



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 06:44 pm GMT

PDB ID : 1UK4  
Title : Crystal structure of SARS Coronavirus Main Proteinase (3CLpro) Complexed With An Inhibitor  
Authors : Yang, H.; Yang, M.; Liu, Y.; Bartlam, M.; Ding, Y.; Lou, Z.; Sun, L.; Zhou, Z.; Ye, S.; Anand, K.; Pang, H.; Gao, G.F.; Hilgenfeld, R.; Rao, Z.  
Deposited on : 2003-08-14  
Resolution : 2.50 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

<http://wwpdb.org/validation/2016/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)	:	1.9-1692
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

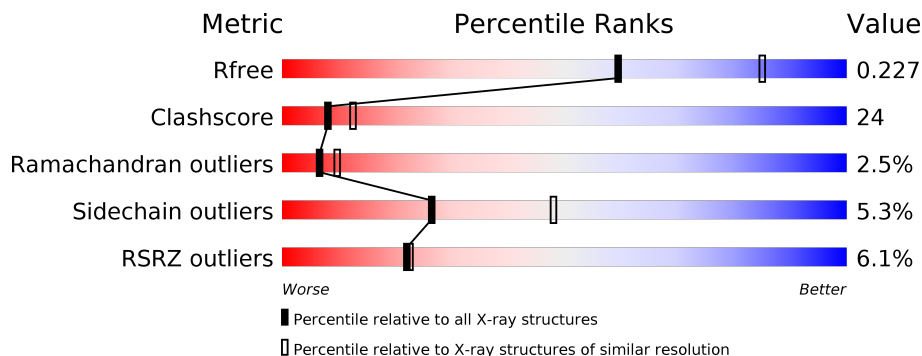
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.50 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))
$R_{free}$	100719	3846 (2.50-2.50)
Clashscore	112137	4554 (2.50-2.50)
Ramachandran outliers	110173	4463 (2.50-2.50)
Sidechain outliers	110143	4465 (2.50-2.50)
RSRZ outliers	101464	3876 (2.50-2.50)

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. 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	306	<div> <div>4%</div> <div>62%</div> <div>33%</div> <div>• •</div> </div>
1	B	306	<div> <div>5%</div> <div>64%</div> <div>31%</div> <div>• •</div> </div>
2	G	5	<div> <div>100%</div> <div>20%</div> <div>60%</div> <div>20%</div> </div>
2	H	5	<div> <div>40%</div> <div>80%</div> <div>20%</div> </div>
2	K	5	<div> <div>60%</div> <div>20%</div> <div>80%</div> </div>

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	ATO	A	1006	-	-	X	-

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 4875 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.

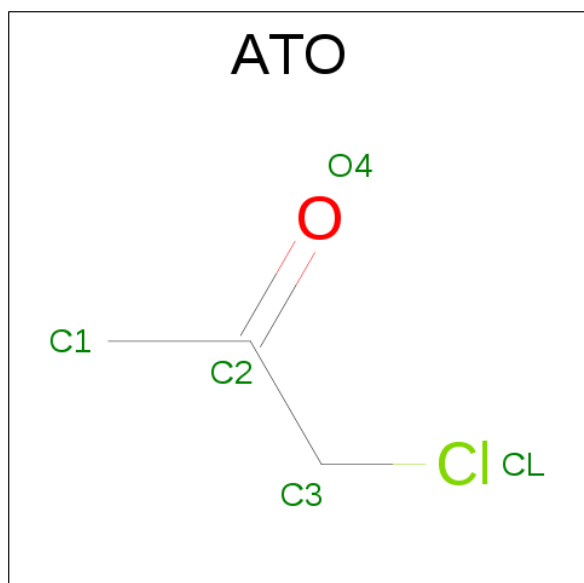
- Molecule 1 is a protein called 3C-like proteinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	301	Total	C	N	O	S	0	0	0
			2333	1476	399	436	22			
1	B	302	Total	C	N	O	S	0	0	0
			2336	1476	400	438	22			

- Molecule 2 is a protein called 5-mer peptide of inhibitor.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	G	5	Total	C	N	O	0	0	0
			39	22	7	10			
2	H	5	Total	C	N	O	0	0	0
			39	22	7	10			
2	K	5	Total	C	N	O	0	0	0
			39	22	7	10			

- Molecule 3 is CHLOROACETONE (three-letter code: ATO) (formula: C<sub>3</sub>H<sub>5</sub>ClO).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C 1 1	0	0
3	B	1	Total C 1 1	0	0

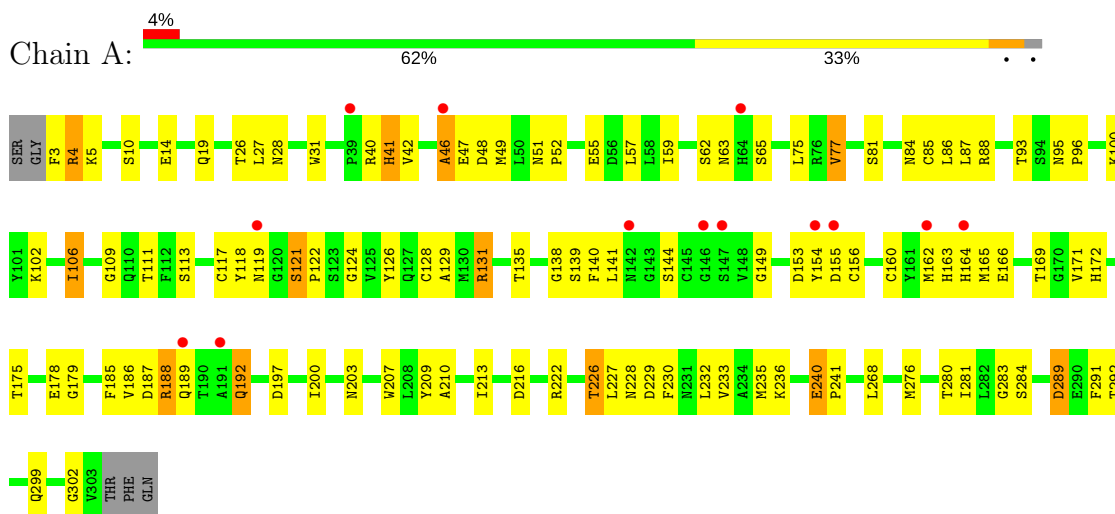
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	43	Total O 43 43	0	0
4	B	42	Total O 42 42	0	0
4	G	1	Total O 1 1	0	0
4	K	1	Total O 1 1	0	0

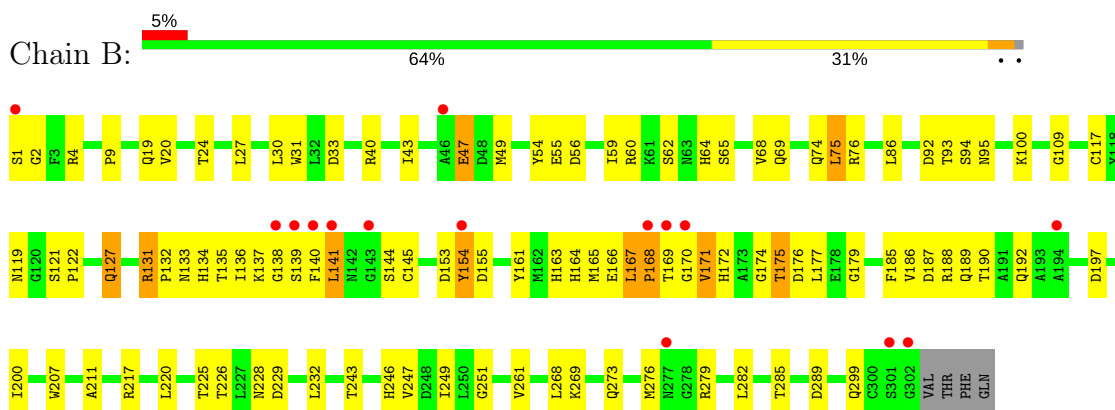
### 3 Residue-property plots

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

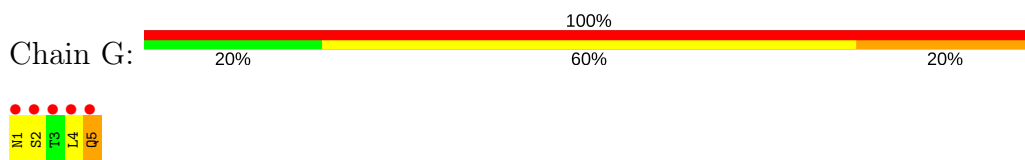
#### • Molecule 1: 3C-like proteinase



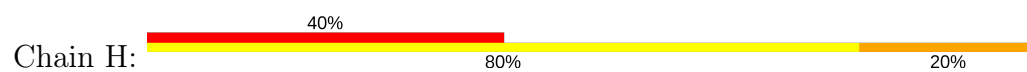
#### • Molecule 1: 3C-like proteinase



#### • Molecule 2: 5-mer peptide of inhibitor



#### • Molecule 2: 5-mer peptide of inhibitor



- Molecule 2: 5-mer peptide of inhibitor



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	52.63Å 97.73Å 67.93Å 90.00° 102.79° 90.00°	Depositor
Resolution (Å)	25.00 – 2.50 25.07 – 2.21	Depositor EDS
% Data completeness (in resolution range)	(Not available) (25.00-2.50) 88.2 (25.07-2.21)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.43 (at 2.22Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, $R_{free}$	0.213 , 0.231 0.210 , 0.227	Depositor DCC
$R_{free}$ test set	1039 reflections (5.03%)	DCC
Wilson B-factor (Å <sup>2</sup> )	29.1	Xtriage
Anisotropy	0.558	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 46.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	4875	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.73% 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: ATO

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	0.39	0/2385	0.65	0/3241
1	B	0.38	0/2388	0.69	0/3244
2	G	0.56	0/38	1.63	1/49 (2.0%)
2	H	0.69	0/38	0.94	0/49
2	K	0.67	0/38	0.74	0/49
All	All	0.39	0/4887	0.68	1/6632 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	5	GLN	CA-C-O	9.14	139.29	120.10

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	2333	0	2286	106	0
1	B	2336	0	2288	113	0
2	G	39	0	39	16	0
2	H	39	0	39	21	0
2	K	39	0	39	6	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	1	0	0	4	0
3	B	1	0	0	2	0
4	A	43	0	0	1	0
4	B	42	0	0	0	0
4	G	1	0	0	1	0
4	K	1	0	0	0	0
All	All	4875	0	4691	221	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 24.

All (221) 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:226:THR:HG22	1:A:229:ASP:H	1.19	1.03
1:A:144:SER:N	2:G:5:GLN:O	1.92	1.01
1:B:189:GLN:HE21	2:H:4:LEU:HD12	1.26	0.99
3:A:1006:ATO:C1	2:G:5:GLN:HB3	1.92	0.97
1:B:168:PRO:HD3	2:H:2:SER:HA	1.50	0.94
2:K:1:ASN:CG	2:K:2:SER:H	1.68	0.92
1:A:169:THR:HB	1:A:171:VAL:HG23	1.54	0.89
1:A:171:VAL:HG12	1:A:172:HIS:H	1.38	0.87
1:A:299:GLN:HB3	1:B:140:PHE:HE2	1.40	0.86
2:H:1:ASN:HD22	2:H:1:ASN:H1	1.22	0.86
3:A:1006:ATO:C1	2:G:5:GLN:CB	2.53	0.85
1:B:168:PRO:CD	2:H:2:SER:HA	2.07	0.85
3:A:1006:ATO:C1	2:G:5:GLN:CA	2.55	0.84
1:B:145:CYS:CB	2:H:5:GLN:OXT	2.25	0.84
1:B:145:CYS:HB2	2:H:5:GLN:OXT	1.80	0.81
2:H:1:ASN:HD22	2:H:1:ASN:N	1.75	0.81
1:A:19:GLN:NE2	1:A:119:ASN:HB3	1.96	0.81
1:A:291:PHE:HE2	1:A:299:GLN:HE22	1.26	0.79
1:A:3:PHE:HA	1:B:140:PHE:HE1	1.49	0.78
1:B:186:VAL:H	1:B:192:GLN:HE22	1.32	0.76
1:B:167:LEU:HD21	1:B:185:PHE:CE1	2.21	0.75
1:A:141:LEU:HD11	1:B:299:GLN:O	1.85	0.75
1:B:189:GLN:HE22	2:H:3:THR:HA	1.53	0.73
1:B:131:ARG:HB2	1:B:135:THR:O	1.87	0.73
1:B:145:CYS:SG	2:H:5:GLN:OXT	2.46	0.73
1:B:76:ARG:HB3	1:B:92:ASP:OD2	1.88	0.73
1:B:69:GLN:HG2	1:B:74:GLN:HG2	1.71	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:1:ASN:ND2	2:H:1:ASN:N	2.38	0.72
3:B:2006:ATO:C1	2:H:5:GLN:CA	2.67	0.72
1:A:10:SER:O	1:A:14:GLU:HG3	1.90	0.71
1:A:118:TYR:CE1	1:A:144:SER:HB3	2.25	0.71
1:B:225:THR:HG22	1:B:226:THR:O	1.91	0.71
1:A:232:LEU:O	1:A:236:LYS:HE2	1.91	0.71
1:B:243:THR:H	1:B:246:HIS:HD2	1.39	0.70
1:A:3:PHE:HA	1:B:140:PHE:CE1	2.27	0.70
1:A:131:ARG:HB2	1:A:135:THR:O	1.92	0.69
2:K:1:ASN:CG	2:K:2:SER:N	2.43	0.69
1:B:49:MET:HA	1:B:49:MET:HE3	1.73	0.69
1:B:86:LEU:HG	1:B:179:GLY:HA2	1.75	0.68
1:A:226:THR:HG22	1:A:229:ASP:N	2.01	0.68
1:A:226:THR:CG2	1:A:229:ASP:H	2.03	0.67
1:B:56:ASP:O	1:B:59:ILE:HG22	1.94	0.67
1:B:170:GLY:O	1:B:171:VAL:HG22	1.95	0.66
1:B:127:GLN:HA	1:B:127:GLN:HE21	1.60	0.66
1:B:269:LYS:O	1:B:273:GLN:HG3	1.94	0.66
1:A:126:TYR:CD1	1:B:4:ARG:HG2	2.31	0.66
1:B:165:MET:HG2	2:H:4:LEU:HD23	1.78	0.66
1:B:163:HIS:HE1	1:B:172:HIS:HB3	1.61	0.66
1:B:243:THR:H	1:B:246:HIS:CD2	2.14	0.65
1:A:100:LYS:HD2	1:A:155:ASP:OD2	1.97	0.65
1:A:171:VAL:HG12	1:A:172:HIS:N	2.11	0.65
1:A:189:GLN:OE1	2:G:1:ASN:HB3	1.96	0.65
1:A:166:GLU:HB2	2:G:5:GLN:OE1	1.96	0.64
1:A:139:SER:HA	1:B:1:SER:O	1.97	0.64
1:A:19:GLN:HE21	1:A:119:ASN:HB3	1.59	0.64
1:B:19:GLN:NE2	1:B:119:ASN:HB3	2.13	0.64
1:A:299:GLN:HB3	1:B:140:PHE:CE2	2.28	0.63
1:B:140:PHE:O	1:B:141:LEU:HG	1.98	0.63
1:A:135:THR:OG1	1:A:171:VAL:HG11	2.00	0.61
1:A:189:GLN:HA	1:A:189:GLN:OE1	1.99	0.61
1:A:117:CYS:O	1:A:144:SER:HA	2.00	0.61
1:B:161:TYR:CE1	1:B:174:GLY:HA3	2.35	0.61
1:B:175:THR:CG2	1:B:176:ASP:O	2.49	0.61
1:B:247:VAL:HG13	1:B:261:VAL:HG11	1.82	0.60
1:A:291:PHE:HE2	1:A:299:GLN:NE2	1.98	0.60
1:B:168:PRO:HG2	2:H:1:ASN:C	2.21	0.60
1:A:276:MET:HE3	1:A:281:ILE:HG13	1.83	0.59
3:B:2006:ATO:C1	2:H:5:GLN:N	2.66	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:41:HIS:HB2	1:A:49:MET:HE3	1.85	0.58
1:A:189:GLN:NE2	2:G:1:ASN:O	2.36	0.58
1:B:47:GLU:CD	1:B:47:GLU:H	2.05	0.58
1:B:141:LEU:HD13	1:B:163:HIS:ND1	2.19	0.58
1:B:164:HIS:CD2	1:B:175:THR:HB	2.39	0.58
1:B:131:ARG:HD3	1:B:132:PRO:HD2	1.86	0.57
1:B:68:VAL:HG23	1:B:75:LEU:HB2	1.86	0.57
1:B:64:HIS:NE2	2:K:4:LEU:HG	2.19	0.57
1:A:186:VAL:H	1:A:192:GLN:HE22	1.52	0.57
1:B:33:ASP:O	1:B:94:SER:HA	2.05	0.57
1:A:111:THR:HG22	1:A:129:ALA:HB2	1.86	0.57
1:A:240:GLU:OE2	1:A:241:PRO:HD2	2.05	0.56
1:A:140:PHE:HB3	1:A:144:SER:OG	2.05	0.56
1:B:131:ARG:HH22	1:B:289:ASP:CG	2.09	0.56
1:B:109:GLY:HA2	1:B:200:ILE:HD13	1.88	0.56
1:B:49:MET:HA	1:B:49:MET:CE	2.35	0.56
1:B:49:MET:SD	2:H:4:LEU:HD13	2.45	0.56
1:A:222:ARG:HG2	1:A:222:ARG:O	2.05	0.56
1:B:175:THR:HG22	1:B:176:ASP:O	2.06	0.55
1:B:211:ALA:HA	1:B:282:LEU:HD13	1.88	0.55
1:B:189:GLN:NE2	2:H:3:THR:HA	2.21	0.55
1:A:49:MET:C	1:A:52:PRO:HD3	2.27	0.55
1:A:140:PHE:HD1	1:A:172:HIS:CD2	2.24	0.55
1:A:113:SER:O	1:A:149:GLY:HA2	2.07	0.55
1:A:5:LYS:HG2	1:A:291:PHE:CE1	2.42	0.55
1:A:63:ASN:HB3	1:A:77:VAL:HG22	1.89	0.55
1:A:226:THR:HG23	1:A:228:ASN:H	1.72	0.55
1:B:31:TRP:CE2	1:B:95:ASN:HB2	2.42	0.54
1:A:109:GLY:HA2	1:A:200:ILE:HD13	1.90	0.54
1:B:163:HIS:CE1	1:B:172:HIS:HB3	2.41	0.54
1:B:165:MET:HG2	2:H:4:LEU:CD2	2.38	0.54
1:B:217:ARG:HG2	1:B:220:LEU:CD1	2.38	0.54
1:B:141:LEU:HD13	1:B:163:HIS:CE1	2.43	0.54
1:A:163:HIS:HE1	1:A:172:HIS:HB3	1.72	0.54
1:B:207:TRP:NE1	1:B:282:LEU:HD23	2.24	0.53
1:B:145:CYS:N	2:H:5:GLN:OXT	2.37	0.53
1:A:5:LYS:HG2	1:A:291:PHE:CZ	2.43	0.53
1:B:40:ARG:O	1:B:43:ILE:HG12	2.09	0.53
1:B:49:MET:HB3	1:B:189:GLN:CG	2.39	0.53
1:A:163:HIS:CE1	1:A:172:HIS:HB3	2.43	0.53
2:K:1:ASN:OD1	2:K:2:SER:N	2.42	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:64:HIS:CD2	2:K:4:LEU:HG	2.44	0.52
1:A:31:TRP:CE2	1:A:75:LEU:HD11	2.45	0.52
2:G:5:GLN:OXT	2:G:5:GLN:HG2	2.10	0.52
1:A:102:LYS:HG3	1:A:156:CYS:SG	2.51	0.51
1:B:225:THR:CG2	1:B:226:THR:N	2.74	0.51
1:A:31:TRP:CZ2	1:A:75:LEU:HD11	2.46	0.50
1:A:3:PHE:CA	1:B:140:PHE:HE1	2.21	0.50
1:A:55:GLU:O	1:A:59:ILE:HG12	2.12	0.50
1:B:188:ARG:NE	1:B:190:THR:HG21	2.26	0.50
1:A:85:CYS:HB2	1:A:179:GLY:O	2.11	0.50
1:B:131:ARG:HD3	1:B:132:PRO:CD	2.42	0.50
1:B:166:GLU:O	1:B:167:LEU:C	2.50	0.50
1:B:131:ARG:NH2	1:B:289:ASP:OD2	2.45	0.50
1:A:111:THR:HG22	1:A:129:ALA:CB	2.42	0.50
1:A:166:GLU:HB2	2:G:5:GLN:CD	2.32	0.50
1:A:48:ASP:HB3	1:A:52:PRO:HB3	1.94	0.49
1:B:225:THR:HG21	1:B:269:LYS:NZ	2.26	0.49
1:A:144:SER:CA	2:G:5:GLN:O	2.59	0.49
1:A:86:LEU:HG	1:A:179:GLY:HA2	1.93	0.49
1:A:166:GLU:HG3	2:G:5:GLN:HE22	1.78	0.49
1:A:138:GLY:O	1:B:2:GLY:HA3	2.13	0.49
1:A:135:THR:CB	1:A:171:VAL:HG11	2.42	0.49
1:A:131:ARG:NH2	1:A:289:ASP:OD2	2.46	0.49
1:A:46:ALA:O	1:A:49:MET:HB2	2.13	0.49
1:B:138:GLY:O	1:B:140:PHE:N	2.45	0.49
1:B:217:ARG:HG2	1:B:220:LEU:HD11	1.94	0.49
1:B:249:ILE:C	1:B:251:GLY:H	2.16	0.49
1:B:92:ASP:OD1	1:B:93:THR:N	2.46	0.49
1:A:75:LEU:HD21	1:A:93:THR:HB	1.95	0.48
1:B:24:THR:HG22	1:B:24:THR:O	2.13	0.48
2:G:2:SER:HA	4:G:51:HOH:O	2.13	0.48
1:A:81:SER:O	1:A:87:LEU:HD12	2.13	0.48
1:B:49:MET:HB3	1:B:189:GLN:HG2	1.94	0.48
1:A:4:ARG:NH1	1:B:137:LYS:O	2.46	0.48
1:A:19:GLN:HE21	1:A:26:THR:HG21	1.79	0.48
1:B:188:ARG:HG2	1:B:190:THR:HG23	1.96	0.48
1:A:153:ASP:C	1:A:154:TYR:CD1	2.87	0.48
1:A:280:THR:HB	1:A:283:GLY:O	2.14	0.48
1:A:49:MET:O	1:A:52:PRO:HD3	2.14	0.48
1:A:222:ARG:HH11	1:A:222:ARG:HG2	1.79	0.47
1:A:135:THR:HB	1:A:171:VAL:CG1	2.44	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:228:ASN:HA	4:A:1026:HOH:O	2.15	0.47
1:A:207:TRP:O	1:A:210:ALA:HB3	2.15	0.47
1:B:47:GLU:N	1:B:47:GLU:OE2	2.45	0.47
1:A:124:GLY:HA2	1:B:9:PRO:HD3	1.96	0.47
1:B:68:VAL:CG2	1:B:75:LEU:HB2	2.44	0.47
1:A:154:TYR:HB2	1:A:155:ASP:H	1.47	0.47
1:B:138:GLY:C	1:B:140:PHE:N	2.68	0.47
1:B:131:ARG:HD2	1:B:197:ASP:OD1	2.14	0.47
1:B:20:VAL:HG22	1:B:68:VAL:HG12	1.98	0.46
1:A:203:ASN:OD1	1:A:292:THR:HA	2.15	0.46
1:B:153:ASP:O	1:B:154:TYR:HB2	2.16	0.46
1:B:225:THR:HG22	1:B:226:THR:N	2.31	0.46
1:A:169:THR:CB	1:A:171:VAL:HG23	2.37	0.46
1:A:88:ARG:NH1	1:A:88:ARG:HB3	2.30	0.46
1:B:127:GLN:NE2	1:B:127:GLN:HA	2.30	0.46
1:B:100:LYS:HD2	1:B:155:ASP:OD2	2.16	0.46
1:A:165:MET:HE3	2:G:1:ASN:OD1	2.16	0.46
1:A:139:SER:OG	1:B:299:GLN:NE2	2.45	0.45
1:A:40:ARG:C	1:A:42:VAL:H	2.19	0.45
1:B:175:THR:HG23	1:B:176:ASP:O	2.17	0.45
1:A:131:ARG:HH22	1:A:289:ASP:CG	2.19	0.45
1:A:88:ARG:CB	1:A:88:ARG:HH11	2.30	0.45
1:A:106:ILE:HD13	1:A:160:CYS:CB	2.46	0.45
1:B:56:ASP:OD1	1:B:60:ARG:NH1	2.50	0.44
1:A:86:LEU:HD21	1:A:162:MET:SD	2.58	0.44
1:B:31:TRP:CD2	1:B:95:ASN:HB2	2.53	0.44
1:B:185:PHE:HA	1:B:192:GLN:NE2	2.33	0.43
1:B:186:VAL:H	1:B:192:GLN:NE2	2.10	0.43
1:A:111:THR:HB	1:A:128:CYS:O	2.18	0.43
1:A:291:PHE:CE2	1:A:299:GLN:NE2	2.83	0.43
1:B:117:CYS:SG	1:B:122:PRO:HA	2.59	0.43
1:B:138:GLY:O	1:B:172:HIS:CE1	2.72	0.43
1:B:54:TYR:OH	1:B:187:ASP:HB3	2.19	0.43
1:A:48:ASP:O	1:A:52:PRO:HD3	2.18	0.43
1:A:62:SER:O	1:A:65:SER:HB2	2.19	0.42
1:B:225:THR:HG23	1:B:229:ASP:HB2	2.00	0.42
1:B:229:ASP:HA	1:B:232:LEU:HD12	2.01	0.42
1:B:145:CYS:SG	2:H:5:GLN:CA	3.07	0.42
1:A:230:PHE:CZ	1:A:268:LEU:HD23	2.54	0.42
1:B:225:THR:HG21	1:B:269:LYS:HZ1	1.84	0.42
1:A:166:GLU:HB3	2:G:2:SER:OG	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:141:LEU:HA	1:B:144:SER:OG	2.19	0.42
3:A:1006:ATO:C1	2:G:5:GLN:OXT	2.65	0.42
1:A:55:GLU:CD	1:A:55:GLU:H	2.23	0.42
1:A:95:ASN:HA	1:A:96:PRO:HD3	1.90	0.42
1:B:86:LEU:HG	1:B:179:GLY:CA	2.44	0.42
1:A:28:ASN:HD21	1:A:144:SER:HA	1.85	0.42
1:B:138:GLY:C	1:B:140:PHE:H	2.22	0.42
1:A:209:TYR:O	1:A:213:ILE:HG13	2.19	0.42
1:A:84:ASN:ND2	1:A:178:GLU:O	2.45	0.41
1:A:121:SER:HA	1:A:122:PRO:HD3	1.86	0.41
1:B:189:GLN:HE22	2:H:4:LEU:H	1.66	0.41
1:A:153:ASP:O	1:A:154:TYR:CD1	2.74	0.41
1:A:165:MET:HE1	1:A:188:ARG:O	2.20	0.41
1:A:233:VAL:O	1:A:236:LYS:HG2	2.21	0.41
1:B:133:ASN:O	1:B:134:HIS:HB2	2.20	0.41
1:B:164:HIS:NE2	1:B:175:THR:HB	2.35	0.41
1:B:121:SER:HA	1:B:122:PRO:HD3	1.87	0.41
1:B:217:ARG:HG2	1:B:220:LEU:HD12	2.03	0.41
1:B:189:GLN:NE2	2:H:4:LEU:HD12	2.11	0.41
1:A:131:ARG:HD2	1:A:197:ASP:OD1	2.21	0.41
1:A:169:THR:HB	1:A:171:VAL:CG2	2.39	0.41
1:A:165:MET:SD	1:A:187:ASP:HA	2.61	0.41
1:B:93:THR:HG22	1:B:94:SER:N	2.37	0.41
1:B:64:HIS:HD2	2:K:5:GLN:O	2.04	0.40
1:A:166:GLU:CG	2:G:5:GLN:HE22	2.34	0.40
1:A:164:HIS:CD2	1:A:175:THR:HG23	2.56	0.40
1:A:230:PHE:HZ	1:A:268:LEU:HD23	1.87	0.40
1:B:276:MET:O	1:B:279:ARG:HG3	2.21	0.40
1:B:62:SER:O	1:B:65:SER:HB2	2.22	0.40
1:B:228:ASN:O	1:B:232:LEU:HG	2.20	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	299/306 (98%)	271 (91%)	22 (7%)	6 (2%)	9	14
1	B	300/306 (98%)	270 (90%)	22 (7%)	8 (3%)	6	9
2	G	3/5 (60%)	2 (67%)	0	1 (33%)	0	0
2	H	3/5 (60%)	1 (33%)	2 (67%)	0	100	100
2	K	3/5 (60%)	2 (67%)	1 (33%)	0	100	100
All	All	608/627 (97%)	546 (90%)	47 (8%)	15 (2%)	6	10

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	192	GLN
1	B	169	THR
1	B	171	VAL
1	A	185	PHE
1	A	302	GLY
2	G	4	LEU
1	A	46	ALA
1	B	139	SER
1	A	41	HIS
1	B	141	LEU
1	B	154	TYR
1	B	167	LEU
1	B	177	LEU
1	A	284	SER
1	B	168	PRO

### 5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	259/263 (98%)	243 (94%)	16 (6%)	21	39
1	B	259/263 (98%)	248 (96%)	11 (4%)	34	59

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	G	5/5 (100%)	5 (100%)	0	100	100
2	H	5/5 (100%)	4 (80%)	1 (20%)	1	2
2	K	5/5 (100%)	5 (100%)	0	100	100
All	All	533/541 (98%)	505 (95%)	28 (5%)	26	48

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	ARG
1	A	27	LEU
1	A	47	GLU
1	A	51	ASN
1	A	57	LEU
1	A	77	VAL
1	A	106	ILE
1	A	121	SER
1	A	131	ARG
1	A	188	ARG
1	A	216	ASP
1	A	226	THR
1	A	227	LEU
1	A	235	MET
1	A	240	GLU
1	A	289	ASP
1	B	27	LEU
1	B	30	LEU
1	B	47	GLU
1	B	55	GLU
1	B	75	LEU
1	B	127	GLN
1	B	131	ARG
1	B	136	ILE
1	B	175	THR
1	B	268	LEU
1	B	285	THR
2	H	1	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	19	GLN
1	A	41	HIS
1	A	51	ASN
1	A	127	GLN
1	A	192	GLN
1	A	246	HIS
1	A	277	ASN
1	B	19	GLN
1	B	127	GLN
1	B	172	HIS
1	B	189	GLN
1	B	192	GLN
1	B	246	HIS
1	B	256	GLN
1	B	274	ASN
1	B	299	GLN
2	H	1	ASN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

### 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

### 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

### 5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 2 are modelled with single atom - 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](#)

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	301/306 (98%)	-0.00	13 (4%) 36 38	15, 34, 72, 95	0
1	B	302/306 (98%)	-0.03	15 (4%) 30 31	17, 33, 65, 106	0
2	G	5/5 (100%)	4.25	5 (100%) 0 0	99, 102, 106, 119	0
2	H	5/5 (100%)	2.34	2 (40%) 0 0	78, 80, 92, 99	0
2	K	5/5 (100%)	2.68	3 (60%) 0 0	66, 75, 92, 104	0
All	All	618/627 (98%)	0.06	38 (6%) 22 22	15, 34, 77, 119	0

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	G	5	GLN	6.3
2	H	1	ASN	5.3
2	G	3	THR	5.1
2	G	1	ASN	4.7
2	K	2	SER	4.6
1	A	191	ALA	4.4
2	K	1	ASN	4.1
1	B	139	SER	4.0
1	B	154	TYR	3.9
1	A	154	TYR	3.9
1	A	142	ASN	3.9
1	B	1	SER	3.8
1	B	138	GLY	3.8
1	B	194	ALA	3.7
1	B	141	LEU	3.5
1	B	168	PRO	3.2
1	B	46	ALA	3.2
1	B	302	GLY	3.2
2	K	3	THR	3.0
1	A	119	ASN	2.9

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Mol	Chain	Res	Type	RSRZ
2	H	2	SER	2.8
2	G	4	LEU	2.8
1	A	189	GLN	2.8
1	A	155	ASP	2.8
1	B	301	SER	2.7
1	A	146	GLY	2.6
1	B	169	THR	2.4
1	A	46	ALA	2.4
1	B	143	GLY	2.3
2	G	2	SER	2.3
1	B	140	PHE	2.2
1	A	147	SER	2.2
1	B	170	GLY	2.1
1	A	164	HIS	2.1
1	B	277	ASN	2.1
1	A	162	MET	2.0
1	A	64	HIS	2.0
1	A	39	PRO	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

## 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
3	ATO	A	1006	1/5	0.92	0.24	-0.53	100,100,100,100	0
3	ATO	B	2006	1/5	0.93	0.10	-2.31	76,76,76,76	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.