



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 12:48 AM GMT

PDB ID : 2BRY
Title : CRYSTAL STRUCTURE OF THE NATIVE MONOOXYGENASE DOMAIN
OF MICAL AT 1.45 Å RESOLUTION
Authors : Siebold, C.; Berrow, N.; Walter, T.S.; Harlos, K.; Owens, R.J.; Terman, J.R.;
Stuart, D.I.; Kolodkin, A.L.; Pasterkamp, R.J.; Jones, E.Y.
Deposited on : 2005-05-13
Resolution : 1.45 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
<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

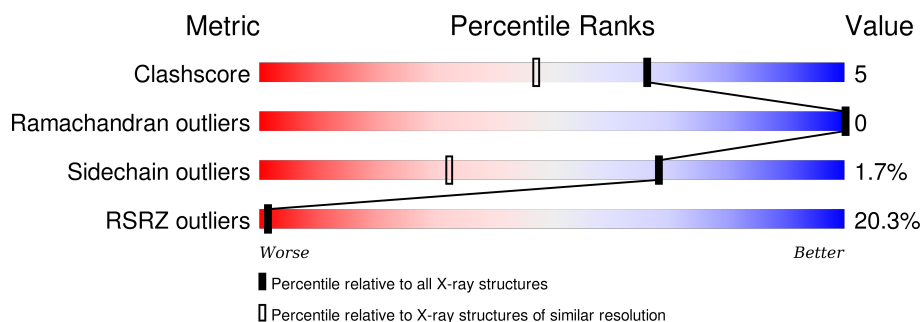
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.45 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	102246	1336 (1.48-1.44)
Ramachandran outliers	100387	1320 (1.48-1.44)
Sidechain outliers	100360	1320 (1.48-1.44)
RSRZ outliers	91569	1279 (1.48-1.44)

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.

Mol	Chain	Length	Quality of chain
1	A	497	<div> <div>31%</div> <div>86%</div> <div>10%</div> <div>•</div> </div>
1	B	497	<div> <div>8%</div> <div>86%</div> <div>10%</div> <div>•</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	GOL	A	1491	-	-	-	X

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 9235 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.

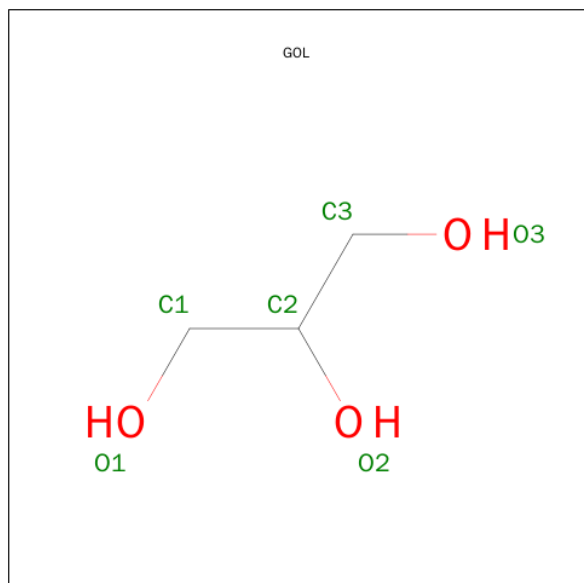
- Molecule 1 is a protein called NEDD9 INTERACTING PROTEIN WITH CALPONIN HO- MOLOGY AND LIM DOMAINS.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	479	Total	C	N	O	S	0	9	0
			3765	2406	669	673	17			
1	B	479	Total	C	N	O	S	0	8	0
			3763	2403	669	674	17			

- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

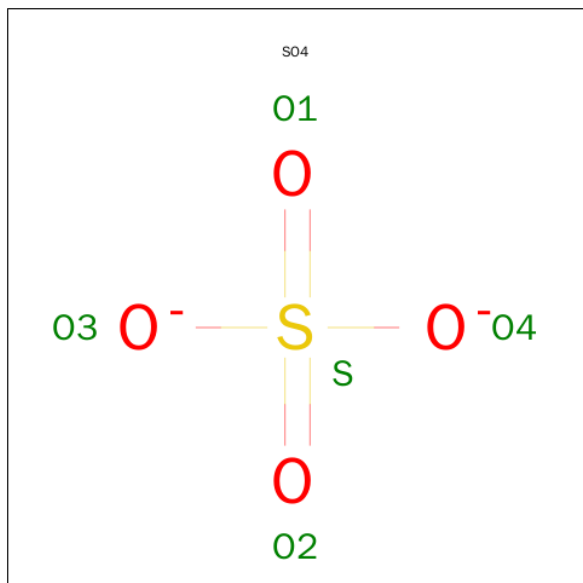
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Cl	0	0
			1	1		
2	A	1	Total	Cl	0	0
			1	1		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



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

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 5 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: C₂₇H₃₃N₉O₁₅P₂).



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

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	757	Total	O	0	0
			757	757		
6	B	826	Total	O	0	0
			826	826		

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	76.59 Å 89.55 Å 83.46 Å 90.00° 114.35° 90.00°	Depositor
Resolution (Å)	76.03 – 1.45 29.08 – 1.45	Depositor EDS
% Data completeness (in resolution range)	88.3 (76.03-1.45) 88.3 (29.08-1.45)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.41 (at 1.45 Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.180 , 0.222 0.230 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	18.0	Xtriage
Anisotropy	0.582	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 76.5	EDS
Estimated twinning fraction	0.023 for h,-k,-h-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	1 of 160456 reflections (0.001%)	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	9235	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.98% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, SO4, FAD, CL

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	0.65	2/3883 (0.1%)	0.69	2/5258 (0.0%)
1	B	0.56	0/3881	0.62	0/5255
All	All	0.61	2/7764 (0.0%)	0.66	2/10513 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	23[A]	CYS	CB-SG	-9.17	1.66	1.82
1	A	23[B]	CYS	CB-SG	-9.17	1.66	1.82

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	146	ARG	NE-CZ-NH2	-7.58	116.51	120.30
1	A	102	GLU	CG-CD-OE1	5.13	128.56	118.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	3765	0	3791	37	1
1	B	3763	0	3787	39	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	6	0	8	0	0
4	A	5	0	0	0	0
4	B	5	0	0	0	0
5	A	53	0	31	0	0
5	B	53	0	31	0	0
6	A	757	0	0	15	0
6	B	826	0	0	21	1
All	All	9235	0	7648	75	2

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 (75) 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:131[B]:THR:HG21	6:B:2318:HOH:O	1.54	1.08
1:B:267:GLN:HG3	1:B:269:PHE:HB3	1.53	0.89
1:A:152:LEU:HG	6:A:2338:HOH:O	1.80	0.81
1:A:254:GLN:HG3	6:A:2463:HOH:O	1.83	0.78
1:B:371:ARG:HB2	1:B:396[B]:VAL:HG12	1.73	0.70
1:A:422:ARG:NH1	1:A:435:GLU:OE1	2.24	0.70
1:B:478:GLN:HG3	6:B:2792:HOH:O	1.99	0.62
1:A:371:ARG:HB2	1:A:396[B]:VAL:HG12	1.81	0.60
1:B:78:GLN:NE2	6:B:2204:HOH:O	2.34	0.58
1:B:152:LEU:HD21	6:B:2508:HOH:O	2.02	0.57
1:B:131[A]:THR:HG21	6:B:2318:HOH:O	2.04	0.57
1:A:216[B]:SER:HB3	6:A:2425:HOH:O	2.06	0.55
1:B:63:LEU:HD12	6:B:2163:HOH:O	2.06	0.54
1:B:231:GLU:OE2	1:B:370:MET:HG3	2.07	0.54
1:B:118:LYS:HE2	6:B:2286:HOH:O	2.07	0.53
1:A:431:GLU:HG2	6:A:2357:HOH:O	2.09	0.53
1:A:59:SER:O	1:A:142:LYS:HE2	2.08	0.52
1:B:188:PRO:HD2	1:B:192:SER:HB3	1.91	0.52
1:B:216[B]:SER:HB3	6:B:2461:HOH:O	2.10	0.52
1:B:86:LYS:HE2	1:B:210:GLU:O	2.10	0.52
1:A:19:GLN:HG3	6:A:2034:HOH:O	2.09	0.52
1:A:371:ARG:CB	1:A:396[B]:VAL:HG12	2.40	0.51
1:A:230:ARG:HD2	1:A:373:GLU:HG3	1.92	0.51
1:A:348:LYS:HD2	6:A:2302:HOH:O	2.09	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:59:SER:O	1:B:142:LYS:HE2	2.10	0.51
1:A:235:LYS:HD2	1:B:225:GLU:OE1	2.11	0.51
1:B:67:LEU:HG	6:B:2163:HOH:O	2.12	0.50
1:A:203:PRO:HB2	6:A:2414:HOH:O	2.10	0.50
1:A:141:LYS:HD3	6:A:2337:HOH:O	2.12	0.50
1:A:146:ARG:HD2	1:A:290:GLU:OE2	2.11	0.49
1:B:477:ASN:HB2	6:B:2792:HOH:O	2.11	0.49
1:A:235:LYS:NZ	6:A:2451:HOH:O	2.44	0.48
1:A:279:ILE:HG22	1:A:281:LEU:HG	1.96	0.48
1:A:267:GLN:HB3	1:A:269:PHE:H	1.78	0.48
1:A:218:ALA:HB3	1:A:222:PHE:CG	2.49	0.48
1:B:133:HIS:HB2	6:B:2125:HOH:O	2.12	0.48
1:A:61:LYS:HE2	1:A:61:LYS:HB2	1.64	0.47
1:B:141:LYS:HE2	1:B:145:GLY:O	2.14	0.47
1:B:141:LYS:HD3	6:B:2335:HOH:O	2.15	0.47
1:A:188:PRO:HD2	1:A:192:SER:CB	2.45	0.46
1:B:397:GLU:HG3	6:B:2687:HOH:O	2.14	0.46
1:B:188:PRO:HD2	1:B:192:SER:CB	2.45	0.46
1:B:315:GLU:HG3	6:B:2595:HOH:O	2.16	0.46
1:B:127:LEU:HB3	1:B:131[B]:THR:HG23	1.99	0.45
1:B:370:MET:HB3	6:B:2665:HOH:O	2.16	0.45
1:B:189:ARG:O	1:B:192:SER:HB2	2.17	0.45
1:A:425:GLU:HG3	6:A:2647:HOH:O	2.16	0.45
1:A:437:GLU:HG2	6:A:2663:HOH:O	2.16	0.45
1:B:24:GLN:NE2	6:B:2039:HOH:O	2.41	0.45
1:A:142:LYS:HD3	6:A:2174:HOH:O	2.17	0.44
1:B:136:ARG:NH1	6:B:2333:HOH:O	2.51	0.44
1:A:77:GLN:HE21	1:A:81:ALA:HA	1.83	0.43
1:B:180:LYS:HE2	6:B:2406:HOH:O	2.18	0.43
1:B:195:ARG:HD3	6:B:2426:HOH:O	2.17	0.43
1:A:202:PRO:HB2	1:A:206:LEU:HB3	1.99	0.43
1:B:151:THR:HB	6:B:2363:HOH:O	2.18	0.43
1:A:188:PRO:HD2	1:A:192:SER:HB3	2.01	0.43
1:B:286:TYR:OH	1:B:292:HIS:CD2	2.72	0.43
1:B:425:GLU:HG3	6:B:2715:HOH:O	2.19	0.42
1:B:252:GLU:OE2	1:B:292:HIS:HE1	2.02	0.42
1:B:95:CYS:SG	1:B:409:GLY:HA3	2.60	0.42
1:B:120[B]:SER:HB2	1:B:355:ALA:HB2	2.01	0.42
1:A:462:PRO:HB2	1:A:469:LEU:HD11	2.00	0.42
1:B:7:THR:HB	1:B:8:ASN:H	1.72	0.42
1:A:57:TYR:CZ	1:A:59:SER:HB3	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:15:GLU:O	1:A:19:GLN:HG2	2.19	0.42
1:A:397:GLU:HB2	6:A:2618:HOH:O	2.20	0.41
1:B:153:ASP:N	1:B:153:ASP:OD1	2.53	0.41
1:A:120:SER:HB2	1:A:355:ALA:HB2	2.01	0.41
1:A:131[B]:THR:HG23	1:A:410:PHE:CD2	2.56	0.41
1:A:109:ARG:NH1	6:A:2251:HOH:O	2.43	0.41
1:B:286:TYR:OH	1:B:292:HIS:HD2	2.04	0.40
1:A:202:PRO:HG2	1:A:207:ALA:HB2	2.02	0.40
1:A:301:GLN:HG2	6:A:2515:HOH:O	2.20	0.40

All (2) 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:A:146:ARG:NH2	1:B:248:ARG:O[1_556]	2.03	0.17
6:B:2039:HOH:O	6:B:2786:HOH:O[2_655]	2.15	0.05

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	484/497 (97%)	481 (99%)	3 (1%)	0	100	100
1	B	484/497 (97%)	477 (99%)	7 (1%)	0	100	100
All	All	968/994 (97%)	958 (99%)	10 (1%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	397/403 (98%)	390 (98%)	7 (2%)	66	30
1	B	397/403 (98%)	389 (98%)	8 (2%)	63	26
All	All	794/806 (98%)	779 (98%)	15 (2%)	68	28

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	70	ARG
1	A	146	ARG
1	A	216[A]	SER
1	A	216[B]	SER
1	A	469	LEU
1	A	477	ASN
1	A	481	ASP
1	B	7	THR
1	B	61	LYS
1	B	216[A]	SER
1	B	216[B]	SER
1	B	453	ARG
1	B	469	LEU
1	B	475	THR
1	B	488	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	77	GLN
1	A	78	GLN
1	A	84	ASN
1	A	133	HIS
1	A	199	GLN
1	A	457	GLN
1	B	84	ASN
1	B	266	ASN
1	B	292	HIS
1	B	382	HIS

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 ⓘ

Of 7 ligands modelled in this entry, 2 are monoatomic - leaving 5 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	GOL	A	1491	-	5,5,5	0.74	0	5,5,5	0.74	0
4	SO4	A	1492	-	4,4,4	0.14	0	6,6,6	0.31	0
5	FAD	A	1493	-	48,58,58	1.30	6 (12%)	54,89,89	2.04	10 (18%)
4	SO4	B	1491	-	4,4,4	0.26	0	6,6,6	0.79	0
5	FAD	B	1492	-	48,58,58	1.30	6 (12%)	54,89,89	2.21	7 (12%)

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	GOL	A	1491	-	-	0/4/4/4	0/0/0/0
4	SO4	A	1492	-	-	0/0/0/0	0/0/0/0
5	FAD	A	1493	-	-	0/30/50/50	0/6/6/6
4	SO4	B	1491	-	-	0/0/0/0	0/0/0/0
5	FAD	B	1492	-	-	0/30/50/50	0/6/6/6

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	1493	FAD	C2A-N1A	2.07	1.37	1.33
5	A	1493	FAD	C5X-N5	2.33	1.39	1.35
5	B	1492	FAD	C2A-N1A	2.53	1.38	1.33
5	A	1493	FAD	C4X-N5	2.75	1.37	1.33
5	B	1492	FAD	C4X-N5	2.79	1.37	1.33
5	A	1493	FAD	C4-N3	2.96	1.38	1.33
5	B	1492	FAD	C4-N3	2.99	1.38	1.33
5	A	1493	FAD	C1'-N10	3.19	1.51	1.48
5	B	1492	FAD	C1'-N10	3.29	1.51	1.48
5	B	1492	FAD	C5X-N5	3.31	1.40	1.35
5	B	1492	FAD	C2A-N3A	4.03	1.39	1.32
5	A	1493	FAD	C2A-N3A	4.16	1.39	1.32

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	1492	FAD	N3A-C2A-N1A	-12.12	119.61	128.89
5	A	1493	FAD	N3A-C2A-N1A	-9.92	121.30	128.89
5	B	1492	FAD	C4X-C4-N3	-3.82	118.36	123.59
5	A	1493	FAD	C4X-C4-N3	-3.48	118.83	123.59
5	A	1493	FAD	C4A-C5A-N7A	-3.02	106.70	109.48
5	B	1492	FAD	C1B-N9A-C4A	-2.55	123.09	126.94
5	A	1493	FAD	C4B-O4B-C1B	-2.43	107.05	109.72
5	A	1493	FAD	C4X-N5-C5X	2.02	119.09	116.76
5	A	1493	FAD	C4-C4X-N5	2.06	121.22	118.72
5	B	1492	FAD	C2A-N1A-C6A	2.10	122.51	118.77
5	A	1493	FAD	C5X-C9A-N10	2.31	119.38	117.62
5	B	1492	FAD	C4X-C10-N10	2.32	121.89	120.52
5	A	1493	FAD	C4X-C10-N10	2.38	121.92	120.52
5	B	1492	FAD	C1'-N10-C9A	3.36	122.63	118.86
5	A	1493	FAD	C1'-N10-C9A	3.94	123.29	118.86
5	A	1493	FAD	C4-N3-C2	5.90	120.34	115.25
5	B	1492	FAD	C4-N3-C2	6.91	121.22	115.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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, 95th 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 ²)	Q<0.9
1	A	479/497 (96%)	1.56	155 (32%) 1 0	15, 20, 31, 45	0
1	B	479/497 (96%)	0.49	39 (8%) 15 15	6, 15, 29, 46	0
All	All	958/994 (96%)	1.02	194 (20%) 1 1	6, 19, 30, 46	0

All (194) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	265	TYR	11.4
1	A	265	TYR	8.5
1	A	151	THR	7.0
1	B	488	LYS	6.8
1	A	488	LYS	6.5
1	A	468	ASN	6.2
1	A	355	ALA	6.1
1	B	266	ASN	6.1
1	B	151	THR	6.1
1	B	7	THR	6.0
1	A	274	LEU	5.8
1	A	410	PHE	5.7
1	A	95	CYS	5.5
1	A	266	ASN	5.5
1	A	215	ILE	5.4
1	A	275	LYS	5.2
1	A	97	LEU	5.2
1	A	271	GLN	5.1
1	A	101	VAL	5.0
1	B	355	ALA	5.0
1	A	88	LEU	4.8
1	A	164	LEU	4.8
1	A	89	VAL	4.6
1	A	406	VAL	4.5

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Mol	Chain	Res	Type	RSRZ
1	A	160	LEU	4.5
1	A	171	LEU	4.5
1	A	100	ALA	4.5
1	A	473	ALA	4.5
1	A	311	GLN	4.4
1	A	391	VAL	4.4
1	A	453	ARG	4.4
1	A	99	ALA	4.3
1	A	276	ALA	4.3
1	A	301	GLN	4.3
1	A	486	MET	4.3
1	A	472	ARG	4.3
1	A	90	VAL	4.2
1	A	173	VAL	4.2
1	A	306	LEU	4.2
1	B	356	ARG	4.2
1	A	169	LEU	4.1
1	A	390	LEU	4.1
1	A	268	LYS	4.1
1	A	413	ALA	4.1
1	A	412	ALA	4.1
1	A	414	PHE	4.1
1	A	157	ILE	4.0
1	B	487	ASP	4.0
1	A	125	LEU	4.0
1	A	302	CYS	3.9
1	A	214	LEU	3.9
1	A	138	LEU	3.8
1	A	96	GLY	3.8
1	A	411	LEU	3.8
1	A	305	ARG	3.8
1	A	170	LEU	3.7
1	A	167	VAL	3.7
1	A	168	ALA	3.6
1	A	103	LEU	3.6
1	A	163	LEU	3.6
1	A	409	GLY	3.6
1	A	204	ALA	3.6
1	A	111	VAL	3.6
1	A	477	ASN	3.6
1	A	110	VAL	3.6
1	A	175	ILE	3.5

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Mol	Chain	Res	Type	RSRZ
1	A	92	ALA	3.5
1	A	416	ALA	3.5
1	A	314	SER	3.5
1	A	203	PRO	3.5
1	A	87	CYS	3.5
1	A	328	GLU	3.4
1	B	149	THR	3.4
1	A	94	PRO	3.4
1	A	400	TRP	3.4
1	A	124	VAL	3.4
1	A	155	ILE	3.4
1	A	112	LEU	3.4
1	A	323	ALA	3.4
1	A	165	LEU	3.3
1	A	304	LEU	3.3
1	B	301	GLN	3.3
1	A	293	TYR	3.3
1	B	148	CYS	3.3
1	B	150	GLY	3.3
1	A	201	ASN	3.3
1	A	485	MET	3.3
1	B	268	LYS	3.2
1	A	278	GLY	3.2
1	B	426	GLY	3.2
1	B	190	LYS	3.2
1	A	315	GLU	3.1
1	A	475	THR	3.1
1	A	131[A]	THR	3.1
1	A	135	LEU	3.1
1	B	152	LEU	3.1
1	A	213	VAL	3.1
1	A	480	GLN	3.1
1	A	313	LEU	3.0
1	A	326	VAL	3.0
1	A	93	GLY	3.0
1	A	143	PHE	3.0
1	A	320	LEU	3.0
1	A	312	ASP	3.0
1	A	481	ASP	3.0
1	B	222	PHE	3.0
1	A	150	GLY	3.0
1	B	453	ARG	3.0

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Mol	Chain	Res	Type	RSRZ
1	B	233	ARG	2.9
1	B	489	GLU	2.9
1	A	190	LYS	2.9
1	A	318	GLN	2.9
1	B	468	ASN	2.9
1	A	104	ALA	2.9
1	A	356	ARG	2.9
1	A	279	ILE	2.9
1	A	294	PHE	2.9
1	A	119	PHE	2.9
1	A	188	PRO	2.9
1	B	481	ASP	2.9
1	B	201	ASN	2.9
1	A	407	ALA	2.8
1	A	191	GLY	2.8
1	B	483	TYR	2.8
1	A	140	ALA	2.8
1	A	332	ARG	2.7
1	A	267	GLN	2.7
1	A	310	ARG	2.7
1	A	308	VAL	2.7
1	A	309	LEU	2.7
1	A	471	LEU	2.7
1	A	479	VAL	2.7
1	B	188	PRO	2.7
1	A	405	GLY	2.7
1	A	382	HIS	2.6
1	A	152	LEU	2.6
1	A	7	THR	2.6
1	A	483	TYR	2.6
1	A	348	LYS	2.6
1	A	489	GLU	2.6
1	B	454	ASN	2.6
1	A	287	TYR	2.6
1	A	162	LEU	2.6
1	A	454	ASN	2.5
1	A	98	ARG	2.5
1	B	147	PHE	2.5
1	A	420	VAL	2.5
1	A	469	LEU	2.5
1	B	477	ASN	2.5
1	A	91	GLY	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	67	LEU	2.4
1	A	137	ALA	2.4
1	A	222	PHE	2.4
1	A	417	ALA	2.4
1	A	139	GLY	2.4
1	A	189	ARG	2.4
1	A	30	PHE	2.4
1	A	270	PHE	2.4
1	A	389	GLY	2.4
1	B	475	THR	2.4
1	A	418	TRP	2.4
1	A	113	VAL	2.4
1	A	295	VAL	2.4
1	A	370	MET	2.4
1	A	17	PHE	2.4
1	A	478	GLN	2.3
1	B	274	LEU	2.3
1	A	144	TYR	2.3
1	A	37	LEU	2.3
1	B	232	MET	2.3
1	A	233	ARG	2.3
1	B	275	LYS	2.3
1	A	325	VAL	2.3
1	A	474	VAL	2.3
1	A	316	THR	2.3
1	A	467	PRO	2.3
1	A	452	HIS	2.2
1	B	486	MET	2.2
1	A	217	ALA	2.2
1	A	149	THR	2.2
1	A	322	LYS	2.2
1	A	23[A]	CYS	2.2
1	A	127	LEU	2.2
1	B	322	LYS	2.2
1	B	276	ALA	2.2
1	B	384	ALA	2.2
1	A	240	ILE	2.2
1	B	305	ARG	2.2
1	A	277	THR	2.1
1	B	223[A]	VAL	2.1
1	A	273	LEU	2.1
1	A	395	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	369	MET	2.0
1	B	485	MET	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

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, 95th 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 ²)	Q<0.9
3	GOL	A	1491	6/6	0.74	0.19	2.96	18,21,22,22	0
2	CL	B	1490	1/1	1.00	0.08	0.38	16,16,16,16	0
5	FAD	B	1492	53/53	0.97	0.07	-0.38	5,8,12,13	0
4	SO4	A	1492	5/5	0.96	0.10	-1.70	30,30,32,32	0
5	FAD	A	1493	53/53	0.98	0.07	-2.29	6,9,10,12	0
2	CL	A	1490	1/1	0.99	0.04	-3.72	12,12,12,12	0
4	SO4	B	1491	5/5	0.97	0.11	-	22,22,26,27	0

6.5 Other polymers [i](#)

There are no such residues in this entry.