



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 05:34 AM UTC

PDB ID : 4MX8 / pdb_00004mx8
Title : Crystal Structure of TroA-like Periplasmic Binding Protein Peripla_BP_2 from Xylanimonas cellulosilytica
Authors : Kim, Y.; Wu, R.; Endres, M.; Joachimiak, A.; Midwest Center for Structural Genomics (MCSG)
Deposited on : 2013-09-26
Resolution : 2.91 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 : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

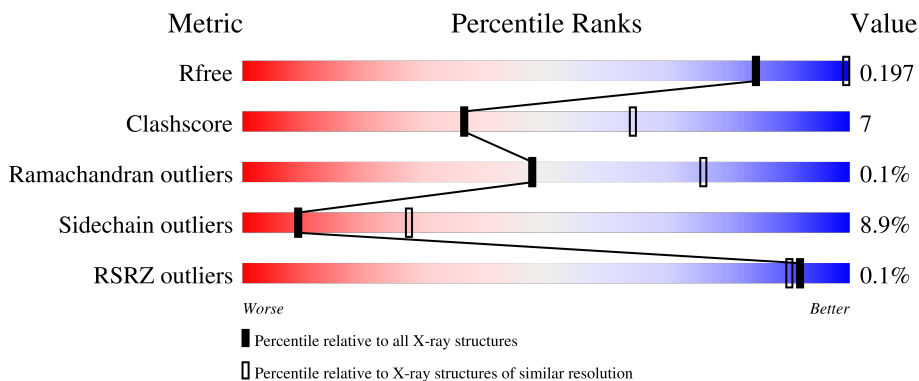
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.91 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}	180053	2995 (2.94-2.90)
Clashscore	190562	3213 (2.94-2.90)
Ramachandran outliers	187476	3128 (2.94-2.90)
Sidechain outliers	187428	3130 (2.94-2.90)
RSRZ outliers	180081	2995 (2.94-2.90)

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 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 $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	315	 79% 16% . .
1	B	315	 77% 17% . .
1	C	315	 73% 19% . .
1	D	315	 78% 16% . .
1	E	315	 73% 19% . .

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Mol	Chain	Length	Quality of chain
1	F	315	 70% 22% . .

2 Entry composition i

There are 2 unique types of molecules in this entry. The entry contains 13799 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 Periplasmic binding protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	Se			
1	A	302	2273	1434	367	469	3	0	2	0
1	B	302	2272	1430	370	469	3	0	2	0
1	C	301	2246	1415	364	464	3	0	0	0
1	D	302	2264	1429	366	466	3	0	1	0
1	E	301	2255	1420	365	467	3	0	1	0
1	F	301	2255	1420	366	466	3	0	1	0

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	SER	-	expression tag	UNP D1BRG9
A	0	ASN	-	expression tag	UNP D1BRG9
A	1	ALA	-	expression tag	UNP D1BRG9
B	-1	SER	-	expression tag	UNP D1BRG9
B	0	ASN	-	expression tag	UNP D1BRG9
B	1	ALA	-	expression tag	UNP D1BRG9
C	-1	SER	-	expression tag	UNP D1BRG9
C	0	ASN	-	expression tag	UNP D1BRG9
C	1	ALA	-	expression tag	UNP D1BRG9
D	-1	SER	-	expression tag	UNP D1BRG9
D	0	ASN	-	expression tag	UNP D1BRG9
D	1	ALA	-	expression tag	UNP D1BRG9
E	-1	SER	-	expression tag	UNP D1BRG9
E	0	ASN	-	expression tag	UNP D1BRG9
E	1	ALA	-	expression tag	UNP D1BRG9
F	-1	SER	-	expression tag	UNP D1BRG9
F	0	ASN	-	expression tag	UNP D1BRG9

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Chain	Residue	Modelled	Actual	Comment	Reference
F	1	ALA	-	expression tag	UNP D1BRG9


- Molecule 2 is water.

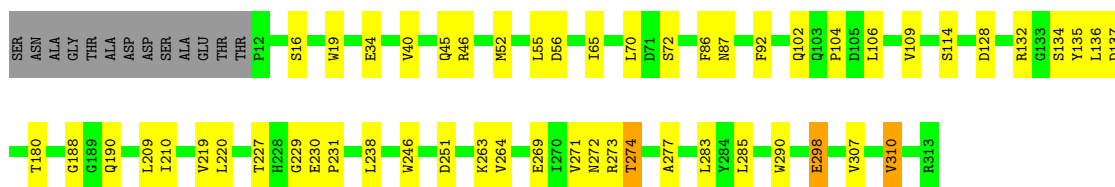
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	73	Total O 73 73	0	0
2	B	56	Total O 56 56	0	0
2	C	18	Total O 18 18	0	0
2	D	56	Total O 56 56	0	0
2	E	18	Total O 18 18	0	0
2	F	13	Total O 13 13	0	0

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.

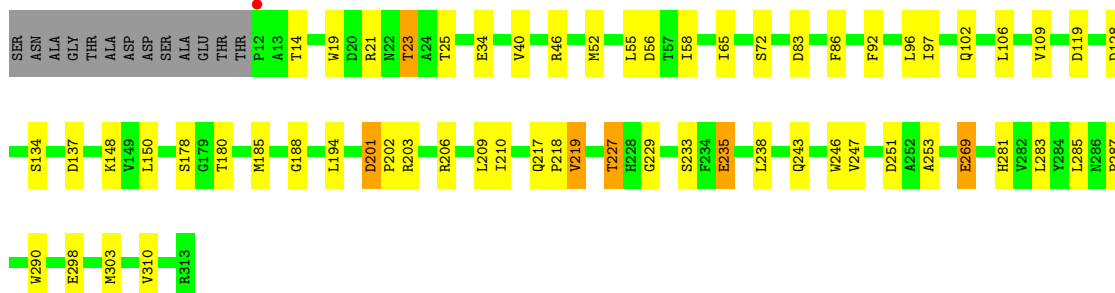
- Molecule 1: Periplasmic binding protein

Chain A:  79% 16% ..



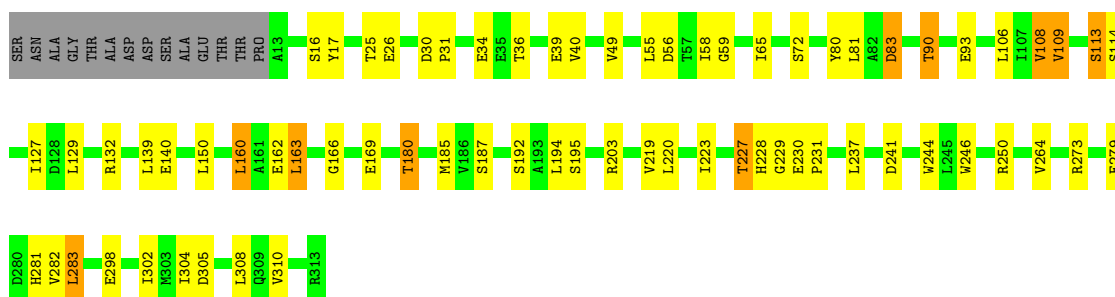
- Molecule 1: Periplasmic binding protein

Chain B:  77% 17% ..




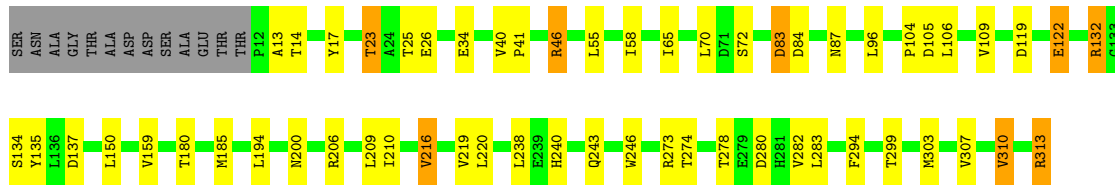
- Molecule 1: Periplasmic binding protein

Chain C:  73% 19% ..



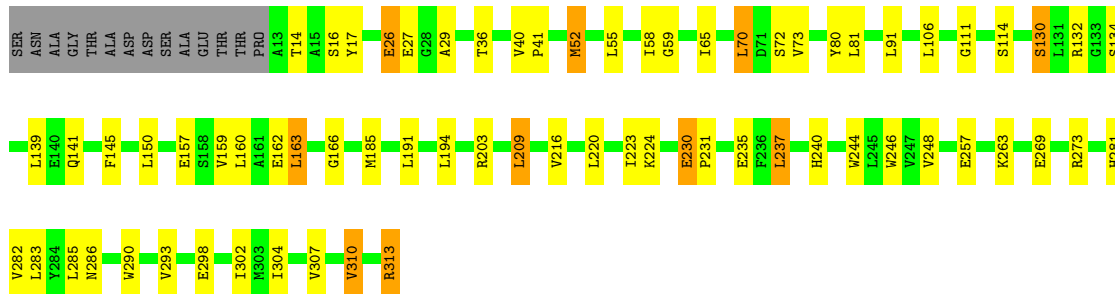
- Molecule 1: Periplasmic binding protein

Chain D:  78% 16%



• Molecule 1: Periplasmic binding protein

Chain E:  73% 19%



• Molecule 1: Periplasmic binding protein

Chain F:  70% 22%



4 Data and refinement statistics i

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	201.66Å 116.55Å 177.16Å 90.00° 112.19° 90.00°	Depositor
Resolution (Å)	49.37 – 2.91 49.37 – 2.91	Depositor EDS
% Data completeness (in resolution range)	96.6 (49.37-2.91) 96.3 (49.37-2.91)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.34 (at 2.91Å)	Xtrriage
Refinement program	PHENIX (phenix.refine: dev_1367)	Depositor
R, R_{free}	0.159 , 0.196 0.161 , 0.197	Depositor DCC
R_{free} test set	4014 reflections (4.76%)	wwPDB-VP
Wilson B-factor (Å ²)	54.1	Xtrriage
Anisotropy	0.053	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 39.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	0.467 for $1/2^*h+3/2^*k, 1/2^*h-1/2^*k, -1/2^*h-1/2^*k-l$ 0.468 for $1/2^*h-3/2^*k, -1/2^*h-1/2^*k, -1/2^*h+1/2^*k-l$	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	13799	wwPDB-VP
Average B, all atoms (Å ²)	59.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.22% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.57	0/2313	0.86	0/3161
1	B	0.58	0/2311	0.87	1/3158 (0.0%)
1	C	0.49	0/2284	0.86	0/3122
1	D	0.58	0/2304	0.89	1/3149 (0.0%)
1	E	0.47	0/2293	0.84	2/3134 (0.1%)
1	F	0.48	0/2293	0.84	2/3134 (0.1%)
All	All	0.53	0/13798	0.86	6/18858 (0.0%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	13	ALA	N-CA-C	5.38	116.86	110.19
1	E	286	ASN	CA-C-N	5.32	124.78	119.24
1	E	286	ASN	C-N-CA	5.32	124.78	119.24
1	F	201	ASP	CA-C-N	5.21	124.83	119.56
1	F	201	ASP	C-N-CA	5.21	124.83	119.56
1	B	247	VAL	N-CA-C	5.04	115.23	108.17

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	2273	0	2185	31	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	2272	0	2187	38	0
1	C	2246	0	2164	35	0
1	D	2264	0	2180	32	0
1	E	2255	0	2169	33	0
1	F	2255	0	2171	40	0
2	A	73	0	0	3	0
2	B	56	0	0	2	0
2	C	18	0	0	1	0
2	D	56	0	0	0	0
2	E	18	0	0	1	0
2	F	13	0	0	1	0
All	All	13799	0	13056	191	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 7.

All (191) 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:E:111:GLY:H	1:E:130:SER:HB3	1.44	0.82
1:C:230:GLU:HG3	1:C:231:PRO:HD2	1.62	0.82
1:C:55:LEU:HG	1:C:65:ILE:HG13	1.66	0.77
1:B:134:SER:HB2	1:B:137:ASP:H	1.51	0.74
1:E:52:MSE:HE2	1:E:91:LEU:HD21	1.70	0.74
1:C:109:VAL:HG22	1:C:113:SER:HB2	1.72	0.72
1:A:271:VAL:O	1:A:274:THR:HB	1.91	0.71
1:D:273:ARG:NH1	1:F:241:ASP:OD2	2.24	0.71
1:E:230:GLU:HG3	1:E:231:PRO:HD2	1.74	0.70
1:F:51:ASP:HB2	1:F:110:GLY:HA3	1.72	0.69
1:A:274:THR:HG22	1:A:277:ALA:H	1.56	0.69
1:E:55:LEU:HG	1:E:65:ILE:HG13	1.75	0.69
1:A:238:LEU:O	1:C:273:ARG:NH1	2.27	0.68
1:B:55:LEU:HG	1:B:65:ILE:HG13	1.75	0.68
1:B:227:THR:HG22	1:B:229:GLY:H	1.59	0.68
1:A:227:THR:HG22	1:A:229:GLY:H	1.59	0.68
1:B:86:PHE:CE1	1:B:102:GLN:HG2	2.29	0.68
1:C:139:LEU:HD21	1:C:163:LEU:HD23	1.76	0.67
1:F:185:MSE:HE2	1:F:250:ARG:HB2	1.75	0.67
1:A:86:PHE:CE2	1:A:102:GLN:HG2	2.30	0.67
1:D:55:LEU:HG	1:D:65:ILE:HG13	1.77	0.66
1:F:139:LEU:HD21	1:F:163:LEU:HD23	1.77	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:238:LEU:O	1:E:273:ARG:NH1	2.29	0.66
1:F:253:ALA:HB2	1:F:287:PRO:HB2	1.76	0.66
1:F:55:LEU:HG	1:F:65:ILE:HG13	1.77	0.66
1:A:273:ARG:NH1	1:C:241:ASP:OD1	2.28	0.65
1:D:83:ASP:N	1:D:83:ASP:OD1	2.28	0.65
1:A:55:LEU:HG	1:A:65:ILE:HG13	1.79	0.63
1:B:58:ILE:HG23	1:B:150:LEU:HD11	1.81	0.63
1:C:83:ASP:N	1:C:83:ASP:OD1	2.22	0.62
1:D:238:LEU:O	1:F:273:ARG:NH1	2.32	0.61
1:A:272:ASN:ND2	2:A:455:HOH:O	2.34	0.61
1:D:70:LEU:HD12	1:D:87:ASN:HB2	1.83	0.61
1:E:220:LEU:HB2	1:E:223:ILE:HG12	1.83	0.60
1:E:282:VAL:O	1:E:313:ARG:NH2	2.35	0.60
1:F:40:VAL:HG13	1:F:106:LEU:HD13	1.83	0.60
1:D:220:LEU:HD21	1:D:240:HIS:CD2	2.38	0.59
1:D:185:MSE:HE3	1:D:194:LEU:HD21	1.85	0.59
1:A:180:THR:OG1	1:A:219:VAL:HA	2.03	0.58
1:D:282:VAL:O	1:D:313:ARG:NH1	2.36	0.58
1:C:227:THR:HG22	1:C:229:GLY:H	1.69	0.58
1:E:185:MSE:HE3	1:E:194:LEU:HD21	1.86	0.58
1:A:92:PHE:HD2	1:B:23:THR:HG23	1.68	0.57
1:F:109:VAL:HG22	1:F:113:SER:HB2	1.86	0.57
1:D:58:ILE:HG23	1:D:150:LEU:HD11	1.86	0.56
1:C:34:GLU:OE2	1:E:72:SER:OG	2.24	0.56
1:C:58:ILE:HG23	1:C:150:LEU:HD11	1.88	0.56
1:A:134:SER:HB3	1:A:137:ASP:H	1.71	0.55
1:B:40:VAL:HG13	1:B:106:LEU:HD13	1.88	0.55
1:D:46:ARG:HG2	1:D:104:PRO:HA	1.88	0.55
1:B:86:PHE:CD1	1:B:102:GLN:HG2	2.42	0.54
1:A:92:PHE:CD2	1:B:23:THR:HG23	2.42	0.54
1:F:107:ILE:HB	1:F:126:THR:HB	1.90	0.54
1:C:49:VAL:HG21	1:C:65:ILE:HD12	1.89	0.54
1:A:40:VAL:HG13	1:A:106:LEU:HD13	1.90	0.54
1:F:300:THR:O	1:F:304:ILE:HG12	2.07	0.54
1:C:40:VAL:HG13	1:C:106:LEU:HD13	1.90	0.54
1:F:58:ILE:HG23	1:F:150:LEU:HD11	1.90	0.54
1:F:17:TYR:HB2	1:F:145:PHE:CZ	2.43	0.54
1:E:58:ILE:HG23	1:E:150:LEU:HD11	1.89	0.54
1:A:219:VAL:HG13	1:A:220:LEU:HG	1.90	0.53
1:B:253:ALA:HB2	1:B:287:PRO:HB2	1.89	0.53
1:F:185:MSE:HE3	1:F:248:VAL:HG22	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:246:TRP:CH2	1:E:283:LEU:HD13	2.45	0.52
1:F:187:SER:OG	1:F:250:ARG:NH2	2.42	0.52
1:B:188:GLY:N	1:B:251:ASP:OD2	2.40	0.52
1:E:40:VAL:HG13	1:E:106:LEU:HD13	1.91	0.52
1:C:169:GLU:HG2	1:C:308:LEU:HD21	1.92	0.51
1:A:70:LEU:HD12	1:A:87:ASN:HB2	1.92	0.51
1:A:188:GLY:N	1:A:251:ASP:OD2	2.44	0.51
1:B:52:MSE:N	2:B:420:HOH:O	2.40	0.51
1:E:17:TYR:OH	1:E:132:ARG:NH2	2.43	0.50
1:E:26[B]:GLU:HG2	1:E:29:ALA:H	1.75	0.50
1:C:185:MSE:HE3	1:C:194:LEU:HD21	1.93	0.50
1:F:121:ASN:HA	1:F:126:THR:HG22	1.93	0.50
1:A:263:LYS:HB2	2:A:441:HOH:O	2.12	0.50
1:C:17:TYR:OH	1:C:132:ARG:NH1	2.39	0.49
2:C:411:HOH:O	1:F:23:THR:HG21	2.12	0.49
1:B:185:MSE:HE3	1:B:194:LEU:HD21	1.94	0.49
1:B:206:ARG:HD2	2:B:413:HOH:O	2.12	0.49
1:E:111:GLY:N	1:E:130:SER:HB3	2.21	0.49
1:F:290:TRP:CE2	1:F:303:MSE:HG2	2.47	0.49
1:A:34:GLU:OE2	1:D:72:SER:OG	2.24	0.49
1:F:246:TRP:CH2	1:F:283:LEU:HD13	2.47	0.49
1:E:285:LEU:HB3	1:E:290:TRP:CD1	2.47	0.48
1:C:230:GLU:HG2	1:F:228:HIS:C	2.38	0.48
1:D:180:THR:OG1	1:D:219:VAL:HA	2.11	0.48
1:A:86:PHE:CD2	1:A:102:GLN:HG2	2.48	0.48
1:B:180:THR:HG23	1:B:243:GLN:H	1.78	0.48
1:C:227:THR:HG22	1:C:229:GLY:N	2.29	0.48
1:A:246:TRP:CH2	1:A:283:LEU:HD13	2.49	0.48
1:D:17:TYR:OH	1:D:132:ARG:NH2	2.47	0.48
1:F:180:THR:OG1	1:F:219:VAL:HA	2.14	0.47
1:D:299:THR:HG22	1:D:303:MSE:HE3	1.95	0.47
1:F:185:MSE:HG3	1:F:194:LEU:HD21	1.96	0.47
1:B:92:PHE:HD1	1:D:23:THR:HG23	1.79	0.47
1:D:209:LEU:HD11	1:D:294[A]:PHE:CE2	2.50	0.47
1:C:26:GLU:CD	1:E:26[B]:GLU:HG3	2.39	0.47
1:A:135:TYR:CG	1:A:209:LEU:HD11	2.50	0.46
1:B:246:TRP:HE3	1:B:285:LEU:HD21	1.81	0.46
1:C:244:TRP:CE3	1:C:281:HIS:HB3	2.51	0.46
1:A:230:GLU:HG2	1:A:231:PRO:HD2	1.96	0.46
1:F:245:LEU:HB3	1:F:282:VAL:HG12	1.97	0.46
1:B:180:THR:OG1	1:B:219:VAL:HA	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:220:LEU:HB2	1:C:223:ILE:HG12	1.96	0.46
1:D:210:ILE:O	1:D:216:VAL:HG13	2.16	0.46
1:E:244:TRP:CE3	1:E:281:HIS:HB3	2.51	0.46
1:D:206:ARG:HA	1:D:294[A]:PHE:CE2	2.51	0.45
1:B:246:TRP:CE3	1:B:285:LEU:HD21	2.52	0.45
1:B:290:TRP:CE2	1:B:303:MSE:HG2	2.52	0.45
1:C:72:SER:OG	1:F:34:GLU:OE1	2.25	0.45
1:F:167:ILE:HG13	1:F:304:ILE:HD11	1.99	0.45
1:B:56:ASP:OD1	1:B:298:GLU:HB2	2.17	0.45
1:B:92:PHE:CD1	1:D:23:THR:HG23	2.52	0.45
1:A:246:TRP:CZ3	1:A:283:LEU:HD13	2.52	0.45
1:E:139:LEU:HD21	1:E:163:LEU:HD23	1.99	0.45
1:F:309:GLN:NE2	2:F:402:HOH:O	2.49	0.45
1:C:140:GLU:HA	1:C:160:LEU:HD11	1.98	0.45
1:D:96:LEU:HD11	1:D:119:ASP:HB3	1.98	0.44
1:E:298:GLU:O	1:E:302:ILE:HG12	2.17	0.44
1:C:298:GLU:O	1:C:302:ILE:HG13	2.18	0.44
1:D:180:THR:HG23	1:D:243:GLN:H	1.82	0.44
1:D:246:TRP:CH2	1:D:283:LEU:HD13	2.52	0.44
1:F:18:THR:HB	1:F:37:THR:OG1	2.18	0.44
1:F:244:TRP:CE3	1:F:281:HIS:HB3	2.52	0.44
1:C:59:GLY:HA3	1:C:80:TYR:CG	2.51	0.44
1:C:246:TRP:CH2	1:C:283:LEU:HD13	2.53	0.44
1:F:112:ARG:NH2	1:F:292:ILE:HA	2.32	0.44
1:B:96:LEU:HD11	1:B:119:ASP:HB3	2.00	0.44
1:E:59:GLY:HA3	1:E:80:TYR:CG	2.53	0.44
1:B:19:TRP:CZ2	1:B:128:ASP:HB3	2.53	0.44
1:B:134:SER:CB	1:B:137:ASP:H	2.26	0.44
1:F:180:THR:HG22	1:F:243[B]:GLN:OE1	2.18	0.44
1:A:136:LEU:HD23	1:A:136:LEU:HA	1.83	0.43
1:E:17:TYR:HB2	1:E:145:PHE:CZ	2.53	0.43
1:C:223:ILE:O	1:C:227:THR:HB	2.18	0.43
1:E:160:LEU:HD23	1:E:160:LEU:HA	1.85	0.43
1:D:135:TYR:CG	1:D:209:LEU:HD21	2.53	0.43
1:C:56:ASP:OD1	1:C:298:GLU:HB2	2.19	0.43
1:F:237:LEU:HD12	1:F:237:LEU:HA	1.80	0.43
1:B:201:ASP:HA	1:B:202:PRO:HD3	1.73	0.43
1:E:220:LEU:HD21	1:E:240:HIS:CD2	2.54	0.42
1:A:285:LEU:HB3	1:A:290:TRP:CD1	2.54	0.42
1:C:49:VAL:HG13	1:C:108:VAL:HG13	2.00	0.42
1:E:52:MSE:HB3	1:E:293:VAL:HG22	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:197:ASN:C	1:F:197:ASN:HD22	2.28	0.42
1:F:307:VAL:O	1:F:310:VAL:HG22	2.18	0.42
1:D:307:VAL:O	1:D:310:VAL:HG22	2.19	0.42
1:A:72:SER:OG	1:B:34:GLU:OE2	2.28	0.42
1:D:280:ASP:O	1:D:313:ARG:HG2	2.20	0.42
1:F:210:ILE:HD12	1:F:210:ILE:HA	1.95	0.42
1:F:217:GLN:HA	1:F:218:PRO:HD3	1.85	0.42
1:B:243:GLN:HE21	1:B:281:HIS:CE1	2.36	0.42
1:D:274:THR:O	1:D:278:THR:HG23	2.19	0.42
1:F:59:GLY:HA3	1:F:80:TYR:CG	2.55	0.42
1:B:285:LEU:HB3	1:B:290:TRP:CD1	2.55	0.42
1:C:90:THR:HG22	1:C:93:GLU:H	1.83	0.42
1:C:127:ILE:HD12	1:C:129:LEU:HD21	2.01	0.42
1:D:41:PRO:HG2	1:D:105:ASP:O	2.19	0.42
1:C:187:SER:OG	1:C:250:ARG:NH2	2.53	0.42
1:E:52:MSE:HE3	1:E:72:SER:O	2.20	0.42
1:F:178:SER:HB2	1:F:179:GLY:H	1.69	0.42
1:F:185:MSE:HE2	1:F:185:MSE:HB3	1.92	0.42
1:A:52:MSE:N	2:A:443:HOH:O	2.50	0.41
1:B:217:GLN:HA	1:B:218:PRO:HD3	1.92	0.41
1:B:233:SER:HB2	1:B:235:GLU:HG2	2.01	0.41
1:E:70:LEU:O	1:E:73:VAL:HG22	2.20	0.41
1:A:46:ARG:HB3	1:A:104:PRO:HA	2.01	0.41
1:D:134:SER:HB2	1:D:137:ASP:H	1.85	0.41
1:A:307:VAL:O	1:A:310:VAL:HG22	2.21	0.41
1:E:40:VAL:HA	1:E:41:PRO:HD3	1.90	0.41
1:F:235:GLU:H	1:F:235:GLU:HG2	1.47	0.41
1:B:72:SER:OG	1:D:34:GLU:OE2	2.35	0.41
1:C:166:GLY:HA3	1:C:304:ILE:HD13	2.02	0.41
1:C:180:THR:OG1	1:C:219:VAL:HA	2.20	0.41
1:D:210:ILE:HD12	1:D:210:ILE:HA	1.85	0.41
1:B:83:ASP:N	1:B:83:ASP:OD1	2.52	0.41
1:B:246:TRP:CH2	1:B:283:LEU:HD13	2.56	0.41
1:F:185:MSE:HE1	1:F:291:TYR:CD1	2.55	0.41
1:E:237:LEU:HD12	1:E:237:LEU:HA	1.86	0.41
1:B:97:ILE:HD11	1:D:122:GLU:CD	2.47	0.40
1:B:210:ILE:HA	1:B:210:ILE:HD12	1.85	0.40
1:C:228:HIS:C	1:E:230:GLU:HG2	2.46	0.40
1:E:209:LEU:HD12	1:E:209:LEU:HA	1.87	0.40
1:E:307:VAL:O	1:E:310:VAL:HG22	2.22	0.40
1:C:30:ASP:HA	1:C:31:PRO:HD2	1.97	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:166:GLY:HA3	1:E:304:ILE:HD13	2.04	0.40
1:B:269:GLU:HG3	2:E:415:HOH:O	2.21	0.40
1:D:40:VAL:HG13	1:D:106:LEU:HD13	2.03	0.40
1:A:19:TRP:CZ2	1:A:128:ASP:HB3	2.57	0.40
1:A:56:ASP:OD1	1:A:298[A]:GLU:HB2	2.20	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	302/315 (96%)	291 (96%)	11 (4%)	0	100	100
1	B	302/315 (96%)	294 (97%)	7 (2%)	1 (0%)	36	64
1	C	299/315 (95%)	289 (97%)	10 (3%)	0	100	100
1	D	301/315 (96%)	292 (97%)	9 (3%)	0	100	100
1	E	300/315 (95%)	285 (95%)	15 (5%)	0	100	100
1	F	300/315 (95%)	285 (95%)	14 (5%)	1 (0%)	36	64
All	All	1804/1890 (95%)	1736 (96%)	66 (4%)	2 (0%)	48	76

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	F	178	SER
1	B	178	SER

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	234/238 (98%)	221 (94%)	13 (6%)	19	48
1	B	234/238 (98%)	219 (94%)	15 (6%)	16	43
1	C	231/238 (97%)	205 (89%)	26 (11%)	5	18
1	D	233/238 (98%)	218 (94%)	15 (6%)	16	43
1	E	232/238 (98%)	201 (87%)	31 (13%)	4	12
1	F	232/238 (98%)	207 (89%)	25 (11%)	6	20
All	All	1396/1428 (98%)	1271 (91%)	125 (9%)	9	27

All (125) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	SER
1	A	45	GLN
1	A	109	VAL
1	A	114	SER
1	A	132	ARG
1	A	190	GLN
1	A	210	ILE
1	A	264	VAL
1	A	269	GLU
1	A	274	THR
1	A	298[A]	GLU
1	A	298[B]	GLU
1	A	310	VAL
1	B	14	THR
1	B	21	ARG
1	B	23	THR
1	B	25	THR
1	B	46	ARG
1	B	109	VAL
1	B	148	LYS
1	B	201	ASP
1	B	203	ARG
1	B	209	LEU
1	B	219	VAL
1	B	227	THR
1	B	235	GLU

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Mol	Chain	Res	Type
1	B	269	GLU
1	B	310	VAL
1	C	16	SER
1	C	25	THR
1	C	36	THR
1	C	39	GLU
1	C	81	LEU
1	C	83	ASP
1	C	90	THR
1	C	108	VAL
1	C	109	VAL
1	C	113	SER
1	C	114	SER
1	C	160	LEU
1	C	162	GLU
1	C	163	LEU
1	C	180	THR
1	C	192	SER
1	C	195	SER
1	C	203	ARG
1	C	227	THR
1	C	237	LEU
1	C	264	VAL
1	C	279	GLU
1	C	282	VAL
1	C	283	LEU
1	C	305	ASP
1	C	310	VAL
1	D	14	THR
1	D	23	THR
1	D	25	THR
1	D	26	GLU
1	D	46	ARG
1	D	83	ASP
1	D	84	ASP
1	D	109	VAL
1	D	122	GLU
1	D	132	ARG
1	D	159	VAL
1	D	200	ASN
1	D	216	VAL
1	D	310	VAL

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Mol	Chain	Res	Type
1	D	313	ARG
1	E	14	THR
1	E	16	SER
1	E	26[A]	GLU
1	E	26[B]	GLU
1	E	27	GLU
1	E	36	THR
1	E	52	MSE
1	E	70	LEU
1	E	81	LEU
1	E	114	SER
1	E	130	SER
1	E	134	SER
1	E	141	GLN
1	E	157	GLU
1	E	159	VAL
1	E	162	GLU
1	E	163	LEU
1	E	191	LEU
1	E	203	ARG
1	E	209	LEU
1	E	216	VAL
1	E	224	LYS
1	E	230	GLU
1	E	235	GLU
1	E	237	LEU
1	E	248	VAL
1	E	257	GLU
1	E	263	LYS
1	E	269	GLU
1	E	310	VAL
1	E	313	ARG
1	F	16	SER
1	F	18	THR
1	F	25	THR
1	F	27	GLU
1	F	32	VAL
1	F	61	LEU
1	F	70	LEU
1	F	81	LEU
1	F	87	ASN
1	F	91	LEU

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Mol	Chain	Res	Type
1	F	109	VAL
1	F	153	GLU
1	F	190	GLN
1	F	223	ILE
1	F	235	GLU
1	F	237	LEU
1	F	248	VAL
1	F	263	LYS
1	F	264	VAL
1	F	265	VAL
1	F	275	THR
1	F	282	VAL
1	F	283	LEU
1	F	310	VAL
1	F	313	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	272	ASN
1	A	281	HIS
1	B	103	GLN
1	B	141	GLN
1	B	200	ASN
1	B	217	GLN
1	B	243	GLN
1	B	309	GLN
1	C	272	ASN
1	D	240	HIS
1	D	281	HIS
1	D	309	GLN
1	E	141	GLN
1	E	240	HIS
1	F	22	ASN
1	F	141	GLN

5.3.3 RNA

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 oligosaccharides 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	299/315 (94%)	-1.29	0 100 100	19, 42, 78, 143	2 (0%)
1	B	299/315 (94%)	-1.31	1 (0%) 90 87	22, 42, 76, 138	2 (0%)
1	C	298/315 (94%)	-0.65	0 100 100	47, 68, 109, 155	0
1	D	299/315 (94%)	-1.31	0 100 100	18, 42, 75, 153	1 (0%)
1	E	298/315 (94%)	-0.69	0 100 100	31, 69, 106, 154	1 (0%)
1	F	298/315 (94%)	-0.72	0 100 100	34, 69, 108, 150	1 (0%)
All	All	1791/1890 (94%)	-1.00	1 (0%) 92 90	18, 58, 98, 155	7 (0%)

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	12	PRO	2.3

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 oligosaccharides 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.