



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 27, 2014 – 12:05 AM GMT

PDB ID : 2VHG  
Title : CRYSTAL STRUCTURE OF THE ISHP608 TRANSPOSASE IN COMPLEX  
WITH RIGHT END 31-MER DNA  
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Deposited on : 2007-11-21  
Resolution : 2.90 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

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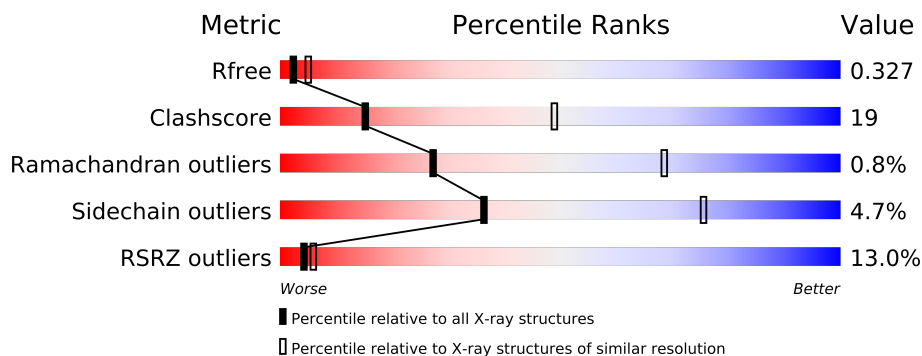
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 2.90 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))
$R_{free}$	66092	1053 (2.90-2.90)
Clashscore	79885	1326 (2.90-2.90)
Ramachandran outliers	78287	1290 (2.90-2.90)
Sidechain outliers	78261	1292 (2.90-2.90)
RSRZ outliers	66119	1054 (2.90-2.90)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	159	
1	B	159	
2	C	31	
2	D	31	

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3032 atoms, of which 0 are hydrogen and 0 are deuterium.

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 TRANSPOSASE ORFA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	126	Total	C	N	O	S	0	0	0
			1032	664	183	179	6			
1	B	127	Total	C	N	O	S	0	0	0
			1041	669	185	181	6			

- Molecule 2 is a DNA chain called RIGHT END 31-MER.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	23	Total	C	N	O	P	0	0	0
			469	224	85	138	22			
2	D	24	Total	C	N	O	P	0	0	0
			489	234	87	145	23			

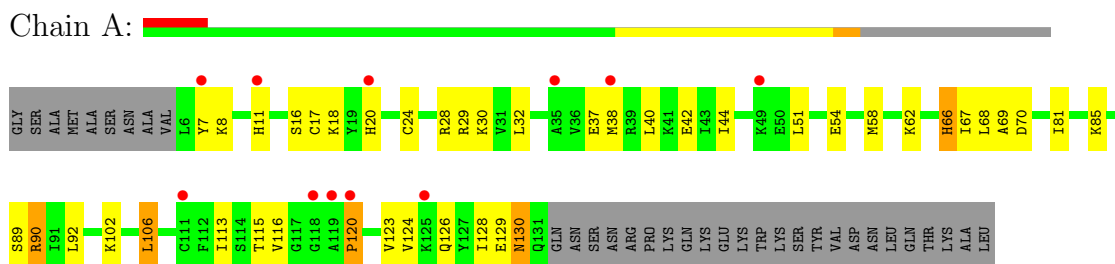
- Molecule 3 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mn	0	0
			1	1		

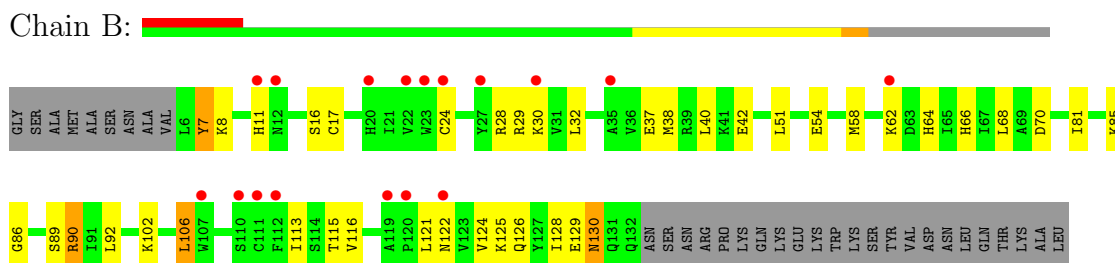
### 3 Residue-property plots

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

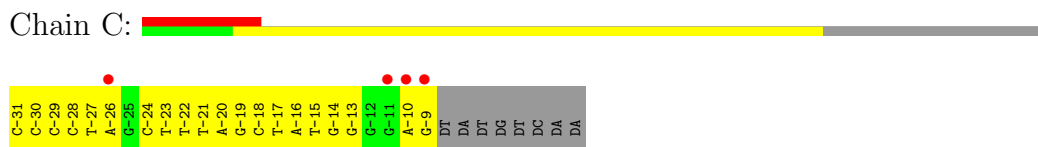
- Molecule 1: TRANSPOSASE ORFA



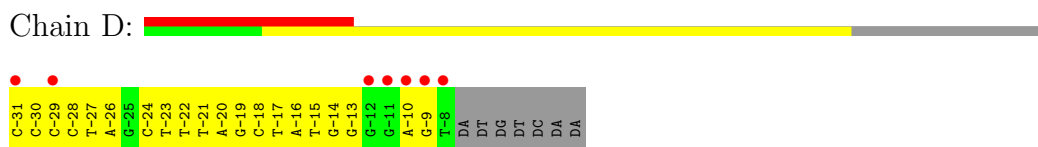
- Molecule 1: TRANSPOSASE ORFA



- Molecule 2: RIGHT END 31-MER



- Molecule 2: RIGHT END 31-MER



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	66.47Å 90.92Å 93.42Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 2.90 29.60 – 2.90	Depositor EDS
% Data completeness (in resolution range)	90.3 (40.00-2.90) 90.3 (29.60-2.90)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	6.54 (at 2.90Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.285 , 0.320 0.303 , 0.327	Depositor DCC
$R_{free}$ test set	843 reflections (7.15%)	DCC
Wilson B-factor (Å <sup>2</sup> )	75.8	Xtriage
Anisotropy	0.355	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.27 , 39.4	EDS
Estimated twinning fraction	0.016 for -h,l,k	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 11787 reflections	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	3032	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	86.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.45% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

## 5 Model quality

### 5.1 Standard geometry

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

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	0.35	0/1053	0.62	0/1419
1	B	0.35	0/1062	0.60	0/1431
2	C	0.35	0/525	0.66	0/809
2	D	0.33	0/547	0.66	0/843
All	All	0.35	0/3187	0.63	0/4502

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
2	C	0	1
2	D	0	1
All	All	0	3

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	7	TYR	Sidechain
2	C	-24	DC	Sidechain
2	D	-24	DC	Sidechain

## 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1032	0	1072	30	0
1	B	1041	0	1080	33	0
2	C	469	0	261	31	0
2	D	489	0	273	28	0
3	A	1	0	0	0	0
All	All	3032	0	2686	108	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 19.

All (108) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:C:-15:DT:H3'	2:C:-14:DG:H5''	1.29	1.15
2:C:-10:DA:H3'	2:C:-9:DG:H5''	1.12	1.12
2:D:-15:DT:H3'	2:D:-14:DG:H5''	1.29	1.09
2:C:-22:DT:H2''	2:C:-21:DT:H5'	1.35	1.06
2:D:-22:DT:H2''	2:D:-21:DT:H5'	1.37	1.03
2:C:-10:DA:H3'	2:C:-9:DG:C5'	1.93	0.99
2:D:-15:DT:H3'	2:D:-14:DG:C5'	1.92	0.97
2:C:-15:DT:H3'	2:C:-14:DG:C5'	1.94	0.96
1:A:81:ILE:HD11	1:A:113:ILE:HD11	1.56	0.88
1:B:81:ILE:HD11	1:B:113:ILE:HD11	1.55	0.88
2:C:-10:DA:H2'	2:C:-9:DG:C8	2.17	0.80
2:C:-10:DA:C3'	2:C:-9:DG:H5''	2.07	0.77
2:C:-14:DG:H2''	2:C:-13:DG:H5'	1.67	0.75
2:D:-14:DG:H2''	2:D:-13:DG:H5'	1.69	0.75
1:A:66:HIS:HE1	1:B:124:VAL:HG11	1.53	0.74
2:D:-30:DC:H2''	2:D:-29:DC:C5	2.25	0.71
1:A:66:HIS:CE1	1:B:124:VAL:HG11	2.30	0.66
1:A:7:TYR:CE2	1:A:123:VAL:HG11	2.32	0.64
2:D:-22:DT:H2''	2:D:-21:DT:C5'	2.22	0.64
1:B:125:LYS:HA	1:B:128:ILE:HD12	1.79	0.63
1:B:54:GLU:HB3	1:B:70:ASP:HB3	1.81	0.62
1:A:89:SER:HA	1:A:106:LEU:HD12	1.80	0.62

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:32:LEU:HD22	1:B:40:LEU:HD22	1.81	0.62
2:C:-22:DT:H2"	2:C:-21:DT:C5'	2.23	0.61
1:A:32:LEU:HD22	1:A:40:LEU:HD22	1.83	0.61
1:A:30:LYS:HB3	1:A:62:LYS:HB3	1.83	0.61
2:C:-30:DC:H2"	2:C:-29:DC:C5	2.35	0.61
1:A:54:GLU:HB3	1:A:70:ASP:HB3	1.83	0.60
1:B:24:CYS:SG	1:B:64:HIS:HB3	2.43	0.59
2:C:-15:DT:C3'	2:C:-14:DG:H5"	2.20	0.58
2:D:-30:DC:H2"	2:D:-29:DC:C6	2.38	0.58
1:B:30:LYS:HB3	1:B:62:LYS:HB3	1.87	0.57
2:D:-18:DC:C6	2:D:-17:DT:H72	2.40	0.57
1:B:89:SER:HA	1:B:106:LEU:HD12	1.86	0.56
2:D:-15:DT:C3'	2:D:-14:DG:H5"	2.21	0.55
2:C:-31:DC:H2"	2:C:-30:DC:C5	2.41	0.55
2:C:-31:DC:H2"	2:C:-30:DC:C6	2.41	0.55
1:B:7:TYR:CD1	1:B:16:SER:HB2	2.42	0.55
1:A:81:ILE:CD1	1:A:113:ILE:HD11	2.33	0.54
1:B:121:LEU:O	1:B:122:ASN:HB3	2.06	0.54
1:A:28:ARG:NH1	1:B:129:GLU:HA	2.23	0.54
1:B:124:VAL:HG12	1:B:128:ILE:HD11	1.91	0.53
2:C:-16:DA:H2"	2:C:-14:DG:H5'	1.92	0.52
2:D:-16:DA:H2"	2:D:-14:DG:H5'	1.91	0.52
2:C:-15:DT:C3'	2:C:-14:DG:C5'	2.80	0.52
1:A:20:HIS:CE1	1:A:68:LEU:HD13	2.44	0.52
2:D:-23:DT:H2"	2:D:-22:DT:H5"	1.90	0.52
2:C:-27:DT:H2"	2:C:-26:DA:H8	1.75	0.52
1:B:92:LEU:HD12	1:B:106:LEU:HD11	1.92	0.51
2:C:-10:DA:H2"	2:C:-9:DG:C5	2.46	0.51
1:B:81:ILE:CD1	1:B:113:ILE:HD11	2.36	0.51
1:A:68:LEU:HD11	1:A:116:VAL:HG11	1.93	0.51
2:C:-27:DT:H2"	2:C:-26:DA:C8	2.45	0.50
1:A:92:LEU:HD12	1:A:106:LEU:HD11	1.94	0.50
2:C:-29:DC:C2	2:C:-28:DC:C5	3.00	0.50
2:C:-30:DC:H2"	2:C:-29:DC:C6	2.47	0.50
1:B:7:TYR:CE1	1:B:16:SER:HB2	2.46	0.49
2:C:-23:DT:H2"	2:C:-22:DT:H5"	1.93	0.49
1:B:124:VAL:O	1:B:128:ILE:HD12	2.13	0.49
1:B:68:LEU:HD11	1:B:116:VAL:HG11	1.95	0.49
1:A:51:LEU:HA	2:D:-15:DT:O4	2.14	0.48
2:D:-15:DT:C3'	2:D:-14:DG:C5'	2.79	0.48
2:D:-18:DC:H2"	2:D:-17:DT:OP2	2.14	0.48
2:C:-18:DC:C6	2:C:-17:DT:H72	2.48	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:C:-10:DA:H2"	2:C:-9:DG:C4	2.49	0.48
2:D:-29:DC:C2	2:D:-28:DC:C5	3.01	0.48
1:A:17:CYS:O	1:A:70:ASP:HA	2.13	0.48
1:A:29:ARG:HH11	1:A:102:LYS:HE2	1.79	0.48
2:C:-14:DG:H2"	2:C:-13:DG:C5'	2.42	0.47
2:D:-27:DT:H2"	2:D:-26:DA:C8	2.50	0.47
2:D:-10:DA:H2"	2:D:-9:DG:O5'	2.14	0.47
2:D:-14:DG:H2"	2:D:-13:DG:C5'	2.44	0.46
1:A:7:TYR:OH	1:A:120:PRO:HG3	2.14	0.46
1:A:38:MET:O	1:A:42:GLU:HG3	2.15	0.46
1:B:89:SER:OG	2:D:-19:DG:H4'	2.15	0.46
1:B:38:MET:O	1:B:42:GLU:HG3	2.15	0.46
1:B:37:GLU:HG3	1:B:38:MET:N	2.31	0.46
2:C:-22:DT:O2	2:C:-20:DA:H8	1.98	0.46
1:B:51:LEU:HA	2:C:-15:DT:O4	2.16	0.45
1:B:7:TYR:CD2	1:B:7:TYR:N	2.84	0.45
2:C:-21:DT:H4'	2:C:-20:DA:O4'	2.16	0.45
2:D:-21:DT:H4'	2:D:-20:DA:O4'	2.17	0.45
1:A:90:ARG:NH1	2:C:-20:DA:H4'	2.32	0.45
1:B:90:ARG:NH1	2:D:-20:DA:H4'	2.32	0.45
1:A:7:TYR:CD1	1:A:16:SER:HB2	2.52	0.44
1:A:17:CYS:HA	1:B:115:THR:OG1	2.17	0.44
1:B:29:ARG:HH11	1:B:102:LYS:HE2	1.82	0.44
1:B:58:MET:HA	1:B:66:HIS:O	2.18	0.44
2:C:-19:DG:C4	2:C:-18:DC:C5	3.06	0.44
2:D:-27:DT:H2"	2:D:-26:DA:H8	1.82	0.44
2:D:-22:DT:O2	2:D:-20:DA:H8	2.01	0.43
1:A:37:GLU:HG3	1:A:38:MET:N	2.33	0.43
1:B:17:CYS:O	1:B:70:ASP:HA	2.19	0.43
2:D:-31:DC:N4	2:D:-10:DA:H61	2.16	0.43
1:B:86:GLY:HA2	2:D:-18:DC:H5'	2.01	0.43
1:A:85:LYS:HB3	2:C:-18:DC:H5"	2.00	0.42
1:A:126:GLN:O	1:A:130:ASN:HB2	2.20	0.42
1:A:44:ILE:HD13	1:A:67:ILE:HD11	2.02	0.42
1:A:124:VAL:O	1:A:128:ILE:HG13	2.20	0.42
2:C:-18:DC:H2"	2:C:-17:DT:OP2	2.19	0.41
1:A:129:GLU:HA	1:B:28:ARG:NH1	2.36	0.41
1:A:115:THR:OG1	1:B:17:CYS:HA	2.20	0.41
2:D:-19:DG:C4	2:D:-18:DC:C5	3.08	0.41
1:A:58:MET:HA	1:A:66:HIS:O	2.20	0.41
1:B:85:LYS:HB3	2:D:-18:DC:H5"	2.02	0.41
2:D:-31:DC:H42	2:D:-10:DA:H61	1.69	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:126:GLN:O	1:B:130:ASN:HB2	2.21	0.40
1:A:18:LYS:HA	1:A:69:ALA:O	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	124/159 (78%)	111 (90%)	12 (10%)	1 (1%)	27	68
1	B	125/159 (79%)	112 (90%)	12 (10%)	1 (1%)	27	68
All	All	249/318 (78%)	223 (90%)	24 (10%)	2 (1%)	27	68

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	11	HIS
1	B	11	HIS

### 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	116/144 (81%)	109 (94%)	7 (6%)	27	63
1	B	117/144 (81%)	113 (97%)	4 (3%)	49	86
All	All	233/288 (81%)	222 (95%)	11 (5%)	36	75

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	LYS
1	A	24	CYS
1	A	66	HIS
1	A	90	ARG
1	A	106	LEU
1	A	120	PRO
1	A	130	ASN
1	B	8	LYS
1	B	90	ARG
1	B	106	LEU
1	B	130	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	66	HIS
1	A	94	GLN
1	B	94	GLN

### 5.3.3 RNA ⓘ

There are no RNA chains 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 1 ligands modelled in this entry, 1 is 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.

## 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	A	126/159 (79%)	0.77	11 (8%) 10 13	40, 79, 116, 131	0
1	B	127/159 (79%)	0.88	17 (13%) 4 5	40, 84, 108, 118	0
2	C	23/31 (74%)	0.68	4 (17%) 2 3	55, 73, 141, 156	0
2	D	24/31 (77%)	1.35	7 (29%) 1 1	51, 87, 164, 169	0
All	All	300/380 (78%)	0.86	39 (13%) 4 6	40, 81, 129, 169	0

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	-8	DT	6.7
1	B	27	TYR	6.3
2	D	-10	DA	5.7
1	B	11	HIS	5.6
1	A	118	GLY	4.6
1	A	7	TYR	4.3
1	B	20	HIS	4.1
1	A	120	PRO	4.0
2	C	-9	DG	3.9
1	B	35	ALA	3.8
1	A	38	MET	3.7
1	A	119	ALA	3.5
1	A	20	HIS	3.4
1	B	120	PRO	3.3
2	D	-9	DG	3.2
2	D	-12	DG	3.2
2	C	-11	DG	3.2
1	A	11	HIS	3.2
1	B	122	ASN	3.1
1	A	35	ALA	3.1
1	B	119	ALA	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	24	CYS	2.8
1	B	23	TRP	2.7
1	B	62	LYS	2.7
2	D	-31	DC	2.6
2	C	-10	DA	2.6
1	A	49	LYS	2.6
1	A	125	LYS	2.6
1	B	22	VAL	2.6
1	B	112	PHE	2.5
1	A	111	CYS	2.5
1	B	12	ASN	2.4
1	B	110	SER	2.4
1	B	30	LYS	2.4
2	D	-11	DG	2.4
1	B	111	CYS	2.4
1	B	107	TRP	2.3
2	D	-29	DC	2.2
2	C	-26	DA	2.1

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

## 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands ⓘ

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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
3	MN	A	1132	1/1	0.13	-2.12	97,97,97,97	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.