



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 12:21 PM GMT

PDB ID : 3QYD  
Title : Crystal structure of a recombinant chimeric trypsin inhibitor  
Authors : Sen, U.; Majumder, S.; Khamrui, S.; Dasgupta, J.  
Deposited on : 2011-03-03  
Resolution : 2.97 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

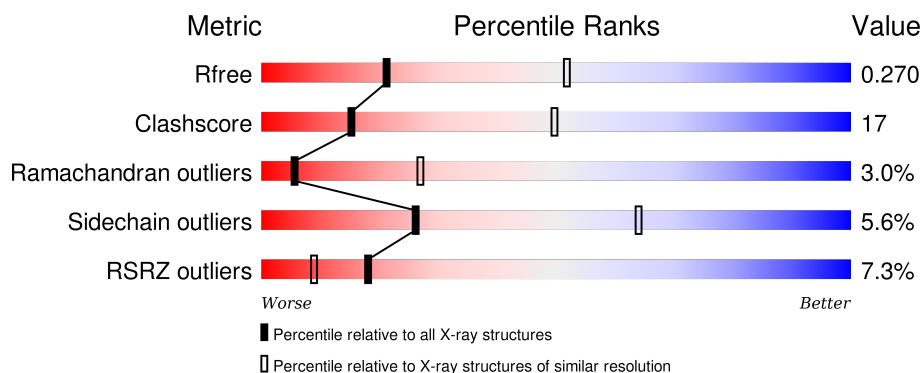
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.97 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	1992 (3.00-2.96)
Clashscore	102246	2349 (3.00-2.96)
Ramachandran outliers	100387	2274 (3.00-2.96)
Sidechain outliers	100360	2277 (3.00-2.96)
RSRZ outliers	91569	2007 (3.00-2.96)

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	181	 3% 62% 33% ..
1	B	181	 10% 59% 36% ...
1	C	181	 8% 56% 39% ...

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4323 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 Chymotrypsin inhibitor 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	177	Total	C	N	O	S	0	0	0
			1393	880	247	262	4			
1	C	177	Total	C	N	O	S	0	0	0
			1393	880	247	262	4			
1	B	177	Total	C	N	O	S	0	0	0
			1393	880	247	262	4			

There are 30 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	EXPRESSION TAG	UNP P10822
A	1	SER	-	EXPRESSION TAG	UNP P10822
A	2	HIS	-	EXPRESSION TAG	UNP P10822
A	3	MET	-	EXPRESSION TAG	UNP P10822
A	66	ARG	GLN	ENGINEERED MUTATION	UNP P10822
A	67	LEU	PHE	ENGINEERED MUTATION	UNP P10822
A	68	ARG	LEU	ENGINEERED MUTATION	UNP P10822
A	70	ALA	LEU	ENGINEERED MUTATION	UNP P10822
A	79	ARG	ALA	ENGINEERED MUTATION	UNP P10822
A	118	TYR	LEU	ENGINEERED MUTATION	UNP P10822
C	0	GLY	-	EXPRESSION TAG	UNP P10822
C	1	SER	-	EXPRESSION TAG	UNP P10822
C	2	HIS	-	EXPRESSION TAG	UNP P10822
C	3	MET	-	EXPRESSION TAG	UNP P10822
C	66	ARG	GLN	ENGINEERED MUTATION	UNP P10822
C	67	LEU	PHE	ENGINEERED MUTATION	UNP P10822
C	68	ARG	LEU	ENGINEERED MUTATION	UNP P10822
C	70	ALA	LEU	ENGINEERED MUTATION	UNP P10822
C	79	ARG	ALA	ENGINEERED MUTATION	UNP P10822
C	118	TYR	LEU	ENGINEERED MUTATION	UNP P10822
B	0	GLY	-	EXPRESSION TAG	UNP P10822
B	1	SER	-	EXPRESSION TAG	UNP P10822
B	2	HIS	-	EXPRESSION TAG	UNP P10822

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Chain	Residue	Modelled	Actual	Comment	Reference
B	3	MET	-	EXPRESSION TAG	UNP P10822
B	66	ARG	GLN	ENGINEERED MUTATION	UNP P10822
B	67	LEU	PHE	ENGINEERED MUTATION	UNP P10822
B	68	ARG	LEU	ENGINEERED MUTATION	UNP P10822
B	70	ALA	LEU	ENGINEERED MUTATION	UNP P10822
B	79	ARG	ALA	ENGINEERED MUTATION	UNP P10822
B	118	TYR	LEU	ENGINEERED MUTATION	UNP P10822

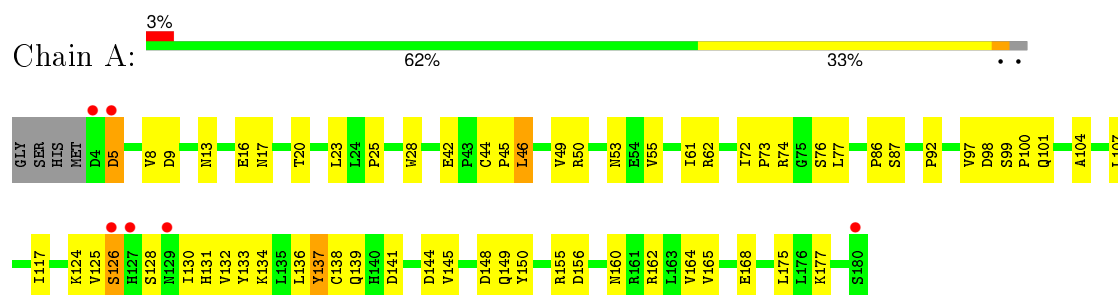
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	64	Total O 64 64	0	0
2	C	44	Total O 44 44	0	0
2	B	36	Total O 36 36	0	0

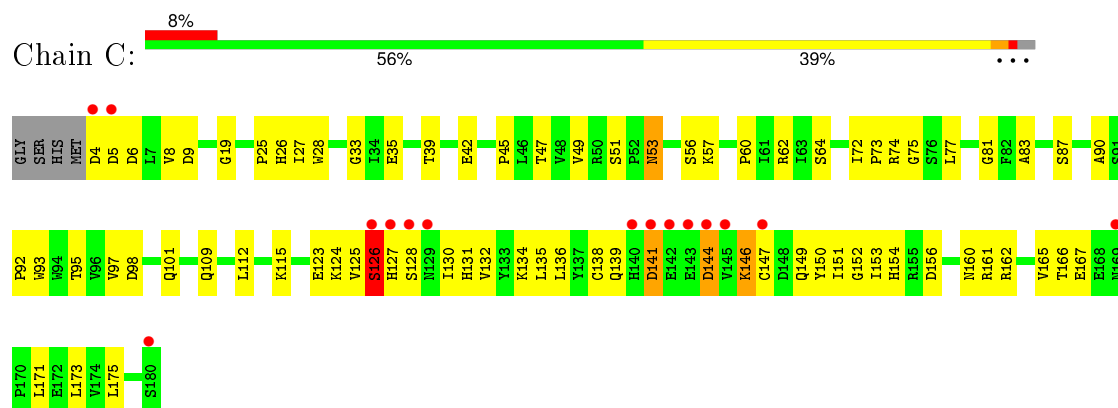
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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 ( $RSRZ > 2$ ). 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.

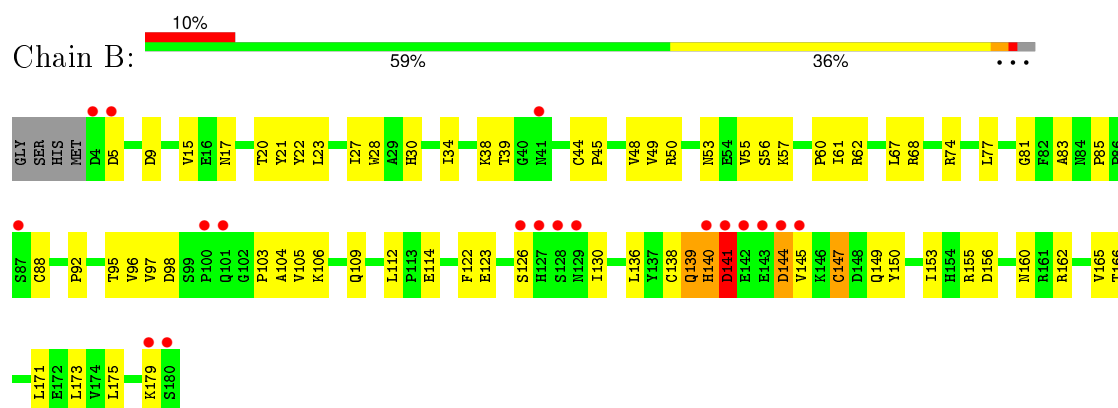
#### • Molecule 1: Chymotrypsin inhibitor 3



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## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	92.53 Å   39.25 Å   94.28 Å 90.00°   118.95°   90.00°	Depositor
Resolution (Å)	30.00 – 2.97 29.93 – 2.97	Depositor EDS
% Data completeness (in resolution range)	91.5 (30.00-2.97) 93.4 (29.93-2.97)	Depositor EDS
$R_{merge}$	0.12	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.90 (at 2.95 Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, $R_{free}$	0.210 , 0.270 0.210 , 0.270	Depositor DCC
$R_{free}$ test set	1148 reflections (9.74%)	DCC
Wilson B-factor (Å <sup>2</sup> )	35.7	Xtriage
Anisotropy	0.362	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 48.6	EDS
Estimated twinning fraction	0.014 for -h-l,k,h 0.014 for l,k,-h-l 0.008 for h,-k,-h-l 0.018 for -h-l,-k,l 0.016 for l,-k,h	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 11790 reflections	Xtriage
$F_o, F_c$ correlation	0.90	EDS
Total number of atoms	4323	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.83% 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 [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	0.38	0/1427	0.77	1/1940 (0.1%)
1	B	0.36	0/1427	0.72	1/1940 (0.1%)
1	C	0.36	0/1427	0.74	0/1940
All	All	0.37	0/4281	0.74	2/5820 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	46	LEU	CA-CB-CG	-6.15	101.16	115.30
1	B	130	ILE	N-CA-C	-5.51	96.11	111.00

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	1393	0	1374	46	0
1	B	1393	0	1374	51	0
1	C	1393	0	1374	50	0
2	A	64	0	0	6	0
2	B	36	0	0	4	0
2	C	44	0	0	5	0
All	All	4323	0	4122	144	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 17.

All (144) 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:C:49:VAL:HG12	1:C:162:ARG:HA	1.53	0.89
1:B:149:GLN:HG2	2:B:205:HOH:O	1.72	0.89
1:B:77:LEU:HD11	1:B:140:HIS:HB2	1.65	0.78
1:B:20:THR:HG22	1:B:62:ARG:HG3	1.68	0.75
1:A:23:LEU:HD13	1:A:61:ILE:HD12	1.69	0.75
1:B:48:VAL:HG21	1:B:105:VAL:HG12	1.69	0.73
1:C:26:HIS:HB3	2:C:276:HOH:O	1.87	0.72
1:A:45:PRO:HB2	1:A:86:PRO:HG3	1.71	0.72
1:B:97:VAL:HG12	1:B:98:ASP:H	1.54	0.72
1:A:137:TYR:O	1:A:138:CYS:HB3	1.88	0.72
1:A:44:CYS:HB2	1:A:45:PRO:CD	2.21	0.70
1:B:53:ASN:HB3	1:B:56:SER:H	1.56	0.69
1:A:20:THR:HG22	1:A:62:ARG:HG2	1.76	0.68
1:B:141:ASP:OD2	1:B:145:VAL:HG12	1.95	0.67
1:B:147:CYS:HB3	2:B:398:HOH:O	1.94	0.66
1:C:135:LEU:HG	1:C:151:ILE:HD12	1.78	0.66
1:C:4:ASP:CG	1:C:5:ASP:H	1.98	0.65
1:B:61:ILE:HG23	1:B:81:GLY:O	1.97	0.64
1:A:117:ILE:HD13	2:A:383:HOH:O	1.98	0.63
1:A:42:GLU:HB3	1:A:46:LEU:HD12	1.81	0.62
1:C:125:VAL:HG22	1:C:134:LYS:HG3	1.82	0.61
1:C:5:ASP:O	1:C:73:PRO:HA	2.01	0.60
1:B:49:VAL:HG12	1:B:162:ARG:HA	1.84	0.60
1:A:125:VAL:CG2	1:A:134:LYS:HG3	2.31	0.59
1:A:5:ASP:O	1:A:73:PRO:HA	2.03	0.59
1:A:155:ARG:HB3	2:A:328:HOH:O	2.03	0.58
1:C:6:ASP:HA	1:C:72:ILE:O	2.03	0.58
1:C:33:GLY:HA3	1:C:57:LYS:O	2.04	0.58
1:A:53:ASN:OD1	1:A:55:VAL:HG22	2.03	0.57
1:A:62:ARG:HH22	1:A:92:PRO:HG3	1.69	0.57
1:B:156:ASP:OD2	1:B:160:ASN:HB2	2.04	0.57
1:A:8:VAL:HG11	1:A:131:HIS:CE1	2.41	0.56
1:B:136:LEU:HD23	1:B:149:GLN:O	2.06	0.56
1:B:97:VAL:O	1:B:103:PRO:HA	2.06	0.55
1:C:101:GLN:HG2	1:C:154:HIS:CD2	2.41	0.55
1:B:138:CYS:O	1:B:139:GLN:HG2	2.07	0.55
1:B:77:LEU:HD11	1:B:140:HIS:CB	2.37	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:60:PRO:HG2	1:C:83:ALA:HB3	1.89	0.54
1:B:85:PRO:HG3	1:B:92:PRO:HB3	1.89	0.54
1:A:100:PRO:HB2	1:A:101:GLN:NE2	2.23	0.53
1:C:27:ILE:HG12	2:C:276:HOH:O	2.08	0.53
1:C:26:HIS:CD2	1:B:67:LEU:HD22	2.43	0.53
1:C:153:ILE:HG21	1:C:161:ARG:HB3	1.90	0.53
1:C:123:GLU:OE1	1:C:136:LEU:HD11	2.09	0.53
1:B:15:VAL:HG13	1:B:21:TYR:CE1	2.44	0.53
1:A:9:ASP:HB2	1:A:175:LEU:O	2.08	0.53
1:A:62:ARG:NH2	1:A:92:PRO:HG3	2.23	0.52
1:B:155:ARG:HA	1:B:160:ASN:O	2.11	0.51
1:C:81:GLY:HA2	1:C:92:PRO:O	2.11	0.51
1:C:128:SER:HB3	2:C:468:HOH:O	2.09	0.51
1:C:9:ASP:HB2	1:C:175:LEU:O	2.10	0.51
1:C:77:LEU:HD21	1:C:141:ASP:OD2	2.11	0.51
1:B:138:CYS:O	1:B:139:GLN:CG	2.59	0.51
1:C:47:THR:HG22	1:C:49:VAL:HG13	1.92	0.51
1:B:34:ILE:HG23	1:B:49:VAL:O	2.11	0.51
1:C:149:GLN:HB3	1:C:165:VAL:CG1	2.41	0.51
1:C:25:PRO:O	1:B:68:ARG:HD2	2.11	0.50
1:A:20:THR:HG22	1:A:62:ARG:CG	2.41	0.50
1:B:22:TYR:CD2	1:B:57:LYS:HB3	2.47	0.50
1:A:104:ALA:HA	1:A:164:VAL:HG12	1.93	0.50
1:C:153:ILE:CG2	1:C:161:ARG:HB3	2.42	0.50
1:A:25:PRO:HB3	1:A:50:ARG:NH2	2.26	0.50
1:A:136:LEU:HD21	1:A:150:TYR:CZ	2.48	0.49
1:B:103:PRO:HB2	1:B:165:VAL:CG2	2.43	0.49
1:A:125:VAL:HG23	1:A:134:LYS:HG3	1.93	0.49
1:B:153:ILE:HG13	1:B:171:LEU:HD21	1.95	0.49
1:A:128:SER:HA	2:A:324:HOH:O	2.12	0.48
1:C:125:VAL:O	1:C:126:SER:O	2.31	0.48
1:A:98:ASP:HB2	2:A:251:HOH:O	2.13	0.48
1:B:171:LEU:HB3	1:B:173:LEU:HD21	1.95	0.48
1:B:96:VAL:HA	1:B:104:ALA:O	2.14	0.48
1:B:9:ASP:HB2	1:B:175:LEU:O	2.13	0.48
1:A:16:GLU:OE1	1:A:177:LYS:HE2	2.14	0.48
1:A:125:VAL:O	1:A:125:VAL:HG12	2.13	0.48
1:B:39:THR:HB	2:B:465:HOH:O	2.13	0.48
1:A:144:ASP:O	1:A:145:VAL:HG23	2.14	0.47
1:C:149:GLN:HB3	1:C:165:VAL:HG11	1.96	0.47
1:B:145:VAL:HG13	1:B:145:VAL:O	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:125:VAL:O	1:A:126:SER:O	2.32	0.47
1:A:156:ASP:OD2	1:A:160:ASN:HB2	2.15	0.47
1:A:17:ASN:HB2	1:A:72:ILE:HG13	1.97	0.47
1:B:123:GLU:OE2	1:B:150:TYR:OH	2.26	0.47
1:A:136:LEU:HD21	1:A:150:TYR:CE1	2.49	0.47
1:B:141:ASP:OD1	1:B:144:ASP:HA	2.14	0.46
1:B:97:VAL:HG12	1:B:98:ASP:N	2.25	0.46
1:C:156:ASP:OD2	1:C:160:ASN:HB2	2.15	0.46
1:C:39:THR:O	1:C:42:GLU:HG3	2.15	0.46
1:C:126:SER:O	1:C:128:SER:N	2.49	0.46
1:A:44:CYS:HB2	1:A:45:PRO:HD2	1.96	0.46
1:C:4:ASP:CG	1:C:5:ASP:N	2.68	0.46
1:C:124:LYS:HE2	2:C:468:HOH:O	2.15	0.46
1:C:87:SER:HB2	2:C:430:HOH:O	2.15	0.46
1:B:62:ARG:HH22	1:B:92:PRO:HG3	1.81	0.46
1:B:88:CYS:O	1:B:109:GLN:HG2	2.16	0.45
1:C:166:THR:OG1	1:C:167:GLU:N	2.49	0.45
1:A:20:THR:HA	1:A:61:ILE:O	2.16	0.45
1:B:105:VAL:O	1:B:106:LYS:HG2	2.16	0.45
1:C:153:ILE:HG22	1:C:154:HIS:N	2.31	0.45
1:A:97:VAL:HG12	1:A:98:ASP:N	2.31	0.45
1:A:46:LEU:HD22	1:A:107:LEU:O	2.17	0.44
1:C:33:GLY:HA3	1:C:57:LYS:C	2.38	0.44
1:B:122:PHE:HB2	2:B:359:HOH:O	2.17	0.44
1:C:35:GLU:CD	1:C:51:SER:HB2	2.37	0.44
1:B:95:THR:HG22	1:B:106:LYS:O	2.18	0.44
1:A:44:CYS:HB2	1:A:45:PRO:HD3	1.96	0.44
1:A:76:SER:HB3	2:A:386:HOH:O	2.18	0.44
1:A:149:GLN:HB3	1:A:165:VAL:CG1	2.48	0.44
1:C:74:ARG:HG3	1:C:75:GLY:N	2.32	0.44
1:B:20:THR:HA	1:B:61:ILE:O	2.17	0.44
1:B:55:VAL:HG13	1:B:56:SER:N	2.31	0.43
1:A:130:ILE:O	1:A:132:VAL:HG23	2.18	0.43
1:C:19:GLY:O	1:C:62:ARG:HA	2.18	0.43
1:A:117:ILE:HG21	2:A:383:HOH:O	2.17	0.43
1:B:103:PRO:HB2	1:B:165:VAL:HG21	2.00	0.43
1:C:125:VAL:CG2	1:C:134:LYS:HG3	2.49	0.43
1:A:77:LEU:HG	1:A:141:ASP:HA	2.00	0.43
1:A:5:ASP:O	1:A:74:ARG:N	2.52	0.43
1:B:153:ILE:HA	1:B:162:ARG:O	2.19	0.43
1:C:53:ASN:HB3	1:C:56:SER:H	1.84	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:27:ILE:HG13	1:B:30:HIS:CE1	2.54	0.43
1:B:60:PRO:HG2	1:B:83:ALA:HB3	2.00	0.42
1:B:145:VAL:CG1	1:B:145:VAL:O	2.68	0.42
1:C:130:ILE:HD12	1:C:131:HIS:H	1.84	0.42
1:C:152:GLY:HA2	1:C:171:LEU:HG	2.01	0.42
1:C:97:VAL:HG12	1:C:98:ASP:N	2.34	0.42
1:C:26:HIS:CD2	1:C:132:VAL:HG21	2.55	0.42
1:A:49:VAL:HG12	1:A:162:ARG:HA	2.00	0.42
1:A:8:VAL:HG13	1:A:13:ASN:O	2.20	0.42
1:C:35:GLU:OE2	1:C:51:SER:HB2	2.19	0.42
1:B:38:LYS:HG3	1:B:44:CYS:C	2.40	0.42
1:C:49:VAL:HG12	1:C:162:ARG:CA	2.36	0.41
1:C:27:ILE:HD13	1:B:67:LEU:HD23	2.02	0.41
1:C:8:VAL:HB	1:C:131:HIS:CD2	2.55	0.41
1:B:48:VAL:CG2	1:B:105:VAL:HG12	2.45	0.41
1:C:62:ARG:HD3	1:C:93:TRP:HH2	1.85	0.41
1:A:23:LEU:HD13	1:A:61:ILE:CD1	2.43	0.41
1:B:95:THR:HB	1:B:112:LEU:CD1	2.51	0.41
1:A:42:GLU:CD	1:A:46:LEU:HB2	2.40	0.41
1:C:90:ALA:HB2	1:C:109:GLN:O	2.21	0.41
1:C:136:LEU:HD21	1:C:150:TYR:CZ	2.56	0.41
1:B:141:ASP:CG	1:B:145:VAL:H	2.24	0.41
1:B:50:ARG:CZ	1:B:153:ILE:HD11	2.52	0.40
1:C:95:THR:HB	1:C:112:LEU:HD13	2.03	0.40
1:A:124:LYS:HA	1:A:133:TYR:CD1	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	175/181 (97%)	159 (91%)	14 (8%)	2 (1%)	17 57

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	175/181 (97%)	155 (89%)	13 (7%)	7 (4%)	4	19
1	C	175/181 (97%)	155 (89%)	13 (7%)	7 (4%)	4	19
All	All	525/543 (97%)	469 (89%)	40 (8%)	16 (3%)	5	27

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	87	SER
1	A	126	SER
1	C	126	SER
1	C	127	HIS
1	C	147	CYS
1	B	5	ASP
1	C	146	LYS
1	B	17	ASN
1	B	126	SER
1	B	144	ASP
1	B	179	LYS
1	C	144	ASP
1	B	141	ASP
1	C	45	PRO
1	C	141	ASP
1	B	45	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	155/158 (98%)	148 (96%)	7 (4%)	34	72
1	B	155/158 (98%)	146 (94%)	9 (6%)	25	62
1	C	155/158 (98%)	145 (94%)	10 (6%)	21	57
All	All	465/474 (98%)	439 (94%)	26 (6%)	26	64

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	ASP
1	A	28	TRP
1	A	99	SER
1	A	137	TYR
1	A	139	GLN
1	A	148	ASP
1	A	168	GLU
1	C	28	TRP
1	C	53	ASN
1	C	64	SER
1	C	115	LYS
1	C	126	SER
1	C	138	CYS
1	C	139	GLN
1	C	144	ASP
1	C	146	LYS
1	C	173	LEU
1	B	23	LEU
1	B	28	TRP
1	B	74	ARG
1	B	114	GLU
1	B	139	GLN
1	B	140	HIS
1	B	141	ASP
1	B	147	CYS
1	B	166	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	41	ASN
1	A	101	GLN
1	A	139	GLN
1	C	26	HIS
1	C	41	ASN
1	C	131	HIS
1	C	139	GLN
1	B	30	HIS
1	B	41	ASN
1	B	139	GLN

### 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 [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	177/181 (97%)	-0.26	6 (3%)	49 28	7, 21, 60, 115	2 (1%)
1	B	177/181 (97%)	0.24	18 (10%)	9 4	11, 32, 98, 139	2 (1%)
1	C	177/181 (97%)	0.22	15 (8%)	13 6	14, 32, 102, 191	2 (1%)
All	All	531/543 (97%)	0.07	39 (7%)	18 9	7, 28, 94, 191	6 (1%)

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	140	HIS	11.7
1	A	180	SER	7.1
1	B	180	SER	6.4
1	C	141	ASP	5.8
1	B	144	ASP	5.6
1	B	142	GLU	5.5
1	B	127	HIS	5.4
1	B	143	GLU	5.2
1	C	144	ASP	5.2
1	C	127	HIS	5.1
1	A	127	HIS	5.1
1	C	180	SER	5.0
1	C	142	GLU	4.6
1	B	4	ASP	4.6
1	C	143	GLU	4.3
1	C	129	ASN	4.3
1	B	140	HIS	4.0
1	B	126	SER	3.9
1	B	129	ASN	3.9
1	B	5	ASP	3.6
1	A	4	ASP	3.6
1	C	128	SER	3.5
1	C	126	SER	3.5

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Mol	Chain	Res	Type	RSRZ
1	A	5	ASP	3.3
1	B	145	VAL	3.3
1	B	128	SER	3.3
1	A	129	ASN	3.2
1	B	87	SER	3.0
1	C	169	ASN	3.0
1	B	100	PRO	2.8
1	C	5	ASP	2.7
1	C	4	ASP	2.7
1	C	147	CYS	2.5
1	A	126	SER	2.5
1	B	41	ASN	2.5
1	C	145	VAL	2.5
1	B	179	LYS	2.3
1	B	141	ASP	2.3
1	B	101	GLN	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.