



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 11:22 PM GMT

PDB ID : 2BXR
Title : HUMAN MONOAMINE OXIDASE A IN COMPLEX WITH CLORGYLINE,
CRYSTAL FORM A
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Deposited on : 2005-07-27
Resolution : 3.00 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

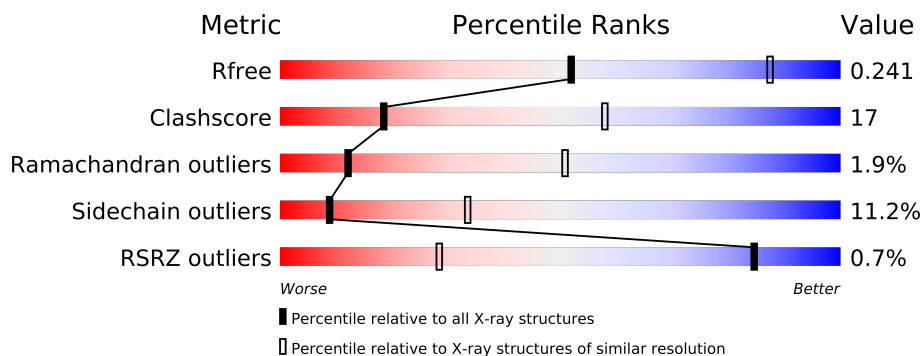
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
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) : stable22683

1 Overall quality at a glance



The reported resolution of this entry is 3.00 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	1216 (3.00-3.00)
Clashscore	79885	1594 (3.00-3.00)
Ramachandran outliers	78287	1537 (3.00-3.00)
Sidechain outliers	78261	1540 (3.00-3.00)
RSRZ outliers	66119	1217 (3.00-3.00)

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. 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	527	
1	B	527	

2 Entry composition i

There are 3 unique types of molecules in this entry. The entry contains 7186 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 AMINE OXIDASE [FLAVIN-CONTAINING] A.

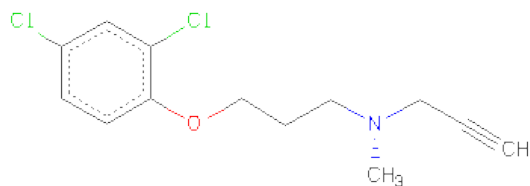
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	445	Total	C	N	O	S	0	0	0
			3523	2253	601	648	21			
1	B	445	Total	C	N	O	S	0	0	0
			3523	2253	601	648	21			

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



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

- Molecule 3 is N-[3-(2,4-DICHLOROPHENOXY)PROPYL]-N-METHYL-N-PROP-2-YNYL AMINE (three-letter code: MLG) (formula: $C_{13}H_{15}Cl_2NO$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	Cl	N	O	0	0
			17	13	2	1	1		
3	B	1	Total	C	Cl	N	O	0	0
			17	13	2	1	1		

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	143.49Å 109.60Å 81.33Å 90.00° 95.18° 90.00°	Depositor
Resolution (Å)	87.04 – 3.00 14.97 – 3.00	Depositor EDS
% Data completeness (in resolution range)	96.5 (87.04-3.00) 97.4 (14.97-3.00)	Depositor EDS
R_{merge}	0.01	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.67 (at 3.01Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.192 , 0.238 0.194 , 0.241	Depositor DCC
R_{free} test set	1245 reflections (5.39%)	DCC
Wilson B-factor (Å ²)	69.1	Xtriage
Anisotropy	0.434	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 26.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 24331 reflections	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	7186	wwPDB-VP
Average B, all atoms (Å ²)	64.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.15% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities 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: MLG, 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	0.75	2/3606 (0.1%)	0.78	3/4886 (0.1%)
1	B	0.74	3/3606 (0.1%)	0.77	3/4886 (0.1%)
All	All	0.75	5/7212 (0.1%)	0.77	6/9772 (0.1%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	395	LYS	CD-CE	7.94	1.71	1.51
1	A	395	LYS	CE-NZ	7.74	1.68	1.49
1	B	395	LYS	CD-CE	7.68	1.70	1.51
1	B	395	LYS	CE-NZ	7.35	1.67	1.49
1	B	116	TRP	CB-CG	5.69	1.60	1.50

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	395	LYS	CD-CE-NZ	7.82	129.68	111.70
1	A	303	VAL	CB-CA-C	-7.35	97.44	111.40
1	B	395	LYS	CD-CE-NZ	7.30	128.49	111.70
1	B	303	VAL	CB-CA-C	-7.06	97.98	111.40
1	A	56	ARG	NE-CZ-NH2	5.65	123.12	120.30
1	B	51	ARG	NE-CZ-NH2	-5.15	117.73	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	3523	0	3466	115	0
1	B	3523	0	3466	111	0
2	A	53	0	29	17	0
2	B	53	0	29	7	0
3	A	17	0	15	9	0
3	B	17	0	15	1	0
All	All	7186	0	7020	236	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 17.

All (236) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:A:600:FAD:C5X	3:A:601:MLG:H16	1.27	1.58
1:A:395:LYS:CE	1:A:395:LYS:NZ	1.68	1.54
2:A:600:FAD:N5	3:A:601:MLG:H16	1.35	1.21
1:A:297:ARG:HG3	1:A:297:ARG:HH11	1.12	1.10
1:B:297:ARG:HG3	1:B:297:ARG:HH11	0.94	1.07
1:B:121:TYR:HA	1:B:124:TYR:HD1	1.23	1.02
1:A:395:LYS:HE2	1:A:400:GLU:OE2	1.59	1.01
1:A:156:THR:HG22	1:A:159:GLU:HG2	1.45	0.97
1:B:121:TYR:HA	1:B:124:TYR:CD1	2.01	0.95
1:A:121:TYR:HA	1:A:124:TYR:HD1	1.30	0.94
2:A:600:FAD:C9A	3:A:601:MLG:H16	1.99	0.92
1:B:156:THR:HG22	1:B:159:GLU:HG2	1.52	0.91
1:A:297:ARG:HG3	1:A:297:ARG:NH1	1.79	0.91
1:B:297:ARG:NH1	1:B:297:ARG:HG3	1.69	0.91
2:A:600:FAD:C4X	3:A:601:MLG:C16	2.49	0.90
1:B:297:ARG:CG	1:B:297:ARG:HH11	1.83	0.89
1:A:121:TYR:HA	1:A:124:TYR:CD1	2.08	0.88
1:A:216:GLU:OE1	1:A:216:GLU:HA	1.74	0.86
2:A:600:FAD:C4X	3:A:601:MLG:H16	2.05	0.85
1:B:275:PRO:HD3	1:B:436:GLU:HG3	1.59	0.85
1:B:324:MET:SD	1:B:380:VAL:HG11	2.17	0.83

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:275:PRO:HD3	1:A:436:GLU:HG3	1.60	0.83
1:B:185:GLU:OE2	1:B:356:ARG:HD2	1.79	0.83
1:B:323:CYS:HA	1:B:336:THR:HG22	1.61	0.81
1:B:216:GLU:HA	1:B:216:GLU:OE1	1.80	0.81
1:A:397:TRP:CE2	2:A:600:FAD:HM82	2.17	0.80
1:A:297:ARG:HH11	1:A:297:ARG:CG	1.94	0.80
1:B:397:TRP:CE2	2:B:600:FAD:HM82	2.17	0.79
1:A:314:PHE:HA	1:A:317:LYS:HE3	1.64	0.78
1:B:314:PHE:HA	1:B:317:LYS:HE3	1.66	0.77
1:A:134:MET:HE1	1:A:160:LEU:HD21	1.67	0.77
1:A:251:SER:O	1:A:253:ASN:N	2.18	0.77
1:B:121:TYR:CA	1:B:124:TYR:HD1	1.99	0.76
1:A:323:CYS:HA	1:A:336:THR:HG22	1.68	0.76
1:A:185:GLU:OE2	1:A:356:ARG:HD2	1.85	0.75
2:B:600:FAD:N1	2:B:600:FAD:H2'	2.00	0.75
1:B:304:ILE:HB	1:B:353:ILE:HG23	1.67	0.74
1:B:134:MET:HE1	1:B:160:LEU:HD21	1.69	0.74
1:B:242:HIS:NE2	1:B:262:GLU:OE2	2.20	0.74
1:A:369:ARG:NH1	1:A:394:GLU:OE2	2.21	0.74
1:A:324:MET:SD	1:A:380:VAL:HG11	2.28	0.73
1:A:400:GLU:HA	1:A:400:GLU:OE1	1.88	0.72
1:A:156:THR:HG22	1:A:159:GLU:H	1.55	0.72
1:B:251:SER:O	1:B:253:ASN:N	2.22	0.72
1:B:156:THR:HG22	1:B:159:GLU:CG	2.20	0.71
1:A:304:ILE:HB	1:A:353:ILE:HG23	1.72	0.71
1:A:395:LYS:CE	1:A:400:GLU:OE2	2.38	0.70
1:A:121:TYR:HA	1:A:124:TYR:HB2	1.74	0.70
1:A:94:SER:HB3	1:A:95:GLU:CD	2.12	0.70
1:B:94:SER:HB3	1:B:95:GLU:CD	2.11	0.70
1:A:94:SER:HB3	1:A:95:GLU:OE1	1.91	0.69
1:B:400:GLU:HA	1:B:400:GLU:OE1	1.91	0.69
1:B:369:ARG:NH1	1:B:394:GLU:OE2	2.26	0.68
1:A:156:THR:HG23	1:A:158:LYS:H	1.57	0.68
1:B:121:TYR:HA	1:B:124:TYR:HB2	1.76	0.67
1:B:94:SER:HB3	1:B:95:GLU:OE1	1.95	0.67
1:A:197:TYR:O	1:A:200:GLN:HG3	1.95	0.66
1:A:156:THR:HG22	1:A:159:GLU:CG	2.22	0.66
1:A:443:GLY:O	2:A:600:FAD:O3'	2.13	0.66
1:A:87:GLU:OE1	1:A:87:GLU:HA	1.96	0.66
1:A:242:HIS:NE2	1:A:262:GLU:OE2	2.30	0.65
1:A:95:GLU:HG2	1:A:319:ASP:OD1	1.97	0.65
1:B:324:MET:H	1:B:336:THR:HG22	1.64	0.63

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:365:HIS:CD2	1:B:367:GLU:H	2.16	0.63
1:A:121:TYR:CA	1:A:124:TYR:HD1	2.07	0.63
1:B:121:TYR:O	1:B:124:TYR:HB2	1.98	0.62
1:B:156:THR:HG23	1:B:158:LYS:H	1.64	0.62
1:B:67:GLY:HA2	2:B:600:FAD:C4X	2.29	0.62
1:B:353:ILE:C	1:B:353:ILE:HD13	2.20	0.62
1:B:324:MET:HG2	1:B:336:THR:HG21	1.82	0.62
1:B:397:TRP:CD1	2:B:600:FAD:HM71	2.34	0.61
1:A:51:ARG:HG2	1:A:406:CYS:SG	2.40	0.61
1:A:124:TYR:HD2	1:A:128:TRP:CH2	2.18	0.61
1:A:253:ASN:HD22	1:A:265:GLU:HB2	1.66	0.61
1:A:324:MET:H	1:A:336:THR:HG22	1.66	0.61
1:A:121:TYR:O	1:A:124:TYR:HB2	2.00	0.61
1:A:406:CYS:HB3	1:A:407:TYR:HA	1.82	0.60
1:B:253:ASN:ND2	1:B:265:GLU:HB2	2.16	0.60
1:A:324:MET:HG2	1:A:336:THR:HG21	1.83	0.60
1:B:156:THR:HG22	1:B:159:GLU:H	1.67	0.60
1:A:365:HIS:CD2	1:A:367:GLU:H	2.20	0.59
1:A:156:THR:CG2	1:A:159:GLU:H	2.15	0.59
1:B:253:ASN:HD22	1:B:265:GLU:HB2	1.67	0.59
1:A:384:GLN:OE1	1:A:384:GLN:HA	2.02	0.59
1:A:397:TRP:CD1	2:A:600:FAD:HM71	2.37	0.59
1:A:324:MET:HG2	1:A:336:THR:CG2	2.32	0.59
1:A:204:THR:O	1:A:207:ILE:HG22	2.03	0.59
1:A:253:ASN:ND2	1:A:265:GLU:HB2	2.17	0.58
1:B:87:GLU:OE1	1:B:87:GLU:HA	2.03	0.58
1:A:353:ILE:HD13	1:A:353:ILE:C	2.24	0.58
1:A:322:GLY:O	1:A:336:THR:CG2	2.52	0.58
1:A:304:ILE:HB	1:A:353:ILE:CG2	2.33	0.58
1:B:395:LYS:HE2	1:B:400:GLU:OE2	2.02	0.58
1:A:422:VAL:HG13	1:A:422:VAL:O	2.04	0.57
1:B:324:MET:N	1:B:336:THR:HG22	2.19	0.57
1:B:324:MET:SD	1:B:380:VAL:CG1	2.91	0.57
1:B:197:TYR:O	1:B:200:GLN:HG3	2.06	0.56
1:A:324:MET:N	1:A:336:THR:HG22	2.20	0.56
1:B:324:MET:HG2	1:B:336:THR:CG2	2.35	0.56
1:B:204:THR:O	1:B:207:ILE:HG22	2.06	0.56
1:B:397:TRP:CE3	1:B:400:GLU:HG3	2.41	0.55
1:A:67:GLY:HA2	2:A:600:FAD:C4X	2.36	0.55
1:A:156:THR:HG23	1:A:158:LYS:N	2.22	0.55
1:A:421:ARG:HH11	1:B:35:TYR:HE1	1.53	0.55
1:B:365:HIS:CD2	1:B:366:LYS:N	2.76	0.54

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:256:ILE:HD11	1:A:269:VAL:HG22	1.89	0.54
1:A:60:VAL:HG21	1:A:63:VAL:HG12	1.89	0.54
1:B:378:ALA:HB2	1:B:387:LEU:HD13	1.89	0.54
1:A:397:TRP:CE3	1:A:400:GLU:HG3	2.43	0.54
1:B:322:GLY:O	1:B:336:THR:CG2	2.56	0.54
1:A:333:ILE:HD13	1:A:353:ILE:HB	1.90	0.54
1:B:95:GLU:HG2	1:B:319:ASP:OD1	2.09	0.53
1:B:117:ASN:HB3	1:B:120:ALA:HB3	1.91	0.53
1:B:365:HIS:HD2	1:B:366:LYS:N	2.06	0.53
1:B:365:HIS:HD2	1:B:367:GLU:H	1.55	0.53
1:B:51:ARG:HG2	1:B:406:CYS:SG	2.49	0.53
1:B:124:TYR:HD2	1:B:128:TRP:CH2	2.27	0.53
1:B:323:CYS:CA	1:B:336:THR:HG22	2.34	0.53
1:A:323:CYS:CA	1:A:336:THR:HG22	2.38	0.53
1:B:156:THR:CG2	1:B:159:GLU:H	2.22	0.52
1:B:397:TRP:O	1:B:400:GLU:HB2	2.10	0.52
1:A:407:TYR:CZ	3:A:601:MLG:C15	2.93	0.52
2:A:600:FAD:C10	3:A:601:MLG:H16	2.40	0.52
1:B:406:CYS:HB3	1:B:407:TYR:HA	1.91	0.51
1:A:365:HIS:CD2	1:A:366:LYS:N	2.78	0.51
1:A:340:THR:HG23	1:A:347:PRO:HA	1.91	0.51
2:A:600:FAD:N10	3:A:601:MLG:H16	2.24	0.51
1:A:92:ASN:ND2	1:A:94:SER:OG	2.33	0.51
1:B:422:VAL:HG13	1:B:422:VAL:O	2.11	0.51
1:B:124:TYR:HE2	1:B:173:PHE:CD2	2.29	0.51
1:B:323:CYS:HB3	1:B:337:LEU:HD23	1.91	0.51
1:A:124:TYR:HE2	1:A:173:PHE:CD2	2.29	0.51
1:A:365:HIS:HD2	1:A:366:LYS:N	2.09	0.51
1:B:14:PHE:O	1:B:266:CYS:HA	2.11	0.50
1:B:400:GLU:CA	1:B:400:GLU:OE1	2.58	0.50
1:A:421:ARG:NH1	1:B:35:TYR:HE1	2.09	0.50
2:B:600:FAD:N10	3:B:601:MLG:H16	2.04	0.50
1:A:294:LEU:HD21	1:A:423:ILE:HG12	1.93	0.50
1:B:384:GLN:OE1	1:B:384:GLN:HA	2.12	0.50
2:A:600:FAD:H2'	2:A:600:FAD:N1	2.26	0.50
1:B:333:ILE:HD13	1:B:353:ILE:HB	1.94	0.50
1:A:43:GLU:OE1	2:A:600:FAD:H1B	2.12	0.50
1:B:395:LYS:CE	1:B:400:GLU:OE2	2.59	0.50
1:B:294:LEU:HD21	1:B:423:ILE:HG12	1.93	0.50
1:B:304:ILE:HB	1:B:353:ILE:CG2	2.39	0.49
1:B:340:THR:HG23	1:B:347:PRO:HA	1.94	0.49
1:B:121:TYR:CA	1:B:124:TYR:HB2	2.43	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:365:HIS:HD2	1:A:367:GLU:H	1.58	0.49
1:B:140:THR:HG22	1:B:199:LYS:CD	2.42	0.49
1:A:322:GLY:O	1:A:336:THR:HG23	2.12	0.49
1:A:323:CYS:HB3	1:A:337:LEU:HD23	1.95	0.48
1:B:322:GLY:O	1:B:336:THR:HG23	2.13	0.48
1:A:45:ARG:NH2	1:A:403:SER:OG	2.42	0.48
2:A:600:FAD:N1	2:A:600:FAD:C2'	2.76	0.48
1:B:445:MET:O	1:B:446:GLU:C	2.52	0.48
1:B:246:HIS:HB2	1:B:257:GLU:HB3	1.95	0.48
1:A:378:ALA:HB2	1:A:387:LEU:HD13	1.95	0.48
1:A:100:TYR:HD2	1:A:326:ILE:HG23	1.79	0.48
1:A:107:PRO:O	1:A:108:PHE:CB	2.61	0.48
1:A:209:SER:HB2	1:A:214:GLY:HA3	1.96	0.47
1:A:406:CYS:CB	1:A:407:TYR:HA	2.42	0.47
1:B:216:GLU:OE1	1:B:216:GLU:CA	2.59	0.47
1:B:140:THR:HG22	1:B:199:LYS:HD3	1.95	0.47
1:B:256:ILE:HD11	1:B:269:VAL:HG22	1.95	0.47
1:A:40:LEU:HD21	1:A:42:LEU:HD21	1.97	0.47
1:A:205:THR:HG22	1:A:206:ARG:N	2.29	0.47
1:A:121:TYR:CA	1:A:124:TYR:HB2	2.43	0.47
1:B:444:TYR:CD1	1:B:444:TYR:N	2.82	0.47
1:B:297:ARG:CG	1:B:297:ARG:NH1	2.51	0.46
1:A:432:PHE:O	1:A:454:ARG:NH2	2.40	0.46
1:B:353:ILE:CD1	1:B:358:ALA:HA	2.46	0.46
1:A:168:LYS:HB2	1:A:168:LYS:HE3	1.71	0.45
1:B:43:GLU:OE1	2:B:600:FAD:H1B	2.17	0.45
2:A:600:FAD:C6	3:A:601:MLG:C16	2.85	0.45
1:B:156:THR:HG23	1:B:158:LYS:N	2.31	0.45
1:A:369:ARG:NH1	1:A:394:GLU:CD	2.70	0.45
1:A:60:VAL:O	1:A:62:TYR:N	2.49	0.45
1:B:209:SER:HB2	1:B:214:GLY:HA3	1.98	0.45
1:B:107:PRO:O	1:B:108:PHE:CB	2.64	0.45
1:A:400:GLU:HB3	1:A:403:SER:HB2	1.98	0.45
1:B:45:ARG:NH2	1:B:403:SER:OG	2.46	0.45
1:B:40:LEU:HD21	1:B:42:LEU:HD21	1.99	0.45
1:A:14:PHE:O	1:A:266:CYS:HA	2.17	0.44
1:B:406:CYS:CB	1:B:407:TYR:HA	2.48	0.44
1:A:124:TYR:CD2	1:A:128:TRP:CH2	3.03	0.44
1:A:19:ILE:O	1:A:272:ALA:HB3	2.17	0.44
1:A:216:GLU:CA	1:A:216:GLU:OE1	2.56	0.44
1:A:106:TYR:HA	1:A:107:PRO:HD3	1.78	0.44
1:A:77:ILE:HD13	1:A:445:MET:HB3	1.99	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:397:TRP:CD2	2:B:600:FAD:HM82	2.52	0.43
1:B:100:TYR:HD2	1:B:326:ILE:HG23	1.83	0.43
1:A:400:GLU:OE1	1:A:400:GLU:CA	2.57	0.43
1:A:397:TRP:CD2	2:A:600:FAD:HM82	2.54	0.43
1:A:16:VAL:HG22	1:A:39:VAL:HG13	1.99	0.43
1:B:323:CYS:HA	1:B:336:THR:CG2	2.42	0.43
1:B:240:LEU:O	1:B:242:HIS:ND1	2.52	0.43
1:A:353:ILE:CD1	1:A:358:ALA:HA	2.48	0.43
1:B:369:ARG:NH1	1:B:394:GLU:CD	2.72	0.43
1:A:269:VAL:HG12	1:A:270:ILE:N	2.33	0.43
1:A:453:GLU:O	1:A:457:ARG:HG3	2.19	0.43
1:A:391:HIS:CG	1:A:392:TYR:N	2.87	0.43
1:A:117:ASN:HB3	1:A:120:ALA:HB3	2.01	0.43
1:A:324:MET:SD	1:A:380:VAL:CG1	3.02	0.42
1:A:241:ASN:C	1:A:243:PRO:HD3	2.39	0.42
1:B:100:TYR:CE1	1:B:103:GLY:O	2.72	0.42
1:A:92:ASN:OD1	1:A:94:SER:HB2	2.18	0.42
1:B:385:GLU:HG2	1:B:385:GLU:H	1.59	0.42
1:B:332:PRO:O	1:B:353:ILE:HG12	2.20	0.42
1:B:205:THR:HG22	1:B:206:ARG:N	2.34	0.42
1:B:307:MET:HB3	1:B:309:TYR:CE2	2.54	0.42
1:B:60:VAL:HG21	1:B:63:VAL:HG12	2.02	0.42
1:B:40:LEU:HG	1:B:41:VAL:N	2.34	0.42
1:A:304:ILE:HD12	1:A:353:ILE:HD12	2.02	0.42
1:A:95:GLU:CG	1:A:319:ASP:OD1	2.67	0.42
1:A:140:THR:HG22	1:A:199:LYS:CD	2.49	0.42
1:A:397:TRP:O	1:A:400:GLU:HB2	2.20	0.41
1:A:121:TYR:HA	1:A:124:TYR:CB	2.47	0.41
1:B:400:GLU:HB3	1:B:403:SER:HB2	2.02	0.41
1:B:420:GLY:O	1:B:423:ILE:HB	2.20	0.41
1:A:408:THR:OG1	1:A:409:ALA:N	2.50	0.41
1:B:292:ASN:O	1:B:296:GLN:HB2	2.20	0.41
1:A:397:TRP:CD2	2:A:600:FAD:C8M	3.03	0.41
1:B:254:ILE:O	1:B:265:GLU:HA	2.20	0.41
1:B:444:TYR:HD1	1:B:444:TYR:N	2.18	0.41
1:B:125:ASN:HA	1:B:128:TRP:HE3	1.85	0.41
1:A:40:LEU:HG	1:A:41:VAL:N	2.35	0.41
1:B:77:ILE:HD13	1:B:445:MET:HB3	2.02	0.41
1:A:100:TYR:CE1	1:A:103:GLY:O	2.73	0.41
1:A:307:MET:SD	1:A:350:MET:HG2	2.61	0.41
1:B:337:LEU:HD23	1:B:337:LEU:HA	1.94	0.40
1:A:369:ARG:HH11	1:A:394:GLU:CD	2.24	0.40

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:332:PRO:O	1:A:353:ILE:HG12	2.22	0.40
1:B:241:ASN:C	1:B:243:PRO:HD3	2.41	0.40
1:B:391:HIS:CG	1:B:392:TYR:N	2.89	0.40
1:B:96:ARG:HE	1:B:96:ARG:HB2	1.57	0.40
1:B:427:VAL:O	1:B:427:VAL:HG23	2.21	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	439/527 (83%)	392 (89%)	38 (9%)	9 (2%)	11	47
1	B	439/527 (83%)	391 (89%)	40 (9%)	8 (2%)	13	53
All	All	878/1054 (83%)	783 (89%)	78 (9%)	17 (2%)	12	51

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	108	PHE
1	A	109	ARG
1	A	215	GLN
1	A	252	ASP
1	B	108	PHE
1	B	109	ARG
1	B	215	GLN
1	B	252	ASP
1	A	118	PRO
1	B	118	PRO
1	A	61	ASP
1	B	61	ASP
1	A	399	GLU
1	A	428	GLY
1	B	107	PRO

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Mol	Chain	Res	Type
1	A	107	PRO
1	B	332	PRO

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	371/450 (82%)	330 (89%)	41 (11%)	9	34
1	B	371/450 (82%)	329 (89%)	42 (11%)	9	33
All	All	742/900 (82%)	659 (89%)	83 (11%)	9	33

All (83) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	HIS
1	A	70	VAL
1	A	77	ILE
1	A	78	LEU
1	A	94	SER
1	A	105	THR
1	A	119	ILE
1	A	121	TYR
1	A	140	THR
1	A	147	GLN
1	A	154	LYS
1	A	156	THR
1	A	176	LEU
1	A	209	SER
1	A	210	VAL
1	A	216	GLU
1	A	252	ASP
1	A	253	ASN
1	A	265	GLU
1	A	296	GLN
1	A	297	ARG
1	A	303	VAL
1	A	317	LYS

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Mol	Chain	Res	Type
1	A	318	LYS
1	A	319	ASP
1	A	321	CYS
1	A	324	MET
1	A	329	GLU
1	A	330	ASP
1	A	336	THR
1	A	353	ILE
1	A	356	ARG
1	A	383	SER
1	A	385	GLU
1	A	387	LEU
1	A	407	TYR
1	A	421	ARG
1	A	422	VAL
1	A	423	ILE
1	A	424	ARG
1	A	436	GLU
1	B	16	VAL
1	B	55	ILE
1	B	70	VAL
1	B	77	ILE
1	B	78	LEU
1	B	94	SER
1	B	105	THR
1	B	118	PRO
1	B	119	ILE
1	B	121	TYR
1	B	147	GLN
1	B	154	LYS
1	B	156	THR
1	B	176	LEU
1	B	209	SER
1	B	210	VAL
1	B	216	GLU
1	B	252	ASP
1	B	253	ASN
1	B	265	GLU
1	B	296	GLN
1	B	297	ARG
1	B	303	VAL
1	B	317	LYS

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Mol	Chain	Res	Type
1	B	318	LYS
1	B	319	ASP
1	B	321	CYS
1	B	329	GLU
1	B	330	ASP
1	B	336	THR
1	B	353	ILE
1	B	356	ARG
1	B	383	SER
1	B	385	GLU
1	B	387	LEU
1	B	407	TYR
1	B	421	ARG
1	B	422	VAL
1	B	423	ILE
1	B	424	ARG
1	B	436	GLU
1	B	444	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	147	GLN
1	A	179	ASN
1	A	237	GLN
1	A	253	ASN
1	A	296	GLN
1	A	365	HIS
1	B	147	GLN
1	B	148	HIS
1	B	237	GLN
1	B	253	ASN
1	B	365	HIS

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 ⓘ

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	600	1,3	58,58,58	2.91	18 (31%)	85,89,89	2.38	27 (31%)
3	MLG	A	601	2	17,17,17	4.03	4 (23%)	21,21,21	12.26	7 (33%)
2	FAD	B	600	1,3	58,58,58	2.75	17 (29%)	85,89,89	2.17	23 (27%)
3	MLG	B	601	2	17,17,17	3.20	3 (17%)	21,21,21	11.63	8 (38%)

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	600	1,3	-	0/34/50/50	0/1/6/6
3	MLG	A	601	2	-	0/9/10/10	0/1/1/1
2	FAD	B	600	1,3	-	0/34/50/50	0/1/6/6
3	MLG	B	601	2	-	0/9/10/10	0/1/1/1

All (42) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	601	MLG	C15-C16	14.85	1.51	1.17
3	B	601	MLG	C15-C16	11.02	1.43	1.17

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	600	FAD	C8M-C8	10.05	1.72	1.51
2	A	600	FAD	C4-C4X	-9.69	1.25	1.41
2	B	600	FAD	C4-C4X	-9.09	1.26	1.41
2	A	600	FAD	C1'-C2'	-9.03	1.43	1.51
2	B	600	FAD	C10-N1	8.96	1.51	1.35
2	B	600	FAD	C1'-N10	7.54	1.56	1.48
3	A	601	MLG	C14-N13	-6.35	1.36	1.46
2	B	600	FAD	C4-N3	6.22	1.47	1.37
2	A	600	FAD	C10-N1	5.88	1.46	1.35
2	A	600	FAD	C9-C9A	5.22	1.51	1.40
2	A	600	FAD	O4-C4	5.16	1.34	1.24
3	B	601	MLG	C14-N13	-5.13	1.38	1.46
2	B	600	FAD	P-O3P	4.61	1.68	1.59
2	A	600	FAD	C9A-N10	-4.32	1.32	1.38
2	B	600	FAD	C9-C9A	4.19	1.49	1.40
3	B	601	MLG	C14-C15	-4.07	1.41	1.47
2	B	600	FAD	C2A-N3A	3.95	1.40	1.32
2	B	600	FAD	C8M-C8	3.94	1.59	1.51
2	B	600	FAD	C5X-N5	3.93	1.41	1.35
2	A	600	FAD	C1'-N10	3.72	1.52	1.48
2	B	600	FAD	C2-N1	-3.43	1.27	1.35
2	A	600	FAD	C5X-N5	3.27	1.40	1.35
2	B	600	FAD	O2'-C2'	3.23	1.50	1.43
2	B	600	FAD	C2A-N1A	2.90	1.39	1.33
2	B	600	FAD	C6-C7	2.83	1.45	1.37
2	A	600	FAD	C4-N3	2.79	1.41	1.37
2	A	600	FAD	C6-C7	2.73	1.45	1.37
2	A	600	FAD	C9A-C5X	-2.67	1.37	1.42
2	B	600	FAD	C2B-C1B	-2.54	1.49	1.53
2	B	600	FAD	P-O5'	2.46	1.70	1.59
2	A	600	FAD	C2A-N3A	2.42	1.37	1.32
2	A	600	FAD	PA-O3P	2.40	1.64	1.59
2	A	600	FAD	C2-N3	2.40	1.42	1.37
2	A	600	FAD	C4X-N5	2.35	1.41	1.36
3	A	601	MLG	C02-CL07	-2.21	1.68	1.73
2	B	600	FAD	C2'-C3'	2.13	1.57	1.53
3	A	601	MLG	C14-C15	-2.06	1.44	1.47
2	A	600	FAD	O3'-C3'	2.03	1.47	1.43
2	B	600	FAD	C10-N10	-2.02	1.34	1.38
2	A	600	FAD	C2A-N1A	2.00	1.37	1.33

All (65) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	601	MLG	C14-C15-C16	-55.01	111.93	177.80
3	B	601	MLG	C14-C15-C16	-51.38	116.28	177.80
2	A	600	FAD	N3A-C2A-N1A	-10.59	119.86	128.71
3	B	601	MLG	C17-N13-C14	10.04	124.77	111.50
2	B	600	FAD	N3A-C2A-N1A	-8.97	121.21	128.71
3	A	601	MLG	C17-N13-C14	8.34	122.52	111.50
2	A	600	FAD	C2'-C1'-N10	7.36	122.22	112.45
3	B	601	MLG	C15-C14-N13	6.39	128.93	113.53
2	B	600	FAD	C4X-C10-N10	-6.26	117.39	120.51
2	A	600	FAD	O5'-P-O1P	5.87	132.35	109.37
2	A	600	FAD	C4X-C10-N1	5.07	127.79	122.73
2	B	600	FAD	C4X-N5-C5X	4.77	122.05	116.69
2	A	600	FAD	C9A-N10-C10	-4.72	117.13	121.77
2	A	600	FAD	P-O3P-PA	-4.68	117.95	131.68
2	B	600	FAD	C5X-C9A-N10	-4.58	112.30	116.80
2	B	600	FAD	N3A-C4A-N9A	4.54	133.62	125.43
2	B	600	FAD	O5'-P-O1P	4.34	126.36	109.37
2	B	600	FAD	P-O3P-PA	-4.17	119.46	131.68
2	A	600	FAD	O2'-C2'-C1'	-3.70	100.52	109.71
3	A	601	MLG	C17-N13-C12	3.62	121.54	110.51
2	A	600	FAD	C9-C8-C7	-3.57	114.16	119.88
3	B	601	MLG	C17-N13-C12	3.50	121.19	110.51
2	B	600	FAD	O2'-C2'-C3'	-3.49	100.36	109.05
2	B	600	FAD	C1'-N10-C9A	-3.48	115.49	118.87
3	B	601	MLG	C05-C04-CL08	3.38	124.91	119.34
3	A	601	MLG	C15-C14-N13	3.37	121.66	113.53
2	A	600	FAD	C10-C4X-N5	-3.31	116.44	120.45
2	A	600	FAD	C6-C5X-N5	3.29	122.81	118.97
2	B	600	FAD	C9-C9A-N10	3.27	128.64	121.59
2	A	600	FAD	C6-C5X-C9A	-3.25	114.51	119.02
2	A	600	FAD	N3A-C4A-N9A	3.18	131.18	125.43
2	A	600	FAD	N6A-C6A-N1A	-3.07	113.35	119.36
3	A	601	MLG	O09-C01-C02	3.02	120.37	116.33
2	B	600	FAD	C4'-C3'-C2'	2.84	119.67	113.25
2	A	600	FAD	C1'-N10-C9A	2.79	121.59	118.87
3	B	601	MLG	C02-C03-C04	2.74	121.83	118.67
2	B	600	FAD	C2-N1-C10	2.71	117.71	114.98
2	B	600	FAD	C9A-C5X-N5	-2.70	118.23	122.37
2	B	600	FAD	C6-C5X-N5	2.69	122.11	118.97
2	B	600	FAD	O3'-C3'-C4'	-2.56	102.27	108.74
3	B	601	MLG	C05-C04-C03	-2.55	118.01	121.54
2	B	600	FAD	C5A-C4A-N3A	-2.55	120.14	125.70
2	A	600	FAD	O2P-P-O1P	-2.52	98.15	112.21

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	600	FAD	O4B-C1B-C2B	-2.52	102.92	106.77
2	B	600	FAD	C4X-C10-N1	2.51	125.23	122.73
2	B	600	FAD	O2B-C2B-C1B	-2.47	103.76	111.23
2	A	600	FAD	C5A-C4A-N3A	-2.47	120.33	125.70
2	B	600	FAD	C10-C4X-N5	-2.45	117.48	120.45
2	A	600	FAD	C4-C4X-N5	2.44	128.63	117.77
2	A	600	FAD	O3'-C3'-C2'	-2.42	102.61	108.74
2	B	600	FAD	O4B-C1B-N9A	2.38	110.65	108.44
2	A	600	FAD	C7-C6-C5X	2.36	124.86	120.91
2	A	600	FAD	C4X-C4-N3	2.24	121.94	115.39
2	A	600	FAD	C4A-C5A-N7A	-2.21	107.63	109.52
2	A	600	FAD	C9A-C5X-N5	-2.19	119.01	122.37
2	A	600	FAD	C4X-N5-C5X	2.17	119.12	116.69
2	B	600	FAD	O4'-C4'-C5'	-2.15	105.71	110.12
3	A	601	MLG	C11-C12-N13	-2.14	107.17	113.93
2	B	600	FAD	C4-C4X-N5	2.14	127.30	117.77
3	A	601	MLG	C05-C04-CL08	2.13	122.86	119.34
2	A	600	FAD	C2A-N3A-C4A	2.10	120.00	114.01
2	A	600	FAD	O2A-PA-O3P	2.07	114.97	105.14
2	A	600	FAD	C8-C9-C9A	2.06	123.97	119.81
2	B	600	FAD	N3-C2-N1	-2.04	116.86	121.19
3	B	601	MLG	O09-C01-C02	2.02	119.03	116.33

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, 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	445/527 (84%)	-0.05	4 (0%) 81 24	62, 64, 67, 70	0
1	B	445/527 (84%)	-0.05	2 (0%) 90 41	62, 64, 67, 70	0
All	All	890/1054 (84%)	-0.05	6 (0%) 84 28	62, 64, 67, 70	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	147	GLN	2.3
1	A	116	TRP	2.2
1	A	464	GLY	2.2
1	B	96	ARG	2.2
1	A	215	GLN	2.1
1	B	147	GLN	2.1

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, 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	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	MLG	B	601	17/17	0.25	0.49	72,79,84,84	0
2	FAD	B	600	53/53	0.23	0.29	56,63,67,70	0
2	FAD	A	600	53/53	0.23	0.12	61,65,69,72	0
3	MLG	A	601	17/17	0.19	-0.58	73,79,84,85	0

6.5 Other polymers ⓘ

There are no such residues in this entry.