



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 17, 2018 – 09:12 am GMT

PDB ID : 1N4W
Title : ATOMIC RESOLUTION STRUCTURE OF CHOLESTEROL OXIDASE @
pH 7.3 (STREPTOMYCES SP. SA-COO)
Authors : Vrielink, A.; Lario, P.I.
Deposited on : 2002-11-01
Resolution : 0.92 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/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 : 1.7.3 (157068), CSD as539be (2018)
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk30686

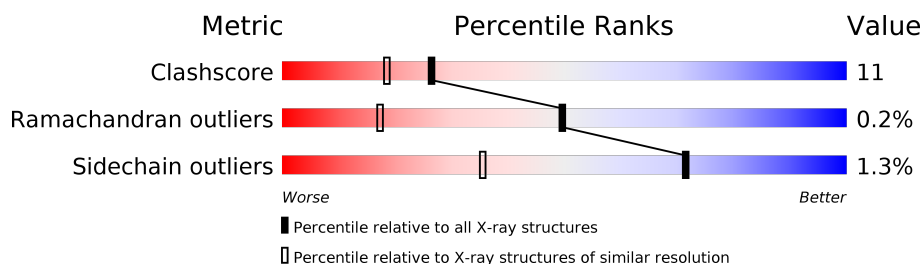
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 0.92 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))
Clashscore	122078	1141 (1.06-0.78)
Ramachandran outliers	120005	1058 (1.06-0.78)
Sidechain outliers	119972	1060 (1.06-0.78)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	504	

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
4	GOL	A	511	-	X	X	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 9520 atoms, of which 4334 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.

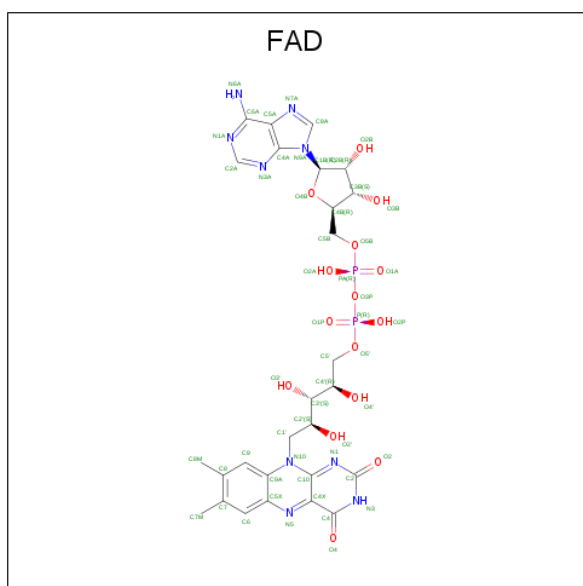
- Molecule 1 is a protein called Cholesterol oxidase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	499	Total	C	H	N	O	S	0	92	1
			8616	2747	4303	725	816	25			

- Molecule 2 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mn	0	0
			1	1		

- Molecule 3 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total	C	H	N	O	P	0	0
			84	27	31	9	15	2		

- 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		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	813	Total	O	0	0
			813	813		

4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	51.34Å 72.96Å 63.05Å 90.00° 105.25° 90.00°	Depositor
Resolution (Å)	31.30 – 0.92	Depositor
% Data completeness (in resolution range)	95.0 (31.30-0.92)	Depositor
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	SHELXL-97	Depositor
R, R_{free}	0.103 , 0.122	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	9520	wwPDB-VP
Average B, all atoms (Å ²)	13.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, MN, 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	0.83	2/4646 (0.0%)	1.29	47/6308 (0.7%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	400[A]	ASN	C-N	-5.18	1.22	1.34
1	A	400[B]	ASN	C-N	-5.18	1.22	1.34

All (47) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	396	ARG	NE-CZ-NH2	-17.37	111.62	120.30
1	A	254	ARG	NE-CZ-NH2	-13.58	113.51	120.30
1	A	283[A]	ARG	NE-CZ-NH1	-12.95	113.82	120.30
1	A	283[B]	ARG	NE-CZ-NH1	-12.95	113.82	120.30
1	A	71	ARG	NE-CZ-NH2	-12.59	114.01	120.30
1	A	254	ARG	NH1-CZ-NH2	10.71	131.18	119.40
1	A	254	ARG	NE-CZ-NH1	-9.99	115.31	120.30
1	A	202[A]	ARG	NE-CZ-NH2	9.73	125.16	120.30
1	A	202[B]	ARG	NE-CZ-NH2	9.73	125.16	120.30
1	A	328	ARG	NE-CZ-NH2	-7.64	116.48	120.30
1	A	283[A]	ARG	NE-CZ-NH2	7.60	124.10	120.30
1	A	283[B]	ARG	NE-CZ-NH2	7.60	124.10	120.30
1	A	328	ARG	NE-CZ-NH1	7.29	123.94	120.30
1	A	202[A]	ARG	NE-CZ-NH1	-6.99	116.81	120.30
1	A	202[B]	ARG	NE-CZ-NH1	-6.99	116.81	120.30
1	A	71	ARG	NE-CZ-NH1	6.87	123.74	120.30
1	A	419	ARG	NE-CZ-NH2	6.75	123.67	120.30
1	A	128	ARG	NE-CZ-NH2	-6.65	116.98	120.30
1	A	122[A]	MET	CG-SD-CE	6.57	110.72	100.20

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	122[B]	MET	CG-SD-CE	6.57	110.72	100.20
1	A	332[A]	MET	CG-SD-CE	-6.40	89.96	100.20
1	A	332[B]	MET	CG-SD-CE	-6.40	89.96	100.20
1	A	332[C]	MET	CG-SD-CE	-6.40	89.96	100.20
1	A	446	TYR	CB-CG-CD1	6.32	124.79	121.00
1	A	59[A]	LEU	CB-CG-CD1	6.21	121.55	111.00
1	A	59[B]	LEU	CB-CG-CD1	6.21	121.55	111.00
1	A	463	ARG	NE-CZ-NH2	-6.13	117.23	120.30
1	A	400[A]	ASN	O-C-N	-6.12	112.90	122.70
1	A	400[B]	ASN	O-C-N	-6.12	112.90	122.70
1	A	396	ARG	NH1-CZ-NH2	6.00	126.00	119.40
1	A	400[A]	ASN	C-N-CA	5.88	136.39	121.70
1	A	400[B]	ASN	C-N-CA	5.88	136.39	121.70
1	A	390	TYR	CB-CG-CD2	-5.59	117.64	121.00
1	A	87	ARG	NE-CZ-NH1	5.57	123.09	120.30
1	A	178	ARG	NE-CZ-NH2	5.46	123.03	120.30
1	A	396	ARG	O-C-N	5.45	131.41	122.70
1	A	361[A]	GLU	CA-CB-CG	5.43	125.35	113.40
1	A	361[B]	GLU	CA-CB-CG	5.43	125.35	113.40
1	A	279[A]	GLU	CG-CD-OE2	5.30	128.90	118.30
1	A	279[B]	GLU	CG-CD-OE2	5.30	128.90	118.30
1	A	219	TYR	CB-CG-CD2	-5.28	117.83	121.00
1	A	367	ALA	C-N-CA	-5.25	111.27	122.30
1	A	388	PHE	CB-CG-CD2	-5.25	117.13	120.80
1	A	390	TYR	CB-CG-CD1	5.23	124.14	121.00
1	A	433[A]	PHE	CB-CG-CD2	5.19	124.43	120.80
1	A	433[B]	PHE	CB-CG-CD2	5.19	124.43	120.80
1	A	64	ARG	NE-CZ-NH1	5.00	122.80	120.30

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	4313	4303	4223	90	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	1	0	0	0	0
3	A	53	31	31	1	0
4	A	6	0	6	7	0
5	A	813	0	0	74	0
All	All	5186	4334	4260	97	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 11.

All (97) 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)
4:A:511:GOL:C2	4:A:511:GOL:C3	1.89	1.48
1:A:243[B]:THR:HG21	5:A:948:HOH:O	1.24	1.35
1:A:365[B]:MET:SD	5:A:678:HOH:O	1.91	1.27
1:A:202[A]:ARG:NE	5:A:1002:HOH:O	1.62	1.26
1:A:359[A]:PHE:CZ	5:A:1325:HOH:O	1.86	1.23
1:A:258[B]:ASP:HA	5:A:1072:HOH:O	1.38	1.21
1:A:400[B]:ASN:ND2	5:A:1163:HOH:O	1.74	1.17
1:A:257[B]:LYS:O	5:A:1072:HOH:O	1.61	1.16
1:A:142[A]:GLU:OE2	5:A:1074:HOH:O	1.67	1.13
1:A:166[A]:GLU:OE2	5:A:1073:HOH:O	1.66	1.12
4:A:511:GOL:C3	4:A:511:GOL:O3	0.81	1.11
4:A:511:GOL:H31	4:A:511:GOL:O3	1.46	1.10
1:A:202[A]:ARG:CZ	5:A:1002:HOH:O	1.88	1.09
4:A:511:GOL:H32	4:A:511:GOL:O3	1.46	1.09
1:A:48[B]:PRO:HD2	5:A:1306:HOH:O	1.55	1.07
1:A:359[A]:PHE:CE2	5:A:1330:HOH:O	2.09	1.02
1:A:150[B]:ARG:NH1	5:A:986:HOH:O	1.94	0.99
1:A:255[B]:GLN:OE1	5:A:1155:HOH:O	1.81	0.98
1:A:359[A]:PHE:CD2	5:A:1324:HOH:O	2.16	0.97
1:A:142[A]:GLU:CG	5:A:1074:HOH:O	2.14	0.95
1:A:207[B]:GLU:OE2	5:A:567:HOH:O	1.85	0.95
1:A:207[B]:GLU:CD	5:A:567:HOH:O	2.05	0.95
1:A:359[A]:PHE:HD2	5:A:1324:HOH:O	1.48	0.93
1:A:166[A]:GLU:CG	5:A:1073:HOH:O	2.17	0.92
1:A:207[B]:GLU:OE1	5:A:567:HOH:O	1.87	0.92
1:A:215[B]:THR:HG21	5:A:1037:HOH:O	1.68	0.92
1:A:359[A]:PHE:CE2	5:A:1325:HOH:O	2.15	0.91
1:A:479[A]:PRO:HG2	5:A:1044:HOH:O	1.68	0.91
1:A:166[A]:GLU:HG2	5:A:1073:HOH:O	1.72	0.89

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:359[A]:PHE:HZ	5:A:1325:HOH:O	1.27	0.89
1:A:479[A]:PRO:O	5:A:1044:HOH:O	1.91	0.88
1:A:332[C]:MET:SD	5:A:698:HOH:O	2.31	0.88
1:A:257[B]:LYS:HB3	5:A:1059:HOH:O	1.74	0.86
1:A:504[B]:GLN:NE2	5:A:877:HOH:O	1.60	0.85
1:A:402[B]:THR:OG1	1:A:404:ASP:OD1	1.92	0.85
1:A:202[A]:ARG:NH2	5:A:1002:HOH:O	1.98	0.83
4:A:511:GOL:C1	4:A:511:GOL:C3	2.56	0.82
1:A:279[A]:GLU:HG2	5:A:1108:HOH:O	1.82	0.80
4:A:511:GOL:C2	4:A:511:GOL:O3	2.31	0.79
1:A:187[A]:GLY:O	5:A:1084:HOH:O	2.02	0.77
1:A:359[A]:PHE:HE2	5:A:1330:HOH:O	1.56	0.77
1:A:259[B]:GLY:N	5:A:1156:HOH:O	2.18	0.75
1:A:215[B]:THR:CG2	5:A:1037:HOH:O	2.30	0.74
1:A:361[B]:GLU:OE2	5:A:1328:HOH:O	2.06	0.72
1:A:485[A]:ASN:OD1	5:A:1330:HOH:O	2.06	0.72
1:A:283[A]:ARG:NE	5:A:922:HOH:O	2.01	0.69
1:A:355:ASP:OD1	5:A:1148:HOH:O	2.11	0.69
1:A:257[B]:LYS:CB	5:A:1059:HOH:O	2.37	0.67
1:A:332[C]:MET:CE	5:A:1240:HOH:O	2.43	0.66
1:A:218[A]:ILE:CG2	5:A:1118:HOH:O	2.44	0.66
1:A:296[A]:GLU:OE1	5:A:1158:HOH:O	2.10	0.64
1:A:332[C]:MET:HE2	5:A:1240:HOH:O	1.97	0.64
4:A:511:GOL:C3	4:A:511:GOL:O2	2.43	0.63
1:A:218[A]:ILE:HG23	5:A:1118:HOH:O	1.99	0.63
1:A:187[A]:GLY:C	5:A:1084:HOH:O	2.38	0.62
1:A:504[B]:GLN:NE2	5:A:788:HOH:O	2.32	0.62
1:A:59[A]:LEU:HD21	5:A:1307:HOH:O	1.99	0.61
1:A:507[B]:THR:N	5:A:1167:HOH:O	2.33	0.60
1:A:283[A]:ARG:CZ	5:A:922:HOH:O	2.46	0.59
1:A:283[A]:ARG:NH2	5:A:922:HOH:O	2.36	0.59
1:A:241:LYS:NZ	5:A:1136:HOH:O	2.37	0.58
1:A:284:TYR:CE1	1:A:506[B]:VAL:HG13	2.40	0.57
1:A:59[A]:LEU:CD2	5:A:1307:HOH:O	2.53	0.57
1:A:398:LYS:HE3	5:A:1035:HOH:O	2.08	0.54
1:A:479[A]:PRO:C	5:A:1044:HOH:O	2.42	0.53
1:A:284:TYR:CZ	1:A:506[B]:VAL:HG13	2.43	0.53
1:A:142[A]:GLU:HG3	5:A:1074:HOH:O	1.96	0.52
1:A:273:LYS:CE	5:A:1061:HOH:O	2.57	0.52
1:A:506[B]:VAL:HG12	1:A:507[B]:THR:N	2.24	0.52
1:A:273:LYS:HD3	5:A:1061:HOH:O	2.11	0.50

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:377[A]:LEU:HD12	5:A:1329:HOH:O	2.12	0.50
1:A:218[A]:ILE:HG22	5:A:545:HOH:O	2.12	0.49
1:A:243[B]:THR:CG2	5:A:948:HOH:O	2.09	0.49
1:A:162:THR:O	1:A:166[A]:GLU:HG2	2.13	0.48
1:A:485[A]:ASN:HB3	3:A:510:FAD:C2	2.43	0.48
1:A:456[B]:LYS:NZ	5:A:883:HOH:O	2.47	0.46
1:A:264[B]:THR:HG21	5:A:663:HOH:O	2.16	0.46
1:A:9[B]:GLY:N	5:A:1195:HOH:O	2.48	0.46
1:A:283[A]:ARG:NH2	5:A:1120:HOH:O	2.50	0.45
1:A:257[B]:LYS:CA	5:A:1059:HOH:O	2.63	0.44
1:A:377[A]:LEU:CD1	5:A:1327:HOH:O	2.66	0.44
1:A:164:TRP:CD1	1:A:201[B]:GLN:HG3	2.53	0.44
1:A:332[C]:MET:HE1	5:A:1240:HOH:O	2.12	0.43
1:A:66[A]:SER:OG	1:A:69[A]:LYS:HB2	2.19	0.43
1:A:389[A]:VAL:HG13	5:A:1164:HOH:O	2.17	0.43
1:A:368:GLY:HA3	5:A:1307:HOH:O	2.18	0.43
1:A:201[B]:GLN:HG2	5:A:1104:HOH:O	2.19	0.43
1:A:170:TRP:CZ3	1:A:339[A]:HIS:HB3	2.53	0.42
1:A:395:ASP:O	1:A:396:ARG:HD2	2.19	0.42
1:A:299[B]:VAL:HG12	1:A:390:TYR:HB2	2.01	0.42
1:A:377[A]:LEU:HD12	5:A:1327:HOH:O	2.18	0.42
1:A:188[A]:THR:HA	1:A:347:GLY:O	2.20	0.42
1:A:34:VAL:HG21	1:A:506[B]:VAL:HG11	2.02	0.42
1:A:258[B]:ASP:CA	5:A:1072:HOH:O	2.23	0.41
1:A:255[B]:GLN:HG2	1:A:256[B]:THR:O	2.21	0.41
1:A:351:TRP:CD2	1:A:357:SER:HB3	2.56	0.41

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	589/504 (117%)	568 (96%)	20 (3%)	1 (0%)	49 17

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	217	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	479/401 (120%)	471 (98%)	8 (2%)	63 27

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	56[A]	CYS
1	A	56[B]	CYS
1	A	59[A]	LEU
1	A	59[B]	LEU
1	A	158	ASN
1	A	279[A]	GLU
1	A	279[B]	GLU
1	A	369	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 1 is monoatomic - leaving 2 for Mogul analysis.

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
3	FAD	A	510	-	50,58,58	1.28	5 (10%)	56,89,89	1.06	2 (3%)
4	GOL	A	511	-	5,5,5	8.25	5 (100%)	5,5,5	3.20	3 (60%)

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
3	FAD	A	510	-	-	0/28/50/50	0/6/6/6
4	GOL	A	511	-	-	0/4/4/4	0/0/0/0

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	511	GOL	O3-C3	-14.37	0.81	1.42
4	A	511	GOL	O2-C2	-3.00	1.34	1.43
3	A	510	FAD	C5A-C4A	-2.60	1.34	1.40
3	A	510	FAD	C10-N1	2.13	1.36	1.33
3	A	510	FAD	C5X-N5	2.44	1.39	1.35
4	A	511	GOL	O1-C1	2.51	1.53	1.42
3	A	510	FAD	C9A-N10	2.92	1.42	1.38
4	A	511	GOL	C1-C2	4.94	1.71	1.52
3	A	510	FAD	C4X-N5	5.16	1.40	1.33
4	A	511	GOL	C3-C2	9.71	1.89	1.52

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	511	GOL	C3-C2-C1	-5.39	90.87	111.63
4	A	511	GOL	O1-C1-C2	-3.78	91.79	110.11
4	A	511	GOL	O2-C2-C3	-2.75	96.41	109.00
3	A	510	FAD	C2-N1-C10	2.83	117.62	114.90
3	A	510	FAD	C1'-N10-C9A	4.50	122.31	118.31

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	510	FAD	1	0
4	A	511	GOL	7	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

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.