



# wwPDB X-ray Structure Validation Summary Report ⓘ

Dec 6, 2023 – 07:04 am GMT

PDB ID : 1OL2  
Title : Cyclin A binding groove inhibitor H-Arg-Arg-Leu-Asn-(p-F-Phe)-NH<sub>2</sub>  
Authors : Kontopidis, G.; Andrews, M.; McInnes, C.; Cowan, A.; Powers, H.; Innes, L.; Plater, A.; Griffiths, G.; Paterson, D.; Zheleva, D.; Lane, D.; Green, S.; Walkinshaw, M.; Fischer, P.  
Deposited on : 2003-08-05  
Resolution : 2.60 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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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 : 1.8.4, CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.36  
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) : 2.36

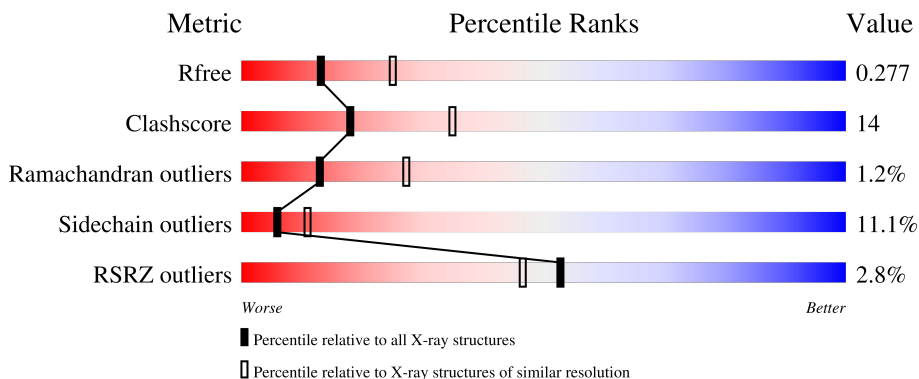
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.60 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	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	298	
1	C	298	
2	B	260	
2	D	260	
3	E	6	

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Mol	Chain	Length	Quality of chain
3	F	6	 <p>A horizontal bar chart representing the quality of chain. The bar is divided into four segments: a red segment on the left labeled '17%', a green segment labeled '33%', a yellow segment labeled '50%', and a red segment on the right labeled '17%'.</p>

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 9386 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 CELL DIVISION PROTEIN KINASE 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	296	Total 2378	C 1547	N 403	O 420	S 8	0	0	0
1	C	296	Total 2378	C 1547	N 403	O 420	S 8	0	0	0

- Molecule 2 is a protein called CYCLIN A2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	258	Total 2083	C 1350	N 339	O 383	S 11	0	0	0
2	D	258	Total 2084	C 1350	N 339	O 384	S 11	0	0	0

- Molecule 3 is a protein called ARG-ARG-LEU-ASN-PFF-NH2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	F	N	O			
3	E	6	Total 51	C 31	F 1	N 13	O 6	0	0	1
3	F	6	Total 51	C 31	F 1	N 13	O 6	0	0	1

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	93	Total 93	O 93	0	0
4	B	75	Total 75	O 75	0	0
4	C	102	Total 102	O 102	0	0
4	D	85	Total 85	O 85	0	0

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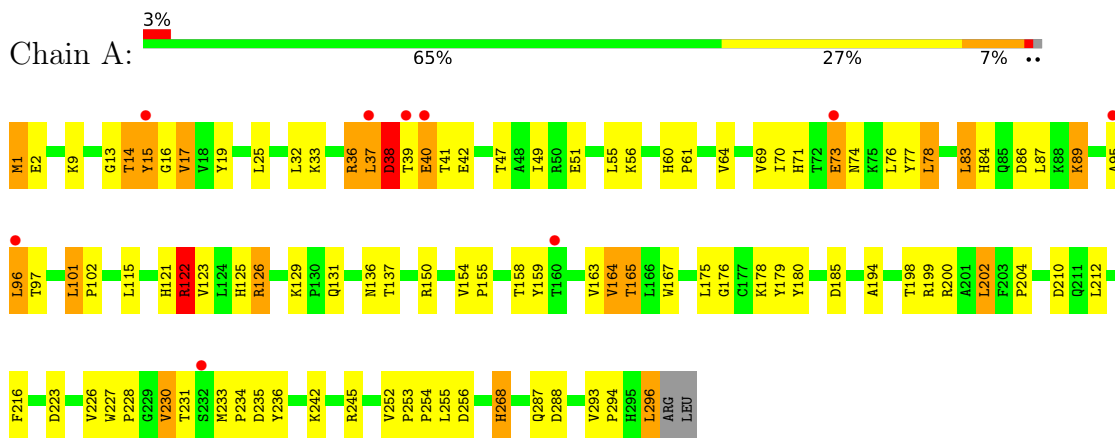
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<b>Mol</b>	<b>Chain</b>	<b>Residues</b>	<b>Atoms</b>	<b>ZeroOcc</b>	<b>AltConf</b>
4	E	3	Total O 3 3	0	0
4	F	3	Total O 3 3	0	0

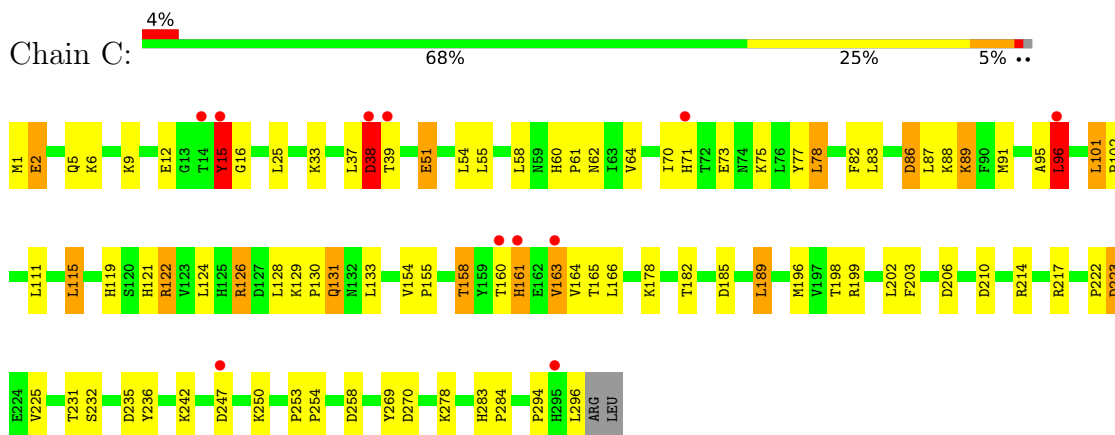
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

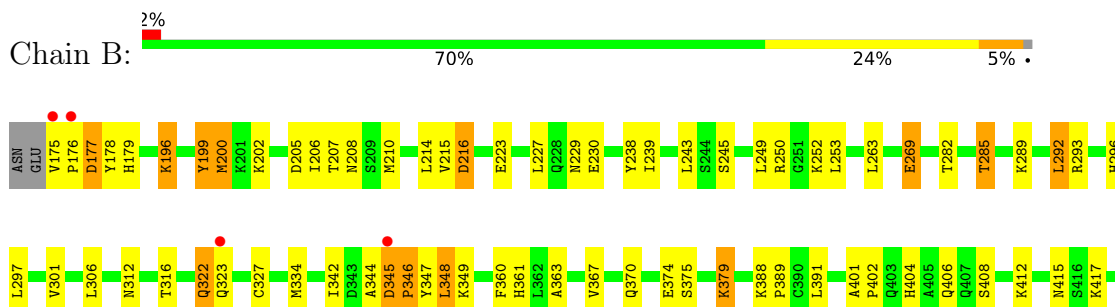
- Molecule 1: CELL DIVISION PROTEIN KINASE 2



- Molecule 1: CELL DIVISION PROTEIN KINASE 2

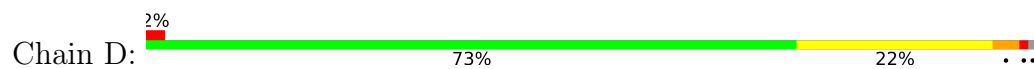


- Molecule 2: CYCLIN A2

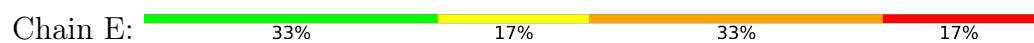




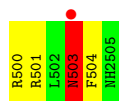
- Molecule 2: CYCLIN A2



- Molecule 3: ARG-ARG-LEU-ASN-PFF-NH2



- Molecule 3: ARG-ARG-LEU-ASN-PFF-NH2



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	73.54Å 112.98Å 153.09Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	12.00 – 2.60 11.98 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.6 (12.00-2.60) 99.6 (11.98-2.60)	Depositor EDS
$R_{merge}$	0.11	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.61 (at 2.59Å)	Xtrriage
Refinement program	REFMAC 5.0	Depositor
R, $R_{free}$	0.191 , 0.290 0.201 , 0.277	Depositor DCC
$R_{free}$ test set	1220 reflections (3.10%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	39.0	Xtrriage
Anisotropy	0.106	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 40.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	9386	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	44.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.53% 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: NH2, PFF

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	0.59	0/2440	0.99	11/3313 (0.3%)
1	C	0.60	0/2440	0.98	8/3313 (0.2%)
2	B	0.63	0/2133	0.95	3/2897 (0.1%)
2	D	0.60	0/2134	0.99	8/2897 (0.3%)
3	E	2.03	3/37 (8.1%)	2.25	2/47 (4.3%)
3	F	2.12	3/37 (8.1%)	1.65	0/47
All	All	0.63	6/9221 (0.1%)	0.99	32/12514 (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	A	0	3
3	E	0	2
All	All	0	5

The worst 5 of 6 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	F	500	ARG	CB-CG	-7.52	1.32	1.52
3	F	501	ARG	CB-CG	-7.01	1.33	1.52
3	E	500	ARG	CB-CG	-6.92	1.33	1.52
3	E	501	ARG	CB-CG	-6.88	1.33	1.52
3	F	503	ASN	CB-CG	-6.74	1.35	1.51

The worst 5 of 32 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	240	ASP	CB-CG-OD2	9.86	127.17	118.30
1	A	223	ASP	CB-CG-OD2	8.16	125.64	118.30
1	C	86	ASP	CB-CG-OD2	7.78	125.30	118.30
1	A	122	ARG	NE-CZ-NH1	-7.70	116.45	120.30
3	E	500	ARG	CA-CB-CG	7.42	129.73	113.40

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	13	GLY	Peptide
1	A	163	VAL	Peptide
1	A	83	LEU	Peptide
3	E	500	ARG	Sidechain
3	E	501	ARG	Sidechain

## 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	2378	0	2426	93	0
1	C	2378	0	2426	73	1
2	B	2083	0	2107	47	1
2	D	2084	0	2107	41	0
3	E	51	0	49	3	0
3	F	51	0	49	3	0
4	A	93	0	0	8	0
4	B	75	0	0	6	0
4	C	102	0	0	11	0
4	D	85	0	0	6	0
4	E	3	0	0	0	0
4	F	3	0	0	0	0
All	All	9386	0	9164	246	1

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 14.

The worst 5 of 246 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:A:37:LEU:HD13	1:A:37:LEU:O	1.10	1.23
1:A:37:LEU:O	1:A:37:LEU:CD1	1.88	1.21
1:A:15:TYR:CE2	1:A:33:LYS:HD3	1.93	1.04
1:A:15:TYR:CD2	1:A:33:LYS:HD3	1.96	1.01
3:F:503:ASN:HD22	3:F:503:ASN:H	1.01	0.97

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:431:ASN:ND2	1:C:210:ASP:OD1[2_664]	1.82	0.38

## 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	294/298 (99%)	273 (93%)	17 (6%)	4 (1%)	11	22
1	C	294/298 (99%)	277 (94%)	14 (5%)	3 (1%)	15	32
2	B	256/260 (98%)	246 (96%)	6 (2%)	4 (2%)	9	19
2	D	256/260 (98%)	247 (96%)	7 (3%)	2 (1%)	19	39
3	E	3/6 (50%)	3 (100%)	0	0	100	100
3	F	3/6 (50%)	3 (100%)	0	0	100	100
All	All	1106/1128 (98%)	1049 (95%)	44 (4%)	13 (1%)	13	27

5 of 13 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	165	THR
1	C	161	HIS
1	A	96	LEU
2	B	424	LEU

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Mol	Chain	Res	Type
1	C	96	LEU

### 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	261/263 (99%)	227 (87%)	34 (13%)	4	7
1	C	261/263 (99%)	231 (88%)	30 (12%)	5	10
2	B	232/234 (99%)	207 (89%)	25 (11%)	6	12
2	D	232/234 (99%)	213 (92%)	19 (8%)	11	22
3	E	4/4 (100%)	3 (75%)	1 (25%)	0	1
3	F	4/4 (100%)	3 (75%)	1 (25%)	0	1
All	All	994/1002 (99%)	884 (89%)	110 (11%)	6	11

5 of 110 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	431	ASN
1	C	88	LYS
3	F	503	ASN
2	D	293	ARG
1	C	2	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 16 such sidechains are listed below:

Mol	Chain	Res	Type
2	D	431	ASN
2	D	404	HIS
2	B	431	ASN
2	D	208	ASN
2	B	419	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](#)

2 non-standard protein/DNA/RNA residues 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
3	PFF	F	504	3	11,12,13	0.76	0	12,15,17	1.16	1 (8%)
3	PFF	E	504	3	11,12,13	0.80	0	12,15,17	1.20	2 (16%)

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
3	PFF	F	504	3	-	2/5/6/8	0/1/1/1
3	PFF	E	504	3	-	0/5/6/8	0/1/1/1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	504	PFF	CB-CA-C	-2.91	106.00	111.47
3	E	504	PFF	CE2-CZ-CE1	-2.41	119.62	122.83
3	E	504	PFF	CD1-CE1-CZ	2.22	120.66	118.36

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	F	504	PFF	N-CA-CB-CG

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Mol	Chain	Res	Type	Atoms
3	F	504	PFF	C-CA-CB-CG

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	E	504	PFF	1	0

## 5.5 Carbohydrates [i](#)

There are no monosaccharides 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, 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	296/298 (99%)	-0.35	9 (3%) 50 43	21, 39, 79, 97	0
1	C	296/298 (99%)	-0.29	11 (3%) 41 34	21, 38, 86, 118	0
2	B	258/260 (99%)	-0.38	4 (1%) 72 68	25, 39, 70, 108	0
2	D	258/260 (99%)	-0.30	6 (2%) 60 54	23, 44, 71, 98	0
3	E	4/6 (66%)	-0.24	0 100 100	48, 49, 61, 79	0
3	F	4/6 (66%)	0.35	1 (25%) 0 0	38, 49, 56, 80	0
All	All	1116/1128 (98%)	-0.33	31 (2%) 53 46	21, 40, 77, 118	0

The worst 5 of 31 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	39	THR	5.1
1	C	14	THR	4.1
1	C	15	TYR	3.6
1	C	163	VAL	3.6
2	B	345	ASP	3.4

### 6.2 Non-standard residues in protein, DNA, RNA chains [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
3	PFF	E	504	12/13	0.86	0.18	34,47,78,87	0
3	PFF	F	504	12/13	0.90	0.15	32,48,52,54	0

### 6.3 Carbohydrates [i](#)

There are no monosaccharides 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.