



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 10, 2026 – 07:20 PM UTC

PDB ID : 6B97 / pdb_00006b97
Title : Crystal structure of PDE2 in complex with complex 9
Authors : Lu, J.
Deposited on : 2017-10-10
Resolution : 1.76 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 ⓘ 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 ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

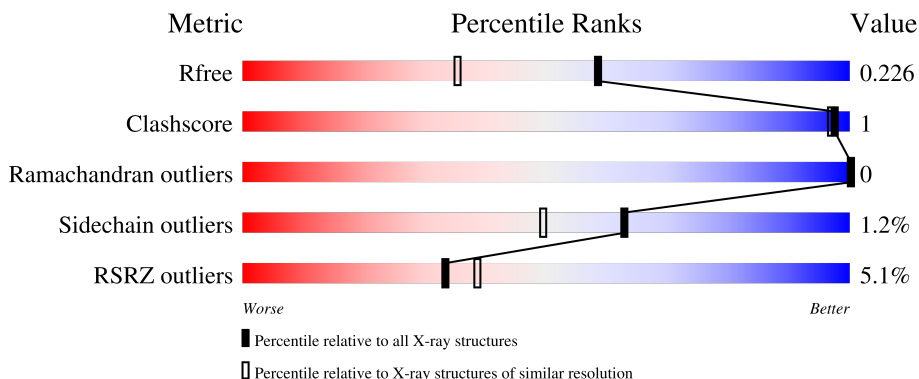
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.76 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}	180053	3183 (1.76-1.76)
Clashscore	190562	3299 (1.76-1.76)
Ramachandran outliers	187476	3274 (1.76-1.76)
Sidechain outliers	187428	3274 (1.76-1.76)
RSRZ outliers	180081	3183 (1.76-1.76)

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 $\leq 5\%$. 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	6% 87% 5% 8%
1	B	373	4% 90% • 7%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6008 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 cGMP-dependent 3',5'-cyclic phosphodiesterase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	343	Total	C	N	O	S	0	4	0
			2752	1755	462	509	26			
1	B	348	Total	C	N	O	S	0	6	0
			2844	1813	484	520	27			

There are 62 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	547	MET	-	initiating methionine	UNP O00408
A	548	HIS	-	expression tag	UNP O00408
A	549	HIS	-	expression tag	UNP O00408
A	550	HIS	-	expression tag	UNP O00408
A	551	HIS	-	expression tag	UNP O00408
A	552	HIS	-	expression tag	UNP O00408
A	553	HIS	-	expression tag	UNP O00408
A	554	GLU	-	expression tag	UNP O00408
A	555	ASN	-	expression tag	UNP O00408
A	556	LEU	-	expression tag	UNP O00408
A	557	TYR	-	expression tag	UNP O00408
A	558	PHE	-	expression tag	UNP O00408
A	559	GLN	-	expression tag	UNP O00408
A	560	GLY	-	expression tag	UNP O00408
A	561	GLU	-	expression tag	UNP O00408
A	562	LEU	-	expression tag	UNP O00408
A	563	SER	-	expression tag	UNP O00408
A	564	THR	-	expression tag	UNP O00408
A	565	SER	-	expression tag	UNP O00408
A	566	LEU	-	expression tag	UNP O00408
A	567	TYR	-	expression tag	UNP O00408
A	568	LYS	-	expression tag	UNP O00408
A	569	LYS	-	expression tag	UNP O00408
A	570	ALA	-	expression tag	UNP O00408
A	571	GLY	-	expression tag	UNP O00408

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Chain	Residue	Modelled	Actual	Comment	Reference
A	572	PHE	-	expression tag	UNP O00408
A	573	ASP	-	expression tag	UNP O00408
A	574	ASP	-	expression tag	UNP O00408
A	575	ASP	-	expression tag	UNP O00408
A	576	ASP	-	expression tag	UNP O00408
A	577	LYS	-	expression tag	UNP O00408
B	547	MET	-	initiating methionine	UNP O00408
B	548	HIS	-	expression tag	UNP O00408
B	549	HIS	-	expression tag	UNP O00408
B	550	HIS	-	expression tag	UNP O00408
B	551	HIS	-	expression tag	UNP O00408
B	552	HIS	-	expression tag	UNP O00408
B	553	HIS	-	expression tag	UNP O00408
B	554	GLU	-	expression tag	UNP O00408
B	555	ASN	-	expression tag	UNP O00408
B	556	LEU	-	expression tag	UNP O00408
B	557	TYR	-	expression tag	UNP O00408
B	558	PHE	-	expression tag	UNP O00408
B	559	GLN	-	expression tag	UNP O00408
B	560	GLY	-	expression tag	UNP O00408
B	561	GLU	-	expression tag	UNP O00408
B	562	LEU	-	expression tag	UNP O00408
B	563	SER	-	expression tag	UNP O00408
B	564	THR	-	expression tag	UNP O00408
B	565	SER	-	expression tag	UNP O00408
B	566	LEU	-	expression tag	UNP O00408
B	567	TYR	-	expression tag	UNP O00408
B	568	LYS	-	expression tag	UNP O00408
B	569	LYS	-	expression tag	UNP O00408
B	570	ALA	-	expression tag	UNP O00408
B	571	GLY	-	expression tag	UNP O00408
B	572	PHE	-	expression tag	UNP O00408
B	573	ASP	-	expression tag	UNP O00408
B	574	ASP	-	expression tag	UNP O00408
B	575	ASP	-	expression tag	UNP O00408
B	576	ASP	-	expression tag	UNP O00408
B	577	LYS	-	expression tag	UNP O00408

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

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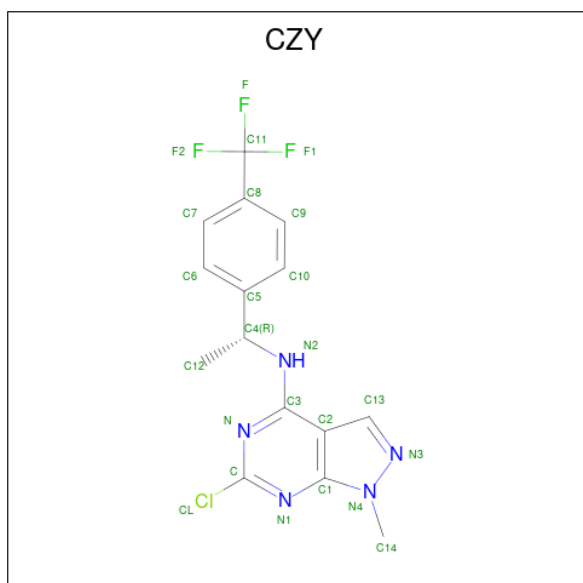
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Zn 1 1	0	0
2	B	1	Total Zn 1 1	0	0

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Mg 1 1	0	0
3	B	1	Total Mg 1 1	0	0

- Molecule 4 is 6-chloro-1-methyl-N-{(1R)-1-[4-(trifluoromethyl)phenyl]ethyl}-1H-pyrazolo[3,4-d]pyrimidin-4-amine (CCD ID: CZY) (formula: C₁₅H₁₃ClF₃N₅).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total 24	C 15	Cl 1	F 3	N 5	0	0
4	B	1	Total 48	C 30	Cl 2	F 6	N 10	0	1

- Molecule 5 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).

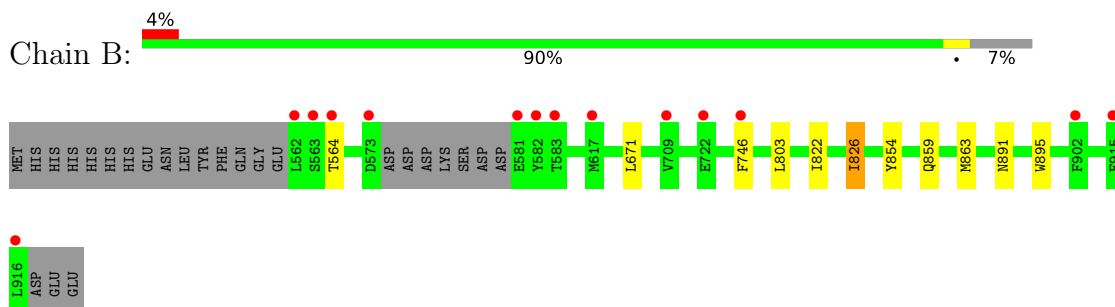


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	155	Total	O	0	0
			155	155		
6	B	173	Total	O	0	0
			173	173		

- Molecule 1: cGMP-dependent 3',5'-cyclic phosphodiesterase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	72.27Å 96.07Å 102.06Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	34.06 – 1.76 34.06 – 1.76	Depositor EDS
% Data completeness (in resolution range)	99.9 (34.06-1.76) 99.9 (34.06-1.76)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.37 (at 1.76Å)	Xtriage
Refinement program	BUSTER 2.11.7	Depositor
R, R_{free}	0.190 , 0.220 0.195 , 0.226	Depositor DCC
R_{free} test set	3584 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	24.4	Xtriage
Anisotropy	0.521	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 32.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6008	wwPDB-VP
Average B, all atoms (Å ²)	31.0	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 20.91 % 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 7.8463e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

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.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CZY, EDO, MG, ZN

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	0.85	0/2818	1.23	4/3816 (0.1%)
1	B	0.85	0/2915	1.17	3/3937 (0.1%)
All	All	0.85	0/5733	1.20	7/7753 (0.1%)

There are no bond length outliers.

The worst 5 of 7 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	854	TYR	CA-C-N	7.46	127.59	120.43
1	B	854	TYR	C-N-CA	7.46	127.59	120.43
1	B	746	PHE	CA-CB-CG	7.44	121.24	113.80
1	A	902	PHE	CA-CB-CG	6.75	120.55	113.80
1	A	845	MET	N-CA-C	-5.76	102.52	110.35

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	2752	0	2610	4	0
1	B	2844	0	2746	5	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	24	0	0	0	0
4	B	48	0	0	0	0
5	A	8	0	12	0	0
6	A	155	0	0	0	0
6	B	173	0	0	0	0
All	All	6008	0	5368	9	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.

The worst 5 of 9 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:817:LYS:HD3	1:A:820:ARG:HH21	1.73	0.54
1:B:822:ILE:HG23	1:B:826:ILE:HD12	1.96	0.47
1:A:859:GLN:HG2	1:A:895:TRP:CE2	2.52	0.45
1:A:645:MET:HG3	1:A:743:CYS:SG	2.59	0.43
1:B:863:MET:HE1	1:B:891:ASN:HB2	2.01	0.43

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	343/373 (92%)	336 (98%)	7 (2%)	0	100	100
1	B	350/373 (94%)	344 (98%)	6 (2%)	0	100	100
All	All	693/746 (93%)	680 (98%)	13 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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	294/337 (87%)	289 (98%)	5 (2%)	53	36
1	B	309/337 (92%)	307 (99%)	2 (1%)	78	71
All	All	603/674 (90%)	596 (99%)	7 (1%)	63	49

5 of 7 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	826	ILE
1	A	898	VAL
1	B	826	ILE
1	B	564	THR
1	A	785	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 8 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	733	GLN
1	B	708	GLN
1	B	591	GLN
1	A	708	GLN
1	B	624	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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 4 are monoatomic - leaving 5 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
5	EDO	A	1005	-	3,3,3	0.54	0	2,2,2	0.53	0
5	EDO	A	1004	-	3,3,3	0.43	0	2,2,2	0.58	0
4	CZY	B	1003[B]	-	26,26,26	0.38	0	33,39,39	0.75	2 (6%)
4	CZY	A	1003	-	26,26,26	0.37	0	33,39,39	0.53	1 (3%)
4	CZY	B	1003[A]	-	26,26,26	0.40	0	33,39,39	0.49	1 (3%)

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	EDO	A	1005	-	-	0/1/1/1	-
5	EDO	A	1004	-	-	0/1/1/1	-
4	CZY	B	1003[B]	-	-	0/14/23/23	0/3/3/3
4	CZY	A	1003	-	-	4/14/23/23	0/3/3/3
4	CZY	B	1003[A]	-	-	5/14/23/23	0/3/3/3

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1003[B]	CZY	C14-N4-C1	3.11	130.49	128.30
4	A	1003	CZY	C14-N4-C1	2.31	129.92	128.30
4	B	1003[B]	CZY	N2-C3-N	2.25	122.88	118.55

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1003[A]	CZY	C14-N4-C1	2.01	129.71	128.30

There are no chirality outliers.

5 of 9 torsion outliers are listed below:

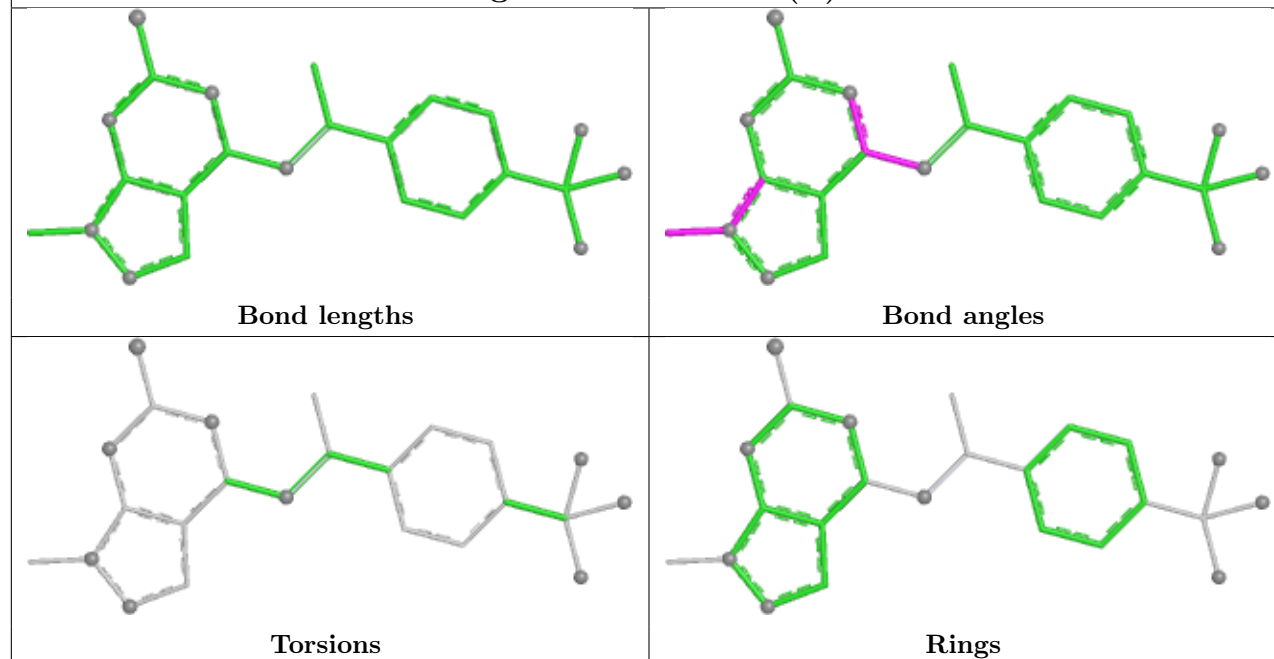
Mol	Chain	Res	Type	Atoms
4	A	1003	CZY	C2-C3-N2-C4
4	B	1003[A]	CZY	C2-C3-N2-C4
4	A	1003	CZY	N-C3-N2-C4
4	B	1003[A]	CZY	N-C3-N2-C4
4	A	1003	CZY	N2-C4-C5-C10

There are no ring outliers.

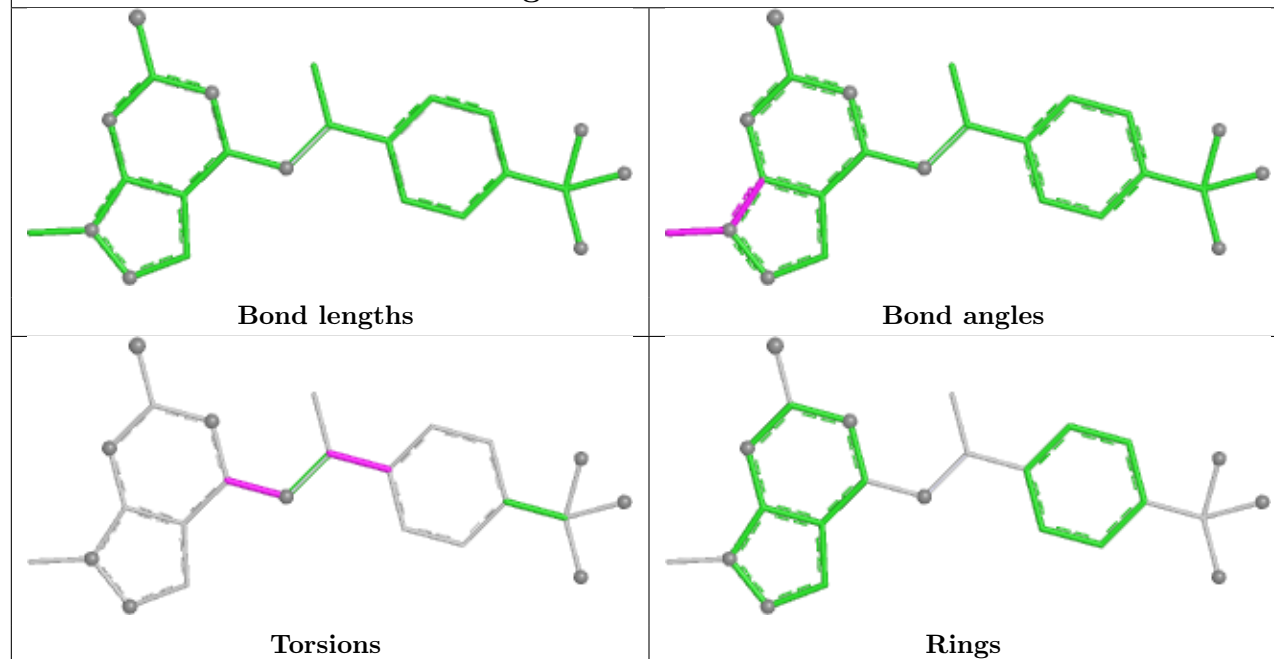
No monomer is involved in short contacts.

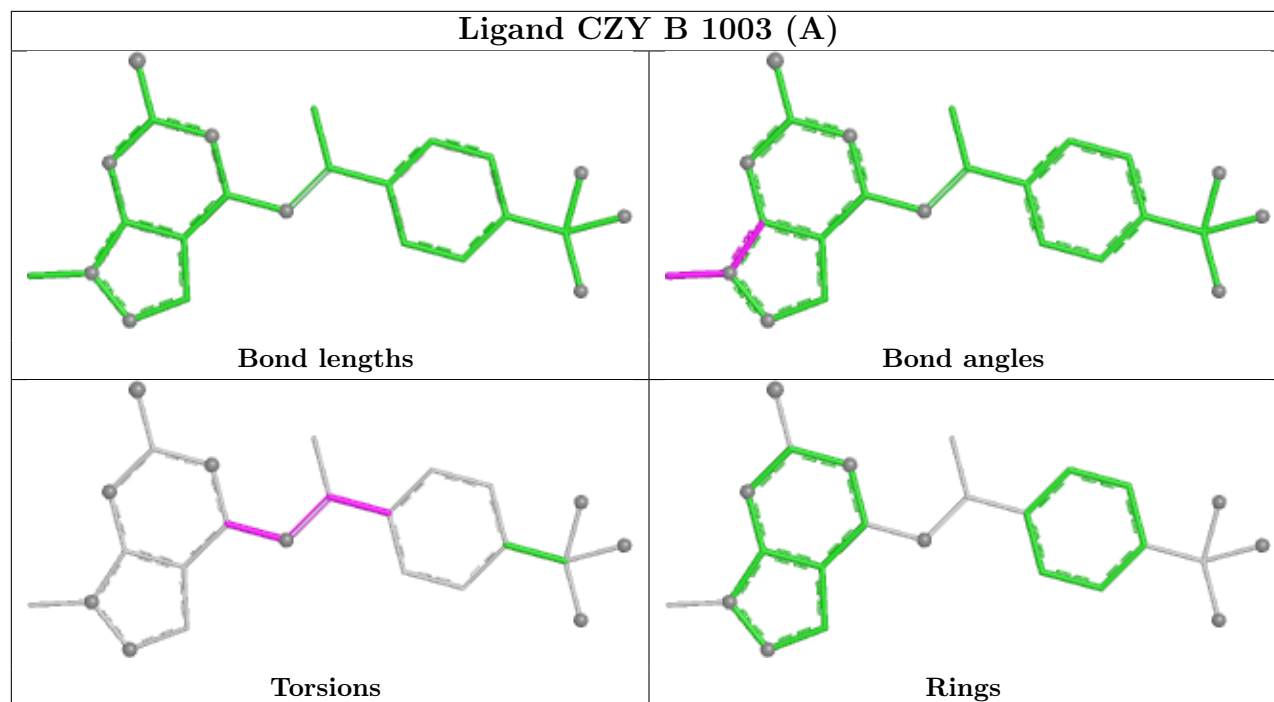
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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 CZY B 1003 (B)



Ligand CZY A 1003





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, 95th 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	A	343/373 (91%)	0.41	21 (6%)	27 31	10, 29, 55, 69	4 (1%)
1	B	348/373 (93%)	0.18	14 (4%)	42 48	10, 28, 48, 78	6 (1%)
All	All	691/746 (92%)	0.30	35 (5%)	33 39	10, 28, 52, 78	10 (1%)

The worst 5 of 35 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	562	LEU	6.3
1	B	916	LEU	5.6
1	A	588	ASP	4.6
1	A	916	LEU	4.5
1	B	563	SER	4.0

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](#)

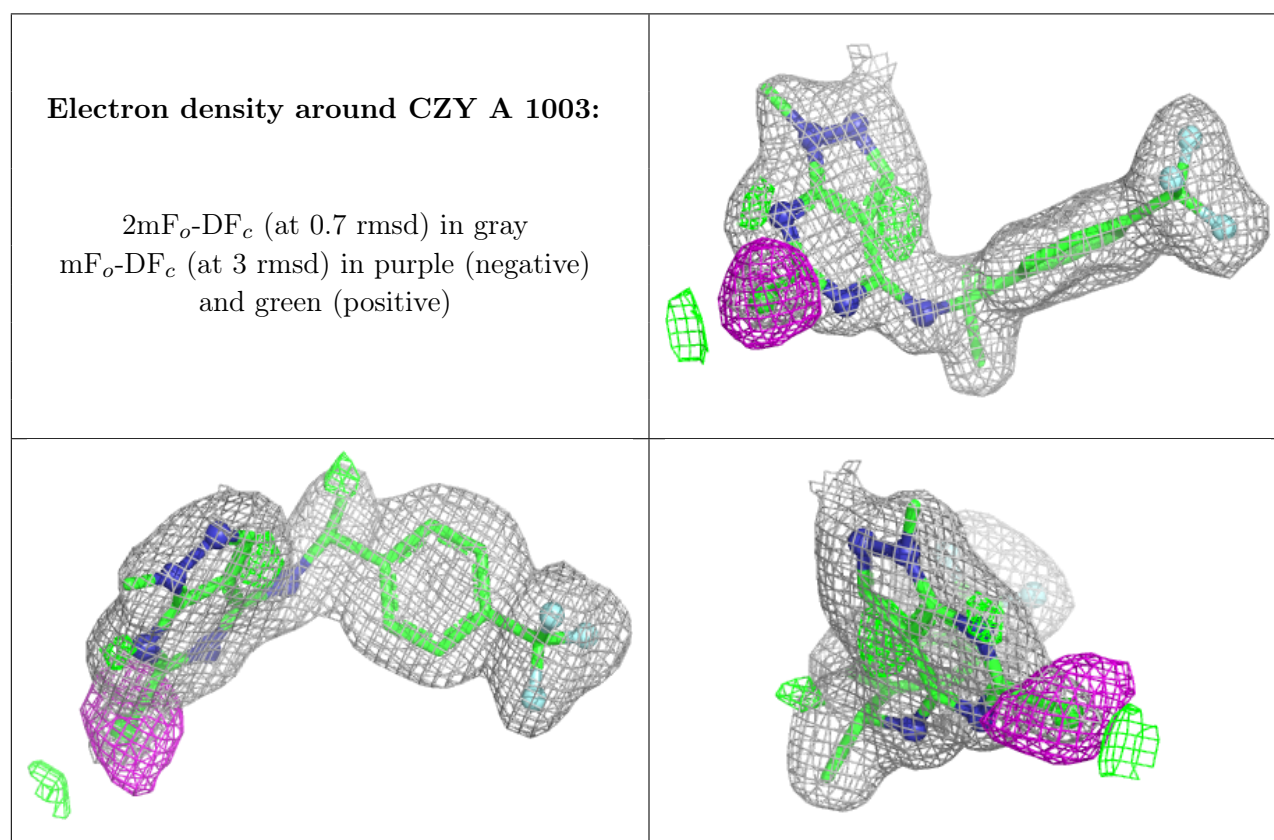
There are no oligosaccharides in this entry.

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, 95th 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.

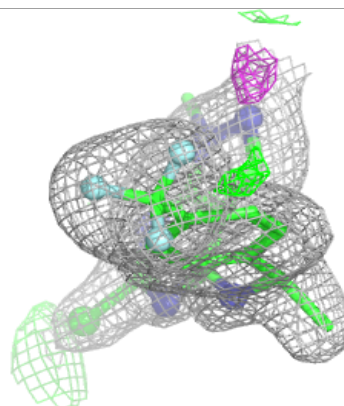
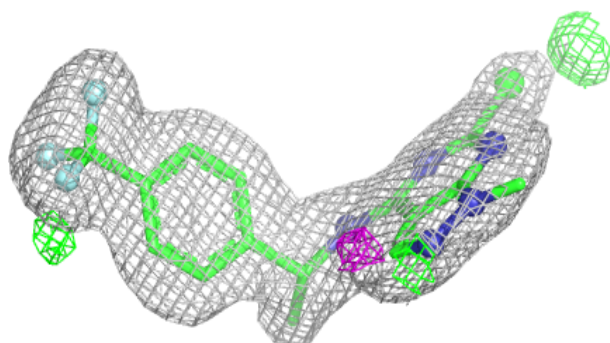
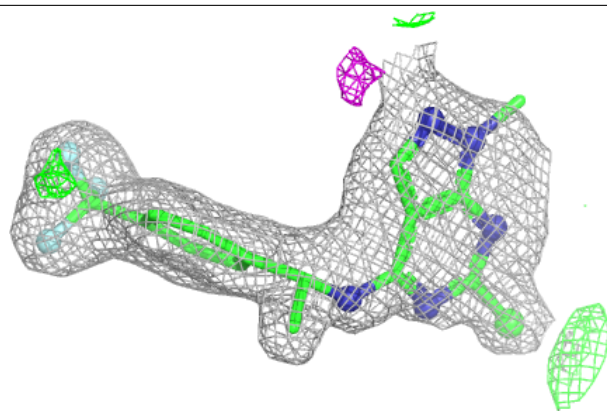
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	EDO	A	1004	4/4	0.81	0.14	43,46,46,47	0
4	CZY	A	1003	24/24	0.86	0.10	22,32,38,56	0
5	EDO	A	1005	4/4	0.89	0.11	38,38,39,39	0
4	CZY	B	1003[A]	24/24	0.90	0.08	20,25,31,35	24
4	CZY	B	1003[B]	24/24	0.90	0.08	31,37,41,43	24
3	MG	A	1002	1/1	0.98	0.03	18,18,18,18	0
3	MG	B	1002	1/1	0.99	0.03	19,19,19,19	0
2	ZN	B	1001	1/1	1.00	0.03	26,26,26,26	0
2	ZN	A	1001	1/1	1.00	0.02	25,25,25,25	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.

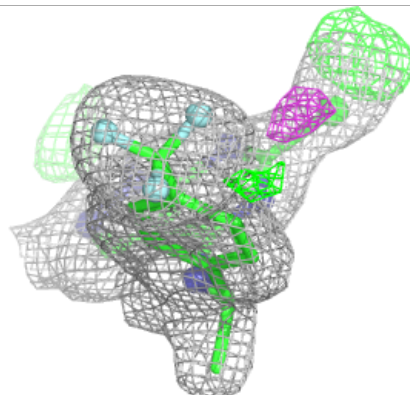
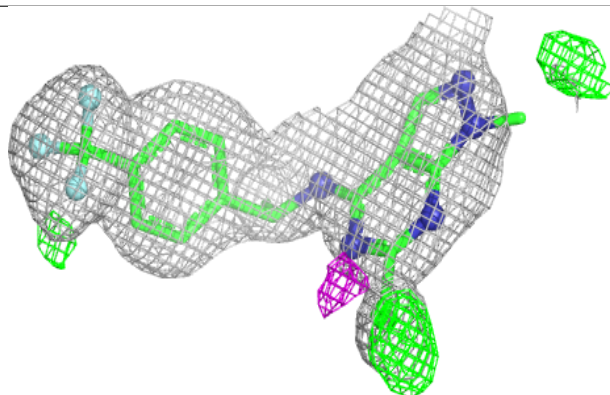
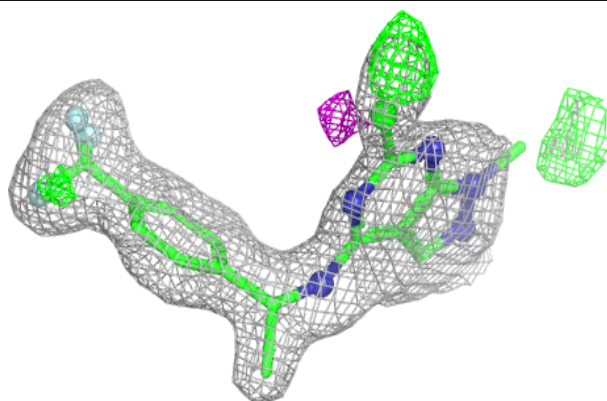


Electron density around CZY B 1003 (A):

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

**Electron density around CZY B 1003 (B):**

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



6.5 Other polymers [i](#)

There are no such residues in this entry.