



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 06:56 AM UTC

PDB ID : 3AQQ / pdb_00003aqq
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

<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>.

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	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
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)	:	2.49

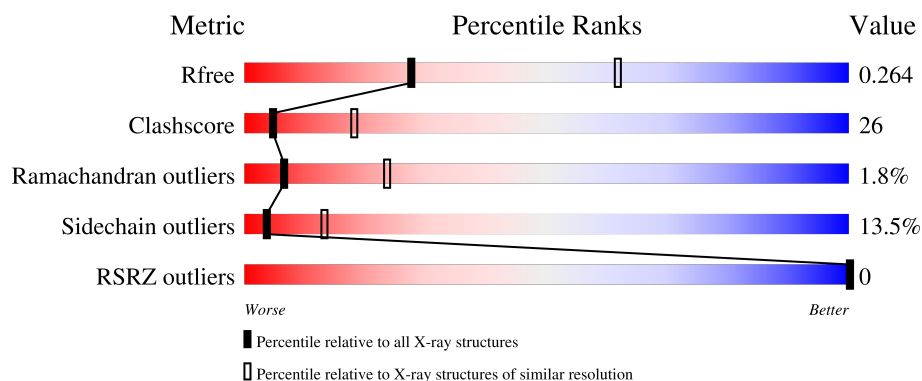
1 Overall quality at a glance

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}	180053	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 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	147	
1	B	147	
1	C	147	
1	D	147	

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			

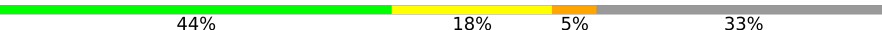
- Molecule 2 is water.

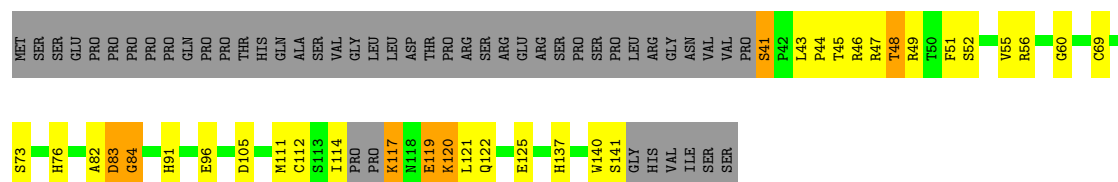
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	4	Total	O	0	0
			4	4		
2	C	8	Total	O	0	0
			8	8		
2	B	12	Total	O	0	0
			12	12		
2	D	8	Total	O	0	0
			8	8		

3 Residue-property plots [i](#)

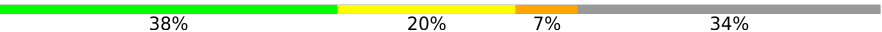
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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.

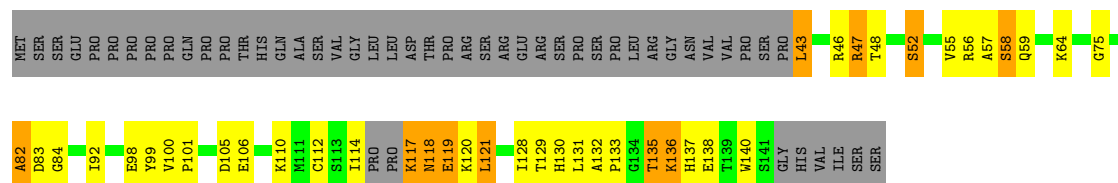
- Molecule 1: Calcium-regulated heat stable protein 1

Chain A: 



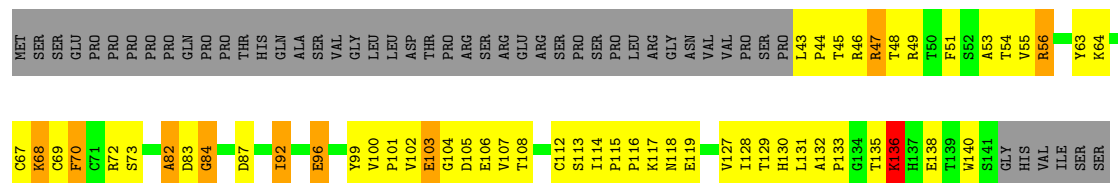
- Molecule 1: Calcium-regulated heat stable protein 1

Chain C: 

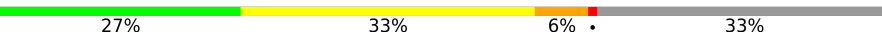


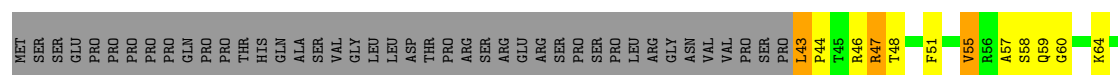
- Molecule 1: Calcium-regulated heat stable protein 1

Chain B: 



- Molecule 1: Calcium-regulated heat stable protein 1

Chain D: 





4 Data and refinement statistics

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α , β , γ	79.86Å 79.86Å 181.98Å 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 2.81Å)	Xtriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, R_{free}	0.229 , 0.272 0.228 , 0.264	Depositor DCC
R_{free} test set	812 reflections (5.03%)	wwPDB-VP
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 , 55.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	0.214 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.96	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, 0.333 respectively for untwinned datasets, and 0.375, 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	0.50	0/780	0.93	4/1056 (0.4%)
1	B	0.73	0/783	1.17	8/1063 (0.8%)
1	C	0.76	0/753	0.96	1/1019 (0.1%)
1	D	0.65	0/773	1.09	9/1049 (0.9%)
All	All	0.67	0/3089	1.04	22/4187 (0.5%)

There are no bond length outliers.

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	43	LEU	CA-C-N	9.33	129.46	120.31
1	B	43	LEU	C-N-CA	9.33	129.46	120.31
1	B	114	ILE	CA-C-N	9.03	129.68	120.38
1	B	114	ILE	C-N-CA	9.03	129.68	120.38
1	D	60	GLY	CA-C-N	6.55	126.52	120.03
1	D	60	GLY	C-N-CA	6.55	126.52	120.03
1	D	132	ALA	CA-C-N	6.51	126.09	119.19
1	D	132	ALA	C-N-CA	6.51	126.09	119.19
1	D	114	ILE	CA-C-N	6.45	124.30	119.66
1	D	114	ILE	C-N-CA	6.45	124.30	119.66
1	A	82	ALA	N-CA-C	-6.45	102.48	110.41
1	D	107	VAL	N-CA-C	5.87	117.13	108.45
1	A	60	GLY	CA-C-N	5.83	125.81	120.21
1	A	60	GLY	C-N-CA	5.83	125.81	120.21
1	B	44	PRO	N-CA-C	5.79	120.37	111.57
1	B	82	ALA	N-CA-C	-5.74	102.91	110.43
1	A	76	HIS	N-CA-C	5.61	116.77	108.86
1	B	53	ALA	N-CA-C	-5.59	105.09	111.07
1	D	85	GLY	CA-C-N	5.48	125.48	119.89
1	D	85	GLY	C-N-CA	5.48	125.48	119.89
1	B	131	LEU	CB-CA-C	5.44	118.72	109.53
1	C	82	ALA	N-CA-C	-5.25	103.10	110.50

There are no chirality outliers.

There are no planarity outliers.

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	762	0	752	24	0
1	B	763	0	755	43	1
1	C	737	0	722	47	1
1	D	754	0	741	49	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	154	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 26.

All (154) 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	1.16
1:C:47:ARG:HG2	1:C:47:ARG:NH1	1.52	1.05
1:C:47:ARG:HH11	1:C:47:ARG:CG	1.65	1.05
1:C:121:LEU:HD23	1:C:121:LEU:N	1.54	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:H	1:C:118:ASN:HD22	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	0.85
1:D:103:GLU:H	1:D:103:GLU:CD	1.84	0.84
1:C:118:ASN:C	1:C:119:GLU:CG	2.49	0.84
1:C:120:LYS:C	1:C:121:LEU:CD2	2.54	0.81

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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: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:ND2	1:C:118:ASN:N	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:52:SER:O	1:C:56:ARG:HG3	1.86	0.76
1:C:117:LYS:O	1:C:117:LYS:CE	2.30	0.75
1:C:117:LYS:O	1:C:117:LYS:CG	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:D:82:ALA:O	1:D:83:ASP:C	2.33	0.71
1:D:83:ASP:CG	1:D:84:GLY:N	2.47	0.69
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: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:B:103:GLU:CD	1:B:103:GLU:H	2.03	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:116:PRO:CD	1:D:118:ASN:N	2.61	0.63
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: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

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:69:CYS:O	1:B:70:PHE:CB	2.53	0.54
1:A:119:GLU:O	1:A:120:LYS:C	2.51	0.53
1:A:83:ASP:O	1:A:84:GLY:O	2.27	0.53
1:C:64:LYS:HE3	1:C:106:GLU:OE2	2.08	0.53
1:C:110:LYS:HE2	1:B:56:ARG:HH12	1.73	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:D:83:ASP:CG	1:D:84:GLY:H	2.17	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:B:115:PRO:HB3	1:B:116:PRO:HA	1.92	0.51
1:D:116:PRO:HG3	1:D:118:ASN:N	2.24	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: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:D:82:ALA:O	1:D:83:ASP:O	2.30	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:51:PHE:C	1:D:51:PHE:CD1	2.90	0.49
1:D:83:ASP:OD1	1:D:84:GLY:O	2.30	0.49
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: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:B:99:TYR:HD1	1:B:138:GLU:O	1.96	0.48
1:D:70:PHE:HA	1:D:77:GLY:HA3	1.96	0.48
1:B:63:TYR:O	1:B:108:THR:HA	2.13	0.47
1:B:102:VAL:HG13	1:B:103:GLU:CD	2.39	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:D:83:ASP:C	1:D:83:ASP:OD1	2.57	0.47
1:A:41:SER:O	1:A:43:LEU:HD23	2.15	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:128:ILE:HG21	1:C:131:LEU:CD2	2.44	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:A:44:PRO:HB2	1:A:122:GLN:HE22	1.80	0.45
1:D:105:ASP:OD1	1:D:137:HIS:CE1	2.69	0.45
1:C:43:LEU:C	1:C:43:LEU:HD22	2.40	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:C:57:ALA:C	1:C:59:GLN:H	2.24	0.45
1:B:102:VAL:CG1	1:B:103:GLU:N	2.80	0.44
1:A:111:MET:SD	1:A:121:LEU:HD13	2.57	0.44
1:C:46:ARG:NH1	1:D:93:SER:O	2.49	0.44
1:C:75:GLY:O	1:C:92:ILE:HG22	2.17	0.44
1:B:104:GLY:C	1:B:132:ALA:HB2	2.42	0.44
1:A:49:ARG:HD3	1:A:112:CYS:SG	2.58	0.44
1:B:56:ARG:HB2	1:B:56:ARG:NH1	2.33	0.44
1:B:136:LYS:HB3	1:B:136:LYS:HE3	1.62	0.44
1:D:64:LYS:HD3	1:D:106:GLU:OE2	2.17	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:C:58:SER:HB2	1:D:127:VAL:HG21	1.99	0.43
1:D:91:HIS:CE1	1:D:92:ILE:HG22	2.54	0.43
1:A:140:TRP:O	1:A:141:SER:OG	2.26	0.42
1:B:135:THR:HG22	1:B:135:THR:O	2.18	0.42
1:D:113:SER:OG	1:D:118:ASN:HB3	2.19	0.42
1:C:132:ALA:HA	1:C:133:PRO:HD3	1.86	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:B:115:PRO:HA	1:B:116:PRO:C	2.45	0.42
1:D:104:GLY:C	1:D:132:ALA:HB2	2.44	0.42
1:B:117:LYS:O	1:B:118:ASN:C	2.61	0.42
1:A:121:LEU:HD23	1:A:121:LEU:HA	1.85	0.42
1:C:131:LEU:HD22	1:C:137:HIS:CE1	2.55	0.42
1:B:96:GLU:H	1:B:96:GLU:HG2	1.69	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:55:VAL:O	1:D:59:GLN:HG3	2.21	0.41
1:D:72:ARG:NH1	1:D:140:TRP:HA	2.36	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: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
1:D:103:GLU:CD	1:D:103:GLU:N	2.63	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 [i](#)

5.3.1 Protein backbone [i](#)

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%)	5	20
1	B	97/147 (66%)	89 (92%)	5 (5%)	3 (3%)	3	12
1	C	93/147 (63%)	82 (88%)	10 (11%)	1 (1%)	11	36
1	D	94/147 (64%)	87 (93%)	6 (6%)	1 (1%)	11	36

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	379/588 (64%)	344 (91%)	28 (7%)	7 (2%)	6	23

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%)	78 (92%)	7 (8%)	10	33
1	B	85/129 (66%)	73 (86%)	12 (14%)	3	12
1	C	80/129 (62%)	67 (84%)	13 (16%)	2	8
1	D	84/129 (65%)	71 (84%)	13 (16%)	2	9
All	All	334/516 (65%)	289 (86%)	45 (14%)	4	13

All (45) 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	117	LYS
1	A	119	GLU
1	C	43	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	C	47	ARG
1	C	48	THR
1	C	52	SER
1	C	58	SER
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	45	THR
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	92	ILE
1	B	96	GLU
1	B	103	GLU
1	B	112	CYS
1	B	136	LYS
1	D	43	LEU
1	D	47	ARG
1	D	48	THR
1	D	55	VAL
1	D	68	LYS
1	D	73	SER
1	D	83	ASP
1	D	96	GLU
1	D	103	GLU
1	D	114	ILE
1	D	119	GLU
1	D	120	LYS
1	D	135	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	76	HIS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	C	59	GLN
1	C	76	HIS
1	C	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 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, 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%)	-1.26	0 100 100	50, 63, 88, 104	0
1	B	99/147 (67%)	-1.37	0 100 100	43, 58, 83, 95	0
1	C	97/147 (65%)	-1.32	0 100 100	45, 58, 84, 112	0
1	D	98/147 (66%)	-1.29	0 100 100	47, 62, 91, 103	0
All	All	393/588 (66%)	-1.31	0 100 100	43, 61, 88, 112	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.