Full wwPDB X-ray Structure Validation Report (i)

Jan 31, 2016 – 09:07 PM GMT

PDB ID	:	1NJK
Title	:	Crystal Structure of YbaW Probable Thioesterase from Escherichia coli
Authors	:	Kim, Y.; Joachimiak, A.; Edwards, A.; Xu, X.; Savchenko, A.; Midwest Center for Structural Genomics (MCSG)
Deposited on	:	2002-12-31
Resolution	:	1.90 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

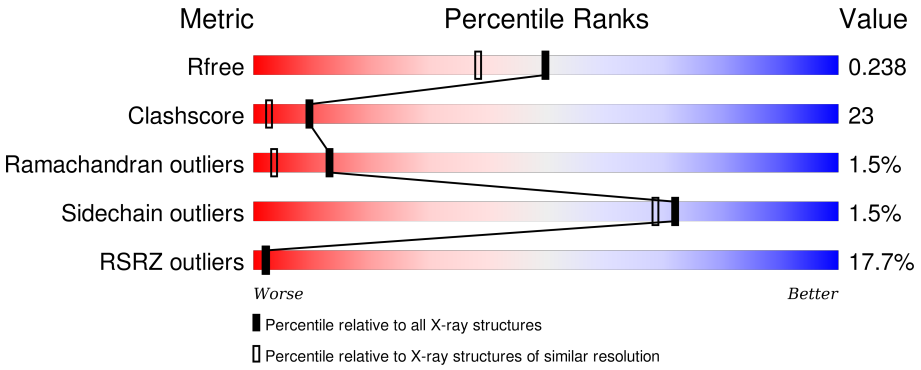
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 1.90 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	4755 (1.90-1.90)
Clashscore	102246	5398 (1.90-1.90)
Ramachandran outliers	100387	5338 (1.90-1.90)
Sidechain outliers	100360	5339 (1.90-1.90)
RSRZ outliers	91569	4766 (1.90-1.90)

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 <=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	156	18% 51% 32% • 15%
1	B	156	18% 55% 29% • 15%
1	C	156	10% 54% 29% • 15%
1	D	156	13% 58% 26% • 15%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4643 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 Hypothetical protein ybaW.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	133	Total	C	N	O	S	Se	0	0	0
			1071	680	186	201	1	3			
1	B	133	Total	C	N	O	S	Se	0	0	0
			1071	680	186	201	1	3			
1	C	133	Total	C	N	O	S	Se	0	0	0
			1071	680	186	201	1	3			
1	D	133	Total	C	N	O	S	Se	0	0	0
			1071	680	186	201	1	3			

There are 104 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-20	GLY	-	CLONING ARTIFACT	UNP P77712
A	-19	SER	-	CLONING ARTIFACT	UNP P77712
A	-18	SER	-	CLONING ARTIFACT	UNP P77712
A	-17	HIS	-	CLONING ARTIFACT	UNP P77712
A	-16	HIS	-	CLONING ARTIFACT	UNP P77712
A	-15	HIS	-	CLONING ARTIFACT	UNP P77712
A	-14	HIS	-	CLONING ARTIFACT	UNP P77712
A	-13	HIS	-	CLONING ARTIFACT	UNP P77712
A	-12	HIS	-	CLONING ARTIFACT	UNP P77712
A	-11	SER	-	CLONING ARTIFACT	UNP P77712
A	-10	SER	-	CLONING ARTIFACT	UNP P77712
A	-9	GLY	-	CLONING ARTIFACT	UNP P77712
A	-8	ARG	-	CLONING ARTIFACT	UNP P77712
A	-7	GLU	-	CLONING ARTIFACT	UNP P77712
A	-6	ASN	-	CLONING ARTIFACT	UNP P77712
A	-5	LEU	-	CLONING ARTIFACT	UNP P77712
A	-4	TYR	-	CLONING ARTIFACT	UNP P77712
A	-3	PHE	-	CLONING ARTIFACT	UNP P77712
A	-2	GLN	-	CLONING ARTIFACT	UNP P77712
A	-1	GLY	-	CLONING ARTIFACT	UNP P77712
A	0	HIS	-	CLONING ARTIFACT	UNP P77712

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Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	MET	MODIFIED RESIDUE	UNP P77712
A	44	MSE	MET	MODIFIED RESIDUE	UNP P77712
A	130	MSE	MET	MODIFIED RESIDUE	UNP P77712
A	133	GLY	-	CLONING ARTIFACT	UNP P77712
A	134	HIS	-	CLONING ARTIFACT	UNP P77712
B	-20	GLY	-	CLONING ARTIFACT	UNP P77712
B	-19	SER	-	CLONING ARTIFACT	UNP P77712
B	-18	SER	-	CLONING ARTIFACT	UNP P77712
B	-17	HIS	-	CLONING ARTIFACT	UNP P77712
B	-16	HIS	-	CLONING ARTIFACT	UNP P77712
B	-15	HIS	-	CLONING ARTIFACT	UNP P77712
B	-14	HIS	-	CLONING ARTIFACT	UNP P77712
B	-13	HIS	-	CLONING ARTIFACT	UNP P77712
B	-12	HIS	-	CLONING ARTIFACT	UNP P77712
B	-11	SER	-	CLONING ARTIFACT	UNP P77712
B	-10	SER	-	CLONING ARTIFACT	UNP P77712
B	-9	GLY	-	CLONING ARTIFACT	UNP P77712
B	-8	ARG	-	CLONING ARTIFACT	UNP P77712
B	-7	GLU	-	CLONING ARTIFACT	UNP P77712
B	-6	ASN	-	CLONING ARTIFACT	UNP P77712
B	-5	LEU	-	CLONING ARTIFACT	UNP P77712
B	-4	TYR	-	CLONING ARTIFACT	UNP P77712
B	-3	PHE	-	CLONING ARTIFACT	UNP P77712
B	-2	GLN	-	CLONING ARTIFACT	UNP P77712
B	-1	GLY	-	CLONING ARTIFACT	UNP P77712
B	0	HIS	-	CLONING ARTIFACT	UNP P77712
B	1	MSE	MET	MODIFIED RESIDUE	UNP P77712
B	44	MSE	MET	MODIFIED RESIDUE	UNP P77712
B	130	MSE	MET	MODIFIED RESIDUE	UNP P77712
B	133	GLY	-	CLONING ARTIFACT	UNP P77712
B	134	HIS	-	CLONING ARTIFACT	UNP P77712
C	-20	GLY	-	CLONING ARTIFACT	UNP P77712
C	-19	SER	-	CLONING ARTIFACT	UNP P77712
C	-18	SER	-	CLONING ARTIFACT	UNP P77712
C	-17	HIS	-	CLONING ARTIFACT	UNP P77712
C	-16	HIS	-	CLONING ARTIFACT	UNP P77712
C	-15	HIS	-	CLONING ARTIFACT	UNP P77712
C	-14	HIS	-	CLONING ARTIFACT	UNP P77712
C	-13	HIS	-	CLONING ARTIFACT	UNP P77712
C	-12	HIS	-	CLONING ARTIFACT	UNP P77712
C	-11	SER	-	CLONING ARTIFACT	UNP P77712
C	-10	SER	-	CLONING ARTIFACT	UNP P77712

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-9	GLY	-	CLONING ARTIFACT	UNP P77712
C	-8	ARG	-	CLONING ARTIFACT	UNP P77712
C	-7	GLU	-	CLONING ARTIFACT	UNP P77712
C	-6	ASN	-	CLONING ARTIFACT	UNP P77712
C	-5	LEU	-	CLONING ARTIFACT	UNP P77712
C	-4	TYR	-	CLONING ARTIFACT	UNP P77712
C	-3	PHE	-	CLONING ARTIFACT	UNP P77712
C	-2	GLN	-	CLONING ARTIFACT	UNP P77712
C	-1	GLY	-	CLONING ARTIFACT	UNP P77712
C	0	HIS	-	CLONING ARTIFACT	UNP P77712
C	1	MSE	MET	MODIFIED RESIDUE	UNP P77712
C	44	MSE	MET	MODIFIED RESIDUE	UNP P77712
C	130	MSE	MET	MODIFIED RESIDUE	UNP P77712
C	133	GLY	-	CLONING ARTIFACT	UNP P77712
C	134	HIS	-	CLONING ARTIFACT	UNP P77712
D	-20	GLY	-	CLONING ARTIFACT	UNP P77712
D	-19	SER	-	CLONING ARTIFACT	UNP P77712
D	-18	SER	-	CLONING ARTIFACT	UNP P77712
D	-17	HIS	-	CLONING ARTIFACT	UNP P77712
D	-16	HIS	-	CLONING ARTIFACT	UNP P77712
D	-15	HIS	-	CLONING ARTIFACT	UNP P77712
D	-14	HIS	-	CLONING ARTIFACT	UNP P77712
D	-13	HIS	-	CLONING ARTIFACT	UNP P77712
D	-12	HIS	-	CLONING ARTIFACT	UNP P77712
D	-11	SER	-	CLONING ARTIFACT	UNP P77712
D	-10	SER	-	CLONING ARTIFACT	UNP P77712
D	-9	GLY	-	CLONING ARTIFACT	UNP P77712
D	-8	ARG	-	CLONING ARTIFACT	UNP P77712
D	-7	GLU	-	CLONING ARTIFACT	UNP P77712
D	-6	ASN	-	CLONING ARTIFACT	UNP P77712
D	-5	LEU	-	CLONING ARTIFACT	UNP P77712
D	-4	TYR	-	CLONING ARTIFACT	UNP P77712
D	-3	PHE	-	CLONING ARTIFACT	UNP P77712
D	-2	GLN	-	CLONING ARTIFACT	UNP P77712
D	-1	GLY	-	CLONING ARTIFACT	UNP P77712
D	0	HIS	-	CLONING ARTIFACT	UNP P77712
D	1	MSE	MET	MODIFIED RESIDUE	UNP P77712
D	44	MSE	MET	MODIFIED RESIDUE	UNP P77712
D	130	MSE	MET	MODIFIED RESIDUE	UNP P77712
D	133	GLY	-	CLONING ARTIFACT	UNP P77712
D	134	HIS	-	CLONING ARTIFACT	UNP P77712

- Molecule 2 is IODIDE ION (three-letter code: IOD) (formula: I).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total I 1 1	0	0
2	D	1	Total I 1 1	0	0
2	C	2	Total I 2 2	0	0

- Molecule 3 is water.

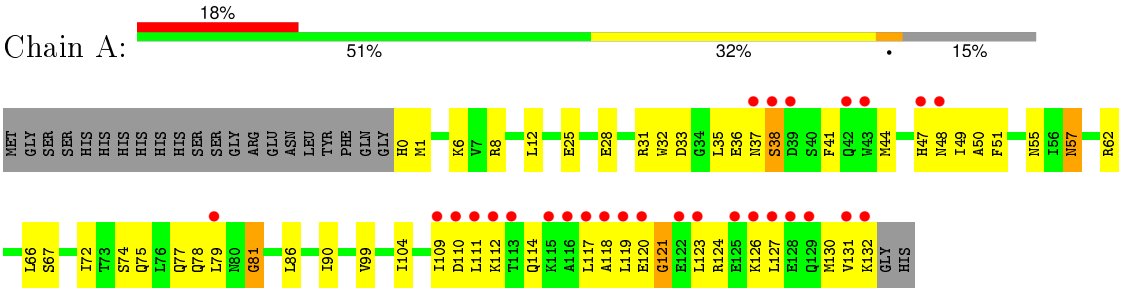
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	89	Total O 89 89	0	0
3	B	80	Total O 80 80	0	0
3	C	92	Total O 92 92	0	0
3	D	94	Total O 94 94	0	0

3 Residue-property plots (i)

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.

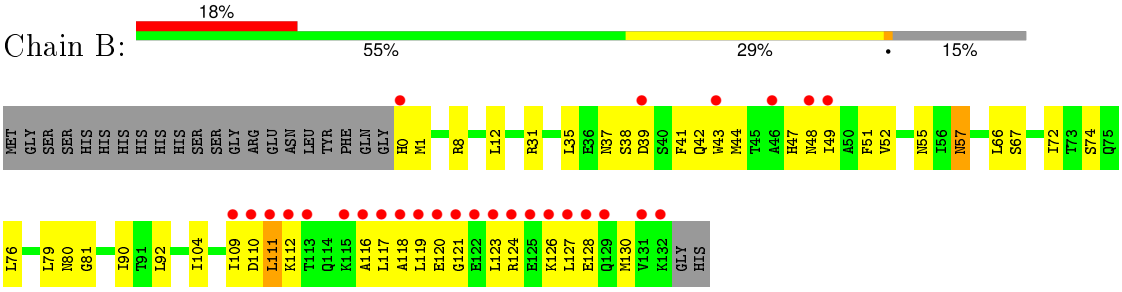
- Molecule 1: Hypothetical protein ybaW

Chain A:



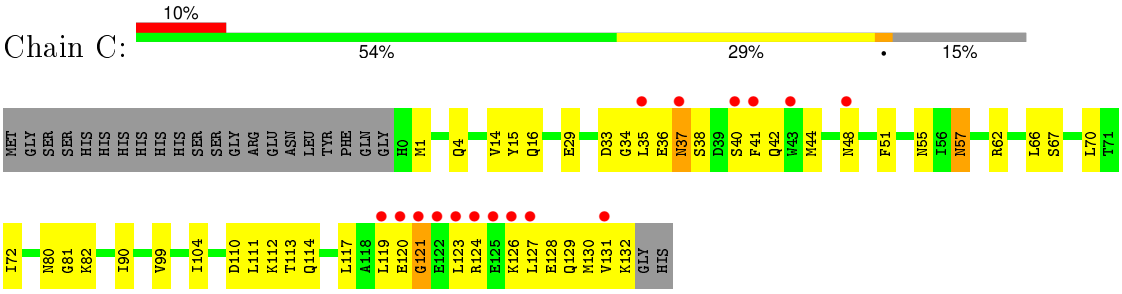
- Molecule 1: Hypothetical protein ybaW

Chain B:



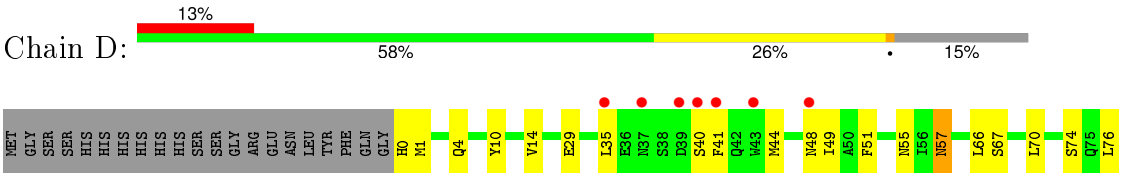
- Molecule 1: Hypothetical protein ybaW

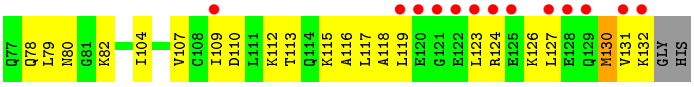
Chain C:



- Molecule 1: Hypothetical protein ybaW

Chain D:




Q77
Q78
L79
N80
G81
K82
I104
V107
C108
I109
D110
L111
K112
T113
Q114
K115
A116
L117
A118
L119
E120
G121
E122
L123
R124
E125
K126
L127
E128
Q129
M130
V131
K132
GLY
HIS

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, α, β, γ	90.13Å 90.13Å 83.37Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	35.35 – 1.90 35.35 – 1.90	Depositor EDS
% Data completeness (in resolution range)	96.6 (35.35-1.90) 96.7 (35.35-1.90)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.10 (at 1.89Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.214 , 0.240 0.212 , 0.238	Depositor DCC
R_{free} test set	5659 reflections (10.87%)	DCC
Wilson B-factor (Å^2)	21.4	Xtriage
Anisotropy	0.041	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 44.5	EDS
Estimated twinning fraction	0.045 for -h,-k,l 0.046 for h,-h-k,-l 0.488 for -k,-h,-l	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.34$	Xtriage
Outliers	0 of 58747 reflections	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	4643	wwPDB-VP
Average B, all atoms (Å^2)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.74% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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

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	0.31	0/1085	0.55	0/1465
1	B	0.32	0/1085	0.56	0/1465
1	C	0.32	0/1085	0.58	0/1465
1	D	0.33	0/1085	0.57	0/1465
All	All	0.32	0/4340	0.56	0/5860

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

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	1071	0	1080	53	0
1	B	1071	0	1080	58	0
1	C	1071	0	1080	48	0
1	D	1071	0	1080	49	0
2	B	1	0	0	0	0
2	C	2	0	0	0	0
2	D	1	0	0	0	0
3	A	89	0	0	0	0
3	B	80	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	92	0	0	0	0
3	D	94	0	0	0	0
All	All	4643	0	4320	194	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 23.

All (194) 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:79:LEU:HD22	1:B:127:LEU:HD13	1.50	0.93
1:B:52:VAL:HG21	1:B:109:ILE:HD12	1.54	0.86
1:A:120:GLU:HG3	1:A:121:GLY:H	1.46	0.80
1:B:123:LEU:O	1:B:126:LYS:HG2	1.83	0.78
1:C:123:LEU:HA	1:C:126:LYS:HE2	1.65	0.78
1:C:35:LEU:HD22	1:C:130:MSE:HE2	1.66	0.77
1:D:110:ASP:HB3	1:D:117:LEU:HD11	1.66	0.77
1:C:44:MSE:HE1	1:C:127:LEU:HD21	1.67	0.77
1:A:123:LEU:HA	1:A:126:LYS:HE2	1.67	0.77
1:A:55:ASN:HD21	1:B:57:ASN:HD21	1.31	0.75
1:C:40:SER:HB3	1:C:126:LYS:HB2	1.67	0.75
1:B:124:ARG:O	1:B:128:GLU:HG2	1.87	0.74
1:B:76:LEU:HD21	1:B:79:LEU:HD11	1.69	0.74
1:B:110:ASP:OD2	1:B:117:LEU:HB2	1.89	0.73
1:B:119:LEU:O	1:B:124:ARG:HG2	1.88	0.72
1:D:119:LEU:O	1:D:124:ARG:HG3	1.91	0.70
1:A:57:ASN:HD21	1:B:55:ASN:HD21	1.41	0.68
1:B:123:LEU:HD22	1:B:126:LYS:NZ	2.10	0.67
1:D:109:ILE:O	1:D:109:ILE:HG13	1.95	0.67
1:D:104:ILE:HD12	1:D:104:ILE:N	2.11	0.66
1:C:104:ILE:HD12	1:C:104:ILE:N	2.10	0.66
1:C:110:ASP:HB3	1:C:113:THR:HG22	1.78	0.65
1:A:55:ASN:HD21	1:B:57:ASN:ND2	1.94	0.65
1:C:48:ASN:ND2	1:C:111:LEU:HB2	2.12	0.65
1:C:110:ASP:HB2	1:C:117:LEU:HD21	1.79	0.65
1:B:48:ASN:HB3	1:B:111:LEU:HB2	1.79	0.64
1:B:119:LEU:HD22	1:B:127:LEU:HD11	1.78	0.64
1:C:41:PHE:HB2	1:C:130:MSE:HE3	1.78	0.64
1:A:47:HIS:HB2	1:A:49:ILE:HG13	1.78	0.64
1:C:33:ASP:O	1:C:36:GLU:HG2	1.97	0.64
1:A:79:LEU:HD22	1:A:119:LEU:HD13	1.79	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:57:ASN:HD21	1:D:55:ASN:HD21	1.47	0.62
1:B:31:ARG:O	1:B:35:LEU:HD23	1.99	0.61
1:B:1:MSE:HE1	1:B:31:ARG:HA	1.81	0.61
1:A:36:GLU:HG3	1:A:37:ASN:ND2	2.15	0.61
1:B:110:ASP:C	1:B:112:LYS:H	2.03	0.61
1:C:1:MSE:HE2	1:C:131:VAL:HA	1.82	0.61
1:A:57:ASN:ND2	1:B:55:ASN:HD21	1.98	0.61
1:D:82:LYS:O	1:D:107:VAL:HG23	2.01	0.61
1:D:51:PHE:HZ	1:D:130:MSE:HE1	1.66	0.60
1:B:76:LEU:HD11	1:B:79:LEU:HD21	1.82	0.60
1:B:38:SER:HB3	1:B:41:PHE:HB2	1.84	0.59
1:D:76:LEU:HD21	1:D:79:LEU:HG	1.83	0.59
1:B:123:LEU:HD22	1:B:126:LYS:HZ2	1.67	0.58
1:C:35:LEU:C	1:C:37:ASN:H	2.07	0.57
1:A:123:LEU:HD23	1:A:126:LYS:HE2	1.85	0.57
1:D:40:SER:HB3	1:D:130:MSE:HB2	1.86	0.57
1:A:104:ILE:HD12	1:A:104:ILE:N	2.21	0.56
1:C:55:ASN:HD21	1:D:57:ASN:HD21	1.53	0.56
1:C:119:LEU:O	1:C:124:ARG:HG3	2.05	0.56
1:A:120:GLU:HG3	1:A:121:GLY:N	2.18	0.55
1:C:57:ASN:ND2	1:D:55:ASN:HD21	2.05	0.55
1:C:126:LYS:O	1:C:130:MSE:HG3	2.06	0.55
1:D:0:HIS:HB3	1:D:132:LYS:HB2	1.88	0.54
1:C:37:ASN:OD1	1:C:132:LYS:HE2	2.08	0.54
1:C:55:ASN:HD21	1:D:57:ASN:ND2	2.05	0.54
1:B:44:MSE:SE	1:B:126:LYS:HD3	2.58	0.53
1:C:38:SER:HB3	1:C:42:GLN:HB2	1.90	0.53
1:D:110:ASP:HB2	1:D:112:LYS:HG3	1.89	0.53
1:B:44:MSE:HE1	1:B:51:PHE:CE2	2.43	0.53
1:B:126:LYS:HG3	1:B:127:LEU:N	2.24	0.53
1:D:49:ILE:HA	1:D:110:ASP:HA	1.91	0.53
1:A:123:LEU:O	1:A:127:LEU:HG	2.09	0.53
1:B:57:ASN:HD22	1:B:57:ASN:C	2.13	0.53
1:B:81:GLY:O	1:B:118:ALA:HA	2.09	0.52
1:C:119:LEU:HD22	1:C:127:LEU:CD1	2.39	0.52
1:D:41:PHE:O	1:D:44:MSE:HB2	2.10	0.52
1:A:44:MSE:HG2	1:A:49:ILE:HB	1.91	0.52
1:B:117:LEU:HB3	1:B:123:LEU:HD11	1.90	0.52
1:B:117:LEU:HD23	1:B:118:ALA:N	2.24	0.52
1:A:35:LEU:HD23	1:A:41:PHE:CD2	2.45	0.52
1:A:31:ARG:O	1:A:35:LEU:HD13	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1:MSE:SE	1:A:131:VAL:HG23	2.59	0.52
1:B:39:ASP:O	1:B:43:TRP:HB2	2.10	0.52
1:B:52:VAL:HG21	1:B:109:ILE:CD1	2.35	0.51
1:D:76:LEU:HD21	1:D:79:LEU:CG	2.41	0.51
1:D:35:LEU:HD12	1:D:130:MSE:HG2	1.93	0.51
1:B:79:LEU:CD2	1:B:127:LEU:HD13	2.31	0.50
1:A:110:ASP:OD1	1:A:112:LYS:HE2	2.11	0.50
1:A:1:MSE:HE1	1:A:31:ARG:HA	1.92	0.50
1:D:123:LEU:HD23	1:D:126:LYS:HE3	1.92	0.50
1:D:48:ASN:HD22	1:D:112:LYS:CG	2.24	0.50
1:A:49:ILE:HG12	1:A:110:ASP:OD2	2.10	0.50
1:A:119:LEU:C	1:A:124:ARG:HB2	2.31	0.50
1:B:80:ASN:O	1:B:119:LEU:HD12	2.12	0.50
1:B:116:ALA:HB1	3:B:449:HOH:O	2.11	0.50
1:B:127:LEU:HA	1:B:130:MSE:HB2	1.93	0.50
1:B:124:ARG:C	1:B:128:GLU:HG2	2.31	0.50
1:C:4:GLN:HA	1:C:70:LEU:O	2.12	0.50
1:D:110:ASP:C	1:D:112:LYS:H	2.15	0.50
1:D:1:MSE:HE3	1:D:74:SER:HB2	1.93	0.50
1:B:1:MSE:HG3	1:B:74:SER:HB2	1.93	0.49
1:B:35:LEU:HD12	1:B:41:PHE:CD2	2.48	0.49
1:A:25:GLU:HG2	1:D:10:TYR:CD1	2.47	0.49
1:C:48:ASN:HD21	1:C:111:LEU:HB2	1.78	0.49
1:B:12:LEU:HD12	1:D:14:VAL:HB	1.95	0.49
1:B:44:MSE:HB3	1:B:49:ILE:O	2.13	0.49
1:C:1:MSE:SE	1:C:34:GLY:HA3	2.63	0.49
1:B:117:LEU:O	1:B:123:LEU:HD12	2.12	0.48
1:D:127:LEU:O	1:D:130:MSE:HB3	2.14	0.48
1:A:110:ASP:CB	1:A:112:LYS:HG2	2.44	0.48
1:A:48:ASN:HD21	1:A:111:LEU:HB2	1.78	0.48
1:A:38:SER:HB3	1:A:130:MSE:HG3	1.96	0.48
1:C:15:TYR:O	1:C:16:GLN:HB2	2.13	0.48
1:B:44:MSE:SE	1:B:126:LYS:HZ3	2.47	0.47
1:D:48:ASN:HD22	1:D:112:LYS:HG3	1.79	0.47
1:C:119:LEU:HD22	1:C:127:LEU:HD11	1.95	0.47
1:A:110:ASP:OD1	1:A:117:LEU:HD11	2.14	0.47
1:D:110:ASP:OD1	1:D:112:LYS:HD2	2.15	0.47
1:D:57:ASN:C	1:D:57:ASN:HD22	2.18	0.47
1:A:12:LEU:HD12	1:C:14:VAL:HB	1.97	0.47
1:B:104:ILE:N	1:B:104:ILE:HD12	2.28	0.47
1:B:0:HIS:ND1	1:B:0:HIS:O	2.48	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:126:LYS:CG	1:B:127:LEU:N	2.77	0.47
1:C:57:ASN:HD22	1:C:57:ASN:C	2.18	0.47
1:A:66:LEU:O	1:A:67:SER:HB2	2.14	0.47
1:C:110:ASP:O	1:C:114:GLN:N	2.45	0.47
1:C:110:ASP:HB3	1:C:113:THR:CG2	2.42	0.47
1:D:110:ASP:C	1:D:112:LYS:N	2.67	0.47
1:B:120:GLU:HA	1:B:124:ARG:HG2	1.96	0.47
1:C:124:ARG:O	1:C:128:GLU:HG3	2.14	0.47
1:C:44:MSE:HE1	1:C:127:LEU:CD2	2.43	0.46
1:A:0:HIS:HB3	1:A:75:GLN:HB3	1.95	0.46
1:A:44:MSE:HE1	1:A:51:PHE:CD2	2.50	0.46
1:C:80:ASN:O	1:C:82:LYS:N	2.49	0.46
1:D:119:LEU:HA	1:D:123:LEU:HD12	1.97	0.46
1:D:35:LEU:N	1:D:35:LEU:HD22	2.31	0.45
1:A:72:ILE:HG12	1:A:90:ILE:HG12	1.99	0.45
1:A:8:ARG:HG3	1:D:29:GLU:OE1	2.16	0.45
1:A:109:ILE:HD11	1:A:114:GLN:C	2.36	0.45
1:B:8:ARG:HG3	1:C:29:GLU:OE1	2.17	0.45
1:D:44:MSE:HE2	1:D:51:PHE:CD2	2.52	0.45
1:D:78:GLN:HG2	1:D:79:LEU:N	2.31	0.45
1:C:66:LEU:O	1:C:67:SER:HB2	2.16	0.45
1:A:0:HIS:HA	1:A:132:LYS:O	2.16	0.45
1:D:104:ILE:CD1	1:D:104:ILE:N	2.80	0.44
1:A:49:ILE:HG22	1:A:50:ALA:N	2.33	0.44
1:D:82:LYS:HG2	1:D:118:ALA:HA	2.00	0.44
1:B:110:ASP:C	1:B:112:LYS:N	2.69	0.44
1:D:40:SER:HB2	1:D:126:LYS:HB2	2.00	0.44
1:A:51:PHE:CZ	1:A:130:MSE:HE1	2.52	0.44
1:B:76:LEU:HD21	1:B:79:LEU:CD1	2.45	0.44
1:B:44:MSE:HG2	1:B:126:LYS:NZ	2.33	0.44
1:A:81:GLY:O	1:A:118:ALA:HA	2.18	0.44
1:D:113:THR:O	1:D:115:LYS:HG3	2.18	0.44
1:A:0:HIS:HB2	1:A:74:SER:O	2.17	0.43
1:A:118:ALA:O	1:A:119:LEU:HD23	2.18	0.43
1:A:77:GLN:HG2	1:A:78:GLN:HG3	2.00	0.43
1:A:110:ASP:HB3	1:A:112:LYS:HG2	2.01	0.43
1:B:120:GLU:HA	1:B:124:ARG:CG	2.48	0.43
1:B:47:HIS:O	1:B:49:ILE:HG13	2.19	0.43
1:C:104:ILE:CD1	1:C:104:ILE:N	2.80	0.43
1:C:48:ASN:O	1:C:111:LEU:HD13	2.18	0.43
1:A:79:LEU:HD13	1:A:124:ARG:HG3	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:35:LEU:O	1:B:38:SER:HB3	2.18	0.43
1:A:110:ASP:CG	1:A:117:LEU:HD11	2.39	0.43
1:A:33:ASP:O	1:A:36:GLU:HG2	2.18	0.43
1:C:51:PHE:HE2	1:C:130:MSE:HE1	1.83	0.43
1:C:120:GLU:O	1:C:121:GLY:C	2.55	0.43
1:D:35:LEU:HB3	1:D:41:PHE:CD2	2.53	0.42
1:C:48:ASN:ND2	1:C:112:LYS:HG3	2.34	0.42
1:B:72:ILE:HG12	1:B:90:ILE:HG12	2.01	0.42
1:A:28:GLU:HG2	1:A:32:TRP:CD1	2.55	0.42
1:D:40:SER:CB	1:D:126:LYS:HB2	2.49	0.42
1:A:35:LEU:HD23	1:A:41:PHE:CE2	2.54	0.42
1:C:123:LEU:HD12	1:C:123:LEU:N	2.35	0.42
1:A:62:ARG:O	1:A:99:VAL:HG13	2.18	0.42
1:A:86:LEU:HD13	1:A:131:VAL:HG21	2.02	0.42
1:B:110:ASP:O	1:B:112:LYS:N	2.53	0.42
1:C:62:ARG:O	1:C:99:VAL:HG13	2.20	0.42
1:B:66:LEU:O	1:B:67:SER:HB2	2.19	0.42
1:C:110:ASP:CB	1:C:113:THR:HG22	2.46	0.41
1:D:66:LEU:O	1:D:67:SER:HB2	2.19	0.41
1:B:42:GLN:C	1:B:44:MSE:H	2.22	0.41
1:D:35:LEU:CD1	1:D:130:MSE:HG2	2.51	0.41
1:C:111:LEU:CD1	1:C:111:LEU:H	2.33	0.41
1:C:126:LYS:HA	1:C:129:GLN:HG2	2.03	0.41
1:D:131:VAL:O	1:D:131:VAL:HG22	2.19	0.41
1:B:47:HIS:O	1:B:112:LYS:HE3	2.21	0.41
1:D:130:MSE:HG2	1:D:130:MSE:O	2.20	0.41
1:C:111:LEU:HD12	1:C:111:LEU:N	2.35	0.41
1:B:76:LEU:HD11	1:B:79:LEU:CD2	2.49	0.41
1:C:1:MSE:CE	1:C:131:VAL:HA	2.49	0.41
1:D:110:ASP:O	1:D:112:LYS:N	2.54	0.41
1:A:57:ASN:HD22	1:A:57:ASN:C	2.24	0.41
1:D:107:VAL:HG21	1:D:116:ALA:CB	2.50	0.41
1:A:6:LYS:HG2	1:A:8:ARG:CZ	2.51	0.41
1:A:110:ASP:HB2	1:A:112:LYS:HG2	2.03	0.41
1:D:80:ASN:C	1:D:124:ARG:HH21	2.25	0.40
1:C:72:ILE:HG12	1:C:90:ILE:HG12	2.02	0.40
1:A:25:GLU:HG2	1:D:10:TYR:CE1	2.56	0.40
1:B:76:LEU:HD21	1:B:79:LEU:HD21	2.04	0.40
1:A:119:LEU:O	1:A:124:ARG:HB2	2.22	0.40
1:D:4:GLN:HA	1:D:70:LEU:O	2.21	0.40

There are no symmetry-related clashes.

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	131/156 (84%)	115 (88%)	13 (10%)	3 (2%)	8 1
1	B	131/156 (84%)	116 (88%)	13 (10%)	2 (2%)	13 3
1	C	131/156 (84%)	120 (92%)	9 (7%)	2 (2%)	13 3
1	D	131/156 (84%)	117 (89%)	13 (10%)	1 (1%)	24 11
All	All	524/624 (84%)	468 (89%)	48 (9%)	8 (2%)	13 3

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	81	GLY
1	A	81	GLY
1	A	121	GLY
1	B	121	GLY
1	C	121	GLY
1	A	38	SER
1	B	111	LEU
1	D	130	MSE

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	118/134 (88%)	117 (99%)	1 (1%)	86 86
1	B	118/134 (88%)	115 (98%)	3 (2%)	55 47

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	C	118/134 (88%)	116 (98%)	2 (2%)	68 64
1	D	118/134 (88%)	117 (99%)	1 (1%)	86 86
All	All	472/536 (88%)	465 (98%)	7 (2%)	72 69

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	57	ASN
1	B	37	ASN
1	B	57	ASN
1	B	92	LEU
1	C	37	ASN
1	C	57	ASN
1	D	57	ASN

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

Mol	Chain	Res	Type
1	A	37	ASN
1	A	48	ASN
1	A	57	ASN
1	A	77	GLN
1	B	17	HIS
1	B	19	ASN
1	B	20	ASN
1	B	37	ASN
1	B	57	ASN
1	B	75	GLN
1	C	19	ASN
1	C	48	ASN
1	C	57	ASN
1	C	129	GLN
1	D	2	GLN
1	D	19	ASN
1	D	37	ASN
1	D	48	ASN
1	D	57	ASN

5.3.3 RNA

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

Of 4 ligands modelled in this entry, 4 are 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

There are no such residues in this entry.

5.8 Polymer linkage issues

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, 95^{th} 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	130/156 (83%)	0.86	28 (21%)	1	1	13, 24, 120, 137	0
1	B	130/156 (83%)	0.74	28 (21%)	1	1	14, 24, 120, 133	0
1	C	130/156 (83%)	0.33	16 (12%)	5	6	12, 24, 105, 120	0
1	D	130/156 (83%)	0.40	20 (15%)	3	3	13, 25, 107, 121	0
All	All	520/624 (83%)	0.58	92 (17%)	2	2	12, 24, 116, 137	0

All (92) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	43	TRP	13.0
1	A	123	LEU	9.9
1	C	131	VAL	9.0
1	A	131	VAL	8.3
1	A	129	GLN	8.0
1	A	119	LEU	7.6
1	B	43	TRP	7.0
1	A	112	LYS	7.0
1	A	113	THR	6.4
1	B	119	LEU	6.2
1	D	43	TRP	6.1
1	B	120	GLU	6.0
1	B	126	LYS	5.9
1	B	112	LYS	5.7
1	B	123	LEU	5.7
1	A	117	LEU	5.6
1	B	48	ASN	5.5
1	B	109	ILE	5.5
1	A	120	GLU	5.3
1	A	122	GLU	5.2
1	B	113	THR	5.0

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Mol	Chain	Res	Type	RSRZ
1	A	118	ALA	5.0
1	A	48	ASN	5.0
1	A	110	ASP	4.9
1	B	121	GLY	4.9
1	C	43	TRP	4.7
1	C	41	PHE	4.7
1	C	121	GLY	4.6
1	A	109	ILE	4.6
1	C	37	ASN	4.4
1	A	132	LYS	4.4
1	A	127	LEU	4.1
1	C	40	SER	4.1
1	B	116	ALA	4.0
1	D	40	SER	4.0
1	D	123	LEU	4.0
1	B	118	ALA	4.0
1	D	131	VAL	4.0
1	D	122	GLU	3.9
1	D	127	LEU	3.8
1	D	124	ARG	3.7
1	C	126	LYS	3.7
1	D	129	GLN	3.7
1	B	39	ASP	3.5
1	D	41	PHE	3.4
1	B	46	ALA	3.4
1	A	111	LEU	3.4
1	B	111	LEU	3.4
1	B	128	GLU	3.4
1	B	127	LEU	3.3
1	D	125	GLU	3.3
1	D	120	GLU	3.2
1	D	132	LYS	3.2
1	B	129	GLN	3.2
1	B	110	ASP	3.2
1	B	117	LEU	3.2
1	B	124	ARG	3.1
1	A	128	GLU	3.0
1	B	125	GLU	3.0
1	B	122	GLU	3.0
1	C	35	LEU	3.0
1	C	123	LEU	2.9
1	D	35	LEU	2.9

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Mol	Chain	Res	Type	RSRZ
1	D	109	ILE	2.9
1	B	132	LYS	2.9
1	D	37	ASN	2.8
1	C	119	LEU	2.7
1	D	128	GLU	2.7
1	A	39	ASP	2.7
1	A	37	ASN	2.7
1	B	131	VAL	2.7
1	C	48	ASN	2.7
1	A	38	SER	2.6
1	D	121	GLY	2.6
1	D	48	ASN	2.6
1	A	126	LYS	2.6
1	C	124	ARG	2.5
1	A	79	LEU	2.5
1	C	127	LEU	2.5
1	C	122	GLU	2.4
1	B	0	HIS	2.4
1	B	49	ILE	2.3
1	A	47	HIS	2.3
1	A	42	GLN	2.2
1	A	125	GLU	2.2
1	D	119	LEU	2.2
1	B	115	LYS	2.2
1	D	39	ASP	2.2
1	C	120	GLU	2.1
1	C	125	GLU	2.1
1	A	116	ALA	2.1
1	A	115	LYS	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)

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95^{th} 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	LLDF	B-factors(Å^2)	Q<0.9
2	IOD	D	403	1/1	0.99	0.06	-0.85	59,59,59,59	1
2	IOD	C	402	1/1	0.99	0.06	-0.85	61,61,61,61	1
2	IOD	C	401	1/1	0.99	0.12	-	25,25,25,25	0
2	IOD	B	404	1/1	0.97	0.21	-	86,86,86,86	1

6.5 Other polymers (i)

There are no such residues in this entry.