



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 09:54 AM GMT

PDB ID : 3K4N
Title : Pyranose 2-oxidase F454A/S455A/Y456A mutant
Authors : Divne, C.; Tan, T.C.
Deposited on : 2009-10-05
Resolution : 2.75 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

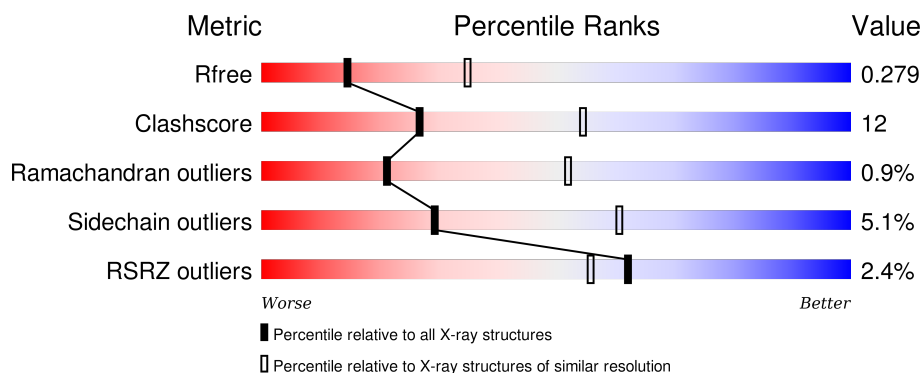
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.75 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	3340 (2.80-2.72)
Clashscore	102246	3829 (2.80-2.72)
Ramachandran outliers	100387	3767 (2.80-2.72)
Sidechain outliers	100360	3770 (2.80-2.72)
RSRZ outliers	91569	3352 (2.80-2.72)

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	623	 2% 70% 20% • 8%
1	B	623	 2% 66% 23% • 8%

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
2	FAD	A	801	X	-	-	-
2	FAD	B	801	X	-	-	-

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 9263 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.

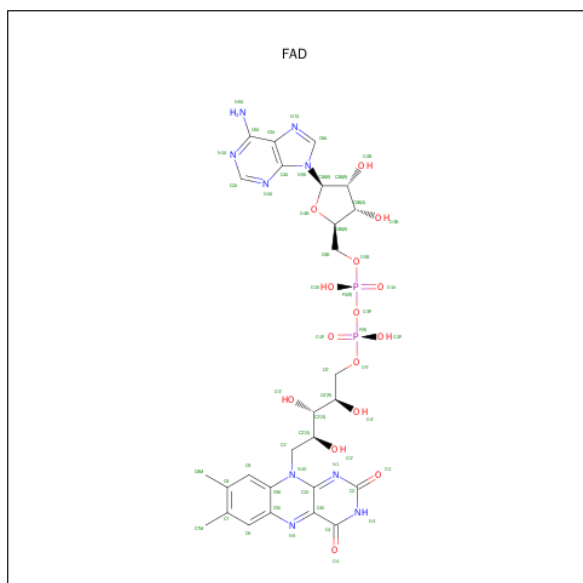
- Molecule 1 is a protein called Pyranose 2-oxidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	573	Total	C	N	O	S	0	0	0
			4504	2841	774	865	24			
1	B	573	Total	C	N	O	S	0	0	0
			4504	2841	774	865	24			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	454	ALA	PHE	ENGINEERED	UNP Q7ZA32
A	455	ALA	SER	ENGINEERED	UNP Q7ZA32
A	456	ALA	TYR	ENGINEERED	UNP Q7ZA32
B	454	ALA	PHE	ENGINEERED	UNP Q7ZA32
B	455	ALA	SER	ENGINEERED	UNP Q7ZA32
B	456	ALA	TYR	ENGINEERED	UNP Q7ZA32

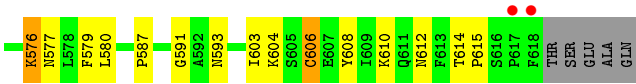
- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



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 water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	94	Total	O	0	0
			94	94		
3	B	55	Total	O	0	0
			55	55		



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	101.58 Å 101.58 Å 250.00 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.75 29.97 – 2.75	Depositor EDS
% Data completeness (in resolution range)	99.8 (30.00-2.75) 99.8 (29.97-2.75)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.19	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.77 (at 2.76 Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.191 , 0.274 0.200 , 0.279	Depositor DCC
R_{free} test set	1931 reflections (5.87%)	DCC
Wilson B-factor (Å ²)	25.8	Xtriage
Anisotropy	0.004	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 42.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Outliers	0 of 34840 reflections	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	9263	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 17.70% 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: 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	0.85	3/4618 (0.1%)	0.84	1/6280 (0.0%)
1	B	0.86	5/4618 (0.1%)	0.85	5/6280 (0.1%)
All	All	0.86	8/9236 (0.1%)	0.85	6/12560 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	228	GLU	CG-CD	8.15	1.64	1.51
1	B	429	GLU	CG-CD	5.49	1.60	1.51
1	A	59	CYS	CB-SG	-5.49	1.72	1.81
1	B	108	GLN	CG-CD	5.41	1.63	1.51
1	A	347	GLU	CB-CG	5.36	1.62	1.52
1	B	606	CYS	CB-SG	-5.35	1.73	1.81
1	A	347	GLU	CG-CD	5.32	1.59	1.51
1	B	228	GLU	CD-OE2	5.09	1.31	1.25

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	139	ARG	NE-CZ-NH2	-5.99	117.30	120.30
1	B	147	ASP	CB-CG-OD1	5.96	123.66	118.30
1	B	159	ARG	NE-CZ-NH1	5.75	123.17	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	517	ASP	CB-CG-OD1	5.62	123.36	118.30
1	A	452	ASP	CB-CG-OD1	5.45	123.20	118.30
1	B	159	ARG	NE-CZ-NH2	-5.05	117.77	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	460	GLN	Peptide

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	4504	0	4356	105	0
1	B	4504	0	4356	120	0
2	A	53	0	27	9	0
2	B	53	0	28	8	0
3	A	94	0	0	1	0
3	B	55	0	0	2	0
All	All	9263	0	8767	215	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 12.

All (215) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:167:HIS:NE2	2:B:801:FAD:HM82	1.21	1.47
1:A:167:HIS:NE2	2:A:801:FAD:HM82	1.04	1.35
1:A:167:HIS:CE1	2:A:801:FAD:HM82	1.63	1.32
1:A:456:ALA:CB	1:A:461:GLN:HE21	1.70	1.04
1:A:167:HIS:NE2	2:A:801:FAD:C8	2.27	0.96
1:B:91:LYS:HD2	1:B:100:ILE:HD11	1.55	0.89
1:A:456:ALA:HB1	1:A:461:GLN:HE21	1.38	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:167:HIS:CE1	2:A:801:FAD:C8M	2.38	0.86
1:A:123:VAL:HG22	1:B:459:VAL:HG22	1.58	0.84
1:A:456:ALA:HB3	1:A:461:GLN:HE21	1.40	0.83
1:B:542:GLU:HG2	1:B:545:LEU:HD13	1.65	0.78
1:A:477:THR:HG23	1:A:511:THR:HG21	1.69	0.75
1:B:104:VAL:HG21	1:B:455:ALA:HB2	1.69	0.73
1:B:167:HIS:CE1	2:B:801:FAD:HM82	2.19	0.73
1:A:123:VAL:HG21	1:B:459:VAL:HG13	1.71	0.73
1:B:343:ALA:O	1:B:345:PRO:HD3	1.89	0.72
1:A:216:SER:O	1:A:220:ASN:ND2	2.21	0.71
1:A:176:ASP:OD1	1:A:179:GLN:NE2	2.21	0.70
1:B:217:ILE:HD12	1:B:427:PRO:CD	2.22	0.69
1:B:126:LEU:HD13	1:B:132:GLN:HG3	1.73	0.69
1:B:217:ILE:HD12	1:B:427:PRO:HD2	1.75	0.68
1:A:465:SER:HA	1:A:468:ILE:HD12	1.75	0.67
1:B:50:VAL:HG13	1:B:313:ALA:HB2	1.76	0.67
1:B:57:ILE:N	1:B:57:ILE:HD13	2.09	0.67
1:A:63:ARG:HD2	1:A:259:VAL:O	1.95	0.66
1:B:461:GLN:H	1:B:461:GLN:HE21	1.43	0.66
1:B:47:TYR:O	1:B:313:ALA:HA	1.96	0.65
1:B:50:VAL:HG13	1:B:313:ALA:CB	2.27	0.65
1:B:468:ILE:HD13	1:B:533:LEU:HD22	1.79	0.64
1:B:549:LEU:HD21	1:B:552:THR:HG21	1.79	0.64
1:B:126:LEU:HD13	1:B:132:GLN:CG	2.27	0.64
1:B:336:GLN:NE2	1:B:344:ASN:O	2.28	0.64
1:B:167:HIS:NE2	2:B:801:FAD:C8	2.61	0.64
1:B:459:VAL:C	1:B:460:GLN:HG2	2.19	0.64
1:B:126:LEU:CD1	1:B:132:GLN:HG3	2.27	0.64
1:A:369:SER:OG	1:A:372:LEU:HD12	1.98	0.63
1:B:184:VAL:HG23	1:B:191:ASP:HA	1.79	0.62
1:B:194:GLU:OE2	1:B:604:LYS:NZ	2.28	0.62
1:B:97:GLN:NE2	1:B:251:VAL:O	2.29	0.62
1:B:456:ALA:HB1	1:B:460:GLN:O	2.00	0.61
1:B:114:VAL:HG22	1:B:157:VAL:HG13	1.82	0.61
1:B:131:TRP:CH2	1:B:133:ALA:HB2	2.35	0.61
1:A:456:ALA:HB1	1:A:461:GLN:NE2	2.12	0.61
1:B:555:MET:HA	1:B:566:VAL:O	2.01	0.61
1:B:608:TYR:O	1:B:612:ASN:ND2	2.31	0.60
1:A:349:LEU:HD11	1:A:572:VAL:HG13	1.81	0.60
1:B:437:THR:O	1:B:437:THR:HG23	2.02	0.60
1:A:299:HIS:NE2	1:A:310:GLU:OE2	2.35	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:230:TYR:O	1:B:231:LYS:C	2.41	0.59
1:A:239:ILE:HG22	1:A:240:PRO:O	2.03	0.58
1:A:167:HIS:CE1	2:A:801:FAD:HM81	2.38	0.57
1:B:321:GLY:O	1:B:325:ASN:HB2	2.05	0.57
1:A:459:VAL:HG11	1:A:534:PRO:HB2	1.87	0.56
1:A:380:MET:HE1	1:A:409:ASN:HB3	1.88	0.56
1:A:458:ALA:HB1	1:B:121:LEU:HD13	1.87	0.56
2:A:801:FAD:O2'	2:A:801:FAD:O4'	2.15	0.55
1:A:167:HIS:C	1:A:167:HIS:CD2	2.79	0.55
1:B:468:ILE:HD13	1:B:533:LEU:CD2	2.37	0.55
1:B:101:ASP:O	1:B:104:VAL:HG12	2.06	0.55
1:B:359:GLN:OE1	1:B:548:HIS:ND1	2.38	0.55
1:A:432:GLU:HG3	1:A:451:ARG:HB2	1.88	0.55
1:B:167:HIS:CD2	2:B:801:FAD:HM82	2.26	0.54
1:A:456:ALA:HB3	1:A:461:GLN:NE2	2.18	0.54
1:B:435:VAL:HG12	1:B:449:ILE:HD12	1.89	0.54
1:B:157:VAL:HG21	1:B:324:HIS:HE1	1.72	0.54
2:B:801:FAD:H8A	3:B:1139:HOH:O	2.07	0.54
1:B:542:GLU:O	1:B:545:LEU:HB2	2.08	0.54
1:A:285:ARG:NH2	1:A:299:HIS:ND1	2.55	0.54
1:A:456:ALA:HB1	1:A:461:GLN:CG	2.38	0.54
1:B:302:ASP:HB3	1:B:305:SER:OG	2.07	0.54
1:A:86:ILE:O	1:A:257:ASN:HB2	2.07	0.53
1:B:64:GLU:OE2	1:B:201:LYS:NZ	2.35	0.53
1:A:380:MET:HE1	1:A:409:ASN:CB	2.39	0.53
1:A:223:LEU:HD13	1:A:223:LEU:O	2.10	0.52
1:A:532:PHE:CE1	1:A:538:PRO:HG3	2.45	0.52
1:A:367:VAL:HG21	1:B:126:LEU:CD2	2.40	0.52
1:A:572:VAL:HG21	1:A:578:LEU:HD23	1.90	0.52
1:B:359:GLN:NE2	1:B:591:GLY:O	2.43	0.52
1:A:282:ALA:HB1	3:A:1045:HOH:O	2.09	0.51
1:B:50:VAL:CG1	1:B:313:ALA:HB2	2.40	0.51
1:A:47:TYR:CE2	1:A:73:ALA:HB2	2.46	0.51
1:A:47:TYR:CD2	1:A:73:ALA:HB2	2.46	0.51
1:B:159:ARG:HA	2:B:801:FAD:O2B	2.11	0.51
1:B:70:TYR:OH	1:B:610:LYS:HA	2.10	0.51
1:A:167:HIS:NE2	2:A:801:FAD:HM81	2.09	0.50
1:B:77:ILE:HG23	1:B:159:ARG:HB3	1.94	0.50
1:A:299:HIS:CE1	1:A:310:GLU:OE2	2.64	0.50
1:A:389:LEU:H	1:A:389:LEU:HD23	1.76	0.50
1:A:437:THR:HG23	1:A:437:THR:O	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:579:PHE:O	1:B:580:LEU:HD23	2.11	0.50
1:A:94:VAL:HG12	1:A:98:LYS:HE2	1.94	0.50
1:A:83:GLY:HA2	1:B:80:ILE:HG22	1.94	0.50
1:A:362:VAL:CG2	1:A:473:PHE:HB2	2.42	0.50
1:B:452:ASP:OD1	1:B:452:ASP:C	2.50	0.49
1:A:163:GLY:HA2	2:A:801:FAD:O3B	2.12	0.49
1:A:300:ILE:CD1	1:A:311:ILE:HG12	2.42	0.49
1:A:300:ILE:HD13	1:A:311:ILE:HG12	1.94	0.49
1:A:131:TRP:CH2	1:A:133:ALA:HB2	2.47	0.49
1:B:437:THR:O	1:B:437:THR:CG2	2.60	0.49
1:B:356:ILE:HG23	1:B:483:ASN:O	2.13	0.49
1:B:373:ILE:HD11	1:B:426:ILE:HD11	1.95	0.49
1:B:284:GLU:O	1:B:331:ASN:ND2	2.40	0.49
1:A:104:VAL:HG21	1:A:455:ALA:HB2	1.94	0.48
1:B:373:ILE:CD1	1:B:426:ILE:HD11	2.43	0.48
1:A:471:TRP:N	1:A:471:TRP:CD1	2.82	0.48
1:A:555:MET:SD	1:A:568:THR:HA	2.54	0.48
1:A:380:MET:HE1	1:A:409:ASN:CA	2.43	0.48
1:B:137:PHE:O	1:B:139:ARG:HD3	2.14	0.48
1:A:98:LYS:HD3	1:B:494:ALA:HB1	1.96	0.48
1:A:532:PHE:CZ	1:A:538:PRO:CG	2.97	0.47
1:B:459:VAL:HG12	1:B:459:VAL:O	2.14	0.47
1:B:110:GLN:HE21	1:B:167:HIS:HD1	1.61	0.47
1:A:187:ASP:HB3	1:A:190:ALA:HB3	1.96	0.47
1:A:456:ALA:CB	1:A:461:GLN:NE2	2.56	0.47
1:A:243:ALA:HB2	1:A:253:TRP:HA	1.96	0.47
1:B:614:THR:HG22	1:B:615:PRO:O	2.14	0.47
1:B:606:CYS:O	1:B:610:LYS:HG3	2.15	0.47
1:A:459:VAL:O	1:A:533:LEU:HD11	2.15	0.47
1:A:169:THR:O	1:A:170:CYS:HB2	2.14	0.47
1:B:104:VAL:HG13	1:B:105:ASN:N	2.31	0.46
1:A:449:ILE:HG12	1:A:471:TRP:CZ3	2.50	0.46
1:A:324:HIS:HD2	1:A:327:GLN:OE1	1.98	0.46
1:B:567:ASN:C	1:B:567:ASN:OD1	2.53	0.46
1:A:75:PHE:CD1	1:A:278:PHE:HB2	2.50	0.46
1:A:608:TYR:CE2	1:A:612:ASN:ND2	2.83	0.46
1:B:50:VAL:HA	1:B:73:ALA:O	2.16	0.46
1:A:101:ASP:O	1:A:104:VAL:HG12	2.16	0.46
1:A:362:VAL:HG23	1:A:473:PHE:HB2	1.96	0.46
1:A:219:HIS:HA	1:A:433:PRO:HB3	1.97	0.46
1:A:123:VAL:CG2	1:B:459:VAL:HG13	2.43	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:576:LYS:HD3	1:B:577:ASN:HB3	1.97	0.46
1:A:459:VAL:HG12	1:A:459:VAL:O	2.16	0.46
1:A:49:VAL:HB	1:A:72:VAL:HG22	1.98	0.46
1:B:58:GLY:HA2	1:B:319:THR:HG21	1.98	0.46
1:B:567:ASN:HB3	1:B:573:PHE:CE1	2.51	0.46
1:B:57:ILE:N	1:B:57:ILE:CD1	2.79	0.45
1:B:461:GLN:HE21	1:B:461:GLN:N	2.13	0.45
1:B:571:ARG:HB2	1:B:579:PHE:CE2	2.51	0.45
1:A:362:VAL:HA	1:A:539:GLN:O	2.16	0.45
1:B:55:GLY:HA3	2:B:801:FAD:O1P	2.16	0.45
1:A:446:HIS:HB3	1:A:474:PHE:HB2	1.99	0.45
1:A:126:LEU:HD12	1:A:132:GLN:HG3	1.98	0.45
1:A:286:VAL:HG23	1:A:332:SER:HB3	1.99	0.45
1:B:264:ASN:ND2	1:B:276:ASN:HD22	2.14	0.45
1:B:496:ASN:O	3:B:1019:HOH:O	2.21	0.45
1:B:554:ARG:HB2	1:B:587:PRO:HB3	1.98	0.45
1:B:163:GLY:HA2	2:B:801:FAD:O3B	2.17	0.45
1:A:176:ASP:O	1:A:180:ARG:HB2	2.16	0.45
1:A:223:LEU:HD13	1:A:223:LEU:C	2.36	0.45
1:A:111:LEU:HD22	1:A:156:ALA:HB3	1.99	0.45
1:B:478:GLU:OE1	1:B:511:THR:OG1	2.21	0.45
1:A:594:PRO:O	1:A:595:THR:C	2.54	0.45
1:A:493:ASP:C	1:A:493:ASP:OD1	2.55	0.44
1:A:261:ASP:C	1:A:261:ASP:OD1	2.55	0.44
1:A:126:LEU:HD23	1:B:534:PRO:HD3	1.98	0.44
1:A:450:HIS:O	1:A:470:ASP:HB2	2.18	0.44
1:A:265:ARG:HA	1:A:266:PRO:C	2.38	0.44
1:B:205:TYR:HB3	1:B:259:VAL:HG13	2.00	0.44
1:B:412:VAL:O	1:B:416:MET:HG3	2.17	0.44
1:A:70:TYR:OH	1:A:613:PHE:O	2.27	0.44
1:A:218:ARG:HD2	1:A:426:ILE:HG21	1.99	0.44
1:B:464:ASP:O	1:B:465:SER:C	2.55	0.43
1:B:131:TRP:HH2	1:B:133:ALA:HB2	1.80	0.43
1:A:131:TRP:CD1	1:B:373:ILE:HG21	2.54	0.43
1:B:119:ASN:OD1	1:B:119:ASN:C	2.57	0.43
1:A:83:GLY:N	1:B:81:ASP:HA	2.33	0.43
1:A:491:ILE:HG22	1:A:492:THR:N	2.33	0.43
1:B:242:ALA:HB1	1:B:254:SER:HB2	2.00	0.43
1:A:336:GLN:HE22	1:A:341:ASN:HB3	1.83	0.43
1:A:80:ILE:O	1:B:83:GLY:HA3	2.19	0.43
1:A:359:GLN:OE1	1:A:548:HIS:ND1	2.33	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:167:HIS:CD2	1:B:167:HIS:C	2.92	0.43
1:B:226:LEU:CD1	1:B:449:ILE:HD13	2.48	0.43
1:B:47:TYR:CD2	1:B:73:ALA:HB2	2.53	0.42
1:A:570:SER:HB2	1:A:605:SER:HB2	2.00	0.42
1:B:452:ASP:O	1:B:453:ALA:C	2.58	0.42
1:B:487:PHE:HB3	1:B:498:PRO:HB2	2.00	0.42
1:B:555:MET:O	1:B:587:PRO:HG3	2.19	0.42
1:A:380:MET:CE	1:A:409:ASN:CB	2.97	0.42
1:B:348:LEU:O	1:B:349:LEU:C	2.58	0.42
1:A:182:LEU:HD13	1:A:185:LYS:HA	2.00	0.42
1:A:94:VAL:CG1	1:A:98:LYS:HE2	2.49	0.42
1:B:181:PRO:HG3	1:B:587:PRO:HD2	2.01	0.42
1:B:163:GLY:O	1:B:166:THR:OG1	2.35	0.42
1:B:532:PHE:CZ	1:B:538:PRO:HG3	2.55	0.42
1:B:415:HIS:NE2	1:B:424:LEU:O	2.44	0.42
1:B:549:LEU:HD21	1:B:552:THR:CG2	2.48	0.42
1:A:380:MET:CE	1:A:409:ASN:HB3	2.48	0.42
1:B:120:THR:HA	1:B:136:PHE:CD1	2.55	0.42
1:A:172:THR:HB	1:A:596:LEU:HD13	2.02	0.42
1:B:230:TYR:CD1	1:B:234:ARG:HG3	2.55	0.42
1:A:126:LEU:HD23	1:B:534:PRO:CD	2.50	0.42
1:B:264:ASN:ND2	1:B:276:ASN:ND2	2.68	0.42
1:A:47:TYR:HA	1:A:71:LYS:HD3	2.02	0.41
1:B:404:HIS:CE1	1:B:408:TRP:CZ2	3.08	0.41
1:B:77:ILE:HG23	1:B:78:GLY:N	2.34	0.41
1:B:439:PHE:CD1	1:B:439:PHE:C	2.93	0.41
1:B:100:ILE:HA	1:B:100:ILE:HD12	1.73	0.41
1:B:263:GLN:O	1:B:265:ARG:NH1	2.53	0.41
1:B:52:VAL:HG12	1:B:320:ALA:HB2	2.01	0.41
1:A:373:ILE:HG23	1:A:425:PRO:HD2	2.02	0.41
1:B:239:ILE:HG22	1:B:240:PRO:O	2.21	0.41
1:A:459:VAL:HA	1:A:462:SER:HB3	2.02	0.41
1:B:603:ILE:O	1:B:606:CYS:HB2	2.21	0.41
1:A:300:ILE:CD1	1:A:311:ILE:CG1	2.99	0.41
1:B:497:MET:HE2	1:B:497:MET:HB3	1.87	0.41
1:B:506:PHE:O	1:B:507:PRO:C	2.57	0.41
1:B:49:VAL:HG22	1:B:315:VAL:HB	2.03	0.41
1:B:219:HIS:HA	1:B:433:PRO:HB3	2.03	0.40
1:A:110:GLN:HE21	1:A:167:HIS:HD1	1.70	0.40
1:B:373:ILE:HD13	1:B:373:ILE:HA	1.92	0.40
1:B:556:GLY:O	1:B:567:ASN:HA	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:167:HIS:CD2	2:A:801:FAD:C8	3.04	0.40
1:A:223:LEU:HD22	1:A:236:PHE:HB3	2.03	0.40
1:A:572:VAL:HG21	1:A:578:LEU:CD2	2.51	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	571/623 (92%)	533 (93%)	35 (6%)	3 (0%)	34	67
1	B	571/623 (92%)	517 (90%)	47 (8%)	7 (1%)	16	43
All	All	1142/1246 (92%)	1050 (92%)	82 (7%)	10 (1%)	21	52

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	231	LYS
1	B	453	ALA
1	A	231	LYS
1	A	453	ALA
1	A	458	ALA
1	B	187	ASP
1	B	264	ASN
1	B	343	ALA
1	B	101	ASP
1	B	456	ALA

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	498/539 (92%)	470 (94%)	28 (6%)	26	56
1	B	498/539 (92%)	475 (95%)	23 (5%)	33	65
All	All	996/1078 (92%)	945 (95%)	51 (5%)	29	61

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	46	LYS
1	A	100	ILE
1	A	112	MET
1	A	165	SER
1	A	167	HIS
1	A	168	TRP
1	A	169	THR
1	A	185	LYS
1	A	204	SER
1	A	223	LEU
1	A	228	GLU
1	A	233	GLN
1	A	279	PRO
1	A	286	VAL
1	A	300	ILE
1	A	347	GLU
1	A	348	LEU
1	A	375	SER
1	A	389	LEU
1	A	450	HIS
1	A	460	GLN
1	A	465	SER
1	A	490	LYS
1	A	496	ASN
1	A	505	ARG
1	A	533	LEU
1	A	561	GLU

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Mol	Chain	Res	Type
1	A	593	ASN
1	B	112	MET
1	B	139	ARG
1	B	160	VAL
1	B	168	TRP
1	B	184	VAL
1	B	206	PHE
1	B	291	LEU
1	B	326	THR
1	B	328	LEU
1	B	344	ASN
1	B	371	GLU
1	B	388	GLU
1	B	408	TRP
1	B	429	GLU
1	B	450	HIS
1	B	452	ASP
1	B	460	GLN
1	B	461	GLN
1	B	462	SER
1	B	545	LEU
1	B	570	SER
1	B	576	LYS
1	B	593	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	105	ASN
1	A	233	GLN
1	A	276	ASN
1	A	324	HIS
1	A	344	ASN
1	A	461	GLN
1	A	612	ASN
1	B	108	GLN
1	B	263	GLN
1	B	264	ASN
1	B	276	ASN
1	B	324	HIS
1	B	461	GLN

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 ⓘ

2 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	801	1	48,58,58	1.28	6 (12%)	54,89,89	3.33	21 (38%)
2	FAD	B	801	1	48,58,58	1.46	7 (14%)	54,89,89	3.20	21 (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	801	1	1/1/9/9	0/30/50/50	0/6/6/6
2	FAD	B	801	1	2/2/9/9	0/30/50/50	0/6/6/6

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	801	FAD	C8M-C8	-4.58	1.41	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	801	FAD	O3B-C3B	-2.53	1.36	1.43
2	A	801	FAD	C2B-C3B	-2.21	1.47	1.53
2	B	801	FAD	C2B-C3B	-2.03	1.47	1.53
2	B	801	FAD	C1'-N10	2.23	1.50	1.48
2	A	801	FAD	C4-N3	2.26	1.37	1.33
2	A	801	FAD	C5'-C4'	2.31	1.55	1.51
2	B	801	FAD	C2A-N1A	2.83	1.39	1.33
2	A	801	FAD	C2A-N1A	2.98	1.39	1.33
2	A	801	FAD	C2A-N3A	3.01	1.37	1.32
2	B	801	FAD	C2A-N3A	3.18	1.37	1.32
2	A	801	FAD	C4-C4X	4.04	1.49	1.41
2	B	801	FAD	C4-C4X	4.21	1.49	1.41

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	801	FAD	N3A-C2A-N1A	-12.94	118.99	128.89
2	B	801	FAD	N3A-C2A-N1A	-11.45	120.13	128.89
2	B	801	FAD	C4-C4X-C10	-8.49	114.51	119.94
2	A	801	FAD	C4X-C4-N3	-7.85	112.86	123.59
2	B	801	FAD	C4B-O4B-C1B	-5.94	103.19	109.72
2	A	801	FAD	C4B-O4B-C1B	-4.62	104.64	109.72
2	A	801	FAD	C4X-C10-N10	-4.45	117.89	120.52
2	B	801	FAD	C4X-C10-N10	-3.86	118.24	120.52
2	B	801	FAD	C1B-N9A-C4A	-3.09	122.27	126.94
2	A	801	FAD	C6-C7-C8	-2.87	114.56	120.04
2	A	801	FAD	C9-C9A-C5X	-2.67	114.88	119.62
2	A	801	FAD	O3P-PA-O5B	-2.64	95.93	102.94
2	A	801	FAD	C1B-N9A-C4A	-2.46	123.23	126.94
2	B	801	FAD	C4X-C4-N3	-2.36	120.36	123.59
2	B	801	FAD	C9A-C5X-N5	-2.17	119.14	122.36
2	A	801	FAD	C9A-C5X-N5	-2.14	119.20	122.36
2	B	801	FAD	P-O3P-PA	-2.13	126.74	132.73
2	A	801	FAD	C7-C6-C5X	-2.11	117.48	120.92
2	B	801	FAD	C4-C4X-N5	2.01	121.16	118.72
2	A	801	FAD	C2A-N1A-C6A	2.48	123.20	118.77
2	B	801	FAD	O5B-C5B-C4B	2.49	118.29	109.12
2	A	801	FAD	O4B-C4B-C3B	2.50	110.19	105.15
2	A	801	FAD	C5X-C9A-N10	2.57	119.57	117.62
2	B	801	FAD	O3B-C3B-C2B	2.73	120.71	111.83
2	A	801	FAD	C5B-C4B-C3B	2.85	126.53	115.21
2	A	801	FAD	O2B-C2B-C3B	2.94	121.39	111.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	801	FAD	C5B-C4B-C3B	3.01	127.16	115.21
2	B	801	FAD	O3P-P-O5'	3.18	111.36	102.94
2	B	801	FAD	C5X-C9A-N10	3.50	120.28	117.62
2	B	801	FAD	O2B-C2B-C3B	3.54	123.35	111.83
2	A	801	FAD	C4X-N5-C5X	3.58	120.89	116.76
2	A	801	FAD	O4B-C4B-C5B	3.83	123.00	109.32
2	B	801	FAD	O4B-C4B-C5B	4.01	123.67	109.32
2	B	801	FAD	C4-N3-C2	4.05	118.74	115.25
2	B	801	FAD	C4X-N5-C5X	4.21	121.61	116.76
2	B	801	FAD	O3B-C3B-C4B	4.94	125.88	111.05
2	A	801	FAD	O4B-C1B-N9A	5.35	119.29	108.10
2	A	801	FAD	O3B-C3B-C4B	6.79	131.43	111.05
2	B	801	FAD	C2B-C1B-N9A	6.95	124.91	114.29
2	A	801	FAD	C2B-C1B-N9A	7.48	125.72	114.29
2	B	801	FAD	O4B-C1B-N9A	7.94	124.73	108.10
2	A	801	FAD	C4-N3-C2	8.11	122.25	115.25

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	B	801	FAD	C4B
2	B	801	FAD	C1B
2	A	801	FAD	C4B

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 17 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	801	FAD	9	0
2	B	801	FAD	8	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	573/623 (91%)	-0.22	12 (2%) 67 61	17, 24, 33, 51	0
1	B	573/623 (91%)	-0.16	15 (2%) 59 53	17, 25, 37, 49	0
All	All	1146/1246 (91%)	-0.19	27 (2%) 62 56	17, 25, 36, 51	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	458	ALA	5.2
1	B	456	ALA	5.2
1	B	459	VAL	5.0
1	A	459	VAL	4.5
1	B	618	PHE	4.5
1	A	456	ALA	4.4
1	B	458	ALA	4.2
1	B	343	ALA	4.1
1	B	457	GLY	3.8
1	B	342	PRO	3.7
1	A	457	GLY	3.7
1	A	618	PHE	3.4
1	B	455	ALA	3.2
1	A	343	ALA	3.1
1	A	341	ASN	2.9
1	B	341	ASN	2.9
1	B	345	PRO	2.9
1	B	617	PRO	2.9
1	A	460	GLN	2.8
1	A	455	ALA	2.7
1	B	344	ASN	2.5
1	A	344	ASN	2.5
1	A	454	ALA	2.4
1	A	617	PRO	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	270	ALA	2.3
1	B	460	GLN	2.2
1	B	271	PRO	2.2

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
2	FAD	B	801	53/53	0.97	0.12	-0.59	16,27,34,35	0
2	FAD	A	801	53/53	0.97	0.12	-0.93	20,25,30,33	0

6.5 Other polymers [i](#)

There are no such residues in this entry.