

wwPDB X-ray Structure Validation Summary Report (i)

Jun 17, 2024 – 01:22 AM EDT

PDB ID : 5GUG

Title : Crystal structure of inositol 1,4,5-trisphosphate receptor large cytosolic domain

with inositol 1,4,5-trisphosphate

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Deposited on : 2016-08-29

Resolution : 7.40 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 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: 4.02b-467

Mogul : 1.8.5 (274361), CSD as541be (2020)

Xtriage (Phenix) : 1.13

EDS : 2.37.1 buster-report : 1.1.7 (2018)

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

 $Refmac \quad : \quad 5.8.0158$

CCP4 : 7.0.044 (Gargrove)

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

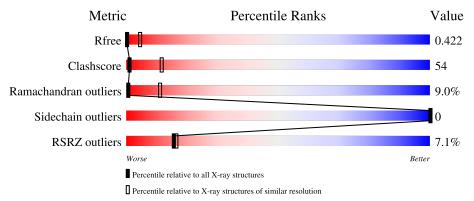
Validation Pipeline (wwPDB-VP) : 2.37.1

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 7.40 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	Similar resolution		
Metric	$(\# \mathrm{Entries})$	$(\# ext{Entries}, ext{ resolution range}(ext{Å}))$		
R_{free}	130704	1004 (10.00-3.90)		
Clashscore	141614	1069 (10.00-3.90)		
Ramachandran outliers	138981	1002 (10.00-3.90)		
Sidechain outliers	138945	1002 (10.00-3.86)		
RSRZ outliers	127900	1004 (9.50-3.80)		

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	2217	43%	30%	•	22%	_	
1	В	2217	43%	30%	5%	22%		

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:



Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	I3P	В	3000	_	_	-	X



2 Entry composition (i)

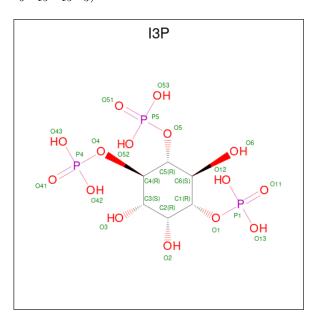
There are 2 unique types of molecules in this entry. The entry contains 25117 atoms, of which 7810 are hydrogens and 0 are deuteriums.

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 Inositol 1,4,5-trisphosphate receptor type 1.

]	Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
	1	A	1721	Total 12529	C 5147	H 3897	N 1746	O 1738	S 1	0	0	0
	1	В	1720	Total 12522	C 5144	H 3895	N 1745	O 1737	S 1	0	0	0

• Molecule 2 is D-MYO-INOSITOL-1,4,5-TRIPHOSPHATE (three-letter code: I3P) (formula: $C_6H_{15}O_{15}P_3$).



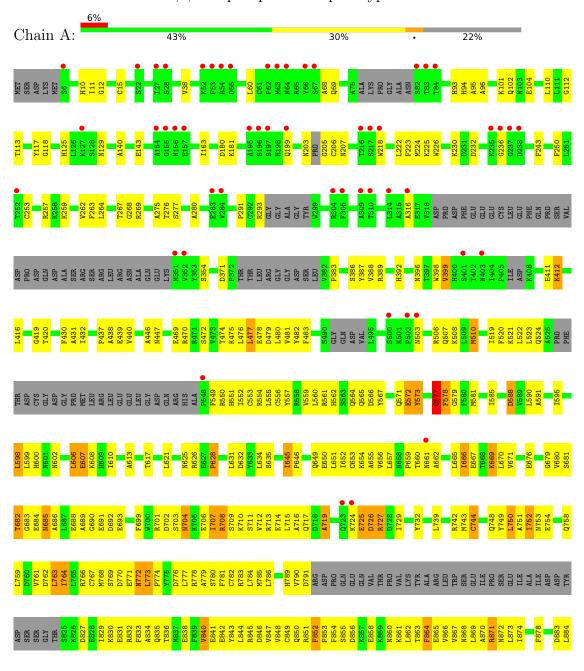
Mol	Chain	Residues	${f Atoms}$					ZeroOcc	AltConf	
9	٨	1	Total	С	Н	О	Р	0	0	
2	A	1	33	6	9	15	3	0	0	
9	D	1	Total	С	Н	О	Р	0	0	
2	Б	1	33	6	9	15	3	0		



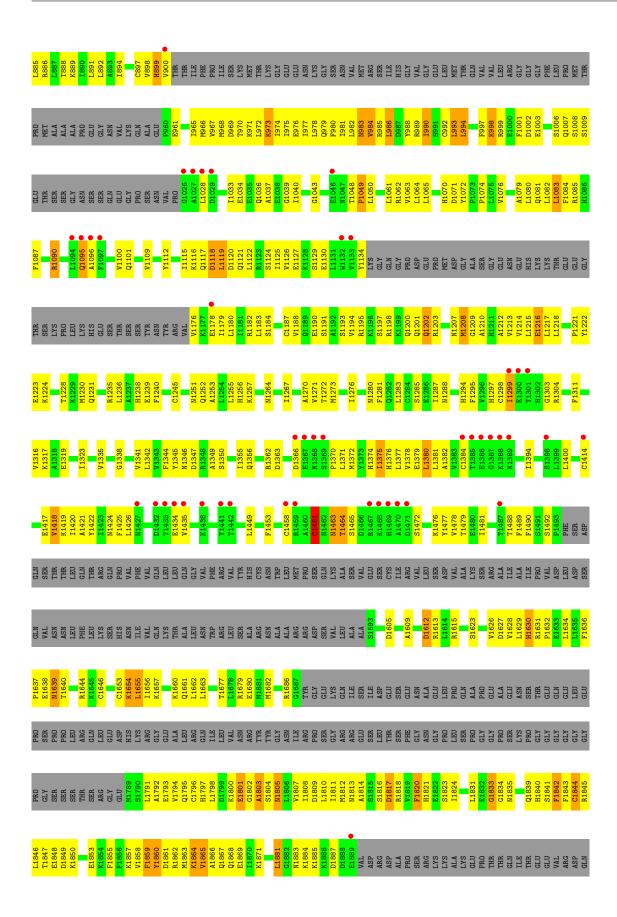
3 Residue-property plots (i)

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.

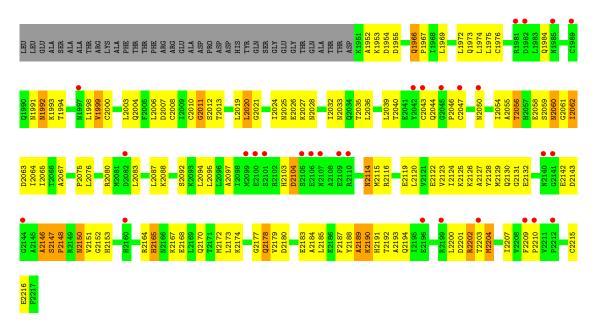
• Molecule 1: Inositol 1,4,5-trisphosphate receptor type 1



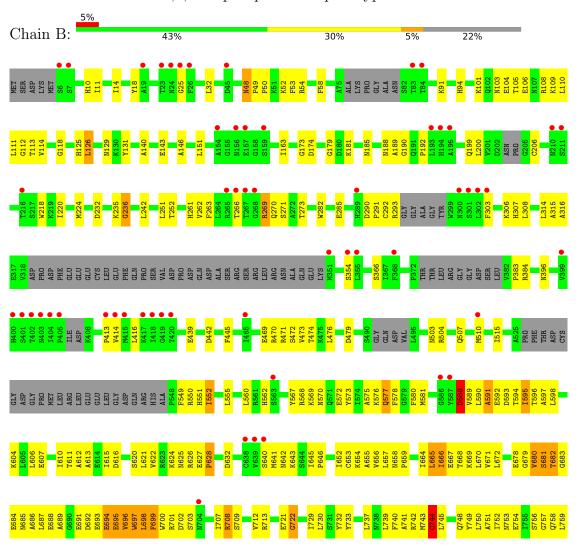




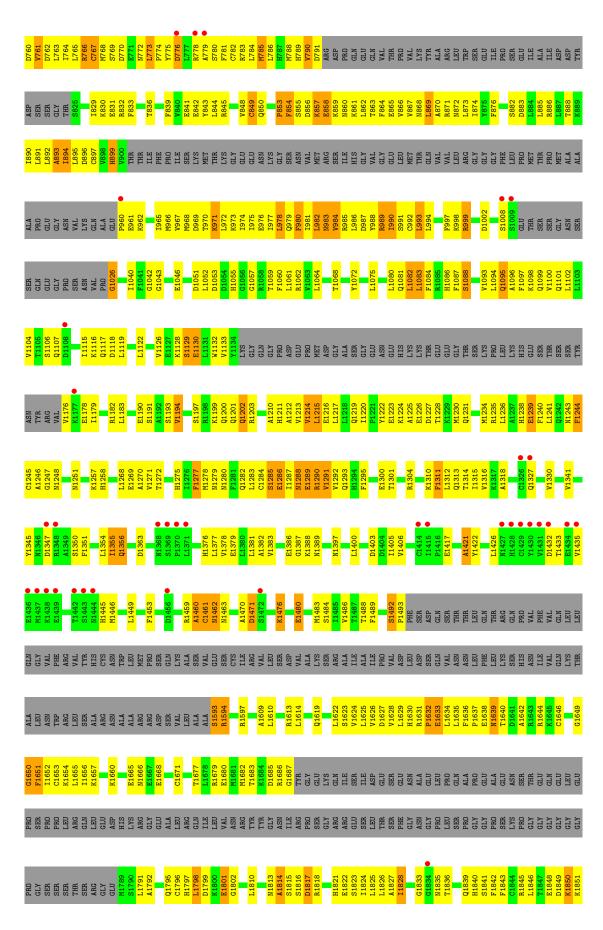




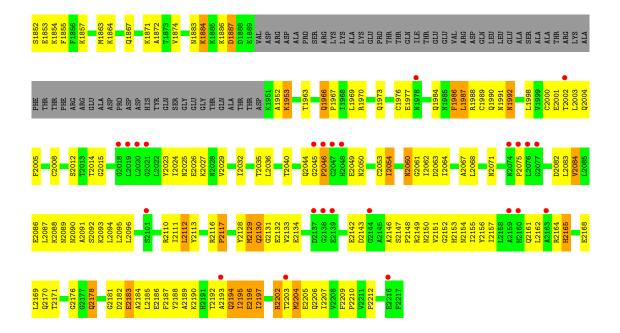
• Molecule 1: Inositol 1,4,5-trisphosphate receptor type 1













4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants	211.99Å 223.49Å 319.87Å	Donositon
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.49 - 7.40	Depositor
Resolution (A)	49.49 - 7.40	EDS
% Data completeness	99.9 (49.49-7.40)	Depositor
(in resolution range)	85.7 (49.49-7.40)	EDS
R_{merge}	0.29	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	1.62 (at 7.37Å)	Xtriage
Refinement program	PHENIX (1.10.1_2155: ???)	Depositor
D D.	0.356 , 0.416	Depositor
R, R_{free}	0.359 , 0.422	DCC
R_{free} test set	526 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	159.4	Xtriage
Anisotropy	0.177	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.26 , 212.2	EDS
L-test for twinning ²	$ < L >=0.23, < L^2>=0.08$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.62	EDS
Total number of atoms	25117	wwPDB-VP
Average B, all atoms (Å ²)	166.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 4.84% 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: I3P

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		ond lengths	Bond angles		
IVIOI	Chain	RMSZ	# Z > 5	RMSZ	# Z >5	
1	A	0.72	9/8617~(0.1%)	0.84	6/11978 (0.1%)	
1	В	0.71	$6/8612 \ (0.1\%)$	0.83	5/11971 (0.0%)	
All	All	0.71	$15/17229 \ (0.1\%)$	0.83	11/23949 (0.0%)	

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 maintenain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	8
1	В	0	9
All	All	0	17

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

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	Observed(A)	$\operatorname{Ideal}(\text{\AA})$
1	В	270	GLN	CD-NE2	-9.15	1.09	1.32
1	В	270	GLN	CD-OE1	-7.90	1.06	1.24
1	A	577	GLN	C-O	7.33	1.37	1.23
1	В	1026	GLY	C-O	-6.82	1.12	1.23
1	A	579	GLY	N-CA	-6.53	1.36	1.46

The worst 5 of 11 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\mathbf{Observed}(^o)$	$\operatorname{Ideal}({}^{o})$
1	A	577	GLN	C-N-CA	9.25	144.83	121.70
1	A	578	PHE	N-CA-CB	8.28	125.51	110.60
1	В	588	ASP	N-CA-C	6.75	129.24	111.00
1	В	588	ASP	N-CA-CB	-6.54	98.82	110.60

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\mathbf{Mol}	Chain	Res	Type	Atoms	${f Z}$	$\operatorname{Observed}(^{o})$	$\operatorname{Ideal}({}^{o})$
1	A	510	MET	CB-CG-SD	-5.91	94.68	112.40

There are no chirality outliers.

5 of 17 planarity outliers are listed below:

Mol	Chain	Res	Type	Group	
1	A	1034	GLU	Peptide	
1	A	1375	ILE	Peptide	
1	A	1380	LEU	Peptide	
1	A	577	GLN	Peptide	
1	A	588	ASP	Peptide	

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	8632	3897	3899	667	0
1	В	8627	3895	3898	687	0
2	A	24	9	9	6	0
2	В	24	9	9	1	0
All	All	17307	7810	7815	1353	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 54.

The worst 5 of 1353 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	$\begin{array}{c} \text{Clash} \\ \text{overlap } (\text{\AA}) \end{array}$
1:A:2020:LEU:O	1:A:2024:ILE:N	1.57	1.38
1:B:863:THR:O	1:B:867:VAL:N	1.72	1.21
1:A:1210:ALA:O	1:A:1214:VAL:CB	1.89	1.20
1:A:1125:ILE:O	1:A:1129:SER:CB	1.91	1.19
1:B:1682:MET:O	1:B:1686:ARG:N	1.78	1.16

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	1689/2217 (76%)	1219 (72%)	330 (20%)	140 (8%)	1 12		
1	В	1688/2217 (76%)	1173 (70%)	350 (21%)	165 (10%)	0 9		
All	All	3377/4434 (76%)	2392 (71%)	680 (20%)	305 (9%)	1 11		

5 of 305 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	598	LEU
1	A	628	PRO
1	A	659	PRO
1	A	666	ILE
1	A	669	LYS

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	26/1980 (1%)	26 (100%)	0	100 100		
1	В	26/1980 (1%)	26 (100%)	0	100 100		
All	All	52/3960~(1%)	52 (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. All (1) such sidechains are listed below:



Mol	Chain	Res	Type
1	В	270	GLN

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

2 ligands are modelled in this entry.

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 T	Type	Chain	Res	Link	Bond lengths			Bond angles		
	туре		rtes	Lilik	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	I3P	A	3000	-	24,24,24	1.13	1 (4%)	36,39,39	1.08	3 (8%)
2	I3P	В	3000	-	24,24,24	1.15	2 (8%)	36,39,39	1.09	4 (11%)

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.

\mathbf{N}	Iol	\mathbf{Type}	Chain	Res	Link	Chirals	Torsions	Rings
	2	I3P	A	3000	-	-	0/15/39/39	0/1/1/1
	2	I3P	В	3000	-	-	0/15/39/39	0/1/1/1

All (3) bond length outliers are listed below:



Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\operatorname{Observed}(\text{\AA})$	$\operatorname{Ideal}(ext{\AA})$
2	В	3000	I3P	P5-O53	-2.23	1.46	1.54
2	A	3000	I3P	P4-O42	-2.15	1.46	1.54
2	В	3000	I3P	P1-O13	-2.12	1.46	1.54

The worst 5 of 7 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	${f Z}$	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$
2	В	3000	I3P	O13-P1-O12	2.51	117.24	107.64
2	В	3000	I3P	O43-P4-O42	2.38	116.74	107.64
2	A	3000	I3P	O53-P5-O52	2.27	116.32	107.64
2	A	3000	I3P	O1-P1-O11	-2.26	100.67	109.39
2	В	3000	I3P	O1-C1-C2	2.12	113.60	108.66

There are no chirality outliers.

There are no torsion outliers.

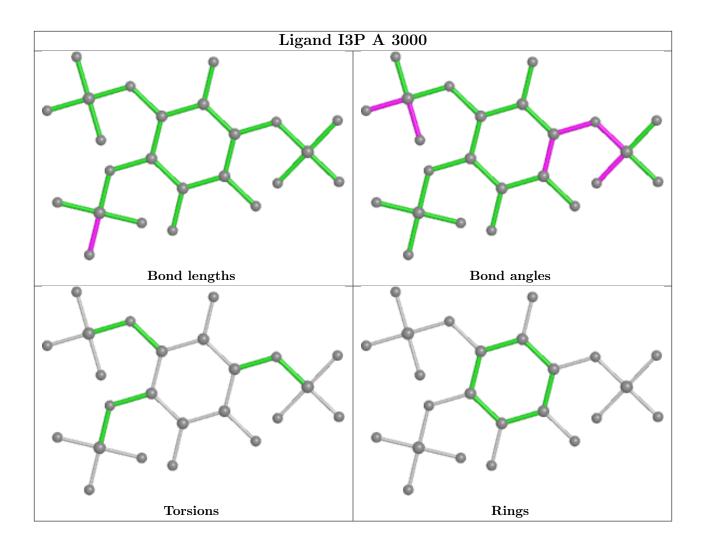
There are no ring outliers.

2 monomers are involved in 7 short contacts:

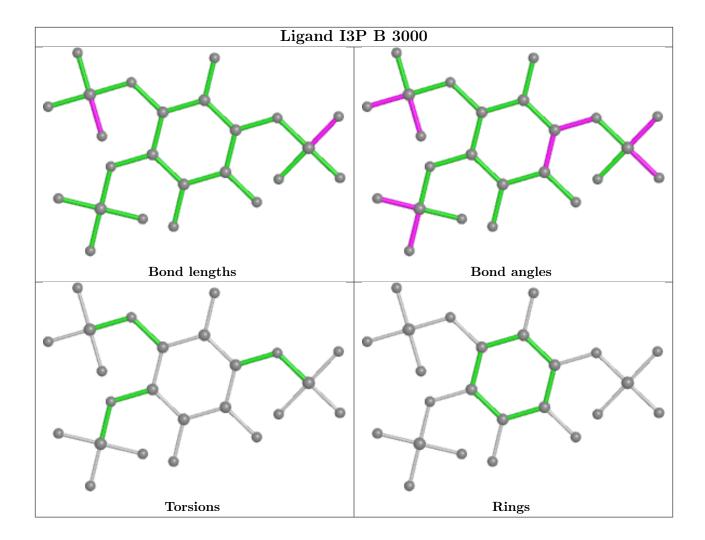
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	3000	I3P	6	0
2	В	3000	I3P	1	0

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.









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 $>$ 2		$OWAB(Å^2)$	Q<0.9
1	A	1721/2217 (77%)	0.15	126 (7%) 15 16	135, 166, 176, 187	0
1	В	$1720/2217 \ (77\%)$	0.10	117 (6%) 17 18	145, 165, 178, 192	0
All	All	3441/4434 (77%)	0.12	243 (7%) 16 16	135, 166, 177, 192	0

The worst 5 of 243 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	В	1436	GLU	7.8
1	В	401	SER	7.8
1	A	1468	LYS	7.3
1	В	1437	MET	6.8
1	A	1368	ASN	6.3

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 monosaccharides in this entry.

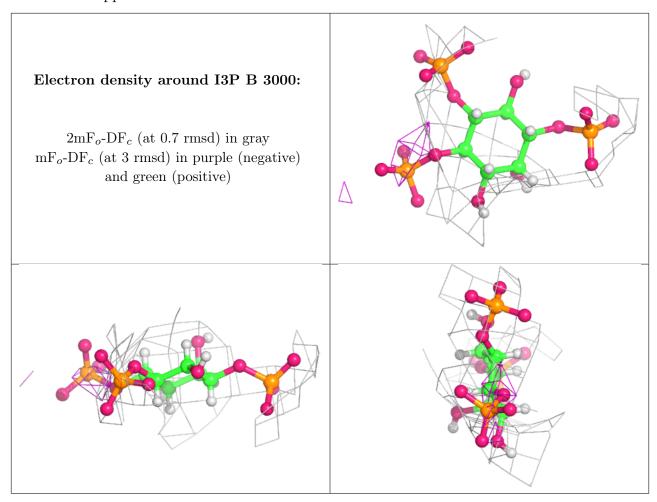
6.4 Ligands (i)

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.

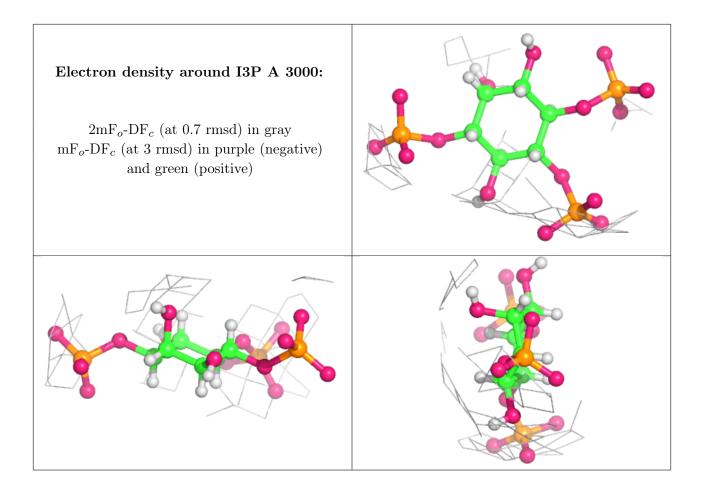


\mathbf{N}	Iol	Type	Chain	Res	Atoms	RSCC	RSR	${f B-factors(\AA^2)}$	Q<0.9
	2	I3P	В	3000	24/24	0.57	0.70	199,199,203,203	0
	2	I3P	A	3000	24/24	0.59	0.35	169,172,172,172	0

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.

