



# Full wwPDB X-ray Structure Validation Report ⓘ

May 29, 2020 – 09:35 am BST

PDB ID : 6BRG  
Title : The SAM domain of mouse SAMHD1 is critical for its activation and regulation  
Authors : Buzovetsky, O.; Tang, C.; Knecht, K.M.; Antonucci, J.M.; Wu, L.; Ji, X.; Xiong, Y.  
Deposited on : 2017-11-30  
Resolution : 3.50 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

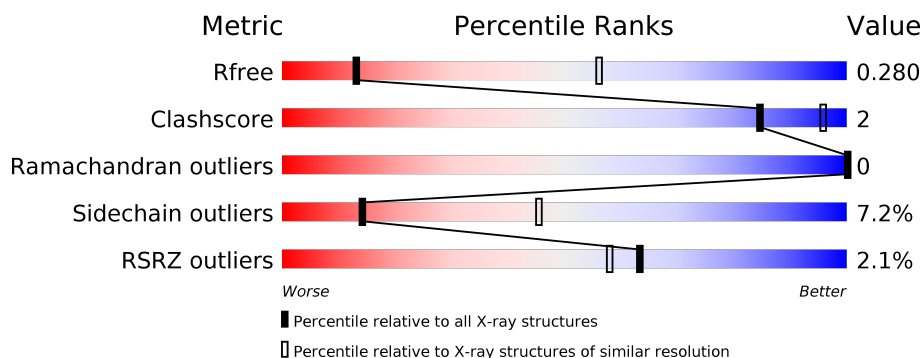
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.50 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	1659 (3.60-3.40)
Clashscore	141614	1036 (3.58-3.42)
Ramachandran outliers	138981	1005 (3.58-3.42)
Sidechain outliers	138945	1006 (3.58-3.42)
RSRZ outliers	127900	1559 (3.60-3.40)

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	672	<div> <div>2%</div> <div> <div></div> <div>64%</div> <div>10%</div> <div>26%</div> </div> </div>
1	B	672	<div> <div></div> <div> <div>66%</div> <div>8%</div> <div>26%</div> </div> </div>
1	C	672	<div> <div>2%</div> <div> <div></div> <div>66%</div> <div>8%</div> <div>26%</div> </div> </div>
1	D	672	<div> <div>2%</div> <div> <div></div> <div>65%</div> <div>9%</div> <div>26%</div> </div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 16571 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 Deoxynucleoside triphosphate triphosphohydrolase SAMHD1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	500	Total	C	N	O	S	0	0	0
			4130	2641	727	741	21			
1	B	500	Total	C	N	O	S	0	0	0
			4130	2641	727	741	21			
1	C	500	Total	C	N	O	S	0	0	0
			4130	2641	727	741	21			
1	D	500	Total	C	N	O	S	0	0	0
			4124	2638	724	741	21			

There are 56 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-13	MET	-	initiating methionine	UNP F8WJE0
A	-12	GLY	-	expression tag	UNP F8WJE0
A	-11	SER	-	expression tag	UNP F8WJE0
A	-10	SER	-	expression tag	UNP F8WJE0
A	-9	HIS	-	expression tag	UNP F8WJE0
A	-8	HIS	-	expression tag	UNP F8WJE0
A	-7	HIS	-	expression tag	UNP F8WJE0
A	-6	HIS	-	expression tag	UNP F8WJE0
A	-5	HIS	-	expression tag	UNP F8WJE0
A	-4	HIS	-	expression tag	UNP F8WJE0
A	-3	SER	-	expression tag	UNP F8WJE0
A	-2	SER	-	expression tag	UNP F8WJE0
A	-1	GLY	-	expression tag	UNP F8WJE0
A	0	LEU	-	expression tag	UNP F8WJE0
B	-13	MET	-	initiating methionine	UNP F8WJE0
B	-12	GLY	-	expression tag	UNP F8WJE0
B	-11	SER	-	expression tag	UNP F8WJE0
B	-10	SER	-	expression tag	UNP F8WJE0
B	-9	HIS	-	expression tag	UNP F8WJE0
B	-8	HIS	-	expression tag	UNP F8WJE0
B	-7	HIS	-	expression tag	UNP F8WJE0

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-6	HIS	-	expression tag	UNP F8WJE0
B	-5	HIS	-	expression tag	UNP F8WJE0
B	-4	HIS	-	expression tag	UNP F8WJE0
B	-3	SER	-	expression tag	UNP F8WJE0
B	-2	SER	-	expression tag	UNP F8WJE0
B	-1	GLY	-	expression tag	UNP F8WJE0
B	0	LEU	-	expression tag	UNP F8WJE0
C	-13	MET	-	initiating methionine	UNP F8WJE0
C	-12	GLY	-	expression tag	UNP F8WJE0
C	-11	SER	-	expression tag	UNP F8WJE0
C	-10	SER	-	expression tag	UNP F8WJE0
C	-9	HIS	-	expression tag	UNP F8WJE0
C	-8	HIS	-	expression tag	UNP F8WJE0
C	-7	HIS	-	expression tag	UNP F8WJE0
C	-6	HIS	-	expression tag	UNP F8WJE0
C	-5	HIS	-	expression tag	UNP F8WJE0
C	-4	HIS	-	expression tag	UNP F8WJE0
C	-3	SER	-	expression tag	UNP F8WJE0
C	-2	SER	-	expression tag	UNP F8WJE0
C	-1	GLY	-	expression tag	UNP F8WJE0
C	0	LEU	-	expression tag	UNP F8WJE0
D	-13	MET	-	initiating methionine	UNP F8WJE0
D	-12	GLY	-	expression tag	UNP F8WJE0
D	-11	SER	-	expression tag	UNP F8WJE0
D	-10	SER	-	expression tag	UNP F8WJE0
D	-9	HIS	-	expression tag	UNP F8WJE0
D	-8	HIS	-	expression tag	UNP F8WJE0
D	-7	HIS	-	expression tag	UNP F8WJE0
D	-6	HIS	-	expression tag	UNP F8WJE0
D	-5	HIS	-	expression tag	UNP F8WJE0
D	-4	HIS	-	expression tag	UNP F8WJE0
D	-3	SER	-	expression tag	UNP F8WJE0
D	-2	SER	-	expression tag	UNP F8WJE0
D	-1	GLY	-	expression tag	UNP F8WJE0
D	0	LEU	-	expression tag	UNP F8WJE0

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

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total Mg 1 1	0	0
2	A	1	Total Mg 1 1	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	D	1	Total 1	Mg 1	0	0
2	C	1	Total 1	Mg 1	0	0

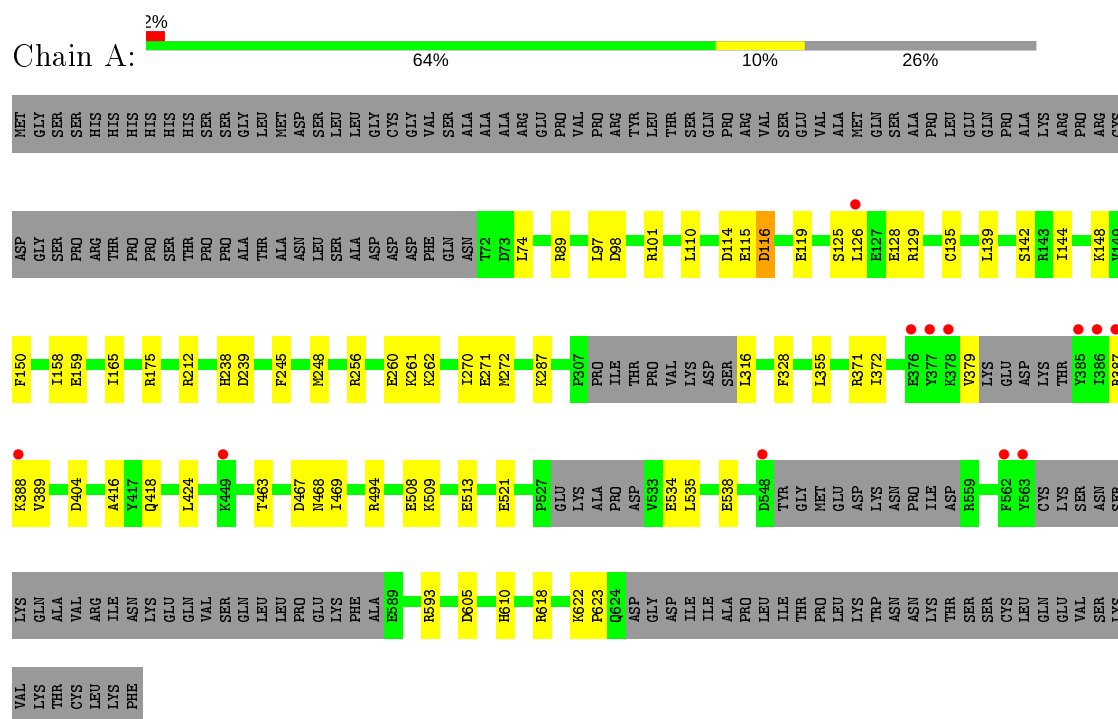
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	14	Total 14	O 14	0	0
3	B	14	Total 14	O 14	0	0
3	C	12	Total 12	O 12	0	0
3	D	13	Total 13	O 13	0	0

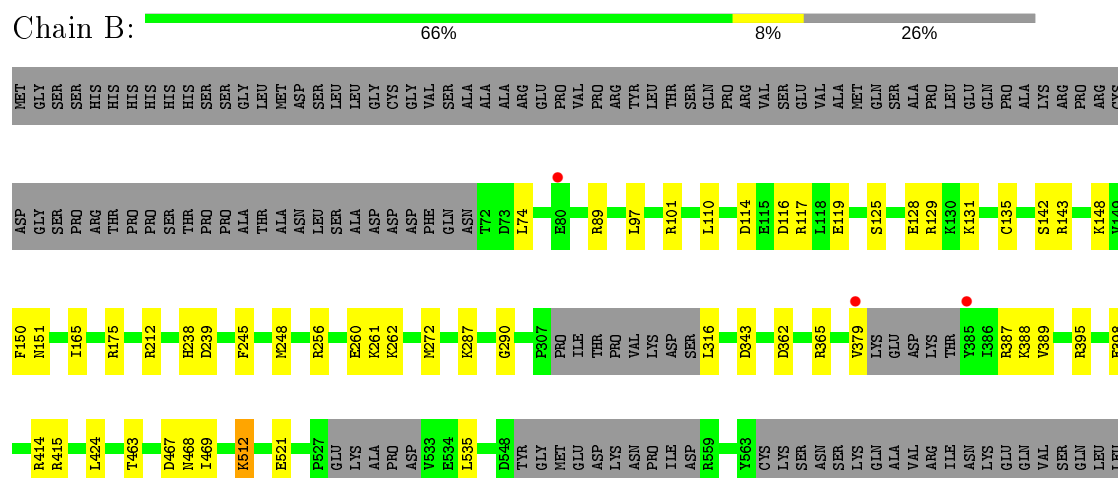
### 3 Residue-property plots [i](#)

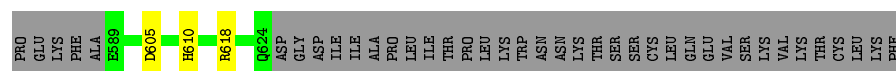
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are 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: Deoxynucleoside triphosphate triphosphohydrolase SAMHD1

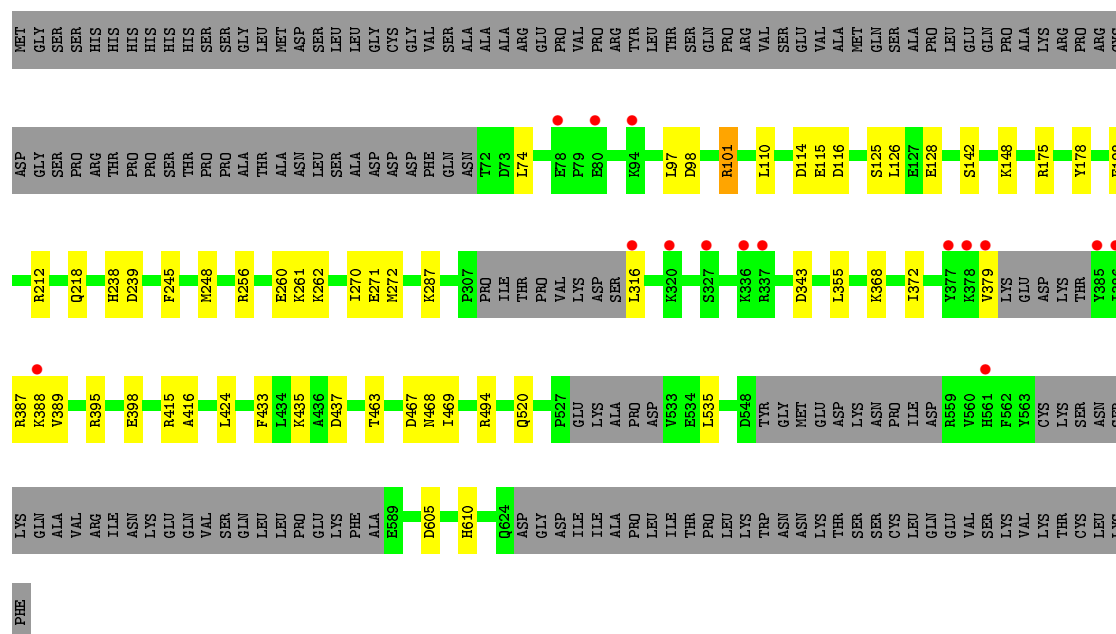


- Molecule 1: Deoxynucleoside triphosphate triphosphohydrolase SAMHD1

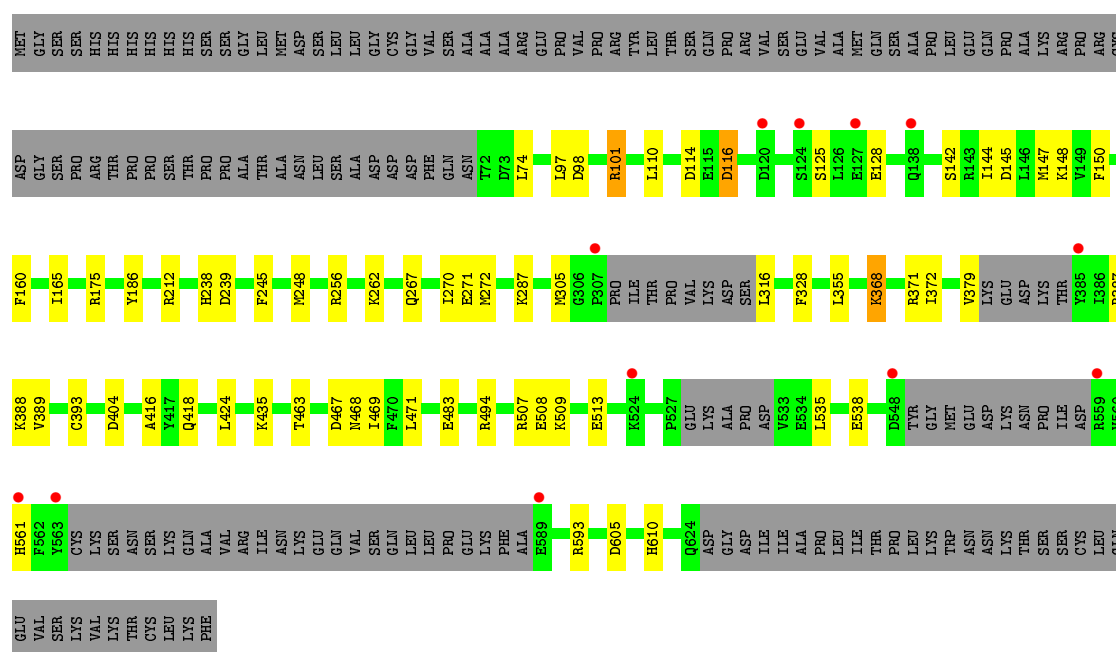




● Molecule 1: Deoxynucleoside triphosphate triphosphohydrolase SAMHD1



● Molecule 1: Deoxynucleoside triphosphate triphosphohydrolase SAMHD1



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	164.58Å 109.92Å 162.09Å 90.00° 105.78° 90.00°	Depositor
Resolution (Å)	20.00 – 3.50 19.98 – 3.50	Depositor EDS
% Data completeness (in resolution range)	98.0 (20.00-3.50) 98.6 (19.98-3.50)	Depositor EDS
$R_{merge}$	0.15	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.37 (at 3.52Å)	Xtriage
Refinement program	REFMAC 5.8.0103	Depositor
R, $R_{free}$	0.244 , 0.278 0.246 , 0.280	Depositor DCC
$R_{free}$ test set	1725 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	74.0	Xtriage
Anisotropy	0.104	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 85.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.89	EDS
Total number of atoms	16571	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	89.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.69% of the height of the origin peak. No significant pseudotranslation is detected.*

<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 [i](#)

### 5.1 Standard geometry [i](#)

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	0.48	0/4221	0.62	0/5674
1	B	0.49	0/4221	0.62	1/5674 (0.0%)
1	C	0.48	0/4221	0.60	0/5674
1	D	0.48	0/4215	0.59	0/5667
All	All	0.48	0/16878	0.61	1/22689 (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	414	ARG	NE-CZ-NH1	-5.67	117.47	120.30

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4130	0	4110	27	0
1	B	4130	0	4110	19	0
1	C	4130	0	4110	15	0
1	D	4124	0	4099	28	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	14	0	0	2	0
3	B	14	0	0	0	0
3	C	12	0	0	1	0
3	D	13	0	0	0	0
All	All	16571	0	16429	76	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 2.

All (76) 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)
3:A:909:HOH:O	1:B:150:PHE:HE2	1.71	0.72
1:B:119:GLU:OE2	1:B:129:ARG:NH1	2.30	0.65
1:B:117:ARG:HD2	1:B:290:GLY:HA2	1.80	0.63
1:A:139:LEU:HD22	3:A:902:HOH:O	1.99	0.61
1:B:148:LYS:HD2	1:B:165:ILE:HG12	1.87	0.56
1:B:260:GLU:HG2	1:B:261:LYS:N	2.21	0.55
1:A:116:ASP:OD1	1:A:116:ASP:N	2.40	0.55
1:A:371:ARG:HH12	1:D:144:ILE:CG2	2.20	0.54
1:A:371:ARG:HH12	1:D:144:ILE:HG22	1.72	0.53
1:C:260:GLU:HG2	1:C:261:LYS:N	2.21	0.53
1:A:125:SER:HB2	1:A:128:GLU:HB3	1.91	0.52
1:B:415:ARG:NH1	1:D:404:ASP:OD1	2.44	0.51
1:B:125:SER:HB2	1:B:128:GLU:HB3	1.93	0.50
1:C:245:PHE:N	1:C:467:ASP:OD1	2.44	0.50
1:A:159:GLU:HG3	1:D:368:LYS:HE2	1.92	0.49
1:D:144:ILE:HA	1:D:147:MET:HG2	1.94	0.49
3:C:912:HOH:O	1:D:150:PHE:HE2	1.97	0.48
1:D:148:LYS:HD2	1:D:165:ILE:HG12	1.96	0.48
1:C:125:SER:HB2	1:C:128:GLU:HB3	1.95	0.48
1:D:125:SER:HB2	1:D:128:GLU:HB3	1.94	0.47
1:A:89:ARG:HG2	1:A:135:CYS:SG	2.54	0.47
1:A:355:LEU:O	1:B:151:ASN:ND2	2.48	0.47
1:B:395:ARG:HB3	1:B:398:GLU:HG2	1.97	0.46
1:A:418:GLN:O	1:A:593:ARG:NH1	2.43	0.45
1:A:521:GLU:OE1	1:A:618:ARG:NH2	2.49	0.45
1:D:245:PHE:N	1:D:467:ASP:OD1	2.49	0.44
1:C:101:ARG:NE	1:C:101:ARG:O	2.50	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:395:ARG:HB3	1:C:398:GLU:HG2	2.00	0.44
1:D:267:GLN:HA	1:D:305:MET:HE1	2.00	0.44
1:D:328:PHE:CG	1:D:372:ILE:HG12	2.52	0.44
1:A:260:GLU:HG2	1:A:261:LYS:N	2.32	0.44
1:D:116:ASP:N	1:D:116:ASP:OD1	2.50	0.44
1:B:238:HIS:NE2	1:B:239:ASP:OD2	2.51	0.43
1:D:355:LEU:HD13	1:D:416:ALA:HB2	2.00	0.43
1:D:101:ARG:O	1:D:101:ARG:NE	2.51	0.43
1:D:160:PHE:HB2	1:D:165:ILE:HD11	1.99	0.43
1:D:270:ILE:HG13	1:D:271:GLU:N	2.33	0.43
1:A:371:ARG:NH1	1:D:144:ILE:HG21	2.33	0.43
1:D:393:CYS:SG	1:D:561:HIS:HB2	2.58	0.43
1:A:270:ILE:HG13	1:A:271:GLU:N	2.34	0.43
1:C:178:TYR:CD1	1:D:471:LEU:HD11	2.54	0.43
1:A:508:GLU:O	1:A:509:LYS:HG3	2.18	0.43
1:A:371:ARG:NH1	1:D:144:ILE:CG2	2.81	0.43
1:C:238:HIS:NE2	1:C:343:ASP:OD1	2.52	0.43
1:B:125:SER:HB2	1:B:128:GLU:CB	2.48	0.42
1:A:144:ILE:CG2	1:D:371:ARG:HH12	2.31	0.42
1:C:238:HIS:NE2	1:C:239:ASP:OD2	2.52	0.42
1:D:238:HIS:NE2	1:D:239:ASP:OD2	2.52	0.42
1:B:362:ASP:OD1	1:B:365:ARG:HG3	2.18	0.42
1:B:89:ARG:HG2	1:B:135:CYS:SG	2.59	0.42
1:D:418:GLN:O	1:D:593:ARG:NH1	2.40	0.42
1:A:404:ASP:OD1	1:C:415:ARG:NH1	2.52	0.42
1:A:144:ILE:HG22	1:D:371:ARG:HH12	1.85	0.42
1:A:148:LYS:HD2	1:A:165:ILE:HG12	2.02	0.42
1:B:238:HIS:NE2	1:B:343:ASP:OD1	2.53	0.42
1:B:245:PHE:N	1:B:467:ASP:OD1	2.51	0.42
1:B:521:GLU:OE1	1:B:618:ARG:NH2	2.52	0.42
1:A:238:HIS:NE2	1:A:239:ASP:OD2	2.53	0.42
1:A:119:GLU:OE2	1:A:129:ARG:NH1	2.53	0.41
1:A:245:PHE:N	1:A:467:ASP:OD1	2.49	0.41
1:C:218:GLN:NE2	1:C:372:ILE:O	2.49	0.41
1:D:435:LYS:HD3	1:D:483:GLU:HB3	2.01	0.41
1:D:508:GLU:O	1:D:509:LYS:HG3	2.20	0.41
1:B:260:GLU:CG	1:B:261:LYS:N	2.83	0.41
1:B:512:LYS:HG2	1:B:512:LYS:H	1.67	0.41
1:C:270:ILE:HG13	1:C:271:GLU:N	2.35	0.41
1:C:433:PHE:O	1:C:437:ASP:N	2.53	0.41
1:A:328:PHE:CG	1:A:372:ILE:HG12	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:117:ARG:CD	1:B:290:GLY:HA2	2.49	0.41
1:A:144:ILE:CG2	1:D:371:ARG:NH1	2.83	0.41
1:A:622:LYS:HA	1:A:623:PRO:HD3	1.99	0.41
1:C:355:LEU:HD13	1:C:416:ALA:HB2	2.02	0.41
1:A:355:LEU:HD13	1:A:416:ALA:HB2	2.02	0.41
1:A:150:PHE:HB2	1:A:158:ILE:HB	2.02	0.41
1:C:260:GLU:CG	1:C:261:LYS:N	2.84	0.41
1:C:198:GLU:OE2	1:D:186:TYR:OH	2.35	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	488/672 (73%)	474 (97%)	14 (3%)	0	100	100
1	B	488/672 (73%)	473 (97%)	15 (3%)	0	100	100
1	C	488/672 (73%)	474 (97%)	14 (3%)	0	100	100
1	D	488/672 (73%)	473 (97%)	15 (3%)	0	100	100
All	All	1952/2688 (73%)	1894 (97%)	58 (3%)	0	100	100

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	449/602 (75%)	416 (93%)	33 (7%)	14	45
1	B	449/602 (75%)	420 (94%)	29 (6%)	17	50
1	C	449/602 (75%)	415 (92%)	34 (8%)	13	43
1	D	448/602 (74%)	415 (93%)	33 (7%)	13	44
All	All	1795/2408 (74%)	1666 (93%)	129 (7%)	14	45

All (129) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	74	LEU
1	A	97	LEU
1	A	98	ASP
1	A	101	ARG
1	A	110	LEU
1	A	114	ASP
1	A	115	GLU
1	A	116	ASP
1	A	126	LEU
1	A	142	SER
1	A	175	ARG
1	A	212	ARG
1	A	248	MET
1	A	256	ARG
1	A	262	LYS
1	A	272	MET
1	A	287	LYS
1	A	316	LEU
1	A	379	VAL
1	A	387	ARG
1	A	388	LYS
1	A	389	VAL
1	A	424	LEU
1	A	463	THR
1	A	468	ASN
1	A	469	ILE
1	A	494	ARG
1	A	513	GLU
1	A	534	GLU
1	A	535	LEU
1	A	538	GLU
1	A	605	ASP
1	A	610	HIS

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Mol	Chain	Res	Type
1	B	74	LEU
1	B	97	LEU
1	B	101	ARG
1	B	110	LEU
1	B	114	ASP
1	B	116	ASP
1	B	131	LYS
1	B	142	SER
1	B	143	ARG
1	B	175	ARG
1	B	212	ARG
1	B	248	MET
1	B	256	ARG
1	B	262	LYS
1	B	272	MET
1	B	287	LYS
1	B	316	LEU
1	B	379	VAL
1	B	387	ARG
1	B	388	LYS
1	B	389	VAL
1	B	424	LEU
1	B	463	THR
1	B	468	ASN
1	B	469	ILE
1	B	512	LYS
1	B	535	LEU
1	B	605	ASP
1	B	610	HIS
1	C	74	LEU
1	C	97	LEU
1	C	98	ASP
1	C	101	ARG
1	C	110	LEU
1	C	114	ASP
1	C	115	GLU
1	C	116	ASP
1	C	126	LEU
1	C	142	SER
1	C	148	LYS
1	C	175	ARG
1	C	212	ARG

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Mol	Chain	Res	Type
1	C	248	MET
1	C	256	ARG
1	C	262	LYS
1	C	272	MET
1	C	287	LYS
1	C	316	LEU
1	C	368	LYS
1	C	379	VAL
1	C	387	ARG
1	C	388	LYS
1	C	389	VAL
1	C	424	LEU
1	C	435	LYS
1	C	463	THR
1	C	468	ASN
1	C	469	ILE
1	C	494	ARG
1	C	520	GLN
1	C	535	LEU
1	C	605	ASP
1	C	610	HIS
1	D	74	LEU
1	D	97	LEU
1	D	98	ASP
1	D	101	ARG
1	D	110	LEU
1	D	114	ASP
1	D	116	ASP
1	D	142	SER
1	D	145	ASP
1	D	175	ARG
1	D	212	ARG
1	D	248	MET
1	D	256	ARG
1	D	262	LYS
1	D	272	MET
1	D	287	LYS
1	D	316	LEU
1	D	368	LYS
1	D	379	VAL
1	D	387	ARG
1	D	388	LYS

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Mol	Chain	Res	Type
1	D	389	VAL
1	D	424	LEU
1	D	463	THR
1	D	468	ASN
1	D	469	ILE
1	D	494	ARG
1	D	507	ARG
1	D	513	GLU
1	D	535	LEU
1	D	538	GLU
1	D	605	ASP
1	D	610	HIS

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

### 5.3.3 RNA ⓘ

There are no RNA 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 4 ligands modelled in this entry, 4 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	500/672 (74%)	-0.20	12 (2%) 59 53	37, 82, 150, 224	0
1	B	500/672 (74%)	-0.20	3 (0%) 89 86	40, 83, 140, 215	0
1	C	500/672 (74%)	-0.12	15 (3%) 50 44	38, 86, 151, 270	0
1	D	500/672 (74%)	-0.24	12 (2%) 59 53	39, 78, 145, 194	0
All	All	2000/2688 (74%)	-0.19	42 (2%) 63 58	37, 82, 148, 270	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	379	VAL	7.7
1	C	385	TYR	4.5
1	C	378	LYS	4.3
1	B	385	TYR	3.8
1	C	377	TYR	3.6
1	D	559	ARG	3.5
1	C	388	LYS	3.4
1	A	377	TYR	3.2
1	A	563	TYR	3.2
1	D	563	TYR	3.2
1	B	80	GLU	3.1
1	A	388	LYS	3.0
1	A	387	ARG	3.0
1	C	386	ILE	2.9
1	D	589	GLU	2.7
1	A	376	GLU	2.7
1	A	386	ILE	2.7
1	D	548	ASP	2.7
1	D	138	GLN	2.6
1	D	307	PRO	2.6
1	C	80	GLU	2.5

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Mol	Chain	Res	Type	RSRZ
1	D	385	TYR	2.4
1	D	524	LYS	2.4
1	C	561	HIS	2.3
1	C	316	LEU	2.3
1	C	336	LYS	2.3
1	C	337	ARG	2.3
1	C	78	GLU	2.2
1	C	320	LYS	2.2
1	D	120	ASP	2.2
1	D	127	GLU	2.2
1	C	327	SER	2.2
1	A	548	ASP	2.2
1	A	385	TYR	2.1
1	A	449	LYS	2.1
1	D	561	HIS	2.1
1	B	379	VAL	2.1
1	C	94	LYS	2.1
1	D	124	SER	2.1
1	A	378	LYS	2.1
1	A	562	PHE	2.0
1	A	126	LEU	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( $\text{\AA}^2$ )	Q<0.9
2	MG	D	800	1/1	0.87	0.41	106,106,106,106	0
2	MG	B	800	1/1	0.93	0.31	126,126,126,126	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	MG	A	800	1/1	0.97	0.10	21,21,21,21	0
2	MG	C	800	1/1	0.97	0.15	34,34,34,34	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.