



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:39 AM GMT

PDB ID : 2XUT  
Title : CRYSTAL STRUCTURE OF A PROTON DEPENDENT OLIGOPEPTIDE (POT) FAMILY TRANSPORTER.  
Authors : Newstead, S.; Drew, D.; Cameron, A.D.; Postis, V.L.; Xia, X.; Fowler, P.W.; Ingram, J.C.; Carpenter, E.P.; Sansom, M.S.P.; McPherson, M.J.; Baldwin, S.A.; Iwata, S.  
Deposited on : 2010-10-21  
Resolution : 3.62 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](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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The following versions of software and data (see [references ⓘ](#)) 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

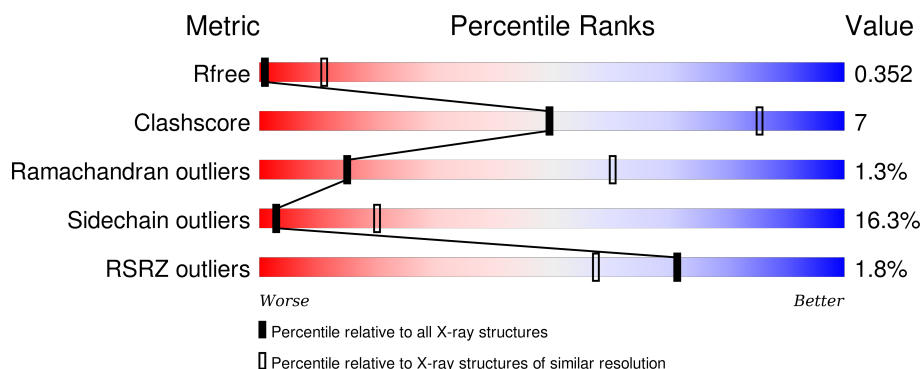
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.62 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	1093 (3.80-3.44)
Clashscore	102246	1043 (3.78-3.46)
Ramachandran outliers	100387	1003 (3.78-3.46)
Sidechain outliers	100360	1003 (3.78-3.46)
RSRZ outliers	91569	1100 (3.80-3.44)

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  $\geq 3$ , 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  $\leq 5\%$ . 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	524	<div> <div> <div>2%</div> <div>62%</div> <div>23%</div> <div>•</div> <div>13%</div> </div> </div>
1	B	524	<div> <div>2%</div> <div>63%</div> <div>23%</div> <div>•</div> <div>13%</div> </div>

## 2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 10533 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 PROTON/PEPTIDE SYMPORTER FAMILY PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	456	Total	C	N	O	S	0	0	0
			3511	2345	553	589	24			
1	B	456	Total	C	N	O	S	0	0	0
			3511	2345	553	589	24			
1	C	456	Total	C	N	O	S	0	0	0
			3511	2345	553	589	24			

There are 30 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	517	GLY	-	EXPRESSION TAG	UNP Q8EKT7
A	518	SER	-	EXPRESSION TAG	UNP Q8EKT7
A	519	GLU	-	EXPRESSION TAG	UNP Q8EKT7
A	520	ASN	-	EXPRESSION TAG	UNP Q8EKT7
A	521	LEU	-	EXPRESSION TAG	UNP Q8EKT7
A	522	TYR	-	EXPRESSION TAG	UNP Q8EKT7
A	523	PHE	-	EXPRESSION TAG	UNP Q8EKT7
A	524	GLN	-	EXPRESSION TAG	UNP Q8EKT7
A	2	ASN	THR	ENGINEERED MUTATION	UNP Q8EKT7
A	3	SER	THR	ENGINEERED MUTATION	UNP Q8EKT7
B	517	GLY	-	EXPRESSION TAG	UNP Q8EKT7
B	518	SER	-	EXPRESSION TAG	UNP Q8EKT7
B	519	GLU	-	EXPRESSION TAG	UNP Q8EKT7
B	520	ASN	-	EXPRESSION TAG	UNP Q8EKT7
B	521	LEU	-	EXPRESSION TAG	UNP Q8EKT7
B	522	TYR	-	EXPRESSION TAG	UNP Q8EKT7
B	523	PHE	-	EXPRESSION TAG	UNP Q8EKT7
B	524	GLN	-	EXPRESSION TAG	UNP Q8EKT7
B	2	ASN	THR	ENGINEERED MUTATION	UNP Q8EKT7
B	3	SER	THR	ENGINEERED MUTATION	UNP Q8EKT7
C	517	GLY	-	EXPRESSION TAG	UNP Q8EKT7
C	518	SER	-	EXPRESSION TAG	UNP Q8EKT7
C	519	GLU	-	EXPRESSION TAG	UNP Q8EKT7

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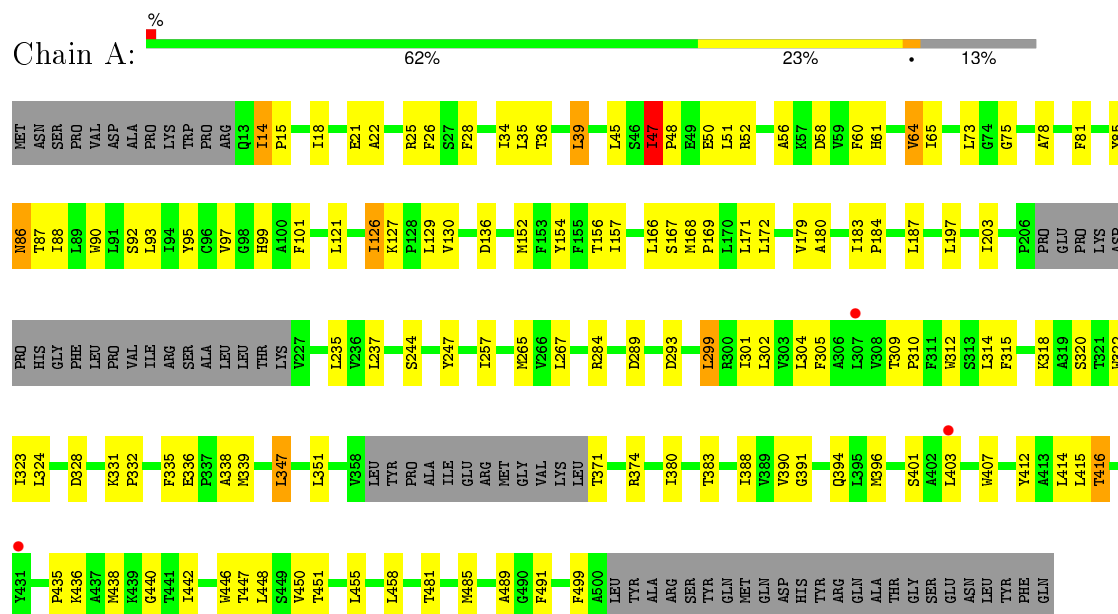
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Chain	Residue	Modelled	Actual	Comment	Reference
C	520	ASN	-	EXPRESSION TAG	UNP Q8EKT7
C	521	LEU	-	EXPRESSION TAG	UNP Q8EKT7
C	522	TYR	-	EXPRESSION TAG	UNP Q8EKT7
C	523	PHE	-	EXPRESSION TAG	UNP Q8EKT7
C	524	GLN	-	EXPRESSION TAG	UNP Q8EKT7
C	2	ASN	THR	ENGINEERED MUTATION	UNP Q8EKT7
C	3	SER	THR	ENGINEERED MUTATION	UNP Q8EKT7

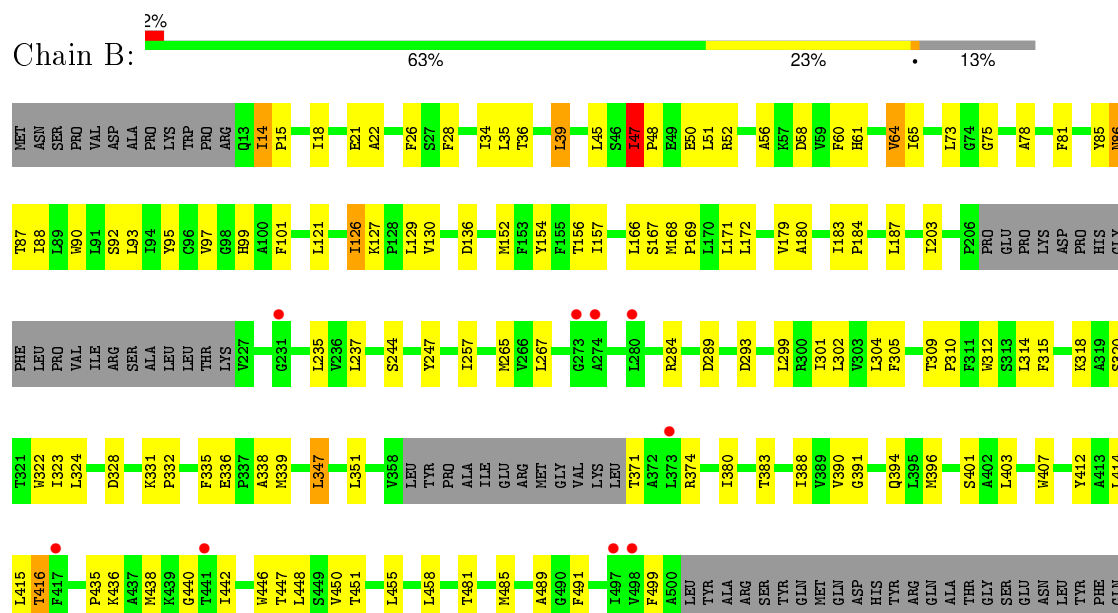
### 3 Residue-property plots

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.

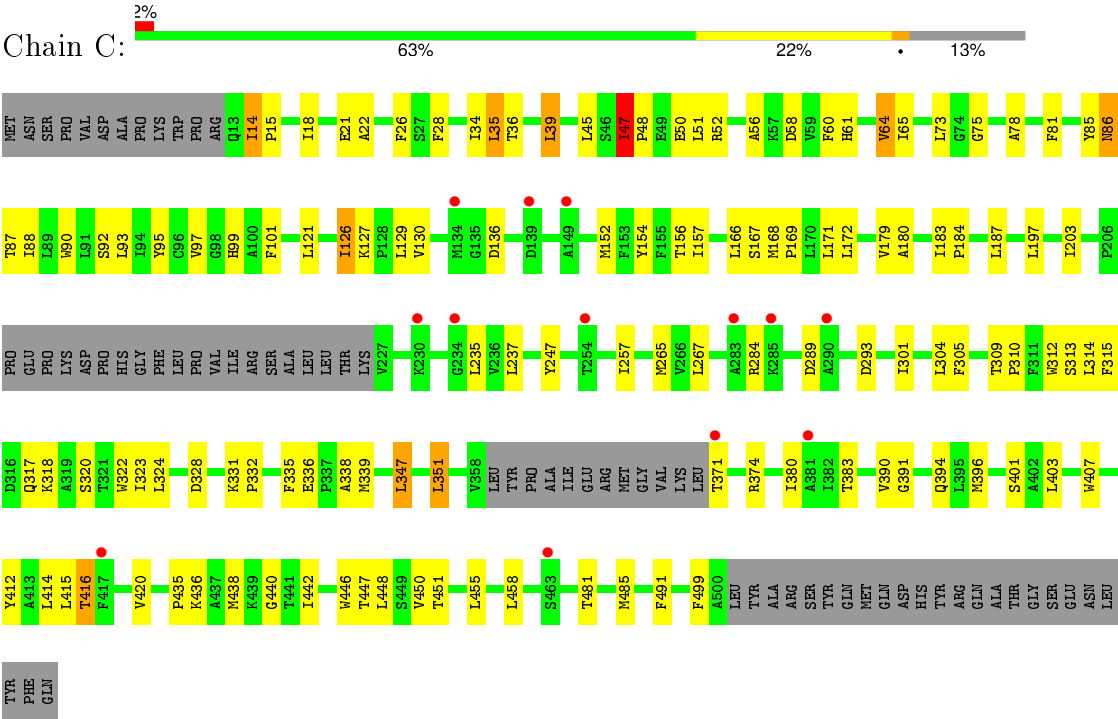
#### • Molecule 1: PROTON/PEPTIDE SYMPORTER FAMILY PROTEIN



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● Molecule 1: PROTON/PEPTIDE SYMPORTER FAMILY PROTEIN



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	159.40Å 159.40Å 152.94Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	35.47 – 3.62 34.51 – 3.62	Depositor EDS
% Data completeness (in resolution range)	(Not available) (35.47-3.62) 58.5 (34.51-3.62)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.39 (at 3.66Å)	Xtriage
Refinement program	BUSTER 2.8.0	Depositor
R, $R_{free}$	0.277 , 0.296 0.325 , 0.352	Depositor DCC
$R_{free}$ test set	1688 reflections (6.20%)	DCC
Wilson B-factor (Å <sup>2</sup> )	118.1	Xtriage
Anisotropy	0.043	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.26 , 273.7	EDS
Estimated twinning fraction	0.110 for -h,-k,l 0.135 for h,-h-k,-l 0.139 for -k,-h,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.39$ , $\langle L^2 \rangle = 0.22$	Xtriage
Outliers	0 of 28894 reflections	Xtriage
$F_o, F_c$ correlation	0.84	EDS
Total number of atoms	10533	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	252.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 15.74% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  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

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	0.49	0/3604	0.69	0/4893
1	B	0.49	0/3604	0.69	0/4893
1	C	0.48	0/3604	0.69	0/4893
All	All	0.48	0/10812	0.69	0/14679

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	3511	0	3592	51	0
1	B	3511	0	3592	51	0
1	C	3511	0	3592	50	0
All	All	10533	0	10776	152	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 (152) 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:14:ILE:H	1:B:15:PRO:HD2	1.39	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:14:ILE:H	1:A:15:PRO:HD2	1.40	0.87
1:C:14:ILE:H	1:C:15:PRO:HD2	1.39	0.87
1:A:48:PRO:HD2	1:A:51:LEU:HB2	1.59	0.84
1:B:48:PRO:HD2	1:B:51:LEU:HB2	1.59	0.84
1:C:48:PRO:HD2	1:C:51:LEU:HB2	1.59	0.82
1:A:318:LYS:HA	1:A:322:TRP:CD1	2.26	0.71
1:C:318:LYS:HA	1:C:322:TRP:CD1	2.26	0.70
1:B:318:LYS:HA	1:B:322:TRP:CD1	2.27	0.70
1:C:14:ILE:H	1:C:15:PRO:CD	2.05	0.69
1:B:14:ILE:H	1:B:15:PRO:CD	2.05	0.68
1:A:14:ILE:H	1:A:15:PRO:CD	2.05	0.68
1:C:435:PRO:HB2	1:C:438:MET:HB3	1.76	0.67
1:B:435:PRO:HB2	1:B:438:MET:HB3	1.77	0.66
1:A:435:PRO:HB2	1:A:438:MET:HB3	1.77	0.65
1:A:347:LEU:HB3	1:A:416:THR:HG21	1.82	0.61
1:B:347:LEU:HB3	1:B:416:THR:HG21	1.82	0.61
1:C:347:LEU:HB3	1:C:416:THR:HG21	1.83	0.61
1:A:56:ALA:O	1:A:60:PHE:HB3	2.00	0.61
1:B:18:ILE:HG21	1:B:152:MET:HB3	1.82	0.61
1:C:56:ALA:O	1:C:60:PHE:HB3	2.01	0.61
1:B:56:ALA:O	1:B:60:PHE:HB3	2.01	0.61
1:A:18:ILE:HG21	1:A:152:MET:HB3	1.83	0.60
1:C:18:ILE:HG21	1:C:152:MET:HB3	1.82	0.60
1:B:28:PHE:CE2	1:B:127:LYS:NZ	2.64	0.60
1:C:47:ILE:HB	1:C:48:PRO:HD3	1.84	0.59
1:A:47:ILE:HB	1:A:48:PRO:HD3	1.84	0.59
1:B:47:ILE:HB	1:B:48:PRO:HD3	1.85	0.57
1:C:28:PHE:CE2	1:C:127:LYS:NZ	2.67	0.56
1:A:171:LEU:HD23	1:A:180:ALA:HA	1.87	0.56
1:B:171:LEU:HD23	1:B:180:ALA:HA	1.87	0.56
1:C:171:LEU:HD23	1:C:180:ALA:HA	1.88	0.56
1:A:28:PHE:CE2	1:A:127:LYS:NZ	2.67	0.56
1:A:136:ASP:HB3	1:A:203:ILE:HD11	1.89	0.55
1:C:47:ILE:HB	1:C:48:PRO:CD	2.38	0.54
1:B:136:ASP:HB3	1:B:203:ILE:HD11	1.89	0.54
1:A:47:ILE:HB	1:A:48:PRO:CD	2.38	0.54
1:C:136:ASP:HB3	1:C:203:ILE:HD11	1.90	0.54
1:B:47:ILE:HB	1:B:48:PRO:CD	2.38	0.54
1:C:184:PRO:HA	1:C:187:LEU:HD12	1.91	0.53
1:B:14:ILE:N	1:B:15:PRO:HD2	2.18	0.53
1:A:184:PRO:HA	1:A:187:LEU:HD12	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:75:GLY:HA2	1:B:78:ALA:HB3	1.91	0.53
1:B:184:PRO:HA	1:B:187:LEU:HD12	1.91	0.53
1:A:126:ILE:O	1:A:129:LEU:HB2	2.09	0.52
1:C:14:ILE:N	1:C:15:PRO:HD2	2.18	0.52
1:A:75:GLY:HA2	1:A:78:ALA:HB3	1.91	0.52
1:C:126:ILE:O	1:C:129:LEU:HB2	2.10	0.52
1:C:34:ILE:HD11	1:C:168:MET:HG3	1.90	0.52
1:C:75:GLY:HA2	1:C:78:ALA:HB3	1.92	0.51
1:A:34:ILE:HD11	1:A:168:MET:HG3	1.91	0.51
1:C:315:PHE:O	1:C:318:LYS:HG2	2.10	0.51
1:B:34:ILE:HD11	1:B:168:MET:HG3	1.91	0.51
1:A:315:PHE:O	1:A:318:LYS:HG2	2.11	0.50
1:B:315:PHE:O	1:B:318:LYS:HG2	2.11	0.50
1:B:126:ILE:O	1:B:129:LEU:HB2	2.11	0.50
1:A:14:ILE:N	1:A:15:PRO:HD2	2.19	0.50
1:C:22:ALA:HA	1:C:156:THR:HG23	1.94	0.50
1:C:86:ASN:ND2	1:C:86:ASN:H	2.10	0.50
1:A:86:ASN:ND2	1:A:86:ASN:H	2.10	0.49
1:B:86:ASN:H	1:B:86:ASN:ND2	2.10	0.49
1:A:22:ALA:HA	1:A:156:THR:HG23	1.95	0.48
1:C:61:HIS:HA	1:C:64:VAL:HG23	1.95	0.48
1:B:22:ALA:HA	1:B:156:THR:HG23	1.95	0.48
1:C:126:ILE:HA	1:C:129:LEU:HD22	1.95	0.48
1:B:320:SER:O	1:B:323:ILE:HG22	2.13	0.48
1:A:126:ILE:HA	1:A:129:LEU:HD22	1.95	0.48
1:B:391:GLY:HA2	1:B:485:MET:HB3	1.96	0.48
1:C:391:GLY:HA2	1:C:485:MET:HB3	1.96	0.48
1:A:391:GLY:HA2	1:A:485:MET:HB3	1.96	0.48
1:A:61:HIS:HA	1:A:64:VAL:HG23	1.96	0.48
1:C:314:LEU:HD12	1:C:383:THR:HB	1.96	0.48
1:A:314:LEU:HD12	1:A:383:THR:HB	1.96	0.47
1:B:314:LEU:HD12	1:B:383:THR:HB	1.96	0.47
1:A:320:SER:O	1:A:323:ILE:HG22	2.14	0.47
1:B:61:HIS:HA	1:B:64:VAL:HG23	1.96	0.47
1:A:336:GLU:HB2	1:A:339:MET:SD	2.55	0.47
1:C:320:SER:O	1:C:323:ILE:HG22	2.14	0.47
1:B:169:PRO:HB3	1:B:338:ALA:HB3	1.97	0.47
1:C:169:PRO:HB3	1:C:338:ALA:HB3	1.97	0.47
1:A:169:PRO:HB3	1:A:338:ALA:HB3	1.97	0.46
1:A:247:TYR:HB3	1:A:267:LEU:HD13	1.97	0.46
1:A:21:GLU:HG3	1:A:156:THR:HG21	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:312:TRP:CE3	1:C:450:VAL:HG13	2.51	0.46
1:C:247:TYR:HB3	1:C:267:LEU:HD13	1.98	0.46
1:A:88:ILE:HG22	1:A:126:ILE:HG12	1.98	0.46
1:A:312:TRP:CE3	1:A:450:VAL:HG13	2.50	0.45
1:B:336:GLU:HB2	1:B:339:MET:SD	2.56	0.45
1:B:21:GLU:HG3	1:B:156:THR:HG21	1.98	0.45
1:C:336:GLU:HB2	1:C:339:MET:SD	2.56	0.45
1:A:99:HIS:CE1	1:A:184:PRO:HB2	2.52	0.45
1:B:312:TRP:CE3	1:B:450:VAL:HG13	2.51	0.45
1:C:21:GLU:HG3	1:C:156:THR:HG21	1.97	0.45
1:B:310:PRO:O	1:B:314:LEU:HG	2.17	0.45
1:C:310:PRO:O	1:C:314:LEU:HG	2.17	0.45
1:A:310:PRO:O	1:A:314:LEU:HG	2.17	0.45
1:B:247:TYR:HB3	1:B:267:LEU:HD13	1.97	0.45
1:C:39:LEU:HD13	1:C:56:ALA:HA	1.99	0.44
1:A:88:ILE:HA	1:A:126:ILE:HD11	1.98	0.44
1:B:371:THR:HB	1:B:374:ARG:HD2	1.99	0.44
1:B:39:LEU:HD13	1:B:56:ALA:HA	1.98	0.44
1:C:371:THR:HB	1:C:374:ARG:HD2	2.00	0.44
1:B:14:ILE:N	1:B:15:PRO:CD	2.79	0.44
1:B:99:HIS:CE1	1:B:184:PRO:HB2	2.52	0.44
1:B:126:ILE:HA	1:B:129:LEU:HD22	1.99	0.44
1:A:39:LEU:HD13	1:A:56:ALA:HA	1.99	0.44
1:C:26:PHE:HB2	1:C:187:LEU:HD13	2.00	0.44
1:C:88:ILE:HG22	1:C:126:ILE:HG12	1.99	0.44
1:A:371:THR:HB	1:A:374:ARG:HD2	1.99	0.43
1:C:99:HIS:CE1	1:C:184:PRO:HB2	2.52	0.43
1:B:88:ILE:HG22	1:B:126:ILE:HG12	2.00	0.43
1:A:86:ASN:HD22	1:A:86:ASN:H	1.66	0.43
1:B:435:PRO:HG2	1:B:442:ILE:HD11	2.01	0.43
1:B:88:ILE:HA	1:B:126:ILE:HD11	2.00	0.43
1:B:154:TYR:HA	1:B:157:ILE:HD12	2.01	0.43
1:C:88:ILE:HA	1:C:126:ILE:HD11	2.01	0.43
1:A:154:TYR:HA	1:A:157:ILE:HD12	2.01	0.43
1:C:154:TYR:HA	1:C:157:ILE:HD12	2.01	0.43
1:B:86:ASN:HD22	1:B:86:ASN:H	1.66	0.42
1:B:39:LEU:HD22	1:B:56:ALA:HA	2.01	0.42
1:A:26:PHE:HB2	1:A:187:LEU:HD13	2.01	0.42
1:C:86:ASN:HD22	1:C:86:ASN:H	1.66	0.42
1:C:305:PHE:CD1	1:C:446:TRP:HD1	2.37	0.42
1:B:305:PHE:CD1	1:B:446:TRP:HD1	2.37	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:318:LYS:O	1:B:322:TRP:HB2	2.20	0.42
1:A:39:LEU:HD22	1:A:56:ALA:HA	2.02	0.42
1:B:332:PRO:HD2	1:B:335:PHE:HB2	2.01	0.42
1:A:305:PHE:CD1	1:A:446:TRP:HD1	2.37	0.42
1:C:351:LEU:HD22	1:C:420:VAL:HG21	2.02	0.42
1:C:39:LEU:HD22	1:C:56:ALA:HA	2.02	0.42
1:A:332:PRO:HD2	1:A:335:PHE:HB2	2.01	0.41
1:C:301:ILE:HA	1:C:304:LEU:HD12	2.01	0.41
1:C:332:PRO:HD2	1:C:335:PHE:HB2	2.01	0.41
1:A:318:LYS:O	1:A:322:TRP:HB2	2.19	0.41
1:C:318:LYS:O	1:C:322:TRP:HB2	2.20	0.41
1:A:301:ILE:HA	1:A:304:LEU:HD12	2.01	0.41
1:A:388:ILE:HB	1:A:489:ALA:HB1	2.02	0.41
1:B:26:PHE:HB2	1:B:187:LEU:HD13	2.02	0.41
1:B:301:ILE:HA	1:B:304:LEU:HD12	2.01	0.41
1:B:244:SER:HA	1:B:247:TYR:CD2	2.56	0.41
1:A:312:TRP:CE2	1:A:450:VAL:HG22	2.56	0.41
1:B:388:ILE:HB	1:B:489:ALA:HB1	2.02	0.41
1:C:312:TRP:CE2	1:C:450:VAL:HG22	2.55	0.41
1:C:435:PRO:HG2	1:C:442:ILE:HD11	2.02	0.40
1:C:313:SER:O	1:C:317:GLN:NE2	2.50	0.40
1:A:435:PRO:HG2	1:A:442:ILE:HD11	2.03	0.40
1:C:35:LEU:O	1:C:39:LEU:HD12	2.22	0.40
1:A:244:SER:HA	1:A:247:TYR:CD2	2.56	0.40
1:A:25:ARG:HD3	1:A:25:ARG:HA	1.95	0.40
1:B:299:LEU:HA	1:B:302:LEU:HD12	2.04	0.40
1:B:312:TRP:CE2	1:B:450:VAL:HG22	2.56	0.40
1:A:299:LEU:HA	1:A:302:LEU:HD12	2.04	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	449/524 (86%)	388 (86%)	55 (12%)	6 (1%)	15	60
1	B	449/524 (86%)	388 (86%)	55 (12%)	6 (1%)	15	60
1	C	449/524 (86%)	389 (87%)	54 (12%)	6 (1%)	15	60
All	All	1347/1572 (86%)	1165 (86%)	164 (12%)	18 (1%)	15	60

All (18) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	14	ILE
1	A	47	ILE
1	B	14	ILE
1	B	47	ILE
1	C	14	ILE
1	C	47	ILE
1	A	436	LYS
1	B	436	LYS
1	C	436	LYS
1	A	401	SER
1	A	403	LEU
1	B	401	SER
1	B	403	LEU
1	C	401	SER
1	C	403	LEU
1	A	440	GLY
1	C	440	GLY
1	B	440	GLY

### 5.3.2 Protein sidechains ⓘ

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	367/427 (86%)	306 (83%)	61 (17%)	3	19
1	B	367/427 (86%)	308 (84%)	59 (16%)	3	21
1	C	367/427 (86%)	307 (84%)	60 (16%)	3	20

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	1101/1281 (86%)	921 (84%)	180 (16%)	3 20

All (180) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	35	LEU
1	A	36	THR
1	A	39	LEU
1	A	45	LEU
1	A	47	ILE
1	A	50	GLU
1	A	52	ARG
1	A	58	ASP
1	A	64	VAL
1	A	65	ILE
1	A	73	LEU
1	A	81	PHE
1	A	85	TYR
1	A	86	ASN
1	A	87	THR
1	A	90	TRP
1	A	92	SER
1	A	93	LEU
1	A	95	TYR
1	A	97	VAL
1	A	101	PHE
1	A	121	LEU
1	A	126	ILE
1	A	130	VAL
1	A	166	LEU
1	A	167	SER
1	A	172	LEU
1	A	179	VAL
1	A	183	ILE
1	A	197	LEU
1	A	235	LEU
1	A	237	LEU
1	A	257	ILE
1	A	265	MET
1	A	284	ARG
1	A	289	ASP
1	A	293	ASP

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Mol	Chain	Res	Type
1	A	299	LEU
1	A	309	THR
1	A	324	LEU
1	A	328	ASP
1	A	331	LYS
1	A	347	LEU
1	A	351	LEU
1	A	380	ILE
1	A	390	VAL
1	A	394	GLN
1	A	396	MET
1	A	407	TRP
1	A	412	TYR
1	A	414	LEU
1	A	415	LEU
1	A	416	THR
1	A	447	THR
1	A	448	LEU
1	A	451	THR
1	A	455	LEU
1	A	458	LEU
1	A	481	THR
1	A	491	PHE
1	A	499	PHE
1	B	35	LEU
1	B	36	THR
1	B	39	LEU
1	B	45	LEU
1	B	47	ILE
1	B	50	GLU
1	B	52	ARG
1	B	58	ASP
1	B	64	VAL
1	B	65	ILE
1	B	73	LEU
1	B	81	PHE
1	B	85	TYR
1	B	86	ASN
1	B	87	THR
1	B	90	TRP
1	B	92	SER
1	B	93	LEU

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Mol	Chain	Res	Type
1	B	95	TYR
1	B	97	VAL
1	B	101	PHE
1	B	121	LEU
1	B	126	ILE
1	B	130	VAL
1	B	166	LEU
1	B	167	SER
1	B	172	LEU
1	B	179	VAL
1	B	183	ILE
1	B	235	LEU
1	B	237	LEU
1	B	257	ILE
1	B	265	MET
1	B	284	ARG
1	B	289	ASP
1	B	293	ASP
1	B	309	THR
1	B	324	LEU
1	B	328	ASP
1	B	331	LYS
1	B	347	LEU
1	B	351	LEU
1	B	380	ILE
1	B	390	VAL
1	B	394	GLN
1	B	396	MET
1	B	407	TRP
1	B	412	TYR
1	B	414	LEU
1	B	415	LEU
1	B	416	THR
1	B	447	THR
1	B	448	LEU
1	B	451	THR
1	B	455	LEU
1	B	458	LEU
1	B	481	THR
1	B	491	PHE
1	B	499	PHE
1	C	35	LEU

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Mol	Chain	Res	Type
1	C	36	THR
1	C	39	LEU
1	C	45	LEU
1	C	47	ILE
1	C	50	GLU
1	C	52	ARG
1	C	58	ASP
1	C	64	VAL
1	C	65	ILE
1	C	73	LEU
1	C	81	PHE
1	C	85	TYR
1	C	86	ASN
1	C	87	THR
1	C	90	TRP
1	C	92	SER
1	C	93	LEU
1	C	95	TYR
1	C	97	VAL
1	C	101	PHE
1	C	121	LEU
1	C	126	ILE
1	C	130	VAL
1	C	166	LEU
1	C	167	SER
1	C	172	LEU
1	C	179	VAL
1	C	183	ILE
1	C	197	LEU
1	C	235	LEU
1	C	237	LEU
1	C	257	ILE
1	C	265	MET
1	C	284	ARG
1	C	289	ASP
1	C	293	ASP
1	C	309	THR
1	C	324	LEU
1	C	328	ASP
1	C	331	LYS
1	C	347	LEU
1	C	351	LEU

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Mol	Chain	Res	Type
1	C	380	ILE
1	C	390	VAL
1	C	394	GLN
1	C	396	MET
1	C	407	TRP
1	C	412	TYR
1	C	414	LEU
1	C	415	LEU
1	C	416	THR
1	C	447	THR
1	C	448	LEU
1	C	451	THR
1	C	455	LEU
1	C	458	LEU
1	C	481	THR
1	C	491	PHE
1	C	499	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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 ⓘ

There are no ligands in this entry.

### 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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	456/524 (87%)	-0.29	3 (0%) 89 81	77, 270, 300, 300	0
1	B	456/524 (87%)	-0.27	9 (1%) 68 54	78, 280, 300, 300	0
1	C	456/524 (87%)	-0.22	13 (2%) 55 40	87, 290, 300, 300	0
All	All	1368/1572 (87%)	-0.26	25 (1%) 71 57	77, 282, 300, 300	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	139	ASP	6.6
1	C	381	ALA	5.8
1	C	149	ALA	4.9
1	B	497	ILE	3.7
1	C	254	THR	3.7
1	C	283	ALA	3.5
1	C	230	LYS	3.4
1	C	417	PHE	3.2
1	C	290	ALA	3.2
1	B	373	LEU	3.1
1	C	234	GLY	3.0
1	B	417	PHE	3.0
1	A	431	TYR	2.8
1	C	134	MET	2.7
1	B	273	GLY	2.6
1	B	441	THR	2.5
1	C	463	SER	2.5
1	B	498	VAL	2.4
1	C	371	THR	2.4
1	B	274	ALA	2.3
1	B	280	LEU	2.3
1	A	403	LEU	2.2
1	B	231	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	285	LYS	2.1
1	A	307	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.