

Full wwPDB X-ray Structure Validation Report (i)

Feb 22, 2024 – 11:00 AM EST

PDB ID	:	4Y86
Title	:	Crystal structure of PDE9 in complex with racemic inhibitor C33
Authors	:	Huang, M.
Deposited on		
Resolution	:	2.01 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:

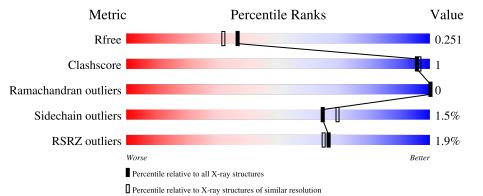
Xtriage (Phenix) EDS buster-report Percentile statistics	: : :	20191225.v01 (using entries in the PDB archive December 25th 2019)
-	:	
CCP4 Ideal geometry (proteins)		7.0.044 (Gargrove) Engh & Huber (2001)
Ideal geometry (DNA, RNA) Validation Pipeline (wwPDB-VP)		

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 2.01 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	$\begin{array}{c} \textbf{Whole archive} \\ \textbf{(\#Entries)} \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
R_{free}	130704	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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				
1	А	533	.% • 58%	•	40%	
1	В	533	% 59%	•	39%	



2 Entry composition (i)

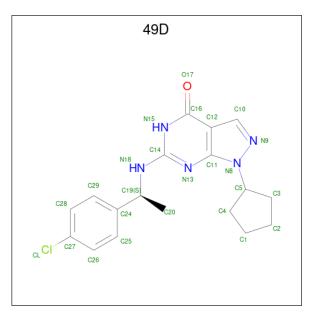
There are 6 unique types of molecules in this entry. The entry contains 5664 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 High affinity cGMP-specific 3',5'-cyclic phosphodiesterase 9A.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	Λ	322	Total	С	Ν	Ο	\mathbf{S}	0	0	0
	Л	522	2661	1704	442	485	30	0	0	0
1	В	326	Total	С	Ν	Ο	S	0	0	0
	D	520	2694	1727	446	491	30	0	U	

• Molecule 2 is $6-\{[(1S)-1-(4-chlorophenyl)ethyl]amino\}-1-cyclopentyl-1,5-dihydro-4H-pyrazol o[3,4-d]pyrimidin-4-one (three-letter code: 49D) (formula: C₁₈H₂₀ClN₅O).$



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf		
9	Δ	1	Total	С	Cl	Ν	0	0	0
	А	1	25	18	1	5	1	0	0

DE

• Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Ator	ns	ZeroOcc	AltConf
3	А	1	Total 1	Zn 1	0	0
				Co	ntinued on r	next page

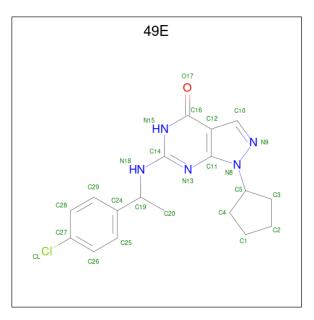
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	В	1	Total Z 1	Zn 1	0	0

• Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mo	Chain	Residues	Atoms	ZeroOcc	AltConf
4	А	1	Total Mg 1 1	0	0
4	В	1	Total Mg 1 1	0	0

• Molecule 5 is 6-{[(1R)-1-(4-chlorophenyl)ethyl]amino}-1-cyclopentyl-1,5-dihydro-4H-pyrazo lo[3,4-d]pyrimidin-4-one (three-letter code: 49E) (formula: $C_{18}H_{20}ClN_5O$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf		
5	В	1	Total 25	C 18	Cl 1	N 5	0 1	0	0

• Molecule 6 is water.

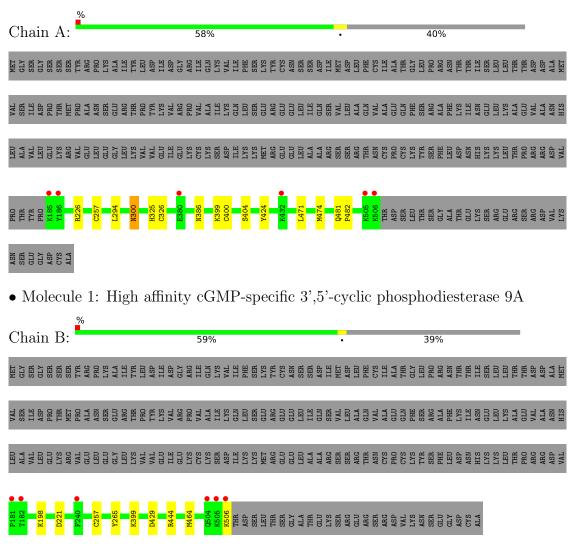
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	А	131	Total O 131 131	0	0
6	В	124	Total O 124 124	0	0



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: High affinity cGMP-specific 3',5'-cyclic phosphodiesterase 9A





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants	104.14Å 104.14Å 269.54Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	97.14 - 2.01	Depositor
Resolution (A)	48.57 - 2.01	EDS
% Data completeness	97.0 (97.14-2.01)	Depositor
(in resolution range)	97.0(48.57-2.01)	EDS
R _{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	$3.84 (at 2.01 \text{\AA})$	Xtriage
Refinement program	REFMAC 5.8.0073	Depositor
D D.	0.223 , 0.247	Depositor
R, R_{free}	0.228 , 0.251	DCC
R_{free} test set	4732 reflections $(4.89%)$	wwPDB-VP
Wilson B-factor $(Å^2)$	28.5	Xtriage
Anisotropy	0.143	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.35 , 32.7	EDS
L-test for twinning ²	$ \langle L \rangle = 0.43, \langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5664	wwPDB-VP
Average B, all atoms $(Å^2)$	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 6.56% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical 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.



¹Intensities 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: MG, 49E, ZN, 49D

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		
INIOI		RMSZ	# Z > 5	RMSZ	# Z > 5	
1	А	0.36	0/2728	0.53	0/3692	
1	В	0.35	0/2764	0.54	0/3743	
All	All	0.36	0/5492	0.53	0/7435	

There are no bond length outliers.

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	А	2661	0	2609	9	0
1	В	2694	0	2640	3	0
2	А	25	0	20	1	0
3	А	1	0	0	0	0
3	В	1	0	0	0	0
4	А	1	0	0	0	0
4	В	1	0	0	0	0
5	В	25	0	20	0	0
6	А	131	0	0	3	0
6	В	124	0	0	0	0
All	All	5664	0	5289	12	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 1.

All (12) 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:294:LEU:O	1:A:325:HIS:HD2	1.85	0.60
1:A:300:ASN:HD22	1:A:300:ASN:C	2.06	0.59
1:A:325:HIS:HE1	6:A:722:HOH:O	1.87	0.56
1:A:226:ARG:NH1	6:A:828:HOH:O	2.33	0.54
1:A:399:LYS:HE3	6:A:831:HOH:O	2.12	0.49
1:B:399:LYS:HE3	1:B:464:MET:HE3	1.97	0.46
1:B:429:ASP:OD2	1:B:444:ARG:NH1	2.49	0.46
1:A:400:CYS:O	1:A:404:SER:HB3	2.17	0.45
1:A:424:TYR:CD2	2:A:601:49D:H16	2.53	0.43
1:B:221:ASP:OD2	1:B:265:TYR:OH	2.31	0.41
1:A:300:ASN:C	1:A:300:ASN:ND2	2.72	0.41
1:A:481:GLN:HB3	1:A:482:PRO:CD	2.51	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 Percen		ntiles
1	А	320/533~(60%)	315~(98%)	5(2%)	0	100	100
1	В	324/533~(61%)	320 (99%)	4 (1%)	0	100	100
All	All	644/1066~(60%)	635~(99%)	9 (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	299/488~(61%)	293~(98%)	6(2%)	55 58		
1	В	303/488~(62%)	300 (99%)	3 (1%)	76 81		
All	All	602/976~(62%)	593~(98%)	9~(2%)	65 69		

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	А	257	CYS
1	А	300	ASN
1	А	326	CYS
1	А	386	ASN
1	А	471	LEU
1	А	474	MET
1	В	198	LYS
1	В	257	CYS
1	В	506	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	А	300	ASN
1	А	304	GLN
1	А	325	HIS
1	А	351	GLN
1	А	381	ASN
1	А	386	ASN
1	В	304	GLN
1	В	343	ASN
1	В	351	GLN

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)

Of 6 ligands modelled in this entry, 4 are 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).

Т	Mol Type		Chain	Dec	Timle	Bo	ond leng	\mathbf{ths}	B	ond ang	les
	101	туре	Chain	nes	Link	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z >2
	5	49E	В	601	-	25,28,28	2.66	5 (20%)	24,40,40	1.50	3 (12%)
	2	49D	А	601	-	25,28,28	2.56	5 (20%)	24,40,40	1.80	8 (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
5	49E	В	601	-	-	0/8/19/19	0/4/4/4
2	49D	А	601	-	-	2/8/19/19	0/4/4/4

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms		Observed(Å)	Ideal(Å)
5	В	601	49E	C12-C16	-7.91	1.34	1.47
2	А	601	49D	C12-C16	-7.79	1.34	1.47
5	В	601	49E	N9-N8	-7.24	1.24	1.36
2	А	601	49D	N9-N8	-5.61	1.27	1.36
2	А	601	49D	C12-C11	-5.13	1.33	1.41
5	В	601	49E	C12-C11	-4.68	1.34	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	А	601	49D	C24-C19	-4.66	1.40	1.52
5	В	601	49E	C24-C19	-4.25	1.41	1.52
2	А	601	49D	C27-CL	3.09	1.81	1.74
5	В	601	49E	C27-CL	2.07	1.79	1.74

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All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Ζ	$Observed(^{o})$	$Ideal(^{o})$
2	А	601	49D	C12-C16-N15	4.06	118.96	114.86
5	В	601	49E	C1-C4-C5	-3.93	95.63	104.32
2	А	601	49D	C2-C3-C5	-3.41	96.79	104.32
5	В	601	49E	C12-C16-N15	3.21	118.10	114.86
2	А	601	49D	C2-C1-C4	-2.92	96.17	105.99
2	А	601	49D	C24-C19-N18	-2.91	104.59	111.31
5	В	601	49E	C10-C12-C11	2.48	106.24	103.80
2	А	601	49D	O17-C16-N15	-2.19	118.07	120.65
2	А	601	49D	C4-C5-C3	2.13	107.15	104.25
2	А	601	49D	C10-C12-C11	2.10	105.87	103.80
2	А	601	49D	C3-C5-N8	-2.09	110.68	113.35

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	А	601	49D	C20-C19-C24-C29
2	А	601	49D	C20-C19-C24-C25

There are no ring outliers.

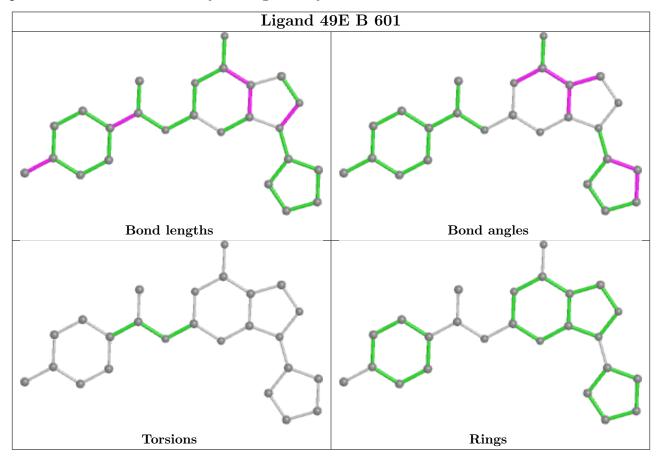
1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes	
2	А	601	49D	1	0	

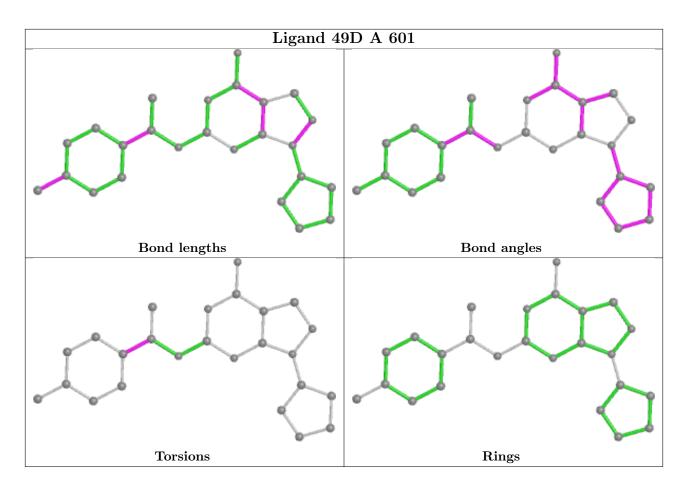
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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 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.







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	$\langle RSRZ \rangle$	#RSRZ>2	$\mathbf{OWAB}(\mathbf{A}^2)$	Q<0.9
1	А	322/533~(60%)	-0.18	6 (1%) 66 65	24, 36, 56, 103	0
1	В	326/533~(61%)	-0.11	6 (1%) 68 66	24, 38, 59, 118	0
All	All	648/1066~(60%)	-0.15	12 (1%) 66 65	24, 38, 57, 118	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	В	506	LYS	7.1
1	В	182	THR	6.9
1	В	505	LYS	5.1
1	А	186	TYR	4.5
1	В	181	PRO	4.3
1	А	505	LYS	2.6
1	А	432	LYS	2.5
1	А	506	LYS	2.2
1	А	380	GLU	2.2
1	А	185	LYS	2.2
1	В	240	PHE	2.1
1	В	504	GLN	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 monosaccharides in this entry.

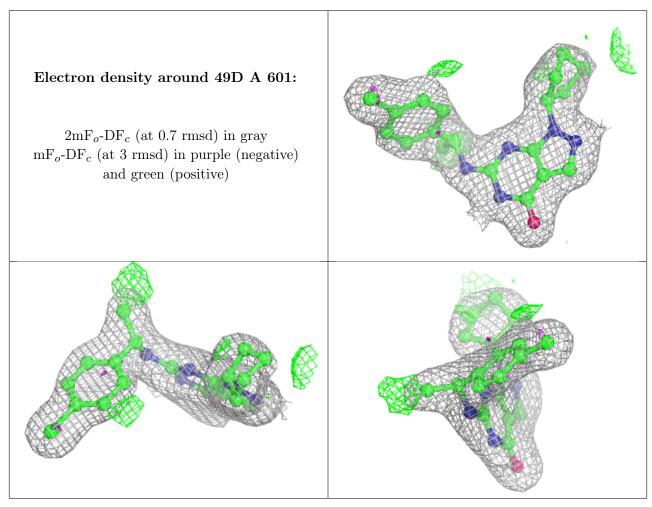


6.4 Ligands (i)

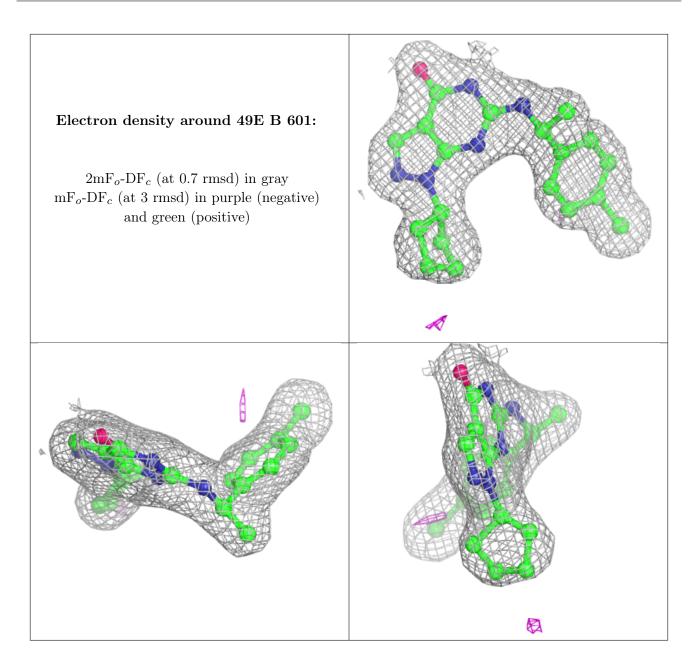
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9'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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-factors}(\mathrm{\AA}^2)$	Q<0.9
2	49D	А	601	25/25	0.95	0.13	$30,\!35,\!45,\!53$	0
4	MG	А	603	1/1	0.97	0.05	30,30,30,30	0
4	MG	В	603	1/1	0.97	0.09	32,32,32,32	0
5	49E	В	601	25/25	0.97	0.10	34,37,40,49	0
3	ZN	В	602	1/1	0.98	0.03	49,49,49,49	0
3	ZN	А	602	1/1	1.00	0.03	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.







6.5 Other polymers (i)

There are no such residues in this entry.

