

# Full wwPDB X-ray Structure Validation Report (i)

#### Dec 19, 2022 – 03:13 pm GMT

PDB ID	:	8B95
Title	:	Crystal structure of PPARG and NCOR2 with BAY-9683, an inverse agonist
Authors	:	Friberg, A.; Orsi, D.L.; Pook, E.; Siegel, S.; Lemke, C.T.; Stellfeld, T.; Puetter,
		V.; Goldstein, J.
Deposited on	:	2022-10-05
Resolution	:	1.72  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 A user guide is available at https://www.wwpdb.org/validation/2017/XrayValidationReportHelp with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (1)) were used in the production of this report:

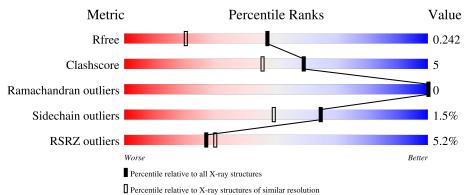
MolProbity		4.02b-467
5		
Mogul	:	1.8.4, CSD as $541be(2020)$
Xtriage (Phenix)	:	1.13
$\mathrm{EDS}$	:	2.31.3
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.31.3

# 1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:  $X\text{-}RAY \, DIFFRACTION$ 

The reported resolution of this entry is 1.72 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	$egin{array}{c} { m Whole \ archive} \ (\#{ m Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
$R_{free}$	130704	5722(1.74-1.70)
Clashscore	141614	6152(1.74-1.70)
Ramachandran outliers	138981	6051 (1.74-1.70)
Sidechain outliers	138945	6051 (1.74-1.70)
RSRZ outliers	127900	5629(1.74-1.70)

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 of the lower bar indicate the fraction of residues that contain outliers for >=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 <=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		C	Quality of chain		
1	А	279	4%		82%	10%	• 7%
1	В	279	5%	7	8%	7% •	14%
2	С	23	4%	39%	9%	52%	
2	D	23	13%	52%		48%	



# 2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 4589 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.

• N	• Molecule 1 is a protein called Peroxisome proliferator-activated receptor gamma.									
Mol	Chain	Residues		At	oms			ZeroOcc	AltConf	Trac
				$\sim$		0	~			

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	А	260	Total 2105	C 1359	N 249	0	S 10	0	3	0
			2105	1998	548	300	10			
1	В	239	Total	$\mathbf{C}$	Ν	Ο	$\mathbf{S}$	0	2	0
1	D	239	1916	1238	315	354	9	0		0

There are 8 discrepancies between the modelled and reference sequences:

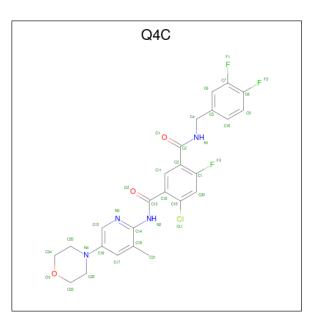
Chain	Residue	Modelled	Actual	Comment	Reference
А	199	GLY	-	expression tag	UNP P37231
А	200	SER	-	expression tag	UNP P37231
А	201	HIS	-	expression tag	UNP P37231
А	202	MET	-	expression tag	UNP P37231
В	199	GLY	-	expression tag	UNP P37231
В	200	SER	-	expression tag	UNP P37231
В	201	HIS	-	expression tag	UNP P37231
В	202	MET	-	expression tag	UNP P37231

• Molecule 2 is a protein called Nuclear receptor corepressor 2.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
2	C	11	Total	С	Ν	Ο	S	0	0	0
	U	11	83	54	15	13	1	0		
2	Л	19	Total	С	Ν	Ο	S	0	0	0
	D	12	87	56	16	14	1	0	0	0

• Molecule 3 is  $\{N\}1-[[3,4-bis(fluoranyl)phenyl]methyl]-4-chloranyl-6-fluoranyl- <math>\{N\}3-(3-methyl-5-morpholin-4-yl-pyridin-2-yl)$ benzene-1,3-dicarboxamide (three-letter code: Q4C) (formula:  $C_{25}H_{22}ClF_3N_4O_3$ ) (labeled as "Ligand of Interest" by depositor).





Mol	Chain	Residues	Atoms				ZeroOcc	AltConf		
2	Λ	1	Total	С	F	Ν	Ο	0	0	
5	A	1	35	25	3	4	3	0	0	
9	D	1	Total	С	F	Ν	0	0	0	
0	D	1	35	25	3	4	3	0	0	

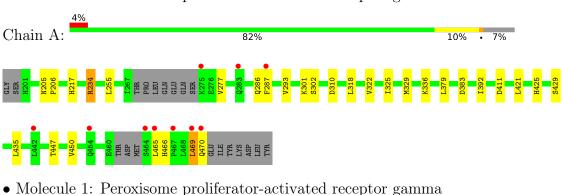
• Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	А	176	Total O 176 176	0	0
4	В	152	Total O 152 152	0	0

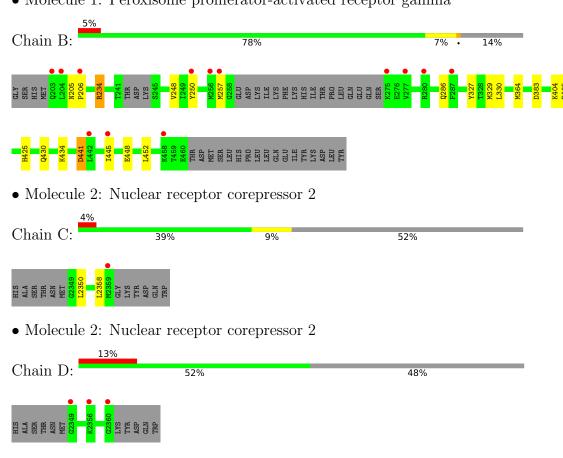


# 3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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: Peroxisome proliferator-activated receptor gamma





## 4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants	60.64Å 86.07Å 121.01Å	Depositor
a, b, c, $\alpha$ , $\beta$ , $\gamma$	$90.00^{\circ}$ $90.00^{\circ}$ $90.00^{\circ}$	Depositor
Resolution (Å)	42.83 - 1.72	Depositor
Resolution (A)	42.83 - 1.72	EDS
% Data completeness	99.8 (42.83-1.72)	Depositor
(in resolution range)	99.8 (42.83-1.72)	EDS
R <sub>merge</sub>	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$< I/\sigma(I) > 1$	$1.12 (at 1.72 \text{\AA})$	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
D D.	0.196 , $0.238$	Depositor
$R, R_{free}$	0.203 , $0.242$	DCC
$R_{free}$ test set	3389 reflections $(5.00%)$	wwPDB-VP
Wilson B-factor $(Å^2)$	34.8	Xtriage
Anisotropy	0.049	Xtriage
Bulk solvent $k_{sol}(e/Å^3), B_{sol}(Å^2)$	0.33 , $42.7$	EDS
L-test for twinning <sup>2</sup>	$ \langle L  \rangle = 0.50, \langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	4589	wwPDB-VP
Average B, all atoms $(Å^2)$	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 5.80% of the height of the origin peak. No significant pseudotranslation is detected.

<sup>&</sup>lt;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.



<sup>&</sup>lt;sup>1</sup>Intensities estimated from amplitudes.

# 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:  $\rm Q4C$ 

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).

Mal	Mol Chain		lengths	Bond angles		
	Chain	RMSZ	# Z  > 5	RMSZ	# Z  > 5	
1	А	0.74	0/2141	0.86	2/2881~(0.1%)	
1	В	0.75	0/1949	0.83	3/2623~(0.1%)	
2	С	0.82	0/82	0.69	0/107	
2	D	0.76	0/86	0.76	0/112	
All	All	0.75	0/4258	0.84	5/5723~(0.1%)	

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\mathbf{Observed}(^{o})$	$\mathbf{Ideal}(^{o})$
1	А	234	ARG	NE-CZ-NH2	-10.97	114.81	120.30
1	А	234	ARG	NE-CZ-NH1	10.03	125.31	120.30
1	В	234	ARG	NE-CZ-NH1	-6.16	117.22	120.30
1	В	234	ARG	NE-CZ-NH2	6.11	123.36	120.30
1	В	327	TYR	CB-CA-C	5.19	120.78	110.40

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	А	2105	0	2157	28	0
1	В	1916	0	1981	12	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes				
2	С	83	0	97	2	0				
2	D	87	0	100	0	0				
3	А	35	0	0	0	0				
3	В	35	0	0	2	0				
4	А	176	0	0	5	0				
4	В	152	0	0	2	0				
All	All	4589	0	4335	40	0				

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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 5.

All (40) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash
1:A:465:LEU:CD2	1:A:469:LEU:HD13	$\frac{2.12}{2.12}$	$\frac{\text{overlap (Å)}}{0.80}$
1:A:405:LE0:CD2	4:A:747:HOH:O	2.12	0.80
1:B:383:ASP:OD2	4:A:747:HOH:0 1:B:425:HIS:HE1	$\frac{2.34}{1.72}$	
			0.71
1:A:234:ARG:HD2	4:A:747:HOH:O	1.93	0.67
1:A:234:ARG:HD3	4:A:747:HOH:O	1.92	0.67
1:A:411[B]:ASP:OD1	4:A:601:HOH:O	2.13	0.65
1:B:452:LEU:HD23	4:B:733:HOH:O	1.99	0.63
1:A:425:HIS:HD2	4:A:756:HOH:O	1.83	0.62
1:A:465:LEU:HG	1:A:469:LEU:HD13	1.82	0.60
1:A:383:ASP:OD2	1:A:425:HIS:HE1	1.85	0.59
1:A:465:LEU:CG	1:A:469:LEU:HD13	2.33	0.58
1:A:379:LEU:HD11	1:A:435:LEU:HD11	1.85	0.58
1:B:330:LEU:HD21	1:B:364:MET:HE1	1.86	0.57
1:B:205:ASN:HB2	1:B:206:PRO:HD2	1.87	0.56
1:A:217:HIS:HE1	1:A:302:SER:O	1.89	0.54
1:A:465:LEU:HD21	1:A:469:LEU:HD13	1.89	0.54
1:A:465:LEU:HD12	1:A:466:HIS:H	1.73	0.53
1:A:286:GLN:HE21	1:A:465:LEU:HD11	1.75	0.52
1:A:447:THR:O	1:A:450:VAL:HG22	2.10	0.51
1:B:430:GLN:O	1:B:434[B]:LYS:HG2	2.11	0.51
1:A:465:LEU:O	1:A:470:GLN:NE2	2.43	0.50
1:A:318:LEU:O	1:A:322:VAL:HG13	2.11	0.49
1:A:255:LEU:HD21	1:A:277:VAL:HG13	1.94	0.49
1:A:287[A]:PHE:CD2	1:A:466:HIS:CD2	3.01	0.49
1:B:441:ASP:O	1:B:445:ILE:HG12	2.13	0.48
1:A:255:LEU:CD2	1:A:277:VAL:HG13	2.45	0.46
1:B:286:GLN:HE22	3:B:501:Q4C:C5	2.29	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:448:GLU:O	1:B:452:LEU:HG	2.16	0.46
1:B:329:MET:HE3	3:B:501:Q4C:C24	2.46	0.45
1:A:325:ILE:HD11	1:A:392:ILE:HG13	2.00	0.44
1:B:404:LYS:HB3	1:B:405:PRO:HD3	2.01	0.42
1:B:425:HIS:HD2	4:B:738:HOH:O	2.02	0.42
1:A:329:MET:HB2	1:A:329:MET:HE3	1.95	0.41
1:A:421:LEU:HD11	1:A:435:LEU:HD22	2.02	0.41
1:A:301:LYS:HE2	2:C:2358:LEU:HA	2.02	0.41
1:A:205:ASN:HB2	1:A:206:PRO:HD2	2.03	0.41
1:A:325:ILE:HD11	1:A:392:ILE:CG1	2.50	0.41
1:B:248:VAL:HG11	1:B:250:TYR:CE2	2.56	0.41
1:A:466:HIS:HB3	1:A:469:LEU:HD12	2.03	0.40
1:A:293:VAL:HG21	2:C:2350:LEU:CD1	2.52	0.40

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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	А	257/279~(92%)	254 (99%)	3~(1%)	0	100 100
1	В	235/279~(84%)	233~(99%)	2(1%)	0	100 100
2	С	9/23~(39%)	9 (100%)	0	0	100 100
2	D	10/23~(44%)	10 (100%)	0	0	100 100
All	All	511/604~(85%)	506 (99%)	5 (1%)	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 side chain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	А	235/251~(94%)	231~(98%)	4 (2%)	60 44
1	В	214/251~(85%)	211 (99%)	3 (1%)	67 52
2	С	8/18 (44%)	8 (100%)	0	100 100
2	D	8/18 (44%)	8 (100%)	0	100 100
All	All	465/538~(86%)	458 (98%)	7~(2%)	65 49

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	$\mathbf{Res}$	Type
1	А	310	ASP
1	А	336	LYS
1	А	429	SER
1	А	469	LEU
1	В	234	ARG
1	В	257	MET
1	В	441	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	А	217	HIS
1	А	266	HIS
1	А	314	GLN
1	А	425	HIS
1	В	286	GLN
1	В	425	HIS

#### 5.3.3 RNA (i)

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 monosaccharides in this entry.

### 5.6 Ligand geometry (i)

2 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	Trune	Chain	Dec	Tinle	Link Bond lengths			В	ond ang	gles
	Type	Chain	$\mathbf{Res}$	LIIIK	Counts	RMSZ	# Z >2	Counts	RMSZ	# Z >2
3	Q4C	В	501	1	38,38,39	0.76	2(5%)	$50,\!53,\!55$	1.40	7 (14%)
3	Q4C	А	501	1	38,38,39	0.77	0	50, 53, 55	1.87	13 (26%)

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	Q4C	В	501	1	-	5/21/29/29	0/4/4/4
3	Q4C	А	501	1	-	5/21/29/29	0/4/4/4

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms		Observed(Å)	$\mathrm{Ideal}(\mathrm{\AA})$
3	В	501	Q4C	C14-N2	-2.47	1.34	1.39
3	В	501	Q4C	C18-C14	-2.06	1.38	1.41

All (20) bond angle outliers are listed below:



Mol	Chain	Res	Type	Atoms	Z	Observed(°)	$Ideal(^{o})$
3	А	501	Q4C	C18-C14-N3	-5.46	117.33	123.54
3	В	501	Q4C	C25-N4-C22	5.04	122.64	111.52
3	А	501	Q4C	C12-C13-N2	-3.90	107.33	115.92
3	А	501	Q4C	C25-N4-C22	3.79	119.88	111.52
3	А	501	Q4C	C17-C16-C15	-3.77	116.50	119.48
3	А	501	Q4C	C20-C19-C12	-3.75	116.42	120.78
3	А	501	Q4C	C15-N3-C14	3.39	122.85	116.05
3	В	501	Q4C	C20-C1-C2	-3.11	119.72	123.11
3	В	501	Q4C	C18-C14-N3	-2.95	120.19	123.54
3	А	501	Q4C	C20-C1-C2	-2.94	119.90	123.11
3	А	501	Q4C	C19-C20-C1	2.70	122.31	119.05
3	В	501	Q4C	C15-N3-C14	2.54	121.15	116.05
3	В	501	Q4C	C16-C17-C18	-2.49	118.83	121.60
3	А	501	Q4C	O2-C13-N2	2.48	129.38	123.71
3	А	501	Q4C	O1-C3-C2	-2.34	116.73	121.01
3	В	501	Q4C	C11-C2-C1	2.30	119.46	116.66
3	В	501	Q4C	C20-C19-C12	-2.28	118.13	120.78
3	А	501	Q4C	C14-N2-C13	-2.16	121.30	126.64
3	А	501	Q4C	C24-C25-N4	2.10	113.90	110.02
3	А	501	Q4C	C19-C12-C11	2.06	121.67	119.24

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	А	501	Q4C	C11-C12-C13-N2
3	А	501	Q4C	C11-C12-C13-O2
3	А	501	Q4C	C19-C12-C13-N2
3	В	501	Q4C	C11-C12-C13-N2
3	В	501	Q4C	C11-C12-C13-O2
3	А	501	Q4C	C19-C12-C13-O2
3	В	501	Q4C	C19-C12-C13-N2
3	В	501	Q4C	C19-C12-C13-O2
3	А	501	Q4C	N3-C14-N2-C13
3	В	501	Q4C	N3-C14-N2-C13

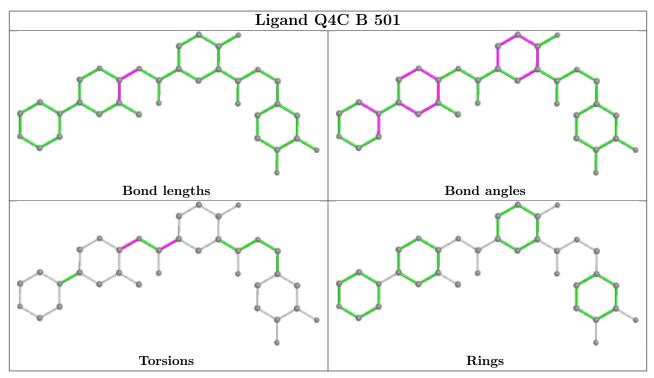
There are no ring outliers.

1 monomer is involved in 2 short contacts:

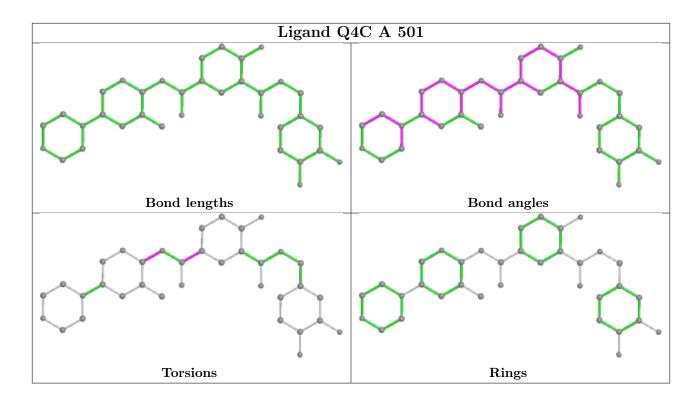
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	В	501	Q4C	2	0



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 then 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 sufficient the outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.







## 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,  $95^{th}$  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	$OWAB(Å^2)$	$\mathbf{Q}{<}0.9$
1	А	260/279~(93%)	0.14	10 (3%) 40 45	26, 39, 76, 93	0
1	В	239/279~(85%)	0.19	13 (5%) 25 28	26, 38, 67, 99	0
2	С	11/23~(47%)	0.51	1 (9%) 9 10	45, 57, 72, 89	0
2	D	12/23~(52%)	0.53	$3\ (25\%)\ 0\ 0$	41, 52, 79, 87	0
All	All	522/604~(86%)	0.18	27 (5%) 27 30	26, 39, 74, 99	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	А	465	LEU	7.1
1	А	464	SER	6.7
1	А	275	LYS	5.9
1	В	256	MET	4.9
1	А	469	LEU	4.6
1	В	287	PHE	4.4
1	В	206	PRO	4.3
1	В	250	TYR	4.0
1	В	204	LEU	3.6
2	С	2359	MET	3.5
1	В	275	LYS	2.9
1	В	458	LYS	2.9
1	А	470	GLN	2.8
1	А	283	GLN	2.7
1	В	203	GLN	2.7
1	А	287[A]	PHE	2.6
1	В	257	MET	2.5
1	А	442	LEU	2.5
2	D	2360	GLY	2.5
1	А	454	GLN	2.5
1	В	280	ARG	2.4

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Mol	Chain	Res	Type	RSRZ
1	В	442	LEU	2.3
1	А	467	PRO	2.3
1	В	445	ILE	2.3
1	В	277	VAL	2.2
2	D	2349	GLY	2.2
2	D	2356	LYS	2.0

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### 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 monosaccharides in this entry.

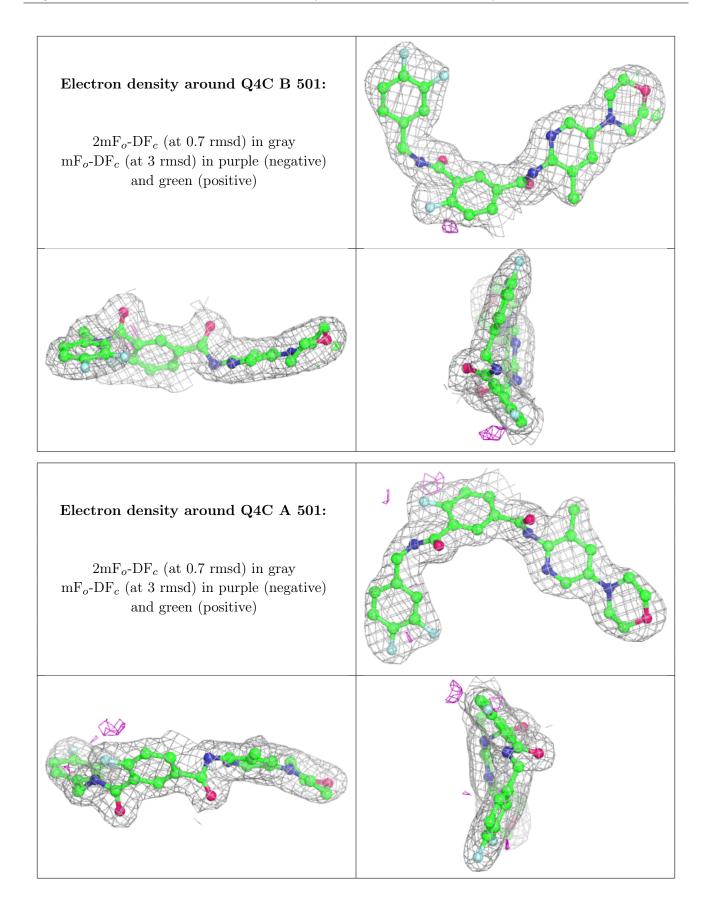
#### 6.4 Ligands (i)

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,  $95^{th}$  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	$B-factors(Å^2)$	Q<0.9
3	Q4C	В	501	35/36	0.90	0.11	$33,\!38,\!59,\!70$	0
3	Q4C	А	501	35/36	0.94	0.09	29,40,53,63	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.







## 6.5 Other polymers (i)

There are no such residues in this entry.

