



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 08:56 AM GMT

PDB ID : 3GMB
Title : Crystal Structure of 2-Methyl-3-hydroxypyridine-5-carboxylic acid Oxygenase
Authors : McCulloch, K.M.; Mukherjee, T.; Begley, T.P.; Ealick, S.E.
Deposited on : 2009-03-13
Resolution : 2.10 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

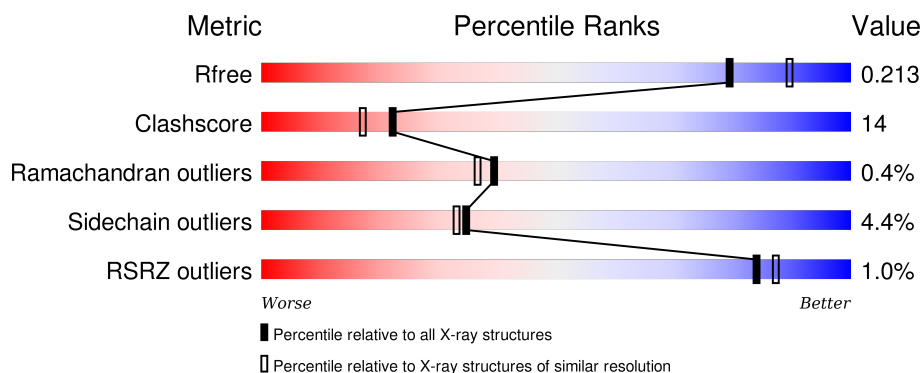
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	3939 (2.10-2.10)
Clashscore	102246	4460 (2.10-2.10)
Ramachandran outliers	100387	4413 (2.10-2.10)
Sidechain outliers	100360	4414 (2.10-2.10)
RSRZ outliers	91569	3948 (2.10-2.10)

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	415	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 18%, green 69%, grey 11%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 69% 18% • 11% </div> </div>
1	B	415	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, green 68%, yellow 19%, orange 1%, grey 11%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 68% 19% • 11% </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6580 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-methyl-3-hydroxypyridine-5-carboxylic acid oxygenase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	369	Total	C	N	O	S	Se	0	0	0
			2865	1807	512	532	5	9			
1	B	369	Total	C	N	O	S	Se	0	0	0
			2865	1807	512	532	5	9			

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-35	MSE	-	EXPRESSION TAG	UNP Q988D3
A	-34	GLY	-	EXPRESSION TAG	UNP Q988D3
A	-33	SER	-	EXPRESSION TAG	UNP Q988D3
A	-32	HIS	-	EXPRESSION TAG	UNP Q988D3
A	-31	HIS	-	EXPRESSION TAG	UNP Q988D3
A	-30	HIS	-	EXPRESSION TAG	UNP Q988D3
A	-29	HIS	-	EXPRESSION TAG	UNP Q988D3
A	-28	HIS	-	EXPRESSION TAG	UNP Q988D3
A	-27	HIS	-	EXPRESSION TAG	UNP Q988D3
A	-26	ASP	-	EXPRESSION TAG	UNP Q988D3
A	-25	ILE	-	EXPRESSION TAG	UNP Q988D3
A	-24	THR	-	EXPRESSION TAG	UNP Q988D3
A	-23	SER	-	EXPRESSION TAG	UNP Q988D3
A	-22	LEU	-	EXPRESSION TAG	UNP Q988D3
A	-21	TYR	-	EXPRESSION TAG	UNP Q988D3
A	-20	LYS	-	EXPRESSION TAG	UNP Q988D3
A	-19	LYS	-	EXPRESSION TAG	UNP Q988D3
A	-18	ALA	-	EXPRESSION TAG	UNP Q988D3
A	-17	GLY	-	EXPRESSION TAG	UNP Q988D3
A	-16	SER	-	EXPRESSION TAG	UNP Q988D3
A	-15	ALA	-	EXPRESSION TAG	UNP Q988D3
A	-14	ALA	-	EXPRESSION TAG	UNP Q988D3
A	-13	ALA	-	EXPRESSION TAG	UNP Q988D3
A	-12	VAL	-	EXPRESSION TAG	UNP Q988D3
A	-11	LEU	-	EXPRESSION TAG	UNP Q988D3

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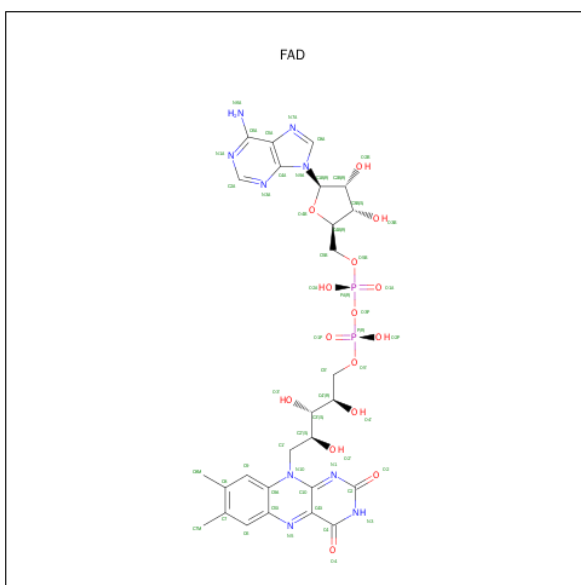
Chain	Residue	Modelled	Actual	Comment	Reference
A	-10	GLU	-	EXPRESSION TAG	UNP Q988D3
A	-9	GLU	-	EXPRESSION TAG	UNP Q988D3
A	-8	ASN	-	EXPRESSION TAG	UNP Q988D3
A	-7	LEU	-	EXPRESSION TAG	UNP Q988D3
A	-6	TYR	-	EXPRESSION TAG	UNP Q988D3
A	-5	PHE	-	EXPRESSION TAG	UNP Q988D3
A	-4	GLY	-	EXPRESSION TAG	UNP Q988D3
A	-3	GLY	-	EXPRESSION TAG	UNP Q988D3
A	-2	SER	-	EXPRESSION TAG	UNP Q988D3
A	-1	PHE	-	EXPRESSION TAG	UNP Q988D3
A	0	THR	-	EXPRESSION TAG	UNP Q988D3
B	-35	MSE	-	EXPRESSION TAG	UNP Q988D3
B	-34	GLY	-	EXPRESSION TAG	UNP Q988D3
B	-33	SER	-	EXPRESSION TAG	UNP Q988D3
B	-32	HIS	-	EXPRESSION TAG	UNP Q988D3
B	-31	HIS	-	EXPRESSION TAG	UNP Q988D3
B	-30	HIS	-	EXPRESSION TAG	UNP Q988D3
B	-29	HIS	-	EXPRESSION TAG	UNP Q988D3
B	-28	HIS	-	EXPRESSION TAG	UNP Q988D3
B	-27	HIS	-	EXPRESSION TAG	UNP Q988D3
B	-26	ASP	-	EXPRESSION TAG	UNP Q988D3
B	-25	ILE	-	EXPRESSION TAG	UNP Q988D3
B	-24	THR	-	EXPRESSION TAG	UNP Q988D3
B	-23	SER	-	EXPRESSION TAG	UNP Q988D3
B	-22	LEU	-	EXPRESSION TAG	UNP Q988D3
B	-21	TYR	-	EXPRESSION TAG	UNP Q988D3
B	-20	LYS	-	EXPRESSION TAG	UNP Q988D3
B	-19	LYS	-	EXPRESSION TAG	UNP Q988D3
B	-18	ALA	-	EXPRESSION TAG	UNP Q988D3
B	-17	GLY	-	EXPRESSION TAG	UNP Q988D3
B	-16	SER	-	EXPRESSION TAG	UNP Q988D3
B	-15	ALA	-	EXPRESSION TAG	UNP Q988D3
B	-14	ALA	-	EXPRESSION TAG	UNP Q988D3
B	-13	ALA	-	EXPRESSION TAG	UNP Q988D3
B	-12	VAL	-	EXPRESSION TAG	UNP Q988D3
B	-11	LEU	-	EXPRESSION TAG	UNP Q988D3
B	-10	GLU	-	EXPRESSION TAG	UNP Q988D3
B	-9	GLU	-	EXPRESSION TAG	UNP Q988D3
B	-8	ASN	-	EXPRESSION TAG	UNP Q988D3
B	-7	LEU	-	EXPRESSION TAG	UNP Q988D3
B	-6	TYR	-	EXPRESSION TAG	UNP Q988D3
B	-5	PHE	-	EXPRESSION TAG	UNP Q988D3

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-4	GLY	-	EXPRESSION TAG	UNP Q988D3
B	-3	GLY	-	EXPRESSION TAG	UNP Q988D3
B	-2	SER	-	EXPRESSION TAG	UNP Q988D3
B	-1	PHE	-	EXPRESSION TAG	UNP Q988D3
B	0	THR	-	EXPRESSION TAG	UNP Q988D3

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

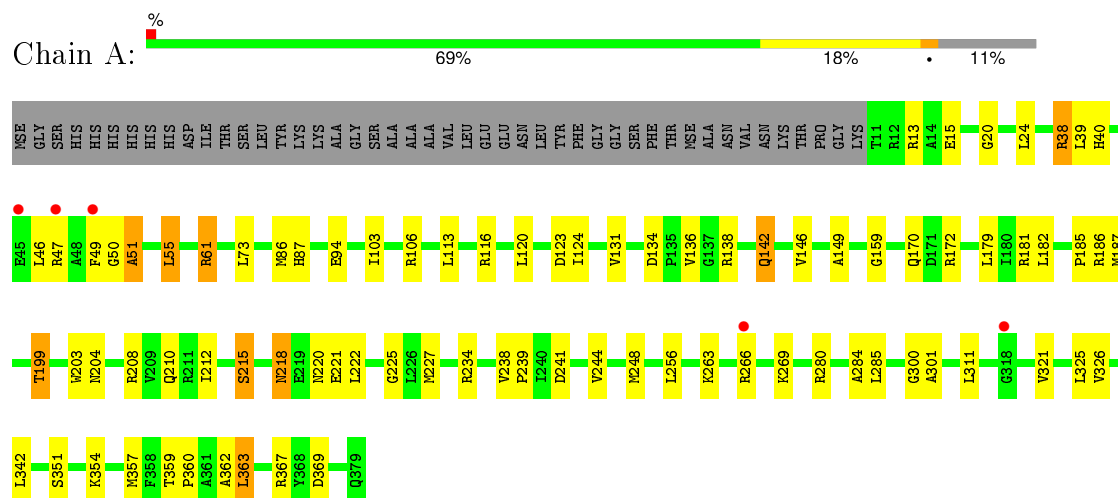
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	364	Total	O	0	0
			364	364		
3	B	380	Total	O	0	0
			380	380		

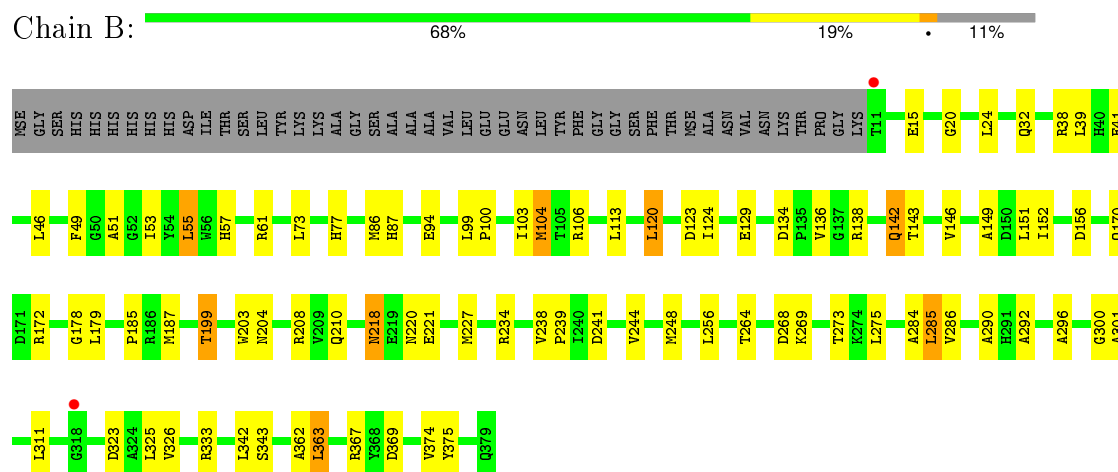
3 Residue-property plots

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: 2-methyl-3-hydroxypyridine-5-carboxylic acid oxygenase



- Molecule 1: 2-methyl-3-hydroxypyridine-5-carboxylic acid oxygenase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	111.28Å 130.20Å 89.54Å 90.00° 122.60° 90.00°	Depositor
Resolution (Å)	29.89 – 2.10 29.89 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.9 (29.89-2.10) 99.3 (29.89-2.10)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.24 (at 2.10Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.193 , 0.224 0.187 , 0.213	Depositor DCC
R_{free} test set	3149 reflections (5.08%)	DCC
Wilson B-factor (Å ²)	17.4	Xtriage
Anisotropy	0.821	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 50.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 62018 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6580	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.96% 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: FAD

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	0.31	0/2922	0.60	1/3954 (0.0%)
1	B	0.31	0/2922	0.61	1/3954 (0.0%)
All	All	0.31	0/5844	0.60	2/7908 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	199	THR	N-CA-C	-5.57	95.95	111.00
1	A	199	THR	N-CA-C	-5.25	96.82	111.00

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	2865	0	2824	77	0
1	B	2865	0	2824	84	0
2	A	53	0	31	4	0
2	B	53	0	31	2	0
3	A	364	0	0	3	1
3	B	380	0	0	10	1
All	All	6580	0	5710	159	2

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 14.

All (159) 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:208:ARG:HD2	1:B:210:GLN:HE21	1.17	1.07
1:A:208:ARG:HD2	1:A:210:GLN:HE21	1.19	1.05
1:B:218:ASN:HD22	1:B:220:ASN:H	1.21	0.89
1:B:208:ARG:HD2	1:B:210:GLN:NE2	1.88	0.88
1:A:218:ASN:HD22	1:A:220:ASN:H	1.26	0.83
1:A:208:ARG:HD2	1:A:210:GLN:NE2	1.93	0.83
1:B:61:ARG:HD3	1:B:369:ASP:OD1	1.79	0.82
1:A:208:ARG:HH11	1:A:210:GLN:NE2	1.81	0.79
1:B:20:GLY:O	1:B:24:LEU:HD23	1.84	0.77
1:A:116:ARG:CZ	1:A:120:LEU:HD11	2.15	0.77
1:B:86:MSE:HE2	1:B:87:HIS:CG	2.21	0.75
1:A:172:ARG:HH11	1:B:170:GLN:HE22	1.35	0.72
1:B:208:ARG:HH11	1:B:210:GLN:NE2	1.88	0.72
1:B:300:GLY:HA3	2:B:400:FAD:H1'2	1.72	0.72
1:B:208:ARG:CD	1:B:210:GLN:HE21	2.02	0.71
1:B:53:ILE:HG23	1:B:104:MSE:HE3	1.72	0.70
1:B:343:SER:HB3	3:B:487:HOH:O	1.90	0.70
1:A:116:ARG:NH1	1:A:120:LEU:HD11	2.06	0.70
1:B:86:MSE:HE2	1:B:87:HIS:CD2	2.27	0.70
1:B:273:THR:HG22	1:B:292:ALA:HB3	1.73	0.69
1:A:218:ASN:ND2	1:A:221:GLU:H	1.92	0.67
1:B:244:VAL:HG21	3:B:390:HOH:O	1.94	0.67
1:B:218:ASN:ND2	1:B:220:ASN:H	1.92	0.67
1:A:300:GLY:HA3	2:A:400:FAD:H1'2	1.76	0.66
1:A:234:ARG:HD2	1:A:241:ASP:OD2	1.97	0.65
1:B:218:ASN:ND2	1:B:221:GLU:H	1.95	0.65
1:A:208:ARG:HH11	1:A:210:GLN:HE22	1.43	0.65
1:B:86:MSE:HE3	1:B:87:HIS:H	1.61	0.64
1:B:273:THR:HB	3:B:667:HOH:O	1.98	0.64
1:A:61:ARG:HD3	1:A:369:ASP:OD1	1.98	0.63
1:A:170:GLN:HE22	1:B:172:ARG:HH11	1.48	0.62
1:A:185:PRO:CB	1:A:187:MSE:HE3	2.29	0.62
1:A:86:MSE:HE3	1:A:87:HIS:H	1.64	0.62
1:A:94:GLU:HB2	3:A:463:HOH:O	2.00	0.61
1:B:55:LEU:HD22	1:B:104:MSE:HE1	1.82	0.61
1:B:208:ARG:HH11	1:B:210:GLN:HE22	1.49	0.61
1:B:273:THR:CG2	1:B:292:ALA:H	2.14	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:238:VAL:HA	1:A:239:PRO:C	2.21	0.61
1:B:138:ARG:HG3	1:B:138:ARG:HH11	1.66	0.60
1:B:208:ARG:HG2	1:B:248:MSE:SE	2.52	0.60
1:A:208:ARG:HG2	1:A:248:MSE:SE	2.51	0.60
1:A:351:SER:HB3	1:A:357:MSE:HE2	1.84	0.60
1:A:185:PRO:HB2	1:A:187:MSE:HE3	1.82	0.59
1:B:46:LEU:HD11	1:B:124:ILE:HG21	1.84	0.59
1:B:238:VAL:HA	1:B:239:PRO:C	2.24	0.58
1:A:86:MSE:HE2	1:A:87:HIS:CE1	2.39	0.57
1:A:86:MSE:HE2	1:A:87:HIS:CG	2.40	0.57
1:B:284:ALA:HB2	1:B:325:LEU:HD13	1.86	0.57
1:B:151:LEU:C	1:B:152:ILE:HD12	2.26	0.56
1:A:86:MSE:HE3	1:A:87:HIS:N	2.21	0.56
1:B:86:MSE:HE1	1:B:204:ASN:HA	1.86	0.56
1:B:210:GLN:HE22	1:B:244:VAL:HG22	1.69	0.56
1:A:204:ASN:HB2	1:A:210:GLN:O	2.06	0.56
1:B:134:ASP:OD1	1:B:136:VAL:HG12	2.05	0.56
1:A:47:ARG:NH1	1:A:47:ARG:HB2	2.21	0.56
1:A:116:ARG:O	1:A:120:LEU:HD13	2.06	0.56
1:A:218:ASN:ND2	1:A:220:ASN:H	1.98	0.55
1:B:86:MSE:HE3	1:B:87:HIS:N	2.21	0.55
1:A:363:LEU:O	1:A:367:ARG:HG3	2.07	0.55
1:B:269:LYS:HD3	3:B:490:HOH:O	2.06	0.55
1:B:204:ASN:HB2	1:B:210:GLN:O	2.07	0.55
1:A:187:MSE:HE2	1:A:187:MSE:HA	1.90	0.54
1:B:134:ASP:HB3	1:B:138:ARG:HB2	1.89	0.54
1:B:264:THR:O	1:B:264:THR:HG22	2.07	0.54
1:A:39:LEU:C	1:A:39:LEU:HD23	2.28	0.54
1:B:55:LEU:HD13	1:B:104:MSE:HE1	1.90	0.54
1:A:46:LEU:HD11	1:A:124:ILE:HG21	1.90	0.53
1:A:134:ASP:HB3	1:A:138:ARG:HB2	1.91	0.53
1:B:296:ALA:N	3:B:487:HOH:O	2.42	0.52
1:A:134:ASP:OD1	1:A:136:VAL:HG12	2.09	0.52
1:B:323:ASP:O	1:B:326:VAL:HG22	2.08	0.52
1:B:234:ARG:HD2	1:B:241:ASP:OD2	2.09	0.51
1:A:20:GLY:O	1:A:24:LEU:HD23	2.09	0.51
1:B:311:LEU:HD11	1:B:325:LEU:HD22	1.93	0.51
1:B:218:ASN:HD21	1:B:221:GLU:H	1.59	0.51
1:A:266:ARG:HH11	1:A:266:ARG:HG3	1.76	0.51
1:A:47:ARG:HB2	1:A:47:ARG:CZ	2.40	0.51
1:A:86:MSE:HE1	1:A:204:ASN:HA	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:49:PHE:O	1:A:106:ARG:HD2	2.11	0.50
1:A:86:MSE:HE2	1:A:87:HIS:CD2	2.47	0.49
1:B:55:LEU:HD22	1:B:104:MSE:CE	2.41	0.49
1:A:342:LEU:HD23	1:A:362:ALA:HA	1.95	0.49
1:A:179:LEU:HD23	1:A:179:LEU:N	2.28	0.49
1:B:49:PHE:O	1:B:106:ARG:HD2	2.13	0.49
1:A:244:VAL:HG11	3:A:420:HOH:O	2.12	0.48
1:A:218:ASN:HD21	1:A:221:GLU:H	1.58	0.48
1:B:301:ALA:HB3	2:B:400:FAD:O2	2.13	0.48
1:A:269:LYS:HD3	3:B:601:HOH:O	2.14	0.48
1:B:286:VAL:HA	3:B:564:HOH:O	2.12	0.48
1:B:152:ILE:N	1:B:152:ILE:HD12	2.29	0.48
1:A:311:LEU:HD11	1:A:325:LEU:HD22	1.95	0.47
1:A:208:ARG:CD	1:A:210:GLN:HE21	2.07	0.47
1:B:273:THR:HG22	1:B:292:ALA:CB	2.43	0.47
1:B:55:LEU:HB2	1:B:104:MSE:HE2	1.96	0.47
1:A:134:ASP:CG	1:A:136:VAL:HG12	2.35	0.47
1:A:103:ILE:HD13	1:A:215:SER:HB3	1.97	0.47
1:B:363:LEU:O	1:B:367:ARG:HG3	2.15	0.46
1:A:86:MSE:HE3	1:A:203:TRP:O	2.16	0.46
1:B:342:LEU:HD23	1:B:362:ALA:HA	1.97	0.46
1:B:326:VAL:HG23	3:B:663:HOH:O	2.16	0.46
1:B:185:PRO:CB	1:B:187:MSE:HE3	2.45	0.46
1:A:39:LEU:HD23	1:A:40:HIS:N	2.31	0.46
1:B:156:ASP:OD1	1:B:156:ASP:N	2.48	0.46
1:B:178:GLY:HA3	1:B:268:ASP:O	2.15	0.46
1:A:179:LEU:HD21	2:A:400:FAD:C7M	2.46	0.45
1:B:311:LEU:HD11	1:B:325:LEU:CD2	2.46	0.45
1:B:187:MSE:HE2	1:B:187:MSE:HA	1.97	0.45
1:A:185:PRO:HB3	1:A:187:MSE:HE3	1.96	0.45
1:B:185:PRO:HB2	1:B:187:MSE:HE3	1.97	0.45
1:B:41:GLU:HG2	1:B:46:LEU:HD23	1.99	0.45
1:A:159:GLY:HA2	3:A:627:HOH:O	2.16	0.45
1:A:301:ALA:HB3	2:A:400:FAD:O2	2.17	0.45
1:A:186:ARG:HD2	1:A:222:LEU:HB2	1.98	0.45
1:A:131:VAL:HG12	1:A:131:VAL:O	2.17	0.45
1:B:94:GLU:HB2	3:B:426:HOH:O	2.17	0.44
1:A:210:GLN:HE22	1:A:244:VAL:CG1	2.30	0.44
1:B:86:MSE:CE	1:B:87:HIS:N	2.80	0.44
1:B:275:LEU:O	1:B:333:ARG:NH2	2.43	0.44
1:A:50:GLY:O	1:A:51:ALA:HB3	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:239:PRO:HG3	1:A:263:LYS:HA	1.99	0.43
1:B:57:HIS:HB2	1:B:100:PRO:HD2	1.99	0.43
1:B:99:LEU:HD13	1:B:100:PRO:O	2.18	0.43
1:B:38:ARG:HD3	1:B:123:ASP:OD2	2.19	0.43
1:A:218:ASN:HD22	1:A:220:ASN:N	2.05	0.43
1:A:181:ARG:O	1:A:182:LEU:HD23	2.19	0.42
1:A:321:VAL:O	1:A:325:LEU:HG	2.19	0.42
1:B:39:LEU:HD23	1:B:39:LEU:C	2.39	0.42
1:A:86:MSE:HE2	1:A:87:HIS:ND1	2.33	0.42
1:A:212:ILE:HA	1:A:225:GLY:O	2.19	0.42
1:A:55:LEU:HD12	1:A:55:LEU:HA	1.90	0.42
1:A:38:ARG:HG3	1:A:123:ASP:OD2	2.19	0.42
1:B:77:HIS:HD2	3:B:606:HOH:O	2.02	0.42
1:B:218:ASN:HD22	1:B:220:ASN:N	2.01	0.42
1:B:15:GLU:HG2	1:B:149:ALA:HB2	2.00	0.42
1:B:86:MSE:HE3	1:B:203:TRP:O	2.20	0.42
1:A:142:GLN:HE21	1:A:142:GLN:HB3	1.69	0.42
1:B:374:VAL:HG22	1:B:375:TYR:N	2.34	0.42
1:A:179:LEU:HB3	1:A:227:MSE:HG2	2.02	0.41
1:B:86:MSE:HE1	1:B:204:ASN:CA	2.48	0.41
1:A:15:GLU:HG2	1:A:149:ALA:HB2	2.01	0.41
1:A:181:ARG:HD3	2:A:400:FAD:HM72	2.03	0.41
1:B:134:ASP:CG	1:B:136:VAL:HG12	2.40	0.41
1:B:285:LEU:HB3	1:B:290:ALA:HB2	2.01	0.41
1:B:273:THR:HG23	1:B:292:ALA:H	1.85	0.41
1:A:38:ARG:HD2	1:A:40:HIS:CE1	2.55	0.41
1:B:179:LEU:HB2	1:B:227:MSE:HG3	2.03	0.41
1:A:311:LEU:HD11	1:A:325:LEU:CD2	2.51	0.41
1:A:359:THR:HB	1:A:360:PRO:CD	2.51	0.41
1:B:32:GLN:HG3	1:B:120:LEU:HD21	2.03	0.41
1:B:146:VAL:O	1:B:146:VAL:HG13	2.21	0.41
1:A:146:VAL:HG23	1:A:146:VAL:O	2.21	0.41
1:B:86:MSE:HA	1:B:86:MSE:HE3	2.02	0.40
1:A:234:ARG:HB3	1:A:244:VAL:HG21	2.04	0.40
1:B:103:ILE:C	1:B:104:MSE:HG3	2.41	0.40
1:B:129:GLU:HB3	1:B:142:GLN:HB2	2.03	0.40
1:B:86:MSE:CE	1:B:204:ASN:HA	2.51	0.40
1:A:284:ALA:HB2	1:A:325:LEU:HD13	2.03	0.40
1:A:354:LYS:HB2	1:A:354:LYS:NZ	2.35	0.40
1:B:138:ARG:HG3	1:B:138:ARG:NH1	2.35	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the sym-

metry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:526:HOH:O	3:A:526:HOH:O[2_656]	1.47	0.73
3:B:616:HOH:O	3:B:616:HOH:O[2_554]	2.12	0.08

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	367/415 (88%)	353 (96%)	13 (4%)	1 (0%)	46	45
1	B	367/415 (88%)	354 (96%)	11 (3%)	2 (0%)	34	30
All	All	734/830 (88%)	707 (96%)	24 (3%)	3 (0%)	39	37

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	143	THR
1	A	51	ALA
1	B	51	ALA

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	297/322 (92%)	282 (95%)	15 (5%)	29	26
1	B	297/322 (92%)	286 (96%)	11 (4%)	41	41

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	594/644 (92%)	568 (96%)	26 (4%)	35	33

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	ARG
1	A	38	ARG
1	A	55	LEU
1	A	61	ARG
1	A	73	LEU
1	A	113	LEU
1	A	142	GLN
1	A	199	THR
1	A	215	SER
1	A	218	ASN
1	A	256	LEU
1	A	280	ARG
1	A	285	LEU
1	A	326	VAL
1	A	363	LEU
1	B	55	LEU
1	B	73	LEU
1	B	104	MSE
1	B	113	LEU
1	B	120	LEU
1	B	142	GLN
1	B	199	THR
1	B	218	ASN
1	B	256	LEU
1	B	285	LEU
1	B	363	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	57	HIS
1	A	97	ASN
1	A	110	HIS
1	A	142	GLN
1	A	170	GLN
1	A	210	GLN

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Mol	Chain	Res	Type
1	A	218	ASN
1	A	340	GLN
1	A	356	ASN
1	B	77	HIS
1	B	97	ASN
1	B	142	GLN
1	B	170	GLN
1	B	210	GLN
1	B	218	ASN
1	B	340	GLN
1	B	356	ASN

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](#)

2 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	FAD	A	400	-	48,58,58	2.33	15 (31%)	54,89,89	2.81	14 (25%)
2	FAD	B	400	-	48,58,58	2.32	15 (31%)	54,89,89	2.86	15 (27%)

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	FAD	A	400	-	-	0/30/50/50	0/6/6/6
2	FAD	B	400	-	-	0/30/50/50	0/6/6/6

All (30) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	400	FAD	C9-C8	2.05	1.43	1.37
2	B	400	FAD	C4X-C10	2.12	1.45	1.41
2	B	400	FAD	C9-C8	2.19	1.43	1.37
2	A	400	FAD	C4X-C10	2.23	1.45	1.41
2	B	400	FAD	C9-C9A	2.34	1.45	1.40
2	A	400	FAD	C8-C7	2.35	1.47	1.41
2	B	400	FAD	C8-C7	2.35	1.47	1.41
2	A	400	FAD	C9-C9A	2.49	1.46	1.40
2	A	400	FAD	C6-C7	2.55	1.44	1.37
2	B	400	FAD	C6-C7	2.68	1.45	1.37
2	B	400	FAD	C10-N1	3.30	1.41	1.35
2	A	400	FAD	C4A-N3A	3.37	1.40	1.35
2	B	400	FAD	C4A-N3A	3.51	1.40	1.35
2	A	400	FAD	C10-N1	3.55	1.41	1.35
2	B	400	FAD	C1'-N10	3.96	1.52	1.48
2	A	400	FAD	C1'-N10	3.97	1.52	1.48
2	A	400	FAD	C2A-N1A	4.06	1.41	1.33
2	B	400	FAD	C2A-N1A	4.06	1.41	1.33
2	A	400	FAD	C10-N10	4.07	1.43	1.39
2	B	400	FAD	C10-N10	4.19	1.44	1.39
2	A	400	FAD	C5X-N5	4.44	1.42	1.35
2	A	400	FAD	C4-N3	4.45	1.41	1.33
2	B	400	FAD	C4-N3	4.46	1.41	1.33
2	B	400	FAD	C5X-N5	4.52	1.42	1.35
2	A	400	FAD	C2A-N3A	5.21	1.41	1.32
2	B	400	FAD	C2A-N3A	5.32	1.41	1.32
2	B	400	FAD	C9A-N10	5.57	1.46	1.38
2	B	400	FAD	C4X-N5	5.65	1.42	1.33
2	A	400	FAD	C9A-N10	5.79	1.46	1.38
2	A	400	FAD	C4X-N5	5.88	1.42	1.33

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	400	FAD	N3A-C2A-N1A	-13.89	118.26	128.89
2	A	400	FAD	N3A-C2A-N1A	-13.74	118.38	128.89
2	B	400	FAD	O2A-PA-O3P	-4.18	86.11	105.09
2	A	400	FAD	O2A-PA-O3P	-4.08	86.60	105.09
2	A	400	FAD	O2A-PA-O1A	-3.48	93.68	112.53
2	B	400	FAD	O3P-PA-O5B	-3.40	93.93	102.94
2	B	400	FAD	O2A-PA-O1A	-3.30	94.66	112.53
2	A	400	FAD	O3P-PA-O5B	-2.80	95.52	102.94
2	B	400	FAD	C4-C4X-C10	-2.71	118.20	119.94
2	B	400	FAD	C4X-C4-N3	-2.70	119.89	123.59
2	A	400	FAD	C4-C4X-C10	-2.60	118.28	119.94
2	A	400	FAD	C4X-C4-N3	-2.54	120.12	123.59
2	B	400	FAD	C1B-N9A-C4A	-2.27	123.52	126.94
2	A	400	FAD	C9A-C5X-N5	-2.10	119.25	122.36
2	A	400	FAD	C1B-N9A-C4A	-2.09	123.79	126.94
2	B	400	FAD	P-O3P-PA	-2.07	126.91	132.73
2	B	400	FAD	C9A-C5X-N5	-2.02	119.37	122.36
2	A	400	FAD	C2B-C3B-C4B	2.28	107.30	102.61
2	B	400	FAD	C2B-C3B-C4B	2.32	107.39	102.61
2	A	400	FAD	C4-C4X-N5	2.56	121.82	118.72
2	B	400	FAD	C4-C4X-N5	2.70	122.00	118.72
2	A	400	FAD	C4B-O4B-C1B	2.73	112.72	109.72
2	B	400	FAD	C4B-O4B-C1B	2.79	112.78	109.72
2	B	400	FAD	C4X-N5-C5X	4.59	122.05	116.76
2	A	400	FAD	C4X-N5-C5X	4.63	122.08	116.76
2	A	400	FAD	C4-N3-C2	6.91	121.22	115.25
2	B	400	FAD	C4-N3-C2	7.01	121.30	115.25
2	A	400	FAD	O2A-PA-O5B	7.71	147.34	108.46
2	B	400	FAD	O2A-PA-O5B	7.93	148.46	108.46

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	400	FAD	4	0
2	B	400	FAD	2	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 [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	360/415 (86%)	-0.22	5 (1%) 78 82	9, 16, 30, 44	0
1	B	360/415 (86%)	-0.21	2 (0%) 90 92	7, 16, 30, 43	0
All	All	720/830 (86%)	-0.21	7 (0%) 84 87	7, 16, 30, 44	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	49	PHE	2.9
1	A	318	GLY	2.7
1	A	45	GLU	2.7
1	A	266	ARG	2.6
1	A	47	ARG	2.1
1	B	318	GLY	2.1
1	B	11	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(\AA^2)	Q<0.9
2	FAD	B	400	53/53	0.91	0.15	0.67	18,21,30,35	0
2	FAD	A	400	53/53	0.93	0.14	0.47	19,24,32,37	0

6.5 Other polymers [i](#)

There are no such residues in this entry.