



# Full wwPDB X-ray Structure Validation Report ⓘ

May 29, 2020 – 04:57 am BST

PDB ID : 4G3H  
Title : Crystal structure of helicobacter pylori arginase  
Authors : Zhang, J.; Zhang, X.; Li, D.; Hu, Y.; Zou, Q.; Wang, D.  
Deposited on : 2012-07-13  
Resolution : 2.20 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](mailto:validation@mail.wwpdb.org)

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

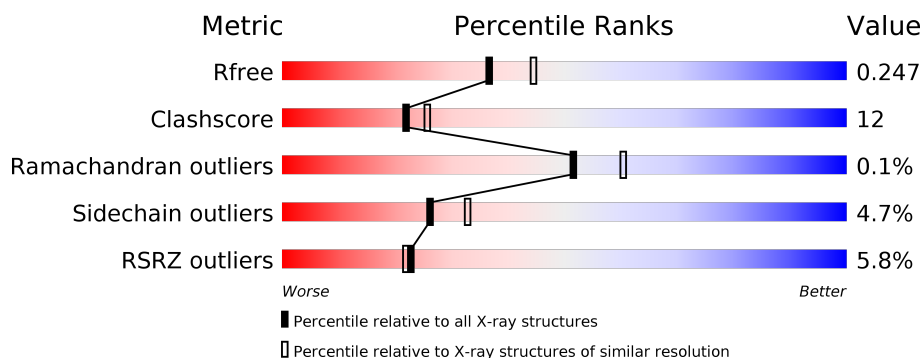
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.20 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))
$R_{free}$	130704	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. 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	330	<div> <div>5%</div> <div> <div></div> <div>65%</div> <div>22%</div> <div>•</div> <div>9%</div> </div> </div>
1	B	330	<div> <div>7%</div> <div> <div></div> <div>65%</div> <div>24%</div> <div>•</div> <div>8%</div> </div> </div>
1	C	330	<div> <div>5%</div> <div> <div></div> <div>70%</div> <div>21%</div> <div>•</div> <div>6%</div> </div> </div>
1	D	330	<div> <div>4%</div> <div> <div></div> <div>66%</div> <div>22%</div> <div>•</div> <div>9%</div> </div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10439 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 Arginase (RocF).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	301	Total	C	N	O	S	0	0	0
			2419	1539	411	454	15			
1	B	305	Total	C	N	O	S	0	0	0
			2447	1551	417	465	14			
1	C	310	Total	C	N	O	S	6	0	0
			2491	1581	424	471	15			
1	D	300	Total	C	N	O	S	0	0	0
			2399	1524	407	453	15			

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	323	LEU	-	EXPRESSION TAG	UNP O25949
A	324	GLU	-	EXPRESSION TAG	UNP O25949
A	325	HIS	-	EXPRESSION TAG	UNP O25949
A	326	HIS	-	EXPRESSION TAG	UNP O25949
A	327	HIS	-	EXPRESSION TAG	UNP O25949
A	328	HIS	-	EXPRESSION TAG	UNP O25949
A	329	HIS	-	EXPRESSION TAG	UNP O25949
A	330	HIS	-	EXPRESSION TAG	UNP O25949
B	323	LEU	-	EXPRESSION TAG	UNP O25949
B	324	GLU	-	EXPRESSION TAG	UNP O25949
B	325	HIS	-	EXPRESSION TAG	UNP O25949
B	326	HIS	-	EXPRESSION TAG	UNP O25949
B	327	HIS	-	EXPRESSION TAG	UNP O25949
B	328	HIS	-	EXPRESSION TAG	UNP O25949
B	329	HIS	-	EXPRESSION TAG	UNP O25949
B	330	HIS	-	EXPRESSION TAG	UNP O25949
C	323	LEU	-	EXPRESSION TAG	UNP O25949
C	324	GLU	-	EXPRESSION TAG	UNP O25949
C	325	HIS	-	EXPRESSION TAG	UNP O25949
C	326	HIS	-	EXPRESSION TAG	UNP O25949
C	327	HIS	-	EXPRESSION TAG	UNP O25949

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Chain	Residue	Modelled	Actual	Comment	Reference
C	328	HIS	-	EXPRESSION TAG	UNP O25949
C	329	HIS	-	EXPRESSION TAG	UNP O25949
C	330	HIS	-	EXPRESSION TAG	UNP O25949
D	323	LEU	-	EXPRESSION TAG	UNP O25949
D	324	GLU	-	EXPRESSION TAG	UNP O25949
D	325	HIS	-	EXPRESSION TAG	UNP O25949
D	326	HIS	-	EXPRESSION TAG	UNP O25949
D	327	HIS	-	EXPRESSION TAG	UNP O25949
D	328	HIS	-	EXPRESSION TAG	UNP O25949
D	329	HIS	-	EXPRESSION TAG	UNP O25949
D	330	HIS	-	EXPRESSION TAG	UNP O25949

- Molecule 2 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	2	Total 2	Mn 2	0	0
2	A	2	Total 2	Mn 2	0	0
2	D	2	Total 2	Mn 2	0	0
2	C	2	Total 2	Mn 2	0	0

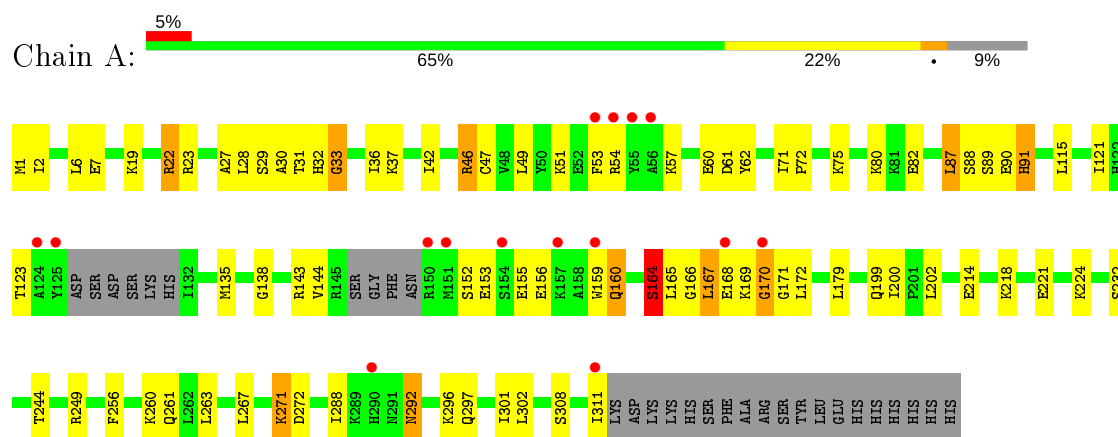
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	174	Total 174	O 174	0	0
3	B	153	Total 153	O 153	0	0
3	C	168	Total 168	O 168	0	0
3	D	180	Total 180	O 180	0	0

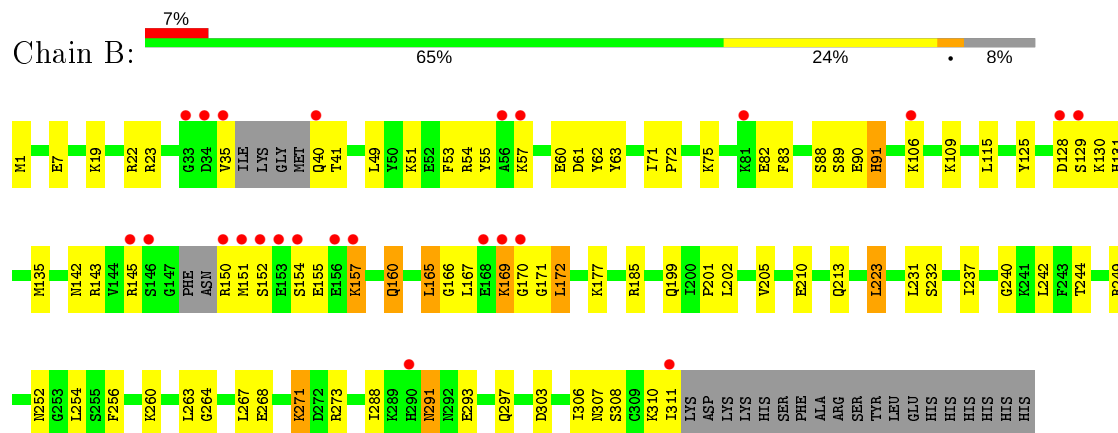
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

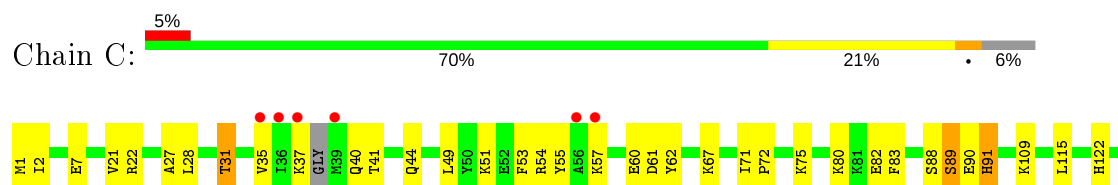
#### • Molecule 1: Arginase (RocF)

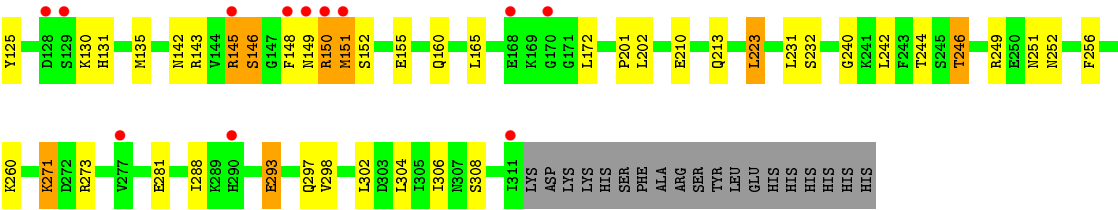


#### • Molecule 1: Arginase (RocF)

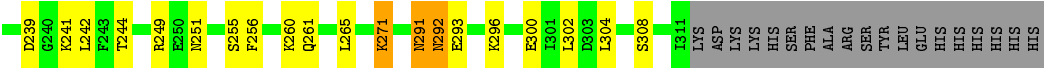
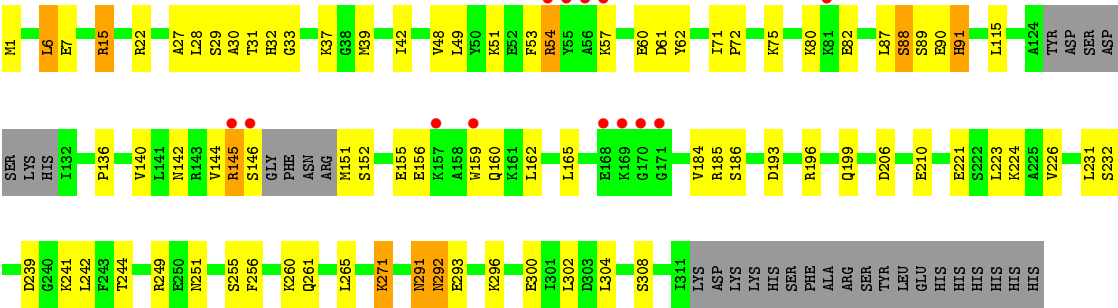


#### • Molecule 1: Arginase (RocF)





● Molecule 1: Arginase (RocF)



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	94.69Å 102.24Å 148.61Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	34.90 – 2.20 34.92 – 2.20	Depositor EDS
% Data completeness (in resolution range)	(Not available) (34.90-2.20) 99.9 (34.92-2.20)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.62 (at 2.20Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, $R_{free}$	0.218 , 0.230 0.208 , 0.247	Depositor DCC
$R_{free}$ test set	7448 reflections (10.10%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	23.1	Xtriage
Anisotropy	1.069	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 55.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.55$ , $\langle L^2 \rangle = 0.39$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	10439	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	36.0	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 52.93 % 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 4.5344e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<sup>1</sup> Intensities estimated from amplitudes.

<sup>2</sup> Theoretical 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: MN

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	1.22	35/2453 (1.4%)	0.78	10/3289 (0.3%)
1	B	0.82	14/2482 (0.6%)	0.67	5/3329 (0.2%)
1	C	0.67	8/2528 (0.3%)	0.63	1/3391 (0.0%)
1	D	1.26	31/2432 (1.3%)	0.75	5/3261 (0.2%)
All	All	1.02	88/9895 (0.9%)	0.71	21/13270 (0.2%)

All (88) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	89	SER	CB-OG	-16.67	1.20	1.42
1	D	29	SER	CB-OG	-16.50	1.20	1.42
1	D	184	VAL	CB-CG2	-15.99	1.19	1.52
1	B	89	SER	CB-OG	-15.55	1.22	1.42
1	D	89	SER	C-O	-15.13	0.94	1.23
1	A	89	SER	C-O	-14.95	0.94	1.23
1	A	29	SER	CB-OG	-14.79	1.23	1.42
1	C	89	SER	CB-OG	-14.72	1.23	1.42
1	D	88	SER	C-O	-13.75	0.97	1.23
1	A	88	SER	CB-OG	-13.72	1.24	1.42
1	D	186	SER	CB-OG	-12.95	1.25	1.42
1	C	88	SER	CB-OG	-12.76	1.25	1.42
1	D	184	VAL	C-O	-12.57	0.99	1.23
1	A	33	GLY	C-O	-12.34	1.03	1.23
1	D	88	SER	CB-OG	-12.31	1.26	1.42
1	D	185	ARG	C-O	-12.06	1.00	1.23
1	B	89	SER	C-O	-11.24	1.01	1.23
1	B	88	SER	CB-OG	-11.23	1.27	1.42
1	A	89	SER	CB-OG	-11.22	1.27	1.42
1	A	88	SER	C-O	-11.20	1.02	1.23
1	D	28	LEU	C-O	-10.99	1.02	1.23
1	A	29	SER	C-O	-10.85	1.02	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	88	SER	C-O	-10.85	1.02	1.23
1	D	30	ALA	C-O	-10.67	1.03	1.23
1	A	90	GLU	CD-OE2	-10.17	1.14	1.25
1	B	90	GLU	CB-CG	-10.16	1.32	1.52
1	B	88	SER	C-O	-10.03	1.04	1.23
1	C	89	SER	C-O	-9.95	1.04	1.23
1	D	185	ARG	CZ-NH1	-9.81	1.20	1.33
1	B	165	LEU	C-O	-9.72	1.04	1.23
1	A	171	GLY	C-O	-9.41	1.08	1.23
1	C	90	GLU	CD-OE2	-9.22	1.15	1.25
1	A	27	ALA	C-O	-9.03	1.06	1.23
1	A	87	LEU	C-O	-9.03	1.06	1.23
1	D	90	GLU	CB-CG	-8.92	1.35	1.52
1	A	166	GLY	C-O	-8.90	1.09	1.23
1	A	30	ALA	C-O	-8.85	1.06	1.23
1	A	164	SER	CB-OG	-8.80	1.30	1.42
1	A	28	LEU	C-O	-8.66	1.06	1.23
1	D	87	LEU	CG-CD1	-8.12	1.21	1.51
1	D	31	THR	CB-CG2	-8.11	1.25	1.52
1	A	172	LEU	C-O	-7.89	1.08	1.23
1	D	185	ARG	CG-CD	-7.81	1.32	1.51
1	D	87	LEU	C-O	-7.78	1.08	1.23
1	D	87	LEU	CG-CD2	-7.71	1.23	1.51
1	D	29	SER	C-O	-7.65	1.08	1.23
1	A	90	GLU	C-O	-7.64	1.08	1.23
1	A	165	LEU	CG-CD1	-7.56	1.23	1.51
1	B	166	GLY	C-O	-7.42	1.11	1.23
1	A	31	THR	CB-CG2	-7.29	1.28	1.52
1	D	185	ARG	CZ-NH2	-7.26	1.23	1.33
1	D	27	ALA	C-O	-7.24	1.09	1.23
1	A	167	LEU	CG-CD2	-7.09	1.25	1.51
1	A	87	LEU	CG-CD1	-7.05	1.25	1.51
1	D	32	HIS	C-O	-7.05	1.09	1.23
1	C	90	GLU	CD-OE1	-7.03	1.18	1.25
1	D	28	LEU	CG-CD2	-6.96	1.26	1.51
1	A	165	LEU	C-O	-6.88	1.10	1.23
1	B	167	LEU	C-O	-6.80	1.10	1.23
1	A	87	LEU	CG-CD2	-6.74	1.26	1.51
1	A	164	SER	C-O	-6.72	1.10	1.23
1	A	32	HIS	C-O	-6.64	1.10	1.23
1	D	90	GLU	CD-OE1	-6.64	1.18	1.25
1	D	29	SER	N-CA	-6.62	1.33	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	166	GLY	N-CA	-6.57	1.36	1.46
1	A	90	GLU	CB-CG	-6.52	1.39	1.52
1	B	90	GLU	C-O	-6.33	1.11	1.23
1	D	29	SER	CA-CB	-6.25	1.43	1.52
1	A	169	LYS	N-CA	-6.12	1.34	1.46
1	A	31	THR	C-O	-6.02	1.11	1.23
1	C	90	GLU	C-O	-5.99	1.11	1.23
1	D	186	SER	N-CA	-5.90	1.34	1.46
1	D	30	ALA	CA-CB	-5.88	1.40	1.52
1	B	171	GLY	C-O	-5.74	1.14	1.23
1	D	31	THR	C-O	-5.62	1.12	1.23
1	A	29	SER	N-CA	-5.53	1.35	1.46
1	D	186	SER	C-O	-5.53	1.12	1.23
1	B	90	GLU	CD-OE2	-5.40	1.19	1.25
1	A	171	GLY	CA-C	-5.32	1.43	1.51
1	D	28	LEU	CG-CD1	-5.23	1.32	1.51
1	A	33	GLY	N-CA	-5.22	1.38	1.46
1	A	172	LEU	C-N	-5.22	1.22	1.34
1	A	30	ALA	N-CA	-5.17	1.36	1.46
1	A	29	SER	CA-CB	-5.13	1.45	1.52
1	B	169	LYS	N-CA	-5.08	1.36	1.46
1	A	28	LEU	CG-CD2	-5.07	1.33	1.51
1	B	165	LEU	CG-CD1	-5.05	1.33	1.51
1	C	89	SER	N-CA	-5.02	1.36	1.46

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	54	ARG	NE-CZ-NH2	-10.39	115.11	120.30
1	D	54	ARG	NE-CZ-NH1	9.65	125.13	120.30
1	B	171	GLY	N-CA-C	8.63	134.67	113.10
1	B	128	ASP	N-CA-C	-7.55	90.62	111.00
1	B	165	LEU	CB-CG-CD1	-7.07	98.97	111.00
1	A	167	LEU	CA-CB-CG	6.79	130.91	115.30
1	C	149	ASN	N-CA-C	-6.31	93.97	111.00
1	D	185	ARG	NE-CZ-NH1	-6.07	117.27	120.30
1	D	185	ARG	NE-CZ-NH2	6.05	123.33	120.30
1	A	167	LEU	O-C-N	-6.05	113.02	122.70
1	A	168	GLU	N-CA-CB	6.00	121.40	110.60
1	A	172	LEU	CB-CG-CD1	5.73	120.74	111.00
1	A	87	LEU	CB-CG-CD1	5.72	120.73	111.00
1	A	165	LEU	CB-CG-CD2	-5.66	101.38	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	169	LYS	CD-CE-NZ	5.62	124.64	111.70
1	D	87	LEU	CB-CG-CD1	5.55	120.43	111.00
1	B	170	GLY	N-CA-C	5.45	126.72	113.10
1	A	170	GLY	N-CA-C	5.16	126.01	113.10
1	A	171	GLY	N-CA-C	-5.12	100.30	113.10
1	A	46	ARG	CB-CG-CD	5.09	124.83	111.60
1	B	169	LYS	CB-CA-C	-5.07	100.25	110.40

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	2419	0	2472	57	0
1	B	2447	0	2482	66	0
1	C	2491	0	2530	62	0
1	D	2399	0	2446	56	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
3	A	174	0	0	4	0
3	B	153	0	0	7	0
3	C	168	0	0	8	0
3	D	180	0	0	4	0
All	All	10439	0	9930	233	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 12.

All (233) 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:A:123:THR:H	1:A:135:MET:HE2	1.26	1.00

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:244:THR:HA	1:C:249:ARG:HD3	1.40	0.98
1:B:244:THR:HA	1:B:249:ARG:HD3	1.44	0.97
1:A:244:THR:HA	1:A:249:ARG:HD3	1.53	0.91
1:B:1:MET:HA	1:B:82:GLU:HG2	1.55	0.88
1:D:244:THR:HA	1:D:249:ARG:HD3	1.56	0.88
1:C:1:MET:HA	1:C:82:GLU:HG2	1.62	0.81
1:D:151:MET:CE	1:D:156:GLU:HA	2.13	0.78
1:A:143:ARG:HD3	3:A:718:HOH:O	1.84	0.77
1:A:311:ILE:HD12	1:A:311:ILE:H	1.50	0.76
1:C:151:MET:CB	1:C:155:GLU:HB2	2.18	0.74
1:D:156:GLU:O	1:D:160:GLN:HG2	1.90	0.72
1:A:22:ARG:HD3	1:A:23:ARG:N	2.06	0.70
1:C:145:ARG:HB3	1:C:151:MET:HE2	1.74	0.70
1:B:106:LYS:HD3	3:B:695:HOH:O	1.93	0.68
1:D:151:MET:HE3	1:D:156:GLU:HA	1.76	0.68
1:A:123:THR:N	1:A:135:MET:HE2	2.07	0.67
1:B:1:MET:HB2	1:B:83:PHE:O	1.94	0.67
1:C:27:ALA:O	1:C:31:THR:HG23	1.95	0.66
1:C:1:MET:HB2	1:C:83:PHE:O	1.95	0.66
1:C:151:MET:HB2	1:C:155:GLU:HB2	1.77	0.66
1:C:151:MET:HB3	1:C:155:GLU:HB2	1.78	0.65
1:C:21:VAL:HG21	1:C:89:SER:HB2	1.79	0.65
1:A:292:ASN:HB3	1:A:296:LYS:HZ3	1.62	0.64
1:A:167:LEU:O	1:A:170:GLY:N	2.30	0.64
1:B:152:SER:OG	1:B:155:GLU:HG3	1.98	0.63
1:D:80:LYS:HE2	3:D:755:HOH:O	1.98	0.63
1:D:145:ARG:O	1:D:146:SER:CB	2.47	0.62
1:B:154:SER:O	1:B:157:LYS:HG2	1.99	0.62
1:C:2:ILE:HD12	1:C:44:GLN:HB3	1.79	0.62
1:B:240:GLY:O	1:B:249:ARG:HD2	2.00	0.62
1:B:271:LYS:HD3	3:B:720:HOH:O	1.99	0.61
1:D:291:ASN:C	1:D:291:ASN:HD22	2.03	0.61
1:D:54:ARG:HB2	1:D:57:LYS:HB2	1.81	0.61
1:A:297:GLN:O	1:A:301:ILE:HG12	2.01	0.60
1:B:157:LYS:NZ	1:B:157:LYS:HB3	2.17	0.60
1:B:154:SER:HA	1:B:157:LYS:HE2	1.84	0.60
1:D:53:PHE:HB2	1:D:62:TYR:CE1	2.37	0.60
1:A:311:ILE:HD12	1:A:311:ILE:N	2.18	0.59
1:B:256:PHE:CE2	1:B:260:LYS:HD2	2.37	0.59
1:C:246:THR:HG22	1:C:281:GLU:O	2.02	0.59
1:B:151:MET:HB3	1:B:155:GLU:HB2	1.85	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:55:TYR:HE2	1:B:135:MET:HE2	1.69	0.58
1:C:223:LEU:O	1:C:273:ARG:NH2	2.36	0.58
1:C:256:PHE:CZ	1:C:260:LYS:HD2	2.38	0.58
1:A:271:LYS:HE3	1:A:308:SER:O	2.04	0.58
1:A:53:PHE:HB2	1:A:62:TYR:CE1	2.39	0.58
1:C:35:VAL:O	1:C:35:VAL:HG13	2.03	0.57
1:A:22:ARG:NH2	1:A:23:ARG:HG2	2.19	0.57
1:A:46:ARG:HG2	1:A:47:CYS:N	2.18	0.57
1:A:91:HIS:HE1	1:A:232:SER:OG	1.88	0.57
1:A:292:ASN:HB3	1:A:296:LYS:NZ	2.19	0.57
1:D:151:MET:HE2	1:D:156:GLU:HA	1.86	0.57
1:D:241:LYS:HE3	1:D:251:ASN:HD21	1.69	0.57
1:C:55:TYR:HE2	1:C:135:MET:HE2	1.70	0.57
1:C:91:HIS:HE1	1:C:232:SER:OG	1.87	0.57
1:A:152:SER:HB2	1:A:155:GLU:HG3	1.87	0.56
1:B:145:ARG:NH1	1:B:160:GLN:HE22	2.03	0.56
1:A:256:PHE:CZ	1:A:260:LYS:HD2	2.41	0.56
1:D:292:ASN:HD22	1:D:292:ASN:H	1.53	0.56
1:C:151:MET:HB3	1:C:155:GLU:OE1	2.06	0.56
1:D:292:ASN:HD22	1:D:292:ASN:N	2.03	0.55
1:B:271:LYS:NZ	1:B:271:LYS:HB2	2.21	0.55
1:C:1:MET:HE3	1:C:306:ILE:HG23	1.87	0.55
1:A:22:ARG:HH21	1:A:23:ARG:HG2	1.72	0.55
1:B:256:PHE:CZ	1:B:260:LYS:HD2	2.41	0.55
1:D:115:LEU:HD12	1:D:115:LEU:N	2.21	0.55
1:D:33:GLY:HA2	1:D:42:ILE:HB	1.88	0.55
1:D:54:ARG:CB	1:D:57:LYS:HB2	2.36	0.55
1:B:264:GLY:O	1:B:268:GLU:HG3	2.07	0.55
1:C:143:ARG:HG2	1:C:143:ARG:HH11	1.72	0.55
1:C:298:VAL:O	1:C:302:LEU:HD13	2.07	0.54
1:A:214:GLU:HG2	1:A:218:LYS:HE3	1.90	0.54
1:A:80:LYS:HE2	3:A:737:HOH:O	2.07	0.54
1:B:23:ARG:HH21	1:B:23:ARG:HG2	1.72	0.54
1:A:115:LEU:N	1:A:115:LEU:HD12	2.22	0.54
1:D:152:SER:OG	1:D:155:GLU:HG3	2.08	0.54
1:D:91:HIS:HE1	1:D:232:SER:OG	1.91	0.54
1:C:53:PHE:HB2	1:C:62:TYR:CE1	2.43	0.53
1:C:271:LYS:HE3	1:C:308:SER:O	2.09	0.53
1:B:1:MET:N	1:B:41:THR:O	2.41	0.53
1:B:53:PHE:HB2	1:B:62:TYR:CE1	2.43	0.53
1:B:291:ASN:H	1:B:291:ASN:HD22	1.56	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:153:GLU:O	1:A:156:GLU:HG2	2.09	0.53
1:A:54:ARG:HB2	1:A:57:LYS:HB2	1.91	0.53
1:D:145:ARG:O	1:D:145:ARG:HD3	2.09	0.53
1:A:160:GLN:O	1:A:164:SER:HB2	2.09	0.53
1:C:28:LEU:HD11	1:C:302:LEU:HD23	1.91	0.53
1:D:292:ASN:ND2	1:D:292:ASN:H	2.06	0.53
1:A:123:THR:H	1:A:135:MET:CE	2.10	0.52
1:A:51:LYS:HE2	1:A:61:ASP:HB3	1.91	0.52
1:D:6:LEU:HG	1:D:88:SER:HA	1.90	0.52
1:B:291:ASN:HD22	1:B:291:ASN:N	2.05	0.52
1:C:51:LYS:HE2	1:C:61:ASP:HB3	1.90	0.52
1:D:293:GLU:HA	1:D:296:LYS:HZ3	1.75	0.52
1:B:41:THR:HG21	1:B:306:ILE:CG2	2.40	0.52
1:C:145:ARG:HB3	1:C:151:MET:CE	2.39	0.51
1:A:221:GLU:O	1:A:224:LYS:HG2	2.11	0.51
1:B:51:LYS:HE2	1:B:61:ASP:HB3	1.91	0.51
1:C:271:LYS:HD3	3:C:760:HOH:O	2.09	0.51
1:B:19:LYS:NZ	3:B:724:HOH:O	2.43	0.51
1:B:54:ARG:CB	1:B:57:LYS:HB2	2.40	0.51
1:C:115:LEU:N	1:C:115:LEU:HD12	2.25	0.51
1:C:54:ARG:CB	1:C:57:LYS:HB2	2.40	0.51
1:D:145:ARG:O	1:D:146:SER:HB2	2.11	0.50
1:B:22:ARG:HH21	1:B:22:ARG:HG2	1.76	0.50
1:C:54:ARG:HB2	1:C:57:LYS:HB2	1.92	0.50
1:A:54:ARG:CB	1:A:57:LYS:HB2	2.42	0.50
1:B:223:LEU:O	1:B:273:ARG:NH2	2.45	0.50
1:C:240:GLY:O	1:C:249:ARG:HD2	2.12	0.50
1:B:54:ARG:HB2	1:B:57:LYS:HB2	1.92	0.50
1:D:51:LYS:HE2	1:D:61:ASP:HB3	1.93	0.50
1:B:115:LEU:HD12	1:B:115:LEU:N	2.26	0.50
1:C:54:ARG:HD3	3:C:637:HOH:O	2.12	0.50
1:A:202:LEU:HD23	1:A:202:LEU:C	2.32	0.49
1:A:244:THR:HG22	1:A:249:ARG:NH1	2.27	0.49
1:C:80:LYS:HE2	3:C:745:HOH:O	2.12	0.49
1:D:15:ARG:HD2	3:D:648:HOH:O	2.13	0.49
1:C:210:GLU:HG3	3:C:750:HOH:O	2.13	0.49
1:B:150:ARG:HA	1:B:150:ARG:HE	1.78	0.49
1:A:261:GLN:HE21	1:D:261:GLN:HE21	1.61	0.49
1:D:39:MET:HE2	3:D:728:HOH:O	2.12	0.49
1:C:67:LYS:HB2	1:C:165:LEU:HD22	1.96	0.48
1:D:1:MET:CA	1:D:82:GLU:HG2	2.42	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:19:LYS:HA	1:A:22:ARG:HG3	1.94	0.48
1:B:271:LYS:HE3	1:B:308:SER:O	2.13	0.48
1:B:291:ASN:ND2	1:B:291:ASN:N	2.61	0.48
1:C:35:VAL:HA	1:C:40:GLN:HB2	1.96	0.48
1:A:144:VAL:HG11	1:A:159:TRP:CD2	2.50	0.47
1:D:1:MET:HA	1:D:82:GLU:HG2	1.96	0.47
1:B:91:HIS:HE1	1:B:232:SER:OG	1.98	0.47
1:D:256:PHE:CZ	1:D:260:LYS:HD2	2.50	0.47
1:D:136:PRO:O	1:D:140:VAL:HG23	2.13	0.47
1:D:162:LEU:HD12	1:D:165:LEU:HD12	1.96	0.47
1:A:36:ILE:HD12	1:B:199:GLN:HE22	1.81	0.46
1:D:271:LYS:HE3	1:D:308:SER:O	2.15	0.46
1:C:231:LEU:HD23	1:C:231:LEU:C	2.35	0.46
1:B:293:GLU:O	1:B:297:GLN:HG3	2.15	0.46
1:B:63:TYR:CE1	1:B:165:LEU:HD21	2.51	0.46
1:B:109:LYS:HE2	3:B:706:HOH:O	2.14	0.46
1:B:242:LEU:HD23	3:B:701:HOH:O	2.16	0.46
1:B:268:GLU:HG2	3:B:718:HOH:O	2.16	0.46
1:B:125:TYR:CD1	1:B:131:HIS:HA	2.51	0.46
1:C:293:GLU:O	1:C:297:GLN:HG3	2.16	0.46
1:B:35:VAL:HA	1:B:40:GLN:HB2	1.98	0.46
1:B:157:LYS:HZ2	1:B:157:LYS:HB3	1.80	0.45
3:A:709:HOH:O	1:B:177:LYS:HB3	2.16	0.45
1:A:311:ILE:CD1	1:A:311:ILE:H	2.22	0.45
1:D:7:GLU:O	1:D:49:LEU:HA	2.16	0.45
1:C:125:TYR:CD1	1:C:131:HIS:HA	2.51	0.45
1:A:7:GLU:O	1:A:49:LEU:HA	2.16	0.45
1:B:151:MET:HA	1:B:155:GLU:OE1	2.17	0.45
1:B:71:ILE:HB	1:B:72:PRO:HD3	1.99	0.45
1:C:67:LYS:HB2	1:C:165:LEU:CD2	2.46	0.45
1:B:244:THR:HG21	1:B:288:ILE:HG12	1.99	0.45
1:D:292:ASN:ND2	1:D:292:ASN:N	2.65	0.45
1:A:1:MET:CA	1:A:82:GLU:HG2	2.46	0.45
1:B:310:LYS:O	1:B:311:ILE:HB	2.17	0.45
1:A:2:ILE:HG12	1:A:82:GLU:HB3	1.97	0.45
1:B:213:GLN:H	1:B:213:GLN:CD	2.20	0.45
1:B:71:ILE:O	1:B:75:LYS:HG3	2.16	0.45
1:C:7:GLU:O	1:C:49:LEU:HA	2.17	0.45
1:D:71:ILE:HB	1:D:72:PRO:HD3	1.98	0.45
1:B:202:LEU:HD23	1:B:202:LEU:C	2.37	0.44
1:B:205:VAL:HG13	1:B:254:LEU:HD23	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:244:THR:HG22	1:B:249:ARG:NH1	2.32	0.44
1:D:223:LEU:O	1:D:226:VAL:HG22	2.17	0.44
1:C:213:GLN:H	1:C:213:GLN:CD	2.21	0.44
1:D:145:ARG:HB3	1:D:151:MET:SD	2.58	0.44
1:A:261:GLN:HG3	1:D:261:GLN:CG	2.48	0.44
1:B:185:ARG:NH2	1:B:237:ILE:O	2.40	0.43
1:D:144:VAL:HG21	1:D:159:TRP:CE3	2.53	0.43
1:C:22:ARG:HH21	1:C:22:ARG:HG2	1.84	0.43
1:A:271:LYS:NZ	1:A:271:LYS:HB2	2.33	0.43
1:D:193:ASP:OD1	1:D:196:ARG:NH2	2.52	0.43
1:D:71:ILE:O	1:D:75:LYS:HG3	2.18	0.43
1:C:151:MET:HB3	1:C:155:GLU:CB	2.46	0.43
1:A:121:ILE:O	1:A:138:GLY:HA3	2.18	0.43
1:C:246:THR:CG2	1:C:281:GLU:O	2.65	0.43
1:C:71:ILE:HB	1:C:72:PRO:HD3	2.00	0.43
1:A:33:GLY:HA2	1:A:42:ILE:HB	1.99	0.43
1:A:71:ILE:HB	1:A:72:PRO:HD3	1.99	0.43
1:A:36:ILE:HD12	1:B:199:GLN:NE2	2.33	0.43
1:D:304:LEU:HD23	1:D:304:LEU:C	2.38	0.43
1:C:244:THR:HG21	1:C:288:ILE:HG12	2.01	0.43
1:C:60:GLU:OE1	1:C:60:GLU:HA	2.19	0.43
1:C:256:PHE:CE2	1:C:260:LYS:HD2	2.54	0.43
1:D:144:VAL:HG11	1:D:159:TRP:CD2	2.54	0.43
1:A:263:LEU:O	1:A:267:LEU:HG	2.19	0.42
1:C:1:MET:N	1:C:41:THR:O	2.49	0.42
1:D:6:LEU:HD22	1:D:48:VAL:HB	2.01	0.42
1:C:122:HIS:CE1	1:C:146:SER:HB3	2.55	0.42
1:D:241:LYS:HG2	1:D:251:ASN:ND2	2.34	0.42
1:A:160:GLN:NE2	1:A:160:GLN:C	2.73	0.42
1:B:172:LEU:HD12	1:B:172:LEU:HA	1.87	0.42
1:C:202:LEU:HD23	1:C:202:LEU:C	2.40	0.42
1:D:231:LEU:HD23	1:D:231:LEU:C	2.39	0.42
1:D:242:LEU:HD23	1:D:255:SER:HA	2.01	0.42
1:D:199:GLN:NE2	3:D:661:HOH:O	2.51	0.42
1:A:71:ILE:O	1:A:75:LYS:HG3	2.19	0.42
1:D:296:LYS:O	1:D:300:GLU:HG3	2.20	0.42
1:A:244:THR:HG22	1:A:249:ARG:HH11	1.84	0.42
1:C:41:THR:HG21	1:C:306:ILE:CG2	2.50	0.42
1:B:23:ARG:NH2	1:B:23:ARG:HG2	2.33	0.42
1:B:60:GLU:OE1	1:B:60:GLU:HA	2.20	0.42
1:D:60:GLU:HA	1:D:60:GLU:OE1	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:251:ASN:ND2	3:C:685:HOH:O	2.53	0.41
1:C:304:LEU:HD23	1:C:304:LEU:O	2.19	0.41
1:C:71:ILE:O	1:C:75:LYS:HG3	2.20	0.41
1:A:160:GLN:NE2	3:A:747:HOH:O	2.52	0.41
1:B:231:LEU:HD23	1:B:231:LEU:C	2.41	0.41
1:B:263:LEU:O	1:B:267:LEU:HG	2.20	0.41
1:B:7:GLU:O	1:B:49:LEU:HA	2.20	0.41
1:C:242:LEU:HD23	3:C:666:HOH:O	2.19	0.41
1:B:307:ASN:O	1:B:310:LYS:HE2	2.21	0.41
1:C:201:PRO:HB3	1:D:37:LYS:HA	2.01	0.41
1:A:244:THR:HG21	1:A:288:ILE:HG12	2.01	0.41
1:B:145:ARG:CZ	1:B:160:GLN:HE22	2.33	0.41
1:B:210:GLU:HG3	3:B:702:HOH:O	2.21	0.41
1:D:206:ASP:O	1:D:210:GLU:HG3	2.20	0.41
1:C:109:LYS:HE2	3:C:668:HOH:O	2.20	0.41
1:D:239:ASP:HB3	1:D:242:LEU:HB2	2.01	0.41
1:A:22:ARG:HH21	1:A:23:ARG:CG	2.33	0.41
1:A:60:GLU:OE1	1:A:60:GLU:HA	2.21	0.41
1:C:304:LEU:HD23	1:C:304:LEU:C	2.41	0.41
1:A:179:LEU:HD23	1:A:200:ILE:HD13	2.02	0.40
1:C:145:ARG:CB	1:C:151:MET:HE2	2.46	0.40
1:D:221:GLU:O	1:D:224:LYS:HG2	2.21	0.40
1:A:1:MET:HA	1:A:82:GLU:HG2	2.02	0.40
1:B:252:ASN:HB2	1:D:199:GLN:HE22	1.87	0.40
1:A:199:GLN:HE22	1:C:252:ASN:HB2	1.87	0.40
1:A:37:LYS:HA	1:B:201:PRO:HB3	2.02	0.40
1:C:273:ARG:HD3	3:C:732:HOH:O	2.21	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	295/330 (89%)	281 (95%)	14 (5%)	0	100	100
1	B	299/330 (91%)	287 (96%)	12 (4%)	0	100	100
1	C	306/330 (93%)	285 (93%)	20 (6%)	1 (0%)	41	46
1	D	294/330 (89%)	280 (95%)	14 (5%)	0	100	100
All	All	1194/1320 (90%)	1133 (95%)	60 (5%)	1 (0%)	51	60

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	150	ARG

### 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	269/296 (91%)	259 (96%)	10 (4%)	34	43
1	B	273/296 (92%)	260 (95%)	13 (5%)	25	32
1	C	278/296 (94%)	261 (94%)	17 (6%)	18	21
1	D	267/296 (90%)	256 (96%)	11 (4%)	30	39
All	All	1087/1184 (92%)	1036 (95%)	51 (5%)	26	33

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	LEU
1	A	22	ARG
1	A	87	LEU
1	A	91	HIS
1	A	160	GLN
1	A	164	SER
1	A	271	LYS
1	A	272	ASP
1	A	292	ASN
1	A	302	LEU

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Mol	Chain	Res	Type
1	B	91	HIS
1	B	129	SER
1	B	130	LYS
1	B	142	ASN
1	B	143	ARG
1	B	157	LYS
1	B	160	GLN
1	B	169	LYS
1	B	172	LEU
1	B	223	LEU
1	B	271	LYS
1	B	291	ASN
1	B	303	ASP
1	C	31	THR
1	C	37	LYS
1	C	91	HIS
1	C	130	LYS
1	C	142	ASN
1	C	145	ARG
1	C	146	SER
1	C	148	PHE
1	C	150	ARG
1	C	151	MET
1	C	152	SER
1	C	160	GLN
1	C	172	LEU
1	C	223	LEU
1	C	246	THR
1	C	271	LYS
1	C	293	GLU
1	D	6	LEU
1	D	15	ARG
1	D	22	ARG
1	D	91	HIS
1	D	142	ASN
1	D	145	ARG
1	D	265	LEU
1	D	271	LYS
1	D	291	ASN
1	D	292	ASN
1	D	302	LEU

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

sidechains are listed below:

Mol	Chain	Res	Type
1	A	40	GLN
1	A	44	GLN
1	A	91	HIS
1	A	160	GLN
1	A	199	GLN
1	A	251	ASN
1	A	261	GLN
1	A	292	ASN
1	B	40	GLN
1	B	44	GLN
1	B	91	HIS
1	B	199	GLN
1	B	217	GLN
1	B	251	ASN
1	B	252	ASN
1	B	291	ASN
1	C	44	GLN
1	C	91	HIS
1	C	160	GLN
1	C	251	ASN
1	C	252	ASN
1	D	40	GLN
1	D	44	GLN
1	D	91	HIS
1	D	199	GLN
1	D	251	ASN
1	D	291	ASN
1	D	292	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry

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

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	301/330 (91%)	0.01	15 (4%) 28 27	15, 28, 69, 127	0
1	B	305/330 (92%)	0.19	24 (7%) 12 11	15, 31, 83, 118	1 (0%)
1	C	310/330 (93%)	0.13	18 (5%) 23 22	14, 28, 80, 130	1 (0%)
1	D	300/330 (90%)	-0.16	13 (4%) 35 33	14, 27, 70, 117	0
All	All	1216/1320 (92%)	0.04	70 (5%) 23 22	14, 28, 78, 130	2 (0%)

All (70) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	148	PHE	8.2
1	A	125	TYR	5.7
1	D	145	ARG	5.7
1	D	55	TYR	5.6
1	A	55	TYR	4.8
1	B	129	SER	4.7
1	C	290	HIS	4.7
1	B	290	HIS	4.3
1	C	37	LYS	4.2
1	C	36	ILE	4.1
1	C	129	SER	4.0
1	C	149	ASN	3.9
1	A	124	ALA	3.9
1	B	145	ARG	3.9
1	C	170	GLY	3.7
1	B	152	SER	3.6
1	A	168	GLU	3.6
1	C	39	MET	3.6
1	B	33	GLY	3.5
1	B	40	GLN	3.5
1	C	128	ASP	3.4

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Mol	Chain	Res	Type	RSRZ
1	B	157	LYS	3.4
1	B	150	ARG	3.4
1	D	146	SER	3.4
1	B	154	SER	3.3
1	C	150	ARG	3.3
1	C	311	ILE	3.2
1	B	35	VAL	3.1
1	C	57	LYS	3.1
1	B	170	GLY	3.1
1	B	57	LYS	3.0
1	B	153	GLU	3.0
1	B	151	MET	3.0
1	C	145	ARG	2.9
1	B	128	ASP	2.9
1	D	54	ARG	2.9
1	C	168	GLU	2.8
1	D	57	LYS	2.7
1	D	81	LYS	2.7
1	B	156	GLU	2.6
1	B	311	ILE	2.6
1	D	169	LYS	2.6
1	B	168	GLU	2.5
1	C	151	MET	2.5
1	C	35	VAL	2.5
1	A	54	ARG	2.5
1	D	56	ALA	2.4
1	D	170	GLY	2.4
1	A	154	SER	2.4
1	B	169	LYS	2.4
1	D	159	TRP	2.4
1	A	151	MET	2.4
1	A	290	HIS	2.4
1	A	150	ARG	2.4
1	A	53	PHE	2.3
1	A	159	TRP	2.3
1	B	81	LYS	2.3
1	D	168	GLU	2.3
1	B	34	ASP	2.3
1	D	157	LYS	2.3
1	B	56	ALA	2.3
1	B	146	SER	2.2
1	D	171	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	157	LYS	2.1
1	A	56	ALA	2.1
1	A	170	GLY	2.1
1	A	311	ILE	2.1
1	C	277	VAL	2.1
1	C	56	ALA	2.0
1	B	106	LYS	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. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	B-factors(Å <sup>2</sup> )	Q<0.9
2	MN	D	500	1/1	0.97	0.06	34,34,34,34	0
2	MN	B	500	1/1	0.97	0.07	40,40,40,40	0
2	MN	C	500	1/1	0.98	0.08	37,37,37,37	0
2	MN	A	500	1/1	0.98	0.05	32,32,32,32	0
2	MN	C	501	1/1	0.99	0.18	25,25,25,25	0
2	MN	B	501	1/1	0.99	0.16	26,26,26,26	0
2	MN	D	501	1/1	0.99	0.10	24,24,24,24	0
2	MN	A	501	1/1	1.00	0.13	28,28,28,28	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.