



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 15, 2017 – 01:20 am GMT

PDB ID : 2XYO
Title : STRUCTURAL BASIS FOR A NEW TETRACYCLINE RESISTANCE
MECHANISM RELYING ON THE TETX MONOOXYGENASE
Authors : Volkers, G.; Palm, G.J.; Weiss, M.S.; Hinrichs, W.
Deposited on : 2010-11-18
Resolution : 3.00 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	:	trunk28620
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)	:	recalc28949

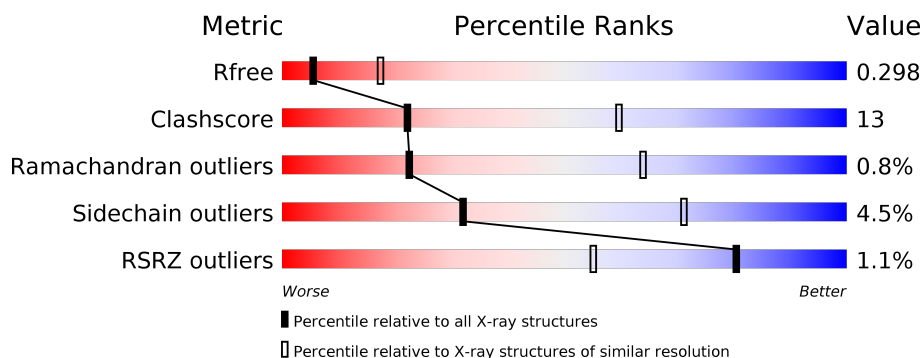
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	1692 (3.00-3.00)
Clashscore	112137	2037 (3.00-3.00)
Ramachandran outliers	110173	1973 (3.00-3.00)
Sidechain outliers	110143	1976 (3.00-3.00)
RSRZ outliers	101464	1716 (3.00-3.00)

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	399	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 24%, green 66%, grey 8%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 66% 24% • 8% </div> </div>
1	B	399	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 24%, green 67%, grey 8%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 67% 24% • 8% </div> </div>
1	C	399	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 22%, green 69%, grey 8%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 69% 22% • 8% </div> </div>
1	D	399	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 22%, green 69%, grey 8%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 69% 22% • 8% </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	SO4	B	1386	-	-	-	X
3	SO4	D	1386	-	-	-	X
4	GOL	B	1387	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 11747 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 TETX2.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	367	Total	C	N	O	S	Se	0	3	0
			2880	1824	491	553	2	10			
1	B	367	Total	C	N	O	S	Se	0	3	0
			2882	1825	492	553	2	10			
1	C	366	Total	C	N	O	S	Se	0	2	0
			2865	1815	488	550	2	10			
1	D	366	Total	C	N	O	S	Se	0	2	0
			2862	1814	486	550	2	10			

There are 84 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-10	SER	-	EXPRESSION TAG	UNP Q93L51
A	-9	MSE	-	EXPRESSION TAG	UNP Q93L51
A	-8	GLY	-	EXPRESSION TAG	UNP Q93L51
A	-7	SER	-	EXPRESSION TAG	UNP Q93L51
A	-6	SER	-	EXPRESSION TAG	UNP Q93L51
A	-5	HIS	-	EXPRESSION TAG	UNP Q93L51
A	-4	HIS	-	EXPRESSION TAG	UNP Q93L51
A	-3	HIS	-	EXPRESSION TAG	UNP Q93L51
A	-2	HIS	-	EXPRESSION TAG	UNP Q93L51
A	-1	HIS	-	EXPRESSION TAG	UNP Q93L51
A	0	HIS	-	EXPRESSION TAG	UNP Q93L51
A	1	SER	-	EXPRESSION TAG	UNP Q93L51
A	2	SER	-	EXPRESSION TAG	UNP Q93L51
A	3	GLY	-	EXPRESSION TAG	UNP Q93L51
A	4	LEU	-	EXPRESSION TAG	UNP Q93L51
A	5	VAL	-	EXPRESSION TAG	UNP Q93L51
A	6	PRO	-	EXPRESSION TAG	UNP Q93L51
A	7	ARG	-	EXPRESSION TAG	UNP Q93L51
A	8	GLY	-	EXPRESSION TAG	UNP Q93L51
A	9	SER	-	EXPRESSION TAG	UNP Q93L51
A	10	HIS	-	EXPRESSION TAG	UNP Q93L51

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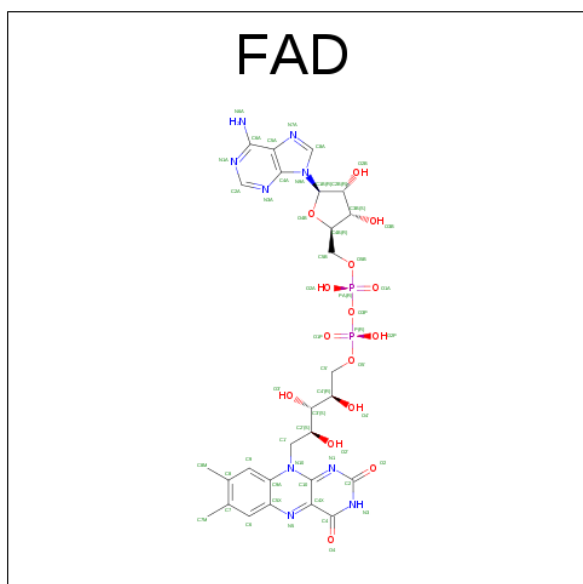
Chain	Residue	Modelled	Actual	Comment	Reference
B	-10	SER	-	EXPRESSION TAG	UNP Q93L51
B	-9	MSE	-	EXPRESSION TAG	UNP Q93L51
B	-8	GLY	-	EXPRESSION TAG	UNP Q93L51
B	-7	SER	-	EXPRESSION TAG	UNP Q93L51
B	-6	SER	-	EXPRESSION TAG	UNP Q93L51
B	-5	HIS	-	EXPRESSION TAG	UNP Q93L51
B	-4	HIS	-	EXPRESSION TAG	UNP Q93L51
B	-3	HIS	-	EXPRESSION TAG	UNP Q93L51
B	-2	HIS	-	EXPRESSION TAG	UNP Q93L51
B	-1	HIS	-	EXPRESSION TAG	UNP Q93L51
B	0	HIS	-	EXPRESSION TAG	UNP Q93L51
B	1	SER	-	EXPRESSION TAG	UNP Q93L51
B	2	SER	-	EXPRESSION TAG	UNP Q93L51
B	3	GLY	-	EXPRESSION TAG	UNP Q93L51
B	4	LEU	-	EXPRESSION TAG	UNP Q93L51
B	5	VAL	-	EXPRESSION TAG	UNP Q93L51
B	6	PRO	-	EXPRESSION TAG	UNP Q93L51
B	7	ARG	-	EXPRESSION TAG	UNP Q93L51
B	8	GLY	-	EXPRESSION TAG	UNP Q93L51
B	9	SER	-	EXPRESSION TAG	UNP Q93L51
B	10	HIS	-	EXPRESSION TAG	UNP Q93L51
C	-10	SER	-	EXPRESSION TAG	UNP Q93L51
C	-9	MSE	-	EXPRESSION TAG	UNP Q93L51
C	-8	GLY	-	EXPRESSION TAG	UNP Q93L51
C	-7	SER	-	EXPRESSION TAG	UNP Q93L51
C	-6	SER	-	EXPRESSION TAG	UNP Q93L51
C	-5	HIS	-	EXPRESSION TAG	UNP Q93L51
C	-4	HIS	-	EXPRESSION TAG	UNP Q93L51
C	-3	HIS	-	EXPRESSION TAG	UNP Q93L51
C	-2	HIS	-	EXPRESSION TAG	UNP Q93L51
C	-1	HIS	-	EXPRESSION TAG	UNP Q93L51
C	0	HIS	-	EXPRESSION TAG	UNP Q93L51
C	1	SER	-	EXPRESSION TAG	UNP Q93L51
C	2	SER	-	EXPRESSION TAG	UNP Q93L51
C	3	GLY	-	EXPRESSION TAG	UNP Q93L51
C	4	LEU	-	EXPRESSION TAG	UNP Q93L51
C	5	VAL	-	EXPRESSION TAG	UNP Q93L51
C	6	PRO	-	EXPRESSION TAG	UNP Q93L51
C	7	ARG	-	EXPRESSION TAG	UNP Q93L51
C	8	GLY	-	EXPRESSION TAG	UNP Q93L51
C	9	SER	-	EXPRESSION TAG	UNP Q93L51
C	10	HIS	-	EXPRESSION TAG	UNP Q93L51

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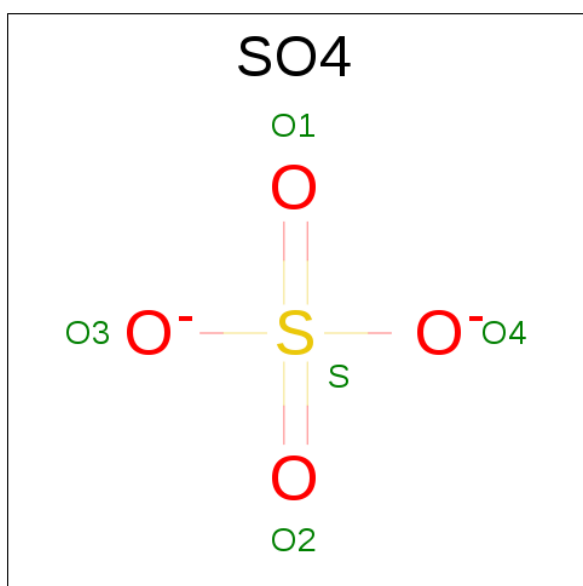
Chain	Residue	Modelled	Actual	Comment	Reference
D	-10	SER	-	EXPRESSION TAG	UNP Q93L51
D	-9	MSE	-	EXPRESSION TAG	UNP Q93L51
D	-8	GLY	-	EXPRESSION TAG	UNP Q93L51
D	-7	SER	-	EXPRESSION TAG	UNP Q93L51
D	-6	SER	-	EXPRESSION TAG	UNP Q93L51
D	-5	HIS	-	EXPRESSION TAG	UNP Q93L51
D	-4	HIS	-	EXPRESSION TAG	UNP Q93L51
D	-3	HIS	-	EXPRESSION TAG	UNP Q93L51
D	-2	HIS	-	EXPRESSION TAG	UNP Q93L51
D	-1	HIS	-	EXPRESSION TAG	UNP Q93L51
D	0	HIS	-	EXPRESSION TAG	UNP Q93L51
D	1	SER	-	EXPRESSION TAG	UNP Q93L51
D	2	SER	-	EXPRESSION TAG	UNP Q93L51
D	3	GLY	-	EXPRESSION TAG	UNP Q93L51
D	4	LEU	-	EXPRESSION TAG	UNP Q93L51
D	5	VAL	-	EXPRESSION TAG	UNP Q93L51
D	6	PRO	-	EXPRESSION TAG	UNP Q93L51
D	7	ARG	-	EXPRESSION TAG	UNP Q93L51
D	8	GLY	-	EXPRESSION TAG	UNP Q93L51
D	9	SER	-	EXPRESSION TAG	UNP Q93L51
D	10	HIS	-	EXPRESSION TAG	UNP Q93L51

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



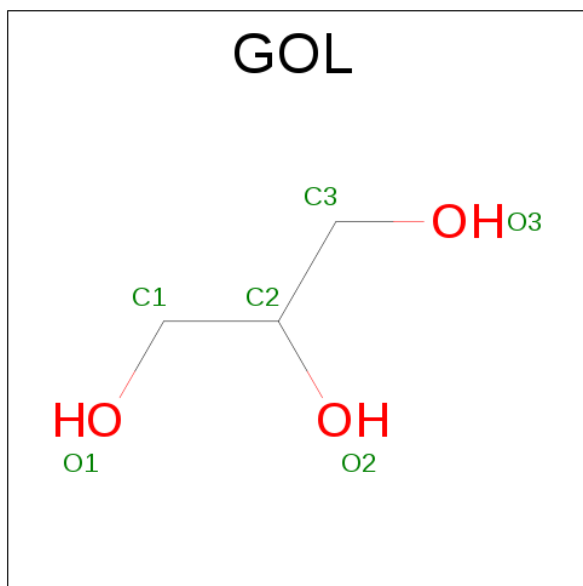
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	C	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	D	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).

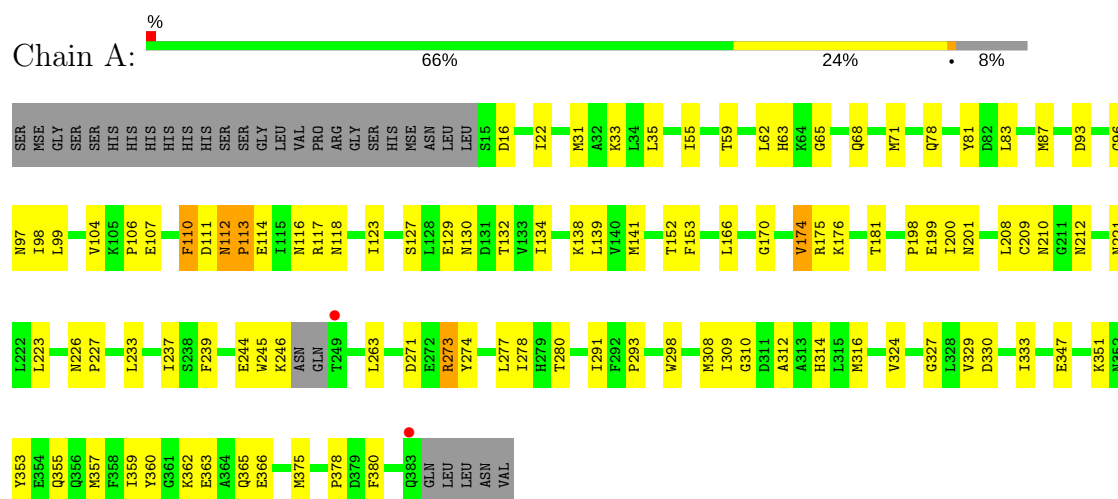


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			6	3	3		

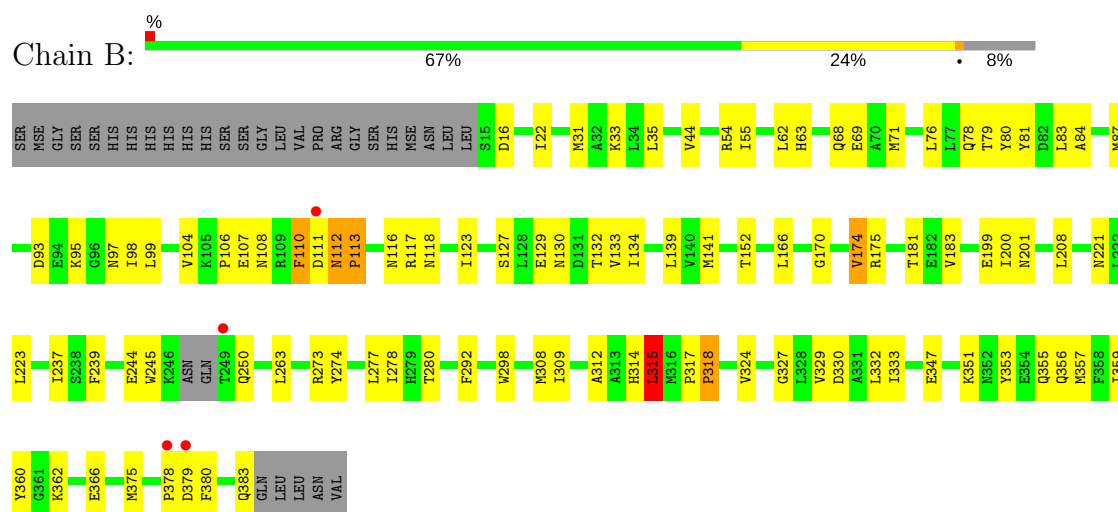
3 Residue-property plots

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

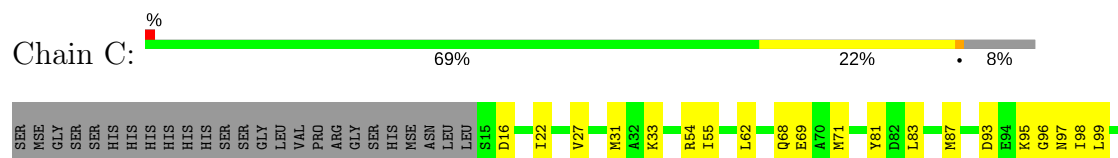
• Molecule 1: TETX2



• Molecule 1: TETX2

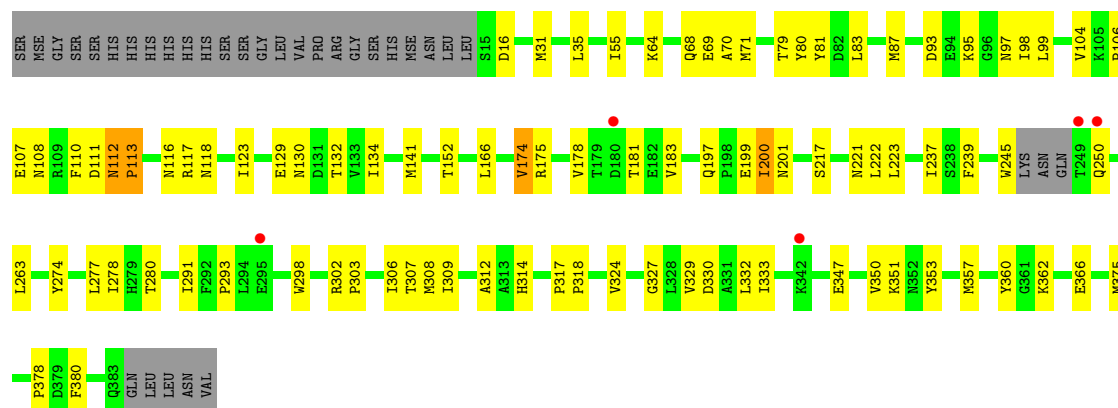


• Molecule 1: TETX2





- Molecule 1: TETX2



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	87.34Å 68.66Å 152.48Å 90.00° 101.08° 90.00°	Depositor
Resolution (Å)	149.64 – 3.00 29.80 – 3.00	Depositor EDS
% Data completeness (in resolution range)	99.4 (149.64-3.00) 99.5 (29.80-3.00)	Depositor EDS
R_{merge}	0.22	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.23 (at 3.00Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.249 , 0.295 0.249 , 0.298	Depositor DCC
R_{free} test set	1783 reflections (5.25%)	DCC
Wilson B-factor (Å ²)	41.7	Xtriage
Anisotropy	0.409	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 37.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	11747	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 59.16 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 1.8384e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: GOL, SO4, 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	0.68	0/2937	0.74	0/3962
1	B	0.68	0/2939	0.77	1/3964 (0.0%)
1	C	0.62	0/2919	0.70	0/3938
1	D	0.61	0/2916	0.72	0/3934
All	All	0.65	0/11711	0.73	1/15798 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	315	LEU	CB-CG-CD2	6.31	121.73	111.00

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	2880	0	2780	81	0
1	B	2882	0	2787	96	0
1	C	2865	0	2768	78	0
1	D	2862	0	2760	63	0
2	A	53	0	31	3	0
2	B	53	0	31	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	53	0	31	7	0
2	D	53	0	31	3	0
3	A	10	0	0	0	0
3	B	10	0	0	0	0
3	C	10	0	0	0	0
3	D	10	0	0	0	0
4	B	6	0	8	4	0
All	All	11747	0	11227	304	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 13.

All (304) 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:B:104:VAL:HG12	1:B:106:PRO:HD2	1.46	0.98
1:D:107:GLU:O	1:D:110:PHE:HB2	1.65	0.96
1:B:107:GLU:O	1:B:110:PHE:HB2	1.64	0.96
1:C:107:GLU:O	1:C:110:PHE:HB2	1.67	0.94
1:A:107:GLU:O	1:A:110:PHE:HB2	1.66	0.94
1:A:175:ARG:HD3	1:A:308:MSE:HE1	1.49	0.92
1:A:104:VAL:HG12	1:A:106:PRO:HD2	1.50	0.92
1:D:104:VAL:HG12	1:D:106:PRO:HD2	1.53	0.91
1:C:175:ARG:HD3	1:C:308:MSE:HE1	1.55	0.89
1:C:104:VAL:HG12	1:C:106:PRO:HD2	1.56	0.88
1:D:31:MSE:HG2	1:D:309:ILE:CD1	2.06	0.86
1:D:175:ARG:HD3	1:D:308:MSE:HE1	1.57	0.86
1:B:112:ASN:H	1:B:113:PRO:CD	1.94	0.80
1:C:117:ARG:NH2	2:C:1384:FAD:O2'	2.15	0.80
1:D:112:ASN:H	1:D:113:PRO:CD	1.95	0.79
1:A:175:ARG:NH1	1:A:312:ALA:O	2.16	0.79
1:B:208:LEU:HG	1:B:273[B]:ARG:HH11	1.45	0.78
1:A:65:GLY:HA3	1:B:379:ASP:CB	2.13	0.78
1:A:112:ASN:H	1:A:113:PRO:CD	1.96	0.78
1:C:175:ARG:NH1	1:C:312:ALA:O	2.17	0.78
1:C:112:ASN:H	1:C:113:PRO:CD	1.98	0.77
1:A:175:ARG:CD	1:A:308:MSE:HE1	2.15	0.75
1:D:175:ARG:NH1	1:D:312:ALA:O	2.19	0.74
1:A:117:ARG:NH2	2:A:1384:FAD:O2'	2.21	0.74
1:A:208:LEU:HG	1:A:273[B]:ARG:HH11	1.53	0.74
1:A:174:VAL:HG13	1:A:174:VAL:O	1.87	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:175:ARG:CD	1:C:308:MSE:HE1	2.19	0.72
2:B:1384:FAD:H2'	2:B:1384:FAD:N1	2.03	0.72
1:B:117:ARG:NH2	2:B:1384:FAD:O2'	2.16	0.72
1:B:175:ARG:HD3	1:B:308:MSE:HE1	1.72	0.71
1:B:208:LEU:HG	1:B:273[B]:ARG:NH1	2.06	0.70
1:B:31:MSE:HG2	1:B:309:ILE:CD1	2.21	0.70
1:D:166:LEU:HB2	1:D:308:MSE:HG2	1.71	0.70
1:C:31:MSE:HG2	1:C:309:ILE:CD1	2.22	0.69
1:D:175:ARG:CD	1:D:308:MSE:HE1	2.22	0.69
1:A:244:GLU:CD	1:A:244:GLU:H	1.95	0.68
1:A:359:ILE:HG23	1:B:98:ILE:CG2	2.22	0.68
1:B:78:GLN:H	4:B:1387:GOL:H12	1.59	0.67
1:B:112:ASN:O	1:B:113:PRO:C	2.32	0.67
1:B:78:GLN:N	4:B:1387:GOL:H12	2.09	0.67
1:A:363:GLU:HG3	1:B:98:ILE:HD11	1.75	0.67
1:D:55:ILE:H	1:D:118:ASN:HD21	1.43	0.66
1:C:174:VAL:HG13	1:C:174:VAL:O	1.95	0.66
1:C:27:VAL:HB	2:C:1384:FAD:O1P	1.96	0.66
1:B:175:ARG:NH1	1:B:312:ALA:O	2.28	0.66
1:B:166:LEU:HB2	1:B:308:MSE:HG2	1.78	0.66
1:B:298:TRP:CH2	1:B:357:MSE:HE1	2.31	0.65
2:A:1384:FAD:N1	2:A:1384:FAD:H2'	2.12	0.65
1:D:117:ARG:NH2	2:D:1384:FAD:O2'	2.29	0.64
1:A:112:ASN:O	1:A:113:PRO:C	2.35	0.64
1:C:298:TRP:CH2	1:C:357:MSE:HE1	2.32	0.64
1:A:65:GLY:HA3	1:B:379:ASP:CG	2.19	0.63
1:B:175:ARG:HG2	1:B:308:MSE:HE1	1.79	0.63
1:A:359:ILE:HG23	1:B:98:ILE:HG22	1.80	0.62
1:C:129:GLU:O	1:C:132:THR:OG1	2.17	0.62
1:B:174:VAL:HG13	1:B:174:VAL:O	1.98	0.61
1:B:110:PHE:O	1:B:112:ASN:N	2.34	0.61
1:A:110:PHE:O	1:A:111:ASP:C	2.39	0.61
1:B:201:ASN:HB3	1:B:280:THR:HG21	1.82	0.61
1:A:166:LEU:HB2	1:A:308:MSE:HG2	1.81	0.61
1:A:31:MSE:HG2	1:A:309:ILE:CD1	2.30	0.60
1:C:55:ILE:H	1:C:118:ASN:HD21	1.49	0.60
1:A:308:MSE:HE2	1:A:312:ALA:HB1	1.83	0.60
2:C:1384:FAD:N1	2:C:1384:FAD:H2'	2.16	0.60
1:C:112:ASN:O	1:C:113:PRO:C	2.38	0.60
1:B:110:PHE:O	1:B:111:ASP:C	2.41	0.59
1:B:329:VAL:HG12	1:B:333:ILE:HD12	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:112:ASN:H	1:D:113:PRO:HD3	1.67	0.59
1:C:175:ARG:HG2	1:C:308:MSE:HE1	1.84	0.59
1:A:110:PHE:O	1:A:112:ASN:N	2.36	0.59
1:B:175:ARG:CD	1:B:308:MSE:HE1	2.31	0.59
1:A:201:ASN:HB3	1:A:280:THR:HG21	1.84	0.59
1:B:314:HIS:NE2	1:B:330:ASP:OD2	2.30	0.59
1:B:54[A]:ARG:HH21	1:C:54[A]:ARG:HG3	1.67	0.59
1:C:112:ASN:H	1:C:113:PRO:HD3	1.66	0.59
1:B:175:ARG:HG2	1:B:308:MSE:CE	2.32	0.59
1:C:308:MSE:HE2	1:C:312:ALA:HB1	1.85	0.59
1:C:314:HIS:NE2	1:C:330:ASP:OD2	2.35	0.58
1:D:175:ARG:HG2	1:D:308:MSE:HE1	1.85	0.58
1:A:298:TRP:CH2	1:A:357:MSE:HE1	2.38	0.58
1:D:347:GLU:O	1:D:351:LYS:HG2	2.03	0.58
1:A:112:ASN:H	1:A:113:PRO:HD3	1.67	0.58
1:C:175:ARG:CG	1:C:308:MSE:HE1	2.34	0.58
1:A:327:GLY:HA2	1:A:330:ASP:OD2	2.04	0.58
1:D:174:VAL:O	1:D:174:VAL:HG13	2.02	0.58
1:C:68:GLN:NE2	1:C:81:TYR:OH	2.37	0.57
1:B:68:GLN:NE2	1:B:81:TYR:OH	2.37	0.57
1:D:107:GLU:O	1:D:110:PHE:CB	2.45	0.57
1:A:329:VAL:HG12	1:A:333:ILE:HD12	1.85	0.57
1:C:166:LEU:HB2	1:C:308:MSE:HG2	1.85	0.57
1:D:112:ASN:O	1:D:113:PRO:C	2.41	0.57
1:B:107:GLU:O	1:B:110:PHE:CB	2.47	0.57
1:B:112:ASN:H	1:B:113:PRO:HD2	1.66	0.57
1:B:54[A]:ARG:HG3	1:C:54[A]:ARG:HH21	1.69	0.57
1:A:359:ILE:HB	1:B:383:GLN:CB	2.34	0.57
1:D:110:PHE:O	1:D:112:ASN:N	2.38	0.57
1:C:117:ARG:HH22	2:C:1384:FAD:HO2'	1.49	0.56
1:C:223:LEU:HD12	1:C:237:ILE:HG12	1.87	0.56
1:D:329:VAL:HG12	1:D:333:ILE:HD12	1.87	0.56
1:D:298:TRP:CH2	1:D:357:MSE:HE1	2.38	0.56
1:C:263:LEU:HB3	1:C:278:ILE:HD13	1.87	0.56
1:A:223:LEU:HD12	1:A:237:ILE:HG12	1.88	0.56
1:B:175:ARG:CG	1:B:308:MSE:HE1	2.34	0.56
1:D:175:ARG:CG	1:D:308:MSE:HE1	2.36	0.56
1:A:175:ARG:CG	1:A:308:MSE:HE1	2.36	0.56
1:C:175:ARG:HG2	1:C:308:MSE:CE	2.36	0.56
1:A:55:ILE:H	1:A:118:ASN:HD21	1.53	0.55
1:A:68:GLN:NE2	1:A:81:TYR:OH	2.39	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:201:ASN:HB3	1:D:280:THR:HG21	1.86	0.55
1:B:356:GLN:HA	1:B:359:ILE:HD12	1.88	0.55
1:C:107:GLU:O	1:C:110:PHE:CB	2.48	0.55
1:D:83:LEU:HD22	1:D:123:ILE:HD11	1.88	0.55
1:B:54[A]:ARG:HH21	1:C:54[A]:ARG:NE	2.05	0.54
1:C:201:ASN:HB3	1:C:280:THR:HG21	1.88	0.54
1:D:31:MSE:CG	1:D:309:ILE:CD1	2.83	0.54
1:C:324:VAL:HG12	2:C:1384:FAD:O2	2.06	0.54
1:C:329:VAL:HG12	1:C:333:ILE:HD12	1.89	0.54
1:A:112:ASN:N	1:A:113:PRO:CD	2.68	0.54
1:A:347:GLU:O	1:A:351:LYS:HG2	2.08	0.54
1:C:110:PHE:O	1:C:111:ASP:C	2.45	0.54
1:D:68:GLN:NE2	1:D:81:TYR:OH	2.41	0.54
1:A:107:GLU:O	1:A:110:PHE:CB	2.49	0.54
1:D:223:LEU:HD12	1:D:237:ILE:HG12	1.90	0.53
1:B:83:LEU:HD22	1:B:123:ILE:HD11	1.88	0.53
1:C:112:ASN:N	1:C:113:PRO:CD	2.71	0.53
1:B:308:MSE:HE2	1:B:312:ALA:HB1	1.91	0.53
1:A:175:ARG:HG2	1:A:308:MSE:HE1	1.91	0.52
1:B:54[A]:ARG:NH2	1:C:54[A]:ARG:CZ	2.72	0.52
1:C:347:GLU:O	1:C:351:LYS:HG2	2.09	0.52
1:D:175:ARG:HG2	1:D:308:MSE:CE	2.39	0.52
1:A:78:GLN:HG2	1:D:64:LYS:HB3	1.90	0.52
1:A:174:VAL:CG1	1:A:174:VAL:O	2.53	0.52
1:A:170:GLY:O	1:A:175:ARG:NH2	2.42	0.52
1:D:110:PHE:O	1:D:111:ASP:C	2.46	0.52
1:A:22:ILE:HG13	1:A:139:LEU:HD22	1.92	0.51
1:A:175:ARG:HG2	1:A:308:MSE:CE	2.39	0.51
1:D:362:LYS:O	1:D:366:GLU:HG2	2.11	0.51
1:C:291:ILE:O	1:C:293:PRO:HD3	2.10	0.51
1:D:314:HIS:NE2	1:D:330:ASP:OD2	2.35	0.51
1:D:357:MSE:HA	1:D:360:TYR:CZ	2.46	0.51
1:A:129:GLU:O	1:A:132:THR:OG1	2.22	0.50
1:C:110:PHE:O	1:C:112:ASN:N	2.45	0.50
1:B:223:LEU:HD12	1:B:237:ILE:HG12	1.93	0.50
1:D:324:VAL:HG12	2:D:1384:FAD:O2	2.12	0.50
1:D:31:MSE:CG	1:D:309:ILE:HD13	2.41	0.50
1:D:69:GLU:HG3	1:D:332:LEU:HD21	1.93	0.50
1:B:112:ASN:N	1:B:113:PRO:CD	2.66	0.50
1:D:31:MSE:HG2	1:D:309:ILE:HD13	1.87	0.50
1:D:221:ASN:HD22	1:D:239:PHE:HB3	1.77	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:87:MSE:H	1:A:112:ASN:HD21	1.60	0.50
1:C:274:TYR:O	1:C:277:LEU:HB3	2.12	0.50
1:B:79:THR:HG1	4:B:1387:GOL:HO1	1.59	0.49
1:B:55:ILE:H	1:B:118:ASN:HD21	1.60	0.49
1:C:83:LEU:HD22	1:C:123:ILE:HD11	1.94	0.49
1:A:209:CYS:HB3	1:A:227:PRO:HG3	1.95	0.49
1:B:263:LEU:HB3	1:B:278:ILE:HD13	1.95	0.49
1:B:274:TYR:O	1:B:277:LEU:HB3	2.11	0.49
1:B:129:GLU:O	1:B:132:THR:OG1	2.21	0.49
1:B:54[A]:ARG:CZ	1:C:54[A]:ARG:NH2	2.76	0.49
1:D:112:ASN:N	1:D:113:PRO:CD	2.66	0.49
1:D:93:ASP:HB3	1:D:98:ILE:HD11	1.95	0.49
1:B:375:MSE:HA	1:B:380:PHE:CG	2.48	0.49
1:B:347:GLU:O	1:B:351:LYS:HG2	2.13	0.48
1:D:263:LEU:HB3	1:D:278:ILE:HD13	1.94	0.48
1:B:362:LYS:O	1:B:366:GLU:HG2	2.13	0.48
1:C:108:ASN:C	1:C:110:PHE:H	2.16	0.48
1:A:87:MSE:HG3	1:A:114:GLU:HG3	1.96	0.47
1:A:59:THR:OG1	1:A:114:GLU:OE1	2.24	0.47
1:A:221:ASN:HD22	1:A:239:PHE:HB3	1.79	0.47
1:A:314:HIS:NE2	1:A:330:ASP:OD2	2.38	0.47
1:D:129:GLU:O	1:D:132:THR:OG1	2.24	0.47
1:D:308:MSE:HE2	1:D:312:ALA:HB1	1.96	0.47
1:B:54[A]:ARG:NE	1:C:54[A]:ARG:HH21	2.13	0.47
1:C:174:VAL:CG1	1:C:174:VAL:O	2.63	0.47
1:B:54[A]:ARG:NH2	1:C:54[A]:ARG:HG3	2.29	0.47
1:B:22:ILE:HG13	1:B:139:LEU:HD22	1.97	0.47
1:B:357:MSE:HA	1:B:360:TYR:CZ	2.50	0.47
1:B:93:ASP:HB3	1:B:98:ILE:HD11	1.97	0.47
1:A:138:LYS:O	1:A:153:PHE:HA	2.14	0.46
1:A:274:TYR:O	1:A:277:LEU:HB3	2.15	0.46
1:D:108:ASN:C	1:D:110:PHE:H	2.18	0.46
1:B:329:VAL:HG12	1:B:333:ILE:CD1	2.46	0.46
1:B:84:ALA:HB1	1:B:113:PRO:CB	2.46	0.46
1:A:362:LYS:O	1:A:366:GLU:HG2	2.15	0.46
1:A:83:LEU:HD22	1:A:123:ILE:HD11	1.98	0.46
1:A:360:TYR:CD2	1:B:383:GLN:CB	2.98	0.46
1:C:22:ILE:HD11	1:C:151:LEU:HD23	1.98	0.46
1:A:245:TRP:CD1	1:A:245:TRP:N	2.83	0.46
1:A:360:TYR:HD2	1:B:383:GLN:CB	2.29	0.46
1:C:31:MSE:SE	1:C:331:ALA:HA	2.65	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:87:MSE:HG3	1:C:114:GLU:HG3	1.98	0.46
1:B:221:ASN:HD22	1:B:239:PHE:HB3	1.81	0.46
1:C:141:MSE:HB3	1:C:152:THR:HB	1.98	0.46
1:C:362:LYS:O	1:C:366:GLU:HG2	2.16	0.46
1:A:263:LEU:HB3	1:A:278:ILE:HD13	1.97	0.46
1:D:245:TRP:HA	1:D:250:GLN:HG2	1.97	0.46
1:A:87:MSE:HE1	1:A:227:PRO:HD2	1.97	0.46
1:A:93:ASP:HB3	1:A:98:ILE:HD11	1.98	0.46
1:B:84:ALA:CB	1:B:113:PRO:HB3	2.45	0.46
1:B:324:VAL:HG12	2:B:1384:FAD:O2	2.16	0.46
1:C:318:PRO:HB3	2:C:1384:FAD:C5X	2.46	0.46
1:C:87:MSE:H	1:C:112:ASN:HD21	1.63	0.45
1:D:112:ASN:H	1:D:113:PRO:HD2	1.78	0.45
2:B:1384:FAD:O2'	2:B:1384:FAD:O4'	2.30	0.45
1:C:245:TRP:HA	1:C:250:GLN:HG2	1.97	0.45
1:C:357:MSE:HA	1:C:360:TYR:CZ	2.52	0.45
1:C:317:PRO:HA	1:C:318:PRO:HD3	1.77	0.45
1:B:54[A]:ARG:NH2	1:C:54[A]:ARG:NE	2.64	0.45
1:B:174:VAL:O	1:B:174:VAL:CG1	2.65	0.45
1:B:62:LEU:HD13	1:B:71:MSE:HG3	1.99	0.45
1:C:93:ASP:HB3	1:C:98:ILE:HD11	1.99	0.45
1:A:363:GLU:HG3	1:B:93:ASP:HB3	1.99	0.45
1:B:87:MSE:H	1:B:112:ASN:HD21	1.65	0.44
1:C:208:LEU:HD23	1:C:208:LEU:HA	1.83	0.44
1:C:307:THR:HB	1:C:350:VAL:CG1	2.47	0.44
1:A:141:MSE:HB3	1:A:152:THR:HB	1.98	0.44
1:A:271:ASP:CG	1:A:273[A]:ARG:HH11	2.20	0.44
1:D:307:THR:HB	1:D:350:VAL:CG1	2.47	0.44
1:D:87:MSE:H	1:D:112:ASN:HD21	1.64	0.44
1:B:170:GLY:O	1:B:175:ARG:NH2	2.50	0.44
1:A:316:MSE:HE2	1:A:365:GLN:HG2	2.00	0.44
1:B:263:LEU:HD23	1:B:263:LEU:HA	1.88	0.44
1:D:291:ILE:O	1:D:293:PRO:HD3	2.17	0.44
1:A:271:ASP:OD1	1:A:273[A]:ARG:HD3	2.17	0.43
1:B:292:PHE:H	1:B:315:LEU:HB3	1.83	0.43
1:D:327:GLY:HA2	1:D:330:ASP:OD2	2.18	0.43
1:B:112:ASN:H	1:B:113:PRO:HD3	1.76	0.43
1:B:375:MSE:HA	1:B:380:PHE:CD1	2.54	0.43
1:B:71:MSE:HE2	1:B:80:TYR:HB2	2.01	0.43
1:A:31:MSE:HE3	1:A:35:LEU:HG	2.00	0.43
1:A:291:ILE:O	1:A:293:PRO:HD3	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:55:ILE:HG21	1:A:55:ILE:HD13	1.81	0.43
1:B:141:MSE:HB3	1:B:152:THR:HB	2.00	0.43
1:C:245:TRP:CD1	1:C:245:TRP:N	2.85	0.43
1:A:62:LEU:HD13	1:A:71:MSE:HG3	2.01	0.43
1:A:93:ASP:OD1	1:A:96:GLY:N	2.52	0.43
1:B:93:ASP:C	1:B:95:LYS:H	2.22	0.43
1:C:93:ASP:OD1	1:C:96:GLY:N	2.50	0.43
1:B:84:ALA:HB3	1:B:113:PRO:HB3	1.99	0.43
1:D:71:MSE:HE2	1:D:80:TYR:HB2	2.01	0.43
1:C:197:GLN:HB3	1:C:200:ILE:HD13	2.01	0.43
1:B:327:GLY:HA2	1:B:330:ASP:OD2	2.18	0.42
1:C:327:GLY:HA2	1:C:330:ASP:OD2	2.18	0.42
1:A:87:MSE:CE	1:A:227:PRO:HD2	2.49	0.42
1:B:31:MSE:HE3	1:B:35:LEU:HG	2.00	0.42
1:C:307:THR:HB	1:C:350:VAL:HG13	2.01	0.42
1:A:63:HIS:HA	1:A:111:ASP:CB	2.50	0.42
1:B:69:GLU:HG3	1:B:332:LEU:HD21	2.01	0.42
2:D:1384:FAD:H9	2:D:1384:FAD:H1'1	1.86	0.42
1:D:31:MSE:HE3	1:D:35:LEU:HG	2.01	0.42
1:A:310:GLY:HA2	1:A:327:GLY:HA3	2.02	0.42
1:C:221:ASN:HD22	1:C:239:PHE:HB3	1.84	0.42
1:D:274:TYR:O	1:D:277:LEU:HB3	2.18	0.42
1:A:226:ASN:C	1:A:226:ASN:OD1	2.58	0.42
1:C:142:LEU:HA	1:C:142:LEU:HD12	1.91	0.42
1:B:245:TRP:HA	1:B:250:GLN:HG2	2.02	0.42
1:D:141:MSE:HB3	1:D:152:THR:HB	2.01	0.42
1:A:245:TRP:O	1:A:246:LYS:CB	2.67	0.42
1:D:302:ARG:HA	1:D:303:PRO:HD3	1.92	0.42
1:A:198:PRO:HB3	1:A:233:LEU:HB2	2.02	0.42
1:B:84:ALA:CB	1:B:113:PRO:CB	2.98	0.42
1:B:166:LEU:HD12	1:B:308:MSE:HG2	2.01	0.42
1:B:244:GLU:CD	1:B:244:GLU:H	2.22	0.42
1:A:359:ILE:HD12	1:B:383:GLN:HA	2.02	0.41
1:A:359:ILE:HG23	1:B:98:ILE:HG21	1.97	0.41
1:C:197:GLN:HB3	1:C:200:ILE:CD1	2.50	0.41
1:C:351:LYS:HD3	1:C:351:LYS:HA	1.94	0.41
1:D:93:ASP:C	1:D:95:LYS:H	2.24	0.41
1:C:318:PRO:HB3	2:C:1384:FAD:C9A	2.50	0.41
1:C:93:ASP:C	1:C:95:LYS:H	2.23	0.41
1:B:33:LYS:HD3	1:B:127:SER:OG	2.20	0.41
1:C:170:GLY:O	1:C:175:ARG:NH2	2.53	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:245:TRP:CD1	1:D:245:TRP:N	2.88	0.41
1:B:245:TRP:N	1:B:245:TRP:CD1	2.87	0.41
1:D:217:SER:HB3	1:D:222:LEU:HG	2.03	0.41
2:B:1384:FAD:H9	2:B:1384:FAD:H1'1	1.91	0.41
1:B:317:PRO:HA	1:B:318:PRO:HD3	1.84	0.41
1:D:70:ALA:HA	1:D:332:LEU:HD13	2.02	0.41
1:A:324:VAL:HG12	2:A:1384:FAD:O2	2.20	0.41
1:B:76:LEU:CA	4:B:1387:GOL:H11	2.51	0.41
1:C:31:MSE:CG	1:C:309:ILE:CD1	2.95	0.41
1:C:33:LYS:HD3	1:C:127:SER:OG	2.21	0.41
1:D:197:GLN:HB3	1:D:200:ILE:CD1	2.51	0.41
1:B:44:VAL:HB	1:B:133:VAL:HG22	2.02	0.41
1:C:22:ILE:HD11	1:C:151:LEU:CD2	2.50	0.41
1:C:31:MSE:CG	1:C:309:ILE:HD13	2.51	0.41
1:D:178:VAL:HG22	1:D:306:ILE:HG23	2.03	0.41
1:B:62:LEU:O	1:B:111:ASP:HA	2.21	0.41
1:D:375:MSE:HA	1:D:380:PHE:CG	2.56	0.41
1:A:175:ARG:O	1:A:176:LYS:C	2.60	0.41
1:A:375:MSE:HA	1:A:380:PHE:CG	2.56	0.40
1:D:317:PRO:HA	1:D:318:PRO:HD3	1.82	0.40
1:A:33:LYS:HD3	1:A:127:SER:OG	2.21	0.40
1:B:63:HIS:HA	1:B:111:ASP:CB	2.51	0.40
1:C:69:GLU:HG3	1:C:332:LEU:HD21	2.04	0.40
1:A:166:LEU:HD12	1:A:308:MSE:HG2	2.03	0.40
1:B:166:LEU:HD12	1:B:308:MSE:HE3	2.02	0.40
1:B:55:ILE:HG21	1:B:55:ILE:HD13	1.88	0.40
1:C:175:ARG:O	1:C:176:LYS:C	2.60	0.40
1:D:197:GLN:HB3	1:D:200:ILE:HD13	2.04	0.40
1:D:307:THR:HB	1:D:350:VAL:HG13	2.03	0.40
1:A:210:ASN:C	1:A:212:ASN:H	2.24	0.40
1:B:108:ASN:C	1:B:110:PHE:H	2.24	0.40
1:C:62:LEU:HD13	1:C:71:MSE:HG3	2.03	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	366/399 (92%)	332 (91%)	31 (8%)	3 (1%)	22	64
1	B	366/399 (92%)	336 (92%)	27 (7%)	3 (1%)	22	64
1	C	364/399 (91%)	335 (92%)	26 (7%)	3 (1%)	22	64
1	D	364/399 (91%)	337 (93%)	24 (7%)	3 (1%)	22	64
All	All	1460/1596 (92%)	1340 (92%)	108 (7%)	12 (1%)	22	64

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	112	ASN
1	B	112	ASN
1	C	112	ASN
1	D	112	ASN
1	B	378	PRO
1	B	113	PRO
1	D	113	PRO
1	A	113	PRO
1	C	113	PRO
1	C	378	PRO
1	D	378	PRO
1	A	378	PRO

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	306/334 (92%)	291 (95%)	15 (5%)	29	68
1	B	307/334 (92%)	291 (95%)	16 (5%)	27	65
1	C	305/334 (91%)	293 (96%)	12 (4%)	37	75
1	D	303/334 (91%)	290 (96%)	13 (4%)	33	72

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	1221/1336 (91%)	1165 (95%)	56 (5%)	32 70

All (56) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	ASP
1	A	97	ASN
1	A	99	LEU
1	A	110	PHE
1	A	116	ASN
1	A	130	ASN
1	A	134	ILE
1	A	174	VAL
1	A	181	THR
1	A	199	GLU
1	A	200	ILE
1	A	273[A]	ARG
1	A	273[B]	ARG
1	A	353	TYR
1	A	355	GLN
1	B	16	ASP
1	B	97	ASN
1	B	99	LEU
1	B	110	PHE
1	B	116	ASN
1	B	130	ASN
1	B	134	ILE
1	B	174	VAL
1	B	181	THR
1	B	183	VAL
1	B	199	GLU
1	B	200	ILE
1	B	315	LEU
1	B	318	PRO
1	B	353	TYR
1	B	355	GLN
1	C	16	ASP
1	C	97	ASN
1	C	99	LEU
1	C	116	ASN
1	C	130	ASN
1	C	134	ILE

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Mol	Chain	Res	Type
1	C	174	VAL
1	C	181	THR
1	C	183	VAL
1	C	199	GLU
1	C	200	ILE
1	C	353	TYR
1	D	16	ASP
1	D	79	THR
1	D	97	ASN
1	D	99	LEU
1	D	116	ASN
1	D	130	ASN
1	D	134	ILE
1	D	174	VAL
1	D	181	THR
1	D	183	VAL
1	D	199	GLU
1	D	200	ILE
1	D	353	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	68	GLN
1	A	112	ASN
1	A	118	ASN
1	A	190	ASN
1	A	201	ASN
1	A	221	ASN
1	A	250	GLN
1	B	68	GLN
1	B	112	ASN
1	B	118	ASN
1	B	190	ASN
1	B	201	ASN
1	B	221	ASN
1	B	228	ASN
1	C	68	GLN
1	C	112	ASN
1	C	118	ASN
1	C	190	ASN
1	C	201	ASN

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Mol	Chain	Res	Type
1	C	221	ASN
1	D	68	GLN
1	D	112	ASN
1	D	118	ASN
1	D	190	ASN
1	D	201	ASN
1	D	221	ASN

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](#)

13 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	1384	-	51,58,58	1.46	6 (11%)	54,89,89	2.77	11 (20%)
3	SO4	A	1385	-	4,4,4	0.17	0	6,6,6	0.29	0
3	SO4	A	1386	-	4,4,4	0.35	0	6,6,6	0.54	0
2	FAD	B	1384	-	51,58,58	1.33	5 (9%)	54,89,89	2.52	11 (20%)
3	SO4	B	1385	-	4,4,4	0.20	0	6,6,6	0.33	0
3	SO4	B	1386	-	4,4,4	0.24	0	6,6,6	0.25	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	GOL	B	1387	-	5,5,5	0.45	0	5,5,5	0.90	0
2	FAD	C	1384	-	51,58,58	1.46	8 (15%)	54,89,89	2.31	7 (12%)
3	SO4	C	1385	-	4,4,4	0.20	0	6,6,6	0.30	0
3	SO4	C	1386	-	4,4,4	0.27	0	6,6,6	0.44	0
2	FAD	D	1384	-	51,58,58	1.44	7 (13%)	54,89,89	2.06	6 (11%)
3	SO4	D	1385	-	4,4,4	0.18	0	6,6,6	0.13	0
3	SO4	D	1386	-	4,4,4	0.18	0	6,6,6	0.12	0

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	1384	-	-	0/28/50/50	0/6/6/6
3	SO4	A	1385	-	-	0/0/0/0	0/0/0/0
3	SO4	A	1386	-	-	0/0/0/0	0/0/0/0
2	FAD	B	1384	-	-	0/28/50/50	0/6/6/6
3	SO4	B	1385	-	-	0/0/0/0	0/0/0/0
3	SO4	B	1386	-	-	0/0/0/0	0/0/0/0
4	GOL	B	1387	-	-	0/4/4/4	0/0/0/0
2	FAD	C	1384	-	-	0/28/50/50	0/6/6/6
3	SO4	C	1385	-	-	0/0/0/0	0/0/0/0
3	SO4	C	1386	-	-	0/0/0/0	0/0/0/0
2	FAD	D	1384	-	-	0/28/50/50	0/6/6/6
3	SO4	D	1385	-	-	0/0/0/0	0/0/0/0
3	SO4	D	1386	-	-	0/0/0/0	0/0/0/0

All (26) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1384	FAD	C4-C4X	-3.72	1.34	1.41
2	C	1384	FAD	C2B-C1B	-2.47	1.49	1.53
2	C	1384	FAD	C4-C4X	-2.46	1.36	1.41
2	C	1384	FAD	C5X-N5	2.12	1.38	1.35
2	D	1384	FAD	C5X-N5	2.21	1.38	1.35
2	C	1384	FAD	C2A-N1A	2.47	1.38	1.33
2	D	1384	FAD	C1'-N10	2.52	1.51	1.48
2	B	1384	FAD	C4-N3	2.62	1.37	1.33
2	D	1384	FAD	C2A-N1A	2.63	1.38	1.33
2	A	1384	FAD	C5X-N5	2.65	1.39	1.35
2	B	1384	FAD	C10-N1	2.66	1.37	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1384	FAD	C1'-N10	2.92	1.51	1.48
2	A	1384	FAD	C2A-N3A	2.96	1.37	1.32
2	A	1384	FAD	C4-N3	3.00	1.38	1.33
2	C	1384	FAD	C4X-N5	3.26	1.38	1.33
2	C	1384	FAD	C2A-N3A	3.27	1.37	1.32
2	A	1384	FAD	C10-N1	3.44	1.38	1.33
2	B	1384	FAD	C2A-N3A	3.53	1.38	1.32
2	D	1384	FAD	C2A-N3A	3.56	1.38	1.32
2	D	1384	FAD	C4-N3	3.80	1.39	1.33
2	C	1384	FAD	C4-N3	3.85	1.40	1.33
2	D	1384	FAD	C10-N1	3.97	1.38	1.33
2	A	1384	FAD	C4X-N5	4.34	1.39	1.33
2	D	1384	FAD	C4X-N5	4.42	1.39	1.33
2	B	1384	FAD	C4X-N5	4.47	1.39	1.33
2	C	1384	FAD	C10-N1	4.50	1.39	1.33

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1384	FAD	N3A-C2A-N1A	-12.45	118.01	128.86
2	B	1384	FAD	N3A-C2A-N1A	-11.77	118.61	128.86
2	A	1384	FAD	N3A-C2A-N1A	-11.52	118.82	128.86
2	D	1384	FAD	N3A-C2A-N1A	-10.09	120.07	128.86
2	A	1384	FAD	C1'-N10-C10	-6.76	111.57	118.50
2	A	1384	FAD	C4-C4X-C10	-4.74	116.12	119.96
2	B	1384	FAD	C1B-N9A-C4A	-3.77	120.13	126.64
2	B	1384	FAD	C4X-C4-N3	-2.92	119.33	123.48
2	C	1384	FAD	C1'-N10-C10	-2.74	115.69	118.50
2	D	1384	FAD	C4X-C4-N3	-2.60	119.78	123.48
2	A	1384	FAD	C1B-N9A-C4A	-2.55	122.22	126.64
2	A	1384	FAD	C4A-C5A-N7A	-2.45	107.04	109.41
2	C	1384	FAD	C1B-N9A-C4A	-2.30	122.67	126.64
2	C	1384	FAD	O3B-C3B-C4B	-2.29	104.40	111.09
2	A	1384	FAD	C4B-O4B-C1B	-2.20	107.42	109.77
2	B	1384	FAD	C4-C4X-C10	-2.10	118.26	119.96
2	A	1384	FAD	C4X-C10-N10	-2.02	119.11	120.52
2	B	1384	FAD	C4-C4X-N5	2.12	121.00	118.68
2	A	1384	FAD	C5X-C9A-N10	2.13	119.24	117.66
2	D	1384	FAD	C4X-N5-C5X	2.39	119.29	116.76
2	B	1384	FAD	C4X-N5-C5X	2.40	119.29	116.76
2	B	1384	FAD	O4'-C4'-C5'	2.42	115.39	110.00
2	A	1384	FAD	C10-C4X-N5	2.56	123.54	120.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1384	FAD	C5X-C9A-N10	2.68	119.65	117.66
2	B	1384	FAD	O3'-C3'-C2'	2.89	115.97	108.82
2	D	1384	FAD	C5X-C9A-N10	3.96	120.60	117.66
2	C	1384	FAD	C4-N3-C2	4.06	118.71	115.16
2	B	1384	FAD	C5X-C9A-N10	4.31	120.86	117.66
2	D	1384	FAD	C4-N3-C2	4.94	119.48	115.16
2	D	1384	FAD	C1'-N10-C9A	5.04	122.96	118.35
2	B	1384	FAD	C1'-N10-C9A	6.60	124.40	118.35
2	A	1384	FAD	C4-N3-C2	6.74	121.06	115.16
2	C	1384	FAD	C1'-N10-C9A	6.88	124.65	118.35
2	B	1384	FAD	C4-N3-C2	7.44	121.67	115.16
2	A	1384	FAD	C1'-N10-C9A	10.17	127.67	118.35

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 22 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1384	FAD	3	0
2	B	1384	FAD	5	0
4	B	1387	GOL	4	0
2	C	1384	FAD	7	0
2	D	1384	FAD	3	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	357/399 (89%)	-0.27	2 (0%) 89 71	9, 30, 65, 128	0
1	B	357/399 (89%)	-0.27	4 (1%) 80 55	10, 29, 62, 133	0
1	C	356/399 (89%)	-0.17	4 (1%) 80 55	17, 43, 78, 108	0
1	D	356/399 (89%)	-0.14	5 (1%) 75 49	17, 45, 82, 119	0
All	All	1426/1596 (89%)	-0.21	15 (1%) 80 55	9, 37, 74, 133	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	378	PRO	4.6
1	B	379	ASP	4.5
1	B	111	ASP	3.4
1	D	249	THR	3.2
1	B	249	THR	2.8
1	C	340	ASP	2.7
1	A	249	THR	2.7
1	C	252	ASP	2.5
1	D	295	GLU	2.4
1	C	249	THR	2.3
1	A	383	GLN	2.2
1	C	327	GLY	2.2
1	D	342	LYS	2.1
1	D	180	ASP	2.1
1	D	250	GLN	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(\AA^2)	Q<0.9
3	SO4	D	1386	5/5	0.93	0.31	2.54	88,88,88,88	0
3	SO4	B	1386	5/5	0.93	0.25	2.22	75,75,75,75	0
3	SO4	C	1386	5/5	0.88	0.25	1.86	71,71,72,72	0
3	SO4	A	1386	5/5	0.91	0.19	0.46	58,59,60,60	0
2	FAD	B	1384	53/53	0.96	0.17	-0.25	13,23,31,34	0
2	FAD	A	1384	53/53	0.95	0.17	-0.34	12,19,31,32	0
4	GOL	B	1387	6/6	0.85	0.16	-0.57	23,30,32,32	0
2	FAD	C	1384	53/53	0.95	0.16	-0.73	22,27,31,32	0
2	FAD	D	1384	53/53	0.97	0.14	-0.92	20,25,33,37	0
3	SO4	A	1385	5/5	0.90	0.28	-	74,74,75,75	0
3	SO4	B	1385	5/5	0.91	0.29	-	61,62,63,63	0
3	SO4	C	1385	5/5	0.97	0.24	-	52,52,53,53	0
3	SO4	D	1385	5/5	0.95	0.23	-	67,68,68,69	0

6.5 Other polymers [i](#)

There are no such residues in this entry.