



## wwPDB EM Map/Model Validation Report ⓘ

Sep 20, 2016 – 08:54 PM EDT

PDB ID : 5I68  
EMDB ID: : EMD-8072  
Title : Alcohol oxidase from Pichia pastoris  
Authors : Vonck, J.; Mills, D.J.; Parcej, D.N.  
Deposited on : 2016-02-16  
Resolution : 3.37 Å(reported)  
Based on PDB ID : 3NNE, 3FIM, 1GAL

This is a wwPDB EM Map/Model Validation Report for a publicly released PDB/EMDB entry.  
For rigid body fitted models, validation errors reported here could stem from errors in the original structure(s) used in the fitting.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/EMValidationReportHelp>

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MolProbity : 4.02b-467  
Mogul : 1.7.1 (RC1), CSD as537be (2016)  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : rb-20027939

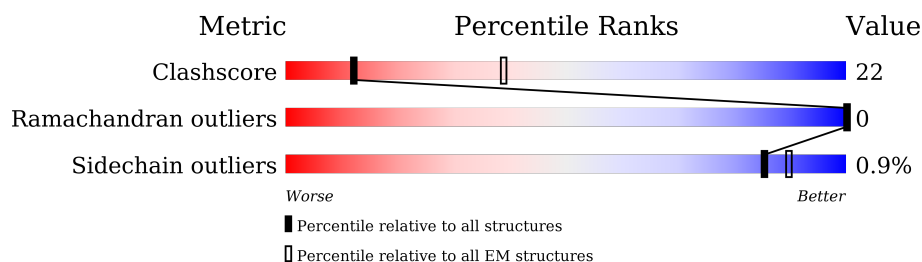
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*ELECTRON MICROSCOPY*

The reported resolution of this entry is 3.37 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)	EM structures (#Entries)
Clashscore	114402	924
Ramachandran outliers	111179	726
Sidechain outliers	111093	686

The table below summarises the geometric issues observed across the polymeric chains. The red, orange, yellow and green segments on the bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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  $\leq 5\%$ .

Mol	Chain	Length	Quality of chain
1	A	663	

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 5252 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, 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.

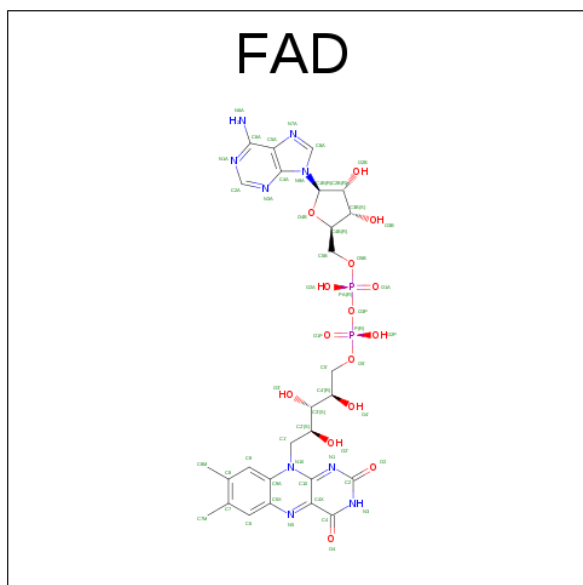
- Molecule 1 is a protein called Alcohol oxidase 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	662	5198	3286	900	986	26	0	0

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		AltConf
2	A	1	Total	Mg	0
			1	1	

- Molecule 3 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: C<sub>27</sub>H<sub>33</sub>N<sub>9</sub>O<sub>15</sub>P<sub>2</sub>).

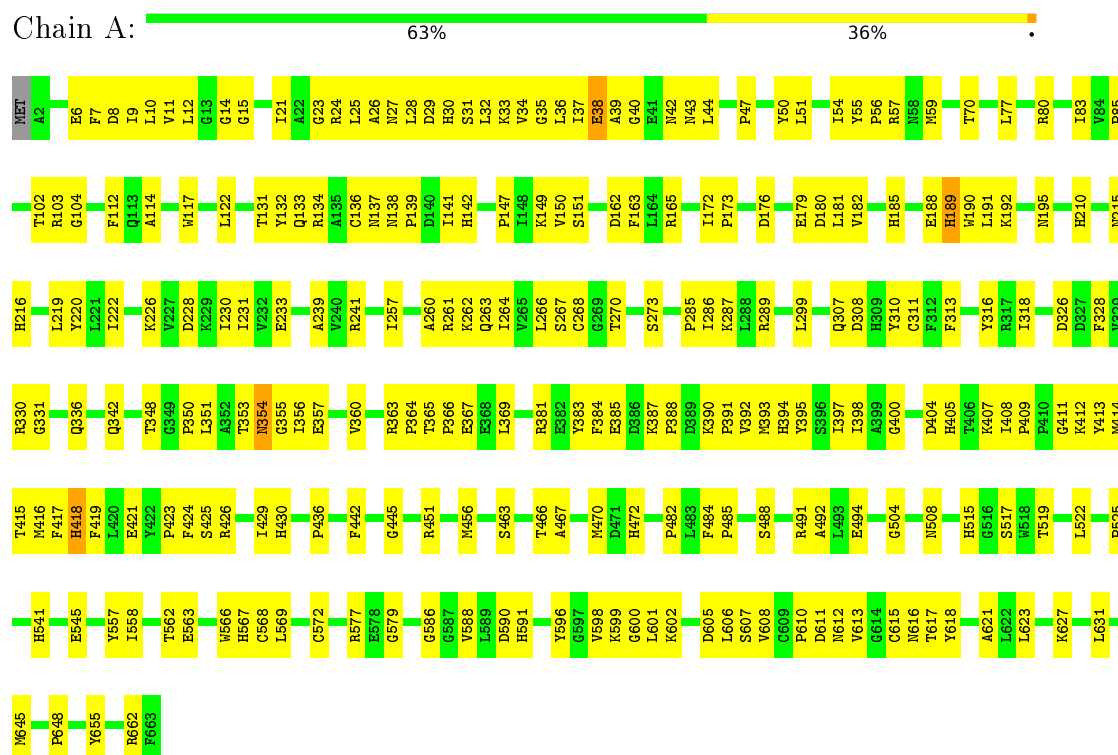


Mol	Chain	Residues	Atoms					AltConf
			Total	C	N	O	P	
3	A	1	53	27	9	15	2	0

### 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. 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. 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: Alcohol oxidase 1



## 4 Experimental information

Property	Value	Source
Reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, D	Depositor
Number of particles used	49559	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	Not provided	Depositor
Microscope	JEOL 3200FSC	Depositor
Voltage (kV)	300	Depositor
Electron dose ( $e^-/\text{\AA}^2$ )	Not provided	Depositor
Minimum defocus (nm)	Not provided	Depositor
Maximum defocus (nm)	Not provided	Depositor
Magnification	30000	Depositor
Image detector	Not provided	Depositor

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MG, 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	$\# Z  > 2$	RMSZ	$\# Z  > 2$
1	A	0.42	0/5340	0.58	1/7245 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	A	189	HIS	N-CA-C	-5.72	95.56	111.00

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	5198	0	5011	231	0
2	A	1	0	0	0	0
3	A	53	0	31	5	0
All	All	5252	0	5042	231	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 22.

All (231) 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:A:103:ARG:CD	1:A:181:LEU:HD11	1.70	1.21
1:A:134:ARG:NH2	1:A:188:GLU:OE2	1.81	1.12
1:A:397:ILE:HD11	1:A:414:MET:HE3	1.31	1.11
1:A:397:ILE:HD11	1:A:414:MET:CE	1.83	1.08
1:A:267:SER:HA	1:A:605:ASP:OD1	1.57	1.04
1:A:397:ILE:CD1	1:A:414:MET:HE3	1.91	1.00
1:A:103:ARG:HD3	1:A:181:LEU:HD11	1.44	1.00
1:A:103:ARG:HD2	1:A:181:LEU:HD11	1.42	0.96
1:A:8:ASP:OD2	1:A:32:LEU:HD22	1.69	0.93
1:A:397:ILE:CD1	1:A:414:MET:CE	2.47	0.92
1:A:103:ARG:HD2	1:A:181:LEU:CD1	2.01	0.91
1:A:59:MET:CE	1:A:59:MET:HA	2.01	0.90
1:A:7:PHE:HE1	1:A:220:TYR:CD2	1.89	0.89
1:A:8:ASP:HB3	1:A:33:LYS:O	1.76	0.85
1:A:59:MET:HE2	1:A:59:MET:HA	1.60	0.84
1:A:264:ILE:HG22	1:A:600:GLY:O	1.79	0.81
1:A:318:ILE:HG22	1:A:412:LYS:O	1.82	0.78
1:A:616:ASN:HB3	3:A:702:FAD:O2	1.85	0.76
1:A:363:ARG:NH2	1:A:384:PHE:O	2.19	0.76
1:A:14:GLY:HA3	1:A:38:GLU:HG3	1.67	0.74
1:A:267:SER:CA	1:A:605:ASP:OD1	2.36	0.74
1:A:397:ILE:HD11	1:A:414:MET:HE2	1.68	0.74
1:A:7:PHE:CE1	1:A:220:TYR:CD2	2.75	0.74
1:A:381:ARG:O	1:A:385:GLU:HB3	1.87	0.74
1:A:356:ILE:H	1:A:356:ILE:HD12	1.54	0.73
1:A:579:GLY:HA2	1:A:586:GLY:HA2	1.73	0.71
1:A:122:LEU:HD23	1:A:623:LEU:HD21	1.72	0.70
1:A:180:ASP:O	1:A:181:LEU:HB2	1.91	0.70
1:A:567:HIS:HD2	1:A:617:THR:H	1.38	0.69
1:A:70:THR:HG22	1:A:83:ILE:HG12	1.75	0.69
1:A:137:ASN:O	1:A:139:PRO:HD3	1.93	0.68
1:A:28:LEU:HD23	1:A:645:MET:SD	2.34	0.68
1:A:400:GLY:O	1:A:415:THR:HG22	1.94	0.68
1:A:356:ILE:HG13	1:A:398:ILE:HG12	1.76	0.68
1:A:30:HIS:CE1	1:A:216:HIS:NE2	2.62	0.67
1:A:132:TYR:CE1	1:A:142:HIS:CD2	2.82	0.67
1:A:504:GLY:HA3	1:A:508:ASN:HD22	1.59	0.67
1:A:172:ILE:HD13	1:A:360:VAL:HG11	1.77	0.66
1:A:394:HIS:HB2	1:A:421:GLU:OE2	1.95	0.66
1:A:150:VAL:HG12	1:A:191:LEU:HD22	1.78	0.66
1:A:21:ILE:HG22	1:A:25:LEU:HD12	1.80	0.64
1:A:264:ILE:CG2	1:A:601:LEU:HA	2.27	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:38:GLU:OE1	1:A:40:GLY:N	2.31	0.64
1:A:266:LEU:O	1:A:605:ASP:OD1	2.15	0.63
1:A:567:HIS:CD2	1:A:617:THR:H	2.15	0.63
1:A:311:CYS:HA	1:A:418:HIS:O	1.98	0.63
1:A:102:THR:CG2	1:A:616:ASN:H	2.12	0.62
1:A:231:ILE:HB	1:A:239:ALA:HB3	1.82	0.61
1:A:408:ILE:HG12	1:A:409:PRO:HD2	1.81	0.61
1:A:318:ILE:HG13	1:A:472:HIS:HB3	1.81	0.61
1:A:605:ASP:O	1:A:608:VAL:HG23	2.01	0.60
1:A:162:ASP:OD1	1:A:165:ARG:NH2	2.34	0.60
1:A:602:LYS:NZ	1:A:631:LEU:HB3	2.16	0.60
1:A:287:LYS:HB3	1:A:430:HIS:CE1	2.37	0.60
1:A:11:VAL:O	1:A:36:LEU:HD12	2.02	0.60
1:A:14:GLY:HA3	1:A:38:GLU:CG	2.32	0.59
1:A:588:VAL:O	1:A:596:TYR:HD2	1.84	0.59
1:A:264:ILE:HG23	1:A:601:LEU:HD13	1.83	0.59
1:A:577:ARG:HB2	1:A:596:TYR:CE1	2.37	0.59
1:A:182:VAL:HG23	1:A:182:VAL:O	2.02	0.59
1:A:307:GLN:HE21	1:A:426:ARG:NE	2.01	0.58
1:A:354:ASN:OD1	1:A:356:ILE:HD12	2.04	0.57
1:A:381:ARG:HA	1:A:385:GLU:HB2	1.84	0.57
1:A:394:HIS:CD2	1:A:395:TYR:N	2.72	0.57
1:A:397:ILE:CD1	1:A:414:MET:HE2	2.29	0.57
1:A:363:ARG:NH1	1:A:390:LYS:O	2.37	0.57
1:A:611:ASP:OD1	1:A:612:ASN:N	2.37	0.57
1:A:414:MET:CE	1:A:470:MET:SD	2.93	0.56
1:A:26:ALA:HB2	1:A:219:LEU:HD13	1.87	0.56
1:A:394:HIS:CB	1:A:421:GLU:OE2	2.54	0.56
1:A:286:ILE:HG23	1:A:287:LYS:N	2.21	0.56
1:A:7:PHE:CE1	1:A:220:TYR:HD2	2.22	0.55
1:A:29:ASP:O	1:A:30:HIS:HB2	2.05	0.55
1:A:132:TYR:CD2	1:A:136:CYS:HB2	2.42	0.55
1:A:326:ASP:OD1	1:A:405:HIS:CD2	2.60	0.54
1:A:24:ARG:NH2	1:A:147:PRO:HG2	2.22	0.54
1:A:416:MET:CE	1:A:467:ALA:HB2	2.37	0.54
1:A:162:ASP:OD2	1:A:472:HIS:CD2	2.61	0.54
1:A:264:ILE:CG2	1:A:601:LEU:HD13	2.39	0.54
1:A:114:ALA:HB3	1:A:117:TRP:HB2	1.90	0.53
1:A:188:GLU:HG2	1:A:189:HIS:N	2.23	0.53
1:A:416:MET:HE3	1:A:467:ALA:HB2	1.89	0.53
1:A:414:MET:HE1	1:A:470:MET:SD	2.48	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:397:ILE:HD12	1:A:414:MET:CE	2.34	0.53
1:A:7:PHE:HE1	1:A:220:TYR:CG	2.27	0.53
1:A:408:ILE:CG1	1:A:409:PRO:HD2	2.39	0.53
1:A:567:HIS:HD2	1:A:617:THR:N	2.07	0.52
1:A:15:GLY:HA3	3:A:702:FAD:O5B	2.08	0.52
1:A:133:GLN:HB2	1:A:151:SER:HB3	1.90	0.52
1:A:602:LYS:HZ1	1:A:631:LEU:HB3	1.73	0.52
1:A:463:SER:HA	1:A:466:THR:HG22	1.91	0.52
1:A:569:LEU:HB3	1:A:606:LEU:HD23	1.92	0.51
1:A:601:LEU:HD12	1:A:602:LYS:N	2.25	0.51
1:A:425:SER:OG	1:A:445:GLY:O	2.23	0.51
1:A:39:ALA:HB2	1:A:226:LYS:HA	1.93	0.51
1:A:241:ARG:HB3	1:A:257:ILE:HG12	1.92	0.51
1:A:85:PRO:HB2	1:A:566:TRP:HE1	1.75	0.51
1:A:10:LEU:HD12	1:A:260:ALA:HB2	1.93	0.51
1:A:47:PRO:O	1:A:51:LEU:HG	2.10	0.51
1:A:577:ARG:HB2	1:A:596:TYR:CD1	2.46	0.51
1:A:102:THR:HG22	1:A:616:ASN:HD22	1.76	0.51
1:A:163:PHE:HE2	1:A:360:VAL:HG23	1.75	0.51
1:A:409:PRO:O	1:A:413:TYR:OH	2.23	0.51
1:A:228:ASP:OD1	1:A:436:PRO:HG3	2.10	0.51
1:A:192:LYS:NZ	1:A:357:GLU:OE2	2.40	0.51
1:A:215:ASN:OD1	1:A:648:PRO:CB	2.59	0.51
1:A:414:MET:HG3	1:A:414:MET:O	2.10	0.50
1:A:318:ILE:HG22	1:A:412:LYS:C	2.32	0.50
1:A:522:LEU:HD21	1:A:541:HIS:CE1	2.47	0.50
1:A:613:VAL:O	1:A:613:VAL:HG23	2.11	0.50
1:A:342:GLN:NE2	1:A:348:THR:O	2.45	0.50
1:A:51:LEU:HD11	1:A:54:ILE:HD12	1.94	0.50
1:A:273:SER:HB2	1:A:605:ASP:OD2	2.11	0.50
1:A:179:GLU:O	1:A:181:LEU:HD12	2.12	0.49
1:A:264:ILE:HG23	1:A:601:LEU:CD1	2.41	0.49
1:A:132:TYR:CE2	1:A:136:CYS:HB2	2.47	0.49
1:A:215:ASN:OD1	1:A:648:PRO:CG	2.60	0.49
1:A:215:ASN:ND2	1:A:215:ASN:N	2.60	0.49
1:A:12:LEU:HD12	1:A:266:LEU:HD23	1.93	0.49
1:A:28:LEU:HD12	1:A:29:ASP:N	2.28	0.49
1:A:366:PRO:HA	1:A:369:LEU:HB3	1.94	0.49
1:A:456:MET:HB3	1:A:558:ILE:HD13	1.93	0.49
1:A:354:ASN:OD1	1:A:355:GLY:N	2.45	0.48
1:A:590:ASP:OD1	1:A:591:HIS:N	2.46	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:350:PRO:O	1:A:353:THR:HG22	2.13	0.48
1:A:393:MET:HA	1:A:421:GLU:HG2	1.96	0.48
1:A:606:LEU:HD11	1:A:617:THR:HB	1.96	0.48
1:A:102:THR:HG23	1:A:616:ASN:H	1.79	0.48
1:A:616:ASN:HB3	3:A:702:FAD:C2	2.44	0.48
1:A:134:ARG:NH1	1:A:176:ASP:O	2.47	0.47
1:A:567:HIS:CD2	1:A:616:ASN:HA	2.49	0.47
1:A:28:LEU:HB3	1:A:645:MET:HG2	1.97	0.47
1:A:7:PHE:CE1	1:A:220:TYR:HB2	2.50	0.47
1:A:38:GLU:OE1	1:A:39:ALA:C	2.53	0.47
1:A:11:VAL:O	1:A:11:VAL:HG12	2.14	0.47
1:A:482:PRO:HG2	1:A:484:PHE:CE2	2.48	0.47
1:A:131:THR:CG2	1:A:149:LYS:HG2	2.45	0.47
1:A:12:LEU:HD12	1:A:266:LEU:CD2	2.44	0.47
1:A:330:ARG:NH2	1:A:411:GLY:O	2.48	0.47
1:A:285:PRO:O	1:A:289:ARG:HG3	2.14	0.47
1:A:30:HIS:NE2	1:A:216:HIS:NE2	2.62	0.47
1:A:54:ILE:O	1:A:54:ILE:HG22	2.15	0.47
1:A:572:CYS:HG	1:A:607:SER:HG	1.54	0.47
1:A:6:GLU:OE2	1:A:261:ARG:HG2	2.15	0.46
1:A:424:PHE:HD2	1:A:451:ARG:HH11	1.62	0.46
1:A:7:PHE:CD1	1:A:35:GLY:HA3	2.49	0.46
1:A:103:ARG:HH12	1:A:141:ILE:HG22	1.80	0.46
1:A:264:ILE:HG22	1:A:601:LEU:HA	1.96	0.46
1:A:29:ASP:O	1:A:31:SER:N	2.43	0.46
1:A:162:ASP:OD2	1:A:472:HIS:NE2	2.49	0.46
1:A:525:PRO:HG3	1:A:541:HIS:CE1	2.51	0.46
1:A:230:ILE:HD11	1:A:266:LEU:HD21	1.98	0.46
1:A:429:ILE:HG13	1:A:442:PHE:HB3	1.98	0.46
1:A:44:LEU:HB2	1:A:210:HIS:CD2	2.51	0.46
1:A:567:HIS:CD2	1:A:617:THR:HG23	2.50	0.45
1:A:173:PRO:HD2	1:A:185:HIS:HB3	1.99	0.45
1:A:313:PHE:CE1	1:A:417:PHE:CD1	3.05	0.45
1:A:286:ILE:CG2	1:A:287:LYS:N	2.79	0.45
1:A:601:LEU:HD12	1:A:602:LYS:H	1.82	0.45
1:A:215:ASN:HD22	1:A:215:ASN:N	2.13	0.45
1:A:383:TYR:HD2	1:A:384:PHE:CE2	2.34	0.45
1:A:263:GLN:HG3	1:A:264:ILE:N	2.32	0.45
1:A:517:SER:O	1:A:517:SER:OG	2.33	0.45
1:A:104:GLY:HA2	1:A:615:CYS:HB3	1.98	0.45
1:A:307:GLN:O	1:A:568:CYS:HA	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:328:PHE:CE1	1:A:336:GLN:HG3	2.52	0.44
1:A:408:ILE:HG12	1:A:409:PRO:CD	2.45	0.44
1:A:112:PHE:CD1	1:A:610:PRO:HD2	2.52	0.44
1:A:131:THR:HG22	1:A:149:LYS:HG2	1.98	0.44
1:A:618:TYR:CD1	1:A:618:TYR:O	2.70	0.44
1:A:567:HIS:HB3	1:A:617:THR:HG23	1.99	0.44
1:A:491:ARG:HG2	1:A:492:ALA:N	2.33	0.44
1:A:299:LEU:HD22	1:A:588:VAL:HG12	1.99	0.44
1:A:414:MET:HE3	1:A:470:MET:SD	2.57	0.44
1:A:7:PHE:CE2	1:A:222:ILE:HD11	2.53	0.44
1:A:558:ILE:O	1:A:562:THR:HG23	2.18	0.44
1:A:47:PRO:HA	1:A:50:TYR:CE2	2.53	0.43
1:A:494:GLU:HG3	1:A:519:THR:HB	1.99	0.43
1:A:54:ILE:CG2	1:A:57:ARG:HD3	2.48	0.43
1:A:308:ASP:O	1:A:423:PRO:HG3	2.18	0.43
1:A:424:PHE:O	1:A:426:ARG:NH1	2.51	0.43
1:A:606:LEU:HD12	1:A:621:ALA:HB2	1.99	0.43
1:A:404:ASP:OD2	1:A:407:LYS:HD3	2.17	0.43
1:A:363:ARG:NH2	1:A:387:LYS:O	2.52	0.43
1:A:485:PRO:HG2	1:A:488:SER:HB2	2.01	0.43
1:A:390:LYS:HA	1:A:391:PRO:HD2	1.82	0.42
1:A:545:GLU:OE1	1:A:545:GLU:N	2.51	0.42
1:A:588:VAL:O	1:A:596:TYR:CD2	2.69	0.42
1:A:270:THR:HA	1:A:607:SER:HB3	2.00	0.42
1:A:419:PHE:CD1	1:A:419:PHE:O	2.72	0.42
1:A:598:VAL:HG12	1:A:599:LYS:N	2.34	0.42
1:A:8:ASP:OD1	1:A:262:LYS:HE3	2.19	0.42
1:A:23:GLY:O	1:A:27:ASN:OD1	2.38	0.42
1:A:33:LYS:HD3	1:A:220:TYR:HE2	1.84	0.42
1:A:508:ASN:OD1	1:A:508:ASN:O	2.37	0.42
1:A:215:ASN:OD1	1:A:648:PRO:HG3	2.19	0.42
1:A:15:GLY:CA	3:A:702:FAD:O5B	2.67	0.42
1:A:149:LYS:HD2	1:A:195:ASN:ND2	2.33	0.42
1:A:316:TYR:HE2	1:A:416:MET:HG3	1.83	0.42
1:A:264:ILE:HG23	1:A:601:LEU:HA	2.01	0.42
1:A:655:TYR:CG	1:A:662:ARG:HD3	2.55	0.42
1:A:365:THR:HG22	1:A:367:GLU:H	1.85	0.42
1:A:387:LYS:HA	1:A:388:PRO:HD3	1.76	0.42
1:A:9:ILE:HB	1:A:34:VAL:HG22	2.02	0.42
1:A:310:TYR:CE2	1:A:423:PRO:HB3	2.54	0.42
1:A:51:LEU:HD13	1:A:351:LEU:HD11	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:37:ILE:HG12	1:A:222:ILE:HD12	2.02	0.41
1:A:364:PRO:HD3	1:A:392:VAL:HG21	2.01	0.41
1:A:356:ILE:H	1:A:356:ILE:CD1	2.29	0.41
1:A:364:PRO:HG2	1:A:369:LEU:HD12	2.03	0.41
1:A:268:CYS:C	3:A:702:FAD:H52A	2.41	0.41
1:A:51:LEU:O	1:A:51:LEU:HD12	2.20	0.41
1:A:215:ASN:OD1	1:A:648:PRO:HB2	2.21	0.41
1:A:136:CYS:SG	1:A:138:ASN:O	2.79	0.41
1:A:414:MET:CG	1:A:414:MET:O	2.69	0.41
1:A:627:LYS:HD2	1:A:627:LYS:HA	1.92	0.41
1:A:77:LEU:HB3	1:A:80:ARG:HB3	2.02	0.41
1:A:233:GLU:OE1	1:A:233:GLU:N	2.54	0.41
1:A:563:GLU:OE1	1:A:563:GLU:N	2.54	0.41
1:A:188:GLU:CG	1:A:189:HIS:N	2.84	0.41
1:A:54:ILE:HG22	1:A:57:ARG:HD3	2.03	0.41
1:A:132:TYR:OH	1:A:134:ARG:HD3	2.21	0.40
1:A:330:ARG:HH22	1:A:411:GLY:H	1.69	0.40
1:A:42:ASN:OD1	1:A:43:ASN:N	2.54	0.40
1:A:163:PHE:CE2	1:A:360:VAL:HG23	2.56	0.40
1:A:55:TYR:N	1:A:56:PRO:HD2	2.36	0.40
1:A:104:GLY:CA	1:A:615:CYS:HB3	2.51	0.40
1:A:482:PRO:HB3	1:A:557:TYR:HB2	2.03	0.40
1:A:262:LYS:HB3	1:A:262:LYS:HE2	1.94	0.40
1:A:330:ARG:N	1:A:331:GLY:HA2	2.37	0.40
1:A:567:HIS:NE2	1:A:616:ASN:HA	2.36	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 PDB entries followed by that with respect to all EM entries.

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	660/663 (100%)	626 (95%)	34 (5%)	0	100	100

There are no Ramachandran outliers to report.

### 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 PDB entries followed by that with respect to all EM entries.

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	560/561 (100%)	555 (99%)	5 (1%)	84 93

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	38	GLU
1	A	190	TRP
1	A	354	ASN
1	A	418	HIS
1	A	515	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	30	HIS
1	A	142	HIS
1	A	394	HIS
1	A	430	HIS
1	A	508	ASN
1	A	532	HIS
1	A	541	HIS
1	A	616	ASN

### 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 [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
3	FAD	A	702	-	52,58,58	1.21	4 (7%)	52,89,89	2.52	14 (26%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	FAD	A	702	-	-	0/30/50/50	0/6/6/6

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	702	FAD	C5X-N5	2.47	1.39	1.35
3	A	702	FAD	C2A-N3A	2.59	1.36	1.32
3	A	702	FAD	C4-N3	2.71	1.37	1.33
3	A	702	FAD	C4X-N5	3.73	1.39	1.33

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	702	FAD	N3A-C2A-N1A	-10.48	120.64	128.87
3	A	702	FAD	C8M-C8-C9	-5.06	106.03	120.33
3	A	702	FAD	N3-C2-N1	-4.81	119.59	127.69
3	A	702	FAD	C1B-N9A-C4A	-4.30	122.01	126.81
3	A	702	FAD	C4-C4X-C10	-3.17	117.91	119.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	702	FAD	C4X-C4-N3	-2.51	120.25	123.52
3	A	702	FAD	O3B-C3B-C4B	-2.38	103.90	111.01
3	A	702	FAD	N6A-C6A-N1A	-2.21	114.80	118.52
3	A	702	FAD	O3'-C3'-C4'	-2.02	103.48	108.73
3	A	702	FAD	C4-C4X-N5	2.28	121.47	118.70
3	A	702	FAD	C1'-N10-C9A	2.47	121.70	118.83
3	A	702	FAD	C8M-C8-C7	3.75	128.80	120.73
3	A	702	FAD	C5X-C9A-N10	4.87	121.23	117.58
3	A	702	FAD	C4-N3-C2	6.73	120.77	115.16

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	702	FAD	5	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.