



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 06:15 pm GMT

PDB ID : 2WMO  
Title : STRUCTURE OF THE COMPLEX BETWEEN DOCK9 AND CDC42.  
Authors : Yang, J.; Roe, S.M.; Barford, D.  
Deposited on : 2009-07-02  
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

<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  
Mogul : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : 1.9-1692  
EDS : trunk28620  
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) : recalc28949

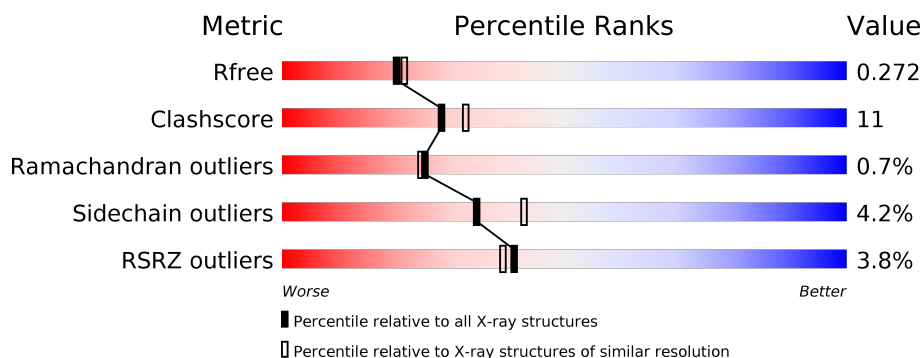
# 1 Overall quality at a glance

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))
$R_{free}$	100719	4002 (2.20-2.20)
Clashscore	112137	4730 (2.20-2.20)
Ramachandran outliers	110173	4656 (2.20-2.20)
Sidechain outliers	110143	4657 (2.20-2.20)
RSRZ outliers	101464	4033 (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. 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	428	<div> <div>5%</div> <div> <div></div> <div>69%</div> <div>24%</div> <div>• 6%</div> </div> </div>
2	B	190	<div> <div>%</div> <div> <div></div> <div>69%</div> <div>22%</div> <div>• 6%</div> </div> </div>

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 4865 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.

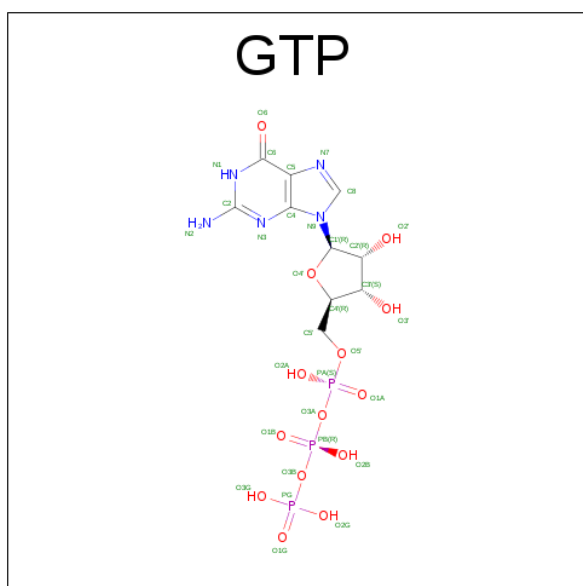
- Molecule 1 is a protein called DEDICATOR OF CYTOKINESIS PROTEIN 9.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	403	Total	C	N	O	S	0	0	0
			3185	2029	546	590	20			

- Molecule 2 is a protein called CELL DIVISION CONTROL PROTEIN 42 HOMOLOG.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	178	Total	C	N	O	S	0	0	0
			1339	863	208	261	7			

- Molecule 3 is GUANOSINE-5'-TRIPHOSPHATE (three-letter code: GTP) (formula:  $C_{10}H_{16}N_5O_{14}P_3$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	B	1	Total	C	N	O	P	0	0
			32	10	5	14	3		

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total 1	Mg 1	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	235	Total 235	O 235	0	0
5	B	73	Total 73	O 73	0	0



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	93.06Å 88.59Å 89.98Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.99 – 2.20 44.99 – 2.20	Depositor EDS
% Data completeness (in resolution range)	98.5 (44.99-2.20) 99.6 (44.99-2.20)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.70 (at 2.20Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, $R_{free}$	0.216 , 0.272 0.210 , 0.272	Depositor DCC
$R_{free}$ test set	1909 reflections (4.99%)	DCC
Wilson B-factor (Å <sup>2</sup> )	38.3	Xtriage
Anisotropy	0.485	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 58.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.018 for l,-k,h 0.022 for -k,-h,l 0.017 for -h,l,k 0.007 for k,l,h 0.007 for l,h,k	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	4865	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.47% 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: GTP, MG

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	0.40	1/3246 (0.0%)	0.52	0/4383
2	B	0.32	0/1368	0.51	0/1874
All	All	0.38	1/4614 (0.0%)	0.52	0/6257

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	94	CYS	CB-SG	-5.14	1.73	1.81

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	A	3185	0	3063	69	0
2	B	1339	0	1301	36	0
3	B	32	0	12	3	0
4	B	1	0	0	0	0
5	A	235	0	0	10	0
5	B	73	0	0	0	0
All	All	4865	0	4376	99	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 11.

All (99) 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)
2:B:98:VAL:HG11	2:B:149:LEU:HD13	1.54	0.86
1:A:268:GLN:HA	1:A:274:GLN:HE22	1.41	0.85
1:A:7:PRO:HB3	1:A:40:LEU:HD11	1.59	0.83
1:A:215:PRO:HA	1:A:218:LEU:HD13	1.62	0.81
2:B:138:THR:OG1	2:B:141:THR:HG22	1.83	0.79
1:A:335:LEU:HD11	1:A:389:ALA:HB1	1.72	0.71
1:A:369:ASP:O	1:A:372:VAL:HG23	1.90	0.71
1:A:320:GLU:O	1:A:324:LEU:HD13	1.90	0.70
1:A:418:GLU:O	1:A:422:ILE:HG13	1.92	0.69
1:A:184:LEU:H	2:B:26:ASN:ND2	1.90	0.69
1:A:401:GLN:HB3	5:A:2231:HOH:O	1.92	0.68
1:A:388:GLN:O	1:A:392:VAL:HG23	1.96	0.65
2:B:90:PHE:CE1	2:B:145:LEU:HD12	2.32	0.64
1:A:60:VAL:HG22	5:A:2046:HOH:O	1.97	0.63
2:B:86:SER:O	2:B:89:SER:HB2	1.99	0.62
2:B:16:LYS:HD3	3:B:1178:GTP:O2B	2.00	0.61
1:A:409:LYS:HD3	1:A:413:ARG:NH1	2.16	0.60
2:B:93:VAL:O	2:B:98:VAL:HG13	2.04	0.58
1:A:154:THR:HG23	1:A:180:LYS:O	2.03	0.58
2:B:110:PHE:CE1	2:B:151:ALA:HB2	2.38	0.57
2:B:93:VAL:HG11	2:B:112:LEU:HD11	1.86	0.57
2:B:11:ASP:O	2:B:14:VAL:HG13	2.04	0.57
1:A:106:LEU:HD11	1:A:250:HIS:CE1	2.38	0.57
1:A:372:VAL:HA	5:A:2226:HOH:O	2.04	0.57
2:B:90:PHE:CZ	2:B:145:LEU:HD12	2.40	0.57
1:A:358:PHE:O	1:A:359:LEU:HD23	2.04	0.56
1:A:171:GLY:HA3	5:A:2138:HOH:O	2.05	0.56
1:A:368:PRO:O	1:A:369:ASP:CB	2.54	0.55
1:A:368:PRO:O	1:A:369:ASP:HB3	2.07	0.55
1:A:209:ASP:OD1	1:A:267:ARG:NH1	2.38	0.55
1:A:423:MET:O	1:A:425:GLU:HG2	2.07	0.54
2:B:6:CYS:SG	2:B:79:LEU:HD22	2.48	0.54
1:A:58:PHE:HA	5:A:2046:HOH:O	2.07	0.53
1:A:305:ASN:O	1:A:309:VAL:HG23	2.09	0.53
1:A:409:LYS:HD3	1:A:413:ARG:HH12	1.74	0.53
1:A:161:PHE:HB2	1:A:224:TYR:HB2	1.89	0.53
2:B:68:ARG:HB3	2:B:69:PRO:HD3	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:393:ASN:HD22	1:A:405:GLN:NE2	2.07	0.52
2:B:82:PHE:HB2	2:B:89:SER:HB3	1.91	0.52
1:A:285:HIS:HB2	1:A:294:ILE:HD13	1.91	0.52
1:A:15:ASP:O	1:A:19:ARG:HG3	2.10	0.52
1:A:205:LYS:HB3	5:A:2155:HOH:O	2.09	0.51
1:A:101:ALA:O	1:A:102:GLU:HB2	2.10	0.51
2:B:98:VAL:HA	2:B:101:ILE:HG12	1.93	0.51
1:A:222:TYR:HB2	1:A:224:TYR:CE2	2.46	0.50
1:A:320:GLU:OE1	1:A:338:LYS:HE3	2.12	0.50
1:A:274:GLN:NE2	2:B:34:PRO:HG2	2.26	0.50
1:A:337:LEU:HG	2:B:39:ASN:HB3	1.94	0.49
1:A:269:GLY:H	1:A:274:GLN:NE2	2.10	0.49
1:A:352:LEU:O	1:A:356:ARG:HG3	2.13	0.49
2:B:79:LEU:HG	2:B:113:VAL:HG21	1.95	0.49
1:A:314:MET:O	1:A:318:VAL:HG22	2.12	0.49
1:A:280:ILE:HD12	1:A:280:ILE:N	2.28	0.48
1:A:331:ASP:OD2	2:B:3:THR:OG1	2.26	0.47
1:A:213:VAL:CG2	1:A:218:LEU:HD11	2.44	0.47
1:A:301:HIS:HE1	5:A:2207:HOH:O	1.96	0.47
1:A:210:SER:HB2	1:A:259:MET:SD	2.55	0.47
1:A:351:PRO:HG3	2:B:37:PHE:CE1	2.49	0.47
1:A:371:LYS:HA	5:A:2224:HOH:O	2.16	0.46
2:B:83:SER:HB3	2:B:86:SER:HB3	1.98	0.46
1:A:65:ILE:HA	1:A:65:ILE:HD12	1.82	0.46
2:B:77:VAL:HB	2:B:176:ALA:HB2	1.98	0.46
1:A:379:PHE:O	1:A:382:PHE:HB3	2.16	0.45
1:A:103:ARG:CZ	1:A:288:PRO:HG3	2.47	0.45
2:B:98:VAL:O	2:B:102:THR:HG23	2.16	0.45
1:A:236:GLU:HA	1:A:239:LEU:HD12	1.99	0.45
2:B:25:THR:O	2:B:26:ASN:HB2	2.17	0.44
2:B:101:ILE:HG13	2:B:102:THR:N	2.33	0.44
1:A:130:LEU:HD12	1:A:131:TYR:N	2.32	0.44
1:A:218:LEU:HD12	1:A:218:LEU:N	2.33	0.44
1:A:415:MET:HG3	1:A:416:ALA:N	2.32	0.43
1:A:318:VAL:HG12	1:A:382:PHE:HA	2.00	0.43
1:A:284:ILE:CG2	1:A:297:MET:HB2	2.49	0.43
2:B:1:MET:SD	2:B:50:PRO:HB2	2.59	0.43
1:A:213:VAL:HG21	1:A:218:LEU:HD11	2.01	0.43
1:A:141:VAL:O	1:A:145:MET:HG3	2.18	0.43
1:A:390:LEU:HD11	1:A:412:TYR:CG	2.54	0.43
1:A:172:LYS:HA	5:A:2141:HOH:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:182:THR:HG23	1:A:186:GLU:HB3	2.02	0.42
1:A:181:LEU:HD12	2:B:45:MET:HG2	2.02	0.42
1:A:235:ASP:HA	5:A:2169:HOH:O	2.19	0.42
1:A:256:MET:HG2	1:A:257:PHE:N	2.35	0.42
2:B:49:GLU:HA	2:B:50:PRO:HD3	1.91	0.42
2:B:116:GLN:HG2	3:B:1178:GTP:C6	2.55	0.41
2:B:70:LEU:O	2:B:73:PRO:HD2	2.20	0.41
1:A:315:SER:HA	1:A:318:VAL:CG2	2.51	0.41
1:A:182:THR:HA	1:A:183:PRO:HD3	1.91	0.41
1:A:157:ARG:HD2	1:A:157:ARG:C	2.41	0.41
1:A:168:ASP:OD1	1:A:168:ASP:N	2.52	0.41
2:B:8:VAL:HG21	2:B:20:LEU:HD11	2.03	0.41
2:B:116:GLN:HG2	3:B:1178:GTP:C5	2.56	0.41
1:A:277:ARG:HB2	1:A:277:ARG:HE	1.50	0.41
1:A:191:LEU:O	1:A:195:TYR:HD1	2.03	0.41
1:A:237:LYS:O	1:A:241:GLU:HG3	2.21	0.40
2:B:108:THR:HA	2:B:109:PRO:HD3	1.92	0.40
2:B:138:THR:H	2:B:141:THR:CG2	2.34	0.40
2:B:98:VAL:CG1	2:B:149:LEU:HD13	2.38	0.40
2:B:161:THR:OG1	2:B:163:LYS:HG3	2.21	0.40
1:A:376:LYS:O	1:A:380:ARG:HG3	2.21	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	A	395/428 (92%)	363 (92%)	28 (7%)	4 (1%)	18	16
2	B	176/190 (93%)	170 (97%)	6 (3%)	0	100	100
All	All	571/618 (92%)	533 (93%)	34 (6%)	4 (1%)	25	24

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	369	ASP
1	A	55	CYS
1	A	266	LYS
1	A	372	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	326/381 (86%)	314 (96%)	12 (4%)	39	49
2	B	145/170 (85%)	137 (94%)	8 (6%)	25	29
All	All	471/551 (86%)	451 (96%)	20 (4%)	34	43

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	THR
1	A	16	SER
1	A	53	GLN
1	A	202	GLU
1	A	213	VAL
1	A	243	LYS
1	A	256	MET
1	A	268	GLN
1	A	292	LYS
1	A	304	LEU
1	A	318	VAL
1	A	344	SER
2	B	16	LYS
2	B	52	THR
2	B	63	ASP
2	B	79	LEU
2	B	89	SER
2	B	98	VAL
2	B	141	THR

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Mol	Chain	Res	Type
2	B	145	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	53	GLN
1	A	263	GLN
1	A	274	GLN
1	A	285	HIS
1	A	300	HIS
1	A	405	GLN
1	A	411	ASN
2	B	2	GLN
2	B	26	ASN
2	B	132	ASN
2	B	162	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 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 for Mogul analysis.

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	GTP	B	1178	4	27,34,34	0.96	1 (3%)	27,54,54	1.84	5 (18%)

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	GTP	B	1178	4	-	0/18/38/38	0/3/3/3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1178	GTP	C6-N1	3.04	1.38	1.33

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1178	GTP	N3-C2-N1	-5.36	119.63	127.46
3	B	1178	GTP	C5-C6-N1	-3.11	119.06	123.48
3	B	1178	GTP	N2-C2-N1	2.06	120.53	117.24
3	B	1178	GTP	C6-N1-C2	3.24	120.72	116.06
3	B	1178	GTP	C2-N3-C4	4.26	120.13	115.16

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1178	GTP	3	0

## 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	403/428 (94%)	0.24	20 (4%) 30 28	20, 44, 73, 86	0
2	B	178/190 (93%)	-0.02	2 (1%) 80 79	36, 51, 62, 66	0
All	All	581/618 (94%)	0.16	22 (3%) 41 39	20, 47, 69, 86	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	327	SER	4.2
1	A	378	VAL	3.9
1	A	328	ALA	3.5
1	A	222	TYR	3.4
1	A	204	VAL	3.3
1	A	423	MET	3.1
1	A	360	ASP	2.8
1	A	422	ILE	2.8
1	A	419	LEU	2.7
1	A	355	ALA	2.6
1	A	424	HIS	2.6
1	A	265	GLY	2.5
1	A	354	TYR	2.4
1	A	374	LEU	2.4
1	A	215	PRO	2.3
2	B	90	PHE	2.2
1	A	394	GLU	2.2
1	A	356	ARG	2.2
1	A	166	PHE	2.1
1	A	420	SER	2.1
1	A	164	GLY	2.1
2	B	74	GLN	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. 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( $\text{\AA}^2$ )	Q<0.9
3	GTP	B	1178	32/32	0.95	0.13	0.23	38,42,47,49	0
4	MG	B	1179	1/1	0.92	0.14	-	46,46,46,46	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.