Full wwPDB X-ray Structure Validation Report (i)

Aug 30, 2023 – 10:33 AM EDT

PDB ID	:	3NXZ
Title	:	Crystal Structure of UreE from Helicobacter pylori (Cu2+ bound form)
Authors	:	Shi, R.; Munger, C.; Assinas, A.; Matte, A.; Cygler, M.; Montreal-Kingston Bacterial Structural Genomics Initiative (BSGI)
Deposited on	:	2010-07-14
Resolution	:	2.70 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*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) 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 (1)) were used in the production of this report:

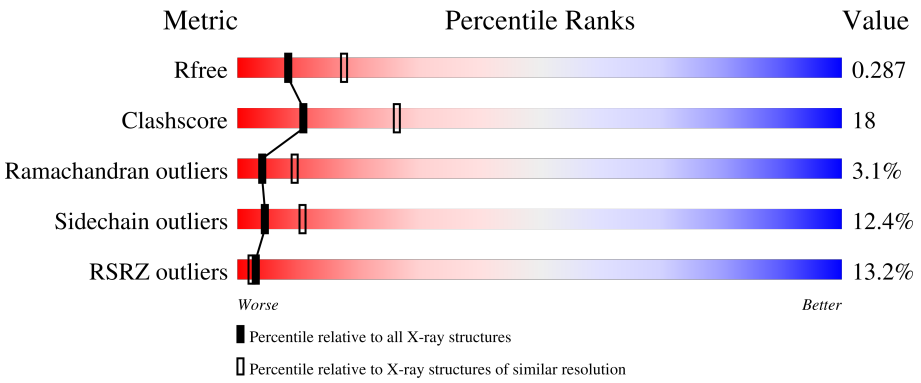
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.35
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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}	130704	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (2.70-2.70)

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 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 <=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	170	5% 51% 32% • • 12%
1	B	170	7% 51% 31% • • 14%
1	C	170	16% 51% 29% 7% • 13%
1	D	170	18% 54% 29% • • 12%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 4751 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 Urease accessory protein ureE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	149	1194	764	208	220	2	0	0	0
			Total	C	N	O	S			
1	B	147	1178	754	206	217	1	0	0	0
			Total	C	N	O	S			
1	C	148	1186	759	207	219	1	0	0	0
			Total	C	N	O	S			
1	D	149	1191	762	208	220	1	0	0	0

- Molecule 2 is COPPER (II) ION (three-letter code: CU) (formula: Cu).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Cu		
2	A	1	1	1	0	0

- Molecule 3 is water.

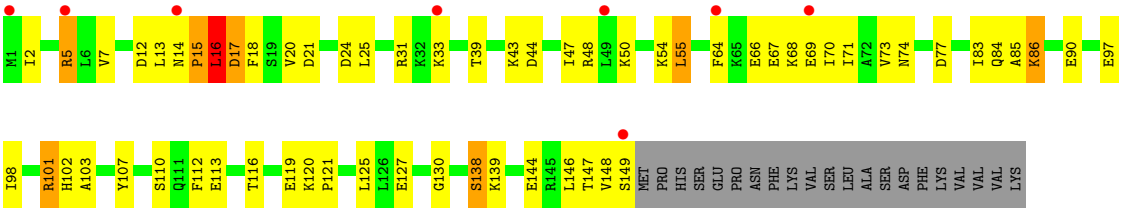
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	C	1	1	1	0	0

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.

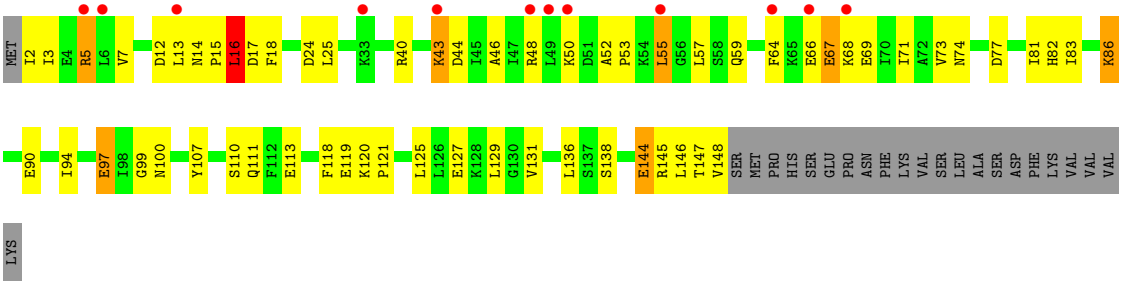
• Molecule 1: Urease accessory protein ureE

Chain A:



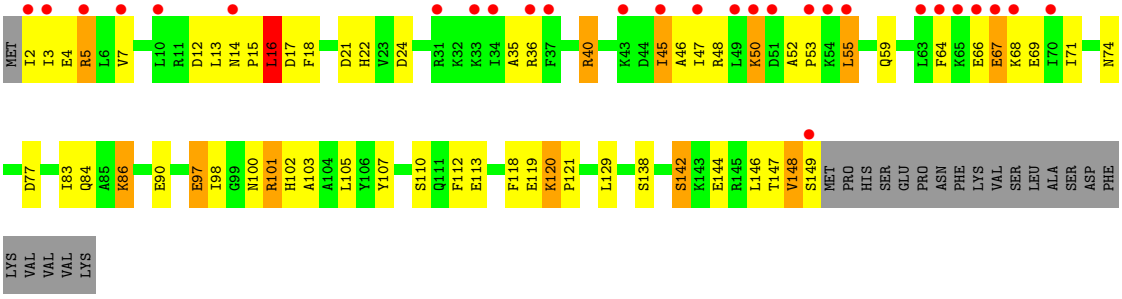
• Molecule 1: Urease accessory protein ureE

Chain B:



• Molecule 1: Urease accessory protein ureE

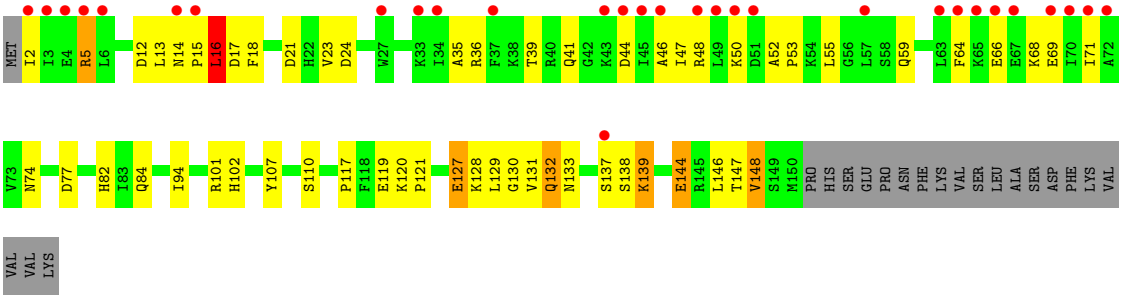
Chain C:



• Molecule 1: Urease accessory protein ureE

Chain D:

18%
54%
29%
12%


MET I2 I3 E4 R5 L6 D12 L13 N14 P15 L16 D17 F18 D21 H22 V23 D24 W27 K33 I34 A35 R36 F37 K38 T39 R40 Q41 G42 K43 D44 I45 A46 I47 R48 L49 K50 D51 A52 P53 K54 L55 G56 L57 S58 Q59 L63 F64 K65 E66 E67 K68 E69 I70 I71 A72
V73 N74 D77 H82 I83 Q84 I94 R101 H102 Y107 S110 P117 F118 E119 K120 P121 E127 K128 L129 G130 V131 Q132 N133 S137 S138 K139 E144 R145 L146 T147 V148 S149 M150 PRO HIS SER GLU PRO ASN PHE LYS VAL SER LEU ALA SER ASP PHE LYS VAL
VAL VAL LYS

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α, β, γ	88.83Å 88.83Å 203.63Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.17 – 2.70 44.17 – 2.70	Depositor EDS
% Data completeness (in resolution range)	95.9 (44.17-2.70) 95.9 (44.17-2.70)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.94 (at 2.69Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.256 , 0.294 0.246 , 0.287	Depositor DCC
R_{free} test set	1142 reflections (5.15%)	wwPDB-VP
Wilson B-factor (Å^2)	70.9	Xtriage
Anisotropy	0.743	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.30 , 72.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	4751	wwPDB-VP
Average B, all atoms (Å^2)	120.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.55% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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)

Bond lengths and bond angles in the following residue types are not validated in this section: CU

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	0.70	0/1212	0.83	0/1629
1	B	0.68	0/1196	0.79	0/1608
1	C	0.64	0/1204	0.82	1/1619 (0.1%)
1	D	0.64	0/1209	0.77	0/1626
All	All	0.67	0/4821	0.80	1/6482 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	129	LEU	CA-CB-CG	5.15	127.14	115.30

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	1194	0	1236	42	1
1	B	1178	0	1212	48	1
1	C	1186	0	1224	48	1
1	D	1191	0	1226	46	0
2	A	1	0	0	0	0
3	C	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	4751	0	4898	176	2

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 18.

All (176) 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:5:ARG:HH21	1:C:50:LYS:HB2	1.33	0.91
1:D:16:LEU:HA	1:D:17:ASP:HB3	1.54	0.90
1:D:84:GLN:HB3	1:D:132:GLN:HE21	1.34	0.90
1:C:24:ASP:OD1	1:C:74:ASN:HB3	1.71	0.88
1:C:16:LEU:HA	1:C:17:ASP:HB3	1.55	0.88
1:D:16:LEU:HA	1:D:17:ASP:CB	2.04	0.86
1:D:24:ASP:OD1	1:D:74:ASN:HB3	1.75	0.85
1:B:59:GLN:HE22	1:B:74:ASN:HA	1.40	0.84
1:C:59:GLN:HE22	1:C:74:ASN:HA	1.42	0.83
1:B:16:LEU:HA	1:B:17:ASP:HB3	1.64	0.78
1:C:3:ILE:HD13	1:C:45:ILE:HD11	1.68	0.75
1:D:2:ILE:CG2	1:D:48:ARG:HG3	2.16	0.75
1:D:16:LEU:CA	1:D:17:ASP:CB	2.65	0.74
1:C:16:LEU:HA	1:C:17:ASP:CB	2.16	0.74
1:A:24:ASP:OD1	1:A:74:ASN:HB3	1.87	0.74
1:C:16:LEU:CA	1:C:17:ASP:CB	2.65	0.73
1:B:16:LEU:HA	1:B:17:ASP:CB	2.19	0.71
1:D:84:GLN:CB	1:D:132:GLN:HE21	2.02	0.71
1:B:66:GLU:O	1:B:69:GLU:HB3	1.90	0.71
1:C:5:ARG:NH2	1:C:50:LYS:HB2	2.06	0.70
1:B:2:ILE:CG2	1:B:48:ARG:HG3	2.22	0.70
1:A:86:LYS:HG2	1:A:90:GLU:OE2	1.92	0.70
1:C:16:LEU:CA	1:C:17:ASP:HB3	2.21	0.70
1:A:16:LEU:N	1:A:17:ASP:HB2	2.09	0.68
1:B:97:GLU:OE2	1:C:148:VAL:HA	1.93	0.68
1:B:16:LEU:CA	1:B:17:ASP:CB	2.71	0.68
1:B:5:ARG:HH21	1:B:50:LYS:HB2	1.58	0.68
1:B:18:PHE:HD1	1:B:68:LYS:HA	1.57	0.68
1:A:64:PHE:HB3	1:A:71:ILE:HB	1.76	0.67
1:B:24:ASP:OD1	1:B:74:ASN:HB3	1.94	0.67
1:A:5:ARG:HH21	1:A:50:LYS:HB2	1.60	0.66
1:A:2:ILE:HG21	1:A:48:ARG:HG3	1.75	0.66
1:C:13:LEU:HA	1:C:14:ASN:HB2	1.78	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:18:PHE:HD1	1:D:68:LYS:HA	1.62	0.65
1:A:2:ILE:CG2	1:A:48:ARG:HG3	2.26	0.65
1:C:55:LEU:O	1:C:142:SER:HB3	1.96	0.65
1:D:5:ARG:HH21	1:D:50:LYS:HB2	1.60	0.65
1:A:97:GLU:HG2	1:A:125:LEU:HD11	1.78	0.64
1:A:16:LEU:HA	1:A:17:ASP:CB	2.28	0.64
1:B:25:LEU:HD12	1:B:73:VAL:CG1	2.28	0.64
1:A:149:SER:HB3	1:D:128:LYS:NZ	2.14	0.63
1:A:18:PHE:HD1	1:A:68:LYS:HA	1.64	0.62
1:B:97:GLU:OE2	1:C:149:SER:N	2.28	0.62
1:C:21:ASP:O	1:C:71:ILE:HA	1.99	0.62
1:C:2:ILE:CG2	1:C:48:ARG:HG3	2.30	0.62
1:A:16:LEU:CA	1:A:17:ASP:CB	2.78	0.61
1:D:107:TYR:CZ	1:D:146:LEU:HD21	2.35	0.61
1:D:13:LEU:HA	1:D:14:ASN:HB2	1.80	0.61
1:D:2:ILE:HG21	1:D:48:ARG:HG3	1.83	0.60
1:D:64:PHE:HB3	1:D:71:ILE:HB	1.83	0.60
1:B:13:LEU:HA	1:B:14:ASN:HB2	1.83	0.60
1:C:18:PHE:HD1	1:C:68:LYS:HA	1.65	0.60
1:C:16:LEU:N	1:C:17:ASP:CB	2.65	0.60
1:B:2:ILE:HD13	1:B:46:ALA:HB3	1.84	0.59
1:A:55:LEU:H	1:A:55:LEU:HD12	1.69	0.57
1:B:16:LEU:CA	1:B:17:ASP:HB3	2.35	0.57
1:B:64:PHE:HB3	1:B:71:ILE:HB	1.87	0.56
1:C:3:ILE:HD13	1:C:45:ILE:CD1	2.34	0.56
1:C:16:LEU:N	1:C:17:ASP:HB2	2.20	0.56
1:D:16:LEU:N	1:D:17:ASP:HB2	2.20	0.56
1:A:33:LYS:O	1:A:47:ILE:HD12	2.06	0.56
1:A:97:GLU:OE2	1:D:148:VAL:HA	2.06	0.56
1:B:18:PHE:HB3	1:B:68:LYS:O	2.05	0.56
1:C:66:GLU:O	1:C:69:GLU:HB3	2.05	0.55
1:D:94:ILE:HD11	1:D:131:VAL:HG11	1.87	0.55
1:B:16:LEU:N	1:B:17:ASP:HB2	2.22	0.55
1:B:146:LEU:N	1:B:146:LEU:HD23	2.21	0.55
1:B:18:PHE:CD1	1:B:68:LYS:HA	2.41	0.55
1:A:18:PHE:CD1	1:A:68:LYS:HA	2.41	0.55
1:C:84:GLN:HE22	1:C:112:PHE:HB2	1.72	0.55
1:A:16:LEU:CA	1:A:17:ASP:HB2	2.38	0.54
1:C:98:ILE:HD13	1:C:105:LEU:HD13	1.90	0.53
1:D:107:TYR:O	1:D:144:GLU:HG2	2.09	0.53
1:A:13:LEU:HA	1:A:14:ASN:HB2	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:66:GLU:O	1:A:69:GLU:HB3	2.09	0.53
1:D:39:THR:HG21	1:D:71:ILE:HD13	1.90	0.52
1:D:16:LEU:CA	1:D:17:ASP:HB2	2.40	0.52
1:A:16:LEU:HA	1:A:17:ASP:HB3	1.92	0.52
1:B:83:ILE:O	1:B:113:GLU:HA	2.08	0.52
1:C:64:PHE:HB3	1:C:71:ILE:HB	1.91	0.52
1:C:83:ILE:O	1:C:113:GLU:HA	2.10	0.52
1:B:16:LEU:N	1:B:17:ASP:CB	2.73	0.52
1:B:94:ILE:HD11	1:B:131:VAL:HG11	1.91	0.52
1:A:21:ASP:O	1:A:71:ILE:HA	2.09	0.52
1:D:35:ALA:O	1:D:47:ILE:HG13	2.11	0.51
1:C:16:LEU:N	1:C:17:ASP:HB3	2.25	0.51
1:D:21:ASP:O	1:D:71:ILE:HA	2.10	0.51
1:D:59:GLN:HE22	1:D:74:ASN:HA	1.76	0.51
1:D:82:HIS:O	1:D:133:ASN:HA	2.10	0.51
1:C:59:GLN:NE2	1:C:74:ASN:HA	2.19	0.51
1:D:18:PHE:CD1	1:D:68:LYS:HA	2.43	0.51
1:D:127:GLU:O	1:D:130:GLY:N	2.40	0.50
1:B:59:GLN:NE2	1:B:74:ASN:HA	2.19	0.50
1:D:84:GLN:HB3	1:D:132:GLN:NE2	2.16	0.50
1:A:98:ILE:HG23	1:A:103:ALA:HB3	1.94	0.50
1:C:118:PHE:C	1:C:118:PHE:CD1	2.85	0.50
1:D:59:GLN:OE1	1:D:139:LYS:HD3	2.12	0.49
1:B:55:LEU:H	1:B:55:LEU:HD12	1.75	0.49
1:C:86:LYS:HG2	1:C:90:GLU:OE2	2.12	0.49
1:A:84:GLN:HE22	1:A:112:PHE:HB2	1.77	0.49
1:B:52:ALA:O	1:B:53:PRO:C	2.50	0.49
1:B:118:PHE:C	1:B:118:PHE:CD1	2.85	0.49
1:D:16:LEU:CA	1:D:17:ASP:HB3	2.33	0.49
1:B:148:VAL:HA	1:C:97:GLU:OE2	2.13	0.48
1:A:18:PHE:HB3	1:A:68:LYS:O	2.12	0.48
1:B:25:LEU:HD12	1:B:73:VAL:HG11	1.95	0.48
1:A:138:SER:OG	1:A:139:LYS:N	2.43	0.48
1:B:16:LEU:HA	1:B:18:PHE:N	2.29	0.48
1:D:66:GLU:O	1:D:69:GLU:HB3	2.13	0.48
1:A:83:ILE:O	1:A:113:GLU:HA	2.14	0.48
1:D:17:ASP:HB3	1:D:18:PHE:CD2	2.48	0.48
1:C:17:ASP:HB3	1:C:18:PHE:CD2	2.49	0.48
1:C:52:ALA:O	1:C:53:PRO:C	2.52	0.48
1:D:120:LYS:N	1:D:121:PRO:HD2	2.29	0.48
1:C:4:GLU:OE1	1:C:50:LYS:HG2	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:120:LYS:N	1:C:121:PRO:HD2	2.29	0.47
1:D:16:LEU:N	1:D:17:ASP:CB	2.78	0.47
1:B:120:LYS:N	1:B:121:PRO:HD2	2.30	0.46
1:A:2:ILE:HG21	1:A:48:ARG:CG	2.44	0.46
1:D:101:ARG:O	1:D:102:HIS:HB2	2.14	0.46
1:C:36:ARG:NH2	1:C:46:ALA:HB2	2.30	0.46
1:A:16:LEU:N	1:A:17:ASP:CB	2.78	0.46
1:A:85:ALA:O	1:A:112:PHE:HB3	2.16	0.46
1:D:52:ALA:O	1:D:53:PRO:C	2.53	0.45
1:D:77:ASP:OD1	1:D:139:LYS:HE2	2.17	0.45
1:B:57:LEU:HD22	1:B:73:VAL:HG21	1.99	0.45
1:D:36:ARG:NH2	1:D:46:ALA:HB2	2.32	0.45
1:B:25:LEU:HD12	1:B:73:VAL:HG13	1.98	0.45
1:C:3:ILE:CD1	1:C:45:ILE:HD11	2.44	0.45
1:C:98:ILE:HG23	1:C:103:ALA:HB3	1.99	0.45
1:A:120:LYS:N	1:A:121:PRO:HD2	2.33	0.44
1:C:67:GLU:HB2	1:C:68:LYS:H	1.55	0.44
1:D:107:TYR:CZ	1:D:146:LEU:CD2	2.99	0.44
1:C:64:PHE:CE2	1:C:66:GLU:HB2	2.53	0.44
1:B:107:TYR:CE2	1:B:146:LEU:CD2	3.01	0.44
1:B:97:GLU:HG2	1:B:125:LEU:HD11	1.99	0.44
1:D:107:TYR:CE2	1:D:146:LEU:CD2	3.00	0.44
1:D:14:ASN:N	1:D:15:PRO:HD3	2.33	0.43
1:A:101:ARG:O	1:A:102:HIS:HB2	2.18	0.43
1:D:21:ASP:HB3	1:D:71:ILE:HG12	2.00	0.43
1:B:111:GLN:HG2	1:B:111:GLN:O	2.18	0.43
1:B:2:ILE:HD13	1:B:46:ALA:CB	2.47	0.43
1:A:39:THR:OG1	1:A:43:LYS:HB2	2.18	0.43
1:B:2:ILE:HG21	1:B:48:ARG:HG3	1.99	0.43
1:D:21:ASP:OD1	1:D:41:GLN:HG3	2.18	0.43
1:D:107:TYR:CE2	1:D:146:LEU:HD21	2.54	0.43
1:A:15:PRO:C	1:A:17:ASP:HB2	2.38	0.43
1:A:5:ARG:NH2	1:A:50:LYS:HB2	2.32	0.42
1:D:2:ILE:HD13	1:D:46:ALA:HB3	2.01	0.42
1:A:127:GLU:O	1:A:130:GLY:N	2.46	0.42
1:C:84:GLN:NE2	1:C:112:PHE:HB2	2.35	0.42
1:A:25:LEU:HD12	1:A:73:VAL:HG13	2.02	0.42
1:A:17:ASP:HB3	1:A:18:PHE:CD2	2.55	0.42
1:B:16:LEU:HA	1:B:18:PHE:H	1.85	0.42
1:C:101:ARG:O	1:C:102:HIS:HB2	2.18	0.42
1:C:107:TYR:CZ	1:C:146:LEU:CD2	3.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:77:ASP:OD1	1:A:139:LYS:HE2	2.20	0.42
1:A:107:TYR:CE2	1:A:146:LEU:CD2	3.02	0.42
1:A:16:LEU:HA	1:A:18:PHE:N	2.35	0.42
1:B:99:GLY:O	1:C:100:ASN:HA	2.19	0.42
1:B:144:GLU:OE2	1:B:144:GLU:HA	2.19	0.42
1:C:14:ASN:N	1:C:15:PRO:HD3	2.35	0.42
1:C:35:ALA:O	1:C:47:ILE:HG13	2.20	0.42
1:A:20:VAL:HG13	1:A:70:ILE:HG22	2.02	0.41
1:B:86:LYS:HG2	1:B:90:GLU:OE2	2.20	0.41
1:C:22:HIS:O	1:C:40:ARG:HB2	2.20	0.41
1:B:81:ILE:HG13	1:B:118:PHE:HA	2.03	0.41
1:B:82:HIS:CE1	1:B:136:LEU:HD12	2.56	0.41
1:C:16:LEU:HA	1:C:18:PHE:N	2.35	0.41
1:D:23:VAL:HG22	1:D:39:THR:HG22	2.01	0.41
1:B:15:PRO:C	1:B:16:LEU:HD12	2.41	0.40
1:B:67:GLU:HB2	1:B:68:LYS:H	1.49	0.40
1:B:100:ASN:OD1	1:C:102:HIS:ND1	2.49	0.40
1:A:149:SER:HB3	1:D:128:LYS:HZ3	1.86	0.40
1:B:82:HIS:HD2	1:B:113:GLU:OE1	2.04	0.40
1:C:2:ILE:HG21	1:C:48:ARG:HG3	2.04	0.40

All (2) 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:B:77:ASP:OD2	1:C:77:ASP:OD2[4_454]	1.90	0.30
1:A:77:ASP:OD2	1:A:77:ASP:OD2[7_555]	2.17	0.03

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	147/170 (86%)	130 (88%)	12 (8%)	5 (3%)	3 8
1	B	145/170 (85%)	132 (91%)	8 (6%)	5 (3%)	3 8
1	C	146/170 (86%)	130 (89%)	11 (8%)	5 (3%)	3 8
1	D	147/170 (86%)	131 (89%)	13 (9%)	3 (2%)	7 19
All	All	585/680 (86%)	523 (89%)	44 (8%)	18 (3%)	4 9

All (18) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	147	THR
1	D	148	VAL
1	A	16	LEU
1	B	16	LEU
1	C	16	LEU
1	C	147	THR
1	D	16	LEU
1	A	147	THR
1	B	147	THR
1	C	148	VAL
1	A	67	GLU
1	A	148	VAL
1	B	67	GLU
1	B	43	LYS
1	C	67	GLU
1	B	7	VAL
1	A	7	VAL
1	C	7	VAL

5.3.2 Protein sidechains (i)

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	130/152 (86%)	114 (88%)	16 (12%)	4 11
1	B	127/152 (84%)	110 (87%)	17 (13%)	4 9

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	129/152 (85%)	113 (88%)	16 (12%)	4	11
1	D	129/152 (85%)	114 (88%)	15 (12%)	5	12
All	All	515/608 (85%)	451 (88%)	64 (12%)	4	11

All (64) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	ARG
1	A	12	ASP
1	A	15	PRO
1	A	16	LEU
1	A	17	ASP
1	A	31	ARG
1	A	44	ASP
1	A	54	LYS
1	A	55	LEU
1	A	86	LYS
1	A	101	ARG
1	A	110	SER
1	A	116	THR
1	A	119	GLU
1	A	138	SER
1	A	144	GLU
1	B	3	ILE
1	B	5	ARG
1	B	12	ASP
1	B	16	LEU
1	B	40	ARG
1	B	43	LYS
1	B	44	ASP
1	B	55	LEU
1	B	86	LYS
1	B	97	GLU
1	B	110	SER
1	B	119	GLU
1	B	127	GLU
1	B	129	LEU
1	B	138	SER
1	B	144	GLU
1	B	145	ARG
1	C	5	ARG

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Mol	Chain	Res	Type
1	C	12	ASP
1	C	16	LEU
1	C	40	ARG
1	C	45	ILE
1	C	50	LYS
1	C	55	LEU
1	C	86	LYS
1	C	97	GLU
1	C	101	ARG
1	C	110	SER
1	C	119	GLU
1	C	120	LYS
1	C	138	SER
1	C	142	SER
1	C	144	GLU
1	D	5	ARG
1	D	12	ASP
1	D	16	LEU
1	D	44	ASP
1	D	55	LEU
1	D	110	SER
1	D	117	PRO
1	D	119	GLU
1	D	127	GLU
1	D	129	LEU
1	D	132	GLN
1	D	137	SER
1	D	138	SER
1	D	139	LYS
1	D	144	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	14	ASN
1	A	82	HIS
1	A	84	GLN
1	B	82	HIS
1	B	84	GLN
1	C	82	HIS
1	C	84	GLN
1	D	14	ASN

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Mol	Chain	Res	Type
1	D	82	HIS
1	D	84	GLN
1	D	132	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	149/170 (87%)	0.49	8 (5%)	25	24	54, 96, 159, 187	0
1	B	147/170 (86%)	0.57	12 (8%)	11	9	59, 100, 182, 211	0
1	C	148/170 (87%)	0.89	28 (18%)	1	0	59, 103, 208, 240	0
1	D	149/170 (87%)	1.14	30 (20%)	1	0	62, 119, 221, 241	0
All	All	593/680 (87%)	0.77	78 (13%)	3	2	54, 104, 208, 241	0

All (78) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	70	ILE	10.6
1	D	45	ILE	7.4
1	D	64	PHE	7.1
1	C	63	LEU	6.8
1	D	44	ASP	6.3
1	C	37	PHE	5.7
1	D	37	PHE	5.6
1	C	33	LYS	5.1
1	D	2	ILE	5.1
1	D	46	ALA	4.9
1	D	3	ILE	4.8
1	C	14	ASN	4.7
1	D	6	LEU	4.5
1	C	36	ARG	4.4
1	C	50	LYS	4.3
1	C	43	LYS	4.3
1	D	14	ASN	4.1
1	C	68	LYS	4.0
1	D	72	ALA	4.0
1	D	71	ILE	3.9
1	C	64	PHE	3.8

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Mol	Chain	Res	Type	RSRZ
1	C	54	LYS	3.8
1	B	66	GLU	3.7
1	D	48	ARG	3.6
1	C	70	ILE	3.5
1	C	7	VAL	3.4
1	D	43	LYS	3.4
1	C	3	ILE	3.3
1	A	149	SER	3.2
1	D	63	LEU	3.2
1	D	33	LYS	3.2
1	C	51	ASP	3.2
1	C	49	LEU	3.2
1	D	5	ARG	3.1
1	D	50	LYS	3.1
1	C	66	GLU	3.1
1	D	34	ILE	3.0
1	D	69	GLU	3.0
1	C	10	LEU	3.0
1	B	48	ARG	3.0
1	B	6	LEU	2.9
1	C	45	ILE	2.9
1	A	1	MET	2.9
1	C	55	LEU	2.9
1	D	65	LYS	2.8
1	B	33	LYS	2.8
1	C	2	ILE	2.8
1	D	51	ASP	2.8
1	B	43	LYS	2.7
1	B	64	PHE	2.7
1	D	49	LEU	2.7
1	B	50	LYS	2.7
1	C	53	PRO	2.7
1	B	49	LEU	2.7
1	D	66	GLU	2.7
1	A	14	ASN	2.6
1	D	67	GLU	2.6
1	B	68	LYS	2.6
1	D	137	SER	2.5
1	B	13	LEU	2.5
1	A	33	LYS	2.5
1	D	15	PRO	2.5
1	A	69	GLU	2.5

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Mol	Chain	Res	Type	RSRZ
1	C	47	ILE	2.4
1	A	64	PHE	2.4
1	C	5	ARG	2.4
1	C	31	ARG	2.4
1	B	55	LEU	2.4
1	C	149	SER	2.4
1	C	34	ILE	2.3
1	B	5	ARG	2.2
1	A	5	ARG	2.2
1	C	65	LYS	2.2
1	A	49	LEU	2.2
1	D	4	GLU	2.2
1	D	27	TRP	2.1
1	D	57	LEU	2.0
1	C	67	GLU	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 monosaccharides in this entry.

6.4 Ligands (i)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	CU	A	200	1/1	0.98	0.02	68,68,68,68	0

6.5 Other polymers (i)

There are no such residues in this entry.