



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 27, 2014 – 12:13 PM GMT

PDB ID : 3N0C
Title : TM0449 mutant crystal grown by hanging drop method
Authors : Mathews, I.I.
Deposited on : 2010-05-13
Resolution : 2.30 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

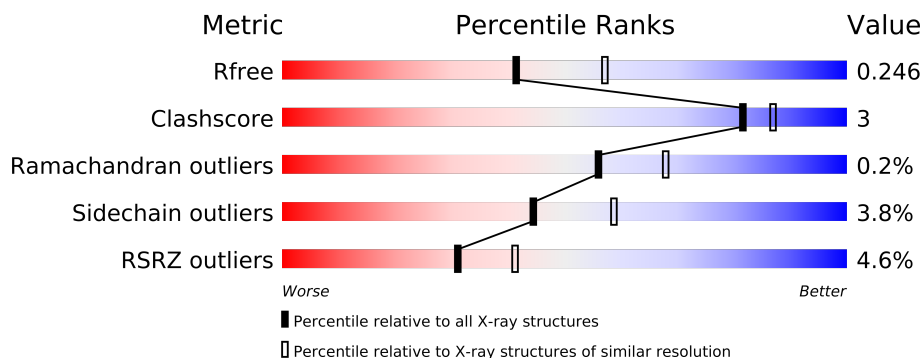
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.30 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}	66092	2929 (2.30-2.30)
Clashscore	79885	3679 (2.30-2.30)
Ramachandran outliers	78287	3642 (2.30-2.30)
Sidechain outliers	78261	3641 (2.30-2.30)
RSRZ outliers	66119	2930 (2.30-2.30)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	232	
1	B	232	
1	C	232	
1	D	232	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7471 atoms, of which 0 are hydrogen and 0 are deuterium.

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	0	0
			1778	1153	307	312	6			
1	B	220	Total	C	N	O	S	0	0	0
			1821	1180	313	322	6			
1	C	216	Total	C	N	O	S	0	0	0
			1793	1162	308	318	5			
1	D	210	Total	C	N	O	S	0	0	0
			1741	1128	300	308	5			

There are 56 discrepancies between the modelled and reference sequences:

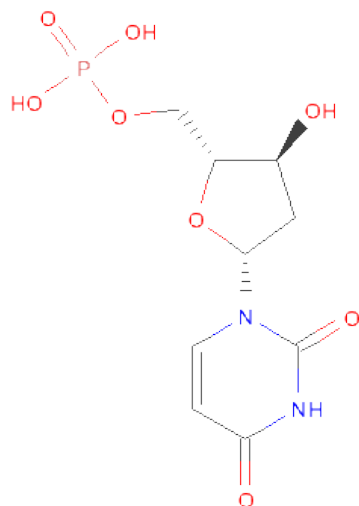
Chain	Residue	Modelled	Actual	Comment	Reference
A	-11	MET	-	EXPRESSION TAG	UNP Q9WYT0
A	-10	GLY	-	EXPRESSION TAG	UNP Q9WYT0
A	-9	SER	-	EXPRESSION TAG	UNP Q9WYT0
A	-8	ASP	-	EXPRESSION TAG	UNP Q9WYT0
A	-7	LYS	-	EXPRESSION TAG	UNP Q9WYT0
A	-6	ILE	-	EXPRESSION TAG	UNP Q9WYT0
A	-5	HIS	-	EXPRESSION TAG	UNP 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	158	ALA	PHE	ENGINEERED MUTATION	UNP Q9WYT0
A	160	ALA	TRP	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

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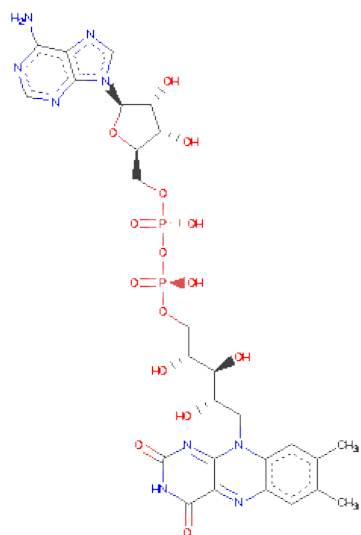
Chain	Residue	Modelled	Actual	Comment	Reference
B	-4	HIS	-	EXPRESSION TAG	UNP Q9WYT0
B	-3	HIS	-	EXPRESSION TAG	UNP Q9WYT0
B	-2	HIS	-	EXPRESSION TAG	UNP Q9WYT0
B	-1	HIS	-	EXPRESSION TAG	UNP Q9WYT0
B	0	HIS	-	EXPRESSION TAG	UNP Q9WYT0
B	158	ALA	PHE	ENGINEERED MUTATION	UNP Q9WYT0
B	160	ALA	TRP	ENGINEERED 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	158	ALA	PHE	ENGINEERED MUTATION	UNP Q9WYT0
C	160	ALA	TRP	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	158	ALA	PHE	ENGINEERED MUTATION	UNP Q9WYT0
D	160	ALA	TRP	ENGINEERED MUTATION	UNP Q9WYT0

- Molecule 2 is 2'-DEOXYURIDINE 5'-MONOPHOSPHATE (three-letter code: UMP) (formula: $C_9H_{13}N_2O_8P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			20	9	2	8	1		
2	B	1	Total	C	N	O	P	0	0
			20	9	2	8	1		
2	C	1	Total	C	N	O	P	0	0
			20	9	2	8	1		
2	D	1	Total	C	N	O	P	0	0
			20	9	2	8	1		

- Molecule 3 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



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

- Molecule 4 is water.

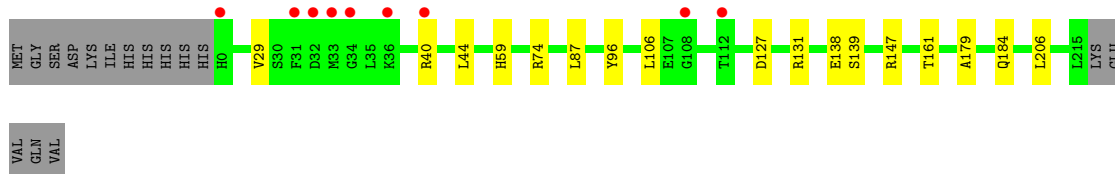
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	11	Total	O	0	0
			11	11		
4	B	13	Total	O	0	0
			13	13		
4	C	11	Total	O	0	0
			11	11		
4	D	11	Total	O	0	0
			11	11		

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.

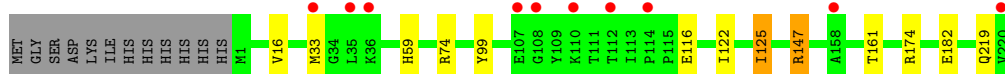
- Molecule 1: Thymidylate synthase thyX

Chain A: 



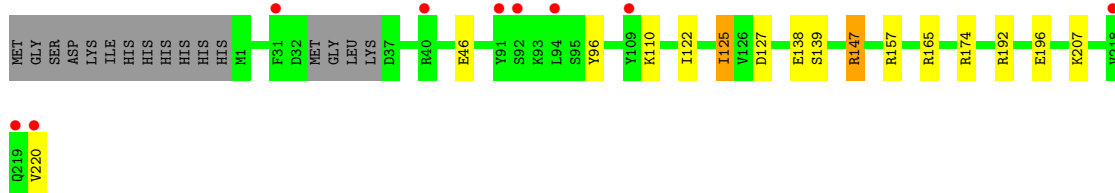
- Molecule 1: Thymidylate synthase thyX

Chain B: 



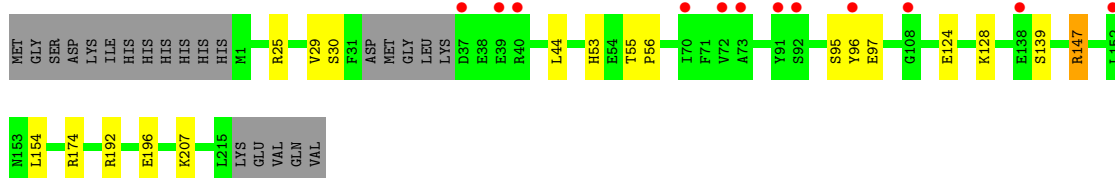
- Molecule 1: Thymidylate synthase thyX

Chain C: 



- Molecule 1: Thymidylate synthase thyX

Chain D: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	54.55Å 116.72Å 141.20Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.30 29.99 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.3 (30.00-2.30) 99.4 (29.99-2.30)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.73 (at 2.31Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.193 , 0.248 0.191 , 0.246	Depositor DCC
R_{free} test set	2037 reflections (5.28%)	DCC
Wilson B-factor (Å ²)	34.9	Xtriage
Anisotropy	0.030	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 25.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 40619 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7471	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.66% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: UMP, 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	0.87	1/1824 (0.1%)	0.83	1/2465 (0.0%)
1	B	0.84	0/1867	0.79	1/2522 (0.0%)
1	C	0.87	0/1838	0.82	1/2483 (0.0%)
1	D	0.85	0/1785	0.77	0/2414
All	All	0.86	1/7314 (0.0%)	0.80	3/9884 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	138	GLU	CG-CD	5.32	1.59	1.51

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	74	ARG	NE-CZ-NH2	-9.29	115.65	120.30
1	C	147	ARG	NE-CZ-NH2	-5.87	117.36	120.30
1	B	74	ARG	NE-CZ-NH2	-5.32	117.64	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1778	0	0	4	0
1	B	1821	0	0	6	0
1	C	1793	0	0	7	0
1	D	1741	0	0	9	0
2	A	20	0	11	1	0
2	B	20	0	11	1	0
2	C	20	0	11	1	0
2	D	20	0	11	1	0
3	A	53	0	31	2	0
3	B	53	0	31	1	0
3	C	53	0	31	0	0
3	D	53	0	31	0	0
4	A	11	0	0	0	0
4	B	13	0	0	2	0
4	C	11	0	0	0	0
4	D	11	0	0	0	0
All	All	7471	0	168	24	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 3.

All (24) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:95:SER:O	1:D:97:GLU:N	2.27	0.67
1:A:127:ASP:OD2	1:A:131:ARG:NH1	2.33	0.62
1:B:122:ILE:O	1:B:125:ILE:CD1	2.50	0.59
1:C:46:GLU:OE2	1:C:207:LYS:NZ	2.39	0.56
1:D:124:GLU:OE2	1:D:128:LYS:NZ	2.38	0.56
3:B:310:FAD:N3A	4:B:226:HOH:O	2.34	0.54
1:C:157:ARG:NH1	1:D:29:VAL:O	2.41	0.53
2:B:308:UMP:HN3	1:C:174:ARG:NH1	2.06	0.53
1:C:122:ILE:O	1:C:125:ILE:CD1	2.59	0.51
1:A:179:ALA:O	1:A:184:GLN:NE2	2.44	0.51
1:D:147:ARG:NH1	2:D:318:UMP:OP1	2.46	0.49
1:B:174:ARG:NH1	2:C:313:UMP:HN3	2.12	0.48
1:C:192:ARG:CZ	1:C:220:VAL:CG2	2.92	0.47
2:A:303:UMP:HN3	1:D:174:ARG:NH1	2.12	0.47
1:A:59:HIS:ND1	1:B:161:THR:OG1	2.48	0.46
3:A:315:FAD:H2B	1:C:165:ARG:CZ	2.45	0.46
1:A:161:THR:OG1	1:B:59:HIS:ND1	2.50	0.45
1:D:192:ARG:NH1	1:D:196:GLU:OE1	2.49	0.45
1:B:147:ARG:NH1	4:B:228:HOH:O	2.50	0.44
1:D:147:ARG:C	1:D:147:ARG:CD	2.86	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:A:315:FAD:O3B	1:C:165:ARG:NH2	2.52	0.43
1:B:99:TYR:OH	1:B:182:GLU:OE2	2.36	0.43
1:D:55:THR:N	1:D:56:PRO:CD	2.82	0.43
1:D:25:ARG:O	1:D:29:VAL:N	2.52	0.42

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	214/232 (92%)	207 (97%)	7 (3%)	0	100	100
1	B	218/232 (94%)	210 (96%)	8 (4%)	0	100	100
1	C	212/232 (91%)	205 (97%)	6 (3%)	1 (0%)	38	45
1	D	206/232 (89%)	202 (98%)	3 (2%)	1 (0%)	38	45
All	All	850/928 (92%)	824 (97%)	24 (3%)	2 (0%)	56	68

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	96	TYR
1	D	96	TYR

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	187/205 (91%)	178 (95%)	9 (5%)	35	46

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	193/205 (94%)	187 (97%)	6 (3%)	52	68
1	C	190/205 (93%)	183 (96%)	7 (4%)	45	60
1	D	184/205 (90%)	177 (96%)	7 (4%)	44	59
All	All	754/820 (92%)	725 (96%)	29 (4%)	44	59

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	29	VAL
1	A	40	ARG
1	A	44	LEU
1	A	87	LEU
1	A	96	TYR
1	A	106	LEU
1	A	139	SER
1	A	147	ARG
1	A	206	LEU
1	B	16	VAL
1	B	33	MET
1	B	116	GLU
1	B	125	ILE
1	B	147	ARG
1	B	219	GLN
1	C	110	LYS
1	C	125	ILE
1	C	127	ASP
1	C	138	GLU
1	C	139	SER
1	C	147	ARG
1	C	196	GLU
1	D	30	SER
1	D	44	LEU
1	D	53	HIS
1	D	139	SER
1	D	147	ARG
1	D	154	LEU
1	D	207	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

8 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	UMP	A	303	-	21,21,21	1.12	1 (4%)	26,31,31	1.07	1 (3%)
3	FAD	A	315	-	58,58,58	3.09	17 (29%)	85,89,89	2.84	23 (27%)
2	UMP	B	308	-	21,21,21	0.98	2 (9%)	26,31,31	1.50	2 (7%)
3	FAD	B	310	-	58,58,58	2.83	17 (29%)	85,89,89	2.70	18 (21%)
3	FAD	C	305	-	58,58,58	2.76	15 (25%)	85,89,89	2.76	19 (22%)
2	UMP	C	313	-	21,21,21	0.91	1 (4%)	26,31,31	1.47	2 (7%)
3	FAD	D	300	-	58,58,58	2.86	18 (31%)	85,89,89	2.73	20 (23%)
2	UMP	D	318	-	21,21,21	1.51	5 (23%)	26,31,31	1.18	3 (11%)

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	UMP	A	303	-	-	0/7/22/22	0/2/2/2
3	FAD	A	315	-	-	0/34/50/50	0/1/6/6
2	UMP	B	308	-	-	0/7/22/22	0/2/2/2
3	FAD	B	310	-	-	0/34/50/50	0/1/6/6
3	FAD	C	305	-	-	0/34/50/50	0/1/6/6
2	UMP	C	313	-	-	0/7/22/22	0/2/2/2
3	FAD	D	300	-	-	0/34/50/50	0/1/6/6
2	UMP	D	318	-	-	0/7/22/22	0/2/2/2

All (76) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	315	FAD	C6-C5X	10.50	1.54	1.41
3	D	300	FAD	C6-C5X	9.81	1.53	1.41
3	B	310	FAD	C6-C5X	9.22	1.52	1.41
3	C	305	FAD	C6-C5X	8.70	1.52	1.41
3	A	315	FAD	C4X-C10	8.06	1.54	1.40
3	C	305	FAD	C4X-C10	7.75	1.54	1.40
3	B	310	FAD	C4X-C10	7.49	1.53	1.40
3	A	315	FAD	C5X-N5	7.28	1.46	1.35
3	D	300	FAD	C4X-C10	6.84	1.52	1.40
3	B	310	FAD	C9A-N10	-6.58	1.29	1.38
3	A	315	FAD	C4-C4X	6.51	1.51	1.41
3	A	315	FAD	C10-N1	6.46	1.47	1.35
3	D	300	FAD	C5X-N5	6.29	1.45	1.35
3	A	315	FAD	C9A-N10	-6.26	1.29	1.38
3	C	305	FAD	C9A-N10	-6.08	1.29	1.38
3	D	300	FAD	C10-N1	6.03	1.46	1.35
3	A	315	FAD	C6-C7	5.83	1.54	1.37
3	C	305	FAD	C10-N1	5.80	1.46	1.35
3	B	310	FAD	C5X-N5	5.75	1.44	1.35
3	C	305	FAD	C5X-N5	5.73	1.44	1.35
3	D	300	FAD	C9A-N10	-5.72	1.30	1.38
3	C	305	FAD	C4-C4X	5.71	1.50	1.41
3	B	310	FAD	C10-N1	5.63	1.45	1.35
3	B	310	FAD	C4-C4X	5.62	1.50	1.41
3	D	300	FAD	C6-C7	5.54	1.53	1.37
3	B	310	FAD	C6-C7	5.35	1.52	1.37
3	C	305	FAD	C6-C7	5.10	1.51	1.37
3	D	300	FAD	C4-C4X	5.09	1.49	1.41
3	A	315	FAD	C9A-C5X	4.85	1.52	1.42
3	D	300	FAD	C9A-C5X	4.76	1.52	1.42
3	D	300	FAD	C9-C9A	4.56	1.49	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	310	FAD	C10-N10	4.55	1.48	1.38
3	A	315	FAD	C4X-N5	4.49	1.45	1.36
2	D	318	UMP	C2-N1	4.47	1.43	1.38
3	B	310	FAD	C9A-C5X	4.46	1.51	1.42
3	D	300	FAD	C10-N10	4.43	1.48	1.38
3	C	305	FAD	C5A-C4A	4.31	1.50	1.40
3	D	300	FAD	C2-N1	-4.23	1.25	1.35
3	C	305	FAD	C8-C7	4.21	1.53	1.40
3	C	305	FAD	C9A-C5X	4.18	1.51	1.42
3	C	305	FAD	C10-N10	4.11	1.47	1.38
3	A	315	FAD	C10-N10	4.06	1.47	1.38
3	A	315	FAD	C5A-C4A	3.96	1.49	1.40
3	A	315	FAD	C8-C7	3.95	1.52	1.40
3	B	310	FAD	C8-C7	3.73	1.51	1.40
3	D	300	FAD	C8-C7	3.62	1.51	1.40
3	D	300	FAD	C4A-N9A	-3.60	1.32	1.37
3	A	315	FAD	C4A-N9A	-3.59	1.32	1.37
3	B	310	FAD	C2-N1	-3.56	1.26	1.35
3	C	305	FAD	C9-C9A	3.49	1.47	1.40
3	B	310	FAD	C4X-N5	3.48	1.43	1.36
3	B	310	FAD	C9-C9A	3.36	1.47	1.40
3	B	310	FAD	C4A-N9A	-3.28	1.33	1.37
3	D	300	FAD	C5A-C4A	3.23	1.47	1.40
3	B	310	FAD	C5A-C4A	3.18	1.47	1.40
3	A	315	FAD	C2-N1	-3.13	1.28	1.35
3	A	315	FAD	C9-C9A	3.13	1.46	1.40
3	C	305	FAD	C4X-N5	3.11	1.42	1.36
3	D	300	FAD	C4X-N5	3.10	1.42	1.36
3	D	300	FAD	C1'-N10	-2.87	1.44	1.48
3	C	305	FAD	C2-N1	-2.85	1.28	1.35
3	C	305	FAD	C4A-N9A	-2.74	1.33	1.37
2	D	318	UMP	C4-N3	-2.70	1.32	1.37
3	A	315	FAD	C2A-N3A	2.63	1.37	1.32
2	B	308	UMP	C4-N3	-2.58	1.32	1.37
2	B	308	UMP	C6-C5	2.56	1.40	1.36
2	C	313	UMP	C6-C5	2.55	1.40	1.36
3	A	315	FAD	C1'-C2'	2.45	1.53	1.51
2	D	318	UMP	O2-C2	2.44	1.28	1.23
2	D	318	UMP	C2-N3	-2.34	1.32	1.37
3	B	310	FAD	PA-O3P	2.19	1.63	1.59
2	A	303	UMP	O2-C2	2.15	1.27	1.23
3	D	300	FAD	C1'-C2'	2.14	1.53	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	318	UMP	C6-C5	2.14	1.39	1.36
3	D	300	FAD	C4'-C3'	-2.13	1.49	1.53
3	B	310	FAD	C8A-N9A	-2.06	1.33	1.36

All (88) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	315	FAD	C1'-N10-C9A	11.22	129.79	118.87
3	A	315	FAD	C2-N1-C10	11.11	126.18	114.98
3	C	305	FAD	C4X-C10-N1	-11.02	111.72	122.73
3	A	315	FAD	C4X-C10-N1	-10.55	112.19	122.73
3	B	310	FAD	C1'-N10-C9A	10.50	129.09	118.87
3	C	305	FAD	C1'-N10-C9A	10.15	128.75	118.87
3	B	310	FAD	C4X-C10-N1	-9.85	112.89	122.73
3	B	310	FAD	C2-N1-C10	9.61	124.66	114.98
3	D	300	FAD	C2-N1-C10	9.52	124.58	114.98
3	D	300	FAD	C1'-N10-C9A	9.52	128.13	118.87
3	C	305	FAD	C2-N1-C10	9.51	124.57	114.98
3	D	300	FAD	C4X-C10-N1	-9.24	113.50	122.73
3	D	300	FAD	C1'-N10-C10	-7.85	108.03	119.17
3	D	300	FAD	N3A-C2A-N1A	-7.55	122.40	128.71
3	A	315	FAD	C1'-N10-C10	-7.34	108.76	119.17
3	B	310	FAD	N3A-C2A-N1A	-7.19	122.70	128.71
3	B	310	FAD	C1'-N10-C10	-7.11	109.07	119.17
3	C	305	FAD	C1'-N10-C10	-6.74	109.60	119.17
3	D	300	FAD	C4X-C10-N10	-6.39	117.32	120.51
3	C	305	FAD	N3A-C2A-N1A	-6.20	123.52	128.71
3	A	315	FAD	N3A-C2A-N1A	-5.96	123.73	128.71
2	B	308	UMP	N3-C2-N1	5.64	120.68	115.97
3	D	300	FAD	N3A-C4A-N9A	5.38	135.14	125.43
2	C	313	UMP	N3-C2-N1	5.34	120.43	115.97
3	B	310	FAD	N3A-C4A-N9A	5.23	134.88	125.43
3	A	315	FAD	N3A-C4A-N9A	5.22	134.87	125.43
3	C	305	FAD	N3A-C4A-N9A	5.16	134.75	125.43
3	C	305	FAD	C6-C5X-N5	-4.99	113.16	118.97
3	A	315	FAD	C5X-C9A-N10	4.67	121.40	116.80
3	D	300	FAD	C6-C5X-N5	-4.51	113.72	118.97
3	D	300	FAD	C7-C6-C5X	-4.48	113.38	120.91
3	A	315	FAD	P-O3P-PA	-4.35	118.94	131.68
3	C	305	FAD	C7-C6-C5X	-4.27	113.74	120.91
3	B	310	FAD	C7-C6-C5X	-4.26	113.76	120.91
3	D	300	FAD	N1-C10-N10	4.03	126.57	115.97

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	315	FAD	C7-C6-C5X	-3.98	114.22	120.91
3	C	305	FAD	N1-C10-N10	3.95	126.37	115.97
3	C	305	FAD	C2B-C1B-N9A	3.87	123.19	113.27
3	C	305	FAD	C4A-C5A-N7A	-3.87	106.21	109.52
3	B	310	FAD	C5X-C9A-N10	3.81	120.55	116.80
3	C	305	FAD	C5A-C4A-N3A	-3.77	117.49	125.70
3	A	315	FAD	C4A-C5A-N7A	-3.72	106.33	109.52
3	B	310	FAD	C6-C5X-N5	-3.63	114.74	118.97
3	B	310	FAD	N1-C10-N10	3.58	125.39	115.97
3	B	310	FAD	C3B-C2B-C1B	3.56	106.48	100.91
3	C	305	FAD	O4B-C1B-N9A	-3.55	105.14	108.44
3	B	310	FAD	C4X-C10-N10	-3.55	118.74	120.51
3	A	315	FAD	C10-C4X-N5	-3.55	116.15	120.45
3	C	305	FAD	C5X-C9A-N10	3.54	120.28	116.80
3	B	310	FAD	C2B-C1B-N9A	3.38	121.94	113.27
3	A	315	FAD	N1-C10-N10	3.37	124.83	115.97
3	D	300	FAD	C8A-N9A-C4A	3.13	109.29	106.90
3	D	300	FAD	C5A-C4A-N3A	-3.05	119.05	125.70
3	A	315	FAD	C5A-C4A-N3A	-3.05	119.06	125.70
3	A	315	FAD	C9-C8-C7	2.97	124.63	119.88
3	B	310	FAD	C5A-C4A-N3A	-2.94	119.30	125.70
3	C	305	FAD	C10-C4X-N5	-2.87	116.97	120.45
3	D	300	FAD	C3B-C2B-C1B	2.71	105.14	100.91
3	C	305	FAD	C3B-C2B-C1B	2.60	104.97	100.91
3	B	310	FAD	C4A-C5A-N7A	-2.52	107.36	109.52
3	A	315	FAD	C4B-O4B-C1B	2.52	112.48	109.75
3	A	315	FAD	C6-C5X-N5	-2.49	116.07	118.97
3	D	300	FAD	C4A-C5A-N7A	-2.49	107.39	109.52
3	B	310	FAD	C10-C4X-N5	-2.49	117.43	120.45
2	A	303	UMP	N3-C2-N1	2.49	118.05	115.97
3	A	315	FAD	C8A-N9A-C4A	2.47	108.79	106.90
2	D	318	UMP	OP3-P-OP1	2.41	118.32	110.44
2	C	313	UMP	C6-N1-C1'	2.37	124.84	119.35
2	B	308	UMP	OP3-P-OP1	2.28	117.89	110.44
3	D	300	FAD	O3'-C3'-C4'	-2.27	102.99	108.74
3	D	300	FAD	C5'-C4'-C3'	-2.26	107.80	112.06
3	C	305	FAD	C7M-C7-C6	-2.25	114.95	120.38
3	D	300	FAD	C8M-C8-C7	-2.24	115.57	120.74
3	C	305	FAD	C4X-C10-N10	-2.23	119.40	120.51
3	A	315	FAD	O3'-C3'-C4'	-2.21	103.14	108.74
2	D	318	UMP	C5-C4-N3	2.20	120.84	116.70
3	D	300	FAD	C2A-N3A-C4A	2.19	120.25	114.01

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	315	FAD	C2B-C1B-N9A	2.17	118.84	113.27
3	A	315	FAD	C9A-C5X-N5	-2.16	119.05	122.37
3	B	310	FAD	O4B-C1B-N9A	-2.14	106.45	108.44
3	A	315	FAD	O2A-PA-O1A	2.13	124.10	112.21
3	B	310	FAD	C2'-C1'-N10	-2.09	109.68	112.45
3	D	300	FAD	O2A-PA-O1A	2.08	123.81	112.21
3	A	315	FAD	C6-C7-C8	-2.05	116.60	119.88
3	D	300	FAD	O2B-C2B-C1B	-2.04	105.05	111.23
3	C	305	FAD	O2A-PA-O1A	2.03	123.58	112.21
2	D	318	UMP	C5-C6-N1	-2.02	118.93	121.21
3	A	315	FAD	O3P-PA-O5B	-2.00	94.45	103.41

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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%)	-0.07	9 (4%) 35 45	18, 29, 60, 66	0
1	B	220/232 (94%)	0.07	10 (4%) 32 42	17, 31, 60, 72	0
1	C	216/232 (93%)	-0.12	9 (4%) 35 45	16, 27, 53, 62	0
1	D	210/232 (90%)	0.08	12 (5%) 23 32	17, 29, 56, 65	0
All	All	862/928 (92%)	-0.01	40 (4%) 31 41	16, 29, 58, 72	0

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	33	MET	4.1
1	D	96	TYR	4.0
1	D	91	TYR	3.7
1	C	219	GLN	3.7
1	B	35	LEU	3.7
1	D	40	ARG	3.7
1	A	40	ARG	3.5
1	C	109	TYR	3.2
1	D	92	SER	3.1
1	B	112	THR	3.1
1	A	34	GLY	3.0
1	B	108	GLY	2.8
1	C	220	VAL	2.8
1	A	112	THR	2.8
1	A	32	ASP	2.7
1	C	40	ARG	2.6
1	D	37	ASP	2.6
1	D	70	ILE	2.6
1	A	36	LYS	2.6
1	D	108	GLY	2.6
1	C	94	LEU	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	158	ALA	2.5
1	C	31	PHE	2.5
1	D	138	GLU	2.4
1	B	220	VAL	2.4
1	A	31	PHE	2.4
1	C	91	TYR	2.4
1	A	108	GLY	2.4
1	C	92	SER	2.3
1	B	33	MET	2.3
1	D	73	ALA	2.2
1	C	218	VAL	2.2
1	D	39	GLU	2.2
1	B	114	PRO	2.2
1	D	152	LEU	2.2
1	B	110	LYS	2.2
1	B	107	GLU	2.1
1	A	0	HIS	2.1
1	D	72	VAL	2.1
1	B	36	LYS	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	FAD	A	315	53/53	0.18	1.43	32,45,51,52	0
3	FAD	B	310	53/53	0.16	1.07	33,46,53,54	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	FAD	D	300	53/53	0.14	1.06	31,42,49,50	0
3	FAD	C	305	53/53	0.14	0.39	30,40,48,49	0
2	UMP	C	313	20/20	0.12	-0.29	31,35,41,41	0
2	UMP	D	318	20/20	0.13	-0.45	42,46,48,49	0
2	UMP	A	303	20/20	0.11	-0.49	25,27,30,31	0
2	UMP	B	308	20/20	0.11	-0.69	26,31,35,38	0

6.5 Other polymers ⓘ

There are no such residues in this entry.