



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:14 AM GMT

PDB ID : 2CDU  
Title : THE CRYSTAL STRUCTURE OF WATER-FORMING NAD(P)H OXIDASE FROM LACTOBACILLUS SANFRANCISCENSIS  
Authors : Lountos, G.T.; Jiang, R.; Wellborn, W.B.; Thaler, T.L.; Bommarius, A.S.; Orville, A.M.  
Deposited on : 2006-01-28  
Resolution : 1.80 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.

---

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

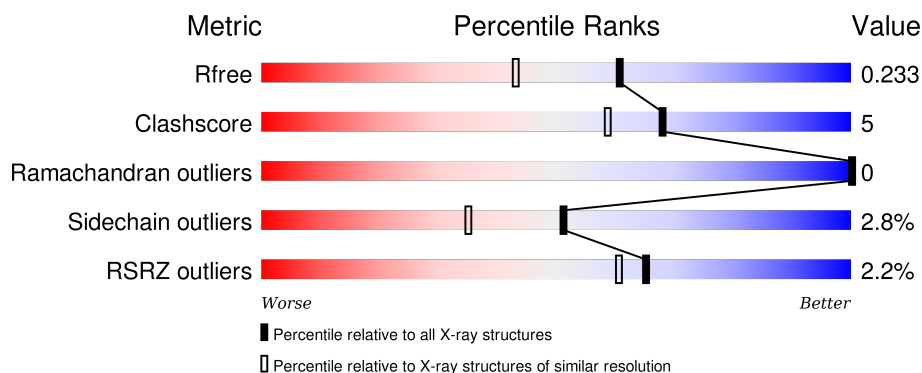
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*



The reported resolution of this entry is 1.80 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	4533 (1.80-1.80)
Clashscore	102246	5383 (1.80-1.80)
Ramachandran outliers	100387	5320 (1.80-1.80)
Sidechain outliers	100360	5319 (1.80-1.80)
RSRZ outliers	91569	4547 (1.80-1.80)

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	452	
1	B	452	

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
2	FAD	A	500[*]	-	-	-	X
2	FAD	B	500[*]	-	-	-	X
3	ADP	B	501	-	-	-	X

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 7846 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 NADPH OXIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	451	Total	C	N	O	S	0	4	0
			3493	2197	587	692	17			
1	B	449	Total	C	N	O	S	0	8	0
			3485	2191	585	691	18			

- 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	1
			53	27	9	15	2		
2	B	1	Total	C	N	O	P	0	1
			53	27	9	15	2		

- Molecule 3 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula:  $C_{10}H_{15}N_5O_{10}P_2$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	B	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

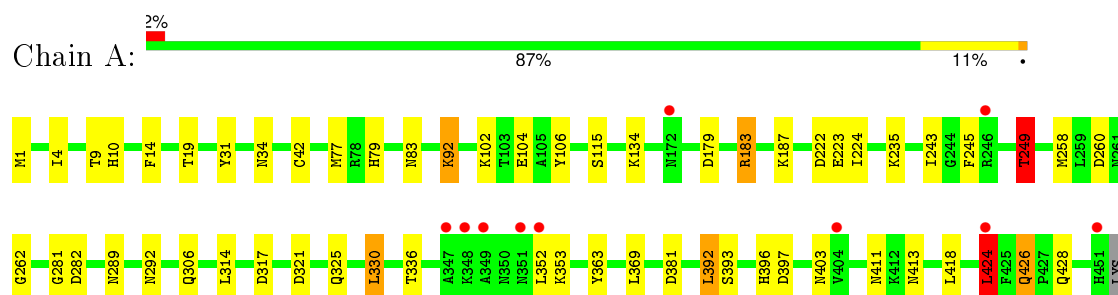
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	382	Total	O	0	0
			382	382		
4	B	326	Total	O	0	0
			326	326		

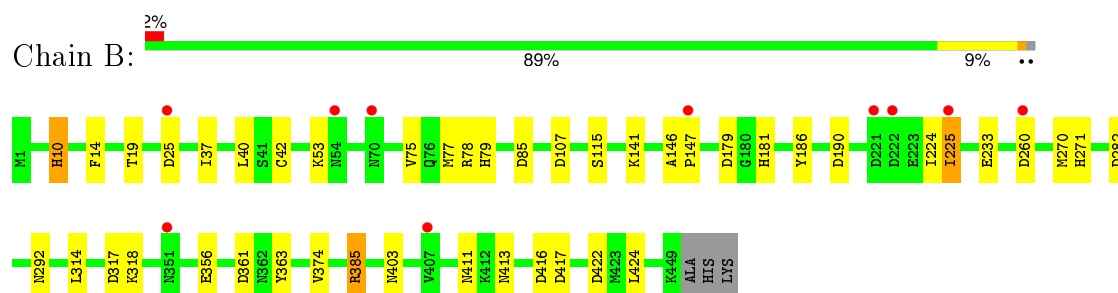
### 3 Residue-property plots

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: NADPH OXIDASE



#### • Molecule 1: NADPH OXIDASE



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	59.64Å 92.64Å 163.47Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	81.65 – 1.80 40.30 – 1.80	Depositor EDS
% Data completeness (in resolution range)	96.1 (81.65-1.80) 96.0 (40.30-1.80)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.56 (at 1.79Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, $R_{free}$	0.178 , 0.223 0.189 , 0.233	Depositor DCC
$R_{free}$ test set	4053 reflections (5.24%)	DCC
Wilson B-factor (Å <sup>2</sup> )	23.8	Xtriage
Anisotropy	0.613	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 39.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	1 of 81424 reflections (0.001%)	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	7846	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.65% of the height of the origin peak. No significant pseudotranslation is detected.*

<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: CSX, ADP, 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	0.66	0/3555	0.89	14/4818 (0.3%)
1	B	0.64	0/3572	0.84	10/4840 (0.2%)
All	All	0.65	0/7127	0.86	24/9658 (0.2%)

There are no bond length outliers.

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	183	ARG	NE-CZ-NH2	-9.39	115.60	120.30
1	A	183	ARG	NE-CZ-NH1	9.32	124.96	120.30
1	B	385	ARG	NE-CZ-NH2	-9.31	115.65	120.30
1	A	392	LEU	CA-CB-CG	8.66	135.21	115.30
1	A	330	LEU	CA-CB-CG	8.40	134.61	115.30
1	A	282	ASP	CB-CG-OD2	8.03	125.53	118.30
1	B	416	ASP	CB-CG-OD2	6.91	124.52	118.30
1	A	381	ASP	CB-CG-OD2	6.76	124.38	118.30
1	B	385	ARG	NE-CZ-NH1	6.65	123.62	120.30
1	B	179	ASP	CB-CG-OD2	5.94	123.65	118.30
1	B	417	ASP	CB-CG-OD2	5.77	123.49	118.30
1	A	260	ASP	CB-CG-OD2	5.73	123.46	118.30
1	A	424	LEU	CA-CB-CG	5.61	128.19	115.30
1	A	249	THR	N-CA-CB	-5.59	99.67	110.30
1	A	317	ASP	CB-CG-OD2	5.44	123.19	118.30
1	A	222	ASP	CB-CG-OD2	5.42	123.17	118.30
1	B	260	ASP	CB-CG-OD2	5.39	123.15	118.30
1	B	85	ASP	CB-CG-OD2	5.35	123.11	118.30
1	A	392	LEU	CB-CG-CD2	5.23	119.89	111.00
1	A	179	ASP	CB-CG-OD2	5.22	123.00	118.30
1	B	361	ASP	CB-CG-OD2	5.09	122.88	118.30
1	B	107	ASP	CB-CG-OD2	5.09	122.88	118.30

*Continued on next page...*



*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	78	ARG	NE-CZ-NH1	5.04	122.82	120.30
1	A	321	ASP	CB-CG-OD1	5.03	122.82	118.30

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	3493	0	3464	43	0
1	B	3485	0	3458	30	0
2	A	53	0	8	0	0
2	B	53	0	8	0	0
3	A	27	0	12	1	0
3	B	27	0	12	0	0
4	A	382	0	0	12	1
4	B	326	0	0	6	1
All	All	7846	0	6962	67	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 5.

All (67) 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:289:ASN:HD22	1:A:325:GLN:HE22	1.09	0.99
1:A:249:THR:HG21	1:A:262:GLY:O	1.68	0.93
1:A:411:ASN:HD21	1:B:413:ASN:HD21	1.20	0.85
1:A:363:TYR:CZ	1:A:396[B]:HIS:HD2	1.96	0.84
1:A:289:ASN:HD22	1:A:325:GLN:NE2	1.80	0.80
1:A:413:ASN:HD21	1:B:411:ASN:HD21	1.29	0.78
1:B:25:ASP:OD1	4:B:2014:HOH:O	2.02	0.77
1:B:318:LYS:NZ	4:B:2224:HOH:O	2.21	0.74
1:A:363:TYR:CE1	1:A:396[B]:HIS:HD2	2.06	0.73
1:A:223:GLU:OE2	1:A:235:LYS:NZ	2.21	0.73

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:37:ILE:HG23	1:B:77[A]:MET:HE2	1.70	0.72
1:B:270:MET:SD	4:B:2194:HOH:O	2.48	0.72
1:B:10[A]:HIS:ND1	4:B:2006:HOH:O	2.23	0.71
1:A:10[A]:HIS:ND1	4:A:2006:HOH:O	2.24	0.69
1:A:403:ASN:HD21	1:B:424:LEU:H	1.38	0.68
1:A:10[A]:HIS:CE1	4:A:2006:HOH:O	2.46	0.67
1:A:428[B]:GLN:OE1	4:A:2352:HOH:O	2.14	0.66
1:A:281:GLY:H	1:A:306:GLN:HE21	1.43	0.65
1:A:34:ASN:ND2	4:A:2024:HOH:O	2.30	0.64
1:A:187:LYS:NZ	4:A:2173:HOH:O	2.29	0.64
1:B:225:ILE:HD12	1:B:233:GLU:HB3	1.81	0.61
1:A:31:TYR:OH	1:A:104:GLU:OE2	2.09	0.60
1:A:363:TYR:CE1	1:A:396[B]:HIS:CD2	2.88	0.59
1:A:396[B]:HIS:ND1	1:A:397:ASP:N	2.50	0.59
1:B:53:LYS:NZ	4:B:2037:HOH:O	2.37	0.57
1:A:424:LEU:H	1:B:403:ASN:HD21	1.53	0.56
1:A:34:ASN:CG	4:A:2026:HOH:O	2.44	0.56
1:A:183:ARG:HD3	4:A:2168:HOH:O	2.06	0.55
1:A:281:GLY:H	1:A:306:GLN:NE2	2.05	0.54
1:A:249:THR:HG23	1:A:258:MET:CE	2.38	0.53
1:B:40:LEU:HD22	1:B:42[A]:CSX:OD	2.08	0.53
1:A:363:TYR:CZ	1:A:396[B]:HIS:CD2	2.87	0.53
1:A:83:ASN:HB3	1:A:92:LYS:HB2	1.90	0.53
1:B:37:ILE:HG23	1:B:77[A]:MET:CE	2.39	0.52
1:A:79:HIS:HE1	4:A:2093:HOH:O	1.92	0.51
1:B:146:ALA:HB3	1:B:147:PRO:HD3	1.92	0.51
1:B:19:THR:HG23	1:B:314:LEU:HD22	1.95	0.49
1:B:271:HIS:HE1	1:B:317:ASP:OD1	1.96	0.48
1:A:134:LYS:HG3	1:A:245:PHE:CE2	2.48	0.48
1:A:249:THR:HG22	4:A:2227:HOH:O	2.13	0.47
1:A:77:MET:O	1:A:79:HIS:HD2	1.98	0.47
1:B:75:VAL:CG1	1:B:77[B]:MET:HE3	2.45	0.46
1:A:243:ILE:HA	3:A:501:ADP:H5'2	1.98	0.45
1:B:190:ASP:OD2	1:B:356:GLU:OE2	2.35	0.45
1:A:9:THR:HG22	4:A:2034:HOH:O	2.16	0.45
1:A:10[B]:HIS:CE1	4:A:2373:HOH:O	2.70	0.45
1:A:4:ILE:HD12	1:A:106:TYR:HB3	1.99	0.45
1:A:411:ASN:ND2	1:B:413:ASN:HD21	2.00	0.44
1:B:225:ILE:CD1	1:B:233:GLU:HB3	2.48	0.44
1:A:249:THR:HG23	1:A:258:MET:HE1	1.99	0.43
1:A:115:SER:HB2	1:A:245:PHE:HB3	1.99	0.43

*Continued on next page...*

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:10[A]:HIS:CD2	4:B:2005:HOH:O	2.72	0.43
1:B:115:SER:HB3	1:B:282:ASP:HB3	2.00	0.43
1:A:336:THR:O	1:A:393:SER:HA	2.19	0.42
1:B:363:TYR:HB3	1:B:374[B]:VAL:CG2	2.49	0.42
1:B:75:VAL:CG1	1:B:77[B]:MET:CE	2.98	0.41
1:A:325:GLN:HB2	1:B:422:ASP:OD2	2.20	0.41
1:A:19:THR:HG23	1:A:314:LEU:HD22	2.02	0.41
1:B:19:THR:HG23	1:B:314:LEU:CD2	2.51	0.41
1:A:292:ASN:HB2	4:A:2256:HOH:O	2.22	0.40
1:B:181:HIS:CD2	1:B:186:TYR:CD2	3.10	0.40
1:A:289:ASN:HB2	1:A:325:GLN:HE21	1.86	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)
4:A:2121:HOH:O	4:B:2074:HOH:O[2_554]	2.11	0.09

## 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	451/452 (100%)	437 (97%)	14 (3%)	0	100	100
1	B	453/452 (100%)	441 (97%)	12 (3%)	0	100	100
All	All	904/904 (100%)	878 (97%)	26 (3%)	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	384/383 (100%)	370 (96%)	14 (4%)	42	24
1	B	387/383 (101%)	379 (98%)	8 (2%)	61	47
All	All	771/766 (101%)	749 (97%)	22 (3%)	51	34

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

Mol	Chain	Res	Type
1	A	1	MET
1	A	14	PHE
1	A	92	LYS
1	A	102	LYS
1	A	224	ILE
1	A	249	THR
1	A	330	LEU
1	A	352	LEU
1	A	353	LYS
1	A	369	LEU
1	A	392	LEU
1	A	418	LEU
1	A	424	LEU
1	A	426	GLN
1	B	10[A]	HIS
1	B	10[B]	HIS
1	B	14	PHE
1	B	141	LYS
1	B	224	ILE
1	B	225	ILE
1	B	292	ASN
1	B	385	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	34	ASN
1	A	79	HIS
1	A	271	HIS
1	A	306	GLN
1	A	325	GLN
1	A	403	ASN
1	A	411	ASN
1	A	426	GLN
1	A	451	HIS
1	B	79	HIS
1	B	271	HIS
1	B	292	ASN
1	B	403	ASN
1	B	411	ASN
1	B	428	GLN
1	B	446	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

4 non-standard protein/DNA/RNA residues 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	$\# Z  > 2$
1	CSX	A	42[A]	-	3,6,7	0.58	0	3,6,8	1.00	0
1	CSX	A	42[B]	-	3,6,7	0.58	0	3,6,8	1.00	0
1	CSX	B	42[A]	-	3,6,7	0.53	0	3,6,8	1.15	0
1	CSX	B	42[B]	-	3,6,7	0.53	0	3,6,8	1.15	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means

no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	CSX	A	42[A]	-	-	0/1/5/7	0/0/0/0
1	CSX	A	42[B]	-	-	0/1/5/7	0/0/0/0
1	CSX	B	42[A]	-	-	0/1/5/7	0/0/0/0
1	CSX	B	42[B]	-	-	0/1/5/7	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	42[B]	CSX	1	0
1	B	42[A]	CSX	1	0
1	B	42[B]	CSX	2	0

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

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	500[*]	-	48,58,58	1.35	5 (10%)	54,89,89	2.20	11 (20%)
3	ADP	A	501	-	22,29,29	1.21	2 (9%)	27,45,45	2.48	5 (18%)
2	FAD	B	500[*]	-	48,58,58	1.67	9 (18%)	54,89,89	2.10	6 (11%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	ADP	B	501	-	22,29,29	1.20	3 (13%)	27,45,45	2.57	3 (11%)

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	500[*]	-	-	0/30/50/50	0/6/6/6
3	ADP	A	501	-	-	0/12/32/32	0/3/3/3
2	FAD	B	500[*]	-	-	0/30/50/50	0/6/6/6
3	ADP	B	501	-	-	0/12/32/32	0/3/3/3

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	500[*]	FAD	O5'-C5'	-2.18	1.35	1.44
2	B	500[*]	FAD	C7M-C7	2.01	1.55	1.51
2	B	500[*]	FAD	P-O5'	2.19	1.69	1.59
2	A	500[*]	FAD	C9A-N10	2.28	1.41	1.38
3	B	501	ADP	C2-N1	2.32	1.38	1.33
3	B	501	ADP	PB-O3B	2.48	1.63	1.54
2	A	500[*]	FAD	C5X-N5	2.48	1.39	1.35
2	B	500[*]	FAD	C5'-C4'	2.56	1.55	1.51
3	A	501	ADP	PB-O3B	2.58	1.64	1.54
2	B	500[*]	FAD	C10-N1	2.65	1.40	1.35
2	A	500[*]	FAD	C4-N3	3.03	1.38	1.33
3	A	501	ADP	C2-N3	3.25	1.37	1.32
2	A	500[*]	FAD	O4B-C1B	3.33	1.45	1.41
3	B	501	ADP	C2-N3	3.35	1.38	1.32
2	B	500[*]	FAD	C4-N3	3.92	1.40	1.33
2	B	500[*]	FAD	C4X-N5	4.24	1.40	1.33
2	B	500[*]	FAD	C1'-N10	4.29	1.52	1.48
2	B	500[*]	FAD	C5X-N5	4.34	1.42	1.35
2	A	500[*]	FAD	C4X-N5	4.34	1.40	1.33

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	501	ADP	N3-C2-N1	-11.86	119.81	128.89
3	A	501	ADP	N3-C2-N1	-10.93	120.53	128.89
2	B	500[*]	FAD	N3A-C2A-N1A	-9.76	121.42	128.89

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	500[*]	FAD	N3A-C2A-N1A	-8.87	122.11	128.89
2	A	500[*]	FAD	C4B-O4B-C1B	-4.63	104.64	109.72
2	B	500[*]	FAD	C4X-C4-N3	-3.74	118.48	123.59
3	B	501	ADP	C1'-N9-C4	-3.50	121.66	126.94
3	A	501	ADP	C1'-N9-C4	-3.29	121.98	126.94
2	A	500[*]	FAD	C4X-C4-N3	-3.22	119.19	123.59
2	A	500[*]	FAD	C9A-C5X-N5	-2.47	118.70	122.36
2	A	500[*]	FAD	O4B-C1B-N9A	-2.46	102.96	108.10
3	A	501	ADP	PA-O3A-PB	-2.45	124.44	132.67
3	B	501	ADP	PA-O3A-PB	-2.29	124.99	132.67
2	A	500[*]	FAD	C4A-C5A-N7A	-2.16	107.49	109.48
2	A	500[*]	FAD	O2A-PA-O3P	2.01	114.20	105.09
3	A	501	ADP	C2'-C3'-C4'	2.01	106.74	102.61
2	B	500[*]	FAD	O2P-P-O1P	2.07	123.74	112.53
2	B	500[*]	FAD	C4X-N5-C5X	2.24	119.34	116.76
3	A	501	ADP	O2B-PB-O3A	2.26	115.34	105.09
2	A	500[*]	FAD	C4-C4X-N5	2.66	121.95	118.72
2	B	500[*]	FAD	C5X-C9A-N10	2.83	119.77	117.62
2	A	500[*]	FAD	C4X-N5-C5X	3.78	121.11	116.76
2	A	500[*]	FAD	C5X-C9A-N10	3.89	120.57	117.62
2	A	500[*]	FAD	C4-N3-C2	7.90	122.08	115.25
2	B	500[*]	FAD	C4-N3-C2	8.17	122.31	115.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	501	ADP	1	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

### 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	450/452 (99%)	-0.13	10 (2%) 65 60	16, 24, 37, 56	0
1	B	448/452 (99%)	0.01	10 (2%) 65 60	15, 28, 42, 55	0
All	All	898/904 (99%)	-0.06	20 (2%) 65 60	15, 26, 41, 56	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	451	HIS	5.4
1	B	221	ASP	5.3
1	B	25	ASP	4.0
1	B	147	PRO	3.5
1	A	348	LYS	3.4
1	A	351	ASN	3.1
1	B	222	ASP	2.9
1	A	349	ALA	2.9
1	A	347	ALA	2.7
1	B	351	ASN	2.7
1	B	260	ASP	2.5
1	A	172	ASN	2.4
1	A	246	ARG	2.4
1	B	54	ASN	2.3
1	B	70	ASN	2.3
1	A	424	LEU	2.2
1	A	404	VAL	2.2
1	B	407	VAL	2.2
1	A	352	LEU	2.2
1	B	225	ILE	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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
1	CSX	A	42[A]	7/8	0.98	0.12	-	18,19,20,22	1
1	CSX	B	42[A]	7/8	0.95	0.13	-	18,25,25,26	1
1	CSX	B	42[B]	7/8	0.95	0.13	-	21,25,25,26	1
1	CSX	A	42[B]	7/8	0.98	0.12	-	18,19,21,22	1

## 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	RSCC	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	FAD	B	500[*]	53/53	0.87	0.15	3.88	20,24,27,31	9
3	ADP	B	501	27/27	0.85	0.17	3.05	37,43,59,61	0
2	FAD	A	500[*]	53/53	0.90	0.15	1.95	15,20,23,23	9
3	ADP	A	501	27/27	0.91	0.09	-0.16	22,30,48,49	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.