



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:03 AM GMT

PDB ID : 2ZAF
Title : Mechanistic and Structural Analyses of the Roles of Arg409 and Asp402 in the Reaction of the Flavoprotein Nitroalkane Oxidase
Authors : Fitzpatrick, P.F.; Bozinovski, D.M.; Heroux, A.; Shaw, P.G.; Valley, M.P.; Orville, A.M.
Deposited on : 2007-10-05
Resolution : 2.50 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

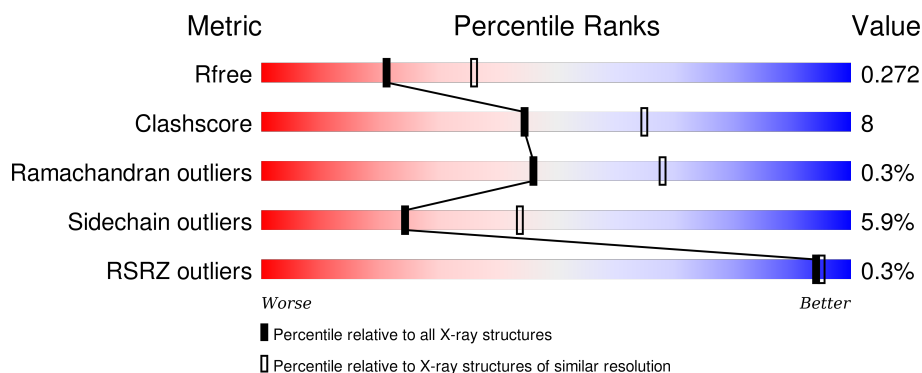
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)

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	439	<div> <div>83%</div> <div>12%</div> <div>• •</div> </div>
1	B	439	<div> <div>81%</div> <div>16%</div> <div>• •</div> </div>
1	C	439	<div> <div>%</div> <div>78%</div> <div>17%</div> <div>• •</div> </div>
1	D	439	<div> <div>79%</div> <div>16%</div> <div>• •</div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 13440 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 Nitroalkane oxidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	430	Total	C	N	O	S	0	1	0
			3307	2096	565	625	21			
1	B	430	Total	C	N	O	S	0	1	0
			3307	2096	565	625	21			
1	C	430	Total	C	N	O	S	0	1	0
			3307	2096	565	625	21			
1	D	430	Total	C	N	O	S	0	1	0
			3307	2096	565	625	21			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	409	LYS	ARG	ENGINEERED	UNP Q8X1D8
B	409	LYS	ARG	ENGINEERED	UNP Q8X1D8
C	409	LYS	ARG	ENGINEERED	UNP Q8X1D8
D	409	LYS	ARG	ENGINEERED	UNP Q8X1D8

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

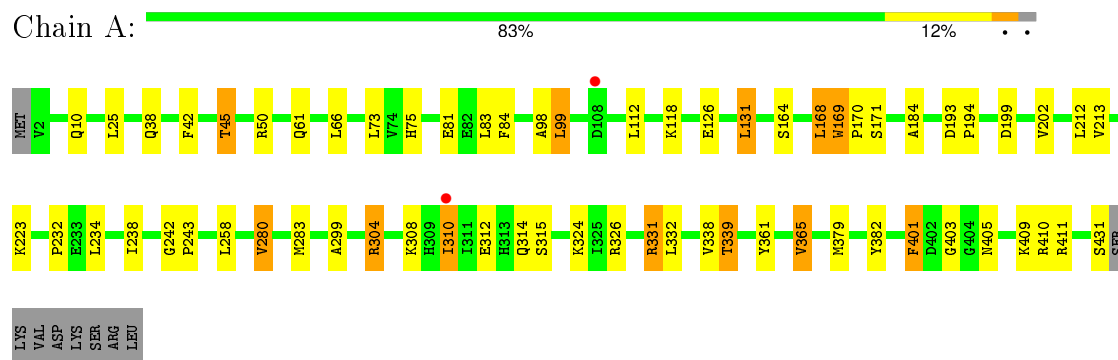


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	C	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	D	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

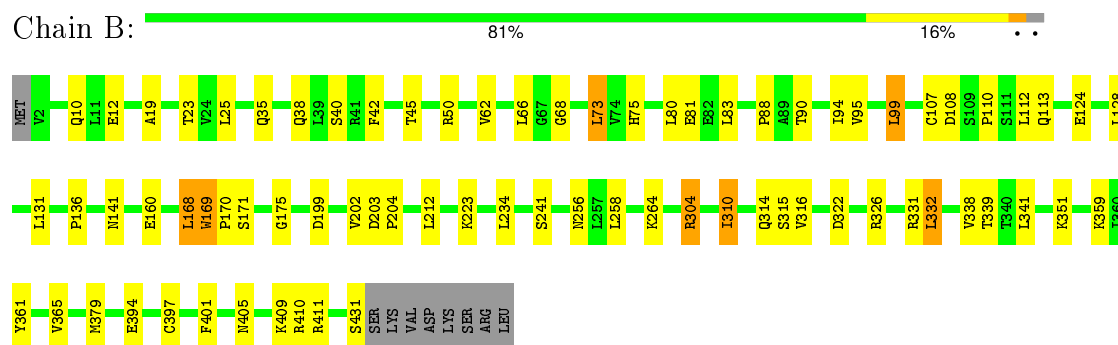
3 Residue-property plots

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.

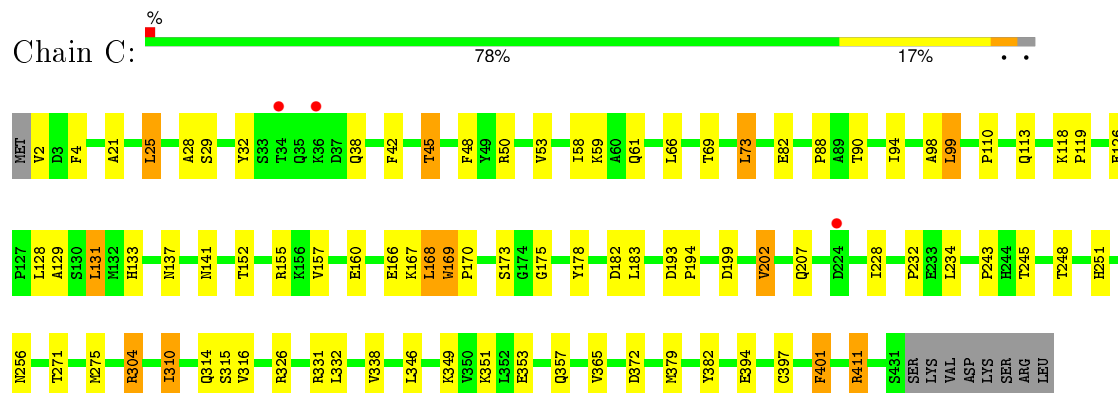
• Molecule 1: Nitroalkane oxidase



• Molecule 1: Nitroalkane oxidase



• Molecule 1: Nitroalkane oxidase



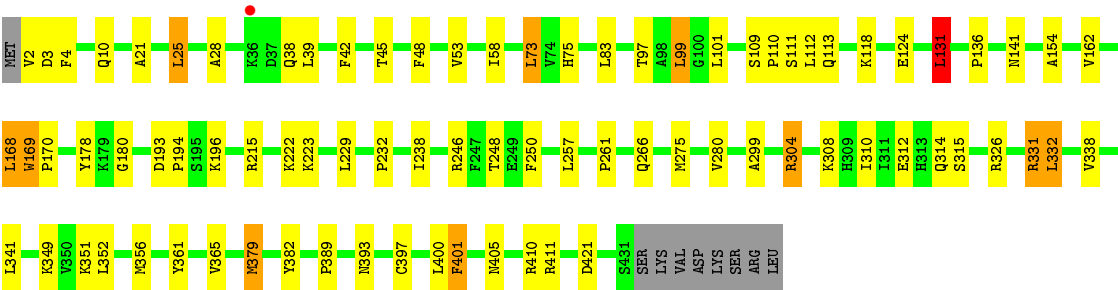
● Molecule 1: Nitroalkane oxidase

Chain D:

79%

16%

••



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	108.32Å 108.32Å 340.01Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.00 – 2.50 72.27 – 2.50	Depositor EDS
% Data completeness (in resolution range)	91.0 (50.00-2.50) 91.0 (72.27-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.13	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.69 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.228 , 0.277 0.221 , 0.272	Depositor DCC
R_{free} test set	3776 reflections (5.39%)	DCC
Wilson B-factor (Å ²)	46.6	Xtriage
Anisotropy	0.088	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 21.6	EDS
Estimated twinning fraction	0.024 for -h,-k,l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 73886 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	13440	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.69% 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	0.47	0/3382	0.63	1/4589 (0.0%)
1	B	0.48	0/3382	0.62	0/4589
1	C	0.50	0/3382	0.63	1/4589 (0.0%)
1	D	0.52	0/3382	0.64	2/4589 (0.0%)
All	All	0.49	0/13528	0.63	4/18356 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	331	ARG	NE-CZ-NH2	-6.62	116.99	120.30
1	D	331	ARG	NE-CZ-NH2	-6.03	117.28	120.30
1	A	331	ARG	NE-CZ-NH2	-5.86	117.37	120.30
1	D	131	LEU	CA-CB-CG	5.33	127.55	115.30

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	3307	0	3320	53	0
1	B	3307	0	3320	53	0
1	C	3307	0	3320	71	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	3307	0	3320	70	0
2	A	53	0	31	2	0
2	B	53	0	31	12	0
2	C	53	0	31	15	0
2	D	53	0	31	8	0
All	All	13440	0	13404	213	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 8.

All (213) 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:379:MET:CE	2:B:500:FAD:HM81	1.70	1.20
2:C:500:FAD:H9	1:D:379:MET:HE3	1.17	1.17
1:C:379:MET:HE2	2:D:500:FAD:H3'	1.26	1.13
1:C:379:MET:CE	1:D:401:PHE:HA	1.83	1.08
1:C:379:MET:HE3	1:D:401:PHE:HA	1.09	1.05
1:A:379:MET:HE2	2:B:500:FAD:HM81	1.41	1.01
1:A:401:PHE:HA	1:B:379:MET:CE	1.92	0.99
1:A:379:MET:CE	2:B:500:FAD:C8M	2.43	0.95
2:C:500:FAD:C9	1:D:379:MET:HE3	1.97	0.94
1:D:141:ASN:HD21	2:D:500:FAD:H61A	1.12	0.94
2:C:500:FAD:H9	1:D:379:MET:CE	1.97	0.93
1:B:141:ASN:HD21	2:B:500:FAD:H61A	1.19	0.91
1:C:379:MET:HE3	1:D:401:PHE:CA	2.00	0.90
2:C:500:FAD:C9	1:D:379:MET:CE	2.51	0.89
1:A:401:PHE:CA	1:B:379:MET:HE1	2.04	0.88
1:D:405:ASN:O	1:D:410:ARG:HG3	1.77	0.84
1:C:326:ARG:NH2	1:C:372:ASP:OD2	2.12	0.83
1:B:168:LEU:C	1:B:170:PRO:HD3	1.99	0.82
1:C:379:MET:HE2	2:D:500:FAD:C3'	2.09	0.80
1:C:379:MET:CE	2:D:500:FAD:H3'	2.09	0.76
1:D:73:LEU:HB3	1:D:338:VAL:HB	1.67	0.76
1:A:73:LEU:HB3	1:A:338:VAL:HB	1.68	0.76
1:C:199:ASP:HB3	1:C:202:VAL:CG1	2.16	0.75
2:C:500:FAD:H2'	1:D:379:MET:HE2	1.69	0.74
1:D:168:LEU:C	1:D:170:PRO:HD3	2.08	0.74
1:D:141:ASN:ND2	2:D:500:FAD:H61A	1.85	0.74
1:A:401:PHE:CA	1:B:379:MET:CE	2.62	0.73
1:D:42:PHE:O	1:D:45:THR:HB	1.89	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:38:GLN:NE2	1:B:234:LEU:H	1.87	0.71
2:C:500:FAD:C8A	1:D:310:ILE:HD11	2.21	0.71
1:A:379:MET:HE1	2:B:500:FAD:C8M	2.20	0.70
1:C:304:ARG:HG2	2:D:500:FAD:O5B	1.91	0.70
1:C:379:MET:HE1	1:D:400:LEU:O	1.90	0.70
1:A:315:SER:H	1:D:314:GLN:HE21	1.41	0.69
1:B:128:LEU:HD12	1:B:175:GLY:HA2	1.76	0.68
1:A:168:LEU:C	1:A:170:PRO:HD2	2.15	0.67
1:A:326:ARG:HB3	1:A:365:VAL:HG13	1.76	0.67
1:C:128:LEU:HD12	1:C:175:GLY:HA2	1.76	0.67
1:B:315:SER:H	1:C:314:GLN:HE21	1.42	0.66
2:C:500:FAD:O5B	1:D:304:ARG:HG2	1.96	0.66
1:D:99:LEU:HD22	1:D:131:LEU:HD23	1.78	0.65
1:B:314:GLN:HB2	1:C:314:GLN:HB2	1.78	0.65
1:C:379:MET:HE1	1:D:400:LEU:C	2.19	0.64
1:B:73:LEU:HB3	1:B:338:VAL:HB	1.80	0.64
1:B:304:ARG:HB2	1:B:310:ILE:HD13	1.80	0.63
1:A:379:MET:HE3	2:B:500:FAD:HM81	1.71	0.63
1:A:99:LEU:HD13	1:A:131:LEU:HB3	1.80	0.63
1:A:401:PHE:CB	1:B:379:MET:HE1	2.29	0.62
1:B:35:GLN:HE21	1:B:40:SER:HB3	1.65	0.62
1:B:38:GLN:HE21	1:B:234:LEU:H	1.48	0.61
1:D:97:THR:O	1:D:101:LEU:HG	2.01	0.60
1:B:42:PHE:O	1:B:45:THR:HB	2.02	0.60
1:A:112:LEU:HD23	1:A:258:LEU:HD12	1.83	0.60
1:A:379:MET:HE2	2:B:500:FAD:H9	1.83	0.59
1:A:42:PHE:O	1:A:45:THR:HB	2.03	0.59
1:C:379:MET:CE	1:D:401:PHE:CA	2.71	0.59
1:A:315:SER:H	1:D:314:GLN:NE2	2.01	0.59
1:A:401:PHE:CB	1:B:379:MET:CE	2.81	0.58
1:D:326:ARG:HB3	1:D:365:VAL:HG13	1.83	0.58
1:A:401:PHE:HA	1:B:379:MET:HE3	1.85	0.58
1:D:28:ALA:HA	1:D:48:PHE:CZ	2.38	0.58
1:C:99:LEU:HD22	1:C:131:LEU:HD23	1.85	0.57
1:C:73:LEU:HB3	1:C:338:VAL:HB	1.85	0.57
1:B:10:GLN:HG3	1:B:75:HIS:CE1	2.40	0.57
2:C:500:FAD:C8M	1:D:379:MET:HE1	2.34	0.56
1:C:199:ASP:HB3	1:C:202:VAL:HG12	1.85	0.56
1:C:168:LEU:C	1:C:170:PRO:HD2	2.26	0.56
1:C:38:GLN:NE2	1:C:234:LEU:H	2.03	0.56
1:C:173:SER:OG	1:C:243:PRO:HD2	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:500:FAD:HM81	1:B:379:MET:SD	2.45	0.56
1:B:199:ASP:HB3	1:B:202:VAL:HG12	1.87	0.56
2:C:500:FAD:C9	1:D:379:MET:HE1	2.35	0.56
1:D:352:LEU:O	1:D:356:MET:HG2	2.05	0.56
1:D:38:GLN:HE22	1:D:238:ILE:HA	1.70	0.55
1:A:401:PHE:N	1:B:379:MET:HE1	2.21	0.55
1:C:310:ILE:HG23	1:C:316:VAL:HG11	1.89	0.55
1:D:21:ALA:HA	1:D:25:LEU:HB2	1.88	0.55
1:D:361:TYR:O	1:D:365:VAL:HB	2.07	0.55
1:B:110:PRO:HA	1:B:113:GLN:HE21	1.72	0.55
1:B:81:GLU:OE2	1:B:331:ARG:HD3	2.07	0.54
1:D:275:MET:HG2	1:D:341:LEU:HD13	1.88	0.54
1:C:82:GLU:HA	1:C:82:GLU:OE1	2.07	0.54
1:D:308:LYS:HD2	1:D:312:GLU:HB2	1.88	0.54
1:C:310:ILE:CG2	1:C:316:VAL:HG11	2.38	0.54
1:C:401:PHE:HB2	1:D:379:MET:CE	2.38	0.54
1:D:308:LYS:HB2	1:D:312:GLU:HG3	1.89	0.54
1:C:50:ARG:HD2	1:C:126:GLU:OE1	2.08	0.53
1:A:61:GLN:HE21	1:A:98:ALA:HB2	1.73	0.52
1:C:42:PHE:O	1:C:45:THR:HB	2.10	0.52
1:D:109:SER:OG	1:D:112:LEU:HB2	2.10	0.52
1:B:141:ASN:ND2	2:B:500:FAD:H61A	1.99	0.52
1:C:379:MET:CE	2:D:500:FAD:C3'	2.78	0.52
1:A:314:GLN:HE21	1:D:315:SER:H	1.58	0.51
1:A:361:TYR:O	1:A:365:VAL:HB	2.10	0.51
1:C:178:TYR:CG	1:C:228:ILE:HD12	2.44	0.51
1:D:168:LEU:O	1:D:169:TRP:HB2	2.10	0.51
1:D:162:VAL:HA	1:D:250:PHE:O	2.10	0.51
1:B:361:TYR:O	1:B:365:VAL:HB	2.10	0.51
1:A:379:MET:CE	2:B:500:FAD:H9	2.40	0.50
1:C:59:LYS:HE2	1:C:69:THR:HG23	1.92	0.50
1:C:141:ASN:HD21	2:C:500:FAD:H61A	1.58	0.50
1:D:299:ALA:HA	1:D:310:ILE:HG22	1.94	0.50
1:C:53:VAL:HG22	1:C:58:ILE:HG13	1.93	0.50
1:D:389:PRO:O	1:D:393:ASN:ND2	2.43	0.50
1:C:401:PHE:HB2	1:D:379:MET:HE1	1.94	0.50
1:C:168:LEU:C	1:C:170:PRO:CD	2.80	0.50
1:B:405:ASN:O	1:B:410:ARG:HG3	2.12	0.49
2:C:500:FAD:C8	1:D:379:MET:HE1	2.42	0.49
1:C:88:PRO:HG2	1:C:394:GLU:OE2	2.12	0.49
1:A:38:GLN:HG3	1:A:232:PRO:O	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:160:GLU:OE2	1:C:251:HIS:ND1	2.38	0.49
1:A:50:ARG:HD2	1:A:126:GLU:OE1	2.13	0.49
1:A:403:GLY:HA3	1:A:409:LYS:NZ	2.26	0.49
1:B:310:ILE:HG23	1:B:316:VAL:HG11	1.95	0.49
1:C:45:THR:HG23	1:C:90:THR:OG1	2.13	0.49
1:C:207:GLN:NE2	1:C:207:GLN:HA	2.29	0.48
1:C:353:GLU:O	1:C:357:GLN:HG3	2.13	0.48
1:C:169:TRP:N	1:C:170:PRO:HD2	2.27	0.48
1:C:168:LEU:O	1:C:170:PRO:HD3	2.13	0.48
1:C:61:GLN:HE21	1:C:98:ALA:HB2	1.79	0.48
1:C:21:ALA:HA	1:C:25:LEU:HB2	1.95	0.48
1:C:29:SER:HA	1:C:32:TYR:CE1	2.48	0.48
1:B:314:GLN:HE21	1:C:315:SER:H	1.61	0.48
1:B:171:SER:HA	1:B:241:SER:O	2.13	0.48
1:A:38:GLN:NE2	1:A:234:LEU:H	2.11	0.48
1:B:112:LEU:CD2	1:B:258:LEU:HD12	2.45	0.47
1:B:169:TRP:N	1:B:170:PRO:HD3	2.28	0.47
1:D:168:LEU:O	1:D:170:PRO:HD3	2.14	0.47
1:D:38:GLN:HG3	1:D:232:PRO:O	2.13	0.47
1:A:405:ASN:O	1:A:410:ARG:HG3	2.13	0.47
1:B:326:ARG:HB3	1:B:365:VAL:HG13	1.96	0.47
1:D:136:PRO:HA	1:D:168:LEU:HD12	1.97	0.47
1:A:304:ARG:HG2	2:B:500:FAD:O5B	2.15	0.47
1:C:346:LEU:O	1:C:351:LYS:HE3	2.15	0.47
2:C:500:FAD:C2'	1:D:379:MET:HE2	2.42	0.46
1:C:411:ARG:HA	1:C:411:ARG:HE	1.80	0.46
1:D:193:ASP:HA	1:D:194:PRO:HD2	1.77	0.46
1:B:332:LEU:HD12	1:B:332:LEU:HA	1.77	0.46
1:A:401:PHE:HA	1:B:379:MET:HE1	1.66	0.46
1:C:271:THR:O	1:C:275:MET:HG3	2.16	0.46
1:A:199:ASP:HB3	1:A:202:VAL:HG12	1.97	0.46
1:B:19:ALA:O	1:B:23:THR:OG1	2.25	0.45
1:A:299:ALA:HA	1:A:310:ILE:HG22	1.98	0.45
1:A:308:LYS:HD2	1:A:312:GLU:HB2	1.98	0.45
1:A:38:GLN:HE22	1:A:238:ILE:HA	1.81	0.45
1:B:12:GLU:HA	1:B:12:GLU:OE2	2.17	0.45
1:C:38:GLN:HG3	1:C:232:PRO:O	2.17	0.45
1:D:421:ASP:N	1:D:421:ASP:OD1	2.50	0.45
1:D:169:TRP:N	1:D:170:PRO:HD3	2.32	0.45
1:A:84:PHE:CE1	1:A:283:MET:HG2	2.52	0.45
1:A:314:GLN:HB2	1:D:314:GLN:HB2	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:137:ASN:N	1:C:137:ASN:HD22	2.13	0.44
1:C:155:ARG:HE	1:C:157:VAL:CG2	2.30	0.44
1:C:304:ARG:HD3	2:D:500:FAD:H51A	1.98	0.44
1:D:53:VAL:HG22	1:D:58:ILE:HG13	1.98	0.44
1:D:110:PRO:HA	1:D:113:GLN:HE21	1.82	0.44
1:A:339:THR:HG21	1:C:4:PHE:HD1	1.82	0.44
1:A:184:ALA:HB3	1:A:213:VAL:HB	1.99	0.44
1:D:154:ALA:HA	1:D:162:VAL:O	2.18	0.44
1:B:62:VAL:HB	1:B:68:GLY:HA3	1.99	0.44
1:D:341:LEU:HD23	1:D:351:LYS:HB3	2.00	0.44
1:D:73:LEU:HA	1:D:73:LEU:HD12	1.86	0.44
1:A:379:MET:CE	2:B:500:FAD:C9	2.96	0.43
1:C:133:HIS:HB3	2:C:500:FAD:O2	2.18	0.43
1:A:403:GLY:HA3	1:A:409:LYS:HZ1	1.82	0.43
1:B:88:PRO:HG2	1:B:394:GLU:OE2	2.18	0.43
1:C:166:GLU:HA	1:C:245:THR:O	2.18	0.43
1:A:193:ASP:HA	1:A:194:PRO:HD2	1.86	0.43
2:A:500:FAD:H3'	1:B:379:MET:HE2	2.00	0.43
1:C:118:LYS:HB3	1:C:119:PRO:HD3	2.01	0.43
1:C:382:TYR:CZ	1:D:397:CYS:HB2	2.53	0.43
1:A:409:LYS:HA	1:A:409:LYS:HD3	1.90	0.43
1:D:180:GLY:O	1:D:215:ARG:NH1	2.44	0.43
1:D:10:GLN:HG3	1:D:75:HIS:CE1	2.53	0.43
1:A:379:MET:HE1	2:B:500:FAD:C8	2.49	0.42
1:D:257:LEU:HD21	1:D:261:PRO:HD3	2.01	0.42
1:C:141:ASN:ND2	2:C:500:FAD:H61A	2.17	0.42
1:A:10:GLN:HG3	1:A:75:HIS:CE1	2.53	0.42
1:A:81:GLU:OE2	1:A:331:ARG:HD3	2.19	0.42
1:B:136:PRO:HA	1:B:168:LEU:HD12	2.00	0.42
1:A:169:TRP:N	1:A:170:PRO:HD2	2.34	0.42
2:C:500:FAD:HM81	1:D:379:MET:CE	2.49	0.42
1:B:341:LEU:HD23	1:B:351:LYS:HB3	2.00	0.42
1:A:382:TYR:CZ	1:B:397:CYS:HB2	2.54	0.42
1:A:401:PHE:HB2	1:B:379:MET:HE3	2.01	0.42
1:A:324:LYS:HG3	1:C:357:GLN:NE2	2.34	0.42
1:C:193:ASP:HA	1:C:194:PRO:HD2	1.89	0.42
1:B:203:ASP:HA	1:B:204:PRO:HD3	1.95	0.41
1:D:39:LEU:HD22	1:D:178:TYR:CE2	2.55	0.41
1:C:129:ALA:HA	1:C:183:LEU:O	2.19	0.41
1:C:110:PRO:HA	1:C:113:GLN:HE21	1.85	0.41
1:A:242:GLY:N	1:A:243:PRO:HD3	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:152:THR:OG1	1:C:167:LYS:HD3	2.20	0.41
1:B:314:GLN:NE2	1:C:315:SER:H	2.17	0.41
1:D:229:LEU:HD11	1:D:246:ARG:HB2	2.03	0.41
1:D:331:ARG:O	1:D:332:LEU:C	2.59	0.41
1:B:90:THR:O	1:B:94:ILE:HG13	2.21	0.41
1:B:322:ASP:O	1:B:326:ARG:HG3	2.21	0.41
1:B:339:THR:HG21	1:D:4:PHE:HD1	1.85	0.41
1:C:397:CYS:HB2	1:D:382:TYR:CZ	2.56	0.41
1:C:401:PHE:HA	1:D:379:MET:HE2	2.03	0.41
1:A:401:PHE:CB	1:B:379:MET:HE3	2.49	0.41
1:D:168:LEU:O	1:D:169:TRP:CB	2.69	0.41
1:C:90:THR:O	1:C:94:ILE:HG13	2.21	0.41
1:B:80:LEU:HD22	1:B:95:VAL:CG1	2.51	0.41
1:B:99:LEU:HD22	1:B:131:LEU:HD23	2.03	0.41
1:C:310:ILE:HD12	1:C:310:ILE:HA	1.67	0.40
1:B:128:LEU:CD1	1:B:175:GLY:HA2	2.50	0.40
1:C:169:TRP:N	1:C:170:PRO:CD	2.84	0.40
1:B:359:LYS:HD2	1:B:409:LYS:HG3	2.03	0.40
1:C:28:ALA:HA	1:C:48:PHE:CZ	2.57	0.40
1:D:304:ARG:HB2	1:D:310:ILE:HD13	2.03	0.40
1:C:401:PHE:CB	1:D:379:MET:CE	2.98	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	429/439 (98%)	414 (96%)	13 (3%)	2 (0%)	34	55
1	B	429/439 (98%)	414 (96%)	14 (3%)	1 (0%)	52	75
1	C	429/439 (98%)	412 (96%)	16 (4%)	1 (0%)	52	75
1	D	429/439 (98%)	409 (95%)	19 (4%)	1 (0%)	52	75

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	1716/1756 (98%)	1649 (96%)	62 (4%)	5 (0%)	46 68

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	169	TRP
1	B	169	TRP
1	C	169	TRP
1	D	169	TRP
1	A	280	VAL

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	356/364 (98%)	335 (94%)	21 (6%)	24 44
1	B	356/364 (98%)	335 (94%)	21 (6%)	24 44
1	C	356/364 (98%)	337 (95%)	19 (5%)	28 50
1	D	356/364 (98%)	333 (94%)	23 (6%)	21 39
All	All	1424/1456 (98%)	1340 (94%)	84 (6%)	24 44

All (84) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	LEU
1	A	45	THR
1	A	66	LEU
1	A	83	LEU
1	A	99	LEU
1	A	118	LYS
1	A	131	LEU
1	A	164	SER
1	A	168	LEU
1	A	171	SER

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Mol	Chain	Res	Type
1	A	212	LEU
1	A	223	LYS
1	A	280	VAL
1	A	304	ARG
1	A	310	ILE
1	A	332	LEU
1	A	339	THR
1	A	365	VAL
1	A	401	PHE
1	A	411	ARG
1	A	431	SER
1	B	25	LEU
1	B	50	ARG
1	B	66	LEU
1	B	73	LEU
1	B	83	LEU
1	B	99	LEU
1	B	107	CYS
1	B	108	ASP
1	B	124	GLU
1	B	160	GLU
1	B	168	LEU
1	B	212	LEU
1	B	223	LYS
1	B	256	ASN
1	B	264	LYS
1	B	304	ARG
1	B	310	ILE
1	B	332	LEU
1	B	401	PHE
1	B	411	ARG
1	B	431	SER
1	C	2	VAL
1	C	25	LEU
1	C	45	THR
1	C	66	LEU
1	C	73	LEU
1	C	99	LEU
1	C	131	LEU
1	C	168	LEU
1	C	182	ASP
1	C	202	VAL

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Mol	Chain	Res	Type
1	C	248	THR
1	C	256	ASN
1	C	304	ARG
1	C	310	ILE
1	C	332	LEU
1	C	349	LYS
1	C	365	VAL
1	C	401	PHE
1	C	411	ARG
1	D	2	VAL
1	D	3	ASP
1	D	25	LEU
1	D	73	LEU
1	D	83	LEU
1	D	99	LEU
1	D	111	SER
1	D	118	LYS
1	D	124	GLU
1	D	131	LEU
1	D	168	LEU
1	D	196	LYS
1	D	222	LYS
1	D	223	LYS
1	D	248	THR
1	D	266	GLN
1	D	280	VAL
1	D	304	ARG
1	D	332	LEU
1	D	349	LYS
1	D	379	MET
1	D	401	PHE
1	D	411	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (54) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	22	ASN
1	A	38	GLN
1	A	43	GLN
1	A	61	GLN
1	A	113	GLN
1	A	137	ASN

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Mol	Chain	Res	Type
1	A	141	ASN
1	A	159	ASN
1	A	256	ASN
1	A	266	GLN
1	A	314	GLN
1	B	18	GLN
1	B	22	ASN
1	B	35	GLN
1	B	38	GLN
1	B	43	GLN
1	B	113	GLN
1	B	137	ASN
1	B	141	ASN
1	B	198	GLN
1	B	256	ASN
1	B	313	HIS
1	B	314	GLN
1	C	18	GLN
1	C	22	ASN
1	C	35	GLN
1	C	38	GLN
1	C	43	GLN
1	C	61	GLN
1	C	113	GLN
1	C	137	ASN
1	C	141	ASN
1	C	144	GLN
1	C	198	GLN
1	C	207	GLN
1	C	244	HIS
1	C	256	ASN
1	C	313	HIS
1	C	314	GLN
1	D	22	ASN
1	D	35	GLN
1	D	38	GLN
1	D	43	GLN
1	D	61	GLN
1	D	113	GLN
1	D	137	ASN
1	D	141	ASN
1	D	144	GLN

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Mol	Chain	Res	Type
1	D	198	GLN
1	D	256	ASN
1	D	266	GLN
1	D	313	HIS
1	D	314	GLN
1	D	357	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 ⓘ

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	-	48,58,58	1.32	5 (10%)	54,89,89	2.27	7 (12%)
2	FAD	B	500	-	48,58,58	1.30	5 (10%)	54,89,89	2.04	6 (11%)
2	FAD	C	500	-	48,58,58	1.21	5 (10%)	54,89,89	2.12	7 (12%)
2	FAD	D	500	-	48,58,58	1.27	5 (10%)	54,89,89	2.10	9 (16%)

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	-	-	0/30/50/50	0/6/6/6
2	FAD	B	500	-	-	0/30/50/50	0/6/6/6
2	FAD	C	500	-	-	0/30/50/50	0/6/6/6
2	FAD	D	500	-	-	0/30/50/50	0/6/6/6

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	500	FAD	C2A-N1A	2.40	1.38	1.33
2	C	500	FAD	C2A-N1A	2.47	1.38	1.33
2	B	500	FAD	C2A-N1A	2.50	1.38	1.33
2	A	500	FAD	C4-N3	2.53	1.37	1.33
2	D	500	FAD	C2A-N1A	2.54	1.38	1.33
2	C	500	FAD	C4-N3	2.80	1.38	1.33
2	B	500	FAD	C4-N3	2.81	1.38	1.33
2	D	500	FAD	C4-N3	2.83	1.38	1.33
2	C	500	FAD	C4X-N5	2.89	1.37	1.33
2	B	500	FAD	C4X-N5	2.96	1.38	1.33
2	D	500	FAD	C4X-N5	3.11	1.38	1.33
2	D	500	FAD	C1'-N10	3.17	1.51	1.48
2	C	500	FAD	C1'-N10	3.20	1.51	1.48
2	B	500	FAD	C2A-N3A	3.24	1.37	1.32
2	C	500	FAD	C2A-N3A	3.38	1.38	1.32
2	A	500	FAD	C2A-N3A	3.42	1.38	1.32
2	A	500	FAD	C4X-N5	3.45	1.38	1.33
2	B	500	FAD	C1'-N10	3.65	1.52	1.48
2	D	500	FAD	C2A-N3A	3.68	1.38	1.32
2	A	500	FAD	C1'-N10	3.70	1.52	1.48

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	500	FAD	N3A-C2A-N1A	-13.18	118.80	128.89
2	B	500	FAD	N3A-C2A-N1A	-11.95	119.74	128.89
2	C	500	FAD	N3A-C2A-N1A	-11.51	120.08	128.89
2	D	500	FAD	N3A-C2A-N1A	-10.84	120.59	128.89
2	D	500	FAD	C2B-C1B-N9A	-4.32	107.69	114.29
2	A	500	FAD	P-O3P-PA	-3.39	123.22	132.73
2	C	500	FAD	C4X-C4-N3	-3.05	119.42	123.59
2	D	500	FAD	P-O3P-PA	-2.81	124.84	132.73

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	500	FAD	C2B-C1B-N9A	-2.75	110.08	114.29
2	A	500	FAD	C4X-C4-N3	-2.70	119.90	123.59
2	C	500	FAD	P-O3P-PA	-2.62	125.38	132.73
2	D	500	FAD	C4X-C4-N3	-2.47	120.21	123.59
2	B	500	FAD	C4X-C4-N3	-2.12	120.69	123.59
2	D	500	FAD	C4A-C5A-N7A	-2.08	107.56	109.48
2	C	500	FAD	O2A-PA-O3P	2.09	114.57	105.09
2	D	500	FAD	C5X-C9A-N10	2.44	119.48	117.62
2	C	500	FAD	C5X-C9A-N10	2.76	119.72	117.62
2	D	500	FAD	O3P-P-O5'	2.88	110.58	102.94
2	B	500	FAD	O2A-PA-O3P	2.91	118.27	105.09
2	A	500	FAD	C5X-C9A-N10	2.93	119.85	117.62
2	D	500	FAD	C4X-N5-C5X	2.94	120.14	116.76
2	B	500	FAD	C5X-C9A-N10	3.02	119.91	117.62
2	B	500	FAD	C4X-N5-C5X	3.45	120.73	116.76
2	C	500	FAD	C4X-N5-C5X	3.64	120.95	116.76
2	A	500	FAD	C4X-N5-C5X	3.71	121.03	116.76
2	B	500	FAD	C4-N3-C2	4.32	118.98	115.25
2	D	500	FAD	C4-N3-C2	5.05	119.61	115.25
2	C	500	FAD	C4-N3-C2	5.14	119.69	115.25
2	A	500	FAD	C4-N3-C2	5.15	119.69	115.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 37 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	500	FAD	2	0
2	B	500	FAD	12	0
2	C	500	FAD	15	0
2	D	500	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	430/439 (97%)	0.03	2 (0%) 91 92	26, 39, 51, 62	0
1	B	430/439 (97%)	-0.05	0 100 100	28, 39, 51, 60	0
1	C	430/439 (97%)	0.02	3 (0%) 89 90	27, 40, 52, 61	0
1	D	430/439 (97%)	-0.12	1 (0%) 95 96	28, 40, 52, 62	0
All	All	1720/1756 (97%)	-0.03	6 (0%) 94 95	26, 39, 52, 62	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	224	ASP	2.8
1	A	108	ASP	2.6
1	D	36	LYS	2.4
1	C	36	LYS	2.4
1	A	310	ILE	2.2
1	C	34	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(Å ²)	Q<0.9
2	FAD	D	500	53/53	0.96	0.14	0.07	34,37,45,45	0
2	FAD	A	500	53/53	0.97	0.14	-0.07	25,29,39,39	0
2	FAD	C	500	53/53	0.96	0.15	-0.15	29,37,41,41	0
2	FAD	B	500	53/53	0.97	0.14	-0.33	26,30,35,35	0

6.5 Other polymers

There are no such residues in this entry.