



# Full wwPDB X-ray Structure Validation Report ⓘ

May 18, 2020 – 11:25 am BST

PDB ID : 5RAM  
Title : PanDDA analysis group deposition – Crystal Structure of JMJD1B in complex with XS038544d  
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Deposited on : 2020-03-16  
Resolution : 2.05 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](mailto: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.

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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

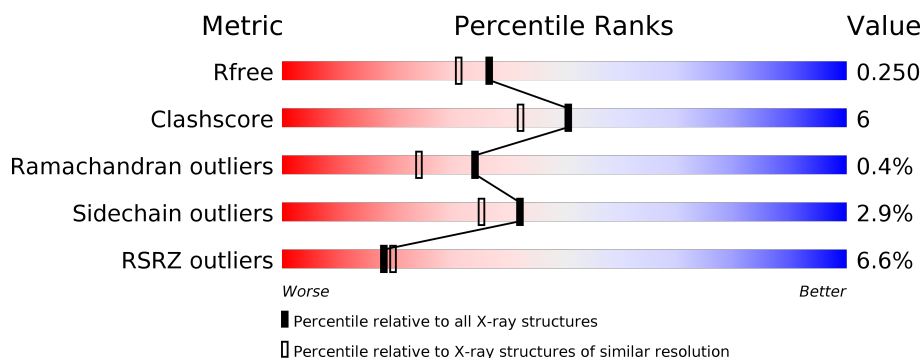
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.05 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	1692 (2.04-2.04)
Clashscore	141614	1773 (2.04-2.04)
Ramachandran outliers	138981	1752 (2.04-2.04)
Sidechain outliers	138945	1752 (2.04-2.04)
RSRZ outliers	127900	1672 (2.04-2.04)

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  $\geq 3$ , 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	<div> <div>9%</div> <div> <div></div> <div>81%</div> <div>10%</div> <div>•</div> <div>9%</div> </div> </div>
1	B	372	<div> <div>3%</div> <div> <div></div> <div>80%</div> <div>11%</div> <div>•</div> <div>8%</div> </div> </div>

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:

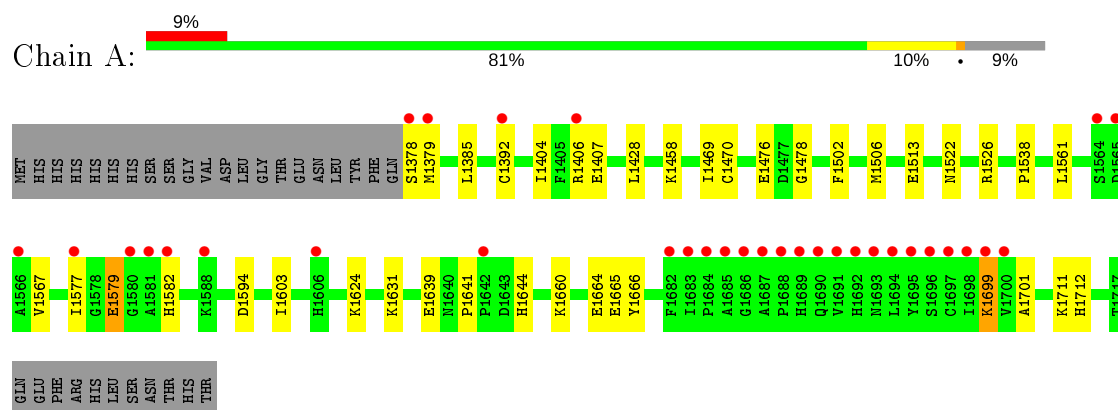
Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	CL	A	1804	-	-	-	X

ENTRY-COMPOSITION INFOmissingINFO

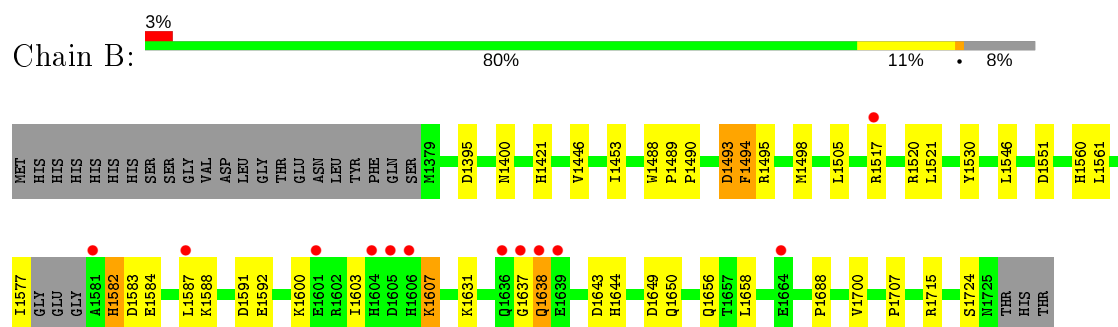
## 2 Residue-property plots [i](#)

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



### 3 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	57.03 Å 93.50 Å 92.92 Å 90.00° 108.05° 90.00°	Depositor
Resolution (Å)	88.35 – 2.05 88.35 – 2.05	Depositor EDS
% Data completeness (in resolution range)	88.6 (88.35-2.05) 88.6 (88.35-2.05)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.68 (at 2.05 Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, $R_{free}$	0.175 , 0.235 0.197 , 0.250	Depositor DCC
$R_{free}$ test set	2634 reflections (5.11%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	31.1	Xtriage
Anisotropy	0.101	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 58.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.035 for h,-k,-h-l	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	6293	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.49% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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.

## 4 Model quality [i](#)

### 4.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MN, S3V, CL

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	1/2841 (0.0%)	0.91	0/3854
1	B	0.79	0/2872	0.88	0/3895
All	All	0.80	1/5713 (0.0%)	0.89	0/7749

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1478	GLY	C-O	5.16	1.31	1.23

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 4.2 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	29	0
1	B	2798	0	2685	33	0
2	A	9	0	0	1	0
2	B	9	0	0	0	0
3	A	4	0	0	1	0
3	B	1	0	0	1	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	384	0	0	13	2
5	B	318	0	0	9	1
All	All	6293	0	5335	62	2

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 (62) 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:1476:GLU:CB	5:A:2225:HOH:O	2.15	0.93
1:A:1711:LYS:CB	5:A:2239:HOH:O	2.16	0.93
1:A:1522:ASN:O	1:A:1526:ARG:NH1	2.02	0.92
1:A:1407:GLU:HG2	5:A:2221:HOH:O	1.73	0.87
1:A:1406[A]:ARG:NH1	5:A:1903:HOH:O	2.10	0.83
1:A:1639:GLU:O	5:A:1902:HOH:O	1.96	0.82
1:B:1421:HIS:HB2	5:B:2168:HOH:O	1.81	0.80
1:A:1513:GLU:HG2	5:A:2175:HOH:O	1.81	0.80
1:A:1406[A]:ARG:NH2	5:A:1906:HOH:O	2.15	0.78
1:B:1517[B]:ARG:HH11	1:B:1517[B]:ARG:HG3	1.51	0.75
1:A:1385:LEU:HD12	1:A:1392[B]:CYS:SG	2.27	0.74
1:B:1517[B]:ARG:HH11	1:B:1517[B]:ARG:CG	2.01	0.73
1:B:1644:HIS:NE2	1:B:1649:ASP:OD2	2.23	0.69
1:B:1395:ASP:OD1	5:B:1901:HOH:O	2.10	0.69
1:A:1428:LEU:O	5:A:1904:HOH:O	2.11	0.68
1:B:1517[B]:ARG:NH1	1:B:1517[B]:ARG:HG3	2.06	0.68
1:B:1650:GLN:OE1	5:B:1902:HOH:O	2.10	0.68
1:A:1470:CYS:O	5:A:1905:HOH:O	2.12	0.67
1:A:1385:LEU:CD1	1:A:1392[B]:CYS:SG	2.86	0.64
1:B:1520:ARG:NE	5:B:1909:HOH:O	2.32	0.61
1:A:1631:LYS:HE2	1:A:1666:TYR:OH	2.04	0.57
1:B:1715:ARG:HD3	5:B:2184:HOH:O	2.05	0.57
1:B:1490:PRO:O	1:B:1493:ASP:HB2	2.05	0.56
1:A:1631:LYS:NZ	1:A:1665:GLU:OE1	2.36	0.53
1:A:1469:ILE:HD12	1:A:1579:GLU:HB2	1.91	0.53
1:A:1699:LYS:NZ	2:A:1801:S3V:O	2.41	0.52
1:B:1587:LEU:HD13	1:B:1603:ILE:HD11	1.91	0.52
1:B:1488:TRP:HA	1:B:1489:PRO:C	2.29	0.52
1:B:1591:ASP:O	5:B:1903:HOH:O	2.17	0.52
1:B:1607:LYS:HA	1:B:1607:LYS:NZ	2.26	0.51
1:B:1505:LEU:HD22	1:B:1700:VAL:HG11	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1446:VAL:HG22	1:B:1453:ILE:HD13	1.92	0.51
1:B:1631:LYS:HG2	5:B:2100:HOH:O	2.12	0.50
1:B:1560:HIS:O	1:B:1688:PRO:HA	2.12	0.49
1:B:1494:PHE:CE1	1:B:1498:MET:HE3	2.47	0.49
1:A:1506:MET:HE1	1:A:1538:PRO:HD3	1.93	0.49
1:A:1385:LEU:HG	1:A:1392[B]:CYS:SG	2.53	0.48
1:A:1624:LYS:NZ	5:A:1901:HOH:O	1.95	0.48
1:B:1498:MET:HA	3:B:1802:CL:CL	2.51	0.47
1:B:1521:LEU:O	1:B:1707:PRO:HD3	2.15	0.47
1:B:1715:ARG:HG3	5:B:2130:HOH:O	2.15	0.47
1:A:1458:LYS:HG2	5:A:2197:HOH:O	2.15	0.46
1:A:1378:SER:HA	5:A:1959:HOH:O	2.14	0.46
1:A:1582:HIS:O	1:A:1582:HIS:CG	2.69	0.46
1:B:1637:GLY:O	1:B:1638:GLN:C	2.54	0.45
1:B:1546:LEU:HB3	1:B:1551:ASP:HB3	1.98	0.45
1:A:1660:LYS:HE2	1:A:1664:GLU:OE2	2.17	0.45
1:B:1582:HIS:C	1:B:1584:GLU:H	2.18	0.45
1:A:1641:PRO:CD	5:A:2249:HOH:O	2.65	0.44
1:B:1607:LYS:HA	1:B:1607:LYS:CE	2.48	0.44
1:A:1502:PHE:CE2	1:A:1506:MET:HE2	2.53	0.43
1:B:1591:ASP:HA	1:B:1600:LYS:HE2	1.99	0.43
1:B:1517[B]:ARG:CB	1:B:1517[B]:ARG:HH11	2.30	0.43
1:B:1561:LEU:C	1:B:1561:LEU:HD12	2.38	0.43
1:B:1658:LEU:HD23	1:B:1658:LEU:HA	1.89	0.42
1:A:1385:LEU:CG	1:A:1392[B]:CYS:SG	3.08	0.41
1:B:1517[B]:ARG:HB2	1:B:1517[B]:ARG:HH11	1.86	0.41
1:B:1592:GLU:HG3	1:B:1656:GLN:CD	2.40	0.41
1:A:1567:VAL:O	1:A:1701:ALA:HA	2.20	0.41
1:A:1561:LEU:HD12	1:A:1561:LEU:C	2.42	0.41
1:A:1712:HIS:NE2	3:A:1805:CL:CL	2.73	0.40
1:B:1400:ASN:OD1	5:B:1904:HOH:O	2.22	0.40

All (2) 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)
5:A:2063:HOH:O	5:B:1909:HOH:O[2_557]	2.08	0.12
5:A:2197:HOH:O	5:A:2234:HOH:O[2_648]	2.17	0.03



## 4.3 Torsion angles

### 4.3.1 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	343/372 (92%)	334 (97%)	8 (2%)	1 (0%)	41	31
1	B	342/372 (92%)	325 (95%)	15 (4%)	2 (1%)	25	15
All	All	685/744 (92%)	659 (96%)	23 (3%)	3 (0%)	34	24

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	1583	ASP
1	B	1638	GLN
1	A	1594	ASP

### 4.3.2 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	299/331 (90%)	291 (97%)	8 (3%)	44	38
1	B	302/331 (91%)	292 (97%)	10 (3%)	38	31
All	All	601/662 (91%)	583 (97%)	18 (3%)	42	34

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1379[A]	MET
1	A	1379[B]	MET
1	A	1404	ILE

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Mol	Chain	Res	Type
1	A	1577	ILE
1	A	1579	GLU
1	A	1603	ILE
1	A	1644	HIS
1	A	1699	LYS
1	B	1493	ASP
1	B	1494	PHE
1	B	1495	ARG
1	B	1530	TYR
1	B	1577	ILE
1	B	1582	HIS
1	B	1588	LYS
1	B	1607	LYS
1	B	1643	ASP
1	B	1724	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

#### 4.3.3 RNA ⓘ

There are no RNA molecules in this entry.

#### 4.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

#### 4.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

#### 4.6 Ligand geometry ⓘ

Of 9 ligands modelled in this entry, 2 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.

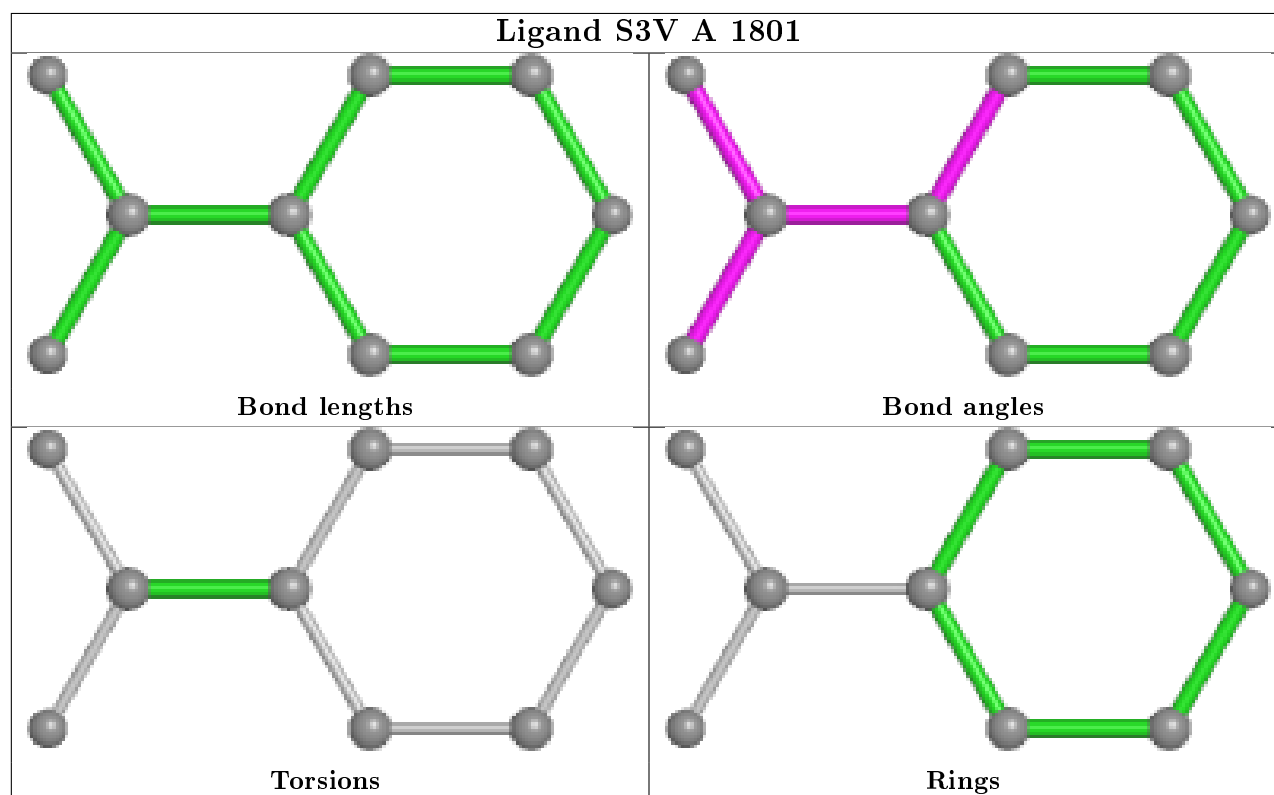
There are no chirality outliers.

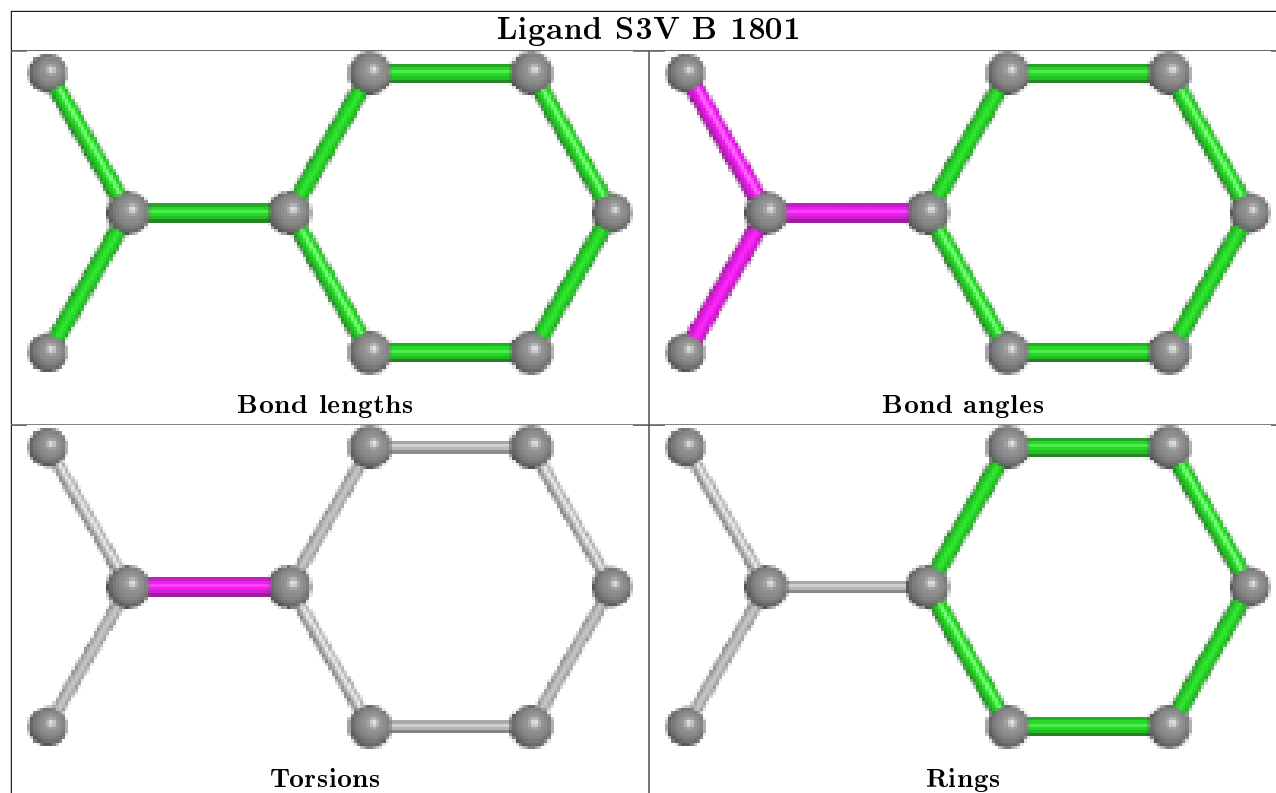
There are no torsion outliers.

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.





#### 4.7 Other polymers [i](#)

There are no such residues in this entry.

#### 4.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 5 Fit of model and data ⓘ

### 5.1 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	340/372 (91%)	-0.04	33 (9%) 7 8	19, 32, 71, 116	31 (9%)
1	B	344/372 (92%)	-0.27	12 (3%) 44 48	22, 38, 80, 142	6 (1%)
All	All	684/744 (91%)	-0.16	45 (6%) 18 19	19, 35, 79, 142	37 (5%)

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	1605	ASP	12.0
1	B	1581	ALA	7.9
1	A	1392[A]	CYS	7.3
1	A	1379[A]	MET	6.8
1	A	1697	CYS	6.8
1	A	1581	ALA	6.7
1	B	1604	HIS	6.2
1	A	1694	LEU	6.1
1	B	1637	GLY	6.1
1	A	1566	ALA	6.0
1	A	1698	ILE	5.9
1	A	1691	VAL	5.8
1	A	1580	GLY	5.6
1	A	1687	ALA	5.5
1	A	1683	ILE	5.4
1	A	1695	TYR	5.4
1	A	1682	PHE	5.3
1	A	1700	VAL	5.2
1	A	1685	ALA	5.1
1	A	1688	PRO	5.1
1	A	1582	HIS	5.0
1	A	1696	SER	4.5
1	A	1689	HIS	4.4
1	A	1684	PRO	4.3

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Mol	Chain	Res	Type	RSRZ
1	A	1588[A]	LYS	4.3
1	B	1587	LEU	4.3
1	B	1517[A]	ARG	4.3
1	A	1692	HIS	4.3
1	A	1686	GLY	4.2
1	B	1606	HIS	4.2
1	A	1693	ASN	4.2
1	A	1564[A]	SER	4.2
1	B	1638	GLN	4.2
1	A	1406[A]	ARG	4.0
1	B	1636	GLN	3.7
1	A	1699	LYS	3.5
1	A	1565	ASP	3.5
1	A	1577	ILE	3.4
1	B	1601	GLU	3.2
1	B	1664[A]	GLU	3.1
1	A	1690	GLN	2.9
1	A	1378	SER	2.9
1	B	1639	GLU	2.8
1	A	1642	PRO	2.7
1	A	1606	HIS	2.7

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

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

## 5.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.4 Ligands [i](#)

LIGAND-RSR INFOmissingINFO

## 5.5 Other polymers [i](#)

There are no such residues in this entry.