Full wwPDB X-ray Structure Validation Report (i)

Feb 19, 2016 – 08:07 PM GMT

PDB ID	:	4XQ6
Title	:	CRYSTAL STRUCTURE OF DIHYDROOROTATE DEHYDROGENSE from MYCOBACTERIUM TUBERCULOSIS
Authors	:	Kishor, C.; Addlagatta, A.
Deposited on	:	2015-01-19
Resolution	:	2.00 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

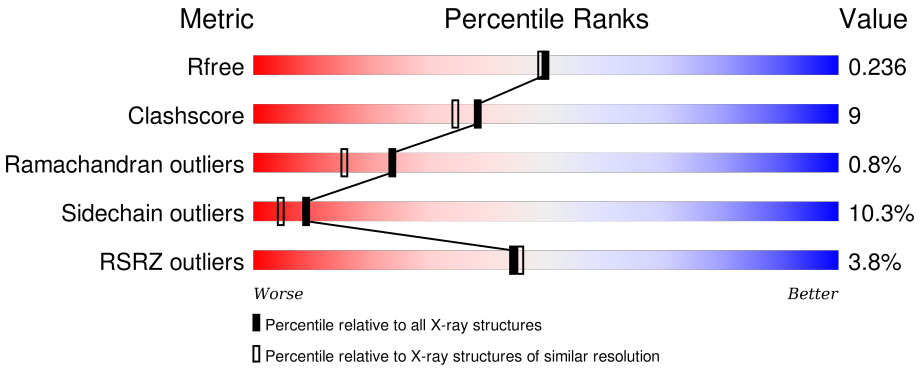
MolProbity	:	4.02b-467
Mogul	:	1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026982
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20026982

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 2.00 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	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (2.00-2.00)

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 <=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	344	3% 76% 13% • 8%
1	B	344	4% 72% 17% • 7%

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 5133 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 Dihydroorotate dehydrogenase (quinone).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	317	Total	C	N	O	S	0	7	0
			2420	1519	458	440	3			
1	B	320	Total	C	N	O	S	0	7	0
			2432	1526	451	451	4			

There are 42 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	10	MET	-	expression tag	UNP A5U4G5
A	11	GLY	-	expression tag	UNP A5U4G5
A	12	SER	-	expression tag	UNP A5U4G5
A	13	SER	-	expression tag	UNP A5U4G5
A	14	HIS	-	expression tag	UNP A5U4G5
A	15	HIS	-	expression tag	UNP A5U4G5
A	16	HIS	-	expression tag	UNP A5U4G5
A	17	HIS	-	expression tag	UNP A5U4G5
A	18	HIS	-	expression tag	UNP A5U4G5
A	19	HIS	-	expression tag	UNP A5U4G5
A	20	SER	-	expression tag	UNP A5U4G5
A	21	SER	-	expression tag	UNP A5U4G5
A	22	GLY	-	expression tag	UNP A5U4G5
A	23	LEU	-	expression tag	UNP A5U4G5
A	24	VAL	-	expression tag	UNP A5U4G5
A	25	PRO	-	expression tag	UNP A5U4G5
A	26	ARG	-	expression tag	UNP A5U4G5
A	27	GLY	-	expression tag	UNP A5U4G5
A	28	SER	-	expression tag	UNP A5U4G5
A	29	HIS	-	expression tag	UNP A5U4G5
A	30	MET	-	expression tag	UNP A5U4G5
B	10	MET	-	expression tag	UNP A5U4G5
B	11	GLY	-	expression tag	UNP A5U4G5
B	12	SER	-	expression tag	UNP A5U4G5
B	13	SER	-	expression tag	UNP A5U4G5

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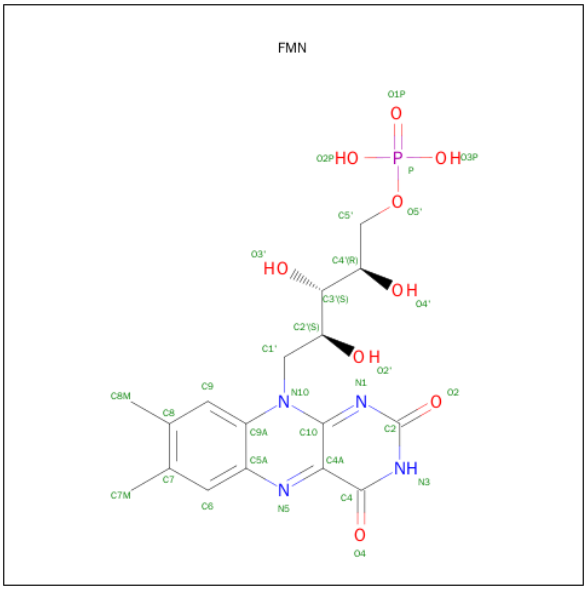
Chain	Residue	Modelled	Actual	Comment	Reference
B	14	HIS	-	expression tag	UNP A5U4G5
B	15	HIS	-	expression tag	UNP A5U4G5
B	16	HIS	-	expression tag	UNP A5U4G5
B	17	HIS	-	expression tag	UNP A5U4G5
B	18	HIS	-	expression tag	UNP A5U4G5
B	19	HIS	-	expression tag	UNP A5U4G5
B	20	SER	-	expression tag	UNP A5U4G5
B	21	SER	-	expression tag	UNP A5U4G5
B	22	GLY	-	expression tag	UNP A5U4G5
B	23	LEU	-	expression tag	UNP A5U4G5
B	24	VAL	-	expression tag	UNP A5U4G5
B	25	PRO	-	expression tag	UNP A5U4G5
B	26	ARG	-	expression tag	UNP A5U4G5
B	27	GLY	-	expression tag	UNP A5U4G5
B	28	SER	-	expression tag	UNP A5U4G5
B	29	HIS	-	expression tag	UNP A5U4G5
B	30	MET	-	expression tag	UNP A5U4G5

- Molecule 2 is FLAVIN MONONUCLEOTIDE (three-letter code: FMN) (formula: $C_{17}H_{21}N_4O_9P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			31	17	4	9	1		
2	B	1	Total	C	N	O	P	0	0
			31	17	4	9	1		

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	3	Total Cl 3 3	0	0
3	A	2	Total Cl 2 2	0	0

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total Mg 1 1	0	0

- Molecule 5 is water.

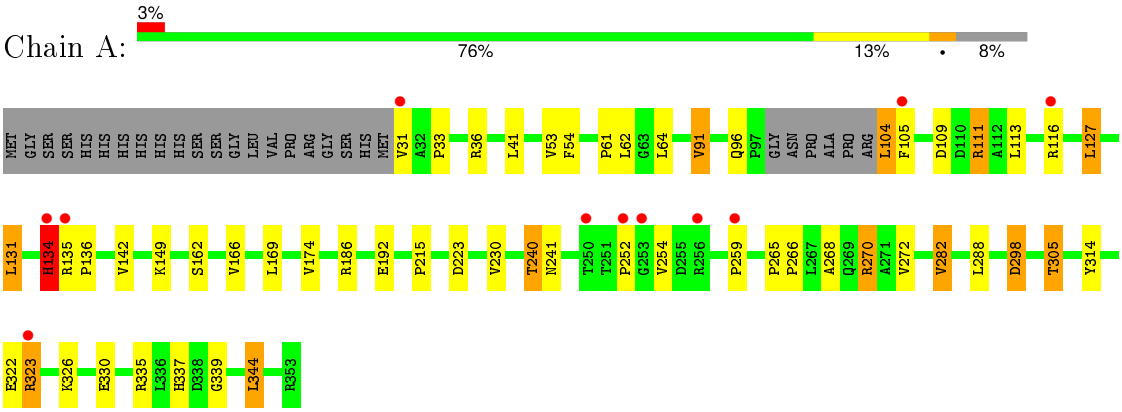
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	110	Total O 110 110	0	0
5	B	103	Total O 103 103	0	0

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

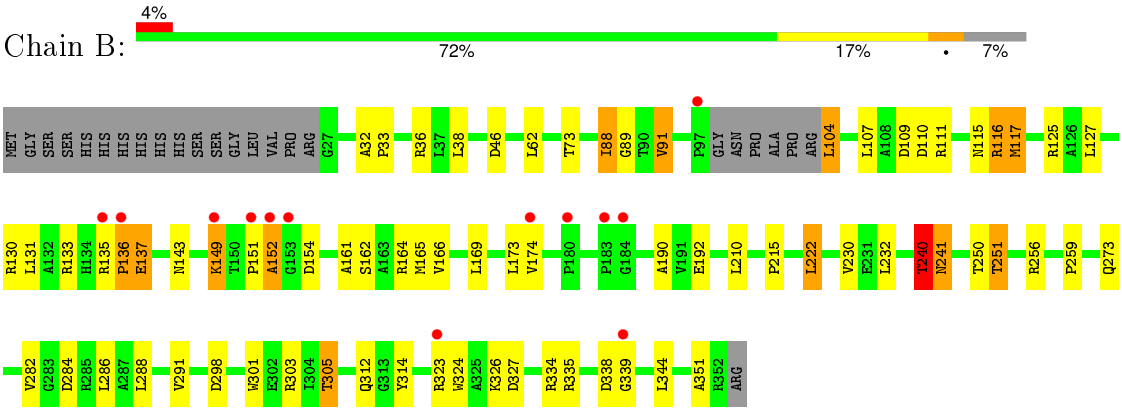
- Molecule 1: Dihydroorotate dehydrogenase (quinone)

Chain A:



- Molecule 1: Dihydroorotate dehydrogenase (quinone)

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	48.34Å 106.92Å 65.09Å 90.00° 106.21° 90.00°	Depositor
Resolution (Å)	23.47 – 2.00 23.47 – 2.00	Depositor EDS
% Data completeness (in resolution range)	92.4 (23.47-2.00) 92.5 (23.47-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.41 (at 2.01Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.176 , 0.231 0.184 , 0.236	Depositor DCC
R_{free} test set	1979 reflections (5.27%)	DCC
Wilson B-factor (Å^2)	24.2	Xtriage
Anisotropy	0.086	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.38 , 46.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.33$	Xtriage
Outliers	0 of 39530 reflections	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	5133	wwPDB-VP
Average B, all atoms (Å^2)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.18% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: FMN, MG, CL

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\#\|Z\| > 5$	RMSZ	$\#\|Z\| > 5$
1	A	0.83	0/2479	1.02	9/3367 (0.3%)
1	B	0.85	0/2486	1.03	10/3380 (0.3%)
All	All	0.84	0/4965	1.02	19/6747 (0.3%)

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	A	0	1
1	B	0	1
All	All	0	2

There are no bond length outliers.

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	186[A]	ARG	NE-CZ-NH2	-9.28	115.66	120.30
1	A	186[B]	ARG	NE-CZ-NH2	-9.28	115.66	120.30
1	B	46	ASP	CB-CG-OD1	6.74	124.36	118.30
1	B	303	ARG	NE-CZ-NH1	6.23	123.41	120.30
1	B	130	ARG	NE-CZ-NH2	-5.88	117.36	120.30
1	B	133	ARG	NE-CZ-NH1	5.78	123.19	120.30
1	B	130	ARG	NE-CZ-NH1	5.67	123.14	120.30
1	A	298	ASP	CB-CG-OD1	5.46	123.22	118.30
1	B	284	ASP	CB-CG-OD1	5.40	123.16	118.30
1	B	91	VAL	CG1-CB-CG2	5.38	119.50	110.90
1	B	303	ARG	NE-CZ-NH2	-5.37	117.61	120.30
1	A	174[A]	VAL	CB-CA-C	-5.29	101.36	111.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	174[B]	VAL	CB-CA-C	-5.29	101.36	111.40
1	A	223	ASP	CB-CG-OD1	5.20	122.98	118.30
1	A	186[A]	ARG	NE-CZ-NH1	5.18	122.89	120.30
1	A	186[B]	ARG	NE-CZ-NH1	5.18	122.89	120.30
1	B	240	THR	N-CA-CB	-5.06	100.69	110.30
1	A	91	VAL	CG1-CB-CG2	5.04	118.97	110.90
1	B	339	GLY	N-CA-C	-5.03	100.52	113.10

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	134	HIS	Peptide
1	B	338	ASP	Peptide

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	2420	0	2481	50	0
1	B	2432	0	2465	43	0
2	A	31	0	19	0	0
2	B	31	0	19	0	0
3	A	2	0	0	0	0
3	B	3	0	0	0	0
4	B	1	0	0	0	0
5	A	110	0	0	4	0
5	B	103	0	0	13	0
All	All	5133	0	4984	90	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 (90) 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:36[B]:ARG:HB3	1:A:36[B]:ARG:NH1	1.61	1.12
1:A:111[A]:ARG:HG2	1:A:111[A]:ARG:HH11	0.97	1.08
1:A:36[B]:ARG:HH11	1:A:36[B]:ARG:HB3	0.92	1.04
1:A:36[B]:ARG:NH1	1:A:36[B]:ARG:CB	2.20	1.02
1:A:36[B]:ARG:HH11	1:A:36[B]:ARG:CB	1.73	1.00
1:A:111[A]:ARG:CG	1:A:111[A]:ARG:HH11	1.77	0.97
1:B:273[A]:GLN:CD	5:B:501:HOH:O	2.09	0.89
1:A:298:ASP:OD1	1:A:335[A]:ARG:NH2	2.06	0.88
1:A:111[A]:ARG:HG2	1:A:111[A]:ARG:NH1	1.79	0.86
1:B:273[A]:GLN:CG	5:B:501:HOH:O	2.26	0.84
1:B:174[B]:VAL:HG12	1:B:210:LEU:HB2	1.58	0.83
1:A:305:THR:HG21	5:A:595:HOH:O	1.82	0.78
1:A:134:HIS:HA	1:A:135:ARG:HB2	1.67	0.75
1:B:149:LYS:HB3	5:B:569:HOH:O	1.85	0.74
1:A:36[B]:ARG:HB2	1:A:36[B]:ARG:NH1	2.04	0.70
1:B:125[B]:ARG:NH2	5:B:502:HOH:O	2.24	0.70
1:B:273[A]:GLN:OE1	5:B:501:HOH:O	2.07	0.67
1:B:240:THR:HG23	1:B:241:ASN:O	1.95	0.67
1:B:305:THR:HG23	1:B:351:ALA:HB2	1.76	0.67
1:A:240:THR:HG23	1:A:241:ASN:O	1.96	0.65
1:A:104:LEU:HD12	1:A:104:LEU:C	2.17	0.65
1:A:134:HIS:HA	1:A:135:ARG:CB	2.26	0.64
1:A:215:PRO:O	1:A:270:ARG:HD3	1.98	0.64
1:A:134:HIS:CG	1:A:135:ARG:HB2	2.34	0.62
1:A:305:THR:HG22	5:A:550:HOH:O	1.98	0.62
1:A:36[B]:ARG:CZ	1:A:36[B]:ARG:CB	2.76	0.60
1:A:116:ARG:NH1	1:A:259:PRO:O	2.34	0.60
1:A:36[B]:ARG:CZ	1:A:36[B]:ARG:HB2	2.32	0.59
1:A:135:ARG:N	1:A:136:PRO:HD3	2.18	0.59
1:A:111[A]:ARG:CG	1:A:111[A]:ARG:NH1	2.44	0.59
1:B:240:THR:CG2	1:B:241:ASN:O	2.51	0.59
1:B:115:ASN:OD1	1:B:117:MET:HG2	2.03	0.59
1:A:305:THR:CG2	5:A:550:HOH:O	2.51	0.58
1:B:104:LEU:N	1:B:117:MET:HE1	2.18	0.58
1:A:192:GLU:HG2	1:B:36:ARG:NH2	2.18	0.58
1:B:110:ASP:OD2	1:B:251:THR:HG23	2.05	0.56
1:B:125[B]:ARG:HB3	5:B:547:HOH:O	2.05	0.56
1:A:230:VAL:HG21	1:A:282:VAL:CG2	2.36	0.55
1:B:151:PRO:O	1:B:152:ALA:CB	2.56	0.54
1:B:190:ALA:HB1	1:B:192:GLU:OE1	2.06	0.54
1:B:136:PRO:O	1:B:137:GLU:CB	2.56	0.53
1:A:298:ASP:OD1	1:A:335[A]:ARG:CZ	2.56	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:127[B]:LEU:HG	1:A:131:LEU:HD22	1.89	0.53
1:A:298:ASP:OD1	1:A:335[A]:ARG:NH1	2.41	0.53
1:B:273[A]:GLN:NE2	5:B:578:HOH:O	0.68	0.53
1:B:169:LEU:CD2	5:B:504:HOH:O	2.57	0.52
1:A:215:PRO:HB3	1:A:240:THR:HG21	1.91	0.52
1:B:161:ALA:O	1:B:165:MET:HG3	2.08	0.52
1:B:116:ARG:HG2	1:B:259:PRO:O	2.10	0.52
1:A:31:VAL:HG12	1:A:33:PRO:HD2	1.92	0.52
1:B:230:VAL:HG21	1:B:282:VAL:HG22	1.93	0.50
1:A:326:LYS:O	1:A:330:GLU:HG3	2.12	0.50
1:A:104:LEU:HD12	1:A:105:PHE:N	2.28	0.49
1:B:117:MET:HE2	5:B:591:HOH:O	2.12	0.49
1:B:149:LYS:CB	5:B:569:HOH:O	2.52	0.48
1:B:215:PRO:HB3	1:B:240:THR:HG21	1.94	0.48
1:B:301:TRP:O	1:B:305:THR:HB	2.13	0.48
1:B:273[A]:GLN:HG2	5:B:501:HOH:O	2.03	0.47
1:A:61:PRO:HB3	1:A:344:LEU:HD11	1.97	0.47
1:A:135:ARG:N	1:A:136:PRO:CD	2.77	0.46
1:A:149:LYS:NZ	5:A:504:HOH:O	2.48	0.46
1:B:222[A]:LEU:HA	1:B:222[A]:LEU:HD13	1.76	0.46
1:B:88[A]:ILE:HD12	1:B:89:GLY:N	2.30	0.46
1:A:192:GLU:HG2	1:B:36:ARG:HH21	1.80	0.46
1:A:230:VAL:CG2	1:A:282:VAL:CG2	2.94	0.45
1:A:53:VAL:HG12	1:A:54:PHE:CD2	2.50	0.45
1:B:154:ASP:OD1	1:B:154:ASP:O	2.34	0.45
1:B:286:LEU:HD12	1:B:286:LEU:C	2.36	0.45
1:B:298:ASP:OD1	1:B:335:ARG:NH2	2.46	0.45
1:A:268:ALA:O	1:A:272:VAL:HG23	2.17	0.45
1:B:125[B]:ARG:NE	5:B:503:HOH:O	2.49	0.44
1:B:169:LEU:HD22	5:B:504:HOH:O	2.15	0.44
1:A:104:LEU:HD11	1:A:113:LEU:HD22	2.00	0.43
1:B:32:ALA:N	1:B:33:PRO:CD	2.81	0.43
1:A:142:VAL:HG11	1:A:166:VAL:HG21	2.00	0.43
1:B:143:ASN:HA	1:B:174[B]:VAL:HG23	2.01	0.43
1:A:337:HIS:O	1:A:339:GLY:O	2.37	0.43
1:B:291:VAL:HB	1:B:312:GLN:HB2	2.00	0.43
1:A:109:ASP:OD2	1:A:252:PRO:HD3	2.18	0.43
1:A:323:ARG:O	1:A:326:LYS:N	2.52	0.42
1:B:109:ASP:HB3	1:B:250:THR:HG22	2.01	0.42
1:B:162:SER:O	1:B:166:VAL:HG22	2.21	0.41
1:A:215:PRO:HB3	1:A:240:THR:CG2	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:104:LEU:HB3	1:B:117:MET:HE3	2.03	0.41
1:A:265:PRO:N	1:A:266:PRO:CD	2.84	0.41
1:A:192:GLU:CG	1:B:36:ARG:HH21	2.32	0.41
1:A:162:SER:O	1:A:166:VAL:HG22	2.21	0.41
1:A:240:THR:CG2	1:A:241:ASN:O	2.66	0.41
1:A:322:GLU:O	1:A:323:ARG:C	2.59	0.40
1:B:143:ASN:HA	1:B:174[B]:VAL:CG2	2.52	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	320/344 (93%)	309 (97%)	10 (3%)	1 (0%)	46 41
1	B	323/344 (94%)	307 (95%)	12 (4%)	4 (1%)	16 8
All	All	643/688 (94%)	616 (96%)	22 (3%)	5 (1%)	24 15

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	136	PRO
1	B	137	GLU
1	B	152	ALA
1	B	324	TRP
1	A	134	HIS

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	249/264 (94%)	227 (91%)	22 (9%)	12 7
1	B	251/264 (95%)	219 (87%)	32 (13%)	5 3
All	All	500/528 (95%)	446 (89%)	54 (11%)	9 4

All (54) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	41	LEU
1	A	62	LEU
1	A	64	LEU
1	A	91	VAL
1	A	96	GLN
1	A	104	LEU
1	A	111[A]	ARG
1	A	111[B]	ARG
1	A	127[A]	LEU
1	A	127[B]	LEU
1	A	131	LEU
1	A	134	HIS
1	A	169	LEU
1	A	240	THR
1	A	254	VAL
1	A	270	ARG
1	A	282	VAL
1	A	288	LEU
1	A	305	THR
1	A	314	TYR
1	A	323	ARG
1	A	344	LEU
1	B	38	LEU
1	B	62	LEU
1	B	73	THR
1	B	88[A]	ILE
1	B	88[B]	ILE
1	B	91	VAL
1	B	104	LEU
1	B	107	LEU
1	B	111	ARG
1	B	116	ARG
1	B	117	MET

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Mol	Chain	Res	Type
1	B	127	LEU
1	B	131	LEU
1	B	135	ARG
1	B	149	LYS
1	B	164	ARG
1	B	173	LEU
1	B	222[A]	LEU
1	B	222[B]	LEU
1	B	232	LEU
1	B	240	THR
1	B	241	ASN
1	B	251	THR
1	B	256	ARG
1	B	288	LEU
1	B	305	THR
1	B	314	TYR
1	B	323	ARG
1	B	326	LYS
1	B	327	ASP
1	B	334	ARG
1	B	344	LEU

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

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no carbohydrates in this entry.

5.6 Ligand geometry (i)

Of 8 ligands modelled in this entry, 6 are monoatomic - leaving 2 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	FMN	A	401	-	32,33,33	1.43	6 (18%)	34,50,50	2.84	9 (26%)
2	FMN	B	402	-	32,33,33	1.61	7 (21%)	34,50,50	2.59	9 (26%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	FMN	A	401	-	-	0/18/18/18	0/3/3/3
2	FMN	B	402	-	-	0/18/18/18	0/3/3/3

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	FMN	C1'-N10	-3.44	1.44	1.48
2	B	402	FMN	P-O3P	-2.72	1.45	1.54
2	A	401	FMN	C2-N3	-2.10	1.33	1.38
2	B	402	FMN	C4A-N5	2.06	1.36	1.33
2	A	401	FMN	C8-C7	2.09	1.46	1.41
2	A	401	FMN	C9A-C5A	2.18	1.47	1.42
2	B	402	FMN	C9A-N10	2.76	1.42	1.38
2	B	402	FMN	C4-C4A	2.93	1.47	1.41
2	A	401	FMN	C4A-C10	3.11	1.46	1.40
2	A	401	FMN	C4-C4A	3.29	1.48	1.41
2	B	402	FMN	C9A-C5A	3.39	1.49	1.42
2	B	402	FMN	C8-C7	3.62	1.50	1.41
2	B	402	FMN	C4A-C10	3.67	1.47	1.40

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	401	FMN	C4-C4A-C10	-8.80	114.31	119.94
2	B	402	FMN	C4-C4A-C10	-7.30	115.27	119.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	401	FMN	N3-C2-N1	-3.58	121.66	127.69
2	B	402	FMN	C4A-C4-N3	-3.53	118.91	123.52
2	B	402	FMN	N3-C2-N1	-3.41	121.95	127.69
2	A	401	FMN	C4A-C4-N3	-2.95	119.66	123.52
2	B	402	FMN	O2P-P-O5'	-2.02	100.82	106.72
2	A	401	FMN	C4A-N5-C5A	2.20	119.32	116.72
2	A	401	FMN	O3P-P-O2P	2.99	118.42	107.44
2	A	401	FMN	C5A-C9A-N10	3.17	119.96	117.58
2	B	402	FMN	C5A-C9A-N10	3.20	119.97	117.58
2	B	402	FMN	C4A-N5-C5A	3.35	120.67	116.72
2	B	402	FMN	C4-C4A-N5	3.42	122.86	118.70
2	B	402	FMN	C1'-N10-C9A	3.71	123.13	118.83
2	A	401	FMN	C1'-N10-C9A	4.22	123.72	118.83
2	A	401	FMN	C4-C4A-N5	4.29	123.92	118.70
2	B	402	FMN	C4-N3-C2	8.95	122.63	115.16
2	A	401	FMN	C4-N3-C2	9.46	123.05	115.16

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

6.1 Protein, DNA and RNA chains

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95^{th} 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	317/344 (92%)	-0.14	11 (3%) 48 49	13, 23, 53, 88	0
1	B	320/344 (93%)	-0.10	13 (4%) 41 42	16, 27, 51, 68	0
All	All	637/688 (92%)	-0.12	24 (3%) 44 45	13, 25, 51, 88	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	134	HIS	5.1
1	A	135	ARG	4.8
1	A	256	ARG	4.3
1	A	105	PHE	3.8
1	A	31	VAL	3.7
1	A	323	ARG	3.0
1	B	152	ALA	2.9
1	B	183	PRO	2.8
1	A	253	GLY	2.5
1	A	116	ARG	2.5
1	B	153	GLY	2.5
1	B	135	ARG	2.4
1	B	184	GLY	2.4
1	A	259	PRO	2.3
1	B	136	PRO	2.3
1	B	323	ARG	2.2
1	B	149	LYS	2.2
1	A	250	THR	2.2
1	B	174[A]	VAL	2.1
1	B	180	PRO	2.1
1	B	97	PRO	2.1
1	B	151	PRO	2.1
1	B	339	GLY	2.1
1	A	252	PRO	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, 95^{th} 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
2	FMN	B	402	31/31	0.98	0.12	-0.34	16,22,26,27	0
2	FMN	A	401	31/31	0.97	0.10	-0.52	13,16,19,26	0
3	CL	A	403	1/1	0.99	0.08	-0.73	24,24,24,24	0
3	CL	A	402	1/1	0.98	0.07	-1.06	29,29,29,29	0
3	CL	B	403	1/1	0.98	0.07	-1.13	38,38,38,38	0
3	CL	B	404	1/1	0.98	0.08	-1.44	45,45,45,45	0
4	MG	B	401	1/1	0.98	0.07	-1.51	19,19,19,19	0
3	CL	B	405	1/1	0.98	0.04	-	25,25,25,25	0

6.5 Other polymers (i)

There are no such residues in this entry.