



# Full wwPDB X-ray Structure Validation Report ⓘ

Apr 18, 2014 – 05:19 AM EDT

PDB ID : 3D9G  
Title : Nitroalkane oxidase: wild type crystallized in a trapped state forming a cyanoadduct with FAD  
Authors : Heroux, A.; Bozinovski, D.M.; Valley, M.P.; Fitzpatrick, P.F.; Orville, A.M.  
Deposited on : 2008-05-27  
Resolution : 2.15 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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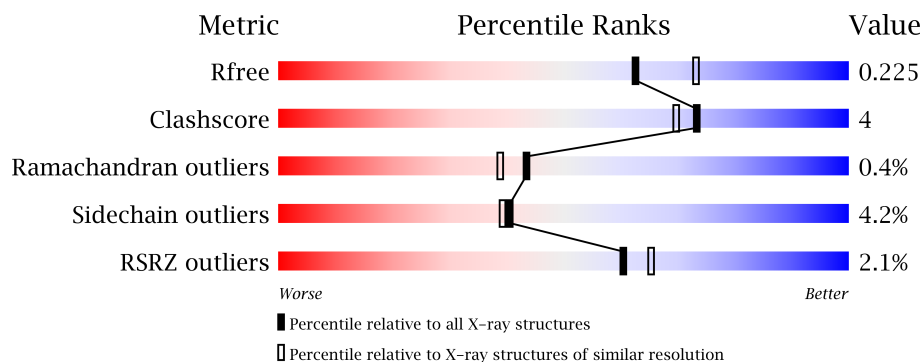
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.16 November 2013  
Xtriage (Phenix) : dev-1439  
EDS : stable22978  
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) : stable22978

# 1 Overall quality at a glance

The reported resolution of this entry is 2.15 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	1094 (2.18-2.14)
Clashscore	79885	1299 (2.18-2.14)
Ramachandran outliers	78287	1272 (2.18-2.14)
Sidechain outliers	78261	1272 (2.18-2.14)
RSRZ outliers	66119	1094 (2.18-2.14)

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	438	
1	B	438	
1	C	438	
1	D	438	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
4	GOL	D	701	-	X

## 2 Entry composition i

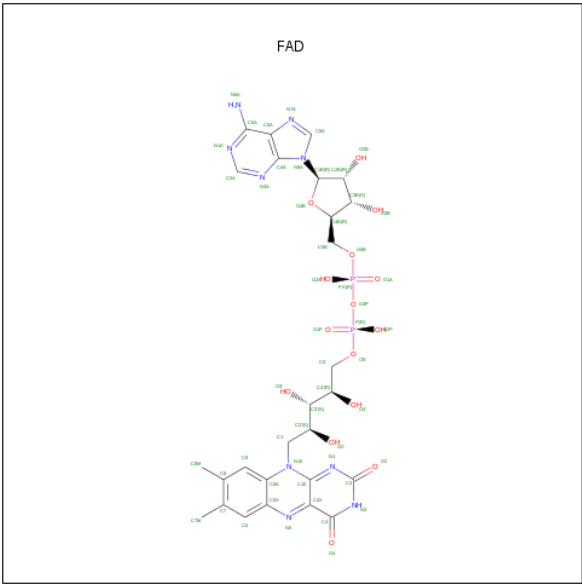
There are 5 unique types of molecules in this entry. The entry contains 14353 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Nitroalkane oxidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	430	Total 3315	2100	568	626	21	0	1	0
1	B	431	Total 3320	2103	569	627	21	0	1	0
1	C	430	Total 3307	2095	567	625	20	0	0	0
1	D	430	Total 3315	2100	568	626	21	0	1	0

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: C<sub>27</sub>H<sub>33</sub>N<sub>9</sub>O<sub>15</sub>P<sub>2</sub>).



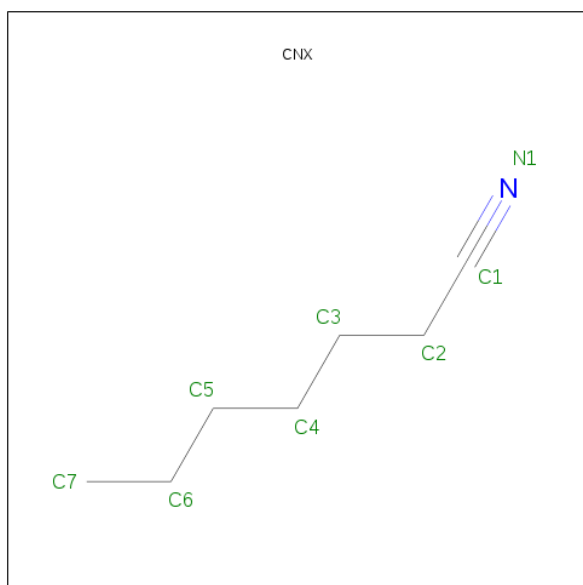
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	A	1	Total 53	27	9	15	2	0	0
2	B	1	Total 53	27	9	15	2	0	0

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

- Molecule 3 is HEPTANENITRILE (three-letter code: CNX) (formula:  $C_7H_{13}N$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	N	0	0
			8	7	1		
3	B	1	Total	C	N	0	0
			8	7	1		
3	C	1	Total	C	N	0	0
			8	7	1		
3	D	1	Total	C	N	0	0
			8	7	1		

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula:  $C_3H_8O_3$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		
4	D	1	Total	C	O	0	0
			6	3	3		

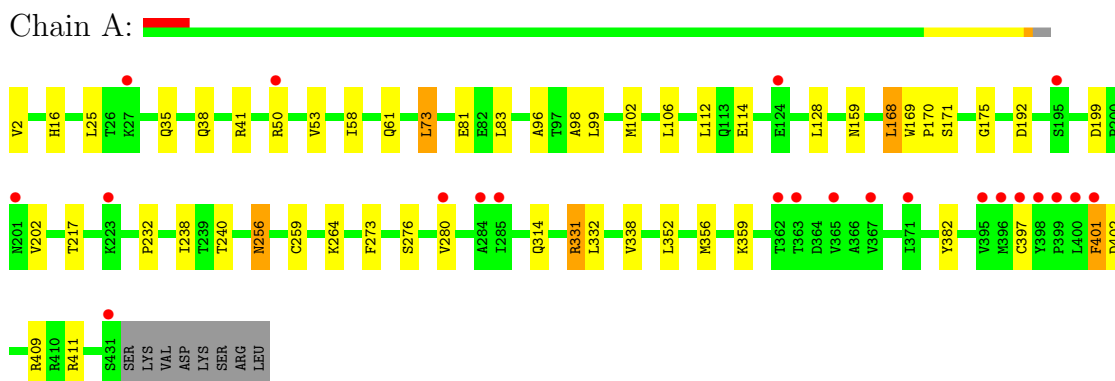
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	204	Total	O	0	0
			204	204		
5	B	194	Total	O	0	0
			194	194		
5	C	219	Total	O	0	0
			219	219		
5	D	223	Total	O	0	0
			223	223		

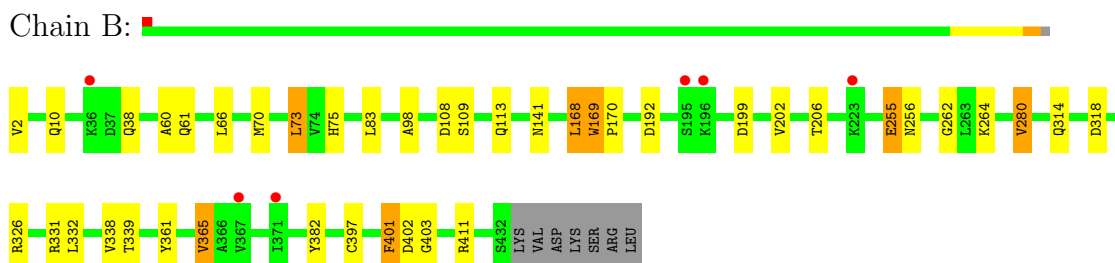
### 3 Residue-property plots

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.

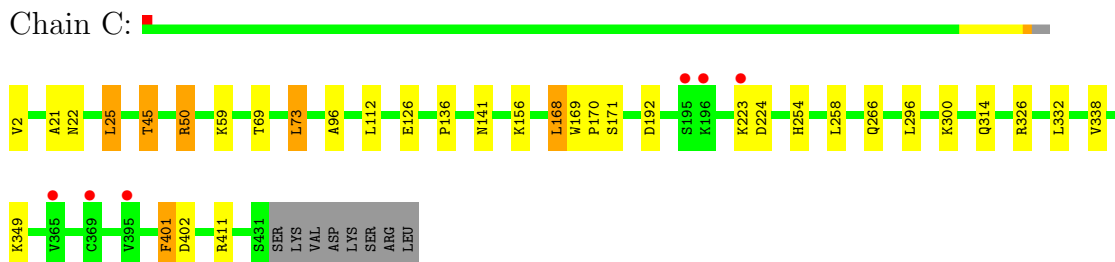
#### • Molecule 1: Nitroalkane oxidase



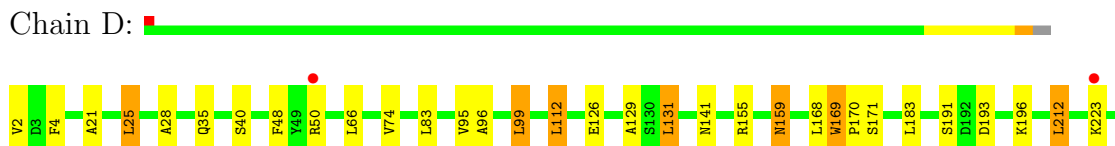
#### • Molecule 1: Nitroalkane oxidase

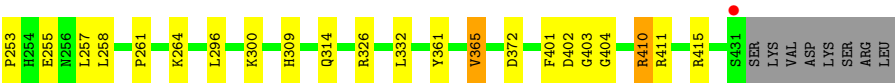


#### • Molecule 1: Nitroalkane oxidase



#### • Molecule 1: Nitroalkane oxidase





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	109.26Å 109.26Å 343.94Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.00 – 2.15 49.32 – 2.15	Depositor EDS
% Data completeness (in resolution range)	94.0 (50.00-2.15) 94.0 (49.32-2.15)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	0.10	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.98 (at 2.16Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.194 , 0.233 0.191 , 0.225	Depositor DCC
$R_{free}$ test set	6137 reflections (5.30%)	DCC
Wilson B-factor (Å <sup>2</sup> )	37.0	Xtriage
Anisotropy	0.061	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 27.6	EDS
Estimated twinning fraction	0.011 for -h,-k,l	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.35$	Xtriage
Outliers	0 of 121858 reflections	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	14353	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.80% 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: GOL, FAD, CNX

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	0.50	0/3384	0.61	1/4592 (0.0%)
1	B	0.50	0/3389	0.62	2/4599 (0.0%)
1	C	0.52	0/3376	0.59	0/4581
1	D	0.51	0/3384	0.63	2/4592 (0.0%)
All	All	0.51	0/13533	0.61	5/18364 (0.0%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	331	ARG	NE-CZ-NH2	-8.92	115.84	120.30
1	A	331	ARG	NE-CZ-NH2	-6.78	116.91	120.30
1	D	410	ARG	NE-CZ-NH1	6.02	123.31	120.30
1	D	410	ARG	NE-CZ-NH2	-5.85	117.37	120.30
1	B	331	ARG	NE-CZ-NH1	5.83	123.22	120.30

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	3315	0	3328	30	0
1	B	3320	0	3330	30	0
1	C	3307	0	3319	23	0
1	D	3315	0	3328	32	0
2	A	53	0	31	3	0
2	B	53	0	31	1	0
2	C	53	0	31	5	0
2	D	53	0	31	3	0
3	A	8	0	12	3	0
3	B	8	0	12	4	0
3	C	8	0	12	4	0
3	D	8	0	12	3	0
4	A	6	0	8	1	0
4	D	6	0	8	2	0
5	A	204	0	0	1	0
5	B	194	0	0	3	0
5	C	219	0	0	5	0
5	D	223	0	0	3	0
All	All	14353	0	13493	114	0

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 4.

All (114) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:159:ASN:HD22	1:D:159:ASN:H	1.21	0.83
1:C:141:ASN:HD21	2:C:500:FAD:H61A	1.24	0.82
1:C:22:ASN:HB3	5:C:697:HOH:O	1.78	0.81
1:B:199:ASP:HB3	1:B:202:VAL:HG12	1.64	0.79
1:B:199:ASP:HB3	1:B:202:VAL:CG1	2.16	0.76
1:C:45:THR:HG23	5:C:685:HOH:O	1.86	0.76
1:C:168:LEU:C	1:C:170:PRO:HD3	2.07	0.75
1:C:402:ASP:HA	3:C:501:CNX:N1	2.04	0.73
1:B:73:LEU:HB3	1:B:338:VAL:HB	1.72	0.72
1:A:168:LEU:C	1:A:170:PRO:HD3	2.11	0.71
1:C:326:ARG:NH2	5:C:606:HOH:O	2.03	0.71
1:A:73:LEU:HB3	1:A:338:VAL:HB	1.73	0.69
1:C:73:LEU:HB3	1:C:338:VAL:HB	1.73	0.69
1:D:50:ARG:HD2	1:D:126:GLU:OE1	1.92	0.69
1:B:168:LEU:C	1:B:170:PRO:HD3	2.14	0.67
1:A:38:GLN:HG3	1:A:232:PRO:O	1.95	0.66
1:D:326:ARG:NH2	1:D:372:ASP:OD2	2.30	0.64
1:B:2:VAL:N	4:D:701:GOL:HO3	1.96	0.64

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:95:VAL:HG22	3:D:501:CNX:H7A	1.79	0.63
1:C:22:ASN:CB	5:C:697:HOH:O	2.40	0.62
1:D:141:ASN:HD21	2:D:500:FAD:H61A	1.46	0.62
1:C:96:ALA:HB1	1:C:171:SER:HB2	1.79	0.62
1:D:253:PRO:HB2	1:D:255:GLU:HG2	1.81	0.62
1:B:326:ARG:HB3	1:B:365:VAL:HG13	1.82	0.61
1:D:35:GLN:HE21	1:D:40:SER:HB3	1.66	0.61
1:B:141:ASN:HD21	2:B:500:FAD:H61A	1.49	0.60
1:D:96:ALA:HB1	1:D:171:SER:HB2	1.83	0.60
1:C:45:THR:CG2	5:C:685:HOH:O	2.48	0.60
1:A:168:LEU:O	1:A:170:PRO:HD3	2.00	0.60
1:A:402:ASP:HA	3:A:501:CNX:N1	2.17	0.60
1:A:16:HIS:HD2	5:A:846:HOH:O	1.85	0.59
1:B:339:THR:HG21	1:D:4:PHE:HD1	1.67	0.58
1:B:61:GLN:HE21	1:B:98:ALA:HB2	1.68	0.58
1:D:402:ASP:HA	3:D:501:CNX:N1	2.19	0.58
1:B:255:GLU:HG3	5:B:682:HOH:O	2.03	0.57
1:B:314:GLN:HB2	1:C:314:GLN:HB2	1.88	0.56
2:C:500:FAD:O4	3:C:501:CNX:H2A	2.05	0.55
1:A:53:VAL:HG22	1:A:58:ILE:HG13	1.88	0.55
1:B:264:LYS:HE3	5:B:659:HOH:O	2.07	0.54
1:A:96:ALA:HB1	1:A:171:SER:HB2	1.90	0.54
1:A:276:SER:HB2	3:A:501:CNX:H7B	1.89	0.53
2:C:500:FAD:H6	3:C:501:CNX:H3	1.91	0.53
1:A:38:GLN:HE22	1:A:238:ILE:HA	1.73	0.52
1:A:382:TYR:CZ	1:B:397:CYS:HB2	2.45	0.52
1:D:326:ARG:HB3	1:D:365:VAL:HG13	1.91	0.52
1:C:141:ASN:ND2	2:C:500:FAD:H61A	2.02	0.51
1:A:259:CYS:HB2	1:A:264:LYS:HD3	1.92	0.51
1:A:314:GLN:HB2	1:D:314:GLN:HB2	1.93	0.51
1:D:99:LEU:HD22	1:D:131:LEU:HD23	1.92	0.51
1:B:280:VAL:HG23	3:B:501:CNX:H7A	1.94	0.50
1:B:402:ASP:HA	3:B:501:CNX:N1	2.27	0.50
1:A:199:ASP:HB3	1:A:202:VAL:HG12	1.94	0.50
1:A:256:ASN:N	1:A:256:ASN:HD22	2.08	0.50
1:C:50:ARG:HG3	1:C:126:GLU:OE1	2.12	0.49
1:C:21:ALA:HA	1:C:25:LEU:HB2	1.93	0.49
1:A:273:PHE:CD1	4:A:700:GOL:H12	2.47	0.49
1:B:168:LEU:O	1:B:170:PRO:HD3	2.13	0.49
1:D:296:LEU:O	1:D:300:LYS:HG3	2.12	0.49
1:A:61:GLN:HE21	1:A:98:ALA:HB2	1.79	0.48
1:C:96:ALA:HB1	1:C:171:SER:CB	2.43	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:401:PHE:CD2	1:A:401:PHE:C	2.87	0.47
1:B:206:THR:O	1:B:262:GLY:HA2	2.13	0.47
1:C:168:LEU:O	1:C:170:PRO:HD3	2.13	0.47
1:A:102[B]:MET:HE3	1:A:106:LEU:HG	1.95	0.47
1:A:199:ASP:HB3	1:A:202:VAL:CG1	2.45	0.47
1:D:21:ALA:HA	1:D:25:LEU:HB2	1.96	0.47
1:D:74:VAL:HB	4:D:701:GOL:H2	1.96	0.47
1:C:296:LEU:O	1:C:300:LYS:HG3	2.15	0.46
1:C:401:PHE:CD2	1:C:401:PHE:C	2.86	0.46
1:D:112:LEU:HD23	1:D:258:LEU:HD12	1.97	0.46
1:D:193:ASP:OD1	1:D:196:LYS:HE2	2.16	0.46
2:A:500:FAD:O4	3:A:501:CNX:H2A	2.16	0.46
1:C:112:LEU:HD23	1:C:258:LEU:HD12	1.97	0.45
1:D:257:LEU:HD21	1:D:261:PRO:HD3	1.98	0.45
1:B:168:LEU:O	1:B:169:TRP:HB2	2.16	0.45
1:C:402:ASP:CA	3:C:501:CNX:N1	2.77	0.45
1:D:159:ASN:H	1:D:159:ASN:ND2	2.01	0.45
1:A:35:GLN:HB2	1:A:41:ARG:HG2	1.99	0.45
1:D:309:HIS:HE1	5:D:909:HOH:O	1.98	0.45
1:A:53:VAL:HG22	1:A:58:ILE:CG1	2.46	0.44
1:A:81:GLU:OE2	1:A:331:ARG:HD3	2.18	0.44
1:D:155:ARG:HB3	1:D:191:SER:O	2.17	0.44
1:A:170:PRO:HA	2:A:500:FAD:C4	2.47	0.44
1:D:169:TRP:N	1:D:170:PRO:HD2	2.33	0.44
1:A:397:CYS:HB2	1:B:382:TYR:CZ	2.52	0.44
1:B:361:TYR:O	1:B:365:VAL:HB	2.18	0.44
1:D:159:ASN:N	1:D:159:ASN:HD22	1.98	0.44
1:D:361:TYR:O	1:D:365:VAL:HB	2.17	0.44
1:B:402:ASP:N	3:B:501:CNX:N1	2.66	0.43
1:A:352:LEU:O	1:A:356:MET:HG2	2.19	0.43
1:A:128:LEU:HD12	1:A:175:GLY:HA2	2.00	0.43
1:D:28:ALA:HA	1:D:48:PHE:CZ	2.54	0.43
2:D:500:FAD:H52A	2:D:500:FAD:O2P	2.18	0.43
1:D:169:TRP:N	1:D:170:PRO:CD	2.82	0.43
1:B:401:PHE:CD2	1:B:401:PHE:C	2.92	0.43
1:B:10:GLN:HG3	1:B:75:HIS:CE1	2.54	0.42
1:D:404:GLY:HA3	5:D:716:HOH:O	2.18	0.42
2:C:500:FAD:H52A	2:C:500:FAD:O2P	2.20	0.42
1:C:136:PRO:HA	1:C:168:LEU:HD12	2.02	0.42
1:D:415:ARG:NH2	5:D:810:HOH:O	2.36	0.42
1:B:109:SER:O	1:B:113:GLN:HG3	2.20	0.41
1:C:59:LYS:HE2	1:C:69:THR:HG23	2.02	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:183:LEU:HD11	1:D:212:LEU:HG	2.01	0.41
1:A:359:LYS:HD2	1:A:409:ARG:HG3	2.02	0.41
1:A:382:TYR:OH	1:B:397:CYS:HB2	2.21	0.41
1:B:108:ASP:HB3	5:B:689:HOH:O	2.20	0.41
2:D:500:FAD:O4	3:D:501:CNX:H2A	2.21	0.41
1:B:318:ASP:HA	1:D:410:ARG:HH21	1.84	0.41
1:B:402:ASP:CA	3:B:501:CNX:N1	2.84	0.41
1:A:240:THR:HG21	2:A:500:FAD:HM72	2.03	0.41
1:C:156:LYS:HE3	1:C:254:HIS:ND1	2.36	0.41
1:B:38:GLN:HE21	1:B:38:GLN:HB2	1.70	0.40
1:B:60:ALA:O	1:B:70:MET:HB2	2.21	0.40
1:D:129:ALA:HA	1:D:183:LEU:O	2.22	0.40

There are no symmetry-related clashes.

## 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	429/438 (98%)	423 (99%)	5 (1%)	1 (0%)	56	55
1	B	430/438 (98%)	422 (98%)	6 (1%)	2 (0%)	38	31
1	C	428/438 (98%)	420 (98%)	7 (2%)	1 (0%)	56	55
1	D	429/438 (98%)	419 (98%)	8 (2%)	2 (0%)	38	31
All	All	1716/1752 (98%)	1684 (98%)	26 (2%)	6 (0%)	43	47

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	169	TRP
1	B	169	TRP
1	C	169	TRP
1	D	169	TRP
1	D	403	GLY
1	B	403	GLY

### 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	356/363 (98%)	339 (95%)	17 (5%)	35	32
1	B	356/363 (98%)	344 (97%)	12 (3%)	49	50
1	C	355/363 (98%)	341 (96%)	14 (4%)	43	43
1	D	356/363 (98%)	340 (96%)	16 (4%)	38	35
All	All	1423/1452 (98%)	1364 (96%)	59 (4%)	40	40

All (59) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	VAL
1	A	25	LEU
1	A	50	ARG
1	A	73	LEU
1	A	83	LEU
1	A	99	LEU
1	A	112	LEU
1	A	114	GLU
1	A	159	ASN
1	A	168	LEU
1	A	192	ASP
1	A	217	THR
1	A	256	ASN
1	A	280	VAL
1	A	332	LEU
1	A	401	PHE
1	A	411	ARG
1	B	66	LEU
1	B	73	LEU
1	B	83	LEU
1	B	168	LEU
1	B	192	ASP
1	B	255	GLU
1	B	256	ASN
1	B	280	VAL
1	B	332	LEU

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Mol	Chain	Res	Type
1	B	365	VAL
1	B	401	PHE
1	B	411	ARG
1	C	2	VAL
1	C	25	LEU
1	C	45	THR
1	C	50	ARG
1	C	73	LEU
1	C	168	LEU
1	C	192	ASP
1	C	223	LYS
1	C	224	ASP
1	C	266	GLN
1	C	332	LEU
1	C	349	LYS
1	C	401	PHE
1	C	411	ARG
1	D	2	VAL
1	D	25	LEU
1	D	66	LEU
1	D	83	LEU
1	D	99	LEU
1	D	112	LEU
1	D	131	LEU
1	D	159	ASN
1	D	168	LEU
1	D	212	LEU
1	D	223	LYS
1	D	264	LYS
1	D	332	LEU
1	D	365	VAL
1	D	401	PHE
1	D	411	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (40) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	22	ASN
1	A	38	GLN
1	A	43	GLN
1	A	61	GLN
1	A	113	GLN

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Mol	Chain	Res	Type
1	A	137	ASN
1	A	159	ASN
1	A	198	GLN
1	A	256	ASN
1	B	22	ASN
1	B	35	GLN
1	B	38	GLN
1	B	113	GLN
1	B	137	ASN
1	B	141	ASN
1	B	198	GLN
1	B	237	HIS
1	B	256	ASN
1	B	266	GLN
1	B	357	GLN
1	C	22	ASN
1	C	43	GLN
1	C	113	GLN
1	C	137	ASN
1	C	141	ASN
1	C	144	GLN
1	C	198	GLN
1	C	357	GLN
1	D	16	HIS
1	D	22	ASN
1	D	35	GLN
1	D	43	GLN
1	D	113	GLN
1	D	137	ASN
1	D	141	ASN
1	D	144	GLN
1	D	159	ASN
1	D	198	GLN
1	D	256	ASN
1	D	266	GLN

### 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 ⓘ

10 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	3	58,58,58	1.35	4 (6%)	85,89,89	1.92	13 (15%)
3	CNX	A	501	2	7,7,7	0.35	0	6,6,6	0.98	0
4	GOL	A	700	-	5,5,5	0.41	0	5,5,5	0.21	0
2	FAD	B	500	3	58,58,58	1.42	6 (10%)	85,89,89	1.88	11 (12%)
3	CNX	B	501	2	7,7,7	0.45	0	6,6,6	1.28	1 (16%)
2	FAD	C	500	3	58,58,58	1.35	6 (10%)	85,89,89	1.79	14 (16%)
3	CNX	C	501	2	7,7,7	0.50	0	6,6,6	1.30	1 (16%)
2	FAD	D	500	3	58,58,58	1.38	5 (8%)	85,89,89	1.90	12 (14%)
3	CNX	D	501	2	7,7,7	0.73	0	6,6,6	1.51	2 (33%)
4	GOL	D	701	-	5,5,5	0.31	0	5,5,5	0.31	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
2	FAD	A	500	3	-	0/34/50/50	0/6/6/6
3	CNX	A	501	2	-	0/5/5/5	0/0/0/0
4	GOL	A	700	-	-	0/4/4/4	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	FAD	B	500	3	-	0/34/50/50	0/6/6/6
3	CNX	B	501	2	-	0/5/5/5	0/0/0/0
2	FAD	C	500	3	-	0/34/50/50	0/6/6/6
3	CNX	C	501	2	-	0/5/5/5	0/0/0/0
2	FAD	D	500	3	-	0/34/50/50	0/6/6/6
3	CNX	D	501	2	-	0/5/5/5	0/0/0/0
4	GOL	D	701	-	-	0/4/4/4	0/0/0/0

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	500	FAD	C4X-N5	6.39	1.44	1.33
2	B	500	FAD	C4X-N5	6.17	1.43	1.33
2	A	500	FAD	C4X-N5	5.88	1.43	1.33
2	C	500	FAD	C4X-N5	5.84	1.43	1.33
2	B	500	FAD	C5X-N5	4.82	1.43	1.35
2	D	500	FAD	C5X-N5	4.63	1.42	1.35
2	C	500	FAD	C5X-N5	4.54	1.42	1.35
2	A	500	FAD	C5X-N5	4.50	1.42	1.35
2	B	500	FAD	O4B-C1B	3.73	1.45	1.41
2	A	500	FAD	O2-C2	3.68	1.26	1.21
2	D	500	FAD	O2-C2	3.02	1.25	1.21
2	B	500	FAD	O2-C2	2.88	1.25	1.21
2	C	500	FAD	C9A-N10	2.86	1.43	1.38
2	D	500	FAD	O4B-C1B	2.85	1.44	1.41
2	A	500	FAD	C9A-N10	2.71	1.42	1.38
2	C	500	FAD	C4A-N9A	-2.54	1.34	1.37
2	D	500	FAD	C9A-N10	2.27	1.42	1.38
2	B	500	FAD	C4-N3	-2.21	1.33	1.36
2	C	500	FAD	O4B-C1B	2.15	1.43	1.41
2	C	500	FAD	O2-C2	2.14	1.24	1.21
2	B	500	FAD	C9A-N10	2.08	1.41	1.38

All (54) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	500	FAD	N3A-C2A-N1A	-9.92	120.17	128.89
2	A	500	FAD	N3A-C2A-N1A	-9.54	120.50	128.89
2	D	500	FAD	N3A-C2A-N1A	-9.44	120.58	128.89
2	C	500	FAD	N3A-C2A-N1A	-8.79	121.15	128.89
2	D	500	FAD	C5A-C4A-N3A	-7.12	119.04	125.98
2	B	500	FAD	C5A-C4A-N3A	-6.99	119.17	125.98

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	500	FAD	C2-N1-C10	6.70	121.40	114.95
2	A	500	FAD	C5A-C4A-N3A	-6.68	119.47	125.98
2	D	500	FAD	C2-N1-C10	6.04	120.77	114.95
2	B	500	FAD	C2-N1-C10	5.58	120.32	114.95
2	C	500	FAD	C5A-C4A-N3A	-5.57	120.55	125.98
2	C	500	FAD	C2-N1-C10	5.04	119.81	114.95
2	C	500	FAD	C9A-N10-C10	-4.80	117.03	121.77
2	D	500	FAD	C9A-N10-C10	-4.74	117.10	121.77
2	D	500	FAD	N3A-C4A-N9A	4.68	133.42	125.39
2	B	500	FAD	C9A-N10-C10	-4.51	117.32	121.77
2	B	500	FAD	N3A-C4A-N9A	4.42	132.97	125.39
2	A	500	FAD	N3A-C4A-N9A	4.38	132.90	125.39
2	A	500	FAD	C9A-N10-C10	-4.17	117.65	121.77
2	C	500	FAD	N3A-C4A-N9A	4.06	132.36	125.39
2	D	500	FAD	C2'-C1'-N10	-3.11	108.34	112.60
3	C	501	CNX	C2-C1-N1	-3.04	161.35	176.69
3	D	501	CNX	C2-C1-N1	-2.93	161.93	176.69
2	D	500	FAD	C4X-N5-C5X	-2.85	113.36	116.68
2	B	500	FAD	C4X-N5-C5X	-2.80	113.41	116.68
2	A	500	FAD	C2A-N3A-C4A	2.78	121.27	113.27
2	B	500	FAD	C2A-N3A-C4A	2.77	121.25	113.27
2	D	500	FAD	C2A-N3A-C4A	2.76	121.23	113.27
2	C	500	FAD	C4X-C10-N1	-2.67	119.23	123.00
2	C	500	FAD	C4X-C10-N10	-2.66	118.92	120.53
2	B	500	FAD	C2'-C1'-N10	-2.66	108.96	112.60
2	A	500	FAD	C4A-C5A-N7A	-2.63	106.87	109.41
2	A	500	FAD	C4X-C10-N1	-2.62	119.30	123.00
3	B	501	CNX	C2-C1-N1	-2.56	163.79	176.69
2	A	500	FAD	C4X-C10-N10	-2.52	119.00	120.53
2	C	500	FAD	C1'-N10-C10	2.43	121.47	118.82
2	A	500	FAD	C3B-C2B-C1B	2.41	104.70	100.92
2	B	500	FAD	C1'-N10-C9A	2.38	121.17	118.67
2	C	500	FAD	C4-N3-C2	-2.36	120.54	125.39
2	C	500	FAD	C4X-N5-C5X	-2.32	113.98	116.68
2	A	500	FAD	C4-N3-C2	-2.31	120.65	125.39
2	D	500	FAD	C4X-C10-N10	-2.28	119.15	120.53
2	C	500	FAD	C1'-N10-C9A	2.25	121.03	118.67
2	D	500	FAD	C4X-C10-N1	-2.20	119.89	123.00
2	A	500	FAD	C1'-N10-C9A	2.19	120.97	118.67
2	B	500	FAD	C4A-C5A-N7A	-2.16	107.32	109.41
2	D	500	FAD	C4-N3-C2	-2.16	120.97	125.39
2	C	500	FAD	C2A-N3A-C4A	2.14	119.43	113.27

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	500	FAD	C2'-C1'-N10	-2.10	109.72	112.60
2	B	500	FAD	C4X-C10-N1	-2.09	120.06	123.00
2	D	500	FAD	C3B-C2B-C1B	2.07	104.17	100.92
2	C	500	FAD	C5X-C9A-N10	2.06	119.15	117.63
3	D	501	CNX	C4-C3-C2	-2.05	105.18	113.33
2	A	500	FAD	C8A-N9A-C4A	2.03	108.61	106.96

There are no chirality outliers.

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	430/438 (98%)	0.22	22 (5%) 27 30	27, 40, 52, 61	0
1	B	431/438 (98%)	0.08	6 (1%) 72 77	26, 37, 50, 58	0
1	C	430/438 (98%)	-0.11	6 (1%) 72 77	27, 35, 48, 59	0
1	D	430/438 (98%)	0.06	3 (0%) 84 89	26, 37, 51, 60	1 (0%)
All	All	1721/1752 (98%)	0.06	37 (2%) 60 65	26, 37, 50, 61	1 (0%)

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	367	VAL	3.8
1	A	395	VAL	3.7
1	A	363	THR	3.1
1	A	400	LEU	3.1
1	A	223	LYS	3.1
1	A	284	ALA	3.0
1	A	365	VAL	3.0
1	B	371	ILE	2.9
1	C	223	LYS	2.9
1	B	196	LYS	2.9
1	A	362	THR	2.9
1	B	36	LYS	2.9
1	A	396	MET	2.8
1	A	397	CYS	2.8
1	B	195	SER	2.7
1	C	365	VAL	2.7
1	C	195	SER	2.6
1	A	431	SER	2.6
1	A	50	ARG	2.5
1	A	124	GLU	2.4
1	A	280	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	367	VAL	2.3
1	D	223	LYS	2.3
1	A	398	TYR	2.2
1	A	401	PHE	2.2
1	B	223	LYS	2.2
1	A	27	LYS	2.2
1	A	285	ILE	2.2
1	A	201	ASN	2.1
1	C	196	LYS	2.1
1	A	371	ILE	2.1
1	A	399	PRO	2.1
1	D	431	SER	2.1
1	C	369	CYS	2.0
1	A	195	SER	2.0
1	C	395	VAL	2.0
1	D	50	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 ⓘ

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
4	GOL	D	701	6/6	0.23	4.11	55,56,57,58	0
3	CNX	B	501	8/8	0.15	1.15	36,40,42,44	0
3	CNX	A	501	8/8	0.19	0.77	39,41,43,44	0
3	CNX	C	501	8/8	0.11	0.19	35,39,42,43	0
3	CNX	D	501	8/8	0.13	0.07	37,39,40,41	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	FAD	D	500	53/53	0.10	-0.40	24,31,34,36	0
2	FAD	A	500	53/53	0.11	-0.46	29,33,36,37	0
2	FAD	C	500	53/53	0.09	-0.60	24,27,32,32	0
2	FAD	B	500	53/53	0.09	-0.62	26,30,32,32	0
4	GOL	A	700	6/6	0.10	-0.90	41,42,43,43	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.