



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 06:58 PM GMT

PDB ID : 1DEU  
Title : CRYSTAL STRUCTURE OF HUMAN PROCATHEPSIN X: A CYSTEINE  
PROTEASE WITH THE PROREGION COVALENTLY LINKED TO THE  
ACTIVE SITE CYSTEINE  
Authors : Sivaraman, J.; Nagler, D.K.; Zhang, R.; Menard, R.; Cygler, M.  
Deposited on : 1999-11-15  
Resolution : 1.70 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  
<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  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
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) : trunk26865

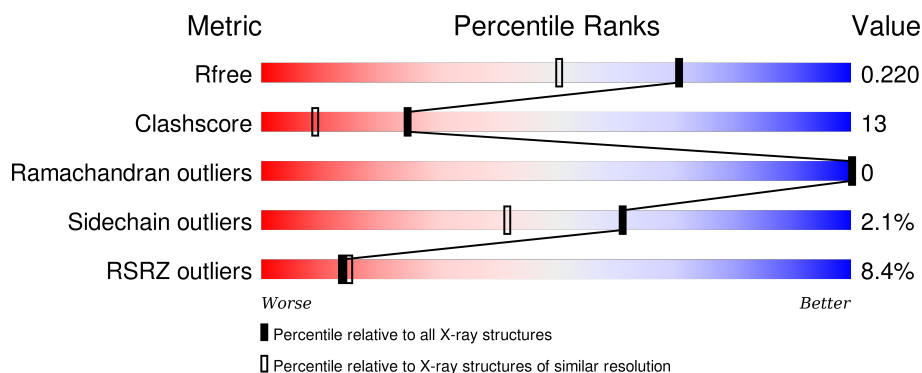
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.70 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))
$R_{free}$	91344	3190 (1.70-1.70)
Clashscore	102246	3585 (1.70-1.70)
Ramachandran outliers	100387	3527 (1.70-1.70)
Sidechain outliers	100360	3527 (1.70-1.70)
RSRZ outliers	91569	3200 (1.70-1.70)

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	277	<div> <div>9%</div> <div>82%</div> <div>16%</div> <div>..</div> </div>
1	B	277	<div> <div>7%</div> <div>79%</div> <div>12%</div> <div>.. 6%</div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4621 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 PROCATHEPSIN X.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	275	Total	C	N	O	S	0	0	0
			2163	1347	385	415	16			
1	B	261	Total	C	N	O	S	0	0	0
			2030	1265	357	392	16			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	25P	THR	SER	CONFLICT	UNP Q9UBR2
A	68	ALA	ARG	SEE REMARK 999	UNP Q9UBR2
A	89	PRO	SER	CONFLICT	UNP Q9UBR2
B	25P	THR	SER	CONFLICT	UNP Q9UBR2
B	68	ALA	ARG	SEE REMARK 999	UNP Q9UBR2
B	89	PRO	SER	CONFLICT	UNP Q9UBR2

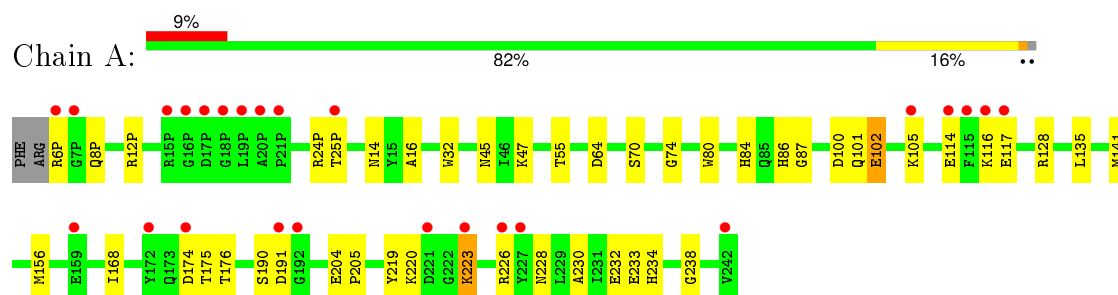
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	222	Total	O	0	0
			222	222		
2	B	206	Total	O	0	0
			206	206		

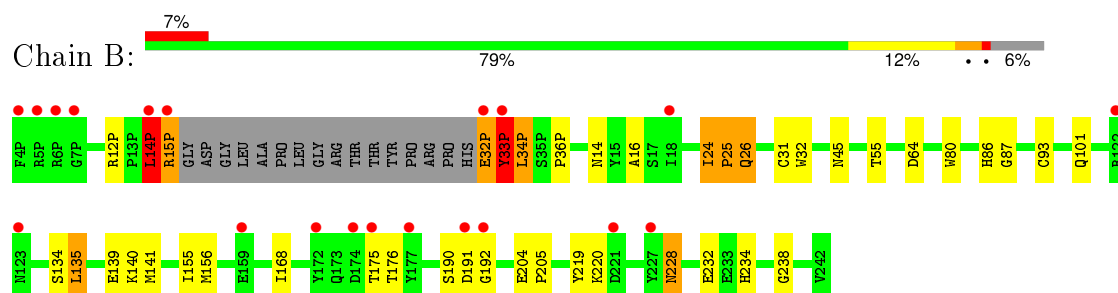
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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 ( $\text{RSRZ} > 2$ ). 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: PROCATHEPSIN X



#### • Molecule 1: PROCATHEPSIN X



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	84.82Å 84.82Å 169.72Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	45.00 – 1.70 30.81 – 1.65	Depositor EDS
% Data completeness (in resolution range)	95.1 (45.00-1.70) 95.6 (30.81-1.65)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.32 (at 1.65Å)	Xtriage
Refinement program	CNS	Depositor
R, $R_{free}$	0.204 , 0.215 0.208 , 0.220	Depositor DCC
$R_{free}$ test set	1900 reflections (2.48%)	DCC
Wilson B-factor (Å <sup>2</sup> )	15.3	Xtriage
Anisotropy	0.426	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.41 , 47.6	EDS
Estimated twinning fraction	0.026 for -h,-k,l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 81942 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	4621	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.27% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

## 5 Model quality

### 5.1 Standard geometry

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	0.30	0/2224	0.58	0/3026
1	B	8.13	19/2084 (0.9%)	1.85	23/2835 (0.8%)
All	All	5.66	19/4308 (0.4%)	1.35	23/5861 (0.4%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	2

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	15(P)	ARG	C-O	279.33	6.54	1.23
1	B	32(P)	GLU	C-O	148.54	4.05	1.23
1	B	32(P)	GLU	CA-CB	97.22	3.67	1.53
1	B	32(P)	GLU	N-CA	78.51	3.03	1.46
1	B	15(P)	ARG	CA-C	74.82	3.47	1.52
1	B	15(P)	ARG	CA-CB	72.54	3.13	1.53
1	B	15(P)	ARG	N-CA	65.22	2.76	1.46
1	B	33(P)	TYR	N-CA	52.75	2.51	1.46
1	B	33(P)	TYR	CA-C	39.20	2.54	1.52
1	B	33(P)	TYR	CA-CB	28.61	2.16	1.53
1	B	34(P)	LEU	N-CA	20.81	1.88	1.46
1	B	14(P)	LEU	N-CA	16.13	1.78	1.46
1	B	33(P)	TYR	C-O	15.82	1.53	1.23
1	B	14(P)	LEU	CA-C	14.78	1.91	1.52
1	B	14(P)	LEU	C-N	13.57	1.65	1.34
1	B	14(P)	LEU	C-O	9.45	1.41	1.23
1	B	33(P)	TYR	C-N	8.24	1.52	1.34
1	B	32(P)	GLU	CA-C	-6.54	1.35	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	25	PRO	CA-C	5.31	1.63	1.52

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	15(P)	ARG	CA-C-O	-51.50	11.96	120.10
1	B	32(P)	GLU	N-CA-CB	32.33	168.79	110.60
1	B	15(P)	ARG	CB-CA-C	24.43	159.26	110.40
1	B	32(P)	GLU	CA-C-O	-23.25	71.28	120.10
1	B	33(P)	TYR	O-C-N	-22.08	87.38	122.70
1	B	15(P)	ARG	N-CA-C	-21.46	53.07	111.00
1	B	33(P)	TYR	CB-CA-C	-21.25	67.89	110.40
1	B	15(P)	ARG	N-CA-CB	20.48	147.46	110.60
1	B	33(P)	TYR	N-CA-C	19.52	163.69	111.00
1	B	32(P)	GLU	CB-CA-C	-17.54	75.33	110.40
1	B	14(P)	LEU	CA-CB-CG	16.63	153.54	115.30
1	B	14(P)	LEU	CA-C-O	-15.89	86.73	120.10
1	B	14(P)	LEU	O-C-N	10.73	139.87	122.70
1	B	33(P)	TYR	CA-C-N	9.42	137.93	117.20
1	B	14(P)	LEU	CB-CG-CD2	-9.20	95.36	111.00
1	B	33(P)	TYR	N-CA-CB	8.02	125.04	110.60
1	B	14(P)	LEU	CA-C-N	7.13	132.88	117.20
1	B	135	LEU	CB-CG-CD1	-6.63	99.73	111.00
1	B	14(P)	LEU	N-CA-CB	-6.54	97.33	110.40
1	B	14(P)	LEU	N-CA-C	-6.43	93.63	111.00
1	B	14(P)	LEU	C-N-CA	6.11	136.97	121.70
1	B	33(P)	TYR	C-N-CA	5.99	136.67	121.70
1	B	135	LEU	CA-CB-CG	5.14	127.13	115.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	24	ILE	Mainchain
1	B	33(P)	TYR	Mainchain

## 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	2163	0	1999	40	0
1	B	2030	0	1841	61	0
2	A	222	0	0	5	0
2	B	206	0	0	7	0
All	All	4621	0	3840	101	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 13.

All (101) 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:B:14(P):LEU:N	1:B:14(P):LEU:CA	1.78	1.43
1:B:14(P):LEU:C	1:B:14(P):LEU:CA	1.91	1.38
1:B:34(P):LEU:CA	1:B:34(P):LEU:N	1.88	1.37
1:B:32(P):GLU:C	1:B:135:LEU:CA	1.90	1.37
1:B:33(P):TYR:CB	1:B:33(P):TYR:CA	2.16	1.24
1:B:32(P):GLU:C	1:B:135:LEU:HA	1.46	1.23
1:B:15(P):ARG:N	2:B:446:HOH:O	1.73	1.21
1:B:32(P):GLU:CB	1:B:33(P):TYR:N	2.09	1.15
1:B:15(P):ARG:CB	2:B:438:HOH:O	2.03	1.03
1:B:15(P):ARG:CB	1:B:32(P):GLU:N	2.24	1.01
1:B:15(P):ARG:O	2:B:438:HOH:O	1.89	0.89
1:B:15(P):ARG:CA	2:B:438:HOH:O	2.24	0.85
1:A:220:LYS:O	1:A:223:LYS:HG3	1.78	0.83
1:A:168:ILE:HD11	1:A:190:SER:HB3	1.61	0.82
1:B:32(P):GLU:C	1:B:135:LEU:CB	2.47	0.81
1:B:32(P):GLU:CA	1:B:135:LEU:HA	2.09	0.81
1:A:102:GLU:CD	1:A:102:GLU:H	1.83	0.81
1:B:15(P):ARG:CB	1:B:134:SER:HB2	2.12	0.80
1:B:14(P):LEU:CB	1:B:14(P):LEU:N	2.47	0.78
1:B:14(P):LEU:O	1:B:14(P):LEU:CA	2.31	0.78
1:B:33(P):TYR:C	1:B:33(P):TYR:CA	2.54	0.76
1:B:32(P):GLU:N	1:B:134:SER:HB2	2.01	0.75
1:B:15(P):ARG:O	1:B:234:HIS:NE2	2.20	0.75
1:B:156:MET:HE1	1:B:176:THR:HA	1.69	0.75
1:A:47:LYS:NZ	1:A:47:LYS:HB2	2.02	0.74
1:B:15(P):ARG:HA	2:B:438:HOH:O	1.84	0.74
1:B:33(P):TYR:HA	1:B:33(P):TYR:N	2.03	0.73
1:B:33(P):TYR:N	1:B:33(P):TYR:CA	2.51	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:32(P):GLU:C	1:B:135:LEU:HB3	2.07	0.73
1:B:15(P):ARG:O	1:B:134:SER:CB	2.39	0.71
1:B:15(P):ARG:CB	1:B:15(P):ARG:HA	2.20	0.70
1:B:15(P):ARG:CA	2:B:446:HOH:O	2.38	0.70
1:B:156:MET:CE	1:B:176:THR:HA	2.22	0.69
1:B:168:ILE:HD11	1:B:190:SER:HB3	1.77	0.67
1:B:15(P):ARG:CB	1:B:134:SER:CB	2.73	0.66
1:B:33(P):TYR:C	1:B:33(P):TYR:CB	2.65	0.65
1:A:176:THR:HG22	1:A:176:THR:O	1.94	0.65
1:A:176:THR:O	1:A:176:THR:CG2	2.47	0.61
1:A:6(P):ARG:C	1:A:8(P):GLN:H	2.04	0.61
1:A:230:ALA:HB1	1:A:233:GLU:HG3	1.82	0.61
1:B:36(P):PRO:HG3	1:B:139:GLU:HG3	1.81	0.60
1:B:14(P):LEU:CB	1:B:14(P):LEU:C	2.69	0.59
1:A:47:LYS:HB2	1:A:47:LYS:HZ3	1.66	0.58
1:B:14(P):LEU:O	1:B:14(P):LEU:N	2.36	0.58
1:A:223:LYS:NZ	1:A:223:LYS:HB3	2.19	0.57
1:B:24:ILE:HA	1:B:25:PRO:C	2.26	0.56
1:B:45:ASN:HD21	1:B:55:THR:H	1.53	0.56
1:B:36(P):PRO:HG3	1:B:139:GLU:CG	2.36	0.55
1:B:14(P):LEU:C	1:B:14(P):LEU:HB2	2.29	0.53
1:A:25(P):THR:HG23	2:A:337:HOH:O	2.08	0.53
1:A:24(P):ARG:HD3	1:A:84:HIS:CE1	2.45	0.51
1:A:174:ASP:HA	1:A:226:ARG:NH1	2.26	0.51
1:A:45:ASN:HD21	1:A:55:THR:H	1.59	0.51
1:B:14:ASN:HD21	1:B:16:ALA:HB3	1.77	0.49
1:A:226:ARG:HA	1:A:226:ARG:HE	1.77	0.49
1:B:24:ILE:HB	1:B:25:PRO:HA	1.94	0.49
1:A:86:HIS:HE1	2:A:394:HOH:O	1.95	0.49
1:B:32(P):GLU:O	1:B:140:LYS:HD3	2.14	0.48
1:A:80:TRP:CE3	1:A:238:GLY:HA3	2.48	0.48
1:A:135:LEU:C	1:A:135:LEU:HD12	2.34	0.48
1:B:15(P):ARG:N	1:B:15(P):ARG:CA	2.76	0.48
1:A:168:ILE:HD11	1:A:190:SER:CB	2.41	0.47
1:A:32:TRP:CD2	1:A:74:GLY:HA3	2.50	0.47
1:B:219:TYR:CD2	1:B:220:LYS:HG2	2.50	0.47
1:B:24:ILE:HA	1:B:26:GLN:N	2.30	0.47
1:A:168:ILE:CD1	1:A:190:SER:HB3	2.39	0.47
1:A:14:ASN:ND2	1:A:16:ALA:H	2.13	0.46
1:B:14(P):LEU:C	1:B:14(P):LEU:N	2.69	0.46
1:B:64:ASP:CG	1:B:101:GLN:HB2	2.36	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:204:GLU:N	1:B:205:PRO:CD	2.79	0.45
1:A:86:HIS:HD2	1:A:87:GLY:O	2.00	0.44
1:B:12(P):ARG:HD2	2:B:447:HOH:O	2.17	0.44
1:A:70:SER:HB2	1:A:100:ASP:OD1	2.17	0.44
1:B:86:HIS:HD2	1:B:87:GLY:O	2.00	0.44
1:B:141:MET:SD	1:B:232:GLU:HG2	2.58	0.44
1:A:14:ASN:HD21	1:A:16:ALA:HB3	1.82	0.44
1:A:47:LYS:HZ2	1:A:47:LYS:HB2	1.82	0.44
1:A:6(P):ARG:C	1:A:8(P):GLN:N	2.71	0.43
1:B:14:ASN:ND2	1:B:16:ALA:H	2.17	0.43
1:B:26:GLN:HB3	1:B:26:GLN:HE21	1.60	0.43
1:A:219:TYR:CD2	1:A:220:LYS:HG2	2.54	0.43
1:B:156:MET:HE2	1:B:176:THR:HA	1.99	0.42
1:A:204:GLU:N	1:A:205:PRO:CD	2.82	0.42
1:B:191:ASP:HB3	1:B:192:GLY:H	1.53	0.42
1:A:114:GLU:HB2	1:A:117:GLU:HB3	2.02	0.42
1:A:64:ASP:CG	1:A:101:GLN:HB2	2.40	0.42
1:B:80:TRP:CE3	1:B:238:GLY:HA3	2.55	0.41
1:A:156:MET:SD	1:A:175:THR:HG23	2.60	0.41
1:A:105:LYS:HD2	2:A:352:HOH:O	2.20	0.41
1:B:228:ASN:C	1:B:228:ASN:HD22	2.24	0.41
1:B:31:CYS:SG	1:B:32:TRP:N	2.93	0.41
1:A:87:GLY:CA	1:A:128:ARG:HD2	2.50	0.41
1:A:219:TYR:CE2	1:A:220:LYS:HG2	2.56	0.41
1:B:156:MET:HE1	1:B:176:THR:HG22	2.03	0.41
1:A:234:HIS:CE1	2:A:371:HOH:O	2.73	0.41
1:B:15(P):ARG:O	1:B:134:SER:HB3	2.20	0.40
1:A:116:LYS:HA	1:A:116:LYS:HD3	1.93	0.40
1:A:223:LYS:HB3	1:A:223:LYS:HZ2	1.84	0.40
1:B:155:ILE:HG12	1:B:156:MET:N	2.36	0.40
1:A:141:MET:SD	1:A:232:GLU:HG2	2.61	0.40
1:A:12(P):ARG:HD2	2:A:430: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	273/277 (99%)	264 (97%)	9 (3%)	0	100	100
1	B	256/277 (92%)	246 (96%)	10 (4%)	0	100	100
All	All	529/554 (96%)	510 (96%)	19 (4%)	0	100	100

There are no Ramachandran outliers to report.

### 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	225/228 (99%)	221 (98%)	4 (2%)	66	49
1	B	208/228 (91%)	203 (98%)	5 (2%)	57	36
All	All	433/456 (95%)	424 (98%)	9 (2%)	61	42

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	102	GLU
1	A	191	ASP
1	A	223	LYS
1	A	228	ASN
1	B	14(P)	LEU
1	B	26	GLN
1	B	93	CYS
1	B	175	THR
1	B	228	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	8(P)	GLN
1	A	14	ASN
1	A	23	HIS
1	A	45	ASN
1	A	61	ASN
1	A	67	ASN
1	A	86	HIS
1	A	113	ASN
1	A	173	GLN
1	A	228	ASN
1	B	14	ASN
1	B	23	HIS
1	B	26	GLN
1	B	45	ASN
1	B	61	ASN
1	B	86	HIS
1	B	113	ASN
1	B	163	ASN
1	B	228	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](#)

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 ⓘ

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	275/277 (99%)	0.36	25 (9%) 11 13	11, 18, 34, 38	0
1	B	261/277 (94%)	0.32	20 (7%) 16 18	11, 19, 33, 38	0
All	All	536/554 (96%)	0.34	45 (8%) 14 15	11, 18, 34, 38	0

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	15(P)	ARG	17.1
1	B	32(P)	GLU	9.7
1	A	20(P)	ALA	9.0
1	A	18(P)	GLY	8.5
1	B	33(P)	TYR	8.4
1	B	4(P)	PHE	5.7
1	B	6(P)	ARG	5.6
1	A	21(P)	PRO	5.6
1	A	115	PHE	5.5
1	A	19(P)	LEU	5.4
1	A	221	ASP	5.2
1	B	14(P)	LEU	5.0
1	A	15(P)	ARG	4.9
1	B	175	THR	4.8
1	A	174	ASP	4.5
1	B	191	ASP	4.5
1	B	174	ASP	4.5
1	A	191	ASP	4.4
1	A	7(P)	GLY	4.3
1	B	122	ARG	4.3
1	B	5(P)	ARG	3.7
1	A	6(P)	ARG	3.7
1	A	116	LYS	3.4
1	A	16(P)	GLY	3.4

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Mol	Chain	Res	Type	RSRZ
1	B	221	ASP	3.4
1	A	242	VAL	3.1
1	A	114	GLU	3.1
1	A	17(P)	ASP	3.1
1	A	223	LYS	3.0
1	A	226	ARG	2.9
1	A	192	GLY	2.7
1	B	177	TYR	2.7
1	B	159	GLU	2.7
1	B	123	ASN	2.7
1	A	172	TYR	2.7
1	A	227	TYR	2.7
1	B	192	GLY	2.6
1	B	7(P)	GLY	2.5
1	A	117	GLU	2.5
1	B	172	TYR	2.4
1	A	25(P)	THR	2.4
1	B	227	TYR	2.3
1	A	159	GLU	2.3
1	B	18	ILE	2.1
1	A	105	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](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.