



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 06:08 AM UTC

PDB ID : 5SDP / pdb_00005sdp
Title : PanDDA analysis group deposition – Crystal Structure of Porphyromonas gingivalis in complex with Z2277255954
Authors : Tham, C.T.; Coker, J.A.; Krojer, T.; Foster, W.R.; Koekemoer, L.; Douangamath, A.; Talon, R.; Fearon, D.; von Delft, F.; Yue, W.W.; Bountra, C.; Bezerra, G.A.
Deposited on : 2022-01-20
Resolution : 2.19 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](#) ①) 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

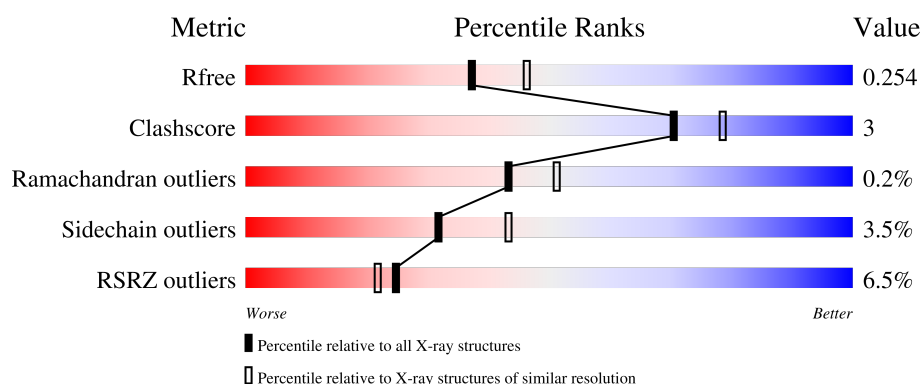
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.19 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	707	7% 87% 10% ..
1	B	707	5% 88% 10% ..

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 11230 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.

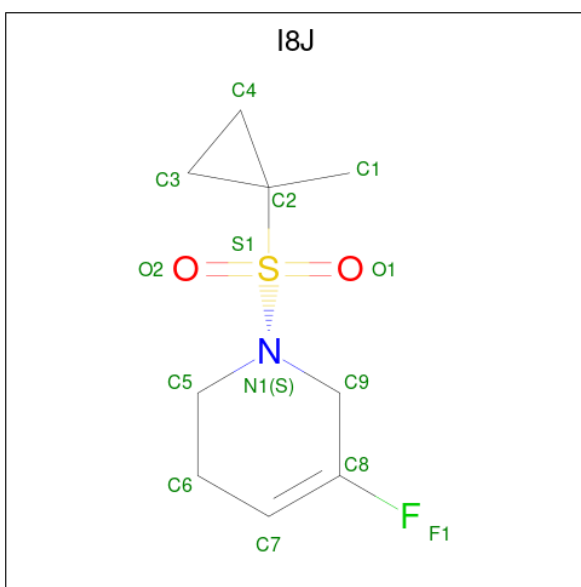
- Molecule 1 is a protein called Asp/Glu-specific dipeptidyl-peptidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	697	Total	C	N	O	S	0	1	0
			5502	3494	950	1031	27			
1	B	695	Total	C	N	O	S	0	1	0
			5469	3471	944	1027	27			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	21	MET	-	initiating methionine	UNP B2RID1
A	721	ALA	-	expression tag	UNP B2RID1
A	722	HIS	-	expression tag	UNP B2RID1
A	723	HIS	-	expression tag	UNP B2RID1
A	724	HIS	-	expression tag	UNP B2RID1
A	725	HIS	-	expression tag	UNP B2RID1
A	726	HIS	-	expression tag	UNP B2RID1
A	727	HIS	-	expression tag	UNP B2RID1
B	21	MET	-	initiating methionine	UNP B2RID1
B	721	ALA	-	expression tag	UNP B2RID1
B	722	HIS	-	expression tag	UNP B2RID1
B	723	HIS	-	expression tag	UNP B2RID1
B	724	HIS	-	expression tag	UNP B2RID1
B	725	HIS	-	expression tag	UNP B2RID1
B	726	HIS	-	expression tag	UNP B2RID1
B	727	HIS	-	expression tag	UNP B2RID1

- Molecule 2 is 5-fluoro-1-(1-methylcyclopropane-1-sulfonyl)-1,2,3,6-tetrahydropyridine (CCD ID: I8J) (formula: C₉H₁₄FN₂O₂S) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	F	N	O	S	0	1
			14	9	1	1	2	1		
2	B	1	Total	C	F	N	O	S	0	1
			14	9	1	1	2	1		

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

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

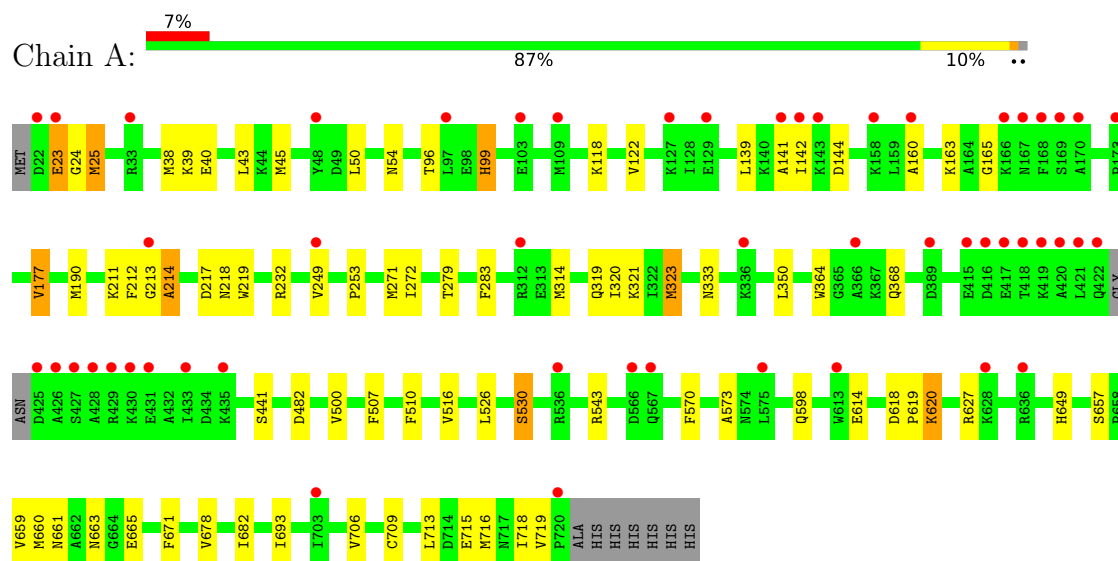
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	117	Total	O	0	0
			117	117		
4	B	111	Total	O	0	0
			111	111		

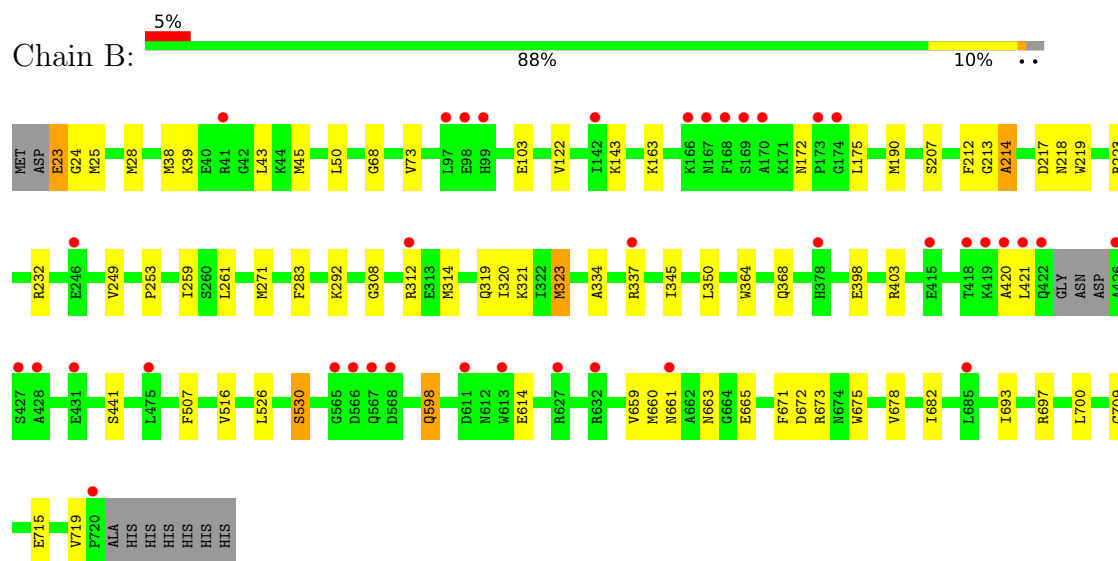
3 Residue-property plots [i](#)

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: Asp/Glu-specific dipeptidyl-peptidase



- Molecule 1: Asp/Glu-specific dipeptidyl-peptidase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	102.11Å 117.24Å 147.93Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	84.03 – 2.19 84.03 – 2.19	Depositor EDS
% Data completeness (in resolution range)	72.7 (84.03-2.19) 72.7 (84.03-2.19)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.55 (at 2.20Å)	Xtriage
Refinement program	BUSTER 2.10.3 (18-SEP-2020)	Depositor
R, R_{free}	0.215 , 0.248 0.222 , 0.254	Depositor DCC
R_{free} test set	3381 reflections (3.71%)	wwPDB-VP
Wilson B-factor (Å ²)	42.4	Xtriage
Anisotropy	0.028	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 28.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	11230	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.83% 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: CL, I8J

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	0.75	3/5626 (0.1%)	1.12	13/7618 (0.2%)
1	B	0.76	3/5592 (0.1%)	1.10	10/7575 (0.1%)
All	All	0.75	6/11218 (0.1%)	1.11	23/15193 (0.2%)

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	323	MET	SD-CE	-6.57	1.63	1.79
1	A	214	ALA	CA-C	5.94	1.58	1.53
1	B	214	ALA	CA-C	5.77	1.58	1.53
1	B	25	MET	SD-CE	-5.76	1.65	1.79
1	B	323	MET	SD-CE	-5.51	1.65	1.79
1	A	25	MET	SD-CE	-5.28	1.66	1.79

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	24	GLY	N-CA-C	7.48	130.91	113.18
1	A	141	ALA	CA-C-N	7.37	135.24	121.97
1	A	141	ALA	C-N-CA	7.37	135.24	121.97
1	B	24	GLY	N-CA-C	7.03	121.41	111.14
1	B	23	GLU	CB-CG-CD	6.24	123.20	112.60
1	A	660	MET	CA-C-N	6.05	131.82	121.29
1	A	660	MET	C-N-CA	6.05	131.82	121.29
1	A	253	PRO	CA-C-N	5.63	127.83	120.28
1	A	253	PRO	C-N-CA	5.63	127.83	120.28
1	A	165	GLY	N-CA-C	5.55	118.02	111.63
1	B	660	MET	CA-C-N	5.48	132.00	121.54
1	B	660	MET	C-N-CA	5.48	132.00	121.54
1	A	144	ASP	CA-CB-CG	5.35	117.95	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	682	ILE	N-CA-C	-5.32	106.98	111.56
1	A	482	ASP	N-CA-C	5.26	116.83	111.14
1	B	122	VAL	N-CA-CB	5.14	119.03	111.52
1	B	28	MET	CA-C-N	5.11	127.64	120.28
1	B	28	MET	C-N-CA	5.11	127.64	120.28
1	A	507	PHE	CA-CB-CG	-5.09	108.71	113.80
1	B	507	PHE	CA-CB-CG	-5.08	108.72	113.80
1	A	682	ILE	N-CA-C	-5.07	107.20	111.56
1	B	671	PHE	CA-CB-CG	5.06	118.86	113.80
1	A	671	PHE	CA-CB-CG	5.01	118.81	113.80

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	5502	0	5294	41	0
1	B	5469	0	5243	33	0
2	A	14	0	0	1	0
2	B	14	0	0	1	0
3	A	2	0	0	0	0
3	B	1	0	0	0	0
4	A	117	0	0	0	0
4	B	111	0	0	0	0
All	All	11230	0	10537	74	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 3.

All (74) 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:223:ARG:NH2	1:B:672:ASP:OD2	2.06	0.87
1:A:713:LEU:HA	1:A:716:MET:HE2	1.56	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:96:THR:HG23	1:A:99:HIS:HE1	1.44	0.83
1:B:319:GLN:HG3	1:B:323:MET:HE2	1.59	0.82
1:A:319:GLN:HG3	1:A:323:MET:HE2	1.60	0.82
1:B:314:MET:HG2	1:B:321:LYS:HA	1.69	0.74
1:A:314:MET:HG2	1:A:321:LYS:HA	1.70	0.73
1:B:213:GLY:HA3	1:B:217:ASP:HB2	1.70	0.72
1:B:50:LEU:HD21	1:B:271:MET:HE1	1.72	0.71
1:B:334:ALA:HA	1:B:337[B]:ARG:HD3	1.74	0.70
1:B:261:LEU:HD21	1:B:700:LEU:HD12	1.73	0.69
1:A:213:GLY:HA3	1:A:217:ASP:HB2	1.73	0.69
1:A:96:THR:HG23	1:A:99:HIS:CE1	2.27	0.68
1:A:25:MET:HE1	1:A:573:ALA:HB2	1.77	0.67
1:B:73:VAL:HG11	1:B:253:PRO:HG3	1.76	0.67
1:A:50:LEU:HD21	1:A:271:MET:HE1	1.77	0.66
1:B:232:ARG:NH2	1:B:715:GLU:OE2	2.31	0.64
1:A:232:ARG:NH2	1:A:715:GLU:OE2	2.32	0.62
1:B:526:LEU:O	1:B:530:SER:HB2	2.00	0.62
1:A:526:LEU:O	1:A:530:SER:HB2	2.00	0.61
1:B:308:GLY:O	1:B:312:ARG:HD2	2.00	0.61
1:A:25:MET:CE	1:A:573:ALA:HB2	2.31	0.60
1:B:398:GLU:O	1:B:403:ARG:HG2	2.01	0.60
1:B:172:ASN:HB3	1:B:175:LEU:HD12	1.85	0.58
1:B:43:LEU:HD11	1:B:45:MET:HG2	1.86	0.58
1:A:620:LYS:HD3	1:A:706:VAL:HA	1.86	0.57
1:A:96:THR:CG2	1:A:99:HIS:HE1	2.15	0.56
1:A:139:LEU:O	1:A:142:ILE:HG13	2.04	0.56
1:B:259:ILE:HG21	1:B:700:LEU:HD11	1.87	0.56
1:B:50:LEU:CD2	1:B:271:MET:HE1	2.36	0.56
1:B:163:LYS:HB3	1:B:190:MET:HE1	1.87	0.56
1:A:25:MET:HE1	1:A:279:THR:CG2	2.36	0.56
1:A:219:TRP:HB2	2:A:801[B]:I8J:C6	2.36	0.55
1:B:232:ARG:HH22	1:B:715:GLU:CD	2.14	0.55
1:A:232:ARG:HH22	1:A:715:GLU:CD	2.15	0.55
1:A:163:LYS:HB3	1:A:190:MET:HE1	1.88	0.55
1:A:50:LEU:CD2	1:A:271:MET:HE1	2.38	0.54
1:A:160:ALA:HB1	1:A:177:VAL:CG1	2.38	0.54
1:A:25:MET:HE2	1:A:570:PHE:CE2	2.43	0.53
1:A:212:PHE:O	1:A:614:GLU:O	2.28	0.52
1:B:212:PHE:O	1:B:614:GLU:O	2.29	0.50
1:A:43:LEU:HD11	1:A:45:MET:HG2	1.93	0.50
1:A:25:MET:CE	1:A:279:THR:HG21	2.42	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:319:GLN:O	1:A:323:MET:HG3	2.13	0.49
1:A:25:MET:HE2	1:A:570:PHE:HE2	1.78	0.49
1:A:214:ALA:HB3	1:A:614:GLU:HB3	1.95	0.48
1:A:716:MET:HE3	1:A:718:ILE:HD11	1.95	0.48
1:B:320:ILE:HD13	1:B:323:MET:HE3	1.96	0.47
1:B:319:GLN:O	1:B:323:MET:HG3	2.15	0.47
1:B:364:TRP:O	1:B:368:GLN:HG2	2.15	0.47
1:A:25:MET:HE1	1:A:279:THR:HG21	1.97	0.46
1:B:214:ALA:HB3	1:B:614:GLU:HB3	1.97	0.45
1:A:364:TRP:O	1:A:368:GLN:HG3	2.16	0.45
1:A:25:MET:HE1	1:A:279:THR:HG23	1.99	0.45
1:A:320:ILE:HD13	1:A:323:MET:HE3	1.98	0.45
1:A:661:ASN:HB3	1:A:663:ASN:H	1.82	0.45
1:B:334:ALA:HA	1:B:337[B]:ARG:CD	2.46	0.44
1:A:38:MET:HE2	1:A:43:LEU:HD22	2.00	0.44
1:B:659:VAL:HG21	1:B:693:ILE:HG12	2.00	0.43
1:A:160:ALA:HB1	1:A:177:VAL:HG12	2.00	0.43
1:A:500:VAL:HG22	1:A:510:PHE:HB2	2.00	0.43
1:B:219:TRP:CZ2	1:B:337[B]:ARG:HG2	2.54	0.42
1:B:675:TRP:HB2	2:B:801[B]:I8J:C8	2.49	0.42
1:B:421:LEU:HB2	1:B:516:VAL:HG12	2.00	0.42
1:A:661:ASN:HD22	1:A:665:GLU:HB2	1.85	0.42
1:B:38:MET:HE2	1:B:43:LEU:HD22	2.02	0.42
1:B:661:ASN:HB3	1:B:663:ASN:H	1.85	0.42
1:A:272:ILE:HG23	1:A:657:SER:HB3	2.02	0.41
1:A:618:ASP:HA	1:A:619:PRO:HD3	1.95	0.41
1:B:661:ASN:HD22	1:B:665:GLU:HB2	1.85	0.41
1:A:659:VAL:HG21	1:A:693:ILE:HG12	2.01	0.41
1:B:598:GLN:HE22	1:B:697:ARG:NH2	2.19	0.41
1:A:23:GLU:O	1:A:649:HIS:HB2	2.20	0.41
1:B:292:LYS:HB2	1:B:345:ILE:HD13	2.03	0.40

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	694/707 (98%)	673 (97%)	21 (3%)	0	100	100
1	B	692/707 (98%)	671 (97%)	18 (3%)	3 (0%)	30	34
All	All	1386/1414 (98%)	1344 (97%)	39 (3%)	3 (0%)	43	51

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	143	LYS
1	B	420	ALA
1	B	68	GLY

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	561/597 (94%)	537 (96%)	24 (4%)	26	35
1	B	555/597 (93%)	540 (97%)	15 (3%)	39	53
All	All	1116/1194 (94%)	1077 (96%)	39 (4%)	32	43

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	23	GLU
1	A	39	LYS
1	A	40	GLU
1	A	54	ASN
1	A	99	HIS
1	A	118	LYS
1	A	122	VAL
1	A	177	VAL
1	A	211	LYS
1	A	218	ASN

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Mol	Chain	Res	Type
1	A	249	VAL
1	A	283	PHE
1	A	333	ASN
1	A	350	LEU
1	A	441	SER
1	A	516	VAL
1	A	530	SER
1	A	543	ARG
1	A	598	GLN
1	A	620	LYS
1	A	627	ARG
1	A	678	VAL
1	A	709	CYS
1	A	719	VAL
1	B	23	GLU
1	B	39	LYS
1	B	103	GLU
1	B	207	SER
1	B	218	ASN
1	B	249	VAL
1	B	283	PHE
1	B	350	LEU
1	B	441	SER
1	B	530	SER
1	B	598	GLN
1	B	673	ARG
1	B	678	VAL
1	B	709	CYS
1	B	719	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	100	ASN
1	A	157	GLN
1	A	186	ASN
1	A	355	GLN
1	A	567	GLN
1	A	583	GLN
1	B	186	ASN
1	B	332	GLN
1	B	355	GLN

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Mol	Chain	Res	Type
1	B	583	GLN
1	B	598	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 5 ligands modelled in this entry, 3 are monoatomic - leaving 2 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	I8J	B	801[B]	-	12,15,15	1.80	2 (16%)	18,24,24	1.85	3 (16%)
2	I8J	A	801[B]	-	12,15,15	1.48	1 (8%)	18,24,24	1.99	7 (38%)

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	I8J	B	801[B]	-	-	5/10/29/29	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	I8J	A	801[B]	-	-	6/10/29/29	0/2/2/2

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	801[B]	I8J	C2-S1	-5.29	1.75	1.81
2	A	801[B]	I8J	C2-S1	-4.23	1.76	1.81
2	B	801[B]	I8J	F1-C8	-2.48	1.34	1.36

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	801[B]	I8J	C6-C5-N1	4.69	113.63	109.19
2	B	801[B]	I8J	O2-S1-N1	-4.68	104.71	109.68
2	B	801[B]	I8J	O1-S1-N1	-3.74	105.72	109.68
2	B	801[B]	I8J	C2-S1-N1	3.24	115.34	106.20
2	A	801[B]	I8J	O1-S1-N1	-3.05	106.45	109.68
2	A	801[B]	I8J	C5-N1-S1	-2.97	111.78	116.78
2	A	801[B]	I8J	C5-C6-C7	-2.79	108.72	111.51
2	A	801[B]	I8J	C9-N1-C5	2.62	117.20	113.34
2	A	801[B]	I8J	O2-S1-N1	-2.11	107.45	109.68
2	A	801[B]	I8J	C8-C9-N1	2.09	114.97	110.67

There are no chirality outliers.

All (11) torsion outliers are listed below:

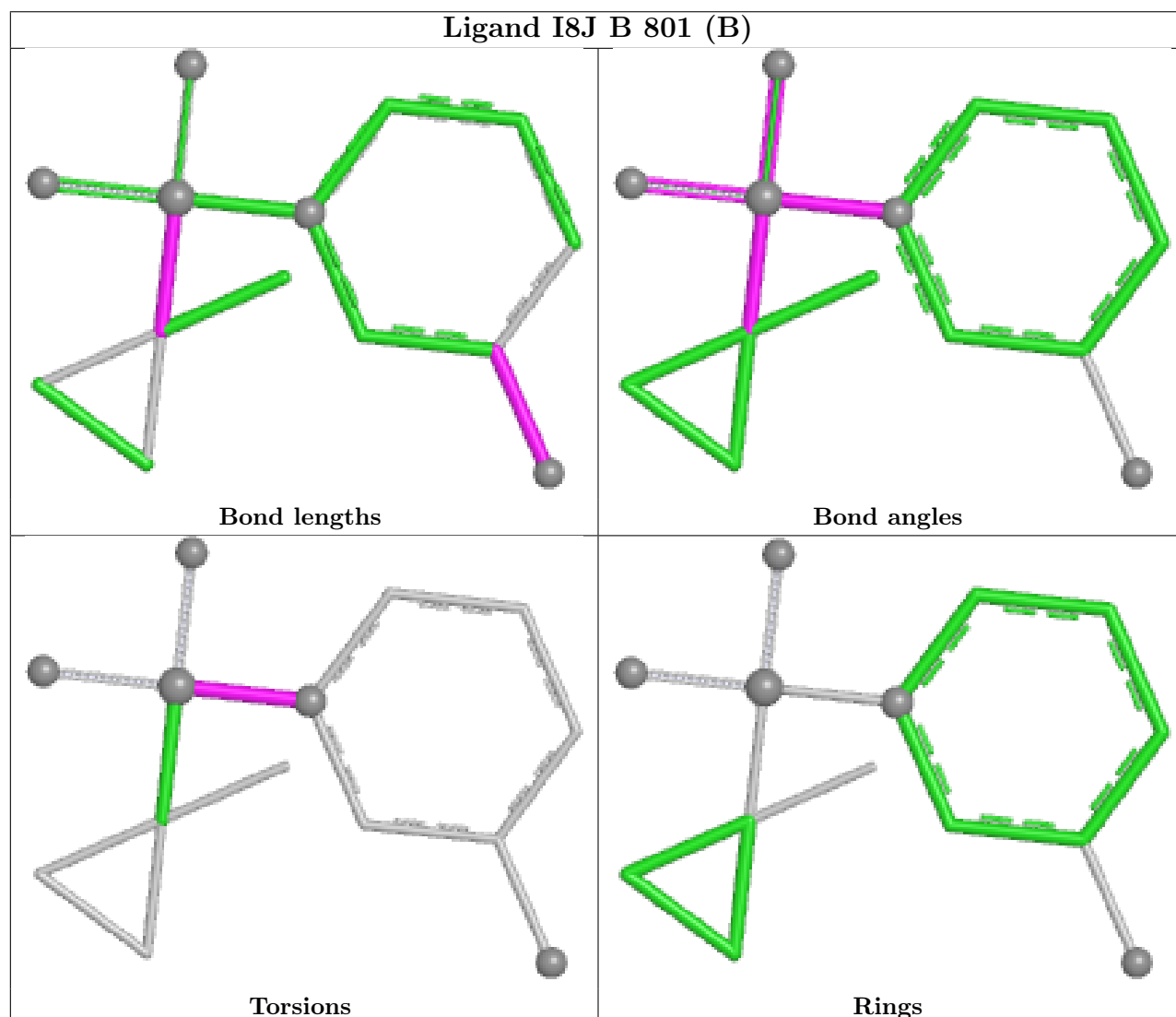
Mol	Chain	Res	Type	Atoms
2	A	801[B]	I8J	C5-N1-S1-O2
2	A	801[B]	I8J	C9-N1-S1-O1
2	A	801[B]	I8J	C4-C2-S1-O1
2	A	801[B]	I8J	C3-C2-S1-O2
2	B	801[B]	I8J	C5-N1-S1-C2
2	B	801[B]	I8J	C5-N1-S1-O2
2	B	801[B]	I8J	C9-N1-S1-C2
2	B	801[B]	I8J	C9-N1-S1-O1
2	B	801[B]	I8J	C9-N1-S1-O2
2	A	801[B]	I8J	C5-N1-S1-C2
2	A	801[B]	I8J	C5-N1-S1-O1

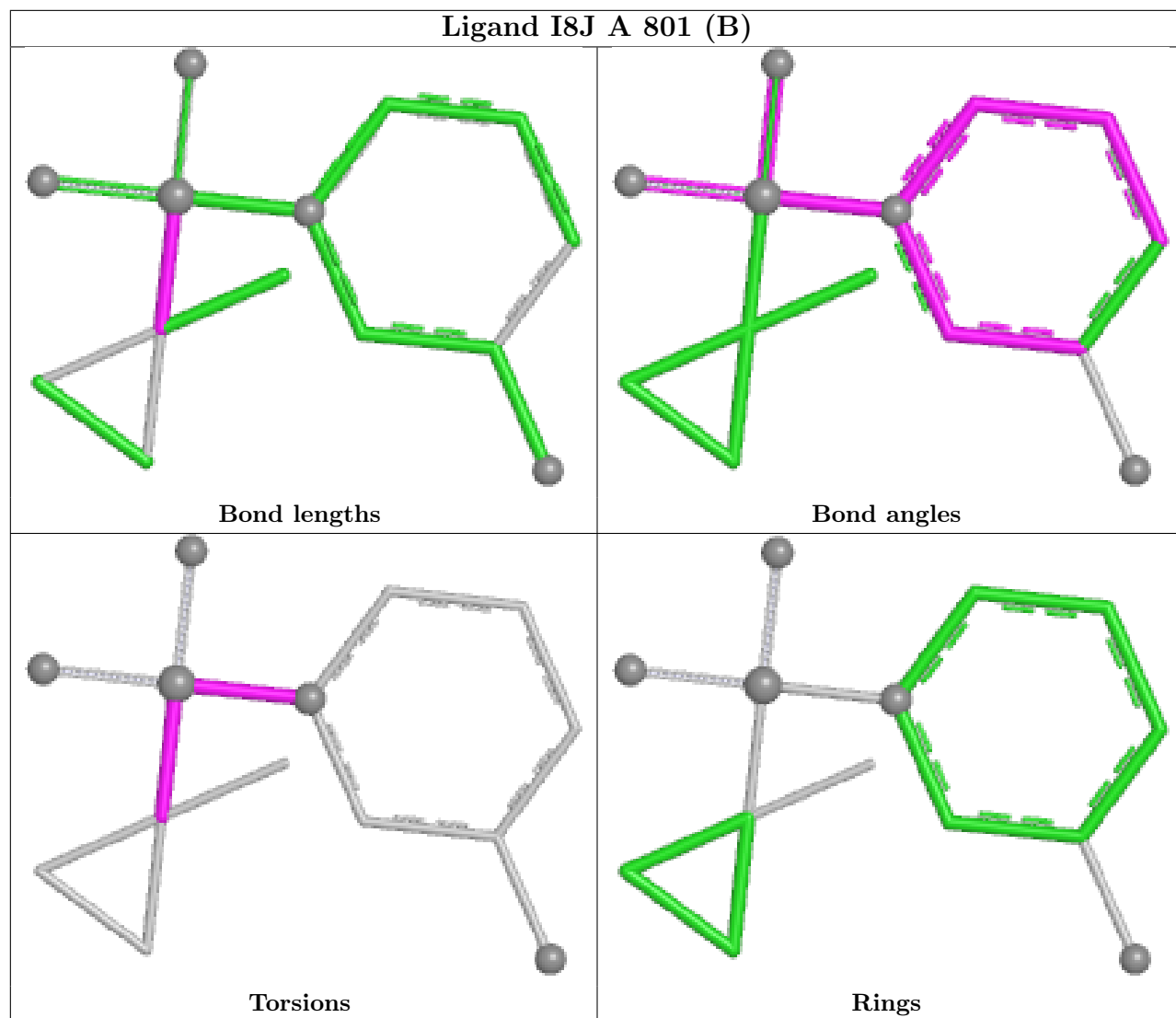
There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	801[B]	I8J	1	0
2	A	801[B]	I8J	1	0

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	697/707 (98%)	0.63	52 (7%)	20 17	19, 48, 69, 85	1 (0%)
1	B	695/707 (98%)	0.55	38 (5%)	30 27	22, 44, 66, 87	1 (0%)
All	All	1392/1414 (98%)	0.59	90 (6%)	25 22	19, 46, 67, 87	2 (0%)

All (90) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	23	GLU	6.4
1	A	142	ILE	6.2
1	B	422	GLN	6.0
1	B	426	ALA	5.8
1	A	426	ALA	5.8
1	A	425	ASP	5.6
1	A	22	ASP	5.5
1	B	720	PRO	5.4
1	A	422	GLN	5.4
1	B	566	ASP	5.3
1	A	566	ASP	5.2
1	A	427	SER	5.0
1	B	428	ALA	4.6
1	A	418	THR	4.6
1	B	173	PRO	4.5
1	B	167	ASN	4.3
1	B	568	ASP	4.2
1	A	167	ASN	4.2
1	B	420	ALA	4.1
1	B	475	LEU	4.1
1	A	169	SER	4.1
1	A	720	PRO	3.9
1	B	415	GLU	3.8
1	B	312	ARG	3.7

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Mol	Chain	Res	Type	RSRZ
1	B	166	LYS	3.6
1	A	613	TRP	3.6
1	A	158	LYS	3.6
1	B	613	TRP	3.5
1	A	428	ALA	3.5
1	A	33	ARG	3.5
1	A	312	ARG	3.5
1	B	419	LYS	3.4
1	B	567	GLN	3.4
1	A	417	GLU	3.4
1	B	418	THR	3.3
1	B	169	SER	3.2
1	B	97	LEU	3.1
1	B	565	GLY	3.1
1	A	575	LEU	3.0
1	B	661	ASN	3.0
1	B	99	HIS	3.0
1	A	429	ARG	2.9
1	B	378	HIS	2.9
1	B	142	ILE	2.8
1	A	416	ASP	2.7
1	A	168	PHE	2.7
1	B	427	SER	2.7
1	A	536	ARG	2.7
1	A	389	ASP	2.7
1	B	98	GLU	2.7
1	A	48	TYR	2.7
1	A	421	LEU	2.6
1	A	419	LYS	2.6
1	A	129	GLU	2.6
1	B	431	GLU	2.6
1	A	97	LEU	2.6
1	A	249	VAL	2.5
1	B	174	GLY	2.5
1	A	143	LYS	2.5
1	A	173	PRO	2.5
1	A	166	LYS	2.5
1	A	433	ILE	2.5
1	B	337[A]	ARG	2.4
1	A	109	MET	2.4
1	A	160	ALA	2.4
1	A	336[A]	LYS	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	415	GLU	2.4
1	A	431	GLU	2.4
1	B	170	ALA	2.3
1	B	421	LEU	2.3
1	B	627	ARG	2.3
1	A	567	GLN	2.3
1	A	636	ARG	2.3
1	B	41	ARG	2.3
1	A	103	GLU	2.3
1	A	435	LYS	2.3
1	A	703	ILE	2.2
1	B	246	GLU	2.2
1	B	168	PHE	2.2
1	A	628	LYS	2.2
1	A	366	ALA	2.2
1	A	430	LYS	2.2
1	B	685	LEU	2.2
1	A	127	LYS	2.2
1	B	611	ASP	2.2
1	A	213	GLY	2.1
1	B	632	ARG	2.1
1	A	420	ALA	2.1
1	A	141	ALA	2.1
1	A	170	ALA	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](#)

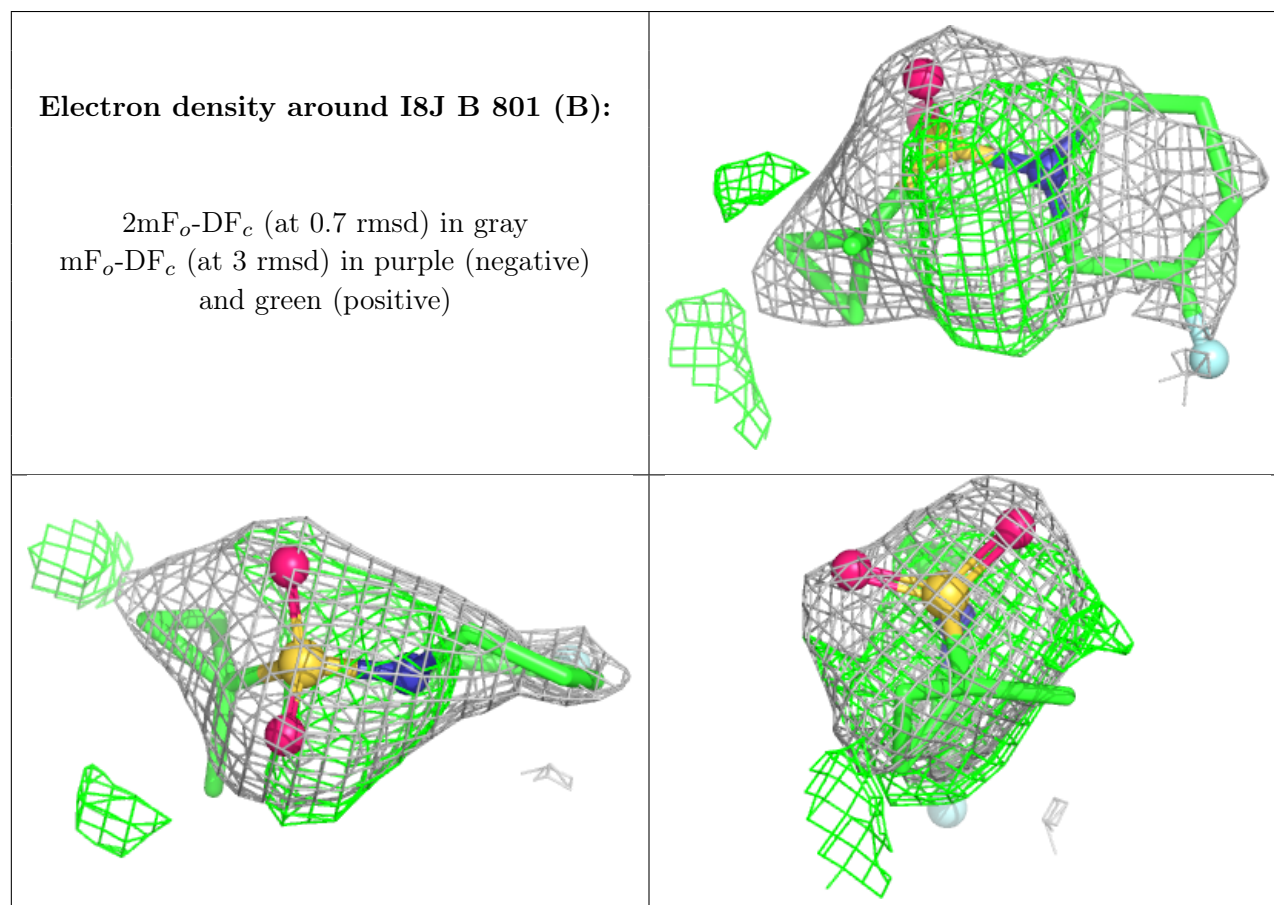
There are no oligosaccharides in this entry.

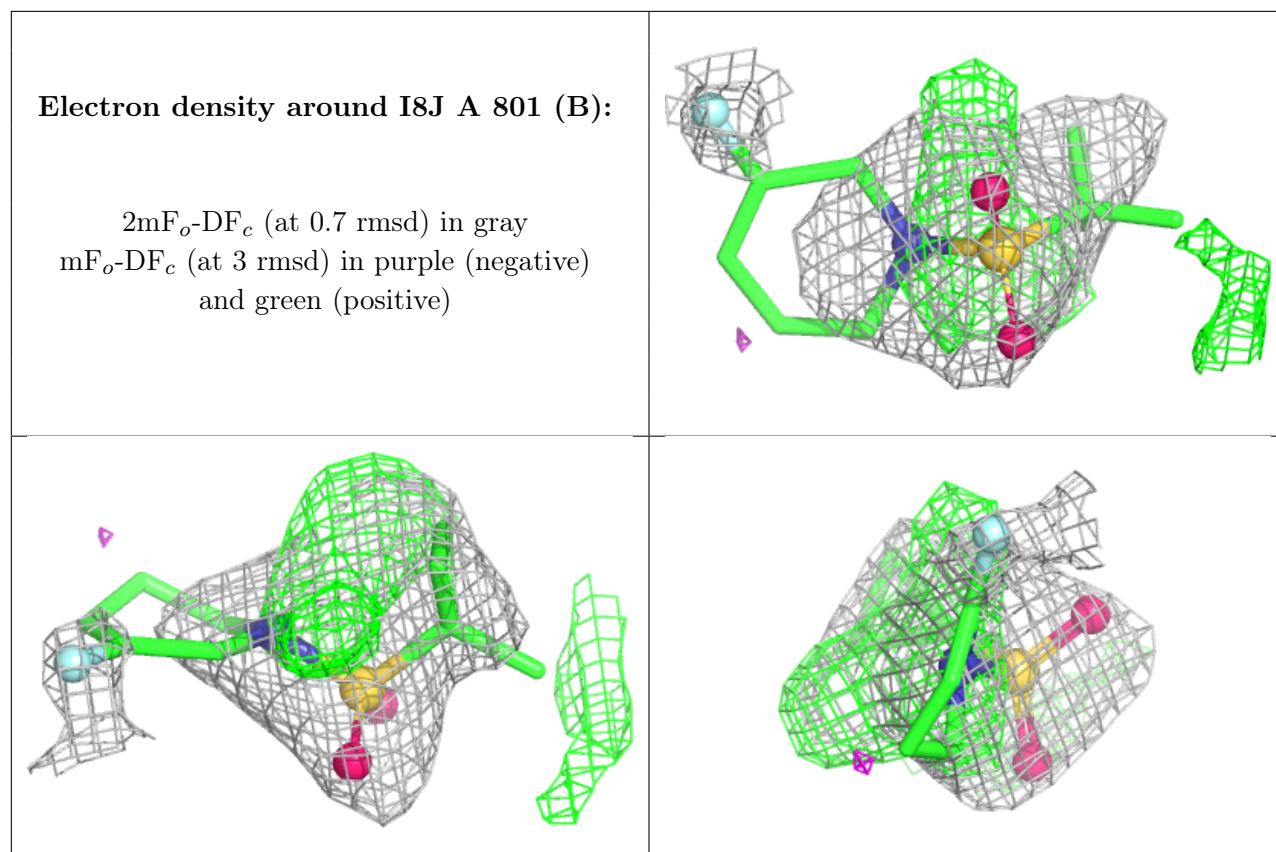
6.4 Ligands [i](#)

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
2	I8J	B	801[B]	14/14	0.79	0.34	41,42,44,44	14
2	I8J	A	801[B]	14/14	0.81	0.28	42,43,44,44	14
3	CL	A	802	1/1	0.82	0.22	82,82,82,82	0
3	CL	A	803	1/1	0.90	0.13	60,60,60,60	0
3	CL	B	802	1/1	0.93	0.24	70,70,70,70	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.