



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 02:56 PM UTC

PDB ID : 6KDX / pdb_00006kdx
Title : Crystal structure of PDE10A in complex with a triazolopyrimidine inhibitor
Authors : Amano, Y.; Honbou, K.
Deposited on : 2019-07-03
Resolution : 2.44 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 ⓘ 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

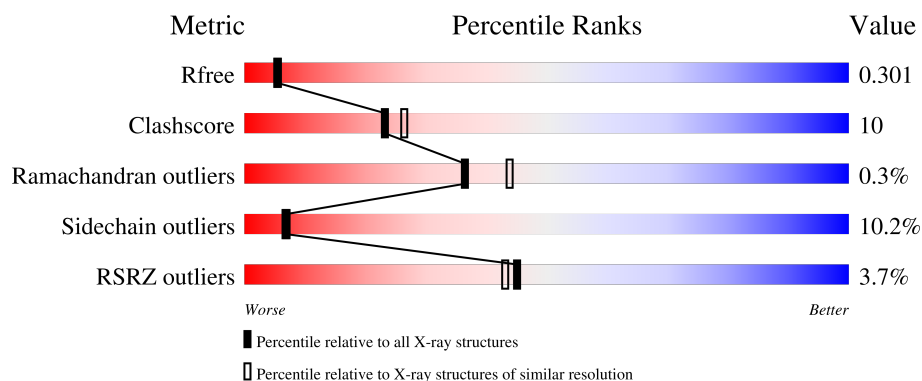
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.44 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	2340 (2.46-2.42)
Clashscore	190562	2400 (2.46-2.42)
Ramachandran outliers	187476	2379 (2.46-2.42)
Sidechain outliers	187428	2379 (2.46-2.42)
RSRZ outliers	180081	2340 (2.46-2.42)

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	345	4% 64% 25% • 6%
1	B	345	3% 67% 23% • 6%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5306 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 cAMP and cAMP-inhibited cGMP 3',5'-cyclic phosphodiesterase 10A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	324	Total	C	N	O	S	0	0	0
			2629	1678	447	479	25			
1	B	324	Total	C	N	O	S	0	0	0
			2629	1678	447	479	25			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	445	GLY	-	expression tag	UNP Q9Y233
A	446	SER	-	expression tag	UNP Q9Y233
A	447	HIS	-	expression tag	UNP Q9Y233
A	448	MET	-	expression tag	UNP Q9Y233
B	445	GLY	-	expression tag	UNP Q9Y233
B	446	SER	-	expression tag	UNP Q9Y233
B	447	HIS	-	expression tag	UNP Q9Y233
B	448	MET	-	expression tag	UNP Q9Y233

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

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

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

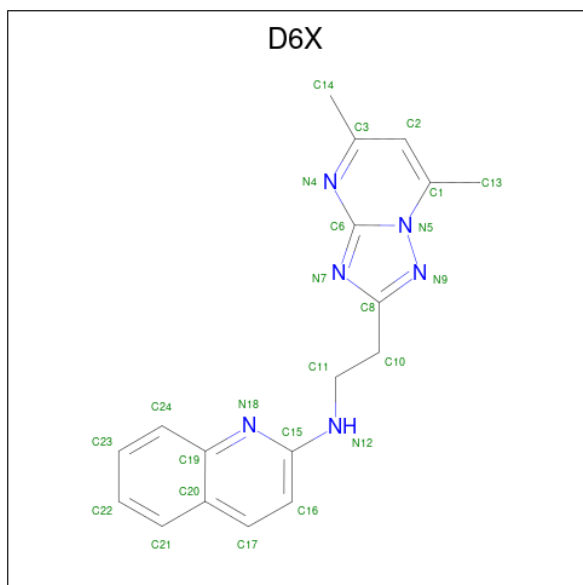
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mg	0	0
			1	1		

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

- Molecule 4 is N-[2-(5,7-dimethyl-[1,2,4]triazolo[1,5-a]pyrimidin-2-yl)ethyl]quinolin-2-amine (CCD ID: D6X) (formula: C₁₈H₁₈N₆) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	N	0	0
			24	18	6		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	11	Total	O	0	0
			11	11		
5	B	9	Total	O	0	0
			9	9		

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	50.42Å 81.44Å 159.98Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	26.66 – 2.44 26.66 – 2.44	Depositor EDS
% Data completeness (in resolution range)	91.1 (26.66-2.44) 91.2 (26.66-2.44)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.14 (at 2.44Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.241 , 0.301 0.241 , 0.301	Depositor DCC
R_{free} test set	1187 reflections (4.68%)	wwPDB-VP
Wilson B-factor (Å ²)	49.7	Xtriage
Anisotropy	0.150	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 27.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5306	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.95% of the height of the origin peak. No significant pseudotranslation is detected.*

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: MG, ZN, D6X

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	1.16	7/2694 (0.3%)	1.24	22/3647 (0.6%)
1	B	1.05	5/2694 (0.2%)	1.16	6/3647 (0.2%)
All	All	1.11	12/5388 (0.2%)	1.20	28/7294 (0.4%)

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	545	HIS	CE1-NE2	24.46	1.57	1.32
1	A	620	GLU	C-O	8.72	1.38	1.24
1	A	580	HIS	CE1-NE2	8.36	1.41	1.32
1	A	545	HIS	CG-ND1	7.52	1.46	1.38
1	A	580	HIS	ND1-CE1	7.28	1.39	1.32
1	B	577	LYS	CE-NZ	5.36	1.65	1.49
1	B	567	HIS	CG-CD2	5.23	1.41	1.35
1	B	563	HIS	CG-ND1	-5.10	1.32	1.38
1	A	579	ASP	CG-OD2	5.09	1.35	1.25
1	B	697	TRP	CD2-CE2	5.06	1.50	1.41
1	B	580	HIS	CG-CD2	5.04	1.41	1.35
1	A	536	CYS	CA-C	5.02	1.59	1.52

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	578	PHE	CB-CA-C	-10.67	93.00	111.35
1	A	621	GLN	N-CA-C	-10.12	99.72	113.18
1	A	484	ASN	CB-CA-C	-9.64	91.23	110.42
1	A	580	HIS	ND1-CE1-NE2	-9.25	99.15	108.40
1	A	578	PHE	CA-C-N	8.16	136.39	121.70
1	A	578	PHE	C-N-CA	8.16	136.39	121.70
1	A	579	ASP	N-CA-C	-7.15	90.99	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	578	PHE	N-CA-C	7.07	125.85	110.80
1	A	467	ARG	NE-CZ-NH2	6.97	125.48	119.20
1	B	579	ASP	N-CA-C	6.72	120.35	111.28
1	A	545	HIS	CE1-NE2-CD2	-6.71	102.29	109.00
1	B	563	HIS	N-CA-C	6.71	119.42	111.71
1	A	769	GLU	N-CA-C	6.00	118.55	110.88
1	B	621	GLN	N-CA-C	-5.74	106.82	113.88
1	A	578	PHE	CA-CB-CG	5.65	119.45	113.80
1	A	746	PRO	N-CA-C	5.44	117.34	110.70
1	A	563	HIS	N-CA-C	5.38	117.90	111.71
1	A	649	TYR	N-CA-C	-5.33	104.90	111.40
1	A	767	ARG	NE-CZ-NH2	5.32	123.99	119.20
1	B	635	LEU	CD1-CG-CD2	-5.28	99.17	110.80
1	A	486	TRP	CA-C-N	-5.28	113.09	118.85
1	A	486	TRP	C-N-CA	-5.28	113.09	118.85
1	A	650	GLN	CA-CB-CG	-5.20	103.71	114.10
1	A	635	LEU	CD1-CG-CD2	-5.17	99.42	110.80
1	A	728	GLY	CA-C-N	5.15	127.14	120.44
1	A	728	GLY	C-N-CA	5.15	127.14	120.44
1	B	767	ARG	NE-CZ-NH2	5.15	123.84	119.20
1	A	580	HIS	CE1-NE2-CD2	5.09	114.09	109.00

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	2629	0	2588	61	1
1	B	2629	0	2588	58	1
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	24	0	0	0	0
5	A	11	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	B	9	0	0	0	0
All	All	5306	0	5176	108	1

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 10.

All (108) 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:580:HIS:ND1	5:A:901:HOH:O	1.68	1.26
1:A:724:GLN:OE1	1:B:683:PRO:HD3	1.33	1.26
1:A:724:GLN:HG3	1:B:683:PRO:HG3	1.30	1.08
1:A:641:ASN:HD22	1:A:664:ARG:HH11	1.09	0.99
1:A:724:GLN:OE1	1:B:682:TRP:HB3	1.64	0.97
1:B:523:PRO:HD2	1:B:695:GLU:HG3	1.48	0.94
1:A:502:CYS:SG	1:A:554:LYS:HD2	2.11	0.91
1:B:472:ILE:HD11	1:B:489:ILE:HG23	1.61	0.83
1:A:724:GLN:HG3	1:B:683:PRO:CG	2.08	0.82
1:B:658:ASN:C	1:B:658:ASN:HD22	1.88	0.81
1:B:573:SER:O	1:B:577:LYS:HB2	1.81	0.81
1:A:655:ASN:ND2	1:A:657:ASN:H	1.79	0.79
1:A:641:ASN:HD22	1:A:664:ARG:NH1	1.80	0.79
1:A:578:PHE:HZ	1:A:703:MET:HG2	1.48	0.78
1:A:677:SER:OG	1:A:688:THR:HG21	1.84	0.78
1:A:573:SER:O	1:A:577:LYS:HB2	1.84	0.77
1:A:730:TYR:HB3	1:A:759:LEU:HD12	1.67	0.74
1:B:641:ASN:HD22	1:B:664:ARG:HH11	1.41	0.67
1:B:655:ASN:ND2	1:B:657:ASN:H	1.93	0.67
1:A:472:ILE:HD11	1:A:489:ILE:HG23	1.78	0.65
1:A:730:TYR:CB	1:A:759:LEU:HD12	2.26	0.65
1:A:634:ASP:HB3	1:A:637:LEU:HD12	1.78	0.64
1:B:654:LEU:HD12	1:B:661:HIS:HD2	1.62	0.64
1:B:706:LEU:O	1:B:708:ILE:HG12	1.98	0.63
1:A:655:ASN:C	1:A:655:ASN:HD22	2.07	0.62
1:A:523:PRO:HD2	1:A:695:GLU:HG3	1.82	0.61
1:A:550:ASP:O	1:A:554:LYS:HG3	2.01	0.60
1:B:658:ASN:C	1:B:658:ASN:ND2	2.54	0.60
1:B:641:ASN:HD22	1:B:664:ARG:NH1	2.00	0.60
1:A:716:ARG:O	1:A:719:LYS:HG3	2.02	0.60
1:B:616:SER:O	1:B:620:GLU:HB2	2.03	0.59
1:B:578:PHE:CZ	1:B:703:MET:HG2	2.38	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:580:HIS:ND1	1:A:581:PRO:HD2	2.19	0.58
1:B:658:ASN:OD1	1:B:661:HIS:ND1	2.37	0.57
1:B:580:HIS:ND1	1:B:581:PRO:HD2	2.20	0.57
1:A:545:HIS:O	1:A:545:HIS:HD2	1.87	0.57
1:A:641:ASN:ND2	1:A:664:ARG:HH11	1.91	0.56
1:A:658:ASN:HD22	1:A:658:ASN:C	2.13	0.56
1:A:448:MET:HE1	1:A:598:GLN:OE1	2.06	0.56
1:A:682:TRP:NE1	1:A:770:GLU:OE2	2.40	0.55
1:A:616:SER:O	1:A:620:GLU:HB2	2.07	0.54
1:B:507:LYS:HG2	1:B:607:GLY:O	2.07	0.54
1:A:527:TRP:O	1:A:530:ALA:HB3	2.07	0.54
1:B:483:GLU:HA	1:B:486:TRP:CE2	2.43	0.54
1:A:624:GLU:OE2	1:A:624:GLU:HA	2.08	0.53
1:A:767:ARG:HD2	1:B:476:HIS:NE2	2.23	0.52
1:B:730:TYR:CB	1:B:759:LEU:HD12	2.39	0.52
1:A:724:GLN:CD	1:B:683:PRO:HD3	2.26	0.52
1:B:654:LEU:HD12	1:B:661:HIS:CD2	2.44	0.52
1:A:756:ARG:NH1	1:B:485:MET:HE3	2.24	0.52
1:B:532:THR:HG22	1:B:676:CYS:SG	2.50	0.52
1:A:655:ASN:ND2	1:A:655:ASN:C	2.67	0.51
1:B:658:ASN:HD22	1:B:659:GLN:N	2.09	0.50
1:A:483:GLU:HA	1:A:486:TRP:CE2	2.46	0.50
1:A:756:ARG:HH22	1:B:484:ASN:HD21	1.60	0.50
1:B:730:TYR:HB3	1:B:759:LEU:CD1	2.41	0.49
1:B:634:ASP:HB3	1:B:637:LEU:HD12	1.94	0.49
1:A:507:LYS:NZ	1:A:612:SER:HB2	2.28	0.48
1:A:578:PHE:CZ	1:A:703:MET:HG2	2.39	0.48
1:B:506:GLU:OE1	1:B:510:ARG:NH2	2.45	0.48
1:B:658:ASN:OD1	1:B:661:HIS:CE1	2.66	0.48
1:B:572:ASN:O	1:B:576:GLN:HG3	2.13	0.48
1:A:506:GLU:OE1	1:A:510:ARG:NH2	2.46	0.48
1:B:700:GLY:HA3	1:B:714:MET:O	2.14	0.48
1:A:756:ARG:HH22	1:B:484:ASN:ND2	2.11	0.47
1:A:766:ILE:HG21	1:B:681:LEU:HD22	1.97	0.47
1:B:644:GLN:O	1:B:648:MET:HG3	2.16	0.46
1:B:655:ASN:ND2	1:B:655:ASN:C	2.72	0.46
1:A:724:GLN:HE21	1:A:724:GLN:HB2	1.54	0.46
1:A:689:ALA:O	1:A:692:ILE:HG22	2.15	0.46
1:A:450:ILE:HG23	1:A:605:LEU:HD13	1.98	0.46
1:B:609:ASN:HB3	1:B:612:SER:OG	2.16	0.46
1:A:658:ASN:C	1:A:658:ASN:ND2	2.74	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:724:GLN:OE1	1:B:683:PRO:CD	2.29	0.45
1:B:603:LEU:HA	1:B:603:LEU:HD23	1.77	0.45
1:A:655:ASN:HD22	1:A:657:ASN:H	1.57	0.45
1:A:545:HIS:O	1:A:545:HIS:CD2	2.68	0.45
1:B:575:LEU:HD11	1:B:591:MET:SD	2.57	0.45
1:B:482:PHE:O	1:B:483:GLU:C	2.60	0.45
1:B:484:ASN:H	1:B:484:ASN:HD22	1.63	0.44
1:B:448:MET:HE1	1:B:598:GLN:OE1	2.17	0.44
1:A:569:GLY:HA2	1:A:695:GLU:OE2	2.18	0.44
1:B:550:ASP:O	1:B:554:LYS:HG3	2.18	0.44
1:B:519:TYR:CZ	1:B:530:ALA:HB2	2.52	0.44
1:B:504:GLU:CD	1:B:507:LYS:HD2	2.43	0.43
1:B:569:GLY:HA2	1:B:695:GLU:OE2	2.18	0.43
1:A:656:LEU:HD11	1:A:665:VAL:HG21	1.99	0.43
1:B:682:TRP:CG	1:B:761:GLN:HE21	2.37	0.43
1:A:682:TRP:HB3	1:A:683:PRO:HD3	2.01	0.43
1:B:578:PHE:HZ	1:B:703:MET:HG2	1.84	0.43
1:A:675:LEU:O	1:A:678:VAL:HG22	2.18	0.43
1:B:648:MET:SD	1:B:661:HIS:CD2	3.12	0.43
1:B:682:TRP:CD1	1:B:761:GLN:HE21	2.37	0.43
1:A:617:SER:O	1:A:621:GLN:HB2	2.19	0.42
1:B:450:ILE:HG23	1:B:605:LEU:HD13	2.01	0.42
1:A:700:GLY:HA3	1:A:714:MET:O	2.20	0.42
1:B:448:MET:HE2	1:B:602:ILE:HG13	2.01	0.42
1:A:568:ARG:N	1:A:568:ARG:HD2	2.34	0.42
1:A:580:HIS:CE1	5:A:901:HOH:O	2.46	0.41
1:A:462:PHE:CD2	1:A:488:GLY:HA3	2.55	0.41
1:A:666:ILE:O	1:A:669:MET:HB2	2.20	0.41
1:A:712:PRO:HB3	1:A:718:LYS:HD3	2.03	0.41
1:B:666:ILE:O	1:B:669:MET:HB2	2.21	0.41
1:A:490:PHE:HA	1:A:493:MET:HE2	2.03	0.41
1:A:735:ILE:HB	1:A:736:PRO:HD3	2.03	0.41
1:B:617:SER:O	1:B:621:GLN:HB2	2.21	0.40
1:A:578:PHE:HA	1:A:579:ASP:OD1	2.22	0.40
1:B:580:HIS:CG	1:B:581:PRO:HD2	2.56	0.40

All (1) 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:449:SER:OG	1:B:614:LEU:O[1_545]	1.95	0.25

5.3 Torsion angles

5.3.1 Protein backbone

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	322/345 (93%)	296 (92%)	25 (8%)	1 (0%)	36	44
1	B	322/345 (93%)	302 (94%)	19 (6%)	1 (0%)	36	44
All	All	644/690 (93%)	598 (93%)	44 (7%)	2 (0%)	36	44

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	707	GLY
1	B	707	GLY

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	293/308 (95%)	261 (89%)	32 (11%)	6	5
1	B	293/308 (95%)	265 (90%)	28 (10%)	8	8
All	All	586/616 (95%)	526 (90%)	60 (10%)	7	7

All (60) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	448	MET
1	A	450	ILE
1	A	453	SER
1	A	463	THR

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Mol	Chain	Res	Type
1	A	466	VAL
1	A	497	SER
1	A	498	CYS
1	A	500	THR
1	A	501	SER
1	A	504	GLU
1	A	510	ARG
1	A	521	ARG
1	A	578	PHE
1	A	579	ASP
1	A	585	LEU
1	A	587	SER
1	A	588	THR
1	A	620	GLU
1	A	624	GLU
1	A	653	SER
1	A	655	ASN
1	A	658	ASN
1	A	662	ARG
1	A	677	SER
1	A	681	LEU
1	A	706	LEU
1	A	708	ILE
1	A	709	GLN
1	A	724	GLN
1	A	760	SER
1	A	767	ARG
1	A	770	GLU
1	B	448	MET
1	B	450	ILE
1	B	453	SER
1	B	454	GLU
1	B	463	THR
1	B	466	VAL
1	B	471	GLU
1	B	472	ILE
1	B	484	ASN
1	B	497	SER
1	B	501	SER
1	B	504	GLU
1	B	510	ARG
1	B	577	LYS

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Mol	Chain	Res	Type
1	B	585	LEU
1	B	587	SER
1	B	588	THR
1	B	606	GLU
1	B	635	LEU
1	B	658	ASN
1	B	660	SER
1	B	677	SER
1	B	706	LEU
1	B	708	ILE
1	B	709	GLN
1	B	724	GLN
1	B	760	SER
1	B	770	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	461	GLN
1	A	476	HIS
1	A	545	HIS
1	A	621	GLN
1	A	641	ASN
1	A	655	ASN
1	A	658	ASN
1	A	661	HIS
1	A	690	ASN
1	A	726	GLN
1	B	461	GLN
1	B	484	ASN
1	B	593	GLN
1	B	641	ASN
1	B	655	ASN
1	B	658	ASN
1	B	661	HIS
1	B	690	ASN
1	B	709	GLN
1	B	724	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 5 ligands modelled in this entry, 4 are monoatomic - leaving 1 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
4	D6X	A	803	-	27,27,27	1.43	4 (14%)	34,38,38	2.63	9 (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
4	D6X	A	803	-	-	3/6/6/6	0/4/4/4

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	803	D6X	C10-C8	-3.55	1.45	1.49
4	A	803	D6X	C2-C3	2.94	1.47	1.41
4	A	803	D6X	C15-N12	2.09	1.39	1.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	803	D6X	C3-N4	-2.09	1.30	1.33

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	803	D6X	C11-N12-C15	-6.61	113.01	123.41
4	A	803	D6X	C1-C2-C3	-6.31	112.51	120.89
4	A	803	D6X	C10-C11-N12	-5.41	100.74	111.86
4	A	803	D6X	C2-C3-N4	5.38	128.00	122.71
4	A	803	D6X	C2-C1-N5	4.58	120.70	115.03
4	A	803	D6X	C16-C15-N12	-3.68	113.32	121.17
4	A	803	D6X	N5-C6-N4	-3.64	119.45	122.61
4	A	803	D6X	C14-C3-C2	-3.34	116.07	120.44
4	A	803	D6X	C10-C8-N7	2.41	125.71	120.27

There are no chirality outliers.

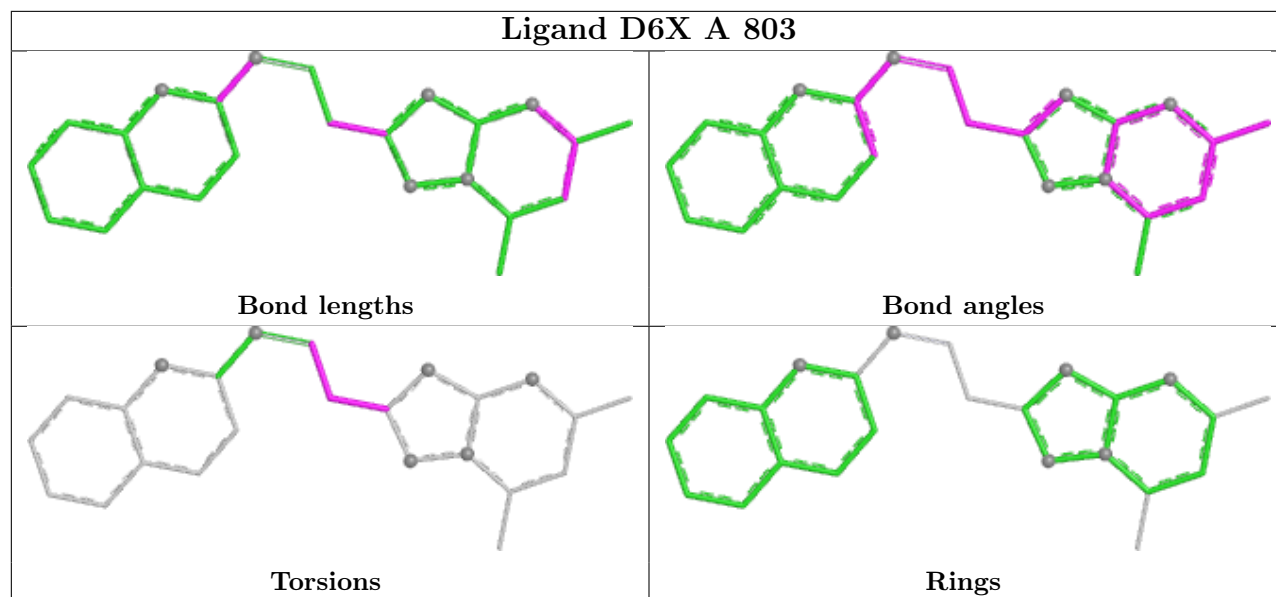
All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	803	D6X	C11-C10-C8-N9
4	A	803	D6X	C8-C10-C11-N12
4	A	803	D6X	C11-C10-C8-N7

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.



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, 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	324/345 (93%)	0.32	13 (4%) 42 39	33, 56, 92, 120	0
1	B	324/345 (93%)	0.22	11 (3%) 48 46	32, 51, 81, 105	0
All	All	648/690 (93%)	0.27	24 (3%) 45 43	32, 53, 88, 120	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	578	PHE	3.2
1	B	498	CYS	3.2
1	A	545	HIS	3.2
1	A	578	PHE	3.1
1	B	447	HIS	3.1
1	B	768	GLY	2.9
1	B	585	LEU	2.8
1	A	585	LEU	2.7
1	A	584	ALA	2.7
1	B	500	THR	2.6
1	B	584	ALA	2.5
1	A	498	CYS	2.4
1	A	613	THR	2.3
1	A	450	ILE	2.3
1	A	720	ASP	2.2
1	B	580	HIS	2.2
1	A	707	GLY	2.2
1	A	470	LYS	2.2
1	A	447	HIS	2.1
1	B	550	ASP	2.1
1	B	462	PHE	2.1
1	B	770	GLU	2.1
1	A	672	ALA	2.1
1	A	586	TYR	2.1

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](#)

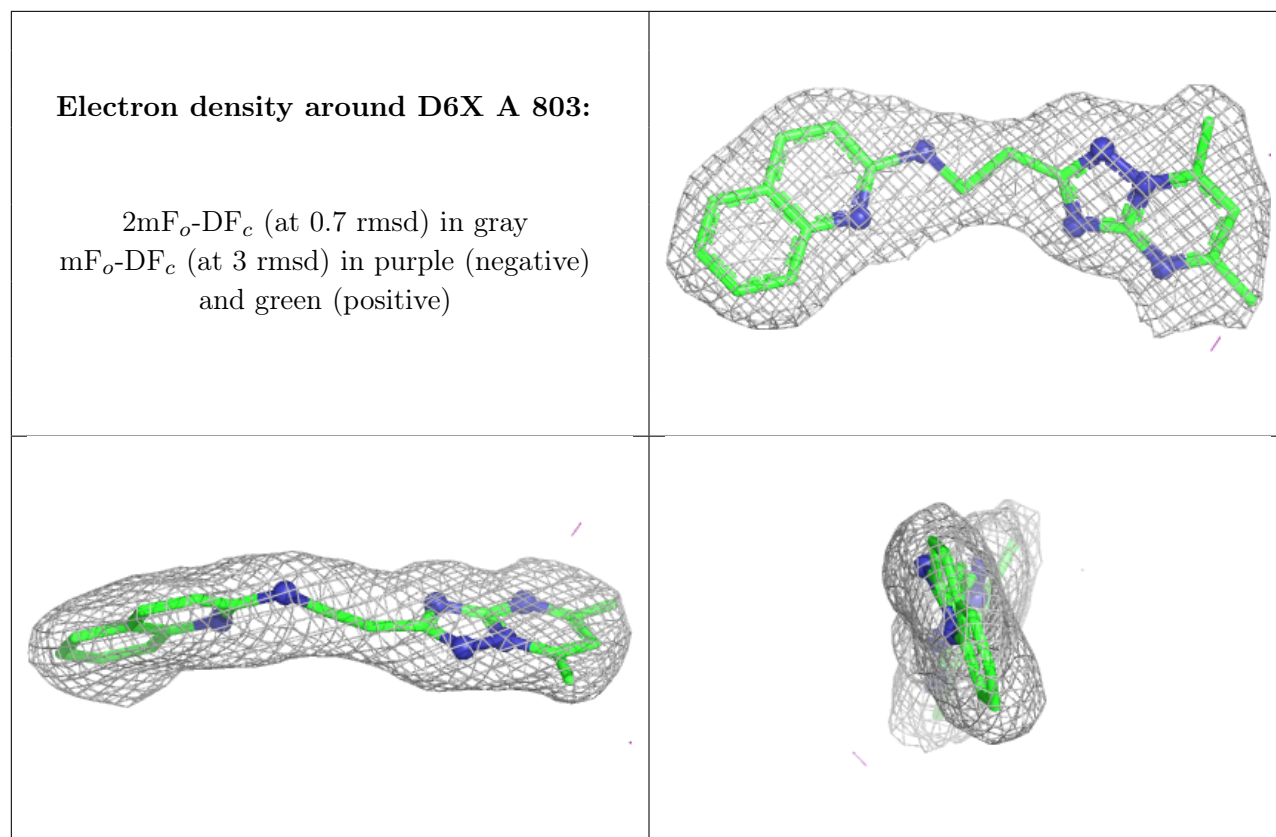
There are no oligosaccharides in this entry.

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
4	D6X	A	803	24/24	0.94	0.07	35,37,41,43	0
3	MG	A	802	1/1	0.95	0.06	43,43,43,43	0
3	MG	B	802	1/1	0.99	0.04	34,34,34,34	0
2	ZN	A	801	1/1	0.99	0.02	51,51,51,51	0
2	ZN	B	801	1/1	1.00	0.02	39,39,39,39	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.