



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 09:49 PM GMT

PDB ID : 4WDX  
Title : 17beta-HSD5 in complex with [4-(2-hydroxyethyl)piperidin-1-yl](5-methyl-1H-indol-2-yl)methanone  
Authors : Amano, Y.; Yamaguchi, T.  
Deposited on : 2014-09-09  
Resolution : 1.64 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 (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

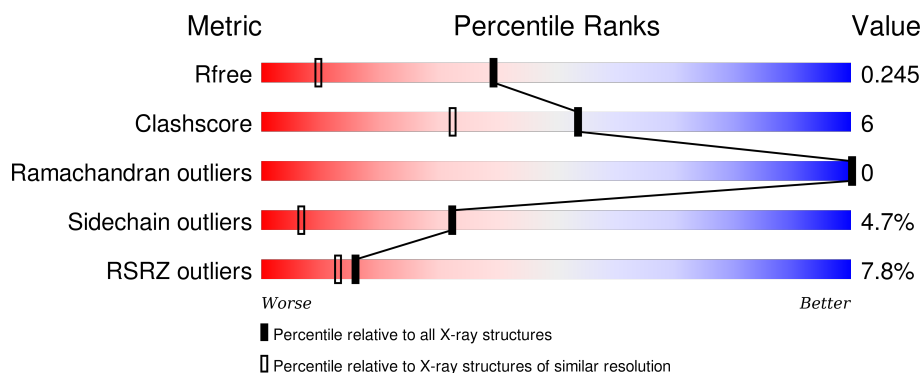
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.64 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	1953 (1.66-1.62)
Clashscore	102246	2091 (1.66-1.62)
Ramachandran outliers	100387	2052 (1.66-1.62)
Sidechain outliers	100360	2052 (1.66-1.62)
RSRZ outliers	91569	1955 (1.66-1.62)

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	331	
1	B	331	

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 5218 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 Aldo-keto reductase family 1 member C3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	298	Total	C	N	O	S	0	0	0
			2388	1526	418	432	12			
1	B	302	Total	C	N	O	S	0	0	0
			2416	1543	423	438	12			

There are 18 discrepancies between the modelled and reference sequences:

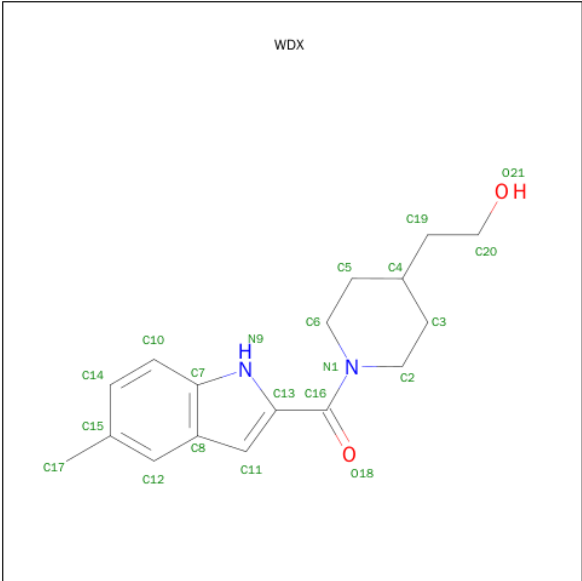
Chain	Residue	Modelled	Actual	Comment	Reference
A	5	GLN	HIS	natural variant	UNP P42330
A	324	LEU	-	expression tag	UNP P42330
A	325	GLU	-	expression tag	UNP P42330
A	326	HIS	-	expression tag	UNP P42330
A	327	HIS	-	expression tag	UNP P42330
A	328	HIS	-	expression tag	UNP P42330
A	329	HIS	-	expression tag	UNP P42330
A	330	HIS	-	expression tag	UNP P42330
A	331	HIS	-	expression tag	UNP P42330
B	5	GLN	HIS	natural variant	UNP P42330
B	324	LEU	-	expression tag	UNP P42330
B	325	GLU	-	expression tag	UNP P42330
B	326	HIS	-	expression tag	UNP P42330
B	327	HIS	-	expression tag	UNP P42330
B	328	HIS	-	expression tag	UNP P42330
B	329	HIS	-	expression tag	UNP P42330
B	330	HIS	-	expression tag	UNP P42330
B	331	HIS	-	expression tag	UNP P42330

- Molecule 2 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NAP) (formula: C<sub>21</sub>H<sub>28</sub>N<sub>7</sub>O<sub>17</sub>P<sub>3</sub>).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	B	1	Total	C	N	O	P	0	0
			48	21	7	17	3		

- Molecule 3 is [4-(2-hydroxyethyl)piperidin-1-yl](5-methyl-1H-indol-2-yl)methanone (three-letter code: WDX) (formula: C<sub>17</sub>H<sub>22</sub>N<sub>2</sub>O<sub>2</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			21	17	2	2		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	C	N	O	0	0
			21	17	2	2		

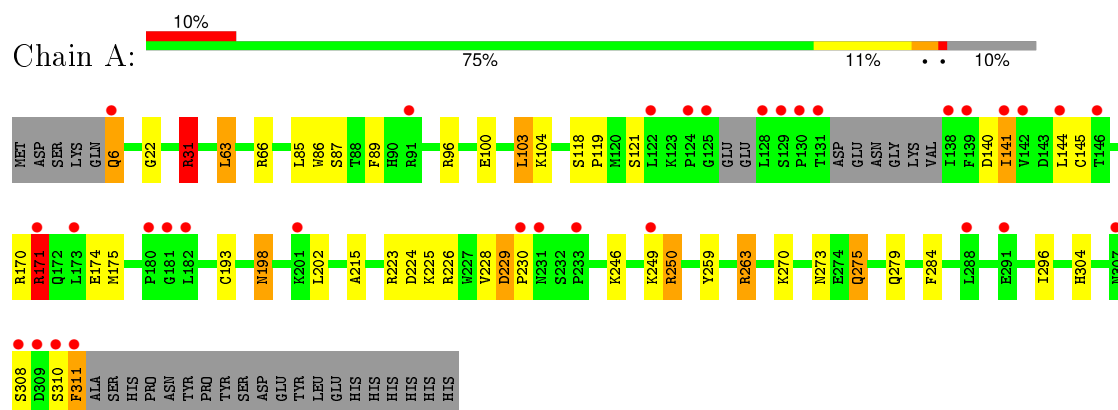
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	122	Total	O	0	0
			122	122		
4	B	154	Total	O	0	0
			154	154		

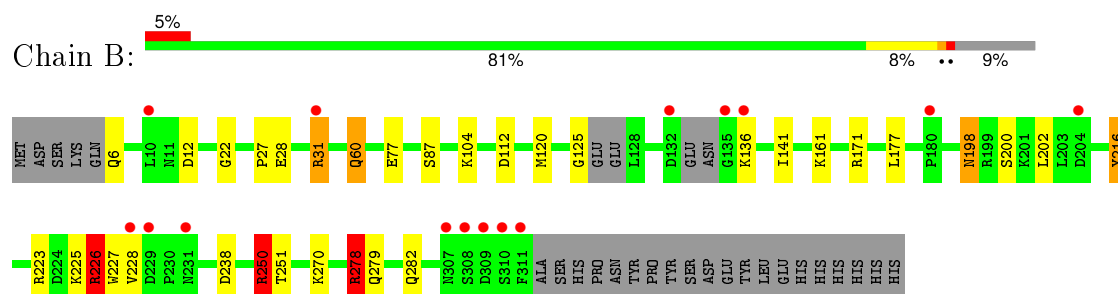
### 3 Residue-property plots [i](#)

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: Aldo-keto reductase family 1 member C3



- Molecule 1: Aldo-keto reductase family 1 member C3



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	60.04Å 77.23Å 72.21Å 90.00° 97.62° 90.00°	Depositor
Resolution (Å)	14.27 – 1.64 14.27 – 1.64	Depositor EDS
% Data completeness (in resolution range)	95.6 (14.27-1.64) 95.8 (14.27-1.64)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.41 (at 1.64Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, $R_{free}$	0.214 , 0.249 0.211 , 0.245	Depositor DCC
$R_{free}$ test set	3855 reflections (5.30%)	DCC
Wilson B-factor (Å <sup>2</sup> )	26.3	Xtriage
Anisotropy	0.021	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.50 , 56.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.35$	Xtriage
Outliers	1 of 76526 reflections (0.001%)	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	5218	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.71% 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, 0.375 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: WDX, NAP

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	1.06	0/2439	1.14	10/3295 (0.3%)
1	B	1.09	1/2467 (0.0%)	1.18	17/3332 (0.5%)
All	All	1.07	1/4906 (0.0%)	1.16	27/6627 (0.4%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	216	TYR	CG-CD2	5.38	1.46	1.39

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	66	ARG	NE-CZ-NH1	-18.73	110.94	120.30
1	B	250	ARG	NE-CZ-NH2	16.35	128.48	120.30
1	A	66	ARG	NE-CZ-NH2	13.35	126.98	120.30
1	A	250	ARG	NE-CZ-NH2	9.73	125.17	120.30
1	B	226	ARG	CG-CD-NE	9.61	131.97	111.80
1	B	278	ARG	NE-CZ-NH2	9.51	125.05	120.30
1	A	103	LEU	CB-CG-CD2	8.30	125.11	111.00
1	A	171	ARG	NE-CZ-NH1	8.17	124.39	120.30
1	B	250	ARG	NE-CZ-NH1	-7.96	116.32	120.30
1	B	31	ARG	NE-CZ-NH1	-7.89	116.36	120.30
1	B	177	LEU	CB-CG-CD1	-7.48	98.28	111.00
1	B	31	ARG	NE-CZ-NH2	7.26	123.93	120.30
1	B	226	ARG	CB-CA-C	7.24	124.88	110.40
1	A	31	ARG	NE-CZ-NH1	-6.75	116.93	120.30
1	B	226	ARG	CA-CB-CG	6.62	127.96	113.40
1	B	238	ASP	CB-CG-OD2	-6.40	112.54	118.30
1	B	12	ASP	CB-CG-OD2	-6.38	112.56	118.30
1	A	63	LEU	CA-CB-CG	6.35	129.90	115.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	112	ASP	CB-CG-OD1	-6.34	112.60	118.30
1	B	278	ARG	NE-CZ-NH1	-6.30	117.15	120.30
1	A	250	ARG	NE-CZ-NH1	-6.29	117.16	120.30
1	B	226	ARG	NE-CZ-NH1	5.89	123.25	120.30
1	A	229	ASP	CB-CG-OD2	-5.88	113.01	118.30
1	B	278	ARG	CD-NE-CZ	5.59	131.42	123.60
1	A	31	ARG	NE-CZ-NH2	5.45	123.03	120.30
1	B	250	ARG	CB-CG-CD	5.43	125.72	111.60
1	B	28	GLU	OE1-CD-OE2	-5.36	116.86	123.30

There are no chirality outliers.

There are no planarity outliers.

## 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	2388	0	2416	34	0
1	B	2416	0	2445	29	0
2	A	48	0	25	3	0
2	B	48	0	25	4	0
3	A	21	0	22	1	0
3	B	21	0	22	1	0
4	A	122	0	0	2	0
4	B	154	0	0	10	0
All	All	5218	0	4955	64	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 6.

All (64) 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:A:311:PHE:O	1:A:311:PHE:HD2	1.22	1.19
1:B:125:GLY:C	4:B:612:HOH:O	1.98	1.01
1:A:311:PHE:O	1:A:311:PHE:CD2	2.14	0.99

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:77:GLU:HG2	4:B:586:HOH:O	1.61	0.99
1:A:31:ARG:NH1	4:A:622:HOH:O	2.00	0.94
1:A:119:PRO:HA	1:A:144:LEU:HD21	1.59	0.83
1:B:104:LYS:CE	4:B:523:HOH:O	2.29	0.79
1:A:6:GLN:HG3	1:A:284:PHE:CE1	2.17	0.79
1:B:31:ARG:HD2	1:B:31:ARG:N	2.01	0.75
1:B:104:LYS:HE2	4:B:523:HOH:O	1.84	0.75
1:A:311:PHE:C	1:A:311:PHE:HD2	1.92	0.73
1:A:279:GLN:NE2	2:A:401:NAP:H62A	1.86	0.73
1:B:31:ARG:NH1	4:B:639:HOH:O	2.24	0.69
1:A:145:CYS:SG	1:A:175:MET:HG2	2.31	0.69
1:B:278:ARG:HH21	1:B:278:ARG:HG2	1.57	0.67
1:A:198:ASN:HD22	1:A:198:ASN:C	1.98	0.67
1:B:250:ARG:HH11	1:B:282:GLN:HE21	1.43	0.67
1:A:311:PHE:C	1:A:311:PHE:CD2	2.66	0.66
1:A:224:ASP:OD2	1:A:226:ARG:HD3	1.96	0.66
1:A:171:ARG:HD2	1:A:171:ARG:C	2.17	0.65
1:A:259:TYR:O	1:A:263:ARG:HG2	1.99	0.63
1:B:279:GLN:NE2	2:B:401:NAP:H62A	1.98	0.62
1:A:273:ASN:OD1	1:A:275:GLN:HG3	2.00	0.61
1:B:104:LYS:HE3	4:B:523:HOH:O	1.99	0.59
1:A:170:ARG:O	1:A:174:GLU:HG3	2.03	0.58
1:B:198:ASN:C	1:B:198:ASN:HD22	2.06	0.57
1:A:121:SER:HB2	1:A:140:ASP:O	2.05	0.56
1:A:171:ARG:HH11	1:A:175:MET:HB2	1.71	0.55
1:A:63:LEU:HD12	1:B:141:ILE:HD12	1.89	0.54
1:B:250:ARG:HD3	4:B:552:HOH:O	2.09	0.53
1:A:171:ARG:HD2	1:A:171:ARG:O	2.08	0.53
1:B:279:GLN:O	1:B:282:GLN:HB2	2.09	0.52
1:B:223:ARG:HA	1:B:228:VAL:HG21	1.93	0.51
1:B:270:LYS:O	2:B:401:NAP:H8A	2.11	0.50
3:A:402:WDX:H10	3:A:402:WDX:C11	2.41	0.50
1:A:270:LYS:O	2:A:401:NAP:H8A	2.11	0.50
1:B:60:GLN:NE2	1:B:60:GLN:H	2.11	0.49
1:B:6:GLN:N	4:B:643:HOH:O	2.45	0.48
1:B:198:ASN:HD22	1:B:200:SER:H	1.61	0.48
1:A:246:LYS:HZ2	1:A:246:LYS:HB3	1.80	0.47
1:B:251:THR:OG1	1:B:279:GLN:NE2	2.48	0.46
1:B:31:ARG:N	1:B:31:ARG:CD	2.66	0.46
1:A:304:HIS:N	1:A:304:HIS:CD2	2.83	0.46
1:A:250:ARG:HD3	4:A:582:HOH:O	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:198:ASN:C	1:A:198:ASN:ND2	2.69	0.44
2:B:401:NAP:C5N	3:B:402:WDX:H8	2.47	0.44
1:A:223:ARG:HA	1:A:228:VAL:HG21	2.00	0.44
1:B:198:ASN:ND2	1:B:200:SER:H	2.15	0.43
1:B:226:ARG:HG2	1:B:227:TRP:CE2	2.53	0.43
1:A:259:TYR:HB2	1:A:296:ILE:HD13	2.00	0.43
1:A:193:CYS:HB3	1:A:215:ALA:CB	2.48	0.43
1:A:141:ILE:H	1:A:141:ILE:HG12	1.61	0.42
1:A:100:GLU:O	1:A:104:LYS:HG3	2.18	0.42
1:A:85:LEU:HD11	1:A:89:PHE:HB2	2.01	0.42
1:B:171:ARG:HD3	1:B:171:ARG:HA	1.90	0.42
1:B:27:PRO:HG2	1:B:226:ARG:NH1	2.35	0.42
1:B:22:GLY:HA3	2:B:401:NAP:H4D	2.02	0.41
1:A:86:TRP:CG	1:A:87:SER:N	2.88	0.41
1:A:22:GLY:HA3	2:A:401:NAP:H4D	2.02	0.41
1:A:229:ASP:HA	1:A:230:PRO:HD2	1.92	0.41
1:B:6:GLN:N	4:B:501:HOH:O	2.53	0.41
1:A:6:GLN:HG3	1:A:284:PHE:CZ	2.55	0.41
1:B:161:LYS:HE2	4:B:645:HOH:O	2.21	0.40
1:B:87:SER:HB3	1:B:120:MET:O	2.20	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	292/331 (88%)	279 (96%)	13 (4%)	0	100	100
1	B	296/331 (89%)	290 (98%)	6 (2%)	0	100	100
All	All	588/662 (89%)	569 (97%)	19 (3%)	0	100	100

There are no Ramachandran outliers to report.

### 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	264/295 (90%)	248 (94%)	16 (6%)	23	3
1	B	267/295 (90%)	258 (97%)	9 (3%)	44	14
All	All	531/590 (90%)	506 (95%)	25 (5%)	32	7

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	GLN
1	A	31	ARG
1	A	96	ARG
1	A	103	LEU
1	A	118	SER
1	A	141	ILE
1	A	171	ARG
1	A	198	ASN
1	A	202	LEU
1	A	225	LYS
1	A	249	LYS
1	A	263	ARG
1	A	275	GLN
1	A	308	SER
1	A	310	SER
1	A	311	PHE
1	B	60	GLN
1	B	136	LYS
1	B	198	ASN
1	B	202	LEU
1	B	216	TYR
1	B	225	LYS
1	B	226	ARG
1	B	250	ARG
1	B	278	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	56	ASN
1	A	60	GLN
1	A	101	ASN
1	A	198	ASN
1	A	279	GLN
1	A	282	GLN
1	A	304	HIS
1	B	6	GLN
1	B	56	ASN
1	B	60	GLN
1	B	101	ASN
1	B	198	ASN
1	B	275	GLN
1	B	279	GLN
1	B	282	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 ⓘ

4 ligands are modelled in this entry.

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
2	NAP	A	401	-	42,52,52	1.27	4 (9%)	54,80,80	1.81	11 (20%)
3	WDX	A	402	-	21,23,23	1.03	1 (4%)	27,32,32	2.24	9 (33%)
2	NAP	B	401	-	42,52,52	1.27	4 (9%)	54,80,80	2.06	10 (18%)
3	WDX	B	402	-	21,23,23	1.06	2 (9%)	27,32,32	1.80	9 (33%)

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
2	NAP	A	401	-	-	0/27/67/67	0/5/5/5
3	WDX	A	402	-	-	0/8/21/21	0/3/3/3
2	NAP	B	401	-	-	0/27/67/67	0/5/5/5
3	WDX	B	402	-	-	0/8/21/21	0/3/3/3

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	402	WDX	C13-C16	-3.34	1.44	1.50
2	A	401	NAP	C3N-C7N	-2.85	1.46	1.50
3	B	402	WDX	C13-C16	-2.76	1.45	1.50
2	A	401	NAP	O4B-C1B	-2.48	1.38	1.41
2	B	401	NAP	C2A-N3A	2.03	1.35	1.32
2	B	401	NAP	C4A-N3A	2.12	1.38	1.35
2	B	401	NAP	O4D-C1D	2.57	1.44	1.41
3	B	402	WDX	C2-N1	2.71	1.51	1.47
2	A	401	NAP	C7N-N7N	3.01	1.39	1.33
2	A	401	NAP	C6A-N6A	3.27	1.45	1.34
2	B	401	NAP	C7N-N7N	4.76	1.42	1.33

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	NAP	C5N-C4N-C3N	-7.36	111.07	120.33
3	A	402	WDX	C14-C10-C7	-6.01	114.35	120.88
2	A	401	NAP	C3N-C7N-N7N	-5.22	112.11	117.82
2	A	401	NAP	C5N-C4N-C3N	-4.79	114.31	120.33
3	A	402	WDX	C15-C12-C8	-4.54	115.89	121.82
2	B	401	NAP	P2B-O2B-C2B	-3.65	112.82	121.56
3	A	402	WDX	C3-C2-N1	-3.36	106.19	111.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	402	WDX	C15-C12-C8	-2.78	118.19	121.82
3	B	402	WDX	C3-C2-N1	-2.72	107.13	111.07
2	B	401	NAP	C4B-O4B-C1B	-2.66	106.79	109.72
3	B	402	WDX	C14-C10-C7	-2.59	118.07	120.88
3	B	402	WDX	C11-C8-C7	-2.42	104.16	106.27
3	A	402	WDX	O18-C16-C13	-2.41	113.39	118.75
2	A	401	NAP	C4A-C5A-N7A	-2.32	107.35	109.48
2	B	401	NAP	O4D-C4D-C3D	-2.27	100.57	105.15
2	A	401	NAP	O3D-C3D-C2D	2.18	118.93	111.83
3	B	402	WDX	C2-N1-C6	2.21	116.66	112.56
2	A	401	NAP	O3X-P2B-O1X	2.28	117.94	110.58
2	B	401	NAP	C2B-C3B-C4B	2.33	107.37	101.85
2	A	401	NAP	PN-O3-PA	2.35	139.32	132.73
2	B	401	NAP	C3N-C2N-N1N	2.41	123.13	120.36
2	A	401	NAP	O5B-PA-O1A	2.42	119.01	109.62
3	A	402	WDX	C5-C6-N1	2.44	114.61	111.07
3	A	402	WDX	C13-N9-C7	2.56	109.75	104.47
3	B	402	WDX	C14-C15-C12	2.60	121.71	118.49
2	A	401	NAP	C4D-O4D-C1D	2.63	112.61	109.72
3	B	402	WDX	C13-N9-C7	2.68	109.99	104.47
3	A	402	WDX	O18-C16-N1	2.85	126.96	122.42
2	A	401	NAP	C3N-C2N-N1N	2.90	123.71	120.36
2	B	401	NAP	O7N-C7N-C3N	2.91	122.76	119.59
2	B	401	NAP	C2N-C3N-C4N	3.25	121.91	118.29
3	A	402	WDX	C11-C8-C7	3.32	109.17	106.27
3	A	402	WDX	C14-C15-C12	3.41	122.72	118.49
2	A	401	NAP	C6N-C5N-C4N	3.55	124.81	119.44
3	B	402	WDX	C13-C11-C8	3.71	111.19	106.55
3	B	402	WDX	C6-C5-C4	3.76	118.09	111.73
2	B	401	NAP	C6N-C5N-C4N	4.86	126.79	119.44
2	A	401	NAP	O7N-C7N-C3N	6.01	126.14	119.59
2	B	401	NAP	N3A-C2A-N1A	7.23	134.43	128.89

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	NAP	3	0
3	A	402	WDX	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	401	NAP	4	0
3	B	402	WDX	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	298/331 (90%)	0.65	32 (10%) <b>8</b> <b>6</b>	15, 27, 51, 67	0
1	B	302/331 (91%)	0.29	15 (4%) <b>32</b> <b>28</b>	15, 23, 39, 58	0
All	All	600/662 (90%)	0.47	47 (7%) <b>16</b> <b>13</b>	15, 25, 47, 67	0

All (47) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	310	SER	9.9
1	A	131	THR	6.9
1	A	125	GLY	6.7
1	B	310	SER	6.6
1	B	135	GLY	6.4
1	A	309	ASP	5.8
1	B	311	PHE	5.4
1	A	311	PHE	5.2
1	B	309	ASP	5.2
1	A	231	ASN	5.1
1	A	141	ILE	4.9
1	A	146	THR	4.9
1	A	128	LEU	4.5
1	A	308	SER	4.0
1	A	138	ILE	3.8
1	A	180	PRO	3.7
1	A	139	PHE	3.4
1	B	136	LYS	3.3
1	A	181	GLY	3.2
1	B	308	SER	3.1
1	A	307	ASN	3.0
1	A	182	LEU	2.8
1	A	130	PRO	2.7
1	A	124	PRO	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	6	GLN	2.6
1	B	180	PRO	2.6
1	A	249	LYS	2.5
1	A	122	LEU	2.5
1	A	144	LEU	2.5
1	A	171	ARG	2.4
1	B	228	VAL	2.4
1	B	132	ASP	2.3
1	B	231	ASN	2.3
1	A	201	LYS	2.3
1	B	307	ASN	2.3
1	A	230	PRO	2.2
1	A	129	SER	2.2
1	B	204	ASP	2.2
1	B	229	ASP	2.2
1	A	291	GLU	2.1
1	A	288	LEU	2.1
1	A	233	PRO	2.1
1	B	31	ARG	2.1
1	A	173	LEU	2.1
1	A	91	ARG	2.1
1	A	142	VAL	2.0
1	B	10	LEU	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	WDX	B	402	21/21	0.93	0.10	-0.22	16,19,23,30	0
3	WDX	A	402	21/21	0.91	0.10	-0.31	22,25,30,35	0
2	NAP	B	401	48/48	0.96	0.08	-0.45	13,17,20,24	0
2	NAP	A	401	48/48	0.95	0.08	-1.11	16,21,26,28	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.