



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:25 PM UTC

PDB ID : 3T61 / pdb_00003t61
Title : Crystal Structure of a gluconokinase from Sinorhizobium meliloti 1021
Authors : Kumaran, D.; Chamala, S.; Evans, B.; Foti, R.; Gizzi, A.; Hillerich, B.; Kar, A.; LaFleur, J.; Seidel, R.; Villigas, G.; Zencheck, W.; Almo, S.C.; Swaminathan, S.; New York Structural Genomics Research Consortium (NYSGRG)
Deposited on : 2011-07-28
Resolution : 2.20 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

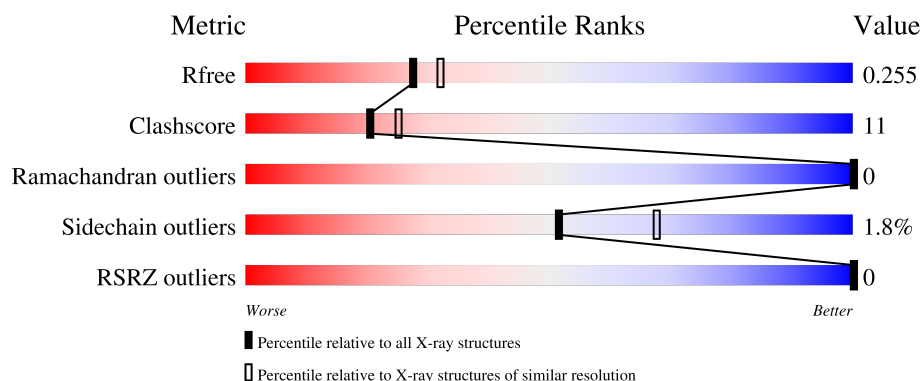
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	202	
1	B	202	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2860 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 Gluconokinase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	176	Total	C	N	O	S	Se	0	0	0
			1353	848	247	252	2	4			
1	B	175	Total	C	N	O	S	Se	0	0	0
			1343	842	244	251	2	4			

There are 48 discrepancies between the modelled and reference sequences:

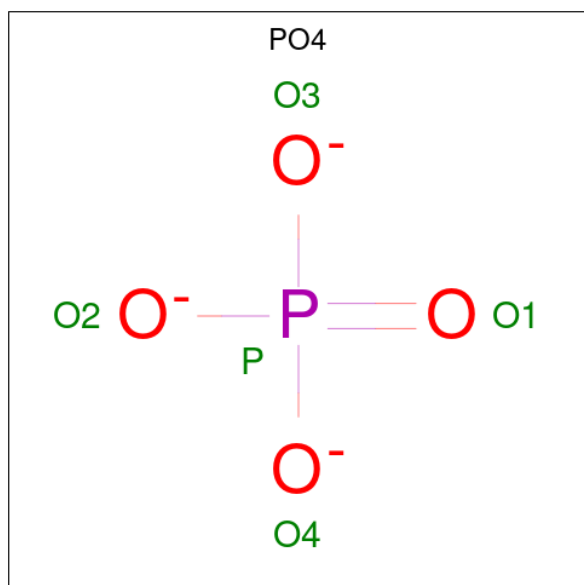
Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MSE	-	expression tag	UNP Q92VK3
A	0	VAL	-	expression tag	UNP Q92VK3
A	179	ALA	-	expression tag	UNP Q92VK3
A	180	GLU	-	expression tag	UNP Q92VK3
A	181	ASN	-	expression tag	UNP Q92VK3
A	182	LEU	-	expression tag	UNP Q92VK3
A	183	TYR	-	expression tag	UNP Q92VK3
A	184	PHE	-	expression tag	UNP Q92VK3
A	185	GLN	-	expression tag	UNP Q92VK3
A	186	SER	-	expression tag	UNP Q92VK3
A	187	HIS	-	expression tag	UNP Q92VK3
A	188	HIS	-	expression tag	UNP Q92VK3
A	189	HIS	-	expression tag	UNP Q92VK3
A	190	HIS	-	expression tag	UNP Q92VK3
A	191	HIS	-	expression tag	UNP Q92VK3
A	192	HIS	-	expression tag	UNP Q92VK3
A	193	TRP	-	expression tag	UNP Q92VK3
A	194	SER	-	expression tag	UNP Q92VK3
A	195	HIS	-	expression tag	UNP Q92VK3
A	196	PRO	-	expression tag	UNP Q92VK3
A	197	GLN	-	expression tag	UNP Q92VK3
A	198	PHE	-	expression tag	UNP Q92VK3
A	199	GLU	-	expression tag	UNP Q92VK3
A	200	LYS	-	expression tag	UNP Q92VK3
B	-1	MSE	-	expression tag	UNP Q92VK3

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Chain	Residue	Modelled	Actual	Comment	Reference
B	0	VAL	-	expression tag	UNP Q92VK3
B	179	ALA	-	expression tag	UNP Q92VK3
B	180	GLU	-	expression tag	UNP Q92VK3
B	181	ASN	-	expression tag	UNP Q92VK3
B	182	LEU	-	expression tag	UNP Q92VK3
B	183	TYR	-	expression tag	UNP Q92VK3
B	184	PHE	-	expression tag	UNP Q92VK3
B	185	GLN	-	expression tag	UNP Q92VK3
B	186	SER	-	expression tag	UNP Q92VK3
B	187	HIS	-	expression tag	UNP Q92VK3
B	188	HIS	-	expression tag	UNP Q92VK3
B	189	HIS	-	expression tag	UNP Q92VK3
B	190	HIS	-	expression tag	UNP Q92VK3
B	191	HIS	-	expression tag	UNP Q92VK3
B	192	HIS	-	expression tag	UNP Q92VK3
B	193	TRP	-	expression tag	UNP Q92VK3
B	194	SER	-	expression tag	UNP Q92VK3
B	195	HIS	-	expression tag	UNP Q92VK3
B	196	PRO	-	expression tag	UNP Q92VK3
B	197	GLN	-	expression tag	UNP Q92VK3
B	198	PHE	-	expression tag	UNP Q92VK3
B	199	GLU	-	expression tag	UNP Q92VK3
B	200	LYS	-	expression tag	UNP Q92VK3

- Molecule 2 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



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

- Molecule 3 is water.

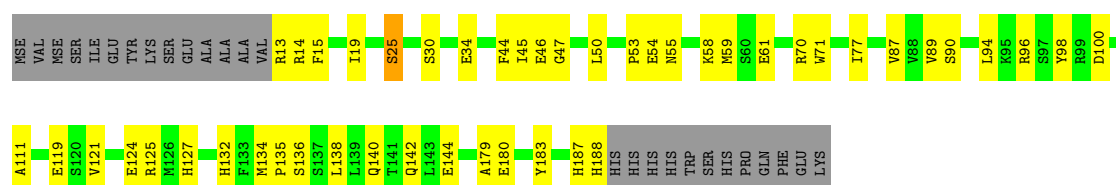
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	78	Total	O	0	0
			78	78		
3	B	76	Total	O	0	0
			76	76		

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.

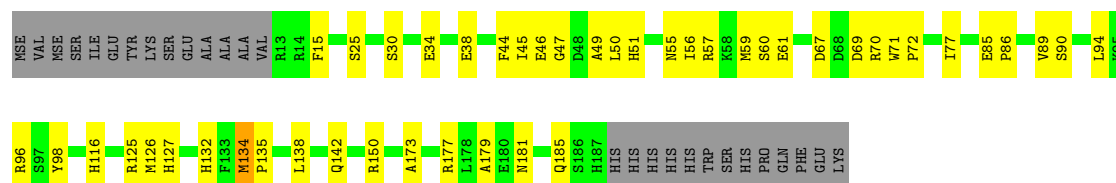
• Molecule 1: Gluconokinase

Chain A: 



• Molecule 1: Gluconokinase

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 43	Depositor
Cell constants a, b, c, α , β , γ	68.20Å 68.20Å 118.62Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.68 – 2.20 44.68 – 2.20	Depositor EDS
% Data completeness (in resolution range)	97.6 (44.68-2.20) 97.6 (44.68-2.20)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	9.63 (at 2.20Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.237 , 0.250 0.237 , 0.255	Depositor DCC
R_{free} test set	1860 reflections (6.71%)	wwPDB-VP
Wilson B-factor (Å ²)	32.1	Xtriage
Anisotropy	0.107	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 23.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.488 for h,-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	2860	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 4.71% 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: PO4

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	0.46	0/1379	0.93	4/1861 (0.2%)
1	B	0.46	0/1368	0.94	3/1846 (0.2%)
All	All	0.46	0/2747	0.94	7/3707 (0.2%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	94	LEU	N-CA-C	8.15	120.16	111.28
1	A	94	LEU	N-CA-C	7.89	119.88	111.28
1	B	70	ARG	N-CA-C	6.52	118.47	111.36
1	A	70	ARG	N-CA-C	6.52	118.47	111.36
1	A	71	TRP	CA-C-N	-5.04	113.85	119.19
1	A	71	TRP	C-N-CA	-5.04	113.85	119.19
1	B	134	MSE	N-CA-C	-5.02	103.45	109.83

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	1353	0	1338	37	0
1	B	1343	0	1331	29	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	5	0	0	0	0
2	B	5	0	0	0	0
3	A	78	0	0	2	0
3	B	76	0	0	1	0
All	All	2860	0	2669	60	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 11.

All (60) 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:13:ARG:HH11	1:A:180:GLU:HA	1.20	1.01
1:A:13:ARG:NH1	1:A:180:GLU:HA	1.76	1.00
1:B:25:SER:H	1:B:142:GLN:HE21	1.23	0.86
1:A:25:SER:H	1:A:142:GLN:HE21	1.18	0.86
1:A:124:GLU:HG2	3:A:226:HOH:O	1.93	0.69
1:B:30:SER:O	1:B:34:GLU:HG3	1.93	0.68
1:B:135:PRO:HD2	1:B:138:LEU:HD12	1.76	0.66
1:B:56:ILE:HA	1:B:59:MSE:HE3	1.77	0.66
1:A:135:PRO:HD2	1:A:138:LEU:HD12	1.79	0.65
1:A:61:GLU:O	1:B:134:MSE:HG2	1.96	0.65
1:B:55:ASN:HB3	1:B:59:MSE:HE2	1.78	0.64
1:A:134:MSE:HG2	1:B:61:GLU:O	1.97	0.64
1:A:13:ARG:NH1	1:A:183:TYR:HB2	2.13	0.63
1:A:25:SER:H	1:A:142:GLN:NE2	1.94	0.62
1:A:136:SER:O	1:A:140:GLN:HG3	2.01	0.60
1:A:179:ALA:O	1:A:183:TYR:HD1	1.84	0.60
1:B:45:ILE:HB	1:B:89:VAL:HG22	1.85	0.59
1:A:140:GLN:O	1:A:144:GLU:HG3	2.04	0.58
1:A:180:GLU:HB3	3:A:221:HOH:O	2.04	0.56
1:A:119:GLU:OE1	1:A:140:GLN:HG2	2.07	0.55
1:B:96:ARG:HG3	1:B:96:ARG:HH11	1.70	0.55
1:B:50:LEU:HD12	1:B:77:ILE:HD13	1.88	0.54
1:B:51:HIS:CG	1:B:59:MSE:HE1	2.43	0.54
1:A:121:VAL:O	1:A:125:ARG:HG3	2.08	0.53
1:B:116:HIS:HE1	3:B:260:HOH:O	1.91	0.53
1:B:34:GLU:HG2	1:B:44:PHE:CD2	2.44	0.52
1:A:30:SER:O	1:A:34:GLU:HG3	2.10	0.51
1:A:50:LEU:HD12	1:A:77:ILE:HD13	1.92	0.51
1:A:183:TYR:O	1:A:187:HIS:HD2	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:173:ALA:O	1:B:177:ARG:HG3	2.10	0.50
1:A:15:PHE:CD2	1:A:19:ILE:HD11	2.47	0.49
1:A:45:ILE:HD12	1:A:87:VAL:HG21	1.94	0.49
1:A:59:MSE:HE2	1:B:60:SER:OG	2.12	0.48
1:B:15:PHE:HA	1:B:179:ALA:HB2	1.96	0.48
1:A:34:GLU:HG2	1:A:44:PHE:CD2	2.49	0.47
1:B:47:GLY:HA3	1:B:90:SER:O	2.14	0.47
1:B:56:ILE:CA	1:B:59:MSE:HE3	2.44	0.47
1:A:187:HIS:O	1:A:188:HIS:C	2.57	0.47
1:A:19:ILE:CD1	1:A:111:ALA:HB3	2.46	0.47
1:A:46:GLU:OE1	1:B:57:ARG:NH1	2.48	0.46
1:A:47:GLY:HA3	1:A:90:SER:O	2.15	0.46
1:A:13:ARG:HH12	1:A:183:TYR:HB2	1.80	0.45
1:A:53:PRO:HG3	1:B:49:ALA:HB2	1.98	0.44
1:A:19:ILE:HD13	1:A:111:ALA:HB3	2.00	0.43
1:B:181:ASN:O	1:B:185:GLN:HG3	2.17	0.43
1:A:45:ILE:HB	1:A:89:VAL:HG22	1.98	0.43
1:A:53:PRO:HB3	1:B:46:GLU:HG2	2.02	0.42
1:A:98:TYR:N	1:A:98:TYR:CD1	2.87	0.42
1:B:85:GLU:HA	1:B:86:PRO:HD3	1.91	0.42
1:B:127:HIS:HA	1:B:132:HIS:O	2.20	0.42
1:A:96:ARG:HD2	1:A:100:ASP:OD2	2.19	0.42
1:B:55:ASN:C	1:B:59:MSE:HE3	2.44	0.42
1:B:69:ASP:O	1:B:72:PRO:HD2	2.20	0.42
1:B:125:ARG:NH1	1:B:126:MSE:CE	2.83	0.42
1:A:55:ASN:O	1:A:59:MSE:HG3	2.20	0.41
1:A:127:HIS:HA	1:A:132:HIS:O	2.20	0.41
1:A:13:ARG:HB3	1:A:14:ARG:H	1.44	0.41
1:A:54:GLU:OE2	1:A:58:LYS:HE3	2.22	0.40
1:B:98:TYR:N	1:B:98:TYR:CD1	2.88	0.40
1:B:56:ILE:N	1:B:59:MSE:HE3	2.37	0.40

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	174/202 (86%)	168 (97%)	6 (3%)	0	100	100
1	B	173/202 (86%)	169 (98%)	4 (2%)	0	100	100
All	All	347/404 (86%)	337 (97%)	10 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	143/160 (89%)	142 (99%)	1 (1%)	76	87
1	B	142/160 (89%)	138 (97%)	4 (3%)	38	52
All	All	285/320 (89%)	280 (98%)	5 (2%)	51	68

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	SER
1	B	38	GLU
1	B	67	ASP
1	B	71	TRP
1	B	150	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	127	HIS
1	A	128	HIS
1	A	142	GLN
1	A	162	GLN
1	A	187	HIS

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Mol	Chain	Res	Type
1	B	116	HIS
1	B	140	GLN
1	B	142	GLN
1	B	162	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 oligosaccharides 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	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	PO4	A	300	-	4,4,4	3.89	3 (75%)	6,6,6	1.05	0
2	PO4	B	300	-	4,4,4	4.00	3 (75%)	6,6,6	0.89	0

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	300	PO4	P-O1	7.08	1.67	1.50
2	A	300	PO4	P-O1	6.78	1.66	1.50
2	A	300	PO4	P-O2	2.88	1.63	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	300	PO4	P-O2	2.55	1.62	1.54
2	B	300	PO4	P-O4	2.53	1.62	1.54
2	A	300	PO4	P-O4	2.37	1.61	1.54

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	172/202 (85%)	-1.29	0 100 100	24, 35, 58, 73	0
1	B	171/202 (84%)	-1.29	0 100 100	23, 35, 58, 71	0
All	All	343/404 (84%)	-1.29	0 100 100	23, 35, 58, 73	0

There are no RSRZ outliers to report.

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q < 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	PO4	A	300	5/5	1.00	0.03	28,31,34,34	0
2	PO4	B	300	5/5	1.00	0.03	26,29,33,33	0

6.5 Other polymers [i](#)

There are no such residues in this entry.