



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 03:30 PM GMT

PDB ID : 4CF6  
Title : Crystal structure of the complex of the P187S variant of human NAD(P)  
H:quinone oxidoreductase with Cibacron blue at 2.7 Å resolution  
Authors : Lienhart, W.D.; Gudipati, V.; Uhl, M.K.; Binter, A.; Pulido, S.; Saf, R.;  
Zanger, K.; Gruber, K.; Macheroux, P.  
Deposited on : 2013-11-13  
Resolution : 2.69 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](mailto: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.

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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 : 1.7 (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

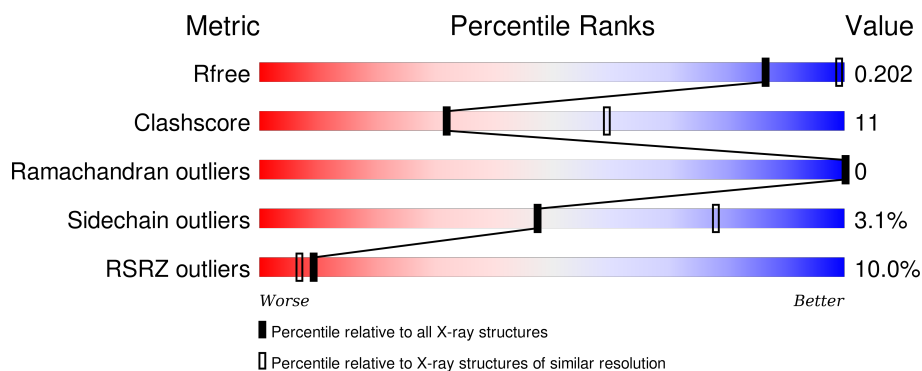
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.69 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	2103 (2.70-2.70)
Clashscore	102246	2422 (2.70-2.70)
Ramachandran outliers	100387	2382 (2.70-2.70)
Sidechain outliers	100360	2382 (2.70-2.70)
RSRZ outliers	91569	2107 (2.70-2.70)

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. 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	294	<div> <div>9%</div> <div>66%</div> <div>26%</div> <div>8%</div> </div>
1	B	294	<div> <div>9%</div> <div>71%</div> <div>20%</div> <div>8%</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
3	CBD	B	1276	-	-	-	X

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4531 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 NAD(P)H DEHYDROGENASE [QUINONE] 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	270	Total	C	N	O	S	0	0	0
			2153	1399	361	386	7			
1	B	271	Total	C	N	O	S	0	0	0
			2163	1405	363	388	7			

There are 42 discrepancies between the modelled and reference sequences:

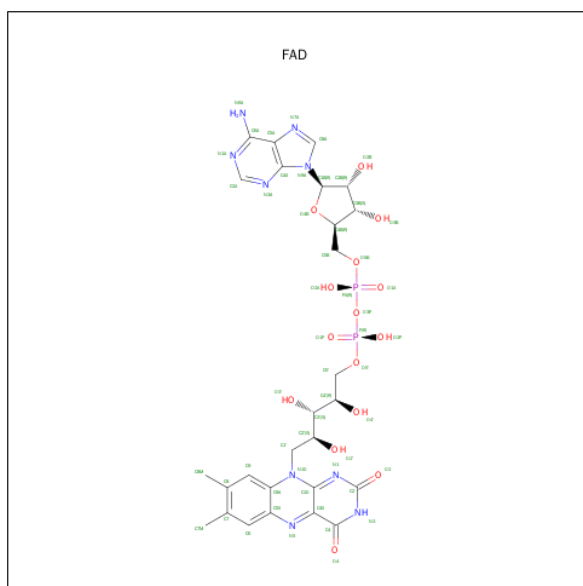
Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	EXPRESSION TAG	UNP P15559
A	-18	GLY	-	EXPRESSION TAG	UNP P15559
A	-17	SER	-	EXPRESSION TAG	UNP P15559
A	-16	SER	-	EXPRESSION TAG	UNP P15559
A	-15	HIS	-	EXPRESSION TAG	UNP P15559
A	-14	HIS	-	EXPRESSION TAG	UNP P15559
A	-13	HIS	-	EXPRESSION TAG	UNP P15559
A	-12	HIS	-	EXPRESSION TAG	UNP P15559
A	-11	HIS	-	EXPRESSION TAG	UNP P15559
A	-10	HIS	-	EXPRESSION TAG	UNP P15559
A	-9	SER	-	EXPRESSION TAG	UNP P15559
A	-8	SER	-	EXPRESSION TAG	UNP P15559
A	-7	GLY	-	EXPRESSION TAG	UNP P15559
A	-6	LEU	-	EXPRESSION TAG	UNP P15559
A	-5	VAL	-	EXPRESSION TAG	UNP P15559
A	-4	PRO	-	EXPRESSION TAG	UNP P15559
A	-3	ARG	-	EXPRESSION TAG	UNP P15559
A	-2	GLY	-	EXPRESSION TAG	UNP P15559
A	-1	SER	-	EXPRESSION TAG	UNP P15559
A	0	HIS	-	EXPRESSION TAG	UNP P15559
A	187	SER	PRO	ENGINEERED MUTATION	UNP P15559
B	-19	MET	-	EXPRESSION TAG	UNP P15559
B	-18	GLY	-	EXPRESSION TAG	UNP P15559
B	-17	SER	-	EXPRESSION TAG	UNP P15559
B	-16	SER	-	EXPRESSION TAG	UNP P15559

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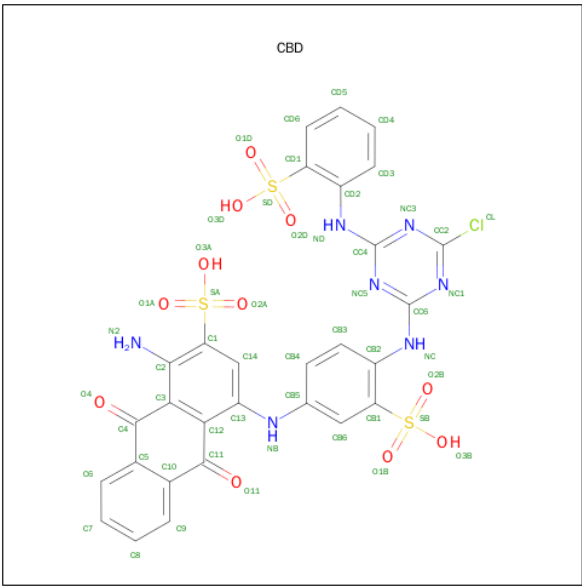
Chain	Residue	Modelled	Actual	Comment	Reference
B	-15	HIS	-	EXPRESSION TAG	UNP P15559
B	-14	HIS	-	EXPRESSION TAG	UNP P15559
B	-13	HIS	-	EXPRESSION TAG	UNP P15559
B	-12	HIS	-	EXPRESSION TAG	UNP P15559
B	-11	HIS	-	EXPRESSION TAG	UNP P15559
B	-10	HIS	-	EXPRESSION TAG	UNP P15559
B	-9	SER	-	EXPRESSION TAG	UNP P15559
B	-8	SER	-	EXPRESSION TAG	UNP P15559
B	-7	GLY	-	EXPRESSION TAG	UNP P15559
B	-6	LEU	-	EXPRESSION TAG	UNP P15559
B	-5	VAL	-	EXPRESSION TAG	UNP P15559
B	-4	PRO	-	EXPRESSION TAG	UNP P15559
B	-3	ARG	-	EXPRESSION TAG	UNP P15559
B	-2	GLY	-	EXPRESSION TAG	UNP P15559
B	-1	SER	-	EXPRESSION TAG	UNP P15559
B	0	HIS	-	EXPRESSION TAG	UNP P15559
B	187	SER	PRO	ENGINEERED MUTATION	UNP P15559

- 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 CIBACRON BLUE (three-letter code: CBD) (formula: C<sub>29</sub>H<sub>20</sub>ClN<sub>7</sub>O<sub>11</sub>S<sub>3</sub>).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total	C	Cl	N	O	S	0	0
			51	29	1	7	11	3		
3	B	1	Total	C	Cl	N	O	S	0	0
			51	29	1	7	11	3		

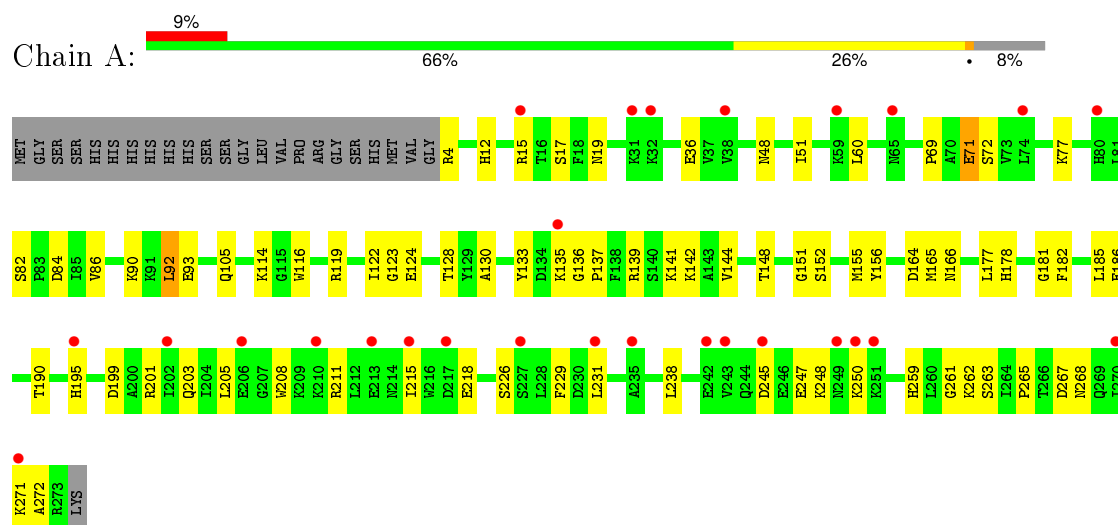
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	4	Total	O	0	0
			4	4		
4	B	3	Total	O	0	0
			3	3		

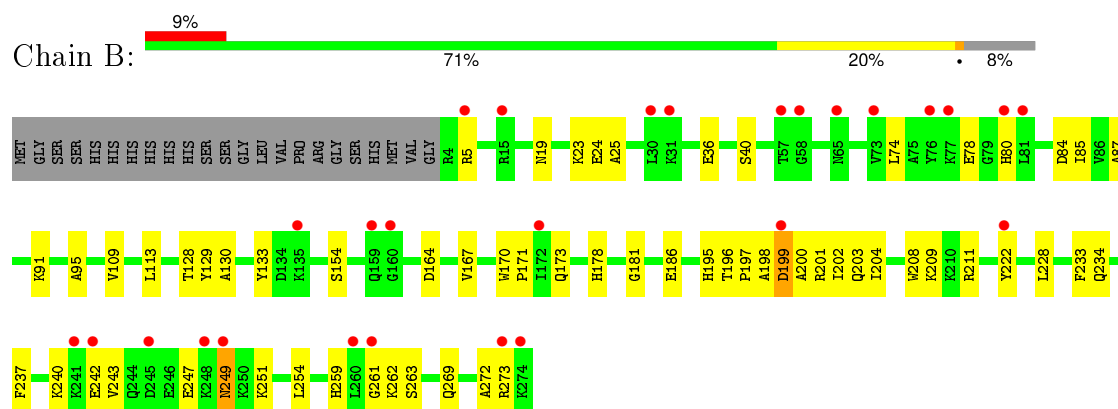
### 3 Residue-property plots [i](#)

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

#### • Molecule 1: NAD(P)H DEHYDROGENASE [QUINONE] 1



#### • Molecule 1: NAD(P)H DEHYDROGENASE [QUINONE] 1



## 4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	104.17Å 104.56Å 118.57Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.22 – 2.69 46.22 – 2.69	Depositor EDS
% Data completeness (in resolution range)	99.4 (46.22-2.69) 99.4 (46.22-2.69)	Depositor EDS
$R_{merge}$	0.16	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.60 (at 2.69Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, $R_{free}$	0.185 , 0.207 0.180 , 0.202	Depositor DCC
$R_{free}$ test set	928 reflections (5.39%)	DCC
Wilson B-factor (Å <sup>2</sup> )	48.0	Xtriage
Anisotropy	0.151	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 37.4	EDS
Estimated twinning fraction	0.390 for -K,-H,-L 0.358 for -k,-h,-l	Xtriage
Reported twinning fraction	0.390 for -K,-H,-L	Depositor
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.39$ , $\langle L^2 \rangle = 0.22$	Xtriage
Outliers	0 of 18137 reflections	Xtriage
$F_o, F_c$ correlation	0.84	EDS
Total number of atoms	4531	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 48.56 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 8.4505e-05.*

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

<sup>2</sup> Theoretical 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, CBD

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	0.24	0/2210	0.37	0/2985
1	B	0.24	0/2220	0.37	0/2996
All	All	0.24	0/4430	0.37	0/5981

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-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 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	2153	0	2149	51	0
1	B	2163	0	2162	41	1
2	A	53	0	31	4	0
2	B	53	0	31	2	0
3	A	51	0	20	8	0
3	B	51	0	20	4	0
4	A	4	0	0	0	0
4	B	3	0	0	0	0
All	All	4531	0	4413	95	1

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 11.

All (95) 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:23:LYS:NZ	1:B:24:GLU:OE2	2.10	0.84
1:B:247:GLU:O	1:B:262:LYS:NZ	2.15	0.80
1:A:92:LEU:O	1:A:141:LYS:NZ	2.20	0.74
1:B:196:THR:O	1:B:201:ARG:NH1	2.20	0.74
1:A:12:HIS:NE2	2:A:1274:FAD:O2P	2.23	0.72
1:A:19:ASN:ND2	2:A:1274:FAD:O1P	2.22	0.71
1:A:77:LYS:HE3	1:A:124:GLU:HG3	1.74	0.68
1:B:19:ASN:ND2	2:B:1275:FAD:O1P	2.25	0.67
1:A:208:TRP:HZ3	1:A:211:ARG:HH11	1.41	0.67
1:B:249:ASN:HD22	1:B:249:ASN:H	1.45	0.65
3:A:1275:CBD:NC1	3:A:1275:CBD:HB3	2.11	0.64
1:A:267:ASP:HB3	1:A:271:LYS:HB2	1.80	0.63
1:A:114:LYS:NZ	1:B:109:VAL:O	2.28	0.62
1:A:128:THR:HG22	1:A:130:ALA:H	1.66	0.60
1:B:259:HIS:CD2	1:B:261:GLY:H	2.19	0.59
1:B:84:ASP:OD1	1:B:85:ILE:N	2.35	0.59
1:A:263:SER:OG	1:B:261:GLY:O	2.20	0.59
1:B:5:ARG:HB2	1:B:95:ALA:HA	1.84	0.59
1:B:222:TYR:HB2	1:B:273:ARG:HG2	1.83	0.58
1:A:247:GLU:O	1:A:262:LYS:NZ	2.37	0.58
1:B:173:GLN:NE2	1:B:269:GLN:OE1	2.36	0.57
1:A:248:LYS:HB2	1:A:248:LYS:NZ	2.21	0.55
1:B:154:SER:OG	3:B:1276:CBD:O3D	2.19	0.54
1:A:259:HIS:CD2	1:A:261:GLY:H	2.26	0.53
1:A:259:HIS:O	1:A:262:LYS:HG2	2.09	0.52
1:A:164:ASP:OD2	1:A:166:ASN:HB2	2.10	0.52
1:A:93:GLU:HA	1:A:141:LYS:NZ	2.25	0.51
1:A:156:TYR:HB3	1:A:165:MET:HB2	1.92	0.51
1:B:133:TYR:O	1:B:181:GLY:HA2	2.09	0.51
3:B:1276:CBD:HB4	3:B:1276:CBD:H14	1.93	0.51
1:A:60:LEU:HD22	1:A:71:GLU:HG2	1.92	0.50
1:A:265:PRO:HB2	1:A:268:ASN:HB2	1.93	0.50
1:A:245:ASP:HA	1:A:248:LYS:NZ	2.28	0.49
1:A:238:LEU:HD21	1:B:154:SER:HB2	1.95	0.49
1:B:25:ALA:HB1	1:B:209:LYS:HD2	1.95	0.49
1:A:259:HIS:HD2	1:A:261:GLY:H	1.61	0.48
1:A:247:GLU:HA	1:A:250:LYS:HG3	1.94	0.48
3:B:1276:CBD:NB	3:B:1276:CBD:O11	2.44	0.48
1:A:84:ASP:N	1:A:84:ASP:OD1	2.42	0.48
1:A:155:MET:HG2	1:B:237:PHE:HB2	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:93:GLU:HA	1:A:141:LYS:HZ1	1.79	0.47
1:A:152:SER:H	1:A:155:MET:HE3	1.80	0.47
1:A:148:THR:HA	1:A:190:THR:HB	1.95	0.47
1:A:226:SER:HA	1:A:229:PHE:HD1	1.80	0.46
1:B:109:VAL:HB	1:B:113:LEU:HD23	1.96	0.46
1:B:186:GLU:O	1:B:211:ARG:NH2	2.36	0.46
1:B:233:PHE:CE1	1:B:237:PHE:HE2	2.33	0.46
1:B:78:GLU:HB2	1:B:80:HIS:HD2	1.80	0.46
1:B:128:THR:HG22	1:B:130:ALA:H	1.81	0.45
1:B:87:ALA:O	1:B:91:LYS:HG3	2.16	0.45
1:B:78:GLU:HB2	1:B:80:HIS:CD2	2.51	0.45
1:B:199:ASP:O	1:B:203:GLN:HG2	2.16	0.45
1:A:186:GLU:OE2	1:A:272:ALA:N	2.39	0.45
1:B:249:ASN:HD22	1:B:249:ASN:N	2.11	0.45
3:A:1275:CBD:NB	3:A:1275:CBD:O11	2.47	0.45
1:A:177:LEU:O	1:A:182:PHE:HB2	2.17	0.45
1:B:200:ALA:O	1:B:204:ILE:HG13	2.18	0.44
1:A:48:ASN:O	1:A:119:ARG:NH1	2.48	0.44
1:A:86:VAL:HG12	1:A:90:LYS:HE3	2.00	0.44
1:B:197:PRO:O	1:B:201:ARG:HG3	2.17	0.43
3:A:1275:CBD:HB4	1:B:129:TYR:CE1	2.54	0.43
3:B:1276:CBD:HB4	3:B:1276:CBD:C14	2.47	0.43
1:A:92:LEU:HD11	1:A:116:TRP:CH2	2.53	0.43
1:A:82:SER:HB3	1:A:84:ASP:OD1	2.19	0.43
1:B:164:ASP:HB3	1:B:167:VAL:HG23	1.99	0.43
1:B:5:ARG:HE	1:B:36:GLU:HB3	1.83	0.43
1:B:74:LEU:HD11	1:B:78:GLU:OE2	2.18	0.43
1:A:195:HIS:HE1	3:A:1275:CBD:CB2	2.32	0.43
1:A:133:TYR:HD1	1:A:178:HIS:CD2	2.38	0.42
3:A:1275:CBD:O1B	3:A:1275:CBD:NC	2.51	0.42
3:A:1275:CBD:ND	3:A:1275:CBD:O2D	2.53	0.42
1:A:133:TYR:O	1:A:181:GLY:HA2	2.19	0.42
1:B:186:GLU:OE2	1:B:272:ALA:N	2.44	0.41
1:B:84:ASP:OD1	1:B:85:ILE:HG13	2.20	0.41
1:B:228:LEU:HB3	1:B:243:VAL:HG11	2.02	0.41
1:B:170:TRP:HB3	1:B:171:PRO:HD3	2.02	0.41
1:A:105:GLN:HA	2:A:1274:FAD:C5X	2.50	0.41
1:A:186:GLU:O	1:A:211:ARG:NH2	2.44	0.41
1:B:208:TRP:HZ3	1:B:211:ARG:HD3	1.85	0.41
1:A:123:GLY:O	1:A:124:GLU:HB2	2.21	0.41
3:A:1275:CBD:HB4	1:B:129:TYR:CZ	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:240:LYS:HB2	1:B:243:VAL:HG23	2.02	0.41
1:A:51:ILE:HG22	1:A:119:ARG:HG2	2.03	0.41
1:B:198:ALA:O	1:B:202:ILE:HG13	2.20	0.41
1:A:136:GLY:HA2	1:A:137:PRO:HD3	1.93	0.41
1:A:142:LYS:HD3	1:A:185:LEU:HD21	2.03	0.41
1:A:199:ASP:O	1:A:203:GLN:HG2	2.21	0.41
1:A:201:ARG:O	1:A:205:LEU:HD12	2.21	0.40
1:A:69:PRO:HG3	2:B:1275:FAD:HM81	2.02	0.40
1:A:17:SER:HB2	2:A:1274:FAD:O2P	2.21	0.40
1:A:151:GLY:HA2	3:A:1275:CBP:HB6	2.02	0.40
1:B:133:TYR:HA	1:B:178:HIS:O	2.21	0.40
1:A:135:LYS:HA	1:A:139:ARG:HD2	2.02	0.40
1:A:215:ILE:HA	1:A:218:GLU:HG3	2.04	0.40
1:A:72:SER:HB2	1:A:122:ILE:HD12	2.03	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	Interatomic distance (Å)	Clash overlap (Å)
1:B:251:LYS:NZ	1:B:263:SER:OG[2_555]	2.04	0.16

## 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	268/294 (91%)	250 (93%)	18 (7%)	0	100	100
1	B	269/294 (92%)	252 (94%)	17 (6%)	0	100	100
All	All	537/588 (91%)	502 (94%)	35 (6%)	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	225/245 (92%)	218 (97%)	7 (3%)	47	78
1	B	226/245 (92%)	219 (97%)	7 (3%)	47	78
All	All	451/490 (92%)	437 (97%)	14 (3%)	47	78

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	ARG
1	A	15	ARG
1	A	36	GLU
1	A	71	GLU
1	A	92	LEU
1	A	144	VAL
1	A	231	LEU
1	B	40	SER
1	B	195	HIS
1	B	199	ASP
1	B	234	GLN
1	B	242	GLU
1	B	249	ASN
1	B	254	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	162	HIS
1	A	249	ASN
1	B	65	ASN
1	B	80	HIS
1	B	105	GLN
1	B	249	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 ⓘ

4 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	1274	-	48,58,58	1.33	5 (10%)	54,89,89	2.16	10 (18%)
3	CBD	A	1275	-	55,56,56	1.72	9 (16%)	76,87,87	2.56	19 (25%)
2	FAD	B	1275	-	48,58,58	1.31	5 (10%)	54,89,89	2.12	12 (22%)
3	CBD	B	1276	-	55,56,56	1.74	11 (20%)	76,87,87	2.46	18 (23%)

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	1274	-	-	0/30/50/50	0/6/6/6
3	CBD	A	1275	-	-	0/30/46/46	0/6/6/6
2	FAD	B	1275	-	-	0/30/50/50	0/6/6/6
3	CBD	B	1276	-	-	0/30/46/46	0/6/6/6

All (30) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1276	CBD	C5-C4	-4.35	1.39	1.48
3	A	1275	CBD	C5-C4	-4.11	1.39	1.48
3	A	1275	CBD	C10-C11	-3.94	1.40	1.48
3	B	1276	CBD	C10-C11	-3.67	1.40	1.48
3	A	1275	CBD	C3-C4	-3.44	1.39	1.47
3	B	1276	CBD	C3-C4	-3.44	1.39	1.47
3	A	1275	CBD	C12-C11	-3.23	1.39	1.47
3	B	1276	CBD	C12-C11	-2.96	1.40	1.47
3	A	1275	CBD	CC2-NC3	2.07	1.35	1.32
3	A	1275	CBD	C2-N2	2.13	1.43	1.37
3	B	1276	CBD	C2-N2	2.17	1.43	1.37
3	A	1275	CBD	CB2-NC	2.42	1.46	1.39
3	B	1276	CBD	CB2-NC	2.47	1.46	1.39
3	B	1276	CBD	CC2-NC3	2.56	1.35	1.32
3	B	1276	CBD	CC2-NC1	2.63	1.35	1.32
3	B	1276	CBD	CB5-NB	2.68	1.46	1.40
2	A	1274	FAD	C8-C7	2.89	1.48	1.41
2	B	1275	FAD	C8-C7	2.92	1.48	1.41
2	B	1275	FAD	C5A-C4A	3.04	1.47	1.40
2	A	1274	FAD	C5A-C4A	3.09	1.47	1.40
2	B	1275	FAD	C9A-C5X	3.36	1.49	1.42
2	A	1274	FAD	C9A-C5X	3.48	1.49	1.42
2	A	1274	FAD	C4-C4X	3.67	1.48	1.41
2	B	1275	FAD	C4-C4X	3.69	1.48	1.41
2	B	1275	FAD	C4X-C10	3.91	1.48	1.41
2	A	1274	FAD	C4X-C10	3.93	1.48	1.41
3	A	1275	CBD	CC4-ND	4.39	1.43	1.36
3	B	1276	CBD	CC4-ND	4.69	1.44	1.36
3	B	1276	CBD	CC6-NC	4.84	1.44	1.36
3	A	1275	CBD	CC6-NC	5.59	1.46	1.36

All (59) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1275	CBD	NC3-CC2-NC1	-7.46	119.28	129.68
2	A	1274	FAD	N3A-C2A-N1A	-7.03	123.51	128.89
2	B	1275	FAD	N3A-C2A-N1A	-6.89	123.62	128.89
3	B	1276	CBD	NC3-CC2-NC1	-6.73	120.29	129.68
2	A	1274	FAD	C4-C4X-C10	-5.67	116.31	119.94
2	B	1275	FAD	C4-C4X-C10	-5.31	116.55	119.94
3	A	1275	CBD	NC5-CC6-NC1	-5.03	118.49	126.22
3	A	1275	CBD	CB5-NB-C13	-4.54	113.88	126.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1276	CBD	NC5-CC6-NC1	-4.11	119.91	126.22
3	B	1276	CBD	NC5-CC4-NC3	-4.07	119.96	126.22
2	A	1274	FAD	C4X-C4-N3	-3.77	118.43	123.59
2	B	1275	FAD	P-O3P-PA	-3.77	122.14	132.73
3	B	1276	CBD	CD2-ND-CC4	-3.75	119.34	129.22
2	B	1275	FAD	C4X-C4-N3	-3.65	118.60	123.59
3	A	1275	CBD	NC5-CC4-NC3	-3.54	120.78	126.22
3	B	1276	CBD	C12-C13-NB	-3.38	117.04	120.47
2	B	1275	FAD	C4A-C5A-N7A	-3.32	106.43	109.48
2	A	1274	FAD	C4A-C5A-N7A	-3.15	106.58	109.48
2	A	1274	FAD	P-O3P-PA	-3.12	123.98	132.73
3	A	1275	CBD	C3-C12-C13	-3.11	117.99	119.88
3	B	1276	CBD	CB3-CB2-NC	-2.39	116.04	121.56
3	A	1275	CBD	CD2-ND-CC4	-2.21	123.39	129.22
2	B	1275	FAD	C2B-C1B-N9A	-2.21	110.92	114.29
2	B	1275	FAD	C4X-C10-N10	-2.19	119.23	120.52
3	A	1275	CBD	CD2-CD1-SD	-2.11	119.00	122.64
3	A	1275	CBD	O1B-SB-CB1	2.03	108.51	106.17
2	B	1275	FAD	O4B-C1B-N9A	2.20	112.70	108.10
3	B	1276	CBD	CB5-CB6-CB1	2.26	121.65	119.27
3	A	1275	CBD	O1D-SD-CD1	2.31	108.84	106.17
3	A	1275	CBD	O2B-SB-CB1	2.40	108.94	106.17
3	A	1275	CBD	CD6-CD1-SD	2.47	121.14	117.51
3	A	1275	CBD	O1A-SA-C1	2.56	109.12	106.17
2	A	1274	FAD	C1'-N10-C9A	2.59	121.77	118.86
2	B	1275	FAD	C4-C4X-N5	2.64	121.92	118.72
3	A	1275	CBD	O2A-SA-C1	2.70	109.29	106.17
3	B	1276	CBD	O1A-SA-C1	2.71	109.30	106.17
2	A	1274	FAD	C5X-C9A-N10	2.74	119.70	117.62
2	B	1275	FAD	C5X-C9A-N10	2.84	119.78	117.62
2	A	1274	FAD	C4-C4X-N5	2.87	122.20	118.72
3	A	1275	CBD	O2D-SD-CD1	2.94	109.56	106.17
3	B	1276	CBD	O1D-SD-CD1	3.00	109.62	106.17
2	A	1274	FAD	C4X-N5-C5X	3.25	120.50	116.76
3	B	1276	CBD	O2B-SB-CB1	3.30	109.98	106.17
3	B	1276	CBD	O2A-SA-C1	3.35	110.03	106.17
3	B	1276	CBD	CL-CC2-NC3	3.38	119.16	115.08
2	B	1275	FAD	C4X-N5-C5X	3.43	120.71	116.76
3	A	1275	CBD	CC6-NC5-CC4	4.06	119.95	113.99
3	A	1275	CBD	CL-CC2-NC1	4.26	120.22	115.08
3	B	1276	CBD	O1B-SB-CB1	4.28	111.10	106.17
3	B	1276	CBD	CC6-NC5-CC4	4.31	120.31	113.99

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1275	CBD	CL-CC2-NC3	4.49	120.50	115.08
3	B	1276	CBD	CL-CC2-NC1	4.53	120.55	115.08
3	B	1276	CBD	O2D-SD-CD1	4.74	111.64	106.17
2	B	1275	FAD	C4-N3-C2	7.91	122.08	115.25
2	A	1274	FAD	C4-N3-C2	8.17	122.31	115.25
3	B	1276	CBD	CC6-NC1-CC2	9.27	119.62	112.49
3	A	1275	CBD	CC4-NC3-CC2	9.63	119.90	112.49
3	B	1276	CBD	CC4-NC3-CC2	9.64	119.91	112.49
3	A	1275	CBD	CC6-NC1-CC2	11.85	121.61	112.49

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 18 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1274	FAD	4	0
3	A	1275	CBD	8	0
2	B	1275	FAD	2	0
3	B	1276	CBD	4	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, 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	270/294 (91%)	0.87	27 (10%) 9 7	28, 40, 58, 68	0
1	B	271/294 (92%)	0.86	27 (9%) 9 7	29, 42, 56, 62	0
All	All	541/588 (92%)	0.87	54 (9%) 9 7	28, 41, 57, 68	0

All (54) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	32	LYS	5.4
1	A	250	LYS	5.3
1	A	210	LYS	5.1
1	B	274	LYS	4.6
1	A	245	ASP	4.5
1	A	251	LYS	4.3
1	A	243	VAL	3.8
1	A	235	ALA	3.7
1	B	241	LYS	3.7
1	A	206	GLU	3.7
1	A	249	ASN	3.6
1	B	261	GLY	3.5
1	A	38	VAL	3.4
1	B	273	ARG	3.4
1	A	231	LEU	3.3
1	B	57	THR	3.3
1	B	81	LEU	3.3
1	B	159	GLN	3.2
1	B	31	LYS	3.2
1	B	245	ASP	3.1
1	B	15	ARG	3.0
1	B	260	LEU	3.0
1	B	222	TYR	2.9
1	A	80	HIS	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	227	SER	2.9
1	B	249	ASN	2.9
1	B	80	HIS	2.7
1	A	202	ILE	2.7
1	A	65	ASN	2.7
1	B	58	GLY	2.7
1	A	31	LYS	2.6
1	B	30	LEU	2.6
1	B	248	LYS	2.6
1	A	270	ILE	2.6
1	A	74	LEU	2.5
1	A	59	LYS	2.5
1	B	242	GLU	2.5
1	A	135	LYS	2.5
1	A	242	GLU	2.3
1	B	160	GLY	2.3
1	A	217	ASP	2.3
1	B	199	ASP	2.3
1	B	65	ASN	2.3
1	B	73	VAL	2.3
1	B	76	TYR	2.2
1	A	195	HIS	2.2
1	B	5	ARG	2.1
1	B	77	LYS	2.1
1	A	215	ILE	2.1
1	A	213	GLU	2.1
1	A	271	LYS	2.0
1	B	172	ILE	2.0
1	B	135	LYS	2.0
1	A	15	ARG	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, 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	RSCC	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
3	CBD	B	1276	51/51	0.65	0.37	2.22	36,54,63,77	51
3	CBD	A	1275	51/51	0.81	0.26	0.37	34,47,64,64	51
2	FAD	B	1275	53/53	0.86	0.23	-0.04	29,37,43,45	0
2	FAD	A	1274	53/53	0.88	0.20	-0.40	31,33,39,41	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.