



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 11:45 AM GMT

PDB ID : 3POP
Title : The crystal structure of GilR, an oxidoreductase that catalyzes the terminal step of gilvocarcin biosynthesis
Authors : Noinaj, N.; Bosserman, M.A.; Schickli, M.A.; Kharel, M.K.; Rohr, J.; Buchanan, S.K.
Deposited on : 2010-11-23
Resolution : 1.65 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

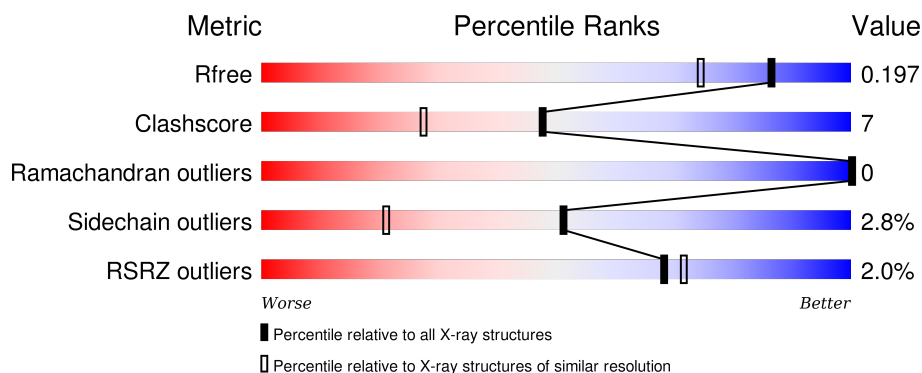
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.65 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	1226 (1.66-1.66)
Clashscore	102246	1323 (1.66-1.66)
Ramachandran outliers	100387	1295 (1.66-1.66)
Sidechain outliers	100360	1295 (1.66-1.66)
RSRZ outliers	91569	1227 (1.66-1.66)

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	501	<div> <div>2%</div> <div>85%</div> <div>12%</div> <div>..</div> </div>
1	B	501	<div> <div>%</div> <div>85%</div> <div>12%</div> <div>..</div> </div>
1	C	501	<div> <div>%</div> <div>85%</div> <div>12%</div> <div>..</div> </div>
1	D	501	<div> <div>3%</div> <div>85%</div> <div>12%</div> <div>..</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 17205 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 GilR oxidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	493	Total	C	N	O	S	0	1	0
			3775	2370	690	701	14			
1	B	493	Total	C	N	O	S	0	4	0
			3795	2380	693	708	14			
1	C	493	Total	C	N	O	S	0	1	0
			3779	2372	690	703	14			
1	D	493	Total	C	N	O	S	0	1	0
			3769	2367	687	701	14			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	EXPRESSION TAG	UNP Q7X2G7
A	-1	SER	-	EXPRESSION TAG	UNP Q7X2G7
A	0	HIS	-	EXPRESSION TAG	UNP Q7X2G7
B	-2	GLY	-	EXPRESSION TAG	UNP Q7X2G7
B	-1	SER	-	EXPRESSION TAG	UNP Q7X2G7
B	0	HIS	-	EXPRESSION TAG	UNP Q7X2G7
C	-2	GLY	-	EXPRESSION TAG	UNP Q7X2G7
C	-1	SER	-	EXPRESSION TAG	UNP Q7X2G7
C	0	HIS	-	EXPRESSION TAG	UNP Q7X2G7
D	-2	GLY	-	EXPRESSION TAG	UNP Q7X2G7
D	-1	SER	-	EXPRESSION TAG	UNP Q7X2G7
D	0	HIS	-	EXPRESSION TAG	UNP Q7X2G7

- 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		

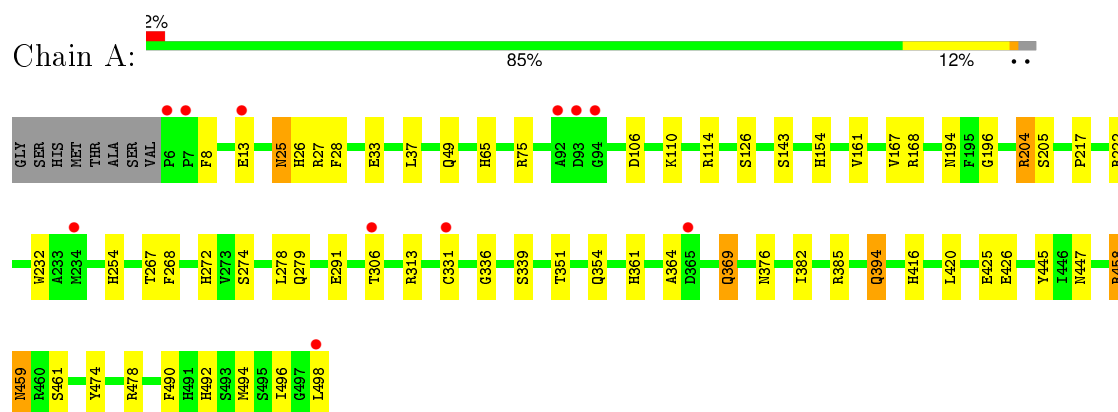
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	468	Total	O	0	0
			468	468		
3	B	472	Total	O	0	0
			472	472		
3	C	512	Total	O	0	0
			512	512		
3	D	423	Total	O	0	0
			423	423		

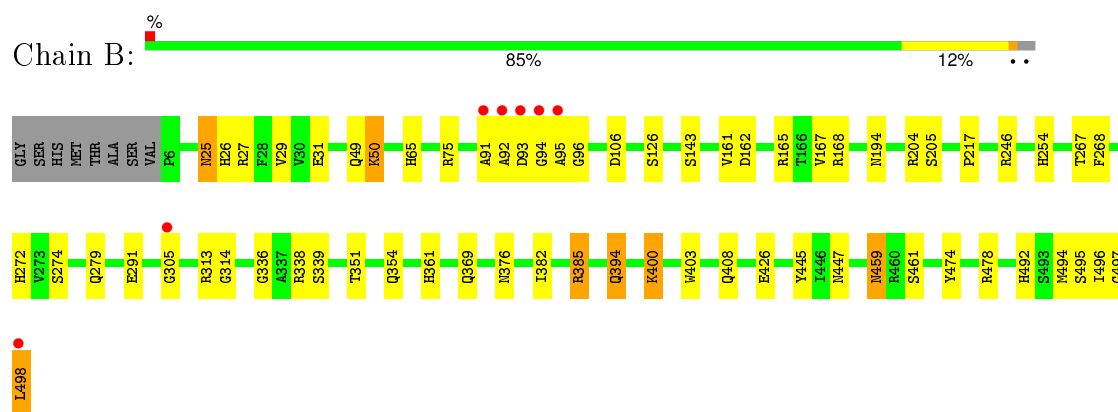
3 Residue-property plots

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.

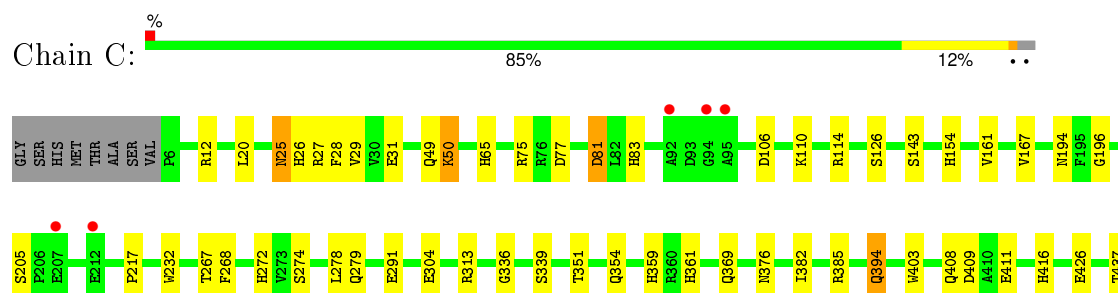
• Molecule 1: GilR oxidase



• Molecule 1: GilR oxidase

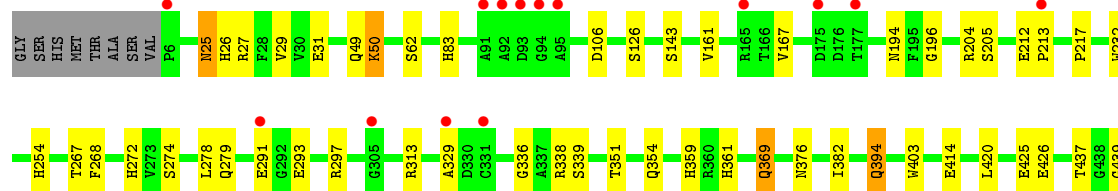
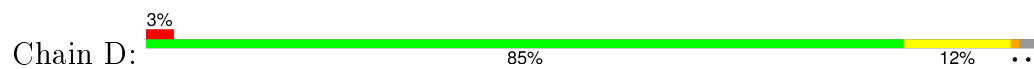


• Molecule 1: GilR oxidase





● Molecule 1: GilR oxidase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	68.08 Å 115.83 Å 290.10 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.97 – 1.65 24.24 – 1.65	Depositor EDS
% Data completeness (in resolution range)	98.0 (19.97-1.65) 97.8 (24.24-1.65)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.41 (at 1.65 Å)	Xtriage
Refinement program	PHENIX (phenix.refine: dev_501)	Depositor
R, R_{free}	0.173 , 0.198 0.174 , 0.197	Depositor DCC
R_{free} test set	2000 reflections (0.75%)	DCC
Wilson B-factor (Å ²)	16.8	Xtriage
Anisotropy	0.200	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 44.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 271402 reflections	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	17205	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.26% 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	0.40	0/3879	0.58	2/5276 (0.0%)
1	B	0.43	0/3899	0.58	1/5304 (0.0%)
1	C	0.41	0/3883	0.56	0/5281
1	D	0.39	0/3873	0.56	0/5269
All	All	0.41	0/15534	0.57	3/21130 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	385	ARG	NE-CZ-NH1	-6.37	117.12	120.30
1	A	385	ARG	NE-CZ-NH1	-6.33	117.13	120.30
1	A	385	ARG	NE-CZ-NH2	6.18	123.39	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3775	0	3592	52	0
1	B	3795	0	3604	57	0
1	C	3779	0	3596	58	0
1	D	3769	0	3581	54	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	53	0	29	1	0
2	B	53	0	29	5	0
2	C	53	0	29	1	0
2	D	53	0	29	1	0
3	A	468	0	0	11	0
3	B	472	0	0	4	0
3	C	512	0	0	16	0
3	D	423	0	0	10	0
All	All	17205	0	14489	219	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 7.

All (219) 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:C:408:GLN:HG2	3:C:2019:HOH:O	1.47	1.15
1:C:154:HIS:HD2	3:C:2012:HOH:O	1.34	1.11
1:B:92:ALA:HA	1:B:93:ASP:CB	1.79	1.11
1:D:361:HIS:HD2	3:D:2036:HOH:O	1.42	0.99
1:C:409:ASP:OD2	1:C:411:GLU:HB2	1.68	0.92
1:C:75:ARG:NH2	1:C:77:ASP:OD1	2.07	0.87
1:B:385:ARG:HD2	3:B:2008:HOH:O	1.74	0.87
1:D:293:GLU:HG2	1:D:297:ARG:NH1	1.91	0.86
1:A:351:THR:H	1:A:354:GLN:HE21	1.24	0.85
1:D:351:THR:H	1:D:354:GLN:HE21	1.28	0.81
1:A:33:GLU:HG3	1:A:75:ARG:NH2	1.95	0.80
1:B:96:GLY:HA3	1:B:204:ARG:HG3	1.62	0.80
1:B:65:HIS:CE1	2:B:499:FAD:C8M	2.62	0.79
1:D:293:GLU:OE1	1:D:297:ARG:NH1	2.16	0.79
1:A:492:HIS:HD2	1:A:494:MET:H	1.32	0.77
1:B:351:THR:H	1:B:354:GLN:HE21	1.29	0.77
1:A:110:LYS:O	1:A:114:ARG:HG3	1.85	0.77
1:D:492:HIS:HD2	1:D:494:MET:H	1.33	0.77
1:D:293:GLU:CD	1:D:297:ARG:HH12	1.88	0.76
1:C:351:THR:H	1:C:354:GLN:HE21	1.34	0.76
1:C:492:HIS:HD2	1:C:494:MET:H	1.33	0.76
1:B:492:HIS:HD2	1:B:494:MET:H	1.32	0.75
1:B:92:ALA:CA	1:B:93:ASP:CB	2.63	0.74
1:C:361:HIS:HD2	3:C:2017:HOH:O	1.71	0.73
1:D:369:GLN:HG3	3:D:2037:HOH:O	1.88	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:293:GLU:CG	1:D:297:ARG:NH1	2.53	0.70
1:B:361:HIS:HD2	3:B:2009:HOH:O	1.74	0.70
1:C:385:ARG:HD2	3:C:2018:HOH:O	1.90	0.70
1:A:267:THR:HG23	1:A:376:ASN:HD21	1.56	0.70
1:B:65:HIS:ND1	2:B:499:FAD:HM81	2.00	0.70
1:A:33:GLU:HG3	1:A:75:ARG:HH22	1.56	0.69
1:C:110:LYS:O	1:C:114:ARG:HG3	1.93	0.68
1:C:267:THR:HG23	1:C:376:ASN:HD21	1.60	0.67
1:B:279:GLN:HE22	1:B:313:ARG:HH22	1.44	0.66
1:D:254:HIS:HE1	3:D:2033:HOH:O	1.79	0.65
1:B:246:ARG:HH12	1:B:305:GLY:H	1.45	0.65
1:B:267:THR:HG23	1:B:376:ASN:HD21	1.61	0.64
1:C:279:GLN:HE22	1:C:313:ARG:HH22	1.43	0.64
1:B:272:HIS:CD2	1:B:274:SER:H	2.16	0.63
1:B:96:GLY:CA	1:B:204:ARG:HG3	2.27	0.63
1:A:254:HIS:HE1	3:A:1785:HOH:O	1.82	0.62
1:B:361:HIS:HE1	1:B:426:GLU:OE1	1.83	0.62
1:D:267:THR:HG23	1:D:376:ASN:HD21	1.63	0.62
1:C:361:HIS:HE1	1:C:426:GLU:OE1	1.82	0.62
1:C:83:HIS:HE1	3:C:1758:HOH:O	1.82	0.62
1:D:361:HIS:HE1	1:D:426:GLU:OE1	1.82	0.61
1:A:106:ASP:HB2	1:A:126:SER:HB2	1.81	0.61
1:D:293:GLU:CD	1:D:297:ARG:NH1	2.52	0.61
1:A:361:HIS:HD2	3:A:1996:HOH:O	1.83	0.61
1:D:279:GLN:HE22	1:D:313:ARG:HH22	1.47	0.61
1:B:106:ASP:HB2	1:B:126:SER:HB2	1.83	0.61
1:D:361:HIS:CD2	3:D:2036:HOH:O	2.29	0.60
1:D:272:HIS:CD2	1:D:274:SER:H	2.18	0.60
1:B:91:ALA:O	1:B:93:ASP:HA	2.02	0.60
1:C:272:HIS:CD2	1:C:274:SER:H	2.19	0.60
1:A:49:GLN:HE22	1:A:167:VAL:H	1.48	0.60
1:A:498:LEU:HD23	3:A:1958:HOH:O	2.01	0.60
1:B:65:HIS:CG	2:B:499:FAD:C8M	2.77	0.60
1:A:26:HIS:HE1	3:D:2035:HOH:O	1.84	0.59
1:A:272:HIS:CD2	1:A:274:SER:H	2.19	0.59
1:B:49:GLN:NE2	1:B:167:VAL:H	2.01	0.59
1:D:359:HIS:HD2	3:D:1663:HOH:O	1.85	0.59
1:A:361:HIS:HE1	1:A:426:GLU:OE1	1.85	0.59
1:A:49:GLN:NE2	1:A:167:VAL:H	2.00	0.59
1:C:359:HIS:HD2	3:C:1457:HOH:O	1.87	0.58
1:C:49:GLN:HE22	1:C:167:VAL:H	1.52	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:194:ASN:ND2	1:B:496:ILE:H	2.00	0.58
1:D:49:GLN:HE22	1:D:167:VAL:H	1.51	0.58
1:B:254:HIS:HE1	3:B:2041:HOH:O	1.86	0.58
1:C:154:HIS:CD2	3:C:2012:HOH:O	2.22	0.57
1:B:49:GLN:HE22	1:B:167:VAL:H	1.50	0.57
1:B:27:ARG:HD3	1:B:339:SER:O	2.04	0.57
1:B:95:ALA:O	1:B:204:ARG:HD2	2.04	0.57
1:A:279:GLN:HE22	1:A:313:ARG:HH22	1.51	0.57
1:D:49:GLN:NE2	1:D:167:VAL:H	2.03	0.56
1:C:114:ARG:HD3	3:C:2011:HOH:O	2.05	0.56
1:A:194:ASN:ND2	1:A:496:ILE:H	2.03	0.56
1:C:466:HIS:HD2	3:C:2021:HOH:O	1.89	0.56
1:A:267:THR:HG23	1:A:376:ASN:ND2	2.21	0.55
1:D:481:LYS:NZ	3:D:1596:HOH:O	2.39	0.54
1:B:93:ASP:CB	1:B:94:GLY:HA2	2.38	0.54
1:C:194:ASN:ND2	1:C:496:ILE:H	2.06	0.54
1:B:65:HIS:ND1	2:B:499:FAD:C8	2.64	0.54
1:C:49:GLN:NE2	1:C:167:VAL:H	2.04	0.54
1:C:408:GLN:CG	3:C:2019:HOH:O	2.23	0.54
1:D:194:ASN:ND2	1:D:496:ILE:H	2.06	0.54
1:C:445:TYR:CZ	1:C:447:ASN:HB2	2.43	0.53
1:C:27:ARG:HD2	3:C:756:HOH:O	2.08	0.53
1:B:161:VAL:HG22	1:B:167:VAL:HG22	1.90	0.53
1:D:351:THR:H	1:D:354:GLN:NE2	2.03	0.53
3:A:1592:HOH:O	1:D:26:HIS:HE1	1.91	0.53
1:D:106:ASP:HB2	1:D:126:SER:HB2	1.91	0.52
1:A:351:THR:H	1:A:354:GLN:NE2	2.00	0.51
1:A:27:ARG:HD3	1:A:339:SER:O	2.10	0.51
1:A:222:ARG:NH1	3:A:1471:HOH:O	2.43	0.51
1:D:293:GLU:CG	1:D:297:ARG:HH12	2.21	0.51
1:B:267:THR:HG23	1:B:376:ASN:ND2	2.25	0.51
1:C:385:ARG:CD	3:C:2018:HOH:O	2.53	0.51
1:C:267:THR:HG23	1:C:376:ASN:ND2	2.25	0.51
1:A:445:TYR:CZ	1:A:447:ASN:HB2	2.46	0.50
1:D:161:VAL:HG22	1:D:167:VAL:HG22	1.93	0.50
1:D:27:ARG:HD3	1:D:339:SER:O	2.11	0.50
1:A:143:SER:HB3	1:A:382:ILE:HD11	1.94	0.50
1:B:445:TYR:CZ	1:B:447:ASN:HB2	2.47	0.50
1:B:25:ASN:C	1:B:25:ASN:HD22	2.14	0.49
1:D:25:ASN:HD22	1:D:25:ASN:C	2.15	0.49
1:C:106:ASP:HB2	1:C:126:SER:HB2	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:25:ASN:HD22	1:A:25:ASN:C	2.15	0.49
1:A:425:GLU:OE1	1:A:458:ARG:HD3	2.12	0.49
1:C:143:SER:HB3	1:C:382:ILE:HD11	1.94	0.49
1:A:205:SER:OG	1:A:217:PRO:HD3	2.12	0.49
1:D:205:SER:OG	1:D:217:PRO:HD3	2.12	0.49
1:B:272:HIS:HD2	1:B:274:SER:H	1.57	0.49
1:D:267:THR:HG23	1:D:376:ASN:ND2	2.27	0.49
1:C:272:HIS:HD2	1:C:274:SER:H	1.61	0.49
1:C:351:THR:H	1:C:354:GLN:NE2	2.07	0.49
1:D:459:ASN:ND2	1:D:461:SER:H	2.11	0.49
1:A:369:GLN:HG3	3:A:1995:HOH:O	2.12	0.49
1:D:272:HIS:HD2	1:D:274:SER:H	1.61	0.48
1:C:279:GLN:NE2	1:C:313:ARG:HH22	2.11	0.48
1:B:27:ARG:HG3	1:B:338:ARG:HB3	1.95	0.48
1:D:445:TYR:CZ	1:D:447:ASN:HB2	2.48	0.48
1:D:439:GLY:HA3	3:D:1597:HOH:O	2.13	0.48
1:A:161:VAL:HG22	1:A:167:VAL:HG22	1.93	0.48
1:B:279:GLN:NE2	1:B:313:ARG:HH22	2.10	0.48
1:D:212:GLU:OE2	1:D:213:PRO:HA	2.13	0.48
1:C:25:ASN:C	1:C:25:ASN:HD22	2.17	0.48
1:D:143:SER:HB3	1:D:382:ILE:HD11	1.96	0.48
1:B:143:SER:HB3	1:B:382:ILE:HD11	1.96	0.48
1:A:13:GLU:HB2	3:A:1547:HOH:O	2.14	0.48
1:C:474:TYR:CE2	1:C:478:ARG:HD2	2.49	0.47
1:B:459:ASN:ND2	1:B:461:SER:H	2.12	0.47
1:A:26:HIS:CE1	1:A:336:GLY:HA2	2.50	0.47
1:B:194:ASN:HD21	1:B:496:ILE:H	1.63	0.47
1:C:459:ASN:ND2	1:C:461:SER:H	2.12	0.47
1:B:394:GLN:HE21	1:B:394:GLN:H	1.63	0.47
1:D:29:VAL:HG22	1:D:31:GLU:HG3	1.97	0.47
1:A:459:ASN:ND2	1:A:461:SER:H	2.12	0.47
1:D:474:TYR:CE2	1:D:478:ARG:HD2	2.50	0.47
1:B:96:GLY:HA3	1:B:204:ARG:CG	2.39	0.47
1:D:425:GLU:OE1	1:D:458:ARG:HD3	2.15	0.46
1:A:272:HIS:HD2	1:A:274:SER:H	1.64	0.46
1:C:50:LYS:HE2	1:C:50:LYS:HB3	1.80	0.46
1:A:154:HIS:HB2	1:A:204:ARG:HB2	1.96	0.46
1:A:474:TYR:CE2	1:A:478:ARG:HD2	2.51	0.46
1:C:161:VAL:HG22	1:C:167:VAL:HG22	1.97	0.46
1:B:459:ASN:C	1:B:459:ASN:HD22	2.20	0.46
1:A:154:HIS:HD2	3:A:1135:HOH:O	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:26:HIS:CE1	1:B:336:GLY:HA2	2.51	0.46
1:C:29:VAL:HG22	1:C:31:GLU:HG3	1.98	0.45
1:C:394:GLN:H	1:C:394:GLN:HE21	1.64	0.45
1:A:25:ASN:HD22	1:A:27:ARG:H	1.65	0.45
1:D:27:ARG:HG3	1:D:338:ARG:HB3	1.99	0.45
1:C:205:SER:OG	1:C:217:PRO:HD3	2.17	0.45
1:C:25:ASN:ND2	1:C:27:ARG:HB2	2.31	0.45
1:A:394:GLN:H	1:A:394:GLN:HE21	1.64	0.45
1:C:25:ASN:HD21	1:C:339:SER:H	1.63	0.45
1:B:497:GLY:O	1:B:498:LEU:O	2.34	0.45
1:C:12:ARG:HD3	3:C:1787:HOH:O	2.16	0.44
1:C:26:HIS:CE1	1:C:336:GLY:HA2	2.52	0.44
1:D:26:HIS:CE1	1:D:336:GLY:HA2	2.52	0.44
1:B:351:THR:H	1:B:354:GLN:NE2	2.07	0.44
1:D:232:TRP:CZ2	1:D:278:LEU:HD21	2.53	0.44
1:A:25:ASN:ND2	1:A:27:ARG:H	2.16	0.44
1:C:28:PHE:O	3:C:1219:HOH:O	2.21	0.44
1:A:232:TRP:CZ2	1:A:278:LEU:HD21	2.53	0.44
1:B:29:VAL:HG22	1:B:31:GLU:HG3	1.98	0.44
1:D:50:LYS:HE2	1:D:50:LYS:HB3	1.78	0.44
1:D:359:HIS:HE1	3:D:711:HOH:O	2.00	0.43
1:B:194:ASN:HD21	1:B:495:SER:HA	1.83	0.43
1:B:314:GLY:N	1:C:408:GLN:OE1	2.41	0.43
1:B:474:TYR:CE2	1:B:478:ARG:HD2	2.53	0.43
1:C:65:HIS:CE1	2:C:499:FAD:HM71	2.53	0.43
1:D:62[A]:SER:HB3	2:D:499:FAD:O1P	2.17	0.43
1:A:272:HIS:HE1	1:A:364:ALA:O	2.01	0.43
1:C:459:ASN:HD22	1:C:459:ASN:C	2.20	0.43
1:B:205:SER:OG	1:B:217:PRO:HD3	2.17	0.43
1:D:279:GLN:NE2	1:D:313:ARG:HH22	2.14	0.43
1:A:8:PHE:CZ	1:A:37:LEU:HD22	2.54	0.43
1:A:65:HIS:CE1	2:A:499:FAD:HM71	2.54	0.43
1:B:400:LYS:HE3	1:B:445:TYR:HB2	2.00	0.43
1:D:459:ASN:HD22	1:D:459:ASN:C	2.21	0.42
1:A:196:GLY:HA2	1:A:490:PHE:CE2	2.54	0.42
1:C:194:ASN:HD21	1:C:496:ILE:H	1.67	0.42
1:B:162:ASP:OD1	1:B:162:ASP:C	2.57	0.42
1:C:196:GLY:HA2	1:C:490:PHE:CE2	2.54	0.42
1:D:194:ASN:HD21	1:D:495:SER:HA	1.84	0.42
1:C:416:HIS:HE1	3:C:1119:HOH:O	2.01	0.42
1:A:331:CYS:SG	1:D:329:ALA:HB3	2.59	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:403:TRP:CD1	1:C:403:TRP:N	2.87	0.42
1:A:28:PHE:O	3:A:615:HOH:O	2.21	0.42
1:B:403:TRP:CD1	1:B:403:TRP:N	2.87	0.42
1:C:25:ASN:ND2	1:C:27:ARG:H	2.18	0.42
1:D:196:GLY:HA2	1:D:490:PHE:CE2	2.54	0.42
1:B:498:LEU:HD12	1:B:498:LEU:HA	1.81	0.42
1:D:25:ASN:HD21	1:D:339:SER:H	1.67	0.41
1:B:65:HIS:CE1	2:B:499:FAD:HM71	2.55	0.41
1:A:459:ASN:C	1:A:459:ASN:HD22	2.22	0.41
1:D:394:GLN:H	1:D:394:GLN:HE21	1.69	0.41
1:D:437:THR:HA	1:D:442:ASP:HA	2.01	0.41
1:B:25:ASN:HD21	1:B:339:SER:H	1.68	0.41
1:B:394:GLN:NE2	1:B:394:GLN:H	2.18	0.41
1:C:25:ASN:HD22	1:C:27:ARG:H	1.67	0.41
1:B:165:ARG:NE	3:B:1880:HOH:O	2.52	0.41
1:A:33:GLU:CG	1:A:75:ARG:NH2	2.78	0.40
1:C:437:THR:HA	1:C:442:ASP:HA	2.04	0.40
1:C:361:HIS:CD2	3:C:2017:HOH:O	2.56	0.40
1:C:232:TRP:CZ2	1:C:278:LEU:HD21	2.56	0.40
1:C:394:GLN:NE2	1:C:394:GLN:H	2.19	0.40
1:D:83:HIS:HE1	3:D:1359:HOH:O	2.03	0.40
1:A:254:HIS:HD2	3:A:562:HOH:O	2.04	0.40
1:B:25:ASN:ND2	1:B:27:ARG:H	2.20	0.40
1:A:37:LEU:HD12	1:A:37:LEU:N	2.37	0.40
1:B:50:LYS:HE2	1:B:50:LYS:HB3	1.76	0.40
1:A:416:HIS:HE1	3:A:1905:HOH:O	2.05	0.40
1:C:20:LEU:HD22	1:C:81:ASP:OD2	2.21	0.40
1:A:351:THR:N	1:A:354:GLN:HE21	2.06	0.40
1:A:394:GLN:H	1:A:394:GLN:NE2	2.19	0.40
1:D:403:TRP:CD1	1:D:403:TRP:N	2.89	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	492/501 (98%)	478 (97%)	14 (3%)	0	100	100
1	B	495/501 (99%)	479 (97%)	16 (3%)	0	100	100
1	C	492/501 (98%)	475 (96%)	17 (4%)	0	100	100
1	D	492/501 (98%)	478 (97%)	14 (3%)	0	100	100
All	All	1971/2004 (98%)	1910 (97%)	61 (3%)	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 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	387/396 (98%)	376 (97%)	11 (3%)	51	21
1	B	390/396 (98%)	378 (97%)	12 (3%)	47	16
1	C	388/396 (98%)	379 (98%)	9 (2%)	58	30
1	D	386/396 (98%)	375 (97%)	11 (3%)	51	21
All	All	1551/1584 (98%)	1508 (97%)	43 (3%)	51	21

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	ASN
1	A	168	ARG
1	A	204	ARG
1	A	268	PHE
1	A	291	GLU
1	A	306	THR
1	A	369	GLN
1	A	394	GLN
1	A	420	LEU
1	A	458	ARG

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Mol	Chain	Res	Type
1	A	459	ASN
1	B	25	ASN
1	B	50	LYS
1	B	75	ARG
1	B	168	ARG
1	B	268	PHE
1	B	291	GLU
1	B	369	GLN
1	B	394	GLN
1	B	400	LYS
1	B	408	GLN
1	B	459	ASN
1	B	498	LEU
1	C	25	ASN
1	C	50	LYS
1	C	81	ASP
1	C	268	PHE
1	C	291	GLU
1	C	304	GLU
1	C	369	GLN
1	C	394	GLN
1	C	459	ASN
1	D	25	ASN
1	D	50	LYS
1	D	204	ARG
1	D	268	PHE
1	D	291	GLU
1	D	369	GLN
1	D	394	GLN
1	D	414	GLU
1	D	420	LEU
1	D	458	ARG
1	D	459	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (71) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	25	ASN
1	A	49	GLN
1	A	68	GLN
1	A	83	HIS
1	A	154	HIS

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Mol	Chain	Res	Type
1	A	194	ASN
1	A	272	HIS
1	A	279	GLN
1	A	354	GLN
1	A	361	HIS
1	A	369	GLN
1	A	376	ASN
1	A	394	GLN
1	A	416	HIS
1	A	459	ASN
1	A	466	HIS
1	A	492	HIS
1	B	25	ASN
1	B	26	HIS
1	B	49	GLN
1	B	68	GLN
1	B	83	HIS
1	B	194	ASN
1	B	254	HIS
1	B	272	HIS
1	B	279	GLN
1	B	354	GLN
1	B	359	HIS
1	B	361	HIS
1	B	369	GLN
1	B	376	ASN
1	B	394	GLN
1	B	416	HIS
1	B	459	ASN
1	B	466	HIS
1	B	492	HIS
1	C	25	ASN
1	C	49	GLN
1	C	68	GLN
1	C	83	HIS
1	C	194	ASN
1	C	272	HIS
1	C	279	GLN
1	C	354	GLN
1	C	359	HIS
1	C	361	HIS
1	C	369	GLN

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Mol	Chain	Res	Type
1	C	376	ASN
1	C	394	GLN
1	C	416	HIS
1	C	459	ASN
1	C	466	HIS
1	C	492	HIS
1	D	25	ASN
1	D	26	HIS
1	D	49	GLN
1	D	68	GLN
1	D	83	HIS
1	D	194	ASN
1	D	272	HIS
1	D	279	GLN
1	D	354	GLN
1	D	359	HIS
1	D	361	HIS
1	D	369	GLN
1	D	376	ASN
1	D	394	GLN
1	D	416	HIS
1	D	459	ASN
1	D	466	HIS
1	D	492	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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](#)

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	499	1	48,58,58	1.69	10 (20%)	54,89,89	2.54	12 (22%)
2	FAD	B	499	1	48,58,58	1.71	13 (27%)	54,89,89	2.85	16 (29%)
2	FAD	C	499	1	48,58,58	1.79	10 (20%)	54,89,89	2.04	15 (27%)
2	FAD	D	499	1	48,58,58	1.75	8 (16%)	54,89,89	2.26	10 (18%)

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	499	1	-	0/30/50/50	0/6/6/6
2	FAD	B	499	1	-	0/30/50/50	0/6/6/6
2	FAD	C	499	1	-	0/30/50/50	0/6/6/6
2	FAD	D	499	1	-	0/30/50/50	0/6/6/6

All (41) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	499	FAD	C5A-C4A	-2.79	1.34	1.40
2	B	499	FAD	C6-C5X	-2.36	1.38	1.41
2	D	499	FAD	O4B-C1B	-2.27	1.38	1.41
2	A	499	FAD	C5A-N7A	-2.05	1.32	1.39
2	B	499	FAD	PA-O1A	-2.05	1.43	1.51
2	C	499	FAD	C6A-N6A	-2.04	1.28	1.34
2	A	499	FAD	C9-C8	2.01	1.43	1.37
2	A	499	FAD	C9A-N10	2.04	1.41	1.38
2	B	499	FAD	O4'-C4'	2.10	1.48	1.43
2	C	499	FAD	C9-C8	2.14	1.43	1.37
2	B	499	FAD	C5'-C4'	2.20	1.55	1.51
2	B	499	FAD	C4'-C3'	2.22	1.58	1.53
2	A	499	FAD	C10-N10	2.27	1.41	1.39
2	B	499	FAD	C10-N10	2.34	1.41	1.39
2	C	499	FAD	O3'-C3'	2.45	1.48	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	499	FAD	C7M-C7	2.46	1.56	1.51
2	C	499	FAD	C7M-C7	2.50	1.56	1.51
2	B	499	FAD	O4B-C1B	2.51	1.44	1.41
2	A	499	FAD	C5X-N5	2.63	1.39	1.35
2	D	499	FAD	C2A-N3A	2.72	1.37	1.32
2	D	499	FAD	C9-C8	2.79	1.45	1.37
2	C	499	FAD	C10-N10	2.83	1.42	1.39
2	D	499	FAD	C2A-N1A	3.03	1.39	1.33
2	B	499	FAD	C8M-C8	3.04	1.57	1.51
2	D	499	FAD	C4-N3	3.28	1.39	1.33
2	B	499	FAD	C7M-C7	3.32	1.57	1.51
2	B	499	FAD	C5X-N5	3.32	1.40	1.35
2	A	499	FAD	C5'-C4'	3.36	1.56	1.51
2	C	499	FAD	C4-N3	3.37	1.39	1.33
2	A	499	FAD	C4-N3	3.60	1.39	1.33
2	B	499	FAD	C4-N3	3.61	1.39	1.33
2	B	499	FAD	C4X-N5	3.66	1.39	1.33
2	C	499	FAD	C5X-N5	3.71	1.41	1.35
2	C	499	FAD	C4X-N5	3.84	1.39	1.33
2	D	499	FAD	C9A-N10	4.23	1.44	1.38
2	A	499	FAD	C2A-N1A	4.36	1.42	1.33
2	D	499	FAD	C10-N1	4.51	1.43	1.35
2	B	499	FAD	C2A-N3A	4.73	1.40	1.32
2	A	499	FAD	C4X-N5	5.66	1.42	1.33
2	D	499	FAD	C4X-N5	5.81	1.42	1.33
2	C	499	FAD	C2A-N3A	6.00	1.42	1.32

All (53) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	499	FAD	N3A-C2A-N1A	-14.76	117.60	128.89
2	A	499	FAD	N3A-C2A-N1A	-13.85	118.29	128.89
2	D	499	FAD	N3A-C2A-N1A	-9.84	121.36	128.89
2	C	499	FAD	N3A-C2A-N1A	-7.32	123.29	128.89
2	C	499	FAD	C4X-C4-N3	-4.83	116.99	123.59
2	D	499	FAD	C4X-C4-N3	-4.55	117.36	123.59
2	B	499	FAD	C4X-C4-N3	-4.38	117.60	123.59
2	B	499	FAD	C2B-C1B-N9A	-4.27	107.77	114.29
2	A	499	FAD	C4X-C4-N3	-3.26	119.13	123.59
2	B	499	FAD	C4X-C10-N10	-3.22	118.62	120.52
2	D	499	FAD	O2'-C2'-C1'	-3.17	102.16	109.94
2	D	499	FAD	O3B-C3B-C2B	-2.84	102.60	111.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	499	FAD	C9A-C5X-N5	-2.78	118.24	122.36
2	A	499	FAD	O2'-C2'-C1'	-2.70	103.30	109.94
2	C	499	FAD	C1B-N9A-C4A	-2.70	122.87	126.94
2	B	499	FAD	C2B-C3B-C4B	-2.56	97.35	102.61
2	A	499	FAD	C6-C5X-C9A	-2.49	115.70	118.98
2	D	499	FAD	C4B-O4B-C1B	-2.49	106.98	109.72
2	C	499	FAD	O3P-P-O5'	-2.45	96.42	102.94
2	B	499	FAD	O2'-C2'-C1'	-2.43	103.98	109.94
2	C	499	FAD	O2'-C2'-C3'	-2.41	102.95	109.02
2	D	499	FAD	C6-C5X-N5	-2.39	115.88	118.96
2	B	499	FAD	C4B-O4B-C1B	-2.27	107.22	109.72
2	A	499	FAD	C1'-C2'-C3'	-2.26	103.36	109.82
2	A	499	FAD	C1B-N9A-C4A	-2.19	123.64	126.94
2	C	499	FAD	C2B-C1B-N9A	-2.12	111.06	114.29
2	B	499	FAD	C1B-N9A-C4A	-2.11	123.75	126.94
2	C	499	FAD	O3'-C3'-C2'	-2.11	103.44	108.75
2	C	499	FAD	O2'-C2'-C1'	-2.01	105.00	109.94
2	A	499	FAD	O4B-C4B-C3B	2.10	109.39	105.15
2	C	499	FAD	C6-C5X-N5	2.11	121.67	118.96
2	D	499	FAD	O2A-PA-O3P	2.14	114.81	105.09
2	B	499	FAD	C9-C8-C7	2.17	124.18	120.04
2	D	499	FAD	O4B-C4B-C3B	2.19	109.55	105.15
2	B	499	FAD	C6-C5X-C9A	2.21	121.89	118.98
2	C	499	FAD	N6A-C6A-N1A	2.27	124.07	119.20
2	C	499	FAD	O4B-C1B-N9A	2.32	112.96	108.10
2	C	499	FAD	O2A-PA-O3P	2.37	115.82	105.09
2	B	499	FAD	C4X-N5-C5X	2.42	119.55	116.76
2	C	499	FAD	C2A-N1A-C6A	2.58	123.38	118.77
2	A	499	FAD	C5X-C9A-N10	2.82	119.76	117.62
2	A	499	FAD	C7-C6-C5X	2.84	125.55	120.92
2	A	499	FAD	C2A-N1A-C6A	2.88	123.91	118.77
2	C	499	FAD	C1'-N10-C9A	3.30	122.56	118.86
2	A	499	FAD	C1'-N10-C9A	3.58	122.88	118.86
2	B	499	FAD	C4-N3-C2	4.28	118.94	115.25
2	B	499	FAD	C1'-N10-C9A	4.38	123.78	118.86
2	B	499	FAD	C2A-N1A-C6A	5.04	127.78	118.77
2	B	499	FAD	C5X-C9A-N10	5.45	121.76	117.62
2	D	499	FAD	C4-N3-C2	5.94	120.38	115.25
2	D	499	FAD	C4X-C10-N10	6.06	124.09	120.52
2	C	499	FAD	C4-N3-C2	6.38	120.76	115.25
2	A	499	FAD	C4-N3-C2	6.42	120.80	115.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	499	FAD	1	0
2	B	499	FAD	5	0
2	C	499	FAD	1	0
2	D	499	FAD	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	493/501 (98%)	-0.21	11 (2%) 65 67	13, 24, 47, 93	0
1	B	493/501 (98%)	-0.30	7 (1%) 78 80	12, 21, 46, 96	0
1	C	493/501 (98%)	-0.31	6 (1%) 81 83	12, 23, 46, 106	0
1	D	493/501 (98%)	-0.17	15 (3%) 54 55	15, 25, 46, 93	0
All	All	1972/2004 (98%)	-0.25	39 (1%) 68 71	12, 23, 47, 106	0

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	94	GLY	7.4
1	D	498	LEU	7.1
1	B	92	ALA	7.0
1	A	92	ALA	5.9
1	A	6	PRO	4.9
1	B	498	LEU	4.9
1	C	92	ALA	4.7
1	D	92	ALA	4.2
1	A	498	LEU	4.1
1	B	94	GLY	3.6
1	D	94	GLY	3.5
1	B	93	ASP	3.4
1	D	177	THR	3.4
1	A	94	GLY	3.4
1	A	331	CYS	3.3
1	C	95	ALA	3.2
1	B	305	GLY	3.2
1	A	7	PRO	3.0
1	D	93	ASP	3.0
1	D	329	ALA	3.0
1	D	175	ASP	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	13	GLU	2.8
1	D	305	GLY	2.8
1	B	95	ALA	2.7
1	D	213	PRO	2.7
1	A	365	ASP	2.6
1	D	291	GLU	2.5
1	D	6	PRO	2.5
1	C	498	LEU	2.4
1	C	207	GLU	2.4
1	D	331	CYS	2.4
1	D	165	ARG	2.4
1	A	93	ASP	2.3
1	D	95	ALA	2.2
1	A	234	MET	2.2
1	C	212	GLU	2.1
1	D	91	ALA	2.1
1	B	91	ALA	2.0
1	A	306	THR	2.0

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	B	499	53/53	0.98	0.07	0.31	9,13,17,17	0
2	FAD	D	499	53/53	0.97	0.07	0.20	11,16,22,24	0
2	FAD	C	499	53/53	0.98	0.07	0.17	10,13,17,17	0
2	FAD	A	499	53/53	0.98	0.06	-0.33	10,13,17,20	0

6.5 Other polymers [i](#)

There are no such residues in this entry.