



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 09:47 AM UTC

PDB ID : 5QSN / pdb_00005qsn
Title : PanDDA analysis group deposition – Crystal Structure of human STAG1 in complex with Z1272480091
Authors : Newman, J.A.; Katis, V.L.; Gavard, A.E.; von Delft, F.; Arrowsmith, C.H.; Edwards, A.; Bountra, C.; Gileadi, O.
Deposited on : 2019-05-25
Resolution : 2.66 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

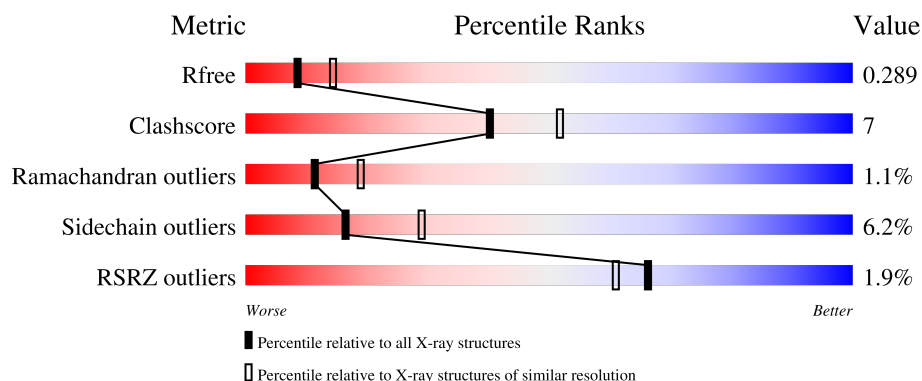
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.66 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	1110 (2.66-2.66)
Clashscore	190562	1141 (2.66-2.66)
Ramachandran outliers	187476	1126 (2.66-2.66)
Sidechain outliers	187428	1126 (2.66-2.66)
RSRZ outliers	180081	1110 (2.66-2.66)

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 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	459	
1	B	459	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 7091 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.

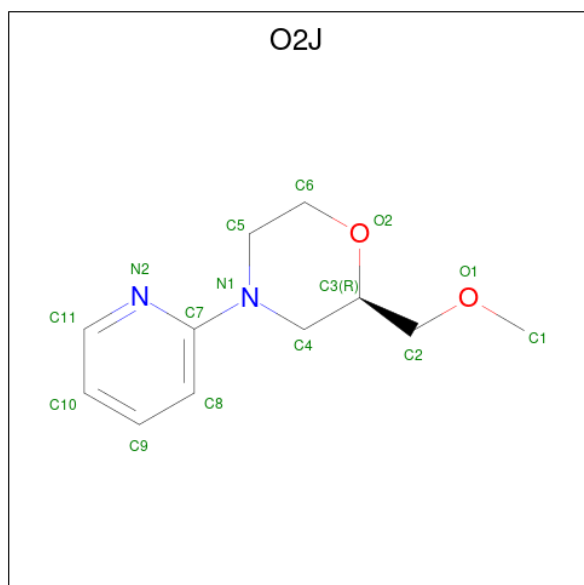
- Molecule 1 is a protein called Cohesin subunit SA-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	435	Total	C	N	O	S	0	0	0
			3493	2225	581	665	22			
1	B	425	Total	C	N	O	S	0	0	0
			3429	2186	570	651	22			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	457	SER	-	expression tag	UNP Q8WVM7
A	458	MET	-	expression tag	UNP Q8WVM7
B	457	SER	-	expression tag	UNP Q8WVM7
B	458	MET	-	expression tag	UNP Q8WVM7

- Molecule 2 is (2R)-2-(methoxymethyl)-4-(pyridin-2-yl)morpholine (CCD ID: O2J) (formula: C₁₁H₁₆N₂O₂) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	B	1	Total	C	N	O	0	0
			15	11	2	2		

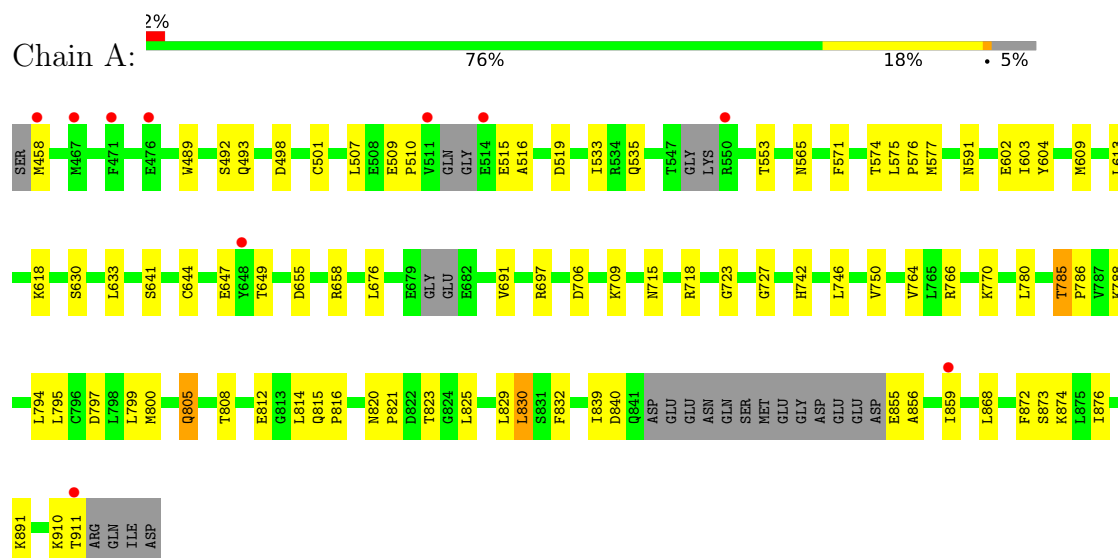
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	113	Total	O	0	0
			113	113		
3	B	41	Total	O	0	0
			41	41		

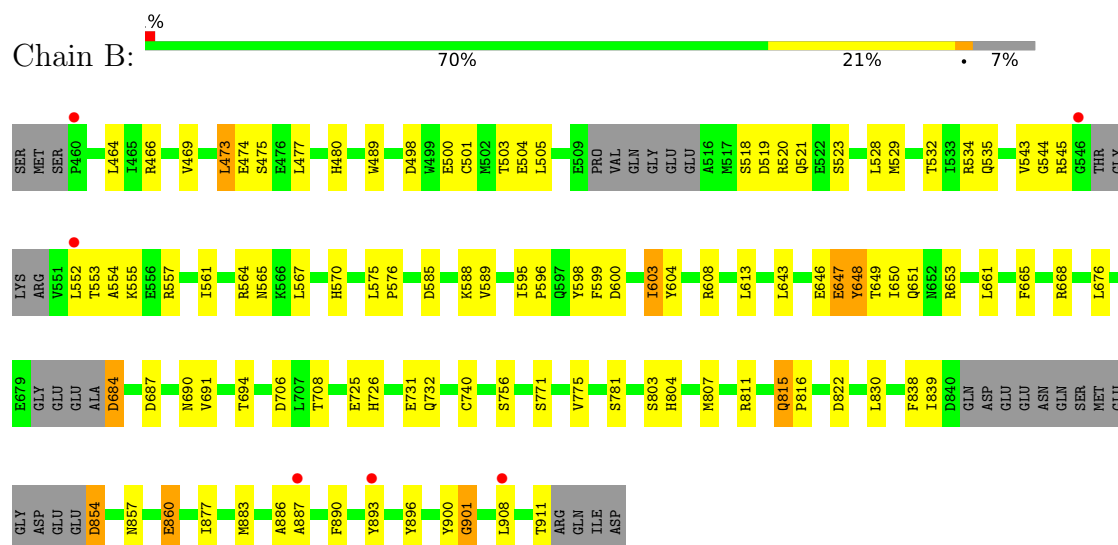
3 Residue-property plots [i](#)

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: Cohesin subunit SA-1



• Molecule 1: Cohesin subunit SA-1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	154.98Å 167.69Å 47.11Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	113.82 – 2.66 113.82 – 2.66	Depositor EDS
% Data completeness (in resolution range)	99.8 (113.82-2.66) 100.0 (113.82-2.66)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.36 (at 2.65Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.217 , 0.288 0.223 , 0.289	Depositor DCC
R_{free} test set	1833 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	63.9	Xtriage
Anisotropy	0.035	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 43.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7091	wwPDB-VP
Average B, all atoms (Å ²)	81.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.81% 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: O2J

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	1.03	0/3554	1.56	1/4803 (0.0%)
1	B	1.05	0/3489	1.56	2/4711 (0.0%)
All	All	1.04	0/7043	1.56	3/9514 (0.0%)

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	B	0	1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	820	ASN	CB-CA-C	6.08	116.78	109.85
1	B	740	CYS	CA-C-N	5.01	127.25	120.38
1	B	740	CYS	C-N-CA	5.01	127.25	120.38

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	599	PHE	Peptide

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	3493	0	3473	41	0
1	B	3429	0	3418	50	0
2	B	15	0	0	0	0
3	A	113	0	0	5	0
3	B	41	0	0	1	0
All	All	7091	0	6891	91	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 7.

All (91) 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:469:VAL:O	1:B:473:LEU:HD23	1.86	0.74
1:A:829:LEU:C	3:A:1001:HOH:O	2.32	0.72
1:A:829:LEU:O	3:A:1001:HOH:O	2.09	0.71
1:B:857:ASN:O	1:B:860:GLU:HB3	1.92	0.69
1:B:830:LEU:CD1	1:B:886:ALA:HA	2.27	0.65
1:A:800:MET:HE3	1:A:874:LYS:HB3	1.79	0.64
1:A:766:ARG:NH1	1:A:816:PRO:O	2.33	0.60
1:A:655:ASP:OD1	1:A:658:ARG:NH2	2.34	0.60
1:B:554:ALA:HA	1:B:557:ARG:HB3	1.83	0.60
1:A:676:LEU:HD22	1:A:691:VAL:HG21	1.86	0.58
1:B:771:SER:O	1:B:775:VAL:HG23	2.05	0.57
1:A:489:TRP:CD2	1:A:535:GLN:HG2	2.41	0.56
1:B:604:TYR:CE1	1:B:613:LEU:HB2	2.41	0.56
1:B:815:GLN:N	1:B:816:PRO:HD2	2.21	0.55
1:B:854:ASP:HB2	1:B:857:ASN:HD22	1.72	0.55
1:A:839:ILE:HD12	1:A:839:ILE:O	2.08	0.54
1:B:676:LEU:HD22	1:B:691:VAL:HG21	1.90	0.54
1:B:732:GLN:NE2	3:B:1102:HOH:O	2.32	0.53
1:A:872:PHE:CE2	1:A:876:ILE:HD11	2.44	0.53
1:B:887:ALA:HA	1:B:890:PHE:CE2	2.44	0.53
1:A:489:TRP:CE2	1:A:535:GLN:HG2	2.44	0.53
1:A:785:THR:N	1:A:786:PRO:CD	2.72	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:529:MET:O	1:B:532:THR:OG1	2.26	0.52
1:A:489:TRP:CE2	1:A:493:GLN:HG3	2.45	0.52
1:A:805:GLN:O	1:A:808:THR:HG23	2.10	0.52
1:A:873:SER:HA	1:A:876:ILE:HD12	1.92	0.52
1:B:650:ILE:O	1:B:653:ARG:N	2.43	0.52
1:A:832:PHE:N	3:A:1001:HOH:O	2.42	0.51
1:A:510:PRO:CG	1:A:516:ALA:HB2	2.40	0.51
1:B:830:LEU:HD13	1:B:886:ALA:HA	1.92	0.51
1:B:585:ASP:O	1:B:589:VAL:HG23	2.11	0.51
1:A:633:LEU:HD13	1:A:697:ARG:HA	1.93	0.49
1:A:821:PRO:CB	1:A:825:LEU:HD23	2.41	0.49
1:B:503:THR:HG21	1:B:570:HIS:NE2	2.27	0.49
1:A:910:LYS:O	1:A:911:THR:HG23	2.13	0.49
1:B:647:GLU:HB3	1:B:648:TYR:CD1	2.48	0.49
1:A:498:ASP:CG	1:A:501:CYS:HB2	2.37	0.49
1:B:561:ILE:O	1:B:565:ASN:ND2	2.45	0.48
1:A:832:PHE:HB3	3:A:1001:HOH:O	2.13	0.48
1:A:604:TYR:CD1	1:A:613:LEU:HB2	2.49	0.48
1:B:553:THR:C	1:B:555:LYS:H	2.22	0.48
1:B:887:ALA:HA	1:B:890:PHE:CD2	2.49	0.48
1:A:575:LEU:N	1:A:576:PRO:CD	2.76	0.47
1:B:661:LEU:HG	1:B:665:PHE:CE1	2.49	0.47
1:A:814:LEU:C	1:A:816:PRO:HD2	2.39	0.47
1:B:564:ARG:HD3	1:B:600:ASP:HB2	1.95	0.47
1:A:815:GLN:N	1:A:816:PRO:CD	2.78	0.47
1:B:803:SER:O	1:B:804:HIS:C	2.58	0.47
1:A:706:ASP:OD1	1:A:706:ASP:C	2.57	0.46
1:B:603:ILE:HG22	1:B:603:ILE:O	2.16	0.46
1:B:661:LEU:HD11	1:B:665:PHE:CZ	2.51	0.46
1:A:780:LEU:O	1:A:788:LYS:HG2	2.16	0.45
1:A:746:LEU:O	1:A:750:VAL:HG23	2.17	0.45
1:B:883:MET:CE	1:B:908:LEU:HD23	2.46	0.45
1:B:567:LEU:C	1:B:567:LEU:HD12	2.42	0.45
1:A:830:LEU:HD13	1:A:830:LEU:HA	1.88	0.45
1:A:715:ASN:OD1	1:A:718:ARG:NH1	2.48	0.44
1:A:458:MET:HA	3:A:1020:HOH:O	2.16	0.44
1:B:706:ASP:OD1	1:B:708:THR:HG23	2.18	0.44
1:B:585:ASP:HB3	1:B:588:LYS:HB2	1.98	0.44
1:B:822:ASP:C	1:B:822:ASP:OD1	2.60	0.44
1:B:646:GLU:HA	1:B:651:GLN:OE1	2.18	0.44
1:B:725:GLU:HG3	1:B:726:HIS:ND1	2.33	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:477:LEU:O	1:B:480:HIS:HB2	2.18	0.44
1:B:604:TYR:CD1	1:B:613:LEU:HB2	2.53	0.43
1:B:528:LEU:O	1:B:532:THR:HG23	2.18	0.43
1:A:742:HIS:CG	1:A:794:LEU:HD22	2.53	0.43
1:B:575:LEU:HB3	1:B:576:PRO:HD3	2.01	0.43
1:B:595:ILE:HB	1:B:596:PRO:HD3	2.00	0.43
1:A:533:ILE:HD11	1:A:571:PHE:CE2	2.53	0.43
1:B:500:GLU:HA	1:B:570:HIS:CE1	2.53	0.43
1:B:519:ASP:O	1:B:521:GLN:N	2.52	0.42
1:B:684:ASP:HB3	1:B:687:ASP:OD2	2.19	0.42
1:B:498:ASP:OD2	1:B:501:CYS:HB2	2.19	0.42
1:B:807:MET:HA	1:B:811:ARG:O	2.19	0.42
1:A:565:ASN:HD22	1:A:603:ILE:HD11	1.85	0.42
1:B:838:PHE:HD1	1:B:896:TYR:CE2	2.38	0.42
1:B:469:VAL:O	1:B:473:LEU:CD2	2.64	0.42
1:A:507:LEU:HD13	1:A:577:MET:HE3	2.02	0.41
1:A:723:GLY:O	1:A:727:GLY:HA2	2.19	0.41
1:A:855:GLU:HG3	1:A:856:ALA:N	2.36	0.41
1:A:814:LEU:C	1:A:816:PRO:CD	2.94	0.41
1:B:690:ASN:O	1:B:694:THR:OG1	2.38	0.41
1:B:900:TYR:O	1:B:901:GLY:C	2.64	0.41
1:A:799:LEU:HD23	1:A:799:LEU:HA	1.96	0.41
1:B:489:TRP:CD2	1:B:535:GLN:HG2	2.55	0.41
1:B:544:GLY:O	1:B:545:ARG:HG2	2.21	0.41
1:A:509:GLU:HA	1:A:509:GLU:OE1	2.21	0.41
1:A:574:THR:HA	1:A:577:MET:HE2	2.02	0.40
1:B:501:CYS:O	1:B:504:GLU:HB2	2.21	0.40
1:B:534:ARG:HG3	1:B:598:TYR:OH	2.22	0.40

There are no symmetry-related clashes.

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	425/459 (93%)	400 (94%)	22 (5%)	3 (1%)	18	30
1	B	415/459 (90%)	373 (90%)	36 (9%)	6 (1%)	9	14
All	All	840/918 (92%)	773 (92%)	58 (7%)	9 (1%)	11	19

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	644	CYS
1	A	647	GLU
1	B	603	ILE
1	B	901	GLY
1	A	515	GLU
1	B	520	ARG
1	B	649	THR
1	B	860	GLU
1	B	543	VAL

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	391/415 (94%)	367 (94%)	24 (6%)	17	30
1	B	385/415 (93%)	361 (94%)	24 (6%)	16	29
All	All	776/830 (94%)	728 (94%)	48 (6%)	16	29

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	492	SER
1	A	519	ASP
1	A	553	THR
1	A	591	ASN
1	A	602	GLU
1	A	609	MET
1	A	618	LYS

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Mol	Chain	Res	Type
1	A	630	SER
1	A	641	SER
1	A	649	THR
1	A	709	LYS
1	A	764	VAL
1	A	770	LYS
1	A	785	THR
1	A	795	LEU
1	A	797	ASP
1	A	805	GLN
1	A	812	GLU
1	A	823	THR
1	A	830	LEU
1	A	840	ASP
1	A	859	ILE
1	A	868	LEU
1	A	891	LYS
1	B	464	LEU
1	B	466	ARG
1	B	473	LEU
1	B	474	GLU
1	B	475	SER
1	B	505	LEU
1	B	518	SER
1	B	523	SER
1	B	552	LEU
1	B	608	ARG
1	B	643	LEU
1	B	647	GLU
1	B	648	TYR
1	B	668	ARG
1	B	684	ASP
1	B	731	GLU
1	B	756	SER
1	B	781	SER
1	B	815	GLN
1	B	839	ILE
1	B	854	ASP
1	B	877	ILE
1	B	893	TYR
1	B	911	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17)

such sidechains are listed below:

Mol	Chain	Res	Type
1	A	463	ASN
1	A	565	ASN
1	A	591	ASN
1	A	678	GLN
1	A	736	GLN
1	A	748	GLN
1	A	778	GLN
1	A	804	HIS
1	A	820	ASN
1	B	480	HIS
1	B	560	GLN
1	B	591	ASN
1	B	619	GLN
1	B	703	ASN
1	B	715	ASN
1	B	736	GLN
1	B	857	ASN

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](#)

1 ligand is 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	$\# Z > 2$
2	O2J	B	1001	-	16,16,16	0.32	0	17,20,20	0.79	1 (5%)

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	O2J	B	1001	-	-	1/7/17/17	0/2/2/2

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1001	O2J	O2-C6-C5	2.30	113.76	110.73

There are no chirality outliers.

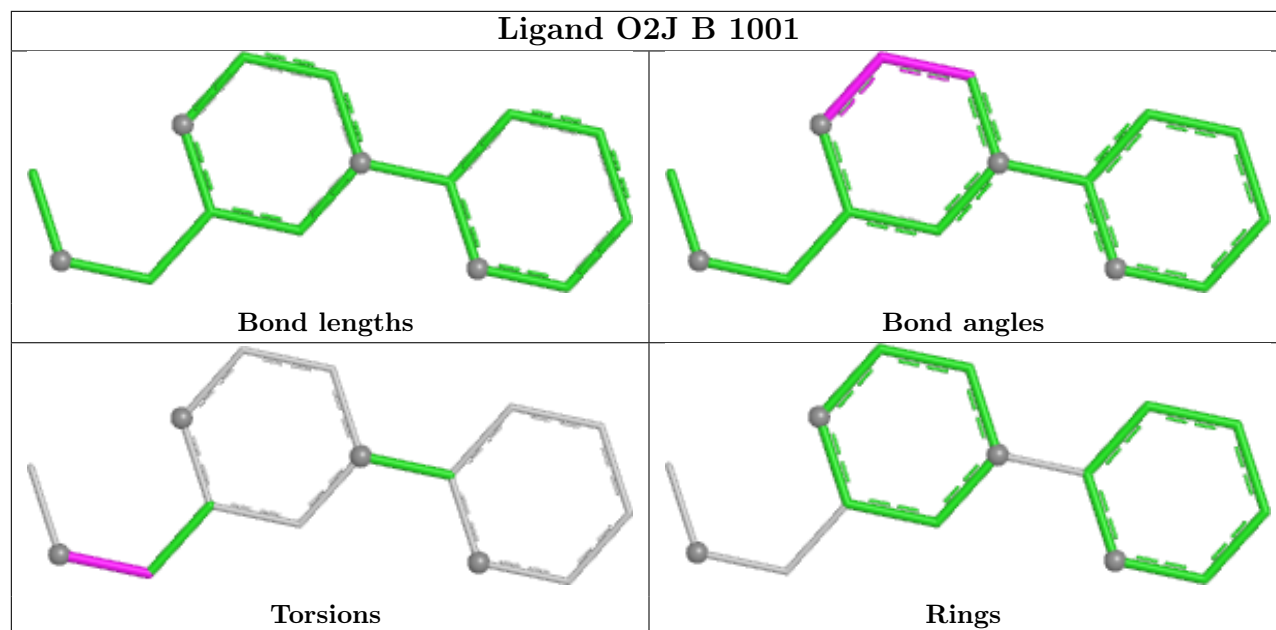
All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	1001	O2J	C3-C2-O1-C1

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	435/459 (94%)	-0.17	10 (2%) 61 54	44, 69, 113, 138	0
1	B	425/459 (92%)	0.11	6 (1%) 73 69	54, 86, 127, 169	0
All	All	860/918 (93%)	-0.03	16 (1%) 66 61	44, 78, 122, 169	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	511	VAL	5.1
1	B	460	PRO	5.0
1	A	458	MET	3.7
1	B	887	ALA	2.9
1	B	893	TYR	2.9
1	B	908	LEU	2.6
1	A	467	MET	2.6
1	A	514	GLU	2.6
1	B	546	GLY	2.4
1	A	911	THR	2.4
1	B	552	LEU	2.3
1	A	859	ILE	2.2
1	A	471	PHE	2.2
1	A	476	GLU	2.1
1	A	648	TYR	2.1
1	A	550	ARG	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](#)

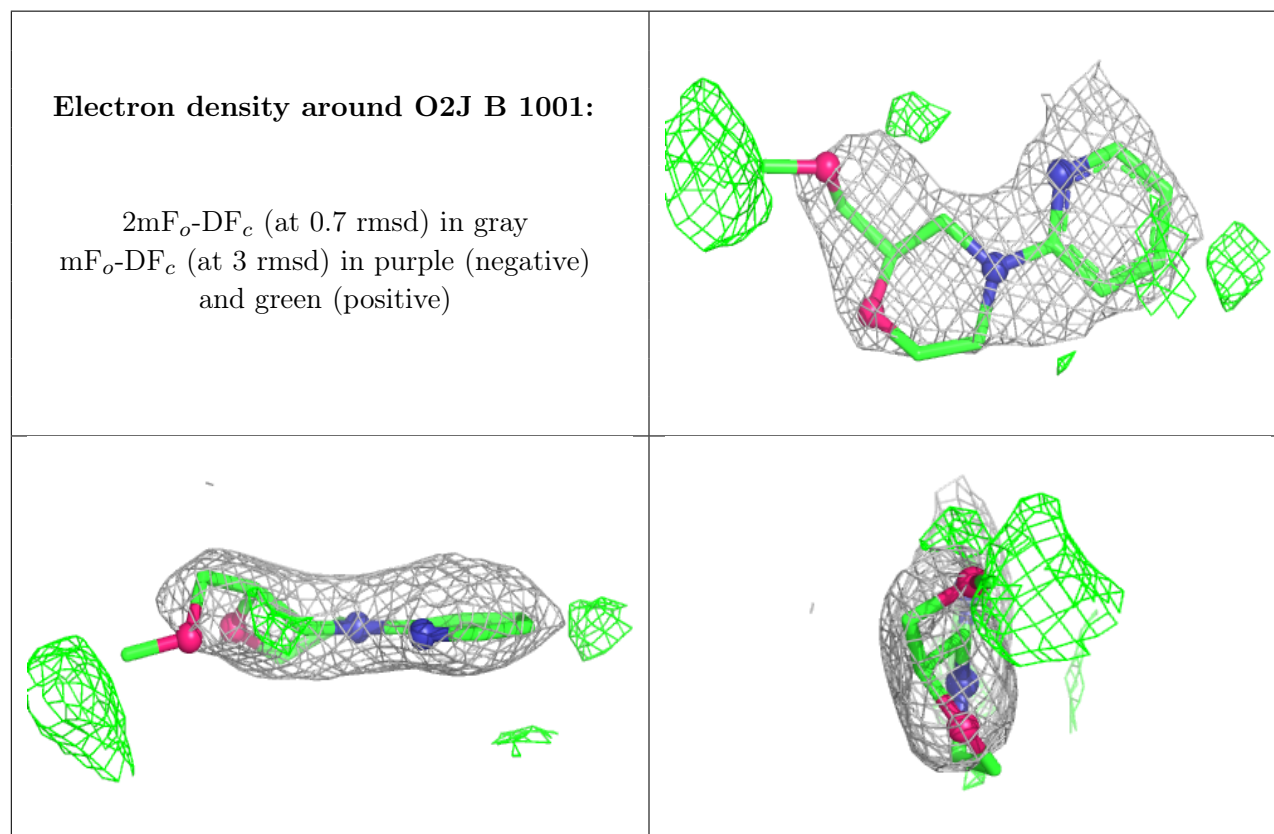
There are no oligosaccharides in this entry.

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	RSR	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	O2J	B	1001	15/15	0.81	0.19	86,92,100,102	15

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.