



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 10:49 PM UTC

PDB ID : 5L9O / pdb_00005l9o
Title : Crystal structure of Agrobacterium tumefaciens C58 strain PBP SocA in complex with glucopine
Authors : Marty, L.; Morera, S.
Deposited on : 2016-06-10
Resolution : 1.84 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

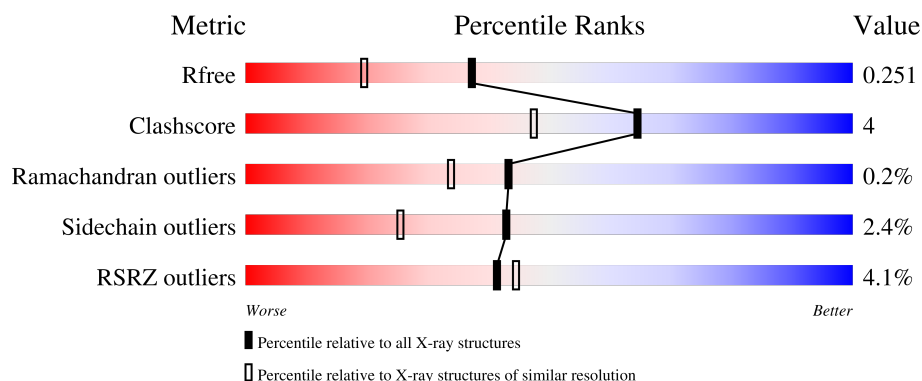
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.84 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	1296 (1.84-1.84)
Clashscore	190562	1329 (1.84-1.84)
Ramachandran outliers	187476	1318 (1.84-1.84)
Sidechain outliers	187428	1318 (1.84-1.84)
RSRZ outliers	180081	1296 (1.84-1.84)

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	268	 82% 8% 10%
1	B	268	 79% 10% 9%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 3944 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 Deoxyfructosyl-amino Acid Transporter Periplasmic Binding Protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	241	Total	C	N	O	S	0	0	0
			1817	1163	303	347	4			
1	B	243	Total	C	N	O	S	0	0	0
			1830	1171	305	349	5			

There are 52 discrepancies between the modelled and reference sequences:

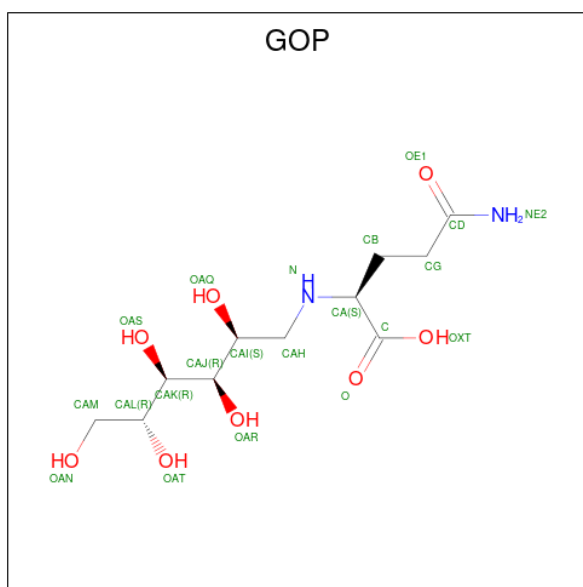
Chain	Residue	Modelled	Actual	Comment	Reference
A	16	GLY	-	expression tag	UNP Q7D447
A	17	SER	-	expression tag	UNP Q7D447
A	18	SER	-	expression tag	UNP Q7D447
A	19	HIS	-	expression tag	UNP Q7D447
A	20	HIS	-	expression tag	UNP Q7D447
A	21	HIS	-	expression tag	UNP Q7D447
A	22	HIS	-	expression tag	UNP Q7D447
A	23	HIS	-	expression tag	UNP Q7D447
A	24	HIS	-	expression tag	UNP Q7D447
A	25	SER	-	expression tag	UNP Q7D447
A	26	SER	-	expression tag	UNP Q7D447
A	27	GLY	-	expression tag	UNP Q7D447
A	28	LEU	-	expression tag	UNP Q7D447
A	29	VAL	-	expression tag	UNP Q7D447
A	30	PRO	-	expression tag	UNP Q7D447
A	31	ARG	-	expression tag	UNP Q7D447
A	32	GLY	-	expression tag	UNP Q7D447
A	33	SER	-	expression tag	UNP Q7D447
A	34	HIS	-	expression tag	UNP Q7D447
A	35	MET	-	expression tag	UNP Q7D447
A	278	HIS	-	expression tag	UNP Q7D447
A	279	HIS	-	expression tag	UNP Q7D447
A	280	HIS	-	expression tag	UNP Q7D447
A	281	HIS	-	expression tag	UNP Q7D447

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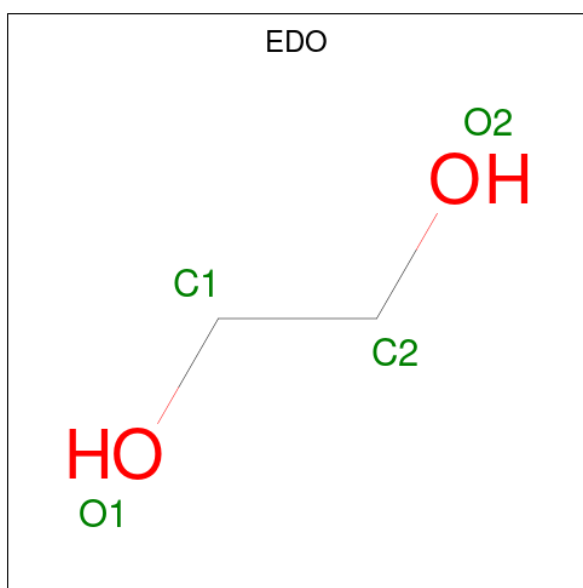
Chain	Residue	Modelled	Actual	Comment	Reference
A	282	HIS	-	expression tag	UNP Q7D447
A	283	HIS	-	expression tag	UNP Q7D447
B	16	GLY	-	expression tag	UNP Q7D447
B	17	SER	-	expression tag	UNP Q7D447
B	18	SER	-	expression tag	UNP Q7D447
B	19	HIS	-	expression tag	UNP Q7D447
B	20	HIS	-	expression tag	UNP Q7D447
B	21	HIS	-	expression tag	UNP Q7D447
B	22	HIS	-	expression tag	UNP Q7D447
B	23	HIS	-	expression tag	UNP Q7D447
B	24	HIS	-	expression tag	UNP Q7D447
B	25	SER	-	expression tag	UNP Q7D447
B	26	SER	-	expression tag	UNP Q7D447
B	27	GLY	-	expression tag	UNP Q7D447
B	28	LEU	-	expression tag	UNP Q7D447
B	29	VAL	-	expression tag	UNP Q7D447
B	30	PRO	-	expression tag	UNP Q7D447
B	31	ARG	-	expression tag	UNP Q7D447
B	32	GLY	-	expression tag	UNP Q7D447
B	33	SER	-	expression tag	UNP Q7D447
B	34	HIS	-	expression tag	UNP Q7D447
B	35	MET	-	expression tag	UNP Q7D447
B	278	HIS	-	expression tag	UNP Q7D447
B	279	HIS	-	expression tag	UNP Q7D447
B	280	HIS	-	expression tag	UNP Q7D447
B	281	HIS	-	expression tag	UNP Q7D447
B	282	HIS	-	expression tag	UNP Q7D447
B	283	HIS	-	expression tag	UNP Q7D447

- Molecule 2 is Glucopine (CCD ID: GOP) (formula: C₁₁H₂₂N₂O₈).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total 21	C 11	N 2	O 8	0	0
2	B	1	Total 21	C 11	N 2	O 8	0	0

- Molecule 3 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $\text{C}_2\text{H}_6\text{O}_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 4 2 2	0	0
3	A	1	Total C O 4 2 2	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 4 2 2	0	0
3	A	1	Total C O 4 2 2	0	0
3	A	1	Total C O 4 2 2	0	0
3	A	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0

- Molecule 4 is CALCIUM ION (CCD ID: CA) (formula: Ca).

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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	126	Total O 126 126	0	0
5	B	87	Total O 87 87	0	0

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.

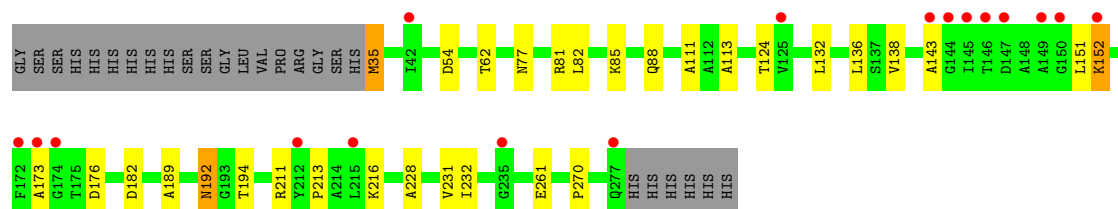
• Molecule 1: Deoxyfructosyl-amino Acid Transporter Periplasmic Binding Protein

Chain A: 



• Molecule 1: Deoxyfructosyl-amino Acid Transporter Periplasmic Binding Protein

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	39.60Å 62.29Å 84.23Å 90.00° 101.40° 90.00°	Depositor
Resolution (Å)	33.83 – 1.84 33.83 – 1.84	Depositor EDS
% Data completeness (in resolution range)	96.5 (33.83-1.84) 96.4 (33.83-1.84)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.98 (at 1.84Å)	Xtriage
Refinement program	BUSTER 2.10.2	Depositor
R, R_{free}	0.218 , 0.252 0.221 , 0.251	Depositor DCC
R_{free} test set	1688 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	28.6	Xtriage
Anisotropy	0.622	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 49.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.000 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3944	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 77.00 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 9.3045e-07. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: GOP, CA, EDO

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	0.77	0/1857	1.23	5/2517 (0.2%)
1	B	0.74	0/1870	1.21	3/2534 (0.1%)
All	All	0.76	0/3727	1.22	8/5051 (0.2%)

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	136	LEU	N-CA-C	-6.00	99.92	109.76
1	B	136	LEU	N-CA-C	-6.00	99.93	109.76
1	A	144	GLY	CA-C-N	5.79	128.39	120.35
1	A	144	GLY	C-N-CA	5.79	128.39	120.35
1	A	172	PHE	CA-CB-CG	5.59	119.39	113.80
1	B	192	ASN	CA-CB-CG	5.57	118.17	112.60
1	B	176	ASP	CA-CB-CG	5.09	117.69	112.60
1	A	99	LEU	N-CA-C	5.06	116.49	110.97

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	A	1817	0	1795	13	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1830	0	1806	18	0
2	A	21	0	0	1	0
2	B	21	0	0	1	0
3	A	24	0	36	3	0
3	B	16	0	24	5	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	126	0	0	0	0
5	B	87	0	0	0	0
All	All	3944	0	3661	29	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 4.

All (29) 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:270:PRO:HD3	3:B:304:EDO:H21	1.53	0.91
1:A:132:LEU:HD12	3:A:306:EDO:H12	1.64	0.78
1:B:143:ALA:HB2	1:B:216:LYS:HE2	1.68	0.74
1:A:77:ASN:HD21	1:A:81:ARG:HH21	1.45	0.65
1:B:77:ASN:HD21	1:B:81:ARG:HH21	1.45	0.64
1:B:35:MET:N	3:B:305:EDO:HO2	1.98	0.60
1:B:211:ARG:O	1:B:213:PRO:HD3	2.02	0.59
1:A:77:ASN:ND2	1:A:81:ARG:HE	2.04	0.56
1:B:132:LEU:HD12	3:B:304:EDO:H11	1.87	0.56
1:B:77:ASN:ND2	1:B:81:ARG:HE	2.05	0.54
1:A:270:PRO:HD3	3:A:306:EDO:H22	1.89	0.53
1:B:261:GLU:HA	3:B:302:EDO:H21	1.92	0.52
1:A:228:ALA:HB2	2:A:301:GOP:OAN	2.10	0.52
1:B:270:PRO:CD	3:B:304:EDO:H21	2.35	0.52
1:A:270:PRO:HD3	3:A:306:EDO:C2	2.40	0.51
1:A:189:ALA:HB1	1:A:194:THR:HB	1.93	0.50
1:A:258:LYS:HG2	1:B:152:LYS:HE2	1.93	0.50
1:B:228:ALA:HB2	2:B:301:GOP:OAN	2.13	0.49
1:B:85:LYS:H	1:B:88:GLN:HE21	1.62	0.48
1:B:111:ALA:HB3	1:B:231:VAL:HB	1.97	0.47
1:A:85:LYS:H	1:A:88:GLN:HE21	1.63	0.46
1:A:255:THR:HG23	1:B:173:ALA:HB2	1.96	0.46
1:A:85:LYS:H	1:A:88:GLN:NE2	2.15	0.45
1:B:85:LYS:H	1:B:88:GLN:NE2	2.14	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:54:ASP:OD2	1:B:182:ASP:HB2	2.18	0.44
1:B:138:VAL:HG11	1:B:151:LEU:HD21	1.98	0.44
1:A:54:ASP:OD2	1:A:182:ASP:HB2	2.18	0.43
1:A:167:TYR:CZ	1:A:221:ILE:HG21	2.54	0.43
1:B:189:ALA:HB1	1:B:194:THR:HB	2.01	0.42

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	239/268 (89%)	235 (98%)	4 (2%)	0	100	100
1	B	241/268 (90%)	234 (97%)	6 (2%)	1 (0%)	30	18
All	All	480/536 (90%)	469 (98%)	10 (2%)	1 (0%)	43	34

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	113	ALA

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.

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	185/209 (88%)	183 (99%)	2 (1%)	65	54
1	B	186/209 (89%)	179 (96%)	7 (4%)	29	12
All	All	371/418 (89%)	362 (98%)	9 (2%)	43	25

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	145	ILE
1	A	232	ILE
1	B	35	MET
1	B	62	THR
1	B	82	LEU
1	B	124	THR
1	B	152	LYS
1	B	192	ASN
1	B	232	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	77	ASN
1	A	88	GLN
1	A	94	GLN
1	A	171	ASN
1	A	183	ASN
1	A	191	ASN
1	A	220	ASN
1	B	77	ASN
1	B	88	GLN
1	B	160	GLN
1	B	171	ASN
1	B	183	ASN
1	B	184	ASN
1	B	191	ASN
1	B	192	ASN
1	B	220	ASN

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 14 ligands modelled in this entry, 2 are monoatomic - leaving 12 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	$\# Z > 2$
3	EDO	B	305	-	3,3,3	0.52	0	2,2,2	0.31	0
3	EDO	B	304	-	3,3,3	0.44	0	2,2,2	0.45	0
3	EDO	A	302	-	3,3,3	0.52	0	2,2,2	0.31	0
2	GOP	A	301	-	20,20,20	0.76	0	24,26,26	1.97	4 (16%)
3	EDO	A	303	-	3,3,3	0.50	0	2,2,2	0.39	0
3	EDO	A	307	-	3,3,3	0.56	0	2,2,2	0.27	0
3	EDO	A	306	-	3,3,3	0.52	0	2,2,2	0.28	0
2	GOP	B	301	-	20,20,20	0.72	0	24,26,26	1.81	3 (12%)
3	EDO	B	302	-	3,3,3	0.41	0	2,2,2	0.51	0
3	EDO	B	303	-	3,3,3	0.54	0	2,2,2	0.32	0
3	EDO	A	305	-	3,3,3	0.54	0	2,2,2	0.34	0
3	EDO	A	304	-	3,3,3	0.51	0	2,2,2	0.39	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
3	EDO	B	305	-	-	0/1/1/1	-
3	EDO	B	304	-	-	1/1/1/1	-
3	EDO	A	302	-	-	0/1/1/1	-
2	GOP	A	301	-	-	2/28/28/28	-
3	EDO	A	303	-	-	0/1/1/1	-
3	EDO	A	307	-	-	0/1/1/1	-
3	EDO	A	306	-	-	1/1/1/1	-
2	GOP	B	301	-	-	2/28/28/28	-
3	EDO	B	302	-	-	1/1/1/1	-
3	EDO	B	303	-	-	0/1/1/1	-
3	EDO	A	305	-	-	0/1/1/1	-
3	EDO	A	304	-	-	1/1/1/1	-

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	301	GOP	CAL-CAK-CAJ	6.86	123.01	112.48
2	B	301	GOP	CAL-CAK-CAJ	6.45	122.39	112.48
2	A	301	GOP	CAM-CAL-CAK	3.39	119.09	112.17
2	B	301	GOP	CAM-CAL-CAK	3.32	118.94	112.17
2	A	301	GOP	OAN-CAM-CAL	2.99	117.44	111.16
2	B	301	GOP	O-C-CA	-2.72	113.49	122.26
2	A	301	GOP	O-C-CA	-2.66	113.67	122.26

There are no chirality outliers.

All (8) torsion outliers are listed below:

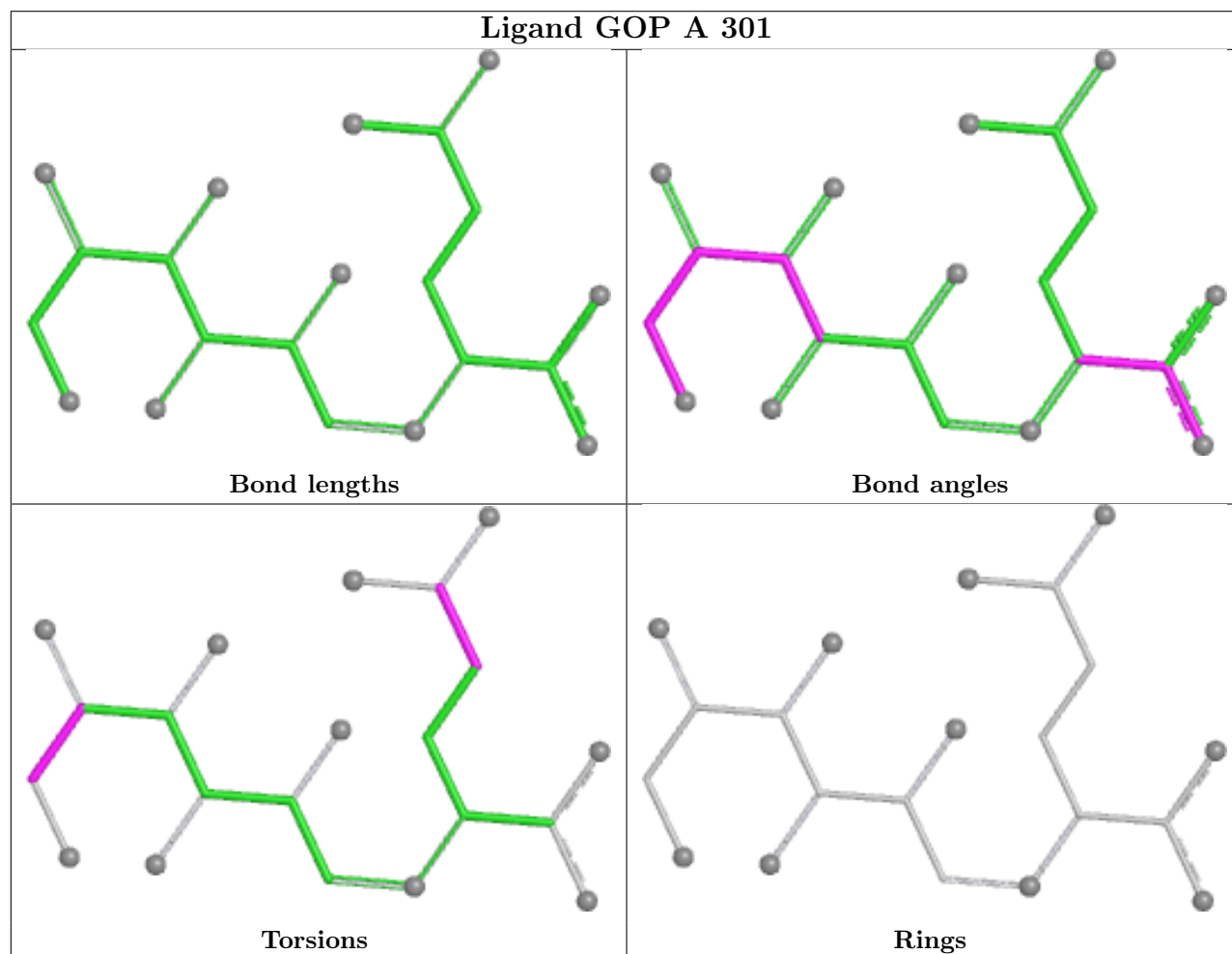
Mol	Chain	Res	Type	Atoms
2	A	301	GOP	CAK-CAL-CAM-OAN
3	B	302	EDO	O1-C1-C2-O2
3	A	304	EDO	O1-C1-C2-O2
2	B	301	GOP	NE2-CD-CG-CB
3	B	304	EDO	O1-C1-C2-O2
2	B	301	GOP	OE1-CD-CG-CB
3	A	306	EDO	O1-C1-C2-O2
2	A	301	GOP	NE2-CD-CG-CB

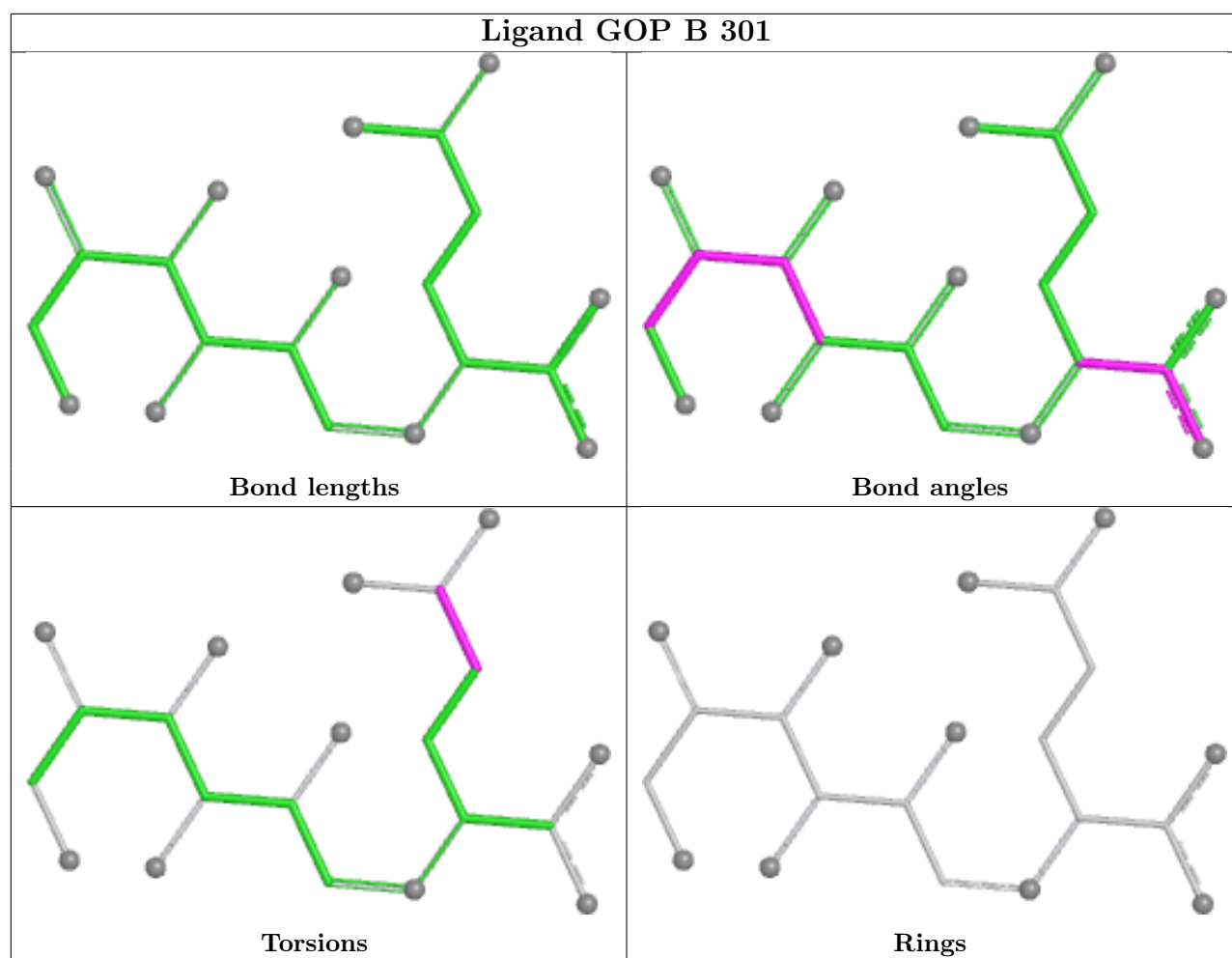
There are no ring outliers.

6 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	305	EDO	1	0
3	B	304	EDO	3	0
2	A	301	GOP	1	0
3	A	306	EDO	3	0
2	B	301	GOP	1	0
3	B	302	EDO	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 [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	241/268 (89%)	0.25	3 (1%) 76 82	20, 36, 58, 77	0
1	B	243/268 (90%)	0.64	17 (6%) 22 23	23, 45, 77, 109	0
All	All	484/536 (90%)	0.45	20 (4%) 41 44	20, 40, 71, 109	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	143	ALA	4.4
1	B	144	GLY	4.1
1	B	145	ILE	3.5
1	B	277	GLN	3.3
1	B	173	ALA	3.1
1	A	145	ILE	2.8
1	B	152	LYS	2.8
1	B	146	THR	2.8
1	B	212	TYR	2.7
1	B	174	GLY	2.6
1	B	149	ALA	2.5
1	B	235	GLY	2.4
1	B	147	ASP	2.4
1	B	172	PHE	2.4
1	A	170	LYS	2.3
1	B	42	ILE	2.2
1	B	150	GLY	2.1
1	A	235	GLY	2.1
1	B	125	VAL	2.1
1	B	215	LEU	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](#)

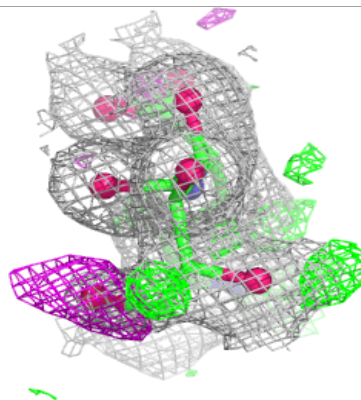
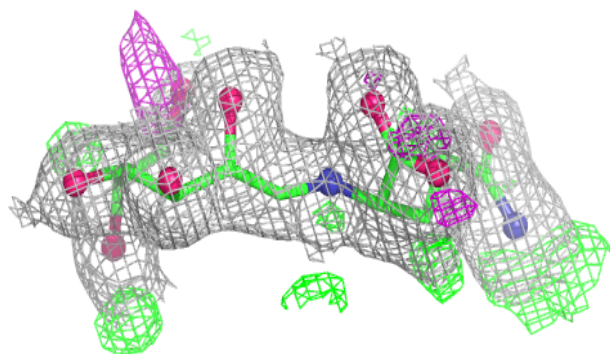
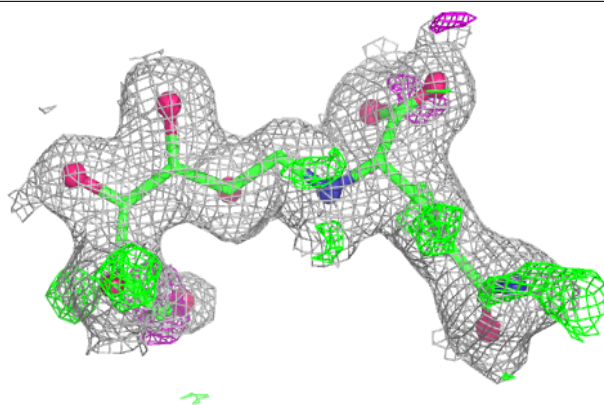
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
3	EDO	A	303	4/4	0.51	0.15	70,71,72,72	0
3	EDO	A	307	4/4	0.69	0.16	67,67,68,68	0
3	EDO	B	305	4/4	0.69	0.13	67,67,69,70	0
3	EDO	A	302	4/4	0.72	0.15	66,67,67,68	0
3	EDO	B	303	4/4	0.74	0.13	65,66,66,66	0
3	EDO	A	305	4/4	0.75	0.14	54,55,56,57	0
4	CA	A	308	1/1	0.84	0.12	99,99,99,99	0
3	EDO	A	304	4/4	0.85	0.14	49,51,51,52	0
3	EDO	B	302	4/4	0.87	0.24	36,38,39,41	0
2	GOP	A	301	21/21	0.88	0.09	18,23,30,31	0
3	EDO	A	306	4/4	0.91	0.15	25,29,31,31	0
3	EDO	B	304	4/4	0.91	0.17	32,37,37,37	0
2	GOP	B	301	21/21	0.92	0.08	23,31,33,34	0
4	CA	B	306	1/1	0.93	0.13	59,59,59,59	0

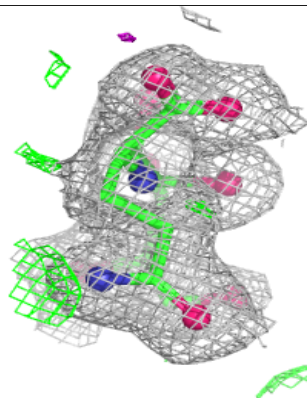
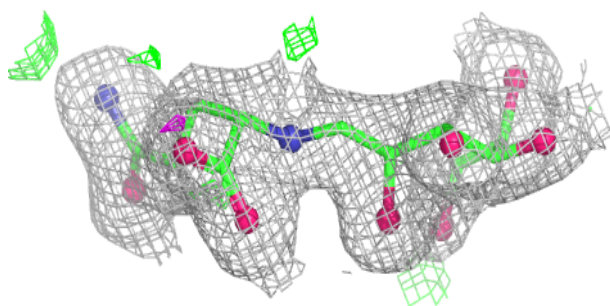
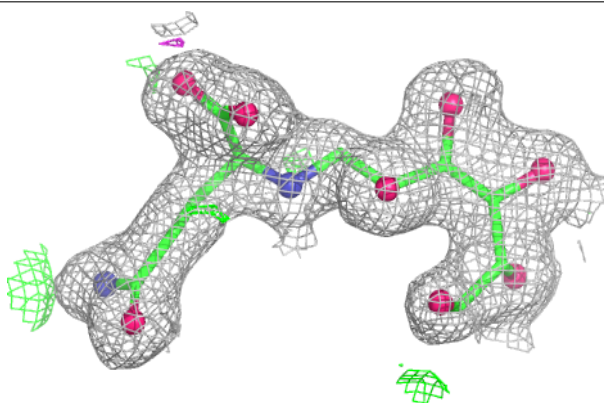
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around GOP A 301:

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

**Electron density around GOP B 301:**

$2mF_o-DF_c$ (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.