



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 22, 2020 – 01:11 PM BST

PDB ID : 5ME7  
Title : Crystal Structure of eiF4E from *C. melo*  
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Deposited on : 2016-11-14  
Resolution : 2.20 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

<https://www.wwpdb.org/validation/2017/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 : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.13.1  
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) : 2.13.1

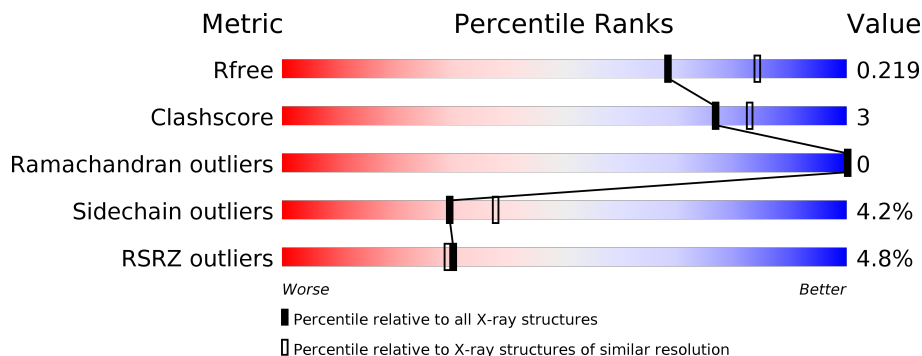
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.20 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	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

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 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	186	 84% 8% • 6%
1	B	186	 2% 82% 8% • 7%
1	C	186	 2% 90% 8% •
1	D	186	 15% 81% 10% •• 7%

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	GOL	C	301	-	-	-	X
2	GOL	C	302	-	-	-	X

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 5880 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 Eukaryotic transcription initiation factor 4E.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	175	1424	915	244	258	7	0	0	0
1	B	173	1415	911	242	255	7	0	0	0
1	C	182	1460	936	249	268	7	0	0	0
1	D	173	1409	906	241	255	7	0	0	0

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	50	SER	-	expression tag	UNP Q00LS8
A	51	ASN	-	expression tag	UNP Q00LS8
A	52	ALA	-	expression tag	UNP Q00LS8
B	50	SER	-	expression tag	UNP Q00LS8
B	51	ASN	-	expression tag	UNP Q00LS8
B	52	ALA	-	expression tag	UNP Q00LS8
C	50	SER	-	expression tag	UNP Q00LS8
C	51	ASN	-	expression tag	UNP Q00LS8
C	52	ALA	-	expression tag	UNP Q00LS8
D	50	SER	-	expression tag	UNP Q00LS8
D	51	ASN	-	expression tag	UNP Q00LS8
D	52	ALA	-	expression tag	UNP Q00LS8

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	C O	0	0
			6	3 3		
2	B	1	Total	C O	0	0
			6	3 3		
2	C	1	Total	C O	0	0
			6	3 3		
2	C	1	Total	C O	0	0
			6	3 3		
2	C	1	Total	C O	0	0
			6	3 3		


- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	42	Total	O	0	0
			42	42		
3	B	46	Total	O	0	0
			46	46		
3	C	46	Total	O	0	0
			46	46		
3	D	8	Total	O	0	0
			8	8		

### 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.

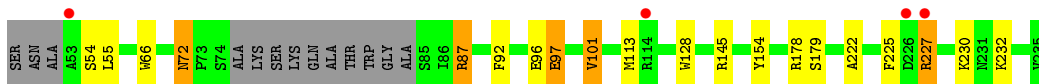
- Molecule 1: Eukaryotic transcription initiation factor 4E

Chain A: 




- Molecule 1: Eukaryotic transcription initiation factor 4E

Chain B: 




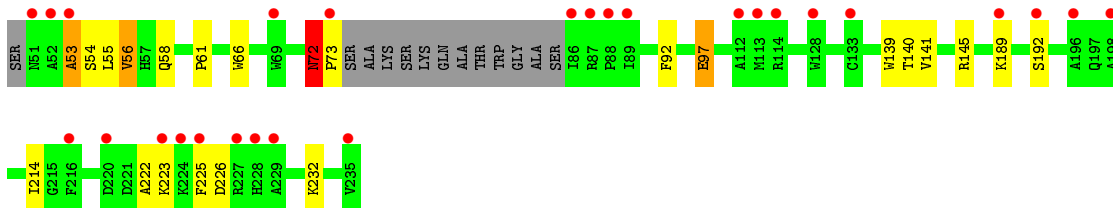
- Molecule 1: Eukaryotic transcription initiation factor 4E

Chain C: 



- Molecule 1: Eukaryotic transcription initiation factor 4E

Chain D: 



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	57.75Å 108.96Å 122.14Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.20 47.08 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.2 (30.00-2.20) 99.3 (47.08-2.20)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.10	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.25 (at 2.20Å)	Xtrriage
Refinement program	REFMAC 5.8.0135	Depositor
R, $R_{free}$	0.187 , 0.214 0.192 , 0.219	Depositor DCC
$R_{free}$ test set	1985 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	39.1	Xtrriage
Anisotropy	0.418	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 37.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	5880	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	50.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.90% 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 i

### 5.1 Standard geometry i

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

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
1	A	1.01	2/1470 (0.1%)	0.96	4/1991 (0.2%)
1	B	1.06	3/1462 (0.2%)	1.06	6/1981 (0.3%)
1	C	0.99	0/1507	0.99	6/2045 (0.3%)
1	D	0.82	0/1456	0.97	5/1974 (0.3%)
All	All	0.98	5/5895 (0.1%)	1.00	21/7991 (0.3%)

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	D	0	1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	97	GLU	CD-OE2	6.02	1.32	1.25
1	A	166	CYS	CB-SG	-5.44	1.73	1.81
1	B	97	GLU	CG-CD	5.12	1.59	1.51
1	B	96	GLU	CG-CD	5.11	1.59	1.51
1	A	100	SER	CB-OG	-5.10	1.35	1.42

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	87	ARG	NE-CZ-NH1	10.28	125.44	120.30
1	A	145	ARG	NE-CZ-NH2	-9.20	115.70	120.30
1	B	145	ARG	NE-CZ-NH2	-9.03	115.79	120.30
1	B	145	ARG	NE-CZ-NH1	8.79	124.69	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	227	ARG	NE-CZ-NH1	8.28	124.44	120.30
1	C	189	LYS	N-CA-C	8.16	133.02	111.00
1	D	145	ARG	NE-CZ-NH1	7.91	124.25	120.30
1	A	145	ARG	NE-CZ-NH1	7.89	124.25	120.30
1	D	145	ARG	NE-CZ-NH2	-7.83	116.39	120.30
1	C	145	ARG	NE-CZ-NH2	-7.72	116.44	120.30
1	C	87	ARG	NE-CZ-NH2	-7.09	116.75	120.30
1	C	145	ARG	NE-CZ-NH1	6.76	123.68	120.30
1	D	97	GLU	N-CA-CB	6.49	122.28	110.60
1	B	154	TYR	CB-CG-CD2	-6.06	117.36	121.00
1	D	189	LYS	CD-CE-NZ	5.99	125.48	111.70
1	A	225	PHE	N-CA-C	-5.72	95.55	111.00
1	B	87	ARG	NE-CZ-NH1	5.48	123.04	120.30
1	A	235	VAL	N-CA-CB	-5.35	99.74	111.50
1	B	227	ARG	NE-CZ-NH2	-5.17	117.72	120.30
1	D	53	ALA	C-N-CA	5.15	134.57	121.70
1	C	178	ARG	NE-CZ-NH1	5.08	122.84	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	72	ASN	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	1424	0	1335	9	0
1	B	1415	0	1328	11	0
1	C	1460	0	1366	5	0
1	D	1409	0	1313	14	0
2	A	6	0	8	0	0
2	B	6	0	8	0	0
2	C	18	0	24	0	0
3	A	42	0	0	0	0
3	B	46	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	46	0	0	0	0
3	D	8	0	0	0	0
All	All	5880	0	5382	33	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 3.

All (33) 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:D:72:ASN:HD22	1:D:73:PRO:HD2	1.36	0.89
1:B:113:MET:HE1	1:B:178:ARG:C	1.98	0.84
1:A:228:HIS:HB3	1:D:58:GLN:HE22	1.54	0.70
1:B:113:MET:HE1	1:B:178:ARG:CA	2.25	0.66
1:D:72:ASN:ND2	1:D:73:PRO:HD2	2.11	0.66
1:A:225:PHE:CE1	1:D:61:PRO:HG3	2.33	0.63
1:D:53:ALA:HA	1:D:55:LEU:H	1.68	0.58
1:D:139:TRP:HB3	1:D:214:ILE:HD11	1.86	0.56
1:A:228:HIS:CB	1:D:58:GLN:HE22	2.19	0.55
1:B:113:MET:HA	1:B:113:MET:HE3	1.92	0.51
1:D:214:ILE:HG23	1:D:214:ILE:O	2.12	0.50
1:D:53:ALA:HB1	1:D:56:VAL:HG23	1.94	0.48
1:B:72:ASN:HD22	1:B:72:ASN:C	2.17	0.48
1:A:52:ALA:HA	1:A:55:LEU:HB2	1.95	0.47
1:D:140:THR:O	1:D:214:ILE:HG13	2.14	0.47
1:B:97:GLU:O	1:B:101:VAL:HG12	2.14	0.46
1:B:97:GLU:HG2	1:C:80:ALA:O	2.16	0.46
1:B:222:ALA:O	1:B:225:PHE:O	2.35	0.45
1:B:113:MET:CE	1:B:179:SER:N	2.80	0.44
1:B:113:MET:HE1	1:B:179:SER:N	2.31	0.44
1:A:66:TRP:HB2	1:A:92:PHE:CZ	2.53	0.44
1:C:165:ASP:O	1:C:166:CYS:HB2	2.19	0.43
1:D:222:ALA:O	1:D:225:PHE:O	2.37	0.42
1:A:123:LYS:HG3	1:C:179:SER:HB3	2.02	0.42
1:A:222:ALA:O	1:A:225:PHE:O	2.39	0.41
1:C:65:SER:OG	1:C:93:SER:HB3	2.21	0.41
1:A:224:LYS:HB3	1:D:61:PRO:HB3	2.02	0.41
1:D:66:TRP:HB2	1:D:92:PHE:CZ	2.56	0.41
1:B:66:TRP:HB2	1:B:92:PHE:CZ	2.56	0.41
1:B:113:MET:CA	1:B:113:MET:HE3	2.49	0.41
1:A:60:HIS:CD2	1:A:96:GLU:OE1	2.74	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:66:TRP:HB2	1:C:92:PHE:CZ	2.57	0.40
1:D:214:ILE:CG2	1:D:214:ILE:O	2.69	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	171/186 (92%)	165 (96%)	6 (4%)	0	100	100
1	B	169/186 (91%)	164 (97%)	5 (3%)	0	100	100
1	C	178/186 (96%)	173 (97%)	5 (3%)	0	100	100
1	D	169/186 (91%)	163 (96%)	6 (4%)	0	100	100
All	All	687/744 (92%)	665 (97%)	22 (3%)	0	100	100

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	149/159 (94%)	144 (97%)	5 (3%)	37	47
1	B	149/159 (94%)	140 (94%)	9 (6%)	19	22
1	C	152/159 (96%)	150 (99%)	2 (1%)	69	81

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	147/159 (92%)	138 (94%)	9 (6%)	18	21
All	All	597/636 (94%)	572 (96%)	25 (4%)	30	38

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	51	ASN
1	A	55	LEU
1	A	87	ARG
1	A	232	LYS
1	A	235	VAL
1	B	54	SER
1	B	55	LEU
1	B	72	ASN
1	B	87	ARG
1	B	101	VAL
1	B	128	TRP
1	B	227	ARG
1	B	230	LYS
1	B	232	LYS
1	C	78	LYS
1	C	232	LYS
1	D	54	SER
1	D	56	VAL
1	D	72	ASN
1	D	97	GLU
1	D	141	VAL
1	D	192	SER
1	D	223	LYS
1	D	226	ASP
1	D	232	LYS

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	60	HIS
1	A	228	HIS
1	B	72	ASN
1	B	219	HIS
1	C	122	HIS
1	D	58	GLN

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Mol	Chain	Res	Type
1	D	72	ASN
1	D	122	HIS
1	D	228	HIS

### 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 monosaccharides in this entry.

### 5.6 Ligand geometry [i](#)

5 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
2	GOL	A	301	-	5,5,5	0.55	0	5,5,5	0.56	0
2	GOL	B	301	-	5,5,5	0.49	0	5,5,5	0.31	0
2	GOL	C	302	-	5,5,5	0.64	0	5,5,5	0.28	0
2	GOL	C	303	-	5,5,5	0.56	0	5,5,5	0.26	0
2	GOL	C	301	-	5,5,5	0.54	0	5,5,5	1.11	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GOL	A	301	-	-	1/4/4/4	-
2	GOL	B	301	-	-	2/4/4/4	-
2	GOL	C	302	-	-	0/4/4/4	-
2	GOL	C	303	-	-	2/4/4/4	-
2	GOL	C	301	-	-	4/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	303	GOL	C1-C2-C3-O3
2	C	301	GOL	C1-C2-C3-O3
2	B	301	GOL	C1-C2-C3-O3
2	C	301	GOL	O2-C2-C3-O3
2	C	303	GOL	O2-C2-C3-O3
2	C	301	GOL	O1-C1-C2-O2
2	A	301	GOL	O2-C2-C3-O3
2	B	301	GOL	O2-C2-C3-O3
2	C	301	GOL	O1-C1-C2-C3

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	175/186 (94%)	-0.11	0 <b>100</b> <b>100</b>	27, 40, 71, 94	0
1	B	173/186 (93%)	-0.20	4 (2%) 60 58	27, 40, 68, 91	0
1	C	182/186 (97%)	-0.25	3 (1%) 72 70	27, 42, 75, 97	0
1	D	173/186 (93%)	0.82	27 (15%) <b>2</b> <b>1</b>	40, 68, 97, 126	0
All	All	703/744 (94%)	0.06	34 (4%) 30 29	27, 46, 88, 126	0

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	224	LYS	6.2
1	D	189	LYS	4.9
1	D	128	TRP	4.1
1	D	196	ALA	4.0
1	D	133	CYS	4.0
1	D	87	ARG	4.0
1	D	86	ILE	3.8
1	D	220	ASP	3.8
1	D	225	PHE	3.8
1	C	53	ALA	3.7
1	D	227	ARG	3.6
1	D	114	ARG	3.4
1	C	229	ALA	3.4
1	D	51	ASN	3.2
1	D	53	ALA	3.2
1	D	235	VAL	3.0
1	D	223	LYS	3.0
1	B	53	ALA	2.7
1	D	89	ILE	2.7
1	D	216	PHE	2.7
1	C	228	HIS	2.6

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Mol	Chain	Res	Type	RSRZ
1	D	192	SER	2.5
1	B	227	ARG	2.5
1	B	114	ARG	2.5
1	D	228	HIS	2.5
1	D	198	ALA	2.5
1	D	69	TRP	2.4
1	D	73	PRO	2.3
1	D	229	ALA	2.3
1	D	113	MET	2.2
1	B	226	ASP	2.2
1	D	112	ALA	2.2
1	D	52	ALA	2.1
1	D	88	PRO	2.1

## 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 monosaccharides in this entry.

## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<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	B-factors(Å <sup>2</sup> )	Q<0.9
2	GOL	B	301	6/6	0.70	0.26	67,93,98,100	0
2	GOL	C	301	6/6	0.71	0.66	75,83,86,86	0
2	GOL	C	302	6/6	0.74	0.44	69,78,85,85	0
2	GOL	C	303	6/6	0.83	0.16	51,70,73,78	0
2	GOL	A	301	6/6	0.89	0.24	42,74,80,89	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.