



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 10:33 PM GMT

PDB ID : 1U5K
Title : Recombinational repair protein RecO
Authors : Makharashvili, N.; Koroleva, O.; Bera, S.; Grandgenett, D.P.; Korolev, S.
Deposited on : 2004-07-27
Resolution : 2.00 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
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

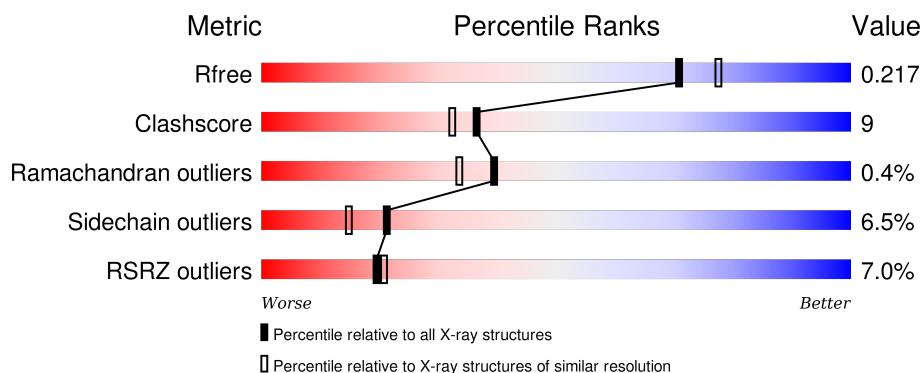
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (2.00-2.00)

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 ≥ 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	244	<div> <div>8%</div> <div>86%</div> <div>9% ...</div> </div>
1	B	244	<div> <div>6%</div> <div>75%</div> <div>15% 5% 5%</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3861 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 hypothetical protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	242	Total	C	N	O	S	0	0	0
			1822	1164	333	319	6			
1	B	232	Total	C	N	O	S	0	0	0
			1764	1125	325	307	7			

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Zn	0	0
			1	1		
2	A	1	Total	Zn	0	0
			1	1		

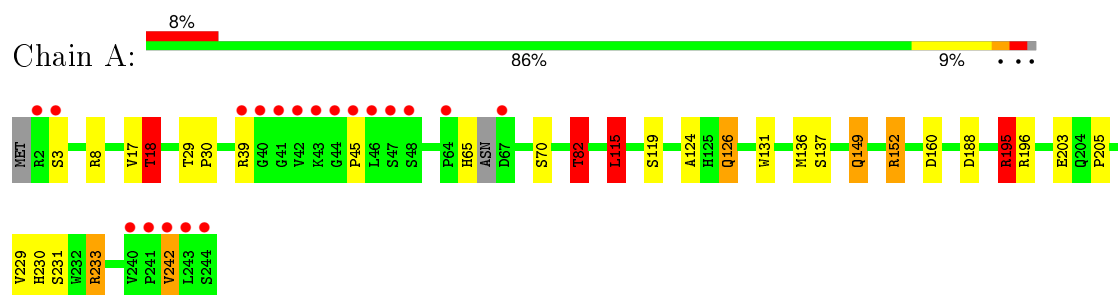
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	165	Total	O	0	0
			165	165		
3	B	108	Total	O	0	0
			108	108		

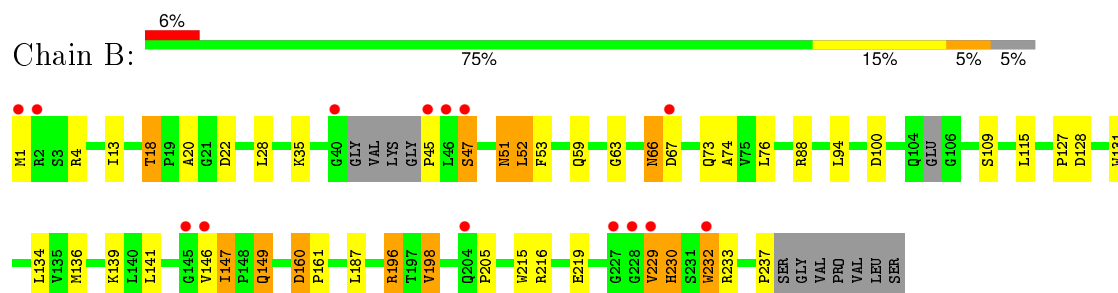
3 Residue-property plots [i](#)

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



• Molecule 1: hypothetical protein



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	134.97Å 52.24Å 100.68Å 90.00° 107.09° 90.00°	Depositor
Resolution (Å)	50.00 – 2.00 33.71 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.5 (50.00-2.00) 99.5 (33.71-2.00)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.58 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.198 , 0.213 0.204 , 0.217	Depositor DCC
R_{free} test set	2296 reflections (5.32%)	DCC
Wilson B-factor (Å ²)	31.9	Xtriage
Anisotropy	0.175	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 46.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 45528 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3861	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.59% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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

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	A	0.37	0/1868	1.04	15/2547 (0.6%)
1	B	0.35	0/1809	0.93	5/2463 (0.2%)
All	All	0.36	0/3677	0.99	20/5010 (0.4%)

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	B	0	1

There are no bond length outliers.

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	8	ARG	NE-CZ-NH2	-13.93	113.34	120.30
1	A	8	ARG	NE-CZ-NH1	10.26	125.43	120.30
1	B	100	ASP	CB-CG-OD2	8.60	126.04	118.30
1	A	233	ARG	NE-CZ-NH1	8.39	124.49	120.30
1	A	160	ASP	CB-CG-OD2	6.72	124.35	118.30
1	A	82	THR	N-CA-CB	-6.68	97.61	110.30
1	B	160	ASP	CB-CG-OD2	6.64	124.28	118.30
1	A	188	ASP	CB-CG-OD2	6.60	124.24	118.30
1	A	152	ARG	NE-CZ-NH1	6.42	123.51	120.30
1	A	233	ARG	NE-CZ-NH2	-6.00	117.30	120.30
1	B	45	PRO	N-CA-CB	5.96	110.45	103.30
1	B	128	ASP	CB-CG-OD1	5.92	123.63	118.30
1	A	195	ARG	NE-CZ-NH1	5.61	123.11	120.30
1	A	18	THR	N-CA-CB	-5.33	100.17	110.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	152	ARG	NE-CZ-NH2	-5.31	117.65	120.30
1	A	8	ARG	CG-CD-NE	-5.22	100.84	111.80
1	A	45	PRO	N-CA-CB	5.16	109.49	103.30
1	A	115	LEU	CB-CG-CD2	5.13	119.73	111.00
1	A	137	SER	CB-CA-C	5.09	119.78	110.10
1	B	88	ARG	NE-CZ-NH1	-5.08	117.76	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	230	HIS	Peptide

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	1822	0	1826	24	0
1	B	1764	0	1763	46	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	165	0	0	8	0
3	B	108	0	0	14	0
All	All	3861	0	3589	68	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 9.

All (68) 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:66:ASN:HB3	3:B:404:HOH:O	1.55	1.05
1:A:18:THR:HG21	3:A:365:HOH:O	1.68	0.94
1:B:141:LEU:HD22	1:B:146:VAL:HG11	1.58	0.86
1:A:149:GLN:HE22	1:A:152:ARG:H	1.28	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:237:PRO:C	3:B:359:HOH:O	2.21	0.79
1:A:126:GLN:HE21	1:A:126:GLN:HA	1.53	0.74
1:A:229:VAL:HG11	3:B:370:HOH:O	1.87	0.73
1:A:82:THR:HG21	1:A:124:ALA:O	1.89	0.72
1:B:230:HIS:HB2	1:B:233:ARG:HG3	1.73	0.71
1:B:115:LEU:HD21	1:B:139:LYS:HB3	1.74	0.70
1:B:18:THR:HG22	1:B:22:ASP:H	1.56	0.70
1:A:229:VAL:HG21	3:B:370:HOH:O	1.92	0.68
1:B:146:VAL:HG13	1:B:147:ILE:N	2.11	0.66
1:A:82:THR:CG2	1:A:124:ALA:O	2.44	0.65
1:A:230:HIS:O	1:A:233:ARG:HD3	1.97	0.65
1:B:66:ASN:CB	3:B:404:HOH:O	2.27	0.65
1:B:229:VAL:O	1:B:230:HIS:CG	2.51	0.63
1:A:242:VAL:HG23	1:B:76:LEU:O	1.98	0.63
1:B:232:TRP:HZ3	3:B:328:HOH:O	1.82	0.61
1:A:152:ARG:HD3	3:A:445:HOH:O	1.99	0.61
1:B:66:ASN:ND2	3:B:404:HOH:O	2.34	0.59
1:B:146:VAL:CG1	1:B:147:ILE:N	2.67	0.57
1:A:82:THR:HB	3:A:307:HOH:O	2.04	0.57
1:B:230:HIS:HB2	1:B:233:ARG:CG	2.35	0.56
1:B:66:ASN:CG	3:B:404:HOH:O	2.44	0.56
1:B:52:LEU:HD13	1:B:53:PHE:CZ	2.43	0.54
1:A:17:VAL:O	1:B:127:PRO:HB3	2.07	0.54
1:B:219:GLU:HG3	1:B:233:ARG:HD2	1.88	0.54
1:B:51:ASN:HD22	1:B:52:LEU:H	1.56	0.54
1:B:59:GLN:NE2	1:B:73:GLN:HE21	2.05	0.53
1:B:136:MET:HE3	1:B:198:VAL:HB	1.91	0.52
1:A:115:LEU:HD22	1:A:136:MET:HE1	1.91	0.52
1:B:18:THR:HG22	1:B:22:ASP:N	2.24	0.52
1:B:149:GLN:HG2	1:B:149:GLN:O	2.10	0.52
1:B:139:LYS:HE2	1:B:196:ARG:O	2.10	0.52
1:B:136:MET:CE	1:B:198:VAL:HB	2.39	0.51
1:A:115:LEU:HD22	1:A:136:MET:CE	2.40	0.51
1:A:131:TRP:CE2	1:A:205:PRO:HA	2.47	0.50
1:B:13:ILE:HD13	1:B:28:LEU:HG	1.94	0.50
1:B:147:ILE:HB	3:B:400:HOH:O	2.12	0.50
1:A:195:ARG:HD3	3:A:420:HOH:O	2.12	0.49
1:B:134:LEU:HG	1:B:215:TRP:CZ2	2.48	0.49
1:B:35:LYS:HB3	3:B:392:HOH:O	2.14	0.47
1:B:147:ILE:CB	3:B:400:HOH:O	2.63	0.47
1:A:29:THR:HB	1:A:30:PRO:HD2	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:18:THR:HG23	1:B:20:ALA:H	1.81	0.46
1:B:109:SER:HA	3:B:390:HOH:O	2.15	0.46
1:B:52:LEU:HD13	1:B:53:PHE:CE1	2.51	0.45
1:B:67:ASP:OD1	1:B:67:ASP:N	2.50	0.45
1:B:229:VAL:O	1:B:230:HIS:ND1	2.50	0.45
1:B:13:ILE:HD13	1:B:28:LEU:CG	2.48	0.44
1:A:82:THR:HG22	1:A:124:ALA:HB1	1.99	0.44
1:B:229:VAL:C	1:B:230:HIS:CG	2.91	0.44
1:B:94:LEU:HD11	1:B:232:TRP:CZ3	2.53	0.43
1:A:39:ARG:CD	3:A:450:HOH:O	2.66	0.43
1:B:147:ILE:HG22	3:B:400:HOH:O	2.18	0.43
1:A:229:VAL:HG23	1:A:231:SER:H	1.84	0.43
1:A:119:SER:CA	1:A:136:MET:HE2	2.49	0.43
1:B:18:THR:HG21	3:B:391:HOH:O	2.17	0.42
1:B:63:GLY:H	1:B:66:ASN:HD21	1.66	0.42
1:A:39:ARG:HD2	3:A:450:HOH:O	2.20	0.42
1:B:139:LYS:CE	1:B:196:ARG:O	2.68	0.41
1:A:65:HIS:C	3:A:412:HOH:O	2.59	0.41
1:A:18:THR:HG23	3:A:424:HOH:O	2.20	0.41
1:B:131:TRP:CE2	1:B:205:PRO:HA	2.56	0.41
1:B:63:GLY:H	1:B:66:ASN:ND2	2.19	0.40
1:B:160:ASP:N	1:B:161:PRO:CD	2.85	0.40
1:B:47:SER:HB3	1:B:74:ALA:O	2.21	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	238/244 (98%)	228 (96%)	9 (4%)	1 (0%)	39	33
1	B	226/244 (93%)	221 (98%)	4 (2%)	1 (0%)	39	33

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	464/488 (95%)	449 (97%)	13 (3%)	2 (0%)	39	33

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	229	VAL
1	A	242	VAL

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	187/195 (96%)	177 (95%)	10 (5%)	28	22
1	B	181/195 (93%)	167 (92%)	14 (8%)	16	10
All	All	368/390 (94%)	344 (94%)	24 (6%)	21	15

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	SER
1	A	18	THR
1	A	70	SER
1	A	82	THR
1	A	115	LEU
1	A	126	GLN
1	A	149	GLN
1	A	195	ARG
1	A	196	ARG
1	A	203	GLU
1	B	1	MET
1	B	4	ARG
1	B	18	THR
1	B	47	SER
1	B	51	ASN
1	B	52	LEU

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Mol	Chain	Res	Type
1	B	66	ASN
1	B	147	ILE
1	B	149	GLN
1	B	187	LEU
1	B	196	ARG
1	B	198	VAL
1	B	216	ARG
1	B	232	TRP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	93	HIS
1	A	126	GLN
1	A	149	GLN
1	B	51	ASN
1	B	59	GLN
1	B	66	ASN
1	B	149	GLN

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 ⓘ

Of 2 ligands modelled in this entry, 2 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, 95th 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 ²)	Q<0.9
1	A	242/244 (99%)	0.21	19 (7%) 15 16	20, 32, 59, 89	0
1	B	232/244 (95%)	0.24	14 (6%) 25 27	21, 35, 57, 66	0
All	All	474/488 (97%)	0.23	33 (6%) 19 21	20, 33, 59, 89	0

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	244	SER	14.5
1	A	243	LEU	10.5
1	B	228	GLY	8.3
1	A	242	VAL	8.2
1	B	227	GLY	5.9
1	B	229	VAL	5.6
1	B	146	VAL	5.5
1	A	44	GLY	5.2
1	A	240	VAL	5.2
1	B	145	GLY	5.2
1	B	46	LEU	5.0
1	B	45	PRO	4.6
1	A	42	VAL	4.6
1	B	40	GLY	4.0
1	A	43	LYS	3.8
1	A	45	PRO	3.8
1	B	67	ASP	3.4
1	A	2	ARG	3.3
1	B	47	SER	3.0
1	B	2	ARG	2.9
1	A	64	PRO	2.8
1	A	241	PRO	2.8
1	A	3	SER	2.7
1	B	204	GLN	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	232	TRP	2.5
1	A	46	LEU	2.5
1	A	67	ASP	2.4
1	A	48	SER	2.4
1	A	39	ARG	2.2
1	A	40	GLY	2.2
1	A	41	GLY	2.1
1	B	1	MET	2.1
1	A	47	SER	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 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, 95th 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(\AA^2)	Q<0.9
2	ZN	B	300	1/1	0.97	0.04	-2.28	39,39,39,39	0
2	ZN	A	300	1/1	1.00	0.05	-3.15	32,32,32,32	0

6.5 Other polymers [i](#)

There are no such residues in this entry.