



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 22, 2026 – 06:21 AM EDT

PDB ID : 11LJ / pdb_000011lj
Title : Human OGA IN COMPLEX WITH LIGAND 24
Authors : Shaffer, P.L.; Cedervall, P.
Deposited on : 2026-03-03
Resolution : 2.33 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

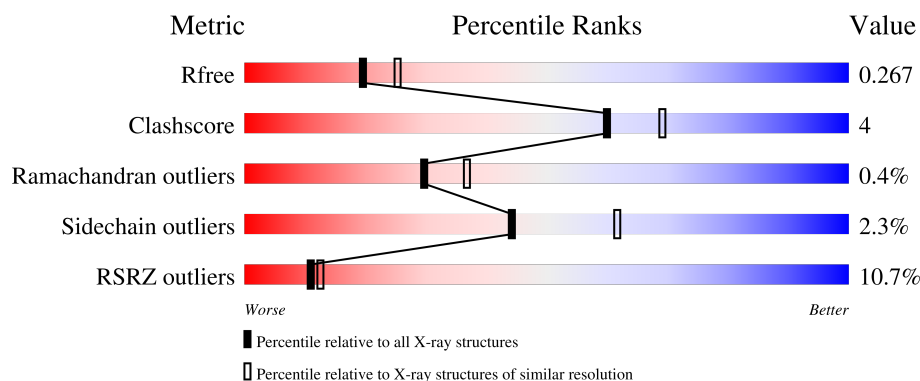
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.33 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	7754 (2.34-2.30)
Clashscore	190562	8383 (2.34-2.30)
Ramachandran outliers	187476	8303 (2.34-2.30)
Sidechain outliers	187428	8303 (2.34-2.30)
RSRZ outliers	180081	7760 (2.34-2.30)

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	386	<div> <div>5%</div> <div> <div>69%</div> <div>10%</div> <div>•</div> <div>21%</div> </div> </div>
1	B	386	<div> <div>6%</div> <div> <div>68%</div> <div>8%</div> <div>•</div> <div>24%</div> </div> </div>
2	C	170	<div> <div>20%</div> <div> <div>56%</div> <div>11%</div> <div>32%</div> </div> </div>
2	D	170	<div> <div>9%</div> <div> <div>64%</div> <div>14%</div> <div>•</div> <div>21%</div> </div> </div>

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 7063 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.

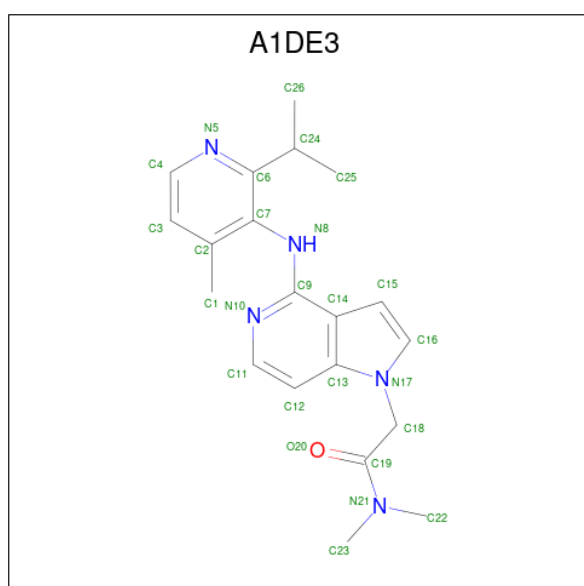
- Molecule 1 is a protein called Protein O-GlcNAcase N-Terminal Fragment.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	305	Total	C	N	O	S	0	0	0
			2501	1621	410	456	14			
1	B	295	Total	C	N	O	S	0	0	0
			2430	1577	401	440	12			

- Molecule 2 is a protein called Protein O-GlcNAcase 535-704 Peptide.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	115	Total	C	N	O	S	0	0	0
			932	603	153	166	10			
2	D	134	Total	C	N	O	S	0	0	0
			1095	710	183	192	10			

- Molecule 3 is N,N-dimethyl-2-(4-{[4-methyl-2-(propan-2-yl)pyridin-3-yl]amino}-1H-pyrrolo[3,2-c]pyridin-1-yl)acetamide (CCD ID: A1DE3) (formula: C₂₀H₂₅N₅O) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			26	20	5	1		
3	B	1	Total	C	N	O	0	0
			26	20	5	1		

- Molecule 4 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Ca	0	0
			1	1		
4	B	1	Total	Ca	0	0
			1	1		

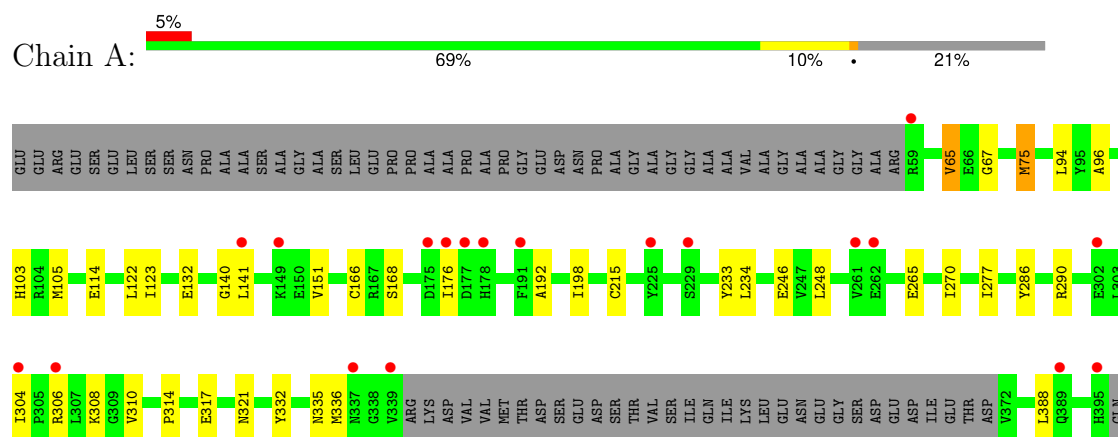
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	27	Total	O	0	0
			27	27		
5	C	4	Total	O	0	0
			4	4		
5	B	15	Total	O	0	0
			15	15		
5	D	5	Total	O	0	0
			5	5		

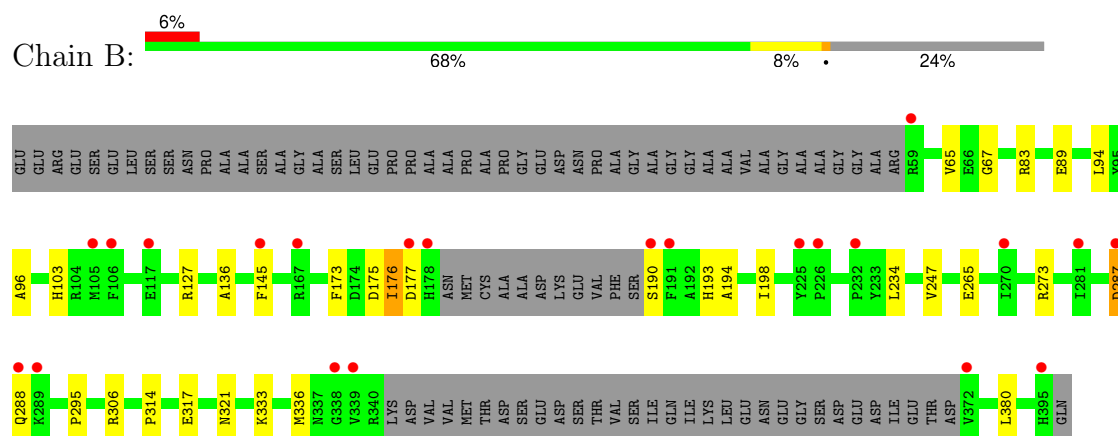
3 Residue-property plots [i](#)

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.

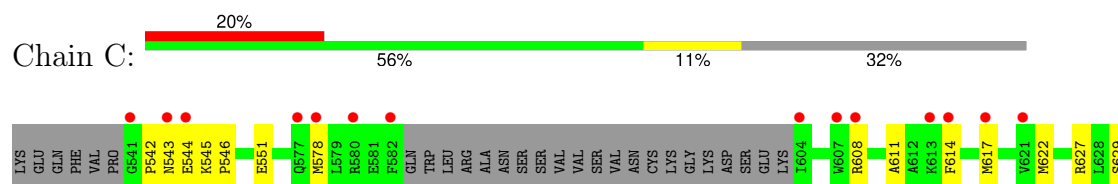
• Molecule 1: Protein O-GlcNAcase N-Terminal Fragment

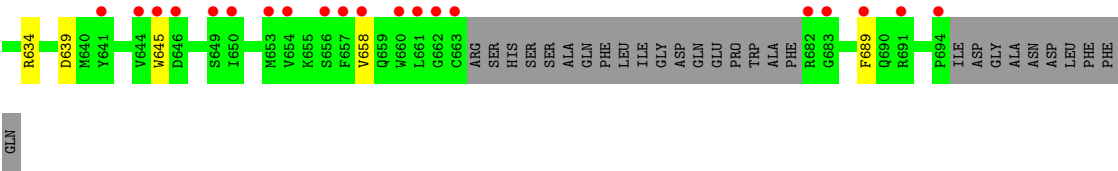


• Molecule 1: Protein O-GlcNAcase N-Terminal Fragment

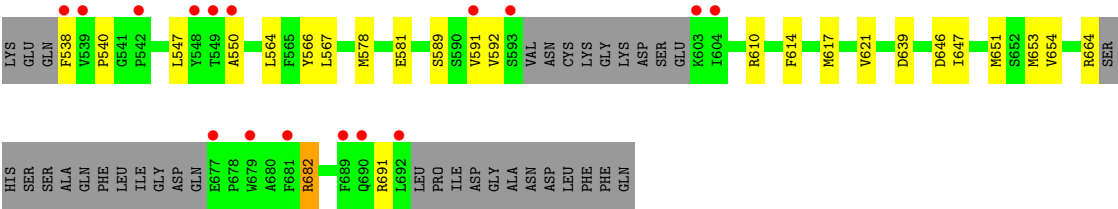


• Molecule 2: Protein O-GlcNAcase 535-704 Peptide





● Molecule 2: Protein O-GlcNAcase 535-704 Peptide



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	96.84Å 96.84Å 257.11Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	60.44 – 2.33 60.44 – 2.33	Depositor EDS
% Data completeness (in resolution range)	93.8 (60.44-2.33) 69.8 (60.44-2.33)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.57 (at 2.32Å)	Xtriage
Refinement program	PHENIX 1.13_2998	Depositor
R, R_{free}	0.223 , 0.268 0.224 , 0.267	Depositor DCC
R_{free} test set	1847 reflections (3.45%)	wwPDB-VP
Wilson B-factor (Å ²)	51.1	Xtriage
Anisotropy	0.053	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 42.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7063	wwPDB-VP
Average B, all atoms (Å ²)	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.41% 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: A1DE3, CA

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	0.09	0/2570	0.29	0/3486
1	B	0.09	0/2497	0.28	0/3386
2	C	0.09	0/954	0.27	0/1284
2	D	0.08	0/1124	0.24	0/1516
All	All	0.09	0/7145	0.27	0/9672

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 [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	2501	0	2454	21	0
1	B	2430	0	2390	17	0
2	C	932	0	918	10	0
2	D	1095	0	1074	16	0
3	A	26	0	0	1	0
3	B	26	0	0	1	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	27	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	B	15	0	0	0	0
5	C	4	0	0	0	0
5	D	5	0	0	0	0
All	All	7063	0	6836	59	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 4.

All (59) 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)
2:D:578:MET:HE2	2:D:617:MET:HB3	1.67	0.77
1:A:286:TYR:O	2:D:682:ARG:NH1	2.17	0.76
2:C:645:TRP:HE1	1:B:287:ASP:N	1.87	0.73
2:C:578:MET:HE2	2:C:617:MET:HB3	1.76	0.68
1:A:265:GLU:OE2	1:A:306:ARG:NH1	2.26	0.67
1:A:141:LEU:HD12	1:A:176:ILE:HD13	1.77	0.66
1:A:105:MET:HE1	2:D:550:ALA:HB2	1.79	0.65
2:D:614:PHE:HA	2:D:617:MET:HE2	1.81	0.63
2:D:589:SER:O	2:D:664:ARG:NH1	2.32	0.58
2:C:608:ARG:HA	2:C:611:ALA:HB3	1.86	0.57
2:C:545:LYS:HD2	2:C:546:PRO:HD2	1.87	0.57
1:A:246:GLU:HG2	1:A:308:LYS:HG2	1.88	0.56
1:B:333:LYS:HG3	1:B:336:MET:HE2	1.89	0.55
2:D:646:ASP:OD1	2:D:682:ARG:NH2	2.40	0.55
2:C:645:TRP:HE1	1:B:287:ASP:H	1.54	0.54
1:A:215:CYS:HB2	1:A:248:LEU:HD12	1.90	0.53
2:D:653:MET:HE1	2:D:682:ARG:HD2	1.91	0.52
2:D:591:VAL:HG23	2:D:592:VAL:HG13	1.92	0.51
3:B:801:A1DE3:C15	3:B:801:A1DE3:C7	2.91	0.49
1:A:317:GLU:HB3	2:C:639:ASP:HB3	1.94	0.49
2:C:551:GLU:HG2	2:C:627:ARG:HG2	1.95	0.49
2:D:621:VAL:HG11	2:D:651:MET:HG3	1.96	0.48
1:A:304:ILE:HD12	1:A:335:ASN:HB3	1.97	0.47
1:B:265:GLU:OE1	1:B:306:ARG:NH2	2.47	0.47
1:B:247:VAL:HG13	1:B:273:ARG:HD2	1.96	0.47
1:A:151:VAL:HG22	1:A:198:ILE:HD13	1.96	0.47
1:B:173:PHE:HZ	1:B:198:ILE:HD11	1.80	0.46
1:A:192:ALA:HB2	1:A:233:TYR:CD2	2.51	0.46
1:A:277:ILE:HB	1:A:310:VAL:HG22	1.98	0.46
1:B:295:PRO:HD3	1:B:380:LEU:HD22	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:317:GLU:HG2	2:D:639:ASP:O	2.15	0.46
2:C:689:PHE:HB3	2:D:654:VAL:HG21	1.97	0.46
1:A:75:MET:HE1	1:A:122:LEU:HB2	1.97	0.46
3:A:801:A1DE3:C7	3:A:801:A1DE3:C15	2.94	0.45
1:A:140:GLY:O	1:A:176:ILE:HG12	2.17	0.45
1:B:190:SER:HB3	1:B:193:HIS:ND1	2.32	0.45
1:B:96:ALA:HB2	1:B:136:ALA:HB3	2.00	0.44
1:B:94:LEU:HD11	1:B:136:ALA:HB2	1.98	0.44
1:A:65:VAL:HB	1:A:94:LEU:HB3	1.98	0.44
2:C:629:SER:O	2:C:634:ARG:NH2	2.37	0.44
1:A:314:PRO:HB2	1:A:321:ASN:CG	2.42	0.44
2:D:581:GLU:OE1	2:D:610:ARG:NH2	2.36	0.44
1:A:332:TYR:CZ	1:A:336:MET:HG3	2.53	0.44
2:D:564:LEU:HD13	2:D:647:ILE:HD13	2.01	0.43
1:A:132:GLU:OE2	1:A:168:SER:OG	2.30	0.43
1:B:89:GLU:OE1	1:B:333:LYS:NZ	2.52	0.42
1:A:314:PRO:HB2	1:A:321:ASN:OD1	2.20	0.42
2:D:538:PHE:HD2	2:D:547:LEU:HD21	1.84	0.42
1:B:67:GLY:HA2	1:B:96:ALA:O	2.20	0.42
1:B:314:PRO:HB2	1:B:321:ASN:OD1	2.19	0.42
1:B:145:PHE:CG	1:B:194:ALA:HB1	2.55	0.41
1:B:127:ARG:HB3	1:B:127:ARG:NH1	2.35	0.41
2:D:540:PRO:HG3	2:D:547:LEU:HG	2.02	0.41
2:C:614:PHE:CE2	2:C:658:VAL:HG21	2.55	0.41
2:D:566:TYR:HB3	2:D:567:LEU:H	1.73	0.41
1:A:67:GLY:HA2	1:A:96:ALA:O	2.21	0.40
1:B:176:ILE:HG13	1:B:177:ASP:H	1.86	0.40
1:A:123:ILE:HD13	1:A:166:CYS:HB2	2.04	0.40
1:A:234:LEU:HB3	1:A:270:ILE:HG21	2.03	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	301/386 (78%)	288 (96%)	13 (4%)	0	100	100
1	B	289/386 (75%)	271 (94%)	17 (6%)	1 (0%)	36	45
2	C	109/170 (64%)	106 (97%)	1 (1%)	2 (2%)	6	5
2	D	128/170 (75%)	125 (98%)	3 (2%)	0	100	100
All	All	827/1112 (74%)	790 (96%)	34 (4%)	3 (0%)	30	37

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	C	544	GLU
2	C	542	PRO
1	B	287	ASP

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	273/330 (83%)	267 (98%)	6 (2%)	45	63
1	B	265/330 (80%)	258 (97%)	7 (3%)	40	57
2	C	99/147 (67%)	97 (98%)	2 (2%)	48	66
2	D	116/147 (79%)	114 (98%)	2 (2%)	53	70
All	All	753/954 (79%)	736 (98%)	17 (2%)	44	62

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	65	VAL
1	A	75	MET
1	A	103	HIS
1	A	114	GLU
1	A	290	ARG
1	A	388	LEU
2	C	543	ASN
2	C	622	MET

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Mol	Chain	Res	Type
1	B	65	VAL
1	B	83	ARG
1	B	103	HIS
1	B	175	ASP
1	B	176	ILE
1	B	234	LEU
1	B	288	GLN
2	D	682	ARG
2	D	691	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	315	ASN
2	C	559	GLN
1	B	288	GLN
1	B	315	ASN
2	D	690	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 4 ligands modelled in this entry, 2 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
3	A1DE3	B	801	-	28,28,28	1.09	3 (10%)	34,40,40	1.45	3 (8%)
3	A1DE3	A	801	-	28,28,28	1.09	3 (10%)	34,40,40	1.42	5 (14%)

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	A1DE3	B	801	-	-	3/16/25/25	0/3/3/3
3	A1DE3	A	801	-	-	2/16/25/25	0/3/3/3

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	801	A1DE3	C18-C19	-2.35	1.49	1.53
3	B	801	A1DE3	C18-C19	-2.35	1.49	1.53
3	A	801	A1DE3	C19-N21	2.19	1.37	1.34
3	B	801	A1DE3	C19-N21	2.18	1.37	1.34
3	A	801	A1DE3	C9-N8	2.13	1.40	1.36
3	B	801	A1DE3	C9-N8	2.10	1.40	1.36

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	801	A1DE3	C18-C19-N21	-4.57	111.55	116.69
3	A	801	A1DE3	C18-C19-N21	-4.30	111.85	116.69
3	B	801	A1DE3	C15-C16-N17	-2.74	107.41	110.54
3	A	801	A1DE3	C15-C16-N17	-2.65	107.52	110.54
3	B	801	A1DE3	O20-C19-N21	2.27	125.49	121.89
3	A	801	A1DE3	O20-C19-N21	2.16	125.31	121.89
3	A	801	A1DE3	C19-C18-N17	-2.03	108.16	110.81
3	A	801	A1DE3	C11-N10-C9	2.00	120.64	116.74

There are no chirality outliers.

All (5) torsion outliers are listed below:

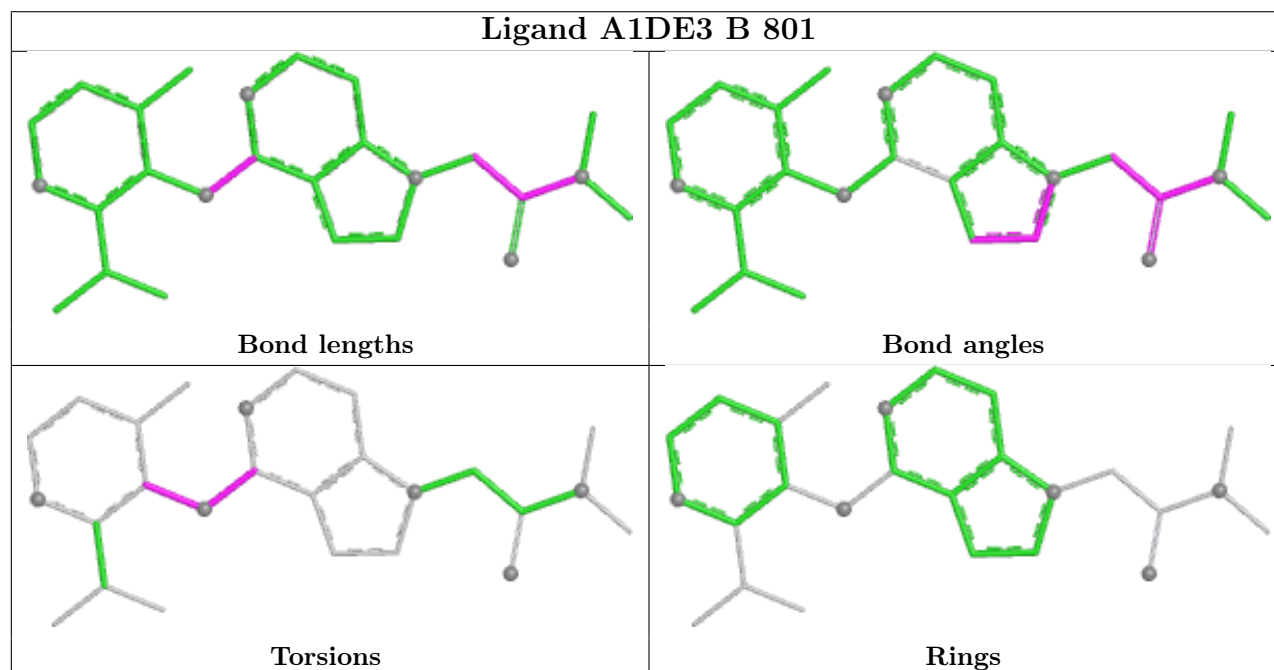
Mol	Chain	Res	Type	Atoms
3	A	801	A1DE3	N10-C9-N8-C7
3	B	801	A1DE3	N10-C9-N8-C7
3	A	801	A1DE3	C14-C9-N8-C7
3	B	801	A1DE3	C14-C9-N8-C7
3	B	801	A1DE3	C6-C7-N8-C9

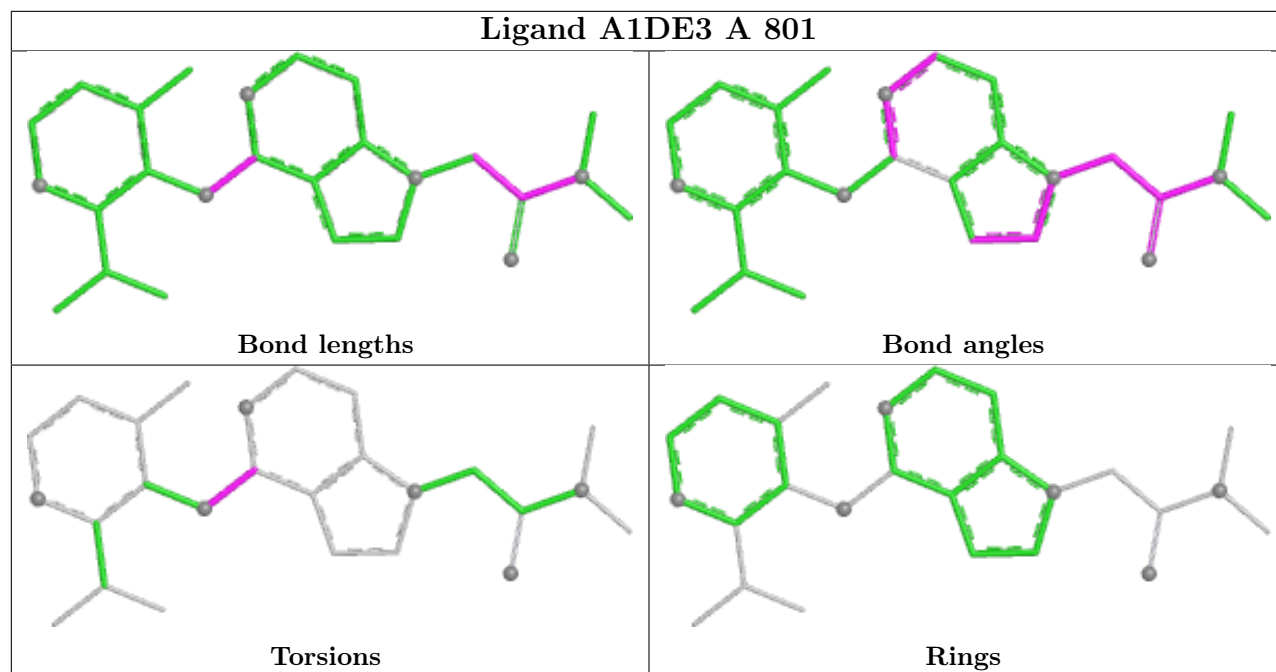
There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	801	A1DE3	1	0
3	A	801	A1DE3	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(Å²)	Q<0.9
1	A	305/386 (79%)	0.51	19 (6%)	26 29	29, 52, 88, 131	0
1	B	295/386 (76%)	0.51	22 (7%)	20 23	27, 51, 88, 120	0
2	C	115/170 (67%)	1.55	34 (29%)	1 1	44, 74, 122, 150	0
2	D	134/170 (78%)	0.79	16 (11%)	9 10	30, 54, 101, 120	0
All	All	849/1112 (76%)	0.70	91 (10%)	11 12	27, 55, 100, 150	0

All (91) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	C	582	PHE	5.8
2	C	691	ARG	5.4
1	B	287	ASP	5.4
2	C	644	VAL	4.9
1	B	395	HIS	4.9
2	C	604	ILE	4.6
2	C	607	TRP	4.6
2	D	603	LYS	4.5
1	A	262	GLU	4.4
2	D	550	ALA	4.3
2	D	593	SER	4.3
1	A	261	VAL	4.1
2	C	650	ILE	4.0
1	B	339	VAL	3.9
2	D	681	PHE	3.7
2	C	645	TRP	3.6
1	B	225	TYR	3.6
2	D	548	TYR	3.6
2	C	661	LEU	3.6
2	C	663	CYS	3.6
1	A	339	VAL	3.6

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Mol	Chain	Res	Type	RSRZ
1	B	178	HIS	3.5
2	C	614	PHE	3.5
2	C	580	ARG	3.4
1	A	306	ARG	3.3
2	C	683	GLY	3.3
1	A	395	HIS	3.2
2	C	543	ASN	3.2
2	D	692	LEU	3.1
1	A	178	HIS	3.0
2	D	549	THR	2.9
2	D	591	VAL	2.9
1	A	176	ILE	2.9
2	C	544	GLU	2.9
2	D	677	GLU	2.9
2	D	604	ILE	2.9
2	C	694	PRO	2.9
2	D	690	GLN	2.9
2	C	646	ASP	2.9
2	C	662	GLY	2.8
1	B	167	ARG	2.8
2	C	660	TRP	2.7
2	C	617	MET	2.7
2	C	613	LYS	2.7
1	B	191	PHE	2.7
1	B	117	GLU	2.7
2	D	539	VAL	2.7
2	D	542	PRO	2.7
1	B	289	LYS	2.6
2	D	538	PHE	2.6
2	C	658	VAL	2.6
2	D	689	PHE	2.6
2	C	608	ARG	2.6
1	B	270	ILE	2.6
1	A	229	SER	2.6
2	C	656	SER	2.5
1	A	175	ASP	2.5
1	B	288	GLN	2.5
1	A	59	ARG	2.5
2	C	682	ARG	2.5
1	A	302	GLU	2.4
1	A	191	PHE	2.4
1	B	145	PHE	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	281	ILE	2.4
2	D	679	TRP	2.4
1	B	106	PHE	2.4
1	A	177	ASP	2.4
1	B	177	ASP	2.3
2	C	653	MET	2.3
2	C	621	VAL	2.3
1	B	226	PRO	2.3
1	B	338	GLY	2.3
2	C	541	GLY	2.3
2	C	654	VAL	2.3
1	B	59	ARG	2.2
1	A	389	GLN	2.2
1	A	149	LYS	2.2
1	B	190	SER	2.2
2	C	578	MET	2.2
1	A	225	TYR	2.1
2	C	657	PHE	2.1
2	C	689	PHE	2.1
1	B	105	MET	2.1
2	C	641	TYR	2.1
1	B	372	VAL	2.1
2	C	649	SER	2.1
1	A	141	LEU	2.1
2	C	577	GLN	2.0
1	A	337	ASN	2.0
1	B	232	PRO	2.0
1	A	304	ILE	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.

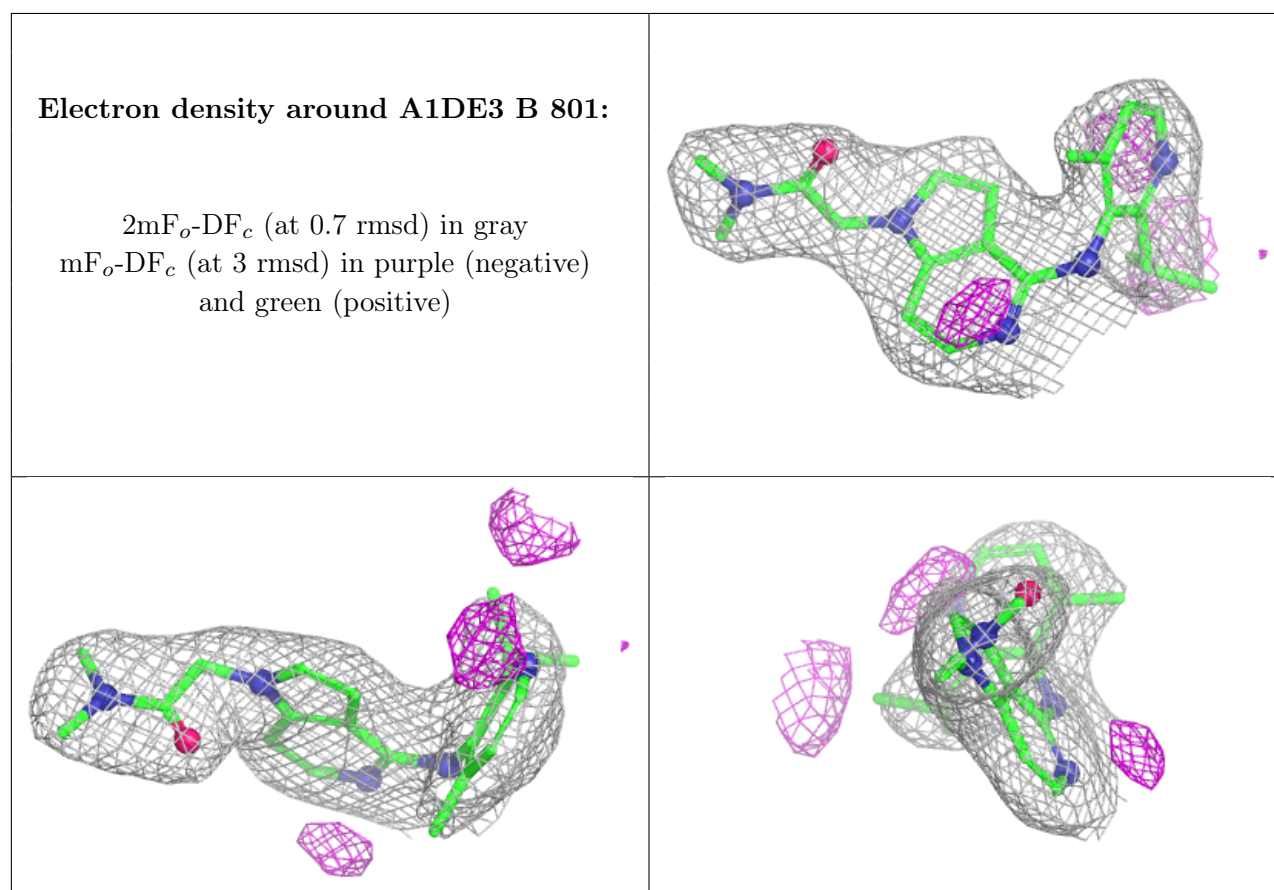
6.4 Ligands [i](#)

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.

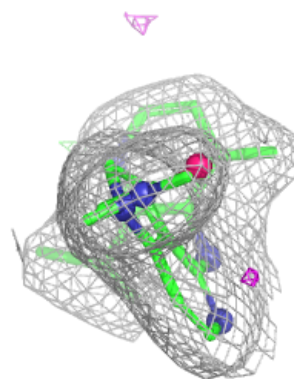
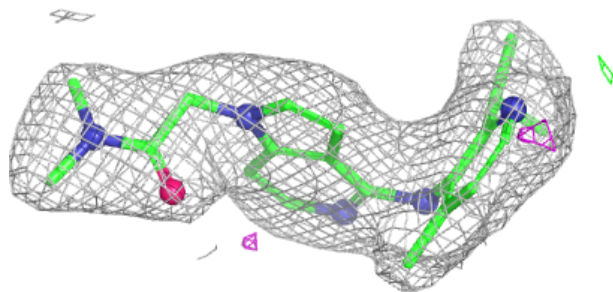
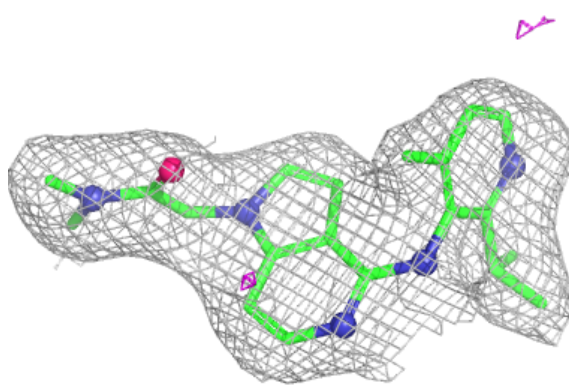
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	CA	A	802	1/1	0.92	0.10	66,66,66,66	0
3	A1DE3	B	801	26/26	0.93	0.10	25,39,52,54	0
4	CA	B	802	1/1	0.94	0.08	64,64,64,64	0
3	A1DE3	A	801	26/26	0.95	0.08	20,35,47,51	0

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 A1DE3 A 801:

$2mF_o-DF_c$ (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.