

wwPDB X-ray Structure Validation Summary Report (i)

May 14, 2020 – 02:30 am BST

PDB ID : 4GMX

Title: Crystal structure of KPT185 in complex with CRM1-Ran-RanBP1

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Deposited on : 2012-08-16

Resolution : 2.10 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 (i) symbol.

The following versions of software and data (see references (1)) 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 : 2.11

117/0

buster-report : 1.1.7 (2018)

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

Refmac: 5.8.0158

CCP4 : 7.0.044 (Gargrove)

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

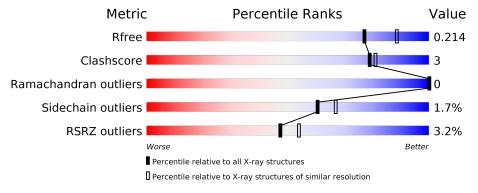
Validation Pipeline (wwPDB-VP) : 2.11

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 2.10 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	$\begin{array}{c} \text{Whole archive} \\ (\#\text{Entries}) \end{array}$	$\begin{array}{c} {\rm Similar \; resolution} \\ (\#{\rm Entries, \; resolution \; range(\AA)}) \end{array}$
R_{free}	130704	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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 on 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	216	6% 87%	8%	5%
2	В	141	7% 9%	• 13	%
3	С	1060	90%	6%	6 •



2 Entry composition (i)

There are 10 unique types of molecules in this entry. The entry contains 12477 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 GTP-binding nuclear protein Ran.

Mol	Chain	Residues		At	oms			ZeroOcc	AltConf	Trace
1	Λ	205	Total	С	N	О	S	0	2	0
1	Α	203	1672	1079	287	300	6	0	J	0

• Molecule 2 is a protein called Ran-specific GTPase-activating protein 1.

\mathbf{Mol}	Chain	Residues		\mathbf{At}	oms			ZeroOcc	AltConf	Trace
2	В	123	Total 1031	C 652	N 179	O 195	S 5	0	3	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
В	61	SER	-	EXPRESSION TAG	UNP P41920

• Molecule 3 is a protein called Exportin-1.

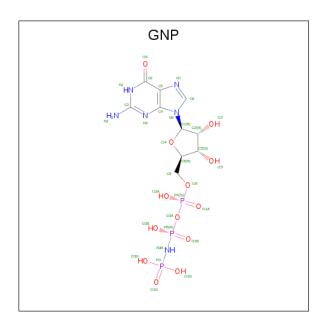
\mathbf{Mol}	Chain	Residues		Α	toms			ZeroOcc	AltConf	Trace
3	С	1015	Total 8439	C 5437	N 1380	O 1579	S 43	0	40	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
С	-1	GLY	-	EXPRESSION TAG	UNP P30822
С	0	ALA	-	EXPRESSION TAG	UNP P30822
С	539	CYS	THR	ENGINEERED MUTATION	UNP P30822
С	1022	CYS	TYR	CONFLICT	UNP P30822

• Molecule 4 is PHOSPHOAMINOPHOSPHONIC ACID-GUANYLATE ESTER (three-letter code: GNP) (formula: C₁₀H₁₇N₆O₁₃P₃).



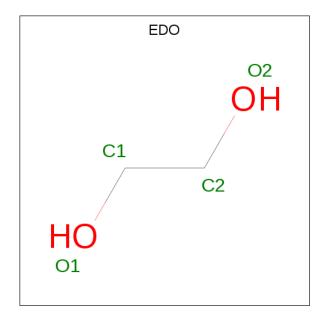


Mol	Chain	Residues		Ato	oms			ZeroOcc	AltConf
4	A	1	Total 32	C 10	_	O 13	P 3	0	0

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

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

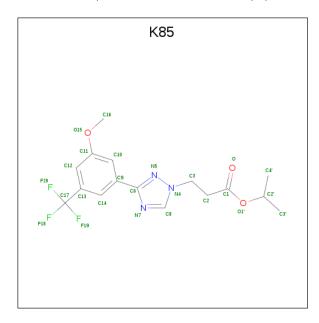
 \bullet Molecule 6 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $\mathrm{C_2H_6O_2}).$





\mathbf{Mol}	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	1	Total C O 4 2 2	0	0
6	С	1	Total C O 4 2 2	0	0

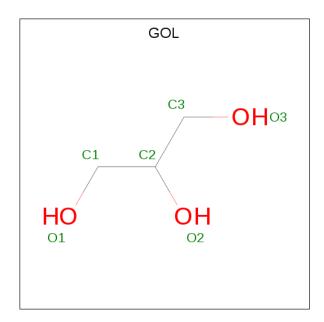
 $\bullet \ \, \text{Molecule 7 is propan-2-yl 3-\{3-[3-methoxy-5-(trifluoromethyl)phenyl]-1H-1,2,4-triazol-1-yl\} propanoate (three-letter code: K85) (formula: $C_{16}H_{18}F_3N_3O_3$). }$



Mol	Chain	Residues		Ato	oms			ZeroOcc	AltConf
7	С	1	Total	С	F	N	О	0	0
'		1	25	16	3	3	3		

• Molecule 8 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).





Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	С	1	Total C O 6 3 3	0	0
8	С	1	Total C O 6 3 3	0	0
8	С	1	Total C O 6 3 3	0	0
8	С	1	Total C O 6 3 3	0	0
8	С	1	Total C O 6 3 3	0	0
8	С	1	Total C O 6 3 3	0	0

• Molecule 9 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	С	3	Total Cl 3 3	0	0

• Molecule 10 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
10	A	205	Total O 205 205	0	0
10	В	65	Total O 65 65	0	0

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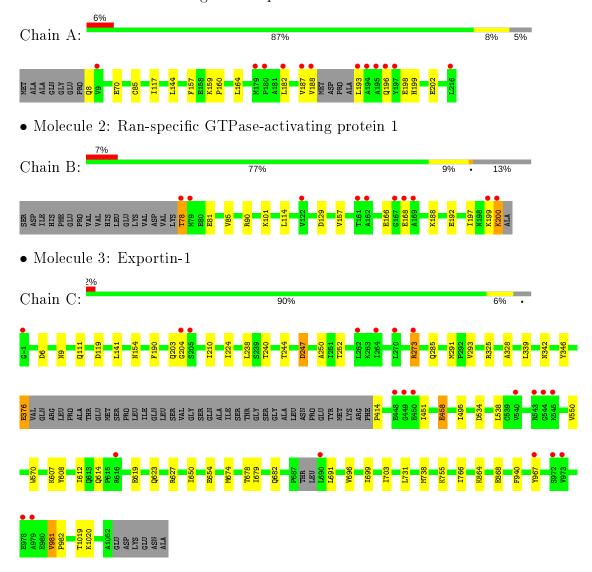
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
10	С	960	Total O 960 960	0	0



3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA 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: GTP-binding nuclear protein Ran





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants	106.49Å 106.49Å 306.80Å	Danagitan
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 - 2.10	Depositor
Resolution (A)	47.57 - 2.10	EDS
% Data completeness	96.0 (50.00-2.10)	Depositor
(in resolution range)	96.0 (47.57-2.10)	EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	1.95 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
D D.	0.148 , 0.197	Depositor
R, R_{free}	0.176 , 0.214	DCC
R_{free} test set	5014 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	39.5	Xtriage
Anisotropy	0.202	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.33, 47.3	EDS
L-test for twinning ²	$ < L > = 0.47, < L^2> = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	12477	wwPDB-VP
Average B, all atoms $(Å^2)$	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 4.16% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical 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.



¹Intensities estimated from amplitudes.

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: GOL, GNP, CL, MG, EDO, K85

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	
MIOI		RMSZ	# Z > 5	RMSZ	# Z >5
1	A	0.45	0/1714	0.56	0/2319
2	В	0.41	0/1050	0.55	0/1399
3	С	0.50	1/8685~(0.0%)	0.50	0/11761
All	All	0.49	1/11449~(0.0%)	0.52	0/15479

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	${ m Observed}({ m \AA})$	$\operatorname{Ideal}(ext{\AA})$
3	С	376	GLU	C-N	26.84	1.85	1.34

There are no bond angle outliers.

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	1672	0	1679	6	0
2	В	1031	0	1031	8	0
3	С	8439	0	8592	49	0
4	A	32	0	13	0	0
5	A	1	0	0	0	0
6	A	4	0	6	0	0
6	С	4	0	6	0	0

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-	110116	picolous	puyc

Mol	Chain	Non-H	$\mathbf{H}(\mathbf{model})$	$\mathbf{H}(\mathbf{added})$	Clashes	Symm-Clashes
7	С	25	0	16	0	0
8	С	36	0	48	0	0
9	С	3	0	0	0	0
10	A	205	0	0	1	0
10	В	65	0	0	0	0
10	С	960	0	0	4	0
All	All	12477	0	11391	63	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.

The worst 5 of 63 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$egin{array}{l} ext{Interatomic} \ ext{distance} \ (ext{Å}) \end{array}$	$egin{array}{c} ext{Clash} \ ext{overlap } (ext{Å}) \end{array}$
3:C:376:GLU:C	3:C:414:PRO:N	1.85	1.28
3:C:627[B]:ARG:HG2	3:C:627[B]:ARG:HH11	1.04	1.07
3:C:627[B]:ARG:HG2	3:C:627[B]:ARG:NH1	1.84	0.90
3:C:627[B]:ARG:HH11	3:C:627[B]:ARG:CG	1.88	0.86
3:C:244:THR:HG22	3:C:285:GLN:OE1	1.87	0.75

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	Perce	ntiles
1	A	$204/216 \ (94\%)$	199 (98%)	5 (2%)	0	100	100
2	В	123/141 (87%)	118 (96%)	5 (4%)	0	100	100
3	С	1051/1060 (99%)	1033 (98%)	18 (2%)	0	100	100
All	All	1378/1417 (97%)	1350 (98%)	28 (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	A	181/185 (98%)	174 (96%)	7 (4%)	32 33
2	В	108/122 (88%)	102 (94%)	6 (6%)	21 18
3	С	966/964 (100%)	956 (99%)	10 (1%)	76 82
All	All	1255/1271 (99%)	1232 (98%)	23 (2%)	60 65

5 of 23 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	В	166	GLU
2	В	200[B]	LYS
3	С	981	VAL
2	В	200[A]	LYS
3	С	203	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 7 such sidechains are listed below:

Mol	Chain	Res	Type
3	С	30	GLN
3	С	434	ASN
3	С	56	GLN
1	A	199	HIS
3	С	233	ASN

5.3.3 RNA (i)

There are no RNA molecules in this entry.



5.4 Non-standard residues in protein, DNA, RNA chains (i)

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

5.5 Carbohydrates (i)

There are no carbohydrates in this entry.

5.6 Ligand geometry (i)

Of 14 ligands modelled in this entry, 4 are monoatomic - leaving 10 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	Tune	Chain	Res	Link	Во	ond leng	ths	В	Bond angles		
MIOI	Type	Chain	nes	Lilik	Counts	RMSZ	# Z > 2	Counts	RMSZ	$\mid \# Z > 2$	
8	GOL	С	1111	-	5,5,5	0.37	0	5,5,5	0.32	0	
6	EDO	A	303	-	3,3,3	0.53	0	2,2,2	0.13	0	
7	K85	С	1101	3	25,26,26	0.96	1 (4%)	32,37,37	1.75	3 (9%)	
8	GOL	С	1106	-	5,5,5	0.37	0	5,5,5	0.32	0	
6	EDO	С	1107	-	3,3,3	0.51	0	2,2,2	0.29	0	
8	GOL	С	1105	-	5,5,5	0.44	0	5,5,5	0.69	0	
8	GOL	С	1102	-	5,5,5	0.35	0	5,5,5	0.25	0	
4	GNP	A	301	5	28,34,34	2.61	7 (25%)	30,54,54	1.21	3 (10%)	
8	GOL	С	1104	-	5,5,5	0.32	0	5,5,5	0.15	0	
8	GOL	С	1103	-	5,5,5	0.35	0	5,5,5	0.16	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
8	GOL	С	1111	-	-	3/4/4/4	-
6	EDO	A	303	-	-	0/1/1/1	-
7	K85	С	1101	3	-	4/21/21/21	0/2/2/2
8	GOL	С	1106	-	-	0/4/4/4	-

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Mol	Type	Chain	${ m Res}$	Link	Chirals	Torsions	Rings
6	EDO	С	1107	-	-	0/1/1/1	1
8	GOL	С	1105	-	-	2/4/4/4	-
8	GOL	С	1102	-	-	3/4/4/4	-
4	GNP	A	301	5	-	3/17/38/38	0/3/3/3
8	GOL	С	1104	_	-	4/4/4/4	-
8	GOL	С	1103	_	-	2/4/4/4	-

The worst 5 of 8 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\operatorname{Observed}(\operatorname{\AA})$	$\operatorname{Ideal}(\text{\AA})$
4	A	301	GNP	C4-N9	-10.93	1.33	1.47
7	С	1101	K85	O1'-C1	4.11	1.45	1.34
4	A	301	GNP	PG-N3B	4.00	1.73	1.63
4	A	301	GNP	PB-N3B	3.91	1.73	1.63
4	A	301	GNP	C8-N9	-2.84	1.35	1.45

The worst 5 of 6 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\mathbf{Observed}(^o)$	$\operatorname{Ideal}({}^o)$
7	С	1101	K85	N5-C6-N7	-7.03	108.89	114.72
7	С	1101	K85	N7-C8-N4	-4.06	106.42	112.21
4	A	301	GNP	C4-C5-N7	3.32	106.86	102.46
4	A	301	GNP	O1G-PG-N3B	-3.17	107.10	111.77
7	С	1101	K85	O1'-C1-C2	3.08	118.14	111.50

There are no chirality outliers.

5 of 21 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
8	С	1111	GOL	O1-C1-C2-C3
8	С	1105	GOL	C1-C2-C3-O3
8	С	1102	GOL	O1-C1-C2-C3
4	A	301	GNP	PG-N3B-PB-O1B
4	A	301	GNP	C2'-C1'-N9-C4

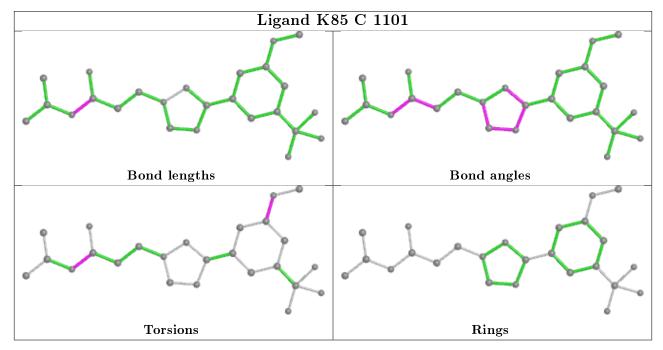
There are no ring outliers.

No monomer is involved in short contacts.

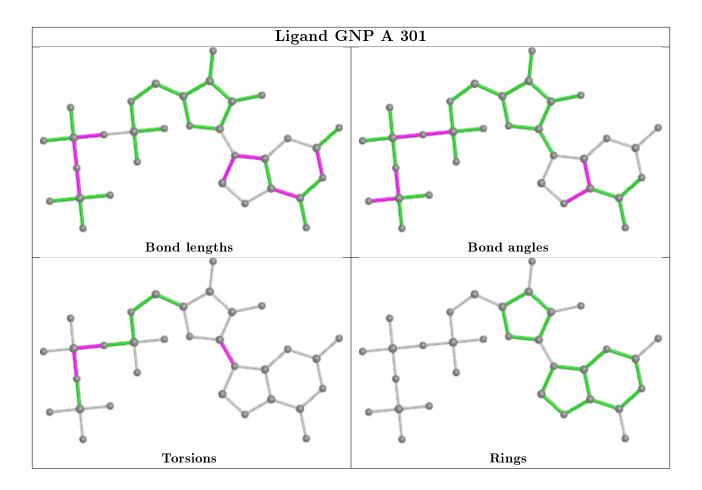
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will



also be included. For torsion angles, if less 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.







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, 95^{th} 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>	$\mathrm{SRZ}{>}$ $\mathrm{\#RSRZ}{>}2$		$OWAB(A^2)$	Q < 0.9
1	A	205/216~(94%)	0.28	12 (5%) 2	2 27	22, 31, 64, 98	0
2	В	123/141 (87%)	0.33	10 (8%) 1	2 15	29, 37, 53, 76	0
3	С	1015/1060 (95%)	-0.02	21 (2%) 6	3 68	22, 33, 58, 85	0
All	All	1343/1417 (94%)	0.06	43 (3%) 4	7 54	22, 34, 58, 98	0

The worst 5 of 43 RSRZ outliers are listed below:

Mol	Chain	81		RSRZ	
2	В	79	MET	7.5	
1	A	197	TYR	5.7	
1	A	194	ALA	5.5	
1	A	193	LEU	5.2	
1	A	196	GLN	4.8	

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 carbohydrates in this entry.

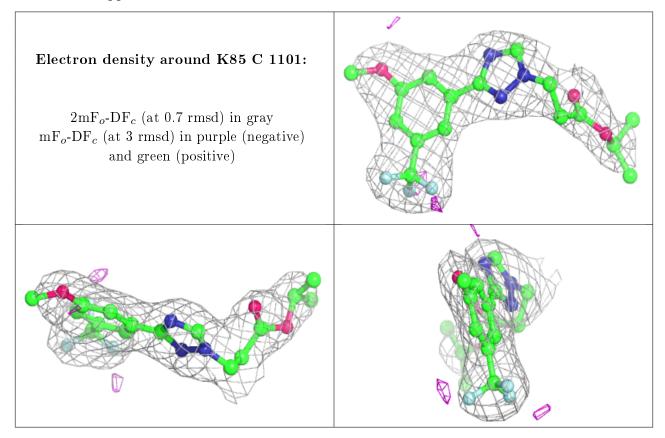
6.4 Ligands (i)

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, 95^{th} 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.

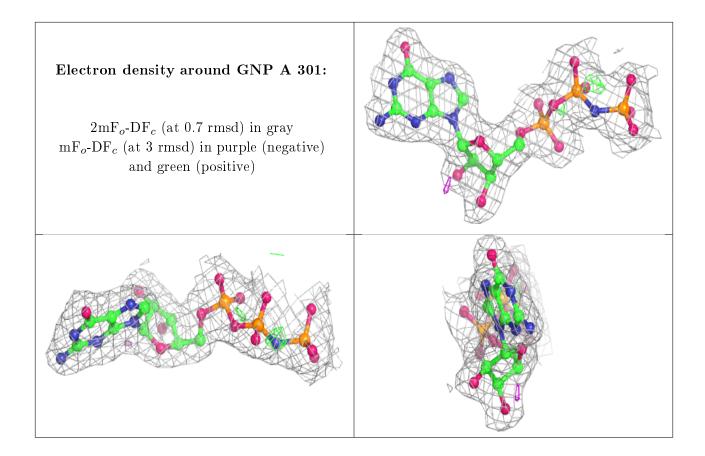


Mol	Type	Chain	Res	Atoms	RSCC	RSR	${f B-factors}({f \AA}^2)$	Q < 0.9
8	GOL	С	1111	6/6	0.68	0.20	56,69,75,77	0
8	GOL	С	1105	6/6	0.72	0.20	68,75,80,81	0
8	GOL	С	1106	6/6	0.81	0.30	58,69,75,76	0
8	GOL	С	1103	6/6	0.83	0.28	72,80,83,89	0
6	EDO	С	1107	4/4	0.88	0.17	58,59,61,70	0
8	GOL	С	1102	6/6	0.90	0.21	62,66,71,77	0
6	EDO	A	303	4/4	0.90	0.30	59,65,66,69	0
8	GOL	С	1104	6/6	0.92	0.15	59,62,63,66	0
7	K85	С	1101	25/25	0.95	0.18	37,42,57,61	0
9	CL	С	1110	1/1	0.96	0.11	89,89,89,89	0
9	CL	С	1108	1/1	0.96	0.09	71,71,71,71	0
9	CL	С	1109	1/1	0.98	0.08	66,66,66,66	0
5	MG	A	302	1/1	0.98	0.07	25,25,25,25	0
4	GNP	A	301	32/32	0.99	0.12	24,27,29,30	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.

