



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 03:15 AM UTC

PDB ID : 8AIY / pdb_00008aiy
Title : STRUCTURE OF THE LECB LECTIN FROM PSEUDOMONAS AERUGINOSA STRAIN PAO1 IN COMPLEX WITH N-(beta-L-Fucopyranosyl)-biphenyl-3-carboxamide (4i)
Authors : Meiers, J.; Mala, P.; Varrot, A.; Siebs, E.; Imberty, A.; Titz, A.
Deposited on : 2022-07-27
Resolution : 1.55 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:

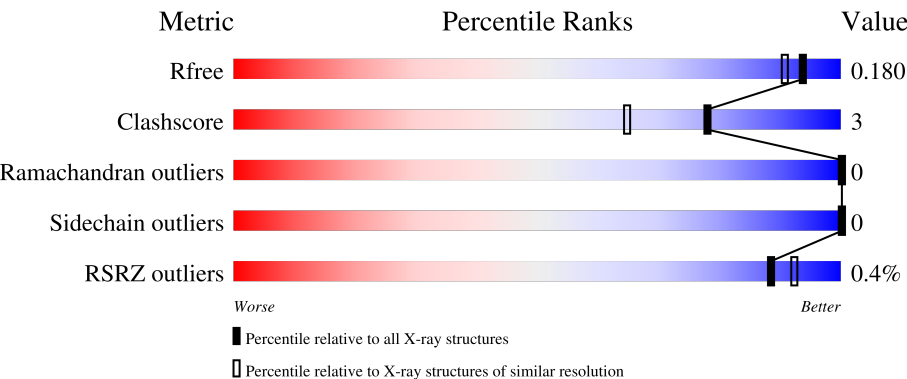
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal 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 i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.55 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	2145 (1.56-1.56)
Clashscore	190562	2189 (1.56-1.56)
Ramachandran outliers	187476	2153 (1.56-1.56)
Sidechain outliers	187428	2150 (1.56-1.56)
RSRZ outliers	180081	2146 (1.56-1.56)

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	AAA	114	93%7%
1	BBB	114	% 94%6%
1	CCC	114	% 96%.
1	DDD	114	90%9%.

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4094 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.

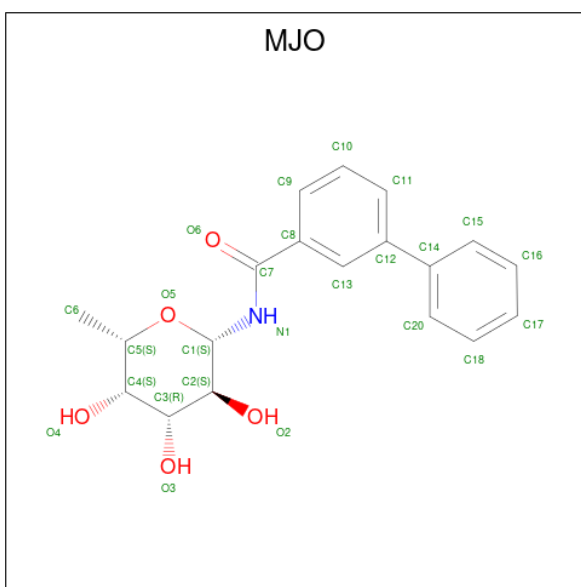
- Molecule 1 is a protein called Fucose-binding lectin PA-IIL.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	AAA	114	Total	C	N	O	0	1	0
			833	511	144	178			
1	BBB	114	Total	C	N	O	0	2	0
			833	513	145	175			
1	CCC	114	Total	C	N	O	0	1	0
			830	511	144	175			
1	DDD	114	Total	C	N	O	0	3	0
			843	517	146	180			

- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca).

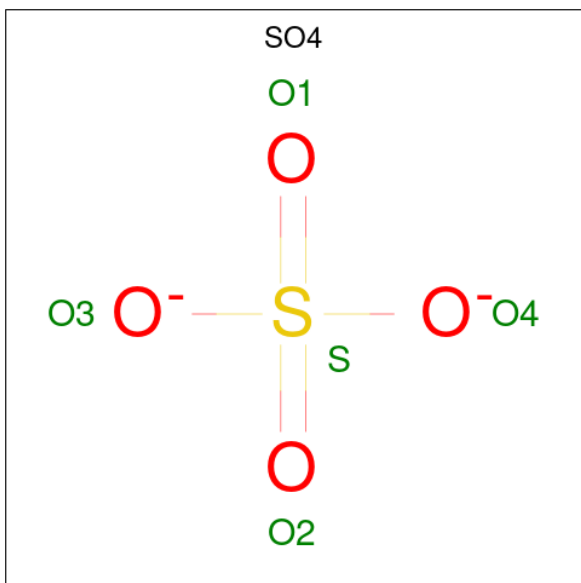
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	AAA	2	Total	Ca	0	0
			2	2		
2	BBB	2	Total	Ca	0	0
			2	2		
2	CCC	2	Total	Ca	0	0
			2	2		
2	DDD	2	Total	Ca	0	0
			2	2		

- Molecule 3 is N-(beta-L-Fucopyranosyl)-biphenyl-3-carboxamide (CCD ID: MJO) (formula: C₁₉H₂₁NO₅) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	AAA	1	Total 25	C 19	N 1	O 5	0	0
3	BBB	1	Total 25	C 19	N 1	O 5	0	0
3	CCC	1	Total 25	C 19	N 1	O 5	0	0
3	DDD	1	Total 25	C 19	N 1	O 5	0	0

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O_4S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	AAA	1	Total	O	S	0	0
			5	4	1		
4	CCC	1	Total	O	S	0	0
			5	4	1		

- Molecule 5 is water.

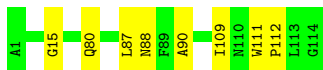
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	AAA	155	Total	O	0	3
			158	158		
5	BBB	150	Total	O	0	1
			151	151		
5	CCC	157	Total	O	0	4
			161	161		
5	DDD	165	Total	O	0	2
			167	167		

3 Residue-property plots [i](#)

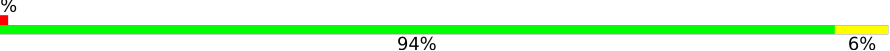
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Fucose-binding lectin PA-IIL

Chain AAA:  93% 7%

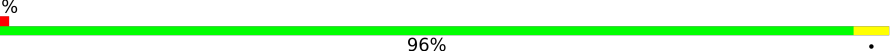


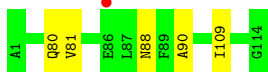
- Molecule 1: Fucose-binding lectin PA-IIL

Chain BBB:  % 94% 6%



- Molecule 1: Fucose-binding lectin PA-IIL

Chain CCC:  % 96% 4%



- Molecule 1: Fucose-binding lectin PA-IIL

Chain DDD:  90% 9% 1%



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	71.91Å 76.50Å 184.31Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.08 – 1.55 46.08 – 1.55	Depositor EDS
% Data completeness (in resolution range)	100.0 (46.08-1.55) 100.0 (46.08-1.55)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.02 (at 1.55Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.142 , 0.166 (Not available) , 0.180	Depositor DCC
R_{free} test set	3662 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	18.9	Xtriage
Anisotropy	0.046	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 41.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	4094	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.30% 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: CA, MJO, SO4

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	AAA	1.07	0/843	1.05	0/1154
1	BBB	1.04	1/843 (0.1%)	1.08	0/1155
1	CCC	0.99	0/840	1.08	0/1151
1	DDD	1.02	1/853 (0.1%)	1.05	1/1168 (0.1%)
All	All	1.03	2/3379 (0.1%)	1.06	1/4628 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	BBB	39	THR	C-O	6.41	1.31	1.23
1	DDD	15	GLY	CA-C	-5.02	1.48	1.52

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	DDD	101	ASP	CA-CB-CG	5.57	118.17	112.60

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	AAA	833	0	804	5	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	BBB	833	0	806	4	0
1	CCC	830	0	804	5	0
1	DDD	843	0	812	9	0
2	AAA	2	0	0	0	0
2	BBB	2	0	0	0	0
2	CCC	2	0	0	0	0
2	DDD	2	0	0	0	0
3	AAA	25	0	0	0	0
3	BBB	25	0	0	0	0
3	CCC	25	0	0	0	0
3	DDD	25	0	0	0	0
4	AAA	5	0	0	0	0
4	CCC	5	0	0	0	0
5	AAA	158	0	0	0	0
5	BBB	151	0	0	0	0
5	CCC	161	0	0	0	0
5	DDD	167	0	0	0	0
All	All	4094	0	3226	19	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 (19) 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:CCC:81[B]:VAL:HG21	1:DDD:81:VAL:HG21	1.30	1.05
1:CCC:81[B]:VAL:HG21	1:DDD:81:VAL:CG2	2.18	0.67
1:DDD:3:GLN:HE21	1:DDD:77:VAL:HA	1.69	0.57
1:CCC:80:GLN:HE21	1:CCC:88:ASN:HD22	1.53	0.56
1:CCC:81[B]:VAL:CG2	1:DDD:81:VAL:HG21	2.21	0.54
1:DDD:29:ASN:OD1	1:DDD:39[B]:THR:HG22	2.09	0.52
1:DDD:29:ASN:OD1	1:DDD:39[A]:THR:HG23	2.12	0.50
1:DDD:80:GLN:HE21	1:DDD:88:ASN:HD22	1.60	0.48
1:AAA:80:GLN:HE21	1:AAA:88:ASN:HD22	1.61	0.48
1:BBB:40:PHE:CD1	1:BBB:51:GLY:HA3	2.50	0.46
1:CCC:90:ALA:HB3	1:CCC:109:ILE:HB	1.97	0.46
1:AAA:90:ALA:HB3	1:AAA:109:ILE:HB	1.99	0.43
1:AAA:111:TRP:HB2	1:AAA:112:PRO:HA	2.00	0.42
1:BBB:90:ALA:HB3	1:BBB:109:ILE:HB	2.01	0.42
1:DDD:103:ASN:OD1	1:DDD:103:ASN:C	2.62	0.42
1:BBB:15:GLY:O	1:BBB:109:ILE:HA	2.20	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AAA:15:GLY:O	1:AAA:109:ILE:HA	2.20	0.41
1:AAA:87:LEU:HD22	1:BBB:93:GLY:HA3	2.02	0.40
1:DDD:15:GLY:O	1:DDD:109:ILE:HA	2.21	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	AAA	113/114 (99%)	111 (98%)	2 (2%)	0	100	100
1	BBB	114/114 (100%)	112 (98%)	2 (2%)	0	100	100
1	CCC	113/114 (99%)	110 (97%)	3 (3%)	0	100	100
1	DDD	115/114 (101%)	112 (97%)	3 (3%)	0	100	100
All	All	455/456 (100%)	445 (98%)	10 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	AAA	94/93 (101%)	94 (100%)	0	100	100
1	BBB	93/93 (100%)	93 (100%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	CCC	93/93 (100%)	93 (100%)	0	100	100
1	DDD	95/93 (102%)	95 (100%)	0	100	100
All	All	375/372 (101%)	375 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

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

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](#)

Of 14 ligands modelled in this entry, 8 are monoatomic - leaving 6 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	$\# Z > 2$
4	SO4	AAA	204	-	4,4,4	0.34	0	6,6,6	0.60	0
3	MJO	CCC	201	2	27,27,27	0.63	1 (3%)	38,38,38	1.16	2 (5%)
3	MJO	AAA	203	2	27,27,27	0.64	0	38,38,38	1.04	1 (2%)
3	MJO	BBB	203	2	27,27,27	0.61	0	38,38,38	0.76	1 (2%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	SO4	CCC	204	-	4,4,4	0.42	0	6,6,6	0.20	0
3	MJO	DDD	203	2	27,27,27	0.72	0	38,38,38	0.85	2 (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
3	MJO	DDD	203	2	-	0/12/32/32	0/3/3/3
3	MJO	CCC	201	2	-	0/12/32/32	0/3/3/3
3	MJO	AAA	203	2	-	6/12/32/32	0/3/3/3
3	MJO	BBB	203	2	-	0/12/32/32	0/3/3/3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	CCC	201	MJO	C1-N1	-2.06	1.40	1.43

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	CCC	201	MJO	O2-C2-C1	3.60	115.69	109.42
3	AAA	203	MJO	O2-C2-C1	3.49	115.51	109.42
3	BBB	203	MJO	O2-C2-C1	3.14	114.90	109.42
3	DDD	203	MJO	O2-C2-C1	2.74	114.20	109.42
3	DDD	203	MJO	C3-C2-C1	-2.17	106.68	109.86
3	CCC	201	MJO	O5-C1-C2	-2.06	107.41	109.72

There are no chirality outliers.

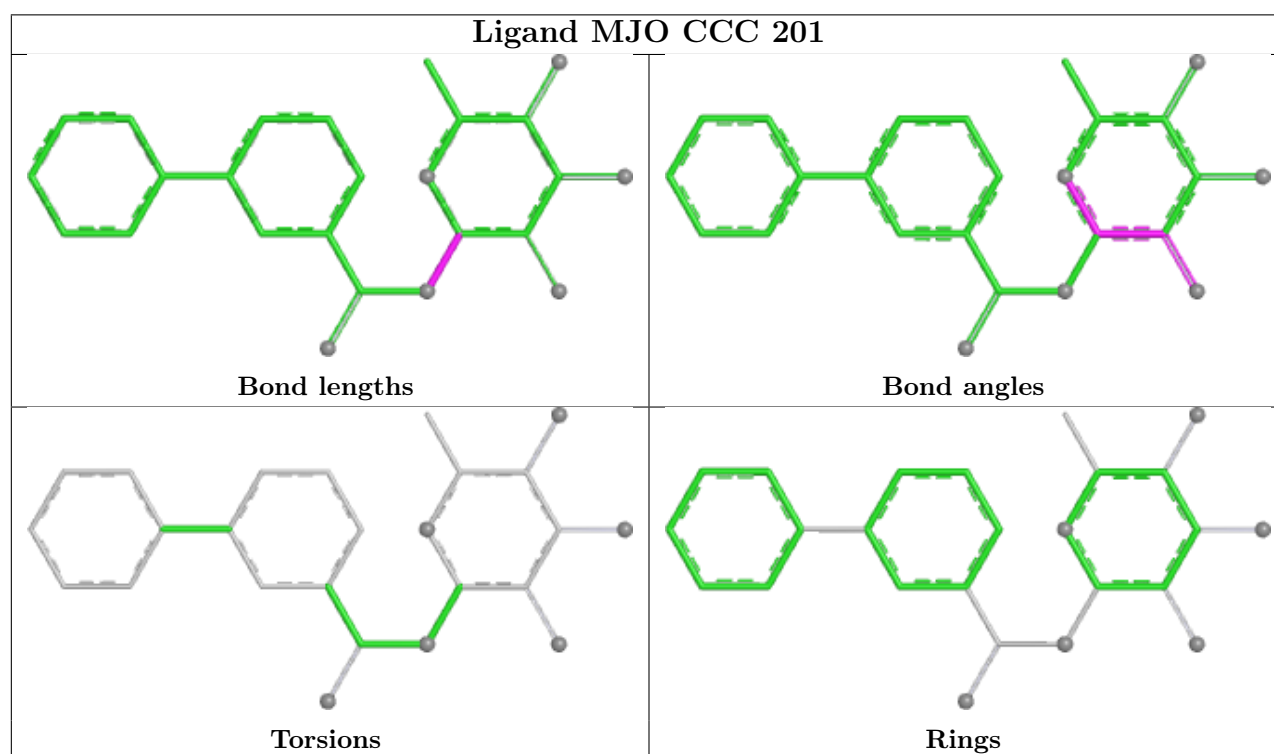
All (6) torsion outliers are listed below:

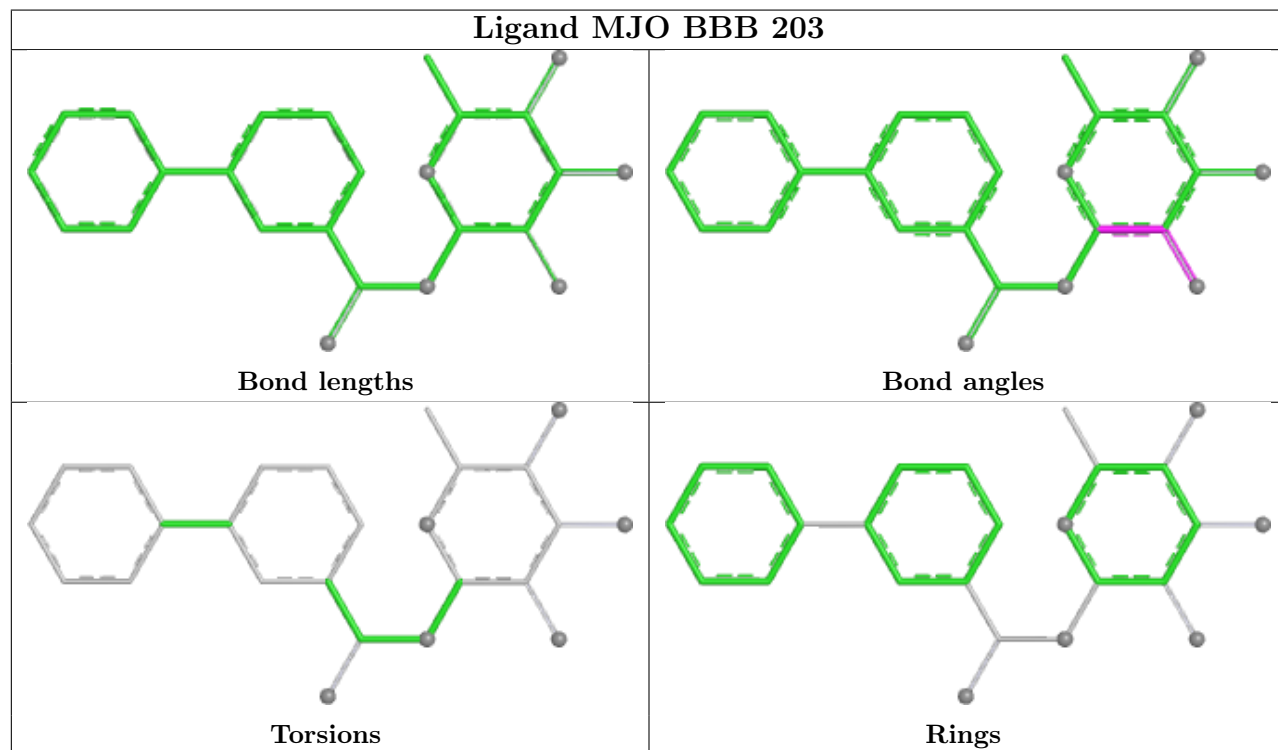
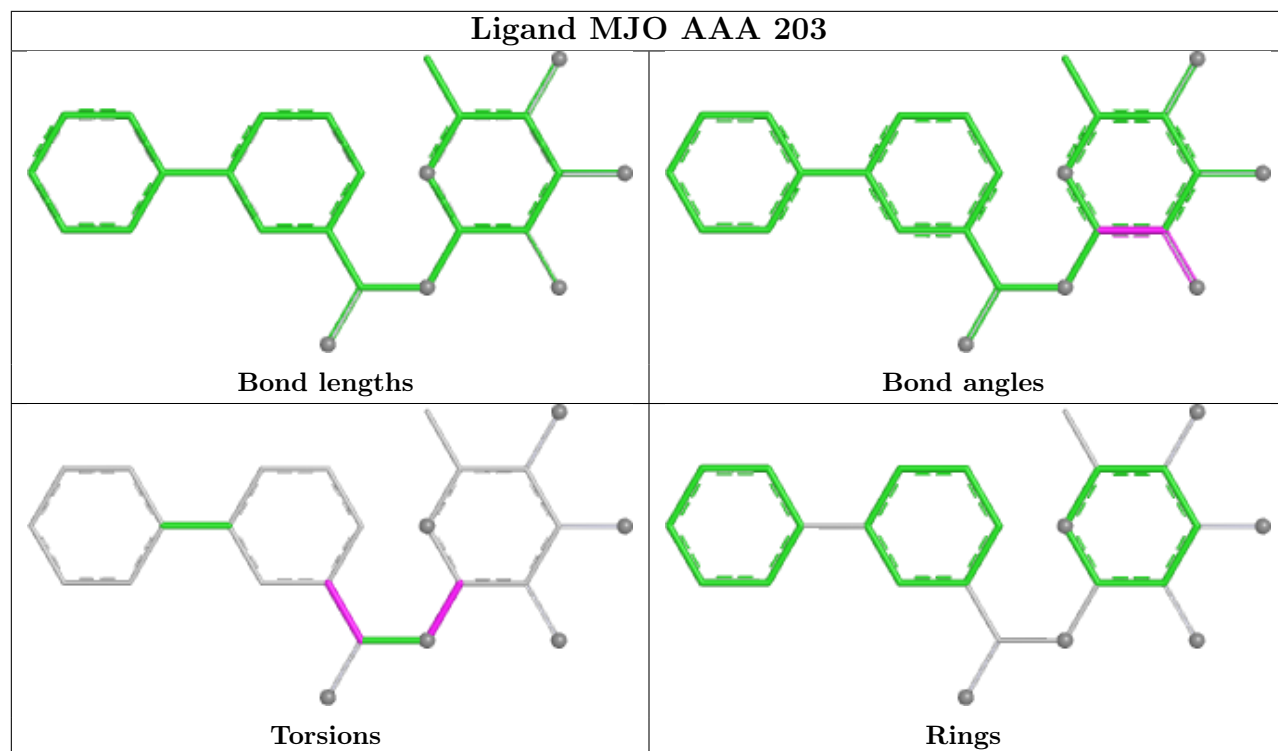
Mol	Chain	Res	Type	Atoms
3	AAA	203	MJO	O5-C1-N1-C7
3	AAA	203	MJO	C2-C1-N1-C7
3	AAA	203	MJO	N1-C7-C8-C13
3	AAA	203	MJO	O6-C7-C8-C13
3	AAA	203	MJO	N1-C7-C8-C9
3	AAA	203	MJO	O6-C7-C8-C9

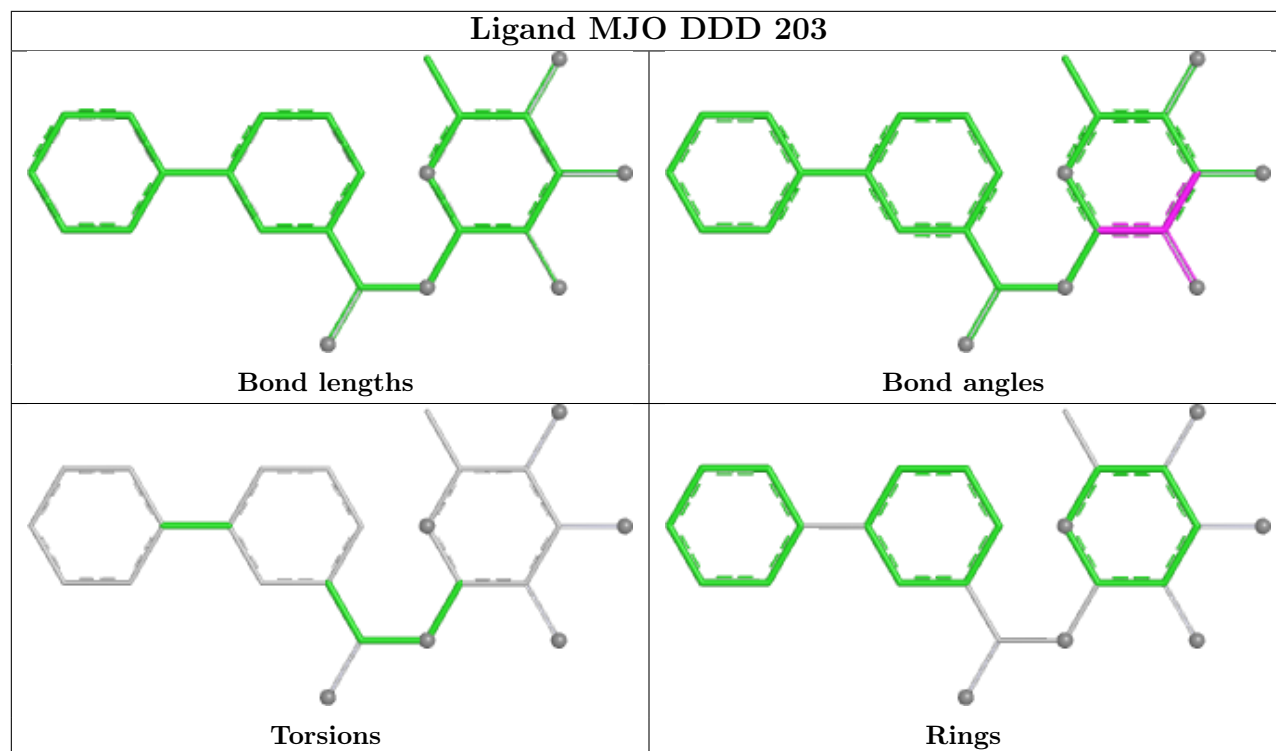
There are no ring outliers.

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	AAA	114/114 (100%)	-0.06	0 100 100	11, 20, 28, 37	1 (0%)
1	BBB	114/114 (100%)	-0.21	1 (0%) 81 86	8, 18, 28, 38	2 (1%)
1	CCC	114/114 (100%)	-0.29	1 (0%) 81 86	9, 18, 25, 29	1 (0%)
1	DDD	114/114 (100%)	-0.28	0 100 100	10, 18, 26, 33	3 (2%)
All	All	456/456 (100%)	-0.21	2 (0%) 88 92	8, 18, 27, 38	7 (1%)

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	BBB	33	ASN	2.2
1	CCC	86	GLU	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](#)

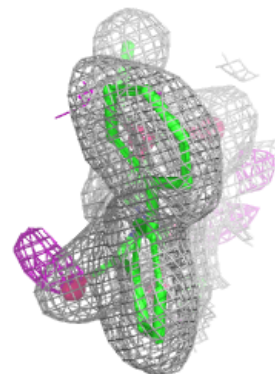
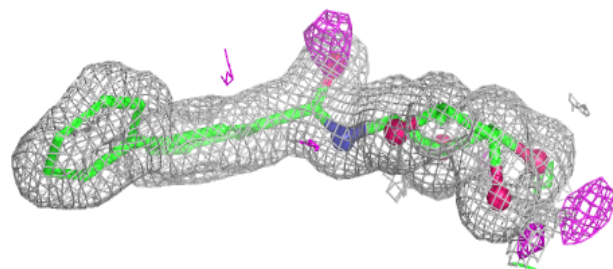
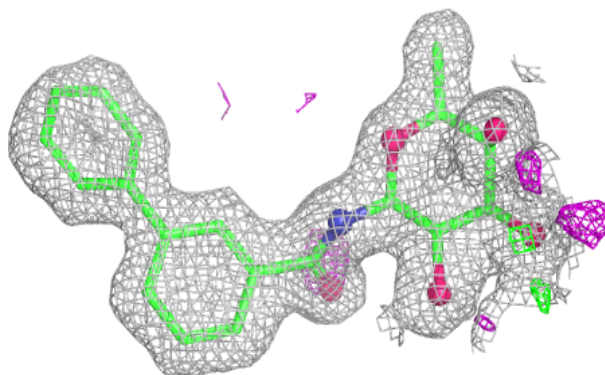
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists 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
4	SO4	AAA	204	5/5	0.84	0.12	20,25,30,32	5
3	MJO	AAA	203	25/25	0.96	0.07	18,20,23,24	0
3	MJO	DDD	203	25/25	0.97	0.06	16,18,26,27	0
3	MJO	BBB	203	25/25	0.97	0.05	16,17,20,21	0
2	CA	AAA	201	1/1	0.98	0.05	18,18,18,18	0
3	MJO	CCC	201	25/25	0.98	0.05	16,18,21,22	0
2	CA	DDD	202	1/1	0.99	0.02	17,17,17,17	0
2	CA	AAA	202	1/1	0.99	0.03	19,19,19,19	0
4	SO4	CCC	204	5/5	0.99	0.04	20,20,21,22	0
2	CA	CCC	202	1/1	1.00	0.02	17,17,17,17	0
2	CA	CCC	203	1/1	1.00	0.02	16,16,16,16	0
2	CA	DDD	201	1/1	1.00	0.02	17,17,17,17	0
2	CA	BBB	201	1/1	1.00	0.01	17,17,17,17	0
2	CA	BBB	202	1/1	1.00	0.02	17,17,17,17	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.

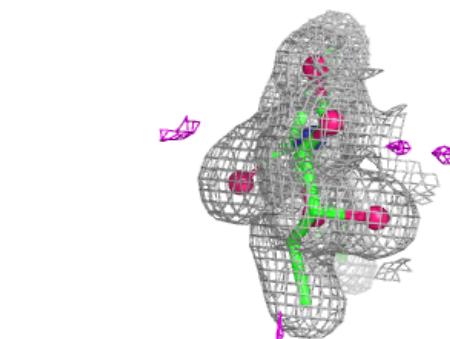
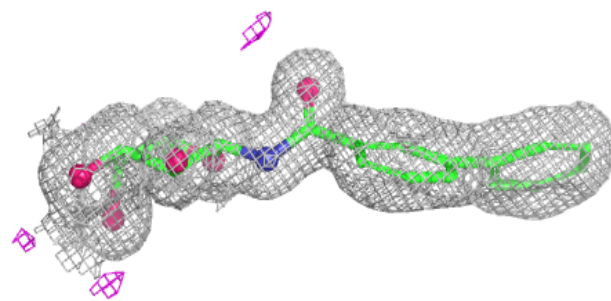
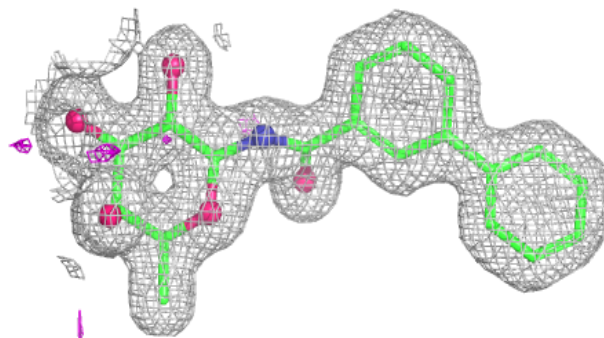
Electron density around MJO AAA 203:

2mF_o-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

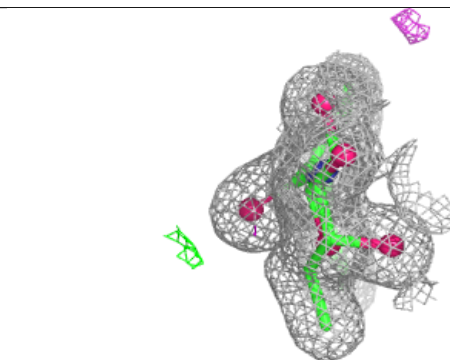
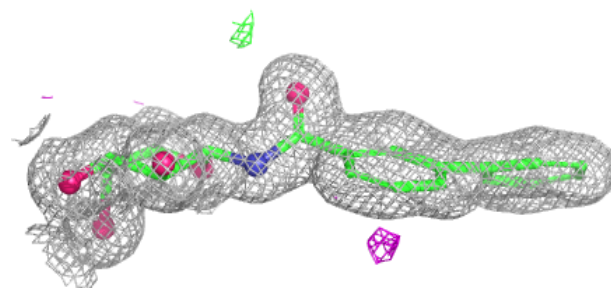
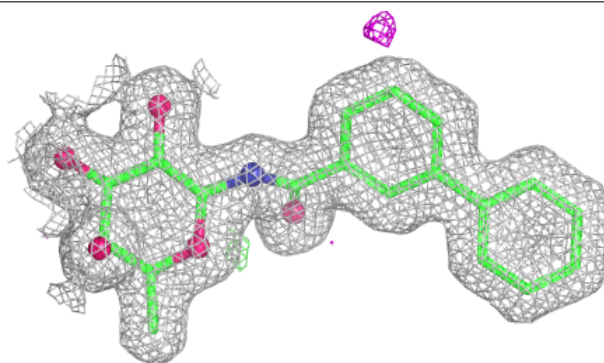


Electron density around MJO DDD 203:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

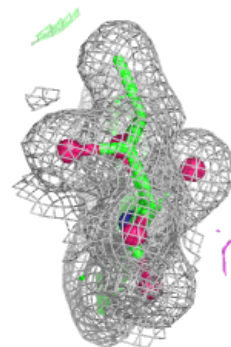
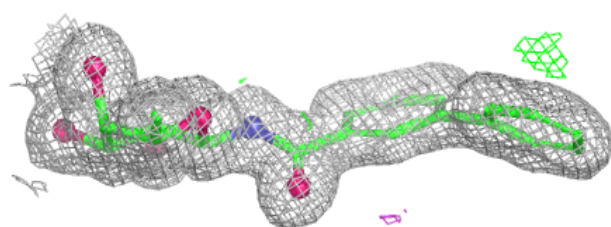
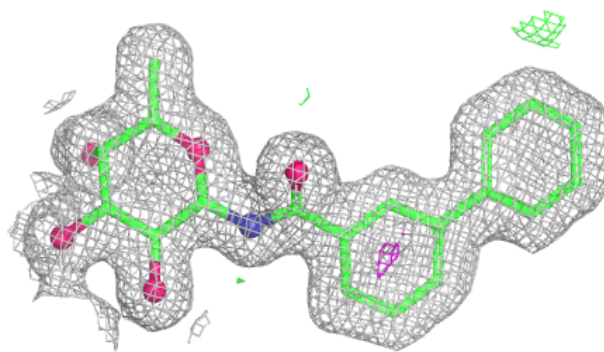
**Electron density around MJO BBB 203:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around MJO CCC 201:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.