



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:08 AM UTC

PDB ID : 6EG3 / pdb_00006eg3
Title : Crystal structure of human BRM in complex with compound 15
Authors : Zhu, X.; Kulathila, R.; Hu, T.; Xie, X.
Deposited on : 2018-08-17
Resolution : 2.84 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

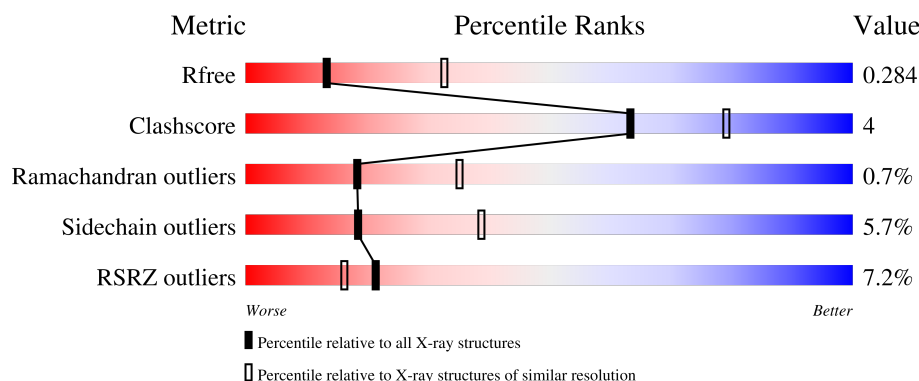
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.84 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	1520 (2.86-2.82)
Clashscore	190562	1559 (2.86-2.82)
Ramachandran outliers	187476	1517 (2.86-2.82)
Sidechain outliers	187428	1518 (2.86-2.82)
RSRZ outliers	180081	1521 (2.86-2.82)

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	621	7% 83% 15% ..

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4931 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.

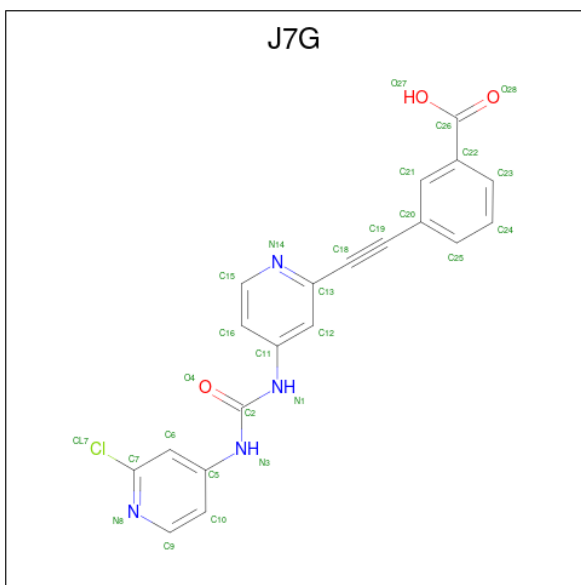
- Molecule 1 is a protein called Maltose/maltodextrin-binding periplasmic protein, Probable global transcription activator SNF2L2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	614	Total	C	N	O	S	0	0	0
			4858	3143	814	886	15			

There are 9 discrepancies between the modelled and reference sequences:

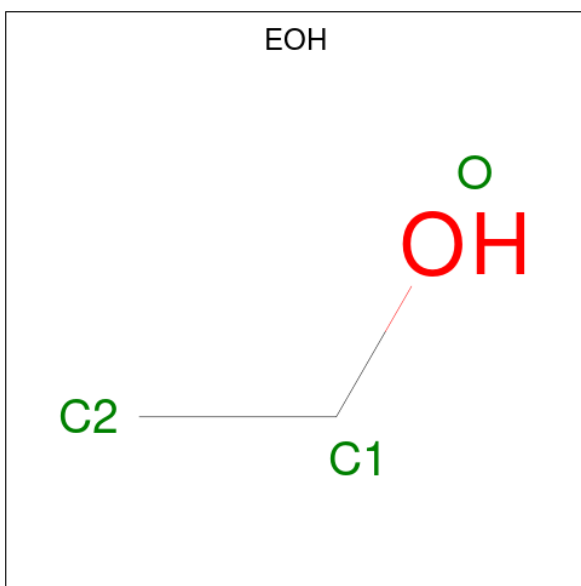
Chain	Residue	Modelled	Actual	Comment	Reference
A	416	ALA	ASP	conflict	UNP P0AEY0
A	417	ALA	LYS	conflict	UNP P0AEY0
A	506	ALA	GLU	conflict	UNP P0AEY0
A	507	ALA	ASN	conflict	UNP P0AEY0
A	573	ALA	LYS	conflict	UNP P0AEY0
A	701	ASN	-	linker	UNP P0AEY0
A	702	ALA	-	linker	UNP P0AEY0
A	703	ALA	-	linker	UNP P0AEY0
A	704	ALA	-	linker	UNP P0AEY0

- Molecule 2 is 3-[(4-{[(2-chloropyridin-4-yl)carbamoyl]amino}pyridin-2-yl)ethynyl]benzoic acid (CCD ID: J7G) (formula: C₂₀H₁₃ClN₄O₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	Cl	N	O	
			28	20	1	4	3	

- Molecule 3 is ETHANOL (CCD ID: EOH) (formula: C₂H₆O).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O		
			3	2	1	0	0
3	A	1	Total	C	O		
			3	2	1	0	0
3	A	1	Total	C	O		
			3	2	1	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			3	2	1		
3	A	1	Total	C	O	0	0
			3	2	1		
3	A	1	Total	C	O	0	0
			3	2	1		

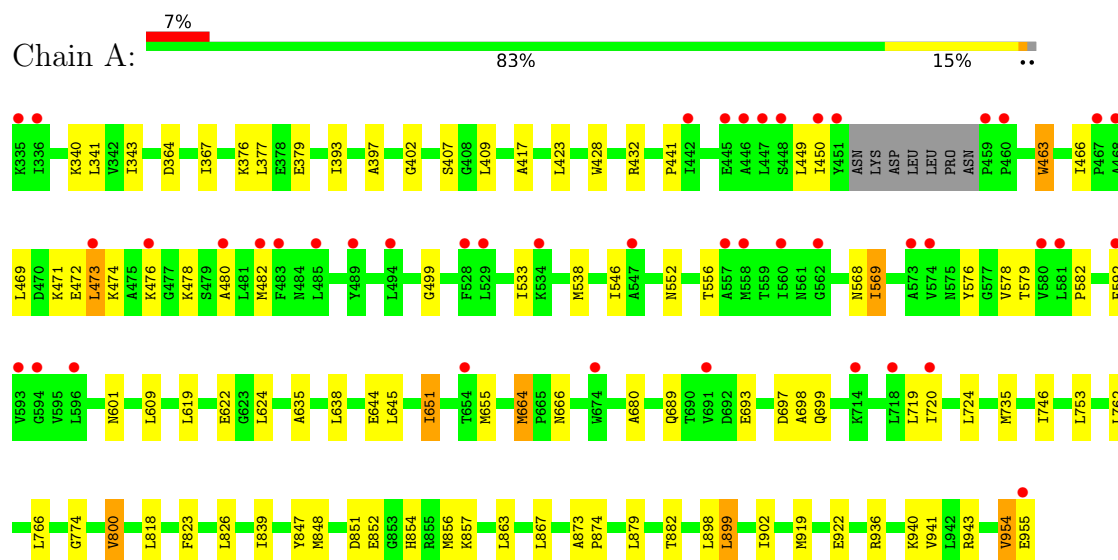
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	27	Total	O	0	0
			27	27		

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: Maltose/maltodextrin-binding periplasmic protein, Probable global transcription activator SNF2L2



4 Data and refinement statistics

Property	Value	Source
Space group	P 42 21 2	Depositor
Cell constants a, b, c, α , β , γ	136.47Å 136.47Å 121.37Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	96.50 – 2.84 96.50 – 2.84	Depositor EDS
% Data completeness (in resolution range)	99.6 (96.50-2.84) 100.0 (96.50-2.84)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.37 (at 2.86Å)	Xtriage
Refinement program	BUSTER 2.11.6	Depositor
R, R_{free}	0.229 , 0.267 0.241 , 0.284	Depositor DCC
R_{free} test set	1323 reflections (4.80%)	wwPDB-VP
Wilson B-factor (Å ²)	72.8	Xtriage
Anisotropy	0.513	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 75.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4931	wwPDB-VP
Average B, all atoms (Å ²)	78.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.16% 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: EOH, J7G

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	0.81	0/4975	1.34	16/6750 (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	579	THR	CA-C-N	5.80	128.41	120.35
1	A	579	THR	C-N-CA	5.80	128.41	120.35
1	A	397	ALA	CA-C-N	5.65	128.63	120.38
1	A	397	ALA	C-N-CA	5.65	128.63	120.38
1	A	364	ASP	CA-CB-CG	5.62	118.22	112.60
1	A	624	LEU	CA-C-N	5.62	128.07	120.38
1	A	624	LEU	C-N-CA	5.62	128.07	120.38
1	A	463	TRP	CA-C-N	5.41	128.28	120.38
1	A	463	TRP	C-N-CA	5.41	128.28	120.38
1	A	697	ASP	CA-CB-CG	5.32	117.92	112.60
1	A	800	VAL	N-CA-CB	5.22	116.95	111.00
1	A	601	ASN	CA-CB-CG	5.18	117.78	112.60
1	A	940	LYS	CA-C-N	5.14	127.71	120.42
1	A	940	LYS	C-N-CA	5.14	127.71	120.42
1	A	417	ALA	CA-C-N	5.11	127.38	120.38
1	A	417	ALA	C-N-CA	5.11	127.38	120.38

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	4858	0	4910	39	0
2	A	28	0	0	0	0
3	A	18	0	36	0	0
4	A	27	0	0	0	0
All	All	4931	0	4946	39	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 (39) 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:954:VAL:HG12	1:A:955:GLU:H	1.27	0.93
1:A:954:VAL:HG12	1:A:955:GLU:N	1.95	0.80
1:A:376:LYS:HD3	1:A:379:GLU:OE1	1.88	0.73
1:A:852:GLU:OE1	1:A:856:MET:HB3	1.89	0.71
1:A:635:ALA:HB1	1:A:655:MET:HE2	1.78	0.64
1:A:449:LEU:HD22	1:A:582:PRO:HD3	1.84	0.60
1:A:552:ASN:ND2	1:A:569:ILE:HG23	2.17	0.59
1:A:848:MET:HE3	1:A:873:ALA:HB3	1.83	0.58
1:A:423:LEU:HD12	1:A:428:TRP:CZ2	2.39	0.57
1:A:954:VAL:CG1	1:A:955:GLU:N	2.67	0.57
1:A:473:LEU:HD13	1:A:480:ALA:HA	1.89	0.55
1:A:569:ILE:HB	1:A:576:TYR:CE2	2.42	0.55
1:A:343:ILE:HG12	1:A:393:ILE:HB	1.89	0.54
1:A:899:LEU:HB3	1:A:902:ILE:HD13	1.90	0.53
1:A:851:ASP:HB2	1:A:879:LEU:HD12	1.92	0.51
1:A:645:LEU:HB3	1:A:651:ILE:HG21	1.92	0.51
1:A:450:ILE:HG12	1:A:578:VAL:HG22	1.93	0.50
1:A:689:GLN:HE21	1:A:693:GLU:HB3	1.77	0.50
1:A:818:LEU:HD21	1:A:826:LEU:HD13	1.95	0.49
1:A:367:ILE:HG21	1:A:609:LEU:HD21	1.94	0.48
1:A:746:ILE:HD11	1:A:898:LEU:HD11	1.96	0.48
1:A:340:LYS:HG2	1:A:341:LEU:N	2.29	0.48
1:A:800:VAL:HG23	1:A:823:PHE:HA	1.95	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:735:MET:HB3	1:A:762:LEU:HD13	1.96	0.46
1:A:402:GLY:HA3	1:A:666:ASN:HB2	1.98	0.46
1:A:423:LEU:HD13	1:A:441:PRO:HG2	1.99	0.45
1:A:407:SER:HB2	1:A:409:LEU:HG	1.98	0.44
1:A:719:LEU:HD21	1:A:724:LEU:HD13	1.99	0.44
1:A:423:LEU:HD23	1:A:638:LEU:HA	1.99	0.44
1:A:680:ALA:HB2	1:A:698:ALA:HB2	1.99	0.43
1:A:592:PHE:HB3	1:A:664:MET:HG2	1.99	0.43
1:A:619:LEU:HD23	1:A:619:LEU:HA	1.94	0.43
1:A:463:TRP:HA	1:A:466:ILE:HD12	2.00	0.43
1:A:482:MET:HE1	1:A:546:ILE:HG22	2.01	0.42
1:A:482:MET:HB2	1:A:556:THR:HG21	2.02	0.42
1:A:473:LEU:HD22	1:A:478:LYS:HB2	2.02	0.42
1:A:854:HIS:O	1:A:882:THR:HG23	2.20	0.41
1:A:533:ILE:HD13	1:A:538:MET:HB3	2.01	0.41
1:A:473:LEU:CD2	1:A:478:LYS:HB2	2.50	0.41

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	610/621 (98%)	576 (94%)	30 (5%)	4 (1%)	18 35

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	954	VAL
1	A	499	GLY
1	A	774	GLY
1	A	874	PRO

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	510/517 (99%)	481 (94%)	29 (6%)	18	39

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	377	LEU
1	A	432	ARG
1	A	469	LEU
1	A	471	LYS
1	A	472	GLU
1	A	473	LEU
1	A	474	LYS
1	A	476	LYS
1	A	568	ASN
1	A	569	ILE
1	A	622	GLU
1	A	644	GLU
1	A	651	ILE
1	A	664	MET
1	A	699	GLN
1	A	720	ILE
1	A	753	LEU
1	A	766	LEU
1	A	839	ILE
1	A	847	TYR
1	A	857	LYS
1	A	863	LEU
1	A	867	LEU
1	A	899	LEU
1	A	919	MET
1	A	922	GLU
1	A	936	ARG
1	A	941	VAL
1	A	943	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	352	ASN
1	A	398	HIS
1	A	484	ASN
1	A	552	ASN
1	A	587	GLN
1	A	659	GLN
1	A	701	ASN
1	A	726	HIS
1	A	728	GLN
1	A	740	ASN
1	A	854	HIS
1	A	865	GLN
1	A	911	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](#)

7 ligands are modelled in this entry.

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	EOH	A	1003	-	2,2,2	0.42	0	1,1,1	0.31	0
3	EOH	A	1002	-	2,2,2	0.44	0	1,1,1	0.35	0
3	EOH	A	1005	-	2,2,2	0.45	0	1,1,1	0.32	0
2	J7G	A	1001	-	30,30,30	0.91	0	39,40,40	1.96	7 (17%)
3	EOH	A	1006	-	2,2,2	0.43	0	1,1,1	0.42	0
3	EOH	A	1007	-	2,2,2	0.40	0	1,1,1	0.33	0
3	EOH	A	1004	-	2,2,2	0.42	0	1,1,1	0.28	0

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	J7G	A	1001	-	-	0/16/17/17	0/3/3/3

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1001	J7G	C6-C7-N8	-5.57	120.74	125.59
2	A	1001	J7G	C15-N14-C13	5.44	123.86	116.51
2	A	1001	J7G	C9-N8-C7	4.47	122.02	115.58
2	A	1001	J7G	C16-C15-N14	-4.12	118.93	123.97
2	A	1001	J7G	C10-C9-N8	-3.66	119.49	123.97
2	A	1001	J7G	C12-C13-N14	-3.17	119.31	123.57
2	A	1001	J7G	C12-C13-C18	2.00	122.90	120.08

There are no chirality outliers.

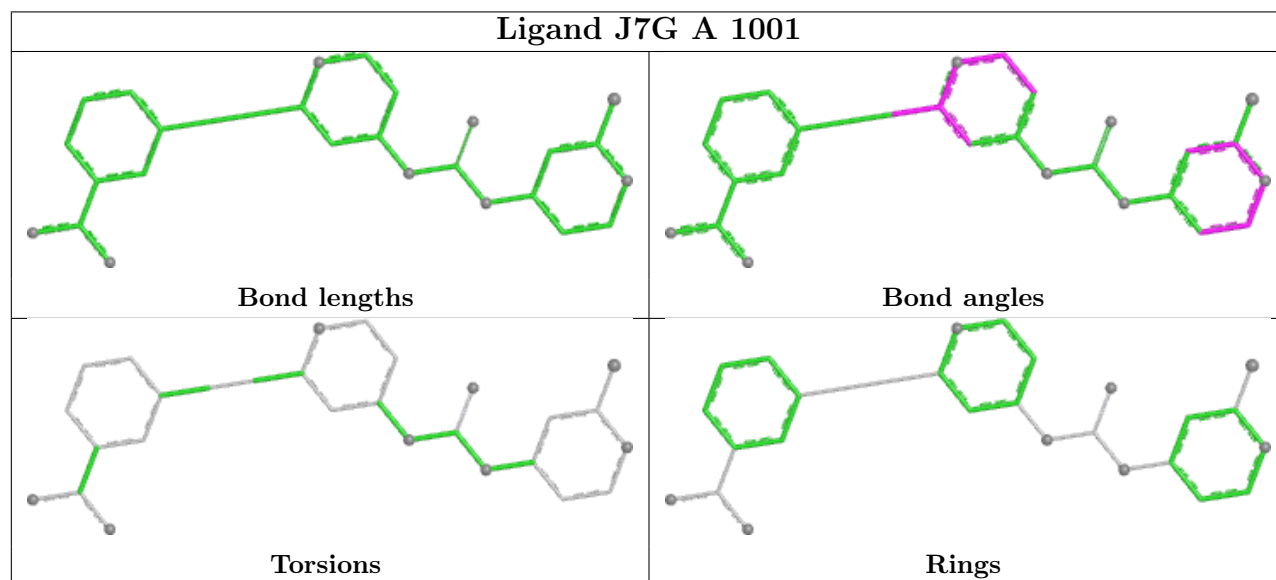
There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the

average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	614/621 (98%)	0.65	44 (7%) 21 16	35, 76, 110, 130	0

All (44) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	451	TYR	4.3
1	A	335	LYS	3.5
1	A	529	LEU	3.5
1	A	336	ILE	3.4
1	A	447	LEU	3.3
1	A	581	LEU	3.2
1	A	445	GLU	3.2
1	A	450	ILE	3.1
1	A	560	ILE	3.1
1	A	448	SER	3.1
1	A	557	ALA	3.0
1	A	718	LEU	3.0
1	A	473	LEU	3.0
1	A	483	PHE	3.0
1	A	489	TYR	3.0
1	A	547	ALA	2.9
1	A	580	VAL	2.9
1	A	467	PRO	2.9
1	A	558	MET	2.8
1	A	720	ILE	2.7
1	A	674	TRP	2.6
1	A	494	LEU	2.6
1	A	459	PRO	2.6
1	A	955	GLU	2.5
1	A	594	GLY	2.5
1	A	446	ALA	2.5
1	A	485	LEU	2.5

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	654	THR	2.4
1	A	534	LYS	2.3
1	A	592	PHE	2.3
1	A	468	ALA	2.3
1	A	442	ILE	2.3
1	A	574	VAL	2.2
1	A	562	GLY	2.2
1	A	593	VAL	2.2
1	A	480	ALA	2.1
1	A	573	ALA	2.1
1	A	482	MET	2.1
1	A	691	VAL	2.1
1	A	476	LYS	2.0
1	A	460	PRO	2.0
1	A	596	LEU	2.0
1	A	714	LYS	2.0
1	A	528	PHE	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.

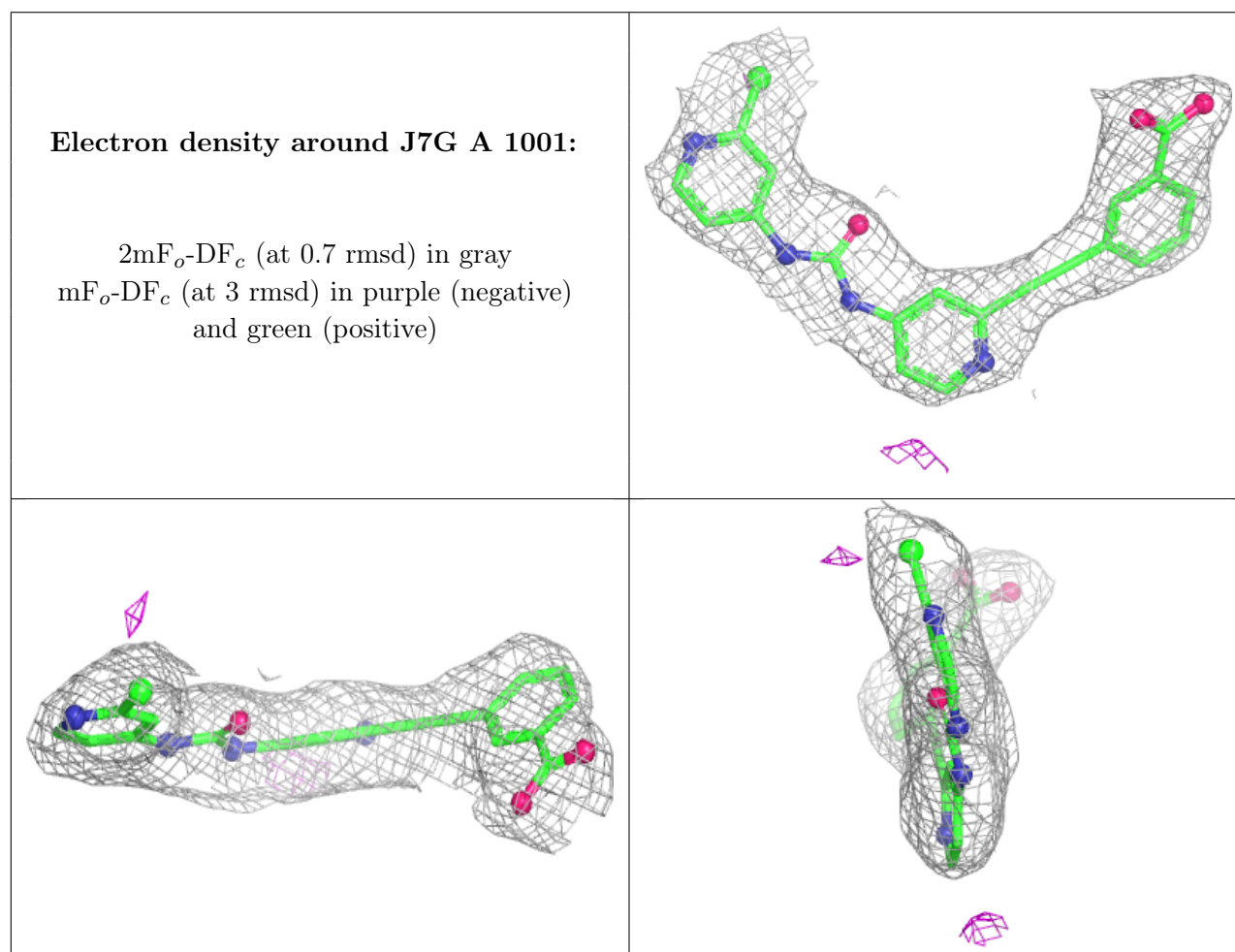
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	EOH	A	1002	3/3	0.73	0.26	66,66,68,70	0
3	EOH	A	1006	3/3	0.73	0.40	72,72,73,74	0
3	EOH	A	1007	3/3	0.80	0.32	67,67,68,69	0
3	EOH	A	1005	3/3	0.87	0.20	60,60,60,61	0
3	EOH	A	1004	3/3	0.88	0.21	55,55,56,57	0
3	EOH	A	1003	3/3	0.91	0.30	55,55,60,63	0

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	J7G	A	1001	28/28	0.96	0.08	53,64,70,72	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



6.5 Other polymers [i](#)

There are no such residues in this entry.