



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 06:16 pm GMT

PDB ID : 1X23
Title : Crystal structure of ubch5c
Authors : Nakanishi, M.; Teshima, N.; Mizushima, T.; Murata, S.; Tanaka, K.; Yamane, T.
Deposited on : 2005-04-19
Resolution : 1.85 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
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

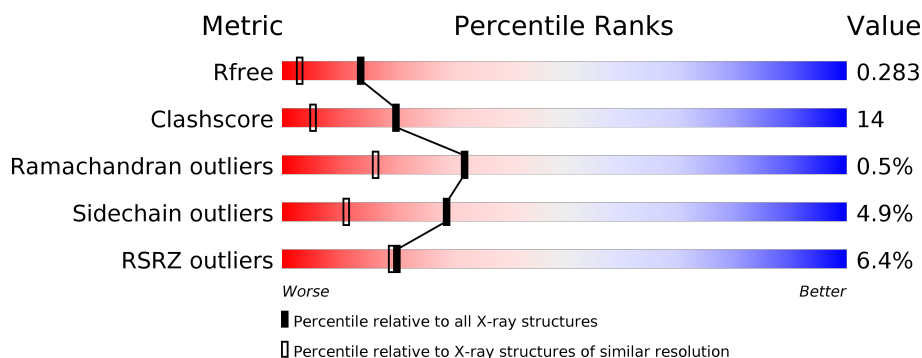
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.85 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	1923 (1.86-1.86)
Clashscore	112137	2083 (1.86-1.86)
Ramachandran outliers	110173	2060 (1.86-1.86)
Sidechain outliers	110143	2060 (1.86-1.86)
RSRZ outliers	101464	1932 (1.86-1.86)

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	155	
1	B	155	
1	C	155	
1	D	155	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5544 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 Ubiquitin-conjugating enzyme E2 D3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	153	Total	C	N	O	S	0	0	0
			1219	781	206	224	8			
1	B	152	Total	C	N	O	S	0	0	0
			1211	775	205	223	8			
1	C	152	Total	C	N	O	S	0	0	0
			1211	775	205	223	8			
1	D	154	Total	C	N	O	S	0	0	0
			1226	786	207	225	8			

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-8	GLY	-	CLONING ARTIFACT	UNP P61077
A	-7	PRO	-	CLONING ARTIFACT	UNP P61077
A	-6	LEU	-	CLONING ARTIFACT	UNP P61077
A	-5	GLY	-	CLONING ARTIFACT	UNP P61077
A	-4	SER	-	CLONING ARTIFACT	UNP P61077
A	-3	PRO	-	CLONING ARTIFACT	UNP P61077
A	-2	GLU	-	CLONING ARTIFACT	UNP P61077
A	-1	PHE	-	CLONING ARTIFACT	UNP P61077
B	-8	GLY	-	CLONING ARTIFACT	UNP P61077
B	-7	PRO	-	CLONING ARTIFACT	UNP P61077
B	-6	LEU	-	CLONING ARTIFACT	UNP P61077
B	-5	GLY	-	CLONING ARTIFACT	UNP P61077
B	-4	SER	-	CLONING ARTIFACT	UNP P61077
B	-3	PRO	-	CLONING ARTIFACT	UNP P61077
B	-2	GLU	-	CLONING ARTIFACT	UNP P61077
B	-1	PHE	-	CLONING ARTIFACT	UNP P61077
C	-8	GLY	-	CLONING ARTIFACT	UNP P61077
C	-7	PRO	-	CLONING ARTIFACT	UNP P61077
C	-6	LEU	-	CLONING ARTIFACT	UNP P61077
C	-5	GLY	-	CLONING ARTIFACT	UNP P61077
C	-4	SER	-	CLONING ARTIFACT	UNP P61077

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-3	PRO	-	CLONING ARTIFACT	UNP P61077
C	-2	GLU	-	CLONING ARTIFACT	UNP P61077
C	-1	PHE	-	CLONING ARTIFACT	UNP P61077
D	-8	GLY	-	CLONING ARTIFACT	UNP P61077
D	-7	PRO	-	CLONING ARTIFACT	UNP P61077
D	-6	LEU	-	CLONING ARTIFACT	UNP P61077
D	-5	GLY	-	CLONING ARTIFACT	UNP P61077
D	-4	SER	-	CLONING ARTIFACT	UNP P61077
D	-3	PRO	-	CLONING ARTIFACT	UNP P61077
D	-2	GLU	-	CLONING ARTIFACT	UNP P61077
D	-1	PHE	-	CLONING ARTIFACT	UNP P61077

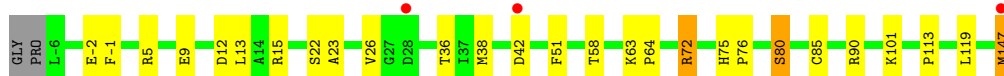
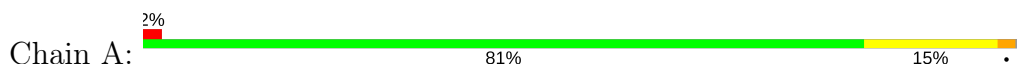
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	177	Total O 177 177	0	0
2	B	183	Total O 183 183	0	0
2	C	173	Total O 173 173	0	0
2	D	144	Total O 144 144	0	0

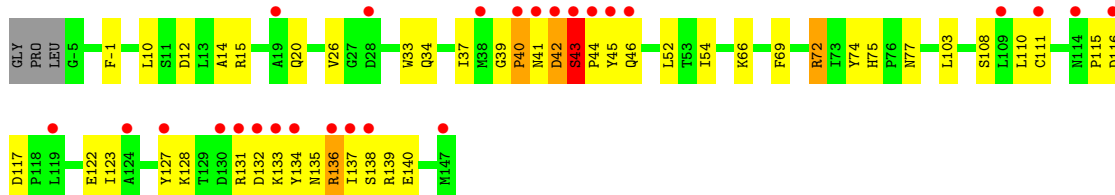
3 Residue-property plots

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.

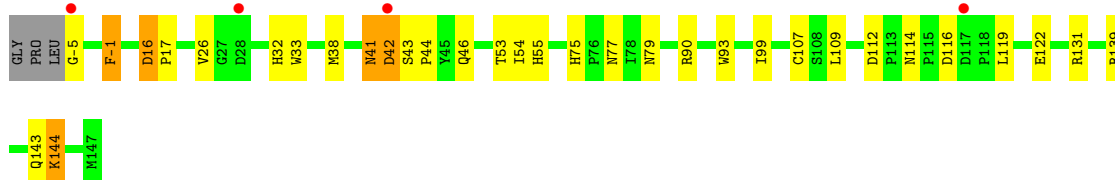
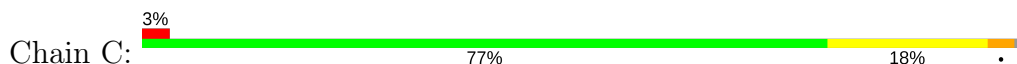
• Molecule 1: Ubiquitin-conjugating enzyme E2 D3



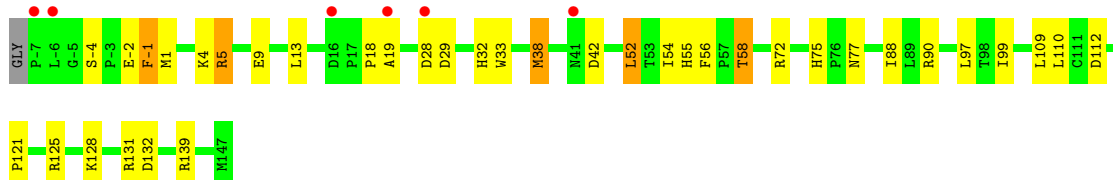
• Molecule 1: Ubiquitin-conjugating enzyme E2 D3



• Molecule 1: Ubiquitin-conjugating enzyme E2 D3



• Molecule 1: Ubiquitin-conjugating enzyme E2 D3



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	62.00Å 44.33Å 97.58Å 90.00° 93.36° 90.00°	Depositor
Resolution (Å)	20.00 – 1.85 19.95 – 1.85	Depositor EDS
% Data completeness (in resolution range)	94.7 (20.00-1.85) 94.7 (19.95-1.85)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.91 (at 1.85Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.199 , 0.282 0.199 , 0.283	Depositor DCC
R_{free} test set	2178 reflections (5.31%)	DCC
Wilson B-factor (Å ²)	23.6	Xtriage
Anisotropy	0.057	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 57.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5544	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 27.86 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 2.0403e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² 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](#)

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	0.71	1/1256 (0.1%)	0.75	0/1711
1	B	0.70	1/1248 (0.1%)	0.73	0/1700
1	C	0.69	1/1248 (0.1%)	0.70	1/1700 (0.1%)
1	D	0.78	1/1264 (0.1%)	0.70	3/1722 (0.2%)
All	All	0.72	4/5016 (0.1%)	0.72	4/6833 (0.1%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	-1	PHE	C-N	18.16	1.75	1.34
1	A	-1	PHE	C-N	13.27	1.64	1.34
1	B	-1	PHE	C-N	12.24	1.62	1.34
1	C	-1	PHE	C-N	12.10	1.61	1.34

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	-1	PHE	O-C-N	-6.60	112.15	122.70
1	D	52	LEU	CA-CB-CG	5.79	128.61	115.30
1	D	5	ARG	NE-CZ-NH1	5.07	122.83	120.30
1	D	5	ARG	NE-CZ-NH2	-5.02	117.79	120.30

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	1219	0	1202	23	0
1	B	1211	0	1191	52	0
1	C	1211	0	1191	32	0
1	D	1226	0	1211	33	0
2	A	177	0	0	11	0
2	B	183	0	0	27	0
2	C	173	0	0	10	0
2	D	144	0	0	17	0
All	All	5544	0	4795	135	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 14.

All (135) 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:D:-1:PHE:C	1:D:1:MET:N	1.75	1.37
1:D:-1:PHE:C	1:D:1:MET:H2	1.38	1.17
1:B:108:SER:HA	2:B:274:HOH:O	1.51	1.10
1:B:137:ILE:O	2:B:284:HOH:O	1.80	0.99
1:B:140:GLU:N	2:B:284:HOH:O	1.95	0.97
1:B:131:ARG:HD2	2:B:216:HOH:O	1.66	0.96
1:B:116:ASP:OD1	2:B:311:HOH:O	1.84	0.94
1:D:-1:PHE:C	1:D:1:MET:H1	1.52	0.94
1:D:75:HIS:HD2	1:D:77:ASN:H	1.17	0.93
1:A:90:ARG:HD3	2:A:285:HOH:O	1.67	0.93
1:A:63:LYS:HG3	1:A:64:PRO:HD2	1.52	0.90
1:A:72:ARG:NH1	2:A:282:HOH:O	2.05	0.89
1:B:136:ARG:O	2:B:285:HOH:O	1.89	0.89
1:B:72:ARG:CG	1:B:72:ARG:HH11	1.92	0.82
1:B:123:ILE:O	2:B:172:HOH:O	1.98	0.82
1:C:38:MET:HB2	2:D:200:HOH:O	1.80	0.80
1:C:107:CYS:SG	2:C:258:HOH:O	2.35	0.80
1:A:38:MET:HG3	2:B:311:HOH:O	1.82	0.79
1:D:-4:SER:OG	1:D:-2:GLU:HG2	1.83	0.78
1:B:140:GLU:OE2	1:D:28:ASP:HB2	1.86	0.76
1:C:75:HIS:HD2	1:C:77:ASN:H	1.32	0.75
1:D:90:ARG:NH2	2:D:213:HOH:O	2.20	0.75
1:B:122:GLU:OE2	1:D:55:HIS:HE1	1.69	0.74
1:B:139:ARG:O	2:B:286:HOH:O	2.06	0.74
1:B:39:GLY:O	2:B:323:HOH:O	2.06	0.72
1:B:140:GLU:HG2	2:D:210:HOH:O	1.88	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:132:ASP:O	2:B:301:HOH:O	2.07	0.71
1:C:139:ARG:O	1:C:143:GLN:HG3	1.91	0.71
1:B:75:HIS:HD2	1:B:77:ASN:H	1.38	0.70
1:C:77:ASN:HD21	1:C:114:ASN:H	1.38	0.69
1:A:26:VAL:HG12	1:D:72:ARG:NH2	2.08	0.68
1:B:72:ARG:HG3	1:B:72:ARG:HH11	1.59	0.68
1:D:55:HIS:HD2	2:D:230:HOH:O	1.77	0.67
1:A:15:ARG:NH1	2:A:180:HOH:O	2.20	0.66
1:C:41:ASN:O	1:C:46:GLN:NE2	2.30	0.65
1:B:42:ASP:N	1:B:43:SER:HA	2.12	0.64
1:D:4:LYS:HE2	2:D:148:HOH:O	1.98	0.63
1:D:1:MET:HG2	2:D:212:HOH:O	1.98	0.63
1:A:147:MET:OXT	2:A:290:HOH:O	2.15	0.63
1:B:39:GLY:HA3	1:B:45:TYR:HB2	1.81	0.63
1:C:38:MET:SD	1:D:128:LYS:HD3	2.39	0.63
1:B:133:LYS:C	2:B:301:HOH:O	2.36	0.62
1:C:77:ASN:ND2	1:C:114:ASN:H	1.97	0.62
1:B:134:TYR:HB3	2:B:216:HOH:O	2.00	0.62
1:A:85:CYS:HB2	1:A:119:LEU:HG	1.82	0.61
1:B:43:SER:H	1:B:44:PRO:CD	2.12	0.61
1:C:79:ASN:ND2	1:C:119:LEU:HB3	2.16	0.60
2:A:273:HOH:O	1:D:121:PRO:HG2	2.02	0.60
1:B:14:ALA:HB2	2:B:196:HOH:O	2.00	0.59
1:C:144:LYS:HA	1:C:144:LYS:HE3	1.82	0.59
1:B:111:CYS:HB2	2:B:274:HOH:O	2.03	0.59
1:B:75:HIS:CD2	1:B:77:ASN:H	2.19	0.59
1:B:40:PRO:HG2	1:B:111:CYS:SG	2.43	0.58
1:D:-1:PHE:C	1:D:1:MET:CA	2.71	0.58
1:C:38:MET:CE	2:D:200:HOH:O	2.52	0.58
1:B:12:ASP:OD2	1:B:15:ARG:NH1	2.37	0.57
1:B:132:ASP:C	2:B:301:HOH:O	2.43	0.56
1:C:90:ARG:NH1	2:C:204:HOH:O	2.39	0.55
1:D:75:HIS:CD2	1:D:77:ASN:H	2.09	0.55
1:A:38:MET:HE3	2:A:246:HOH:O	2.07	0.54
1:B:41:ASN:HA	2:B:246:HOH:O	2.07	0.54
1:B:72:ARG:HH11	1:B:72:ARG:HG2	1.69	0.54
1:B:108:SER:HB3	2:B:273:HOH:O	2.06	0.53
1:C:79:ASN:HD21	1:C:119:LEU:HB3	1.71	0.53
1:C:55:HIS:HD2	2:C:157:HOH:O	1.92	0.53
1:C:107:CYS:HB3	2:C:258:HOH:O	2.10	0.52
1:B:72:ARG:HG3	1:B:72:ARG:NH1	2.22	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:29:ASP:OD2	2:D:206:HOH:O	2.19	0.52
1:D:4:LYS:NZ	2:D:212:HOH:O	2.40	0.52
1:B:72:ARG:NH1	1:B:72:ARG:CG	2.62	0.51
1:C:42:ASP:HA	2:C:265:HOH:O	2.10	0.51
1:D:58:THR:HG23	2:D:154:HOH:O	2.09	0.51
1:B:44:PRO:HA	1:B:139:ARG:HG3	1.91	0.51
1:A:75:HIS:ND1	1:A:76:PRO:HD2	2.26	0.51
1:B:74:TYR:HB3	1:B:138:SER:HA	1.93	0.51
1:A:58:THR:HG22	2:A:314:HOH:O	2.11	0.51
1:A:72:ARG:NH2	2:A:273:HOH:O	2.43	0.51
1:B:135:ASN:N	2:B:301:HOH:O	2.43	0.51
1:B:131:ARG:HA	2:B:216:HOH:O	2.10	0.50
1:D:5:ARG:HD2	1:D:9:GLU:OE1	2.11	0.50
1:D:75:HIS:HE1	1:D:109:LEU:O	1.93	0.50
1:D:125:ARG:NE	2:D:269:HOH:O	2.37	0.50
1:C:75:HIS:CD2	1:C:77:ASN:H	2.21	0.50
1:A:26:VAL:HG12	1:D:72:ARG:HH22	1.76	0.49
1:C:107:CYS:CB	2:C:258:HOH:O	2.56	0.49
1:B:115:PRO:HB2	1:B:128:LYS:HE2	1.95	0.48
1:A:36:THR:HG22	1:A:51:PHE:CD1	2.48	0.48
1:D:139:ARG:NH1	2:D:235:HOH:O	2.45	0.48
1:D:13:LEU:HD23	1:D:18:PRO:HD3	1.96	0.48
1:A:80:SER:HB2	2:A:172:HOH:O	2.14	0.47
1:C:42:ASP:HB2	2:C:240:HOH:O	2.13	0.47
1:B:127:TYR:HA	2:B:216:HOH:O	2.14	0.47
1:B:127:TYR:N	2:B:172:HOH:O	2.47	0.47
1:B:33:TRP:HB2	1:B:54:ILE:HB	1.97	0.47
1:A:12:ASP:OD1	1:A:15:ARG:NH2	2.48	0.47
1:D:38:MET:HE3	2:D:258:HOH:O	2.14	0.47
1:B:128:LYS:CE	2:B:311:HOH:O	2.64	0.46
1:C:77:ASN:HD21	1:C:114:ASN:N	2.09	0.46
1:C:75:HIS:HE1	1:C:109:LEU:O	1.98	0.46
1:C:16:ASP:HA	1:C:17:PRO:HD2	1.77	0.45
1:B:136:ARG:HA	2:B:218:HOH:O	2.16	0.45
1:A:13:LEU:HD21	1:A:23:ALA:HB3	1.98	0.45
1:A:76:PRO:HG2	1:A:113:PRO:HB3	1.99	0.45
1:C:32:HIS:HB2	2:C:169:HOH:O	2.16	0.45
1:D:112:ASP:OD1	2:D:281:HOH:O	2.20	0.45
1:B:66:LYS:NZ	2:B:186:HOH:O	2.50	0.45
1:C:-5:GLY:HA2	1:C:-1:PHE:CE1	2.52	0.44
1:D:42:ASP:N	1:D:42:ASP:OD1	2.31	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:22:SER:OG	1:A:36:THR:OG1	2.35	0.44
1:B:41:ASN:H	1:B:42:ASP:CA	2.31	0.44
1:D:56:PHE:CE1	1:D:99:ILE:HD11	2.51	0.44
1:B:52:LEU:HD23	1:B:69:PHE:HA	1.98	0.44
1:B:42:ASP:HA	1:B:43:SER:OG	2.18	0.44
1:C:93:TRP:HB3	2:C:174:HOH:O	2.17	0.44
1:B:37:ILE:HD12	1:B:52:LEU:HD11	2.00	0.43
1:D:132:ASP:HB3	2:D:261:HOH:O	2.18	0.43
1:A:5:ARG:O	1:A:9:GLU:HG3	2.18	0.43
1:D:18:PRO:O	1:D:19:ALA:C	2.56	0.43
1:A:63:LYS:HE3	2:A:275:HOH:O	2.18	0.43
1:C:38:MET:HE3	2:D:200:HOH:O	2.16	0.43
1:C:33:TRP:HB2	1:C:54:ILE:HB	2.01	0.43
1:D:33:TRP:HB2	1:D:54:ILE:HB	2.00	0.43
1:C:53:THR:HG23	2:C:157:HOH:O	2.19	0.42
1:B:137:ILE:C	2:B:284:HOH:O	2.40	0.42
1:A:101:LYS:HG2	2:A:245:HOH:O	2.19	0.42
1:C:99:ILE:HA	1:C:99:ILE:HD13	1.82	0.42
1:B:26:VAL:CG2	1:B:26:VAL:O	2.67	0.41
1:B:20:GLN:NE2	2:B:206:HOH:O	2.51	0.41
1:C:38:MET:HE2	2:D:200:HOH:O	2.16	0.41
1:A:72:ARG:HD3	1:A:72:ARG:HA	1.67	0.41
1:C:43:SER:HB2	1:C:44:PRO:HD2	2.03	0.40
1:C:41:ASN:HD22	1:C:41:ASN:HA	1.68	0.40
1:B:103:LEU:HD23	1:B:103:LEU:HA	1.91	0.40
1:B:110:LEU:HD23	1:B:110:LEU:HA	1.84	0.40
1:D:88:ILE:HB	1:D:97:LEU:HD13	2.02	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	151/155 (97%)	146 (97%)	5 (3%)	0	100	100
1	B	150/155 (97%)	144 (96%)	3 (2%)	3 (2%)	9	1
1	C	150/155 (97%)	145 (97%)	5 (3%)	0	100	100
1	D	152/155 (98%)	146 (96%)	6 (4%)	0	100	100
All	All	603/620 (97%)	581 (96%)	19 (3%)	3 (0%)	32	16

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	43	SER
1	B	42	ASP
1	B	40	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	137/138 (99%)	132 (96%)	5 (4%)	40	20
1	B	136/138 (99%)	129 (95%)	7 (5%)	28	10
1	C	136/138 (99%)	127 (93%)	9 (7%)	19	5
1	D	138/138 (100%)	132 (96%)	6 (4%)	33	15
All	All	547/552 (99%)	520 (95%)	27 (5%)	29	11

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	-2	GLU
1	A	42	ASP
1	A	72	ARG
1	A	80	SER
1	A	147	MET
1	B	10	LEU
1	B	34	GLN
1	B	43	SER

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Mol	Chain	Res	Type
1	B	46	GLN
1	B	72	ARG
1	B	117	ASP
1	B	136	ARG
1	C	16	ASP
1	C	26	VAL
1	C	41	ASN
1	C	42	ASP
1	C	112	ASP
1	C	116	ASP
1	C	122	GLU
1	C	131	ARG
1	C	144	LYS
1	D	32	HIS
1	D	38	MET
1	D	52	LEU
1	D	58	THR
1	D	110	LEU
1	D	131	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	34	GLN
1	B	34	GLN
1	B	41	ASN
1	B	75	HIS
1	B	135	ASN
1	C	32	HIS
1	C	41	ASN
1	C	55	HIS
1	C	75	HIS
1	C	77	ASN
1	C	79	ASN
1	D	32	HIS
1	D	46	GLN
1	D	55	HIS
1	D	75	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

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, 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	153/155 (98%)	0.21	3 (1%) 65 65	14, 24, 37, 42	0
1	B	152/155 (98%)	0.75	26 (17%) 2 2	16, 27, 40, 45	0
1	C	152/155 (98%)	0.31	4 (2%) 56 55	16, 24, 39, 41	0
1	D	154/155 (99%)	0.31	6 (3%) 40 38	18, 25, 37, 47	0
All	All	611/620 (98%)	0.39	39 (6%) 20 19	14, 25, 39, 47	0

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	-5	GLY	5.5
1	C	42	ASP	4.8
1	B	43	SER	4.8
1	B	28	ASP	3.9
1	B	42	ASP	3.9
1	B	124	ALA	3.8
1	D	28	ASP	3.8
1	B	40	PRO	3.6
1	B	136	ARG	3.6
1	B	137	ILE	3.4
1	C	28	ASP	3.4
1	B	41	ASN	3.3
1	A	147	MET	3.1
1	A	42	ASP	3.0
1	B	134	TYR	2.9
1	D	19	ALA	2.9
1	D	-7	PRO	2.7
1	C	117	ASP	2.7
1	B	130	ASP	2.7
1	D	41	ASN	2.6
1	B	133	LYS	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	28	ASP	2.5
1	B	46	GLN	2.5
1	B	127	TYR	2.5
1	B	109	LEU	2.5
1	B	131	ARG	2.5
1	B	38	MET	2.4
1	B	147	MET	2.4
1	B	119	LEU	2.3
1	B	132	ASP	2.3
1	D	-6	LEU	2.2
1	B	111	CYS	2.2
1	B	114	ASN	2.2
1	B	45	TYR	2.2
1	B	138	SER	2.1
1	D	16	ASP	2.1
1	B	44	PRO	2.0
1	B	116	ASP	2.0
1	B	19	ALA	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.