



Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 01:30 pm BST

PDB ID : 4MX8
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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
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) : 2.11

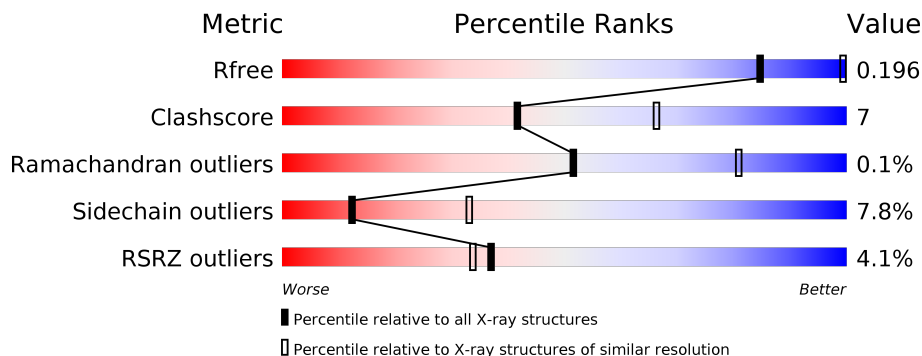
1 Overall quality at a glance

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}	130704	2307 (2.94-2.90)
Clashscore	141614	2531 (2.94-2.90)
Ramachandran outliers	138981	2462 (2.94-2.90)
Sidechain outliers	138945	2464 (2.94-2.90)
RSRZ outliers	127900	2248 (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 on 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	
1	B	315	
1	C	315	
1	D	315	
1	E	315	
1	F	315	

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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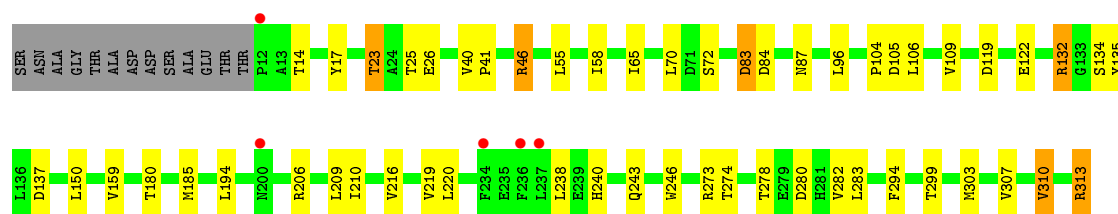
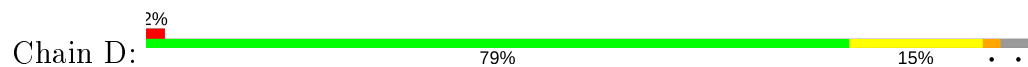
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Chain	Residue	Modelled	Actual	Comment	Reference
F	1	ALA	-	EXPRESSION TAG	UNP D1BRG9

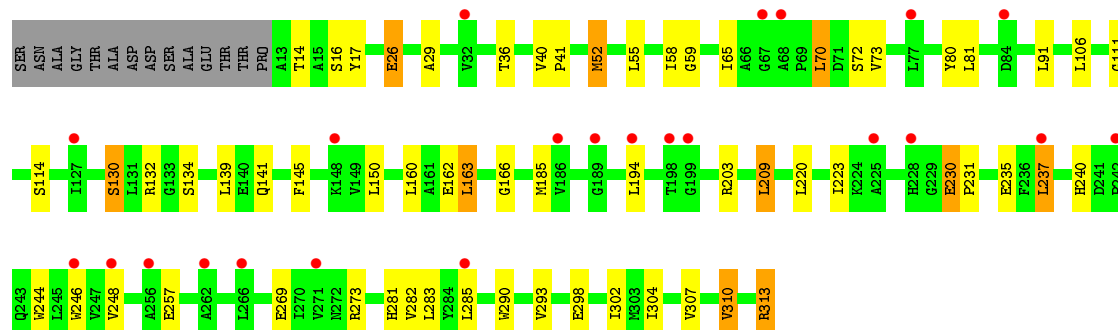
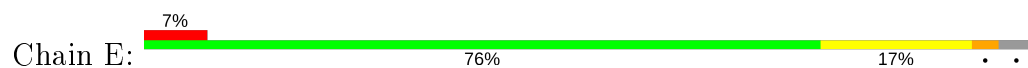
- Molecule 2 is water.

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

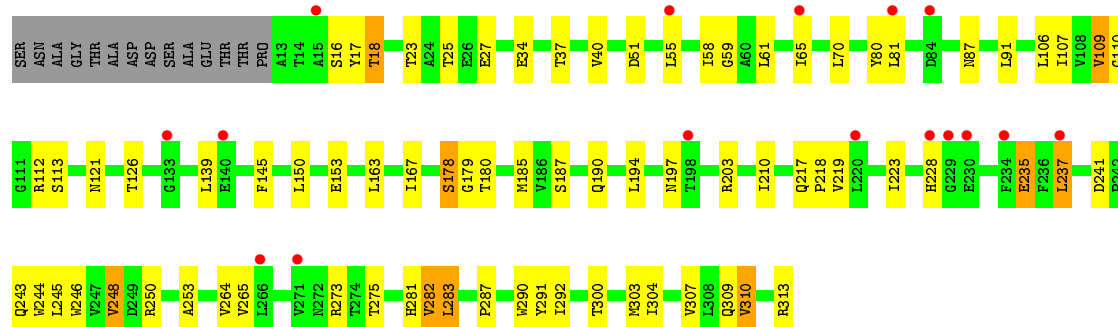
- Molecule 1: Periplasmic binding protein



- Molecule 1: Periplasmic binding protein



- Molecule 1: Periplasmic binding protein



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 50.19 – 2.91	Depositor EDS
% Data completeness (in resolution range)	96.6 (49.37-2.91) 96.3 (50.19-2.91)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.34 (at 2.91Å)	Xtriage
Refinement program	PHENIX (phenix.refine: dev_1367)	Depositor
R, R_{free}	0.159 , 0.196 0.159 , 0.196	Depositor DCC
R_{free} test set	4014 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	54.1	Xtriage
Anisotropy	0.053	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 36.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
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 $1/2^*h-3/2^*k, -1/2^*h-1/2^*k, -1/2^*h+1/2^*k-l$	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	13799	wwPDB-VP
Average B, all atoms (Å ²)	59.0	wwPDB-VP

Xtriage'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	0.45	0/2313	0.61	0/3161
1	B	0.46	0/2311	0.61	0/3158
1	C	0.38	0/2284	0.54	0/3122
1	D	0.46	0/2304	0.62	0/3149
1	E	0.37	0/2293	0.54	0/3134
1	F	0.37	0/2293	0.53	0/3134
All	All	0.42	0/13798	0.58	0/18858

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 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	30	0
1	B	2272	0	2187	35	0
1	C	2246	0	2164	34	0
1	D	2264	0	2180	30	0
1	E	2255	0	2169	32	0
1	F	2255	0	2171	39	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

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	E	18	0	0	1	0
2	F	13	0	0	1	0
All	All	13799	0	13056	186	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 (186) 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
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:OD1	1:C:83:ASP:N	2.22	0.62
1:D:238:LEU:O	1:F:273:ARG:NH1	2.32	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:58:ILE:HG23	1:C:150:LEU:HD11	1.88	0.56
1:C:34:GLU:OE2	1:E:72:SER:OG	2.24	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
1:C:72:SER:OG	1:F:34:GLU:OE1	2.25	0.53
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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: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: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:A:34:GLU:OE2	1:D:72:SER:OG	2.24	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:D:209:LEU:HD11	1:D:294[A]:PHE:CE2	2.50	0.47
1:B:92:PHE:HD1	1:D:23:THR:HG23	1.79	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
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: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:C:230:GLU:HG2	1:F:228:HIS:C	2.38	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:HA	1:A:136:LEU:HD23	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:HA	1:E:160:LEU:HD23	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
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:C:26:GLU:CD	1:E:26[B]:GLU:HG3	2.39	0.42
1:D:280:ASP:O	1:D:313:ARG:HG2	2.20	0.42
1:F:210:ILE:HA	1:F:210:ILE:HD12	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:127:ILE:HD12	1:C:129:LEU:HD21	2.01	0.42
1:C:90:THR:HG22	1:C:93:GLU:H	1.83	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:HB3	1:F:185:MSE:HE2	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:HG2	1:F:235:GLU:H	1.47	0.41
1:C:180:THR:OG1	1:C:219:VAL:HA	2.20	0.41
1:C:166:GLY:HA3	1:C:304:ILE:HD13	2.02	0.41
1:D:210:ILE:HA	1:D:210:ILE:HD12	1.85	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:B:83:ASP:N	1:B:83:ASP:OD1	2.52	0.41
1:E:237:LEU:HA	1:E:237:LEU:HD12	1.86	0.41
1:B:210:ILE:HD12	1:B:210:ILE:HA	1.85	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
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%)	41	70
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%)	41	70
All	All	1804/1890 (95%)	1736 (96%)	66 (4%)	2 (0%)	51	81

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%)	21	50
1	B	234/238 (98%)	221 (94%)	13 (6%)	21	50
1	C	231/238 (97%)	208 (90%)	23 (10%)	7	22
1	D	233/238 (98%)	220 (94%)	13 (6%)	21	50
1	E	232/238 (98%)	208 (90%)	24 (10%)	7	21
1	F	232/238 (98%)	207 (89%)	25 (11%)	6	19
All	All	1396/1428 (98%)	1285 (92%)	111 (8%)	12	32

All (111) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	SER
1	A	46	ARG
1	A	83	ASP
1	A	109	VAL
1	A	114	SER
1	A	132	ARG
1	A	190	GLN
1	A	210	ILE
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	23	THR
1	B	25	THR
1	B	46	ARG
1	B	83	ASP
1	B	109	VAL
1	B	148	LYS
1	B	201	ASP
1	B	209	LEU
1	B	219	VAL
1	B	235	GLU
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	103	GLN
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	192	SER
1	C	195	SER
1	C	203	ARG

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Mol	Chain	Res	Type
1	C	227	THR
1	C	237	LEU
1	C	279	GLU
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	310	VAL
1	D	313	ARG
1	E	14	THR
1	E	16	SER
1	E	26[A]	GLU
1	E	26[B]	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	162	GLU
1	E	163	LEU
1	E	203	ARG
1	E	209	LEU
1	E	230	GLU
1	E	235	GLU
1	E	237	LEU
1	E	248	VAL
1	E	257	GLU
1	E	269	GLU
1	E	310	VAL

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Mol	Chain	Res	Type
1	E	313	ARG
1	F	16	SER
1	F	18	THR
1	F	25	THR
1	F	27	GLU
1	F	61	LEU
1	F	70	LEU
1	F	81	LEU
1	F	87	ASN
1	F	91	LEU
1	F	109	VAL
1	F	153	GLU
1	F	190	GLN
1	F	197	ASN
1	F	203	ARG
1	F	223	ILE
1	F	235	GLU
1	F	237	LEU
1	F	248	VAL
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

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

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

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no carbohydrates 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%)	0.69	4 (1%) 77 77	23, 42, 78, 143	0
1	B	299/315 (94%)	0.66	4 (1%) 77 77	22, 42, 76, 138	0
1	C	298/315 (94%)	0.78	22 (7%) 14 12	47, 68, 109, 155	0
1	D	299/315 (94%)	0.61	5 (1%) 70 70	26, 42, 75, 153	0
1	E	298/315 (94%)	0.80	23 (7%) 13 11	46, 69, 106, 154	0
1	F	298/315 (94%)	0.71	16 (5%) 25 22	45, 69, 108, 150	0
All	All	1791/1890 (94%)	0.71	74 (4%) 37 34	22, 59, 99, 155	0

All (74) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	200	ASN	7.2
1	F	228	HIS	4.3
1	F	81	LEU	4.1
1	F	266	LEU	4.0
1	C	269	GLU	4.0
1	B	228	HIS	4.0
1	E	266	LEU	3.9
1	F	271	VAL	3.7
1	C	99	ILE	3.7
1	F	220	LEU	3.6
1	A	230	GLU	3.4
1	C	228	HIS	3.4
1	E	189	GLY	3.3
1	F	65	ILE	3.3
1	F	234	PHE	3.3
1	C	236	PHE	3.2
1	F	229	GLY	3.1
1	E	285	LEU	3.1
1	C	226	ALA	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	200	ASN	3.1
1	E	271	VAL	3.1
1	E	68	ALA	3.0
1	C	260	GLN	3.0
1	B	200	ASN	2.9
1	E	225	ALA	2.9
1	C	81	LEU	2.8
1	F	237	LEU	2.7
1	F	198	THR	2.7
1	C	284	TYR	2.6
1	E	237	LEU	2.6
1	E	199	GLY	2.6
1	F	140	GLU	2.6
1	C	237	LEU	2.6
1	E	77	LEU	2.6
1	C	297	VAL	2.6
1	C	256	ALA	2.6
1	D	12	PRO	2.5
1	E	32	VAL	2.5
1	E	198	THR	2.4
1	C	229	GLY	2.4
1	E	148	LYS	2.4
1	C	157	GLU	2.4
1	F	230	GLU	2.3
1	D	234	PHE	2.3
1	E	84	ASP	2.3
1	E	228	HIS	2.3
1	F	84	ASP	2.2
1	C	92	PHE	2.2
1	E	248	VAL	2.2
1	C	253	ALA	2.2
1	C	266	LEU	2.2
1	B	225	ALA	2.2
1	F	15	ALA	2.2
1	E	246	TRP	2.2
1	F	133	GLY	2.1
1	A	294[A]	PHE	2.1
1	C	17	TYR	2.1
1	E	67	GLY	2.1
1	F	55	LEU	2.1
1	E	242	PRO	2.1
1	E	194	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	132[A]	ARG	2.1
1	E	256	ALA	2.1
1	C	238	LEU	2.1
1	C	270	ILE	2.1
1	E	186	VAL	2.0
1	D	236	PHE	2.0
1	C	210	ILE	2.0
1	E	127	ILE	2.0
1	A	228	HIS	2.0
1	C	283	LEU	2.0
1	E	262	ALA	2.0
1	C	285	LEU	2.0
1	D	237	LEU	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.