



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 27, 2014 – 05:35 AM GMT

PDB ID : 1VQW
Title : Crystal structure of a protein with similarity to flavin-containing monooxygenases and to mammalian dimethylalanine monooxygenases
Authors : Eswaramoorthy, S.; Swaminathan, S.; Burley, S.K.; New York SGX Research Center for Structural Genomics (NYSGXRC)
Deposited on : 2005-01-05
Resolution : 2.40 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>

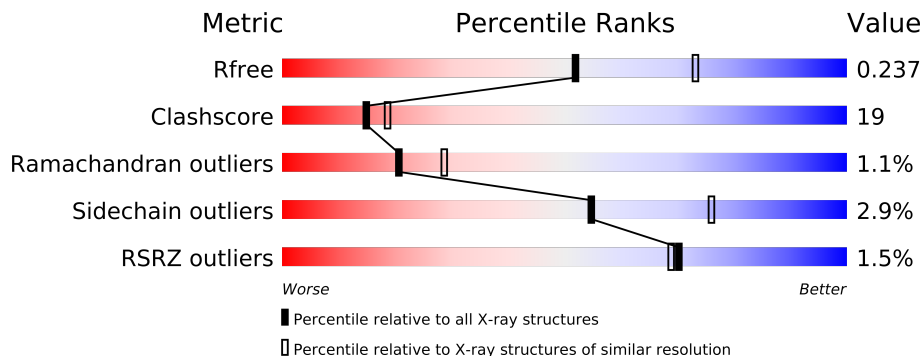
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.40 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	2207 (2.40-2.40)
Clashscore	79885	2789 (2.40-2.40)
Ramachandran outliers	78287	2736 (2.40-2.40)
Sidechain outliers	78261	2737 (2.40-2.40)
RSRZ outliers	66119	2210 (2.40-2.40)

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	457	
1	B	457	

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

Mol	Type	Chain	Res	Geometry	Electron density
2	FAD	A	500	-	X
2	FAD	B	501	-	X
3	EPE	A	502	-	X
3	EPE	A	503	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7492 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 PROTEIN WITH SIMILARITY TO FLAVIN-CONTAINING MONOOXYGENASES AND TO MAMMALIAN DIMETHYLALANINE MONOOXYGENASES.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	442	Total	C	N	O	S	Se	0	0	0
			3490	2261	578	643	4	4			
1	B	442	Total	C	N	O	S	Se	0	0	0
			3490	2261	578	643	4	4			

There are 30 discrepancies between the modelled and reference sequences:

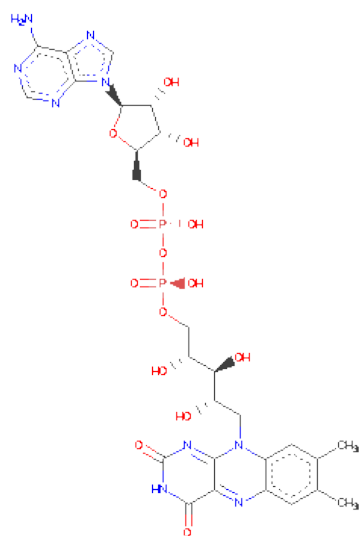
Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MSE	-	CLONING ARTIFACT	UNP Q9HFE4
A	0	SER	-	CLONING ARTIFACT	UNP Q9HFE4
A	1	LEU	-	CLONING ARTIFACT	UNP Q9HFE4
A	97	MSE	MET	MODIFIED RESIDUE	UNP Q9HFE4
A	377	MSE	MET	MODIFIED RESIDUE	UNP Q9HFE4
A	386	MSE	MET	MODIFIED RESIDUE	UNP Q9HFE4
A	433	MSE	MET	MODIFIED RESIDUE	UNP Q9HFE4
A	448	GLU	-	EXPRESSION TAG	UNP Q9HFE4
A	449	GLY	-	EXPRESSION TAG	UNP Q9HFE4
A	450	HIS	-	EXPRESSION TAG	UNP Q9HFE4
A	451	HIS	-	EXPRESSION TAG	UNP Q9HFE4
A	452	HIS	-	EXPRESSION TAG	UNP Q9HFE4
A	453	HIS	-	EXPRESSION TAG	UNP Q9HFE4
A	454	HIS	-	EXPRESSION TAG	UNP Q9HFE4
A	455	HIS	-	EXPRESSION TAG	UNP Q9HFE4
B	-1	MSE	-	CLONING ARTIFACT	UNP Q9HFE4
B	0	SER	-	CLONING ARTIFACT	UNP Q9HFE4
B	1	LEU	-	CLONING ARTIFACT	UNP Q9HFE4
B	97	MSE	MET	MODIFIED RESIDUE	UNP Q9HFE4
B	377	MSE	MET	MODIFIED RESIDUE	UNP Q9HFE4
B	386	MSE	MET	MODIFIED RESIDUE	UNP Q9HFE4
B	433	MSE	MET	MODIFIED RESIDUE	UNP Q9HFE4
B	448	GLU	-	EXPRESSION TAG	UNP Q9HFE4

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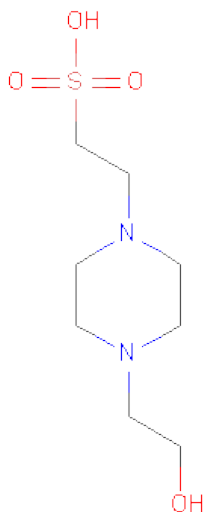
Chain	Residue	Modelled	Actual	Comment	Reference
B	449	GLY	-	EXPRESSION TAG	UNP Q9HFE4
B	450	HIS	-	EXPRESSION TAG	UNP Q9HFE4
B	451	HIS	-	EXPRESSION TAG	UNP Q9HFE4
B	452	HIS	-	EXPRESSION TAG	UNP Q9HFE4
B	453	HIS	-	EXPRESSION TAG	UNP Q9HFE4
B	454	HIS	-	EXPRESSION TAG	UNP Q9HFE4
B	455	HIS	-	EXPRESSION TAG	UNP Q9HFE4

- 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 4-(2-HYDROXYETHYL)-1-PIPERAZINEETHANESULFONIC ACID (three-letter code: EPE) (formula: $C_8H_{18}N_2O_4S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			12	6	2	3	1		
3	A	1	Total	C	N	O	S	0	0
			12	6	2	3	1		

- Molecule 4 is water.

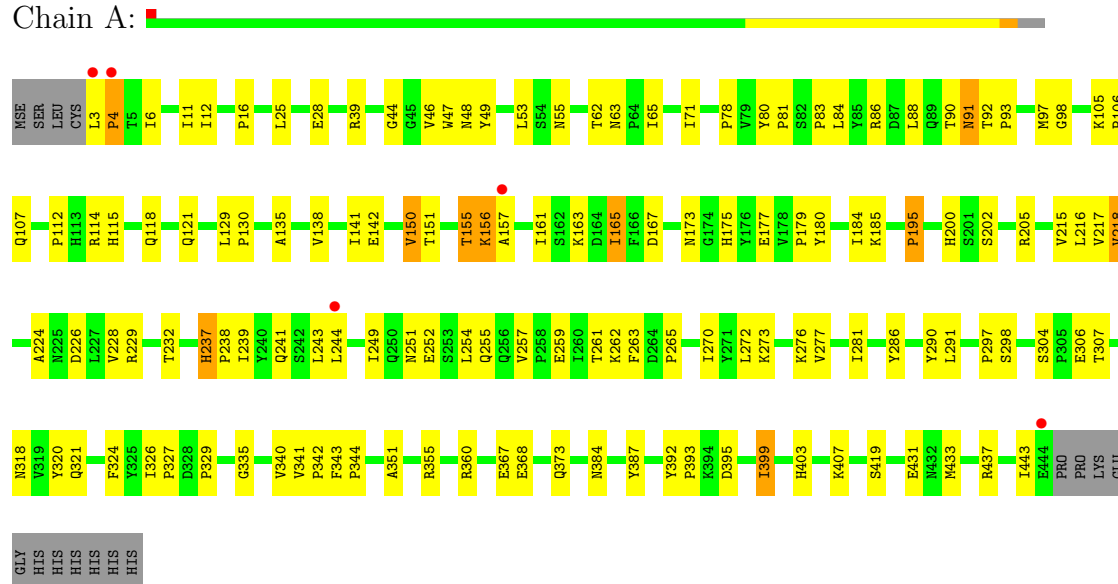
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	200	Total	O	0	0
			200	200		
4	B	182	Total	O	0	0
			182	182		

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

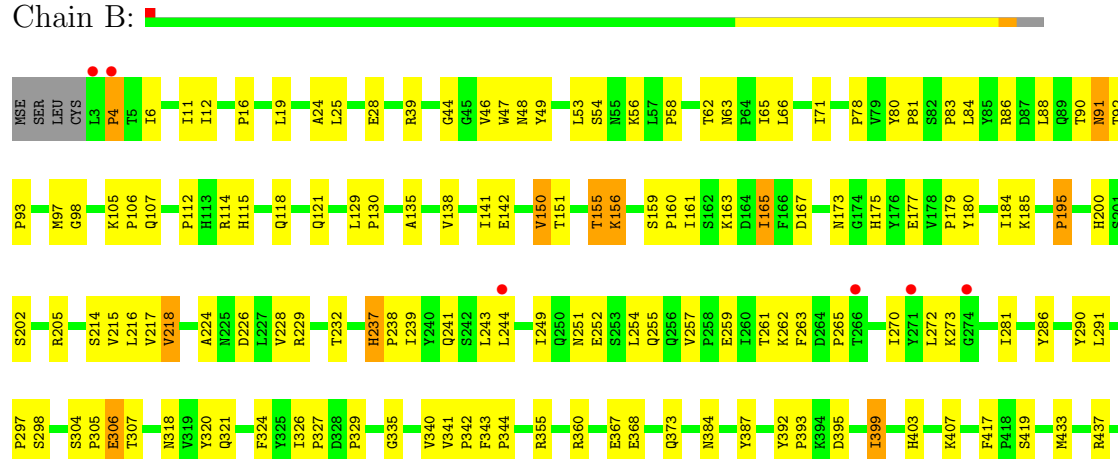
- Molecule 1: PROTEIN WITH SIMILARITY TO FLAVIN-CONTAINING MONOOXYGENASES AND TO MAMMALIAN DIMETHYLALANINE MONOOXYGENASES

Chain A:



- Molecule 1: PROTEIN WITH SIMILARITY TO FLAVIN-CONTAINING MONOOXYGENASES AND TO MAMMALIAN DIMETHYLALANINE MONOOXYGENASES

Chain B:



1443	E449
PRO	
LYS	
GLU	
GLY	
HIS	
HIS	
HIS	
HIS	
HIS	

4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	59.59Å 72.64Å 80.35Å 99.00° 107.09° 102.02°	Depositor
Resolution (Å)	50.00 – 2.40 44.53 – 2.40	Depositor EDS
% Data completeness (in resolution range)	92.3 (50.00-2.40) 92.0 (44.53-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.35 (at 2.39Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.226 , 0.239 0.220 , 0.237	Depositor DCC
R_{free} test set	1335 reflections (3.12%)	DCC
Wilson B-factor (Å ²)	29.5	Xtriage
Anisotropy	0.394	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 22.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 47252 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7492	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.13% 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: EPE, 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	0.37	0/3587	0.64	1/4885 (0.0%)
1	B	0.38	0/3587	0.64	1/4885 (0.0%)
All	All	0.37	0/7174	0.64	2/9770 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	306	GLU	OE1-CD-OE2	6.20	130.74	123.30
1	B	306	GLU	OE1-CD-OE2	5.90	130.38	123.30

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	3490	0	3487	133	0
1	B	3490	0	3487	135	0
2	A	53	0	31	9	0
2	B	53	0	31	8	0
3	A	24	0	24	29	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	200	0	0	14	0
4	B	182	0	0	16	0
All	All	7492	0	7060	270	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 19.

All (270) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:39:ARG:HG2	2:A:500:FAD:N3A	1.78	0.98
1:A:244:LEU:HD12	1:A:244:LEU:H	1.27	0.98
1:B:39:ARG:HG2	2:B:501:FAD:N3A	1.77	0.97
1:B:244:LEU:HD12	1:B:244:LEU:H	1.26	0.96
1:B:114:ARG:HH11	1:B:118:GLN:HE22	1.11	0.95
1:A:114:ARG:HH11	1:A:118:GLN:HE22	1.10	0.94
1:A:48:ASN:HB3	3:A:503:EPE:H102	1.55	0.88
1:A:83:PRO:HG2	2:A:500:FAD:HM82	1.56	0.87
1:B:83:PRO:HG2	2:B:501:FAD:HM82	1.57	0.86
3:A:502:EPE:H102	1:B:48:ASN:HB3	1.58	0.84
1:A:157:ALA:HA	4:A:631:HOH:O	1.75	0.84
3:A:502:EPE:H51	4:B:515:HOH:O	1.79	0.83
1:B:159:SER:HB3	4:B:680:HOH:O	1.80	0.82
1:B:142:GLU:HB3	1:B:298:SER:CB	2.10	0.82
1:A:142:GLU:HB3	1:A:298:SER:CB	2.10	0.82
1:B:155:THR:HG21	4:B:587:HOH:O	1.81	0.80
1:A:218:VAL:HG13	1:A:286:TYR:HA	1.66	0.78
1:A:93:PRO:HD2	1:A:433:MSE:HE2	1.67	0.77
1:B:93:PRO:HD2	1:B:433:MSE:HE2	1.65	0.76
3:A:503:EPE:H51	4:A:509:HOH:O	1.86	0.75
1:B:218:VAL:HG13	1:B:286:TYR:HA	1.68	0.75
1:B:47:TRP:O	1:B:118:GLN:HG2	1.87	0.74
1:B:244:LEU:CD1	1:B:244:LEU:H	2.00	0.74
1:B:155:THR:O	1:B:156:LYS:HB2	1.88	0.73
1:A:244:LEU:H	1:A:244:LEU:CD1	2.00	0.73
1:A:155:THR:O	1:A:156:LYS:HB2	1.88	0.73
1:A:395:ASP:O	1:A:399:ILE:HG23	1.87	0.73
1:A:47:TRP:O	1:A:118:GLN:HG2	1.89	0.71
1:A:91:ASN:HD22	1:A:91:ASN:C	1.91	0.71
1:B:395:ASP:O	1:B:399:ILE:HG23	1.90	0.71
1:B:91:ASN:C	1:B:91:ASN:HD22	1.92	0.71
1:A:39:ARG:HD2	4:A:669:HOH:O	1.91	0.70
1:A:244:LEU:HD12	1:A:244:LEU:N	2.06	0.70

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:318:ASN:H	1:B:373:GLN:HE22	1.39	0.69
1:B:367:GLU:HB3	4:B:642:HOH:O	1.91	0.69
1:A:91:ASN:ND2	1:A:92:THR:HG23	2.08	0.69
1:A:318:ASN:H	1:A:373:GLN:HE22	1.40	0.68
3:A:503:EPE:H61	1:B:114:ARG:HH22	1.59	0.67
1:B:91:ASN:ND2	1:B:92:THR:HG23	2.09	0.67
1:B:244:LEU:HD12	1:B:244:LEU:N	2.06	0.67
1:A:114:ARG:HH22	3:A:502:EPE:H61	1.61	0.66
1:A:175:HIS:HE1	1:A:335:GLY:O	1.77	0.66
1:B:175:HIS:HE1	1:B:335:GLY:O	1.78	0.66
1:A:48:ASN:HD22	3:A:503:EPE:H91	1.60	0.65
1:B:155:THR:O	1:B:156:LYS:CB	2.44	0.65
3:A:502:EPE:H91	1:B:48:ASN:HD22	1.61	0.65
1:A:261:THR:HG22	1:A:262:LYS:HG3	1.79	0.64
1:A:276:LYS:HB3	4:A:597:HOH:O	1.97	0.64
1:A:105:LYS:HE3	4:B:609:HOH:O	1.98	0.64
1:A:342:PRO:HD2	2:A:500:FAD:O2	1.98	0.64
1:A:155:THR:O	1:A:156:LYS:CB	2.45	0.63
1:B:263:PHE:O	1:B:265:PRO:HD3	1.97	0.63
1:A:263:PHE:O	1:A:265:PRO:HD3	1.99	0.63
1:B:261:THR:HG22	1:B:262:LYS:HG3	1.80	0.63
1:B:106:PRO:O	1:B:107:GLN:HB2	1.97	0.63
1:B:355:ARG:NH1	1:B:360:ARG:HH12	1.97	0.62
1:A:224:ALA:O	1:A:228:VAL:HG23	1.99	0.62
1:A:355:ARG:NH1	1:A:360:ARG:HH12	1.97	0.62
1:B:142:GLU:HB3	1:B:298:SER:HB3	1.82	0.62
1:A:63:ASN:OD1	1:A:65:ILE:HG22	1.99	0.62
1:B:86:ARG:HA	1:B:114:ARG:HH21	1.65	0.62
1:B:342:PRO:HD2	2:B:501:FAD:O2	2.00	0.61
1:A:106:PRO:O	1:A:107:GLN:HB2	1.99	0.61
1:B:217:VAL:O	1:B:241:GLN:HA	2.01	0.60
1:A:142:GLU:HB3	1:A:298:SER:HB3	1.82	0.60
1:B:63:ASN:OD1	1:B:65:ILE:HG22	2.02	0.60
1:A:217:VAL:O	1:A:241:GLN:HA	2.02	0.60
1:B:224:ALA:O	1:B:228:VAL:HG23	2.01	0.59
3:A:503:EPE:H61	1:B:114:ARG:NH2	2.17	0.59
1:A:304:SER:HB3	1:A:307:THR:OG1	2.02	0.59
1:B:392:TYR:CD1	1:B:393:PRO:HA	2.37	0.59
1:A:62:THR:HG23	4:A:694:HOH:O	2.03	0.59
1:B:238:PRO:HB3	1:B:252:GLU:O	2.03	0.59
1:A:49:TYR:CE2	3:A:502:EPE:H22	2.38	0.59
1:A:155:THR:HG21	4:A:695:HOH:O	2.02	0.58

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:114:ARG:NH2	3:A:502:EPE:H61	2.18	0.58
1:A:392:TYR:CD1	1:A:393:PRO:HA	2.38	0.58
1:B:304:SER:HB3	1:B:307:THR:OG1	2.04	0.58
1:A:86:ARG:HA	1:A:114:ARG:HH21	1.69	0.58
3:A:503:EPE:H22	1:B:49:TYR:CE2	2.38	0.58
3:A:503:EPE:C6	1:B:114:ARG:HH22	2.17	0.58
1:A:55:ASN:HB3	4:A:565:HOH:O	2.04	0.58
1:A:114:ARG:HH22	3:A:502:EPE:C5	2.17	0.57
1:B:184:ILE:HG22	1:B:185:LYS:N	2.19	0.57
1:B:54:SER:HB3	4:B:536:HOH:O	2.04	0.57
1:A:185:LYS:HG2	1:A:262:LYS:HZ3	1.69	0.57
1:B:259:GLU:OE2	1:B:273:LYS:HD2	2.05	0.57
3:A:503:EPE:C5	1:B:114:ARG:HH22	2.18	0.56
1:A:114:ARG:HH22	3:A:502:EPE:C6	2.17	0.56
1:B:343:PHE:HB2	1:B:344:PRO:HD3	1.87	0.56
1:A:91:ASN:HD22	1:A:92:THR:HG23	1.68	0.56
1:A:184:ILE:HG22	1:A:185:LYS:N	2.20	0.56
1:A:81:PRO:HG2	4:A:607:HOH:O	2.05	0.55
1:B:161:ILE:N	1:B:161:ILE:HD12	2.22	0.55
1:A:238:PRO:HB3	1:A:252:GLU:O	2.06	0.55
1:A:259:GLU:OE2	1:A:273:LYS:HD2	2.06	0.55
1:A:226:ASP:OD1	1:A:229:ARG:NH2	2.41	0.54
1:B:91:ASN:HD22	1:B:92:THR:HG23	1.71	0.54
1:A:343:PHE:HB2	1:A:344:PRO:HD3	1.88	0.54
1:B:184:ILE:HG22	1:B:185:LYS:H	1.71	0.54
1:B:226:ASP:OD1	1:B:229:ARG:NH2	2.40	0.54
1:B:114:ARG:HB2	1:B:118:GLN:HE21	1.73	0.54
1:A:161:ILE:HD12	1:A:161:ILE:N	2.22	0.54
1:A:185:LYS:HB3	1:A:262:LYS:HG2	1.90	0.54
1:B:28:GLU:OE1	1:B:360:ARG:HD3	2.08	0.54
1:B:341:VAL:O	1:B:344:PRO:HD2	2.08	0.54
1:B:185:LYS:HG2	1:B:262:LYS:HZ3	1.73	0.53
1:A:114:ARG:HB2	1:A:118:GLN:HE21	1.72	0.53
1:A:3:LEU:HD23	4:A:602:HOH:O	2.07	0.53
1:B:399:ILE:HD12	1:B:399:ILE:C	2.29	0.53
1:A:106:PRO:HG2	1:B:106:PRO:HG2	1.90	0.53
1:A:341:VAL:O	1:A:344:PRO:HD2	2.09	0.53
1:A:11:ILE:HD11	1:A:25:LEU:HD12	1.91	0.53
1:B:185:LYS:HB3	1:B:262:LYS:HG2	1.90	0.52
1:A:28:GLU:OE1	1:A:360:ARG:HD3	2.08	0.52
1:A:78:PRO:HG2	1:A:155:THR:HG23	1.91	0.52
1:B:11:ILE:HD11	1:B:25:LEU:HD12	1.91	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:83:PRO:CG	2:A:500:FAD:HM82	2.34	0.52
1:A:141:ILE:O	1:A:298:SER:HB2	2.10	0.52
1:A:399:ILE:HD12	1:A:399:ILE:C	2.30	0.52
1:B:78:PRO:HG2	1:B:155:THR:HG23	1.92	0.52
1:A:249:ILE:HG21	1:A:254:LEU:HD23	1.92	0.52
1:B:46:VAL:HG22	2:B:501:FAD:O4'	2.10	0.52
1:A:138:VAL:HG13	1:A:150:VAL:HG23	1.91	0.52
1:B:249:ILE:HG21	1:B:254:LEU:HD23	1.91	0.51
1:B:138:VAL:HG13	1:B:150:VAL:HG23	1.92	0.51
1:B:403:HIS:O	1:B:407:LYS:HG2	2.11	0.51
1:A:84:LEU:HD22	1:A:88:LEU:HD13	1.92	0.51
1:B:98:GLY:HA2	1:B:343:PHE:HD2	1.76	0.51
1:A:46:VAL:HG22	2:A:500:FAD:O4'	2.11	0.51
1:A:184:ILE:HG22	1:A:185:LYS:H	1.75	0.51
1:A:175:HIS:CE1	1:A:335:GLY:O	2.62	0.50
1:B:83:PRO:CG	2:B:501:FAD:HM82	2.35	0.50
1:A:200:HIS:HD2	1:A:202:SER:OG	1.94	0.50
1:A:343:PHE:N	1:A:343:PHE:CD1	2.79	0.50
1:A:48:ASN:HD22	3:A:503:EPE:C9	2.23	0.50
1:B:141:ILE:O	1:B:298:SER:HB2	2.10	0.50
1:A:98:GLY:HA2	1:A:343:PHE:HD2	1.75	0.50
1:A:49:TYR:CD2	3:A:502:EPE:H22	2.47	0.49
1:B:318:ASN:N	1:B:373:GLN:HE22	2.10	0.49
1:B:343:PHE:N	1:B:343:PHE:CD1	2.79	0.49
1:B:200:HIS:HD2	1:B:202:SER:OG	1.94	0.49
1:A:403:HIS:O	1:A:407:LYS:HG2	2.12	0.49
1:B:84:LEU:HD22	1:B:88:LEU:HD13	1.94	0.49
3:A:503:EPE:C5	1:B:115:HIS:NE2	2.76	0.49
3:A:503:EPE:H22	1:B:49:TYR:CD2	2.47	0.49
1:B:91:ASN:C	1:B:91:ASN:ND2	2.64	0.49
1:B:88:LEU:C	1:B:88:LEU:HD23	2.32	0.49
1:A:16:PRO:HG2	2:A:500:FAD:H4'	1.94	0.49
1:A:115:HIS:NE2	3:A:502:EPE:C5	2.76	0.48
1:A:165:ILE:H	1:A:165:ILE:HD13	1.78	0.48
1:A:277:VAL:HG12	4:A:563:HOH:O	2.13	0.48
3:A:502:EPE:C9	1:B:48:ASN:HD22	2.24	0.48
1:B:407:LYS:CD	4:B:595:HOH:O	2.60	0.48
1:A:129:LEU:N	1:A:130:PRO:HD2	2.29	0.48
1:A:97:MSE:HE1	1:A:112:PRO:HD2	1.94	0.48
1:B:251:ASN:HB3	4:B:627:HOH:O	2.12	0.48
1:B:175:HIS:CE1	1:B:335:GLY:O	2.63	0.48
1:A:39:ARG:HD3	1:A:80:TYR:CD1	2.49	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:39:ARG:HD3	1:B:80:TYR:CD1	2.49	0.47
1:B:12:ILE:HD12	1:B:138:VAL:HG21	1.95	0.47
1:B:251:ASN:O	1:B:255:GLN:NE2	2.38	0.47
1:A:185:LYS:HG2	1:A:262:LYS:HG2	1.97	0.47
1:A:12:ILE:HD12	1:A:138:VAL:HG21	1.95	0.47
1:A:88:LEU:C	1:A:88:LEU:HD23	2.34	0.47
1:B:185:LYS:HG2	1:B:262:LYS:HG2	1.96	0.47
1:B:129:LEU:N	1:B:130:PRO:HD2	2.30	0.47
1:A:93:PRO:CD	1:A:433:MSE:HE2	2.43	0.47
1:B:6:ILE:HG23	1:B:167:ASP:HB2	1.96	0.47
1:B:93:PRO:CD	1:B:433:MSE:HE2	2.40	0.47
1:B:105:LYS:NZ	4:B:618:HOH:O	2.47	0.47
1:A:71:ILE:HD12	1:A:71:ILE:N	2.29	0.46
1:A:49:TYR:O	3:A:503:EPE:H21	2.15	0.46
1:A:318:ASN:N	1:A:373:GLN:HE22	2.10	0.46
1:A:232:THR:O	1:A:237:HIS:HE1	1.98	0.46
1:B:165:ILE:HD13	1:B:165:ILE:H	1.80	0.46
1:A:6:ILE:HG23	1:A:167:ASP:HB2	1.97	0.46
1:B:97:MSE:HE1	1:B:112:PRO:HD2	1.98	0.46
1:B:355:ARG:HH11	1:B:360:ARG:HH12	1.62	0.46
1:A:114:ARG:NH1	1:A:118:GLN:HE22	1.94	0.46
1:B:367:GLU:HG3	1:B:368:GLU:N	2.31	0.46
1:B:270:ILE:HD12	1:B:281:ILE:HG13	1.98	0.46
1:A:142:GLU:HB3	1:A:298:SER:OG	2.16	0.46
1:B:407:LYS:HD2	4:B:595:HOH:O	2.16	0.46
1:B:71:ILE:HD12	1:B:71:ILE:N	2.32	0.45
1:B:226:ASP:OD1	1:B:437:ARG:NH2	2.49	0.45
1:B:81:PRO:HG2	4:B:542:HOH:O	2.17	0.45
1:B:173:ASN:HB2	2:B:501:FAD:C8A	2.46	0.45
3:A:503:EPE:O3S	4:A:505:HOH:O	2.20	0.45
3:A:502:EPE:H21	1:B:49:TYR:O	2.16	0.45
1:A:226:ASP:OD1	1:A:437:ARG:NH2	2.49	0.45
1:B:232:THR:O	1:B:237:HIS:HE1	1.99	0.45
1:B:142:GLU:HB3	1:B:298:SER:OG	2.17	0.45
1:B:16:PRO:HG2	2:B:501:FAD:H4'	1.99	0.45
1:A:62:THR:O	1:A:62:THR:HG22	2.17	0.45
1:A:115:HIS:NE2	3:A:502:EPE:H52	2.32	0.45
1:B:62:THR:HG22	1:B:62:THR:O	2.17	0.45
1:B:324:PHE:N	1:B:324:PHE:CD1	2.84	0.45
1:B:56:LYS:N	4:B:536:HOH:O	2.35	0.45
1:B:257:VAL:HB	1:B:272:LEU:HD13	1.99	0.44
3:A:503:EPE:H52	1:B:115:HIS:NE2	2.33	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:367:GLU:HG3	1:A:368:GLU:N	2.32	0.44
1:A:257:VAL:HB	1:A:272:LEU:HD13	1.99	0.44
1:A:249:ILE:HG23	1:A:251:ASN:ND2	2.32	0.44
1:B:180:TYR:CD1	1:B:291:LEU:HD21	2.53	0.44
1:B:135:ALA:C	1:B:155:THR:HB	2.38	0.44
1:B:54:SER:CB	4:B:536:HOH:O	2.62	0.44
1:B:249:ILE:HG23	1:B:251:ASN:ND2	2.32	0.44
1:B:241:GLN:HG2	1:B:243:LEU:HD22	1.99	0.44
1:A:384:ASN:O	1:A:387:TYR:HB2	2.18	0.43
1:A:324:PHE:CD1	1:A:324:PHE:N	2.85	0.43
1:A:44:GLY:HA2	1:A:121:GLN:HE21	1.83	0.43
1:A:48:ASN:CB	3:A:503:EPE:H102	2.39	0.43
1:B:384:ASN:O	1:B:387:TYR:HB2	2.18	0.43
1:A:355:ARG:HH11	1:A:360:ARG:HH12	1.63	0.43
1:B:114:ARG:NH1	1:B:118:GLN:HE22	1.94	0.43
1:B:320:TYR:CZ	1:B:321:GLN:HG2	2.54	0.43
1:B:290:TYR:CD1	1:B:290:TYR:N	2.87	0.43
1:B:90:THR:HA	2:B:501:FAD:O4	2.18	0.43
1:B:251:ASN:CB	4:B:627:HOH:O	2.67	0.43
1:A:326:ILE:HB	1:A:327:PRO:HD3	2.01	0.43
1:B:151:THR:OG1	1:B:163:LYS:HE2	2.19	0.43
1:A:90:THR:HA	2:A:500:FAD:O4	2.19	0.43
1:A:373:GLN:NE2	1:A:387:TYR:OH	2.40	0.43
1:B:355:ARG:HH11	1:B:360:ARG:NH1	2.17	0.43
1:A:232:THR:HA	1:A:239:ILE:HD11	2.01	0.43
1:A:290:TYR:N	1:A:290:TYR:CD1	2.87	0.43
1:A:173:ASN:HB2	2:A:500:FAD:C8A	2.48	0.42
1:A:39:ARG:HD3	1:A:80:TYR:CE1	2.55	0.42
1:A:251:ASN:O	1:A:255:GLN:NE2	2.39	0.42
1:A:351:ALA:HB1	4:A:568:HOH:O	2.18	0.42
1:A:431:GLU:CG	4:A:671:HOH:O	2.68	0.42
1:A:91:ASN:C	1:A:91:ASN:ND2	2.63	0.42
1:B:232:THR:HA	1:B:239:ILE:HD11	2.01	0.42
1:B:373:GLN:NE2	1:B:387:TYR:OH	2.38	0.42
1:B:214:SER:HA	4:B:599:HOH:O	2.20	0.42
1:B:44:GLY:HA2	1:B:121:GLN:HE21	1.83	0.42
1:B:244:LEU:O	1:B:259:GLU:HG2	2.19	0.42
1:A:53:LEU:HD21	1:A:205:ARG:HG3	2.02	0.42
1:A:355:ARG:HH11	1:A:360:ARG:NH1	2.18	0.42
1:B:24:ALA:HB2	1:B:417:PHE:CE2	2.55	0.42
1:B:39:ARG:HD3	1:B:80:TYR:CE1	2.54	0.42
1:B:340:VAL:O	1:B:342:PRO:HD3	2.20	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:241:GLN:HG2	1:A:243:LEU:HD22	2.02	0.41
1:A:270:ILE:HD12	1:A:281:ILE:HG13	2.00	0.41
1:A:39:ARG:HG2	2:A:500:FAD:C2A	2.48	0.41
1:A:135:ALA:C	1:A:155:THR:HB	2.41	0.41
1:A:151:THR:OG1	1:A:163:LYS:HE2	2.21	0.41
1:B:177:GLU:O	1:B:179:PRO:HD3	2.20	0.41
1:A:185:LYS:CB	1:A:262:LYS:HG2	2.51	0.41
1:B:185:LYS:CG	1:B:262:LYS:HG2	2.50	0.41
1:A:244:LEU:O	1:A:259:GLU:HG2	2.21	0.41
1:B:160:PRO:HD2	4:B:680:HOH:O	2.20	0.41
1:A:185:LYS:CG	1:A:262:LYS:HG2	2.51	0.41
1:B:19:LEU:HD12	1:B:121:GLN:HE22	1.86	0.41
1:A:215:VAL:HG12	1:A:216:LEU:N	2.35	0.41
1:A:180:TYR:CD1	1:A:291:LEU:HD21	2.56	0.41
1:A:165:ILE:HD13	4:A:556:HOH:O	2.20	0.41
1:B:53:LEU:HD21	1:B:205:ARG:HG3	2.03	0.40
1:B:215:VAL:HG12	1:B:216:LEU:N	2.35	0.40
1:B:185:LYS:CB	1:B:262:LYS:HG2	2.51	0.40
1:A:320:TYR:CZ	1:A:321:GLN:HG2	2.57	0.40
1:A:177:GLU:O	1:A:179:PRO:HD3	2.22	0.40
3:A:502:EPE:H102	1:B:48:ASN:CB	2.40	0.40
1:B:326:ILE:HB	1:B:327:PRO:HD3	2.02	0.40
1:A:71:ILE:N	1:A:71:ILE:CD1	2.84	0.40
1:A:340:VAL:O	1:A:342:PRO:HD3	2.22	0.40
1:A:272:LEU:HD12	1:A:276:LYS:HG3	2.04	0.40
1:B:58:PRO:HG3	1:B:66:LEU:CD2	2.51	0.40
1:B:305:PRO:HD2	1:B:306:GLU:OE1	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	440/457 (96%)	414 (94%)	21 (5%)	5 (1%)	21 29

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	440/457 (96%)	412 (94%)	23 (5%)	5 (1%)	21	29
All	All	880/914 (96%)	826 (94%)	44 (5%)	10 (1%)	21	29

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	4	PRO
1	A	156	LYS
1	B	4	PRO
1	B	156	LYS
1	A	195	PRO
1	B	195	PRO
1	A	443	ILE
1	B	329	PRO
1	B	443	ILE
1	A	329	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	385/394 (98%)	374 (97%)	11 (3%)	55	76
1	B	385/394 (98%)	374 (97%)	11 (3%)	55	76
All	All	770/788 (98%)	748 (97%)	22 (3%)	55	76

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	PRO
1	A	91	ASN
1	A	150	VAL
1	A	155	THR
1	A	165	ILE
1	A	195	PRO
1	A	218	VAL
1	A	237	HIS

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Mol	Chain	Res	Type
1	A	297	PRO
1	A	399	ILE
1	A	419	SER
1	B	4	PRO
1	B	91	ASN
1	B	150	VAL
1	B	155	THR
1	B	165	ILE
1	B	195	PRO
1	B	218	VAL
1	B	237	HIS
1	B	297	PRO
1	B	399	ILE
1	B	419	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	48	ASN
1	A	91	ASN
1	A	118	GLN
1	A	121	GLN
1	A	200	HIS
1	A	230	HIS
1	A	237	HIS
1	A	251	ASN
1	A	373	GLN
1	B	48	ASN
1	B	91	ASN
1	B	118	GLN
1	B	121	GLN
1	B	200	HIS
1	B	230	HIS
1	B	237	HIS
1	B	251	ASN
1	B	373	GLN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

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	500	-	58,58,58	14.63	17 (29%)	85,89,89	2.88	23 (27%)
3	EPE	A	502	-	12,12,15	1.34	3 (25%)	16,16,20	2.49	5 (31%)
3	EPE	A	503	-	12,12,15	1.22	2 (16%)	16,16,20	2.54	7 (43%)
2	FAD	B	501	-	58,58,58	14.63	17 (29%)	85,89,89	2.88	22 (25%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	FAD	A	500	-	-	0/34/50/50	0/1/6/6
3	EPE	A	502	-	-	0/6/14/19	0/1/1/1
3	EPE	A	503	-	-	0/6/14/19	0/1/1/1
2	FAD	B	501	-	-	0/34/50/50	0/1/6/6

All (39) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	FAD	PA-O2A	110.06	6.46	1.55
2	A	500	FAD	PA-O2A	110.06	6.46	1.55

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	500	FAD	PA-O5B	-6.51	1.29	1.59
2	B	501	FAD	PA-O5B	-6.51	1.29	1.59
2	A	500	FAD	P-O5'	5.80	1.85	1.59
2	B	501	FAD	P-O5'	5.80	1.85	1.59
2	B	501	FAD	C1'-N10	-5.45	1.42	1.48
2	A	500	FAD	C1'-N10	-5.40	1.42	1.48
2	A	500	FAD	C9A-N10	5.17	1.46	1.38
2	B	501	FAD	C9A-N10	5.17	1.46	1.38
2	B	501	FAD	C5A-C4A	-4.45	1.30	1.40
2	A	500	FAD	C5A-C4A	-4.44	1.30	1.40
2	A	500	FAD	C5X-N5	4.00	1.41	1.35
2	B	501	FAD	C5X-N5	3.99	1.41	1.35
2	A	500	FAD	C4A-N3A	3.83	1.41	1.35
2	B	501	FAD	C4A-N3A	3.81	1.41	1.35
2	A	500	FAD	C8A-N9A	3.68	1.42	1.36
2	B	501	FAD	C8A-N9A	3.68	1.42	1.36
2	B	501	FAD	C2A-N3A	3.67	1.39	1.32
2	A	500	FAD	C2A-N3A	3.67	1.39	1.32
2	B	501	FAD	C4X-N5	3.33	1.43	1.36
2	A	500	FAD	C4X-N5	3.27	1.43	1.36
2	A	500	FAD	C4-C4X	-2.61	1.37	1.41
2	B	501	FAD	C4-C4X	-2.60	1.37	1.41
2	A	500	FAD	C5'-C4'	-2.59	1.47	1.51
2	B	501	FAD	C5'-C4'	-2.57	1.47	1.51
2	A	500	FAD	O4B-C1B	2.49	1.45	1.41
2	B	501	FAD	O4B-C1B	2.47	1.45	1.41
2	A	500	FAD	C2B-C3B	-2.43	1.46	1.53
2	B	501	FAD	C2B-C3B	-2.42	1.46	1.53
3	A	502	EPE	C9-N1	2.39	1.53	1.47
2	B	501	FAD	O5B-C5B	-2.32	1.35	1.44
2	A	500	FAD	O5B-C5B	-2.32	1.35	1.44
3	A	503	EPE	C9-N1	2.31	1.52	1.47
3	A	502	EPE	C2-N1	2.26	1.53	1.47
2	A	500	FAD	C8A-N7A	-2.09	1.30	1.34
2	B	501	FAD	C8A-N7A	-2.08	1.30	1.34
3	A	502	EPE	C6-N1	2.05	1.52	1.47
3	A	503	EPE	C6-N1	2.02	1.52	1.47

All (57) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	500	FAD	O2A-PA-O5B	-15.01	32.81	108.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	FAD	O2A-PA-O5B	-15.01	32.85	108.51
2	B	501	FAD	C8A-N9A-C4A	-8.58	100.35	106.90
2	A	500	FAD	C8A-N9A-C4A	-8.57	100.36	106.90
2	A	500	FAD	O5B-C5B-C4B	7.82	137.64	108.94
2	B	501	FAD	O5B-C5B-C4B	7.82	137.63	108.94
2	B	501	FAD	O2A-PA-O1A	-7.02	73.01	112.21
2	A	500	FAD	O2A-PA-O1A	-7.02	73.02	112.21
2	A	500	FAD	O5B-PA-O1A	-6.59	83.56	109.37
2	B	501	FAD	O5B-PA-O1A	-6.59	83.56	109.37
2	B	501	FAD	PA-O5B-C5B	6.45	168.42	122.03
2	A	500	FAD	PA-O5B-C5B	6.45	168.40	122.03
3	A	503	EPE	O1S-S-C10	-6.08	101.60	106.81
2	B	501	FAD	O2A-PA-O3P	5.57	131.58	105.14
2	A	500	FAD	O2A-PA-O3P	5.57	131.57	105.14
3	A	502	EPE	O1S-S-C10	-5.47	102.12	106.81
3	A	502	EPE	C5-N4-C3	4.97	113.59	110.50
3	A	503	EPE	C5-N4-C3	4.49	113.30	110.50
2	A	500	FAD	N3A-C2A-N1A	-3.79	125.54	128.71
2	B	501	FAD	N3A-C2A-N1A	-3.79	125.54	128.71
2	A	500	FAD	N3A-C4A-N9A	-3.70	118.74	125.43
2	B	501	FAD	N3A-C4A-N9A	-3.70	118.75	125.43
2	B	501	FAD	O3P-P-O5'	-3.69	86.89	103.41
2	A	500	FAD	O3P-P-O5'	-3.69	86.91	103.41
3	A	502	EPE	O3S-S-C10	3.41	110.24	105.93
3	A	503	EPE	O3S-S-C10	3.24	110.03	105.93
3	A	502	EPE	C6-N1-C2	3.23	116.87	108.86
3	A	503	EPE	C6-N1-C2	3.22	116.85	108.86
2	A	500	FAD	O5'-P-O1P	3.10	121.51	109.37
2	B	501	FAD	O5'-P-O1P	3.10	121.50	109.37
2	B	501	FAD	C5A-C4A-N9A	3.10	111.64	107.16
2	A	500	FAD	C5A-C4A-N9A	3.08	111.61	107.16
2	A	500	FAD	P-O5'-C5'	3.01	143.69	122.03
2	B	501	FAD	P-O5'-C5'	3.01	143.69	122.03
3	A	503	EPE	C9-N1-C6	-2.73	104.27	111.32
2	A	500	FAD	O4B-C1B-C2B	-2.70	102.64	106.77
2	B	501	FAD	O4B-C1B-C2B	-2.69	102.65	106.77
3	A	502	EPE	C9-N1-C6	-2.59	104.64	111.32
2	A	500	FAD	O3B-C3B-C4B	2.46	118.31	111.08
2	B	501	FAD	O3B-C3B-C4B	2.41	118.17	111.08
2	B	501	FAD	C2'-C1'-N10	2.38	115.61	112.45
2	A	500	FAD	C2'-C1'-N10	2.36	115.58	112.45
2	B	501	FAD	C4-C4X-C10	2.33	120.70	116.95

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	500	FAD	C4-C4X-C10	2.30	120.67	116.95
3	A	503	EPE	O2S-S-C10	2.29	108.77	106.81
2	A	500	FAD	C10-C4X-N5	2.19	123.12	120.45
2	B	501	FAD	C10-C4X-N5	2.18	123.11	120.45
2	B	501	FAD	C1'-C2'-C3'	2.14	115.94	109.82
2	A	500	FAD	C1'-C2'-C3'	2.13	115.92	109.82
2	B	501	FAD	C6-C5X-C9A	-2.09	116.12	119.02
2	A	500	FAD	C6-C5X-C9A	-2.09	116.12	119.02
2	B	501	FAD	C2-N1-C10	2.08	117.07	114.98
2	A	500	FAD	O2B-C2B-C3B	2.05	118.50	111.83
2	B	501	FAD	C8A-N9A-C1B	2.05	130.42	126.38
2	A	500	FAD	C2-N1-C10	2.04	117.04	114.98
2	A	500	FAD	C8A-N9A-C1B	2.04	130.40	126.38
3	A	503	EPE	O3S-S-O1S	2.02	116.15	111.78

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	442/457 (96%)	-0.16	5 (1%) 77 77	13, 31, 46, 56	0
1	B	442/457 (96%)	-0.13	6 (1%) 72 71	14, 32, 47, 57	0
All	All	884/914 (96%)	-0.15	11 (1%) 70 75	13, 32, 47, 57	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	3	LEU	5.5
1	B	3	LEU	4.5
1	B	244	LEU	3.4
1	A	244	LEU	3.3
1	A	4	PRO	3.3
1	B	4	PRO	2.9
1	B	271	TYR	2.6
1	A	444	GLU	2.4
1	B	266	THR	2.3
1	A	157	ALA	2.3
1	B	274	GLY	2.3

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(Å ²)	Q<0.9
3	EPE	A	502	12/15	0.26	5.71	32,34,50,50	0
3	EPE	A	503	12/15	0.24	3.69	34,36,50,51	0
2	FAD	B	501	53/53	0.22	3.05	19,27,33,43	0
2	FAD	A	500	53/53	0.22	2.02	20,27,32,41	0

6.5 Other polymers ⓘ

There are no such residues in this entry.