



# Full wwPDB X-ray Structure Validation Report ⓘ

May 26, 2020 – 07:30 pm BST

PDB ID : 4UND  
Title : HUMAN ARTD1 (PARP1) - CATALYTIC DOMAIN IN COMPLEX WITH INHIBITOR TALAZOPARIB  
Authors : Karlberg, T.; Thorsell, A.G.; Ekblad, T.; Klepsch, M.; Pinto, A.F.; Tre-saugues, L.; Moche, M.; Schuler, H.  
Deposited on : 2014-05-27  
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

<https://www.wwpdb.org/validation/2017/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 : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.11

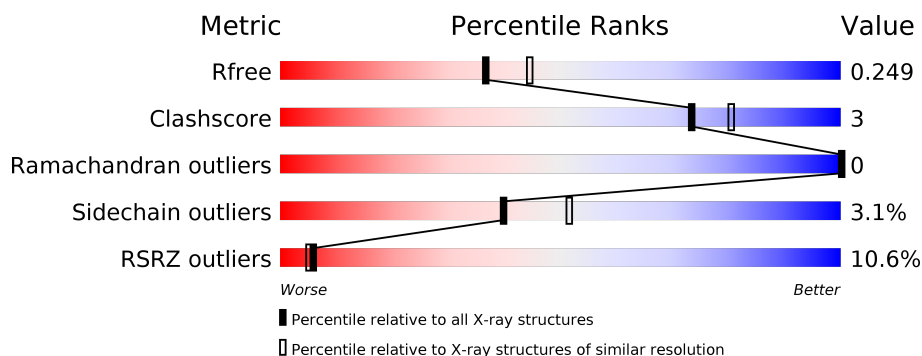
# 1 Overall quality at a glance

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))
$R_{free}$	130704	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (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 respectively. 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	373	<div> <div>8%</div> <div> <div></div> <div>84%</div> <div>10%</div> <div>• 5%</div> </div> </div>
1	B	373	<div> <div>12%</div> <div> <div></div> <div>84%</div> <div>9%</div> <div>• 6%</div> </div> </div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5677 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 POLY [ADP-RIBOSE] POLYMERASE 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	356	Total	C	N	O	S	0	0	0
			2806	1787	473	534	12			
1	B	349	Total	C	N	O	S	0	0	0
			2744	1746	464	523	11			

There are 48 discrepancies between the modelled and reference sequences:

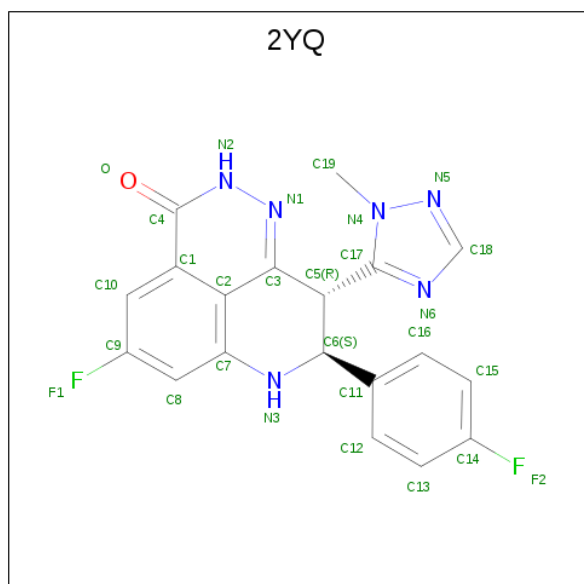
Chain	Residue	Modelled	Actual	Comment	Reference
A	639	MET	-	expression tag	UNP P09874
A	640	HIS	-	expression tag	UNP P09874
A	641	HIS	-	expression tag	UNP P09874
A	642	HIS	-	expression tag	UNP P09874
A	643	HIS	-	expression tag	UNP P09874
A	644	HIS	-	expression tag	UNP P09874
A	645	HIS	-	expression tag	UNP P09874
A	646	SER	-	expression tag	UNP P09874
A	647	SER	-	expression tag	UNP P09874
A	648	GLY	-	expression tag	UNP P09874
A	649	VAL	-	expression tag	UNP P09874
A	650	ASP	-	expression tag	UNP P09874
A	651	LEU	-	expression tag	UNP P09874
A	652	GLY	-	expression tag	UNP P09874
A	653	THR	-	expression tag	UNP P09874
A	654	GLU	-	expression tag	UNP P09874
A	655	ASN	-	expression tag	UNP P09874
A	656	LEU	-	expression tag	UNP P09874
A	657	TYR	-	expression tag	UNP P09874
A	658	PHE	-	expression tag	UNP P09874
A	659	GLN	-	expression tag	UNP P09874
A	660	SER	-	expression tag	UNP P09874
A	661	MET	-	expression tag	UNP P09874
A	762	ALA	VAL	variant	UNP P09874
B	639	MET	-	expression tag	UNP P09874

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Chain	Residue	Modelled	Actual	Comment	Reference
B	640	HIS	-	expression tag	UNP P09874
B	641	HIS	-	expression tag	UNP P09874
B	642	HIS	-	expression tag	UNP P09874
B	643	HIS	-	expression tag	UNP P09874
B	644	HIS	-	expression tag	UNP P09874
B	645	HIS	-	expression tag	UNP P09874
B	646	SER	-	expression tag	UNP P09874
B	647	SER	-	expression tag	UNP P09874
B	648	GLY	-	expression tag	UNP P09874
B	649	VAL	-	expression tag	UNP P09874
B	650	ASP	-	expression tag	UNP P09874
B	651	LEU	-	expression tag	UNP P09874
B	652	GLY	-	expression tag	UNP P09874
B	653	THR	-	expression tag	UNP P09874
B	654	GLU	-	expression tag	UNP P09874
B	655	ASN	-	expression tag	UNP P09874
B	656	LEU	-	expression tag	UNP P09874
B	657	TYR	-	expression tag	UNP P09874
B	658	PHE	-	expression tag	UNP P09874
B	659	GLN	-	expression tag	UNP P09874
B	660	SER	-	expression tag	UNP P09874
B	661	MET	-	expression tag	UNP P09874
B	762	ALA	VAL	variant	UNP P09874

- Molecule 2 is (8S,9R)-5-fluoro-8-(4-fluorophenyl)-9-(1-methyl-1H-1,2,4-triazol-5-yl)-2,7,8,9-tetrahydro-3H-pyrido[4,3,2-de]phthalazin-3-one (three-letter code: 2YQ) (formula: C<sub>19</sub>H<sub>14</sub>F<sub>2</sub>N<sub>6</sub>O).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	F	N	O	0	0
			28	19	2	6	1		
2	B	1	Total	C	F	N	O	0	0
			28	19	2	6	1		

- Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Na	0	0
			1	1		

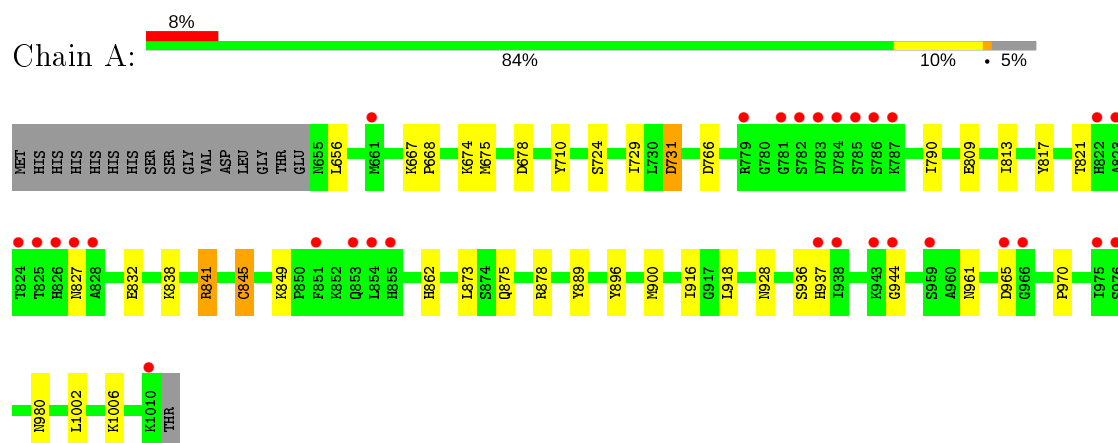
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	55	Total	O	0	0
			55	55		
4	B	15	Total	O	0	0
			15	15		

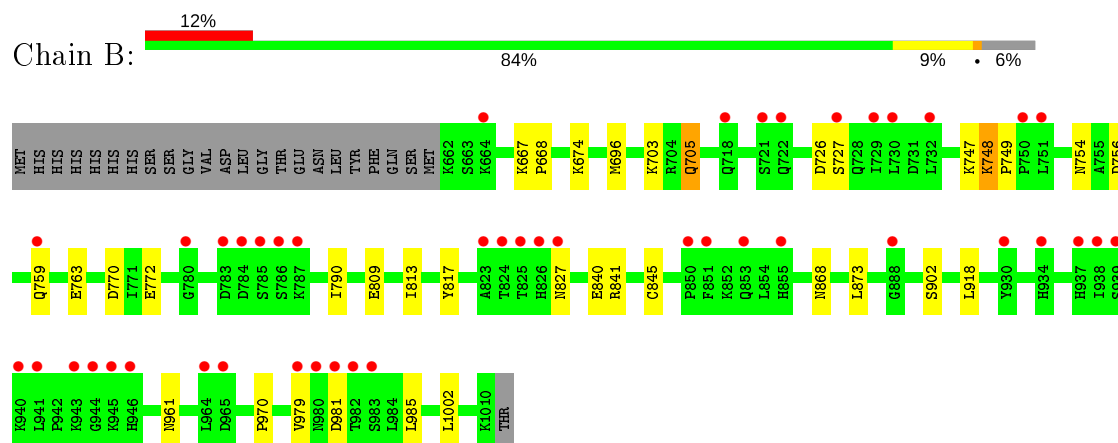
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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: POLY [ADP-RIBOSE] POLYMERASE 1



#### • Molecule 1: POLY [ADP-RIBOSE] POLYMERASE 1



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	103.45Å 103.45Å 168.54Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.27 – 2.20 46.27 – 2.20	Depositor EDS
% Data completeness (in resolution range)	100.0 (46.27-2.20) 100.0 (46.27-2.20)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.57 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, $R_{free}$	0.226 , 0.246 0.230 , 0.249	Depositor DCC
$R_{free}$ test set	2358 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	40.4	Xtriage
Anisotropy	0.218	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 39.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	5677	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.30% 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: NA, 2YQ

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	0.66	1/2860 (0.0%)	0.79	4/3859 (0.1%)
1	B	0.58	0/2796	0.74	0/3773
All	All	0.62	1/5656 (0.0%)	0.77	4/7632 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	731	ASP	CB-CG	-6.38	1.38	1.51

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	841	ARG	NE-CZ-NH2	-8.86	115.87	120.30
1	A	841	ARG	NE-CZ-NH1	6.54	123.57	120.30
1	A	731	ASP	CB-CG-OD2	-6.42	112.52	118.30
1	A	678	ASP	CB-CG-OD1	5.09	122.88	118.30

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	2806	0	2841	18	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	2744	0	2784	15	0
2	A	28	0	13	3	0
2	B	28	0	14	3	0
3	B	1	0	0	0	0
4	A	55	0	0	0	0
4	B	15	0	0	1	0
All	All	5677	0	5652	37	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 3.

All (37) 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:2011:2YQ:F2	2:B:2011:2YQ:C14	1.66	1.34
2:A:2011:2YQ:C14	2:A:2011:2YQ:F2	1.89	1.10
1:B:726:ASP:HB2	4:B:3002:HOH:O	1.88	0.72
1:A:841:ARG:HD2	1:A:873:LEU:O	1.92	0.70
2:B:2011:2YQ:H12	2:B:2011:2YQ:C12	2.28	0.64
1:B:770:ASP:HB2	1:B:868:ASN:OD1	2.01	0.61
1:B:754:ASN:OD1	1:B:756:ASP:N	2.36	0.59
1:A:724:SER:HB2	1:A:729:ILE:HD13	1.84	0.58
1:B:748:LYS:HD3	1:B:749:PRO:HD2	1.89	0.55
1:B:961:ASN:ND2	1:B:970:PRO:HA	2.23	0.53
1:A:961:ASN:ND2	1:A:970:PRO:HA	2.25	0.51
1:A:862:HIS:HE1	2:A:2011:2YQ:C18	2.23	0.51
1:B:759:GLN:O	1:B:763:GLU:HG2	2.11	0.51
1:A:896:TYR:H	2:A:2011:2YQ:C18	2.25	0.50
1:A:928:ASN:HB2	1:A:944:GLY:O	2.13	0.49
1:B:841:ARG:NE	1:B:873:LEU:O	2.46	0.49
1:B:770:ASP:HB3	1:B:868:ASN:HA	1.95	0.48
1:A:845:CYS:O	1:A:849:LYS:HG2	2.14	0.47
1:A:710:TYR:OH	1:A:766:ASP:OD1	2.26	0.47
1:B:667:LYS:HB3	1:B:668:PRO:HD3	1.96	0.46
1:B:703:LYS:HG3	1:B:772:GLU:OE1	2.15	0.46
1:B:696:MET:CE	1:B:705:GLN:OE1	2.65	0.44
1:B:918:LEU:HD22	1:B:1002:LEU:HD21	2.00	0.43
1:A:832:GLU:HB3	1:A:1006:LYS:HB3	1.99	0.43
1:A:809:GLU:O	1:A:813:ILE:HG12	2.18	0.43
1:A:980:ASN:N	1:A:980:ASN:OD1	2.40	0.42
1:A:674:LYS:HE2	1:A:790:ILE:HD11	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:821:THR:HB	1:A:900:MET:HA	2.00	0.42
1:B:809:GLU:O	1:B:813:ILE:HG12	2.20	0.42
1:A:675:MET:SD	1:A:916:ILE:HD11	2.59	0.41
1:B:674:LYS:HE2	1:B:790:ILE:HD11	2.03	0.41
1:A:875:GLN:OE1	1:A:878:ARG:NH1	2.52	0.41
1:A:918:LEU:HD22	1:A:1002:LEU:HD21	2.02	0.41
1:A:889:TYR:O	1:A:937:HIS:CE1	2.73	0.41
1:A:667:LYS:HB3	1:A:668:PRO:HD3	2.02	0.41
2:B:2011:2YQ:H12	2:B:2011:2YQ:C13	2.51	0.41
1:B:747:LYS:O	1:B:748:LYS:C	2.59	0.41

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	354/373 (95%)	339 (96%)	15 (4%)	0	100	100
1	B	347/373 (93%)	331 (95%)	16 (5%)	0	100	100
All	All	701/746 (94%)	670 (96%)	31 (4%)	0	100	100

There are no Ramachandran outliers to report.

### 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	313/328 (95%)	305 (97%)	8 (3%)	46	58
1	B	306/328 (93%)	295 (96%)	11 (4%)	35	45
All	All	619/656 (94%)	600 (97%)	19 (3%)	40	51

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	656	LEU
1	A	731	ASP
1	A	817	TYR
1	A	827	ASN
1	A	838	LYS
1	A	845	CYS
1	A	936	SER
1	A	965	ASP
1	B	705	GLN
1	B	727	SER
1	B	748	LYS
1	B	817	TYR
1	B	827	ASN
1	B	840	GLU
1	B	845	CYS
1	B	902	SER
1	B	979	VAL
1	B	981	ASP
1	B	985	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	717	GLN
1	A	734	ASN
1	A	827	ASN
1	A	906	ASN
1	A	961	ASN
1	B	793	ASN
1	B	827	ASN
1	B	961	ASN

### 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 ⓘ

Of 3 ligands modelled in this entry, 1 is 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	$\# Z  > 2$
2	2YQ	A	2011	-	27,32,32	10.11	11 (40%)	30,48,48	3.31	15 (50%)
2	2YQ	B	2011	-	27,32,32	5.39	5 (18%)	30,48,48	2.61	12 (40%)

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	2YQ	A	2011	-	-	3/4/20/20	0/4/5/5
2	2YQ	B	2011	-	-	1/4/20/20	0/4/5/5

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	2011	2YQ	C3-C5	-44.65	1.34	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	2011	2YQ	C3-C5	-22.67	1.42	1.51
2	A	2011	2YQ	F2-C14	21.95	1.89	1.36
2	A	2011	2YQ	C11-C6	-12.82	1.34	1.52
2	B	2011	2YQ	F2-C14	12.34	1.66	1.36
2	B	2011	2YQ	C11-C6	-7.34	1.42	1.52
2	B	2011	2YQ	C6-N3	5.05	1.51	1.46
2	A	2011	2YQ	F1-C9	4.96	1.48	1.36
2	A	2011	2YQ	C17-C5	4.38	1.57	1.50
2	A	2011	2YQ	C7-N3	-3.53	1.36	1.40
2	A	2011	2YQ	C17-N6	3.21	1.39	1.34
2	A	2011	2YQ	N2-N1	3.03	1.42	1.34
2	A	2011	2YQ	C7-C2	-2.90	1.38	1.42
2	A	2011	2YQ	C4-N2	2.67	1.42	1.31
2	A	2011	2YQ	C6-C5	2.13	1.59	1.55
2	B	2011	2YQ	C6-C5	2.12	1.59	1.55

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	2011	2YQ	C11-C6-N3	8.62	124.53	110.93
2	A	2011	2YQ	C19-N4-C17	7.26	137.45	128.84
2	A	2011	2YQ	C10-C9-C8	-6.56	118.90	124.09
2	A	2011	2YQ	C5-C3-N1	-6.07	111.77	118.18
2	B	2011	2YQ	C11-C6-N3	6.02	120.43	110.93
2	B	2011	2YQ	C10-C9-C8	-4.91	120.21	124.09
2	A	2011	2YQ	C12-C11-C6	-4.41	113.78	120.78
2	B	2011	2YQ	C2-C3-N1	-4.38	118.50	121.50
2	A	2011	2YQ	F1-C9-C8	4.31	124.40	118.25
2	B	2011	2YQ	C15-C14-C13	-4.09	117.38	122.83
2	B	2011	2YQ	C2-C7-N3	4.06	120.11	117.45
2	A	2011	2YQ	C2-C3-N1	3.96	124.22	121.50
2	B	2011	2YQ	C3-N1-N2	3.67	123.41	119.67
2	B	2011	2YQ	C19-N4-C17	3.66	133.18	128.84
2	B	2011	2YQ	C12-C13-C14	3.19	121.66	118.36
2	B	2011	2YQ	F1-C9-C8	3.19	122.81	118.25
2	B	2011	2YQ	C3-C5-C6	2.98	115.08	109.81
2	A	2011	2YQ	C15-C16-C11	-2.94	118.24	121.20
2	A	2011	2YQ	C10-C1-C2	2.82	121.53	118.94
2	A	2011	2YQ	C3-N1-N2	-2.80	116.83	119.67
2	B	2011	2YQ	C16-C15-C14	2.77	121.23	118.36
2	B	2011	2YQ	F2-C14-C15	2.61	122.97	118.54
2	A	2011	2YQ	F2-C14-C15	2.34	122.52	118.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	2011	2YQ	C16-C11-C12	2.33	121.19	118.29
2	A	2011	2YQ	C16-C15-C14	2.31	120.75	118.36
2	A	2011	2YQ	C15-C14-C13	-2.24	119.85	122.83
2	A	2011	2YQ	C9-C10-C1	2.12	120.95	119.43

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	2011	2YQ	C16-C11-C6-C5
2	A	2011	2YQ	C12-C11-C6-N3
2	A	2011	2YQ	C12-C11-C6-C5
2	B	2011	2YQ	C16-C11-C6-N3

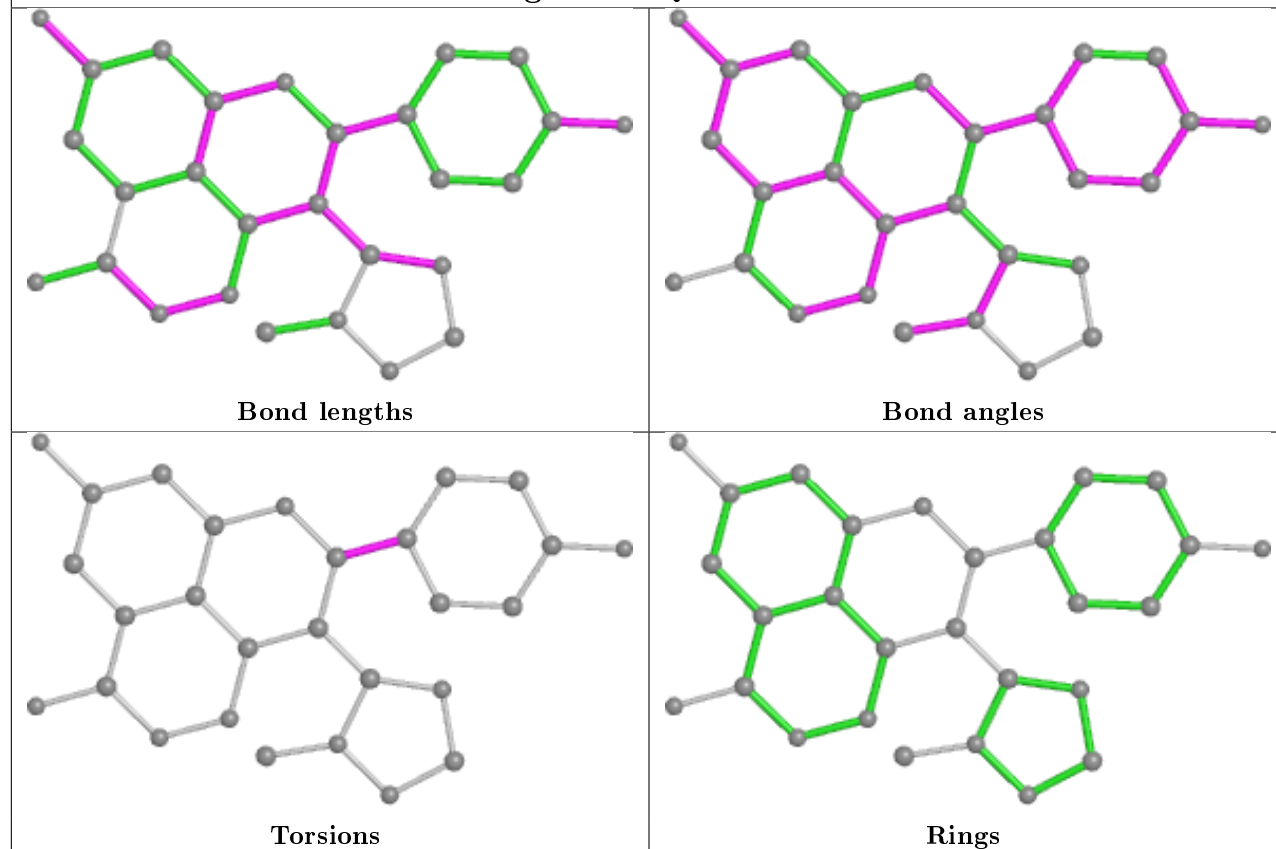
There are no ring outliers.

2 monomers are involved in 6 short contacts:

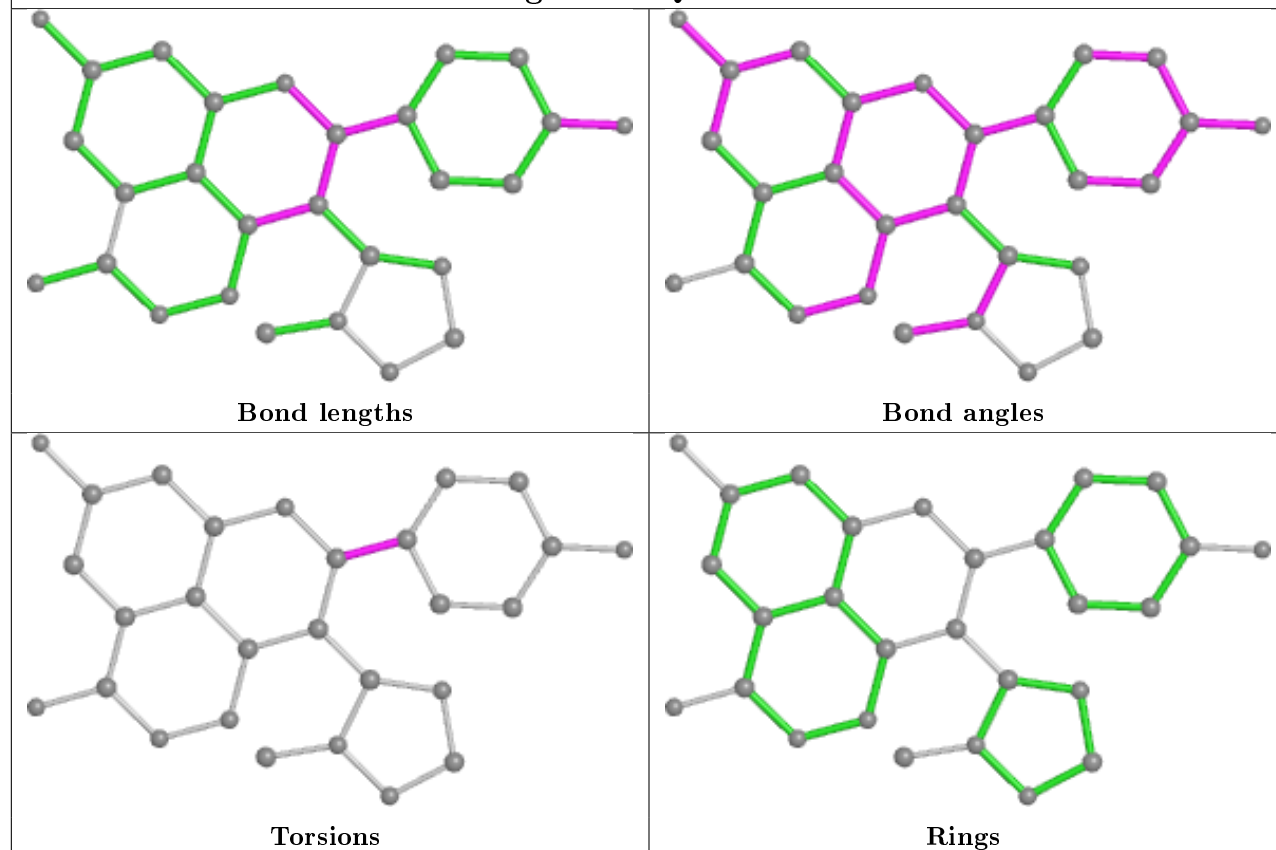
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	2011	2YQ	3	0
2	B	2011	2YQ	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

## Ligand 2YQ A 2011



## Ligand 2YQ B 2011



## 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	356/373 (95%)	0.34	30 (8%)	11 9	24, 47, 99, 164	0
1	B	349/373 (93%)	0.68	45 (12%)	3 3	31, 62, 125, 148	0
All	All	705/746 (94%)	0.51	75 (10%)	6 5	24, 54, 117, 164	0

All (75) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	826	HIS	10.1
1	A	825	THR	8.9
1	B	823	ALA	8.3
1	B	941	LEU	8.3
1	A	785	SER	8.0
1	A	781	GLY	7.7
1	B	721	SER	7.3
1	A	784	ASP	7.2
1	A	824	THR	7.2
1	A	826	HIS	6.6
1	B	980	ASN	6.5
1	B	939	SER	6.4
1	A	823	ALA	5.9
1	B	943	LYS	5.7
1	B	938	ILE	5.4
1	A	966	GLY	5.3
1	B	930	TYR	5.1
1	A	782	SER	5.0
1	A	783	ASP	5.0
1	B	825	THR	4.8
1	A	937	HIS	4.7
1	B	940	LYS	4.6
1	B	944	GLY	4.5
1	A	827	ASN	4.2

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Mol	Chain	Res	Type	RSRZ
1	B	727	SER	4.2
1	A	943	LYS	3.9
1	B	786	SER	3.9
1	B	934	HIS	3.8
1	B	750	PRO	3.8
1	B	785	SER	3.5
1	B	759	GLN	3.4
1	B	982	THR	3.4
1	B	937	HIS	3.4
1	B	732	LEU	3.2
1	A	959	SER	3.1
1	A	855	HIS	3.1
1	B	722	GLN	3.0
1	B	729	ILE	2.9
1	B	946	HIS	2.9
1	B	783	ASP	2.9
1	B	784	ASP	2.9
1	B	718	GLN	2.9
1	A	965	ASP	2.8
1	B	751	LEU	2.8
1	B	945	LYS	2.7
1	A	822	HIS	2.6
1	B	979	VAL	2.6
1	B	855	HIS	2.6
1	A	944	GLY	2.6
1	A	779	ARG	2.6
1	A	787	LYS	2.6
1	B	888	GLY	2.5
1	B	983	SER	2.4
1	A	661	MET	2.4
1	B	824	THR	2.4
1	A	976	SER	2.4
1	A	786	SER	2.4
1	B	664	LYS	2.3
1	B	787	LYS	2.3
1	A	851	PHE	2.3
1	A	853	GLN	2.2
1	B	965	ASP	2.2
1	B	853	GLN	2.2
1	B	780	GLY	2.2
1	A	975	ILE	2.2
1	B	730	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	981	ASP	2.2
1	B	827	ASN	2.1
1	B	964	LEU	2.1
1	A	938	ILE	2.1
1	A	854	LEU	2.1
1	B	850	PRO	2.1
1	A	828	ALA	2.1
1	A	1010	LYS	2.0
1	B	851	PHE	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](#)

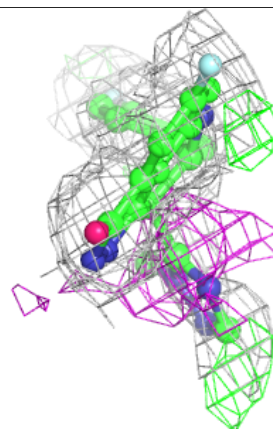
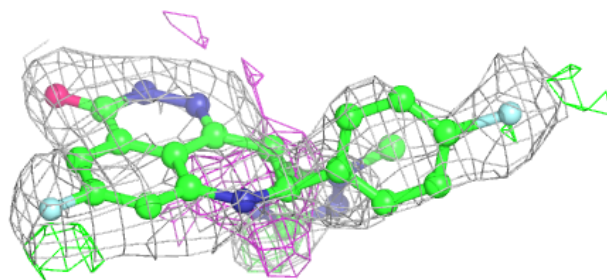
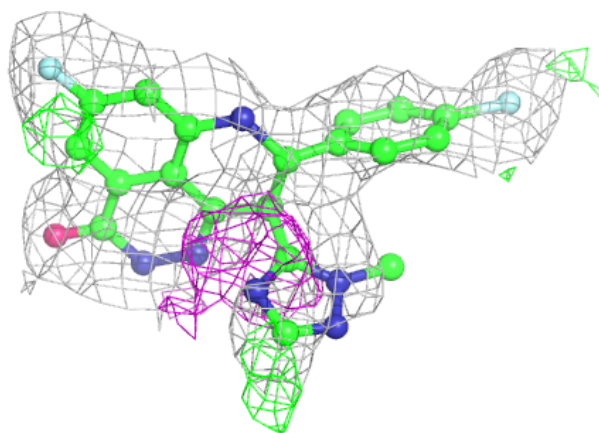
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	2YQ	B	2011	28/28	0.72	0.23	53,77,93,94	0
2	2YQ	A	2011	28/28	0.78	0.18	33,44,65,69	0
3	NA	B	2012	1/1	0.94	0.24	46,46,46,46	0

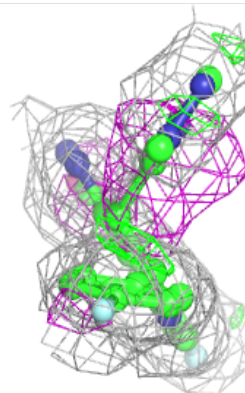
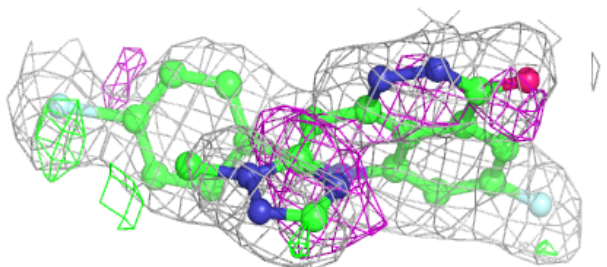
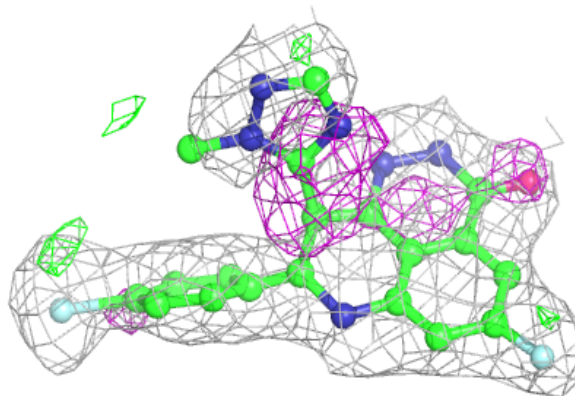
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

**Electron density around 2YQ B 2011:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around 2YQ A 2011:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



## 6.5 Other polymers

There are no such residues in this entry.