



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:25 AM GMT

PDB ID : 3AQQ
Title : Crystal structure of human CRHSP-24
Authors : Hou, H.; Wang, F.; Zhang, W.; Wang, D.; Li, X.; Bartlam, M.; Yao, X.; Rao, Z.
Deposited on : 2010-11-17
Resolution : 2.80 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
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

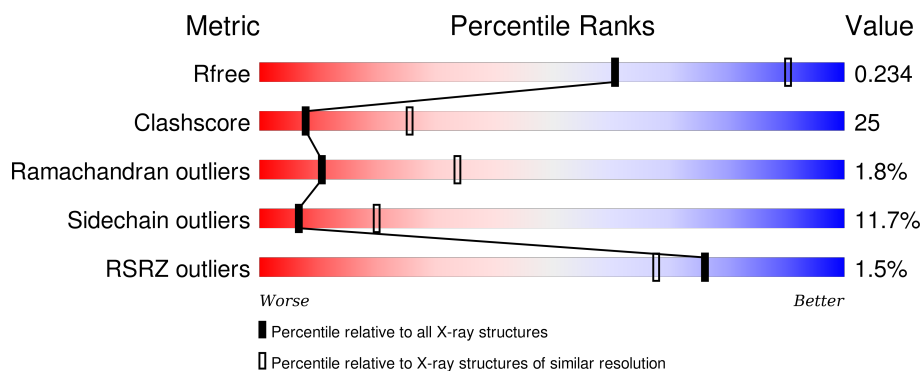
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	2393 (2.80-2.80)
Clashscore	102246	2827 (2.80-2.80)
Ramachandran outliers	100387	2782 (2.80-2.80)
Sidechain outliers	100360	2784 (2.80-2.80)
RSRZ outliers	91569	2404 (2.80-2.80)

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 ≥ 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	147	<div> <div>3%</div> <div>46%</div> <div>16%</div> <div>5%</div> <div>33%</div> </div>
1	B	147	<div> <div>35%</div> <div>27%</div> <div>•</div> <div>•</div> <div>33%</div> </div>
1	C	147	<div> <div>%</div> <div>38%</div> <div>20%</div> <div>7%</div> <div>34%</div> </div>
1	D	147	<div> <div>%</div> <div>31%</div> <div>32%</div> <div>•</div> <div>33%</div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3048 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 Calcium-regulated heat stable protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	99	Total	C	N	O	S	0	0	0
			762	481	133	143	5			
1	C	97	Total	C	N	O	S	0	0	0
			737	463	131	138	5			
1	B	99	Total	C	N	O	S	0	0	0
			763	483	133	142	5			
1	D	98	Total	C	N	O	S	0	0	0
			754	477	131	141	5			

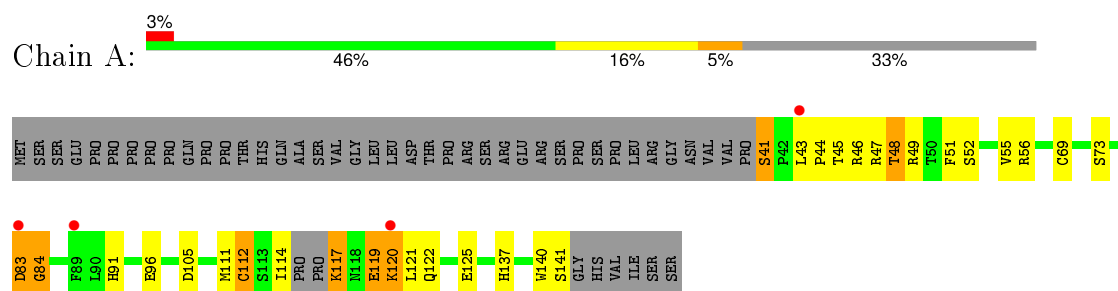
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	4	Total	O	0	0
			4	4		
2	C	8	Total	O	0	0
			8	8		
2	B	12	Total	O	0	0
			12	12		
2	D	8	Total	O	0	0
			8	8		

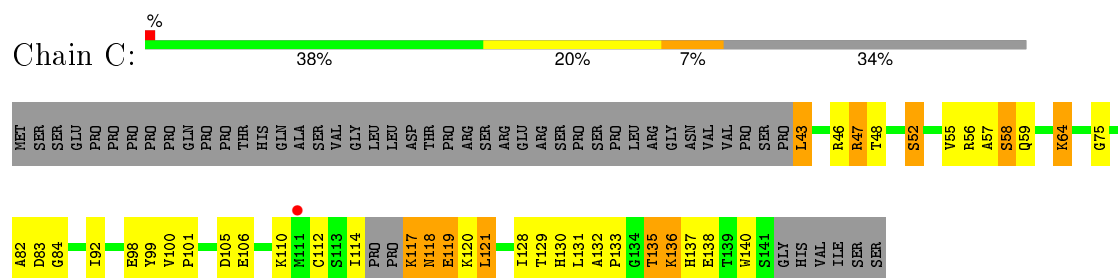
3 Residue-property plots

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

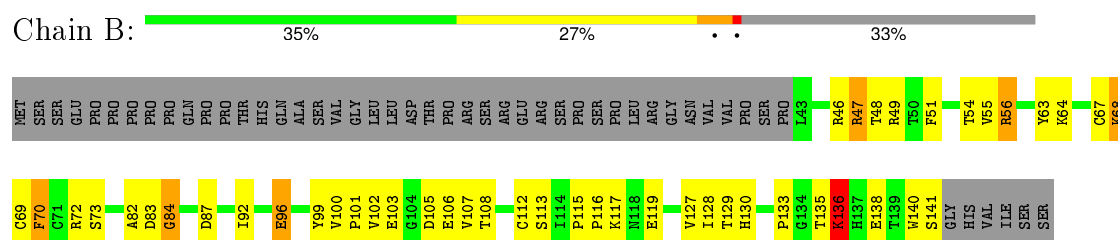
- Molecule 1: Calcium-regulated heat stable protein 1



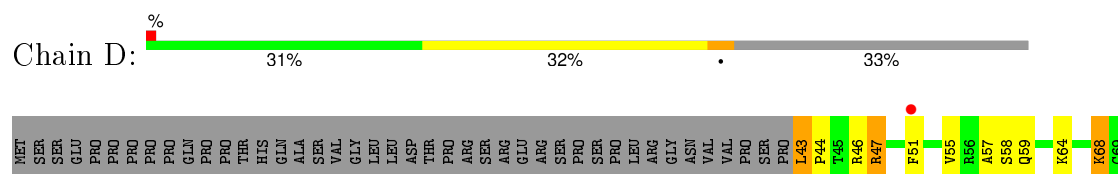
- Molecule 1: Calcium-regulated heat stable protein 1



- Molecule 1: Calcium-regulated heat stable protein 1



- Molecule 1: Calcium-regulated heat stable protein 1



GLY	F70
HIS	C71
VAL	R72
ILE	S73
SER	G77
SER	T80
	P81
	A82
	D83
	G84
	D87
	H91
	I92
	S93
	Y99
	V100
	P101
	V102
	E103
	G104
	D105
	E106
	T108
	V109
	K110
	M111
	C112
	S113
	I114
	P115
	P116
	LVS
	N118
	E119
	K120
	L121
	V124
	V127
	I128
	T129
	H130
	L131
	A132
	P133
	G134
	T135
	K136
	H137
	E138
	T139
	H140
	S141

4 Data and refinement statistics

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α , β , γ	79.86 Å 79.86 Å 181.98 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	45.60 – 2.80 45.60 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.6 (45.60-2.80) 99.6 (45.60-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.06 (at 2.81 Å)	Xtriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, R_{free}	0.229 , 0.272 0.200 , 0.234	Depositor DCC
R_{free} test set	811 reflections (5.02%)	DCC
Wilson B-factor (Å ²)	60.0	Xtriage
Anisotropy	0.482	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 46.3	EDS
Estimated twinning fraction	0.214 for h,-h-k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.26$	Xtriage
Outliers	0 of 16140 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3048	wwPDB-VP
Average B, all atoms (Å ²)	63.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 19.25% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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	0.42	0/780	0.60	0/1056
1	B	0.58	0/783	0.71	0/1063
1	C	0.58	0/753	0.66	0/1019
1	D	0.55	0/773	0.62	0/1049
All	All	0.54	0/3089	0.65	0/4187

There are no bond length outliers.

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	762	0	752	24	0
1	B	763	0	755	40	1
1	C	737	0	722	47	1
1	D	754	0	741	47	0
2	A	4	0	0	0	0
2	B	12	0	0	0	0
2	C	8	0	0	1	0
2	D	8	0	0	0	0
All	All	3048	0	2970	149	1

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 25.

All (149) 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:117:LYS:O	1:C:117:LYS:HE3	1.42	1.18
1:C:120:LYS:C	1:C:121:LEU:HD23	1.70	1.11
1:C:47:ARG:HG2	1:C:47:ARG:NH1	1.52	1.05
1:C:121:LEU:N	1:C:121:LEU:HD23	1.54	1.05
1:C:47:ARG:HH11	1:C:47:ARG:CG	1.65	1.05
1:C:118:ASN:H	1:C:118:ASN:ND2	1.56	0.95
1:C:121:LEU:N	1:C:121:LEU:CD2	2.30	0.91
1:C:118:ASN:HD22	1:C:118:ASN:H	0.99	0.90
1:D:43:LEU:N	1:D:44:PRO:HD3	1.88	0.87
1:C:117:LYS:O	1:C:117:LYS:HG2	1.70	0.86
1:C:118:ASN:O	1:C:119:GLU:HG2	1.74	0.86
1:C:118:ASN:C	1:C:119:GLU:HG3	2.01	0.81
1:C:118:ASN:O	1:C:119:GLU:CG	2.30	0.80
1:C:47:ARG:HG2	1:C:47:ARG:HH11	0.71	0.80
1:D:103:GLU:CD	1:D:103:GLU:H	1.84	0.80
1:C:118:ASN:C	1:C:119:GLU:CG	2.49	0.80
1:B:69:CYS:O	1:B:70:PHE:HB2	1.82	0.79
1:D:116:PRO:HD3	1:D:118:ASN:N	1.99	0.77
1:B:106:GLU:HB2	1:B:130:HIS:HB2	1.66	0.77
1:B:92:ILE:HD11	1:B:140:TRP:CZ3	2.19	0.77
1:C:118:ASN:N	1:C:118:ASN:ND2	2.29	0.77
1:C:118:ASN:HD22	1:C:118:ASN:N	1.77	0.76
1:A:52:SER:O	1:A:56:ARG:HG3	1.85	0.76
1:C:120:LYS:C	1:C:121:LEU:CD2	2.54	0.76
1:C:52:SER:O	1:C:56:ARG:HG3	1.86	0.76
1:C:117:LYS:CE	1:C:117:LYS:O	2.30	0.75
1:C:117:LYS:CG	1:C:117:LYS:O	2.30	0.75
1:C:82:ALA:O	1:C:83:ASP:HB2	1.87	0.74
1:D:111:MET:SD	1:D:121:LEU:HD13	2.29	0.72
1:B:51:PHE:O	1:B:55:VAL:HG23	1.90	0.71
1:B:56:ARG:CZ	1:B:56:ARG:HB2	2.23	0.68
1:D:116:PRO:CG	1:D:118:ASN:N	2.58	0.67
1:C:136:LYS:HD2	1:C:137:HIS:N	2.10	0.67
1:D:82:ALA:O	1:D:83:ASP:C	2.33	0.66
1:A:45:THR:HG22	1:A:91:HIS:CE1	2.31	0.65
1:C:119:GLU:OE1	1:B:117:LYS:NZ	2.20	0.65
1:D:83:ASP:OD1	1:D:84:GLY:N	2.29	0.65
1:D:114:ILE:O	1:D:118:ASN:HA	1.97	0.65
1:D:83:ASP:CG	1:D:84:GLY:N	2.47	0.65
1:D:116:PRO:CD	1:D:118:ASN:N	2.61	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:135:THR:HG22	1:D:136:LYS:H	1.62	0.63
1:C:47:ARG:NH1	1:C:47:ARG:CG	2.35	0.63
1:B:102:VAL:HG13	1:B:103:GLU:OE1	1.99	0.62
1:D:43:LEU:N	1:D:44:PRO:CD	2.61	0.61
1:D:99:TYR:HD1	1:D:138:GLU:O	1.82	0.61
1:B:103:GLU:H	1:B:103:GLU:CD	2.03	0.60
1:A:47:ARG:HD3	1:B:140:TRP:CD1	2.35	0.60
1:B:117:LYS:HB2	1:B:119:GLU:HG3	1.85	0.59
1:A:140:TRP:CD1	1:B:47:ARG:HB3	2.38	0.59
1:A:47:ARG:NH2	1:B:92:ILE:O	2.36	0.59
1:D:68:LYS:HD2	1:D:87:ASP:OD2	2.04	0.57
1:B:70:PHE:CE1	1:B:101:PRO:HD2	2.40	0.56
1:A:51:PHE:O	1:A:55:VAL:HG23	2.05	0.56
1:C:136:LYS:HD2	1:C:137:HIS:H	1.72	0.55
1:C:128:ILE:HG21	1:C:131:LEU:HD23	1.89	0.55
1:D:105:ASP:OD1	1:D:137:HIS:HE1	1.90	0.54
1:B:69:CYS:O	1:B:70:PHE:CB	2.53	0.54
1:A:83:ASP:O	1:A:84:GLY:O	2.27	0.53
1:C:110:LYS:HE2	1:B:56:ARG:HH12	1.73	0.53
1:C:64:LYS:HE3	1:C:106:GLU:OE2	2.08	0.53
1:D:100:VAL:HG22	1:D:140:TRP:CE3	2.44	0.53
1:D:108:THR:HG23	1:D:129:THR:HG21	1.90	0.53
1:B:72:ARG:HH12	1:B:140:TRP:HA	1.73	0.53
1:B:82:ALA:O	1:B:83:ASP:HB2	2.09	0.52
1:A:45:THR:HG22	1:A:91:HIS:NE2	2.24	0.52
1:A:46:ARG:HA	1:A:49:ARG:NH1	2.24	0.52
1:D:116:PRO:HG3	1:D:118:ASN:N	2.24	0.51
1:B:115:PRO:HB3	1:B:116:PRO:HA	1.92	0.51
1:A:46:ARG:HA	1:A:49:ARG:HH12	1.75	0.51
1:D:70:PHE:CE1	1:D:101:PRO:HD2	2.44	0.51
1:A:96:GLU:OE2	1:A:125:GLU:HA	2.11	0.51
1:A:119:GLU:O	1:A:120:LYS:O	2.30	0.50
1:D:72:ARG:HH22	1:D:141:SER:H	1.59	0.50
1:D:82:ALA:O	1:D:83:ASP:O	2.30	0.50
1:C:130:HIS:HB3	2:C:149:HOH:O	2.12	0.49
1:D:135:THR:HG22	1:D:136:LYS:N	2.27	0.49
1:A:140:TRP:O	1:A:141:SER:OG	2.26	0.49
1:B:117:LYS:HB2	1:B:119:GLU:CG	2.43	0.49
1:B:105:ASP:HB3	1:B:128:ILE:HG23	1.95	0.49
1:D:83:ASP:OD1	1:D:84:GLY:O	2.30	0.49
1:A:119:GLU:O	1:A:120:LYS:C	2.51	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:106:GLU:HB3	1:D:130:HIS:HB2	1.95	0.48
1:B:68:LYS:HE2	1:B:87:ASP:OD2	2.14	0.48
1:D:51:PHE:O	1:D:55:VAL:HG23	2.14	0.48
1:B:115:PRO:CB	1:B:116:PRO:HA	2.44	0.48
1:D:83:ASP:CG	1:D:84:GLY:H	2.17	0.48
1:B:64:LYS:HG2	1:B:108:THR:HG22	1.96	0.48
1:C:98:GLU:HA	1:D:47:ARG:O	2.14	0.48
1:D:99:TYR:CD1	1:D:138:GLU:O	2.64	0.48
1:D:70:PHE:HA	1:D:77:GLY:HA3	1.96	0.48
1:B:99:TYR:HD1	1:B:138:GLU:O	1.96	0.48
1:B:63:TYR:O	1:B:108:THR:HA	2.13	0.47
1:C:114:ILE:HG13	1:C:120:LYS:O	2.14	0.47
1:A:105:ASP:OD1	1:A:137:HIS:HE1	1.98	0.47
1:C:128:ILE:HG21	1:C:131:LEU:CD2	2.44	0.47
1:A:41:SER:O	1:A:43:LEU:HD23	2.15	0.47
1:D:58:SER:HA	1:D:110:LYS:HD2	1.98	0.46
1:B:72:ARG:NH1	1:B:140:TRP:HA	2.31	0.46
1:D:57:ALA:HB1	1:D:124:VAL:HG21	1.96	0.46
1:D:132:ALA:HA	1:D:133:PRO:HD3	1.70	0.45
1:D:105:ASP:OD1	1:D:137:HIS:CE1	2.69	0.45
1:A:44:PRO:HB2	1:A:122:GLN:HE22	1.80	0.45
1:A:114:ILE:O	1:A:117:LYS:HA	2.16	0.45
1:B:107:VAL:HA	1:B:127:VAL:O	2.17	0.45
1:B:102:VAL:CG1	1:B:103:GLU:N	2.80	0.44
1:C:75:GLY:O	1:C:92:ILE:HG22	2.17	0.44
1:C:46:ARG:NH1	1:D:93:SER:O	2.49	0.44
1:A:111:MET:SD	1:A:121:LEU:HD13	2.57	0.44
1:B:56:ARG:HB2	1:B:56:ARG:NH1	2.33	0.44
1:A:49:ARG:HD3	1:A:112:CYS:SG	2.58	0.44
1:B:136:LYS:HE3	1:B:136:LYS:HB3	1.62	0.44
1:D:64:LYS:HD3	1:D:106:GLU:OE2	2.17	0.44
1:D:51:PHE:C	1:D:51:PHE:CD1	2.90	0.44
1:B:113:SER:O	1:B:115:PRO:HD3	2.18	0.44
1:B:70:PHE:HE1	1:B:100:VAL:HG13	1.83	0.43
1:C:92:ILE:HD11	1:C:140:TRP:CZ3	2.54	0.43
1:B:49:ARG:HG2	1:B:54:THR:OG1	2.18	0.43
1:C:105:ASP:OD1	1:C:137:HIS:HE1	2.01	0.43
1:B:83:ASP:O	1:B:84:GLY:O	2.37	0.43
1:B:108:THR:HG23	1:B:129:THR:HG21	2.01	0.43
1:B:102:VAL:HG13	1:B:103:GLU:CD	2.39	0.43
1:C:58:SER:HB2	1:D:127:VAL:HG21	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:91:HIS:CE1	1:D:92:ILE:HG22	2.54	0.43
1:D:113:SER:OG	1:D:118:ASN:HB3	2.19	0.42
1:B:135:THR:O	1:B:135:THR:HG22	2.18	0.42
1:C:132:ALA:HA	1:C:133:PRO:HD3	1.86	0.42
1:D:83:ASP:OD1	1:D:83:ASP:C	2.57	0.42
1:A:45:THR:OG1	1:A:48:THR:HB	2.19	0.42
1:C:114:ILE:N	1:C:120:LYS:O	2.50	0.42
1:A:121:LEU:HA	1:A:121:LEU:HD23	1.85	0.42
1:C:131:LEU:HD22	1:C:137:HIS:CE1	2.55	0.42
1:B:96:GLU:HG2	1:B:96:GLU:H	1.69	0.42
1:C:55:VAL:O	1:C:59:GLN:HB2	2.19	0.42
1:C:135:THR:HG22	1:C:136:LYS:H	1.85	0.42
1:D:46:ARG:HE	1:D:46:ARG:HB3	1.52	0.42
1:D:80:THR:HA	1:D:81:PRO:HD2	1.93	0.41
1:A:43:LEU:HA	1:A:44:PRO:HD3	1.79	0.41
1:A:44:PRO:HB2	1:A:122:GLN:NE2	2.35	0.41
1:B:102:VAL:HG12	1:B:103:GLU:N	2.34	0.41
1:D:72:ARG:NH1	1:D:140:TRP:HA	2.36	0.41
1:D:55:VAL:O	1:D:59:GLN:HG3	2.21	0.41
1:C:43:LEU:C	1:C:43:LEU:HD22	2.40	0.41
1:D:44:PRO:HG2	1:D:114:ILE:HG23	2.03	0.40
1:D:70:PHE:CD1	1:D:101:PRO:HD2	2.55	0.40
1:C:57:ALA:C	1:C:59:GLN:H	2.24	0.40
1:C:99:TYR:CE2	1:D:51:PHE:HB2	2.57	0.40
1:B:67:CYS:SG	1:B:69:CYS:O	2.79	0.40
1:C:57:ALA:HB2	1:C:112:CYS:HB2	2.04	0.40
1:C:100:VAL:HA	1:C:101:PRO:HD3	1.81	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:133:PRO:O	1:B:133:PRO:O[3_564]	2.09	0.11

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	95/147 (65%)	86 (90%)	7 (7%)	2 (2%)	9	29
1	B	97/147 (66%)	89 (92%)	5 (5%)	3 (3%)	5	17
1	C	93/147 (63%)	82 (88%)	10 (11%)	1 (1%)	17	50
1	D	94/147 (64%)	87 (93%)	6 (6%)	1 (1%)	17	50
All	All	379/588 (64%)	344 (91%)	28 (7%)	7 (2%)	11	34

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	84	GLY
1	A	120	LYS
1	B	70	PHE
1	C	84	GLY
1	B	84	GLY
1	B	136	LYS
1	D	119	GLU

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	85/129 (66%)	77 (91%)	8 (9%)	11	31
1	B	85/129 (66%)	75 (88%)	10 (12%)	6	19
1	C	80/129 (62%)	66 (82%)	14 (18%)	2	7
1	D	84/129 (65%)	77 (92%)	7 (8%)	14	38
All	All	334/516 (65%)	295 (88%)	39 (12%)	7	20

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	41	SER
1	A	48	THR
1	A	69	CYS
1	A	73	SER
1	A	83	ASP
1	A	112	CYS
1	A	117	LYS
1	A	119	GLU
1	C	43	LEU
1	C	47	ARG
1	C	48	THR
1	C	52	SER
1	C	58	SER
1	C	64	LYS
1	C	117	LYS
1	C	118	ASN
1	C	119	GLU
1	C	121	LEU
1	C	129	THR
1	C	135	THR
1	C	136	LYS
1	C	138	GLU
1	B	46	ARG
1	B	47	ARG
1	B	48	THR
1	B	56	ARG
1	B	68	LYS
1	B	73	SER
1	B	96	GLU
1	B	112	CYS
1	B	136	LYS
1	B	141	SER
1	D	43	LEU
1	D	47	ARG
1	D	68	LYS
1	D	73	SER
1	D	83	ASP
1	D	120	LYS
1	D	135	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	76	HIS
1	C	118	ASN
1	B	118	ASN
1	D	137	HIS

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 [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, 95th 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 ²)	Q<0.9
1	A	99/147 (67%)	0.66	4 (4%) 42 30	50, 63, 88, 104	0
1	B	99/147 (67%)	0.63	0 100 100	43, 58, 83, 95	0
1	C	97/147 (65%)	0.58	1 (1%) 84 77	45, 58, 84, 112	0
1	D	98/147 (66%)	0.59	1 (1%) 84 77	47, 62, 91, 103	0
All	All	393/588 (66%)	0.62	6 (1%) 76 68	43, 61, 88, 112	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	120	LYS	2.3
1	D	51	PHE	2.2
1	A	83	ASP	2.2
1	A	89	PHE	2.1
1	A	43	LEU	2.1
1	C	111	MET	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.