



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 12:50 AM GMT

PDB ID : 2BXR
Title : HUMAN MONOAMINE OXIDASE A IN COMPLEX WITH CLORGYLINE,
CRYSTAL FORM A
Authors : De Colibus, L.; Binda, C.; Edmondson, D.E.; Mattevi, A.
Deposited on : 2005-07-27
Resolution : 3.00 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

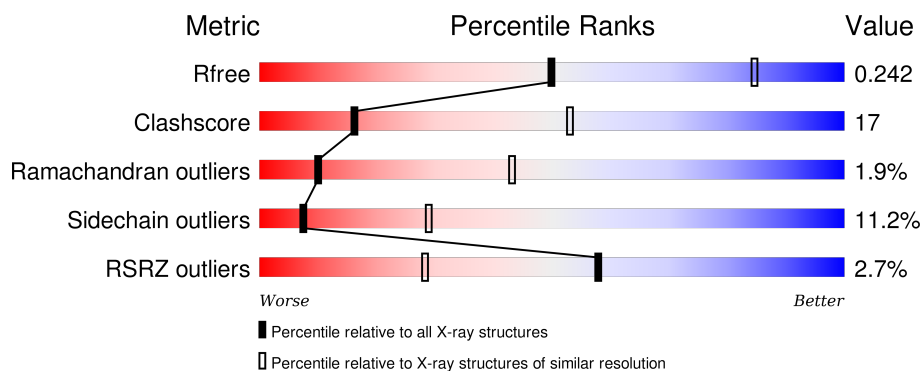
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

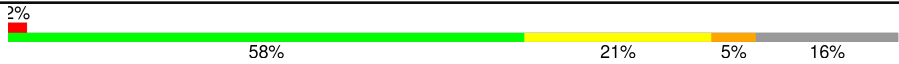

The reported resolution of this entry is 3.00 Å.

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}	91344	1578 (3.00-3.00)
Clashscore	102246	1912 (3.00-3.00)
Ramachandran outliers	100387	1853 (3.00-3.00)
Sidechain outliers	100360	1856 (3.00-3.00)
RSRZ outliers	91569	1592 (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. 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	527	
1	B	527	

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	MLG	A	601	X	-	X	-
3	MLG	B	601	X	-	-	-

2 Entry composition [i](#)

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

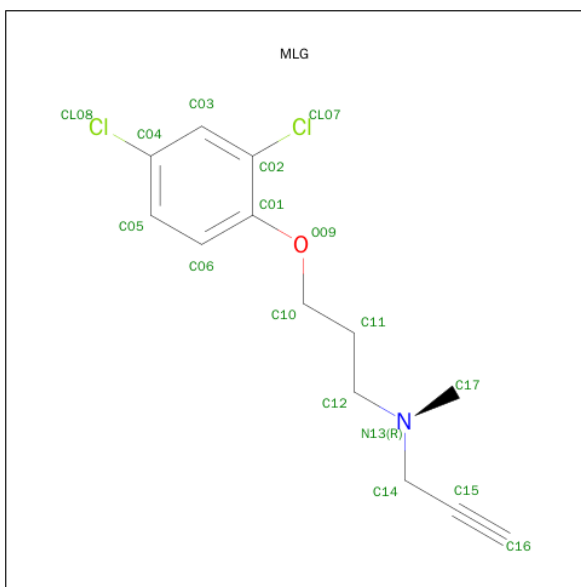
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	445	Total	C	N	O	S	0	0	0
			3523	2253	601	648	21			
1	B	445	Total	C	N	O	S	0	0	0
			3523	2253	601	648	21			

- 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: $C_{13}H_{15}Cl_2NO$).

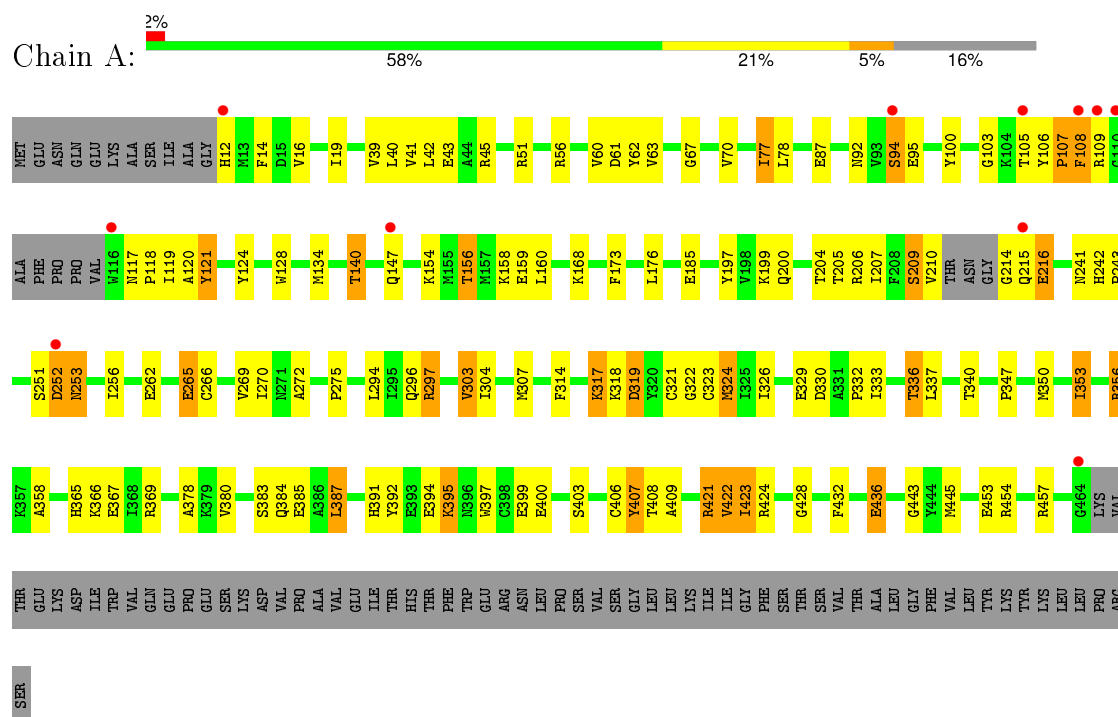


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	Cl	N	O	0	0
			17	13	2	1	1		
3	B	1	Total	C	Cl	N	O	0	0
			17	13	2	1	1		

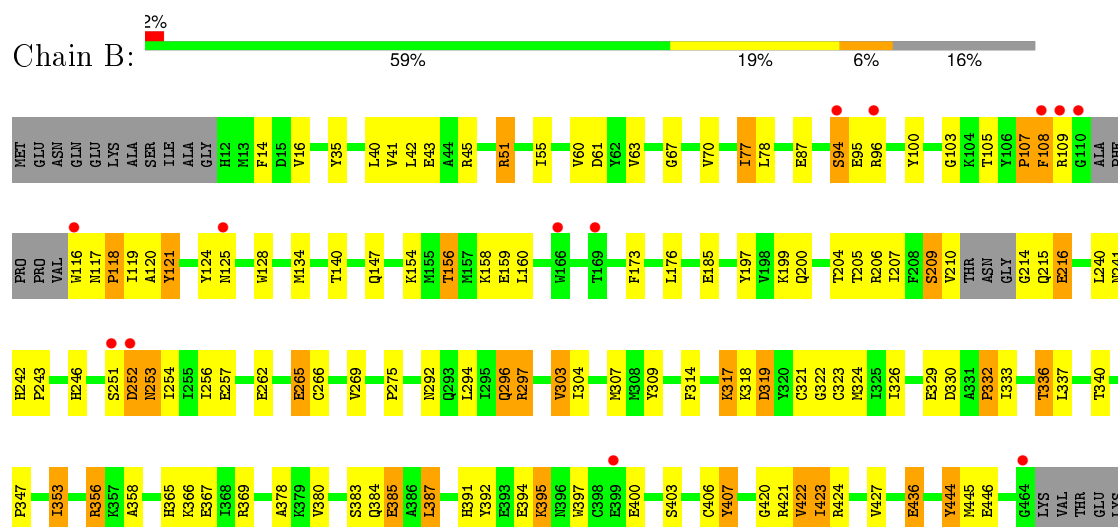
3 Residue-property plots [i](#)

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: AMINE OXIDASE [FLAVIN-CONTAINING] A



• Molecule 1: AMINE OXIDASE [FLAVIN-CONTAINING] A



ASP	ILE	TRP	VAL	GLN	GLU	PRO	GLU	SER	LYS	ASP	VAL	PRO	ALA	VAL	GLU	ILE	THR	HIS	THR	PHE	TRP	GLU	ARG	ASN	LEU	PRO	SER	SER	VAL	SER	GLY	LEU	LEU	LYS	ILE	ILE	GLY	PHE	SER	THR	SER	SER	VAL	THR	ALA	LEU	GLY	PHE	VAL	LEU	TYR	LYS	TYR	LYS	LYS	LEU	LEU	PRO	ARG	SER
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4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	143.49 Å 109.60 Å 81.33 Å 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 3.01 Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.192 , 0.238 0.194 , 0.242	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)	0.31 , 64.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 24331 reflections	Xtriage
F_o, F_c correlation	0.94	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.

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

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 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	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

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 17.

All (236) 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)
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	Interatomic distance (Å)	Clash overlap (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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:324:MET:H	1:B:336:THR:HG22	1.64	0.63
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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 [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	439/527 (83%)	392 (89%)	38 (9%)	9 (2%)	9	40
1	B	439/527 (83%)	391 (89%)	40 (9%)	8 (2%)	11	45
All	All	878/1054 (83%)	783 (89%)	78 (9%)	17 (2%)	10	43

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

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Mol	Chain	Res	Type
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
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%)	8	29
1	B	371/450 (82%)	329 (89%)	42 (11%)	7	28
All	All	742/900 (82%)	659 (89%)	83 (11%)	7	29

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

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Mol	Chain	Res	Type
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
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

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Mol	Chain	Res	Type
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
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 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 ⓘ

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	48,58,58	2.89	15 (31%)	54,89,89	3.15	19 (35%)
3	MLG	A	601	2	16,17,17	3.80	3 (18%)	20,21,21	11.49	6 (30%)
2	FAD	B	600	1,3	48,58,58	2.96	14 (29%)	54,89,89	3.15	16 (29%)
3	MLG	B	601	2	16,17,17	3.06	2 (12%)	20,21,21	10.80	7 (35%)

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	-	0/30/50/50	0/6/6/6
3	MLG	A	601	2	1/1/1/1	0/9/10/10	0/1/1/1
2	FAD	B	600	1,3	-	0/30/50/50	0/6/6/6
3	MLG	B	601	2	1/1/1/1	0/9/10/10	0/1/1/1

All (34) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	600	FAD	C4-C4X	-7.91	1.25	1.41
2	B	600	FAD	C4-C4X	-7.42	1.26	1.41
2	A	600	FAD	C9A-N10	-4.59	1.32	1.38
3	B	601	MLG	C14-C15	-4.55	1.41	1.47
2	B	600	FAD	C10-N10	-4.06	1.34	1.39
2	A	600	FAD	C9A-C5X	-2.75	1.37	1.42
3	A	601	MLG	C02-CL07	-2.31	1.68	1.73
3	A	601	MLG	C14-C15	-2.28	1.44	1.47
2	A	600	FAD	O3'-C3'	2.05	1.47	1.43
2	A	600	FAD	C2A-N1A	2.09	1.37	1.33
2	B	600	FAD	C2'-C3'	2.19	1.57	1.53
2	B	600	FAD	P-O5'	2.47	1.70	1.59
2	A	600	FAD	C2A-N3A	2.69	1.37	1.32
2	A	600	FAD	C6-C7	2.73	1.45	1.37
2	B	600	FAD	C6-C7	2.84	1.45	1.37
2	B	600	FAD	C2A-N1A	3.01	1.39	1.33
2	A	600	FAD	C5X-N5	3.14	1.40	1.35
2	B	600	FAD	O2'-C2'	3.32	1.50	1.43
2	A	600	FAD	C1'-N10	3.77	1.52	1.48
2	B	600	FAD	C5X-N5	3.77	1.41	1.35
2	B	600	FAD	C9-C9A	3.79	1.49	1.40
2	A	600	FAD	O4-C4	4.14	1.34	1.24
2	B	600	FAD	C8M-C8	4.17	1.59	1.51
2	B	600	FAD	C2A-N3A	4.40	1.40	1.32
2	A	600	FAD	C9-C9A	4.73	1.51	1.40
2	A	600	FAD	C4-N3	4.74	1.41	1.33
2	A	600	FAD	C4X-N5	4.93	1.41	1.33
2	A	600	FAD	C10-N1	6.35	1.46	1.35
2	B	600	FAD	C4-N3	7.82	1.47	1.33
2	B	600	FAD	C1'-N10	7.88	1.56	1.48
2	B	600	FAD	C10-N1	9.66	1.51	1.35
2	A	600	FAD	C8M-C8	10.66	1.72	1.51
3	B	601	MLG	C15-C16	10.86	1.43	1.17
3	A	601	MLG	C15-C16	14.65	1.51	1.17

All (48) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	601	MLG	C14-C15-C16	-50.73	111.93	177.76
3	B	601	MLG	C14-C15-C16	-47.38	116.28	177.76
2	A	600	FAD	N3A-C2A-N1A	-11.81	119.86	128.89
2	B	600	FAD	N3A-C2A-N1A	-10.04	121.21	128.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	600	FAD	C4-C4X-C10	-7.99	114.83	119.94
2	B	600	FAD	C4-C4X-C10	-7.38	115.22	119.94
2	B	600	FAD	C5X-C9A-N10	-7.00	112.30	117.62
2	B	600	FAD	C4X-C10-N10	-5.32	117.39	120.52
2	A	600	FAD	P-O3P-PA	-5.26	117.95	132.73
2	B	600	FAD	P-O3P-PA	-4.73	119.46	132.73
2	A	600	FAD	O2'-C2'-C1'	-3.84	100.52	109.94
2	B	600	FAD	O2'-C2'-C3'	-3.44	100.36	109.02
2	A	600	FAD	C6-C5X-C9A	-3.40	114.51	118.98
2	A	600	FAD	C9-C8-C7	-3.08	114.16	120.04
2	B	600	FAD	C1'-N10-C9A	-3.00	115.49	118.86
2	B	600	FAD	C9A-C5X-N5	-2.79	118.23	122.36
2	A	600	FAD	N6A-C6A-N1A	-2.73	113.35	119.20
2	A	600	FAD	O2P-P-O1P	-2.65	98.15	112.53
3	B	601	MLG	C05-C04-C03	-2.63	118.01	121.53
2	B	600	FAD	O3'-C3'-C4'	-2.57	102.27	108.75
2	A	600	FAD	O3'-C3'-C2'	-2.44	102.61	108.75
2	A	600	FAD	C9A-C5X-N5	-2.26	119.01	122.36
3	A	601	MLG	C11-C12-N13	-2.09	107.17	113.90
2	B	600	FAD	O4'-C4'-C5'	-2.06	105.71	110.19
2	B	600	FAD	C4X-C4-N3	-2.02	120.82	123.59
2	A	600	FAD	C4A-C5A-N7A	-2.01	107.63	109.48
3	B	601	MLG	O09-C01-C02	2.04	119.03	116.36
2	A	600	FAD	C4X-N5-C5X	2.05	119.12	116.76
3	A	601	MLG	C05-C04-CL08	2.14	122.86	119.35
2	A	600	FAD	O2A-PA-O3P	2.18	114.97	105.09
2	A	600	FAD	C7-C6-C5X	2.41	124.86	120.92
2	A	600	FAD	C1'-N10-C9A	2.43	121.59	118.86
2	B	600	FAD	C6-C5X-N5	2.45	122.11	118.96
3	B	601	MLG	C02-C03-C04	2.83	121.83	118.69
2	A	600	FAD	C6-C5X-N5	2.99	122.81	118.96
3	A	601	MLG	O09-C01-C02	3.06	120.37	116.36
3	B	601	MLG	C05-C04-CL08	3.39	124.91	119.35
3	B	601	MLG	C17-N13-C12	3.52	121.19	110.51
3	A	601	MLG	C17-N13-C12	3.64	121.54	110.51
2	B	600	FAD	O5'-P-O1P	4.31	126.36	109.62
3	A	601	MLG	C17-N13-C14	4.41	122.52	110.33
2	B	600	FAD	C4X-N5-C5X	4.60	122.05	116.76
3	B	601	MLG	C17-N13-C14	5.23	124.77	110.33
2	A	600	FAD	O5'-P-O1P	5.86	132.35	109.62
2	B	600	FAD	C4-C4X-N5	7.07	127.30	118.72
2	A	600	FAD	C4-C4X-N5	8.16	128.63	118.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	600	FAD	C4-N3-C2	8.75	122.81	115.25
2	B	600	FAD	C4-N3-C2	10.21	124.07	115.25

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	A	601	MLG	N13
3	B	601	MLG	N13

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 25 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	600	FAD	17	0
3	A	601	MLG	9	0
2	B	600	FAD	7	0
3	B	601	MLG	1	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

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%)	-0.34	11 (2%) 61 30	62, 64, 67, 70	0
1	B	445/527 (84%)	-0.28	13 (2%) 55 26	62, 64, 67, 70	0
All	All	890/1054 (84%)	-0.31	24 (2%) 58 28	62, 64, 67, 70	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	94	SER	5.7
1	B	464	GLY	5.3
1	A	116	TRP	4.6
1	A	464	GLY	4.2
1	A	110	GLY	3.8
1	A	252	ASP	3.7
1	B	96	ARG	3.3
1	B	116	TRP	3.2
1	A	108	PHE	3.1
1	B	110	GLY	3.0
1	B	252	ASP	3.0
1	B	125	ASN	2.8
1	A	12	HIS	2.7
1	B	108	PHE	2.7
1	B	109	ARG	2.6
1	B	166	TRP	2.6
1	A	109	ARG	2.5
1	A	105	THR	2.5
1	A	147	GLN	2.4
1	A	215	GLN	2.4
1	A	94	SER	2.3
1	B	251	SER	2.2
1	B	399	GLU	2.2
1	B	169	THR	2.1

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(\AA^2)	Q<0.9
3	MLG	A	601	17/17	0.93	0.23	0.86	73,79,84,85	0
3	MLG	B	601	17/17	0.92	0.24	0.76	72,79,84,84	0
2	FAD	B	600	53/53	0.95	0.14	-0.21	56,63,67,70	0
2	FAD	A	600	53/53	0.96	0.14	-0.57	61,65,69,72	0

6.5 Other polymers [i](#)

There are no such residues in this entry.