



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 03:59 PM UTC

PDB ID : 8IMG / pdb_00008img
Title : Human cGAS catalytic domain bound with C20
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Deposited on : 2023-03-06
Resolution : 1.80 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

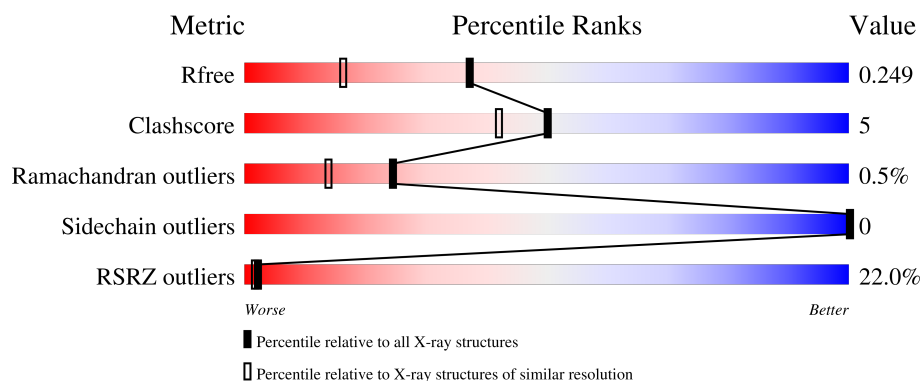
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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	366	21% 81%13%7%
1	B	366	20% 82%12%6%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 5632 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.

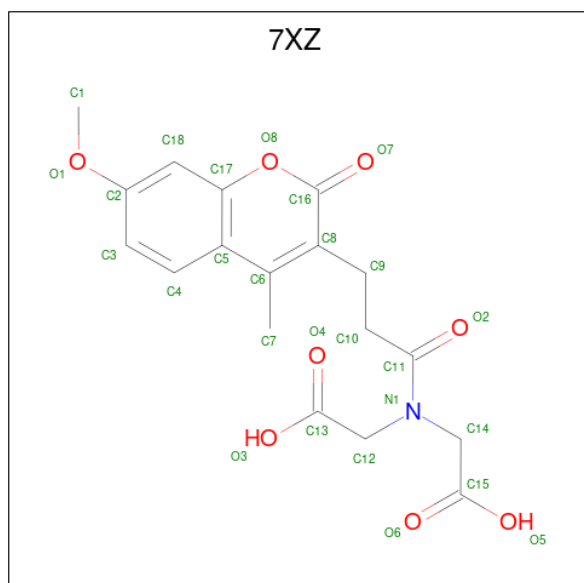
- Molecule 1 is a protein called Cyclic GMP-AMP synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	342	Total	C	N	O	S	0	0	0
			2720	1742	463	500	15			
1	B	343	Total	C	N	O	S	0	0	0
			2749	1760	466	508	15			

- 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 2-[2-hydroxy-2-oxoethyl]-[3-(7-methoxy-4-methyl-2-oxidanylidene-chromen-3-yl)propanoyl]amino]ethanoic acid (CCD ID: 7XZ) (formula: C₁₈H₁₉NO₈) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	C	N	O	0	0
			27	18	1	8		

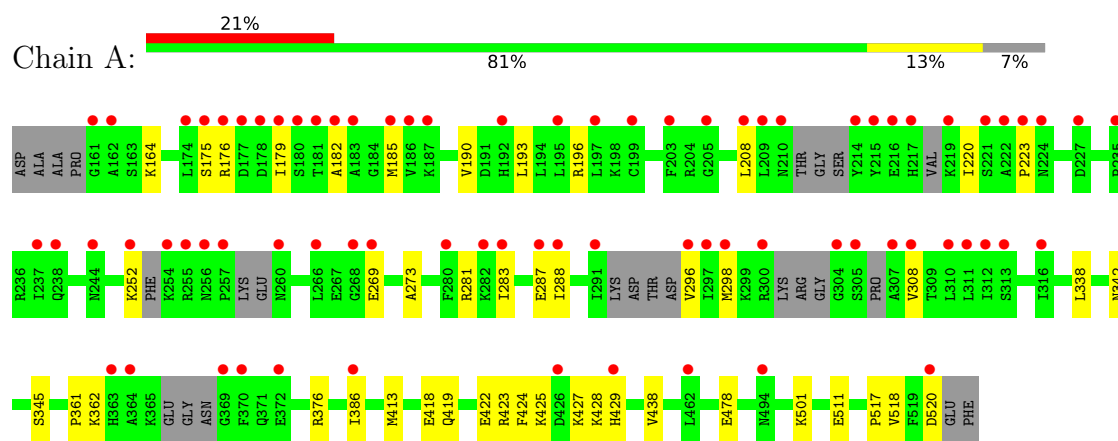
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	65	Total	O	0	0
			65	65		
4	B	69	Total	O	0	0
			69	69		

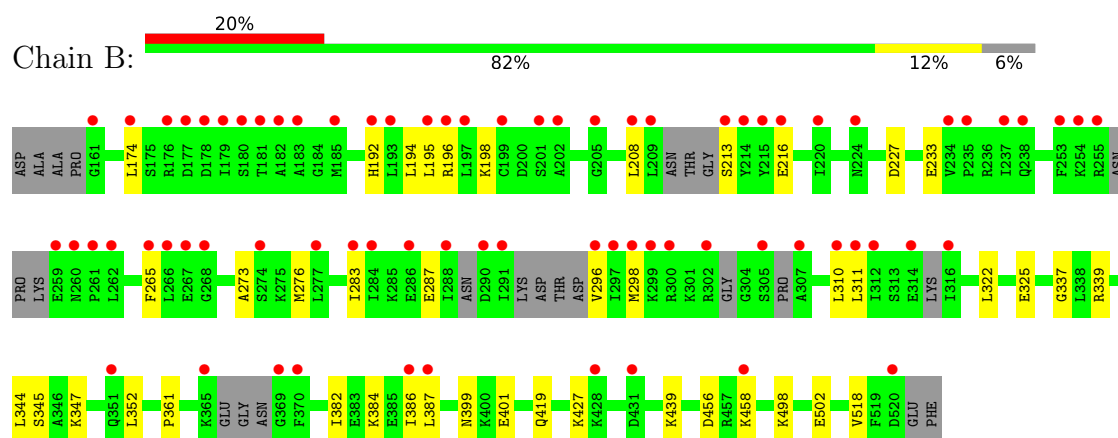
3 Residue-property plots [i](#)

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: Cyclic GMP-AMP synthase



• Molecule 1: Cyclic GMP-AMP synthase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	214.06Å 46.05Å 86.85Å 90.00° 113.94° 90.00°	Depositor
Resolution (Å)	35.68 – 1.80 35.68 – 1.80	Depositor EDS
% Data completeness (in resolution range)	99.2 (35.68-1.80) 99.7 (35.68-1.80)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.73 (at 1.79Å)	Xtriage
Refinement program	PHENIX (1.20.1_4487: ???)	Depositor
R, R_{free}	0.223 , 0.249 0.223 , 0.249	Depositor DCC
R_{free} test set	3556 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å ²)	26.2	Xtriage
Anisotropy	0.113	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 38.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.016 for -h-2*k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5632	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.44% 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: 7XZ, 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	0.28	0/2766	0.48	0/3712
1	B	0.27	0/2796	0.46	0/3749
All	All	0.28	0/5562	0.47	0/7461

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

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	2720	0	2679	28	0
1	B	2749	0	2703	30	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	B	27	0	0	0	0
4	A	65	0	0	1	0
4	B	69	0	0	0	0
All	All	5632	0	5382	58	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 5.

All (58) 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:179:ILE:HG21	1:A:223:PRO:HB2	1.51	0.92
1:B:213:SER:N	1:B:384:LYS:HZ1	1.87	0.73
1:B:276:MET:HE2	1:B:322:LEU:HD21	1.71	0.70
1:B:192:HIS:HA	1:B:195:LEU:HD12	1.75	0.69
1:B:216:GLU:HG2	1:B:227:ASP:HB3	1.74	0.67
1:B:387:LEU:HD12	1:B:439:LYS:HZ1	1.61	0.66
1:A:179:ILE:O	1:A:182:ALA:N	2.37	0.58
1:A:164:LYS:HE3	1:A:520:ASP:OD1	2.04	0.58
1:B:194:LEU:O	1:B:198:LYS:HG3	2.04	0.57
1:A:422:GLU:O	1:A:425:LYS:HG2	2.06	0.55
1:A:196:ARG:HE	1:A:283:ILE:HG23	1.72	0.55
1:A:273:ALA:HB1	1:A:361:PRO:HB3	1.89	0.54
1:A:511:GLU:OE2	1:A:517:PRO:HD2	2.09	0.53
1:A:362:LYS:HD3	1:A:376:ARG:CZ	2.39	0.53
1:B:273:ALA:HB1	1:B:361:PRO:HB3	1.91	0.52
1:A:176:ARG:HA	1:A:179:ILE:HD13	1.93	0.51
1:B:265:PHE:CD2	1:B:276:MET:HG2	2.46	0.51
1:B:298:MET:HG3	1:B:310:LEU:HD23	1.93	0.49
1:A:252:LYS:HG2	1:A:269:GLU:HG2	1.95	0.49
1:B:419:GLN:HB3	1:B:518:VAL:HG22	1.95	0.49
1:A:175:SER:O	1:A:179:ILE:HG23	2.13	0.48
1:A:338:LEU:HD11	1:A:386:ILE:HD11	1.95	0.48
1:A:342:ASN:HB2	4:A:727:HOH:O	2.13	0.48
1:A:298:MET:HE2	1:A:298:MET:HB2	1.76	0.47
1:A:413:MET:HB3	1:A:438:VAL:HG13	1.96	0.47
1:A:478:GLU:HG2	1:A:501:LYS:HE3	1.96	0.46
1:B:192:HIS:O	1:B:196:ARG:HG2	2.15	0.46
1:B:233:GLU:HA	1:B:325:GLU:HB3	1.98	0.46
1:B:498:LYS:HE3	1:B:502:GLU:OE2	2.16	0.46
1:A:419:GLN:HB3	1:A:518:VAL:HG22	1.97	0.46
1:A:220:ILE:HG22	1:A:418:GLU:OE1	2.15	0.46
1:A:427:LYS:C	1:A:429:HIS:H	2.25	0.45
1:B:347:LYS:HB2	1:B:347:LYS:HE2	1.80	0.45
1:B:387:LEU:HD12	1:B:439:LYS:NZ	2.31	0.44
1:A:422:GLU:OE2	1:A:425:LYS:HE3	2.17	0.44
1:B:387:LEU:CD1	1:B:439:LYS:HZ1	2.29	0.44
1:A:423:ARG:HD2	1:A:424:PHE:CE2	2.53	0.44
1:B:296:VAL:HA	1:B:311:LEU:O	2.18	0.44
1:B:344:LEU:CD1	1:B:386:ILE:HD12	2.49	0.43
1:A:362:LYS:HE3	1:A:362:LYS:HB3	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:427:LYS:HB3	1:B:427:LYS:HE3	1.84	0.43
1:B:399:ASN:OD1	1:B:401:GLU:HG2	2.18	0.43
1:B:456:ASP:OD1	1:B:458:LYS:HE2	2.18	0.43
1:A:185:MET:HE3	1:A:185:MET:HA	2.01	0.43
1:B:216:GLU:CG	1:B:227:ASP:HB3	2.45	0.43
1:B:352:LEU:CD1	1:B:386:ILE:HD11	2.48	0.43
1:B:174:LEU:HD23	1:B:174:LEU:HA	1.86	0.42
1:B:208:LEU:HA	1:B:208:LEU:HD23	1.75	0.42
1:B:382:ILE:O	1:B:386:ILE:HG12	2.20	0.42
1:B:283:ILE:O	1:B:287:GLU:HG2	2.19	0.42
1:B:337:GLY:HA2	1:B:339:ARG:HH21	1.85	0.41
1:A:190:VAL:HG11	1:A:208:LEU:HD11	2.03	0.41
1:A:193:LEU:HD23	1:A:287:GLU:HG2	2.01	0.41
1:A:288:ILE:HB	1:A:296:VAL:HB	2.03	0.41
1:B:276:MET:HE2	1:B:276:MET:HB3	1.93	0.41
1:B:298:MET:HE2	1:B:298:MET:HB3	1.75	0.41
1:A:427:LYS:O	1:A:429:HIS:N	2.55	0.40
1:A:281:ARG:HG2	1:A:308:VAL:HG23	2.04	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	324/366 (88%)	316 (98%)	6 (2%)	2 (1%)	21	11
1	B	325/366 (89%)	320 (98%)	4 (1%)	1 (0%)	36	25
All	All	649/732 (89%)	636 (98%)	10 (2%)	3 (0%)	24	14

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	345	SER
1	B	345	SER
1	A	428	LYS

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	295/336 (88%)	295 (100%)	0	100	100
1	B	299/336 (89%)	299 (100%)	0	100	100
All	All	594/672 (88%)	594 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	238	GLN
1	B	192	HIS
1	B	224	ASN
1	B	351	GLN
1	B	429	HIS
1	B	494	ASN
1	B	507	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

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 3 ligands modelled in this entry, 2 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	$\# Z > 2$
3	7XZ	B	601	-	28,28,28	0.85	0	37,39,39	1.20	3 (8%)

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
3	7XZ	B	601	-	-	13/19/19/19	0/2/2/2

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	601	7XZ	C15-C14-N1	3.48	116.75	112.91
3	B	601	7XZ	C12-N1-C14	-2.89	112.57	117.49
3	B	601	7XZ	C10-C9-C8	2.59	119.69	112.53

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	601	7XZ	C10-C11-N1-C14
3	B	601	7XZ	C10-C11-N1-C12
3	B	601	7XZ	O2-C11-N1-C14

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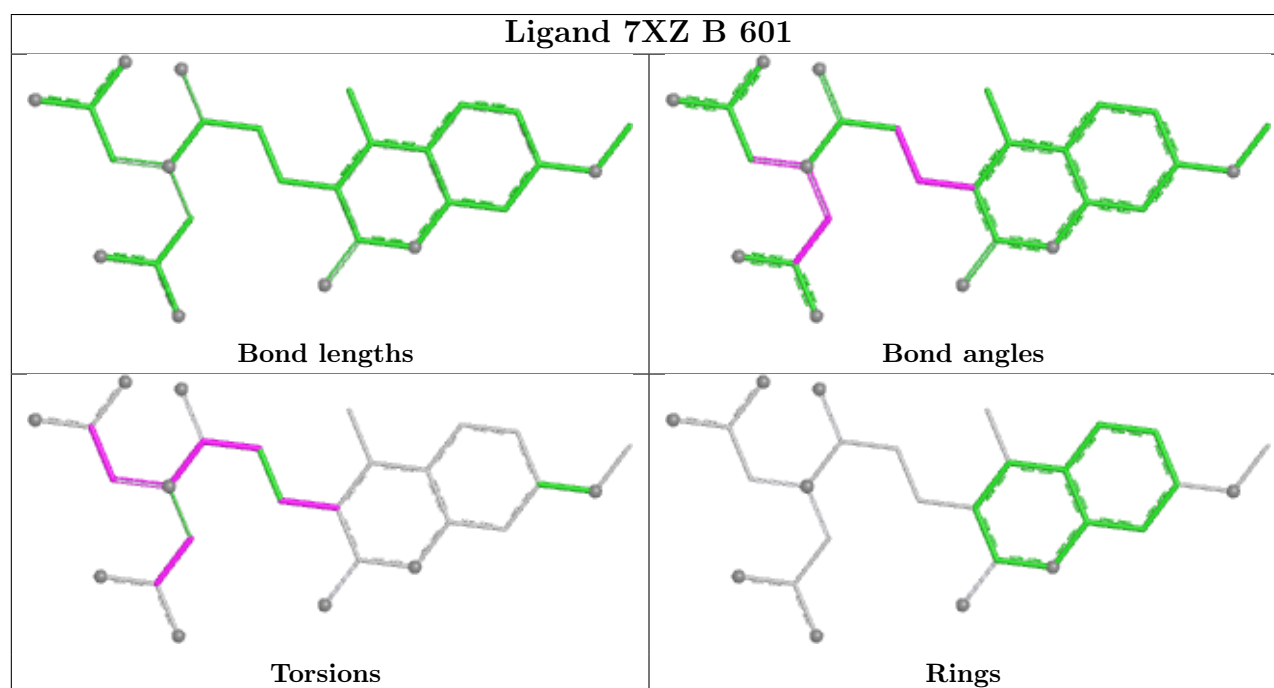
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Mol	Chain	Res	Type	Atoms
3	B	601	7XZ	O2-C11-N1-C12
3	B	601	7XZ	N1-C12-C13-O3
3	B	601	7XZ	C13-C12-N1-C14
3	B	601	7XZ	C13-C12-N1-C11
3	B	601	7XZ	N1-C12-C13-O4
3	B	601	7XZ	C9-C10-C11-N1
3	B	601	7XZ	C9-C10-C11-O2
3	B	601	7XZ	N1-C14-C15-O6
3	B	601	7XZ	N1-C14-C15-O5
3	B	601	7XZ	C16-C8-C9-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.



5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	342/366 (93%)	1.08	77 (22%) 2 2	18, 36, 74, 108	0
1	B	343/366 (93%)	1.08	74 (21%) 2 2	16, 36, 75, 100	0
All	All	685/732 (93%)	1.08	151 (22%) 2 2	16, 36, 75, 108	0

All (151) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	214	TYR	8.0
1	B	291	ILE	6.8
1	B	296	VAL	6.6
1	B	288	ILE	6.0
1	B	214	TYR	5.9
1	A	185	MET	5.8
1	B	520	ASP	5.7
1	A	179	ILE	5.3
1	A	215	TYR	5.2
1	A	291	ILE	5.0
1	A	296	VAL	5.0
1	A	311	LEU	4.8
1	B	311	LEU	4.7
1	B	297	ILE	4.7
1	A	307	ALA	4.6
1	B	179	ILE	4.5
1	B	215	TYR	4.4
1	B	255	ARG	4.4
1	B	370	PHE	4.3
1	A	257	PRO	4.3
1	A	369	GLY	4.1
1	B	181	THR	4.1
1	B	365	LYS	4.0
1	A	300	ARG	3.9

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Mol	Chain	Res	Type	RSRZ
1	B	266	LEU	3.9
1	A	180	SER	3.8
1	A	370	PHE	3.7
1	B	209	LEU	3.7
1	B	302	ARG	3.5
1	A	297	ILE	3.5
1	B	369	GLY	3.5
1	B	180	SER	3.4
1	A	298	MET	3.4
1	A	177	ASP	3.4
1	A	313	SER	3.4
1	B	310	LEU	3.3
1	B	312	ILE	3.3
1	B	237	ILE	3.2
1	A	181	THR	3.2
1	A	305	SER	3.2
1	A	426	ASP	3.2
1	A	429	HIS	3.2
1	A	183	ALA	3.2
1	B	216	GLU	3.2
1	B	284	ILE	3.2
1	A	209	LEU	3.2
1	B	195	LEU	3.2
1	B	298	MET	3.1
1	A	304	GLY	3.1
1	B	193	LEU	3.1
1	B	224	ASN	3.1
1	B	220	ILE	3.1
1	B	253	PHE	3.1
1	A	238	GLN	3.1
1	B	283	ILE	3.0
1	A	308	VAL	3.0
1	B	290	ASP	3.0
1	B	316	ILE	3.0
1	A	162	ALA	3.0
1	A	256	ASN	3.0
1	A	210	ASN	3.0
1	B	299	LYS	2.9
1	B	428	LYS	2.9
1	B	307	ALA	2.9
1	A	316	ILE	2.9
1	B	208	LEU	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	178	ASP	2.9
1	B	300	ARG	2.9
1	B	205	GLY	2.9
1	B	260	ASN	2.9
1	B	262	LEU	2.9
1	A	175	SER	2.9
1	B	314	GLU	2.9
1	A	260	ASN	2.8
1	B	386	ILE	2.8
1	B	213	SER	2.8
1	B	265	PHE	2.8
1	A	199	CYS	2.8
1	B	161	GLY	2.8
1	A	312	ILE	2.8
1	A	520	ASP	2.7
1	A	216	GLU	2.7
1	B	267	GLU	2.7
1	A	186	VAL	2.7
1	A	255	ARG	2.6
1	A	223	PRO	2.6
1	A	161	GLY	2.6
1	B	458	LYS	2.6
1	A	363	HIS	2.5
1	A	252	LYS	2.5
1	B	259	GLU	2.5
1	A	219	LYS	2.5
1	A	208	LEU	2.5
1	A	269	GLU	2.5
1	B	234	VAL	2.5
1	A	217	HIS	2.5
1	A	254	LYS	2.5
1	A	195	LEU	2.4
1	B	238	GLN	2.4
1	B	177	ASP	2.4
1	B	387	LEU	2.4
1	B	192	HIS	2.4
1	B	196	ARG	2.4
1	A	182	ALA	2.3
1	A	364	ALA	2.3
1	A	282	LYS	2.3
1	B	197	LEU	2.3
1	A	237	ILE	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	178	ASP	2.3
1	B	351	GLN	2.3
1	A	197	LEU	2.3
1	B	235	PRO	2.3
1	B	261	PRO	2.3
1	A	288	ILE	2.3
1	A	386	ILE	2.3
1	B	199	CYS	2.3
1	A	268	GLY	2.2
1	B	182	ALA	2.2
1	B	274	SER	2.2
1	A	192	HIS	2.2
1	A	310	LEU	2.2
1	B	183	ALA	2.2
1	A	227	ASP	2.2
1	B	268	GLY	2.2
1	B	277	LEU	2.2
1	B	254	LYS	2.2
1	A	266	LEU	2.1
1	B	286	GLU	2.1
1	A	222	ALA	2.1
1	A	224	ASN	2.1
1	A	244	ASN	2.1
1	A	494	ASN	2.1
1	A	203	PHE	2.1
1	B	431	ASP	2.1
1	A	221	SER	2.1
1	B	201	SER	2.1
1	B	176	ARG	2.1
1	A	187	LYS	2.1
1	A	287	GLU	2.1
1	A	462	LEU	2.1
1	B	174	LEU	2.1
1	B	305	SER	2.1
1	A	174	LEU	2.1
1	B	202	ALA	2.1
1	A	280	PHE	2.0
1	A	283	ILE	2.0
1	A	372	GLU	2.0
1	B	185	MET	2.0
1	A	205	GLY	2.0
1	A	176	ARG	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	235	PRO	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

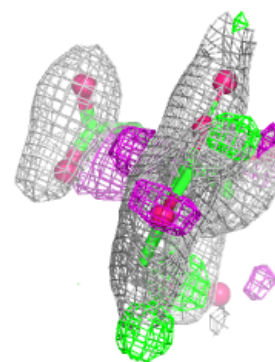
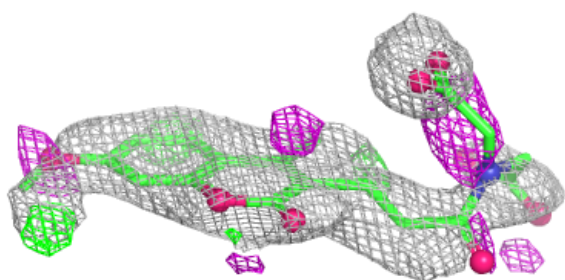
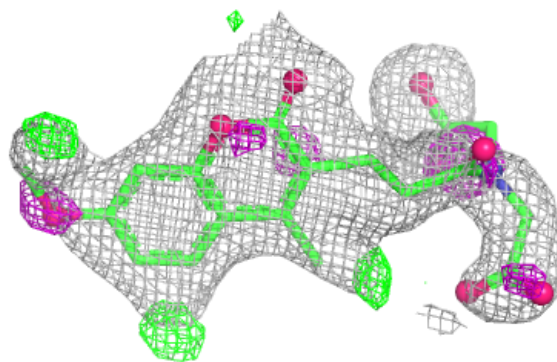
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	7XZ	B	601	27/27	0.68	0.20	47,55,70,70	0
2	ZN	B	602	1/1	0.99	0.05	19,19,19,19	0
2	ZN	A	601	1/1	0.99	0.04	19,19,19,19	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 7XZ B 601:

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



6.5 Other polymers ⓘ

There are no such residues in this entry.