Full wwPDB X-ray Structure Validation Report ⓘ

Feb 18, 2024 – 04:43 PM EST

PDB ID	:	4FO7
Title	:	Pseudomonas aeruginosa MetAP, in Mn form
Authors	:	Ye, Q.Z.; Lu, J.P.
Deposited on	:	2012-06-20
Resolution	:	1.80 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

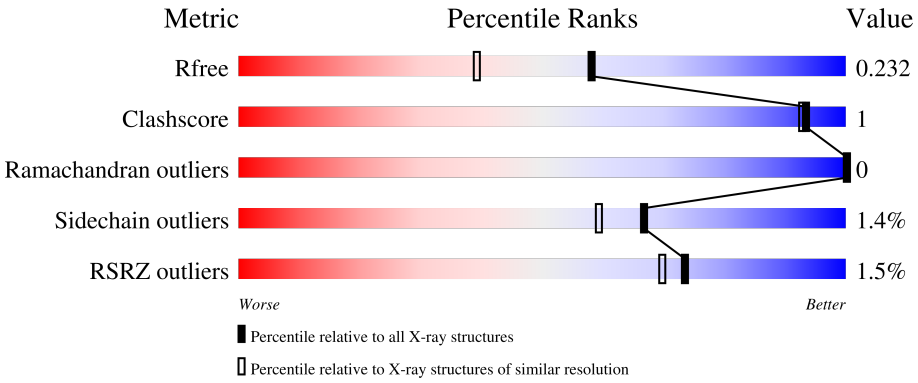
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 1.80 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (1.80-1.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	280	% 86% 6% 8%
1	B	280	84% 8% · 7%
1	C	280	3% 87% 5% 8%
1	D	280	% 85% 6% 8%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 8669 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 Methionine aminopeptidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	258	Total	C	N	O	S	0	1	0
			2036	1288	354	382	12			
1	B	260	Total	C	N	O	S	0	2	0
			2050	1294	357	387	12			
1	C	258	Total	C	N	O	S	0	1	0
			2036	1288	354	382	12			
1	D	258	Total	C	N	O	S	0	2	0
			2037	1285	354	385	13			

There are 76 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	262	GLU	-	expression tag	UNP Q9HXY1
A	263	PHE	-	expression tag	UNP Q9HXY1
A	264	GLU	-	expression tag	UNP Q9HXY1
A	265	LEU	-	expression tag	UNP Q9HXY1
A	266	VAL	-	expression tag	UNP Q9HXY1
A	267	ASP	-	expression tag	UNP Q9HXY1
A	268	LYS	-	expression tag	UNP Q9HXY1
A	269	LEU	-	expression tag	UNP Q9HXY1
A	270	ALA	-	expression tag	UNP Q9HXY1
A	271	ALA	-	expression tag	UNP Q9HXY1
A	272	ALA	-	expression tag	UNP Q9HXY1
A	273	LEU	-	expression tag	UNP Q9HXY1
A	274	GLU	-	expression tag	UNP Q9HXY1
A	275	HIS	-	expression tag	UNP Q9HXY1
A	276	HIS	-	expression tag	UNP Q9HXY1
A	277	HIS	-	expression tag	UNP Q9HXY1
A	278	HIS	-	expression tag	UNP Q9HXY1
A	279	HIS	-	expression tag	UNP Q9HXY1
A	280	HIS	-	expression tag	UNP Q9HXY1
B	262	GLU	-	expression tag	UNP Q9HXY1
B	263	PHE	-	expression tag	UNP Q9HXY1

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Chain	Residue	Modelled	Actual	Comment	Reference
B	264	GLU	-	expression tag	UNP Q9HXY1
B	265	LEU	-	expression tag	UNP Q9HXY1
B	266	VAL	-	expression tag	UNP Q9HXY1
B	267	ASP	-	expression tag	UNP Q9HXY1
B	268	LYS	-	expression tag	UNP Q9HXY1
B	269	LEU	-	expression tag	UNP Q9HXY1
B	270	ALA	-	expression tag	UNP Q9HXY1
B	271	ALA	-	expression tag	UNP Q9HXY1
B	272	ALA	-	expression tag	UNP Q9HXY1
B	273	LEU	-	expression tag	UNP Q9HXY1
B	274	GLU	-	expression tag	UNP Q9HXY1
B	275	HIS	-	expression tag	UNP Q9HXY1
B	276	HIS	-	expression tag	UNP Q9HXY1
B	277	HIS	-	expression tag	UNP Q9HXY1
B	278	HIS	-	expression tag	UNP Q9HXY1
B	279	HIS	-	expression tag	UNP Q9HXY1
B	280	HIS	-	expression tag	UNP Q9HXY1
C	262	GLU	-	expression tag	UNP Q9HXY1
C	263	PHE	-	expression tag	UNP Q9HXY1
C	264	GLU	-	expression tag	UNP Q9HXY1
C	265	LEU	-	expression tag	UNP Q9HXY1
C	266	VAL	-	expression tag	UNP Q9HXY1
C	267	ASP	-	expression tag	UNP Q9HXY1
C	268	LYS	-	expression tag	UNP Q9HXY1
C	269	LEU	-	expression tag	UNP Q9HXY1
C	270	ALA	-	expression tag	UNP Q9HXY1
C	271	ALA	-	expression tag	UNP Q9HXY1
C	272	ALA	-	expression tag	UNP Q9HXY1
C	273	LEU	-	expression tag	UNP Q9HXY1
C	274	GLU	-	expression tag	UNP Q9HXY1
C	275	HIS	-	expression tag	UNP Q9HXY1
C	276	HIS	-	expression tag	UNP Q9HXY1
C	277	HIS	-	expression tag	UNP Q9HXY1
C	278	HIS	-	expression tag	UNP Q9HXY1
C	279	HIS	-	expression tag	UNP Q9HXY1
C	280	HIS	-	expression tag	UNP Q9HXY1
D	262	GLU	-	expression tag	UNP Q9HXY1
D	263	PHE	-	expression tag	UNP Q9HXY1
D	264	GLU	-	expression tag	UNP Q9HXY1
D	265	LEU	-	expression tag	UNP Q9HXY1
D	266	VAL	-	expression tag	UNP Q9HXY1
D	267	ASP	-	expression tag	UNP Q9HXY1

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Chain	Residue	Modelled	Actual	Comment	Reference
D	268	LYS	-	expression tag	UNP Q9HXY1
D	269	LEU	-	expression tag	UNP Q9HXY1
D	270	ALA	-	expression tag	UNP Q9HXY1
D	271	ALA	-	expression tag	UNP Q9HXY1
D	272	ALA	-	expression tag	UNP Q9HXY1
D	273	LEU	-	expression tag	UNP Q9HXY1
D	274	GLU	-	expression tag	UNP Q9HXY1
D	275	HIS	-	expression tag	UNP Q9HXY1
D	276	HIS	-	expression tag	UNP Q9HXY1
D	277	HIS	-	expression tag	UNP Q9HXY1
D	278	HIS	-	expression tag	UNP Q9HXY1
D	279	HIS	-	expression tag	UNP Q9HXY1
D	280	HIS	-	expression tag	UNP Q9HXY1

- Molecule 2 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	4	Total Mn 4 4	0	0
2	B	3	Total Mn 3 3	0	0
2	C	3	Total Mn 3 3	0	0
2	D	4	Total Mn 4 4	0	0

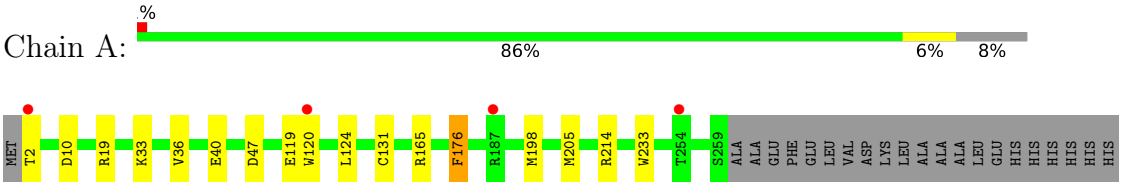
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	130	Total O 130 130	0	0
3	B	148	Total O 148 148	0	0
3	C	102	Total O 102 102	0	0
3	D	116	Total O 116 116	0	0

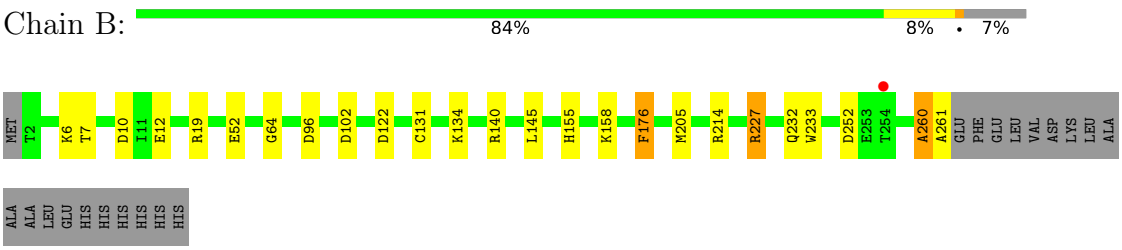
3 Residue-property plots (i)

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.

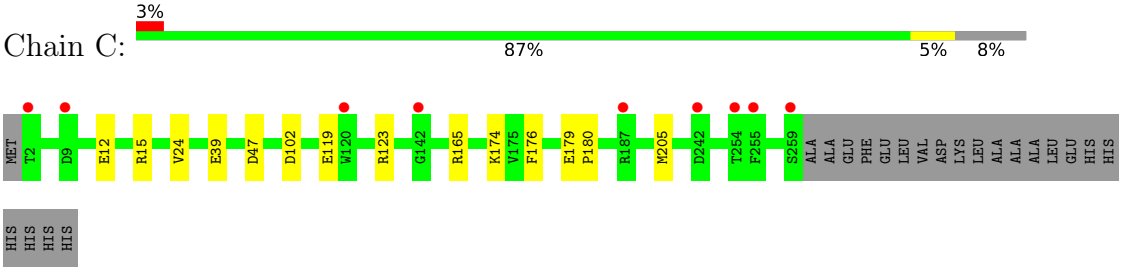
• Molecule 1: Methionine aminopeptidase



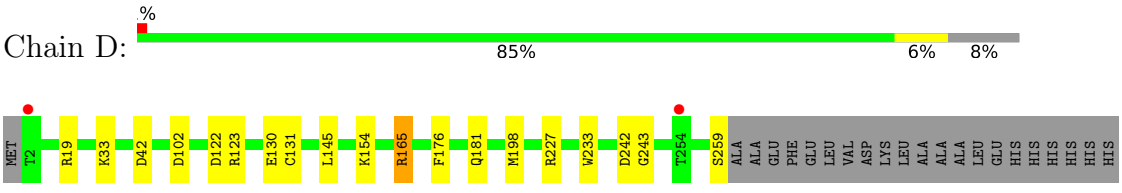
• Molecule 1: Methionine aminopeptidase



• Molecule 1: Methionine aminopeptidase



• Molecule 1: Methionine aminopeptidase



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	85.02Å 111.42Å 140.16Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	35.80 – 1.80 35.80 – 1.80	Depositor EDS
% Data completeness (in resolution range)	92.2 (35.80-1.80) 92.2 (35.80-1.80)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.26 (at 1.81Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.192 , 0.230 0.193 , 0.232	Depositor DCC
R_{free} test set	5723 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å^2)	20.3	Xtriage
Anisotropy	0.073	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.38 , 84.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	8669	wwPDB-VP
Average B, all atoms (Å^2)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.71% 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

5.1 Standard geometry

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

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	1.44	1/2080 (0.0%)	1.25	12/2816 (0.4%)
1	B	1.49	3/2092 (0.1%)	1.26	13/2830 (0.5%)
1	C	1.28	1/2080 (0.0%)	1.20	5/2816 (0.2%)
1	D	1.32	2/2079 (0.1%)	1.26	11/2813 (0.4%)
All	All	1.39	7/8331 (0.1%)	1.24	41/11275 (0.4%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	233	TRP	CB-CG	6.75	1.62	1.50
1	B	131	CYS	CB-SG	6.50	1.93	1.82
1	A	233	TRP	CB-CG	6.19	1.61	1.50
1	C	24	VAL	CB-CG1	5.53	1.64	1.52
1	B	233	TRP	CB-CG	5.39	1.59	1.50
1	B	52	GLU	CB-CG	5.37	1.62	1.52
1	D	131	CYS	CB-SG	5.26	1.91	1.82

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	D	19	ARG	NE-CZ-NH2	-11.13	114.73	120.30
1	B	19	ARG	NE-CZ-NH1	10.58	125.59	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	19	ARG	NE-CZ-NH2	-10.11	115.25	120.30
1	D	19	ARG	NE-CZ-NH1	9.78	125.19	120.30
1	C	165	ARG	NE-CZ-NH1	8.48	124.54	120.30
1	B	176	PHE	CB-CG-CD1	8.25	126.58	120.80
1	A	19	ARG	NE-CZ-NH1	7.77	124.19	120.30
1	B	122	ASP	CB-CG-OD1	7.57	125.11	118.30
1	D	165	ARG	NE-CZ-NH1	-7.42	116.59	120.30
1	B	96	ASP	CB-CG-OD1	7.39	124.95	118.30
1	B	252	ASP	CB-CG-OD1	7.28	124.85	118.30
1	A	165	ARG	NE-CZ-NH2	-7.27	116.67	120.30
1	A	10	ASP	CB-CG-OD1	7.02	124.62	118.30
1	B	140	ARG	NE-CZ-NH1	6.94	123.77	120.30
1	A	47	ASP	CB-CG-OD1	6.63	124.27	118.30
1	C	165	ARG	NE-CZ-NH2	-6.57	117.01	120.30
1	C	205	MET	CG-SD-CE	6.48	110.57	100.20
1	C	47	ASP	CB-CG-OD1	6.41	124.07	118.30
1	D	122	ASP	CB-CG-OD1	6.24	123.92	118.30
1	B	140	ARG	NE-CZ-NH2	-6.13	117.24	120.30
1	B	145	LEU	CA-CB-CG	6.12	129.38	115.30
1	C	102	ASP	CB-CG-OD1	6.09	123.78	118.30
1	A	214	ARG	NE-CZ-NH1	6.05	123.33	120.30
1	A	19	ARG	NE-CZ-NH2	-5.89	117.36	120.30
1	D	130	GLU	OE1-CD-OE2	-5.80	116.34	123.30
1	B	102	ASP	CB-CG-OD1	5.71	123.44	118.30
1	A	198	MET	CA-CB-CG	5.63	122.88	113.30
1	D	154	LYS	CD-CE-NZ	-5.61	98.80	111.70
1	D	102	ASP	CB-CG-OD1	5.60	123.34	118.30
1	A	176	PHE	CB-CG-CD1	5.58	124.71	120.80
1	D	123	ARG	NE-CZ-NH1	5.48	123.04	120.30
1	D	42	ASP	CB-CG-OD2	5.42	123.17	118.30
1	A	131	CYS	CA-CB-SG	5.35	123.64	114.00
1	B	10	ASP	CB-CG-OD1	5.32	123.08	118.30
1	A	205	MET	CG-SD-CE	5.24	108.58	100.20
1	A	124	LEU	CB-CG-CD2	5.21	119.86	111.00
1	B	64	GLY	N-CA-C	5.21	126.13	113.10
1	B	227	ARG	NE-CZ-NH2	5.18	122.89	120.30
1	D	198	MET	CA-CB-CG	5.15	122.06	113.30
1	D	145	LEU	CA-CB-CG	5.15	127.14	115.30
1	A	40	GLU	OE1-CD-OE2	-5.05	117.24	123.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	260	ALA	Peptide

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	2036	0	2031	4	0
1	B	2050	0	2049	13	0
1	C	2036	0	2031	3	0
1	D	2037	0	2031	4	0
2	A	4	0	0	0	0
2	B	3	0	0	0	0
2	C	3	0	0	0	0
2	D	4	0	0	0	0
3	A	130	0	0	0	1
3	B	148	0	0	0	1
3	C	102	0	0	0	0
3	D	116	0	0	1	0
All	All	8669	0	8142	22	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 1.

All (22) 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:6:LYS:H	1:D:181:GLN:HE22	1.10	0.98
1:B:214:ARG:HH21	1:B:227:ARG:NH1	1.73	0.86
1:B:6:LYS:H	1:D:181:GLN:NE2	1.73	0.86
1:C:12:GLU:OE2	1:C:15:ARG:NH1	2.22	0.72
1:B:6:LYS:HD2	1:B:7:THR:H	1.61	0.65
1:B:214:ARG:HH21	1:B:227:ARG:HH12	1.46	0.63
1:D:165:ARG:NH1	3:D:459:HOH:O	2.40	0.54
1:A:120[B]:TRP:N	1:A:120[B]:TRP:CD1	2.77	0.52
1:A:33:LYS:O	1:A:36:VAL:HG12	2.10	0.51
1:B:214:ARG:NH2	1:B:227:ARG:NH1	2.51	0.50
1:C:179:GLU:OE1	1:C:180:PRO:HA	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:119:GLU:HG3	1:A:120[B]:TRP:CD1	2.48	0.48
1:B:205:MET:CE	1:B:232:GLN:NE2	2.78	0.46
1:A:119:GLU:HG3	1:A:120[B]:TRP:HD1	1.80	0.45
1:D:243:GLY:HA3	1:D:259:SER:OG	2.17	0.44
1:B:6:LYS:CD	1:B:7:THR:H	2.29	0.44
1:B:260:ALA:N	1:B:261:ALA:HA	2.33	0.43
1:B:214:ARG:NH2	1:B:227:ARG:HH12	2.14	0.42
1:B:7:THR:HG22	1:B:12[B]:GLU:HG3	2.01	0.41
1:B:205:MET:HE2	1:B:232:GLN:NE2	2.36	0.41
1:B:134:LYS:HE3	1:B:155:HIS:CD2	2.56	0.41
1:C:123:ARG:HA	1:C:123:ARG:HD3	1.84	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)
3:A:514:HOH:O	3:B:542:HOH:O[4_545]	1.50	0.70

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	257/280 (92%)	252 (98%)	5 (2%)	0	100	100
1	B	260/280 (93%)	256 (98%)	4 (2%)	0	100	100
1	C	257/280 (92%)	253 (98%)	4 (2%)	0	100	100
1	D	258/280 (92%)	252 (98%)	6 (2%)	0	100	100
All	All	1032/1120 (92%)	1013 (98%)	19 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	222/238 (93%)	220 (99%)	2 (1%)	78	75
1	B	223/238 (94%)	221 (99%)	2 (1%)	78	75
1	C	222/238 (93%)	218 (98%)	4 (2%)	59	48
1	D	223/238 (94%)	219 (98%)	4 (2%)	59	48
All	All	890/952 (94%)	878 (99%)	12 (1%)	67	62

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	THR
1	A	176	PHE
1	B	158	LYS
1	B	176	PHE
1	C	39	GLU
1	C	119	GLU
1	C	174	LYS
1	C	176	PHE
1	D	33	LYS
1	D	176	PHE
1	D	227	ARG
1	D	242	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	155	HIS
1	A	232	GLN
1	B	46	HIS
1	B	155	HIS
1	B	232	GLN
1	C	155	HIS
1	C	232	GLN

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Mol	Chain	Res	Type
1	D	46	HIS
1	D	181	GLN

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 monosaccharides in this entry.

5.6 Ligand geometry

Of 14 ligands modelled in this entry, 14 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	258/280 (92%)	-0.22	4 (1%)	72	68	9, 18, 33, 41	0
1	B	260/280 (92%)	-0.28	1 (0%)	92	90	9, 17, 33, 40	0
1	C	258/280 (92%)	0.00	9 (3%)	44	38	13, 22, 40, 52	0
1	D	258/280 (92%)	-0.14	2 (0%)	86	84	13, 21, 36, 47	0
All	All	1034/1120 (92%)	-0.16	16 (1%)	73	70	9, 20, 36, 52	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	2	THR	5.8
1	A	120[A]	TRP	4.1
1	C	254	THR	3.6
1	A	2	THR	3.3
1	C	242	ASP	3.0
1	C	2	THR	2.9
1	C	9	ASP	2.8
1	C	259	SER	2.6
1	C	142	GLY	2.5
1	C	120[A]	TRP	2.4
1	A	187	ARG	2.4
1	C	187	ARG	2.3
1	A	254	THR	2.2
1	B	254	THR	2.2
1	D	254	THR	2.2
1	C	255	PHE	2.1

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

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, 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	B-factors(Å^2)	Q<0.9
2	MN	D	303	1/1	0.90	0.13	71,71,71,71	0
2	MN	D	304	1/1	0.98	0.05	24,24,24,24	1
2	MN	C	301	1/1	0.99	0.08	17,17,17,17	0
2	MN	A	303	1/1	0.99	0.05	19,19,19,19	0
2	MN	B	303	1/1	0.99	0.07	21,21,21,21	1
2	MN	B	302	1/1	1.00	0.07	10,10,10,10	0
2	MN	A	302	1/1	1.00	0.08	12,12,12,12	0
2	MN	A	301	1/1	1.00	0.11	13,13,13,13	0
2	MN	C	302	1/1	1.00	0.08	14,14,14,14	0
2	MN	C	303	1/1	1.00	0.06	19,19,19,19	1
2	MN	D	301	1/1	1.00	0.10	16,16,16,16	0
2	MN	D	302	1/1	1.00	0.09	14,14,14,14	0
2	MN	A	304	1/1	1.00	0.06	17,17,17,17	1
2	MN	B	301	1/1	1.00	0.09	11,11,11,11	0

6.5 Other polymers (i)

There are no such residues in this entry.