



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 7, 2017 – 05:04 AM EST

PDB ID : 4GTL
Title : T. Maritima FDTS (R174K mutant) with FAD
Authors : Mathews, I.I.; Lesley, S.A.; Kohen, A.
Deposited on : unknown
Resolution : 2.17 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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20030345
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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) : rb-20030345

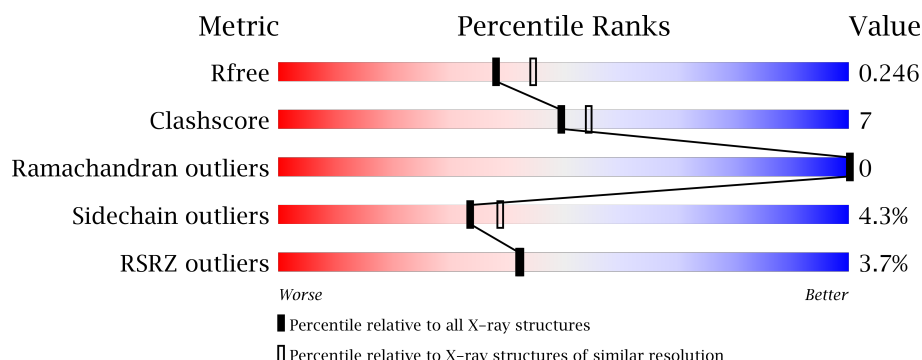
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.17 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}	100719	5526 (2.20-2.16)
Clashscore	112137	6386 (2.20-2.16)
Ramachandran outliers	110173	6282 (2.20-2.16)
Sidechain outliers	110143	6282 (2.20-2.16)
RSRZ outliers	101464	5562 (2.20-2.16)

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	232	<div> <div>5%</div> <div> <div></div> <div>75%</div> <div>16%</div> <div>7%</div> </div> </div>
1	B	232	<div> <div>2%</div> <div> <div></div> <div>75%</div> <div>16%</div> <div>7%</div> </div> </div>
1	C	232	<div> <div>3%</div> <div> <div></div> <div>79%</div> <div>13%</div> <div>8%</div> </div> </div>
1	D	232	<div> <div>4%</div> <div> <div></div> <div>78%</div> <div>15%</div> <div>7%</div> </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7454 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 Thymidylate synthase thyX.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	216	Total	C	N	O	S	0	2	0
			1833	1192	316	319	6			
1	B	215	Total	C	N	O	S	0	0	0
			1804	1175	308	316	5			
1	C	213	Total	C	N	O	S	0	0	0
			1787	1164	303	315	5			
1	D	216	Total	C	N	O	S	0	0	0
			1807	1177	307	318	5			

There are 52 discrepancies between the modelled and reference sequences:

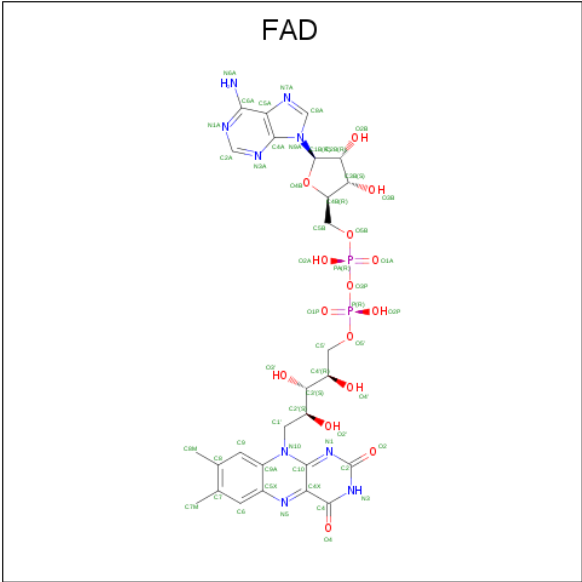
Chain	Residue	Modelled	Actual	Comment	Reference
A	-11	MET	-	EXPRESSION TAG	UNP Q9WYT0
A	-10	GLY	-	EXPRESSION TAG	UNP Q9WYT0
A	-9	SER	-	EXPRESSION TAG	UNP Q9WYT0
A	-8	ASP	-	EXPRESSION TAG	UNP Q9WYT0
A	-7	LYS	-	EXPRESSION TAG	UNP Q9WYT0
A	-6	ILE	-	EXPRESSION TAG	UNP Q9WYT0
A	-5	HIS	-	EXPRESSION TAG	UNP Q9WYT0
A	-4	HIS	-	EXPRESSION TAG	UNP Q9WYT0
A	-3	HIS	-	EXPRESSION TAG	UNP Q9WYT0
A	-2	HIS	-	EXPRESSION TAG	UNP Q9WYT0
A	-1	HIS	-	EXPRESSION TAG	UNP Q9WYT0
A	0	HIS	-	EXPRESSION TAG	UNP Q9WYT0
A	174	LYS	ARG	ENGINEERED MUTATION	UNP Q9WYT0
B	-11	MET	-	EXPRESSION TAG	UNP Q9WYT0
B	-10	GLY	-	EXPRESSION TAG	UNP Q9WYT0
B	-9	SER	-	EXPRESSION TAG	UNP Q9WYT0
B	-8	ASP	-	EXPRESSION TAG	UNP Q9WYT0
B	-7	LYS	-	EXPRESSION TAG	UNP Q9WYT0
B	-6	ILE	-	EXPRESSION TAG	UNP Q9WYT0
B	-5	HIS	-	EXPRESSION TAG	UNP Q9WYT0
B	-4	HIS	-	EXPRESSION TAG	UNP Q9WYT0

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-3	HIS	-	EXPRESSION TAG	UNP Q9WYT0
B	-2	HIS	-	EXPRESSION TAG	UNP Q9WYT0
B	-1	HIS	-	EXPRESSION TAG	UNP Q9WYT0
B	0	HIS	-	EXPRESSION TAG	UNP Q9WYT0
B	174	LYS	ARG	ENGINEERED MUTATION	UNP Q9WYT0
C	-11	MET	-	EXPRESSION TAG	UNP Q9WYT0
C	-10	GLY	-	EXPRESSION TAG	UNP Q9WYT0
C	-9	SER	-	EXPRESSION TAG	UNP Q9WYT0
C	-8	ASP	-	EXPRESSION TAG	UNP Q9WYT0
C	-7	LYS	-	EXPRESSION TAG	UNP Q9WYT0
C	-6	ILE	-	EXPRESSION TAG	UNP Q9WYT0
C	-5	HIS	-	EXPRESSION TAG	UNP Q9WYT0
C	-4	HIS	-	EXPRESSION TAG	UNP Q9WYT0
C	-3	HIS	-	EXPRESSION TAG	UNP Q9WYT0
C	-2	HIS	-	EXPRESSION TAG	UNP Q9WYT0
C	-1	HIS	-	EXPRESSION TAG	UNP Q9WYT0
C	0	HIS	-	EXPRESSION TAG	UNP Q9WYT0
C	174	LYS	ARG	ENGINEERED MUTATION	UNP Q9WYT0
D	-11	MET	-	EXPRESSION TAG	UNP Q9WYT0
D	-10	GLY	-	EXPRESSION TAG	UNP Q9WYT0
D	-9	SER	-	EXPRESSION TAG	UNP Q9WYT0
D	-8	ASP	-	EXPRESSION TAG	UNP Q9WYT0
D	-7	LYS	-	EXPRESSION TAG	UNP Q9WYT0
D	-6	ILE	-	EXPRESSION TAG	UNP Q9WYT0
D	-5	HIS	-	EXPRESSION TAG	UNP Q9WYT0
D	-4	HIS	-	EXPRESSION TAG	UNP Q9WYT0
D	-3	HIS	-	EXPRESSION TAG	UNP Q9WYT0
D	-2	HIS	-	EXPRESSION TAG	UNP Q9WYT0
D	-1	HIS	-	EXPRESSION TAG	UNP Q9WYT0
D	0	HIS	-	EXPRESSION TAG	UNP Q9WYT0
D	174	LYS	ARG	ENGINEERED MUTATION	UNP Q9WYT0

- 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	B	1	Total	C	N	O	P	0	0
			35	15	5	13	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		

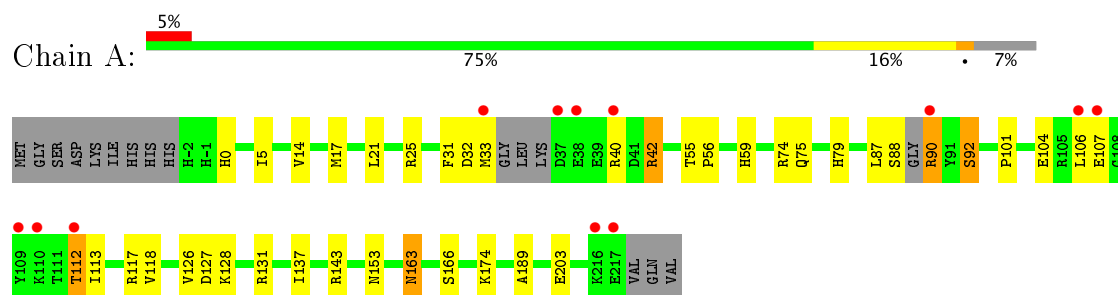
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	21	Total	O	0	1
			22	22		
3	B	19	Total	O	0	0
			19	19		
3	C	25	Total	O	0	0
			25	25		
3	D	16	Total	O	0	0
			16	16		

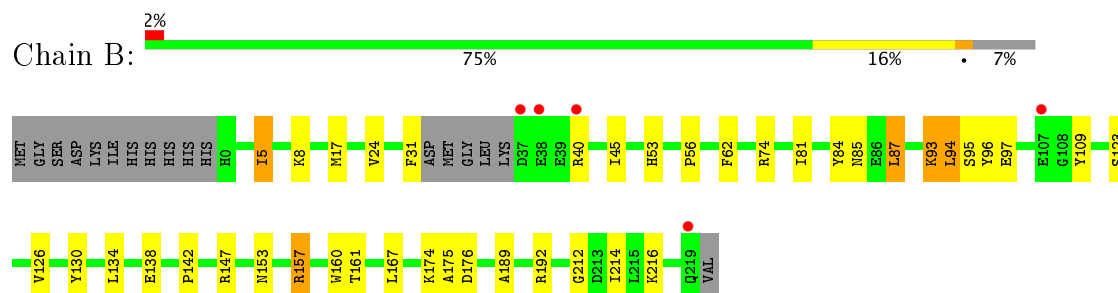
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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.

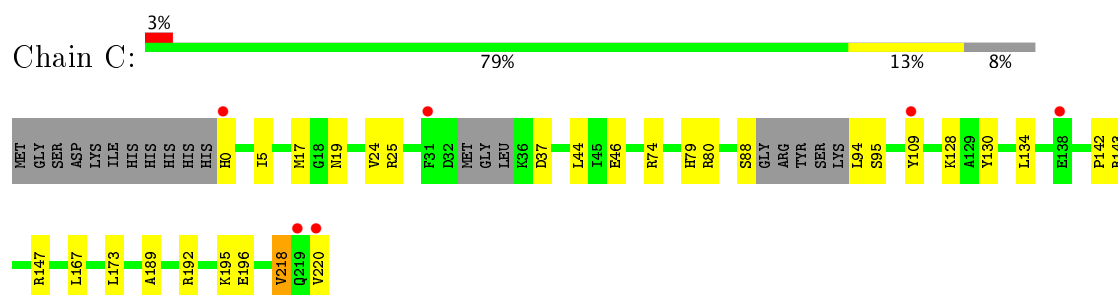
- Molecule 1: Thymidylate synthase thyX



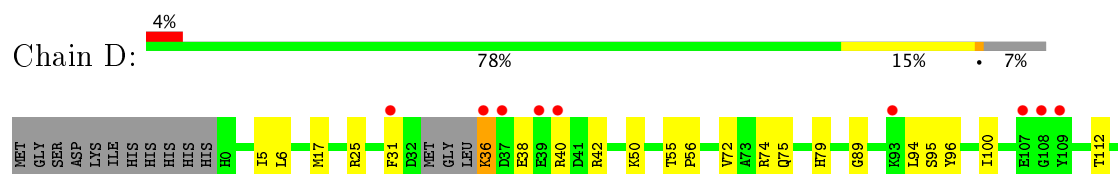
- Molecule 1: Thymidylate synthase thyX



- Molecule 1: Thymidylate synthase thyX



- Molecule 1: Thymidylate synthase thyX



T119	S123	E124	I125	L136	R143	R147	K174	A175	D176	Q180	I183	A189	K207	G212	K216	E217	V218	GLN	VAL
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4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	53.78Å 116.46Å 141.19Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.05 – 2.17 38.05 – 2.17	Depositor EDS
% Data completeness (in resolution range)	99.8 (38.05-2.17) 99.8 (38.05-2.17)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.00 (at 2.18Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.189 , 0.241 0.196 , 0.246	Depositor DCC
R_{free} test set	2415 reflections (5.33%)	DCC
Wilson B-factor (Å ²)	42.3	Xtriage
Anisotropy	0.079	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 46.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7454	wwPDB-VP
Average B, all atoms (Å ²)	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.47% 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, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	0.76	0/1889	0.82	0/2551
1	B	0.73	0/1853	0.81	1/2505 (0.0%)
1	C	0.79	0/1834	0.85	2/2478 (0.1%)
1	D	0.79	0/1855	0.86	2/2507 (0.1%)
All	All	0.77	0/7431	0.83	5/10041 (0.0%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	147	ARG	NE-CZ-NH2	-7.48	116.56	120.30
1	D	147	ARG	NE-CZ-NH1	6.11	123.36	120.30
1	B	157	ARG	NE-CZ-NH1	5.67	123.14	120.30
1	C	80	ARG	NE-CZ-NH1	5.07	122.83	120.30
1	C	37	ASP	CB-CG-OD1	5.03	122.83	118.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	1833	0	1821	30	0
1	B	1804	0	1796	26	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	1787	0	1778	25	1
1	D	1807	0	1800	22	1
2	B	35	0	20	1	0
2	C	53	0	31	3	0
2	D	53	0	31	2	0
3	A	22	0	0	1	0
3	B	19	0	0	1	0
3	C	25	0	0	3	0
3	D	16	0	0	0	0
All	All	7454	0	7277	98	1

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 (98) 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:17:MET:HB2	1:D:17:MET:HB2	1.57	0.86
1:A:163:ASN:HD22	1:A:163:ASN:C	1.82	0.82
1:A:17:MET:HB2	1:B:17:MET:HB2	1.70	0.73
1:A:5:ILE:HD11	1:A:189:ALA:HB2	1.71	0.72
1:D:96:TYR:OH	1:D:143:ARG:NH2	2.22	0.71
1:C:195:LYS:NZ	1:C:220:VAL:HG12	2.05	0.71
1:B:130:TYR:CE2	1:B:134:LEU:HD11	2.26	0.70
1:A:126:VAL:HG21	1:A:153:ASN:HD21	1.59	0.67
1:B:93:LYS:HD2	1:B:93:LYS:N	2.09	0.67
1:D:36:LYS:HZ2	1:D:36:LYS:N	1.95	0.65
1:A:112:THR:HG22	1:A:113:ILE:HG23	1.80	0.63
2:C:301:FAD:N7A	3:C:408:HOH:O	2.31	0.62
1:B:134:LEU:O	1:B:138:GLU:HB2	2.00	0.61
1:D:79:HIS:NE2	1:D:174:LYS:NZ	2.40	0.61
1:C:130:TYR:CE2	1:C:134:LEU:HD11	2.36	0.60
1:C:195:LYS:HZ3	1:C:220:VAL:HG12	1.66	0.60
1:B:87:LEU:HD22	1:B:157:ARG:HB2	1.83	0.60
1:A:163:ASN:ND2	1:A:166:SER:H	2.02	0.58
1:D:72:VAL:HG13	1:D:183:ILE:HD12	1.86	0.58
1:A:163:ASN:ND2	1:A:163:ASN:C	2.55	0.57
1:A:88:SER:C	1:A:90:ARG:N	2.58	0.57
1:C:195:LYS:HD3	1:C:220:VAL:HG12	1.86	0.56
1:A:127:ASP:OD2	1:A:131[A]:ARG:NH1	2.38	0.56
1:B:130:TYR:CZ	1:B:134:LEU:HD11	2.41	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:195:LYS:NZ	1:C:220:VAL:CG1	2.69	0.55
1:B:93:LYS:HD2	1:B:93:LYS:H	1.72	0.55
1:A:25:ARG:HD2	1:A:31:PHE:HA	1.89	0.54
1:A:75:GLN:NE2	3:A:319:HOH:O	2.41	0.54
2:C:301:FAD:O4'	2:C:301:FAD:O2'	2.26	0.54
1:D:112:THR:HG22	1:D:112:THR:O	2.06	0.54
1:C:130:TYR:CZ	1:C:134:LEU:HD11	2.43	0.54
1:D:176:ASP:OD1	1:D:212:GLY:HA2	2.08	0.53
2:B:301:FAD:N7A	3:B:401:HOH:O	2.34	0.52
1:C:195:LYS:CD	1:C:220:VAL:HG12	2.39	0.52
1:A:21:LEU:HD13	1:A:33:MET:HA	1.90	0.51
1:D:5:ILE:HD11	1:D:189:ALA:HB2	1.93	0.51
1:D:25:ARG:HD3	1:D:31:PHE:O	2.11	0.51
1:B:62:PHE:O	1:B:161:THR:HA	2.11	0.51
1:B:8:LYS:HD3	1:B:97:GLU:HB3	1.91	0.51
1:B:94:LEU:HD22	1:B:147:ARG:CZ	2.41	0.51
1:A:113:ILE:HD12	1:A:117:ARG:CB	2.42	0.50
1:D:112:THR:O	1:D:112:THR:CG2	2.59	0.50
1:A:5:ILE:HD11	1:A:189:ALA:CB	2.42	0.49
1:A:137:ILE:HD11	1:A:143:ARG:HA	1.93	0.49
2:D:300:FAD:HO4'	2:D:300:FAD:HO2'	1.56	0.49
1:C:167:LEU:O	1:C:167:LEU:HD23	2.13	0.48
1:C:88:SER:O	1:C:94:LEU:HD21	2.12	0.48
1:C:195:LYS:HZ3	1:C:220:VAL:CG1	2.25	0.48
1:C:218:VAL:O	1:C:218:VAL:HG12	2.14	0.48
1:A:79:HIS:NE2	1:A:174:LYS:NZ	2.38	0.47
1:B:74:ARG:HD3	1:C:74:ARG:HG3	1.96	0.47
1:A:42:ARG:NH2	1:A:203:GLU:OE1	2.47	0.47
1:B:84:TYR:CE1	1:B:160:TRP:CD1	3.03	0.47
1:C:25:ARG:NH1	3:C:423:HOH:O	2.41	0.46
1:A:87:LEU:HD11	1:A:92:SER:HB2	1.97	0.46
1:D:100:ILE:HG22	1:D:119:THR:HG23	1.97	0.46
1:A:59:HIS:NE2	1:B:85:ASN:ND2	2.62	0.46
1:C:192:ARG:HH21	1:C:220:VAL:HG23	1.80	0.46
1:D:89:GLY:HA3	1:D:94:LEU:CD1	2.46	0.46
2:D:300:FAD:H9	2:D:300:FAD:H1'1	1.68	0.46
1:A:128:LYS:HG3	1:D:125:ILE:HD11	1.99	0.45
1:D:5:ILE:HG22	1:D:6:LEU:HG	1.98	0.45
1:B:24:VAL:HG23	1:B:45:ILE:HD11	1.98	0.45
1:B:5:ILE:HD11	1:B:189:ALA:HB2	1.99	0.45
1:D:94:LEU:HD13	1:D:143:ARG:NH2	2.31	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:104:GLU:O	1:A:107:GLU:HG2	2.18	0.44
1:C:5:ILE:HD11	1:C:189:ALA:HB2	1.98	0.44
1:C:143:ARG:O	1:C:147:ARG:HG2	2.18	0.44
1:D:75:GLN:HB2	1:D:183:ILE:HD13	1.99	0.44
1:A:87:LEU:HD12	1:A:88:SER:N	2.32	0.44
1:B:74:ARG:HD3	1:C:74:ARG:CG	2.48	0.44
1:A:113:ILE:HD12	1:A:117:ARG:HB2	2.00	0.43
1:D:125:ILE:HD13	1:D:125:ILE:HA	1.85	0.43
1:A:5:ILE:CD1	1:A:189:ALA:HB2	2.44	0.43
1:C:167:LEU:HD23	1:C:167:LEU:C	2.39	0.43
1:D:55:THR:N	1:D:56:PRO:CD	2.82	0.43
1:A:113:ILE:HD12	1:A:117:ARG:HB3	2.01	0.43
1:C:79:HIS:CE1	1:C:173:LEU:HD12	2.54	0.43
1:B:167:LEU:HD23	1:B:167:LEU:C	2.39	0.42
1:B:126:VAL:HG21	1:B:153:ASN:HD21	1.85	0.42
1:B:95:SER:O	1:B:96:TYR:C	2.58	0.42
1:B:53:HIS:CD2	1:B:56:PRO:HD3	2.55	0.42
1:C:24:VAL:HG13	1:C:44:LEU:HD23	2.02	0.42
1:A:74:ARG:HG2	1:D:74:ARG:CD	2.50	0.41
1:A:0:HIS:CD2	1:B:31:PHE:CZ	3.08	0.41
1:D:207:LYS:HE2	1:D:207:LYS:HA	2.00	0.41
1:B:109:TYR:CE2	1:C:142:PRO:HG3	2.55	0.41
2:C:301:FAD:H1'1	2:C:301:FAD:H9	1.95	0.41
1:B:176:ASP:OD1	1:B:212:GLY:HA2	2.20	0.41
1:C:218:VAL:O	1:C:218:VAL:CG1	2.68	0.41
1:B:94:LEU:HD22	1:B:147:ARG:NH2	2.35	0.41
1:D:180:GLN:O	1:D:183:ILE:HG22	2.21	0.41
1:A:163:ASN:HD21	1:A:166:SER:H	1.65	0.41
1:B:142:PRO:HG3	1:C:109:TYR:CZ	2.55	0.41
1:A:55:THR:N	1:A:56:PRO:CD	2.83	0.41
1:C:19:ASN:HB2	3:C:416:HOH:O	2.20	0.41
1:A:118:VAL:HG13	1:D:136:LEU:HD22	2.02	0.40
1:B:175:ALA:HA	1:B:214:ILE:HD11	2.03	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:46:GLU:OE1	1:D:216:LYS:NZ[1_455]	2.18	0.02

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	212/232 (91%)	203 (96%)	9 (4%)	0	100	100
1	B	211/232 (91%)	203 (96%)	8 (4%)	0	100	100
1	C	207/232 (89%)	200 (97%)	7 (3%)	0	100	100
1	D	212/232 (91%)	204 (96%)	8 (4%)	0	100	100
All	All	842/928 (91%)	810 (96%)	32 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	196/207 (95%)	186 (95%)	10 (5%)	28	30
1	B	192/207 (93%)	182 (95%)	10 (5%)	27	30
1	C	191/207 (92%)	186 (97%)	5 (3%)	51	61
1	D	192/207 (93%)	184 (96%)	8 (4%)	34	40
All	All	771/828 (93%)	738 (96%)	33 (4%)	33	39

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	VAL
1	A	32	ASP

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Mol	Chain	Res	Type
1	A	40	ARG
1	A	42	ARG
1	A	90	ARG
1	A	92	SER
1	A	101	PRO
1	A	106	LEU
1	A	112	THR
1	A	163	ASN
1	B	5	ILE
1	B	40	ARG
1	B	81	ILE
1	B	87	LEU
1	B	93	LYS
1	B	94	LEU
1	B	123	SER
1	B	174	LYS
1	B	192	ARG
1	B	216	LYS
1	C	0	HIS
1	C	95	SER
1	C	128	LYS
1	C	196	GLU
1	C	218	VAL
1	D	36	LYS
1	D	38	GLU
1	D	40	ARG
1	D	42	ARG
1	D	50	LYS
1	D	95	SER
1	D	123	SER
1	D	183	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	163	ASN
1	A	178	HIS
1	B	85	ASN
1	B	219	GLN
1	D	185	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 ⓘ

3 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	B	301	-	31,37,58	1.27	3 (9%)	29,56,89	2.02	6 (20%)
2	FAD	C	301	-	51,58,58	3.30	15 (29%)	54,89,89	2.81	16 (29%)
2	FAD	D	300	-	51,58,58	3.27	16 (31%)	54,89,89	2.51	13 (24%)

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	B	301	-	-	0/26/46/50	0/3/3/6
2	FAD	C	301	-	-	0/28/50/50	0/6/6/6
2	FAD	D	300	-	-	0/28/50/50	0/6/6/6

All (34) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	300	FAD	C9A-N10	-7.45	1.29	1.38
2	C	301	FAD	C9A-N10	-5.45	1.31	1.38
2	C	301	FAD	C2-N1	-4.58	1.29	1.38
2	D	300	FAD	C2-N1	-4.28	1.29	1.38
2	C	301	FAD	PA-O1A	2.02	1.58	1.50
2	B	301	FAD	C5'-C4'	2.12	1.54	1.51
2	C	301	FAD	O4B-C1B	2.17	1.44	1.41
2	D	300	FAD	C4-N3	2.19	1.37	1.33
2	D	300	FAD	C2A-N3A	2.58	1.36	1.32
2	D	300	FAD	O4B-C1B	2.90	1.45	1.41
2	B	301	FAD	PA-O1A	3.23	1.63	1.50
2	C	301	FAD	C5A-C4A	3.53	1.48	1.40
2	D	300	FAD	C9-C9A	3.54	1.48	1.40
2	B	301	FAD	C5A-C4A	3.97	1.49	1.40
2	D	300	FAD	C8-C7	4.16	1.51	1.41
2	D	300	FAD	C5A-C4A	4.24	1.50	1.40
2	C	301	FAD	C8-C7	4.56	1.52	1.41
2	C	301	FAD	C9A-C5X	4.59	1.52	1.42
2	C	301	FAD	C4-C4X	4.69	1.50	1.41
2	C	301	FAD	C9-C9A	4.87	1.51	1.40
2	D	300	FAD	C4-C4X	4.91	1.50	1.41
2	D	300	FAD	C9A-C5X	5.13	1.53	1.42
2	D	300	FAD	C6-C7	5.19	1.51	1.37
2	C	301	FAD	C6-C7	5.53	1.52	1.37
2	D	300	FAD	C6-C5X	5.77	1.50	1.41
2	D	300	FAD	C5X-N5	5.97	1.44	1.35
2	D	300	FAD	C4X-C10	6.17	1.51	1.41
2	C	301	FAD	C4X-C10	6.21	1.51	1.41
2	C	301	FAD	C5X-N5	6.43	1.45	1.35
2	C	301	FAD	C6-C5X	8.10	1.53	1.41
2	D	300	FAD	C4X-N5	8.21	1.45	1.33
2	C	301	FAD	C4X-N5	8.24	1.45	1.33
2	D	300	FAD	C10-N1	11.00	1.48	1.33
2	C	301	FAD	C10-N1	11.01	1.48	1.33

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	301	FAD	N3A-C2A-N1A	-8.64	121.33	128.86
2	B	301	FAD	N3A-C2A-N1A	-7.75	122.11	128.86
2	D	300	FAD	N3A-C2A-N1A	-7.52	122.31	128.86
2	C	301	FAD	C4-C4X-C10	-4.85	116.03	119.96
2	C	301	FAD	C1'-N10-C10	-4.70	113.69	118.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	300	FAD	C4-C4X-C10	-4.52	116.31	119.96
2	C	301	FAD	C4X-C4-N3	-3.94	117.88	123.48
2	D	300	FAD	C4X-C4-N3	-3.77	118.12	123.48
2	D	300	FAD	O2'-C2'-C1'	-3.23	102.33	109.79
2	D	300	FAD	C7-C6-C5X	-2.88	116.63	121.08
2	C	301	FAD	C4A-C5A-N7A	-2.48	107.01	109.41
2	C	301	FAD	O4'-C4'-C3'	-2.41	103.12	109.09
2	C	301	FAD	C4X-C10-N10	-2.39	118.86	120.52
2	C	301	FAD	O2'-C2'-C1'	-2.31	104.45	109.79
2	B	301	FAD	O3'-C3'-C4'	-2.27	103.18	108.82
2	C	301	FAD	C7-C6-C5X	-2.03	117.94	121.08
2	C	301	FAD	C1B-N9A-C4A	-2.00	123.18	126.64
2	D	300	FAD	C1'-N10-C9A	2.04	120.22	118.35
2	D	300	FAD	C2A-N1A-C6A	2.07	122.40	118.77
2	B	301	FAD	O5'-C5'-C4'	2.15	115.09	109.36
2	B	301	FAD	C2A-N1A-C6A	2.27	122.75	118.77
2	D	300	FAD	O4'-C4'-C5'	2.34	115.20	110.00
2	D	300	FAD	O2P-P-O1P	2.44	124.91	112.28
2	C	301	FAD	O2A-PA-O1A	2.48	125.12	112.28
2	B	301	FAD	C4B-O4B-C1B	2.60	112.54	109.77
2	D	300	FAD	N6A-C6A-N1A	2.65	124.02	118.77
2	C	301	FAD	C2A-N1A-C6A	2.77	123.62	118.77
2	B	301	FAD	C4'-C3'-C2'	3.29	117.78	112.46
2	C	301	FAD	C5X-C9A-N10	3.82	120.50	117.66
2	D	300	FAD	C5X-C9A-N10	4.50	121.00	117.66
2	D	300	FAD	C4-C4X-N5	5.44	124.64	118.68
2	C	301	FAD	C4-C4X-N5	5.77	125.00	118.68
2	C	301	FAD	C1'-N10-C9A	6.01	123.85	118.35
2	D	300	FAD	C4-N3-C2	10.37	124.23	115.16
2	C	301	FAD	C4-N3-C2	11.00	124.78	115.16

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	301	FAD	1	0
2	C	301	FAD	3	0
2	D	300	FAD	2	0

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	216/232 (93%)	0.24	12 (5%) 25 26	36, 49, 85, 106	0
1	B	215/232 (92%)	0.07	5 (2%) 61 61	36, 51, 79, 99	0
1	C	213/232 (91%)	-0.04	6 (2%) 53 54	31, 47, 76, 95	0
1	D	216/232 (93%)	0.07	9 (4%) 37 37	33, 50, 81, 98	0
All	All	860/928 (92%)	0.08	32 (3%) 42 42	31, 49, 81, 106	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	220	VAL	5.3
1	A	106	LEU	4.5
1	D	40	ARG	4.0
1	A	40	ARG	3.5
1	D	36	LYS	3.5
1	B	40	ARG	3.5
1	B	37	ASP	3.5
1	A	33	MET	3.3
1	A	38	GLU	3.3
1	C	0	HIS	3.3
1	D	109	TYR	3.3
1	B	38	GLU	3.2
1	D	39	GLU	3.1
1	A	217	GLU	3.1
1	D	93	LYS	3.0
1	A	110	LYS	3.0
1	A	90	ARG	2.8
1	A	112	THR	2.7
1	C	31	PHE	2.7
1	D	37	ASP	2.7
1	C	138	GLU	2.6

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Mol	Chain	Res	Type	RSRZ
1	C	219	GLN	2.4
1	B	107	GLU	2.3
1	B	219	GLN	2.3
1	D	107	GLU	2.3
1	C	109	TYR	2.3
1	D	31	PHE	2.2
1	A	37	ASP	2.2
1	A	109	TYR	2.1
1	A	107	GLU	2.1
1	D	108	GLY	2.1
1	A	216	LYS	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	C	301	53/53	0.90	0.16	0.64	45,61,73,83	0
2	FAD	D	300	53/53	0.95	0.13	-0.63	39,52,61,71	0
2	FAD	B	301	35/53	0.93	0.12	-1.19	39,65,81,84	0

6.5 Other polymers [i](#)

There are no such residues in this entry.