



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2016 – 08:29 PM GMT

PDB ID : 4YNK  
Title : Crystal structure of vitamin D receptor ligand binding domain complexed with a 19-norvitamin D compound  
Authors : Watarai, Y.; Ikura, T.; Ito, N.  
Deposited on : 2015-03-10  
Resolution : 2.30 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 : 1.7.1 (RC1), CSD as537be (2016)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026982  
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) : rb-20026982

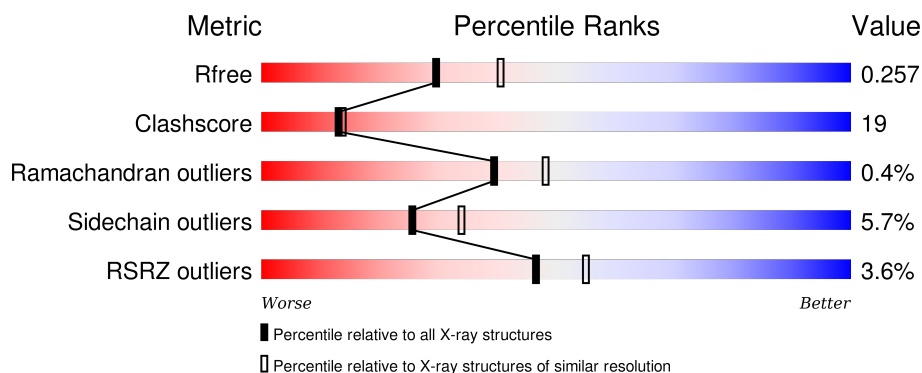
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.30 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	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)

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	271	<div> <div>3%</div> <div> <div></div> <div>54%</div> <div>32%</div> <div>•</div> <div>11%</div> </div> </div>
2	C	13	<div> <div>8%</div> <div> <div></div> <div>31%</div> <div>46%</div> <div>23%</div> </div> </div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 2093 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 Vitamin D3 receptor,Vitamin D3 receptor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	240	Total	C	N	O	S	0	0	0
			1927	1225	329	361	12			

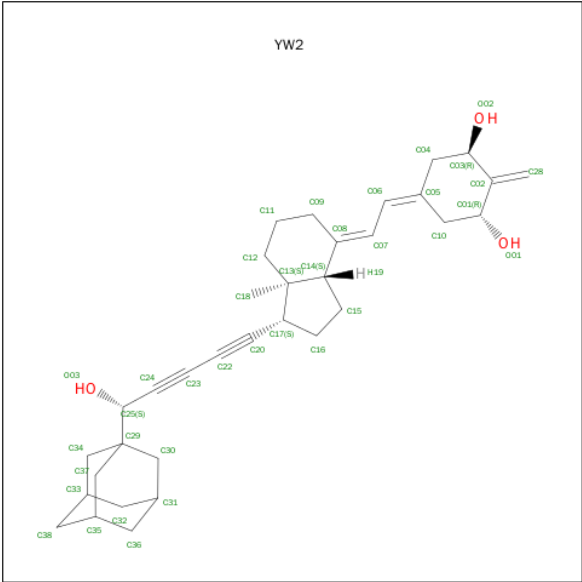
There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	106	GLY	-	expression tag	UNP P13053
A	107	SER	-	expression tag	UNP P13053
A	108	HIS	-	expression tag	UNP P13053
A	109	MET	-	expression tag	UNP P13053
A	110	GLY	-	expression tag	UNP P13053
A	111	SER	-	expression tag	UNP P13053
A	112	PRO	-	expression tag	UNP P13053
A	113	ASN	-	expression tag	UNP P13053
A	114	SER	-	expression tag	UNP P13053
A	115	PRO	-	expression tag	UNP P13053

- Molecule 2 is a protein called Coactivator peptide drip from cDNA FLJ50196, highly similar to Peroxisome proliferator-activated receptor-binding protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	10	Total	C	N	O	S	0	0	0
			82	53	15	12	2			

- Molecule 3 is (1R,3R,7E,17beta)-17-{(5S)-5-hydroxy-5-[(3R,5R,7R)-tricyclo[3.3.1.1<sup>3,7</sup>]dec-1-yl]penta-1,3-dien-1-yl}-2-methylidene-9,10-secoestra-5,7-diene-1,3-diol (three-letter code: YW2) (formula: C<sub>34</sub>H<sub>44</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	C O	0	0
			37	34 3		

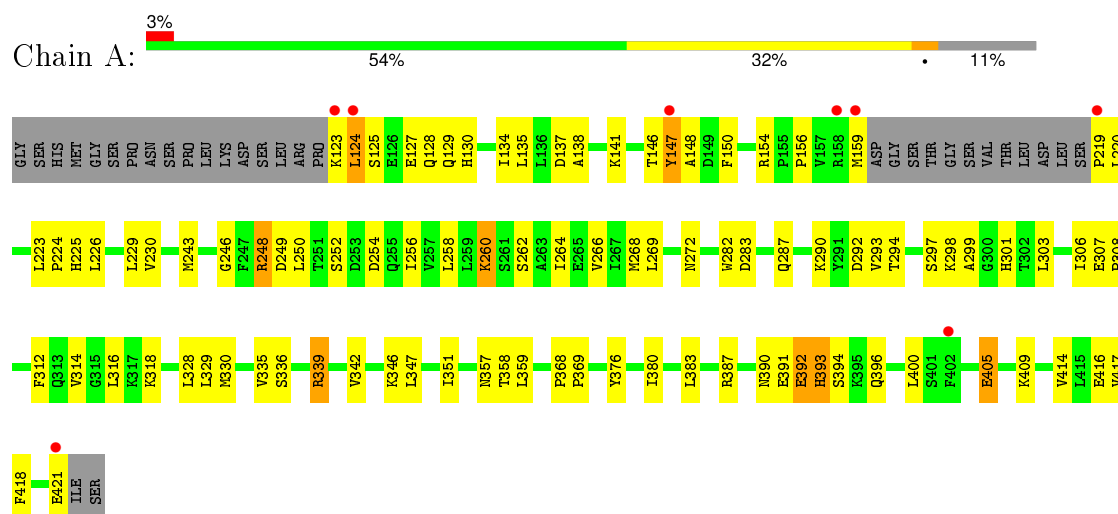
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	46	Total	O	0	0
			46	46		
4	C	1	Total	O	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: Vitamin D3 receptor,Vitamin D3 receptor



- Molecule 2: Coactivator peptide drip from cDNA FLJ50196, highly similar to Peroxisome proliferator-activated receptor-binding protein



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	157.44Å 37.29Å 41.12Å 90.00° 98.47° 90.00°	Depositor
Resolution (Å)	50.00 – 2.30 34.05 – 2.29	Depositor EDS
% Data completeness (in resolution range)	92.8 (50.00-2.30) 92.4 (34.05-2.29)	Depositor EDS
$R_{merge}$	0.13	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.31 (at 2.29Å)	Xtriage
Refinement program	CNS	Depositor
R, $R_{free}$	0.202 , 0.257 0.202 , 0.257	Depositor DCC
$R_{free}$ test set	1039 reflections (10.43%)	DCC
Wilson B-factor (Å <sup>2</sup> )	28.6	Xtriage
Anisotropy	0.734	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.39 , 56.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	1 of 10753 reflections (0.009%)	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	2093	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.74% 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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: YW2

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	0.49	3/1967 (0.2%)	0.58	0/2659
2	C	0.39	0/83	0.51	0/110
All	All	0.49	3/2050 (0.1%)	0.58	0/2769

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	394	SER	CB-OG	-7.86	1.32	1.42
1	A	394	SER	C-O	-7.37	1.09	1.23
1	A	393	HIS	C-O	-5.25	1.13	1.23

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	1927	0	1938	76	0
2	C	82	0	89	4	0
3	A	37	0	44	3	0
4	A	46	0	0	2	0
4	C	1	0	0	0	0
All	All	2093	0	2071	78	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 19.

All (78) 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:376:TYR:CE2	1:A:380:ILE:HD11	2.08	0.89
1:A:138:ALA:HB2	1:A:243:MET:HG2	1.62	0.81
1:A:147:TYR:HA	1:A:150:PHE:HB2	1.61	0.81
1:A:292:ASP:OD2	1:A:294:THR:HG22	1.82	0.78
1:A:124:LEU:HD23	1:A:128:GLN:HB2	1.75	0.69
1:A:124:LEU:HD22	1:A:129:GLN:HG3	1.75	0.69
1:A:252:SER:O	1:A:256:ILE:HG12	1.94	0.69
1:A:130:HIS:O	1:A:134:ILE:HG12	1.94	0.67
1:A:256:ILE:HD12	2:C:630:LEU:HD23	1.77	0.67
1:A:347:LEU:O	1:A:351:ILE:HG12	1.96	0.66
1:A:416:GLU:OE2	2:C:628:PRO:HD2	1.97	0.65
1:A:269:LEU:HD13	1:A:330:MET:HG2	1.78	0.65
1:A:283:ASP:OD2	1:A:290:LYS:HE2	1.98	0.64
1:A:293:VAL:HG13	1:A:306:ILE:HG21	1.80	0.64
1:A:336:SER:HB3	1:A:339:ARG:HH11	1.63	0.63
1:A:336:SER:HB3	1:A:339:ARG:NH1	2.15	0.61
1:A:339:ARG:HG3	4:A:619:HOH:O	1.99	0.61
1:A:347:LEU:O	1:A:347:LEU:HD13	2.01	0.60
1:A:339:ARG:HH11	1:A:339:ARG:HG2	1.66	0.60
1:A:226:LEU:HD12	1:A:299:ALA:HB2	1.84	0.59
1:A:223:LEU:O	1:A:223:LEU:HD23	2.02	0.59
1:A:146:THR:C	1:A:148:ALA:H	2.06	0.58
1:A:346:LYS:HG2	4:A:644:HOH:O	2.04	0.58
1:A:260:LYS:HA	1:A:417:VAL:HG13	1.88	0.56
1:A:396:GLN:O	1:A:400:LEU:HD13	2.05	0.56
1:A:159:MET:HB3	1:A:298:LYS:NZ	2.19	0.56
1:A:124:LEU:HA	1:A:128:GLN:OE1	2.07	0.55
1:A:339:ARG:O	1:A:342:VAL:HG23	2.05	0.55
1:A:220:LEU:HD12	1:A:223:LEU:HD12	1.89	0.55
1:A:339:ARG:HH11	1:A:339:ARG:CG	2.21	0.54
1:A:223:LEU:C	1:A:223:LEU:HD23	2.28	0.54
1:A:262:SER:O	1:A:266:VAL:HG23	2.08	0.54
1:A:128:GLN:HB3	1:A:358:THR:HG23	1.90	0.54
1:A:159:MET:HB3	1:A:298:LYS:HZ1	1.73	0.53
1:A:392:GLU:HA	1:A:392:GLU:OE1	2.07	0.53
1:A:147:TYR:N	1:A:147:TYR:CD2	2.74	0.53
1:A:219:PRO:O	1:A:220:LEU:HB2	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:294:THR:O	1:A:298:LYS:HG3	2.08	0.53
1:A:405:GLU:O	1:A:409:LYS:HE2	2.10	0.52
1:A:387:ARG:O	1:A:391:GLU:HG3	2.10	0.51
1:A:223:LEU:HB3	1:A:224:PRO:HD3	1.92	0.50
1:A:335:VAL:HG12	1:A:335:VAL:O	2.12	0.49
1:A:220:LEU:N	1:A:220:LEU:HD22	2.29	0.48
1:A:123:LYS:HB3	1:A:123:LYS:NZ	2.28	0.48
1:A:329:LEU:HD11	1:A:383:LEU:HG	1.95	0.48
1:A:264:ILE:O	1:A:268:MET:HG3	2.14	0.47
1:A:328:LEU:HD13	1:A:359:LEU:HA	1.97	0.46
1:A:376:TYR:O	1:A:380:ILE:HG13	2.15	0.46
2:C:634:LEU:O	2:C:635:LYS:C	2.54	0.46
1:A:130:HIS:CE1	1:A:134:ILE:HD11	2.51	0.46
1:A:250:LEU:CD2	1:A:347:LEU:HD12	2.46	0.46
1:A:248:ARG:HD2	1:A:249:ASP:N	2.31	0.46
1:A:287:GLN:HE21	1:A:290:LYS:HD3	1.80	0.46
1:A:414:VAL:HG13	1:A:418:PHE:CD1	2.52	0.44
1:A:282:TRP:CE3	3:A:500:YW2:H28	2.52	0.44
1:A:297:SER:HA	1:A:301:HIS:O	2.17	0.44
1:A:312:PHE:CE2	1:A:316:LEU:HD22	2.52	0.43
1:A:307:GLU:HB3	1:A:308:PRO:CD	2.48	0.43
1:A:339:ARG:CG	1:A:339:ARG:NH1	2.82	0.43
2:C:629:MET:O	2:C:633:LEU:HG	2.19	0.43
1:A:314:VAL:O	1:A:318:LYS:HG3	2.19	0.43
1:A:260:LYS:HB2	1:A:417:VAL:O	2.19	0.43
1:A:316:LEU:HA	1:A:316:LEU:HD12	1.87	0.42
1:A:368:PRO:HA	1:A:369:PRO:HD2	1.93	0.42
1:A:124:LEU:HD22	1:A:129:GLN:CG	2.47	0.42
1:A:282:TRP:CD2	3:A:500:YW2:H28	2.55	0.42
1:A:156:PRO:HB3	1:A:225:HIS:CD2	2.55	0.42
1:A:246:GLY:HA3	1:A:351:ILE:HD12	2.02	0.42
1:A:137:ASP:OD1	1:A:141:LYS:HE2	2.20	0.42
1:A:250:LEU:HD23	1:A:347:LEU:HD12	2.01	0.41
1:A:272:ASN:HB2	1:A:316:LEU:HD23	2.02	0.41
1:A:254:ASP:O	1:A:258:LEU:HG	2.19	0.41
1:A:123:LYS:O	1:A:124:LEU:CB	2.68	0.41
1:A:146:THR:C	1:A:148:ALA:N	2.72	0.41
1:A:127:GLU:O	1:A:130:HIS:HB3	2.21	0.40
1:A:135:LEU:HA	1:A:135:LEU:HD23	1.91	0.40
1:A:230:VAL:HG21	3:A:500:YW2:C37	2.52	0.40
1:A:125:SER:OG	1:A:128:GLN:HG3	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	236/271 (87%)	221 (94%)	14 (6%)	1 (0%)	39	48
2	C	8/13 (62%)	7 (88%)	1 (12%)	0	100	100
All	All	244/284 (86%)	228 (93%)	15 (6%)	1 (0%)	39	48

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	124	LEU

### 5.3.2 Protein sidechains [i](#)

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	220/247 (89%)	207 (94%)	13 (6%)	24	32
2	C	10/13 (77%)	10 (100%)	0	100	100
All	All	230/260 (88%)	217 (94%)	13 (6%)	25	34

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	147	TYR
1	A	154	ARG

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Mol	Chain	Res	Type
1	A	229	LEU
1	A	248	ARG
1	A	260	LYS
1	A	303	LEU
1	A	339	ARG
1	A	357	ASN
1	A	390	ASN
1	A	392	GLU
1	A	393	HIS
1	A	405	GLU
1	A	421	GLU

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	130	HIS
1	A	255	GLN
1	A	273	GLN
1	A	287	GLN
1	A	343	GLN
1	A	390	ASN
1	A	396	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 ⓘ

1 ligand is 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	$\# Z  > 2$
3	YW2	A	500	-	34,42,42	2.37	16 (47%)	46,64,64	2.03	7 (15%)

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	YW2	A	500	-	-	0/14/89/89	0/3/6/6

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	500	YW2	C14-C08	2.03	1.56	1.51
3	A	500	YW2	C04-C05	2.28	1.55	1.50
3	A	500	YW2	C36-C35	2.29	1.58	1.52
3	A	500	YW2	C32-C33	2.55	1.59	1.52
3	A	500	YW2	C32-C31	2.58	1.59	1.52
3	A	500	YW2	C38-C35	2.64	1.59	1.52
3	A	500	YW2	C36-C31	2.69	1.59	1.52
3	A	500	YW2	C38-C33	2.71	1.59	1.52
3	A	500	YW2	C01-C02	2.83	1.54	1.51
3	A	500	YW2	C30-C29	2.92	1.61	1.54
3	A	500	YW2	C34-C29	3.31	1.62	1.54
3	A	500	YW2	C37-C29	3.40	1.62	1.54
3	A	500	YW2	C03-C02	3.52	1.54	1.51
3	A	500	YW2	C18-C13	3.61	1.61	1.54
3	A	500	YW2	C10-C05	3.72	1.58	1.50
3	A	500	YW2	C23-C22	6.14	1.51	1.37

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	500	YW2	C22-C23-C24	-8.34	172.34	178.12

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	500	YW2	C23-C22-C20	-7.97	172.59	178.12
3	A	500	YW2	C15-C14-C13	-2.87	101.53	104.19
3	A	500	YW2	C25-C24-C23	-2.20	174.15	177.07
3	A	500	YW2	C12-C13-C17	-2.04	113.76	116.11
3	A	500	YW2	C18-C13-C14	2.25	113.95	110.07
3	A	500	YW2	C06-C07-C08	3.12	129.73	125.82

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	A	500	YW2	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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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	240/271 (88%)	0.29	8 (3%) 50 59	22, 35, 57, 75	0
2	C	10/13 (76%)	1.01	1 (10%) 9 14	43, 48, 57, 60	0
All	All	250/284 (88%)	0.32	9 (3%) 46 55	22, 35, 57, 75	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	159	MET	8.5
1	A	402	PHE	4.6
2	C	626	ASN	4.5
1	A	147	TYR	3.3
1	A	123	LYS	2.9
1	A	158	ARG	2.8
1	A	421	GLU	2.7
1	A	219	PRO	2.6
1	A	124	LEU	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, 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
3	YW2	A	500	37/37	0.95	0.16	0.26	20,29,33,34	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.