



# Full wwPDB X-ray Structure Validation Report ⓘ

May 29, 2020 – 09:07 am BST

PDB ID : 5ZBU  
Title : Crystal Structure of PA-TM-RING E3 ligase RNF13 RING domain in complex with E2 Ub  
Authors : Datta, A.B.; Sarkar, S.  
Deposited on : 2018-02-12  
Resolution : 3.20 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
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

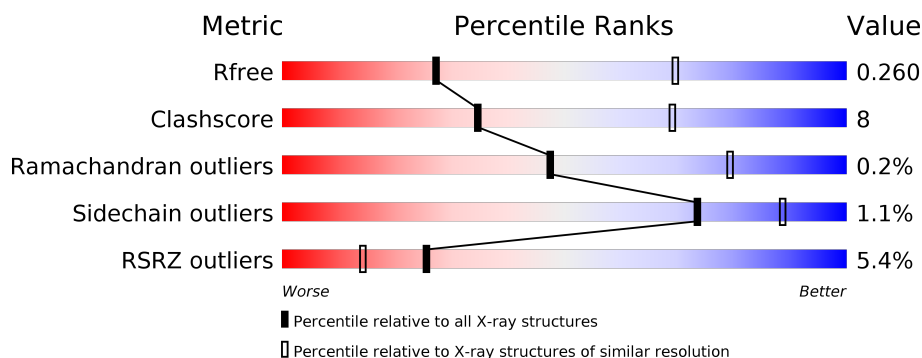
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.20 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	1133 (3.20-3.20)
Clashscore	141614	1253 (3.20-3.20)
Ramachandran outliers	138981	1234 (3.20-3.20)
Sidechain outliers	138945	1233 (3.20-3.20)
RSRZ outliers	127900	1095 (3.20-3.20)

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	B	148	<div> <div>12%</div> <div> <div></div> <div>78%</div> <div>20%</div> <div>..</div> </div> </div>
1	C	148	<div> <div>4%</div> <div> <div></div> <div>84%</div> <div>16%</div> </div> </div>
2	A	76	<div> <div>%</div> <div> <div></div> <div>74%</div> <div>22%</div> <div>..</div> </div> </div>
2	D	76	<div> <div>3%</div> <div> <div></div> <div>68%</div> <div>26%</div> <div>5%</div> </div> </div>
3	E	76	<div> <div>%</div> <div> <div></div> <div>86%</div> <div>13%</div> <div>.</div> </div> </div>

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4163 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 Ubiquitin-conjugating enzyme E2 D2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	148	Total	C	N	O	S	0	0	0
			1185	760	204	214	7			
1	B	147	Total	C	N	O	S	0	0	0
			1176	755	203	212	6			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	0	GLY	-	expression tag	UNP P62837
C	85	LYS	CYS	engineered mutation	UNP P62837
B	0	GLY	-	expression tag	UNP P62837
B	85	LYS	CYS	engineered mutation	UNP P62837

- Molecule 2 is a protein called E3 ubiquitin-protein ligase RNF13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	74	Total	C	N	O	S	0	0	0
			601	381	107	106	7			
2	D	72	Total	C	N	O	S	0	0	0
			587	373	104	102	8			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	215	MET	-	expression tag	UNP O43567
D	215	MET	-	expression tag	UNP O43567

- Molecule 3 is a protein called Ubiquitin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	76	Total	C	N	O	S	0	0	0
			601	378	105	117	1			

- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total 2	Zn 2	0	0
4	D	2	Total 2	Zn 2	0	0

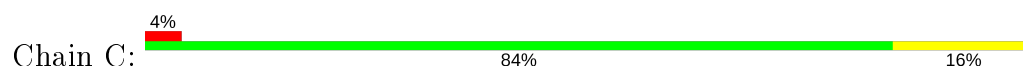
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	C	3	Total 3	O 3	0	0
5	A	3	Total 3	O 3	0	0
5	D	1	Total 1	O 1	0	0
5	E	2	Total 2	O 2	0	0

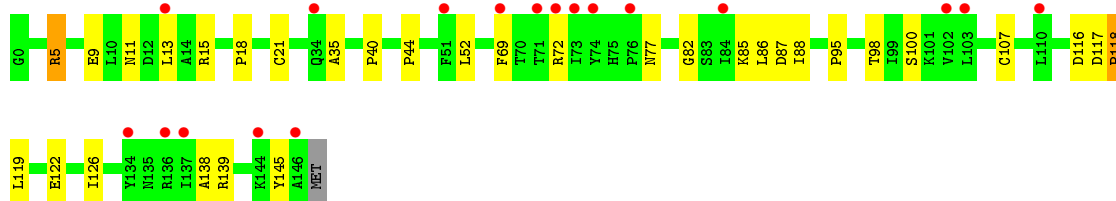
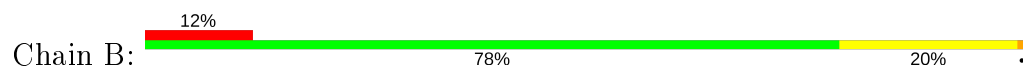
### 3 Residue-property plots [i](#)

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: Ubiquitin-conjugating enzyme E2 D2



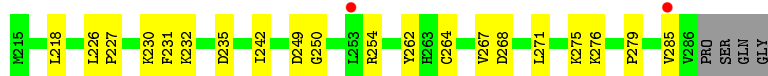
- Molecule 1: Ubiquitin-conjugating enzyme E2 D2



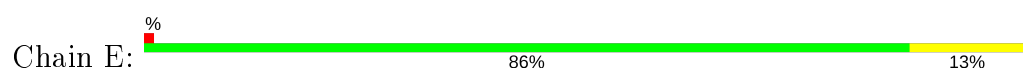
- Molecule 2: E3 ubiquitin-protein ligase RNF13

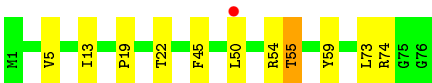


- Molecule 2: E3 ubiquitin-protein ligase RNF13



- Molecule 3: Ubiquitin





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	56.56Å 105.74Å 60.79Å 90.00° 115.49° 90.00°	Depositor
Resolution (Å)	28.26 – 3.20 48.71 – 3.20	Depositor EDS
% Data completeness (in resolution range)	99.7 (28.26-3.20) 99.7 (48.71-3.20)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.50 (at 3.19Å)	Xtriage
Refinement program	PHENIX (1.13_2998: ???)	Depositor
R, $R_{free}$	0.214 , 0.262 0.213 , 0.260	Depositor DCC
$R_{free}$ test set	559 reflections (5.23%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	90.2	Xtriage
Anisotropy	0.412	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 60.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.036 for h,-k,-h-l	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	4163	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	105.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.71% 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: ZN

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	B	0.27	0/1212	0.46	0/1652
1	C	0.26	0/1221	0.44	0/1662
2	A	0.25	0/614	0.45	0/825
2	D	0.26	0/599	0.50	0/803
3	E	0.25	0/607	0.45	0/816
All	All	0.26	0/4253	0.46	0/5758

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1176	0	1164	27	0
1	C	1185	0	1176	15	0
2	A	601	0	614	10	0
2	D	587	0	603	15	0
3	E	601	0	629	10	0
4	A	2	0	0	0	0
4	D	2	0	0	0	0
5	A	3	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	C	3	0	0	0	0
5	D	1	0	0	0	0
5	E	2	0	0	0	0
All	All	4163	0	4186	69	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 8.

All (69) 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:C:79:ASN:OD1	1:C:80:SER:N	2.09	0.86
1:B:5:ARG:NH2	1:B:9:GLU:OE2	2.10	0.85
1:B:85:LYS:HE2	3:E:73:LEU:HD23	1.69	0.75
2:D:226:LEU:HD11	2:D:285:VAL:HG11	1.68	0.73
1:C:5:ARG:NH2	2:D:242:ILE:O	2.25	0.69
1:C:20:GLN:NE2	1:C:38:MET:O	2.25	0.69
1:C:26:VAL:HG22	1:C:34:GLN:HB2	1.79	0.65
1:B:86:LEU:HD12	3:E:74:ARG:HH21	1.65	0.61
1:B:72:ARG:H	1:B:72:ARG:HD2	1.66	0.59
2:A:269:PRO:O	2:A:273:LYS:HB2	2.04	0.58
2:D:230:LYS:HE2	2:D:250:GLY:HA2	1.87	0.56
1:B:117:ASP:OD2	3:E:74:ARG:HG2	2.06	0.56
2:A:254:ARG:NH2	2:A:268:ASP:OD2	2.28	0.55
1:B:86:LEU:HD12	3:E:74:ARG:NH2	2.21	0.55
1:C:85:LYS:HB2	1:C:119:LEU:HG	1.89	0.55
1:B:95:PRO:HB2	2:A:242:ILE:HG23	1.89	0.54
1:B:13:LEU:HD21	1:B:100:SER:HB2	1.89	0.54
1:C:94:SER:OG	2:D:279:PRO:O	2.15	0.54
1:C:5:ARG:NH1	1:C:97:LEU:O	2.41	0.53
1:C:122:GLU:O	1:C:126:ILE:HG12	2.09	0.53
2:A:217:ARG:NH2	5:A:1101:HOH:O	2.42	0.52
1:B:72:ARG:HD3	1:B:145:TYR:CD2	2.45	0.51
1:B:77:ASN:ND2	1:B:117:ASP:O	2.44	0.51
2:A:241:ALA:HB3	2:A:262:TYR:HE2	1.76	0.51
2:D:231:PHE:HD2	2:D:249:ASP:HA	1.75	0.50
1:B:40:PRO:HG3	1:B:107:CYS:SG	2.52	0.50
1:B:69:PHE:O	1:B:82:GLY:HA3	2.11	0.50
1:B:72:ARG:N	1:B:72:ARG:HD2	2.27	0.49
2:D:218:LEU:HG	2:D:276:LYS:HB2	1.93	0.49
1:B:35:ALA:HB3	1:B:52:LEU:HB2	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:85:LYS:HE2	1:C:117:ASP:HB2	1.94	0.49
2:D:271:LEU:HA	2:D:275:LYS:O	2.13	0.48
1:B:116:ASP:O	1:B:118:PRO:HD3	2.13	0.48
1:B:11:ASN:O	1:B:15:ARG:HG2	2.14	0.48
1:B:5:ARG:HH22	1:B:98:THR:HG22	1.80	0.47
1:C:74:TYR:HB2	1:C:141:TRP:CD2	2.50	0.47
2:D:218:LEU:HG	2:D:276:LYS:CB	2.45	0.47
1:C:34:GLN:HE21	1:C:51:PHE:HD2	1.63	0.46
3:E:54:ARG:HB2	3:E:59:TYR:HE1	1.81	0.46
2:A:263:HIS:HB2	2:A:266:CYS:SG	2.55	0.46
1:C:126:ILE:HD12	1:C:133:LYS:HD2	1.98	0.45
1:C:86:LEU:HG	1:C:88:ILE:HG12	1.98	0.45
1:C:79:ASN:HB3	1:C:83:SER:H	1.82	0.44
1:B:69:PHE:N	1:B:82:GLY:O	2.32	0.44
2:A:232:LYS:N	2:A:235:ASP:OD2	2.45	0.44
2:A:271:LEU:HA	2:A:275:LYS:O	2.17	0.44
1:B:87:ASP:OD1	3:E:74:ARG:NH2	2.50	0.44
2:D:271:LEU:HD23	2:D:271:LEU:HA	1.86	0.44
1:B:122:GLU:O	1:B:126:ILE:HG12	2.18	0.44
2:D:262:TYR:HB2	2:D:267:VAL:HB	2.00	0.43
1:B:72:ARG:HD3	1:B:145:TYR:CE2	2.53	0.43
1:C:44:PRO:HB2	1:C:138:ALA:HB3	2.01	0.43
3:E:22:THR:HA	3:E:55:THR:HA	2.01	0.43
2:A:220:LYS:O	2:A:224:LYS:HD3	2.19	0.43
3:E:5:VAL:HB	3:E:13:ILE:HG13	2.01	0.43
1:B:44:PRO:HB2	1:B:138:ALA:HB3	2.01	0.43
2:D:226:LEU:HB2	2:D:254:ARG:HD2	2.01	0.42
3:E:45:PHE:HB3	3:E:50:LEU:HD21	2.00	0.42
1:B:44:PRO:HA	1:B:139:ARG:HD3	2.00	0.42
2:D:267:VAL:HG22	2:D:271:LEU:HG	2.00	0.42
1:B:86:LEU:HG	1:B:88:ILE:HG12	2.01	0.42
1:B:87:ASP:OD1	1:B:87:ASP:N	2.52	0.42
1:B:119:LEU:HA	1:B:119:LEU:HD23	1.69	0.42
1:B:18:PRO:HG2	1:B:21:CYS:O	2.20	0.41
2:A:230:LYS:HB2	2:A:230:LYS:HE2	1.73	0.41
3:E:19:PRO:O	3:E:55:THR:OG1	2.39	0.41
2:D:232:LYS:HB3	2:D:235:ASP:OD2	2.21	0.41
2:D:264:CYS:O	2:D:268:ASP:HB2	2.21	0.41
2:D:226:LEU:HA	2:D:227:PRO:HD3	1.93	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone

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	B	145/148 (98%)	141 (97%)	3 (2%)	1 (1%)	22	61
1	C	146/148 (99%)	142 (97%)	4 (3%)	0	100	100
2	A	72/76 (95%)	69 (96%)	3 (4%)	0	100	100
2	D	70/76 (92%)	64 (91%)	6 (9%)	0	100	100
3	E	74/76 (97%)	74 (100%)	0	0	100	100
All	All	507/524 (97%)	490 (97%)	16 (3%)	1 (0%)	47	79

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	118	PRO

### 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	B	130/131 (99%)	129 (99%)	1 (1%)	81	93
1	C	131/131 (100%)	130 (99%)	1 (1%)	81	93
2	A	70/71 (99%)	68 (97%)	2 (3%)	42	74
2	D	68/71 (96%)	68 (100%)	0	100	100
3	E	68/68 (100%)	67 (98%)	1 (2%)	65	85
All	All	467/472 (99%)	462 (99%)	5 (1%)	73	88

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	81	ASN
1	B	5	ARG
2	A	230	LYS
2	A	237	TYR
3	E	55	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	20	GLN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 4 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

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	B	147/148 (99%)	0.56	18 (12%) 4 2	76, 132, 175, 215	0
1	C	148/148 (100%)	0.11	6 (4%) 37 24	62, 100, 147, 174	0
2	A	74/76 (97%)	-0.04	1 (1%) 75 63	62, 86, 128, 157	0
2	D	72/76 (94%)	0.19	2 (2%) 53 37	59, 86, 114, 123	0
3	E	76/76 (100%)	-0.09	1 (1%) 77 65	54, 79, 106, 149	0
All	All	517/524 (98%)	0.20	28 (5%) 25 14	54, 98, 162, 215	0

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	72	ARG	12.4
1	B	134	TYR	4.5
1	B	74	TYR	4.5
1	B	73	ILE	4.4
1	B	69	PHE	3.7
1	C	123	ILE	3.2
1	B	13	LEU	3.0
1	B	146	ALA	2.9
1	B	110	LEU	2.9
1	C	74	TYR	2.9
1	B	103	LEU	2.9
1	B	84	ILE	2.8
1	C	125	ARG	2.8
1	B	137	ILE	2.8
1	C	134	TYR	2.7
1	B	144	LYS	2.6
2	D	253	LEU	2.6
1	B	136	ARG	2.5
1	B	34	GLN	2.5
1	B	76	PRO	2.4

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Mol	Chain	Res	Type	RSRZ
3	E	50	LEU	2.2
1	C	127	TYR	2.2
1	B	51	PHE	2.1
1	B	102	VAL	2.1
1	B	71	THR	2.1
2	D	285	VAL	2.1
1	C	129	THR	2.1
2	A	230	LYS	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](#)

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( $\text{\AA}^2$ )	Q<0.9
4	ZN	D	1002	1/1	0.96	0.17	88,88,88,88	0
4	ZN	A	1002	1/1	0.98	0.20	70,70,70,70	0
4	ZN	A	1001	1/1	0.99	0.21	74,74,74,74	0
4	ZN	D	1001	1/1	0.99	0.13	71,71,71,71	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.