Full wwPDB X-ray Structure Validation Report (i)

Jun 15, 2024 – 06:23 PM EDT

PDB ID	:	1XS0
Title	:	Structure of the E. coli Ivy protein
Authors	:	Abergel, C.; Monchois, V.; Byrn, D.; Lazzaroni, J.C.; Claverie, J.M.
Deposited on	:	2004-10-18
Resolution	:	1.58 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:

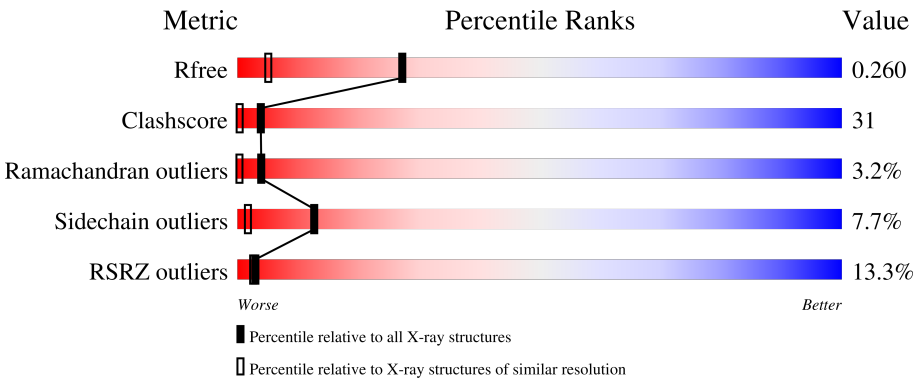
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

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 1.58 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	5534 (1.60-1.56)
Clashscore	141614	5861 (1.60-1.56)
Ramachandran outliers	138981	5708 (1.60-1.56)
Sidechain outliers	138945	5703 (1.60-1.56)
RSRZ outliers	127900	5431 (1.60-1.56)

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	136	12%; 64%; 24%; 6%; 6%
1	B	136	11%; 56%; 28%; 7%; •; 7%
1	C	136	15%; 56%; 30%; 9%; 5%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 3244 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 Inhibitor of vertebrate lysozyme.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	128	978	612	165	194	7	0	0	0
1	B	127	976	611	165	193	7	0	0	0
1	C	129	988	618	168	195	7	0	0	0

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	129	ARG	-	EXPRESSION TAG	UNP P45502
A	130	SER	-	EXPRESSION TAG	UNP P45502
A	131	HIS	-	EXPRESSION TAG	UNP P45502
A	132	HIS	-	EXPRESSION TAG	UNP P45502
A	133	HIS	-	EXPRESSION TAG	UNP P45502
A	134	HIS	-	EXPRESSION TAG	UNP P45502
A	135	HIS	-	EXPRESSION TAG	UNP P45502
A	136	HIS	-	EXPRESSION TAG	UNP P45502
B	129	ARG	-	EXPRESSION TAG	UNP P45502
B	130	SER	-	EXPRESSION TAG	UNP P45502
B	131	HIS	-	EXPRESSION TAG	UNP P45502
B	132	HIS	-	EXPRESSION TAG	UNP P45502
B	133	HIS	-	EXPRESSION TAG	UNP P45502
B	134	HIS	-	EXPRESSION TAG	UNP P45502
B	135	HIS	-	EXPRESSION TAG	UNP P45502
B	136	HIS	-	EXPRESSION TAG	UNP P45502
C	129	ARG	-	EXPRESSION TAG	UNP P45502
C	130	SER	-	EXPRESSION TAG	UNP P45502
C	131	HIS	-	EXPRESSION TAG	UNP P45502
C	132	HIS	-	EXPRESSION TAG	UNP P45502
C	133	HIS	-	EXPRESSION TAG	UNP P45502
C	134	HIS	-	EXPRESSION TAG	UNP P45502
C	135	HIS	-	EXPRESSION TAG	UNP P45502

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Chain	Residue	Modelled	Actual	Comment	Reference
C	136	HIS	-	EXPRESSION TAG	UNP P45502

- Molecule 2 is water.

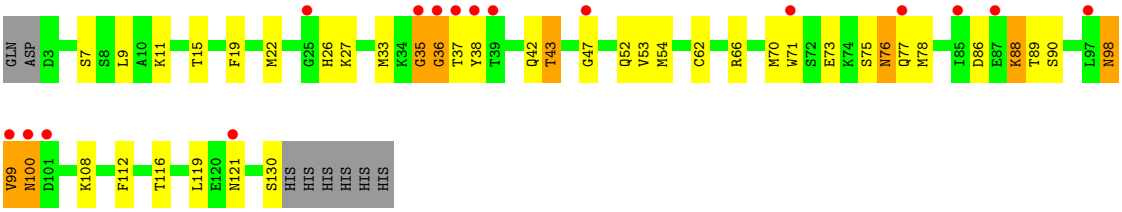
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	104	Total O 104 104	0	0
2	B	101	Total O 101 101	0	0
2	C	97	Total O 97 97	0	0

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.

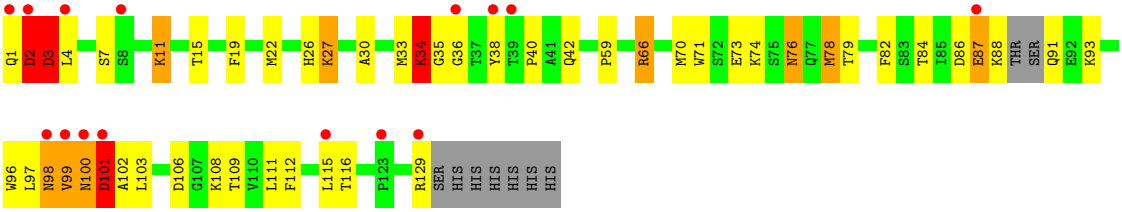
• Molecule 1: Inhibitor of vertebrate lysozyme

Chain A:



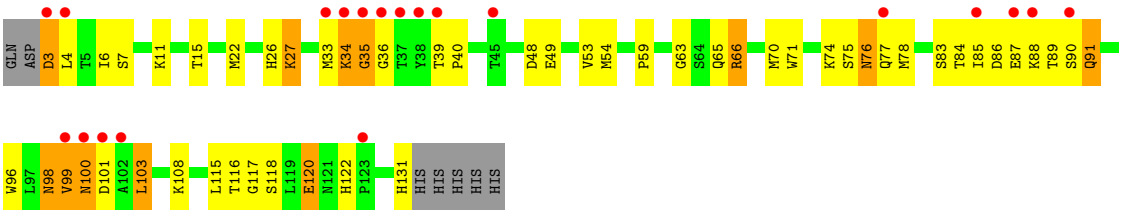
• Molecule 1: Inhibitor of vertebrate lysozyme

Chain B:



• Molecule 1: Inhibitor of vertebrate lysozyme

Chain C:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	81.34Å 46.96Å 88.04Å 90.00° 89.95° 90.00°	Depositor
Resolution (Å)	14.52 – 1.58 19.82 – 1.58	Depositor EDS
% Data completeness (in resolution range)	97.3 (14.52-1.58) 97.2 (19.82-1.58)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.04	Depositor
$< I/\sigma(I) >$ [1]	3.10 (at 1.58Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.225 , 0.260 0.226 , 0.260	Depositor DCC
R_{free} test set	4500 reflections (10.09%)	wwPDB-VP
Wilson B-factor (Å^2)	22.7	Xtriage
Anisotropy	0.273	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 46.9	EDS
L-test for twinning[2]	$< \|L\| > = 0.51$, $< L^2 > = 0.34$	Xtriage
Estimated twinning fraction	0.020 for -1/2*h-3/2*k,-1/2*h+1/2*k,-l 0.022 for -1/2*h+3/2*k,1/2*h+1/2*k,-l 0.477 for 1/2*h-3/2*k,-1/2*h-1/2*k,-l 0.477 for 1/2*h+3/2*k,1/2*h-1/2*k,-l 0.024 for -h,-k,l	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	3244	wwPDB-VP
Average B, all atoms (Å^2)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.28% 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)

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	0.93	0/997	1.01	1/1350 (0.1%)
1	B	0.95	1/994 (0.1%)	1.13	4/1344 (0.3%)
1	C	0.96	2/1008 (0.2%)	1.01	4/1365 (0.3%)
All	All	0.95	3/2999 (0.1%)	1.05	9/4059 (0.2%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	22	MET	SD-CE	-8.61	1.29	1.77
1	B	78	MET	SD-CE	-5.67	1.46	1.77
1	C	120	GLU	CG-CD	5.61	1.60	1.51

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	101	ASP	N-CA-C	-11.20	80.77	111.00
1	B	2	ASP	N-CA-C	7.47	131.18	111.00
1	B	4	LEU	CA-CB-CG	-6.20	101.04	115.30
1	A	108	LYS	CD-CE-NZ	-6.16	97.53	111.70
1	C	78	MET	CA-CB-CG	6.06	123.60	113.30
1	B	78	MET	CG-SD-CE	5.67	109.27	100.20
1	C	78	MET	CB-CG-SD	-5.26	96.63	112.40
1	C	99	VAL	N-CA-C	-5.13	97.15	111.00
1	C	103	LEU	CB-CG-CD2	-5.09	102.36	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	978	0	954	44	0
1	B	976	0	951	85	0
1	C	988	0	961	52	0
2	A	104	0	0	5	0
2	B	101	0	0	7	1
2	C	97	0	0	7	0
All	All	3244	0	2866	178	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 31.

All (178) 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:B:100:ASN:HB2	1:B:103:LEU:N	1.53	1.20
1:B:34:LYS:HD2	1:B:35:GLY:H	1.11	1.09
1:B:100:ASN:CB	1:B:103:LEU:H	1.67	1.07
1:B:34:LYS:CD	1:B:35:GLY:H	1.68	1.04
1:C:99:VAL:O	1:C:101:ASP:N	1.93	1.01
1:B:100:ASN:HB2	1:B:103:LEU:H	0.89	1.01
1:A:7:SER:HB2	1:A:35:GLY:HA3	1.45	0.97
1:B:1:GLN:HA	1:B:38:TYR:HE2	1.31	0.95
1:B:33:MET:HE3	2:B:228:HOH:O	1.65	0.95
1:B:100:ASN:HB3	1:B:103:LEU:CB	1.97	0.94
1:C:34:LYS:HD2	1:C:35:GLY:H	1.33	0.91
1:B:1:GLN:HA	1:B:38:TYR:CE2	2.07	0.90
1:C:34:LYS:CD	1:C:35:GLY:H	1.85	0.89
1:A:88:LYS:NZ	1:A:88:LYS:HB3	1.90	0.87
1:C:99:VAL:C	1:C:101:ASP:H	1.75	0.86
1:B:88:LYS:HE3	1:B:91:GLN:NE2	1.92	0.85
1:B:100:ASN:HB3	1:B:103:LEU:HB3	1.58	0.85
1:B:88:LYS:HE3	1:B:91:GLN:HE21	1.41	0.84
1:B:88:LYS:HB2	1:B:91:GLN:HG3	1.60	0.84
1:B:19:PHE:CD1	1:B:78:MET:HE1	2.11	0.83
1:B:1:GLN:N	2:B:178:HOH:O	2.10	0.83
1:A:37:THR:HB	2:A:218:HOH:O	1.77	0.83
1:C:120:GLU:HG3	2:C:208:HOH:O	1.80	0.81
1:C:26:HIS:O	1:C:27:LYS:HG2	1.80	0.80
1:B:19:PHE:HD1	1:B:78:MET:HE1	1.44	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:33:MET:HE2	2:B:193:HOH:O	1.82	0.80
1:B:34:LYS:CD	1:B:35:GLY:N	2.45	0.79
1:B:7:SER:O	1:B:11:LYS:HE3	1.83	0.79
1:A:88:LYS:HB3	1:A:88:LYS:HZ2	1.45	0.79
1:B:100:ASN:O	1:B:101:ASP:HB3	1.82	0.78
1:B:98:ASN:O	1:B:100:ASN:N	2.16	0.78
1:B:33:MET:HE2	2:B:173:HOH:O	1.83	0.78
1:B:100:ASN:HB3	1:B:103:LEU:HB2	1.65	0.77
1:C:3:ASP:HA	2:C:203:HOH:O	1.85	0.77
1:A:89:THR:O	1:A:90:SER:HB2	1.86	0.74
1:C:34:LYS:CG	1:C:35:GLY:H	1.99	0.73
1:B:98:ASN:C	1:B:100:ASN:H	1.89	0.72
1:C:131:HIS:HB2	2:C:222:HOH:O	1.89	0.72
1:A:71:TRP:CZ3	1:A:73:GLU:HG2	2.25	0.72
1:B:70:MET:HG3	1:B:111:LEU:HD21	1.71	0.72
1:B:100:ASN:OD1	1:B:101:ASP:N	2.23	0.71
1:C:89:THR:O	1:C:90:SER:HB3	1.91	0.69
1:A:7:SER:OG	1:A:11:LYS:HE2	1.93	0.68
1:B:97:LEU:C	1:B:99:VAL:H	1.96	0.68
1:B:100:ASN:CB	1:B:103:LEU:CB	2.72	0.67
1:A:75:SER:OG	1:A:77:GLN:HG3	1.95	0.67
1:A:7:SER:HB2	1:A:35:GLY:CA	2.24	0.67
1:C:39:THR:HG21	1:C:120:GLU:CG	2.26	0.66
1:B:99:VAL:O	1:B:100:ASN:CG	2.34	0.65
1:B:34:LYS:HD2	1:B:35:GLY:N	1.96	0.65
1:B:88:LYS:CB	1:B:91:GLN:HG3	2.26	0.65
1:A:43:THR:OG1	1:A:52:GLN:NE2	2.30	0.64
1:A:88:LYS:CD	1:A:89:THR:HG23	2.27	0.64
1:B:100:ASN:CB	1:B:103:LEU:N	2.40	0.63
1:C:99:VAL:HG12	1:C:103:LEU:HD23	1.81	0.63
1:C:4:LEU:HG	1:C:54:MET:CE	2.27	0.63
1:C:39:THR:HG21	1:C:120:GLU:HG3	1.81	0.62
1:B:71:TRP:CZ3	1:B:73:GLU:HG2	2.35	0.61
1:A:88:LYS:HE2	2:A:212:HOH:O	2.00	0.61
1:A:98:ASN:O	1:A:99:VAL:HG22	2.00	0.61
1:B:88:LYS:CE	1:B:91:GLN:HE21	2.12	0.61
1:C:4:LEU:HG	1:C:54:MET:HE2	1.81	0.60
1:A:53:VAL:HG22	1:A:70:MET:HG2	1.82	0.60
1:A:86:ASP:CG	1:A:88:LYS:HZ3	2.05	0.59
1:B:15:THR:HG22	1:B:76:ASN:HD21	1.67	0.59
1:C:53:VAL:HG22	1:C:70:MET:HG2	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:70:MET:HG3	1:B:111:LEU:CD2	2.31	0.59
1:B:98:ASN:C	1:B:100:ASN:N	2.55	0.58
1:C:99:VAL:O	1:C:99:VAL:HG23	2.03	0.58
1:B:34:LYS:CG	1:B:35:GLY:N	2.66	0.58
1:C:96:TRP:CE2	1:C:108:LYS:HE3	2.38	0.58
1:C:34:LYS:HD2	1:C:35:GLY:N	2.12	0.58
1:C:65:GLN:HG2	1:C:84:THR:HG23	1.86	0.58
1:A:86:ASP:OD2	1:A:88:LYS:HD2	2.04	0.57
1:B:36:GLY:HA2	1:B:59:PRO:HD3	1.85	0.57
1:A:22:MET:CE	1:A:78:MET:CE	2.82	0.57
1:B:96:TRP:O	1:B:99:VAL:HG22	2.04	0.57
1:B:26:HIS:O	1:B:27:LYS:HG2	2.04	0.57
1:B:2:ASP:C	1:B:3:ASP:OD2	2.43	0.57
1:C:34:LYS:CG	1:C:35:GLY:N	2.67	0.57
1:A:47:GLY:HA2	1:A:130:SER:OG	2.06	0.56
1:A:62:CYS:HB3	1:B:76:ASN:HB2	1.87	0.56
1:B:99:VAL:O	1:B:99:VAL:HG12	2.06	0.56
1:A:26:HIS:C	1:A:27:LYS:HG2	2.27	0.56
1:C:15:THR:HG22	1:C:76:ASN:HD21	1.71	0.55
1:B:36:GLY:HA2	1:B:59:PRO:CD	2.36	0.55
1:A:15:THR:HG22	1:A:76:ASN:HD21	1.71	0.55
1:B:86:ASP:OD1	1:B:88:LYS:HE2	2.07	0.55
1:B:22:MET:HA	1:B:98:ASN:ND2	2.21	0.55
1:C:6:ILE:HG13	1:C:54:MET:HG3	1.89	0.55
1:C:108:LYS:NZ	2:C:157:HOH:O	2.40	0.54
1:B:2:ASP:O	1:B:3:ASP:OD2	2.26	0.54
1:B:100:ASN:CB	1:B:103:LEU:HB3	2.35	0.54
1:B:108:LYS:NZ	2:B:198:HOH:O	2.41	0.54
1:B:106:ASP:OD2	1:C:122:HIS:HE1	1.92	0.53
1:B:97:LEU:O	1:B:98:ASN:HB2	2.09	0.53
1:C:99:VAL:C	1:C:101:ASP:N	2.44	0.53
1:B:33:MET:HG3	2:B:193:HOH:O	2.09	0.52
1:A:7:SER:OG	1:A:11:LYS:CE	2.58	0.52
1:A:88:LYS:HD2	1:A:89:THR:HG23	1.92	0.52
1:B:1:GLN:CA	1:B:38:TYR:HE2	2.14	0.51
1:C:75:SER:OG	1:C:77:GLN:HG3	2.11	0.51
1:B:100:ASN:C	1:B:102:ALA:H	1.99	0.51
1:B:100:ASN:CG	1:B:101:ASP:H	2.15	0.51
1:A:121:ASN:ND2	2:A:215:HOH:O	2.42	0.51
1:C:120:GLU:H	1:C:120:GLU:CD	2.14	0.51
1:B:88:LYS:CB	1:B:91:GLN:HE21	2.24	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:79:THR:HG21	1:B:103:LEU:HD13	1.94	0.49
1:A:88:LYS:HG2	1:A:89:THR:HG23	1.94	0.49
1:B:1:GLN:N	1:B:40:PRO:HA	2.27	0.49
1:B:99:VAL:O	1:B:100:ASN:CB	2.60	0.49
1:A:112:PHE:CE2	1:A:116:THR:HG21	2.48	0.49
1:C:122:HIS:HD2	2:C:211:HOH:O	1.96	0.49
1:B:97:LEU:C	1:B:99:VAL:N	2.63	0.49
1:B:30:ALA:O	1:B:34:LYS:HB3	2.12	0.48
1:B:71:TRP:HE1	1:B:76:ASN:ND2	2.12	0.48
1:C:36:GLY:HA2	1:C:59:PRO:HD3	1.95	0.48
1:B:109:THR:OG1	1:C:116:THR:HG21	2.14	0.48
1:C:26:HIS:C	1:C:27:LYS:HG2	2.32	0.48
1:B:100:ASN:CB	1:B:103:LEU:HB2	2.39	0.47
1:B:82:PHE:CE2	1:B:84:THR:HG23	2.49	0.47
1:A:11:LYS:NZ	1:A:33:MET:O	2.47	0.47
1:A:36:GLY:O	2:A:173:HOH:O	2.20	0.47
1:B:88:LYS:HB3	1:B:91:GLN:HE21	1.79	0.47
1:B:112:PHE:O	1:B:116:THR:HG23	2.14	0.47
1:B:100:ASN:CG	1:B:101:ASP:N	2.66	0.47
1:A:19:PHE:HD1	1:A:78:MET:HE1	1.79	0.47
1:A:98:ASN:C	1:A:100:ASN:H	2.18	0.46
1:C:98:ASN:C	1:C:100:ASN:H	2.19	0.46
1:B:42:GLN:HA	1:B:42:GLN:HE21	1.81	0.46
1:C:39:THR:HG21	1:C:120:GLU:HG2	1.96	0.46
1:A:19:PHE:CD1	1:A:78:MET:HE1	2.50	0.46
1:B:112:PHE:CE2	1:B:116:THR:HG21	2.50	0.46
1:C:63:GLY:O	1:C:85:ILE:HD12	2.16	0.46
1:C:117:GLY:C	1:C:120:GLU:OE2	2.54	0.46
1:C:49:GLU:CD	1:C:74:LYS:HZ3	2.19	0.46
1:B:88:LYS:CD	1:B:91:GLN:HE21	2.28	0.45
1:A:42:GLN:HG3	1:A:119:LEU:HG	1.98	0.45
1:A:112:PHE:O	1:A:116:THR:HG23	2.17	0.45
1:B:86:ASP:OD1	1:B:88:LYS:HG3	2.17	0.45
1:A:71:TRP:HE1	1:A:76:ASN:ND2	2.15	0.45
1:C:74:LYS:NZ	2:C:202:HOH:O	2.50	0.45
1:A:22:MET:CE	1:A:78:MET:HE3	2.47	0.44
1:B:86:ASP:O	1:B:87:GLU:HB2	2.16	0.44
1:C:35:GLY:O	1:C:59:PRO:HG2	2.17	0.44
1:B:42:GLN:HA	1:B:42:GLN:NE2	2.33	0.44
1:C:116:THR:HG23	1:C:118:SER:H	1.83	0.44
1:A:27:LYS:HE2	1:A:27:LYS:HB3	1.63	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:74:LYS:NZ	2:B:210:HOH:O	2.50	0.43
1:B:88:LYS:NZ	1:B:93:LYS:NZ	2.66	0.43
1:C:7:SER:O	1:C:11:LYS:HG3	2.17	0.43
1:A:75:SER:CB	1:A:77:GLN:HG3	2.48	0.43
1:C:120:GLU:CG	2:C:208:HOH:O	2.53	0.43
1:C:66:ARG:HG3	1:C:115:LEU:HD13	2.01	0.43
1:C:86:ASP:HB3	1:C:89:THR:OG1	2.18	0.43
1:B:71:TRP:HE1	1:B:76:ASN:HD22	1.66	0.42
1:A:88:LYS:CG	1:A:89:THR:HG23	2.49	0.42
1:A:22:MET:HE2	1:A:78:MET:CE	2.48	0.42
1:C:71:TRP:HE1	1:C:76:ASN:ND2	2.17	0.42
1:B:19:PHE:HA	1:B:78:MET:HE2	2.01	0.42
1:C:48:ASP:OD1	1:C:48:ASP:N	2.48	0.42
1:C:6:ILE:CG1	1:C:54:MET:HG3	2.49	0.41
1:B:19:PHE:CD1	1:B:78:MET:CE	2.95	0.41
1:C:98:ASN:C	1:C:100:ASN:N	2.74	0.41
1:B:79:THR:HG21	1:B:103:LEU:HD22	2.01	0.41
1:B:66:ARG:HG3	1:B:115:LEU:HD13	2.03	0.41
1:C:91:GLN:HE21	1:C:91:GLN:HB2	1.58	0.41
1:A:22:MET:SD	1:A:78:MET:CE	3.09	0.41
1:A:38:TYR:CD1	1:A:38:TYR:C	2.94	0.41
1:B:11:LYS:NZ	1:B:33:MET:O	2.44	0.40
1:C:66:ARG:HG2	1:C:83:SER:HB3	2.03	0.40
1:B:1:GLN:H1	1:B:40:PRO:HA	1.87	0.40
1:A:33:MET:HE2	2:A:198:HOH:O	2.20	0.40
1:C:39:THR:HG23	1:C:40:PRO:HD2	2.04	0.40
1:C:115:LEU:HD12	1:C:115:LEU:C	2.42	0.40
1:A:71:TRP:HE1	1:A:76:ASN:HD22	1.70	0.40
1:A:9:LEU:HD11	1:A:54:MET:CE	2.51	0.40
1:B:1:GLN:HG3	1:B:38:TYR:OH	2.21	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)
2:B:176:HOH:O	2:B:176:HOH:O[2_656]	1.92	0.28

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	126/136 (93%)	120 (95%)	3 (2%)	3 (2%)	6 0
1	B	123/136 (90%)	112 (91%)	4 (3%)	7 (6%)	1 0
1	C	127/136 (93%)	120 (94%)	5 (4%)	2 (2%)	9 1
All	All	376/408 (92%)	352 (94%)	12 (3%)	12 (3%)	4 0

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	99	VAL
1	B	2	ASP
1	B	34	LYS
1	B	87	GLU
1	B	99	VAL
1	B	100	ASN
1	C	100	ASN
1	B	101	ASP
1	C	35	GLY
1	A	36	GLY
1	B	3	ASP
1	A	35	GLY

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.

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	108/116 (93%)	102 (94%)	6 (6%)	21	4
1	B	107/116 (92%)	98 (92%)	9 (8%)	11	1
1	C	109/116 (94%)	99 (91%)	10 (9%)	9	1
All	All	324/348 (93%)	299 (92%)	25 (8%)	13	2

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	43	THR
1	A	66	ARG
1	A	76	ASN
1	A	88	LYS
1	A	98	ASN
1	A	100	ASN
1	B	3	ASP
1	B	11	LYS
1	B	27	LYS
1	B	34	LYS
1	B	66	ARG
1	B	76	ASN
1	B	98	ASN
1	B	101	ASP
1	B	129	ARG
1	C	3	ASP
1	C	27	LYS
1	C	33	MET
1	C	34	LYS
1	C	66	ARG
1	C	76	ASN
1	C	87	GLU
1	C	88	LYS
1	C	91	GLN
1	C	98	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	20	ASN

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Mol	Chain	Res	Type
1	A	52	GLN
1	A	76	ASN
1	A	98	ASN
1	A	100	ASN
1	A	121	ASN
1	B	42	GLN
1	B	76	ASN
1	B	77	GLN
1	B	91	GLN
1	B	98	ASN
1	B	121	ASN
1	C	42	GLN
1	C	76	ASN
1	C	91	GLN
1	C	98	ASN
1	C	100	ASN
1	C	121	ASN
1	C	122	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 monosaccharides 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	128/136 (94%)	0.96	16 (12%)	3	3	15, 26, 44, 48	0
1	B	127/136 (93%)	1.03	15 (11%)	4	4	15, 26, 46, 51	0
1	C	129/136 (94%)	1.03	20 (15%)	2	1	14, 26, 46, 50	0
All	All	384/408 (94%)	1.01	51 (13%)	3	2	14, 26, 46, 51	0

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	36	GLY	6.9
1	B	36	GLY	5.9
1	C	38	TYR	5.4
1	B	129	ARG	4.6
1	C	39	THR	4.4
1	A	38	TYR	4.3
1	C	36	GLY	3.9
1	B	2	ASP	3.7
1	B	39	THR	3.7
1	B	38	TYR	3.7
1	C	100	ASN	3.6
1	A	71	TRP	3.6
1	B	1	GLN	3.5
1	A	39	THR	3.4
1	C	101	ASP	3.4
1	A	97	LEU	3.2
1	C	90	SER	3.2
1	A	100	ASN	3.1
1	B	101	ASP	3.0
1	A	85	ILE	3.0
1	A	121	ASN	3.0
1	B	4	LEU	3.0
1	B	87	GLU	3.0

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Mol	Chain	Res	Type	RSRZ
1	C	45	THR	2.9
1	B	99	VAL	2.9
1	A	35	GLY	2.9
1	C	37	THR	2.8
1	A	99	VAL	2.7
1	C	123	PRO	2.7
1	C	99	VAL	2.6
1	A	87	GLU	2.6
1	C	77	GLN	2.5
1	B	123	PRO	2.5
1	B	115	LEU	2.5
1	C	3	ASP	2.4
1	C	34	LYS	2.4
1	A	25	GLY	2.4
1	C	102	ALA	2.3
1	A	101	ASP	2.3
1	C	33	MET	2.3
1	B	100	ASN	2.3
1	C	4	LEU	2.2
1	C	85	ILE	2.2
1	C	88	LYS	2.1
1	C	87	GLU	2.1
1	A	77	GLN	2.1
1	B	98	ASN	2.1
1	C	35	GLY	2.1
1	B	8	SER	2.0
1	A	47	GLY	2.0
1	A	37	THR	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)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.