



Full wwPDB X-ray Structure Validation Report ⓘ

May 17, 2020 – 09:36 pm BST

PDB ID : 5RAL
Title : PanDDA analysis group deposition – Crystal Structure of JMJD1B in complex with XS039332c
Authors : Snee, M.; Nowak, R.; Johansson, C.; Burgess-Brown, N.A.; Arrowsmith, C.H.; Bountra, C.; Edwards, A.M.; Oppermann, U.
Deposited on : 2020-03-16
Resolution : 1.68 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

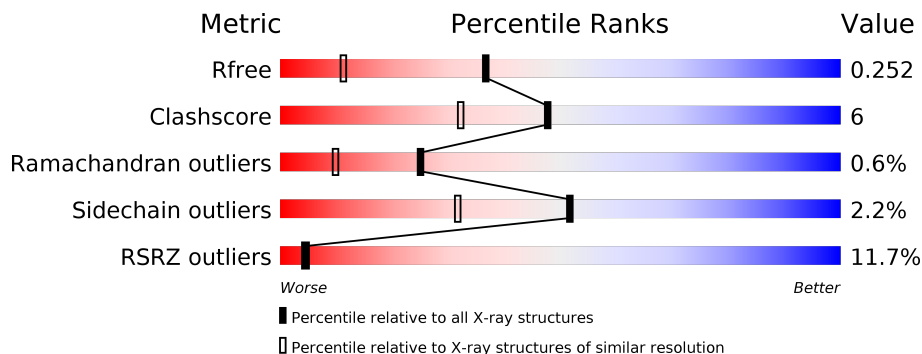
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.68 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}	130704	6780 (1.70-1.66)
Clashscore	141614	7310 (1.70-1.66)
Ramachandran outliers	138981	7173 (1.70-1.66)
Sidechain outliers	138945	7172 (1.70-1.66)
RSRZ outliers	127900	6661 (1.70-1.66)

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 on 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	372	 8% 81% 9% • 9%
1	B	372	 14% 82% 9% • 8%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	SNQ	B	1801	-	-	-	X

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 6251 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 Lysine-specific demethylase 3B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	340	2768	1752	483	516	17	0	5	0
1	B	344	2804	1777	495	517	15	0	2	0

There are 50 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1357	MET	-	initiating methionine	UNP Q7LBC6
A	1358	HIS	-	expression tag	UNP Q7LBC6
A	1359	HIS	-	expression tag	UNP Q7LBC6
A	1360	HIS	-	expression tag	UNP Q7LBC6
A	1361	HIS	-	expression tag	UNP Q7LBC6
A	1362	HIS	-	expression tag	UNP Q7LBC6
A	1363	HIS	-	expression tag	UNP Q7LBC6
A	1364	SER	-	expression tag	UNP Q7LBC6
A	1365	SER	-	expression tag	UNP Q7LBC6
A	1366	GLY	-	expression tag	UNP Q7LBC6
A	1367	VAL	-	expression tag	UNP Q7LBC6
A	1368	ASP	-	expression tag	UNP Q7LBC6
A	1369	LEU	-	expression tag	UNP Q7LBC6
A	1370	GLY	-	expression tag	UNP Q7LBC6
A	1371	THR	-	expression tag	UNP Q7LBC6
A	1372	GLU	-	expression tag	UNP Q7LBC6
A	1373	ASN	-	expression tag	UNP Q7LBC6
A	1374	LEU	-	expression tag	UNP Q7LBC6
A	1375	TYR	-	expression tag	UNP Q7LBC6
A	1376	PHE	-	expression tag	UNP Q7LBC6
A	1377	GLN	-	expression tag	UNP Q7LBC6
A	1378	SER	-	expression tag	UNP Q7LBC6
A	1379	MET	-	expression tag	UNP Q7LBC6
A	1601	GLU	GLN	conflict	UNP Q7LBC6
A	1606	HIS	GLY	conflict	UNP Q7LBC6

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Chain	Residue	Modelled	Actual	Comment	Reference
B	1357	MET	-	initiating methionine	UNP Q7LBC6
B	1358	HIS	-	expression tag	UNP Q7LBC6
B	1359	HIS	-	expression tag	UNP Q7LBC6
B	1360	HIS	-	expression tag	UNP Q7LBC6
B	1361	HIS	-	expression tag	UNP Q7LBC6
B	1362	HIS	-	expression tag	UNP Q7LBC6
B	1363	HIS	-	expression tag	UNP Q7LBC6
B	1364	SER	-	expression tag	UNP Q7LBC6
B	1365	SER	-	expression tag	UNP Q7LBC6
B	1366	GLY	-	expression tag	UNP Q7LBC6
B	1367	VAL	-	expression tag	UNP Q7LBC6
B	1368	ASP	-	expression tag	UNP Q7LBC6
B	1369	LEU	-	expression tag	UNP Q7LBC6
B	1370	GLY	-	expression tag	UNP Q7LBC6
B	1371	THR	-	expression tag	UNP Q7LBC6
B	1372	GLU	-	expression tag	UNP Q7LBC6
B	1373	ASN	-	expression tag	UNP Q7LBC6
B	1374	LEU	-	expression tag	UNP Q7LBC6
B	1375	TYR	-	expression tag	UNP Q7LBC6
B	1376	PHE	-	expression tag	UNP Q7LBC6
B	1377	GLN	-	expression tag	UNP Q7LBC6
B	1378	SER	-	expression tag	UNP Q7LBC6
B	1379	MET	-	expression tag	UNP Q7LBC6
B	1601	GLU	GLN	conflict	UNP Q7LBC6
B	1606	HIS	GLY	conflict	UNP Q7LBC6

- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

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

- Molecule 3 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total Mn 1 1	0	0
3	A	1	Total Mn 1 1	0	0

- Molecule 4 is a ligand with the chemical component id SNQ but its atom names do not match the existing wwPDB Chemical Component Dictionary definition for SNQ. ERROR THIS SHOULD NOT HAPPEN FOLLOWING ANNOTATION.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	N	0	0
			14	11	3		

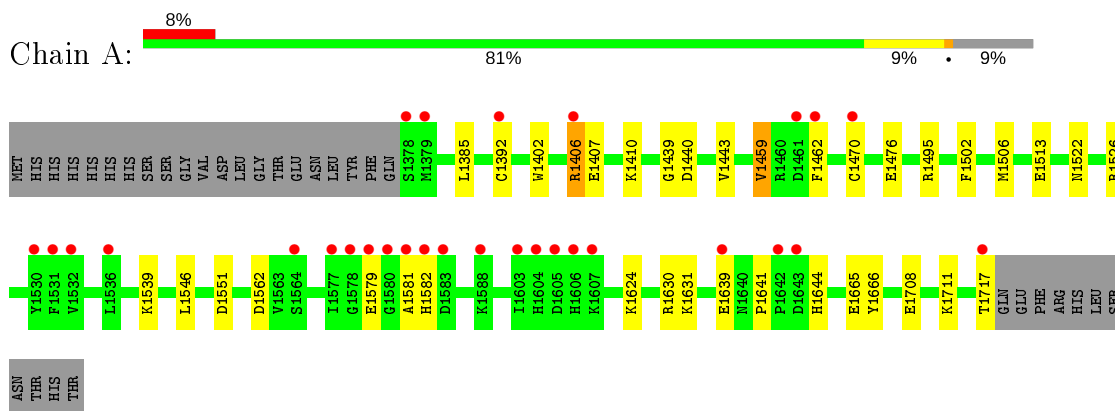
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	374	Total	O	0	0
			374	374		
5	B	284	Total	O	0	0
			284	284		

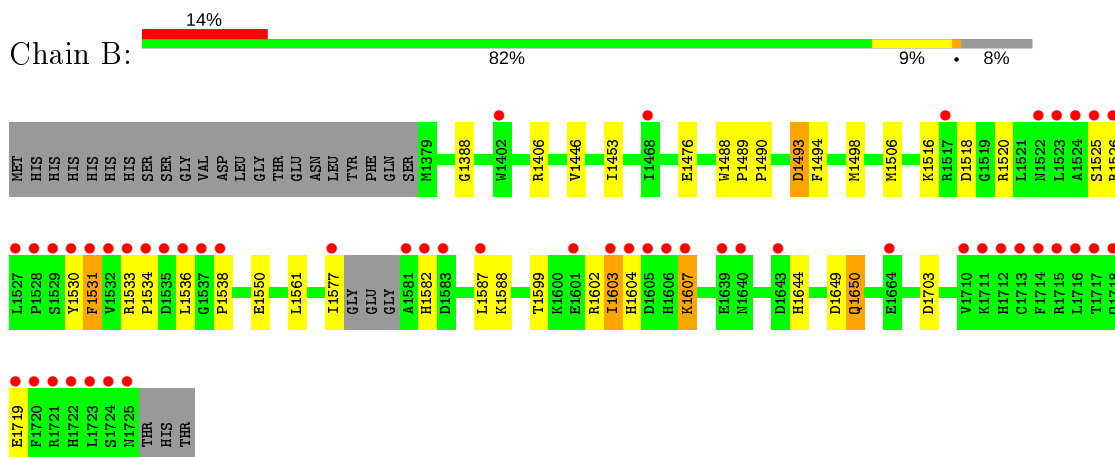
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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: Lysine-specific demethylase 3B



- Molecule 1: Lysine-specific demethylase 3B



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	57.60 Å 93.69 Å 93.43 Å 90.00° 108.03° 90.00°	Depositor
Resolution (Å)	88.85 – 1.68 88.85 – 1.68	Depositor EDS
% Data completeness (in resolution range)	98.9 (88.85-1.68) 98.9 (88.85-1.68)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.46 (at 1.68 Å)	Xtrriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.189 , 0.221 0.221 , 0.252	Depositor DCC
R_{free} test set	5376 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	28.0	Xtrriage
Anisotropy	0.056	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 49.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.024 for h,-k,-h-l	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6251	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.91% 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: CL, MN, SNQ

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	0.87	2/2841 (0.1%)	0.91	2/3854 (0.1%)
1	B	0.79	0/2878	0.81	0/3902
All	All	0.83	2/5719 (0.0%)	0.86	2/7756 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1462	PHE	C-O	7.29	1.37	1.23
1	A	1562	ASP	CG-OD2	5.37	1.37	1.25

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1562	ASP	CB-CG-OD2	6.65	124.29	118.30
1	A	1562	ASP	CB-CA-C	5.04	120.48	110.40

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	2768	0	2650	28	1
1	B	2804	0	2696	39	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	4	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	B	14	0	0	0	0
5	A	374	0	0	16	2
5	B	284	0	0	22	1
All	All	6251	0	5346	67	3

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 6.

All (67) 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:1526:ARG:NH2	5:B:1901:HOH:O	1.61	1.34
1:B:1526:ARG:N	5:B:1902:HOH:O	1.88	1.04
1:B:1520:ARG:HD2	5:B:1907:HOH:O	1.69	0.92
1:B:1533:ARG:NH1	5:B:1906:HOH:O	2.03	0.89
1:B:1476:GLU:OE2	5:B:1903:HOH:O	1.90	0.89
1:B:1520:ARG:NE	5:B:1907:HOH:O	2.09	0.85
1:B:1650:GLN:OE1	5:B:1905:HOH:O	1.93	0.85
1:B:1520:ARG:CD	5:B:1907:HOH:O	2.24	0.76
1:A:1470:CYS:C	5:A:1908:HOH:O	2.26	0.74
1:B:1506:MET:CE	1:B:1538:PRO:HG3	2.20	0.72
1:A:1406[A]:ARG:NH1	5:A:1903:HOH:O	2.23	0.71
1:A:1631:LYS:HE2	1:A:1666:TYR:OH	1.90	0.71
1:A:1624:LYS:NZ	5:A:1901:HOH:O	2.10	0.70
1:B:1644:HIS:NE2	1:B:1649:ASP:OD2	2.25	0.69
1:A:1495:ARG:NH2	5:A:1906:HOH:O	2.26	0.67
1:A:1522:ASN:O	1:A:1526:ARG:NH1	2.31	0.64
1:B:1526:ARG:CZ	5:B:1901:HOH:O	2.24	0.64
1:A:1406[A]:ARG:NH2	5:A:1909:HOH:O	2.31	0.63
1:A:1407:GLU:HG2	5:A:2225:HOH:O	1.99	0.63
1:A:1639:GLU:O	5:A:1902:HOH:O	2.17	0.59
1:B:1446:VAL:HG22	1:B:1453:ILE:HD13	1.84	0.59
1:B:1494:PHE:CE2	1:B:1498:MET:HE3	2.38	0.58
1:B:1530:TYR:O	1:B:1531:PHE:CD2	2.57	0.57
1:B:1587:LEU:HD13	1:B:1603:ILE:HD11	1.86	0.57
1:A:1443:VAL:HG11	1:A:1459:VAL:HG23	1.86	0.57
1:A:1711:LYS:CB	5:A:2109:HOH:O	2.54	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1506:MET:HE1	1:B:1538:PRO:HG3	1.86	0.55
1:A:1407:GLU:O	1:A:1410:LYS:HG2	2.07	0.55
1:B:1506:MET:HE2	1:B:1538:PRO:HG3	1.88	0.54
1:A:1407:GLU:CG	5:A:2225:HOH:O	2.56	0.54
1:B:1550:GLU:HG3	5:B:2079:HOH:O	2.08	0.53
1:A:1502:PHE:CE2	1:A:1506:MET:HE2	2.43	0.53
1:B:1506:MET:O	1:B:1516:LYS:HE3	2.08	0.53
1:B:1388:GLY:N	5:B:1909:HOH:O	2.29	0.53
1:B:1526:ARG:HG3	5:B:1902:HOH:O	2.08	0.52
1:A:1631:LYS:NZ	1:A:1665:GLU:OE1	2.38	0.52
1:A:1385:LEU:HD12	1:A:1392[B]:CYS:SG	2.50	0.52
1:B:1518:ASP:HB3	5:B:2109:HOH:O	2.10	0.52
1:B:1536:LEU:CD1	5:B:1992:HOH:O	2.57	0.52
1:B:1536:LEU:HD12	5:B:2003:HOH:O	2.09	0.51
1:A:1539:LYS:HD3	5:A:2105:HOH:O	2.12	0.50
1:A:1439:GLY:HA2	1:A:1459:VAL:HG12	1.94	0.49
1:B:1494:PHE:CZ	1:B:1498:MET:HE3	2.49	0.48
1:A:1476:GLU:CB	5:A:2126:HOH:O	2.61	0.48
1:B:1490:PRO:O	1:B:1493:ASP:HB2	2.14	0.48
1:B:1561:LEU:C	1:B:1561:LEU:HD12	2.35	0.47
1:B:1536:LEU:HD22	1:B:1703:ASP:HB2	1.98	0.46
1:A:1513:GLU:HG2	5:A:2077:HOH:O	2.16	0.45
1:A:1624:LYS:HE2	5:A:2213:HOH:O	2.17	0.45
1:B:1526:ARG:NE	5:B:1901:HOH:O	2.46	0.45
1:A:1443:VAL:CG1	1:A:1459:VAL:HG23	2.46	0.45
1:B:1536:LEU:HD13	5:B:1992:HOH:O	2.17	0.44
1:B:1406:ARG:CB	5:B:2137:HOH:O	2.66	0.44
1:A:1440:ASP:HB2	5:A:2118:HOH:O	2.17	0.44
1:B:1526:ARG:NH2	5:B:1922:HOH:O	2.51	0.44
1:A:1582:HIS:CE1	5:A:2034:HOH:O	2.70	0.44
1:A:1546:LEU:HB3	1:A:1551:ASP:HB3	2.01	0.43
1:B:1599:THR:O	1:B:1602:ARG:HB3	2.18	0.42
1:B:1607:LYS:NZ	1:B:1607:LYS:HA	2.33	0.42
1:A:1402:TRP:O	1:A:1406[A]:ARG:HG2	2.19	0.42
1:B:1530:TYR:O	1:B:1531:PHE:CB	2.67	0.42
1:A:1582:HIS:O	1:A:1582:HIS:CG	2.73	0.42
1:B:1525:SER:N	5:B:1904:HOH:O	2.52	0.42
1:B:1488:TRP:HA	1:B:1489:PRO:C	2.41	0.41
1:B:1719:GLU:HG2	5:B:1933:HOH:O	2.20	0.41
1:B:1536:LEU:HD12	5:B:1992:HOH:O	2.17	0.41
1:A:1630:ARG:NH2	5:A:1914:HOH:O	2.42	0.40

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1708:GLU:OE2	1:B:1526:ARG:NH1[2_557]	2.00	0.20
5:A:2211:HOH:O	5:B:1907:HOH:O[2_557]	2.12	0.08
5:A:2222:HOH:O	5:A:2244:HOH:O[2_648]	2.13	0.07

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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	343/372 (92%)	332 (97%)	10 (3%)	1 (0%)	41	23
1	B	342/372 (92%)	324 (95%)	15 (4%)	3 (1%)	17	4
All	All	685/744 (92%)	656 (96%)	25 (4%)	4 (1%)	25	10

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1581	ALA
1	B	1531	PHE
1	B	1534	PRO
1	B	1603	ILE

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	299/331 (90%)	292 (98%)	7 (2%)	50	30
1	B	303/331 (92%)	296 (98%)	7 (2%)	50	30
All	All	602/662 (91%)	588 (98%)	14 (2%)	52	30

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1406[A]	ARG
1	A	1406[B]	ARG
1	A	1459	VAL
1	A	1579	GLU
1	A	1641	PRO
1	A	1644	HIS
1	A	1717	THR
1	B	1493	ASP
1	B	1577	ILE
1	B	1582	HIS
1	B	1588	LYS
1	B	1604	HIS
1	B	1607	LYS
1	B	1650	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1640	ASN
1	A	1650	GLN
1	B	1656	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 carbohydrates in this entry.

5.6 Ligand geometry

Of 8 ligands modelled in this entry, 1 could not be matched to an existing wwPDB Chemical Component Dictionary definition at this stage and 7 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

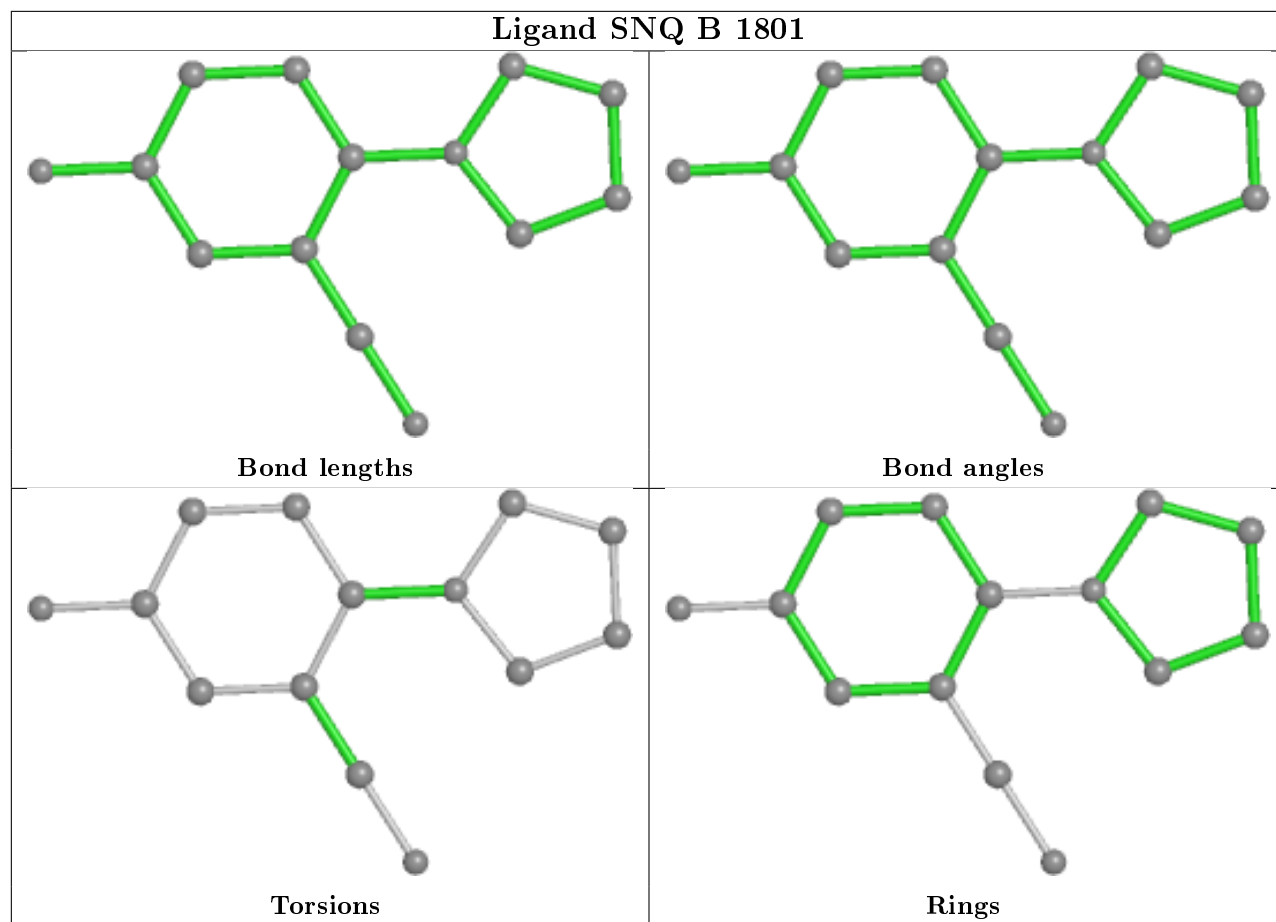
There are no chirality outliers.

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	340/372 (91%)	0.17	29 (8%) 10 11	19, 31, 68, 126	10 (2%)
1	B	344/372 (92%)	1.00	51 (14%) 2 2	21, 36, 70, 141	35 (10%)
All	All	684/744 (91%)	0.58	80 (11%) 4 4	19, 33, 68, 141	45 (6%)

All (80) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	1713	CYS	15.8
1	B	1581	ALA	15.7
1	B	1714	PHE	14.7
1	B	1527	LEU	14.6
1	B	1536	LEU	14.2
1	B	1716	LEU	14.0
1	B	1723	LEU	13.3
1	B	1523	LEU	13.3
1	B	1720	PHE	12.7
1	B	1710	VAL	12.7
1	B	1605	ASP	12.7
1	B	1531	PHE	11.7
1	B	1532	VAL	11.0
1	B	1524	ALA	11.0
1	B	1525	SER	10.8
1	B	1717	THR	10.6
1	A	1581	ALA	9.8
1	B	1537	GLY	9.7
1	B	1533	ARG	9.6
1	A	1577	ILE	9.5
1	B	1538	PRO	9.3
1	A	1462	PHE	9.3
1	B	1530	TYR	9.0
1	B	1712	HIS	9.0

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Mol	Chain	Res	Type	RSRZ
1	A	1531	PHE	8.8
1	A	1392[A]	CYS	8.7
1	B	1522	ASN	8.5
1	B	1711	LYS	8.4
1	B	1604	HIS	8.0
1	A	1530	TYR	8.0
1	B	1517[A]	ARG	7.9
1	B	1535	ASP	7.9
1	B	1526	ARG	7.7
1	A	1578	GLY	7.6
1	B	1725	ASN	7.6
1	B	1534	PRO	7.2
1	B	1715	ARG	7.2
1	A	1379[A]	MET	7.1
1	B	1606	HIS	6.8
1	B	1718	GLN	6.6
1	B	1577	ILE	6.4
1	A	1642	PRO	6.0
1	A	1532	VAL	5.9
1	B	1721	ARG	5.9
1	A	1582	HIS	5.7
1	B	1719	GLU	5.3
1	B	1724	SER	5.2
1	B	1722	HIS	5.1
1	B	1528	PRO	5.0
1	A	1406[A]	ARG	5.0
1	A	1588[A]	LYS	4.8
1	A	1564[A]	SER	4.7
1	A	1580	GLY	4.6
1	B	1529	SER	4.4
1	B	1603	ILE	4.3
1	B	1582	HIS	4.0
1	A	1605	ASP	4.0
1	A	1461	ASP	3.7
1	B	1664[A]	GLU	3.5
1	A	1579	GLU	3.4
1	B	1601	GLU	3.4
1	B	1587	LEU	3.3
1	A	1470	CYS	3.1
1	A	1603	ILE	2.9
1	A	1606	HIS	2.8
1	A	1583	ASP	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	1717	THR	2.7
1	B	1640	ASN	2.6
1	A	1639	GLU	2.5
1	B	1607	LYS	2.5
1	B	1643	ASP	2.4
1	B	1583	ASP	2.3
1	B	1402	TRP	2.3
1	A	1604	HIS	2.2
1	A	1607	LYS	2.1
1	B	1468	ILE	2.1
1	A	1378	SER	2.1
1	A	1536	LEU	2.1
1	A	1643	ASP	2.1
1	B	1639	GLU	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](#)

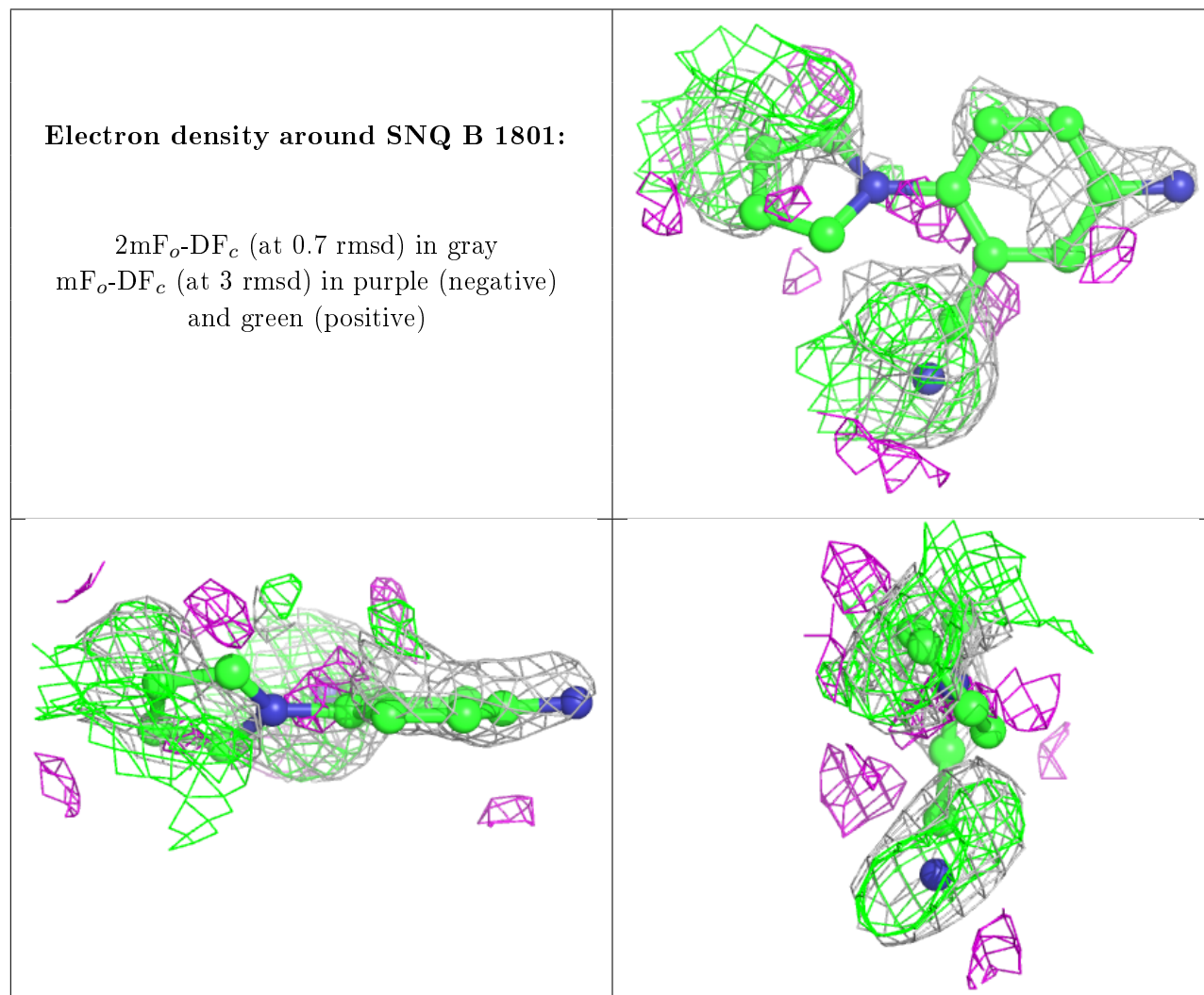
There are no carbohydrates in this entry.

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	SNQ	B	1801	14/14	0.28	0.53	25,27,30,31	14
2	CL	A	1804	1/1	0.97	0.05	57,57,57,57	0
2	CL	A	1803	1/1	0.98	0.06	37,37,37,37	0
3	MN	A	1805	1/1	0.99	0.11	22,22,22,22	0
3	MN	B	1803	1/1	0.99	0.13	24,24,24,24	0
2	CL	A	1802	1/1	0.99	0.10	39,39,39,39	0
2	CL	B	1802	1/1	0.99	0.03	38,38,38,38	0
2	CL	A	1801	1/1	1.00	0.10	26,26,26,26	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.