



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 06:00 PM UTC

PDB ID : 6G3Y / pdb_00006g3y
Title : Structure of the mouse 8-oxoguanine DNA Glycosylase mOGG1 in complex with ligand TH5675
Authors : Masuyer, G.; Helleday, T.; Stenmark, P.
Deposited on : 2018-03-26
Resolution : 2.51 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

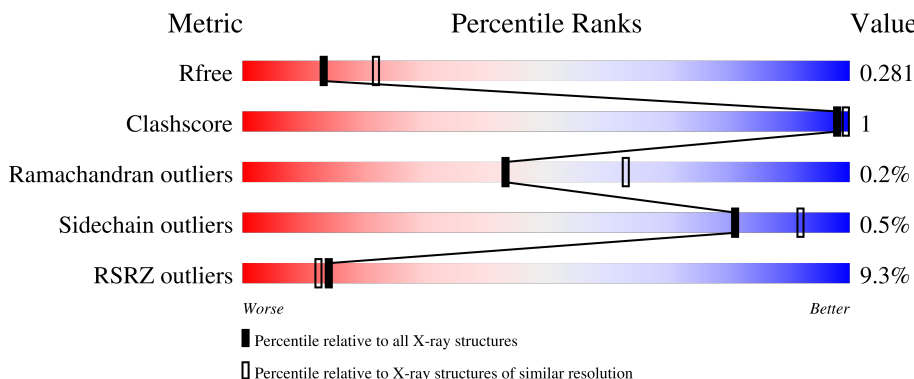
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.51 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	318	3% 96% ..
1	B	318	7% 97% ..
1	C	318	17% 97% ..

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 7750 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.

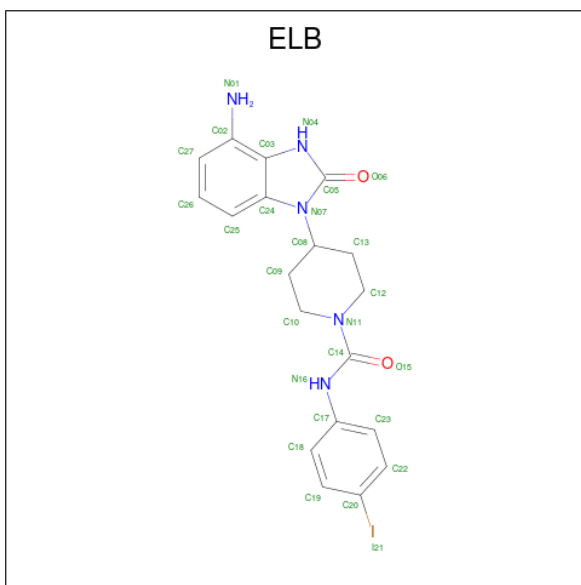
- Molecule 1 is a protein called N-glycosylase/DNA lyase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	316	Total	C	N	O	S	0	1	0
			2526	1605	459	451	11			
1	B	316	Total	C	N	O	S	0	0	0
			2515	1599	455	450	11			
1	C	315	Total	C	N	O	S	0	0	0
			2501	1591	452	447	11			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	8	GLY	-	expression tag	UNP O08760
A	10	HIS	SER	conflict	UNP O08760
B	8	GLY	-	expression tag	UNP O08760
B	10	HIS	SER	conflict	UNP O08760
C	8	GLY	-	expression tag	UNP O08760
C	10	HIS	SER	conflict	UNP O08760

- Molecule 2 is 4-(4-azanyl-2-oxidanylidene-3 {H}-benzimidazol-1-yl)- {N}-(4-iodophenyl)piperidine-1-carboxamide (CCD ID: ELB) (formula: C₁₉H₂₀IN₅O₂).

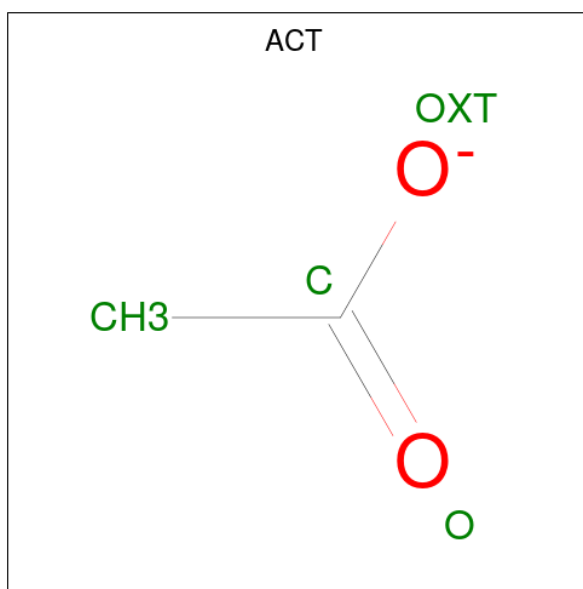


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	I	N	O	0	0
			27	19	1	5	2		
2	B	1	Total	C	I	N	O	0	0
			27	19	1	5	2		
2	C	1	Total	C	I	N	O	0	0
			27	19	1	5	2		

- Molecule 3 is NICKEL (II) ION (CCD ID: NI) (formula: Ni).

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

- Molecule 4 is ACETATE ION (CCD ID: ACT) (formula: C₂H₃O₂).



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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	50	Total	O	0	0
			50	50		
5	B	51	Total	O	0	0
			51	51		
5	C	21	Total	O	0	0
			21	21		

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.

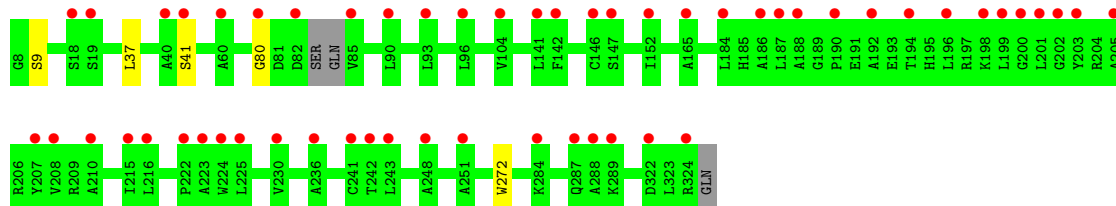
- Molecule 1: N-glycosylase/DNA lyase



- Molecule 1: N-glycosylase/DNA lyase



- Molecule 1: N-glycosylase/DNA lyase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	80.82Å 81.36Å 169.16Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	84.58 – 2.51 84.58 – 2.51	Depositor EDS
% Data completeness (in resolution range)	99.5 (84.58-2.51) 99.4 (84.58-2.51)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.61 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.8.0189	Depositor
R, R_{free}	0.234 , 0.282 0.236 , 0.281	Depositor DCC
R_{free} test set	1862 reflections (4.83%)	wwPDB-VP
Wilson B-factor (Å ²)	58.4	Xtriage
Anisotropy	0.069	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 46.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.014 for k,h,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7750	wwPDB-VP
Average B, all atoms (Å ²)	75.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.80% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ELB, NI, ACT

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	0.54	0/2597	0.79	0/3534
1	B	0.55	0/2586	0.82	2/3520 (0.1%)
1	C	0.55	0/2571	0.78	0/3498
All	All	0.55	0/7754	0.80	2/10552 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	19	SER	CA-C-N	5.18	124.97	119.28
1	B	19	SER	C-N-CA	5.18	124.97	119.28

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2526	0	2470	7	0
1	B	2515	0	2457	3	0
1	C	2501	0	2444	1	0
2	A	27	0	0	1	0
2	B	27	0	0	0	0
2	C	27	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
4	B	4	0	3	0	0
5	A	50	0	0	0	0
5	B	51	0	0	0	0
5	C	21	0	0	0	0
All	All	7750	0	7374	10	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.

All (10) 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:278:ASP:HA	1:B:273:GLN:HE22	1.74	0.53
1:A:34:ARG:HA	1:A:68:GLN:HE22	1.76	0.50
1:A:146:CYS:HA	1:A:204:ARG:HD3	1.95	0.48
1:A:249:LYS:NZ	2:A:401:ELB:O15	2.47	0.47
1:A:80:GLY:HA2	1:A:171:ILE:HD12	1.98	0.46
1:C:37:LEU:O	1:C:41:SER:HB2	2.16	0.46
1:B:80:GLY:N	1:B:171:ILE:HD11	2.33	0.44
1:A:80:GLY:HA2	1:A:171:ILE:HB	2.01	0.41
1:A:265:VAL:HG13	1:A:315:GLN:OE1	2.21	0.40
1:B:36:ASP:HA	1:B:55:TRP:CZ2	2.56	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	315/318 (99%)	305 (97%)	9 (3%)	1 (0%)	36 55
1	B	314/318 (99%)	302 (96%)	12 (4%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	311/318 (98%)	302 (97%)	8 (3%)	1 (0%)	36	55
All	All	940/954 (98%)	909 (97%)	29 (3%)	2 (0%)	43	63

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	82	ASP
1	C	80	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	266/266 (100%)	265 (100%)	1 (0%)	84	93
1	B	265/266 (100%)	264 (100%)	1 (0%)	84	93
1	C	263/266 (99%)	261 (99%)	2 (1%)	73	88
All	All	794/798 (100%)	790 (100%)	4 (0%)	81	92

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	58	VAL
1	B	272	TRP
1	C	9	SER
1	C	272	TRP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	50	GLN
1	A	68	GLN
1	A	164	GLN
1	A	226	GLN

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Mol	Chain	Res	Type
1	A	270	HIS
1	B	68	GLN
1	B	119	HIS
1	B	164	GLN
1	B	270	HIS
1	B	273	GLN
1	C	68	GLN
1	C	112	HIS
1	C	125	GLN
1	C	135	GLN
1	C	150	ASN
1	C	287	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, 1 is monoatomic - leaving 4 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	ACT	B	402	-	3,3,3	0.83	0	3,3,3	0.59	0
2	ELB	A	401	-	30,30,30	1.89	7 (23%)	42,43,43	2.68	14 (33%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	ELB	B	401	-	30,30,30	2.00	7 (23%)	42,43,43	2.58	13 (30%)
2	ELB	C	401	-	30,30,30	2.00	7 (23%)	42,43,43	2.69	16 (38%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	ELB	A	401	-	-	0/12/22/22	0/4/4/4
2	ELB	B	401	-	-	0/12/22/22	0/4/4/4
2	ELB	C	401	-	-	2/12/22/22	0/4/4/4

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	401	ELB	C14-N16	4.81	1.45	1.37
2	C	401	ELB	C14-N11	4.60	1.44	1.36
2	B	401	ELB	C14-N11	4.48	1.44	1.36
2	A	401	ELB	C14-N16	4.45	1.45	1.37
2	C	401	ELB	C05-N04	4.43	1.41	1.37
2	B	401	ELB	C05-N04	4.42	1.41	1.37
2	A	401	ELB	C14-N11	4.16	1.43	1.36
2	C	401	ELB	C14-N16	4.08	1.44	1.37
2	A	401	ELB	C05-N04	4.03	1.41	1.37
2	C	401	ELB	C24-N07	3.60	1.47	1.40
2	C	401	ELB	C02-C03	3.47	1.43	1.40
2	B	401	ELB	C02-C03	3.29	1.43	1.40
2	A	401	ELB	C24-N07	3.25	1.46	1.40
2	B	401	ELB	C24-N07	3.17	1.46	1.40
2	A	401	ELB	C02-C03	2.98	1.43	1.40
2	C	401	ELB	C25-C24	2.84	1.44	1.39
2	A	401	ELB	C03-N04	2.79	1.44	1.38
2	C	401	ELB	C03-N04	2.71	1.44	1.38
2	B	401	ELB	C25-C24	2.69	1.43	1.39
2	B	401	ELB	C03-N04	2.61	1.44	1.38
2	A	401	ELB	C25-C24	2.52	1.43	1.39

All (43) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	ELB	C13-C08-N07	-8.37	91.49	112.27
2	C	401	ELB	C13-C08-N07	-8.31	91.63	112.27
2	B	401	ELB	C13-C08-N07	-7.83	92.82	112.27
2	A	401	ELB	C09-C08-N07	7.70	131.39	112.27
2	C	401	ELB	C09-C08-N07	7.38	130.59	112.27
2	B	401	ELB	C09-C08-N07	6.64	128.76	112.27
2	A	401	ELB	C09-C10-N11	-6.08	98.76	110.66
2	B	401	ELB	C09-C10-N11	-5.46	99.97	110.66
2	C	401	ELB	C24-N07-C05	-5.17	105.76	109.24
2	C	401	ELB	C09-C10-N11	-5.10	100.67	110.66
2	B	401	ELB	C24-N07-C05	-4.85	105.98	109.24
2	A	401	ELB	C24-N07-C05	-4.60	106.15	109.24
2	C	401	ELB	O06-C05-N04	-4.14	123.91	127.49
2	A	401	ELB	C13-C12-N11	-4.12	102.59	110.66
2	B	401	ELB	N16-C14-N11	4.09	120.71	115.91
2	A	401	ELB	O06-C05-N04	-4.02	124.02	127.49
2	B	401	ELB	C03-N04-C05	-3.74	107.80	109.82
2	C	401	ELB	C03-N04-C05	-3.63	107.86	109.82
2	B	401	ELB	O06-C05-N04	-3.54	124.43	127.49
2	B	401	ELB	C13-C12-N11	-3.49	103.83	110.66
2	A	401	ELB	C03-N04-C05	-3.48	107.94	109.82
2	C	401	ELB	C13-C12-N11	-3.22	104.36	110.66
2	B	401	ELB	C17-N16-C14	-3.20	119.57	126.04
2	B	401	ELB	C09-C08-C13	2.97	117.89	111.17
2	C	401	ELB	C17-N16-C14	-2.94	120.09	126.04
2	B	401	ELB	C03-C24-N07	2.86	108.59	106.52
2	C	401	ELB	C03-C24-N07	2.81	108.56	106.52
2	B	401	ELB	O15-C14-N11	-2.79	117.77	121.67
2	A	401	ELB	N16-C14-N11	2.72	119.11	115.91
2	C	401	ELB	C09-C08-C13	2.67	117.21	111.17
2	A	401	ELB	C17-N16-C14	-2.66	120.66	126.04
2	A	401	ELB	C03-C24-N07	2.66	108.45	106.52
2	C	401	ELB	N16-C14-N11	2.65	119.03	115.91
2	C	401	ELB	C19-C20-I21	-2.55	115.88	119.69
2	C	401	ELB	N04-C05-N07	2.53	109.74	106.69
2	C	401	ELB	C18-C17-N16	-2.46	112.16	120.41
2	B	401	ELB	N04-C05-N07	2.31	109.48	106.69
2	A	401	ELB	C09-C08-C13	2.30	116.37	111.17
2	A	401	ELB	N04-C05-N07	2.25	109.40	106.69
2	C	401	ELB	C22-C20-I21	2.20	122.97	119.69
2	A	401	ELB	C23-C17-N16	-2.18	113.08	120.41
2	A	401	ELB	O15-C14-N11	-2.17	118.64	121.67
2	C	401	ELB	C25-C24-C03	-2.13	118.31	121.06

There are no chirality outliers.

All (2) torsion outliers are listed below:

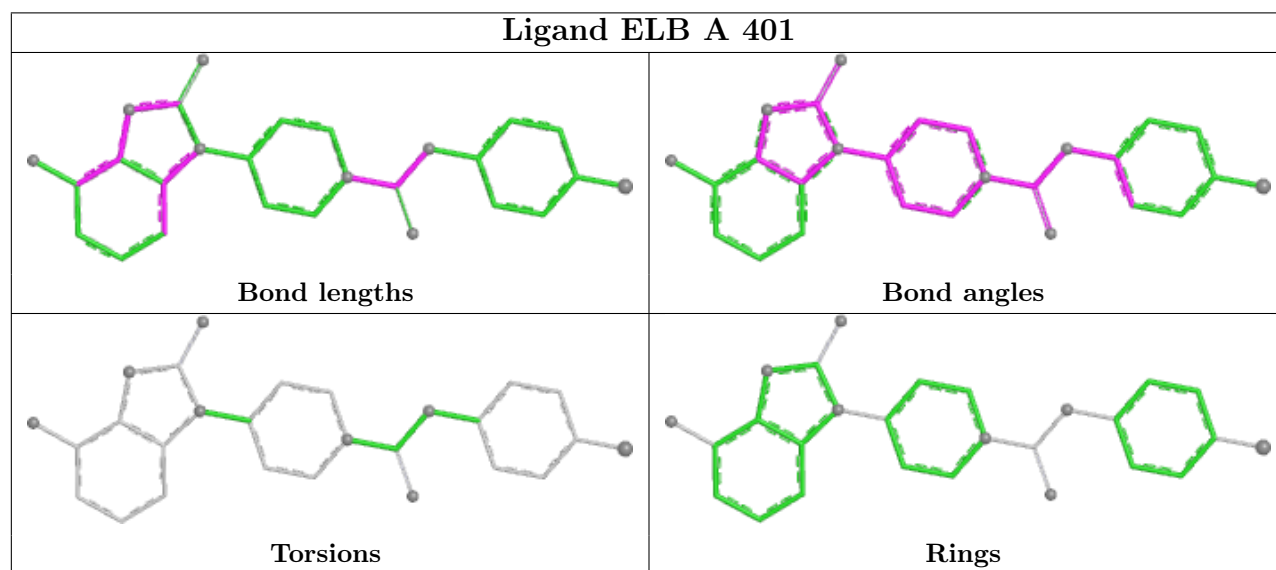
Mol	Chain	Res	Type	Atoms
2	C	401	ELB	C18-C17-N16-C14
2	C	401	ELB	C23-C17-N16-C14

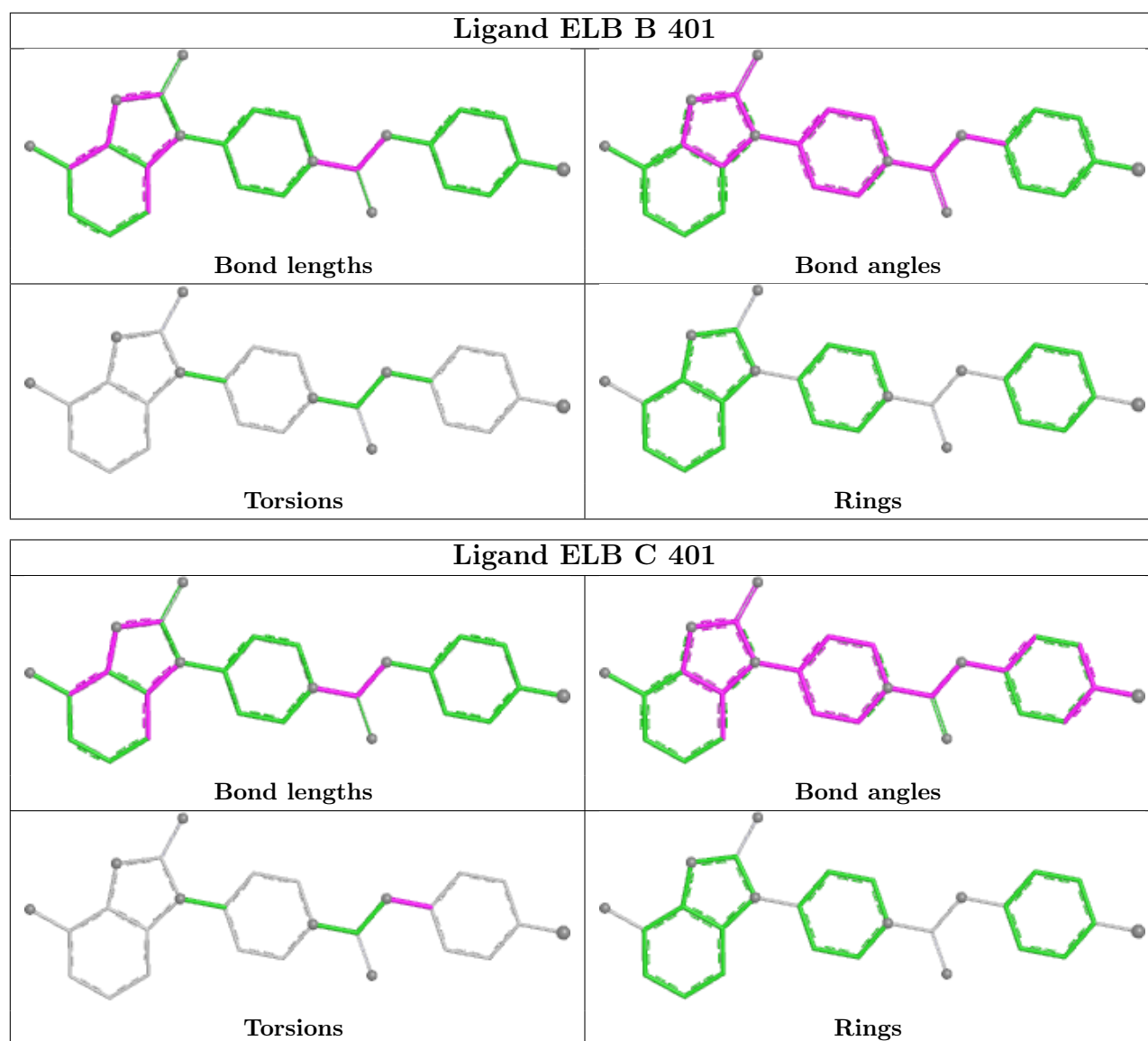
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	ELB	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

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	316/318 (99%)	0.24	10 (3%) 50 46	23, 54, 100, 145	2 (0%)
1	B	316/318 (99%)	0.84	23 (7%) 21 19	43, 74, 114, 141	1 (0%)
1	C	315/318 (99%)	1.12	55 (17%) 4 3	43, 84, 156, 181	1 (0%)
All	All	947/954 (99%)	0.73	88 (9%) 14 12	23, 71, 127, 181	4 (0%)

All (88) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	289	LYS	7.4
1	C	289	LYS	6.9
1	A	289	LYS	6.7
1	B	16	LEU	4.2
1	B	322	ASP	3.9
1	C	142	PHE	3.9
1	A	285	THR	3.9
1	B	85	VAL	3.9
1	A	288	ALA	3.7
1	C	205	ALA	3.5
1	C	85	VAL	3.4
1	C	199	LEU	3.3
1	C	141	LEU	3.2
1	C	190	PRO	3.2
1	B	285	THR	3.2
1	C	152	ILE	3.2
1	C	200	GLY	3.2
1	C	202	GLY	3.1
1	A	290	GLY	3.1
1	C	208	VAL	3.1
1	C	203	TYR	3.0
1	B	288	ALA	2.9
1	C	90	LEU	2.9

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Mol	Chain	Res	Type	RSRZ
1	C	147	SER	2.8
1	C	251	ALA	2.8
1	A	20	PRO	2.8
1	C	222	PRO	2.8
1	A	59	LEU	2.8
1	C	225	LEU	2.8
1	C	96	LEU	2.8
1	B	186	ALA	2.7
1	A	11	MET	2.7
1	C	80	GLY	2.7
1	B	11	MET	2.7
1	C	324	ARG	2.7
1	B	103	ASP	2.6
1	B	187	LEU	2.6
1	B	321	ALA	2.6
1	C	165	ALA	2.6
1	B	171	ILE	2.6
1	C	284	LYS	2.6
1	A	291	PRO	2.6
1	C	82	ASP	2.6
1	C	18	SER	2.5
1	B	207	TYR	2.5
1	C	194	THR	2.5
1	C	243	LEU	2.5
1	C	248	ALA	2.5
1	C	216	LEU	2.4
1	C	104	VAL	2.4
1	C	41	SER	2.4
1	C	186	ALA	2.4
1	A	10	HIS	2.3
1	C	196	LEU	2.3
1	C	201	LEU	2.3
1	C	40	ALA	2.3
1	C	223	ALA	2.3
1	B	147	SER	2.3
1	C	187	LEU	2.3
1	C	322	ASP	2.3
1	C	287	GLN	2.3
1	B	26	ILE	2.3
1	B	323	LEU	2.2
1	C	224	TRP	2.2
1	C	210	ALA	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	146	CYS	2.2
1	C	236	ALA	2.2
1	C	242	THR	2.2
1	C	184	LEU	2.2
1	C	230	VAL	2.1
1	C	192	ALA	2.1
1	C	215	ILE	2.1
1	B	73	LEU	2.1
1	B	324	ARG	2.1
1	C	198	LYS	2.1
1	C	188	ALA	2.1
1	C	19	SER	2.1
1	C	241	CYS	2.1
1	C	93	LEU	2.1
1	B	71	ASP	2.1
1	C	207	TYR	2.1
1	C	60	ALA	2.0
1	C	288	ALA	2.0
1	B	93	LEU	2.0
1	B	272	TRP	2.0
1	B	222	PRO	2.0
1	B	170	LEU	2.0
1	A	83	SER	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.

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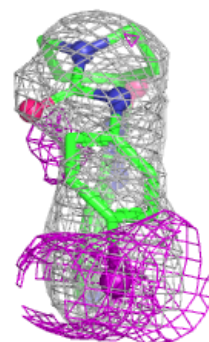
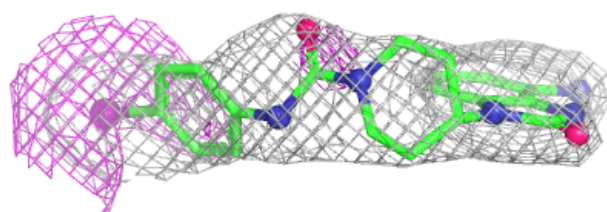
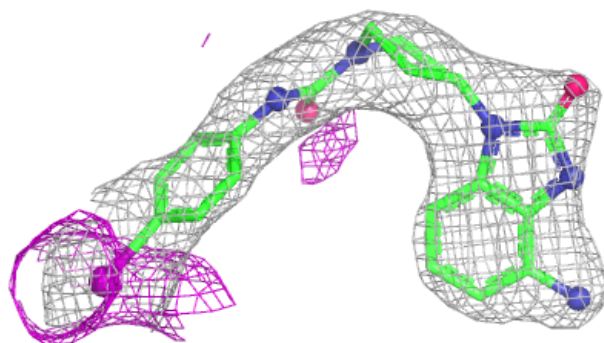
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	ACT	B	402	4/4	0.79	0.15	58,61,61,68	0
2	ELB	C	401	27/27	0.89	0.13	72,80,100,119	0
2	ELB	B	401	27/27	0.92	0.14	59,63,72,88	0
2	ELB	A	401	27/27	0.95	0.12	57,59,64,73	0
3	NI	A	402	1/1	1.00	0.03	40,40,40,40	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.

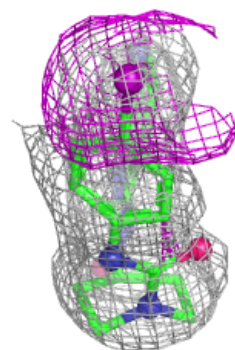
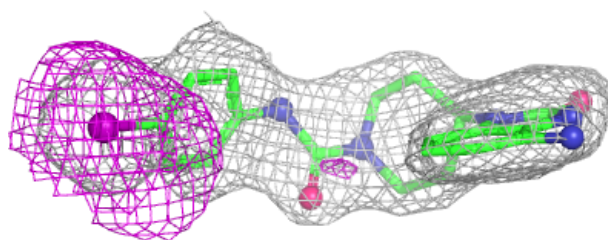
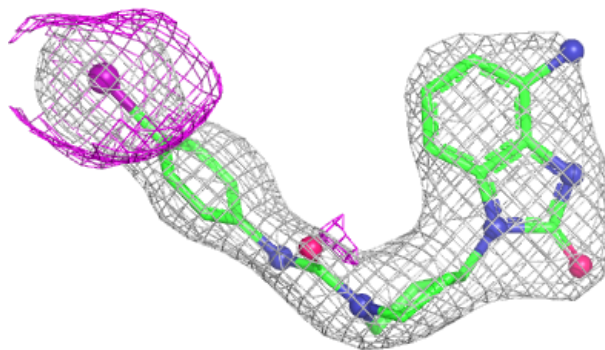
Electron density around ELB C 401:

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

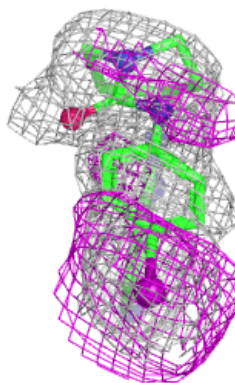
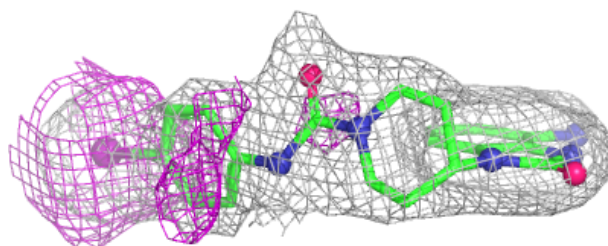
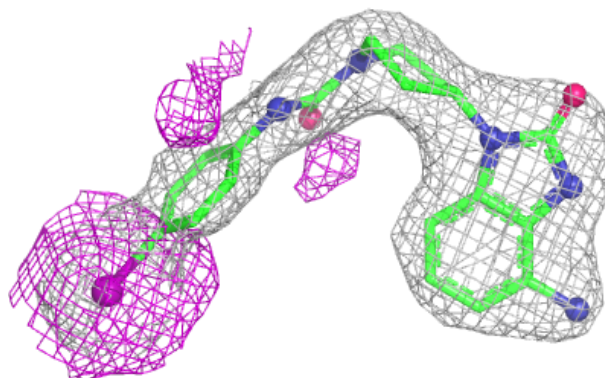


Electron density around ELB B 401:

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

**Electron density around ELB A 401:**

$2mF_o-DF_c$ (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.