



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 09:34 am GMT

PDB ID : 1Z5A  
Title : Topoisomerase VI-B, ADP-bound dimer form  
Authors : Corbett, K.D.; Berger, J.M.  
Deposited on : 2005-03-17  
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

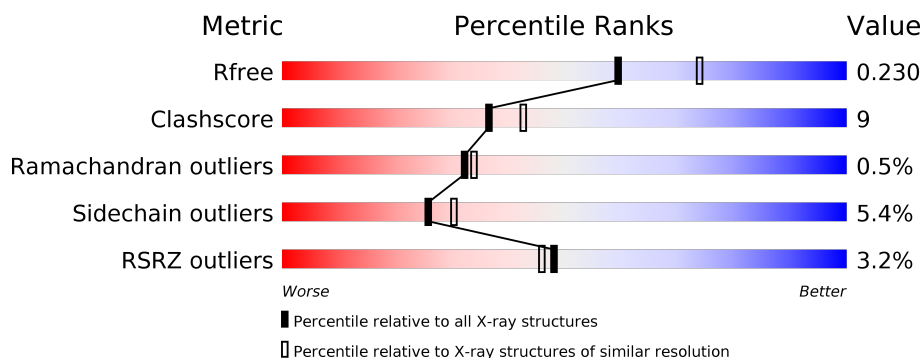
# 1 Overall quality at a glance

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	469	<div> <div>2%</div> <div> <div></div> <div>76%</div> <div>22%</div> <div>..</div> </div> </div>
1	B	469	<div> <div>4%</div> <div> <div></div> <div>79%</div> <div>16%</div> <div>..</div> </div> </div>

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 criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	MG	A	1001	-	-	-	X

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7716 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 Type II DNA topoisomerase VI subunit B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	466	Total	C	N	O	S	0	0	0
			3748	2410	630	702	6			
1	B	460	Total	C	N	O	S	0	0	0
			3704	2383	619	696	6			

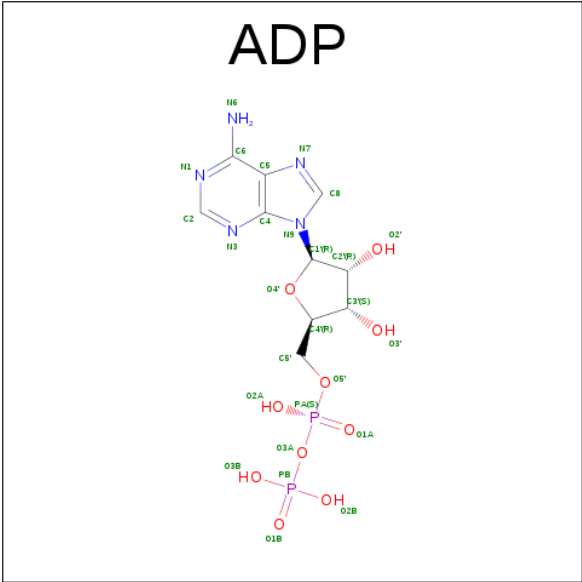
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	303	TYR	ASP	SEE REMARK 999	UNP O05207
A	435	ASP	ASN	SEE REMARK 999	UNP O05207
B	303	TYR	ASP	SEE REMARK 999	UNP O05207
B	435	ASP	ASN	SEE REMARK 999	UNP O05207

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 3 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: C<sub>10</sub>H<sub>15</sub>N<sub>5</sub>O<sub>10</sub>P<sub>2</sub>).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	B	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

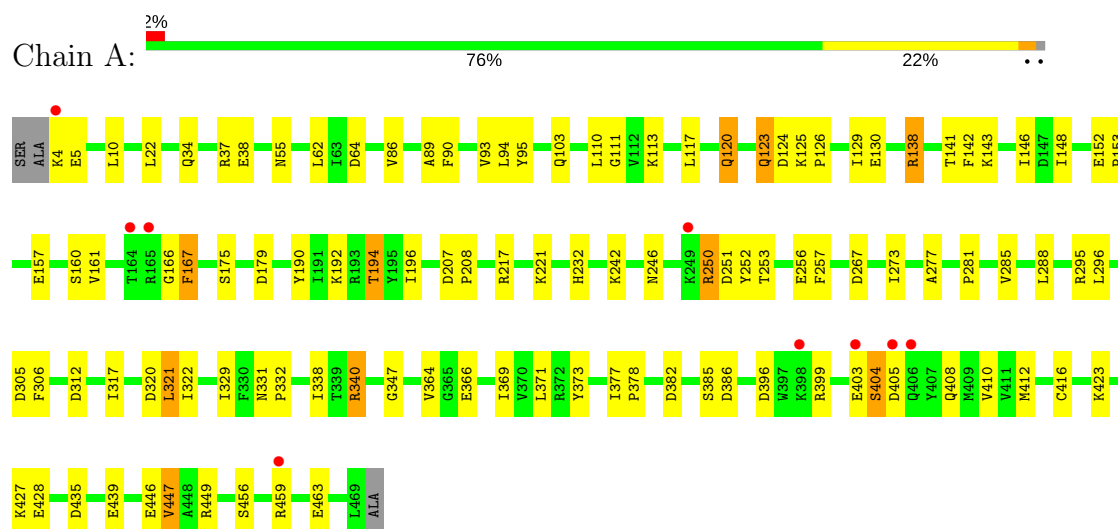
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	87	Total	O	0	0
			87	87		
4	B	121	Total	O	0	0
			121	121		

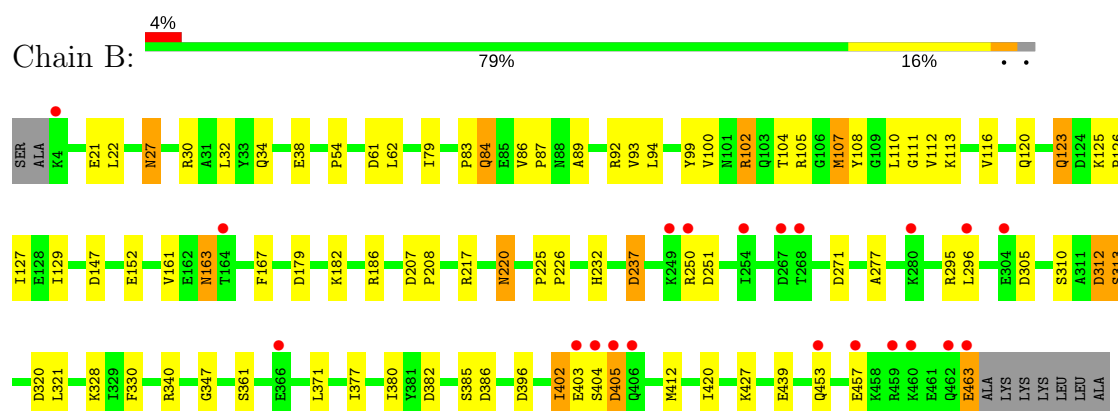
### 3 Residue-property plots

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: Type II DNA topoisomerase VI subunit B



#### • Molecule 1: Type II DNA topoisomerase VI subunit B



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 32 1 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	74.57Å 74.57Å 345.56Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.20 19.94 – 2.20	Depositor EDS
% Data completeness (in resolution range)	93.2 (20.00-2.20) 93.2 (19.94-2.20)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.08	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	5.86 (at 2.19Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, $R_{free}$	0.195 , 0.231 0.197 , 0.230	Depositor DCC
$R_{free}$ test set	2633 reflections (5.26%)	DCC
Wilson B-factor (Å <sup>2</sup> )	27.1	Xtriage
Anisotropy	0.044	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.40 , 47.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.057 for -h,-k,l	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	7716	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.84% 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

### 5.1 Standard geometry

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

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	0.64	0/3829	0.72	9/5178 (0.2%)
1	B	0.66	0/3785	0.72	11/5121 (0.2%)
All	All	0.65	0/7614	0.72	20/10299 (0.2%)

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( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	A	64	ASP	CB-CG-OD2	6.53	124.18	118.30
1	A	386	ASP	CB-CG-OD2	5.71	123.44	118.30
1	A	382	ASP	CB-CG-OD2	5.70	123.43	118.30
1	B	251	ASP	CB-CG-OD2	5.57	123.31	118.30
1	B	61	ASP	CB-CG-OD2	5.55	123.29	118.30
1	A	305	ASP	CB-CG-OD2	5.53	123.27	118.30
1	A	179	ASP	CB-CG-OD2	5.47	123.22	118.30
1	A	320	ASP	CB-CG-OD2	5.46	123.21	118.30
1	B	405	ASP	CB-CG-OD2	5.34	123.10	118.30
1	B	312	ASP	CB-CG-OD2	5.28	123.06	118.30
1	B	320	ASP	CB-CG-OD2	5.26	123.04	118.30
1	A	405	ASP	CB-CG-OD2	5.23	123.01	118.30
1	B	147	ASP	CB-CG-OD2	5.21	122.99	118.30
1	B	271	ASP	CB-CG-OD2	5.18	122.96	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	435	ASP	CB-CG-OD2	5.13	122.91	118.30
1	B	305	ASP	CB-CG-OD2	5.13	122.91	118.30
1	B	237	ASP	CB-CG-OD2	5.11	122.89	118.30
1	A	340	ARG	NE-CZ-NH2	-5.05	117.78	120.30
1	B	396	ASP	CB-CG-OD2	5.05	122.84	118.30
1	B	386	ASP	CB-CG-OD2	5.01	122.81	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	313	SER	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	3748	0	3814	83	0
1	B	3704	0	3766	58	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	27	0	12	0	0
3	B	27	0	12	1	0
4	A	87	0	0	2	0
4	B	121	0	0	5	0
All	All	7716	0	7604	136	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 (136) 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:412:MET:HG2	4:B:2109:HOH:O	1.39	1.17
1:A:4:LYS:HA	4:A:2028:HOH:O	1.51	1.07

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:4:LYS:HG3	1:A:5:GLU:H	1.01	1.06
1:A:4:LYS:CG	1:A:5:GLU:H	1.74	0.99
1:A:4:LYS:HG3	1:A:5:GLU:N	1.77	0.97
1:A:120:GLN:HE22	1:A:146:ILE:N	1.67	0.91
1:B:403:GLU:HG2	1:B:404:SER:H	1.37	0.88
1:B:463:GLU:HA	4:B:2100:HOH:O	1.74	0.86
1:A:120:GLN:HE22	1:A:146:ILE:H	0.88	0.86
1:A:250:ARG:HE	1:A:252:TYR:HE1	1.23	0.86
1:B:277:ALA:O	1:B:295:ARG:NH1	2.10	0.84
1:A:120:GLN:NE2	1:A:146:ILE:H	1.73	0.83
1:A:253:THR:HG23	1:A:256:GLU:H	1.43	0.83
1:A:4:LYS:CG	1:A:5:GLU:N	2.36	0.82
1:B:163:ASN:HD21	1:B:167:PHE:H	1.28	0.81
1:A:123:GLN:HG2	1:A:124:ASP:H	1.47	0.79
1:B:186:ARG:HH11	1:B:380:ILE:CG2	2.00	0.74
1:A:123:GLN:HG2	1:A:124:ASP:N	2.02	0.73
1:A:403:GLU:HG2	1:A:404:SER:OG	1.86	0.73
1:B:340:ARG:HD3	1:B:439:GLU:OE2	1.90	0.72
1:A:403:GLU:HB3	1:A:459:ARG:NH2	2.04	0.72
1:A:403:GLU:HB3	1:A:459:ARG:HH21	1.58	0.69
1:A:340:ARG:HD3	1:A:439:GLU:OE2	1.92	0.68
1:A:123:GLN:NE2	1:A:175:SER:HB2	2.08	0.68
1:B:403:GLU:CG	1:B:404:SER:H	2.03	0.68
1:A:322:ILE:HG12	1:A:412:MET:HE1	1.77	0.67
1:A:89:ALA:O	1:A:113:LYS:HE3	1.95	0.66
1:A:55:ASN:ND2	1:A:217:ARG:HH22	1.94	0.65
1:B:403:GLU:HG2	1:B:404:SER:N	2.11	0.65
1:A:403:GLU:HG2	1:A:404:SER:N	2.12	0.64
1:B:102:ARG:HD3	1:B:104:THR:CG2	2.28	0.63
1:A:207:ASP:HB2	1:A:208:PRO:CD	2.29	0.62
1:A:253:THR:OG1	1:A:281:PRO:O	2.17	0.62
1:B:107:MET:HE1	1:B:420:ILE:H	1.64	0.62
1:A:385:SER:HB3	1:B:385:SER:HB3	1.84	0.60
1:A:90:PHE:CD2	1:A:129:ILE:HD13	2.37	0.60
1:B:207:ASP:HB2	1:B:208:PRO:CD	2.31	0.60
1:A:123:GLN:NE2	1:A:125:LYS:HB2	2.17	0.59
1:B:27:ASN:ND2	1:B:30:ARG:H	2.00	0.59
1:B:123:GLN:NE2	1:B:125:LYS:H	2.01	0.59
1:A:322:ILE:HG12	1:A:412:MET:CE	2.34	0.57
1:A:4:LYS:HE3	1:B:79:ILE:HD12	1.84	0.57
1:B:123:GLN:HE22	1:B:125:LYS:H	1.52	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:84:GLN:NE2	1:B:84:GLN:H	2.04	0.55
1:A:403:GLU:HG2	1:A:404:SER:H	1.69	0.55
1:A:130:GLU:HG2	1:A:141:THR:HG22	1.89	0.54
1:A:446:GLU:HG3	1:A:449:ARG:NH2	2.23	0.54
1:B:463:GLU:CA	4:B:2100:HOH:O	2.44	0.54
1:A:221:LYS:O	1:A:221:LYS:HD3	2.08	0.53
1:B:89:ALA:O	1:B:113:LYS:HE2	2.09	0.53
1:B:186:ARG:HH11	1:B:380:ILE:HG23	1.72	0.53
1:A:55:ASN:HD21	1:A:217:ARG:HH22	1.55	0.52
1:A:117:LEU:CD2	1:A:148:ILE:CD1	2.88	0.52
1:A:253:THR:HG22	1:A:256:GLU:CD	2.30	0.52
1:A:37:ARG:NH1	1:A:428:GLU:OE2	2.43	0.52
1:B:89:ALA:O	1:B:113:LYS:CE	2.58	0.52
1:A:190:TYR:O	1:A:194:THR:CG2	2.59	0.51
1:A:190:TYR:O	1:A:194:THR:HG23	2.11	0.51
1:A:253:THR:HG22	1:A:256:GLU:CG	2.41	0.51
1:B:163:ASN:ND2	1:B:167:PHE:H	2.05	0.50
1:B:127:ILE:HD12	1:B:129:ILE:HD11	1.94	0.49
1:A:10:LEU:HG	1:B:99:TYR:CE2	2.47	0.49
1:A:364:VAL:HG22	1:A:408:GLN:HG2	1.94	0.49
1:B:123:GLN:HE22	1:B:125:LYS:HB2	1.77	0.49
1:B:105:ARG:HB3	1:B:377:ILE:HD11	1.95	0.49
1:B:453:GLN:O	1:B:457:GLU:HB2	2.11	0.49
1:A:143:LYS:HB2	1:A:157:GLU:HB2	1.95	0.48
1:A:138:ARG:HD3	1:A:160:SER:OG	2.13	0.48
1:A:192:LYS:O	1:A:196:ILE:HG12	2.13	0.48
1:A:322:ILE:HG23	1:A:412:MET:CE	2.44	0.48
1:B:310:SER:OG	1:B:312:ASP:OD1	2.30	0.48
1:B:92:ARG:O	1:B:113:LYS:HD3	2.14	0.47
1:B:54:PRO:O	1:B:217:ARG:NH2	2.48	0.47
1:A:89:ALA:O	1:A:113:LYS:CE	2.61	0.47
1:B:83:PRO:HD2	1:B:84:GLN:NE2	2.30	0.47
1:A:123:GLN:CG	1:A:124:ASP:N	2.77	0.46
1:A:331:ASN:N	1:A:332:PRO:HD3	2.30	0.46
1:B:102:ARG:HD3	1:B:104:THR:HG23	1.97	0.46
1:B:179:ASP:OD2	1:B:182:LYS:HE3	2.16	0.46
1:A:166:GLY:O	1:A:167:PHE:C	2.55	0.45
1:A:166:GLY:O	1:A:167:PHE:O	2.34	0.45
1:B:21:GLU:H	1:B:21:GLU:CD	2.19	0.45
1:B:220:ASN:C	1:B:220:ASN:HD22	2.20	0.45
1:A:103:GLN:OE1	1:A:416:CYS:HB2	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:207:ASP:HB2	1:A:208:PRO:HD2	1.98	0.45
1:A:232:HIS:O	1:A:347:GLY:HA2	2.16	0.45
1:B:186:ARG:NH1	1:B:380:ILE:CG2	2.76	0.45
1:A:369:ILE:HB	1:A:410:VAL:HG22	1.98	0.45
1:B:186:ARG:NH1	1:B:380:ILE:HG23	2.31	0.45
1:A:257:PHE:CG	1:A:285:VAL:HG21	2.52	0.45
1:B:116:VAL:HG22	1:B:127:ILE:HG21	1.99	0.44
1:A:340:ARG:CD	1:A:439:GLU:OE2	2.64	0.44
1:B:330:PHE:O	1:B:361:SER:HB2	2.18	0.44
1:A:117:LEU:HD22	1:A:148:ILE:HD12	1.99	0.44
1:B:27:ASN:HD21	1:B:30:ARG:H	1.64	0.44
1:A:152:GLU:HB2	1:A:153:PRO:HD2	1.99	0.44
1:A:338:ILE:HG22	1:A:447:VAL:HG13	1.99	0.44
1:A:86:VAL:HG11	1:A:142:PHE:CD2	2.52	0.44
1:B:403:GLU:CG	1:B:404:SER:N	2.76	0.44
1:B:207:ASP:HB2	1:B:208:PRO:HD2	2.00	0.44
1:B:225:PRO:HA	1:B:226:PRO:HD3	1.91	0.44
1:A:93:VAL:HG12	1:A:94:LEU:HG	1.99	0.43
1:B:220:ASN:C	1:B:220:ASN:ND2	2.71	0.43
1:A:364:VAL:HG22	1:A:408:GLN:CG	2.48	0.43
1:B:84:GLN:H	1:B:84:GLN:HE21	1.65	0.43
1:B:237:ASP:C	1:B:237:ASP:OD1	2.57	0.43
1:B:186:ARG:HH11	1:B:380:ILE:HG22	1.78	0.43
1:A:120:GLN:HG2	4:A:2022:HOH:O	2.18	0.43
1:B:34:GLN:O	1:B:38:GLU:HG2	2.19	0.43
1:A:221:LYS:HD3	1:A:221:LYS:C	2.40	0.42
1:B:108:TYR:HB2	4:B:2003:HOH:O	2.18	0.42
1:A:196:ILE:CD1	1:A:329:ILE:HD11	2.49	0.42
1:A:396:ASP:HB3	1:A:399:ARG:HD3	2.02	0.42
1:A:4:LYS:CE	1:B:79:ILE:HD12	2.47	0.42
1:A:95:TYR:O	1:A:113:LYS:NZ	2.53	0.42
1:B:112:VAL:HG12	3:B:2002:ADP:O1A	2.20	0.42
1:A:277:ALA:O	1:A:295:ARG:NH1	2.50	0.42
1:A:322:ILE:HG23	1:A:412:MET:HE1	2.01	0.42
1:A:207:ASP:CB	1:A:208:PRO:CD	2.95	0.41
1:A:242:LYS:O	1:A:246:ASN:ND2	2.53	0.41
1:A:273:ILE:HD11	1:A:306:PHE:CE1	2.55	0.41
1:B:93:VAL:HG12	1:B:94:LEU:HG	2.01	0.41
1:A:194:THR:HG22	1:A:378:PRO:HG2	2.02	0.41
1:B:120:GLN:HG3	4:B:2051:HOH:O	2.19	0.41
1:B:232:HIS:O	1:B:347:GLY:HA2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:125:LYS:HA	1:A:126:PRO:HD3	1.93	0.41
1:A:288:LEU:HA	1:A:288:LEU:HD23	1.87	0.41
1:B:125:LYS:HA	1:B:126:PRO:HD3	1.88	0.41
1:B:86:VAL:HB	1:B:87:PRO:HD3	2.03	0.41
1:A:373:TYR:HA	1:A:377:ILE:O	2.21	0.41
1:A:423:LYS:NZ	1:B:382:ASP:OD1	2.54	0.41
1:A:317:ILE:HG22	1:A:321:LEU:HB3	2.03	0.41
1:A:117:LEU:CD2	1:A:148:ILE:HD12	2.51	0.40
1:A:34:GLN:O	1:A:38:GLU:HG2	2.21	0.40
1:A:253:THR:CG2	1:A:256:GLU:H	2.22	0.40
1:A:90:PHE:CD2	1:A:129:ILE:CD1	3.05	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	464/469 (99%)	450 (97%)	12 (3%)	2 (0%)	38	41
1	B	458/469 (98%)	445 (97%)	10 (2%)	3 (1%)	25	24
All	All	922/938 (98%)	895 (97%)	22 (2%)	5 (0%)	32	34

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	167	PHE
1	A	111	GLY
1	B	111	GLY
1	B	107	MET
1	B	402	ILE

### 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	412/416 (99%)	391 (95%)	21 (5%)	28	33
1	B	409/416 (98%)	386 (94%)	23 (6%)	25	29
All	All	821/832 (99%)	777 (95%)	44 (5%)	26	30

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	LEU
1	A	62	LEU
1	A	110	LEU
1	A	120	GLN
1	A	123	GLN
1	A	138	ARG
1	A	161	VAL
1	A	194	THR
1	A	250	ARG
1	A	251	ASP
1	A	267	ASP
1	A	296	LEU
1	A	312	ASP
1	A	321	LEU
1	A	366	GLU
1	A	371	LEU
1	A	404	SER
1	A	427	LYS
1	A	447	VAL
1	A	456	SER
1	A	463	GLU
1	B	22	LEU
1	B	27	ASN
1	B	32	LEU
1	B	62	LEU
1	B	84	GLN
1	B	100	VAL

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Mol	Chain	Res	Type
1	B	102	ARG
1	B	110	LEU
1	B	123	GLN
1	B	152	GLU
1	B	161	VAL
1	B	163	ASN
1	B	220	ASN
1	B	250	ARG
1	B	296	LEU
1	B	313	SER
1	B	321	LEU
1	B	328	LYS
1	B	371	LEU
1	B	402	ILE
1	B	405	ASP
1	B	427	LYS
1	B	463	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	55	ASN
1	A	68	GLN
1	A	120	GLN
1	A	123	GLN
1	A	246	ASN
1	A	247	ASN
1	A	408	GLN
1	B	27	ASN
1	B	84	GLN
1	B	101	ASN
1	B	123	GLN
1	B	163	ASN
1	B	211	ASN
1	B	220	ASN
1	B	247	ASN
1	B	408	GLN

### 5.3.3 RNA ⓘ

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, 2 are monoatomic - leaving 2 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	ADP	A	2001	2	25,29,29	1.19	2 (8%)	24,45,45	1.63	2 (8%)
3	ADP	B	2002	2	25,29,29	1.04	1 (4%)	24,45,45	1.46	2 (8%)

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	ADP	A	2001	2	-	0/12/32/32	0/3/3/3
3	ADP	B	2002	2	-	0/12/32/32	0/3/3/3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	2001	ADP	C2-N3	2.53	1.36	1.32
3	B	2002	ADP	C5-C4	2.89	1.47	1.40
3	A	2001	ADP	C5-C4	3.40	1.48	1.40

All (4) bond angle outliers are listed below:



Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
3	A	2001	ADP	N3-C2-N1	-6.30	123.37	128.86
3	B	2002	ADP	N3-C2-N1	-5.56	124.02	128.86
3	A	2001	ADP	C4-C5-N7	-2.64	106.86	109.41
3	B	2002	ADP	C4-C5-N7	-2.55	106.95	109.41

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	2002	ADP	1	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	466/469 (99%)	-0.27	9 (1%) 67 65	12, 23, 42, 54	0
1	B	460/469 (98%)	-0.12	21 (4%) 33 32	10, 22, 46, 78	0
All	All	926/938 (98%)	-0.19	30 (3%) 48 46	10, 22, 45, 78	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	463	GLU	6.1
1	B	405	ASP	4.5
1	B	403	GLU	4.1
1	B	404	SER	3.7
1	A	4	LYS	3.7
1	A	405	ASP	3.5
1	B	406	GLN	3.5
1	B	267	ASP	3.4
1	A	459	ARG	3.3
1	B	366	GLU	3.2
1	A	249	LYS	3.2
1	B	457	GLU	3.2
1	B	250	ARG	3.1
1	B	462	GLN	3.0
1	B	268	THR	2.8
1	B	304	GLU	2.7
1	B	254	ILE	2.6
1	B	4	LYS	2.5
1	B	164	THR	2.4
1	B	296	LEU	2.4
1	B	453	GLN	2.3
1	A	406	GLN	2.3
1	A	164	THR	2.3
1	B	460	LYS	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	165	ARG	2.2
1	B	280	LYS	2.2
1	B	459	ARG	2.1
1	B	249	LYS	2.0
1	A	403	GLU	2.0
1	A	398	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. 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(Å <sup>2</sup> )	Q<0.9
2	MG	A	1001	1/1	0.88	0.19	4.60	63,63,63,63	0
3	ADP	B	2002	27/27	0.98	0.15	1.76	32,34,36,37	0
3	ADP	A	2001	27/27	0.96	0.14	1.43	35,40,41,42	0
2	MG	B	1002	1/1	0.89	0.12	0.75	54,54,54,54	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.