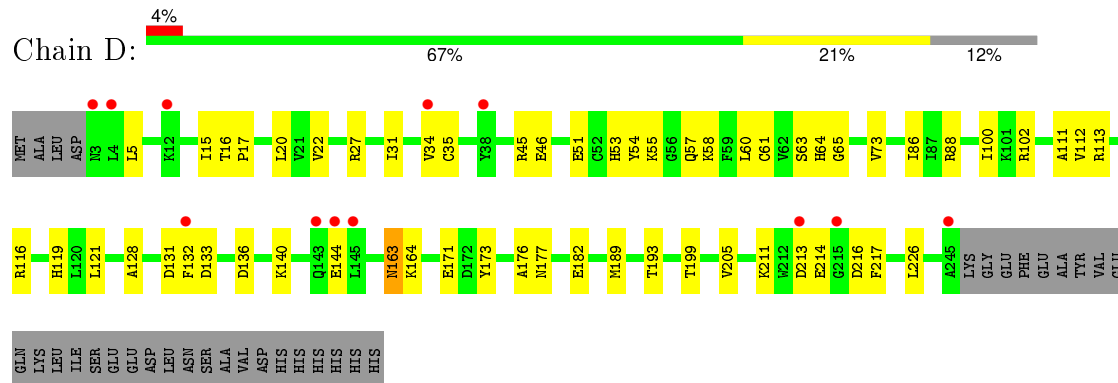
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	D	1	Total	C	O	0	0
			4	3	1		
4	B	1	Total	C	O	0	0
			4	3	1		
4	D	1	Total	C	O	0	0
			4	3	1		

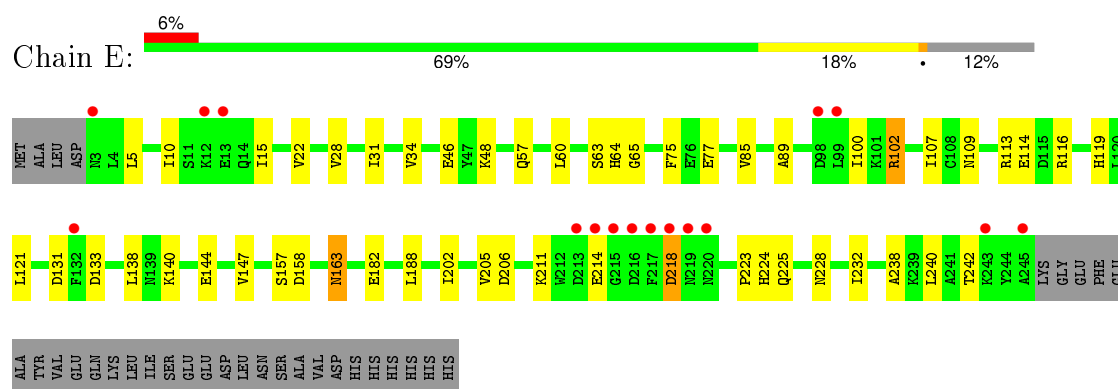
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	56	Total	O	0	0
			56	56		
5	B	45	Total	O	0	0
			45	45		
5	C	51	Total	O	0	0
			51	51		
5	D	61	Total	O	0	0
			61	61		
5	E	65	Total	O	0	0
			65	65		
5	F	79	Total	O	0	0
			79	79		

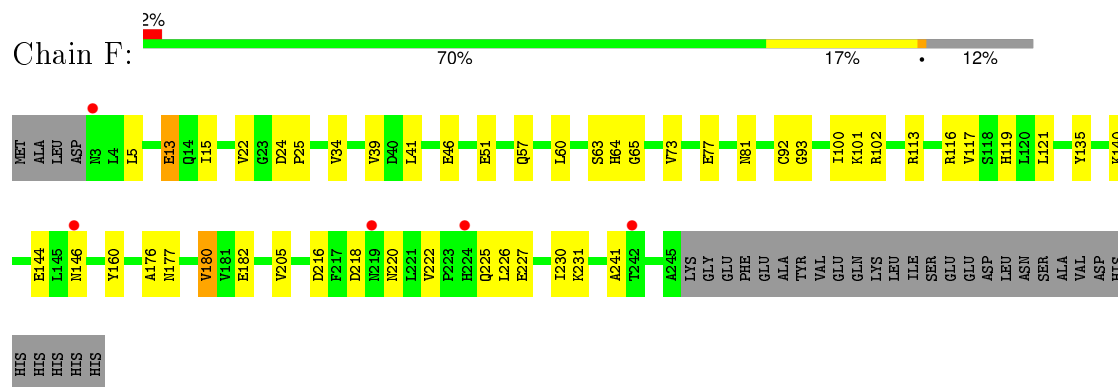
- Molecule 1: Uridine phosphorylase putative



- Molecule 1: Uridine phosphorylase putative



- Molecule 1: Uridine phosphorylase putative



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	86.44Å 91.73Å 238.94Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.02 45.87 – 2.02	Depositor EDS
% Data completeness (in resolution range)	88.8 (30.00-2.02) 91.5 (45.87-2.02)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.97 (at 2.03Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.211 , 0.242 0.215 , 0.246	Depositor DCC
R_{free} test set	11400 reflections (10.01%)	DCC
Wilson B-factor (Å ²)	23.0	Xtriage
Anisotropy	0.478	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 53.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 113901 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	11789	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.44% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: MTI, IPA, 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	0.32	0/1893	0.58	0/2561
1	B	0.30	0/1893	0.56	0/2561
1	C	0.31	0/1893	0.57	0/2561
1	D	0.31	0/1893	0.59	0/2561
1	E	0.33	0/1893	0.59	0/2561
1	F	0.33	0/1893	0.59	0/2561
All	All	0.32	0/11358	0.58	0/15366

There are no bond length outliers.

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	1861	0	1882	35	0
1	B	1861	0	1882	38	0
1	C	1861	0	1882	35	0
1	D	1861	0	1882	41	0
1	E	1861	0	1882	38	0
1	F	1861	0	1882	40	0
2	A	20	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	20	0	0	0	0
2	C	15	0	0	0	0
2	D	20	0	0	1	0
2	E	15	0	0	0	0
2	F	20	0	0	0	0
3	A	20	0	17	0	0
3	B	20	0	17	2	0
3	C	20	0	17	1	0
3	D	20	0	17	0	0
3	E	20	0	17	2	0
3	F	20	0	17	1	0
4	A	4	0	8	0	0
4	B	12	0	24	1	0
4	D	8	0	16	1	0
4	E	4	0	8	0	0
4	F	8	0	16	1	0
5	A	56	0	0	2	0
5	B	45	0	0	1	0
5	C	51	0	0	1	0
5	D	61	0	0	3	0
5	E	65	0	0	3	0
5	F	79	0	0	2	0
All	All	11789	0	11466	215	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 9.

All (215) 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:C:5:LEU:HD11	1:C:15:ILE:HD11	1.49	0.94
1:D:163:ASN:HD22	1:D:164:LYS:H	1.20	0.89
1:E:102:ARG:HH11	1:E:102:ARG:HB2	1.37	0.86
1:A:133:ASP:O	1:A:137:THR:HG23	1.76	0.84
1:B:5:LEU:HD11	1:B:15:ILE:HD11	1.61	0.83
1:F:92:CYS:SG	1:F:180:VAL:HG21	2.18	0.82
1:B:11:SER:HB2	1:B:13:GLU:OE2	1.80	0.80
1:B:117:VAL:O	1:B:121:LEU:HD13	1.82	0.79
1:A:5:LEU:HD11	1:A:15:ILE:HD11	1.67	0.77
1:A:46:GLU:HB3	1:B:46:GLU:HB3	1.70	0.74
5:E:828:HOH:O	1:F:117:VAL:HG23	1.88	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:93:GLY:O	1:F:180:VAL:HG22	1.90	0.72
1:D:64:HIS:HD2	1:D:65:GLY:O	1.73	0.71
1:B:141:CYS:O	1:B:145:LEU:HD13	1.92	0.70
1:E:158:ASP:HB3	1:F:117:VAL:HG22	1.73	0.70
1:C:53:HIS:HE1	1:C:58:LYS:HE2	1.56	0.70
1:F:5:LEU:HD11	1:F:15:ILE:HD11	1.72	0.69
1:B:13:GLU:H	1:B:13:GLU:CD	1.94	0.69
1:E:5:LEU:HD11	1:E:15:ILE:HD11	1.75	0.68
1:F:140:LYS:O	1:F:144:GLU:HG3	1.93	0.67
1:B:102:ARG:NH2	1:B:216:ASP:HA	2.08	0.67
1:A:163:ASN:HB3	5:A:691:HOH:O	1.93	0.67
1:B:3:ASN:N	1:B:3:ASN:HD22	1.92	0.66
1:F:102:ARG:HD2	1:F:216:ASP:OD1	1.94	0.66
1:D:31:ILE:O	1:D:34:VAL:HG12	1.95	0.66
1:E:31:ILE:O	1:E:34:VAL:HG22	1.95	0.66
1:E:163:ASN:HD22	1:E:163:ASN:C	1.98	0.66
1:E:46:GLU:HB3	1:F:46:GLU:HB3	1.76	0.66
1:B:11:SER:OG	1:B:14:GLN:HG3	1.95	0.65
1:A:92:CYS:HB2	1:A:180:VAL:HG13	1.77	0.65
1:C:46:GLU:HB3	1:D:46:GLU:HB3	1.79	0.65
1:C:64:HIS:HD2	1:C:65:GLY:O	1.80	0.65
1:C:13:GLU:CD	1:C:13:GLU:H	1.98	0.65
1:F:64:HIS:HD2	1:F:65:GLY:O	1.81	0.64
1:C:218:ASP:HB3	1:C:220:ASN:O	1.99	0.63
1:B:22:VAL:HG11	1:B:27:ARG:HG2	1.79	0.63
1:D:211:LYS:HE3	1:D:216:ASP:OD2	2.00	0.62
1:A:131:ASP:OD1	1:A:133:ASP:HB2	2.00	0.62
1:C:113:ARG:O	1:C:119:HIS:HE1	1.83	0.62
1:F:92:CYS:SG	1:F:180:VAL:CG2	2.88	0.62
1:B:64:HIS:HD2	1:B:65:GLY:O	1.83	0.61
1:B:163:ASN:HD22	1:B:163:ASN:H	1.48	0.61
1:D:163:ASN:ND2	1:D:164:LYS:H	1.95	0.60
1:A:64:HIS:HD2	1:A:65:GLY:O	1.84	0.60
1:F:146:ASN:HB3	5:F:954:HOH:O	2.01	0.60
1:E:102:ARG:CB	1:E:102:ARG:HH11	2.13	0.58
1:C:14:GLN:NE2	1:C:41:LEU:HD22	2.19	0.58
1:C:53:HIS:CE1	1:C:58:LYS:HE2	2.37	0.58
1:B:3:ASN:N	1:B:3:ASN:ND2	2.50	0.58
1:D:15:ILE:HA	1:D:60:LEU:HD11	1.86	0.58
1:E:140:LYS:O	1:E:144:GLU:HG3	2.04	0.57
1:B:227:GLU:HG2	1:B:231:LYS:HZ2	1.70	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:180:VAL:CG1	1:A:181:VAL:N	2.68	0.56
1:A:15:ILE:HA	1:A:60:LEU:HD11	1.86	0.56
1:D:113:ARG:O	1:D:119:HIS:HE1	1.89	0.56
3:E:305:MTI:HS3	1:F:73:VAL:HG21	1.88	0.55
1:E:85:VAL:HG11	1:E:240:LEU:HD13	1.87	0.55
1:D:51:GLU:OE2	1:D:58:LYS:HE3	2.07	0.55
1:D:22:VAL:O	1:D:63:SER:HA	2.07	0.55
1:C:100:ILE:HG22	1:C:205:VAL:HG21	1.88	0.55
1:A:85:VAL:HG11	1:A:240:LEU:HD13	1.88	0.54
1:D:189:MET:O	1:D:193:THR:HG23	2.05	0.54
1:C:116:ARG:HB2	1:D:116:ARG:HB2	1.88	0.54
1:A:116:ARG:HB2	1:B:116:ARG:HB2	1.90	0.54
1:C:85:VAL:HG11	1:C:240:LEU:HD13	1.88	0.54
1:E:64:HIS:HD2	1:E:65:GLY:O	1.90	0.54
1:D:5:LEU:HD11	1:D:15:ILE:HD11	1.88	0.54
1:A:57:GLN:NE2	1:A:57:GLN:HA	2.22	0.54
1:E:206:ASP:O	1:E:218:ASP:HB2	2.07	0.53
1:E:113:ARG:O	1:E:119:HIS:HE1	1.92	0.53
1:C:97:PRO:O	1:C:101:LYS:HE2	2.08	0.53
1:B:22:VAL:HG11	1:B:27:ARG:CG	2.38	0.53
1:E:5:LEU:HD13	1:E:77:GLU:HB3	1.90	0.53
1:D:116:ARG:HE	4:D:509:IPA:H13	1.74	0.53
1:D:213:ASP:OD2	1:D:214:GLU:HG3	2.09	0.53
1:E:163:ASN:C	1:E:163:ASN:ND2	2.62	0.53
1:C:102:ARG:HD2	1:C:218:ASP:OD2	2.09	0.52
1:C:140:LYS:O	1:C:144:GLU:HG3	2.10	0.52
1:B:113:ARG:O	1:B:119:HIS:HE1	1.93	0.52
1:E:211:LYS:HD3	1:E:214:GLU:OE1	2.09	0.52
1:B:163:ASN:ND2	1:B:163:ASN:H	2.07	0.51
1:F:116:ARG:HE	4:F:504:IPA:H13	1.76	0.51
1:A:222:VAL:HB	1:A:225:GLN:HB2	1.93	0.51
1:F:113:ARG:O	1:F:119:HIS:HE1	1.94	0.51
1:E:22:VAL:O	1:E:63:SER:HA	2.11	0.51
1:A:92:CYS:HB2	1:A:180:VAL:CG1	2.40	0.50
1:D:171:GLU:HG2	5:D:786:HOH:O	2.11	0.50
1:A:46:GLU:CB	1:B:46:GLU:HB3	2.41	0.50
1:B:22:VAL:HG12	5:B:780:HOH:O	2.12	0.50
1:C:73:VAL:O	1:C:77:GLU:HG3	2.12	0.49
1:A:180:VAL:HG13	1:A:181:VAL:N	2.27	0.49
1:B:105:ILE:HB	1:B:149:VAL:HG12	1.93	0.49
1:D:57:GLN:NE2	5:D:725:HOH:O	2.44	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:218:ASP:HB3	1:F:220:ASN:O	2.13	0.49
1:A:22:VAL:O	1:A:63:SER:HA	2.13	0.49
1:B:102:ARG:HH21	1:B:216:ASP:HA	1.75	0.48
1:A:35:CYS:HB3	1:A:53:HIS:O	2.13	0.48
1:A:102:ARG:HD2	1:A:216:ASP:OD1	2.12	0.48
1:D:140:LYS:O	1:D:144:GLU:HG3	2.13	0.48
1:F:5:LEU:HD22	1:F:81:ASN:ND2	2.28	0.48
1:C:176:ALA:O	1:C:177:ASN:HB2	2.13	0.48
1:E:116:ARG:HB2	1:F:116:ARG:HB2	1.94	0.48
1:E:158:ASP:HB3	1:F:117:VAL:CG2	2.43	0.48
1:B:211:LYS:HG2	1:B:214:GLU:OE1	2.13	0.48
1:B:227:GLU:HG2	1:B:231:LYS:NZ	2.29	0.48
1:A:25:PRO:HA	1:A:63:SER:HB3	1.95	0.48
1:D:136:ASP:OD2	1:D:140:LYS:HE3	2.14	0.48
1:F:93:GLY:O	1:F:180:VAL:CG2	2.61	0.47
1:F:100:ILE:HG22	1:F:205:VAL:HG21	1.97	0.47
1:B:14:GLN:NE2	1:B:41:LEU:HD22	2.30	0.47
1:D:64:HIS:CD2	1:D:65:GLY:O	2.61	0.47
1:B:222:VAL:HB	1:B:225:GLN:HB2	1.96	0.47
1:D:102:ARG:HD3	2:D:412:SO4:O3	2.14	0.47
1:F:34:VAL:O	1:F:34:VAL:HG12	2.14	0.47
1:D:34:VAL:O	1:D:34:VAL:HG22	2.14	0.47
1:D:176:ALA:O	1:D:177:ASN:HB2	2.15	0.47
1:A:132:PHE:CG	1:F:135:TYR:HE2	2.32	0.47
1:E:131:ASP:OD1	1:E:133:ASP:HB2	2.14	0.47
1:C:29:ASP:HA	1:C:32:LYS:HE2	1.97	0.47
1:E:100:ILE:HG22	1:E:205:VAL:HG21	1.97	0.47
1:D:16:THR:HB	1:D:17:PRO:HD2	1.97	0.47
1:A:113:ARG:O	1:A:119:HIS:HE1	1.97	0.47
1:B:112:VAL:HG11	1:B:173:TYR:CZ	2.50	0.47
1:C:51:GLU:HA	1:C:60:LEU:HD23	1.97	0.47
1:D:54:TYR:CE2	1:D:55:LYS:HE2	2.50	0.46
1:D:132:PHE:CE1	1:E:109:ASN:HB2	2.50	0.46
1:C:64:HIS:HE1	1:C:88:ARG:HH11	1.63	0.46
1:B:176:ALA:O	1:B:177:ASN:HB2	2.15	0.46
1:C:136:ASP:OD2	1:C:140:LYS:HE3	2.15	0.46
1:B:14:GLN:O	1:B:60:LEU:HD11	2.16	0.46
1:A:139:ASN:O	1:A:143:GLN:HG3	2.16	0.46
1:D:102:ARG:HD2	1:D:216:ASP:OD1	2.14	0.46
1:B:5:LEU:HD22	1:B:81:ASN:ND2	2.31	0.46
1:E:107:ILE:HD13	1:E:138:LEU:HB3	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:22:VAL:O	1:C:63:SER:HA	2.15	0.46
1:F:176:ALA:O	1:F:177:ASN:HB2	2.16	0.46
1:E:224:HIS:CE1	1:E:228:ASN:HD21	2.33	0.46
1:E:57:GLN:OE1	1:E:242:THR:HG22	2.15	0.46
1:F:22:VAL:O	1:F:63:SER:HA	2.15	0.45
1:C:5:LEU:HD13	1:C:77:GLU:HB3	1.99	0.45
1:C:101:LYS:HD2	1:C:101:LYS:N	2.31	0.45
1:D:100:ILE:HG22	1:D:205:VAL:HG21	1.98	0.45
1:F:5:LEU:HD13	1:F:77:GLU:HB3	1.98	0.45
1:E:46:GLU:HB3	1:F:46:GLU:CB	2.43	0.45
3:C:303:MTI:HS3	1:D:73:VAL:HG21	1.98	0.45
1:D:64:HIS:HE1	1:D:88:ARG:HH11	1.63	0.45
1:E:238:ALA:O	1:E:242:THR:HG23	2.16	0.45
1:E:211:LYS:HB3	1:E:214:GLU:HG2	1.99	0.45
1:A:132:PHE:CZ	1:A:133:ASP:OD2	2.71	0.44
1:A:57:GLN:HE21	1:A:57:GLN:HA	1.82	0.44
1:C:11:SER:HB2	1:C:13:GLU:OE2	2.18	0.44
1:C:6:ARG:HD3	1:D:217:PHE:CZ	2.52	0.44
1:D:51:GLU:CD	1:D:58:LYS:HE3	2.37	0.44
1:F:13:GLU:CD	1:F:13:GLU:H	2.20	0.44
1:A:92:CYS:SG	1:A:180:VAL:HG11	2.58	0.44
1:E:114:GLU:HB3	1:E:157:SER:HA	2.00	0.44
1:C:163:ASN:H	1:C:163:ASN:ND2	2.16	0.44
1:D:20:LEU:O	1:D:61:CYS:HA	2.18	0.44
1:C:20:LEU:O	1:C:61:CYS:HA	2.18	0.43
1:F:227:GLU:HG2	1:F:231:LYS:HE3	1.99	0.43
1:F:64:HIS:CD2	1:F:65:GLY:O	2.67	0.43
1:F:51:GLU:HA	1:F:60:LEU:HD23	1.99	0.43
1:E:46:GLU:CB	1:F:46:GLU:HB3	2.46	0.43
1:D:45:ARG:HB3	1:D:46:GLU:OE2	2.18	0.43
1:D:86:ILE:O	1:D:199:THR:HA	2.19	0.43
1:A:5:LEU:HG	1:A:10:ILE:O	2.19	0.43
1:B:22:VAL:HG12	1:B:23:GLY:N	2.34	0.43
1:B:116:ARG:HE	4:B:501:IPA:H13	1.84	0.43
1:A:140:LYS:O	1:A:143:GLN:HB2	2.19	0.43
1:E:228:ASN:O	1:E:232:ILE:HG13	2.18	0.43
1:D:112:VAL:HG11	1:D:173:TYR:CZ	2.54	0.43
1:E:15:ILE:HA	1:E:60:LEU:HD11	2.00	0.43
3:E:305:MTI:H3	5:E:884:HOH:O	2.18	0.43
1:F:57:GLN:HG3	1:F:241:ALA:HB3	2.01	0.43
1:D:35:CYS:HB3	1:D:53:HIS:O	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:131:ASP:OD1	1:D:133:ASP:HB2	2.19	0.42
1:A:176:ALA:O	1:A:177:ASN:HB2	2.19	0.42
1:E:89:ALA:HA	1:E:202:ILE:O	2.18	0.42
1:E:119:HIS:HD2	5:E:833:HOH:O	2.01	0.42
1:B:64:HIS:HE1	1:B:88:ARG:HH11	1.68	0.42
1:F:160:TYR:CD1	3:F:306:MTI:HS1	2.54	0.42
1:E:28:VAL:HG21	1:E:48:LYS:HG2	2.01	0.42
1:E:75:PHE:CE1	1:E:188:LEU:HB2	2.54	0.42
1:A:75:PHE:CE1	1:A:188:LEU:HB2	2.54	0.42
1:C:211:LYS:HE3	1:C:216:ASP:OD2	2.19	0.42
1:F:41:LEU:HD11	1:F:51:GLU:HB2	2.01	0.41
1:E:147:VAL:HG11	1:E:232:ILE:HD11	2.01	0.41
1:C:8:LEU:HG	1:C:77:GLU:OE1	2.20	0.41
1:A:101:LYS:HB3	2:A:413:SO4:O3	2.21	0.41
1:F:24:ASP:HA	1:F:25:PRO:HD2	1.96	0.41
1:C:100:ILE:CG2	1:C:205:VAL:HG21	2.50	0.41
1:A:195:ARG:O	1:A:196:LYS:HB2	2.20	0.41
1:A:132:PHE:CG	1:A:133:ASP:N	2.88	0.41
1:A:34:VAL:O	1:A:34:VAL:HG12	2.21	0.41
1:C:64:HIS:CE1	1:C:88:ARG:HH11	2.38	0.41
1:F:119:HIS:HD2	5:F:836:HOH:O	2.04	0.41
1:E:225:GLN:OE1	1:E:225:GLN:HA	2.20	0.41
1:B:30:LYS:NZ	1:B:30:LYS:HB3	2.36	0.41
1:E:5:LEU:HG	1:E:10:ILE:O	2.21	0.41
1:D:193:THR:HG21	5:D:649:HOH:O	2.21	0.41
1:A:119:HIS:HD2	5:A:834:HOH:O	2.02	0.41
1:F:57:GLN:CG	1:F:241:ALA:HB3	2.50	0.41
3:B:302:MTI:N3	3:B:302:MTI:H1	2.36	0.41
1:C:17:PRO:HG2	5:C:667:HOH:O	2.21	0.41
1:B:23:GLY:HA2	1:B:64:HIS:CD2	2.56	0.40
1:B:66:VAL:HG11	3:B:302:MTI:HS3	2.03	0.40
1:B:204:ILE:HG21	1:B:221:LEU:HD13	2.04	0.40
1:C:131:ASP:OD1	1:C:133:ASP:HB2	2.21	0.40
1:C:112:VAL:HB	1:C:155:VAL:HA	2.03	0.40
1:F:226:LEU:O	1:F:230:ILE:HG13	2.21	0.40
1:F:222:VAL:HB	1:F:225:GLN:HB2	2.03	0.40
1:F:39:VAL:O	1:F:41:LEU:HD12	2.21	0.40
1:B:29:ASP:O	1:B:32:LYS:HG2	2.21	0.40
1:D:111:ALA:O	1:D:128:ALA:HB3	2.22	0.40
1:D:27:ARG:NH1	1:D:226:LEU:HD21	2.37	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	241/276 (87%)	231 (96%)	10 (4%)	0	100	100
1	B	241/276 (87%)	234 (97%)	6 (2%)	1 (0%)	39	32
1	C	241/276 (87%)	228 (95%)	12 (5%)	1 (0%)	39	32
1	D	241/276 (87%)	230 (95%)	11 (5%)	0	100	100
1	E	241/276 (87%)	234 (97%)	6 (2%)	1 (0%)	39	32
1	F	241/276 (87%)	233 (97%)	8 (3%)	0	100	100
All	All	1446/1656 (87%)	1390 (96%)	53 (4%)	3 (0%)	52	47

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	223	PRO
1	E	223	PRO
1	C	223	PRO

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	206/235 (88%)	203 (98%)	3 (2%)	72	74
1	B	206/235 (88%)	201 (98%)	5 (2%)	57	56
1	C	206/235 (88%)	202 (98%)	4 (2%)	65	66
1	D	206/235 (88%)	203 (98%)	3 (2%)	72	74

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	E	206/235 (88%)	201 (98%)	5 (2%)	57	56
1	F	206/235 (88%)	201 (98%)	5 (2%)	57	56
All	All	1236/1410 (88%)	1211 (98%)	25 (2%)	63	64

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	121	LEU
1	A	132	PHE
1	A	182	GLU
1	B	3	ASN
1	B	30	LYS
1	B	57	GLN
1	B	163	ASN
1	B	182	GLU
1	C	101	LYS
1	C	121	LEU
1	C	163	ASN
1	C	182	GLU
1	D	121	LEU
1	D	163	ASN
1	D	182	GLU
1	E	102	ARG
1	E	121	LEU
1	E	163	ASN
1	E	182	GLU
1	E	218	ASP
1	F	13	GLU
1	F	101	LYS
1	F	121	LEU
1	F	180	VAL
1	F	182	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (30) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	53	HIS
1	A	64	HIS
1	A	80	GLN
1	A	119	HIS

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Mol	Chain	Res	Type
1	A	151	ASN
1	B	64	HIS
1	B	119	HIS
1	B	139	ASN
1	C	44	ASN
1	C	53	HIS
1	C	57	GLN
1	C	64	HIS
1	C	119	HIS
1	C	228	ASN
1	D	44	ASN
1	D	57	GLN
1	D	64	HIS
1	D	119	HIS
1	D	151	ASN
1	D	163	ASN
1	E	44	ASN
1	E	53	HIS
1	E	64	HIS
1	E	119	HIS
1	E	163	ASN
1	E	228	ASN
1	F	53	HIS
1	F	64	HIS
1	F	119	HIS
1	F	151	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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

37 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
3	MTI	A	301	-	21,22,22	2.70	10 (47%)	11,32,32	3.66	4 (36%)
2	SO4	A	401	-	4,4,4	3.22	2 (50%)	6,6,6	0.95	0
2	SO4	A	413	-	4,4,4	3.25	2 (50%)	6,6,6	0.96	0
2	SO4	A	414	-	4,4,4	3.26	2 (50%)	6,6,6	0.95	0
2	SO4	A	418	-	4,4,4	3.28	2 (50%)	6,6,6	0.97	0
4	IPA	A	503	-	3,3,3	0.27	0	3,3,3	0.36	0
3	MTI	B	302	-	21,22,22	2.67	10 (47%)	11,32,32	3.76	4 (36%)
2	SO4	B	402	-	4,4,4	3.20	2 (50%)	6,6,6	1.01	0
2	SO4	B	407	-	4,4,4	3.16	2 (50%)	6,6,6	1.01	0
2	SO4	B	410	-	4,4,4	3.26	2 (50%)	6,6,6	0.97	0
2	SO4	B	417	-	4,4,4	3.24	2 (50%)	6,6,6	0.96	0
4	IPA	B	501	-	3,3,3	0.26	0	3,3,3	0.34	0
4	IPA	B	502	-	3,3,3	0.29	0	3,3,3	0.34	0
4	IPA	B	508	-	3,3,3	0.32	0	3,3,3	0.35	0
3	MTI	C	303	-	21,22,22	2.57	10 (47%)	11,32,32	3.85	4 (36%)
2	SO4	C	403	-	4,4,4	3.21	2 (50%)	6,6,6	0.94	0
2	SO4	C	416	-	4,4,4	3.21	2 (50%)	6,6,6	0.91	0
2	SO4	C	422	-	4,4,4	3.27	2 (50%)	6,6,6	0.94	0
3	MTI	D	304	-	21,22,22	2.58	9 (42%)	11,32,32	3.81	4 (36%)
2	SO4	D	404	-	4,4,4	3.25	2 (50%)	6,6,6	0.94	0
2	SO4	D	408	-	4,4,4	3.08	2 (50%)	6,6,6	0.95	0
2	SO4	D	412	-	4,4,4	3.26	2 (50%)	6,6,6	0.95	0
2	SO4	D	421	-	4,4,4	3.32	2 (50%)	6,6,6	0.96	0
4	IPA	D	507	-	3,3,3	0.26	0	3,3,3	0.33	0
4	IPA	D	509	-	3,3,3	0.28	0	3,3,3	0.33	0
3	MTI	E	305	-	21,22,22	2.63	10 (47%)	11,32,32	3.70	4 (36%)
2	SO4	E	405	-	4,4,4	3.23	2 (50%)	6,6,6	0.97	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	SO4	E	415	-	4,4,4	3.16	2 (50%)	6,6,6	0.95	0
2	SO4	E	420	-	4,4,4	3.27	2 (50%)	6,6,6	0.97	0
4	IPA	E	506	-	3,3,3	0.21	0	3,3,3	0.32	0
3	MTI	F	306	-	21,22,22	2.57	10 (47%)	11,32,32	3.76	4 (36%)
2	SO4	F	406	-	4,4,4	3.17	2 (50%)	6,6,6	0.91	0
2	SO4	F	409	-	4,4,4	3.20	2 (50%)	6,6,6	0.98	0
2	SO4	F	411	-	4,4,4	3.17	2 (50%)	6,6,6	0.95	0
2	SO4	F	419	-	4,4,4	3.28	2 (50%)	6,6,6	0.99	0
4	IPA	F	504	-	3,3,3	0.26	0	3,3,3	0.31	0
4	IPA	F	505	-	3,3,3	0.29	0	3,3,3	0.34	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	MTI	A	301	-	-	0/3/23/23	0/3/3/3
2	SO4	A	401	-	-	0/0/0/0	0/0/0/0
2	SO4	A	413	-	-	0/0/0/0	0/0/0/0
2	SO4	A	414	-	-	0/0/0/0	0/0/0/0
2	SO4	A	418	-	-	0/0/0/0	0/0/0/0
4	IPA	A	503	-	-	0/0/0/0	0/0/0/0
3	MTI	B	302	-	-	0/3/23/23	0/3/3/3
2	SO4	B	402	-	-	0/0/0/0	0/0/0/0
2	SO4	B	407	-	-	0/0/0/0	0/0/0/0
2	SO4	B	410	-	-	0/0/0/0	0/0/0/0
2	SO4	B	417	-	-	0/0/0/0	0/0/0/0
4	IPA	B	501	-	-	0/0/0/0	0/0/0/0
4	IPA	B	502	-	-	0/0/0/0	0/0/0/0
4	IPA	B	508	-	-	0/0/0/0	0/0/0/0
3	MTI	C	303	-	-	0/3/23/23	0/3/3/3
2	SO4	C	403	-	-	0/0/0/0	0/0/0/0
2	SO4	C	416	-	-	0/0/0/0	0/0/0/0
2	SO4	C	422	-	-	0/0/0/0	0/0/0/0
3	MTI	D	304	-	-	0/3/23/23	0/3/3/3
2	SO4	D	404	-	-	0/0/0/0	0/0/0/0
2	SO4	D	408	-	-	0/0/0/0	0/0/0/0
2	SO4	D	412	-	-	0/0/0/0	0/0/0/0
2	SO4	D	421	-	-	0/0/0/0	0/0/0/0
4	IPA	D	507	-	-	0/0/0/0	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	IPA	D	509	-	-	0/0/0/0	0/0/0/0
3	MTI	E	305	-	-	0/3/23/23	0/3/3/3
2	SO4	E	405	-	-	0/0/0/0	0/0/0/0
2	SO4	E	415	-	-	0/0/0/0	0/0/0/0
2	SO4	E	420	-	-	0/0/0/0	0/0/0/0
4	IPA	E	506	-	-	0/0/0/0	0/0/0/0
3	MTI	F	306	-	-	0/3/23/23	0/3/3/3
2	SO4	F	406	-	-	0/0/0/0	0/0/0/0
2	SO4	F	409	-	-	0/0/0/0	0/0/0/0
2	SO4	F	411	-	-	0/0/0/0	0/0/0/0
2	SO4	F	419	-	-	0/0/0/0	0/0/0/0
4	IPA	F	504	-	-	0/0/0/0	0/0/0/0
4	IPA	F	505	-	-	0/0/0/0	0/0/0/0

All (103) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	412	SO4	O3-S	-4.60	1.30	1.47
2	B	402	SO4	O3-S	-4.60	1.30	1.47
2	A	414	SO4	O3-S	-4.59	1.30	1.47
2	C	422	SO4	O3-S	-4.57	1.31	1.47
2	F	411	SO4	O3-S	-4.56	1.31	1.47
2	D	404	SO4	O3-S	-4.55	1.31	1.47
2	D	421	SO4	O3-S	-4.54	1.31	1.47
2	A	401	SO4	O3-S	-4.54	1.31	1.47
2	C	403	SO4	O3-S	-4.53	1.31	1.47
2	B	417	SO4	O3-S	-4.52	1.31	1.47
2	B	410	SO4	O3-S	-4.52	1.31	1.47
2	F	419	SO4	O3-S	-4.50	1.31	1.47
2	E	420	SO4	O3-S	-4.49	1.31	1.47
2	F	406	SO4	O3-S	-4.49	1.31	1.47
2	A	418	SO4	O3-S	-4.48	1.31	1.47
2	E	415	SO4	O3-S	-4.46	1.31	1.47
2	A	413	SO4	O3-S	-4.44	1.31	1.47
2	F	409	SO4	O3-S	-4.44	1.31	1.47
2	E	405	SO4	O3-S	-4.44	1.31	1.47
2	C	416	SO4	O3-S	-4.31	1.31	1.47
2	B	407	SO4	O3-S	-4.22	1.32	1.47
2	D	408	SO4	O3-S	-4.15	1.32	1.47
3	A	301	MTI	C8-N7	-2.63	1.31	1.36
3	C	303	MTI	C8-N7	-2.48	1.31	1.36
3	D	304	MTI	C8-N7	-2.45	1.31	1.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	302	MTI	C8-N7	-2.31	1.31	1.36
3	E	305	MTI	C8-N7	-2.30	1.31	1.36
3	F	306	MTI	C8-N7	-2.21	1.32	1.36
3	F	306	MTI	O3'-C3'	2.16	1.48	1.43
3	F	306	MTI	C6-C5	2.22	1.45	1.41
3	C	303	MTI	C6-C5	2.25	1.45	1.41
3	A	301	MTI	C6-C5	2.25	1.45	1.41
3	C	303	MTI	C3'-C2'	2.34	1.59	1.53
3	F	306	MTI	C3'-C2'	2.39	1.59	1.53
3	F	306	MTI	C2'-C1'	2.40	1.57	1.53
3	C	303	MTI	C2'-C1'	2.44	1.57	1.53
3	D	304	MTI	C3'-C2'	2.46	1.60	1.53
3	A	301	MTI	C3'-C2'	2.53	1.60	1.53
3	D	304	MTI	O3'-C3'	2.54	1.49	1.43
3	C	303	MTI	O3'-C3'	2.57	1.49	1.43
3	A	301	MTI	C2'-C1'	2.58	1.57	1.53
3	E	305	MTI	C3'-C2'	2.59	1.60	1.53
3	E	305	MTI	C2'-C1'	2.60	1.57	1.53
3	B	302	MTI	C3'-C2'	2.64	1.60	1.53
3	B	302	MTI	C6-C5	2.65	1.46	1.41
3	A	301	MTI	O3'-C3'	2.67	1.49	1.43
3	E	305	MTI	C6-C5	2.68	1.46	1.41
3	D	304	MTI	C2'-C1'	2.75	1.57	1.53
3	E	305	MTI	O3'-C3'	2.82	1.49	1.43
3	B	302	MTI	C2'-C1'	2.88	1.57	1.53
3	B	302	MTI	O3'-C3'	2.90	1.49	1.43
3	E	305	MTI	C4-N3	3.35	1.41	1.37
3	B	302	MTI	C2-N3	3.74	1.38	1.32
3	D	304	MTI	C2-N3	3.75	1.38	1.32
3	B	302	MTI	C4-N3	3.82	1.42	1.37
3	A	301	MTI	C2-N3	3.85	1.39	1.32
3	A	301	MTI	C4-N3	3.87	1.42	1.37
3	E	305	MTI	C2-N3	3.88	1.39	1.32
3	F	306	MTI	C2-N3	3.90	1.39	1.32
3	F	306	MTI	C4-N3	3.93	1.42	1.37
3	A	301	MTI	C2-N1	3.97	1.41	1.33
3	C	303	MTI	C2-N3	3.98	1.39	1.32
3	D	304	MTI	C4-N3	4.02	1.42	1.37
3	C	303	MTI	C4-N3	4.05	1.42	1.37
3	C	303	MTI	C2-N1	4.09	1.41	1.33
3	D	304	MTI	C2-N1	4.14	1.41	1.33
3	E	305	MTI	C2-N1	4.17	1.41	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	302	MTI	C2-N1	4.20	1.41	1.33
3	F	306	MTI	C2-N1	4.20	1.41	1.33
2	F	411	SO4	O1-S	4.31	1.61	1.47
2	B	402	SO4	O1-S	4.37	1.62	1.47
2	F	406	SO4	O1-S	4.40	1.62	1.47
2	E	415	SO4	O1-S	4.44	1.62	1.47
2	C	403	SO4	O1-S	4.46	1.62	1.47
2	A	401	SO4	O1-S	4.48	1.62	1.47
2	D	408	SO4	O1-S	4.50	1.62	1.47
2	D	412	SO4	O1-S	4.52	1.62	1.47
3	C	303	MTI	C9-C1'	4.54	1.56	1.52
2	D	404	SO4	O1-S	4.55	1.62	1.47
2	F	409	SO4	O1-S	4.55	1.62	1.47
2	B	417	SO4	O1-S	4.56	1.62	1.47
2	E	405	SO4	O1-S	4.57	1.62	1.47
2	C	422	SO4	O1-S	4.57	1.62	1.47
2	A	414	SO4	O1-S	4.59	1.62	1.47
2	B	410	SO4	O1-S	4.63	1.63	1.47
2	B	407	SO4	O1-S	4.65	1.63	1.47
2	A	413	SO4	O1-S	4.65	1.63	1.47
2	F	419	SO4	O1-S	4.68	1.63	1.47
2	E	420	SO4	O1-S	4.68	1.63	1.47
2	C	416	SO4	O1-S	4.69	1.63	1.47
2	A	418	SO4	O1-S	4.72	1.63	1.47
3	D	304	MTI	C9-C1'	4.72	1.56	1.52
2	D	421	SO4	O1-S	4.73	1.63	1.47
3	B	302	MTI	C9-C1'	5.02	1.56	1.52
3	F	306	MTI	C9-C1'	5.31	1.56	1.52
3	E	305	MTI	C9-C1'	5.39	1.56	1.52
3	E	305	MTI	C6-N1	5.41	1.43	1.33
3	F	306	MTI	C6-N1	5.46	1.43	1.33
3	D	304	MTI	C6-N1	5.66	1.43	1.33
3	C	303	MTI	C6-N1	5.70	1.43	1.33
3	A	301	MTI	C6-N1	5.72	1.43	1.33
3	B	302	MTI	C6-N1	5.74	1.43	1.33
3	A	301	MTI	C9-C1'	5.89	1.57	1.52

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	303	MTI	N3-C2-N1	-9.79	121.40	128.89
3	D	304	MTI	N3-C2-N1	-9.69	121.47	128.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	306	MTI	N3-C2-N1	-9.62	121.53	128.89
3	B	302	MTI	N3-C2-N1	-9.58	121.56	128.89
3	A	301	MTI	N3-C2-N1	-9.41	121.69	128.89
3	E	305	MTI	N3-C2-N1	-9.30	121.78	128.89
3	D	304	MTI	C9-C1'-N4'	-4.53	106.64	113.38
3	C	303	MTI	C9-C1'-N4'	-4.34	106.92	113.38
3	B	302	MTI	C9-C1'-N4'	-4.31	106.97	113.38
3	E	305	MTI	C9-C1'-N4'	-4.25	107.05	113.38
3	F	306	MTI	C9-C1'-N4'	-4.07	107.33	113.38
3	A	301	MTI	C9-C1'-N4'	-3.76	107.78	113.38
3	B	302	MTI	C9-C8-N7	3.42	115.31	108.79
3	D	304	MTI	C9-C8-N7	3.50	115.48	108.79
3	F	306	MTI	C9-C8-N7	3.50	115.48	108.79
3	E	305	MTI	C9-C8-N7	3.53	115.52	108.79
3	C	303	MTI	C9-C8-N7	3.58	115.62	108.79
3	A	301	MTI	C9-C8-N7	3.58	115.62	108.79
3	A	301	MTI	C2-N1-C6	5.42	124.25	116.04
3	E	305	MTI	C2-N1-C6	5.45	124.30	116.04
3	B	302	MTI	C2-N1-C6	5.52	124.40	116.04
3	D	304	MTI	C2-N1-C6	5.53	124.42	116.04
3	F	306	MTI	C2-N1-C6	5.59	124.50	116.04
3	C	303	MTI	C2-N1-C6	5.71	124.69	116.04

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

9 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	413	SO4	1	0
3	B	302	MTI	2	0
4	B	501	IPA	1	0
3	C	303	MTI	1	0
2	D	412	SO4	1	0
4	D	509	IPA	1	0
3	E	305	MTI	2	0
3	F	306	MTI	1	0
4	F	504	IPA	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	243/276 (88%)	0.28	14 (5%)	26 28	14, 29, 47, 57	0
1	B	243/276 (88%)	0.64	24 (9%)	9 10	14, 33, 51, 62	0
1	C	243/276 (88%)	0.53	17 (6%)	19 21	12, 29, 48, 62	0
1	D	243/276 (88%)	0.42	12 (4%)	33 35	11, 28, 47, 58	0
1	E	243/276 (88%)	0.26	16 (6%)	22 23	10, 24, 45, 63	0
1	F	243/276 (88%)	0.25	5 (2%)	67 68	11, 24, 41, 51	0
All	All	1458/1656 (88%)	0.39	88 (6%)	25 27	10, 27, 48, 63	0

All (88) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	3	ASN	5.5
1	C	3	ASN	5.4
1	A	132	PHE	5.4
1	E	213	ASP	5.2
1	C	215	GLY	5.0
1	C	219	ASN	5.0
1	E	219	ASN	4.9
1	C	217	PHE	4.6
1	E	3	ASN	4.5
1	D	3	ASN	4.5
1	B	214	GLU	4.5
1	C	221	LEU	4.5
1	A	3	ASN	4.4
1	E	217	PHE	4.3
1	B	132	PHE	4.2
1	D	4	LEU	4.2
1	F	3	ASN	4.1
1	C	218	ASP	4.1
1	B	219	ASN	4.0

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Mol	Chain	Res	Type	RSRZ
1	B	218	ASP	4.0
1	B	215	GLY	4.0
1	B	222	VAL	3.9
1	B	220	ASN	3.9
1	B	146	ASN	3.8
1	D	245	ALA	3.7
1	C	220	ASN	3.7
1	B	213	ASP	3.7
1	B	217	PHE	3.6
1	B	221	LEU	3.6
1	B	216	ASP	3.6
1	C	214	GLU	3.5
1	E	216	ASP	3.5
1	B	212	TRP	3.5
1	C	222	VAL	3.4
1	A	4	LEU	3.4
1	D	34	VAL	3.4
1	C	34	VAL	3.4
1	B	224	HIS	3.3
1	A	213	ASP	3.3
1	D	213	ASP	3.2
1	E	214	GLU	3.2
1	A	215	GLY	3.2
1	A	146	ASN	3.0
1	F	219	ASN	3.0
1	C	146	ASN	3.0
1	F	146	ASN	2.9
1	C	213	ASP	2.9
1	A	13	GLU	2.9
1	E	99	LEU	2.9
1	A	15	ILE	2.8
1	B	4	LEU	2.8
1	C	39	VAL	2.8
1	B	58	LYS	2.7
1	E	245	ALA	2.7
1	E	243	LYS	2.6
1	E	218	ASP	2.6
1	B	33	VAL	2.6
1	C	224	HIS	2.6
1	A	218	ASP	2.5
1	B	13	GLU	2.5
1	B	145	LEU	2.5

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Mol	Chain	Res	Type	RSRZ
1	C	132	PHE	2.5
1	D	145	LEU	2.5
1	A	39	VAL	2.5
1	B	211	LYS	2.5
1	B	12	LYS	2.4
1	D	12	LYS	2.3
1	E	215	GLY	2.3
1	C	216	ASP	2.3
1	D	144	GLU	2.3
1	F	224	HIS	2.3
1	D	132	PHE	2.3
1	D	215	GLY	2.2
1	D	38	TYR	2.2
1	E	132	PHE	2.2
1	E	98	ASP	2.2
1	B	238	ALA	2.2
1	C	4	LEU	2.1
1	E	220	ASN	2.1
1	D	143	GLN	2.1
1	E	13	GLU	2.1
1	A	219	ASN	2.1
1	A	56	GLY	2.0
1	A	245	ALA	2.0
1	B	245	ALA	2.0
1	A	38	TYR	2.0
1	E	12	LYS	2.0
1	F	242	THR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains.

The B-factors column lists the minimum, median, 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	LLDF	B-factors(Å ²)	Q<0.9
4	IPA	A	503	4/4	0.85	0.15	15.28	23,24,25,26	0
4	IPA	B	502	4/4	0.82	0.20	13.02	21,21,22,24	0
4	IPA	F	505	4/4	0.66	0.28	11.81	24,26,27,28	0
4	IPA	B	508	4/4	0.73	0.23	7.09	24,24,25,26	0
4	IPA	E	506	4/4	0.83	0.20	6.72	23,24,26,26	0
4	IPA	D	507	4/4	0.88	0.16	3.92	19,21,22,22	0
3	MTI	D	304	20/20	0.95	0.15	0.50	19,20,33,36	0
3	MTI	F	306	20/20	0.95	0.14	0.31	18,20,32,34	0
3	MTI	C	303	20/20	0.92	0.15	0.23	25,30,41,44	0
3	MTI	E	305	20/20	0.93	0.14	0.19	26,28,37,42	0
3	MTI	A	301	20/20	0.94	0.12	0.17	19,22,34,36	0
2	SO4	F	411	5/5	0.96	0.13	0.09	40,41,44,44	0
2	SO4	E	420	5/5	0.95	0.14	0.07	44,46,47,48	0
2	SO4	D	412	5/5	0.95	0.16	0.02	52,52,54,54	0
2	SO4	D	421	5/5	0.95	0.19	-0.01	56,57,58,58	0
2	SO4	F	419	5/5	0.92	0.18	-0.01	48,51,52,52	0
2	SO4	B	417	5/5	0.93	0.16	-0.03	69,70,70,71	0
3	MTI	B	302	20/20	0.92	0.13	-0.07	27,34,46,48	0
2	SO4	C	403	5/5	0.97	0.13	-0.18	30,31,34,35	0
2	SO4	A	418	5/5	0.94	0.16	-0.39	59,59,59,62	0
2	SO4	A	401	5/5	0.97	0.12	-0.57	32,32,35,35	0
2	SO4	C	422	5/5	0.95	0.10	-0.70	44,45,48,48	0
2	SO4	B	407	5/5	0.98	0.10	-0.72	25,27,29,31	0
2	SO4	E	415	5/5	0.98	0.09	-0.90	20,22,26,27	0
2	SO4	E	405	5/5	0.98	0.11	-1.04	26,27,27,28	0
2	SO4	D	404	5/5	0.97	0.12	-1.06	29,29,30,30	0
2	SO4	F	406	5/5	0.96	0.10	-1.40	23,24,26,27	0
2	SO4	C	416	5/5	0.98	0.07	-1.45	25,26,27,29	0
2	SO4	B	402	5/5	0.96	0.11	-1.57	40,40,42,42	0
2	SO4	D	408	5/5	0.97	0.09	-1.91	18,25,25,27	0
2	SO4	A	414	5/5	0.98	0.08	-2.06	22,26,28,29	0
2	SO4	F	409	5/5	0.98	0.06	-2.61	20,23,29,30	0
2	SO4	B	410	5/5	0.90	0.25	-	69,70,70,70	0
4	IPA	D	509	4/4	0.62	0.27	-	24,24,24,27	0
4	IPA	F	504	4/4	0.79	0.25	-	15,18,21,23	0
2	SO4	A	413	5/5	0.91	0.33	-	65,65,66,66	0
4	IPA	B	501	4/4	0.49	0.30	-	29,30,30,31	0

6.5 Other polymers [i](#)

There are no such residues in this entry.