

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

### Nov 25, 2024 - 07:50 PM EST

PDB ID	:	9EGF
Title	:	Crystal structure of Danio rerio histone deacetylase 6 catalytic domain 2 com-
		plexed with TO-589
Authors	:	Erdogan, F.; Seo, HS.; Dhe-Paganon, S.
Deposited on	:	2024-11-21
Resolution	:	1.25  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:

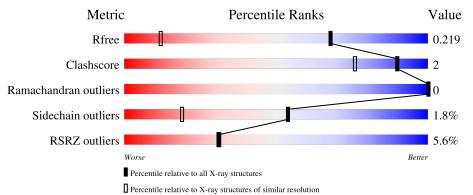
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.21
$\mathrm{EDS}$	:	3.0
buster-report		
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.004 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.40

# 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 1.25 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	$\begin{array}{c} \textbf{Whole archive} \\ (\#\textbf{Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
$R_{free}$	164625	1447 (1.28-1.24)
Clashscore	180529	1571 (1.28-1.24)
Ramachandran outliers	177936	1538 (1.28-1.24)
Sidechain outliers	177891	1537 (1.28-1.24)
RSRZ outliers	164620	1447 (1.28-1.24)

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	Quality of chain					
1	А	357	8%	•••				
1	В	357	<b>4%</b> 95%	5%				



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# 2 Entry composition (i)

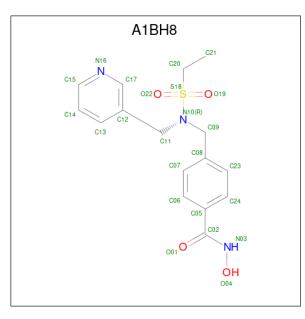
There are 5 unique types of molecules in this entry. The entry contains 6575 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Hdac6 protein.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	А	357	Total 2857	C 1792	N 517	O 529	S 19	0	9	0
1	В	357	Total 2862	C 1795	N 519	O 529	S 19	0	9	0

• Molecule 2 is 4-({(ethanesulfonyl)[(pyridin-3-yl)methyl]amino}methyl)-N-hydroxybenzami de (three-letter code: A1BH8) (formula:  $C_{16}H_{19}N_3O_4S$ ) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
2	А	1	Total 24	C 16			S 1	0	0
2	D	1	Total		0	-	I S	0	0
	Б	1	24	16	3	4	1	U	U

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



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

• Molecule 4 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	А	1	Total K 1 1	0	0
4	В	1	Total K 1 1	0	0

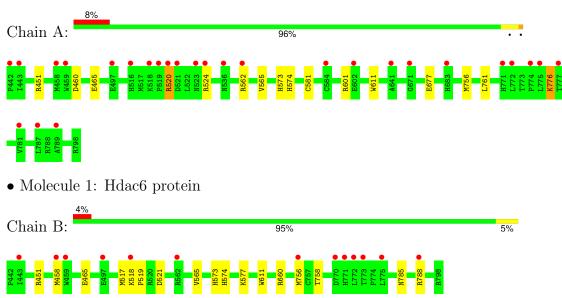
• Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	А	400	Total O 400 400	0	0
5	В	404	Total         O           404         404	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: Hdac6 protein



# 4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants	75.37Å 91.82Å 96.63Å	Depositor
a, b, c, $\alpha$ , $\beta$ , $\gamma$	$90.00^{\circ}$ $90.00^{\circ}$ $90.00^{\circ}$	Depositor
Resolution (Å)	66.56 - 1.25	Depositor
Resolution (A)	66.56 - 1.25	EDS
% Data completeness	$91.4\ (66.56\text{-}1.25)$	Depositor
(in resolution range)	$91.4\ (66.56\text{-}1.25)$	EDS
R <sub>merge</sub>	0.14	Depositor
$\mathrm{R}_{sym}$	(Not available)	Depositor
$< I/\sigma(I) > 1$	$1.18 (at 1.25 \text{\AA})$	Xtriage
Refinement program	PHENIX 1.20_4459	Depositor
R, $R_{free}$	0.200 , $0.221$	Depositor
n, n <sub>free</sub>	0.200 , $0.219$	DCC
$R_{free}$ test set	23710 reflections $(5.03%)$	wwPDB-VP
Wilson B-factor $(Å^2)$	12.3	Xtriage
Anisotropy	0.468	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$ , $B_{sol}(Å^2)$	0.30 , 24.8	EDS
L-test for twinning <sup>2</sup>	$< L >=0.49, < L^2>=0.32$	Xtriage
Estimated twinning fraction	0.000 for -h,l,k	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	6575	wwPDB-VP
Average B, all atoms $(Å^2)$	17.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The analyses of the Patterson function reveals a significant off-origin peak that is 48.13 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 8.9935e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.

<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: A1BH8, ZN, K

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		
	Unain	RMSZ	# Z  > 5	RMSZ	# Z  > 5	
1	А	0.33	0/2930	0.56	0/3970	
1	В	0.34	0/2936	0.56	0/3978	
All	All	0.34	0/5866	0.56	0/7948	

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	2857	0	2778	10	1
1	В	2862	0	2783	9	1
2	А	24	0	0	1	0
2	В	24	0	0	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	А	400	0	0	5	1
5	В	404	0	0	3	1
All	All	6575	0	5561	19	2



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 2.

All (19) 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:581:CYS:SG	5:A:1162:HOH:O	2.49	0.70
1:B:451[A]:ARG:NH2	5:B:902:HOH:O	2.26	0.67
1:A:677:GLU:OE2	1:A:776:LYS:HD2	1.95	0.66
1:A:776:LYS:HE2	5:A:994:HOH:O	2.02	0.60
1:A:520:ARG:NH1	5:A:906:HOH:O	2.36	0.58
1:B:518:LYS:NZ	5:B:907:HOH:O	2.35	0.57
1:A:601:ARG:NH2	5:A:907:HOH:O	2.37	0.56
1:A:451[A]:ARG:NH2	5:A:903:HOH:O	2.29	0.56
1:B:451[A]:ARG:NH1	5:B:910:HOH:O	2.36	0.56
1:B:785:ASN:OD1	1:B:788:ARG:NH2	2.41	0.54
1:A:574:HIS:NE2	2:A:801:A1BH8:N03	2.57	0.53
1:A:565[B]:VAL:HG23	1:A:761:LEU:HD12	1.90	0.53
1:B:574:HIS:NE2	2:B:801:A1BH8:N03	2.61	0.49
1:B:577:LYS:HE2	1:B:577:LYS:HB2	1.59	0.48
1:B:565[B]:VAL:HG21	1:B:758:THR:OG1	2.16	0.46
1:A:520:ARG:NH2	1:A:524:ARG:HH12	2.15	0.45
1:B:458:MET:HE1	1:B:519:PRO:HB3	2.00	0.43
1:A:565[B]:VAL:HG23	1:A:761:LEU:CD1	2.50	0.41
1:B:517:MET:HB3	1:B:521:ASP:HB2	2.04	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:460:ASP:OD2	1:B:660[B]:ARG:NH2[2_654]	2.05	0.15
5:A:1151:HOH:O	5:B:1127:HOH:O[4_455]	2.19	0.01

## 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



Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percer	ntiles
1	А	364/357~(102%)	355~(98%)	9(2%)	0	100	100
1	В	364/357~(102%)	355~(98%)	9 (2%)	0	100	100
All	All	728/714 (102%)	710 (98%)	18 (2%)	0	100	100

analysed, and the total number of residues.

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent side chain 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	А	309/301~(103%)	301~(97%)	8(3%)	41 7		
1	В	310/301~(103%)	305~(98%)	5(2%)	58 23		
All	All	619/602~(103%)	606~(98%)	13 (2%)	54 14		

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

Mol	Chain	Res	Type
1	А	465	GLU
1	А	520	ARG
1	А	562	ARG
1	А	573	HIS
1	А	611	TRP
1	А	756[A]	MET
1	А	756[B]	MET
1	А	776	LYS
1	В	465	GLU
1	В	573	HIS
1	В	611	TRP
1	В	756[A]	MET
1	В	756[B]	MET

Sometimes side chains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such side chains are listed below:



Mol	Chain	Res	Type
1	В	560	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 oligosaccharides 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).

Mol	Turne	Chain	Res	Link	Bo	ond leng	ths	В	ond ang	les
IVIOI	Type	Unam	nes	LIIIK	Counts	RMSZ	# Z >2	Counts	RMSZ	# Z >2
2	A1BH8	А	801	3	$25,\!25,\!25$	2.73	6 (24%)	33,34,34	<mark>3.11</mark>	7 (21%)
2	A1BH8	В	801	3	$25,\!25,\!25$	2.87	6 (24%)	33,34,34	<mark>3.39</mark>	7 (21%)

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	A1BH8	А	801	3	-	2/23/23/23	0/2/2/2
2	A1BH8	В	801	3	-	5/23/23/23	0/2/2/2



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Mol	Chain	Res	Type	Atoms	Ζ	Observed(Å)	Ideal(Å)
2	А	801	A1BH8	C02-N03	9.95	1.45	1.32
2	В	801	A1BH8	C02-N03	9.66	1.45	1.32
2	В	801	A1BH8	C20-S18	7.09	1.82	1.76
2	А	801	A1BH8	C20-S18	5.98	1.81	1.76
2	В	801	A1BH8	S18-N10	5.68	1.71	1.63
2	А	801	A1BH8	S18-N10	4.43	1.69	1.63
2	В	801	A1BH8	O19-S18	3.12	1.46	1.43
2	А	801	A1BH8	O19-S18	2.84	1.46	1.43
2	А	801	A1BH8	O22-S18	2.71	1.46	1.43
2	А	801	A1BH8	O01-C02	-2.57	1.17	1.23
2	В	801	A1BH8	O22-S18	2.43	1.46	1.43
2	В	801	A1BH8	O01-C02	-2.42	1.17	1.23

All (12) bond length outliers are listed below:

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
2	В	801	A1BH8	O22-S18-O19	-16.98	101.89	118.99
2	А	801	A1BH8	O22-S18-O19	-15.14	103.74	118.99
2	В	801	A1BH8	O19-S18-N10	4.72	111.88	107.28
2	А	801	A1BH8	O22-S18-N10	4.17	111.34	107.28
2	В	801	A1BH8	C15-N16-C17	3.40	122.81	116.85
2	А	801	A1BH8	O19-S18-N10	3.34	110.53	107.28
2	А	801	A1BH8	C15-N16-C17	3.31	122.64	116.85
2	В	801	A1BH8	C09-N10-C11	3.01	122.44	116.27
2	А	801	A1BH8	O01-C02-N03	-2.91	117.54	122.90
2	В	801	A1BH8	O01-C02-N03	-2.81	117.72	122.90
2	В	801	A1BH8	C05-C02-N03	2.54	120.17	116.26
2	В	801	A1BH8	O19-S18-C20	2.51	112.09	108.27
2	А	801	A1BH8	C05-C02-N03	2.43	120.01	116.26
2	А	801	A1BH8	C09-N10-C11	2.21	120.80	116.27

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	В	801	A1BH8	C21-C20-S18-N10
2	В	801	A1BH8	C21-C20-S18-O19
2	В	801	A1BH8	C21-C20-S18-O22
2	А	801	A1BH8	C12-C11-N10-C09
2	В	801	A1BH8	C12-C11-N10-C09
2	А	801	A1BH8	C12-C11-N10-S18

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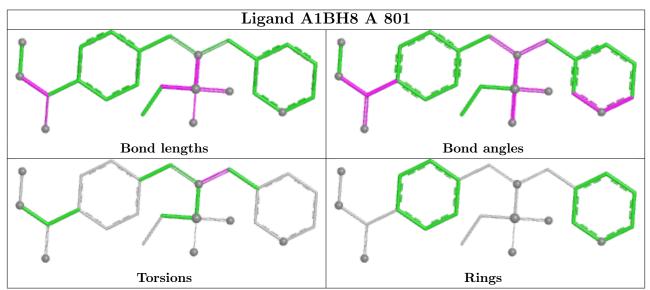
Mol	Chain	Res	Type	Atoms
2	В	801	A1BH8	C12-C11-N10-S18

There are no ring outliers.

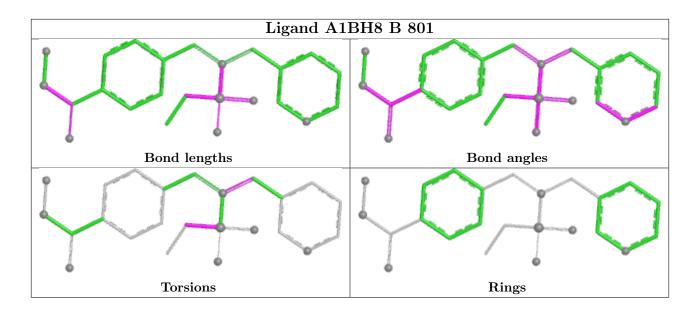
2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	А	801	A1BH8	1	0
2	В	801	A1BH8	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	<RSRZ $>$	# RSRZ > 2	$OWAB(Å^2)$	Q<0.9
1	А	357/357~(100%)	0.52	27 (7%) 21 21	5, 14, 28, 38	11 (3%)
1	В	357/357~(100%)	0.37	13 (3%) 46 46	5, 14, 24, 33	12 (3%)
All	All	714/714~(100%)	0.45	40 (5%) 31 31	5, 14, 25, 38	23 (3%)

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	В	772	LEU	4.9
1	А	459	TRP	4.2
1	А	519	PRO	3.9
1	В	459	TRP	3.6
1	А	772	LEU	3.4
1	В	771	HIS	3.3
1	А	442	PRO	3.1
1	А	518	LYS	3.1
1	А	520	ARG	3.1
1	А	775	LEU	3.0
1	В	497	GLU	3.0
1	А	521	ASP	3.0
1	А	443	ILE	3.0
1	В	775	LEU	2.9
1	А	777	THR	2.9
1	В	770	ASP	2.9
1	А	771	HIS	2.9
1	В	773	THR	2.8
1	В	756[A]	MET	2.7
1	А	787	LEU	2.7
1	А	774	PRO	2.7
1	В	443	ILE	2.6
1	А	458	MET	2.5
1	А	671	GLY	2.5

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Mol	Chain	Res	Type	RSRZ
1	А	683	HIS	2.5
1	В	458	MET	2.5
1	А	536	ASN	2.4
1	В	788	ARG	2.4
1	А	524	ARG	2.3
1	А	523	ASN	2.3
1	А	602	GLU	2.3
1	В	562	ARG	2.2
1	А	584	CYS	2.2
1	А	641	ALA	2.1
1	А	516	HIS	2.1
1	А	497	GLU	2.1
1	А	789	ALA	2.1
1	В	518	LYS	2.1
1	А	562	ARG	2.0
1	А	781	VAL	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.

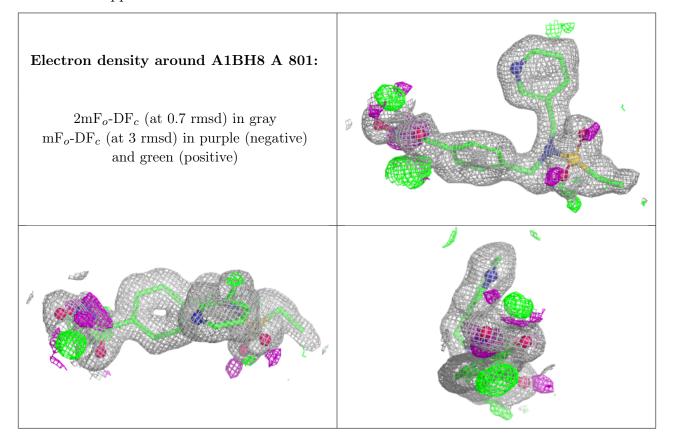
## 6.4 Ligands (i)

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.

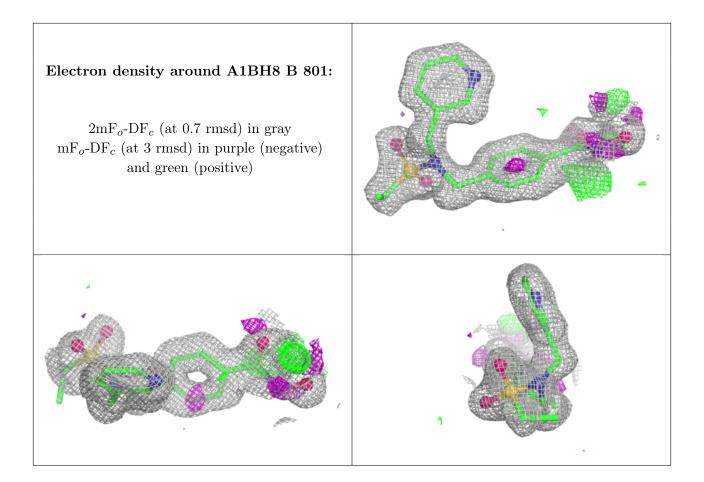
Mol	Type	Chain	Res	Atoms	RSCC	RSR	$B-factors(Å^2)$	Q < 0.9
2	A1BH8	А	801	24/24	0.93	0.10	$13,\!18,\!23,\!25$	0
2	A1BH8	В	801	24/24	0.93	0.10	13,17,25,31	0
4	Κ	А	803	1/1	0.99	0.04	$14,\!14,\!14,\!14$	0
4	К	В	803	1/1	0.99	0.03	11,11,11,11	1
3	ZN	А	802	1/1	1.00	0.05	$15,\!15,\!15,\!15$	0
3	ZN	В	802	1/1	1.00	0.05	$15,\!15,\!15,\!15$	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.

