



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 06:37 PM UTC

PDB ID : 5RAE / pdb_00005rae
Title : PanDDA analysis group deposition – Crystal Structure of JMJD1B in complex with FM001558a
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 : 1.88 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

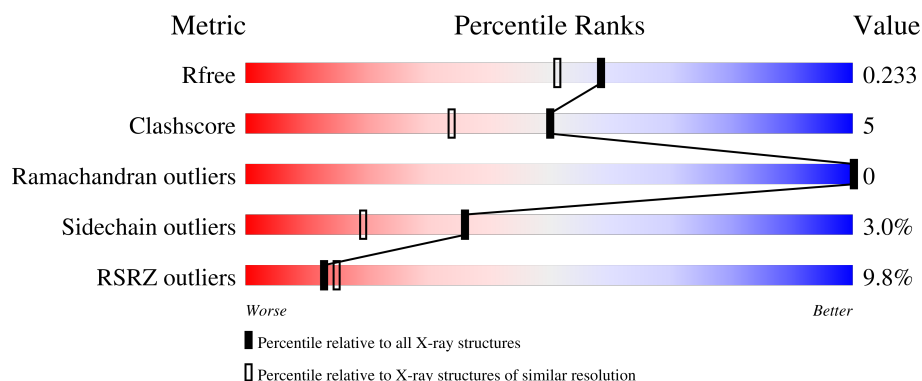
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.88 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	1220 (1.88-1.88)
Clashscore	190562	1234 (1.88-1.88)
Ramachandran outliers	187476	1222 (1.88-1.88)
Sidechain outliers	187428	1222 (1.88-1.88)
RSRZ outliers	180081	1220 (1.88-1.88)

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	372	11% 81% 9% .. 9%
1	B	372	7% 81% 12% 8%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6284 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
1	A	340	Total	C	N	O	S	0	5	0
			2768	1752	483	516	17			
1	B	344	Total	C	N	O	S	0	2	0
			2798	1774	492	517	15			

There are 50 discrepancies between the modelled and reference sequences:

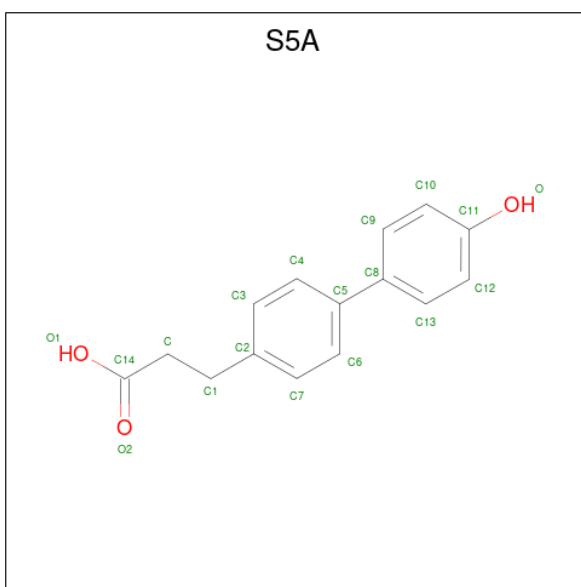
Chain	Residue	Modelled	Actual	Comment	Reference
A	1357	MET	-	initiating methionine	UNP Q7LBC6
A	1358	HIS	-	expression tag	UNP Q7LBC6
A	1359	HIS	-	expression tag	UNP Q7LBC6
A	1360	HIS	-	expression tag	UNP Q7LBC6
A	1361	HIS	-	expression tag	UNP Q7LBC6
A	1362	HIS	-	expression tag	UNP Q7LBC6
A	1363	HIS	-	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 3-[4-(4-hydroxyphenyl)phenyl]propanoic acid (CCD ID: S5A) (formula: C₁₅H₁₄O₃) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			18	15	3		

- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

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

- Molecule 4 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

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

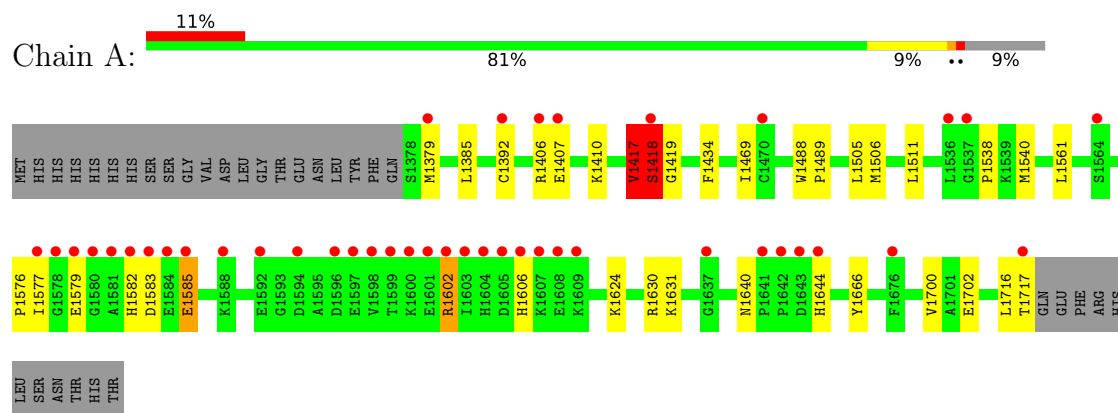
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	375	Total	O	0	0
			375	375		
5	B	318	Total	O	0	0
			318	318		

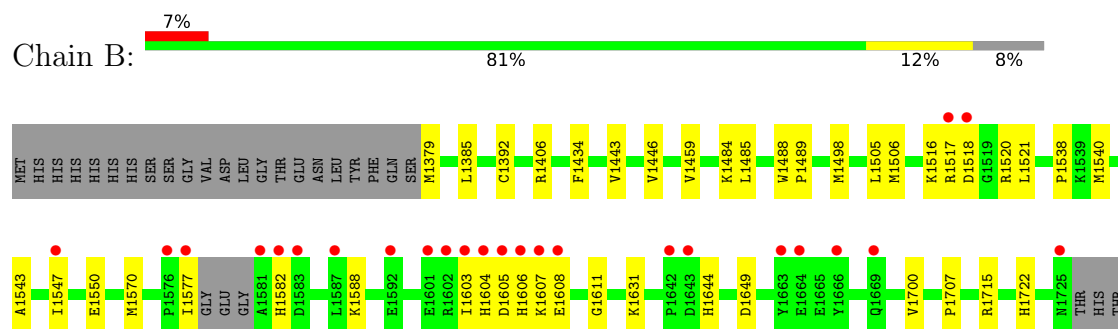
3 Residue-property plots [i](#)

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: 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	57.27Å 93.43Å 92.86Å 90.00° 107.55° 90.00°	Depositor
Resolution (Å)	54.43 – 1.88 54.43 – 1.88	Depositor EDS
% Data completeness (in resolution range)	98.4 (54.43-1.88) 98.4 (54.43-1.88)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.44 (at 1.88Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.180 , 0.222 0.194 , 0.233	Depositor DCC
R_{free} test set	3779 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	28.5	Xtriage
Anisotropy	0.102	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 48.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.032 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6284	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.94% 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: MN, CL, S5A

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	1.36	8/2841 (0.3%)	1.37	13/3854 (0.3%)
1	B	1.10	1/2872 (0.0%)	1.26	0/3895
All	All	1.24	9/5713 (0.2%)	1.32	13/7749 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1418	SER	CB-OG	25.79	1.93	1.42
1	A	1418	SER	N-CA	17.28	1.68	1.46
1	A	1418	SER	CA-CB	13.11	1.75	1.53
1	A	1417	VAL	C-N	10.13	1.47	1.33
1	A	1602	ARG	CZ-NH2	-9.59	1.21	1.33
1	A	1417	VAL	C-O	7.21	1.32	1.24
1	A	1602	ARG	NE-CZ	-6.49	1.25	1.33
1	A	1511	LEU	N-CA	5.84	1.50	1.46
1	B	1547	ILE	N-CA	5.66	1.51	1.46

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1602	ARG	NE-CZ-NH2	-13.34	107.20	119.20
1	A	1602	ARG	NE-CZ-NH1	12.67	134.17	121.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1417	VAL	CA-C-N	-10.66	101.17	121.54
1	A	1417	VAL	C-N-CA	-10.66	101.17	121.54
1	A	1418	SER	N-CA-CB	-10.11	93.40	110.49
1	A	1417	VAL	CA-C-O	-7.33	111.16	120.69
1	A	1602	ARG	CD-NE-CZ	-6.64	115.11	124.40
1	A	1576	PRO	CA-C-N	5.72	130.00	122.51
1	A	1576	PRO	C-N-CA	5.72	130.00	122.51
1	A	1469	ILE	CA-C-O	-5.15	115.39	120.85
1	A	1418	SER	N-CA-C	5.12	121.69	110.80
1	A	1716	LEU	CA-C-N	5.03	130.75	121.70
1	A	1716	LEU	C-N-CA	5.03	130.75	121.70

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1417	VAL	Mainchain
1	A	1418	SER	Mainchain
1	A	1602	ARG	Sidechain

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	2768	0	2650	29	0
1	B	2798	0	2685	29	0
2	A	18	0	0	0	0
3	A	4	0	0	0	0
3	B	1	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	375	0	0	7	2
5	B	318	0	0	10	2
All	All	6284	0	5335	58	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 5.

All (58) 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:1418:SER:CA	1:A:1418:SER:CB	1.75	1.63
1:A:1418:SER:CA	1:A:1418:SER:N	1.68	1.56
1:A:1418:SER:CB	1:A:1418:SER:OG	1.93	1.16
1:A:1417:VAL:C	1:A:1418:SER:CA	2.44	0.89
1:A:1624:LYS:NZ	5:A:1902:HOH:O	2.15	0.79
1:A:1640:ASN:ND2	5:A:1901:HOH:O	2.09	0.79
1:A:1406[A]:ARG:NH2	5:A:1904:HOH:O	2.20	0.75
1:A:1418:SER:CB	1:A:1418:SER:N	2.50	0.74
1:A:1406[A]:ARG:NH1	5:A:1903:HOH:O	2.19	0.74
1:B:1518:ASP:HB3	5:B:2127:HOH:O	1.88	0.73
1:B:1517[B]:ARG:HG3	1:B:1517[B]:ARG:HH11	1.54	0.73
1:B:1517[B]:ARG:HH11	1:B:1517[B]:ARG:CG	2.03	0.72
1:A:1418:SER:CB	1:A:1418:SER:C	2.61	0.71
1:A:1385:LEU:HD12	1:A:1392[B]:CYS:SG	2.33	0.69
1:B:1517[B]:ARG:HG3	1:B:1517[B]:ARG:NH1	2.06	0.68
1:B:1489:PRO:HG2	1:B:1498:MET:HE2	1.73	0.68
1:B:1406:ARG:CB	5:B:2173:HOH:O	2.44	0.65
1:B:1506:MET:CE	1:B:1538:PRO:HG3	2.28	0.63
1:B:1505:LEU:HD22	1:B:1700:VAL:HG11	1.80	0.63
1:A:1418:SER:CB	1:A:1419:GLY:N	2.66	0.58
1:A:1506:MET:HE2	1:A:1538:PRO:HD3	1.87	0.56
1:A:1434:PHE:CZ	1:A:1540:MET:HE1	2.41	0.55
1:B:1520:ARG:NE	5:B:1909:HOH:O	2.41	0.54
1:A:1418:SER:N	1:A:1418:SER:OG	2.41	0.54
1:B:1506:MET:HE3	1:B:1538:PRO:HD3	1.91	0.53
1:A:1631:LYS:HE2	1:A:1666:TYR:OH	2.10	0.51
1:B:1631:LYS:HG2	5:B:2122:HOH:O	2.11	0.50
1:B:1722:HIS:ND1	5:B:1906:HOH:O	2.34	0.49
1:B:1570:MET:HE3	1:B:1611:GLY:HA3	1.94	0.49
1:A:1561:LEU:C	1:A:1561:LEU:HD12	2.37	0.49
1:B:1484:LYS:HB2	1:B:1543:ALA:HB2	1.95	0.48
1:A:1630:ARG:NH2	5:A:1913:HOH:O	2.43	0.48
1:B:1715:ARG:HG3	5:B:2134:HOH:O	2.13	0.47
1:B:1484:LYS:NZ	5:B:1917:HOH:O	2.48	0.47
1:B:1516:LYS:HE3	5:B:1920:HOH:O	2.14	0.47
1:B:1644:HIS:CE1	1:B:1649:ASP:OD2	2.67	0.47
1:B:1385:LEU:HD12	1:B:1392:CYS:SG	2.55	0.47
1:B:1506:MET:HE2	1:B:1538:PRO:HG3	1.96	0.47
1:B:1550:GLU:HG3	5:B:2057:HOH:O	2.15	0.46
1:A:1417:VAL:O	1:A:1418:SER:CA	2.63	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1418:SER:HB3	5:A:2022:HOH:O	2.14	0.46
1:B:1488:TRP:HA	1:B:1489:PRO:C	2.41	0.46
1:B:1506:MET:CE	1:B:1538:PRO:CG	2.94	0.45
1:A:1582:HIS:NE2	1:A:1585:GLU:HG3	2.32	0.45
1:A:1506:MET:HE1	1:A:1702:GLU:CD	2.41	0.45
1:B:1443:VAL:HG21	1:B:1485:LEU:HD11	1.99	0.44
1:A:1505:LEU:HD22	1:A:1700:VAL:HG11	2.00	0.44
1:A:1417:VAL:O	1:A:1418:SER:HA	2.18	0.44
1:A:1385:LEU:CD1	1:A:1392[B]:CYS:SG	3.04	0.43
1:B:1446:VAL:HB	1:B:1484:LYS:HG2	2.01	0.42
1:B:1434:PHE:CZ	1:B:1540:MET:HE1	2.55	0.42
1:B:1506:MET:HE3	1:B:1538:PRO:HG3	2.01	0.41
1:B:1379:MET:N	5:B:1924:HOH:O	2.53	0.41
1:A:1418:SER:HB2	1:A:1419:GLY:N	2.36	0.41
1:B:1521:LEU:O	1:B:1707:PRO:HD3	2.20	0.41
1:A:1379[A]:MET:HE1	5:A:2244:HOH:O	2.20	0.41
1:A:1488:TRP:HA	1:A:1489:PRO:C	2.45	0.40
1:A:1407:GLU:HG3	1:A:1410:LYS:HE2	2.03	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:2206:HOH:O	5:B:1904:HOH:O[2_657]	1.98	0.22
5:A:2163:HOH:O	5:B:1909:HOH:O[2_557]	2.19	0.01

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%)	333 (97%)	10 (3%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	342/372 (92%)	328 (96%)	14 (4%)	0	100	100
All	All	685/744 (92%)	661 (96%)	24 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	299/331 (90%)	291 (97%)	8 (3%)	39	23
1	B	302/331 (91%)	292 (97%)	10 (3%)	33	16
All	All	601/662 (91%)	583 (97%)	18 (3%)	36	19

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

Mol	Chain	Res	Type
1	A	1418	SER
1	A	1577	ILE
1	A	1579	GLU
1	A	1583	ASP
1	A	1585	GLU
1	A	1606	HIS
1	A	1644	HIS
1	A	1717	THR
1	B	1459	VAL
1	B	1577	ILE
1	B	1582	HIS
1	B	1588	LYS
1	B	1603	ILE
1	B	1604	HIS
1	B	1605	ASP
1	B	1606	HIS
1	B	1607	LYS
1	B	1608	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1648	HIS
1	A	1650	GLN
1	B	1479	GLN
1	B	1648	HIS
1	B	1650	GLN
1	B	1656	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 8 ligands modelled in this entry, 7 are monoatomic - leaving 1 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	$\# Z > 2$
2	S5A	A	1801	-	19,19,19	1.25	1 (5%)	25,25,25	0.86	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	S5A	A	1801	-	-	3/9/9/9	0/2/2/2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1801	S5A	C-C14	3.53	1.58	1.50

There are no bond angle outliers.

There are no chirality outliers.

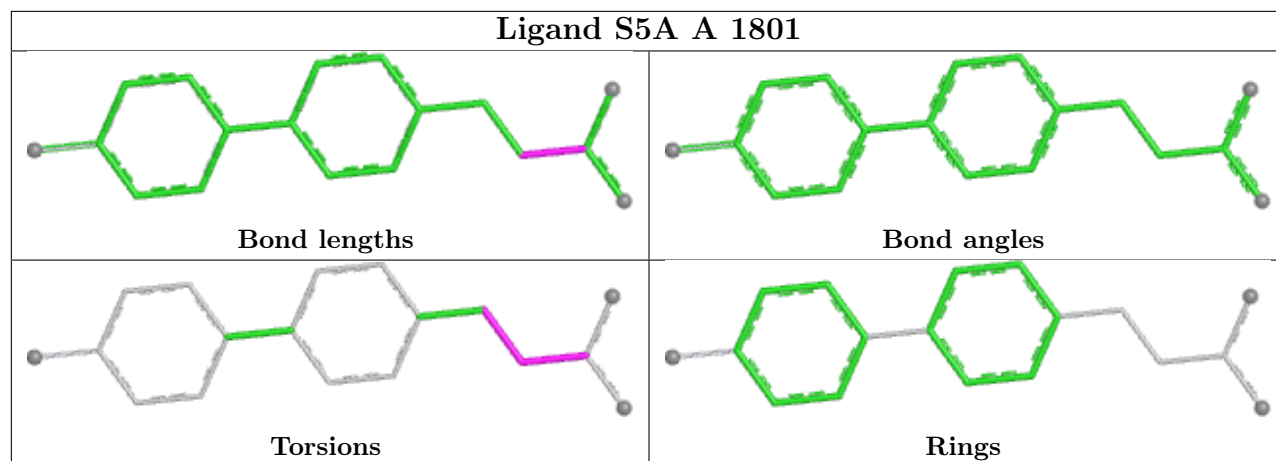
All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1801	S5A	C14-C-C1-C2
2	A	1801	S5A	C1-C-C14-O2
2	A	1801	S5A	C1-C-C14-O1

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%)	0.69	42 (12%) 8 9	5, 29, 55, 102	26 (7%)
1	B	344/372 (92%)	0.48	25 (7%) 21 24	9, 35, 68, 152	10 (2%)
All	All	684/744 (91%)	0.58	67 (9%) 13 15	5, 32, 63, 152	36 (5%)

All (67) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1581	ALA	12.8
1	A	1379[A]	MET	12.2
1	A	1582	HIS	10.8
1	A	1406[A]	ARG	10.7
1	B	1603	ILE	10.2
1	A	1598	VAL	10.1
1	A	1583	ASP	10.0
1	A	1580	GLY	9.6
1	A	1392[A]	CYS	9.5
1	A	1599	THR	9.4
1	B	1664[A]	GLU	9.1
1	A	1577	ILE	9.0
1	A	1588[A]	LYS	9.0
1	A	1676	PHE	8.9
1	B	1607	LYS	8.7
1	B	1602	ARG	8.7
1	B	1517[A]	ARG	8.4
1	A	1602	ARG	8.1
1	A	1584	GLU	8.1
1	A	1579	GLU	8.0
1	A	1564[A]	SER	7.9
1	B	1606	HIS	7.9
1	A	1596	ASP	7.3
1	A	1600	LYS	7.3

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Mol	Chain	Res	Type	RSRZ
1	A	1607	LYS	7.1
1	A	1605	ASP	6.9
1	A	1606	HIS	6.7
1	A	1603	ILE	6.5
1	A	1608	GLU	6.5
1	B	1604	HIS	6.5
1	B	1581	ALA	6.4
1	A	1609	LYS	6.2
1	A	1717	THR	6.2
1	B	1577	ILE	6.0
1	B	1608	GLU	5.8
1	A	1597	GLU	5.7
1	A	1604	HIS	5.4
1	A	1601	GLU	5.2
1	B	1601	GLU	4.9
1	B	1605	ASP	4.7
1	A	1578	GLY	4.4
1	B	1582	HIS	4.2
1	B	1587	LEU	3.8
1	A	1642	PRO	3.0
1	B	1583	ASP	2.8
1	A	1418	SER	2.8
1	A	1536	LEU	2.7
1	B	1642	PRO	2.6
1	B	1725	ASN	2.6
1	A	1644	HIS	2.5
1	A	1585	GLU	2.4
1	B	1663	TYR	2.4
1	A	1643	ASP	2.3
1	A	1537	GLY	2.3
1	A	1641	PRO	2.3
1	A	1637	GLY	2.3
1	A	1470	CYS	2.3
1	A	1592	GLU	2.2
1	B	1669	GLN	2.2
1	B	1666	TYR	2.2
1	B	1547	ILE	2.2
1	A	1407	GLU	2.2
1	A	1594	ASP	2.2
1	B	1576	PRO	2.2
1	B	1518	ASP	2.1
1	B	1592	GLU	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	1643	ASP	2.1

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](#)

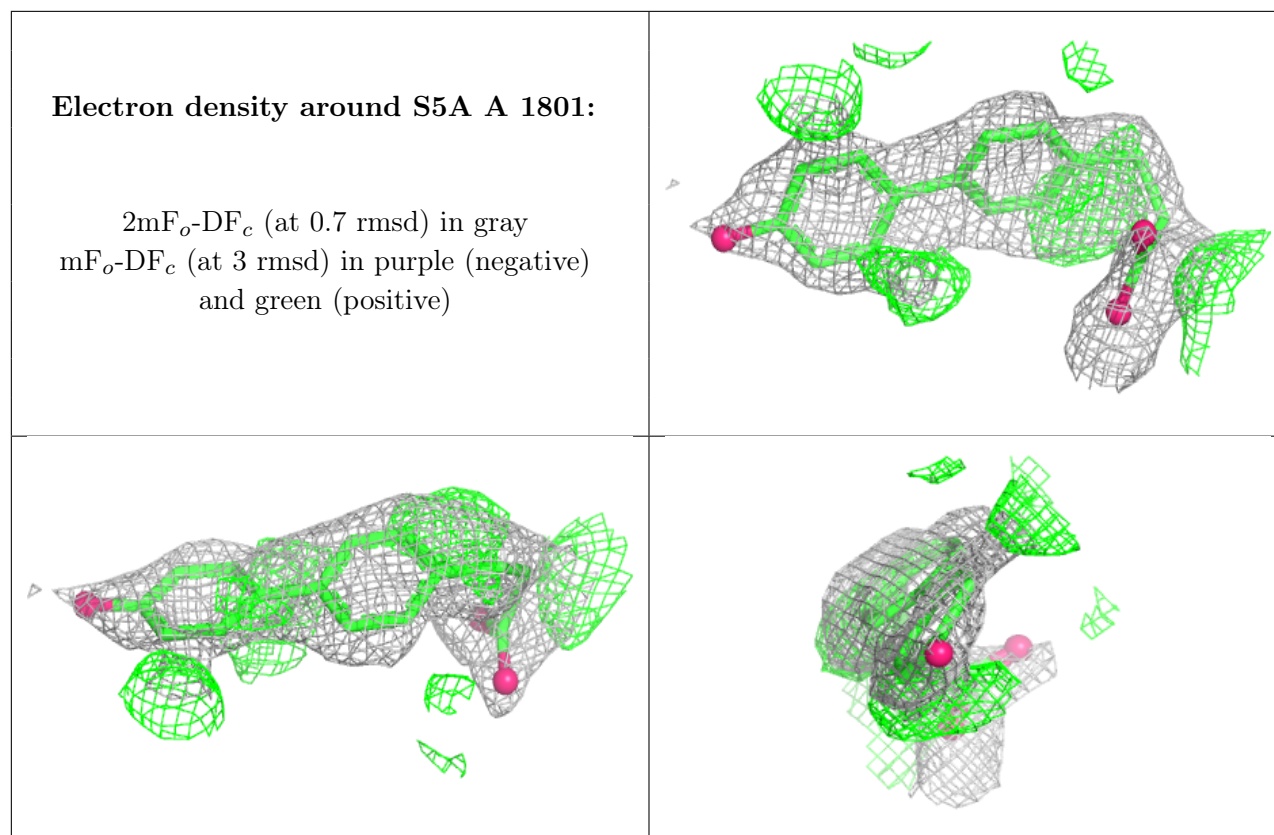
There are no oligosaccharides in this entry.

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
2	S5A	A	1801	18/18	0.72	0.23	34,36,44,45	18
3	CL	A	1805	1/1	0.89	0.13	53,53,53,53	0
4	MN	A	1806	1/1	0.95	0.06	58,58,58,58	0
4	MN	B	1802	1/1	0.97	0.05	41,41,41,41	0
3	CL	A	1804	1/1	0.98	0.08	36,36,36,36	0
3	CL	B	1801	1/1	0.98	0.06	40,40,40,40	0
3	CL	A	1803	1/1	0.99	0.04	41,41,41,41	0
3	CL	A	1802	1/1	1.00	0.03	25,25,25,25	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.