



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 17, 2022 – 02:37 PM EDT

PDB ID : 4G61
Title : Crystal structure of IMPase/NADP phosphatase complexed with Mg²⁺ and phosphate
Authors : Bhattacharyya, S.; Dutta, D.; Ghosh, A.K.; Das, A.K.
Deposited on : 2012-07-18
Resolution : 2.30 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.29
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) : 2.29

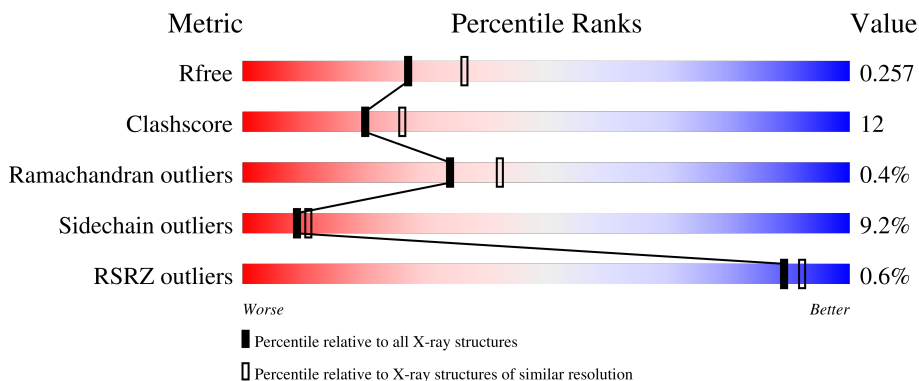
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION


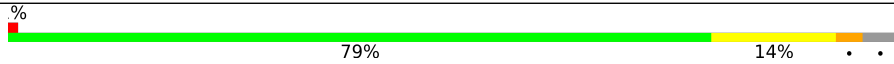
The reported resolution of this entry is 2.30 Å.

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}	130704	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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	271	
1	B	271	

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
4	GOL	B	308	-	-	X	-
6	DTO	B	307[A]	-	X	X	-
6	DTO	B	307[B]	-	-	X	-
7	P33	B	309	-	-	X	X

2 Entry composition i

There are 9 unique types of molecules in this entry. The entry contains 4548 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 Inositol monophosphatase family protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	264	2118	1363	350	392	13	0	2	0
1	B	261	2086	1346	345	381	14	0	1	0

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	HIS	-	expression tag	UNP Q6G709
A	-4	HIS	-	expression tag	UNP Q6G709
A	-3	HIS	-	expression tag	UNP Q6G709
A	-2	HIS	-	expression tag	UNP Q6G709
A	-1	HIS	-	expression tag	UNP Q6G709
A	0	HIS	-	expression tag	UNP Q6G709
B	-5	HIS	-	expression tag	UNP Q6G709
B	-4	HIS	-	expression tag	UNP Q6G709
B	-3	HIS	-	expression tag	UNP Q6G709
B	-2	HIS	-	expression tag	UNP Q6G709
B	-1	HIS	-	expression tag	UNP Q6G709
B	0	HIS	-	expression tag	UNP Q6G709

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	3	Total	Mg	0	0
			3	3		
2	B	3	Total	Mg	0	0
			3	3		

- Molecule 3 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O P 5 4 1	0	0
3	B	1	Total O P 5 4 1	0	0

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 6 3 3	0	0
4	A	1	Total C O 6 3 3	0	0

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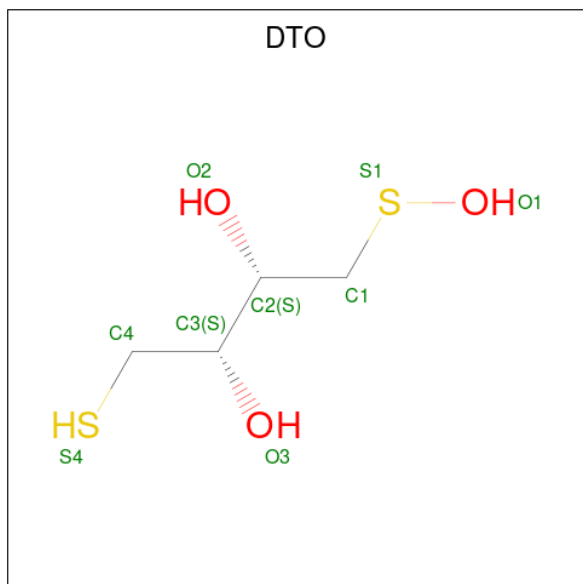
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

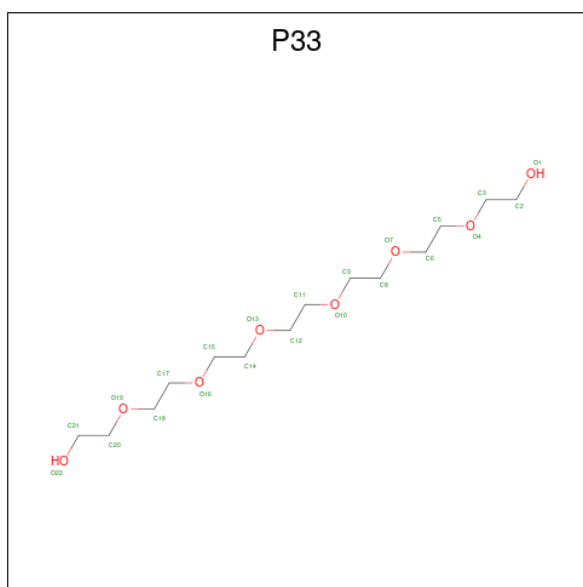
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Cl	0	0
			1	1		
5	B	1	Total	Cl	0	0
			1	1		

- Molecule 6 is 1-HYDROXYSULFANYL-4-MERCAPTO-BUTANE-2,3-DIOL (three-letter code: DTO) (formula: C₄H₁₀O₃S₂).



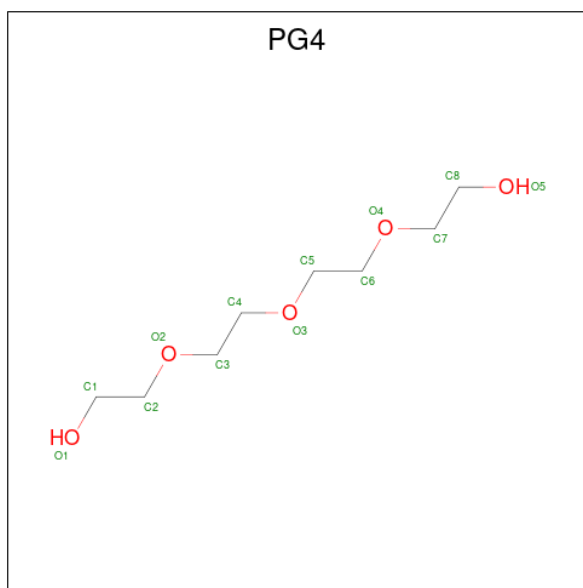
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	B	1	Total	C	O	S	0	1
			16	8	4	4		

- Molecule 7 is 3,6,9,12,15,18-HEXA-OXAICOSANE-1,20-DIOL (three-letter code: P33) (formula: C₁₄H₃₀O₈).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	B	1	Total	C	O	0	0
			22	14	8		

- Molecule 8 is TETRAETHYLENE GLYCOL (three-letter code: PG4) (formula: C₈H₁₈O₅).



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

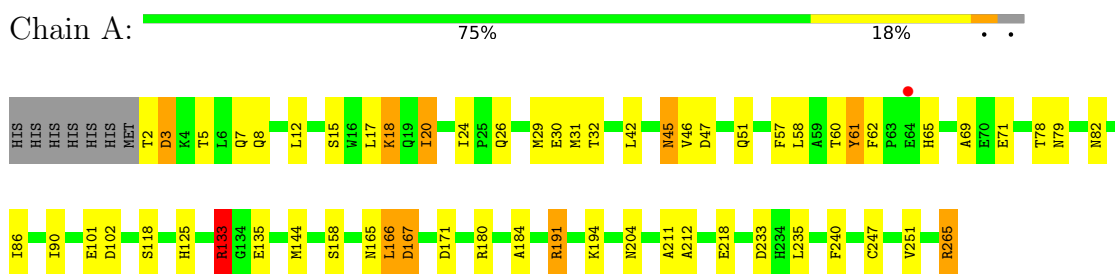
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	148	Total 148	O 148	0	0
9	B	103	Total 103	O 103	0	0

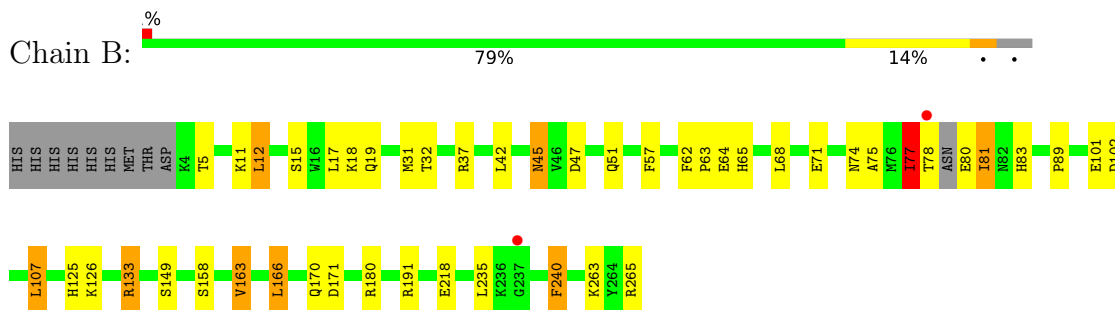
3 Residue-property plots [i](#)

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: Inositol monophosphatase family protein



- Molecule 1: Inositol monophosphatase family protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 2 21	Depositor
Cell constants a, b, c, α , β , γ	60.71Å 63.05Å 141.62Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.67 – 2.30 19.67 – 2.30	Depositor EDS
% Data completeness (in resolution range)	98.2 (19.67-2.30) 98.5 (19.67-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.04 (at 2.30Å)	Xtrriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.176 , 0.255 0.176 , 0.257	Depositor DCC
R_{free} test set	1235 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	22.3	Xtrriage
Anisotropy	0.502	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 37.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.020 for k,h,-l	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4548	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.94% 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: GOL, MG, DTO, PO4, CL, P33, PG4

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	0.92	2/2170 (0.1%)	0.95	7/2935 (0.2%)
1	B	0.88	0/2131	0.97	9/2881 (0.3%)
All	All	0.90	2/4301 (0.0%)	0.96	16/5816 (0.3%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	61	TYR	CD2-CE2	-6.45	1.29	1.39
1	A	46	VAL	CB-CG2	5.39	1.64	1.52

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	180	ARG	NE-CZ-NH2	-13.69	113.45	120.30
1	B	180	ARG	NE-CZ-NH2	-12.27	114.16	120.30
1	B	133	ARG	NE-CZ-NH2	-11.86	114.37	120.30
1	A	180	ARG	NE-CZ-NH1	10.62	125.61	120.30
1	A	133	ARG	NE-CZ-NH2	-9.80	115.40	120.30
1	B	133	ARG	NE-CZ-NH1	9.53	125.06	120.30
1	B	180	ARG	NE-CZ-NH1	8.39	124.49	120.30
1	A	265	ARG	NE-CZ-NH1	7.15	123.88	120.30
1	B	107	LEU	CA-CB-CG	5.84	128.74	115.30
1	A	133	ARG	NE-CZ-NH1	5.70	123.15	120.30
1	B	180	ARG	CG-CD-NE	-5.59	100.06	111.80
1	B	77	ILE	N-CA-C	5.45	125.71	111.00
1	B	37	ARG	NE-CZ-NH2	-5.41	117.59	120.30
1	A	180	ARG	CD-NE-CZ	5.17	130.84	123.60
1	B	107	LEU	CB-CG-CD1	-5.04	102.42	111.00
1	A	191	ARG	NE-CZ-NH1	5.02	122.81	120.30

There are no chirality outliers.

There are no planarity outliers.

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	2118	0	2099	60	0
1	B	2086	0	2061	43	0
2	A	3	0	0	0	0
2	B	3	0	0	0	0
3	A	5	0	0	1	0
3	B	5	0	0	1	0
4	A	12	0	16	4	0
4	B	12	0	16	5	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
6	B	16	0	18	18	0
7	B	22	0	30	14	0
8	B	13	0	18	4	0
9	A	148	0	0	7	0
9	B	103	0	0	2	0
All	All	4548	0	4258	107	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 12.

All (107) 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:64:GLU:HB2	7:B:309:P33:H92	1.42	0.98
1:A:171:ASP:OD1	4:A:307:GOL:H11	1.64	0.97
1:A:57:PHE:O	1:A:61:TYR:HD2	1.46	0.97
1:B:63:PRO:HD2	7:B:309:P33:H111	1.48	0.96
1:A:191:ARG:NH2	6:B:307[B]:DTO:H42	1.82	0.92
1:A:125:HIS:CE1	6:B:307[A]:DTO:H11	2.08	0.88
1:A:184:ALA:HA	4:A:305:GOL:H12	1.62	0.81
1:A:60:THR:HB	1:A:61:TYR:CD2	2.17	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:31:MET:HA	1:A:45:ASN:HD21	1.48	0.79
1:A:51:GLN:HE22	1:A:71:GLU:H	1.29	0.79
1:B:171:ASP:OD1	4:B:308:GOL:C3	2.31	0.78
1:A:57:PHE:O	1:A:61:TYR:CD2	2.36	0.75
1:A:60:THR:HB	1:A:61:TYR:CE2	2.20	0.75
1:A:60:THR:HA	9:A:428:HOH:O	1.85	0.75
1:B:133:ARG:NH2	1:B:218:GLU:OE2	2.21	0.74
1:A:167:ASP:OD1	4:A:307:GOL:H32	1.88	0.73
1:B:51:GLN:HE22	1:B:71:GLU:H	1.34	0.73
1:A:5:THR:H	1:A:8:GLN:HE21	1.37	0.71
1:A:102:ASP:OD1	6:B:307[A]:DTO:S1	2.50	0.69
1:A:194:LYS:HB3	8:B:310:PG4:H52	1.75	0.69
1:A:165[A]:ASN:ND2	1:A:167:ASP:HB2	2.08	0.69
1:A:191:ARG:HH22	6:B:307[B]:DTO:H42	1.58	0.68
1:B:125:HIS:CE1	6:B:307[B]:DTO:S1	2.90	0.65
1:B:101:GLU:OE2	6:B:307[A]:DTO:O3	2.15	0.64
1:B:68:LEU:HD12	1:B:74:ASN:HB3	1.80	0.64
1:B:45:ASN:H	1:B:45:ASN:HD22	1.46	0.64
1:A:133:ARG:NH2	1:A:218:GLU:OE2	2.28	0.64
1:B:81:ILE:HD12	1:B:81:ILE:H	1.63	0.62
1:B:77:ILE:HD12	1:B:80:GLU:HA	1.81	0.62
1:B:77:ILE:O	1:B:81:ILE:HD11	2.00	0.62
3:B:301:PO4:P	9:B:406:HOH:O	2.58	0.62
1:B:64:GLU:HB2	7:B:309:P33:C9	2.24	0.61
1:A:101:GLU:OE2	6:B:307[B]:DTO:H12	2.01	0.60
1:A:7[A]:GLN:OE1	1:A:135:GLU:OE2	2.19	0.60
1:B:62:PHE:HB3	1:B:65:HIS:CD2	2.37	0.59
1:B:171:ASP:OD1	4:B:308:GOL:C2	2.49	0.59
1:A:2:THR:O	1:A:3:ASP:HB2	2.03	0.58
1:A:32:THR:H	1:A:45:ASN:ND2	2.00	0.58
1:B:163:VAL:HG13	9:B:431:HOH:O	2.02	0.58
1:B:191:ARG:NH2	6:B:307[A]:DTO:S4	2.76	0.58
1:B:64:GLU:HG2	7:B:309:P33:H152	1.85	0.58
1:A:118:SER:HB3	1:A:212:ALA:HB1	1.87	0.56
1:A:60:THR:CA	9:A:428:HOH:O	2.47	0.56
1:A:144:MET:HA	8:B:310:PG4:H22	1.86	0.56
1:A:20:ILE:HG23	1:A:24:ILE:CD1	2.35	0.55
1:A:82:ASN:HB2	9:A:546:HOH:O	2.07	0.54
1:A:233:ASP:HB3	9:A:504:HOH:O	2.06	0.54
1:B:51:GLN:HE21	1:B:89:PRO:HG3	1.72	0.54
1:A:20:ILE:CG2	1:A:24:ILE:HD11	2.37	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:191:ARG:NH2	6:B:307[A]:DTO:S1	2.80	0.54
1:A:18:LYS:HE2	9:A:416:HOH:O	2.07	0.53
1:B:45:ASN:HD22	1:B:45:ASN:N	2.05	0.53
1:B:171:ASP:OD1	4:B:308:GOL:H31	2.07	0.53
1:A:165[A]:ASN:ND2	1:A:167:ASP:H	2.06	0.53
1:A:125:HIS:CE1	6:B:307[B]:DTO:O2	2.62	0.53
1:A:20:ILE:HG23	1:A:24:ILE:HD11	1.90	0.53
1:A:101:GLU:OE2	6:B:307[A]:DTO:S1	2.67	0.52
1:B:64:GLU:OE2	7:B:309:P33:H181	2.09	0.52
1:B:63:PRO:HB2	7:B:309:P33:H182	1.92	0.52
1:A:20:ILE:CG2	1:A:24:ILE:CD1	2.89	0.51
1:A:31:MET:HA	1:A:45:ASN:ND2	2.23	0.51
1:A:45:ASN:HD22	1:A:45:ASN:H	1.57	0.51
1:A:165[B]:ASN:HD22	1:A:167:ASP:H	1.58	0.51
1:A:171:ASP:CG	4:A:307:GOL:H11	2.31	0.51
1:A:5:THR:H	1:A:8:GLN:NE2	2.07	0.51
3:A:303:PO4:P	9:A:403:HOH:O	2.71	0.49
1:A:204:ASN:ND2	9:A:525:HOH:O	2.33	0.49
1:A:62:PHE:HB3	1:A:65:HIS:CG	2.48	0.49
1:B:78:THR:C	1:B:80:GLU:CB	2.81	0.48
1:A:165[A]:ASN:HD22	1:A:167:ASP:H	1.61	0.48
1:A:51:GLN:NE2	1:A:69:ALA:HB1	2.28	0.48
7:B:309:P33:H22A	7:B:309:P33:H52	1.66	0.47
1:A:60:THR:CB	1:A:61:TYR:CE2	2.96	0.47
1:A:125:HIS:NE2	6:B:307[A]:DTO:H11	2.30	0.46
1:A:165[A]:ASN:HD21	1:A:167:ASP:HB2	1.79	0.46
1:B:83:HIS:CE1	7:B:309:P33:H51	2.51	0.46
1:A:118:SER:HB3	1:A:212:ALA:CB	2.45	0.46
1:B:64:GLU:CB	7:B:309:P33:H92	2.30	0.46
1:B:65:HIS:HE2	7:B:309:P33:C5	2.29	0.45
1:A:194:LYS:HG2	8:B:310:PG4:H31	1.99	0.45
1:B:171:ASP:OD1	4:B:308:GOL:O3	2.31	0.45
1:A:165[A]:ASN:HD22	1:A:167:ASP:N	2.16	0.44
1:A:191:ARG:HH21	6:B:307[B]:DTO:H42	1.72	0.44
1:A:61:TYR:CD2	1:A:61:TYR:N	2.85	0.43
1:A:47:ASP:OD2	1:A:90:ILE:HB	2.19	0.43
1:B:63:PRO:HG2	7:B:309:P33:H122	2.00	0.43
1:B:45:ASN:N	1:B:45:ASN:ND2	2.67	0.42
1:B:31:MET:HA	1:B:45:ASN:HD21	1.84	0.42
1:B:240:PHE:CD2	1:B:240:PHE:C	2.92	0.42
8:B:310:PG4:H72	8:B:310:PG4:H51	1.47	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:26:GLN:O	1:A:30:GLU:HG2	2.20	0.42
1:A:191:ARG:HH22	6:B:307[B]:DTO:C4	2.27	0.42
1:B:65:HIS:HE2	7:B:309:P33:H51	1.85	0.42
1:A:166:LEU:HD12	1:A:166:LEU:HA	1.88	0.41
1:B:166:LEU:HD12	1:B:166:LEU:HA	1.95	0.41
1:A:191:ARG:HH22	6:B:307[A]:DTO:C4	2.33	0.41
1:A:86:ILE:HG21	1:A:211:ALA:HB2	2.02	0.41
1:B:171:ASP:OD2	4:B:308:GOL:H31	2.21	0.41
1:B:12:LEU:HB3	1:B:57:PHE:HZ	1.84	0.41
1:A:247:CYS:O	1:A:251:VAL:HG23	2.21	0.41
1:B:102:ASP:OD1	6:B:307[A]:DTO:C4	2.69	0.41
1:B:63:PRO:HD2	7:B:309:P33:C11	2.34	0.40
1:B:64:GLU:HG2	7:B:309:P33:H121	2.03	0.40
1:A:191:ARG:HH22	6:B:307[A]:DTO:H42	1.86	0.40
1:B:62:PHE:HB3	1:B:65:HIS:CG	2.57	0.40
1:B:191:ARG:NH2	6:B:307[B]:DTO:S4	2.94	0.40
1:B:32:THR:H	1:B:45:ASN:ND2	2.20	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	A	264/271 (97%)	258 (98%)	5 (2%)	1 (0%)	34	42
1	B	258/271 (95%)	253 (98%)	4 (2%)	1 (0%)	34	42
All	All	522/542 (96%)	511 (98%)	9 (2%)	2 (0%)	34	42

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	75	ALA

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Mol	Chain	Res	Type
1	A	3	ASP

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	228/238 (96%)	210 (92%)	18 (8%)	12	15
1	B	221/238 (93%)	198 (90%)	23 (10%)	7	8
All	All	449/476 (94%)	408 (91%)	41 (9%)	9	11

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	LEU
1	A	15	SER
1	A	17	LEU
1	A	18	LYS
1	A	20	ILE
1	A	29	MET
1	A	42	LEU
1	A	45	ASN
1	A	58	LEU
1	A	78	THR
1	A	79	ASN
1	A	133	ARG
1	A	158	SER
1	A	166	LEU
1	A	167	ASP
1	A	235	LEU
1	A	240	PHE
1	A	265	ARG
1	B	5	THR
1	B	11	LYS
1	B	12	LEU
1	B	15	SER

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Mol	Chain	Res	Type
1	B	17	LEU
1	B	18	LYS
1	B	19	GLN
1	B	42	LEU
1	B	45	ASN
1	B	47	ASP
1	B	77	ILE
1	B	81	ILE
1	B	107	LEU
1	B	126	LYS
1	B	149	SER
1	B	158	SER
1	B	163	VAL
1	B	166	LEU
1	B	170	GLN
1	B	235	LEU
1	B	240	PHE
1	B	263	LYS
1	B	265	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	8	GLN
1	A	45	ASN
1	A	51	GLN
1	A	82	ASN
1	A	125	HIS
1	A	170	GLN
1	A	248	HIS
1	B	45	ASN
1	B	51	GLN
1	B	53	GLN
1	B	82	ASN
1	B	125	HIS
1	B	204	ASN
1	B	248	HIS

5.3.3 RNA

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](#)

Of 18 ligands modelled in this entry, 8 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	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	PO4	B	301	2	4,4,4	1.20	0	6,6,6	1.19	2 (33%)
3	PO4	A	303	2	4,4,4	1.48	0	6,6,6	1.08	0
4	GOL	A	307	-	5,5,5	0.58	0	5,5,5	0.73	0
6	DTO	B	307[A]	-	7,7,8	0.52	0	4,8,9	2.02	2 (50%)
4	GOL	B	305	-	5,5,5	0.24	0	5,5,5	0.74	0
4	GOL	B	308	-	5,5,5	0.27	0	5,5,5	0.75	0
6	DTO	B	307[B]	-	7,7,8	0.57	0	4,8,9	1.05	0
7	P33	B	309	-	21,21,21	0.83	0	20,20,20	0.63	0
8	PG4	B	310	-	12,12,12	1.00	0	11,11,11	0.62	0
4	GOL	A	305	-	5,5,5	0.46	0	5,5,5	1.13	1 (20%)

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	GOL	A	307	-	-	4/4/4/4	-
6	DTO	B	307[A]	-	-	8/8/8/9	-
4	GOL	B	305	-	-	0/4/4/4	-
4	GOL	B	308	-	-	4/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	DTO	B	307[B]	-	-	8/8/8/9	-
7	P33	B	309	-	-	11/19/19/19	-
8	PG4	B	310	-	-	7/10/10/10	-
4	GOL	A	305	-	-	2/4/4/4	-

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	B	307[A]	DTO	O2-C2-C3	2.75	115.37	109.72
6	B	307[A]	DTO	C2-C1-S1	-2.54	107.10	114.47
4	A	305	GOL	O3-C3-C2	-2.30	99.15	110.20
3	B	301	PO4	O4-P-O1	-2.03	103.46	110.89
3	B	301	PO4	O4-P-O2	2.03	114.48	107.97

There are no chirality outliers.

All (44) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	307	GOL	O1-C1-C2-O2
4	B	308	GOL	C1-C2-C3-O3
6	B	307[A]	DTO	S1-C1-C2-O2
6	B	307[A]	DTO	S1-C1-C2-C3
6	B	307[A]	DTO	C1-C2-C3-O3
6	B	307[A]	DTO	C1-C2-C3-C4
6	B	307[A]	DTO	O2-C2-C3-O3
6	B	307[A]	DTO	O2-C2-C3-C4
6	B	307[A]	DTO	C2-C3-C4-S4
6	B	307[A]	DTO	O3-C3-C4-S4
6	B	307[B]	DTO	S1-C1-C2-O2
6	B	307[B]	DTO	S1-C1-C2-C3
6	B	307[B]	DTO	C1-C2-C3-O3
6	B	307[B]	DTO	C1-C2-C3-C4
6	B	307[B]	DTO	O2-C2-C3-C4
6	B	307[B]	DTO	C2-C3-C4-S4
6	B	307[B]	DTO	O3-C3-C4-S4
8	B	310	PG4	C5-C6-O4-C7
7	B	309	P33	C2-C3-O4-C5
7	B	309	P33	C18-C17-O16-C15
7	B	309	P33	O13-C14-C15-O16

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Mol	Chain	Res	Type	Atoms
7	B	309	P33	O4-C5-C6-O7
8	B	310	PG4	O3-C5-C6-O4
7	B	309	P33	O7-C8-C9-O10
7	B	309	P33	O16-C17-C18-O19
8	B	310	PG4	O4-C7-C8-O5
8	B	310	PG4	O2-C3-C4-O3
4	A	305	GOL	O1-C1-C2-C3
4	A	307	GOL	O1-C1-C2-C3
4	A	307	GOL	C1-C2-C3-O3
4	B	308	GOL	O1-C1-C2-C3
7	B	309	P33	C15-C14-O13-C12
4	B	308	GOL	O2-C2-C3-O3
7	B	309	P33	O10-C11-C12-O13
7	B	309	P33	O1-C2-C3-O4
6	B	307[B]	DTO	O2-C2-C3-O3
8	B	310	PG4	C3-C4-O3-C5
7	B	309	P33	C14-C15-O16-C17
8	B	310	PG4	C4-C3-O2-C2
7	B	309	P33	C8-C9-O10-C11
8	B	310	PG4	C8-C7-O4-C6
4	A	305	GOL	O1-C1-C2-O2
4	B	308	GOL	O1-C1-C2-O2
4	A	307	GOL	O2-C2-C3-O3

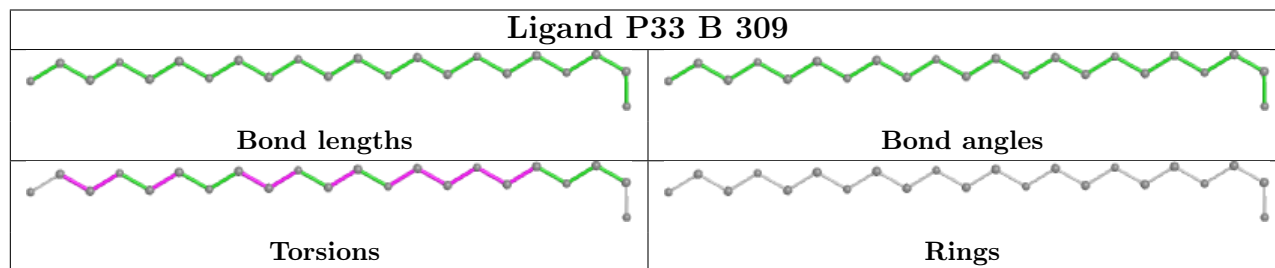
There are no ring outliers.

9 monomers are involved in 47 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	301	PO4	1	0
3	A	303	PO4	1	0
4	A	307	GOL	3	0
6	B	307[A]	DTO	10	0
4	B	308	GOL	5	0
6	B	307[B]	DTO	8	0
7	B	309	P33	14	0
8	B	310	PG4	4	0
4	A	305	GOL	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	264/271 (97%)	-0.52	1 (0%) 92 95	7, 19, 35, 51	0
1	B	261/271 (96%)	-0.41	2 (0%) 86 89	9, 21, 39, 65	0
All	All	525/542 (96%)	-0.46	3 (0%) 89 92	7, 20, 37, 65	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	78	THR	3.1
1	B	237	GLY	2.4
1	A	64	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 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, 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.

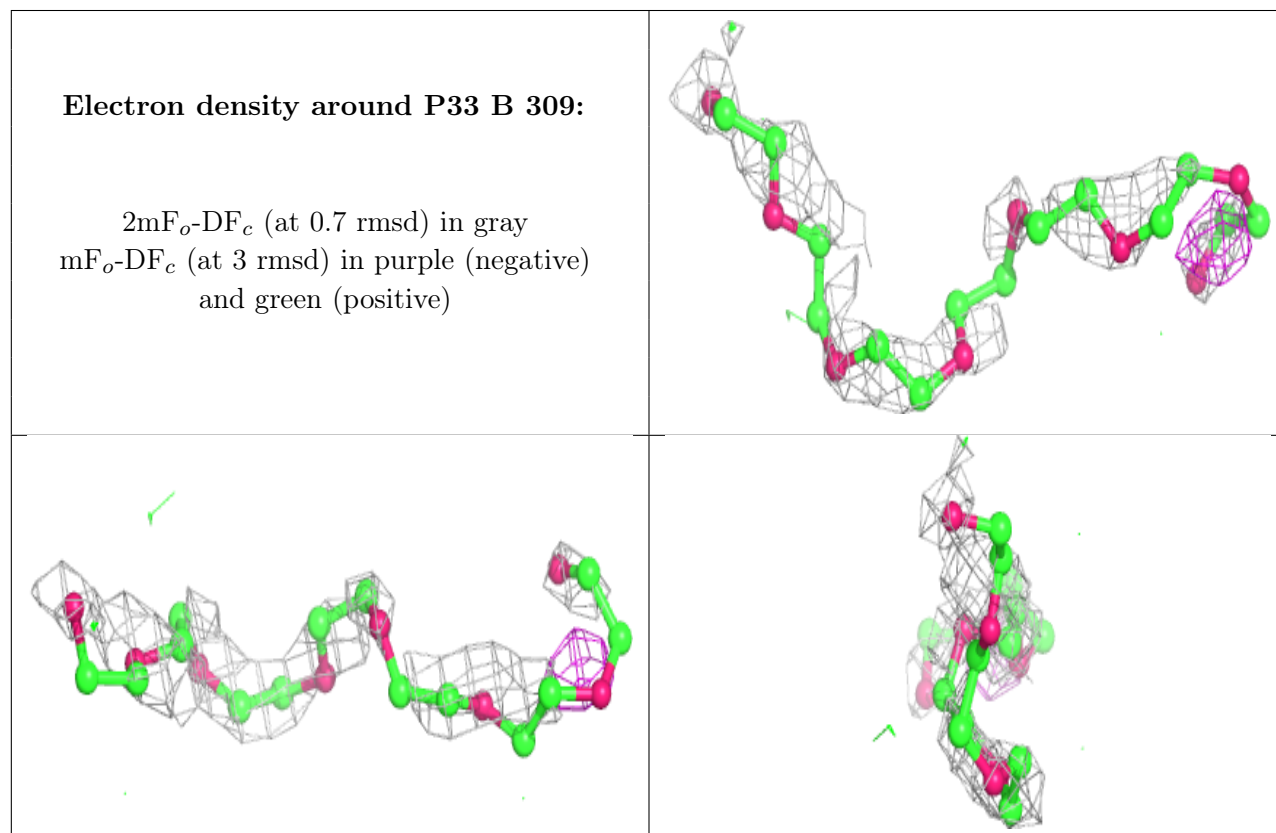
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
7	P33	B	309	22/22	0.65	0.43	70,76,78,78	0
8	PG4	B	310	13/13	0.68	0.31	45,58,63,63	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	DTO	B	307[A]	8/9	0.83	0.34	19,22,27,34	8
6	DTO	B	307[B]	8/9	0.83	0.34	41,45,45,48	8
4	GOL	A	307	6/6	0.84	0.22	40,43,44,44	0
4	GOL	B	308	6/6	0.88	0.19	29,31,33,33	0
4	GOL	B	305	6/6	0.91	0.16	28,38,38,39	0
4	GOL	A	305	6/6	0.93	0.16	33,35,37,40	0
2	MG	B	303	1/1	0.97	0.06	19,19,19,19	0
2	MG	A	301	1/1	0.98	0.07	22,22,22,22	0
5	CL	A	306	1/1	0.98	0.07	22,22,22,22	0
2	MG	B	304	1/1	0.99	0.09	22,22,22,22	0
3	PO4	A	303	5/5	0.99	0.07	14,15,19,20	0
5	CL	B	306	1/1	0.99	0.04	22,22,22,22	0
3	PO4	B	301	5/5	0.99	0.06	20,22,24,25	0
2	MG	A	304	1/1	0.99	0.04	16,16,16,16	0
2	MG	B	302	1/1	0.99	0.04	24,24,24,24	0
2	MG	A	302	1/1	0.99	0.07	15,15,15,15	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.