



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 15, 2017 – 12:15 AM EDT

PDB ID : 4ORM
Title : Crystal structure of Plasmodium falciparum dihydroorotate dehydrogenase bound with Inhibitor DSM338 (N-[3,5-difluoro-4-(trifluoromethyl)phenyl]-5-methyl-2-(trifluoromethyl)[1,2,4]triazolo[1,5-a]pyrimidin-7-amine)
Authors : Deng, X.; Phillips, M.A.
Deposited on : unknown
Resolution : 2.07 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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20029824
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-20029824

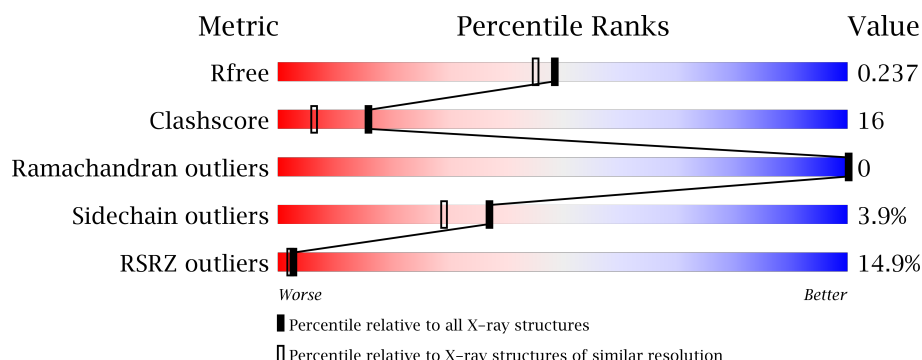
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.07 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	2028 (2.08-2.04)
Clashscore	112137	2143 (2.08-2.04)
Ramachandran outliers	110173	2126 (2.08-2.04)
Sidechain outliers	110143	2126 (2.08-2.04)
RSRZ outliers	101464	2035 (2.08-2.04)

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	401	<div> <div>14%</div> <div>73%</div> <div>19%</div> <div>6%</div> </div>

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
6	SO4	A	1005	-	-	-	X
6	SO4	A	1006	-	-	-	X

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 3194 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.

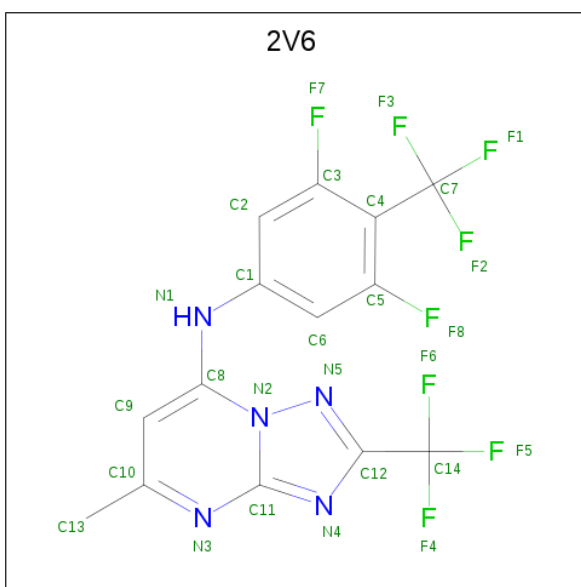
- Molecule 1 is a protein called Dihydroorotate dehydrogenase (quinone), mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	375	Total	C	N	O	S	0	3	0
			2992	1911	500	564	17			

There are 19 discrepancies between the modelled and reference sequences:

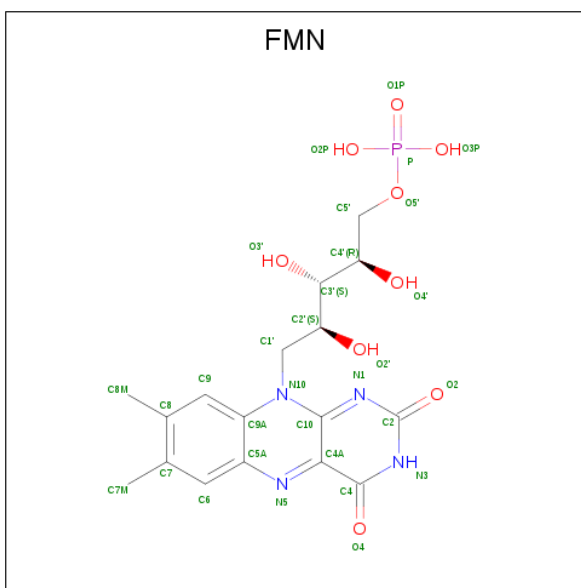
Chain	Residue	Modelled	Actual	Comment	Reference
A	139	MET	-	EXPRESSION TAG	UNP Q08210
A	140	GLY	-	EXPRESSION TAG	UNP Q08210
A	141	HIS	-	EXPRESSION TAG	UNP Q08210
A	142	HIS	-	EXPRESSION TAG	UNP Q08210
A	143	HIS	-	EXPRESSION TAG	UNP Q08210
A	144	HIS	-	EXPRESSION TAG	UNP Q08210
A	145	HIS	-	EXPRESSION TAG	UNP Q08210
A	146	HIS	-	EXPRESSION TAG	UNP Q08210
A	147	ALA	-	EXPRESSION TAG	UNP Q08210
A	148	GLU	-	EXPRESSION TAG	UNP Q08210
A	149	ASN	-	EXPRESSION TAG	UNP Q08210
A	150	LEU	-	EXPRESSION TAG	UNP Q08210
A	151	TYR	-	EXPRESSION TAG	UNP Q08210
A	152	PHE	-	EXPRESSION TAG	UNP Q08210
A	153	GLN	-	EXPRESSION TAG	UNP Q08210
A	154	GLY	-	EXPRESSION TAG	UNP Q08210
A	155	ALA	-	EXPRESSION TAG	UNP Q08210
A	156	ASP	-	EXPRESSION TAG	UNP Q08210
A	157	PRO	-	EXPRESSION TAG	UNP Q08210

- Molecule 2 is N-[3,5-difluoro-4-(trifluoromethyl)phenyl]-5-methyl-2-(trifluoromethyl)[1,2,4]triazolo[1,5-a]pyrimidin-7-amine (three-letter code: 2V6) (formula: C₁₄H₇F₈N₅).



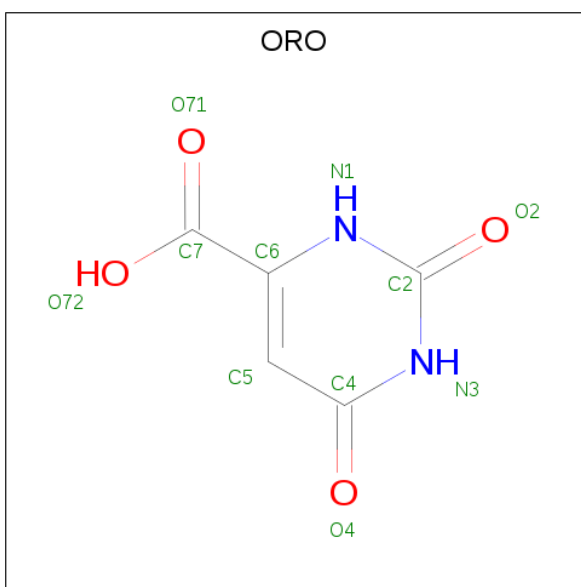
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	F	N	0	0
			27	14	8	5		

- Molecule 3 is FLAVIN MONONUCLEOTIDE (three-letter code: FMN) (formula: $C_{17}H_{21}N_4O_9P$).



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

- Molecule 4 is OROTIC ACID (three-letter code: ORO) (formula: $C_5H_4N_2O_4$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			11	5	2	4		

- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			6	3	3		

- Molecule 6 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



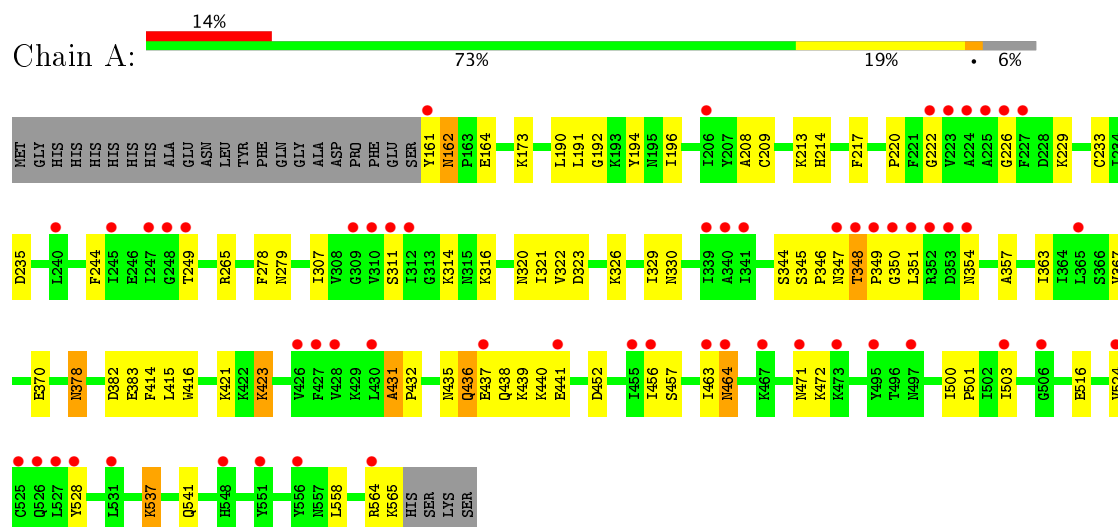
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	O	S	0	0
			5	4	1		
6	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	117	Total	O	0	0
			117	117		

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- Molecule 1: Dihydroorotate dehydrogenase (quinone), mitochondrial



4 Data and refinement statistics

Property	Value	Source
Space group	P 64	Depositor
Cell constants a, b, c, α , β , γ	85.47Å 85.47Å 138.28Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	28.00 – 2.07 27.98 – 2.03	Depositor EDS
% Data completeness (in resolution range)	99.0 (28.00-2.07) 99.0 (27.98-2.03)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.35 (at 2.03Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.186 , 0.240 0.200 , 0.237	Depositor DCC
R_{free} test set	1741 reflections (5.32%)	DCC
Wilson B-factor (Å ²)	44.3	Xtriage
Anisotropy	0.332	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 61.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.047 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3194	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.85% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: FMN, GOL, SO4, 2V6, ORO

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	1.06	2/3042 (0.1%)	0.93	4/4096 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	349	PRO	N-CD	5.30	1.55	1.47
1	A	432	PRO	N-CD	5.12	1.55	1.47

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	500	ILE	C-N-CD	5.87	140.73	128.40
1	A	431	ALA	C-N-CD	5.67	140.31	128.40
1	A	162	ASN	C-N-CD	5.63	140.22	128.40
1	A	348	THR	C-N-CD	5.45	139.85	128.40

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	2992	0	3031	100	0
2	A	27	0	7	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	31	0	19	1	0
4	A	11	0	3	0	0
5	A	6	0	8	1	0
6	A	10	0	0	1	0
7	A	117	0	0	5	0
All	All	3194	0	3068	100	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 16.

All (100) 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:326:LYS:CD	1:A:370:GLU:HG3	1.86	1.04
1:A:321:ILE:HG21	1:A:354:ASN:HD21	1.25	1.02
1:A:326:LYS:HD2	1:A:370:GLU:HG3	1.03	1.01
1:A:326:LYS:HE3	1:A:370:GLU:HG2	1.44	0.97
1:A:326:LYS:HD2	1:A:370:GLU:CG	1.94	0.97
1:A:320:ASN:ND2	1:A:323:ASP:HB2	1.83	0.92
1:A:435:ASN:HD22	1:A:438:GLN:HG3	1.41	0.86
1:A:278:PHE:CD2	1:A:348:THR:HG23	2.10	0.86
1:A:278:PHE:CZ	1:A:345:SER:HB2	2.11	0.85
1:A:383:GLU:OE1	1:A:414:PHE:CZ	2.31	0.84
1:A:278:PHE:HD2	1:A:348:THR:HG23	1.44	0.82
1:A:320:ASN:HD22	1:A:323:ASP:HB2	1.43	0.81
1:A:278:PHE:CZ	1:A:345:SER:CB	2.65	0.80
1:A:321:ILE:HG13	1:A:322:VAL:H	1.45	0.80
1:A:321:ILE:CD1	1:A:363:ILE:HG12	2.11	0.80
1:A:278:PHE:HZ	1:A:345:SER:HB2	1.47	0.76
1:A:321:ILE:HG21	1:A:354:ASN:ND2	2.00	0.76
1:A:326:LYS:HE3	1:A:370:GLU:CG	2.16	0.76
1:A:191:LEU:HD22	1:A:196:ILE:HD11	1.70	0.74
1:A:346:PRO:HG2	1:A:347:ASN:CG	2.07	0.73
1:A:321:ILE:CG1	1:A:322:VAL:N	2.52	0.73
1:A:321:ILE:HG13	1:A:322:VAL:N	2.04	0.73
1:A:345:SER:OG	1:A:346:PRO:HD2	1.89	0.71
1:A:278:PHE:CD2	1:A:348:THR:CG2	2.75	0.70
1:A:321:ILE:HD11	1:A:322:VAL:HG23	1.73	0.69
1:A:217:PHE:CD1	1:A:307:ILE:HD12	2.29	0.67
1:A:435:ASN:O	1:A:439:LYS:HG3	1.92	0.67
1:A:383:GLU:OE1	1:A:414:PHE:CE1	2.47	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:357:ALA:HB2	1:A:441:GLU:HG3	1.77	0.67
1:A:326:LYS:CE	1:A:370:GLU:CG	2.73	0.66
1:A:346:PRO:HD2	1:A:347:ASN:H	1.59	0.66
1:A:321:ILE:HD12	1:A:322:VAL:N	2.10	0.65
1:A:321:ILE:HD11	1:A:363:ILE:HG12	1.78	0.65
1:A:346:PRO:CD	1:A:347:ASN:H	2.09	0.65
1:A:278:PHE:CE2	1:A:345:SER:CB	2.81	0.64
1:A:326:LYS:CD	1:A:370:GLU:CG	2.66	0.63
1:A:314:LYS:HG2	1:A:351:LEU:HD13	1.81	0.61
1:A:435:ASN:HD22	1:A:438:GLN:CG	2.10	0.61
1:A:278:PHE:HD2	1:A:348:THR:CG2	2.13	0.61
1:A:321:ILE:HD13	1:A:363:ILE:HG12	1.82	0.60
1:A:537:LYS:HD2	1:A:541:GLN:NE2	2.16	0.60
1:A:321:ILE:CD1	1:A:322:VAL:N	2.65	0.59
1:A:320:ASN:HB3	1:A:323:ASP:HB3	1.85	0.59
1:A:351:LEU:O	1:A:354:ASN:HB2	2.03	0.58
1:A:209[A]:CYS:SG	7:A:1185:HOH:O	2.57	0.57
1:A:278:PHE:HZ	1:A:345:SER:CB	2.11	0.57
1:A:164:GLU:OE2	6:A:1005:SO4:O2	2.23	0.56
1:A:346:PRO:CD	1:A:347:ASN:N	2.68	0.56
1:A:213:LYS:HB3	7:A:1134:HOH:O	2.05	0.56
1:A:322:VAL:O	1:A:326:LYS:HG2	2.06	0.56
1:A:321:ILE:HD11	1:A:363:ILE:CG1	2.36	0.55
1:A:435:ASN:HB2	1:A:438:GLN:HG3	1.90	0.54
1:A:357:ALA:HB2	1:A:441:GLU:CG	2.38	0.54
1:A:226:GLY:HA3	3:A:1002:FMN:N5	2.23	0.54
1:A:435:ASN:HB2	1:A:438:GLN:H	1.73	0.53
1:A:161:TYR:N	7:A:1178:HOH:O	2.42	0.52
1:A:471:ASN:CB	1:A:472:LYS:HD3	2.40	0.52
1:A:378:ASN:OD1	1:A:415:LEU:HD13	2.10	0.52
1:A:452:ASP:O	1:A:501:PRO:HD2	2.10	0.51
1:A:537:LYS:HD2	1:A:541:GLN:HE22	1.76	0.50
1:A:357:ALA:CB	1:A:441:GLU:HG3	2.42	0.50
1:A:321:ILE:CD1	1:A:363:ILE:CG1	2.87	0.50
1:A:346:PRO:HG2	1:A:347:ASN:ND2	2.26	0.49
1:A:278:PHE:CE2	1:A:345:SER:HB3	2.46	0.49
1:A:346:PRO:HG2	1:A:347:ASN:OD1	2.13	0.49
1:A:463:ILE:C	1:A:464:ASN:HD22	2.16	0.49
1:A:344:SER:HB2	1:A:431:ALA:HB2	1.93	0.49
1:A:350:GLY:O	1:A:351:LEU:HB3	2.13	0.49
1:A:516:GLU:OE1	7:A:1214:HOH:O	2.20	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:435:ASN:ND2	1:A:438:GLN:HG3	2.20	0.48
1:A:436:GLN:O	1:A:440:LYS:HG3	2.14	0.47
1:A:329:ILE:HD11	1:A:367:VAL:HG13	1.95	0.47
1:A:190:LEU:CD1	1:A:194:TYR:CE2	2.97	0.47
1:A:249:THR:HA	1:A:311:SER:HB3	1.95	0.47
1:A:321:ILE:HD12	1:A:322:VAL:CA	2.44	0.47
1:A:382:ASP:CG	5:A:1004:GOL:H32	2.36	0.47
1:A:416:TRP:CZ2	1:A:423:LYS:HB3	2.51	0.46
1:A:190:LEU:HD11	1:A:194:TYR:CE2	2.50	0.46
1:A:278:PHE:CE2	1:A:348:THR:HG23	2.50	0.46
1:A:278:PHE:CZ	1:A:345:SER:HB3	2.49	0.46
1:A:321:ILE:CD1	1:A:322:VAL:HG23	2.44	0.45
1:A:378:ASN:HB2	7:A:1126:HOH:O	2.16	0.45
1:A:190:LEU:CD1	1:A:194:TYR:HE2	2.30	0.44
1:A:330:ASN:HD22	1:A:330:ASN:HA	1.61	0.44
1:A:314:LYS:CD	1:A:351:LEU:HD13	2.49	0.43
1:A:192:GLY:HA3	1:A:233:CYS:SG	2.58	0.43
1:A:314:LYS:HE3	1:A:351:LEU:HB2	2.00	0.43
1:A:314:LYS:CG	1:A:351:LEU:HD13	2.47	0.43
1:A:222:GLY:HA3	1:A:244:PHE:CE1	2.54	0.43
1:A:321:ILE:HD12	1:A:322:VAL:HA	2.00	0.42
1:A:421:LYS:HA	1:A:421:LYS:HD2	1.83	0.42
1:A:435:ASN:ND2	1:A:438:GLN:CD	2.73	0.42
1:A:456:ILE:HA	1:A:457:SER:HA	1.88	0.42
1:A:229:LYS:HD3	1:A:279:ASN:O	2.19	0.42
1:A:503:ILE:HG13	1:A:524:VAL:HG12	2.02	0.42
1:A:208:ALA:O	1:A:220:PRO:HD3	2.19	0.41
1:A:213:LYS:HB3	1:A:214:HIS:H	1.62	0.41
1:A:437:GLU:HA	1:A:440:LYS:HD2	2.02	0.41
1:A:471:ASN:HB3	1:A:472:LYS:HD3	2.02	0.41
1:A:558:LEU:C	1:A:558:LEU:HD23	2.42	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	376/401 (94%)	367 (98%)	9 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	335/354 (95%)	322 (96%)	13 (4%)	37	29

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	162	ASN
1	A	173	LYS
1	A	235	ASP
1	A	265	ARG
1	A	316	LYS
1	A	378	ASN
1	A	423	LYS
1	A	436	GLN
1	A	464	ASN
1	A	528	TYR
1	A	537	LYS
1	A	564	ARG
1	A	565	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	195	ASN
1	A	320	ASN
1	A	330	ASN

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Mol	Chain	Res	Type
1	A	354	ASN
1	A	377	ASN
1	A	435	ASN
1	A	436	GLN
1	A	450	ASN
1	A	464	ASN
1	A	541	GLN
1	A	548	HIS

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](#)

6 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	2V6	A	1001	-	23,29,29	1.61	5 (21%)	32,46,46	3.40	7 (21%)
3	FMN	A	1002	-	31,33,33	1.88	10 (32%)	38,50,50	2.61	13 (34%)
4	ORO	A	1003	-	6,11,11	1.24	1 (16%)	4,15,15	3.68	4 (100%)
5	GOL	A	1004	-	5,5,5	0.71	0	5,5,5	1.45	1 (20%)
6	SO4	A	1005	-	4,4,4	0.31	0	6,6,6	0.43	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	SO4	A	1006	-	4,4,4	0.42	0	6,6,6	0.55	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
2	2V6	A	1001	-	-	0/16/16/16	0/3/3/3
3	FMN	A	1002	-	-	0/16/18/18	0/3/3/3
4	ORO	A	1003	-	-	0/0/4/4	0/1/1/1
5	GOL	A	1004	-	-	0/4/4/4	0/0/0/0
6	SO4	A	1005	-	-	0/0/0/0	0/0/0/0
6	SO4	A	1006	-	-	0/0/0/0	0/0/0/0

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1002	FMN	C2-N3	-2.71	1.32	1.38
3	A	1002	FMN	C1'-N10	-2.68	1.45	1.48
2	A	1001	2V6	C1-N1	-2.49	1.35	1.40
2	A	1001	2V6	C13-C10	2.01	1.55	1.50
4	A	1003	ORO	C4-N3	2.03	1.36	1.33
2	A	1001	2V6	C10-N3	2.08	1.37	1.33
3	A	1002	FMN	C4'-C3'	2.11	1.57	1.53
3	A	1002	FMN	P-O5'	2.13	1.67	1.60
3	A	1002	FMN	C4-N3	2.18	1.37	1.33
3	A	1002	FMN	C8-C7	2.54	1.47	1.41
3	A	1002	FMN	C9A-C5A	2.54	1.47	1.42
2	A	1001	2V6	C9-C10	2.86	1.44	1.38
3	A	1002	FMN	C9A-N10	3.79	1.43	1.38
2	A	1001	2V6	C4-C5	3.84	1.44	1.39
3	A	1002	FMN	C4A-C10	3.96	1.48	1.41
3	A	1002	FMN	C4-C4A	4.14	1.49	1.41

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1001	2V6	C9-C8-N1	-6.73	113.02	123.38
3	A	1002	FMN	C4A-C4-N3	-4.70	116.79	123.48
2	A	1001	2V6	C9-C10-N3	-3.78	117.94	122.61
3	A	1002	FMN	C4-C4A-C10	-3.23	117.35	119.96

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1003	ORO	C5-C4-N3	-2.79	120.76	124.06
4	A	1003	ORO	C5-C6-N1	-2.55	119.31	122.38
3	A	1002	FMN	C9A-C5A-N5	-2.17	119.00	122.24
2	A	1001	2V6	C2-C1-N1	-2.16	113.44	120.36
3	A	1002	FMN	O3'-C3'-C4'	-2.10	103.61	108.82
5	A	1004	GOL	O1-C1-C2	2.18	121.06	110.07
3	A	1002	FMN	O2'-C2'-C3'	2.22	114.60	109.09
3	A	1002	FMN	O3P-P-O2P	2.28	116.81	107.61
3	A	1002	FMN	O3'-C3'-C2'	2.47	114.94	108.82
2	A	1001	2V6	F5-C14-F6	2.70	115.69	105.70
2	A	1001	2V6	C6-C1-C2	2.71	123.40	119.57
2	A	1001	2V6	C8-C9-C10	2.79	119.36	117.10
3	A	1002	FMN	P-O5'-C5'	2.88	126.22	118.30
3	A	1002	FMN	C4-C4A-N5	3.06	122.03	118.68
4	A	1003	ORO	C4-C5-C6	4.12	119.39	116.73
4	A	1003	ORO	C4-N3-C2	4.77	118.23	114.13
3	A	1002	FMN	C5A-C9A-N10	4.80	121.22	117.66
3	A	1002	FMN	C4A-N5-C5A	4.85	121.88	116.76
3	A	1002	FMN	C1'-N10-C9A	5.51	123.39	118.35
3	A	1002	FMN	C4-N3-C2	9.39	123.38	115.16
2	A	1001	2V6	N1-C8-N2	15.82	127.66	113.84

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1002	FMN	1	0
5	A	1004	GOL	1	0
6	A	1005	SO4	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	375/401 (93%)	0.63	56 (14%) 3 2	30, 50, 86, 113	0

All (56) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	161	TYR	7.4
1	A	349	PRO	5.7
1	A	556	TYR	5.2
1	A	441	GLU	5.0
1	A	353	ASP	4.8
1	A	455	ILE	4.6
1	A	467	LYS	4.4
1	A	350	GLY	4.3
1	A	564	ARG	4.2
1	A	223	VAL	4.1
1	A	341	ILE	4.0
1	A	351	LEU	3.9
1	A	310	VAL	3.9
1	A	224	ALA	3.9
1	A	473	LYS	3.7
1	A	471	ASN	3.7
1	A	551	TYR	3.6
1	A	352	ARG	3.5
1	A	354	ASN	3.5
1	A	464	ASN	3.4
1	A	348	THR	3.4
1	A	245	ILE	3.2
1	A	463	ILE	3.1
1	A	227	PHE	3.0
1	A	225	ALA	2.9
1	A	249	THR	2.9
1	A	527	LEU	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	528	TYR	2.8
1	A	309	GLY	2.8
1	A	339	ILE	2.7
1	A	427	PHE	2.7
1	A	206	ILE	2.7
1	A	311	SER	2.6
1	A	340	ALA	2.6
1	A	503	ILE	2.5
1	A	312	ILE	2.5
1	A	525	CYS	2.4
1	A	531	LEU	2.4
1	A	524	VAL	2.4
1	A	226	GLY	2.4
1	A	365	LEU	2.3
1	A	506	GLY	2.3
1	A	526	GLN	2.3
1	A	456	ILE	2.3
1	A	437	GLU	2.3
1	A	347	ASN	2.2
1	A	240	LEU	2.2
1	A	430	LEU	2.1
1	A	497	ASN	2.1
1	A	247	ILE	2.1
1	A	428	VAL	2.1
1	A	426	VAL	2.1
1	A	495	TYR	2.0
1	A	548	HIS	2.0
1	A	222	GLY	2.0
1	A	248	GLY	2.0

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 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
6	SO4	A	1006	5/5	0.83	0.39	5.84	30,30,30,30	0
6	SO4	A	1005	5/5	0.94	0.26	3.43	30,30,30,30	0
3	FMN	A	1002	31/31	0.97	0.23	0.48	33,38,42,44	0
5	GOL	A	1004	6/6	0.95	0.11	0.18	46,48,53,53	0
4	ORO	A	1003	11/11	0.97	0.15	-0.06	39,41,45,46	0
2	2V6	A	1001	27/27	0.97	0.14	-0.41	30,40,46,46	0

6.5 Other polymers [i](#)

There are no such residues in this entry.