WORLDWIDE wwPDB PROTEIN DATA BANK

wwPDB X-ray Structure Validation Summary Report (i)

Nov 18, 2024 – 04:31 PM JST

PDB ID	:	9J6N
Title	:	beta-D-galactofuranosidase ORF1110 from Streptomyces sp. JHA19, D-iminogalactitol complex
Authors	:	Fujio, N.; Yamada, C.; Takegawa, K.; Fushinobu, S.
Deposited on	:	2024-08-16
Resolution	:	1.70 Å(reported)

The following versions of software and data (see references (i)) were used in the production of this report:

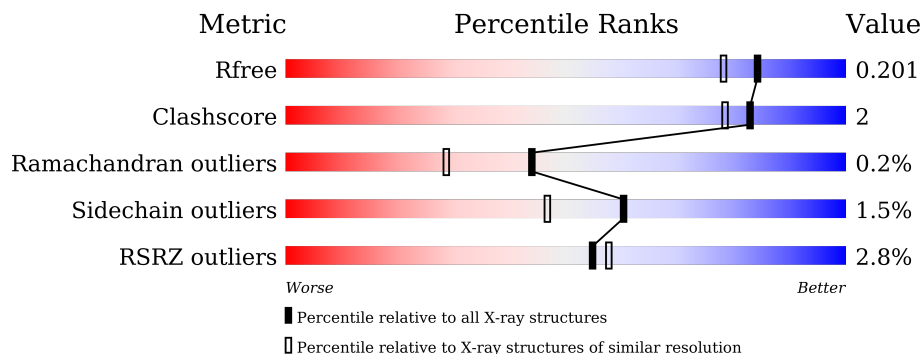|		
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	3.0
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.39

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 1.70 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	Percentile Ranks	Value
Rfree		0.201
Clashscore		2
Ramachandran outliers		0.2%
Sidechain outliers		1.5%
RSRZ outliers		2.8%

Worse   *Better*

▌ Percentile relative to all X-ray structures

▯ Percentile relative to X-ray structures of similar resolution

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	164625	5161 (1.70-1.70)
Clashscore	180529	5671 (1.70-1.70)
Ramachandran outliers	177936	5594 (1.70-1.70)
Sidechain outliers	177891	5594 (1.70-1.70)
RSRZ outliers	164620	5159 (1.70-1.70)

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 <=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	612	89% 5% • 5%
1	B	612	5% 86% 7% • 6%

2 Entry composition ⓘ

There are 7 unique types of molecules in this entry. The entry contains 9889 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 Beta-D-galactofuranosidase 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	581	Total 4524	C 2852	N 812	O 850	S 10	0	3	0
1	B	576	Total 4488	C 2830	N 810	O 839	S 9	0	1	0

There are 42 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	24	MET	-	initiating methionine	UNP A0A0K2SRV3
A	25	GLY	-	expression tag	UNP A0A0K2SRV3
A	26	SER	-	expression tag	UNP A0A0K2SRV3
A	27	SER	-	expression tag	UNP A0A0K2SRV3
A	28	HIS	-	expression tag	UNP A0A0K2SRV3
A	29	HIS	-	expression tag	UNP A0A0K2SRV3
A	30	HIS	-	expression tag	UNP A0A0K2SRV3
A	31	HIS	-	expression tag	UNP A0A0K2SRV3
A	32	HIS	-	expression tag	UNP A0A0K2SRV3
A	33	HIS	-	expression tag	UNP A0A0K2SRV3
A	34	SER	-	expression tag	UNP A0A0K2SRV3
A	35	ALA	-	expression tag	UNP A0A0K2SRV3
A	36	ALA	-	expression tag	UNP A0A0K2SRV3
A	37	LEU	-	expression tag	UNP A0A0K2SRV3
A	38	GLU	-	expression tag	UNP A0A0K2SRV3
A	39	VAL	-	expression tag	UNP A0A0K2SRV3
A	40	LEU	-	expression tag	UNP A0A0K2SRV3
A	41	PHE	-	expression tag	UNP A0A0K2SRV3
A	42	GLN	-	expression tag	UNP A0A0K2SRV3
A	43	GLY	-	expression tag	UNP A0A0K2SRV3
A	44	PRO	-	expression tag	UNP A0A0K2SRV3
B	24	MET	-	initiating methionine	UNP A0A0K2SRV3
B	25	GLY	-	expression tag	UNP A0A0K2SRV3
B	26	SER	-	expression tag	UNP A0A0K2SRV3
B	27	SER	-	expression tag	UNP A0A0K2SRV3

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Chain	Residue	Modelled	Actual	Comment	Reference
B	28	HIS	-	expression tag	UNP A0A0K2SRV3
B	29	HIS	-	expression tag	UNP A0A0K2SRV3
B	30	HIS	-	expression tag	UNP A0A0K2SRV3
B	31	HIS	-	expression tag	UNP A0A0K2SRV3
B	32	HIS	-	expression tag	UNP A0A0K2SRV3
B	33	HIS	-	expression tag	UNP A0A0K2SRV3
B	34	SER	-	expression tag	UNP A0A0K2SRV3
B	35	ALA	-	expression tag	UNP A0A0K2SRV3
B	36	ALA	-	expression tag	UNP A0A0K2SRV3
B	37	LEU	-	expression tag	UNP A0A0K2SRV3
B	38	GLU	-	expression tag	UNP A0A0K2SRV3
B	39	VAL	-	expression tag	UNP A0A0K2SRV3
B	40	LEU	-	expression tag	UNP A0A0K2SRV3
B	41	PHE	-	expression tag	UNP A0A0K2SRV3
B	42	GLN	-	expression tag	UNP A0A0K2SRV3
B	43	GLY	-	expression tag	UNP A0A0K2SRV3
B	44	PRO	-	expression tag	UNP A0A0K2SRV3

- Molecule 2 is HEXAETHYLENE GLYCOL (three-letter code: P6G) (formula: $C_{12}H_{26}O_7$).
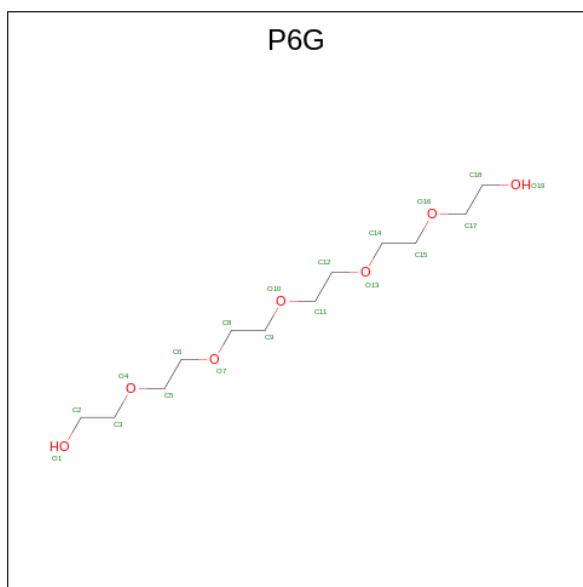
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total 19	C 12	O 7	0	0

- Molecule 3 is (4S)-2-METHYL-2,4-PENTANEDIOL (three-letter code: MPD) (formula: $C_6H_{14}O_2$).

MPD

O4
OH

C1

C5

C4(S)

C2

CM

C3

OH
O2
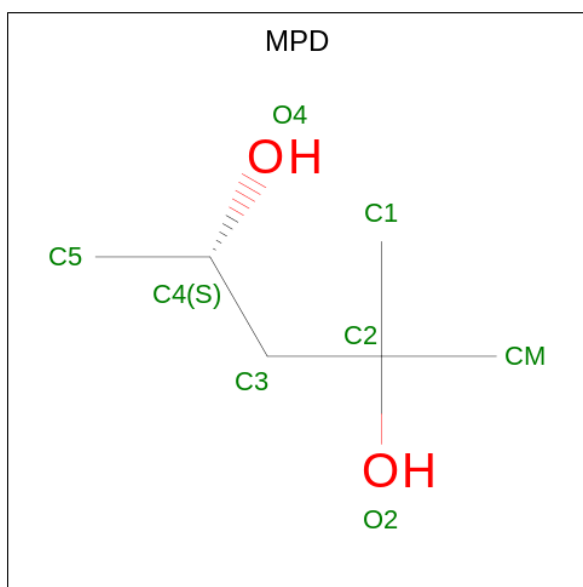
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total 8	C 6	O 2	0	0
3	A	1	Total 8	C 6	O 2	0	0
3	B	1	Total 8	C 6	O 2	0	0

- Molecule 4 is (2 {S},3 {S},4 {S})-2-[(1 {S})-1,2-bis(oxidanyl)ethyl]pyrrolidine-3,4-diol (three-letter code: A1L3Z) (formula: $C_6H_{13}NO_4$) (labeled as "Ligand of Interest" by depositor).
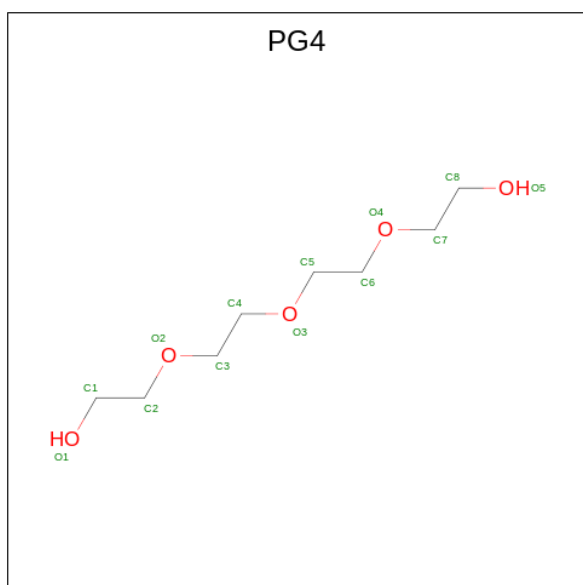
A1L3Z

O2
OH

N1
H
N

C4

O1
OH

C2(S)

C3(S)

C1

C6(S)

C5(S)

HO

OH

O3

O4
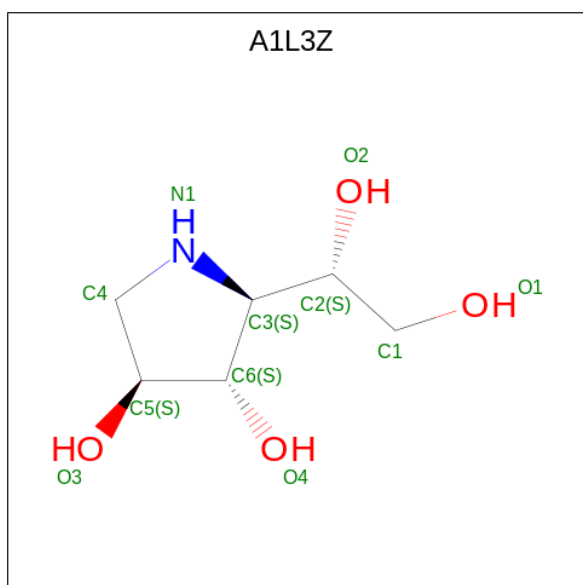
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total 11	C 6	N 1	O 4	0	0
4	B	1	Total 11	C 6	N 1	O 4	0	0

- Molecule 5 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total 1	Mg 1	0	0
5	B	1	Total 1	Mg 1	0	0

- Molecule 6 is TETRAETHYLENE GLYCOL (three-letter code: PG4) (formula: $C_8H_{18}O_5$).

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	B	1	Total 13	C 8	O 5	0	0

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	446	Total 446	O 446	0	0
7	B	351	Total 351	O 351	0	0

3 Residue-property plots ⓘ

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.
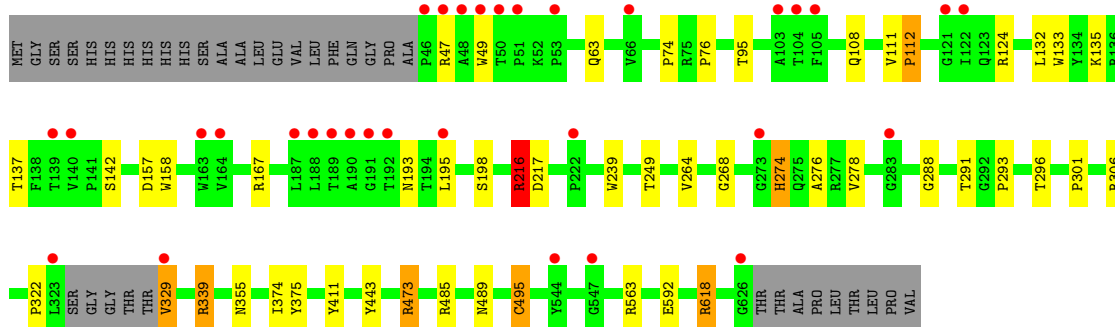
- Molecule 1: Beta-D-galactofuranosidase 2

Chain A:
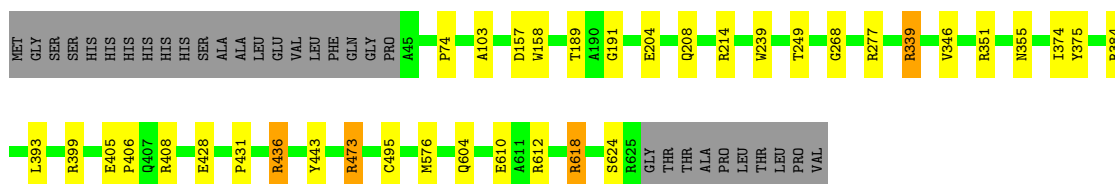
- Molecule 1: Beta-D-galactofuranosidase 2

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	82.36Å 57.35Å 109.73Å 90.00° 94.93° 90.00°	Depositor
Resolution (Å)	47.46 – 1.70 47.46 – 1.70	Depositor EDS
% Data completeness (in resolution range)	98.0 (47.46-1.70) 98.0 (47.46-1.70)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.30 (at 1.70Å)	Xtriage
Refinement program	REFMAC 5.8.0425	Depositor
R, R_{free}	0.153 , 0.189 0.167 , 0.201	Depositor DCC
R_{free} test set	5535 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å^2)	13.4	Xtriage
Anisotropy	0.020	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 46.0	EDS
L-test for twinning[2]	$< |L| > = 0.49, < L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	9889	wwPDB-VP
Average B, all atoms (Å^2)	16.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.51% of the height of the origin peak. No significant pseudotranslation is detected.*

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[1]Intensities estimated from amplitudes.
[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: A1L3Z, MPD, PG4, P6G, MG

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	0.50	1/4660 (0.0%)	0.91	6/6382 (0.1%)
1	B	0.47	1/4617 (0.0%)	0.91	9/6319 (0.1%)
All	All	0.48	2/9277 (0.0%)	0.91	15/12701 (0.1%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3
1	B	0	5
All	All	0	8

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	495	CYS	C-N	10.04	1.57	1.34
1	B	495	CYS	C-N	8.60	1.53	1.34

The worst 5 of 15 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	473	ARG	NE-CZ-NH2	-9.95	115.33	120.30
1	A	473	ARG	NE-CZ-NH2	-8.13	116.23	120.30
1	A	408	ARG	NE-CZ-NH1	7.04	123.82	120.30
1	A	436	ARG	NE-CZ-NH2	-6.59	117.00	120.30
1	A	473	ARG	NE-CZ-NH1	6.52	123.56	120.30

There are no chirality outliers.

5 of 8 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	384	ARG	Sidechain
1	A	399	ARG	Sidechain
1	A	473	ARG	Sidechain
1	B	216	ARG	Sidechain
1	B	47	ARG	Sidechain

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4524	0	4371	20	0
1	B	4488	0	4335	22	0
2	A	19	0	26	1	0
3	A	16	0	28	0	0
3	B	8	0	14	0	0
4	A	11	0	0	0	0
4	B	11	0	0	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
6	B	13	0	18	2	0
7	A	446	0	0	5	0
7	B	351	0	0	0	0
All	All	9889	0	8792	41	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 2.

The worst 5 of 41 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:610[A]:GLU:OE1	1:B:618[A]:ARG:NH2	2.26	0.65
1:A:604:GLN:CG	7:A:850:HOH:O	2.46	0.63
1:A:576[A]:MET:SD	1:A:624:SER:HB3	2.41	0.61
1:A:604:GLN:CD	7:A:850:HOH:O	2.41	0.59
1:A:103:ALA:HB1	2:A:701:P6G:H22	1.86	0.57

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	582/612 (95%)	567 (97%)	14 (2%)	1 (0%)	44	29
1	B	573/612 (94%)	558 (97%)	14 (2%)	1 (0%)	44	29
All	All	1155/1224 (94%)	1125 (97%)	28 (2%)	2 (0%)	44	29

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	374	ILE
1	B	374	ILE

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	476/498 (96%)	472 (99%)	4 (1%)	79	71
1	B	471/498 (95%)	461 (98%)	10 (2%)	48	32
All	All	947/996 (95%)	933 (98%)	14 (2%)	60	47

5 of 14 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	274	HIS

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Mol	Chain	Res	Type
1	B	329	VAL
1	B	489	ASN
1	B	443	TYR
1	B	485	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	63	GLN
1	B	63	GLN
1	B	98	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 9 ligands modelled in this entry, 2 are monoatomic - leaving 7 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	A1L3Z	A	704	-	11,11,11	0.80	0	7,15,15	0.90	1 (14%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#|Z| > 2$	Counts	RMSZ	$\#|Z| > 2$
6	PG4	B	701	-	12,12,12	0.14	0	11,11,11	0.26	0
4	A1L3Z	B	703	-	11,11,11	0.50	0	7,15,15	0.80	0
3	MPD	A	703	-	7,7,7	0.18	0	9,10,10	0.96	1 (11%)
2	P6G	A	701	-	18,18,18	0.17	0	17,17,17	0.22	0
3	MPD	A	702	-	7,7,7	0.19	0	9,10,10	0.90	0
3	MPD	B	702	-	7,7,7	0.25	0	9,10,10	0.51	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
4	A1L3Z	A	704	-	-	0/6/19/19	0/1/1/1
6	PG4	B	701	-	-	4/10/10/10	-
4	A1L3Z	B	703	-	-	0/6/19/19	0/1/1/1
3	MPD	A	703	-	-	0/5/5/5	-
2	P6G	A	701	-	-	9/16/16/16	-
3	MPD	A	702	-	-	3/5/5/5	-
3	MPD	B	702	-	-	3/5/5/5	-

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	703	MPD	O2-C2-C1	-2.24	100.89	108.08
4	A	704	A1L3Z	O1-C1-C2	2.09	115.63	111.07

There are no chirality outliers.
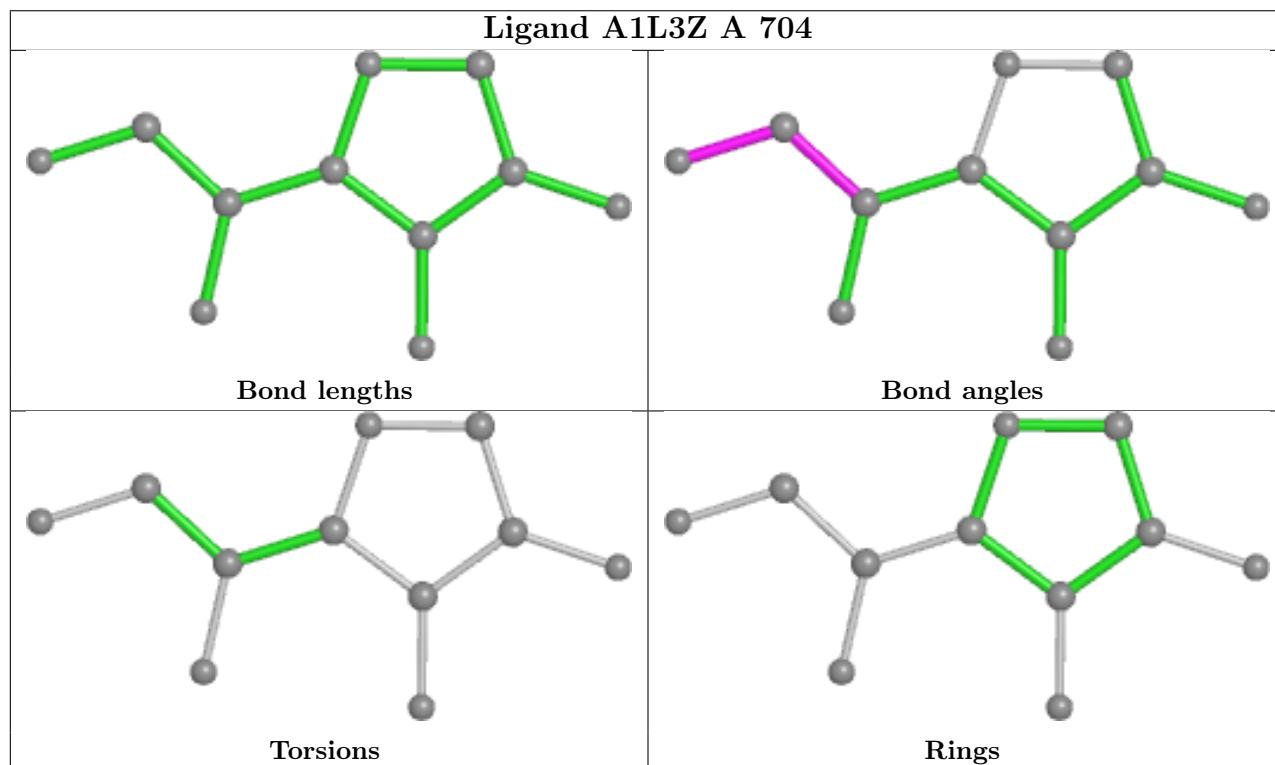
5 of 19 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	702	MPD	C1-C2-C3-C4
3	A	702	MPD	O2-C2-C3-C4
3	B	702	MPD	O2-C2-C3-C4
3	B	702	MPD	CM-C2-C3-C4
2	A	701	P6G	C15-C14-O13-C12

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	B	701	PG4	2	0
2	A	701	P6G	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 then 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.

Ligand A1L3Z A 704

Ligand A1L3Z B 703

Bond lengths

Bond angles

Torsions

Rings

Ligand P6G A 701

Bond lengths

Bond angles

Torsions

Rings
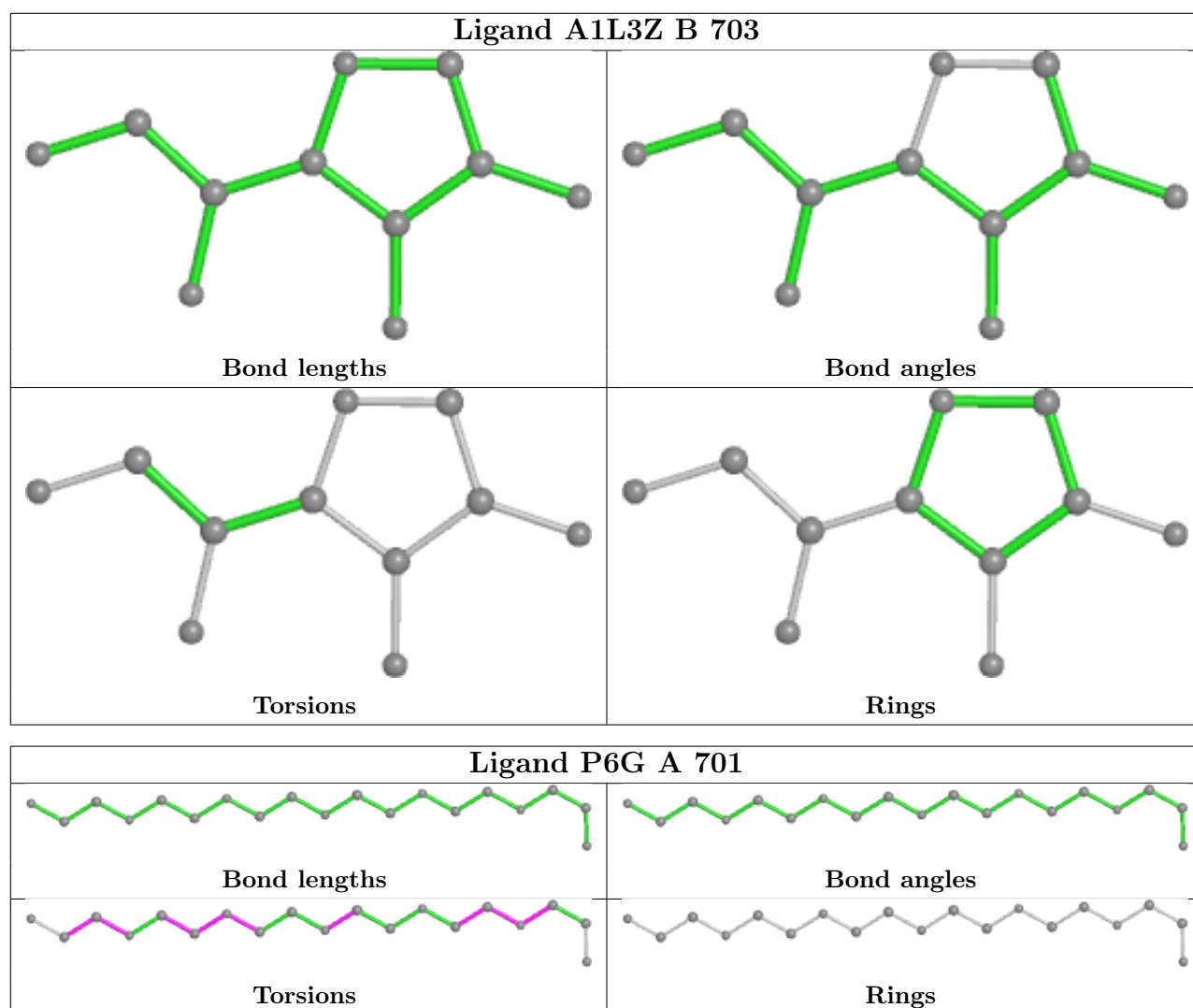
5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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(Å²)	Q<0.9
1	A	581/612 (94%)	-0.62	0	100	100	7, 11, 20, 34	3 (0%)
1	B	576/612 (94%)	0.13	32 (5%)	31	33	8, 18, 39, 65	1 (0%)
All	All	1157/1224 (94%)	-0.25	32 (2%)	55	58	7, 13, 32, 65	4 (0%)

The worst 5 of 32 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	46	PRO	4.9
1	B	329	VAL	4.5
1	B	192	THR	3.8
1	B	626	GLY	3.8
1	B	273	GLY	3.6

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

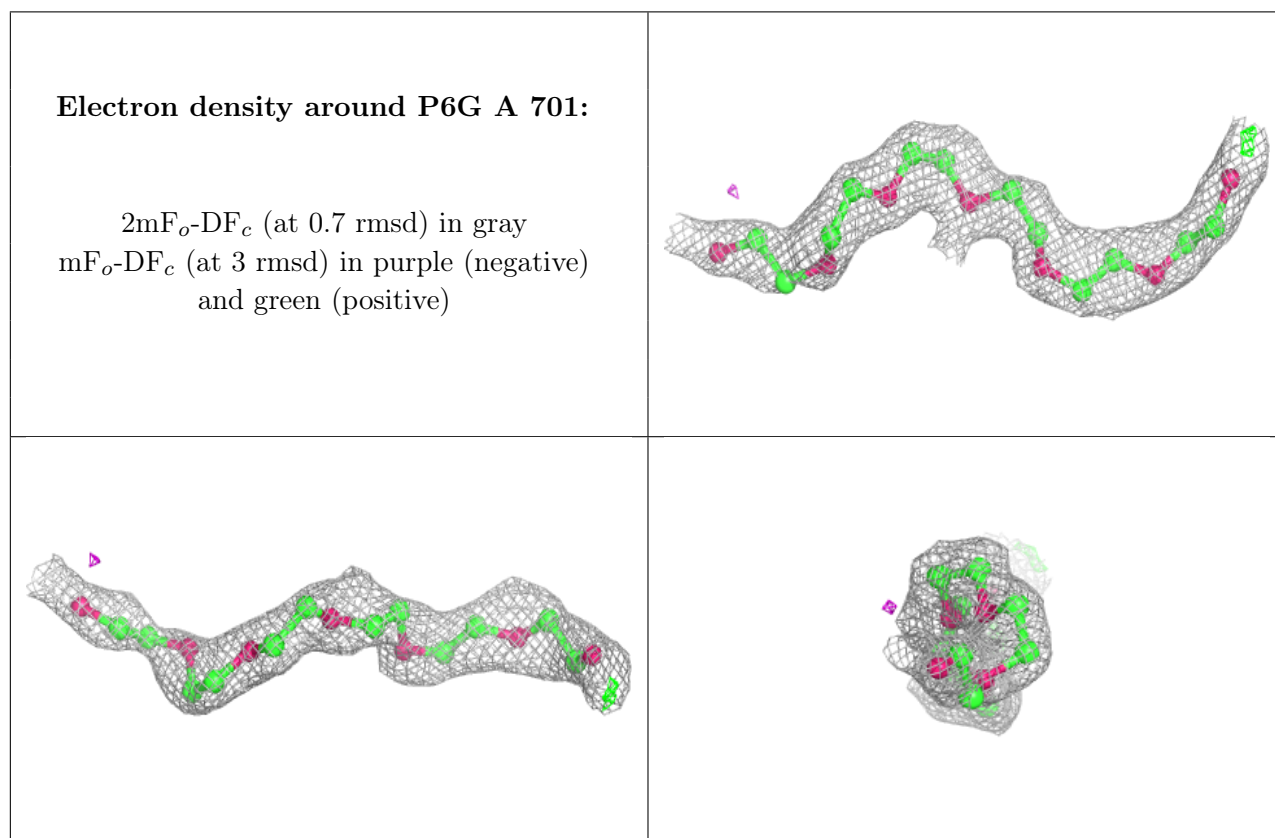There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no monosaccharides in this entry.

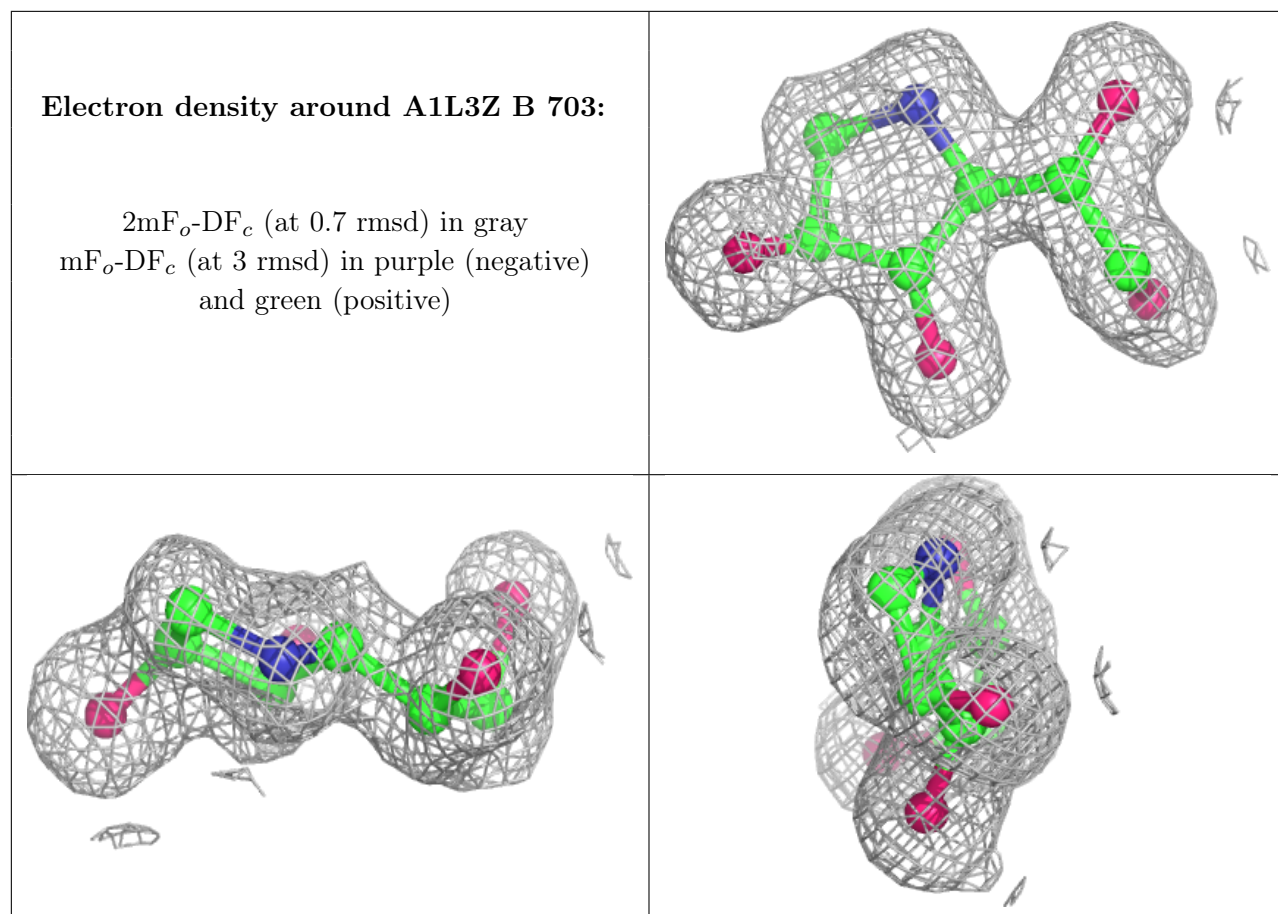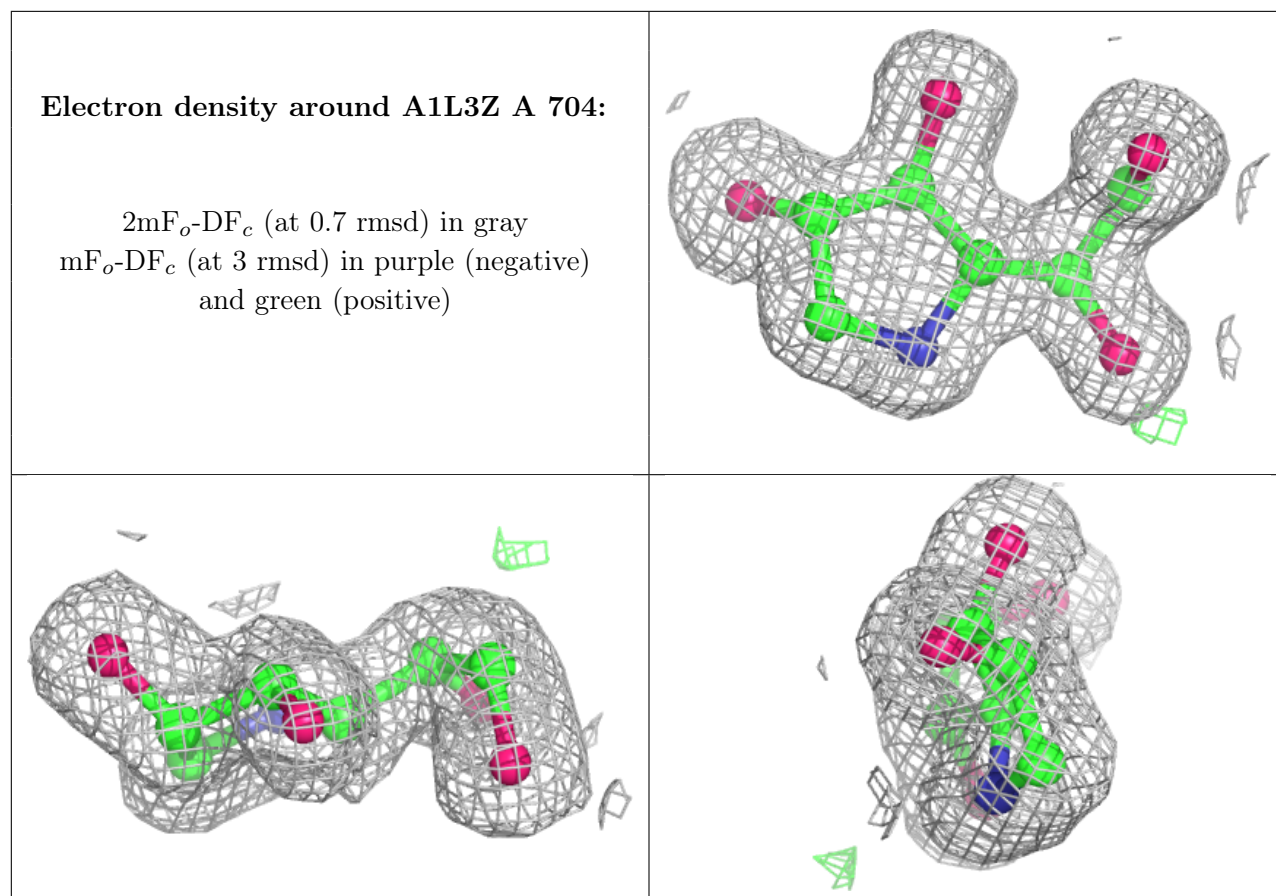## 6.4 Ligands ⓘ

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.

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å²)	Q<0.9
2	P6G	A	701	19/19	0.86	0.13	31,38,43,45	0
3	MPD	A	702	8/8	0.86	0.12	20,23,30,30	0
6	PG4	B	701	13/13	0.88	0.13	32,36,42,48	0
3	MPD	B	702	8/8	0.90	0.12	22,26,29,30	0
3	MPD	A	703	8/8	0.92	0.09	16,20,21,22	0
4	A1L3Z	B	703	11/11	0.96	0.05	11,11,12,12	0
4	A1L3Z	A	704	11/11	0.98	0.04	8,9,10,10	0
5	MG	B	704	1/1	0.99	0.02	14,14,14,14	0
5	MG	A	705	1/1	0.99	0.02	8,8,8,8	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 P6G A 701:

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

Electron density around A1L3Z B 703:

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

Electron density around A1L3Z A 704:

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

6.5 Other polymers (i)

There are no such residues in this entry.