



# Full wwPDB X-ray Structure Validation Report i

Feb 26, 2014 – 08:09 PM GMT

PDB ID : 3C4A  
Title : Crystal structure of vioD hydroxylase in complex with FAD from Chromobacterium violaceum. Northeast Structural Genomics Consortium Target CvR158  
Authors : Forouhar, F.; Neely, H.; Seetharaman, J.; Janjua, H.; Xiao, R.; Maglaqui, M.; Wang, H.; Baran, M.C.; Acton, T.B.; Montelione, G.T.; Hunt, J.F.; Tong, L.; Northeast Structural Genomics Consortium (NESG)  
Deposited on : 2008-01-29  
Resolution : 2.30 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

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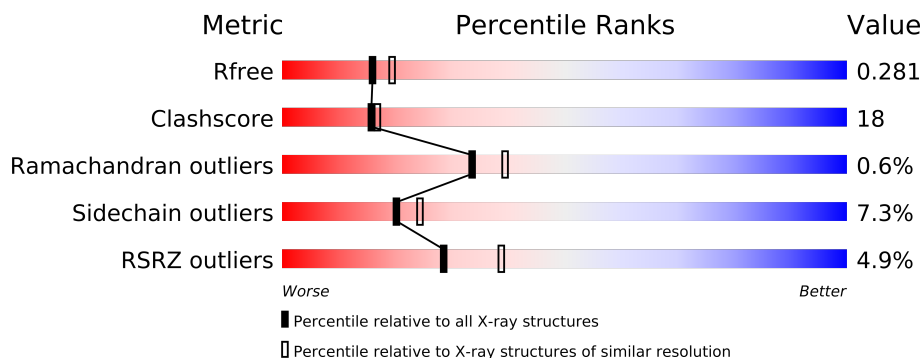
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 2.30 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))
$R_{free}$	66092	2929 (2.30-2.30)
Clashscore	79885	3679 (2.30-2.30)
Ramachandran outliers	78287	3642 (2.30-2.30)
Sidechain outliers	78261	3641 (2.30-2.30)
RSRZ outliers	66119	2930 (2.30-2.30)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	381	

## 2 Entry composition i

There are 3 unique types of molecules in this entry. The entry contains 3009 atoms, of which 0 are hydrogen and 0 are deuterium.

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.

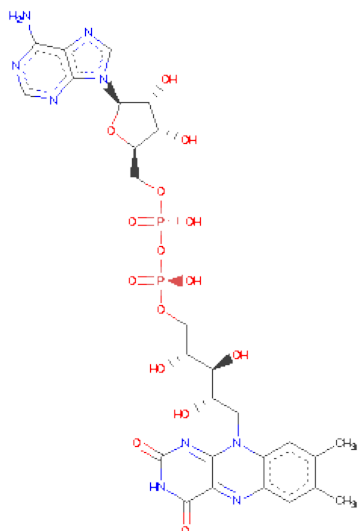
- Molecule 1 is a protein called Probable tryptophan hydroxylase vioD.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	365	2870	1828	507	522	5	8	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	374	LEU	-	EXPRESSION TAG	UNP Q9S3U8
A	375	GLU	-	EXPRESSION TAG	UNP Q9S3U8
A	376	HIS	-	EXPRESSION TAG	UNP Q9S3U8
A	377	HIS	-	EXPRESSION TAG	UNP Q9S3U8
A	378	HIS	-	EXPRESSION TAG	UNP Q9S3U8
A	379	HIS	-	EXPRESSION TAG	UNP Q9S3U8
A	380	HIS	-	EXPRESSION TAG	UNP Q9S3U8
A	381	HIS	-	EXPRESSION TAG	UNP Q9S3U8

- 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		

- Molecule 3 is water.

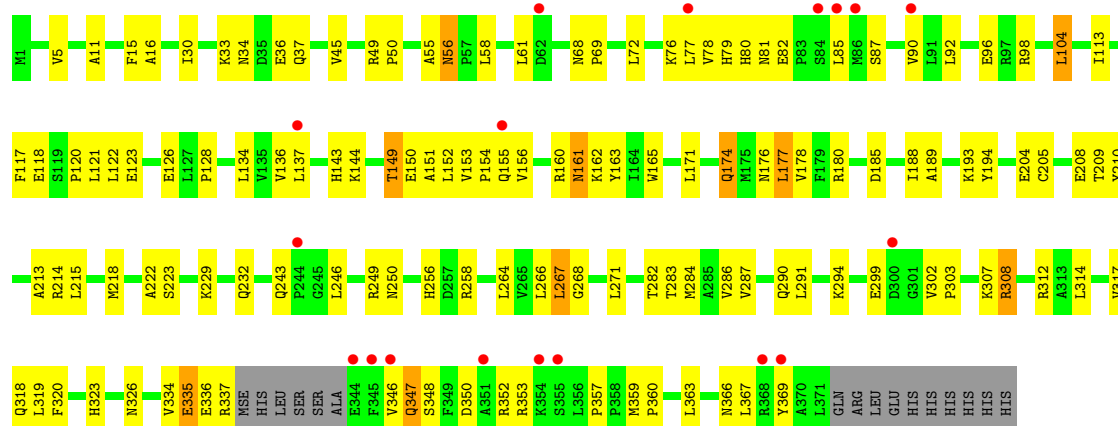
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	86	Total	O	0	0
			86	86		

### 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: Probable tryptophan hydroxylase *vioD*

Chain A: 



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	89.18Å 97.72Å 40.60Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.88 – 2.30 29.47 – 2.30	Depositor EDS
% Data completeness (in resolution range)	84.9 (19.88-2.30) 97.0 (29.47-2.30)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	0.08	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	0.97 (at 2.29Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, $R_{free}$	0.202 , 0.262 0.223 , 0.281	Depositor DCC
$R_{free}$ test set	1578 reflections (9.90%)	DCC
Wilson B-factor (Å <sup>2</sup> )	28.0	Xtriage
Anisotropy	0.400	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 20.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 30443 reflections	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	3009	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.55% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

## 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	0.40	0/2931	0.59	0/3961

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2870	0	2838	107	0
2	A	53	0	31	1	0
3	A	86	0	0	4	1
All	All	3009	0	2869	107	1

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 18.

All (107) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:161:ASN:HB2	1:A:205:CYS:O	1.64	0.96

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:165:TRP:HB3	1:A:246:LEU:HD13	1.55	0.88
1:A:218:MSE:HE3	1:A:223:SER:HA	1.56	0.87
1:A:323:HIS:HD2	1:A:353:ARG:HD3	1.40	0.86
1:A:290:GLN:HE22	1:A:367:LEU:HD23	1.41	0.86
1:A:76:LYS:HB3	1:A:176:ASN:HB3	1.59	0.84
1:A:323:HIS:CD2	1:A:353:ARG:HD3	2.19	0.78
1:A:61:LEU:HD21	1:A:104:LEU:HD12	1.68	0.75
1:A:229:LYS:HA	1:A:232:GLN:HG3	1.68	0.75
1:A:308:ARG:NH2	1:A:369:TYR:HE1	1.86	0.74
1:A:171:LEU:HD22	1:A:193:LYS:HD2	1.69	0.72
1:A:56:ASN:ND2	3:A:431:HOH:O	2.24	0.71
1:A:302:VAL:HB	1:A:303:PRO:HD3	1.72	0.70
1:A:33:LYS:HG3	1:A:120:PRO:HB3	1.73	0.70
1:A:5:VAL:HB	1:A:30:ILE:HG12	1.75	0.68
1:A:264:LEU:HD12	1:A:264:LEU:N	2.11	0.65
1:A:188:ILE:HD11	1:A:334:VAL:HG21	1.78	0.65
1:A:33:LYS:HD2	1:A:34:ASN:ND2	2.12	0.64
1:A:210:TYR:HA	1:A:215:LEU:HD23	1.80	0.64
1:A:78:VAL:HB	1:A:178:VAL:HG13	1.80	0.64
1:A:77:LEU:HD23	1:A:346:VAL:HG13	1.81	0.63
1:A:143:HIS:CD2	1:A:154:PRO:HG3	2.33	0.63
1:A:96:GLU:HG3	1:A:194:TYR:CE2	2.36	0.61
1:A:359:MSE:N	1:A:360:PRO:HD2	2.16	0.60
1:A:162:LYS:NZ	1:A:250:ASN:HD21	1.98	0.60
1:A:314:LEU:O	1:A:318:GLN:HG3	2.02	0.60
1:A:290:GLN:NE2	1:A:367:LEU:HD23	2.15	0.59
1:A:45:VAL:HG11	1:A:92:LEU:HD23	1.85	0.59
1:A:319:LEU:HD11	1:A:357:PRO:HG2	1.85	0.59
1:A:123:GLU:HB2	1:A:126:GLU:HG3	1.84	0.59
1:A:284:MSE:HG3	1:A:320:PHE:CD2	2.38	0.59
1:A:294:LYS:HG3	1:A:308:ARG:HH12	1.68	0.59
1:A:185:ASP:HB3	1:A:209:THR:HG23	1.85	0.58
1:A:314:LEU:HA	1:A:317:VAL:HG22	1.85	0.58
1:A:290:GLN:HE22	1:A:367:LEU:CD2	2.14	0.58
1:A:16:ALA:HB1	1:A:113:ILE:HD13	1.85	0.57
1:A:77:LEU:HD13	1:A:177:LEU:CD1	2.35	0.56
1:A:122:LEU:HD22	1:A:122:LEU:N	2.20	0.56
1:A:81:ASN:HA	1:A:180:ARG:HH22	1.69	0.56
1:A:177:LEU:O	1:A:177:LEU:HD12	2.06	0.55
1:A:243:GLN:HB2	1:A:246:LEU:HG	1.89	0.55
1:A:294:LYS:CG	1:A:308:ARG:HH12	2.19	0.55
1:A:176:ASN:N	1:A:176:ASN:OD1	2.39	0.54

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:160:ARG:HG3	1:A:160:ARG:HH11	1.74	0.53
1:A:33:LYS:CG	1:A:120:PRO:HB3	2.37	0.53
1:A:96:GLU:HG3	1:A:194:TYR:CD2	2.44	0.52
1:A:208:GLU:CD	1:A:208:GLU:H	2.13	0.52
1:A:319:LEU:CD1	1:A:357:PRO:HG2	2.40	0.51
1:A:151:ALA:HB1	1:A:258:ARG:NH1	2.24	0.51
1:A:185:ASP:HB3	1:A:209:THR:CG2	2.41	0.51
1:A:287:VAL:O	1:A:291:LEU:HD13	2.11	0.51
1:A:137:LEU:HD12	1:A:137:LEU:N	2.26	0.50
1:A:161:ASN:ND2	3:A:476:HOH:O	2.44	0.50
1:A:33:LYS:O	1:A:118:GLU:HA	2.11	0.50
1:A:56:ASN:OD1	1:A:58:LEU:N	2.45	0.50
1:A:150:GLU:O	1:A:153:VAL:HG12	2.11	0.50
1:A:326:ASN:HB3	1:A:352:ARG:O	2.13	0.49
1:A:213:ALA:HB3	1:A:215:LEU:CD2	2.43	0.49
1:A:294:LYS:HG3	1:A:308:ARG:NH1	2.27	0.49
1:A:79:HIS:O	1:A:80:HIS:HB2	2.13	0.49
1:A:79:HIS:O	1:A:82:GLU:HG3	2.13	0.49
1:A:55:ALA:HB1	1:A:283:THR:HG23	1.94	0.49
1:A:177:LEU:HA	1:A:189:ALA:O	2.13	0.49
1:A:308:ARG:CZ	1:A:369:TYR:OH	2.61	0.48
1:A:162:LYS:HZ1	1:A:250:ASN:HD21	1.59	0.48
1:A:266:LEU:HB2	1:A:271:LEU:HD13	1.95	0.48
1:A:291:LEU:HD11	1:A:366:ASN:HB3	1.95	0.48
1:A:335:GLU:HG3	1:A:336:GLU:N	2.30	0.47
1:A:210:TYR:CD1	1:A:215:LEU:HB2	2.49	0.47
1:A:68:ASN:N	1:A:69:PRO:HD3	2.29	0.47
1:A:308:ARG:NH2	1:A:369:TYR:CE1	2.75	0.47
1:A:90:VAL:HG23	1:A:92:LEU:HD11	1.97	0.46
1:A:287:VAL:HG21	1:A:363:LEU:HD22	1.97	0.46
1:A:36:GLU:HG2	1:A:117:PHE:CE2	2.50	0.46
1:A:218:MSE:HE3	1:A:223:SER:CA	2.39	0.45
1:A:334:VAL:O	1:A:337:ARG:HB2	2.17	0.45
1:A:213:ALA:HB3	1:A:215:LEU:HD22	1.99	0.44
1:A:312:ARG:HB3	1:A:366:ASN:ND2	2.32	0.44
1:A:149:THR:HG23	1:A:152:LEU:HD12	2.00	0.44
1:A:174:GLN:HE21	1:A:174:GLN:HB3	1.50	0.44
1:A:163:TYR:CE2	2:A:401:FAD:HM71	2.54	0.43
1:A:215:LEU:HA	1:A:218:MSE:HE2	2.00	0.43
1:A:76:LYS:O	1:A:176:ASN:HA	2.19	0.43
1:A:36:GLU:HG2	1:A:117:PHE:CD2	2.55	0.42
1:A:149:THR:CG2	1:A:152:LEU:HD12	2.49	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:267:LEU:HD23	1:A:268:GLY:H	1.84	0.42
1:A:215:LEU:H	1:A:215:LEU:HD22	1.85	0.42
1:A:218:MSE:HG2	1:A:222:ALA:HB3	2.02	0.42
1:A:353:ARG:HD2	3:A:487:HOH:O	2.20	0.42
1:A:90:VAL:HG23	1:A:92:LEU:CD1	2.50	0.42
1:A:347:GLN:HG3	1:A:348:SER:N	2.35	0.42
1:A:81:ASN:OD1	1:A:180:ARG:CZ	2.68	0.42
1:A:264:LEU:CD1	1:A:264:LEU:N	2.82	0.41
1:A:96:GLU:HG3	1:A:194:TYR:HE2	1.80	0.41
1:A:137:LEU:HD11	1:A:264:LEU:HB3	2.02	0.41
1:A:171:LEU:HB3	1:A:193:LYS:HD3	2.01	0.41
1:A:126:GLU:O	1:A:128:PRO:HD3	2.20	0.41
1:A:162:LYS:O	1:A:204:GLU:HA	2.21	0.41
1:A:11:ALA:HB1	1:A:267:LEU:HD22	2.02	0.41
1:A:49:ARG:HG3	1:A:50:PRO:HD2	2.03	0.41
1:A:208:GLU:CD	1:A:208:GLU:N	2.74	0.41
1:A:282:THR:O	1:A:286:VAL:HG23	2.21	0.41
1:A:16:ALA:HB2	1:A:30:ILE:HD11	2.02	0.40
1:A:256:HIS:HB2	3:A:405:HOH:O	2.20	0.40
1:A:299:GLU:HA	1:A:299:GLU:OE2	2.22	0.40
1:A:15:PHE:CE2	1:A:136:VAL:HG21	2.56	0.40
1:A:308:ARG:HH22	1:A:369:TYR:HE1	1.63	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:A:453:HOH:O	3:A:457:HOH:O[2_665]	2.01	0.19

## 5.3 Torsion angles

### 5.3.1 Protein backbone ⓘ

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	361/381 (95%)	333 (92%)	26 (7%)	2 (1%)	33	39

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	87	SER
1	A	156	VAL

### 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	303/309 (98%)	281 (93%)	22 (7%)	20	24

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	GLN
1	A	56	ASN
1	A	72	LEU
1	A	85	LEU
1	A	98	ARG
1	A	104	LEU
1	A	121	LEU
1	A	134	LEU
1	A	144	LYS
1	A	149	THR
1	A	155	GLN
1	A	161	ASN
1	A	174	GLN
1	A	177	LEU
1	A	214	ARG
1	A	249	ARG
1	A	267	LEU
1	A	307	LYS
1	A	308	ARG
1	A	335	GLU
1	A	347	GLN
1	A	350	ASP

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	18	GLN
1	A	34	ASN
1	A	37	GLN
1	A	102	HIS
1	A	155	GLN
1	A	161	ASN
1	A	174	GLN
1	A	250	ASN
1	A	323	HIS
1	A	347	GLN
1	A	366	ASN

### 5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

1 ligand is 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	401	-	58,58,58	3.50	17 (29%)	85,89,89	3.45	27 (31%)

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	401	-	2/2/9/9	0/34/50/50	0/1/6/6

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	FAD	C1'-N10	-21.70	1.23	1.48
2	A	401	FAD	C2A-N1A	-6.08	1.22	1.33
2	A	401	FAD	C2A-N3A	5.06	1.42	1.32
2	A	401	FAD	C6-C5X	-4.57	1.36	1.41
2	A	401	FAD	C1'-C2'	-4.18	1.47	1.51
2	A	401	FAD	C4A-N3A	4.10	1.41	1.35
2	A	401	FAD	C8A-N7A	3.63	1.41	1.34
2	A	401	FAD	C6A-N6A	-3.35	1.24	1.35
2	A	401	FAD	C2B-C1B	-3.24	1.48	1.53
2	A	401	FAD	C9-C9A	-3.23	1.34	1.40
2	A	401	FAD	C5A-N7A	-3.16	1.28	1.40
2	A	401	FAD	C8M-C8	-2.75	1.45	1.51
2	A	401	FAD	C4-N3	2.68	1.41	1.37
2	A	401	FAD	C8A-N9A	-2.53	1.32	1.36
2	A	401	FAD	C4-C4X	2.43	1.45	1.41
2	A	401	FAD	C10-N1	2.28	1.39	1.35
2	A	401	FAD	C5A-C4A	-2.01	1.36	1.40

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	FAD	C2'-C1'-N10	13.31	130.10	112.45
2	A	401	FAD	C2-N1-C10	12.43	127.50	114.98
2	A	401	FAD	C4A-C5A-N7A	12.39	120.14	109.52
2	A	401	FAD	C4X-C10-N1	-8.86	113.88	122.73
2	A	401	FAD	N3A-C2A-N1A	-8.45	121.64	128.71
2	A	401	FAD	N3A-C4A-N9A	6.15	136.54	125.43
2	A	401	FAD	C5A-C4A-N9A	-5.95	98.56	107.16
2	A	401	FAD	C1'-N10-C9A	5.74	124.46	118.87
2	A	401	FAD	O4B-C1B-N9A	-5.04	103.75	108.44
2	A	401	FAD	C1'-C2'-C3'	4.92	123.89	109.82
2	A	401	FAD	C9A-N10-C10	-4.76	117.09	121.77
2	A	401	FAD	C4-C4X-C10	4.72	124.56	116.95
2	A	401	FAD	C5X-C9A-N10	3.90	120.64	116.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	FAD	C8A-N9A-C1B	-3.51	119.46	126.38
2	A	401	FAD	N1-C10-N10	3.44	125.00	115.97
2	A	401	FAD	N3-C2-N1	-3.35	114.06	121.19
2	A	401	FAD	C2A-N1A-C6A	3.22	124.58	118.77
2	A	401	FAD	C8A-N9A-C4A	3.17	109.32	106.90
2	A	401	FAD	O3'-C3'-C2'	2.89	116.04	108.74
2	A	401	FAD	O2'-C2'-C3'	2.88	116.21	109.05
2	A	401	FAD	C1B-N9A-C4A	-2.85	121.70	126.64
2	A	401	FAD	C8M-C8-C9	-2.58	114.15	120.38
2	A	401	FAD	P-O3P-PA	2.31	138.47	131.68
2	A	401	FAD	C6-C5X-N5	-2.22	116.38	118.97
2	A	401	FAD	N6A-C6A-N1A	-2.15	115.15	119.36
2	A	401	FAD	C4B-O4B-C1B	-2.08	107.49	109.75
2	A	401	FAD	O4B-C1B-C2B	-2.04	103.64	106.77

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	401	FAD	C4'
2	A	401	FAD	C2'

There are no torsion outliers.

There are no ring outliers.

## 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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	365/381 (95%)	0.26	18 (4%)	28 39	7, 27, 55, 80	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	85	LEU	7.5
1	A	84	SER	5.4
1	A	86	MSE	4.4
1	A	355	SER	3.5
1	A	77	LEU	3.0
1	A	244	PRO	2.8
1	A	354	LYS	2.8
1	A	369	TYR	2.7
1	A	90	VAL	2.6
1	A	345	PHE	2.6
1	A	300	ASP	2.5
1	A	351	ALA	2.5
1	A	368	ARG	2.5
1	A	346	VAL	2.2
1	A	62	ASP	2.2
1	A	344	GLU	2.2
1	A	155	GLN	2.1
1	A	137	LEU	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

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<sup>th</sup> 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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	FAD	A	401	53/53	0.17	1.22	4,21,29,33	0

### 6.5 Other polymers

There are no such residues in this entry.