



# Full wwPDB X-ray Structure Validation Report ⓘ

Nov 24, 2025 – 02:23 PM JST

PDB ID : 9KU0 / pdb\_00009ku0  
Title : Crystal structure of CtpA F105R mutant from Helicobacter pylori  
Authors : Sun, K.; Au, S.W.N.  
Deposited on : 2024-12-03  
Resolution : 3.70 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.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.46

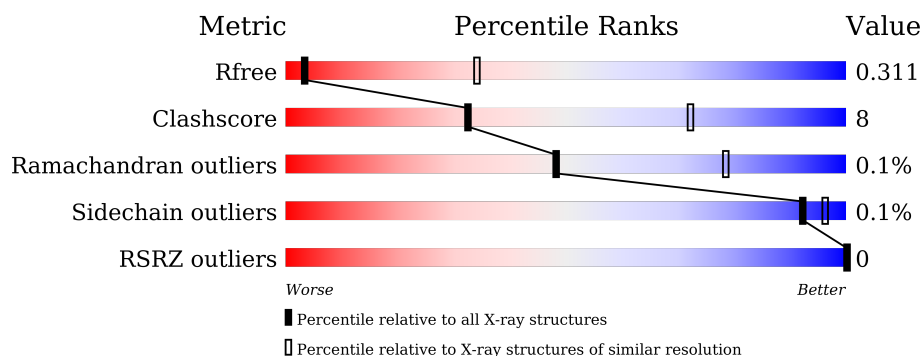
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.70 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))
$R_{free}$	164625	1017 (3.80-3.60)
Clashscore	180529	1074 (3.80-3.60)
Ramachandran outliers	177936	1055 (3.80-3.60)
Sidechain outliers	177891	1052 (3.80-3.60)
RSRZ outliers	164620	1017 (3.80-3.60)

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 of 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	416	<div> <div>68%</div> <div>18%</div> <div>14%</div> </div>
1	B	416	<div> <div>66%</div> <div>20%</div> <div>13%</div> </div>
1	C	416	<div> <div>72%</div> <div>14%</div> <div>13%</div> </div>
1	D	416	<div> <div>69%</div> <div>16%</div> <div>14%</div> </div>

## 2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 11012 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 Carboxyl-terminal protease.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	356	Total	C	N	O	S	0	0	0
			2733	1744	466	516	7			
1	B	360	Total	C	N	O	S	0	0	0
			2769	1770	474	518	7			
1	C	361	Total	C	N	O	S	0	0	0
			2774	1768	473	526	7			
1	D	356	Total	C	N	O	S	0	0	0
			2736	1745	468	516	7			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	39	MET	-	initiating methionine	UNP B5Z8Z7
A	40	HIS	-	expression tag	UNP B5Z8Z7
A	41	HIS	-	expression tag	UNP B5Z8Z7
A	42	HIS	-	expression tag	UNP B5Z8Z7
A	43	HIS	-	expression tag	UNP B5Z8Z7
A	44	HIS	-	expression tag	UNP B5Z8Z7
A	45	HIS	-	expression tag	UNP B5Z8Z7
A	46	MET	-	expression tag	UNP B5Z8Z7
A	105	ARG	PHE	engineered mutation	UNP B5Z8Z7
B	39	MET	-	initiating methionine	UNP B5Z8Z7
B	40	HIS	-	expression tag	UNP B5Z8Z7
B	41	HIS	-	expression tag	UNP B5Z8Z7
B	42	HIS	-	expression tag	UNP B5Z8Z7
B	43	HIS	-	expression tag	UNP B5Z8Z7
B	44	HIS	-	expression tag	UNP B5Z8Z7
B	45	HIS	-	expression tag	UNP B5Z8Z7
B	46	MET	-	expression tag	UNP B5Z8Z7
B	105	ARG	PHE	engineered mutation	UNP B5Z8Z7
C	39	MET	-	initiating methionine	UNP B5Z8Z7
C	40	HIS	-	expression tag	UNP B5Z8Z7
C	41	HIS	-	expression tag	UNP B5Z8Z7

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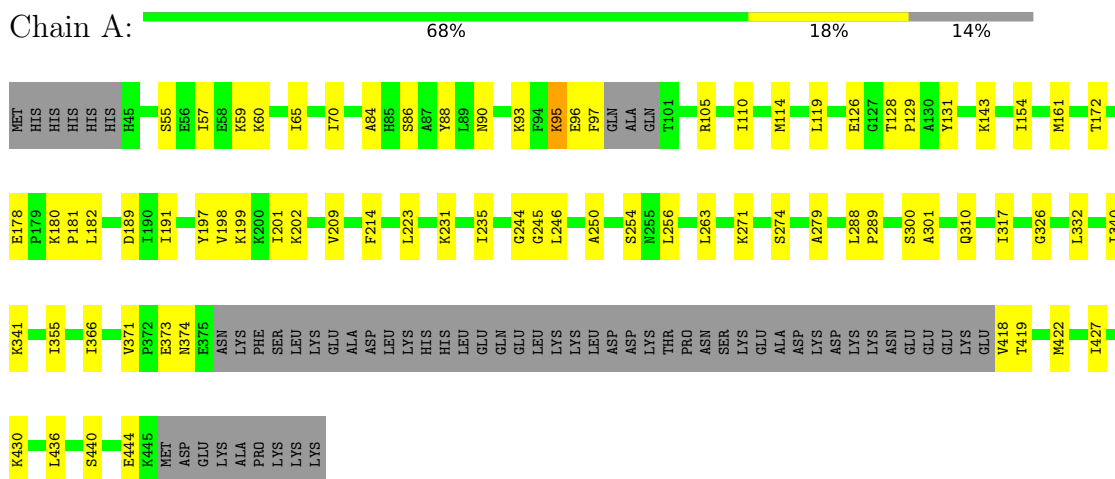
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Chain	Residue	Modelled	Actual	Comment	Reference
C	42	HIS	-	expression tag	UNP B5Z8Z7
C	43	HIS	-	expression tag	UNP B5Z8Z7
C	44	HIS	-	expression tag	UNP B5Z8Z7
C	45	HIS	-	expression tag	UNP B5Z8Z7
C	46	MET	-	expression tag	UNP B5Z8Z7
C	105	ARG	PHE	engineered mutation	UNP B5Z8Z7
D	39	MET	-	initiating methionine	UNP B5Z8Z7
D	40	HIS	-	expression tag	UNP B5Z8Z7
D	41	HIS	-	expression tag	UNP B5Z8Z7
D	42	HIS	-	expression tag	UNP B5Z8Z7
D	43	HIS	-	expression tag	UNP B5Z8Z7
D	44	HIS	-	expression tag	UNP B5Z8Z7
D	45	HIS	-	expression tag	UNP B5Z8Z7
D	46	MET	-	expression tag	UNP B5Z8Z7
D	105	ARG	PHE	engineered mutation	UNP B5Z8Z7

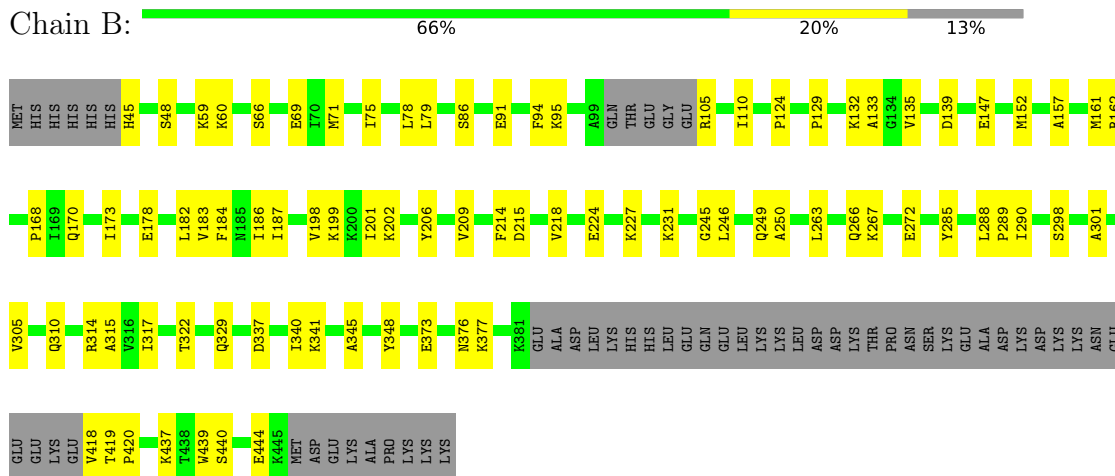
### 3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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: Carboxyl-terminal protease



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## 4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	139.33Å 139.33Å 252.09Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	28.01 – 3.70 28.01 – 3.70	Depositor EDS
% Data completeness (in resolution range)	100.0 (28.01-3.70) 99.7 (28.01-3.70)	Depositor EDS
$R_{merge}$	0.13	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.64 (at 3.75Å)	Xtriage
Refinement program	PHENIX (dev_5526: ???)	Depositor
R, $R_{free}$	0.242 , 0.312 0.242 , 0.311	Depositor DCC
$R_{free}$ test set	1030 reflections (5.28%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	123.6	Xtriage
Anisotropy	0.285	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.28 , 89.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.45$ , $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	0.065 for h,-h-k,-l	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	11012	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	142.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 15.87% 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](#)

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	0.23	0/2772	0.50	0/3737
1	B	0.28	1/2809 (0.0%)	0.48	0/3785
1	C	0.22	0/2813	0.51	0/3791
1	D	0.21	0/2775	0.50	0/3741
All	All	0.24	1/11169 (0.0%)	0.49	0/15054

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	263	LEU	C-N	9.75	1.39	1.33

There are no bond angle outliers.

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	2733	0	2875	46	0
1	B	2769	0	2923	48	0
1	C	2774	0	2913	39	0
1	D	2736	0	2878	43	0
All	All	11012	0	11589	176	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 8.



All (176) 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:93:LYS:HA	1:A:96:GLU:HG3	1.57	0.85
1:C:59:LYS:HG3	1:C:60:LYS:HG3	1.66	0.77
1:B:59:LYS:HG3	1:B:60:LYS:HG3	1.66	0.76
1:C:199:LYS:HD2	1:C:417:GLU:HA	1.65	0.76
1:B:91:GLU:HA	1:B:337:ASP:O	1.87	0.74
1:A:214:PHE:HB2	1:A:245:GLY:HA3	1.70	0.74
1:D:125:LEU:HD21	1:D:193:LEU:HD13	1.72	0.70
1:D:334:VAL:H	1:D:339:ALA:HA	1.57	0.69
1:C:178:GLU:HG3	1:C:182:LEU:HD11	1.75	0.69
1:B:86:SER:HA	1:B:341:LYS:O	1.93	0.69
1:B:373:GLU:HB3	1:B:376:ASN:HB3	1.75	0.68
1:D:65:ILE:HD12	1:D:70:ILE:HD11	1.75	0.68
1:D:202:LYS:HB2	1:D:419:THR:HB	1.75	0.68
1:A:366:ILE:HD11	1:A:427:ILE:HG12	1.76	0.67
1:A:263:LEU:HD12	1:A:279:ALA:HB2	1.76	0.67
1:C:86:SER:HA	1:C:341:LYS:O	1.96	0.65
1:A:59:LYS:HG3	1:A:60:LYS:HG3	1.78	0.65
1:D:362:PRO:HG2	1:D:365:VAL:HG22	1.78	0.65
1:B:214:PHE:HB2	1:B:245:GLY:HA3	1.78	0.64
1:A:65:ILE:HD12	1:A:70:ILE:HD11	1.80	0.64
1:B:178:GLU:HG3	1:B:182:LEU:HD11	1.80	0.64
1:D:193:LEU:HD11	1:D:212:SER:HB2	1.81	0.63
1:B:214:PHE:HA	1:B:218:VAL:HG21	1.80	0.62
1:B:285:TYR:HB3	1:B:288:LEU:HD22	1.80	0.62
1:A:373:GLU:HG2	1:A:374:ASN:H	1.65	0.62
1:D:285:TYR:HB3	1:D:288:LEU:HD22	1.83	0.61
1:D:141:ILE:HG22	1:D:149:THR:HG21	1.81	0.61
1:C:86:SER:HB2	1:C:342:ILE:HG22	1.83	0.61
1:A:244:GLY:HA3	1:A:300:SER:HB3	1.82	0.60
1:B:184:PHE:HB3	1:B:186:ILE:HD11	1.83	0.60
1:B:199:LYS:HD2	1:B:418:VAL:HG23	1.84	0.60
1:A:57:ILE:HG22	1:A:65:ILE:HD11	1.84	0.59
1:C:289:PRO:HB3	1:C:439:TRP:CD2	2.36	0.59
1:A:90:ASN:HB2	1:A:93:LYS:HG2	1.84	0.59
1:C:149:THR:HA	1:C:152:MET:HE3	1.85	0.59
1:B:198:VAL:HG12	1:B:209:VAL:HG22	1.85	0.59
1:C:106:GLY:O	1:C:162:ARG:HG3	2.03	0.59
1:D:59:LYS:HG3	1:D:60:LYS:HG3	1.84	0.58
1:D:216:LYS:HG2	1:D:217:ASN:H	1.68	0.57
1:B:94:PHE:O	1:B:95:LYS:HG2	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:112:VAL:HG23	1:D:154:ILE:HG23	1.84	0.57
1:B:310:GLN:HB2	1:B:317:ILE:HG13	1.86	0.57
1:C:214:PHE:HB2	1:C:245:GLY:HA3	1.87	0.56
1:D:214:PHE:HB2	1:D:245:GLY:HA3	1.87	0.56
1:C:333:PRO:HA	1:C:339:ALA:HA	1.88	0.55
1:A:129:PRO:HG3	1:A:189:ASP:HB2	1.89	0.55
1:B:105:ARG:HH22	1:B:162:ARG:HH11	1.55	0.54
1:A:114:MET:HB2	1:A:154:ILE:HD11	1.88	0.54
1:A:178:GLU:HG3	1:A:182:LEU:HD11	1.88	0.53
1:A:371:VAL:HG21	1:A:430:LYS:HD2	1.90	0.53
1:D:244:GLY:HA3	1:D:300:SER:HB3	1.89	0.53
1:C:112:VAL:HG23	1:C:154:ILE:HG23	1.90	0.53
1:C:285:TYR:HB3	1:C:288:LEU:HD22	1.90	0.53
1:B:75:ILE:HA	1:B:78:LEU:HG	1.90	0.53
1:B:152:MET:HE3	1:B:157:ALA:HB2	1.91	0.53
1:A:143:LYS:HB2	1:A:172:THR:HB	1.89	0.52
1:A:88:TYR:CZ	1:A:90:ASN:HA	2.44	0.52
1:A:288:LEU:HD12	1:A:289:PRO:HD2	1.92	0.52
1:B:71:MET:O	1:B:75:ILE:HG22	2.10	0.52
1:B:224:GLU:HA	1:B:227:LYS:HG2	1.90	0.51
1:B:214:PHE:O	1:B:246:LEU:HD23	2.10	0.51
1:D:89:LEU:HD22	1:D:93:LYS:HB3	1.92	0.51
1:D:329:GLN:HB3	1:D:341:LYS:HD3	1.92	0.51
1:A:202:LYS:HE3	1:A:418:VAL:HG22	1.94	0.50
1:A:326:GLY:HA3	1:A:355:ILE:HB	1.92	0.50
1:C:54:VAL:HG13	1:C:70:ILE:HG21	1.94	0.50
1:D:254:SER:HB2	1:D:263:LEU:HD11	1.93	0.50
1:A:201:ILE:HA	1:A:419:THR:OG1	2.12	0.50
1:D:105:ARG:NH2	1:D:162:ARG:HD2	2.27	0.50
1:C:84:ALA:HB1	1:C:114:MET:HG3	1.93	0.50
1:A:198:VAL:HG12	1:A:209:VAL:HG22	1.92	0.50
1:B:202:LYS:HD2	1:B:420:PRO:HG3	1.94	0.50
1:A:86:SER:HA	1:A:341:LYS:O	2.13	0.49
1:A:90:ASN:H	1:A:93:LYS:HB2	1.78	0.49
1:D:105:ARG:NE	1:D:105:ARG:HA	2.28	0.49
1:A:84:ALA:HB1	1:A:114:MET:HB3	1.95	0.48
1:B:289:PRO:HB3	1:B:439:TRP:CD2	2.48	0.48
1:D:298:SER:HB2	1:D:322:THR:HG22	1.95	0.48
1:A:332:LEU:HD12	1:A:340:ILE:HD11	1.95	0.48
1:D:165:PRO:HB3	1:D:189:ASP:HA	1.95	0.48
1:A:110:ILE:HD11	1:A:161:MET:SD	2.54	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:91:GLU:O	1:C:92:LYS:HB3	2.12	0.48
1:D:119:LEU:HD11	1:D:152:MET:O	2.14	0.48
1:B:201:ILE:HA	1:B:419:THR:OG1	2.14	0.48
1:B:206:TYR:OH	1:B:437:LYS:HA	2.13	0.48
1:C:79:LEU:HD12	1:C:88:TYR:HB2	1.96	0.47
1:B:139:ASP:HB3	1:B:173:ILE:HG23	1.95	0.47
1:B:440:SER:O	1:B:444:GLU:HG2	2.14	0.47
1:B:79:LEU:HD11	1:B:340:ILE:HG22	1.95	0.47
1:C:285:TYR:HB3	1:C:288:LEU:CD2	2.45	0.47
1:B:105:ARG:NH2	1:B:162:ARG:HD2	2.30	0.47
1:D:105:ARG:HH22	1:D:162:ARG:HD2	1.78	0.47
1:D:193:LEU:HG	1:D:195:SER:HA	1.97	0.47
1:B:215:ASP:O	1:B:249:GLN:HG3	2.15	0.47
1:C:330:MET:HE2	1:C:344:THR:HG21	1.96	0.47
1:B:301:ALA:O	1:B:305:VAL:HG12	2.15	0.47
1:C:88:TYR:HD1	1:C:340:ILE:HG22	1.79	0.46
1:A:96:GLU:O	1:A:97:PHE:C	2.57	0.46
1:B:133:ALA:HB2	1:B:186:ILE:HG12	1.97	0.46
1:C:201:ILE:HA	1:C:419:THR:HG21	1.97	0.46
1:D:334:VAL:CG2	1:D:340:ILE:HG23	2.45	0.46
1:A:254:SER:HB2	1:A:263:LEU:HD21	1.96	0.46
1:D:89:LEU:HB2	1:D:339:ALA:O	2.16	0.46
1:D:270:ASN:ND2	1:D:273:GLU:HB2	2.30	0.46
1:D:88:TYR:HD1	1:D:340:ILE:HG22	1.80	0.46
1:D:216:LYS:HG2	1:D:217:ASN:N	2.30	0.46
1:A:129:PRO:HD3	1:A:189:ASP:HB2	1.98	0.46
1:B:147:GLU:OE2	1:B:152:MET:HB2	2.15	0.45
1:A:419:THR:HA	1:A:422:MET:HE3	1.97	0.45
1:D:206:TYR:OH	1:D:437:LYS:HA	2.15	0.45
1:A:223:LEU:HD13	1:A:256:LEU:HD11	1.97	0.45
1:C:289:PRO:HB3	1:C:439:TRP:CE3	2.51	0.45
1:B:168:PRO:HA	1:B:187:ILE:HA	1.99	0.45
1:D:55:SER:O	1:D:59:LYS:HG2	2.16	0.45
1:C:419:THR:H	1:C:420:PRO:CD	2.30	0.45
1:C:298:SER:HB2	1:C:322:THR:HG22	1.98	0.45
1:A:231:LYS:HB2	1:A:231:LYS:HE2	1.81	0.44
1:B:129:PRO:HA	1:B:132:LYS:HD2	1.99	0.44
1:C:79:LEU:CD1	1:C:88:TYR:HB2	2.47	0.44
1:D:334:VAL:HG21	1:D:340:ILE:HG23	1.98	0.44
1:B:267:LYS:HB2	1:B:348:TYR:CE1	2.53	0.44
1:B:170:GLN:NE2	1:B:183:VAL:HG13	2.33	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:250:ALA:HB2	1:C:301:ALA:HB1	1.99	0.44
1:C:329:GLN:HA	1:C:343:THR:HA	1.98	0.44
1:A:55:SER:O	1:A:59:LYS:HG2	2.18	0.43
1:A:197:TYR:HE1	1:A:199:LYS:HE2	1.82	0.43
1:D:442:VAL:O	1:D:445:LYS:HE3	2.18	0.43
1:D:120:THR:HG22	1:D:140:ASN:OD1	2.18	0.43
1:A:180:LYS:HG2	1:A:181:PRO:HD2	2.01	0.43
1:A:271:LYS:HA	1:A:274:SER:OG	2.19	0.43
1:A:440:SER:O	1:A:444:GLU:HG2	2.19	0.43
1:B:250:ALA:HB2	1:B:301:ALA:HB1	1.99	0.43
1:A:105:ARG:HD3	1:A:105:ARG:H	1.83	0.43
1:C:90:ASN:H	1:C:93:LYS:HB2	1.82	0.43
1:B:290:ILE:HG23	1:B:315:ALA:HB2	2.01	0.43
1:A:310:GLN:HB2	1:A:317:ILE:HG13	2.01	0.42
1:A:246:LEU:H	1:A:246:LEU:HD23	1.84	0.42
1:C:141:ILE:HG22	1:C:149:THR:HG21	2.01	0.42
1:C:71:MET:O	1:C:75:ILE:HG22	2.20	0.42
1:D:72:THR:HA	1:D:75:ILE:HG22	2.02	0.42
1:C:205:PRO:HB2	1:C:233:LYS:HG2	2.01	0.42
1:C:295:ASN:O	1:C:322:THR:HA	2.19	0.42
1:C:419:THR:H	1:C:420:PRO:HD2	1.84	0.42
1:B:66:SER:HB2	1:B:69:GLU:OE2	2.20	0.42
1:A:235:ILE:O	1:A:436:LEU:HD11	2.20	0.42
1:C:190:ILE:HG13	1:C:191:ILE:N	2.35	0.42
1:C:164:LYS:O	1:C:188:ARG:HB2	2.19	0.42
1:D:275:LEU:HD23	1:D:276:GLU:N	2.35	0.42
1:D:371:VAL:HG21	1:D:430:LYS:HD3	2.02	0.41
1:B:298:SER:HB2	1:B:322:THR:HG22	2.02	0.41
1:D:269:LYS:HD2	1:D:269:LYS:HA	1.75	0.41
1:D:271:LYS:HD2	1:D:271:LYS:HA	1.74	0.41
1:D:355:ILE:HG22	1:D:355:ILE:O	2.21	0.41
1:A:119:LEU:CD1	1:A:154:ILE:HG13	2.50	0.41
1:B:288:LEU:HD23	1:B:314:ARG:HH22	1.85	0.41
1:C:166:LYS:HA	1:C:166:LYS:HD3	1.50	0.41
1:C:190:ILE:HG13	1:C:191:ILE:H	1.85	0.41
1:B:124:PRO:HG3	1:B:135:VAL:O	2.21	0.41
1:B:272:GLU:H	1:B:272:GLU:HG3	1.66	0.41
1:D:92:LYS:O	1:D:96:GLU:HB2	2.21	0.41
1:D:235:ILE:HD11	1:D:288:LEU:HD11	2.02	0.41
1:D:258:ILE:HD13	1:D:312:HIS:ND1	2.35	0.41
1:C:421:LYS:HB3	1:C:421:LYS:HE3	1.76	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:377:LYS:HD3	1:B:377:LYS:HA	1.90	0.41
1:C:445:LYS:HA	1:C:445:LYS:HD3	1.89	0.41
1:A:95:LYS:HZ2	1:A:95:LYS:HG2	1.72	0.41
1:B:110:ILE:HD11	1:B:161:MET:SD	2.61	0.41
1:B:329:GLN:HB3	1:B:341:LYS:HG3	2.03	0.41
1:A:128:THR:HG22	1:A:191:ILE:HG22	2.03	0.41
1:D:289:PRO:HG3	1:D:439:TRP:CD1	2.56	0.41
1:C:61:TYR:CE2	1:C:63:ASP:HB2	2.56	0.40
1:B:266:GLN:HB3	1:B:345:ALA:HB1	2.04	0.40
1:A:250:ALA:HB2	1:A:301:ALA:HB1	2.04	0.40
1:B:45:HIS:HB2	1:B:48:SER:HB3	2.03	0.40
1:B:231:LYS:HD3	1:B:231:LYS:HA	1.97	0.40
1:A:126:GLU:HA	1:A:131:TYR:CD2	2.57	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	350/416 (84%)	330 (94%)	20 (6%)	0	100	100
1	B	354/416 (85%)	334 (94%)	20 (6%)	0	100	100
1	C	355/416 (85%)	338 (95%)	15 (4%)	2 (1%)	22	54
1	D	350/416 (84%)	324 (93%)	26 (7%)	0	100	100
All	All	1409/1664 (85%)	1326 (94%)	81 (6%)	2 (0%)	48	78

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	419	THR
1	C	217	ASN

### 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	304/360 (84%)	303 (100%)	1 (0%)	91	94
1	B	308/360 (86%)	308 (100%)	0	100	100
1	C	308/360 (86%)	308 (100%)	0	100	100
1	D	304/360 (84%)	304 (100%)	0	100	100
All	All	1224/1440 (85%)	1223 (100%)	1 (0%)	92	96

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	95	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	81	ASN
1	A	266	GLN
1	A	424	ASN
1	B	81	ASN
1	B	85	HIS
1	B	255	ASN
1	B	287	ASN
1	C	249	GLN
1	C	424	ASN
1	D	98	GLN
1	D	170	GLN
1	D	255	ASN
1	D	280	ASN
1	D	287	ASN

### 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 oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

There are no ligands in this entry.

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	356/416 (85%)	-0.38	0 100 100	88, 141, 196, 232	0
1	B	360/416 (86%)	-0.38	0 100 100	84, 143, 205, 242	0
1	C	361/416 (86%)	-0.44	0 100 100	98, 139, 190, 219	0
1	D	356/416 (85%)	-0.42	0 100 100	84, 141, 189, 212	0
All	All	1433/1664 (86%)	-0.41	0 100 100	84, 141, 198, 242	0

There are no RSRZ outliers to report.

### 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 oligosaccharides in this entry.

### 6.4 Ligands [i](#)

There are no ligands in this entry.

### 6.5 Other polymers [i](#)

There are no such residues in this entry.