



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 31, 2017 – 05:21 PM EDT

PDB ID : 3MCM
Title : Crystal Structure of the 6-hydroxymethyl-7,8-dihydropterin pyrophosphokinase dihydropteroate synthase bifunctional enzyme from *Francisella tularensis*
Authors : Pemble IV, C.W.; Mehta, P.K.; Mehra, S.; Li, Z.; Lee, R.E.; White, S.W.
Deposited on : unknown
Resolution : 2.20 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
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20030345
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
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)	:	rb-20030345

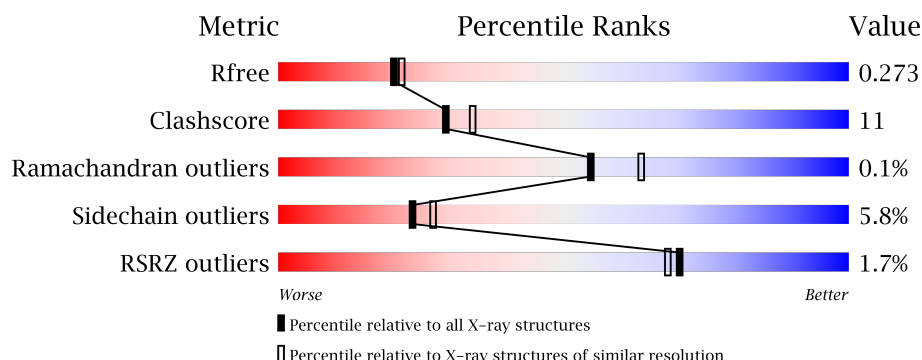
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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}	100719	4002 (2.20-2.20)
Clashscore	112137	4730 (2.20-2.20)
Ramachandran outliers	110173	4656 (2.20-2.20)
Sidechain outliers	110143	4657 (2.20-2.20)
RSRZ outliers	101464	4033 (2.20-2.20)

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	442	<div> <div>2%</div> <div>60%</div> <div>20%</div> <div>17%</div> </div>
1	B	442	<div> <div>%</div> <div>65%</div> <div>19%</div> <div>15%</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6176 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 2-amino-4-hydroxy-6-hydroxymethyldihydropteridine pyrophosphokinase/dihydropteroate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	368	Total	C	N	O	S	0	1	0
			2979	1907	517	549	6			
1	B	377	Total	C	N	O	S	0	2	0
			3064	1963	530	565	6			

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	EXPRESSION TAG	UNP Q2A2W3
A	-18	GLY	-	EXPRESSION TAG	UNP Q2A2W3
A	-17	SER	-	EXPRESSION TAG	UNP Q2A2W3
A	-16	SER	-	EXPRESSION TAG	UNP Q2A2W3
A	-15	HIS	-	EXPRESSION TAG	UNP Q2A2W3
A	-14	HIS	-	EXPRESSION TAG	UNP Q2A2W3
A	-13	HIS	-	EXPRESSION TAG	UNP Q2A2W3
A	-12	HIS	-	EXPRESSION TAG	UNP Q2A2W3
A	-11	HIS	-	EXPRESSION TAG	UNP Q2A2W3
A	-10	HIS	-	EXPRESSION TAG	UNP Q2A2W3
A	-9	SER	-	EXPRESSION TAG	UNP Q2A2W3
A	-8	SER	-	EXPRESSION TAG	UNP Q2A2W3
A	-7	GLY	-	EXPRESSION TAG	UNP Q2A2W3
A	-6	LEU	-	EXPRESSION TAG	UNP Q2A2W3
A	-5	VAL	-	EXPRESSION TAG	UNP Q2A2W3
A	-4	PRO	-	EXPRESSION TAG	UNP Q2A2W3
A	-3	ARG	-	EXPRESSION TAG	UNP Q2A2W3
A	-2	GLY	-	EXPRESSION TAG	UNP Q2A2W3
A	-1	SER	-	EXPRESSION TAG	UNP Q2A2W3
A	0	HIS	-	EXPRESSION TAG	UNP Q2A2W3
A	1	MET	-	EXPRESSION TAG	UNP Q2A2W3
A	2	VAL	-	EXPRESSION TAG	UNP Q2A2W3
B	-19	MET	-	EXPRESSION TAG	UNP Q2A2W3
B	-18	GLY	-	EXPRESSION TAG	UNP Q2A2W3

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-17	SER	-	EXPRESSION TAG	UNP Q2A2W3
B	-16	SER	-	EXPRESSION TAG	UNP Q2A2W3
B	-15	HIS	-	EXPRESSION TAG	UNP Q2A2W3
B	-14	HIS	-	EXPRESSION TAG	UNP Q2A2W3
B	-13	HIS	-	EXPRESSION TAG	UNP Q2A2W3
B	-12	HIS	-	EXPRESSION TAG	UNP Q2A2W3
B	-11	HIS	-	EXPRESSION TAG	UNP Q2A2W3
B	-10	HIS	-	EXPRESSION TAG	UNP Q2A2W3
B	-9	SER	-	EXPRESSION TAG	UNP Q2A2W3
B	-8	SER	-	EXPRESSION TAG	UNP Q2A2W3
B	-7	GLY	-	EXPRESSION TAG	UNP Q2A2W3
B	-6	LEU	-	EXPRESSION TAG	UNP Q2A2W3
B	-5	VAL	-	EXPRESSION TAG	UNP Q2A2W3
B	-4	PRO	-	EXPRESSION TAG	UNP Q2A2W3
B	-3	ARG	-	EXPRESSION TAG	UNP Q2A2W3
B	-2	GLY	-	EXPRESSION TAG	UNP Q2A2W3
B	-1	SER	-	EXPRESSION TAG	UNP Q2A2W3
B	0	HIS	-	EXPRESSION TAG	UNP Q2A2W3
B	1	MET	-	EXPRESSION TAG	UNP Q2A2W3
B	2	VAL	-	EXPRESSION TAG	UNP Q2A2W3

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Mg 1 1	0	0

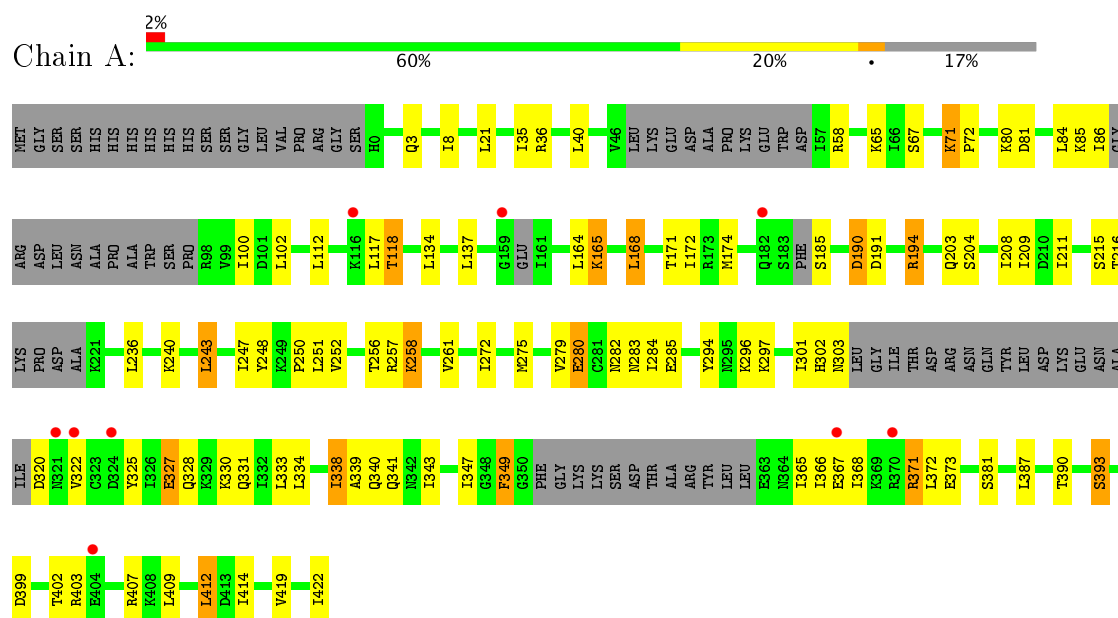
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	63	Total O 63 63	0	0
3	B	69	Total O 69 69	0	0

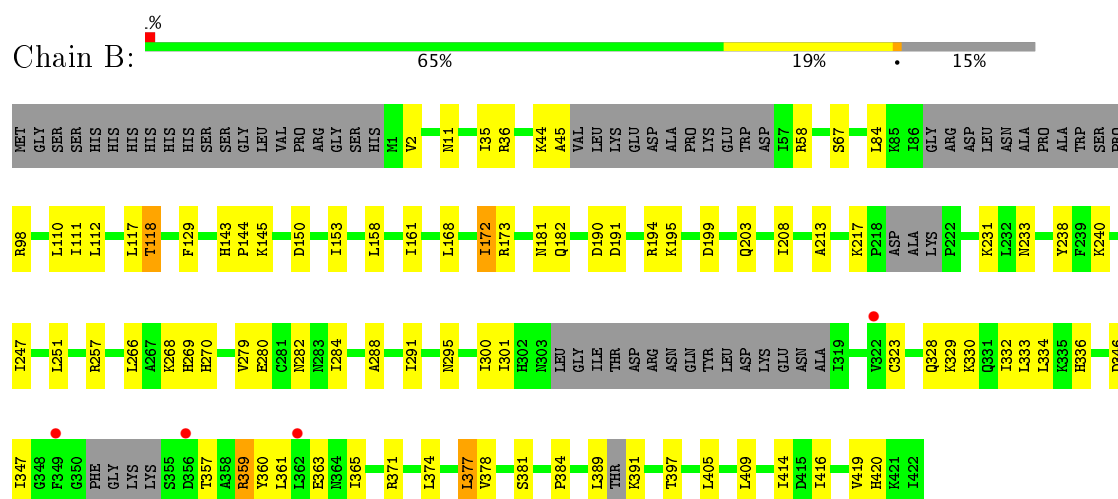
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: 2-amino-4-hydroxy-6-hydroxymethyldihydropteridine pyrophosphokinase/dihydrophosphate synthase



- Molecule 1: 2-amino-4-hydroxy-6-hydroxymethyldihydropteridine pyrophosphokinase/dihydrophosphate synthase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	42.82Å 57.78Å 105.13Å 91.13° 99.39° 111.64°	Depositor
Resolution (Å)	38.44 – 2.20 38.44 – 2.20	Depositor EDS
% Data completeness (in resolution range)	94.2 (38.44-2.20) 92.6 (38.44-2.20)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.21 (at 2.20Å)	Xtriage
Refinement program	PHENIX 1.5_2	Depositor
R, R_{free}	0.215 , 0.278 0.211 , 0.273	Depositor DCC
R_{free} test set	2225 reflections (5.05%)	DCC
Wilson B-factor (Å ²)	33.9	Xtriage
Anisotropy	0.062	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 31.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.085 for h,-h-k,-h-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6176	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.85% 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, 0.333 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: MG

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	0.40	0/3017	0.57	0/4055
1	B	0.40	0/3108	0.57	0/4177
All	All	0.40	0/6125	0.57	0/8232

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

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	2979	0	3119	70	0
1	B	3064	0	3209	63	0
2	A	1	0	0	0	0
3	A	63	0	0	1	0
3	B	69	0	0	4	0
All	All	6176	0	6328	133	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 11.

All (133) 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:B:330:LYS:HG3	1:B:374:LEU:HD11	1.56	0.88
1:A:284:ILE:HD11	1:A:333:LEU:HD21	1.54	0.87
1:A:366:ILE:HD11	1:A:412:LEU:HB3	1.57	0.86
1:B:143:HIS:CE1	1:B:145:LYS:HB2	2.22	0.75
1:A:165:LYS:HD3	1:A:165:LYS:N	2.05	0.72
1:A:257:ARG:HB2	1:A:280:GLU:HB2	1.75	0.68
1:A:340:GLN:NE2	1:A:373:GLU:HB3	2.09	0.67
1:A:372:LEU:O	1:A:373:GLU:HB2	1.96	0.66
1:A:257:ARG:HD2	1:A:280:GLU:HG3	1.78	0.64
1:B:110:LEU:HD12	1:B:111:ILE:N	2.13	0.64
1:B:182:GLN:HE21	1:B:217:LYS:HE3	1.62	0.63
1:B:330:LYS:O	1:B:334:LEU:HG	1.99	0.63
1:A:338:ILE:HD11	1:A:343:ILE:HG13	1.81	0.63
1:A:340:GLN:HE22	1:A:373:GLU:HB3	1.63	0.62
1:A:194:ARG:N	1:A:194:ARG:HD3	2.14	0.62
1:A:204:SER:O	1:A:422:ILE:HG21	1.99	0.62
1:B:58:ARG:NH1	1:B:247:ILE:HG22	2.15	0.61
1:B:279:VAL:HG12	3:B:468:HOH:O	1.99	0.61
1:B:240:LYS:HE3	1:B:269:HIS:HD2	1.66	0.60
1:B:300:ILE:HG23	1:B:329:LYS:HE3	1.83	0.60
1:B:240:LYS:CE	1:B:269:HIS:HD2	2.15	0.60
1:A:275:MET:HB2	1:A:297:LYS:HB2	1.82	0.60
1:B:208:ILE:HG12	1:B:251:LEU:HB2	1.83	0.59
1:A:8:ILE:HG12	1:A:102:LEU:CD2	2.32	0.59
1:B:301:ILE:HG13	1:B:346:ASP:HB3	1.85	0.59
1:A:381:SER:HB2	1:A:419:VAL:HA	1.85	0.58
1:B:233:ASN:OD1	1:B:269:HIS:HE1	1.86	0.58
1:B:328:GLN:O	1:B:332:ILE:HD12	2.03	0.58
1:B:359:ARG:O	1:B:363:GLU:HG3	2.03	0.57
1:B:112:LEU:HD23	1:B:117:LEU:HD11	1.86	0.57
1:B:284:ILE:HD11	1:B:333:LEU:HD21	1.87	0.56
1:A:3:GLN:OE1	1:A:65:LYS:HD3	2.06	0.56
1:A:367:GLU:O	1:A:371:ARG:HG2	2.07	0.55
1:A:58:ARG:HD3	1:A:247:ILE:HA	1.88	0.55
1:A:8:ILE:HG12	1:A:102:LEU:HD22	1.89	0.54
1:A:208:ILE:HG12	1:A:251:LEU:HB2	1.88	0.54
1:A:112:LEU:O	1:A:118:THR:HA	2.07	0.54
1:A:302:HIS:CD2	1:A:322:VAL:HG13	2.44	0.53
1:B:377:LEU:C	1:B:377:LEU:HD12	2.29	0.53
1:A:256:THR:HG21	1:A:261:VAL:HG22	1.89	0.53
1:A:279:VAL:CG1	1:A:301:ILE:HD12	2.39	0.52
1:B:194[B]:ARG:NH2	3:B:467:HOH:O	2.42	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:266:LEU:O	1:B:270:HIS:HB3	2.10	0.52
1:B:117:LEU:C	1:B:117:LEU:HD12	2.30	0.52
1:A:81:ASP:O	1:A:85:LYS:HG3	2.10	0.52
1:A:21:LEU:HB3	1:A:86:ILE:HD12	1.92	0.51
1:A:403:ARG:O	1:A:407:ARG:HG2	2.11	0.51
1:B:35:ILE:HD11	1:B:67:SER:HB2	1.93	0.51
1:B:389:LEU:HD21	1:B:397:THR:HG22	1.93	0.51
1:A:302:HIS:HA	1:A:325:TYR:CE1	2.45	0.51
1:A:211:ILE:HD12	1:A:252:VAL:HG13	1.93	0.51
1:A:409:LEU:HB3	1:A:414:ILE:HG13	1.92	0.51
1:A:171:THR:HG23	1:A:414:ILE:O	2.10	0.51
1:A:347:ILE:HG21	1:A:349:PHE:CE2	2.46	0.51
1:B:381:SER:OG	1:B:420:HIS:HD2	1.93	0.50
1:A:80:LYS:O	1:A:84:LEU:HG	2.12	0.50
1:B:112:LEU:O	1:B:118:THR:HA	2.12	0.50
1:B:182:GLN:NE2	1:B:217:LYS:HG3	2.28	0.49
1:B:58:ARG:HH11	1:B:247:ILE:HG22	1.77	0.49
1:B:257:ARG:HD3	1:B:280:GLU:CG	2.42	0.49
1:B:384:PRO:HG2	1:B:391:LYS:HG3	1.94	0.49
1:B:199:ASP:O	1:B:203:GLN:HG2	2.12	0.49
1:B:172:ILE:HG12	1:B:416:ILE:HG12	1.93	0.49
1:A:215:SER:OG	1:A:216:THR:N	2.42	0.49
1:A:330:LYS:O	1:A:334:LEU:HG	2.14	0.48
1:B:257:ARG:HB2	1:B:280:GLU:CG	2.43	0.48
1:A:399:ASP:O	1:A:402:THR:HG22	2.13	0.48
1:A:279:VAL:HG12	1:A:301:ILE:HD12	1.94	0.48
1:A:35:ILE:HD11	1:A:67:SER:HB2	1.95	0.48
1:B:381:SER:HB2	1:B:419:VAL:HA	1.94	0.48
1:A:236:LEU:HD22	1:A:272:ILE:HD11	1.94	0.48
1:A:134:LEU:HD23	1:A:134:LEU:C	2.35	0.47
1:A:240:LYS:HA	1:A:243:LEU:HD13	1.96	0.47
1:A:409:LEU:O	1:A:414:ILE:HG12	2.14	0.47
1:A:194:ARG:HD3	1:A:194:ARG:H	1.80	0.47
1:B:129:PHE:HA	1:B:161:ILE:CD1	2.45	0.47
1:B:377:LEU:HD12	1:B:378:VAL:N	2.29	0.47
1:B:153:ILE:HD12	1:B:153:ILE:N	2.29	0.47
1:A:248:TYR:CE2	1:A:250:PRO:HA	2.50	0.47
1:A:338:ILE:HD11	1:A:343:ILE:CG1	2.45	0.47
1:B:150:ASP:HB3	1:B:153:ILE:HD13	1.96	0.47
1:A:368:ILE:O	1:A:372:LEU:HG	2.14	0.46
1:B:58:ARG:HD2	3:B:482:HOH:O	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:190:ASP:C	1:A:190:ASP:OD1	2.53	0.46
1:A:40:LEU:HD21	1:A:168:LEU:HD13	1.98	0.46
1:A:412:LEU:N	1:A:412:LEU:HD23	2.31	0.45
1:B:173:ARG:HD3	1:B:173:ARG:N	2.32	0.45
1:B:112:LEU:CD2	1:B:117:LEU:HD11	2.47	0.45
1:A:191:ASP:OD2	1:A:191:ASP:N	2.50	0.45
1:A:294:TYR:O	1:A:296:LYS:HE2	2.17	0.45
1:B:329:LYS:NZ	3:B:476:HOH:O	2.49	0.44
1:A:339:ALA:HB1	1:A:341:GLN:NE2	2.32	0.44
1:A:284:ILE:HG23	1:A:285:GLU:N	2.31	0.44
1:A:390:THR:O	1:A:393:SER:OG	2.35	0.44
1:B:158:LEU:HA	1:B:158:LEU:HD23	1.71	0.44
1:B:240:LYS:NZ	1:B:269:HIS:HD2	2.16	0.44
1:B:257:ARG:HB2	1:B:280:GLU:HB2	2.00	0.44
1:A:21:LEU:HD23	1:A:21:LEU:HA	1.85	0.44
1:B:191:ASP:N	1:B:191:ASP:OD2	2.50	0.43
1:A:327:GLU:O	1:A:331:GLN:HG2	2.18	0.43
1:B:172:ILE:CD1	1:B:208:ILE:HD11	2.49	0.43
1:A:368:ILE:HG22	1:A:372:LEU:CD1	2.49	0.43
1:B:357:THR:HA	1:B:360:TYR:HB3	2.01	0.43
1:B:110:LEU:HD12	1:B:111:ILE:H	1.81	0.43
1:B:347:ILE:HD13	1:B:361:LEU:HB3	2.00	0.43
1:A:164:LEU:C	1:A:165:LYS:HD3	2.39	0.42
1:B:384:PRO:HB3	1:B:389:LEU:O	2.18	0.42
1:B:213:ALA:HB1	1:B:231:LYS:HD2	2.01	0.42
1:B:257:ARG:HD3	1:B:280:GLU:HG3	2.01	0.42
1:B:291:ILE:O	1:B:295:ASN:N	2.52	0.42
1:B:288:ALA:HB1	1:B:336:HIS:HB2	2.00	0.42
1:A:168:LEU:HB2	3:A:475:HOH:O	2.19	0.42
1:A:258:LYS:CB	1:A:261:VAL:HG13	2.50	0.42
1:A:257:ARG:CB	1:A:280:GLU:HB2	2.46	0.42
1:A:257:ARG:HB2	1:A:280:GLU:CB	2.48	0.42
1:A:327:GLU:O	1:A:330:LYS:HB3	2.20	0.41
1:A:168:LEU:HD12	1:A:168:LEU:HA	1.83	0.41
1:A:282:ASN:O	1:A:283:ASN:HB2	2.21	0.41
1:B:381:SER:CB	1:B:420:HIS:HD2	2.33	0.41
1:A:71:LYS:O	1:A:72:PRO:C	2.56	0.41
1:A:203:GLN:NE2	1:A:247:ILE:HD13	2.35	0.41
1:B:117:LEU:O	1:B:117:LEU:HD12	2.20	0.41
1:A:399:ASP:HA	1:A:402:THR:HG22	2.02	0.41
1:B:405:LEU:O	1:B:409:LEU:HG	2.19	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:365:ILE:HG21	1:B:414:ILE:HD11	2.02	0.41
1:A:36:ARG:HD2	1:A:137:LEU:HD23	2.02	0.41
1:B:300:ILE:HG23	1:B:329:LYS:CE	2.48	0.41
1:B:44:LYS:O	1:B:45:ALA:HB2	2.21	0.40
1:A:174:MET:HA	1:A:208:ILE:O	2.21	0.40
1:B:143:HIS:HA	1:B:144:PRO:HD2	1.91	0.40
1:A:209:ILE:HD12	1:A:211:ILE:HD11	2.04	0.40
1:B:195:LYS:HD3	1:B:238:TYR:CE1	2.56	0.40
1:B:58:ARG:HD3	1:B:247:ILE:HA	2.04	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	353/442 (80%)	337 (96%)	15 (4%)	1 (0%)	44	49
1	B	365/442 (83%)	353 (97%)	12 (3%)	0	100	100
All	All	718/884 (81%)	690 (96%)	27 (4%)	1 (0%)	55	63

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	349	PHE

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was

analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	335/396 (85%)	312 (93%)	23 (7%)	18	19
1	B	344/396 (87%)	328 (95%)	16 (5%)	30	37
All	All	679/792 (86%)	640 (94%)	39 (6%)	23	28

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	71	LYS
1	A	100	ILE
1	A	117	LEU
1	A	118	THR
1	A	165	LYS
1	A	168	LEU
1	A	172	ILE
1	A	185	SER
1	A	190	ASP
1	A	194	ARG
1	A	243	LEU
1	A	258	LYS
1	A	280	GLU
1	A	303	ASN
1	A	320	ASP
1	A	327	GLU
1	A	328	GLN
1	A	338	ILE
1	A	365	ILE
1	A	371	ARG
1	A	387	LEU
1	A	393	SER
1	A	412	LEU
1	B	2	VAL
1	B	11	ASN
1	B	36	ARG
1	B	84	LEU
1	B	98	ARG
1	B	118	THR
1	B	168	LEU
1	B	172	ILE
1	B	181	ASN
1	B	190	ASP
1	B	268	LYS

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Mol	Chain	Res	Type
1	B	282	ASN
1	B	323	CYS
1	B	359	ARG
1	B	371	ARG
1	B	377	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	203	GLN
1	A	295	ASN
1	A	302	HIS
1	A	303	ASN
1	A	341	GLN
1	A	342	ASN
1	B	11	ASN
1	B	182	GLN
1	B	188	ASN
1	B	242	GLN
1	B	269	HIS
1	B	282	ASN
1	B	295	ASN
1	B	302	HIS
1	B	342	ASN
1	B	420	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 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	368/442 (83%)	-0.02	9 (2%) 59 57	19, 39, 72, 85	0
1	B	377/442 (85%)	-0.09	4 (1%) 80 79	20, 37, 67, 82	0
All	All	745/884 (84%)	-0.05	13 (1%) 70 68	19, 38, 70, 85	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	159	GLY	4.3
1	A	324	ASP	3.0
1	B	356	ASP	3.0
1	B	362	LEU	2.6
1	A	404	GLU	2.5
1	A	321	ASN	2.4
1	A	367	GLU	2.2
1	B	349	PHE	2.2
1	B	322	VAL	2.1
1	A	116	LYS	2.1
1	A	182	GLN	2.0
1	A	370	ARG	2.0
1	A	322	VAL	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(\AA^2)	Q<0.9
2	MG	A	423	1/1	0.88	0.12	-	58,58,58,58	0

6.5 Other polymers [i](#)

There are no such residues in this entry.