



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:07 PM GMT

PDB ID : 3W8W
Title : The crystal structure of EncM
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Deposited on : 2013-03-22
Resolution : 1.95 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

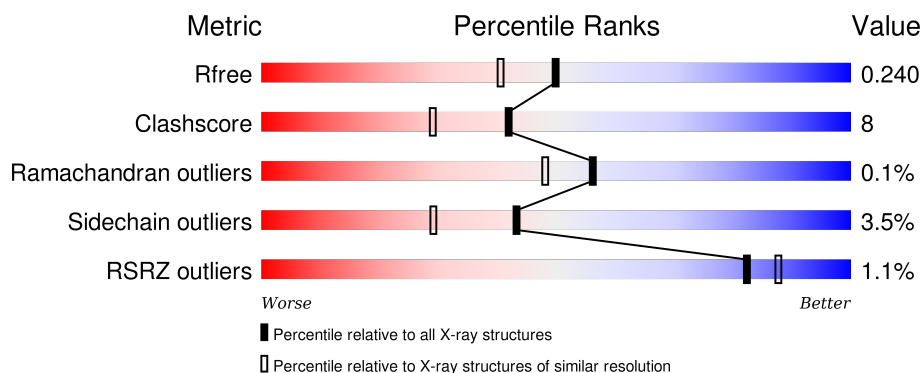
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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	1833 (1.96-1.96)
Clashscore	102246	1953 (1.96-1.96)
Ramachandran outliers	100387	1936 (1.96-1.96)
Sidechain outliers	100360	1936 (1.96-1.96)
RSRZ outliers	91569	1835 (1.96-1.96)

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	468	 82% 15% ..
1	B	468	 81% 16% ..

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7744 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 Putative FAD-dependent oxygenase EncM.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	460	Total	C	N	O	S	0	0	0
			3501	2213	625	650	13			
1	B	461	Total	C	N	O	S	0	0	0
			3507	2216	626	652	13			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	GLY	-	EXPRESSION TAG	UNP Q9KHK2
A	-2	SER	-	EXPRESSION TAG	UNP Q9KHK2
A	-1	HIS	-	EXPRESSION TAG	UNP Q9KHK2
A	0	GLY	-	EXPRESSION TAG	UNP Q9KHK2
B	-3	GLY	-	EXPRESSION TAG	UNP Q9KHK2
B	-2	SER	-	EXPRESSION TAG	UNP Q9KHK2
B	-1	HIS	-	EXPRESSION TAG	UNP Q9KHK2
B	0	GLY	-	EXPRESSION TAG	UNP Q9KHK2

- 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		

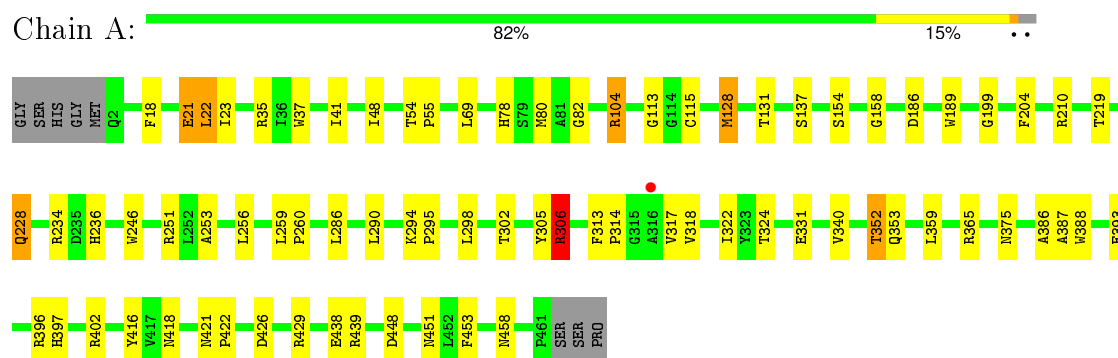
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	307	Total	O	0	0
			307	307		
3	B	323	Total	O	0	0
			323	323		

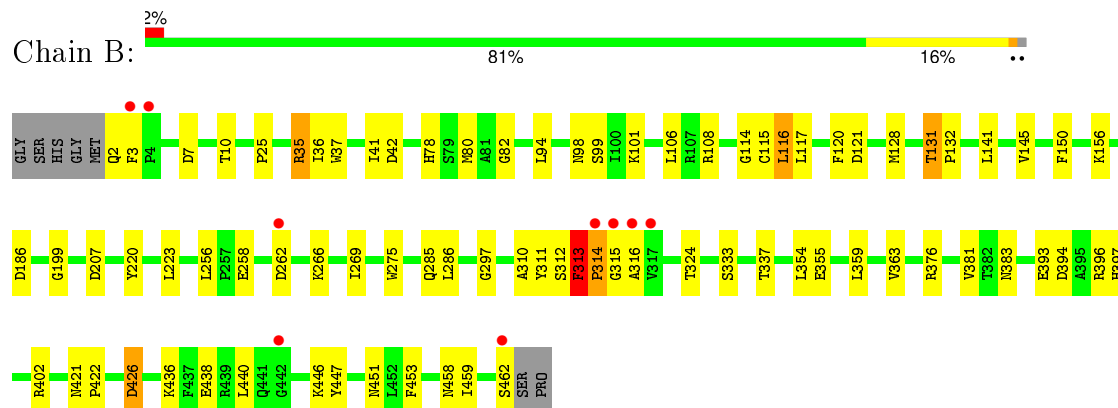
3 Residue-property plots [i](#)

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 ($\text{RSRZ} > 2$). 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: Putative FAD-dependent oxygenase EncM



- Molecule 1: Putative FAD-dependent oxygenase EncM



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	71.41Å 85.28Å 79.67Å 90.00° 99.84° 90.00°	Depositor
Resolution (Å)	47.72 – 1.95 47.68 – 1.95	Depositor EDS
% Data completeness (in resolution range)	99.9 (47.72-1.95) 99.9 (47.68-1.95)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.89 (at 1.95Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.191 , 0.243 0.189 , 0.240	Depositor DCC
R_{free} test set	3442 reflections (5.28%)	DCC
Wilson B-factor (Å ²)	15.7	Xtriage
Anisotropy	0.477	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 50.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 68626 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7744	wwPDB-VP
Average B, all atoms (Å ²)	17.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 16.58% 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

5.1 Standard geometry

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	1.01	4/3592 (0.1%)	1.10	12/4896 (0.2%)
1	B	0.96	2/3598 (0.1%)	1.00	11/4904 (0.2%)
All	All	0.99	6/7190 (0.1%)	1.05	23/9800 (0.2%)

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	B	0	1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	388	TRP	CB-CG	6.99	1.62	1.50
1	A	375	ASN	CG-OD1	5.89	1.36	1.24
1	A	246	TRP	CB-CG	-5.78	1.39	1.50
1	B	121	ASP	CB-CG	5.77	1.63	1.51
1	A	305	TYR	CE1-CZ	5.34	1.45	1.38
1	B	186	ASP	CB-CG	5.16	1.62	1.51

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	306	ARG	NE-CZ-NH2	-24.80	107.90	120.30
1	A	306	ARG	NE-CZ-NH1	18.11	129.36	120.30
1	B	313	PHE	C-N-CD	-12.24	93.68	120.60
1	B	402	ARG	NE-CZ-NH2	-8.64	115.98	120.30
1	A	306	ARG	CD-NE-CZ	8.24	135.14	123.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	186	ASP	CB-CG-OD1	7.83	125.34	118.30
1	B	394	ASP	CB-CG-OD1	6.37	124.03	118.30
1	B	396	ARG	NE-CZ-NH2	-6.33	117.14	120.30
1	A	306	ARG	CG-CD-NE	-6.21	98.76	111.80
1	A	365	ARG	NE-CZ-NH2	-6.19	117.21	120.30
1	A	104	ARG	NE-CZ-NH1	6.10	123.35	120.30
1	B	186	ASP	CB-CG-OD1	5.98	123.68	118.30
1	B	121	ASP	CB-CG-OD2	5.71	123.44	118.30
1	A	186	ASP	CB-CG-OD2	-5.59	113.27	118.30
1	A	128	MET	CG-SD-CE	-5.42	91.52	100.20
1	B	116	LEU	CB-CG-CD1	5.40	120.18	111.00
1	B	117	LEU	CB-CG-CD2	-5.17	102.22	111.00
1	A	402	ARG	NE-CZ-NH2	-5.13	117.74	120.30
1	B	42	ASP	CB-CG-OD1	5.12	122.91	118.30
1	B	426	ASP	CB-CG-OD2	-5.12	113.69	118.30
1	A	448	ASP	CB-CG-OD1	5.08	122.87	118.30
1	B	99	SER	CB-CA-C	-5.03	100.55	110.10
1	A	306	ARG	CB-CG-CD	5.02	124.66	111.60

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	313	PHE	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	3501	0	3413	58	0
1	B	3507	0	3418	67	0
2	A	53	0	31	6	0
2	B	53	0	31	6	0
3	A	307	0	0	12	0
3	B	323	0	0	15	0
All	All	7744	0	6893	117	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 8.

All (117) 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:78:HIS:ND1	2:A:600:FAD:HM83	1.20	1.49
1:B:78:HIS:ND1	2:B:600:FAD:HM83	1.21	1.45
1:A:78:HIS:ND1	2:A:600:FAD:C8M	2.00	1.24
1:B:78:HIS:ND1	2:B:600:FAD:C8M	2.01	1.21
1:A:78:HIS:CE1	2:A:600:FAD:HM83	1.95	0.99
1:A:228:GLN:H	1:A:228:GLN:HE21	1.14	0.96
1:B:78:HIS:CE1	2:B:600:FAD:HM83	2.05	0.92
1:B:35:ARG:HH21	1:B:35:ARG:CG	1.85	0.89
1:B:451:ASN:HD21	1:B:458:ASN:HD21	1.18	0.86
1:A:451:ASN:HD21	1:A:458:ASN:HD21	1.27	0.82
1:B:7:ASP:HB3	1:B:10:THR:HG22	1.59	0.82
1:A:115:CYS:HB3	3:A:920:HOH:O	1.83	0.78
1:A:352:THR:HG21	3:A:900:HOH:O	1.84	0.78
1:B:7:ASP:HB3	1:B:10:THR:CG2	2.14	0.76
1:A:396:ARG:HD3	3:A:874:HOH:O	1.85	0.76
1:B:115:CYS:HB3	3:B:919:HOH:O	1.86	0.75
1:B:35:ARG:HG2	1:B:35:ARG:HH21	1.53	0.73
1:A:352:THR:HB	1:A:386:ALA:HA	1.71	0.73
1:B:128:MET:CE	3:B:875:HOH:O	2.37	0.72
1:B:438:GLU:HB2	3:B:866:HOH:O	1.90	0.71
1:A:228:GLN:N	1:A:228:GLN:HE21	1.87	0.71
1:A:21:GLU:OE1	1:B:101:LYS:NZ	2.23	0.69
1:B:359:LEU:HD11	3:B:779:HOH:O	1.93	0.69
1:B:223:LEU:H	1:B:266:LYS:HZ1	1.40	0.68
1:B:35:ARG:HH21	1:B:35:ARG:HG3	1.58	0.68
1:B:128:MET:HE3	3:B:875:HOH:O	1.93	0.68
1:A:69:LEU:O	3:A:704:HOH:O	2.12	0.68
1:A:128:MET:HE3	1:B:315:GLY:O	1.94	0.67
1:A:128:MET:HE2	3:B:840:HOH:O	1.95	0.67
1:A:306:ARG:HD2	1:B:310:ALA:O	1.95	0.67
1:B:311:TYR:O	1:B:314:PRO:HD3	1.98	0.64
1:A:352:THR:HG22	1:A:387:ALA:H	1.62	0.64
1:A:318:VAL:HG23	1:B:128:MET:HE2	1.80	0.63
1:A:228:GLN:H	1:A:228:GLN:NE2	1.91	0.63
1:B:7:ASP:CB	1:B:10:THR:HG22	2.28	0.62
1:B:421:ASN:HB2	1:B:422:PRO:CD	2.29	0.62
1:A:306:ARG:NH2	3:A:986:HOH:O	2.27	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:236:HIS:HE1	3:A:769:HOH:O	1.81	0.61
1:A:104:ARG:CD	3:A:706:HOH:O	2.49	0.60
1:B:275:TRP:CZ3	1:B:285:GLN:HG3	2.37	0.60
1:A:219:THR:CG2	1:A:298:LEU:HD13	2.31	0.59
1:A:318:VAL:HG23	3:B:875:HOH:O	2.03	0.59
1:B:37:TRP:CH2	2:B:600:FAD:HM82	2.38	0.58
1:B:131:THR:HB	1:B:132:PRO:HD2	1.86	0.58
1:A:18:PHE:CE2	1:A:22:LEU:HD13	2.38	0.58
1:A:294:LYS:HD2	1:A:295:PRO:HD2	1.85	0.57
1:B:333:SER:O	1:B:337:THR:HG23	2.06	0.56
1:B:3:PHE:HE2	3:B:740:HOH:O	1.89	0.56
1:B:35:ARG:NH2	1:B:35:ARG:HG2	2.20	0.56
1:B:421:ASN:HB2	1:B:422:PRO:HD2	1.88	0.56
1:B:128:MET:HE2	3:B:875:HOH:O	2.02	0.55
1:A:219:THR:HG22	1:A:298:LEU:HD13	1.86	0.55
1:B:41:ILE:HG13	1:B:82:GLY:HA3	1.88	0.55
1:B:131:THR:HB	1:B:132:PRO:CD	2.36	0.55
1:B:37:TRP:CZ3	2:B:600:FAD:HM82	2.42	0.55
1:B:223:LEU:H	1:B:266:LYS:NZ	2.05	0.55
1:A:219:THR:HG22	1:A:298:LEU:CD1	2.36	0.55
1:B:35:ARG:NH2	1:B:36:ILE:O	2.41	0.54
1:B:98:ASN:OD1	1:B:115:CYS:HA	2.07	0.54
1:A:35:ARG:HD2	3:A:711:HOH:O	2.07	0.54
1:B:363:VAL:O	1:B:376:ARG:HD3	2.07	0.53
1:B:451:ASN:HD21	1:B:458:ASN:ND2	1.99	0.53
1:A:353:GLN:NE2	3:A:990:HOH:O	2.36	0.52
1:A:37:TRP:CZ3	2:A:600:FAD:HM82	2.45	0.52
1:A:78:HIS:ND1	2:A:600:FAD:C8	2.70	0.52
1:A:41:ILE:HG13	1:A:82:GLY:HA3	1.91	0.52
1:A:104:ARG:HD2	3:A:706:HOH:O	2.09	0.52
1:A:189:TRP:CH2	1:A:439:ARG:HD3	2.45	0.52
1:A:421:ASN:HB2	1:A:422:PRO:CD	2.41	0.51
1:A:104:ARG:HH12	1:B:35:ARG:HD3	1.77	0.50
1:B:78:HIS:ND1	2:B:600:FAD:C8	2.72	0.50
1:B:150:PHE:CE1	3:B:779:HOH:O	2.62	0.49
1:B:440:LEU:HB3	1:B:459:ILE:HD13	1.93	0.49
1:B:393:GLU:HB3	1:B:397:HIS:CD2	2.48	0.49
1:A:128:MET:CE	1:B:315:GLY:O	2.59	0.49
1:B:7:ASP:OD2	1:B:10:THR:HG22	2.12	0.49
1:A:256:LEU:HB2	1:A:259:LEU:CD1	2.42	0.49
1:B:355:GLU:CG	1:B:383:ASN:HB3	2.42	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:286:LEU:O	1:A:290:LEU:HG	2.13	0.49
1:A:23:ILE:HB	1:A:48:ILE:HB	1.96	0.48
1:A:210:ARG:NH2	1:B:316:ALA:O	2.46	0.47
1:B:80:MET:SD	1:B:324:THR:HG21	2.55	0.47
1:B:312:SER:HA	1:B:313:PHE:HA	1.59	0.46
1:A:251:ARG:HD2	1:A:353:GLN:HE21	1.80	0.46
1:B:131:THR:CB	1:B:132:PRO:CD	2.93	0.46
1:A:104:ARG:HD3	3:A:706:HOH:O	2.13	0.46
1:A:322:ILE:HD12	1:A:322:ILE:C	2.35	0.46
1:B:156:LYS:HE2	3:B:931:HOH:O	2.15	0.45
1:A:80:MET:SD	1:A:324:THR:HG21	2.56	0.45
1:B:359:LEU:HD13	1:B:381:VAL:HG23	1.98	0.44
1:A:154:SER:HA	1:A:158:GLY:O	2.18	0.44
1:B:94:LEU:O	1:B:114:GLY:HA3	2.18	0.44
1:A:313:PHE:CG	1:A:314:PRO:HD2	2.53	0.43
1:B:120:PHE:HB2	3:B:919:HOH:O	2.17	0.43
1:B:436:LYS:HB2	3:B:818:HOH:O	2.17	0.43
1:A:37:TRP:CH2	2:A:600:FAD:HM82	2.54	0.43
1:A:199:GLY:HA2	1:A:453:PHE:CE2	2.53	0.42
1:B:220:TYR:HA	1:B:269:ILE:O	2.19	0.42
1:A:393:GLU:HB3	1:A:397:HIS:CD2	2.54	0.42
1:B:258:GLU:O	1:B:297:GLY:HA2	2.18	0.42
1:B:199:GLY:HA2	1:B:453:PHE:CE2	2.55	0.42
1:A:253:ALA:HB1	1:A:259:LEU:CD2	2.50	0.42
1:A:302:THR:HG21	1:B:256:LEU:HD22	2.00	0.41
1:A:416:TYR:CZ	1:A:418:ASN:HB2	2.56	0.41
1:B:141:LEU:O	1:B:145:VAL:HG23	2.20	0.41
1:A:113:GLY:HA2	1:A:204:PHE:HE2	1.86	0.41
1:B:2:GLN:HE21	1:B:25:PRO:HB3	1.86	0.41
1:B:120:PHE:N	3:B:919:HOH:O	2.53	0.41
1:B:128:MET:O	1:B:128:MET:HG3	2.21	0.41
1:B:131:THR:CG2	1:B:207:ASP:O	2.69	0.41
1:B:446:LYS:HD3	1:B:447:TYR:CE1	2.56	0.41
1:A:429:ARG:HD3	3:A:729:HOH:O	2.21	0.41
1:B:376:ARG:HG3	3:B:885:HOH:O	2.20	0.41
1:A:421:ASN:HB2	1:A:422:PRO:HD2	2.03	0.41
1:A:54:THR:HB	1:A:55:PRO:HD3	2.03	0.41
1:B:393:GLU:O	1:B:397:HIS:HD2	2.04	0.40
1:A:451:ASN:HD21	1:A:458:ASN:ND2	2.06	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	458/468 (98%)	448 (98%)	10 (2%)	0	100	100
1	B	459/468 (98%)	444 (97%)	14 (3%)	1 (0%)	52	43
All	All	917/936 (98%)	892 (97%)	24 (3%)	1 (0%)	56	48

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	314	PRO

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	355/361 (98%)	340 (96%)	15 (4%)	36	21
1	B	356/361 (99%)	346 (97%)	10 (3%)	51	39
All	All	711/722 (98%)	686 (96%)	25 (4%)	43	29

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

Mol	Chain	Res	Type
1	A	21	GLU
1	A	22	LEU
1	A	131	THR
1	A	137	SER
1	A	228	GLN

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Mol	Chain	Res	Type
1	A	234	ARG
1	A	260	PRO
1	A	306	ARG
1	A	317	VAL
1	A	331	GLU
1	A	340	VAL
1	A	352	THR
1	A	359	LEU
1	A	426	ASP
1	A	438	GLU
1	B	35	ARG
1	B	106	LEU
1	B	108	ARG
1	B	116	LEU
1	B	131	THR
1	B	262	ASP
1	B	286	LEU
1	B	354	LEU
1	B	426	ASP
1	B	462	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	125	GLN
1	A	228	GLN
1	A	236	HIS
1	A	330	ASN
1	A	353	GLN
1	A	375	ASN
1	A	418	ASN
1	A	458	ASN
1	B	2	GLN
1	B	125	GLN
1	B	163	ASN
1	B	330	ASN
1	B	441	GLN
1	B	458	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

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	600	-	48,58,58	1.55	8 (16%)	54,89,89	2.55	11 (20%)
2	FAD	B	600	-	48,58,58	1.38	6 (12%)	54,89,89	2.90	12 (22%)

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	600	-	-	0/30/50/50	0/6/6/6
2	FAD	B	600	-	-	0/30/50/50	0/6/6/6

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	600	FAD	C1'-N10	-2.65	1.45	1.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	600	FAD	P-O2P	-2.26	1.45	1.54
2	A	600	FAD	C6-C5X	-2.22	1.38	1.41
2	A	600	FAD	C10-N10	2.01	1.41	1.39
2	A	600	FAD	C5X-N5	2.16	1.38	1.35
2	B	600	FAD	C8M-C8	2.46	1.56	1.51
2	B	600	FAD	C8A-N7A	2.47	1.39	1.34
2	B	600	FAD	C2A-N3A	2.70	1.37	1.32
2	A	600	FAD	C8A-N7A	3.40	1.41	1.34
2	B	600	FAD	C4X-N5	3.76	1.39	1.33
2	A	600	FAD	O4B-C1B	4.10	1.46	1.41
2	A	600	FAD	C4X-N5	4.16	1.39	1.33
2	B	600	FAD	C4X-C10	4.63	1.49	1.41
2	A	600	FAD	C4X-C10	4.92	1.50	1.41

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	600	FAD	C4X-C10-N10	-11.97	113.47	120.52
2	A	600	FAD	N3A-C2A-N1A	-9.18	121.86	128.89
2	B	600	FAD	N3A-C2A-N1A	-9.09	121.94	128.89
2	A	600	FAD	C4X-C10-N10	-8.22	115.67	120.52
2	B	600	FAD	C4-C4X-C10	-7.00	115.46	119.94
2	A	600	FAD	C4X-C4-N3	-4.72	117.14	123.59
2	B	600	FAD	C4X-C4-N3	-3.70	118.53	123.59
2	A	600	FAD	O4B-C1B-N9A	-3.34	101.10	108.10
2	B	600	FAD	O3B-C3B-C4B	-3.12	101.70	111.05
2	A	600	FAD	C1B-N9A-C4A	-2.96	122.47	126.94
2	A	600	FAD	C4-C4X-C10	-2.56	118.30	119.94
2	B	600	FAD	O4B-C1B-N9A	-2.50	102.86	108.10
2	B	600	FAD	O4'-C4'-C3'	-2.23	103.40	109.02
2	B	600	FAD	O3P-P-O5'	-2.20	97.11	102.94
2	A	600	FAD	C4X-N5-C5X	2.01	119.07	116.76
2	A	600	FAD	C2A-N1A-C6A	2.11	122.53	118.77
2	A	600	FAD	C2B-C1B-N9A	2.13	117.54	114.29
2	B	600	FAD	C2A-N1A-C6A	2.59	123.40	118.77
2	B	600	FAD	C4X-N5-C5X	2.65	119.81	116.76
2	A	600	FAD	C1'-N10-C9A	2.73	121.92	118.86
2	B	600	FAD	C4-C4X-N5	3.41	122.86	118.72
2	B	600	FAD	C4-N3-C2	8.68	122.75	115.25
2	A	600	FAD	C4-N3-C2	9.96	123.86	115.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	600	FAD	6	0
2	B	600	FAD	6	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	460/468 (98%)	0.09	1 (0%) 95 97	8, 15, 27, 44	0
1	B	461/468 (98%)	0.17	9 (1%) 68 77	8, 16, 31, 46	0
All	All	921/936 (98%)	0.13	10 (1%) 82 88	8, 15, 29, 46	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	316	ALA	3.5
1	B	316	ALA	3.2
1	B	3	PHE	3.0
1	B	315	GLY	2.9
1	B	4	PRO	2.8
1	B	262	ASP	2.7
1	B	442	GLY	2.7
1	B	314	PRO	2.2
1	B	317	VAL	2.2
1	B	462	SER	2.1

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(\AA^2)	Q<0.9
2	FAD	B	600	53/53	0.96	0.11	-0.20	8,9,11,11	0
2	FAD	A	600	53/53	0.97	0.09	-0.35	8,9,10,10	0

6.5 Other polymers [i](#)

There are no such residues in this entry.