



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 17, 2017 – 09:27 AM EDT

PDB ID : 5MLF  
Title : Structure of Psb29 at 1.55Å  
Authors : Murray, J.W.; Kozlo, A.  
Deposited on : unknown  
Resolution : 3.64 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
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20029824
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
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)	:	rb-20029824

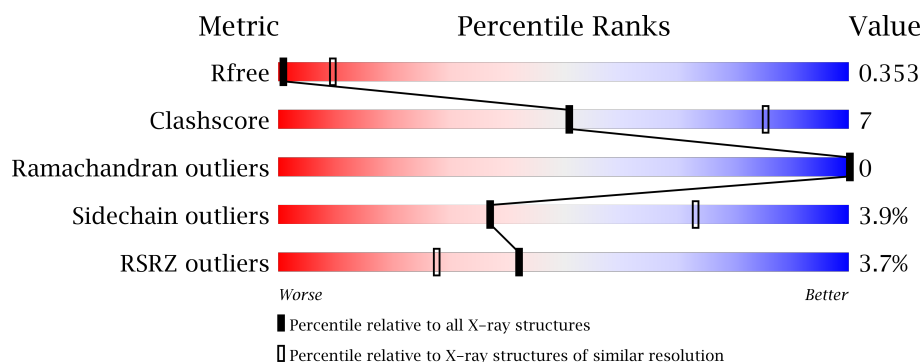
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.64 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))
$R_{free}$	100719	1143 (3.80-3.48)
Clashscore	112137	1092 (3.78-3.50)
Ramachandran outliers	110173	1051 (3.78-3.50)
Sidechain outliers	110143	1051 (3.78-3.50)
RSRZ outliers	101464	1000 (3.78-3.50)

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	239	<div> <div>3%</div> <div>69%</div> <div>15%</div> <div>15%</div> </div>
1	B	239	<div> <div>5%</div> <div>69%</div> <div>14%</div> <div>15%</div> </div>
1	C	239	<div> <div>2%</div> <div>69%</div> <div>14%</div> <div>15%</div> </div>
1	D	239	<div> <div>0%</div> <div>70%</div> <div>13%</div> <div>15%</div> </div>
1	E	239	<div> <div>4%</div> <div>69%</div> <div>15%</div> <div>15%</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	239	
1	G	239	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	HG	D	301	-	-	-	X

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 11277 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 Protein Thf1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	203	Total	C	N	O	S	0	2	0
			1610	1021	282	303	4			
1	B	203	Total	C	N	O	S	0	2	0
			1610	1021	282	303	4			
1	C	203	Total	C	N	O	S	0	2	0
			1610	1021	282	303	4			
1	D	203	Total	C	N	O	S	0	2	0
			1610	1021	282	303	4			
1	E	203	Total	C	N	O	S	0	2	0
			1610	1021	282	303	4			
1	G	203	Total	C	N	O	S	0	2	0
			1610	1021	282	303	4			
1	F	203	Total	C	N	O	S	0	2	0
			1610	1021	282	303	4			

There are 119 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-16	MET	-	initiating methionine	UNP Q8DJT8
A	-15	ARG	-	expression tag	UNP Q8DJT8
A	-14	GLY	-	expression tag	UNP Q8DJT8
A	-13	SER	-	expression tag	UNP Q8DJT8
A	-12	HIS	-	expression tag	UNP Q8DJT8
A	-11	HIS	-	expression tag	UNP Q8DJT8
A	-10	HIS	-	expression tag	UNP Q8DJT8
A	-9	HIS	-	expression tag	UNP Q8DJT8
A	-8	HIS	-	expression tag	UNP Q8DJT8
A	-7	HIS	-	expression tag	UNP Q8DJT8
A	-6	GLY	-	expression tag	UNP Q8DJT8
A	-5	LEU	-	expression tag	UNP Q8DJT8
A	-4	VAL	-	expression tag	UNP Q8DJT8
A	-3	PRO	-	expression tag	UNP Q8DJT8
A	-2	ARG	-	expression tag	UNP Q8DJT8

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP Q8DJT8
A	0	SER	-	expression tag	UNP Q8DJT8
B	-16	MET	-	initiating methionine	UNP Q8DJT8
B	-15	ARG	-	expression tag	UNP Q8DJT8
B	-14	GLY	-	expression tag	UNP Q8DJT8
B	-13	SER	-	expression tag	UNP Q8DJT8
B	-12	HIS	-	expression tag	UNP Q8DJT8
B	-11	HIS	-	expression tag	UNP Q8DJT8
B	-10	HIS	-	expression tag	UNP Q8DJT8
B	-9	HIS	-	expression tag	UNP Q8DJT8
B	-8	HIS	-	expression tag	UNP Q8DJT8
B	-7	HIS	-	expression tag	UNP Q8DJT8
B	-6	GLY	-	expression tag	UNP Q8DJT8
B	-5	LEU	-	expression tag	UNP Q8DJT8
B	-4	VAL	-	expression tag	UNP Q8DJT8
B	-3	PRO	-	expression tag	UNP Q8DJT8
B	-2	ARG	-	expression tag	UNP Q8DJT8
B	-1	GLY	-	expression tag	UNP Q8DJT8
B	0	SER	-	expression tag	UNP Q8DJT8
C	-16	MET	-	initiating methionine	UNP Q8DJT8
C	-15	ARG	-	expression tag	UNP Q8DJT8
C	-14	GLY	-	expression tag	UNP Q8DJT8
C	-13	SER	-	expression tag	UNP Q8DJT8
C	-12	HIS	-	expression tag	UNP Q8DJT8
C	-11	HIS	-	expression tag	UNP Q8DJT8
C	-10	HIS	-	expression tag	UNP Q8DJT8
C	-9	HIS	-	expression tag	UNP Q8DJT8
C	-8	HIS	-	expression tag	UNP Q8DJT8
C	-7	HIS	-	expression tag	UNP Q8DJT8
C	-6	GLY	-	expression tag	UNP Q8DJT8
C	-5	LEU	-	expression tag	UNP Q8DJT8
C	-4	VAL	-	expression tag	UNP Q8DJT8
C	-3	PRO	-	expression tag	UNP Q8DJT8
C	-2	ARG	-	expression tag	UNP Q8DJT8
C	-1	GLY	-	expression tag	UNP Q8DJT8
C	0	SER	-	expression tag	UNP Q8DJT8
D	-16	MET	-	initiating methionine	UNP Q8DJT8
D	-15	ARG	-	expression tag	UNP Q8DJT8
D	-14	GLY	-	expression tag	UNP Q8DJT8
D	-13	SER	-	expression tag	UNP Q8DJT8
D	-12	HIS	-	expression tag	UNP Q8DJT8
D	-11	HIS	-	expression tag	UNP Q8DJT8

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-10	HIS	-	expression tag	UNP Q8DJT8
D	-9	HIS	-	expression tag	UNP Q8DJT8
D	-8	HIS	-	expression tag	UNP Q8DJT8
D	-7	HIS	-	expression tag	UNP Q8DJT8
D	-6	GLY	-	expression tag	UNP Q8DJT8
D	-5	LEU	-	expression tag	UNP Q8DJT8
D	-4	VAL	-	expression tag	UNP Q8DJT8
D	-3	PRO	-	expression tag	UNP Q8DJT8
D	-2	ARG	-	expression tag	UNP Q8DJT8
D	-1	GLY	-	expression tag	UNP Q8DJT8
D	0	SER	-	expression tag	UNP Q8DJT8
E	-16	MET	-	initiating methionine	UNP Q8DJT8
E	-15	ARG	-	expression tag	UNP Q8DJT8
E	-14	GLY	-	expression tag	UNP Q8DJT8
E	-13	SER	-	expression tag	UNP Q8DJT8
E	-12	HIS	-	expression tag	UNP Q8DJT8
E	-11	HIS	-	expression tag	UNP Q8DJT8
E	-10	HIS	-	expression tag	UNP Q8DJT8
E	-9	HIS	-	expression tag	UNP Q8DJT8
E	-8	HIS	-	expression tag	UNP Q8DJT8
E	-7	HIS	-	expression tag	UNP Q8DJT8
E	-6	GLY	-	expression tag	UNP Q8DJT8
E	-5	LEU	-	expression tag	UNP Q8DJT8
E	-4	VAL	-	expression tag	UNP Q8DJT8
E	-3	PRO	-	expression tag	UNP Q8DJT8
E	-2	ARG	-	expression tag	UNP Q8DJT8
E	-1	GLY	-	expression tag	UNP Q8DJT8
E	0	SER	-	expression tag	UNP Q8DJT8
G	-16	MET	-	initiating methionine	UNP Q8DJT8
G	-15	ARG	-	expression tag	UNP Q8DJT8
G	-14	GLY	-	expression tag	UNP Q8DJT8
G	-13	SER	-	expression tag	UNP Q8DJT8
G	-12	HIS	-	expression tag	UNP Q8DJT8
G	-11	HIS	-	expression tag	UNP Q8DJT8
G	-10	HIS	-	expression tag	UNP Q8DJT8
G	-9	HIS	-	expression tag	UNP Q8DJT8
G	-8	HIS	-	expression tag	UNP Q8DJT8
G	-7	HIS	-	expression tag	UNP Q8DJT8
G	-6	GLY	-	expression tag	UNP Q8DJT8
G	-5	LEU	-	expression tag	UNP Q8DJT8
G	-4	VAL	-	expression tag	UNP Q8DJT8
G	-3	PRO	-	expression tag	UNP Q8DJT8

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Chain	Residue	Modelled	Actual	Comment	Reference
G	-2	ARG	-	expression tag	UNP Q8DJT8
G	-1	GLY	-	expression tag	UNP Q8DJT8
G	0	SER	-	expression tag	UNP Q8DJT8
F	-16	MET	-	initiating methionine	UNP Q8DJT8
F	-15	ARG	-	expression tag	UNP Q8DJT8
F	-14	GLY	-	expression tag	UNP Q8DJT8
F	-13	SER	-	expression tag	UNP Q8DJT8
F	-12	HIS	-	expression tag	UNP Q8DJT8
F	-11	HIS	-	expression tag	UNP Q8DJT8
F	-10	HIS	-	expression tag	UNP Q8DJT8
F	-9	HIS	-	expression tag	UNP Q8DJT8
F	-8	HIS	-	expression tag	UNP Q8DJT8
F	-7	HIS	-	expression tag	UNP Q8DJT8
F	-6	GLY	-	expression tag	UNP Q8DJT8
F	-5	LEU	-	expression tag	UNP Q8DJT8
F	-4	VAL	-	expression tag	UNP Q8DJT8
F	-3	PRO	-	expression tag	UNP Q8DJT8
F	-2	ARG	-	expression tag	UNP Q8DJT8
F	-1	GLY	-	expression tag	UNP Q8DJT8
F	0	SER	-	expression tag	UNP Q8DJT8

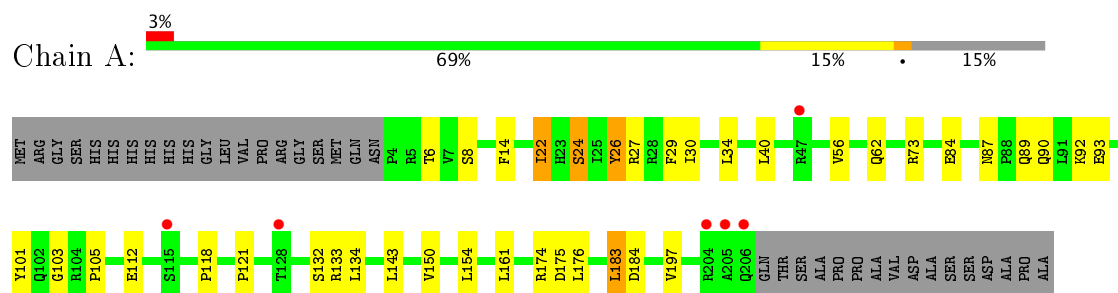
- Molecule 2 is MERCURY (II) ION (three-letter code: HG) (formula: Hg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	G	1	Total Hg 1 1	0	0
2	D	1	Total Hg 1 1	0	0
2	E	1	Total Hg 1 1	0	0
2	B	1	Total Hg 1 1	0	0
2	C	1	Total Hg 1 1	0	0
2	A	1	Total Hg 1 1	0	0
2	F	1	Total Hg 1 1	0	0

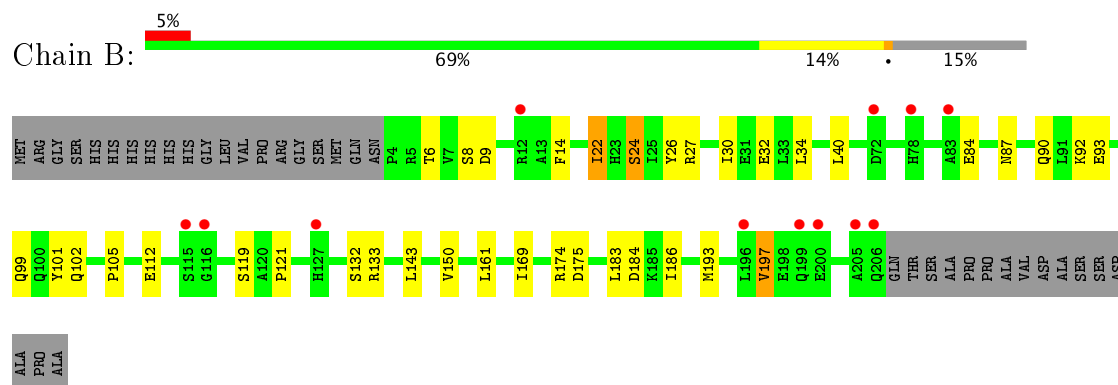
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

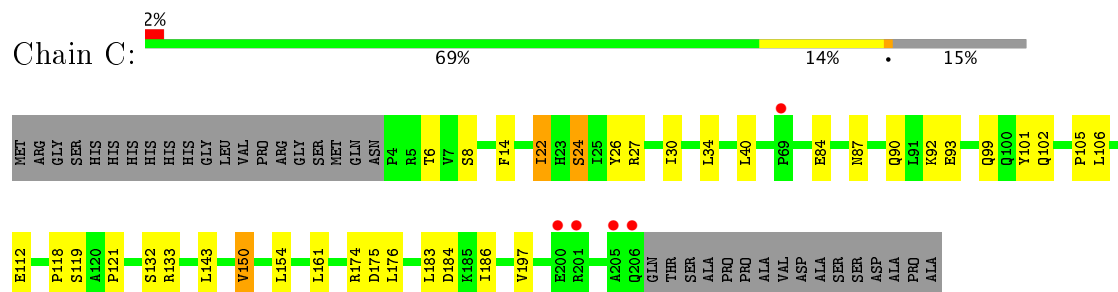
#### • Molecule 1: Protein Thf1



#### • Molecule 1: Protein Thf1

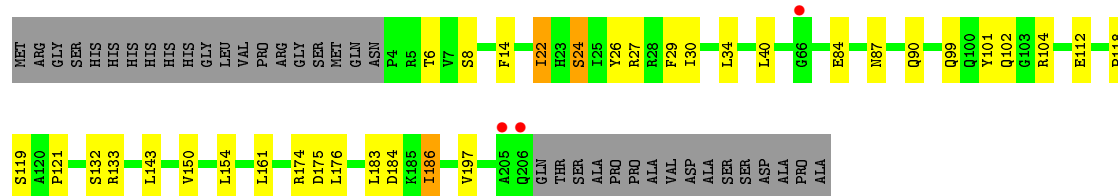


#### • Molecule 1: Protein Thf1

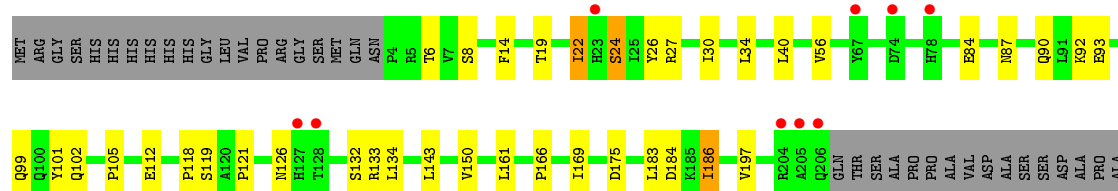


#### • Molecule 1: Protein Thf1

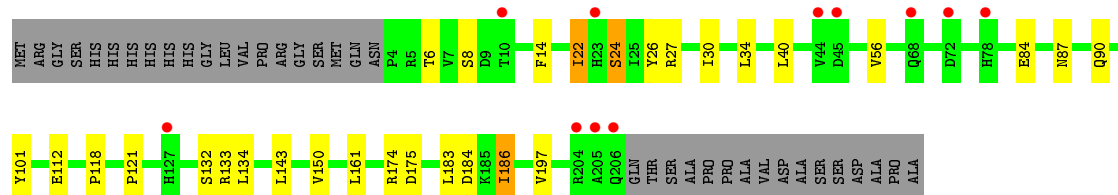
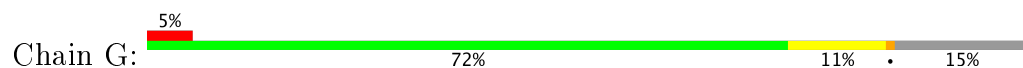




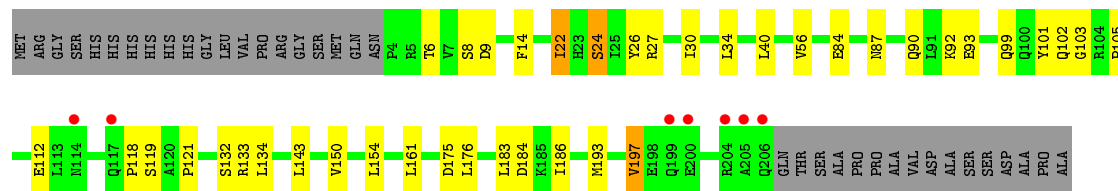
• Molecule 1: Protein Thf1



• Molecule 1: Protein Thf1



• Molecule 1: Protein Thf1



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 63 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	138.91Å 138.91Å 305.91Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	55.98 – 3.64 55.98 – 3.64	Depositor EDS
% Data completeness (in resolution range)	99.8 (55.98-3.64) 99.9 (55.98-3.64)	Depositor EDS
$R_{merge}$	0.39	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.25 (at 3.67Å)	Xtriage
Refinement program	PHENIX (1.11.1 _2575: ???)	Depositor
R, $R_{free}$	0.311 , 0.357 0.308 , 0.353	Depositor DCC
$R_{free}$ test set	1046 reflections (5.13%)	DCC
Wilson B-factor (Å <sup>2</sup> )	53.2	Xtriage
Anisotropy	0.432	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 75.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.41$ , $\langle L^2 \rangle = 0.23$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.79	EDS
Total number of atoms	11277	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	92.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.26% 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, 0.333 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: HG

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	0.24	0/1648	0.37	0/2240
1	B	0.23	0/1648	0.37	0/2240
1	C	0.23	0/1648	0.37	0/2240
1	D	0.23	0/1648	0.37	0/2240
1	E	0.23	0/1648	0.37	0/2240
1	F	0.23	0/1648	0.37	0/2240
1	G	0.23	0/1648	0.37	0/2240
All	All	0.23	0/11536	0.37	0/15680

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	A	0	1
1	B	0	1
1	C	0	1
1	D	0	1
1	E	0	1
1	F	0	1
1	G	0	1
All	All	0	7

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	24	SER	Peptide
1	B	24	SER	Peptide
1	C	24	SER	Peptide
1	D	24	SER	Peptide
1	E	24	SER	Peptide
1	F	24	SER	Peptide
1	G	24	SER	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	1610	0	1552	28	0
1	B	1610	0	1552	26	0
1	C	1610	0	1552	28	0
1	D	1610	0	1552	25	0
1	E	1610	0	1552	22	0
1	F	1610	0	1552	24	0
1	G	1610	0	1552	19	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
2	G	1	0	0	0	0
All	All	11277	0	10864	145	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 (145) 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:174:ARG:HH22	1:C:105:PRO:HD3	1.53	0.72
1:A:24:SER:HB3	1:A:27:ARG:HE	1.57	0.69
1:C:6:THR:HG22	1:C:8:SER:H	1.59	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:6:THR:HG22	1:G:8:SER:H	1.60	0.67
1:D:87:ASN:HB3	1:D:90:GLN:HB3	1.77	0.67
1:E:87:ASN:HB3	1:E:90:GLN:HB3	1.77	0.67
1:G:87:ASN:HB3	1:G:90:GLN:HB3	1.77	0.66
1:A:6:THR:HG22	1:A:8:SER:H	1.61	0.65
1:B:87:ASN:HB3	1:B:90:GLN:HB3	1.77	0.65
1:D:40:LEU:HD21	1:D:132:SER:HB2	1.78	0.65
1:A:87:ASN:HB3	1:A:90:GLN:HB3	1.78	0.65
1:A:40:LEU:HD21	1:A:132:SER:HB2	1.80	0.64
1:E:40:LEU:HD21	1:E:132:SER:HB2	1.79	0.64
1:A:174:ARG:HH22	1:E:105:PRO:HD3	1.62	0.64
1:C:87:ASN:HB3	1:C:90:GLN:HB3	1.78	0.64
1:F:87:ASN:HB3	1:F:90:GLN:HB3	1.79	0.64
1:C:112:GLU:OE2	1:C:119:SER:OG	2.13	0.63
1:A:22:ILE:HD12	1:A:27:ARG:HH12	1.63	0.63
1:G:40:LEU:HD21	1:G:132:SER:HB2	1.79	0.63
1:B:105:PRO:HD3	1:G:174:ARG:HH22	1.63	0.63
1:D:22:ILE:HD12	1:D:27:ARG:HH12	1.63	0.63
1:A:6:THR:HG23	1:E:93:GLU:HG2	1.80	0.63
1:D:6:THR:HG22	1:D:8:SER:H	1.63	0.62
1:F:22:ILE:HD12	1:F:27:ARG:HH12	1.64	0.62
1:E:6:THR:HG22	1:E:8:SER:H	1.65	0.61
1:C:22:ILE:HD12	1:C:27:ARG:HH12	1.65	0.61
1:G:24:SER:HB3	1:G:27:ARG:HE	1.66	0.61
1:B:24:SER:HB3	1:B:27:ARG:HE	1.66	0.61
1:E:22:ILE:HD12	1:E:27:ARG:HH12	1.66	0.61
1:B:40:LEU:HD21	1:B:132:SER:HB2	1.82	0.61
1:E:24:SER:HB3	1:E:27:ARG:HE	1.65	0.61
1:G:22:ILE:HD12	1:G:27:ARG:HH12	1.65	0.61
1:A:103:GLY:O	1:D:174:ARG:NH2	2.34	0.60
1:B:22:ILE:HD12	1:B:27:ARG:HH12	1.66	0.59
1:D:24:SER:HB3	1:D:27:ARG:HE	1.67	0.59
1:F:24:SER:HB3	1:F:27:ARG:HE	1.66	0.59
1:C:6:THR:HG23	1:F:93:GLU:HG2	1.85	0.58
1:B:6:THR:HG23	1:C:93:GLU:HG2	1.84	0.58
1:D:99:GLN:HG3	1:D:102:GLN:HG3	1.84	0.58
1:F:112:GLU:OE2	1:F:119:SER:OG	2.18	0.58
1:F:40:LEU:HD21	1:F:132:SER:HB2	1.84	0.58
1:B:93:GLU:HG2	1:G:6:THR:HG23	1.84	0.57
1:C:24:SER:HB3	1:C:27:ARG:HE	1.68	0.57
1:A:133:ARG:HG3	1:A:175:ASP:HB3	1.87	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:133:ARG:HG3	1:B:175:ASP:HB3	1.88	0.56
1:B:14:PHE:HE2	1:B:22:ILE:HD13	1.69	0.56
1:G:133:ARG:HG3	1:G:175:ASP:HB3	1.88	0.56
1:E:14:PHE:HE2	1:E:22:ILE:HD13	1.71	0.55
1:B:6:THR:HG22	1:B:8:SER:H	1.70	0.55
1:C:174:ARG:HH22	1:F:105:PRO:HD3	1.72	0.55
1:C:40:LEU:HD21	1:C:132:SER:HB2	1.88	0.55
1:C:14:PHE:HE2	1:C:22:ILE:HD13	1.72	0.55
1:D:14:PHE:HE2	1:D:22:ILE:HD13	1.72	0.55
1:D:112:GLU:OE2	1:D:119:SER:OG	2.14	0.54
1:A:14:PHE:HE2	1:A:22:ILE:HD13	1.72	0.54
1:G:14:PHE:HE2	1:G:22:ILE:HD13	1.73	0.54
1:D:133:ARG:HG3	1:D:175:ASP:HB3	1.90	0.53
1:E:133:ARG:HG3	1:E:175:ASP:HB3	1.90	0.53
1:F:14:PHE:HE2	1:F:22:ILE:HD13	1.74	0.53
1:C:133:ARG:HG3	1:C:175:ASP:HB3	1.91	0.53
1:E:112:GLU:OE2	1:E:119:SER:OG	2.20	0.53
1:C:99:GLN:HG3	1:C:102:GLN:HG3	1.91	0.52
1:E:99:GLN:HG3	1:E:102:GLN:HG3	1.91	0.52
1:F:133:ARG:HG3	1:F:175:ASP:HB3	1.93	0.51
1:C:112:GLU:HG2	1:C:118:PRO:HA	1.92	0.51
1:B:99:GLN:HG3	1:B:102:GLN:HG3	1.93	0.50
1:D:112:GLU:HG2	1:D:118:PRO:HA	1.93	0.50
1:A:6:THR:HG22	1:A:8:SER:N	2.26	0.50
1:A:174:ARG:NH2	1:E:105:PRO:HD3	2.27	0.49
1:B:6:THR:HG22	1:B:8:SER:N	2.27	0.49
1:D:143:LEU:HD21	1:D:161:LEU:HD11	1.94	0.49
1:G:101:TYR:HE2	1:G:121:PRO:HD2	1.78	0.48
1:F:6:THR:HB	1:F:9:ASP:OD2	2.12	0.48
1:F:34:LEU:HD23	1:F:84:GLU:OE2	2.13	0.48
1:E:34:LEU:HD23	1:E:84:GLU:OE2	2.13	0.48
1:B:105:PRO:HD3	1:G:174:ARG:NH2	2.29	0.48
1:F:99:GLN:HG3	1:F:102:GLN:HG3	1.96	0.48
1:B:34:LEU:HD23	1:B:84:GLU:OE2	2.13	0.47
1:E:112:GLU:HG2	1:E:118:PRO:HA	1.95	0.47
1:A:34:LEU:HD23	1:A:84:GLU:OE2	2.15	0.47
1:C:174:ARG:NH2	1:F:103:GLY:O	2.49	0.46
1:B:174:ARG:NH2	1:C:105:PRO:HD3	2.26	0.46
1:G:143:LEU:HD21	1:G:161:LEU:HD11	1.97	0.46
1:B:26:TYR:O	1:B:30:ILE:HG13	2.16	0.46
1:G:26:TYR:O	1:G:30:ILE:HG13	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:105:PRO:HD3	1:D:174:ARG:HH22	1.81	0.46
1:C:101:TYR:HE2	1:C:121:PRO:HD2	1.80	0.46
1:E:26:TYR:O	1:E:30:ILE:HG13	2.16	0.46
1:D:101:TYR:HE2	1:D:121:PRO:HD2	1.81	0.45
1:D:34:LEU:HD23	1:D:84:GLU:OE2	2.16	0.45
1:A:101:TYR:HE2	1:A:121:PRO:HD2	1.82	0.45
1:C:26:TYR:O	1:C:30:ILE:HG13	2.15	0.45
1:D:26:TYR:O	1:D:30:ILE:HG13	2.16	0.45
1:A:143:LEU:HD21	1:A:161:LEU:HD11	1.97	0.45
1:E:101:TYR:HE2	1:E:121:PRO:HD2	1.82	0.45
1:F:101:TYR:HE2	1:F:121:PRO:HD2	1.81	0.45
1:B:112:GLU:OE2	1:B:119:SER:OG	2.19	0.45
1:D:186:ILE:H	1:D:186:ILE:HG13	1.61	0.45
1:F:143:LEU:HD21	1:F:161:LEU:HD11	1.97	0.45
1:E:56:VAL:HG22	1:E:134:LEU:HD13	1.97	0.45
1:E:126:ASN:HA	1:E:166:PRO:HG3	1.98	0.45
1:B:101:TYR:HE2	1:B:121:PRO:HD2	1.81	0.45
1:C:6:THR:HG22	1:C:8:SER:N	2.31	0.45
1:G:34:LEU:HD23	1:G:84:GLU:OE2	2.17	0.44
1:F:26:TYR:O	1:F:30:ILE:HG13	2.16	0.44
1:A:56:VAL:HG22	1:A:134:LEU:HD13	1.99	0.44
1:A:26:TYR:O	1:A:30:ILE:HG13	2.17	0.44
1:A:154:LEU:HD21	1:A:176:LEU:HD13	1.99	0.44
1:B:6:THR:HB	1:B:9:ASP:OD2	2.18	0.44
1:C:154:LEU:HD21	1:C:176:LEU:HD13	1.99	0.44
1:D:154:LEU:HD21	1:D:176:LEU:HD13	1.99	0.44
1:G:6:THR:HG22	1:G:8:SER:N	2.30	0.44
1:F:193:MET:O	1:F:197:VAL:N	2.45	0.43
1:G:112:GLU:HG2	1:G:118:PRO:HA	2.00	0.43
1:B:92:LYS:HE2	1:G:8:SER:OG	2.19	0.43
1:A:8:SER:OG	1:E:92:LYS:HE2	2.18	0.43
1:C:8:SER:OG	1:F:92:LYS:HE2	2.19	0.43
1:A:105:PRO:HD3	1:D:174:ARG:NH2	2.33	0.43
1:C:34:LEU:HD23	1:C:84:GLU:OE2	2.18	0.43
1:C:174:ARG:NH2	1:F:105:PRO:HD3	2.33	0.43
1:E:143:LEU:HD21	1:E:161:LEU:HD11	2.01	0.42
1:G:56:VAL:HG22	1:G:134:LEU:HD13	2.00	0.42
1:F:56:VAL:HG22	1:F:134:LEU:HD13	2.00	0.42
1:F:154:LEU:HD21	1:F:176:LEU:HD13	2.02	0.42
1:F:6:THR:HG22	1:F:8:SER:N	2.34	0.42
1:D:101:TYR:HA	1:D:104:ARG:HG3	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:105:PRO:HD3	1:D:174:ARG:HH12	1.85	0.42
1:B:193:MET:O	1:B:197:VAL:N	2.48	0.42
1:A:112:GLU:HG2	1:A:118:PRO:HA	2.01	0.42
1:A:93:GLU:HG2	1:D:6:THR:HG23	2.02	0.42
1:B:169:ILE:HD13	1:B:169:ILE:HA	1.94	0.42
1:C:143:LEU:HD21	1:C:161:LEU:HD11	2.01	0.42
1:A:92:LYS:HE2	1:D:8:SER:OG	2.20	0.42
1:B:143:LEU:HD21	1:B:161:LEU:HD11	2.01	0.41
1:A:62:GLN:HB3	1:A:183:LEU:HD11	2.03	0.41
1:A:89:GLN:HB3	1:D:6:THR:HG21	2.02	0.41
1:G:186:ILE:H	1:G:186:ILE:HG13	1.62	0.41
1:A:105:PRO:HD3	1:D:174:ARG:NH1	2.36	0.41
1:F:112:GLU:HG2	1:F:118:PRO:HA	2.03	0.40
1:E:169:ILE:HA	1:E:169:ILE:HD13	1.95	0.40
1:E:186:ILE:H	1:E:186:ILE:HG13	1.60	0.40
1:C:174:ARG:HH12	1:F:105:PRO:HD3	1.86	0.40
1:B:32:GLU:HA	1:C:99:GLN:HE22	1.87	0.40
1:C:106:LEU:HD22	1:C:150:VAL:HG11	2.04	0.40
1:B:8:SER:OG	1:C:92:LYS:HE2	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	203/239 (85%)	198 (98%)	5 (2%)	0	100	100
1	B	203/239 (85%)	196 (97%)	7 (3%)	0	100	100
1	C	203/239 (85%)	196 (97%)	7 (3%)	0	100	100
1	D	203/239 (85%)	197 (97%)	6 (3%)	0	100	100
1	E	203/239 (85%)	197 (97%)	6 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	F	203/239 (85%)	197 (97%)	6 (3%)	0	100	100
1	G	203/239 (85%)	196 (97%)	7 (3%)	0	100	100
All	All	1421/1673 (85%)	1377 (97%)	44 (3%)	0	100	100

There are no Ramachandran outliers to report.

### 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	169/207 (82%)	161 (95%)	8 (5%)	30	68
1	B	169/207 (82%)	163 (96%)	6 (4%)	40	75
1	C	169/207 (82%)	163 (96%)	6 (4%)	40	75
1	D	169/207 (82%)	162 (96%)	7 (4%)	35	72
1	E	169/207 (82%)	162 (96%)	7 (4%)	35	72
1	F	169/207 (82%)	163 (96%)	6 (4%)	40	75
1	G	169/207 (82%)	163 (96%)	6 (4%)	40	75
All	All	1183/1449 (82%)	1137 (96%)	46 (4%)	37	73

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	ILE
1	A	26	TYR
1	A	29	PHE
1	A	73	ARG
1	A	150	VAL
1	A	183	LEU
1	A	184	ASP
1	A	197	VAL
1	B	22	ILE
1	B	150	VAL

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Mol	Chain	Res	Type
1	B	183	LEU
1	B	184	ASP
1	B	186	ILE
1	B	197	VAL
1	C	22	ILE
1	C	150	VAL
1	C	183	LEU
1	C	184	ASP
1	C	186	ILE
1	C	197	VAL
1	D	22	ILE
1	D	29	PHE
1	D	150	VAL
1	D	183	LEU
1	D	184	ASP
1	D	186	ILE
1	D	197	VAL
1	E	19	THR
1	E	22	ILE
1	E	150	VAL
1	E	183	LEU
1	E	184	ASP
1	E	186	ILE
1	E	197	VAL
1	G	22	ILE
1	G	150	VAL
1	G	183	LEU
1	G	184	ASP
1	G	186	ILE
1	G	197	VAL
1	F	22	ILE
1	F	150	VAL
1	F	183	LEU
1	F	184	ASP
1	F	186	ILE
1	F	197	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	155	ASN
1	G	155	ASN

### 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](#)

Of 7 ligands modelled in this entry, 7 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 [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

### 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	203/239 (84%)	0.27	6 (2%)	51	34	20, 87, 160, 252	0
1	B	203/239 (84%)	0.27	12 (5%)	23	14	35, 91, 191, 244	0
1	C	203/239 (84%)	0.02	5 (2%)	58	40	20, 69, 163, 209	0
1	D	203/239 (84%)	0.08	3 (1%)	74	58	24, 77, 157, 196	0
1	E	203/239 (84%)	0.22	9 (4%)	35	22	32, 97, 175, 232	0
1	F	203/239 (84%)	0.11	7 (3%)	46	30	26, 76, 145, 203	0
1	G	203/239 (84%)	0.44	11 (5%)	26	17	39, 104, 189, 240	0
All	All	1421/1673 (84%)	0.20	53 (3%)	42	28	20, 87, 173, 252	0

All (53) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	205	ALA	5.7
1	A	204	ARG	4.2
1	E	206	GLN	4.2
1	D	206	GLN	4.1
1	D	205	ALA	3.9
1	G	206	GLN	3.6
1	C	200	GLU	3.5
1	G	204	ARG	3.4
1	B	205	ALA	3.4
1	B	115	SER	3.3
1	G	205	ALA	3.2
1	B	206	GLN	3.1
1	B	127	HIS	3.1
1	A	206	GLN	3.1
1	F	205	ALA	3.1
1	E	127	HIS	3.1
1	B	78	HIS	3.0

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Mol	Chain	Res	Type	RSRZ
1	F	204	ARG	2.9
1	B	199	GLN	2.8
1	G	78	HIS	2.8
1	B	72	ASP	2.7
1	F	199	GLN	2.7
1	E	67	TYR	2.7
1	G	68	GLN	2.6
1	F	114	ASN	2.5
1	A	47	ARG	2.5
1	E	205	ALA	2.5
1	G	44	VAL	2.5
1	F	200	GLU	2.5
1	B	83	ALA	2.5
1	A	128	THR	2.4
1	E	128	THR	2.4
1	D	66	GLY	2.4
1	G	23	HIS	2.4
1	E	23	HIS	2.3
1	C	69	PRO	2.3
1	C	206	GLN	2.3
1	G	45	ASP	2.3
1	A	115	SER	2.3
1	C	201	ARG	2.2
1	B	196	LEU	2.2
1	C	205	ALA	2.2
1	F	117	GLN	2.2
1	B	116	GLY	2.2
1	G	127	HIS	2.2
1	G	72	ASP	2.2
1	E	78	HIS	2.2
1	E	74	ASP	2.1
1	F	206	GLN	2.1
1	E	204	ARG	2.1
1	B	12	ARG	2.0
1	G	10	THR	2.0
1	B	200	GLU	2.0

## 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 carbohydrates in this entry.

### 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	HG	D	301	1/1	0.84	0.63	6.59	400,400,400,400	0
2	HG	C	301	1/1	0.97	0.05	-3.25	96,96,96,96	0
2	HG	G	301	1/1	0.96	0.06	-3.40	130,130,130,130	0
2	HG	B	301	1/1	0.96	0.06	-3.41	102,102,102,102	0
2	HG	F	301	1/1	0.96	0.05	-3.58	94,94,94,94	0
2	HG	A	301	1/1	0.97	0.07	-4.97	96,96,96,96	0
2	HG	E	301	1/1	0.97	0.06	-5.46	94,94,94,94	0

### 6.5 Other polymers [i](#)

There are no such residues in this entry.