



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 18, 2026 – 03:47 PM UTC

PDB ID : 3WS8 / pdb_00003ws8
Title : Crystal structure of PDE10A in complex with a benzimidazole inhibitor
Authors : Amano, Y.; Honbou, K.
Deposited on : 2014-03-04
Resolution : 2.60 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

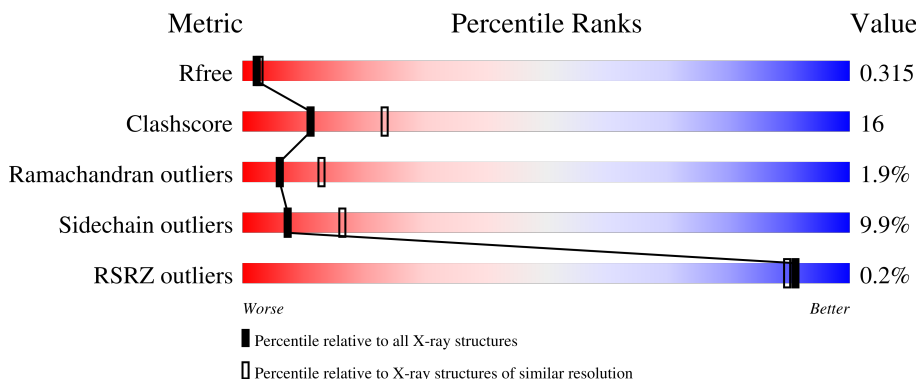
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	
1	B	345	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5337 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).

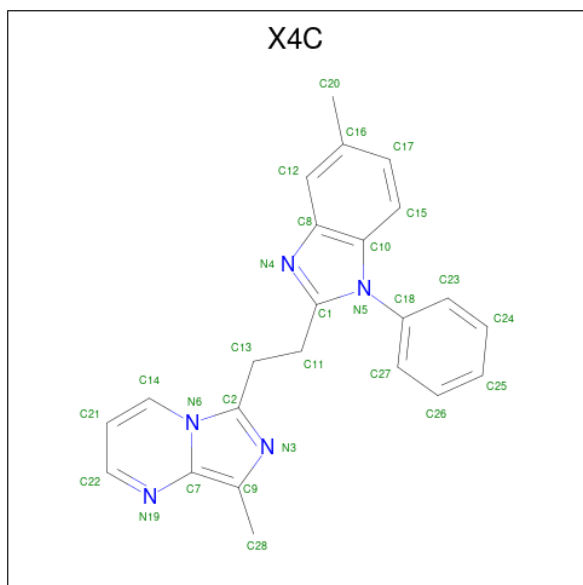
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mg	0	0
			1	1		

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

- Molecule 4 is 8-methyl-6-[2-(5-methyl-1-phenyl-1H-benzimidazol-2-yl)ethyl]imidazo[1,5-a]pyrimidine (CCD ID: X4C) (formula: C₂₃H₂₁N₅).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	N	0	0
			28	23	5		
4	B	1	Total	C	N	0	0
			28	23	5		

- Molecule 5 is water.

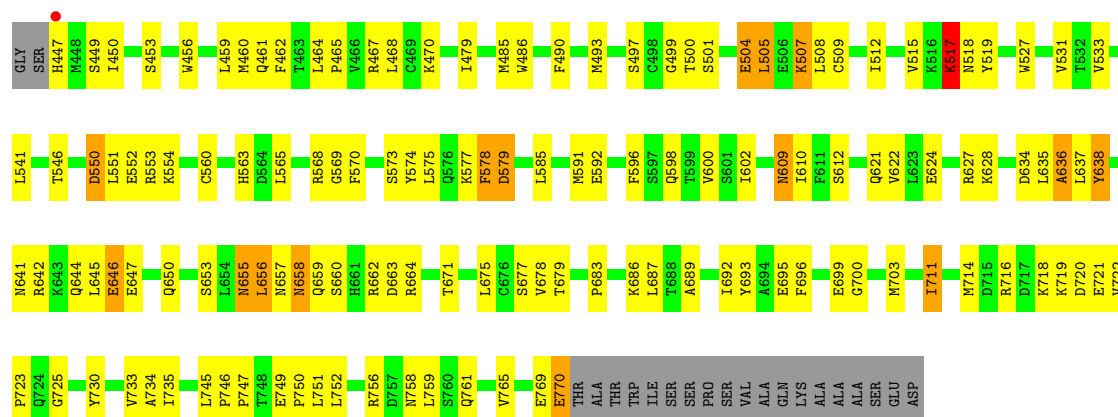
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	9	Total	O	0	0
			9	9		
5	B	10	Total	O	0	0
			10	10		

3 Residue-property plots

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.

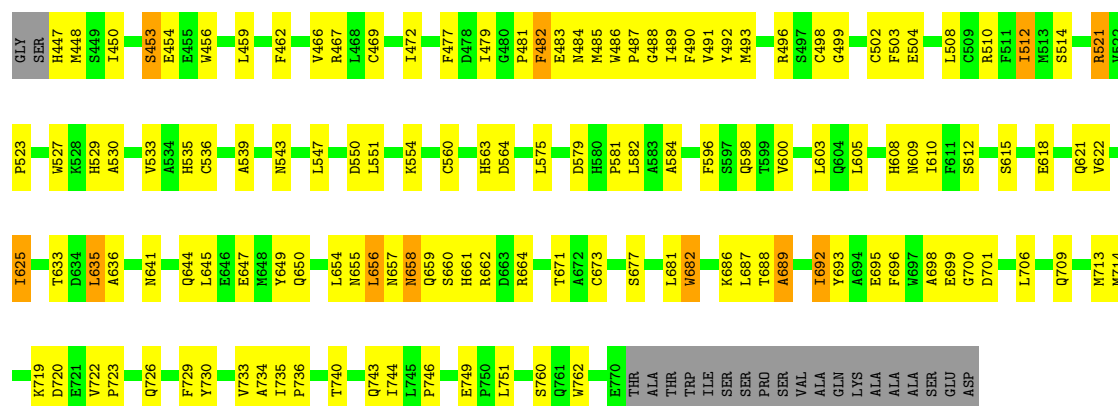
- Molecule 1: cAMP and cAMP-inhibited cGMP 3',5'-cyclic phosphodiesterase 10A

Chain A: 



- Molecule 1: cAMP and cAMP-inhibited cGMP 3',5'-cyclic phosphodiesterase 10A

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	49.31Å 81.11Å 160.47Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.96 – 2.60 19.96 – 2.60	Depositor EDS
% Data completeness (in resolution range)	89.3 (19.96-2.60) 89.3 (19.96-2.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.80 (at 2.59Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.219 , 0.314 0.219 , 0.315	Depositor DCC
R_{free} test set	935 reflections (4.57%)	wwPDB-VP
Wilson B-factor (Å ²)	44.8	Xtriage
Anisotropy	0.230	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 35.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.53$, $\langle L^2 \rangle = 0.37$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5337	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.34% 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: X4C, 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	0.76	0/2694	1.14	6/3647 (0.2%)
1	B	0.83	0/2694	1.12	10/3647 (0.3%)
All	All	0.80	0/5388	1.13	16/7294 (0.2%)

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	745	LEU	CA-C-N	-8.50	111.62	120.38
1	A	745	LEU	C-N-CA	-8.50	111.62	120.38
1	B	625	ILE	N-CA-C	-7.80	103.26	110.82
1	B	682	TRP	CA-C-N	-7.31	112.17	119.56
1	B	682	TRP	C-N-CA	-7.31	112.17	119.56
1	B	536	CYS	N-CA-C	-5.75	105.15	111.82
1	B	722	VAL	N-CA-CB	5.49	114.73	110.45
1	B	746	PRO	N-CA-C	5.34	117.22	110.70
1	B	560	CYS	N-CA-C	-5.32	104.38	112.04
1	A	711	ILE	CA-C-N	-5.26	113.61	119.19
1	A	711	ILE	C-N-CA	-5.26	113.61	119.19
1	A	655	ASN	N-CA-C	5.22	117.51	107.44
1	B	504	GLU	CA-C-N	5.22	127.53	120.38
1	B	504	GLU	C-N-CA	5.22	127.53	120.38
1	A	733	VAL	N-CA-C	5.20	115.73	111.62
1	B	550	ASP	N-CA-C	-5.16	104.92	111.11

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	85	0
1	B	2629	0	2588	87	0
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	28	0	21	1	0
4	B	28	0	21	3	0
5	A	9	0	0	2	0
5	B	10	0	0	0	0
All	All	5337	0	5218	172	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 16.

All (172) 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:B:655:ASN:H	1:B:661:HIS:HD2	1.17	0.92
1:B:656:LEU:H	1:B:656:LEU:HD22	1.37	0.88
1:A:450:ILE:O	1:A:517:LYS:HD2	1.75	0.87
1:A:675:LEU:O	1:A:678:VAL:HG22	1.74	0.87
1:B:658:ASN:HD21	1:B:660:SER:HB3	1.40	0.87
1:A:658:ASN:HD22	1:A:658:ASN:C	1.88	0.82
1:B:641:ASN:HD22	1:B:664:ARG:HH11	1.28	0.82
1:B:658:ASN:C	1:B:658:ASN:HD22	1.89	0.81
1:B:521:ARG:HH11	1:B:521:ARG:HG3	1.46	0.81
1:A:609:ASN:O	1:A:612:SER:OG	1.98	0.78
1:A:769:GLU:O	1:A:770:GLU:HB2	1.82	0.77
1:A:655:ASN:C	1:A:655:ASN:HD22	1.95	0.74
1:B:645:LEU:HD11	1:B:664:ARG:O	1.89	0.72
1:B:735:ILE:HB	1:B:736:PRO:HD3	1.71	0.71
1:A:658:ASN:HD22	1:A:659:GLN:N	1.89	0.70
1:B:658:ASN:C	1:B:658:ASN:ND2	2.49	0.70
1:A:490:PHE:HA	1:A:493:MET:HE2	1.75	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:656:LEU:HD22	1:A:656:LEU:H	1.58	0.69
1:A:655:ASN:ND2	1:A:657:ASN:H	1.91	0.68
1:A:577:LYS:O	1:A:578:PHE:HB3	1.92	0.67
1:A:689:ALA:O	1:A:692:ILE:HG22	1.94	0.67
1:B:533:VAL:HG13	1:B:673:CYS:HB3	1.77	0.66
1:B:615:SER:OG	1:B:618:GLU:HG3	1.96	0.66
1:A:761:GLN:O	1:A:765:VAL:HG23	1.96	0.66
1:A:515:VAL:HG13	1:A:565:LEU:HD21	1.77	0.66
1:B:521:ARG:HH11	1:B:521:ARG:CG	2.09	0.65
1:A:638:TYR:CD1	1:A:671:THR:HG21	2.32	0.64
1:B:523:PRO:HD2	1:B:695:GLU:HG3	1.80	0.63
1:B:693:TYR:OH	4:B:803:X4C:N4	2.28	0.63
1:A:552:GLU:OE1	1:A:662:ARG:NH2	2.32	0.62
1:A:752:LEU:HD21	1:A:756:ARG:NH2	2.15	0.61
1:B:692:ILE:HD11	1:B:696:PHE:CZ	2.36	0.61
1:B:448:MET:HE1	1:B:598:GLN:OE1	2.00	0.61
1:A:624:GLU:OE2	1:A:624:GLU:HA	2.01	0.60
1:A:730:TYR:HA	1:A:734:ALA:HB3	1.83	0.60
1:B:453:SER:HA	1:B:456:TRP:CD1	2.36	0.60
1:B:450:ILE:HG12	1:B:605:LEU:HD13	1.84	0.59
1:A:563:HIS:CD2	5:A:904:HOH:O	2.56	0.58
1:A:634:ASP:HB3	1:A:637:LEU:HD12	1.85	0.58
1:B:498:CYS:HB2	1:B:502:CYS:SG	2.45	0.57
1:A:568:ARG:HD2	1:A:591:MET:HE1	1.87	0.56
1:B:700:GLY:HA3	1:B:714:MET:O	2.06	0.56
1:A:541:LEU:HD22	1:A:553:ARG:HG2	1.87	0.56
1:B:596:PHE:O	1:B:600:VAL:HG23	2.05	0.55
1:B:481:PRO:HB2	1:B:482:PHE:CE1	2.41	0.55
1:B:472:ILE:HD11	1:B:489:ILE:HG23	1.89	0.55
1:B:649:TYR:CZ	1:B:743:GLN:HB3	2.42	0.55
1:B:655:ASN:H	1:B:661:HIS:CD2	2.09	0.55
1:B:563:HIS:HB2	1:B:633:THR:HG21	1.88	0.54
1:B:490:PHE:HA	1:B:493:MET:CE	2.38	0.54
1:B:490:PHE:HA	1:B:493:MET:HE2	1.90	0.54
1:B:482:PHE:HB3	1:B:485:MET:HE3	1.89	0.54
1:B:641:ASN:HD22	1:B:664:ARG:NH1	2.02	0.54
1:A:749:GLU:N	1:A:750:PRO:CD	2.71	0.54
1:B:508:LEU:O	1:B:512:ILE:HD12	2.08	0.54
1:A:460:MET:HG2	1:A:505:LEU:HG	1.89	0.54
1:A:459:LEU:HD23	1:A:509:CYS:HA	1.90	0.53
1:B:655:ASN:N	1:B:661:HIS:HD2	1.96	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:490:PHE:HA	1:A:493:MET:CE	2.38	0.53
1:A:722:VAL:HB	1:A:723:PRO:HD3	1.90	0.53
1:B:681:LEU:O	1:B:682:TRP:C	2.50	0.52
1:B:641:ASN:ND2	1:B:664:ARG:HH11	2.03	0.52
1:A:655:ASN:HD22	1:A:657:ASN:H	1.57	0.52
1:B:462:PHE:CD2	1:B:488:GLY:HA3	2.45	0.52
1:B:644:GLN:HG3	1:B:664:ARG:CZ	2.39	0.52
1:B:483:GLU:HA	1:B:486:TRP:CE2	2.45	0.52
1:A:479:ILE:HD12	1:A:486:TRP:CD1	2.45	0.52
1:B:448:MET:HB2	1:B:605:LEU:HD11	1.92	0.51
1:A:518:ASN:HD22	1:A:602:ILE:HD11	1.75	0.51
1:A:700:GLY:HA3	1:A:714:MET:O	2.11	0.51
1:A:569:GLY:HA2	1:A:695:GLU:OE2	2.10	0.51
1:A:577:LYS:O	1:A:578:PHE:CB	2.59	0.51
1:B:726:GLN:HG3	1:B:762:TRP:CE2	2.45	0.50
1:A:641:ASN:HD22	1:A:664:ARG:HH11	1.58	0.50
1:B:622:VAL:HA	1:B:625:ILE:HD12	1.94	0.50
1:A:628:LYS:NZ	1:A:663:ASP:OD2	2.29	0.50
1:A:550:ASP:O	1:A:554:LYS:HG2	2.11	0.50
1:B:677:SER:HB2	1:B:688:THR:HG21	1.95	0.49
1:A:519:TYR:CE1	1:A:565:LEU:HD23	2.48	0.49
1:A:541:LEU:HB3	1:A:553:ARG:HD2	1.95	0.49
1:B:521:ARG:CG	1:B:521:ARG:NH1	2.74	0.49
1:B:485:MET:O	1:B:489:ILE:HG13	2.13	0.48
1:B:499:GLY:O	1:B:502:CYS:HB2	2.13	0.48
1:B:720:ASP:O	1:B:723:PRO:HD2	2.12	0.48
1:A:504:GLU:O	1:A:505:LEU:C	2.55	0.48
1:A:696:PHE:HB3	1:A:714:MET:HG2	1.94	0.48
1:B:491:VAL:CG2	1:B:512:ILE:HD13	2.44	0.48
1:B:656:LEU:H	1:B:656:LEU:CD2	2.13	0.48
1:B:658:ASN:HD22	1:B:659:GLN:N	2.11	0.48
1:A:533:VAL:HG12	1:A:560:CYS:HB3	1.95	0.47
1:A:467:ARG:NH2	1:A:468:LEU:HD21	2.29	0.47
1:B:698:ALA:O	1:B:701:ASP:HB2	2.14	0.47
1:B:581:PRO:O	1:B:584:ALA:HB3	2.15	0.47
1:B:539:ALA:O	1:B:543:ASN:ND2	2.41	0.47
1:A:758:ASN:O	1:A:761:GLN:HB2	2.15	0.46
1:B:735:ILE:HB	1:B:736:PRO:CD	2.43	0.46
1:A:459:LEU:HD23	1:A:509:CYS:CA	2.46	0.46
1:A:485:MET:O	1:A:486:TRP:C	2.59	0.46
1:B:493:MET:SD	1:B:535:HIS:HA	2.55	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:641:ASN:HA	1:B:664:ARG:NH1	2.31	0.46
1:A:655:ASN:HD22	1:A:656:LEU:N	2.13	0.45
1:B:491:VAL:HG22	1:B:512:ILE:HD13	1.99	0.45
1:B:527:TRP:O	1:B:530:ALA:HB3	2.17	0.45
1:A:533:VAL:HG21	1:A:563:HIS:CE1	2.51	0.45
1:A:573:SER:O	1:A:577:LYS:HB2	2.17	0.45
1:B:459:LEU:HD13	1:B:487:PRO:HB2	1.99	0.45
1:A:592:GLU:HG3	5:A:901:HOH:O	2.17	0.44
1:A:479:ILE:HD12	1:A:486:TRP:NE1	2.32	0.44
1:A:679:THR:HG21	1:A:751:LEU:HD22	1.98	0.44
1:B:498:CYS:CB	1:B:502:CYS:SG	3.05	0.44
1:A:508:LEU:HG	1:A:512:ILE:CD1	2.47	0.44
1:B:603:LEU:HD23	1:B:603:LEU:HA	1.86	0.44
1:B:469:CYS:O	1:B:472:ILE:HG22	2.17	0.44
1:B:496:ARG:HA	1:B:496:ARG:HD2	1.79	0.44
1:A:479:ILE:HD13	1:A:531:VAL:HG12	2.00	0.44
1:B:605:LEU:O	1:B:608:HIS:HB2	2.18	0.43
1:B:635:LEU:HD23	1:B:635:LEU:HA	1.74	0.43
1:A:570:PHE:HA	1:A:699:GLU:OE1	2.19	0.43
1:A:683:PRO:O	1:A:687:LEU:HG	2.18	0.43
1:A:574:TYR:HD1	1:A:703:MET:HE3	1.83	0.43
1:B:654:LEU:HA	1:B:661:HIS:CD2	2.53	0.43
1:B:644:GLN:O	1:B:647:GLU:HG3	2.18	0.43
1:A:449:SER:HB2	1:A:518:ASN:OD1	2.19	0.43
1:A:459:LEU:HD23	1:A:509:CYS:CB	2.49	0.43
1:A:459:LEU:HD23	1:A:509:CYS:HB3	2.00	0.43
1:A:645:LEU:O	1:A:646:GLU:C	2.61	0.43
1:B:686:LYS:O	1:B:689:ALA:HB3	2.19	0.43
1:A:486:TRP:HE3	1:A:527:TRP:CD1	2.37	0.43
1:B:486:TRP:HE3	1:B:527:TRP:CD1	2.36	0.43
1:A:769:GLU:O	1:A:770:GLU:CB	2.61	0.43
1:A:658:ASN:C	1:A:658:ASN:ND2	2.60	0.43
1:B:529:HIS:CD2	1:B:564:ASP:OD2	2.71	0.43
1:A:462:PHE:CE1	1:A:485:MET:HG2	2.53	0.42
1:A:716:ARG:C	1:A:718:LYS:H	2.26	0.42
1:B:730:TYR:HA	1:B:734:ALA:HB3	2.01	0.42
1:A:578:PHE:HA	1:A:579:ASP:OD2	2.19	0.42
1:B:655:ASN:HD21	1:B:657:ASN:HB2	1.83	0.42
1:A:507:LYS:HB3	1:A:610:ILE:HG22	2.02	0.42
1:A:598:GLN:O	1:A:602:ILE:HG13	2.19	0.42
1:A:722:VAL:O	1:A:723:PRO:C	2.62	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:503:PHE:HZ	1:B:554:LYS:HB3	1.85	0.42
1:B:551:LEU:HD12	1:B:551:LEU:HA	1.75	0.42
1:A:658:ASN:ND2	1:A:660:SER:H	2.18	0.42
1:A:504:GLU:H	1:A:504:GLU:HG3	1.58	0.42
1:A:644:GLN:O	1:A:647:GLU:HB3	2.19	0.42
1:A:596:PHE:CE2	1:A:627:ARG:HG3	2.55	0.42
1:B:635:LEU:O	1:B:636:ALA:C	2.63	0.42
1:A:656:LEU:H	1:A:656:LEU:CD2	2.31	0.41
1:B:477:PHE:HB3	1:B:535:HIS:CE1	2.54	0.41
1:A:508:LEU:HG	1:A:512:ILE:HD12	2.02	0.41
1:A:575:LEU:HD11	1:A:591:MET:HE2	2.03	0.41
1:A:693:TYR:CZ	1:A:722:VAL:HG13	2.55	0.41
1:A:725:GLY:O	4:A:803:X4C:H6	2.19	0.41
1:B:609:ASN:OD1	1:B:612:SER:N	2.54	0.41
1:A:635:LEU:O	1:A:636:ALA:C	2.63	0.41
1:B:656:LEU:HD22	1:B:656:LEU:N	2.19	0.41
1:B:492:TYR:CZ	1:B:496:ARG:HG3	2.56	0.41
1:A:655:ASN:C	1:A:655:ASN:ND2	2.65	0.41
1:B:649:TYR:CE2	1:B:743:GLN:HB3	2.55	0.41
1:B:699:GLU:O	1:B:700:GLY:C	2.63	0.41
1:B:740:THR:O	1:B:744:ILE:HG13	2.21	0.41
1:A:456:TRP:O	1:A:459:LEU:HB3	2.21	0.40
1:A:553:ARG:HG3	1:A:553:ARG:HH11	1.86	0.40
1:B:510:ARG:HB3	1:B:608:HIS:CE1	2.57	0.40
1:B:729:PHE:CD1	4:B:803:X4C:H18	2.56	0.40
1:A:746:PRO:N	1:A:747:PRO:CD	2.85	0.40
1:B:751:LEU:HD23	1:B:751:LEU:HA	1.99	0.40
1:B:514:SER:OG	1:B:608:HIS:NE2	2.46	0.40
1:B:641:ASN:HD21	1:B:664:ARG:HB3	1.85	0.40
1:B:656:LEU:O	1:B:662:ARG:HD2	2.21	0.40
1:A:464:LEU:HA	1:A:465:PRO:HD3	1.94	0.40
1:B:713:MET:HG3	4:B:803:X4C:C8	2.51	0.40

There are no symmetry-related clashes.

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%)	291 (90%)	23 (7%)	8 (2%)	4	8
1	B	322/345 (93%)	290 (90%)	28 (9%)	4 (1%)	10	23
All	All	644/690 (93%)	581 (90%)	51 (8%)	12 (2%)	6	13

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	497	SER
1	A	505	LEU
1	A	609	ASN
1	B	579	ASP
1	B	689	ALA
1	A	517	LYS
1	A	578	PHE
1	A	621	GLN
1	B	719	LYS
1	A	636	ALA
1	B	733	VAL
1	A	499	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	13
1	B	293/308 (95%)	267 (91%)	26 (9%)	9	20
All	All	586/616 (95%)	528 (90%)	58 (10%)	7	16

All (58) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	447	HIS
1	A	453	SER
1	A	461	GLN
1	A	470	LYS
1	A	500	THR
1	A	501	SER
1	A	504	GLU
1	A	507	LYS
1	A	517	LYS
1	A	546	THR
1	A	550	ASP
1	A	551	LEU
1	A	579	ASP
1	A	585	LEU
1	A	600	VAL
1	A	622	VAL
1	A	638	TYR
1	A	642	ARG
1	A	646	GLU
1	A	650	GLN
1	A	653	SER
1	A	656	LEU
1	A	658	ASN
1	A	677	SER
1	A	686	LYS
1	A	711	ILE
1	A	719	LYS
1	A	720	ASP
1	A	721	GLU
1	A	735	ILE
1	A	759	LEU
1	A	770	GLU
1	B	447	HIS
1	B	453	SER
1	B	454	GLU
1	B	466	VAL
1	B	467	ARG
1	B	479	ILE
1	B	482	PHE
1	B	484	ASN
1	B	512	ILE
1	B	521	ARG
1	B	547	LEU

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Mol	Chain	Res	Type
1	B	575	LEU
1	B	582	LEU
1	B	610	ILE
1	B	621	GLN
1	B	635	LEU
1	B	650	GLN
1	B	656	LEU
1	B	658	ASN
1	B	671	THR
1	B	687	LEU
1	B	692	ILE
1	B	706	LEU
1	B	709	GLN
1	B	749	GLU
1	B	760	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	447	HIS
1	A	518	ASN
1	A	598	GLN
1	A	641	ASN
1	A	650	GLN
1	A	655	ASN
1	A	658	ASN
1	A	690	ASN
1	A	709	GLN
1	A	724	GLN
1	B	484	ASN
1	B	545	HIS
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	726	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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

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	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	X4C	B	803	-	31,32,32	1.27	4 (12%)	39,46,46	1.90	11 (28%)
4	X4C	A	803	-	31,32,32	1.36	4 (12%)	39,46,46	1.44	8 (20%)

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	X4C	B	803	-	-	2/9/9/9	0/5/5/5
4	X4C	A	803	-	-	0/9/9/9	0/5/5/5

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	803	X4C	C21-C22	4.10	1.49	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	803	X4C	C21-C22	3.97	1.48	1.40
4	A	803	X4C	C18-N5	-3.14	1.40	1.44
4	B	803	X4C	C18-N5	-3.09	1.40	1.44
4	A	803	X4C	C13-C2	-2.67	1.46	1.49
4	B	803	X4C	C10-N5	-2.06	1.37	1.40
4	A	803	X4C	C10-N5	-2.03	1.37	1.40
4	B	803	X4C	C7-N6	-2.02	1.36	1.39

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	803	X4C	C14-C21-C22	-4.73	113.39	118.87
4	A	803	X4C	C14-C21-C22	-3.87	114.39	118.87
4	B	803	X4C	C21-C22-N19	3.79	129.10	124.25
4	A	803	X4C	C8-C12-C16	-3.75	115.04	121.23
4	B	803	X4C	C20-C16-C12	-3.50	116.17	120.92
4	B	803	X4C	C22-N19-C7	-3.15	113.10	115.55
4	B	803	X4C	C17-C16-C12	3.14	122.59	117.96
4	B	803	X4C	C8-C12-C16	-3.14	116.04	121.23
4	A	803	X4C	C17-C15-C10	-3.12	112.84	119.10
4	B	803	X4C	C13-C11-C1	2.89	114.35	111.62
4	A	803	X4C	C17-C16-C12	2.82	122.11	117.96
4	B	803	X4C	C10-N5-C1	2.72	109.25	106.56
4	B	803	X4C	C8-C10-N5	-2.54	103.37	105.25
4	B	803	X4C	C17-C15-C10	-2.43	114.23	119.10
4	B	803	X4C	N5-C1-N4	-2.35	109.76	112.48
4	A	803	X4C	C15-C17-C16	2.24	123.96	121.37
4	A	803	X4C	C13-C11-C1	-2.08	109.65	111.62
4	A	803	X4C	C12-C8-C10	2.07	122.06	120.16
4	A	803	X4C	C21-C22-N19	2.00	126.81	124.25

There are no chirality outliers.

All (2) torsion outliers are listed below:

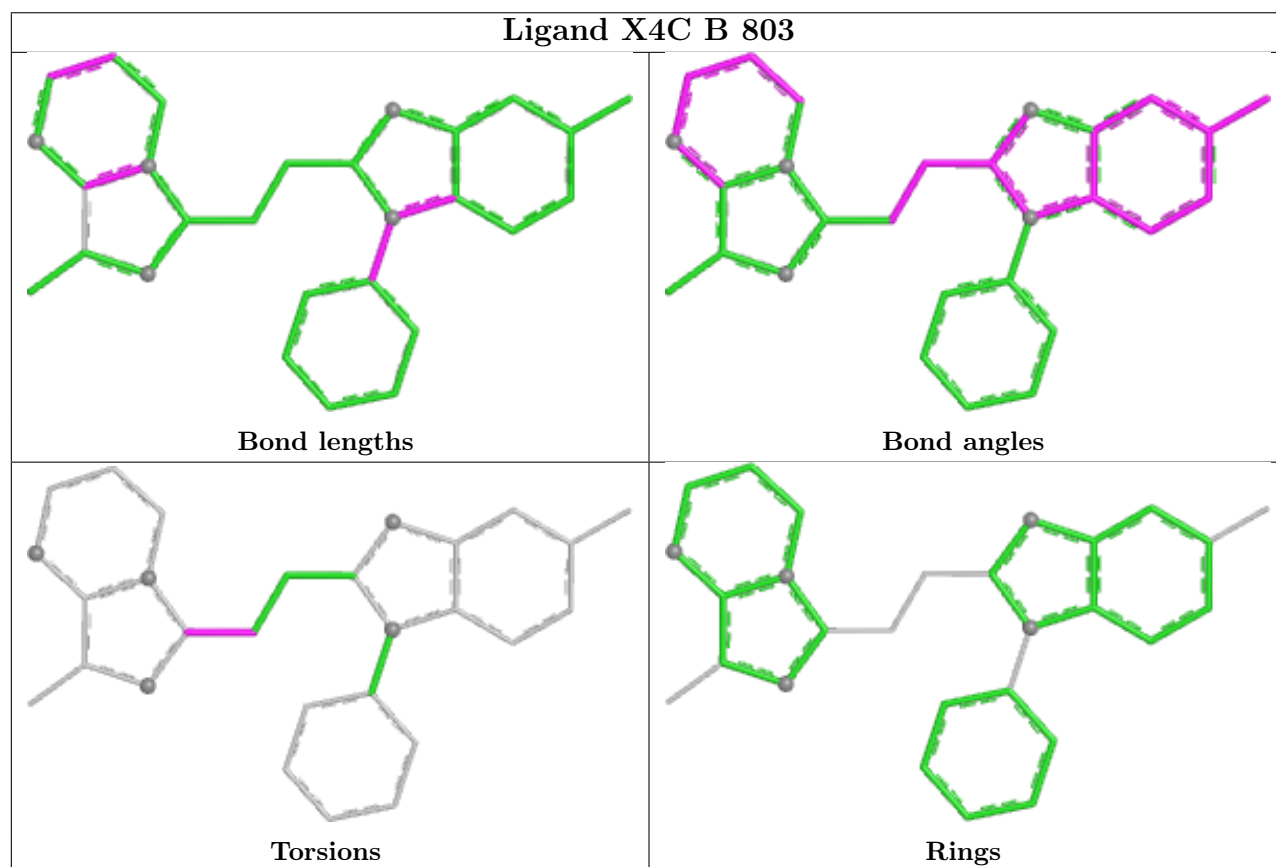
Mol	Chain	Res	Type	Atoms
4	B	803	X4C	C11-C13-C2-N3
4	B	803	X4C	C11-C13-C2-N6

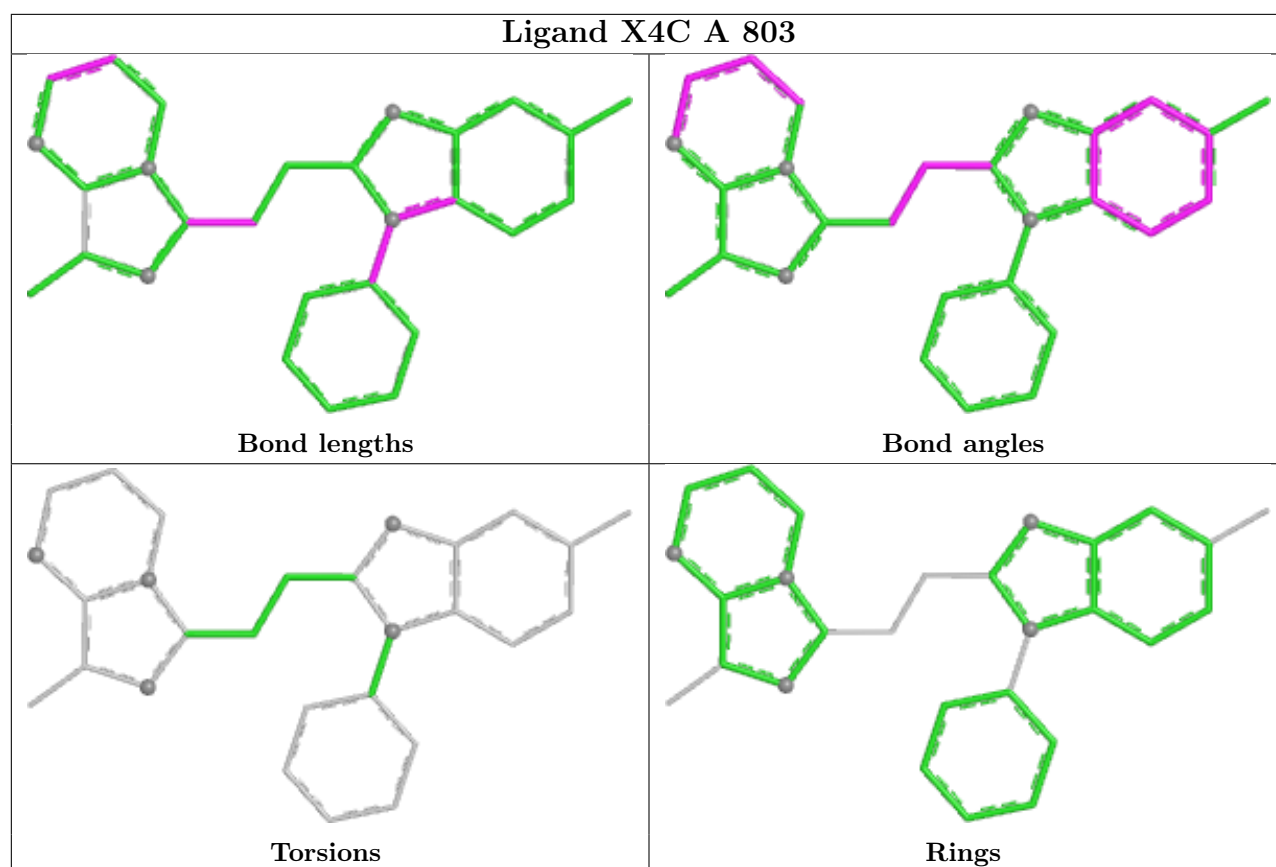
There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	803	X4C	3	0
4	A	803	X4C	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 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 [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	324/345 (93%)	-0.37	1 (0%) 90 88	24, 46, 73, 106	0
1	B	324/345 (93%)	-0.40	0 100 100	27, 45, 71, 85	0
All	All	648/690 (93%)	-0.38	1 (0%) 91 90	24, 46, 72, 106	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	447	HIS	2.4

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(Å ²)	Q<0.9
4	X4C	B	803	28/28	0.92	0.08	35,41,45,46	0
4	X4C	A	803	28/28	0.93	0.08	37,41,46,49	0
3	MG	B	802	1/1	0.96	0.08	25,25,25,25	0
3	MG	A	802	1/1	0.99	0.02	26,26,26,26	0

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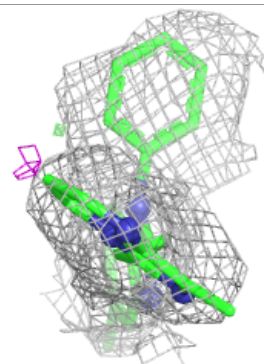
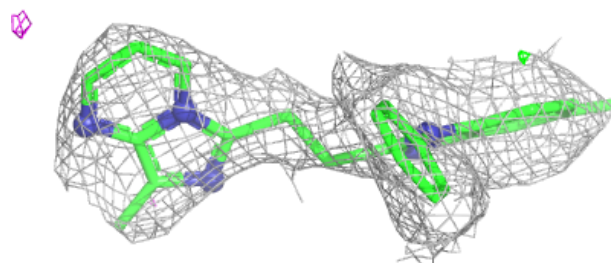
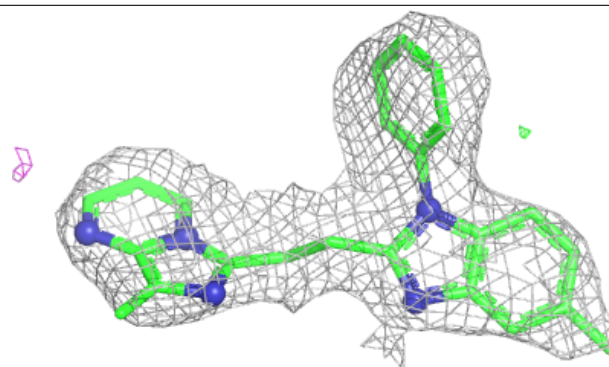
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	ZN	B	801	1/1	0.99	0.03	43,43,43,43	0
2	ZN	A	801	1/1	1.00	0.03	41,41,41,41	0

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.

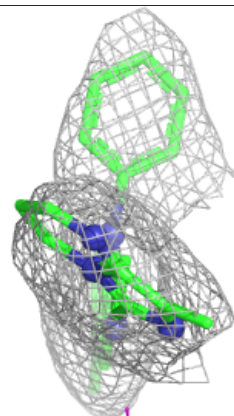
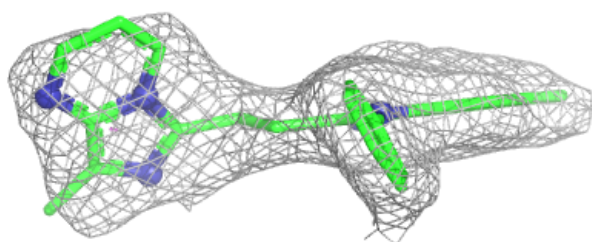
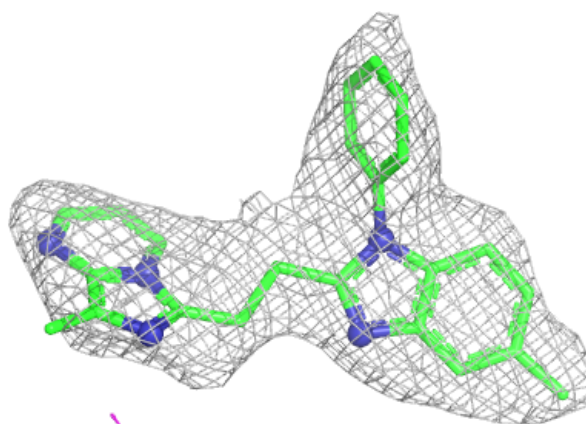
Electron density around X4C B 803:

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 X4C A 803:

$2mF_o-DF_c$ (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.