



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 15, 2017 – 04:05 am GMT

PDB ID : 5MGY
Title : Crystal structure of Pseudomonas stutzeri flavinyl transferase ApbE, apo form
Authors : Zhang, L.; Trncik, C.; Andrade, S.L.A.; Einsle, O.
Deposited on : 2016-11-22
Resolution : 2.60 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
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

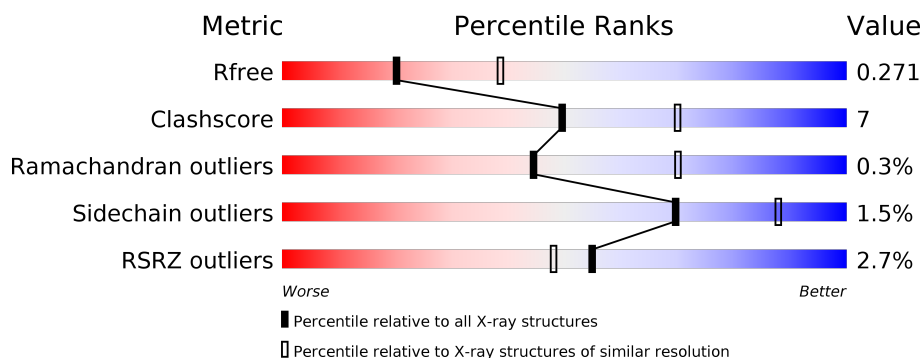
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	2542 (2.60-2.60)
Clashscore	112137	2895 (2.60-2.60)
Ramachandran outliers	110173	2848 (2.60-2.60)
Sidechain outliers	110143	2848 (2.60-2.60)
RSRZ outliers	101464	2550 (2.60-2.60)

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	332	<div> <div>3%</div> <div> <div></div> <div>80%</div> <div>11%</div> <div>8%</div> </div> </div>
1	B	332	<div> <div>2%</div> <div> <div></div> <div>79%</div> <div>11%</div> <div>9%</div> </div> </div>
1	C	332	<div> <div>2%</div> <div> <div></div> <div>83%</div> <div>10%</div> <div>7%</div> </div> </div>
1	D	332	<div> <div>0%</div> <div> <div></div> <div>80%</div> <div>11%</div> <div>7%</div> </div> </div>
1	E	332	<div> <div>2%</div> <div> <div></div> <div>83%</div> <div>11%</div> <div>5%</div> </div> </div>
1	F	332	<div> <div>2%</div> <div> <div></div> <div>79%</div> <div>15%</div> <div>5%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
1	G	332	<div><div></div><div>5%</div><div>72%</div><div>16%</div><div>•</div><div>10%</div></div>
1	H	332	<div><div></div><div>3%</div><div>83%</div><div>12%</div><div>•</div><div>5%</div></div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 18725 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 FAD:protein FMN transferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	H	316	Total	C	N	O	S	0	0	0
			2388	1495	415	470	8			
1	B	303	Total	C	N	O	S	0	0	0
			2288	1431	392	457	8			
1	G	299	Total	C	N	O	S	0	0	0
			2261	1422	385	446	8			
1	A	305	Total	C	N	O	S	0	0	0
			2306	1444	395	459	8			
1	C	310	Total	C	N	O	S	0	0	0
			2334	1466	400	460	8			
1	D	309	Total	C	N	O	S	0	0	0
			2330	1464	399	459	8			
1	E	316	Total	C	N	O	S	0	0	0
			2388	1495	415	470	8			
1	F	315	Total	C	N	O	S	0	0	0
			2377	1489	411	469	8			

There are 200 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
H	27	MET	-	initiating methionine	UNP H7EX88
H	28	ALA	-	expression tag	UNP H7EX88
H	29	MET	-	expression tag	UNP H7EX88
H	30	ASP	-	expression tag	UNP H7EX88
H	31	LEU	-	expression tag	UNP H7EX88
H	32	PHE	-	expression tag	UNP H7EX88
H	33	GLN	-	expression tag	UNP H7EX88
H	34	ASP	-	expression tag	UNP H7EX88
H	35	LYS	-	expression tag	UNP H7EX88
H	36	VAL	-	expression tag	UNP H7EX88
H	37	GLU	-	expression tag	UNP H7EX88
H	38	ALA	-	expression tag	UNP H7EX88
H	39	PHE	-	expression tag	UNP H7EX88

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Chain	Residue	Modelled	Actual	Comment	Reference
H	40	THR	-	expression tag	UNP H7EX88
H	41	GLY	-	expression tag	UNP H7EX88
H	42	PRO	-	expression tag	UNP H7EX88
H	43	THR	-	expression tag	UNP H7EX88
H	351	VAL	-	expression tag	UNP H7EX88
H	352	GLU	-	expression tag	UNP H7EX88
H	353	HIS	-	expression tag	UNP H7EX88
H	354	HIS	-	expression tag	UNP H7EX88
H	355	HIS	-	expression tag	UNP H7EX88
H	356	HIS	-	expression tag	UNP H7EX88
H	357	HIS	-	expression tag	UNP H7EX88
H	358	HIS	-	expression tag	UNP H7EX88
B	27	MET	-	initiating methionine	UNP H7EX88
B	28	ALA	-	expression tag	UNP H7EX88
B	29	MET	-	expression tag	UNP H7EX88
B	30	ASP	-	expression tag	UNP H7EX88
B	31	LEU	-	expression tag	UNP H7EX88
B	32	PHE	-	expression tag	UNP H7EX88
B	33	GLN	-	expression tag	UNP H7EX88
B	34	ASP	-	expression tag	UNP H7EX88
B	35	LYS	-	expression tag	UNP H7EX88
B	36	VAL	-	expression tag	UNP H7EX88
B	37	GLU	-	expression tag	UNP H7EX88
B	38	ALA	-	expression tag	UNP H7EX88
B	39	PHE	-	expression tag	UNP H7EX88
B	40	THR	-	expression tag	UNP H7EX88
B	41	GLY	-	expression tag	UNP H7EX88
B	42	PRO	-	expression tag	UNP H7EX88
B	43	THR	-	expression tag	UNP H7EX88
B	351	VAL	-	expression tag	UNP H7EX88
B	352	GLU	-	expression tag	UNP H7EX88
B	353	HIS	-	expression tag	UNP H7EX88
B	354	HIS	-	expression tag	UNP H7EX88
B	355	HIS	-	expression tag	UNP H7EX88
B	356	HIS	-	expression tag	UNP H7EX88
B	357	HIS	-	expression tag	UNP H7EX88
B	358	HIS	-	expression tag	UNP H7EX88
G	27	MET	-	initiating methionine	UNP H7EX88
G	28	ALA	-	expression tag	UNP H7EX88
G	29	MET	-	expression tag	UNP H7EX88
G	30	ASP	-	expression tag	UNP H7EX88
G	31	LEU	-	expression tag	UNP H7EX88

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Chain	Residue	Modelled	Actual	Comment	Reference
G	32	PHE	-	expression tag	UNP H7EX88
G	33	GLN	-	expression tag	UNP H7EX88
G	34	ASP	-	expression tag	UNP H7EX88
G	35	LYS	-	expression tag	UNP H7EX88
G	36	VAL	-	expression tag	UNP H7EX88
G	37	GLU	-	expression tag	UNP H7EX88
G	38	ALA	-	expression tag	UNP H7EX88
G	39	PHE	-	expression tag	UNP H7EX88
G	40	THR	-	expression tag	UNP H7EX88
G	41	GLY	-	expression tag	UNP H7EX88
G	42	PRO	-	expression tag	UNP H7EX88
G	43	THR	-	expression tag	UNP H7EX88
G	351	VAL	-	expression tag	UNP H7EX88
G	352	GLU	-	expression tag	UNP H7EX88
G	353	HIS	-	expression tag	UNP H7EX88
G	354	HIS	-	expression tag	UNP H7EX88
G	355	HIS	-	expression tag	UNP H7EX88
G	356	HIS	-	expression tag	UNP H7EX88
G	357	HIS	-	expression tag	UNP H7EX88
G	358	HIS	-	expression tag	UNP H7EX88
A	27	MET	-	initiating methionine	UNP H7EX88
A	28	ALA	-	expression tag	UNP H7EX88
A	29	MET	-	expression tag	UNP H7EX88
A	30	ASP	-	expression tag	UNP H7EX88
A	31	LEU	-	expression tag	UNP H7EX88
A	32	PHE	-	expression tag	UNP H7EX88
A	33	GLN	-	expression tag	UNP H7EX88
A	34	ASP	-	expression tag	UNP H7EX88
A	35	LYS	-	expression tag	UNP H7EX88
A	36	VAL	-	expression tag	UNP H7EX88
A	37	GLU	-	expression tag	UNP H7EX88
A	38	ALA	-	expression tag	UNP H7EX88
A	39	PHE	-	expression tag	UNP H7EX88
A	40	THR	-	expression tag	UNP H7EX88
A	41	GLY	-	expression tag	UNP H7EX88
A	42	PRO	-	expression tag	UNP H7EX88
A	43	THR	-	expression tag	UNP H7EX88
A	351	VAL	-	expression tag	UNP H7EX88
A	352	GLU	-	expression tag	UNP H7EX88
A	353	HIS	-	expression tag	UNP H7EX88
A	354	HIS	-	expression tag	UNP H7EX88
A	355	HIS	-	expression tag	UNP H7EX88

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Chain	Residue	Modelled	Actual	Comment	Reference
A	356	HIS	-	expression tag	UNP H7EX88
A	357	HIS	-	expression tag	UNP H7EX88
A	358	HIS	-	expression tag	UNP H7EX88
C	27	MET	-	initiating methionine	UNP H7EX88
C	28	ALA	-	expression tag	UNP H7EX88
C	29	MET	-	expression tag	UNP H7EX88
C	30	ASP	-	expression tag	UNP H7EX88
C	31	LEU	-	expression tag	UNP H7EX88
C	32	PHE	-	expression tag	UNP H7EX88
C	33	GLN	-	expression tag	UNP H7EX88
C	34	ASP	-	expression tag	UNP H7EX88
C	35	LYS	-	expression tag	UNP H7EX88
C	36	VAL	-	expression tag	UNP H7EX88
C	37	GLU	-	expression tag	UNP H7EX88
C	38	ALA	-	expression tag	UNP H7EX88
C	39	PHE	-	expression tag	UNP H7EX88
C	40	THR	-	expression tag	UNP H7EX88
C	41	GLY	-	expression tag	UNP H7EX88
C	42	PRO	-	expression tag	UNP H7EX88
C	43	THR	-	expression tag	UNP H7EX88
C	351	VAL	-	expression tag	UNP H7EX88
C	352	GLU	-	expression tag	UNP H7EX88
C	353	HIS	-	expression tag	UNP H7EX88
C	354	HIS	-	expression tag	UNP H7EX88
C	355	HIS	-	expression tag	UNP H7EX88
C	356	HIS	-	expression tag	UNP H7EX88
C	357	HIS	-	expression tag	UNP H7EX88
C	358	HIS	-	expression tag	UNP H7EX88
D	27	MET	-	initiating methionine	UNP H7EX88
D	28	ALA	-	expression tag	UNP H7EX88
D	29	MET	-	expression tag	UNP H7EX88
D	30	ASP	-	expression tag	UNP H7EX88
D	31	LEU	-	expression tag	UNP H7EX88
D	32	PHE	-	expression tag	UNP H7EX88
D	33	GLN	-	expression tag	UNP H7EX88
D	34	ASP	-	expression tag	UNP H7EX88
D	35	LYS	-	expression tag	UNP H7EX88
D	36	VAL	-	expression tag	UNP H7EX88
D	37	GLU	-	expression tag	UNP H7EX88
D	38	ALA	-	expression tag	UNP H7EX88
D	39	PHE	-	expression tag	UNP H7EX88
D	40	THR	-	expression tag	UNP H7EX88

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Chain	Residue	Modelled	Actual	Comment	Reference
D	41	GLY	-	expression tag	UNP H7EX88
D	42	PRO	-	expression tag	UNP H7EX88
D	43	THR	-	expression tag	UNP H7EX88
D	351	VAL	-	expression tag	UNP H7EX88
D	352	GLU	-	expression tag	UNP H7EX88
D	353	HIS	-	expression tag	UNP H7EX88
D	354	HIS	-	expression tag	UNP H7EX88
D	355	HIS	-	expression tag	UNP H7EX88
D	356	HIS	-	expression tag	UNP H7EX88
D	357	HIS	-	expression tag	UNP H7EX88
D	358	HIS	-	expression tag	UNP H7EX88
E	27	MET	-	initiating methionine	UNP H7EX88
E	28	ALA	-	expression tag	UNP H7EX88
E	29	MET	-	expression tag	UNP H7EX88
E	30	ASP	-	expression tag	UNP H7EX88
E	31	LEU	-	expression tag	UNP H7EX88
E	32	PHE	-	expression tag	UNP H7EX88
E	33	GLN	-	expression tag	UNP H7EX88
E	34	ASP	-	expression tag	UNP H7EX88
E	35	LYS	-	expression tag	UNP H7EX88
E	36	VAL	-	expression tag	UNP H7EX88
E	37	GLU	-	expression tag	UNP H7EX88
E	38	ALA	-	expression tag	UNP H7EX88
E	39	PHE	-	expression tag	UNP H7EX88
E	40	THR	-	expression tag	UNP H7EX88
E	41	GLY	-	expression tag	UNP H7EX88
E	42	PRO	-	expression tag	UNP H7EX88
E	43	THR	-	expression tag	UNP H7EX88
E	351	VAL	-	expression tag	UNP H7EX88
E	352	GLU	-	expression tag	UNP H7EX88
E	353	HIS	-	expression tag	UNP H7EX88
E	354	HIS	-	expression tag	UNP H7EX88
E	355	HIS	-	expression tag	UNP H7EX88
E	356	HIS	-	expression tag	UNP H7EX88
E	357	HIS	-	expression tag	UNP H7EX88
E	358	HIS	-	expression tag	UNP H7EX88
F	27	MET	-	initiating methionine	UNP H7EX88
F	28	ALA	-	expression tag	UNP H7EX88
F	29	MET	-	expression tag	UNP H7EX88
F	30	ASP	-	expression tag	UNP H7EX88
F	31	LEU	-	expression tag	UNP H7EX88
F	32	PHE	-	expression tag	UNP H7EX88

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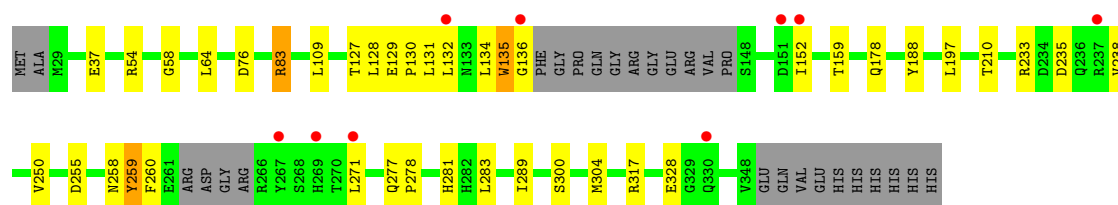
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F	36	VAL	-	expression tag	UNP H7EX88
F	37	GLU	-	expression tag	UNP H7EX88
F	38	ALA	-	expression tag	UNP H7EX88
F	39	PHE	-	expression tag	UNP H7EX88
F	40	THR	-	expression tag	UNP H7EX88
F	41	GLY	-	expression tag	UNP H7EX88
F	42	PRO	-	expression tag	UNP H7EX88
F	43	THR	-	expression tag	UNP H7EX88
F	351	VAL	-	expression tag	UNP H7EX88
F	352	GLU	-	expression tag	UNP H7EX88
F	353	HIS	-	expression tag	UNP H7EX88
F	354	HIS	-	expression tag	UNP H7EX88
F	355	HIS	-	expression tag	UNP H7EX88
F	356	HIS	-	expression tag	UNP H7EX88
F	357	HIS	-	expression tag	UNP H7EX88
F	358	HIS	-	expression tag	UNP H7EX88

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

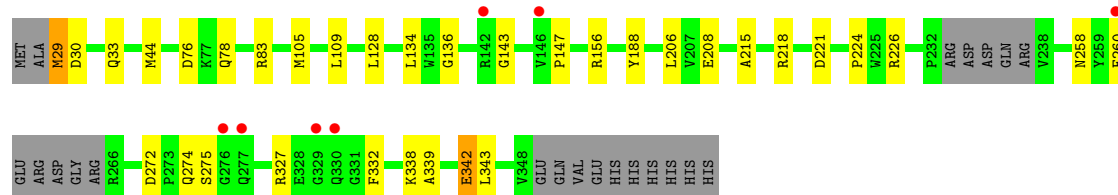
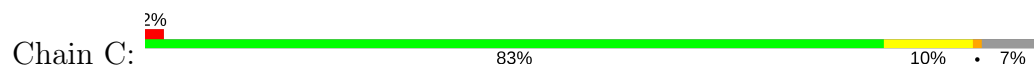
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	G	1	Total Mg 1 1	0	0
2	D	1	Total Mg 1 1	0	0
2	E	1	Total Mg 1 1	0	0
2	H	1	Total Mg 1 1	0	0
2	B	1	Total Mg 1 1	0	0
2	C	1	Total Mg 1 1	0	0
2	A	1	Total Mg 1 1	0	0
2	F	1	Total Mg 1 1	0	0

- Molecule 3 is water.

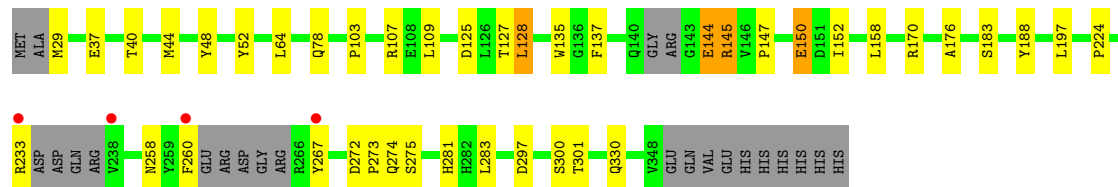
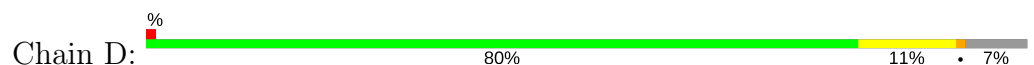
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	H	6	Total 6	O 6	0	0
3	B	9	Total 9	O 9	0	0
3	G	3	Total 3	O 3	0	0
3	A	12	Total 12	O 12	0	0
3	C	3	Total 3	O 3	0	0
3	D	5	Total 5	O 5	0	0
3	E	5	Total 5	O 5	0	0
3	F	2	Total 2	O 2	0	0



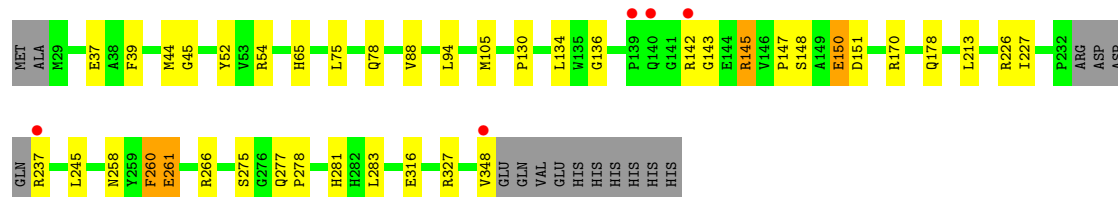
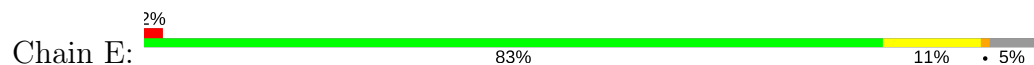
- Molecule 1: FAD:protein FMN transferase



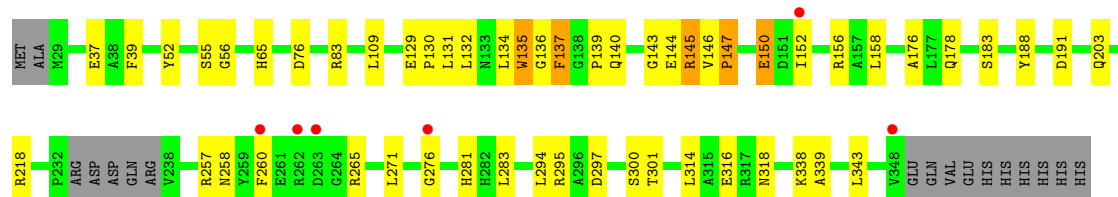
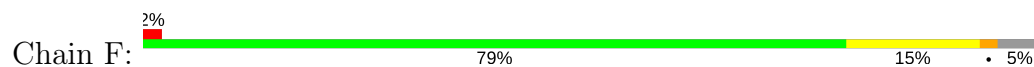
- Molecule 1: FAD:protein FMN transferase



- Molecule 1: FAD:protein FMN transferase



- Molecule 1: FAD:protein FMN transferase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	135.66Å 91.98Å 138.06Å 90.00° 90.41° 90.00°	Depositor
Resolution (Å)	138.06 – 2.60 138.06 – 2.46	Depositor EDS
% Data completeness (in resolution range)	100.0 (138.06-2.60) 100.0 (138.06-2.46)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.12 (at 2.45Å)	Xtriage
Refinement program	REFMAC 5.8.0155	Depositor
R, R_{free}	0.232 , 0.274 0.233 , 0.271	Depositor DCC
R_{free} test set	5134 reflections (5.15%)	DCC
Wilson B-factor (Å ²)	39.9	Xtriage
Anisotropy	0.214	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 21.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	0.000 for l,k,-h 0.000 for h,-k,-l 0.000 for l,-k,h	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	18725	wwPDB-VP
Average B, all atoms (Å ²)	60.0	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 48.34 % 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 8.7302e-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 ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: MG

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	1.04	3/2343 (0.1%)	1.05	6/3171 (0.2%)
1	B	0.95	1/2323 (0.0%)	1.06	12/3143 (0.4%)
1	C	0.88	1/2374 (0.0%)	0.97	5/3214 (0.2%)
1	D	0.91	3/2369 (0.1%)	1.02	8/3206 (0.2%)
1	E	0.87	0/2429	1.04	10/3287 (0.3%)
1	F	0.92	2/2418 (0.1%)	1.10	10/3273 (0.3%)
1	G	1.04	3/2299 (0.1%)	1.08	11/3110 (0.4%)
1	H	0.92	1/2429 (0.0%)	1.05	11/3287 (0.3%)
All	All	0.94	14/18984 (0.1%)	1.05	73/25691 (0.3%)

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	G	281	HIS	C-N	24.09	1.89	1.34
1	A	259	TYR	C-N	18.56	1.76	1.34
1	D	150	GLU	CG-CD	8.63	1.64	1.51
1	C	342	GLU	CD-OE2	-7.88	1.17	1.25
1	B	74	GLN	CG-CD	-7.76	1.33	1.51
1	H	267	TYR	CE1-CZ	7.24	1.48	1.38
1	A	136	GLY	N-CA	7.18	1.56	1.46
1	F	300	SER	CB-OG	-6.96	1.33	1.42
1	A	300	SER	CB-OG	-6.49	1.33	1.42
1	G	135	TRP	CB-CG	6.06	1.61	1.50
1	G	135	TRP	CZ3-CH2	5.48	1.48	1.40
1	D	300	SER	CB-OG	-5.10	1.35	1.42
1	F	150	GLU	CD-OE2	-5.04	1.20	1.25
1	D	137	PHE	CG-CD1	5.03	1.46	1.38

All (73) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	145	ARG	NE-CZ-NH2	-17.41	111.60	120.30
1	G	145	ARG	NE-CZ-NH1	-13.63	113.48	120.30
1	D	145	ARG	NE-CZ-NH2	11.78	126.19	120.30
1	F	146	VAL	C-N-CD	-10.91	96.59	120.60
1	H	76	ASP	CB-CG-OD2	10.39	127.65	118.30
1	G	327	ARG	NE-CZ-NH1	9.95	125.27	120.30
1	A	255	ASP	CB-CG-OD2	8.92	126.33	118.30
1	H	142	ARG	NE-CZ-NH2	8.65	124.63	120.30
1	E	226	ARG	NE-CZ-NH1	-8.11	116.25	120.30
1	H	233	ARG	NE-CZ-NH2	-8.03	116.28	120.30
1	C	218	ARG	NE-CZ-NH1	7.77	124.19	120.30
1	G	142	ARG	NE-CZ-NH2	7.64	124.12	120.30
1	B	295	ARG	NE-CZ-NH2	7.38	123.99	120.30
1	G	148	SER	N-CA-CB	7.29	121.43	110.50
1	G	145	ARG	NE-CZ-NH2	7.26	123.93	120.30
1	H	262	ARG	NE-CZ-NH1	7.03	123.81	120.30
1	B	104	ASP	CB-CG-OD1	7.00	124.61	118.30
1	D	233	ARG	NE-CZ-NH1	-6.98	116.81	120.30
1	H	233	ARG	NE-CZ-NH1	6.97	123.79	120.30
1	F	145	ARG	NH1-CZ-NH2	6.83	126.91	119.40
1	A	255	ASP	CB-CG-OD1	-6.59	112.37	118.30
1	H	208	GLU	OE1-CD-OE2	-6.58	115.41	123.30
1	E	226	ARG	NE-CZ-NH2	6.52	123.56	120.30
1	A	259	TYR	O-C-N	-6.50	112.30	122.70
1	G	327	ARG	NE-CZ-NH2	-6.50	117.05	120.30
1	B	295	ARG	NE-CZ-NH1	-6.45	117.07	120.30
1	F	295	ARG	NE-CZ-NH2	6.38	123.49	120.30
1	G	125	ASP	CB-CG-OD1	6.34	124.00	118.30
1	C	208	GLU	OE1-CD-OE2	-6.28	115.77	123.30
1	H	76	ASP	OD1-CG-OD2	-6.16	111.59	123.30
1	E	260	PHE	CB-CG-CD2	-6.05	116.57	120.80
1	G	77	LYS	CD-CE-NZ	5.99	125.48	111.70
1	A	233	ARG	NE-CZ-NH2	5.99	123.30	120.30
1	H	207	VAL	CG1-CB-CG2	-5.94	101.39	110.90
1	D	145	ARG	NE-CZ-NH1	-5.92	117.34	120.30
1	B	107	ARG	NE-CZ-NH1	5.84	123.22	120.30
1	E	142	ARG	NE-CZ-NH1	5.83	123.21	120.30
1	B	83	ARG	CB-CA-C	-5.82	98.75	110.40
1	F	191	ASP	CB-CG-OD1	5.81	123.53	118.30
1	C	44	MET	CG-SD-CE	5.76	109.42	100.20
1	C	226	ARG	NE-CZ-NH2	-5.76	117.42	120.30
1	E	170	ARG	NE-CZ-NH1	5.75	123.18	120.30
1	B	255	ASP	CB-CG-OD2	5.73	123.45	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	170	ARG	NE-CZ-NH1	5.63	123.11	120.30
1	F	265	ARG	NE-CZ-NH2	5.61	123.11	120.30
1	F	218	ARG	NE-CZ-NH1	5.57	123.09	120.30
1	E	44	MET	CA-CB-CG	5.54	122.72	113.30
1	F	297	ASP	CB-CG-OD2	5.51	123.26	118.30
1	B	34	ASP	CB-CG-OD1	5.46	123.22	118.30
1	H	265	ARG	NE-CZ-NH2	5.41	123.00	120.30
1	F	257	ARG	NE-CZ-NH2	5.38	122.99	120.30
1	D	107	ARG	NE-CZ-NH1	5.37	122.98	120.30
1	E	45	GLY	N-CA-C	5.35	126.48	113.10
1	A	83	ARG	NE-CZ-NH2	-5.30	117.65	120.30
1	G	125	ASP	CB-CG-OD2	-5.26	113.57	118.30
1	A	135	TRP	CA-C-N	5.23	126.65	116.20
1	F	137	PHE	CB-CG-CD2	5.21	124.45	120.80
1	B	75	LEU	CD1-CG-CD2	-5.20	94.91	110.50
1	D	44	MET	CG-SD-CE	5.19	108.50	100.20
1	B	290	ASP	CB-CG-OD1	-5.18	113.64	118.30
1	G	44	MET	CG-SD-CE	5.17	108.48	100.20
1	D	144	GLU	OE1-CD-OE2	-5.16	117.11	123.30
1	C	156	ARG	NE-CZ-NH1	5.15	122.88	120.30
1	E	54	ARG	NE-CZ-NH1	5.13	122.87	120.30
1	B	170	ARG	NE-CZ-NH2	5.13	122.87	120.30
1	H	290	ASP	CB-CG-OD1	-5.11	113.70	118.30
1	H	290	ASP	CB-CG-OD2	5.10	122.89	118.30
1	B	44	MET	CG-SD-CE	5.09	108.35	100.20
1	E	237	ARG	CG-CD-NE	-5.09	101.11	111.80
1	D	297	ASP	CB-CG-OD2	5.09	122.88	118.30
1	B	74	GLN	CB-CA-C	-5.00	100.39	110.40
1	G	54	ARG	NE-CZ-NH2	5.00	122.80	120.30
1	E	88	VAL	CG1-CB-CG2	-5.00	102.90	110.90

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2306	0	2261	31	1
1	B	2288	0	2250	26	1
1	C	2334	0	2295	19	0
1	D	2330	0	2292	21	0
1	E	2388	0	2349	30	0
1	F	2377	0	2336	53	0
1	G	2261	0	2220	61	0
1	H	2388	0	2349	34	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
2	G	1	0	0	0	0
2	H	1	0	0	0	0
3	A	12	0	0	1	0
3	B	9	0	0	0	0
3	C	3	0	0	0	0
3	D	5	0	0	0	0
3	E	5	0	0	0	0
3	F	2	0	0	0	0
3	G	3	0	0	0	0
3	H	6	0	0	0	0
All	All	18725	0	18352	248	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 (248) 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:F:134:LEU:HD11	1:F:147:PRO:CB	1.27	1.54
1:A:259:TYR:C	1:A:260:PHE:N	1.76	1.38
1:F:134:LEU:CD1	1:F:147:PRO:CB	2.02	1.37
1:G:281:HIS:C	1:G:282:HIS:N	1.89	1.24
1:F:134:LEU:CD1	1:F:147:PRO:CG	2.20	1.18
1:F:134:LEU:HD12	1:F:147:PRO:HG3	1.23	1.17
1:F:134:LEU:O	1:F:147:PRO:HG3	1.46	1.13
1:F:134:LEU:HD12	1:F:147:PRO:CG	1.82	1.09
1:H:142:ARG:HH22	1:H:145:ARG:HD2	1.14	1.08
1:H:142:ARG:NH2	1:H:145:ARG:CD	2.17	1.06

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:144:GLU:OE2	1:D:267:TYR:OH	1.71	1.05
1:H:142:ARG:NH2	1:H:145:ARG:HD3	1.75	1.01
1:H:142:ARG:HH22	1:H:145:ARG:CD	1.72	1.00
1:H:224:PRO:CD	1:E:260:PHE:CE2	2.46	0.98
1:G:272:ASP:OD2	1:G:274:GLN:HG2	1.63	0.98
1:F:134:LEU:CD1	1:F:147:PRO:HB3	1.91	0.98
1:G:258:ASN:HB2	1:G:260:PHE:CZ	2.00	0.96
1:G:259:TYR:OH	1:G:282:HIS:HA	1.64	0.96
1:F:134:LEU:HD11	1:F:147:PRO:HB3	1.47	0.96
1:H:224:PRO:HD3	1:E:260:PHE:CE2	2.04	0.91
1:F:134:LEU:CD1	1:F:147:PRO:HB2	1.84	0.91
1:F:134:LEU:HD11	1:F:147:PRO:HB2	0.88	0.86
1:G:282:HIS:HB3	1:G:327:ARG:HB3	1.61	0.83
1:F:134:LEU:CD1	1:F:147:PRO:HG3	1.96	0.83
1:G:327:ARG:HA	1:G:332:PHE:CE1	2.13	0.83
1:H:224:PRO:HD2	1:E:260:PHE:CE2	2.14	0.82
1:F:134:LEU:HD11	1:F:147:PRO:CG	1.97	0.82
1:H:142:ARG:NH2	1:H:145:ARG:HD2	1.88	0.81
1:H:142:ARG:HH21	1:H:145:ARG:HD3	1.46	0.79
1:B:260:PHE:CZ	1:C:224:PRO:HD3	2.18	0.79
1:G:281:HIS:CE1	1:G:307:GLY:CA	2.66	0.79
1:G:283:LEU:HD12	1:G:325:VAL:O	1.82	0.79
1:D:272:ASP:O	1:D:275:SER:O	2.01	0.79
1:A:260:PHE:CZ	1:D:224:PRO:HD3	2.19	0.78
1:G:135:TRP:NE1	1:G:276:GLY:O	2.17	0.77
1:B:233:ARG:HB2	1:B:237:ARG:HB3	1.67	0.76
1:G:226:ARG:HH12	1:F:260:PHE:HE2	1.32	0.75
1:G:326:VAL:O	1:G:332:PHE:HD1	1.69	0.75
1:G:256:TYR:CD2	1:G:257:ARG:N	2.56	0.74
1:H:272:ASP:O	1:H:275:SER:O	2.04	0.74
1:G:272:ASP:OD2	1:G:274:GLN:CG	2.36	0.73
1:A:132:LEU:HD23	1:A:271:LEU:HD11	1.70	0.73
1:A:132:LEU:HD23	1:A:271:LEU:CD1	2.19	0.73
1:F:134:LEU:HD21	1:F:152:ILE:HG13	1.69	0.72
1:G:258:ASN:CB	1:G:260:PHE:CZ	2.71	0.72
1:G:259:TYR:CD1	1:G:284:ALA:HB2	2.25	0.72
1:B:233:ARG:HB2	1:B:237:ARG:CB	2.22	0.70
1:G:268:SER:OG	1:G:279:ILE:HG12	1.92	0.69
1:H:64:LEU:HD13	1:H:197:LEU:HD22	1.76	0.68
1:C:272:ASP:O	1:C:275:SER:O	2.11	0.67
1:F:135:TRP:CZ2	1:F:152:ILE:HD11	2.28	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:259:TYR:HD1	1:G:284:ALA:HB2	1.59	0.66
1:F:134:LEU:O	1:F:134:LEU:HD12	1.95	0.66
1:G:327:ARG:CA	1:G:332:PHE:CE1	2.78	0.66
1:F:136:GLY:O	1:F:143:GLY:HA2	1.95	0.66
1:H:223:SER:HB2	1:E:260:PHE:HZ	1.61	0.65
1:F:135:TRP:CH2	1:F:152:ILE:HD11	2.31	0.65
1:E:145:ARG:O	1:E:147:PRO:HD3	1.98	0.64
1:F:134:LEU:O	1:F:147:PRO:CG	2.37	0.63
1:A:131:LEU:O	1:A:135:TRP:HD1	1.80	0.63
1:A:277:GLN:HB3	1:A:278:PRO:CD	2.28	0.63
1:H:224:PRO:HD3	1:E:260:PHE:CZ	2.33	0.62
1:G:326:VAL:O	1:G:332:PHE:CD1	2.53	0.61
1:G:281:HIS:CE1	1:G:307:GLY:HA3	2.34	0.61
1:C:258:ASN:HB3	1:C:260:PHE:CZ	2.36	0.61
1:H:224:PRO:CD	1:E:260:PHE:CZ	2.83	0.61
1:G:188:TYR:CE2	1:G:192:LEU:HD11	2.36	0.61
1:A:129:GLU:HA	1:A:132:LEU:HD12	1.83	0.60
1:B:260:PHE:CZ	1:C:224:PRO:CD	2.85	0.60
1:D:125:ASP:OD2	1:D:127:THR:OG1	2.14	0.60
1:H:223:SER:HA	1:E:260:PHE:CZ	2.37	0.59
1:H:224:PRO:HD3	1:E:260:PHE:CD2	2.37	0.59
1:F:131:LEU:HD11	1:F:156:ARG:CG	2.33	0.58
1:G:268:SER:OG	1:G:279:ILE:CG1	2.51	0.58
1:E:134:LEU:HG	1:E:147:PRO:HG3	1.84	0.58
1:B:258:ASN:HB3	1:B:260:PHE:CZ	2.40	0.57
1:H:223:SER:HA	1:E:260:PHE:CE1	2.39	0.57
1:H:223:SER:HB2	1:E:260:PHE:CZ	2.38	0.57
1:F:132:LEU:HA	1:F:271:LEU:HD11	1.85	0.56
1:B:271:LEU:HD23	1:B:278:PRO:N	2.20	0.56
1:G:281:HIS:CE1	1:G:307:GLY:N	2.74	0.56
1:G:327:ARG:HA	1:G:332:PHE:CD1	2.40	0.56
1:B:271:LEU:HD23	1:B:278:PRO:CA	2.36	0.56
1:G:274:GLN:HG3	1:G:275:SER:N	2.21	0.56
1:A:258:ASN:HB3	1:A:260:PHE:CZ	2.41	0.55
1:F:140:GLN:OE1	1:F:140:GLN:N	2.39	0.55
1:B:75:LEU:HD23	1:B:105:MET:SD	2.47	0.55
1:F:131:LEU:HD11	1:F:156:ARG:HG3	1.88	0.55
1:C:206:LEU:HD23	1:C:215:ALA:HB2	1.89	0.55
1:G:253:SER:HB2	1:G:300:SER:OG	2.06	0.55
1:E:78:GLN:OE1	1:E:105:MET:HB2	2.07	0.54
1:G:256:TYR:HD2	1:G:257:ARG:H	1.56	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:67:GLU:O	1:H:71:ILE:HG13	2.07	0.54
1:B:158:LEU:HD22	1:B:176:ALA:HB3	1.90	0.54
1:B:260:PHE:CE2	1:C:224:PRO:HD3	2.43	0.54
1:F:135:TRP:CD1	1:F:135:TRP:N	2.75	0.54
1:A:277:GLN:HB3	1:A:278:PRO:HD2	1.90	0.54
1:A:37:GLU:OE1	1:A:54:ARG:NH1	2.41	0.54
1:E:213:LEU:HD11	1:E:227:ILE:HG21	1.88	0.54
1:E:136:GLY:HA3	1:E:143:GLY:HA2	1.90	0.53
1:F:139:PRO:HG2	1:F:140:GLN:OE1	2.08	0.53
1:G:256:TYR:HA	1:G:284:ALA:HB1	1.90	0.53
1:H:136:GLY:HA3	1:H:143:GLY:HA2	1.90	0.52
1:H:142:ARG:HB2	1:F:203:GLN:OE1	2.09	0.52
1:A:134:LEU:O	1:A:134:LEU:HG	2.10	0.52
1:C:272:ASP:OD2	1:C:274:GLN:HB3	2.10	0.52
1:B:210:THR:HB	1:C:29:MET:HB2	1.92	0.52
1:G:258:ASN:HB2	1:G:260:PHE:HZ	1.66	0.52
1:D:272:ASP:OD2	1:D:274:GLN:HB3	2.10	0.52
1:C:76:ASP:OD1	1:C:83:ARG:HD3	2.10	0.51
1:B:260:PHE:CE2	1:C:224:PRO:CD	2.93	0.51
1:C:339:ALA:O	1:C:343:LEU:HD23	2.10	0.51
1:F:134:LEU:CG	1:F:147:PRO:CG	2.88	0.51
1:G:256:TYR:HE2	1:G:257:ARG:HE	1.59	0.51
1:G:259:TYR:HH	1:G:282:HIS:HA	1.75	0.51
1:F:134:LEU:HG	1:F:147:PRO:HG2	1.92	0.51
1:H:224:PRO:HD2	1:E:260:PHE:CZ	2.45	0.51
1:H:188:TYR:CE2	1:H:192:LEU:HD11	2.46	0.50
1:F:134:LEU:HD12	1:F:147:PRO:CB	2.07	0.50
1:G:223:SER:CB	1:F:260:PHE:CZ	2.94	0.50
1:G:259:TYR:OH	1:G:282:HIS:CA	2.48	0.50
1:A:135:TRP:HZ2	1:A:152:ILE:HD11	1.77	0.50
1:D:64:LEU:HD22	1:D:197:LEU:CD2	2.41	0.50
1:F:258:ASN:HB3	1:F:260:PHE:CZ	2.46	0.50
1:B:128:LEU:O	1:B:129:GLU:C	2.49	0.50
1:A:135:TRP:CZ2	1:A:152:ILE:HD11	2.47	0.50
1:G:183:SER:HB2	1:G:301:THR:HG21	1.93	0.49
1:G:281:HIS:ND1	1:G:307:GLY:HA3	2.26	0.49
1:D:78:GLN:OE1	1:D:103:PRO:HB2	2.13	0.49
1:G:327:ARG:NE	1:G:332:PHE:CZ	2.79	0.49
1:B:281:HIS:HD2	1:B:283:LEU:H	1.60	0.49
1:A:317:ARG:HD3	3:A:510:HOH:O	2.12	0.49
1:A:132:LEU:HD23	1:A:271:LEU:HD12	1.93	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:145:ARG:O	1:D:147:PRO:HD3	2.13	0.48
1:G:188:TYR:CZ	1:G:192:LEU:HD11	2.48	0.48
1:F:139:PRO:HB2	1:F:140:GLN:OE1	2.13	0.48
1:H:142:ARG:HA	1:F:203:GLN:HE22	1.78	0.48
1:G:259:TYR:CE1	1:G:282:HIS:O	2.66	0.48
1:A:260:PHE:CE2	1:D:224:PRO:HD3	2.48	0.48
1:D:128:LEU:HD11	1:D:273:PRO:HB3	1.94	0.48
1:G:270:THR:HB	1:G:279:ILE:HD13	1.96	0.48
1:B:272:ASP:OD2	1:B:274:GLN:HB3	2.14	0.48
1:E:75:LEU:HD23	1:E:105:MET:SD	2.54	0.48
1:C:109:LEU:HD21	1:C:188:TYR:CG	2.49	0.48
1:C:338:LYS:O	1:C:342:GLU:HG3	2.13	0.48
1:F:37:GLU:HB2	1:F:52:TYR:CE1	2.49	0.47
1:H:327:ARG:NH2	1:H:330:GLN:HA	2.29	0.47
1:E:281:HIS:HD2	1:E:283:LEU:H	1.62	0.47
1:G:67:GLU:OE1	1:G:196:ARG:NE	2.43	0.47
1:A:281:HIS:HD2	1:A:283:LEU:H	1.61	0.47
1:F:39:PHE:CD1	1:F:65:HIS:CD2	3.02	0.47
1:F:130:PRO:HD3	1:F:178:GLN:HB2	1.95	0.47
1:G:79:LEU:O	1:G:88:VAL:HG11	2.14	0.47
1:A:130:PRO:HD3	1:A:178:GLN:HB2	1.95	0.47
1:E:130:PRO:HD3	1:E:178:GLN:HB2	1.97	0.47
1:A:109:LEU:HD21	1:A:188:TYR:CG	2.50	0.47
1:F:131:LEU:HD11	1:F:156:ARG:HG2	1.97	0.47
1:G:296:ALA:O	1:G:300:SER:HB3	2.15	0.47
1:D:183:SER:HB2	1:D:301:THR:HG21	1.95	0.47
1:G:327:ARG:HD3	1:G:327:ARG:HA	1.68	0.47
1:G:281:HIS:HE1	1:G:305:VAL:O	1.98	0.46
1:H:37:GLU:HB2	1:H:52:TYR:CE1	2.50	0.46
1:A:64:LEU:HD22	1:A:197:LEU:HD22	1.98	0.46
1:G:259:TYR:HD1	1:G:284:ALA:CB	2.25	0.46
1:A:132:LEU:CD2	1:A:271:LEU:HD12	2.45	0.46
1:B:128:LEU:O	1:B:131:LEU:N	2.45	0.46
1:F:316:GLU:OE2	1:F:338:LYS:HB2	2.15	0.46
1:G:75:LEU:HD23	1:G:105:MET:SD	2.56	0.46
1:H:134:LEU:HG	1:H:147:PRO:HG3	1.98	0.46
1:B:294:LEU:HD23	1:B:294:LEU:C	2.35	0.46
1:D:258:ASN:HB3	1:D:260:PHE:CZ	2.51	0.46
1:E:150:GLU:CA	1:E:150:GLU:OE2	2.64	0.46
1:A:210:THR:HB	1:D:29:MET:HB2	1.97	0.45
1:G:223:SER:HB2	1:F:260:PHE:CZ	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:261:GLU:HA	1:E:266:ARG:HA	1.97	0.45
1:E:37:GLU:HB2	1:E:52:TYR:CE1	2.50	0.45
1:F:339:ALA:O	1:F:343:LEU:HD23	2.16	0.45
1:E:258:ASN:HB3	1:E:260:PHE:CZ	2.51	0.45
1:G:218:ARG:NH1	1:G:222:GLY:O	2.50	0.45
1:H:78:GLN:OE1	1:H:105:MET:HB2	2.16	0.45
1:B:272:ASP:HB3	1:B:275:SER:OG	2.16	0.45
1:G:109:LEU:HD21	1:G:188:TYR:CG	2.52	0.45
1:B:129:GLU:HG2	1:B:180:ASP:HB2	1.99	0.45
1:F:314:LEU:HG	1:F:318:ASN:HD22	1.82	0.45
1:F:316:GLU:OE2	1:F:338:LYS:CB	2.65	0.45
1:G:142:ARG:HG3	1:G:145:ARG:CZ	2.47	0.44
1:D:281:HIS:HD2	1:D:283:LEU:H	1.65	0.44
1:H:188:TYR:CZ	1:H:192:LEU:HD11	2.53	0.44
1:A:131:LEU:O	1:A:135:TRP:CD1	2.67	0.44
1:F:135:TRP:HB2	1:F:137:PHE:CE1	2.53	0.44
1:B:37:GLU:HB2	1:B:52:TYR:CE1	2.51	0.44
1:D:37:GLU:HB2	1:D:52:TYR:CE1	2.53	0.44
1:F:109:LEU:HD21	1:F:188:TYR:CG	2.53	0.44
1:E:277:GLN:HB3	1:E:278:PRO:CD	2.48	0.44
1:G:64:LEU:HD13	1:G:197:LEU:HD22	2.00	0.44
1:G:37:GLU:HB2	1:G:52:TYR:CE1	2.53	0.43
1:A:127:THR:O	1:A:159:THR:HG22	2.18	0.43
1:B:182:ASN:OD1	1:C:30:ASP:HB3	2.18	0.43
1:F:55:SER:O	1:F:56:GLY:C	2.57	0.43
1:D:109:LEU:HD21	1:D:188:TYR:CG	2.54	0.43
1:G:135:TRP:CZ2	1:G:277:GLN:HA	2.53	0.43
1:A:64:LEU:HD22	1:A:197:LEU:CD2	2.48	0.43
1:C:136:GLY:HA3	1:C:143:GLY:HA2	2.00	0.43
1:D:158:LEU:HD22	1:D:176:ALA:HB3	2.00	0.43
1:B:312:LEU:HD11	1:B:335:THR:HG21	2.00	0.43
1:E:148:SER:OG	1:E:151:ASP:OD1	2.31	0.43
1:G:208:GLU:HA	1:G:213:LEU:HB3	2.01	0.43
1:G:326:VAL:HB	1:G:333:VAL:CG2	2.48	0.43
1:B:78:GLN:OE1	1:B:103:PRO:HB2	2.19	0.43
1:E:227:ILE:HG12	1:E:245:LEU:HD11	2.01	0.42
1:C:78:GLN:OE1	1:C:105:MET:HB2	2.19	0.42
1:E:261:GLU:OE1	1:E:327:ARG:NH1	2.52	0.42
1:G:158:LEU:HD22	1:G:176:ALA:HB3	2.00	0.42
1:C:327:ARG:HD3	1:C:332:PHE:CE1	2.54	0.42
1:A:271:LEU:HA	1:A:278:PRO:HA	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:135:TRP:CZ2	1:D:152:ILE:HD11	2.54	0.42
1:D:258:ASN:HB3	1:D:260:PHE:CE1	2.54	0.42
1:E:134:LEU:HG	1:E:147:PRO:CG	2.49	0.42
1:H:60:ALA:HB3	1:H:63:VAL:HG23	2.02	0.42
1:A:260:PHE:CZ	1:D:224:PRO:CD	2.95	0.42
1:F:136:GLY:HA3	1:F:143:GLY:HA2	2.00	0.42
1:F:183:SER:HB2	1:F:301:THR:HG21	2.02	0.41
1:G:35:LYS:HE2	1:G:36:VAL:O	2.20	0.41
1:A:277:GLN:CB	1:A:278:PRO:CD	2.94	0.41
1:A:76:ASP:OD1	1:A:83:ARG:HD3	2.20	0.41
1:B:272:ASP:OD2	1:B:274:GLN:N	2.53	0.41
1:H:135:TRP:CZ2	1:H:152:ILE:HD11	2.55	0.41
1:E:39:PHE:CD1	1:E:65:HIS:CD2	3.09	0.41
1:F:158:LEU:HD22	1:F:176:ALA:HB3	2.02	0.41
1:F:281:HIS:HD2	1:F:283:LEU:H	1.69	0.41
1:F:294:LEU:C	1:F:294:LEU:HD23	2.41	0.41
1:G:142:ARG:HG3	1:G:145:ARG:NH2	2.36	0.41
1:G:281:HIS:CE1	1:G:307:GLY:HA2	2.52	0.41
1:B:44:MET:HG3	1:C:33:GLN:O	2.21	0.41
1:F:139:PRO:CB	1:F:140:GLN:OE1	2.69	0.41
1:G:258:ASN:CB	1:G:260:PHE:HZ	2.26	0.41
1:H:294:LEU:C	1:H:294:LEU:HD23	2.41	0.41
1:H:253:SER:HB2	1:H:300:SER:OG	2.21	0.41
1:A:283:LEU:HD23	1:A:304:MET:HA	2.02	0.41
1:A:250:VAL:HG22	1:A:289:ILE:HG12	2.03	0.40
1:C:134:LEU:HG	1:C:147:PRO:HG3	2.03	0.40
1:F:76:ASP:OD1	1:F:83:ARG:HD3	2.21	0.40
1:H:326:VAL:HB	1:H:333:VAL:HG23	2.04	0.40
1:B:109:LEU:HD21	1:B:188:TYR:CG	2.57	0.40
1:D:40:THR:HB	1:D:48:TYR:O	2.22	0.40
1:E:261:GLU:H	1:E:261:GLU:HG2	1.79	0.40
1:F:129:GLU:N	1:F:130:PRO:HD2	2.36	0.40
1:G:105:MET:O	1:G:109:LEU:HG	2.21	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:B:84:SER:OG	1:A:328:GLU:OE1[2_646]	2.18	0.02

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	299/332 (90%)	283 (95%)	14 (5%)	2 (1%)	25	49
1	B	297/332 (90%)	279 (94%)	18 (6%)	0	100	100
1	C	304/332 (92%)	289 (95%)	14 (5%)	1 (0%)	44	70
1	D	301/332 (91%)	291 (97%)	9 (3%)	1 (0%)	44	70
1	E	312/332 (94%)	297 (95%)	15 (5%)	0	100	100
1	F	311/332 (94%)	298 (96%)	11 (4%)	2 (1%)	28	53
1	G	291/332 (88%)	275 (94%)	15 (5%)	1 (0%)	44	70
1	H	312/332 (94%)	301 (96%)	11 (4%)	0	100	100
All	All	2427/2656 (91%)	2313 (95%)	107 (4%)	7 (0%)	44	70

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	235	ASP
1	G	148	SER
1	F	276	GLY
1	C	221	ASP
1	D	330	GLN
1	F	147	PRO
1	A	58	GLY

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	243/265 (92%)	241 (99%)	2 (1%)	85	94
1	B	242/265 (91%)	238 (98%)	4 (2%)	66	86
1	C	245/265 (92%)	243 (99%)	2 (1%)	85	94
1	D	245/265 (92%)	243 (99%)	2 (1%)	85	94
1	E	250/265 (94%)	243 (97%)	7 (3%)	49	76
1	F	249/265 (94%)	245 (98%)	4 (2%)	68	87
1	G	239/265 (90%)	235 (98%)	4 (2%)	66	86
1	H	250/265 (94%)	245 (98%)	5 (2%)	60	83
All	All	1963/2120 (93%)	1933 (98%)	30 (2%)	70	88

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	H	94	LEU
1	H	128	LEU
1	H	145	ARG
1	H	146	VAL
1	H	327	ARG
1	B	74	GLN
1	B	128	LEU
1	B	274	GLN
1	B	348	VAL
1	G	128	LEU
1	G	256	TYR
1	G	258	ASN
1	G	327	ARG
1	A	128	LEU
1	A	238	VAL
1	C	29	MET
1	C	128	LEU
1	D	128	LEU
1	D	150	GLU
1	E	94	LEU
1	E	145	ARG
1	E	150	GLU
1	E	261	GLU
1	E	275	SER
1	E	316	GLU
1	E	348	VAL
1	F	135	TRP

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Mol	Chain	Res	Type
1	F	144	GLU
1	F	145	ARG
1	F	150	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	H	33	GLN
1	H	74	GLN
1	H	240	GLN
1	H	281	HIS
1	B	74	GLN
1	B	240	GLN
1	B	281	HIS
1	B	330	GLN
1	G	240	GLN
1	G	281	HIS
1	A	178	GLN
1	A	240	GLN
1	A	281	HIS
1	C	240	GLN
1	C	281	HIS
1	D	65	HIS
1	D	240	GLN
1	D	281	HIS
1	D	330	GLN
1	E	65	HIS
1	E	178	GLN
1	E	240	GLN
1	E	281	HIS
1	F	65	HIS
1	F	240	GLN
1	F	281	HIS
1	F	318	ASN
1	F	330	GLN

5.3.3 RNA ⓘ

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](#)

Of 8 ligands modelled in this entry, 8 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	G	1
1	A	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	G	281:HIS	C	282:HIS	N	1.89
1	A	259:TYR	C	260:PHE	N	1.76

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	305/332 (91%)	-0.19	9 (2%)	51	43	26, 48, 106, 146	0
1	B	303/332 (91%)	-0.21	8 (2%)	56	49	25, 44, 109, 143	0
1	C	310/332 (93%)	-0.21	7 (2%)	61	54	29, 54, 106, 141	0
1	D	309/332 (93%)	-0.22	4 (1%)	77	73	28, 49, 106, 144	0
1	E	316/332 (95%)	-0.17	5 (1%)	72	67	31, 53, 113, 145	0
1	F	315/332 (94%)	-0.17	6 (1%)	67	61	27, 52, 131, 170	0
1	G	299/332 (90%)	0.11	17 (5%)	24	18	29, 69, 116, 154	0
1	H	316/332 (95%)	-0.20	11 (3%)	44	36	26, 50, 107, 154	0
All	All	2473/2656 (93%)	-0.16	67 (2%)	55	48	25, 51, 113, 170	0

All (67) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	331	GLY	9.4
1	G	282	HIS	8.6
1	G	326	VAL	5.2
1	D	260	PHE	4.8
1	E	140	GLN	4.8
1	B	237	ARG	4.7
1	G	260	PHE	4.2
1	A	269	HIS	4.0
1	A	267	TYR	4.0
1	D	238	VAL	3.9
1	G	312	LEU	3.9
1	D	233	ARG	3.8
1	A	136	GLY	3.8
1	G	335	THR	3.7
1	F	263	ASP	3.7
1	F	152	ILE	3.7

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Mol	Chain	Res	Type	RSRZ
1	A	152	ILE	3.7
1	C	329	GLY	3.6
1	B	132	LEU	3.6
1	A	330	GLN	3.6
1	D	267	TYR	3.6
1	F	276	GLY	3.6
1	F	262	ARG	3.6
1	G	257	ARG	3.5
1	C	260	PHE	3.5
1	E	142	ARG	3.4
1	H	346	ALA	3.4
1	B	152	ILE	3.3
1	F	260	PHE	3.2
1	G	333	VAL	3.1
1	H	44	MET	3.0
1	C	142	ARG	2.9
1	G	323	PHE	2.9
1	B	276	GLY	2.9
1	E	237	ARG	2.8
1	G	29	MET	2.7
1	H	233	ARG	2.7
1	G	283	LEU	2.7
1	G	228	ALA	2.7
1	B	233	ARG	2.7
1	F	348	VAL	2.7
1	A	237	ARG	2.6
1	G	308	PRO	2.6
1	A	271	LEU	2.6
1	A	151	ASP	2.6
1	C	276	GLY	2.6
1	B	278	PRO	2.6
1	H	261	GLU	2.5
1	G	279	ILE	2.4
1	G	306	LEU	2.4
1	A	132	LEU	2.4
1	G	332	PHE	2.4
1	H	326	VAL	2.3
1	H	329	GLY	2.3
1	H	238	VAL	2.3
1	C	277	GLN	2.3
1	H	260	PHE	2.3
1	E	139	PRO	2.2

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Mol	Chain	Res	Type	RSRZ
1	E	348	VAL	2.2
1	B	260	PHE	2.2
1	B	269	HIS	2.2
1	H	330	GLN	2.1
1	C	330	GLN	2.1
1	H	332	PHE	2.1
1	H	144	GLU	2.1
1	G	310	LYS	2.1
1	C	146	VAL	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	MG	D	401	1/1	0.95	0.13	0.55	51,51,51,51	0
2	MG	F	401	1/1	0.82	0.07	-1.30	46,46,46,46	0
2	MG	H	401	1/1	0.83	0.11	-1.53	51,51,51,51	0
2	MG	B	401	1/1	0.92	0.11	-2.09	45,45,45,45	0
2	MG	G	401	1/1	0.94	0.08	-2.72	61,61,61,61	0
2	MG	A	401	1/1	0.97	0.08	-2.97	40,40,40,40	0
2	MG	E	401	1/1	0.95	0.11	-3.49	51,51,51,51	0
2	MG	C	401	1/1	0.99	0.05	-4.13	38,38,38,38	0

6.5 Other polymers [i](#)

There are no such residues in this entry.