



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 9, 2026 – 02:14 PM UTC

PDB ID : 6R2N / pdb_00006r2n
Title : Crystal structure of KIGlk1 glucokinase from Kluyveromyces lactis
Authors : Zak, K.; Wator, E.; Grudnik, P.
Deposited on : 2019-03-18
Resolution : 2.60 Å(reported)

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

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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

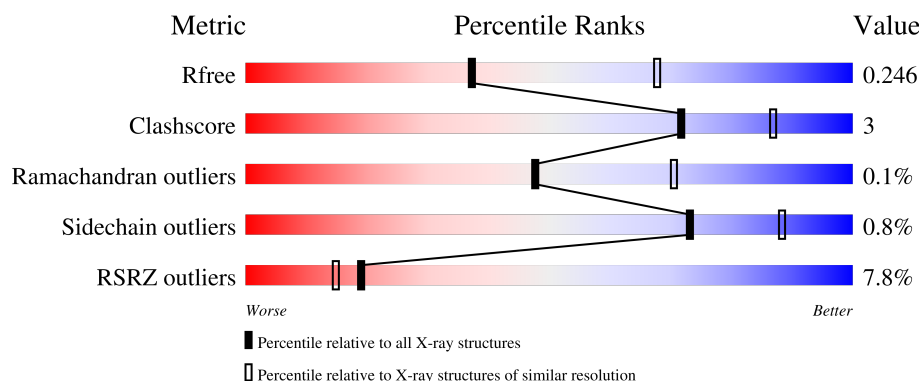
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	475	3% 92% 8%
1	B	475	13% 91% 8%
1	C	475	7% 92% 7%

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	EDO	B	503	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 21946 atoms, of which 10674 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.

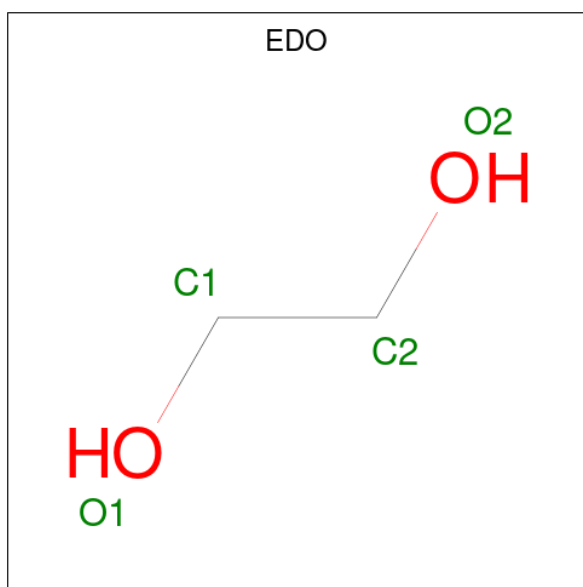
- Molecule 1 is a protein called Glucokinase-1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	475	Total	C	H	N	O	S	0	8	0
			7506	2388	3726	647	724	21			
1	B	474	Total	C	H	N	O	S	12	4	0
			7041	2279	3432	614	696	20			
1	C	475	Total	C	H	N	O	S	0	4	0
			7092	2303	3450	616	703	20			

There are 15 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	?	-	SER	deletion	UNP Q6CUZ3
A	?	-	ASP	deletion	UNP Q6CUZ3
A	?	-	ILE	deletion	UNP Q6CUZ3
A	?	-	SER	deletion	UNP Q6CUZ3
A	?	-	GLN	deletion	UNP Q6CUZ3
B	?	-	SER	deletion	UNP Q6CUZ3
B	?	-	ASP	deletion	UNP Q6CUZ3
B	?	-	ILE	deletion	UNP Q6CUZ3
B	?	-	SER	deletion	UNP Q6CUZ3
B	?	-	GLN	deletion	UNP Q6CUZ3
C	?	-	SER	deletion	UNP Q6CUZ3
C	?	-	ASP	deletion	UNP Q6CUZ3
C	?	-	ILE	deletion	UNP Q6CUZ3
C	?	-	SER	deletion	UNP Q6CUZ3
C	?	-	GLN	deletion	UNP Q6CUZ3

- Molecule 2 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	H	O	0	0
			10	2	6	2		
2	A	1	Total	C	H	O	0	0
			10	2	6	2		
2	A	1	Total	C	H	O	0	0
			10	2	6	2		
2	A	1	Total	C	H	O	0	0
			10	2	6	2		
2	A	1	Total	C	H	O	0	0
			10	2	6	2		
2	A	1	Total	C	H	O	0	0
			10	2	6	2		
2	B	1	Total	C	H	O	0	0
			10	2	6	2		
2	B	1	Total	C	H	O	0	0
			10	2	6	2		
2	B	1	Total	C	H	O	0	0
			10	2	6	2		

- Molecule 3 is BROMIDE ION (CCD ID: BR) (formula: Br).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	1	Total	Br	0	0
			1	1		

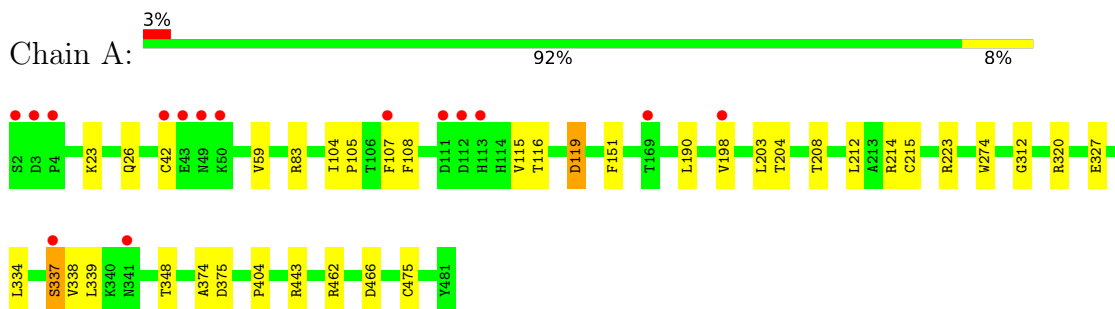
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	127	Total 127	O 127	0	0
4	B	35	Total 35	O 35	0	0
4	C	34	Total 34	O 34	0	0

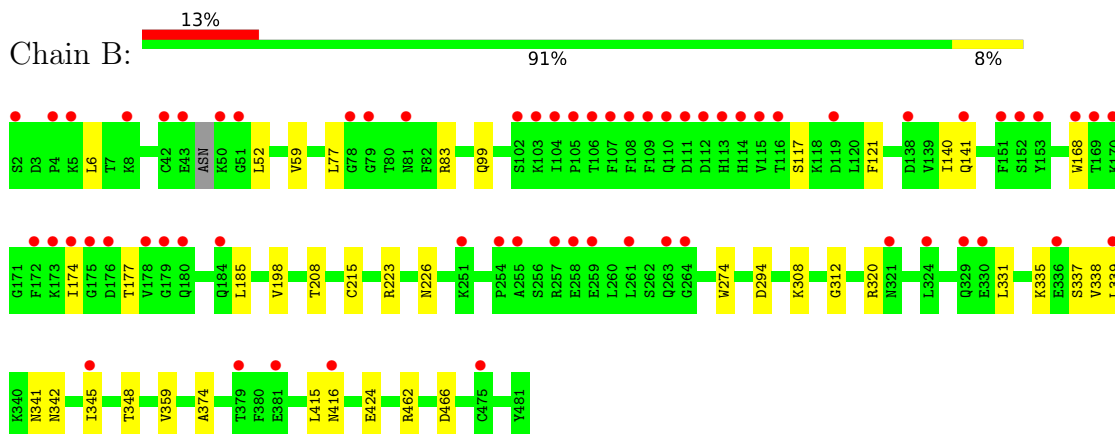
3 Residue-property plots [i](#)

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.

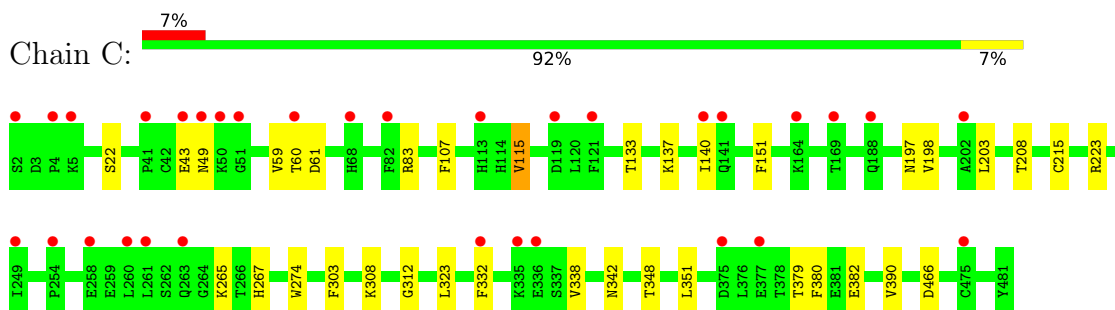
• Molecule 1: Glucokinase-1



• Molecule 1: Glucokinase-1



• Molecule 1: Glucokinase-1



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	102.01Å 122.77Å 360.35Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.08 – 2.60 49.08 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.0 (49.08-2.60) 99.6 (49.08-2.60)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.75 (at 2.61Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.204 , 0.243 0.214 , 0.246	Depositor DCC
R_{free} test set	1099 reflections (1.57%)	wwPDB-VP
Wilson B-factor (Å ²)	59.6	Xtriage
Anisotropy	0.308	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 51.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	21946	wwPDB-VP
Average B, all atoms (Å ²)	83.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.10% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: EDO, BR

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	0.14	0/3863	0.35	0/5227
1	B	0.12	0/3697	0.32	0/5023
1	C	0.11	0/3730	0.30	0/5067
All	All	0.13	0/11290	0.32	0/15317

There are no bond length outliers.

There are no bond angle outliers.

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	3780	3726	3721	27	0
1	B	3609	3432	3421	21	0
1	C	3642	3450	3460	22	0
2	A	32	48	48	1	0
2	B	12	18	18	2	0
3	C	1	0	0	0	0
4	A	127	0	0	1	0
4	B	35	0	0	0	0
4	C	34	0	0	0	0
All	All	11272	10674	10668	71	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 71 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:C:215:CYS:O	1:C:223:ARG:NH2	1.82	1.09
1:A:334:LEU:O	1:A:337:SER:OG	1.81	0.96
1:B:215:CYS:O	1:B:223:ARG:NH2	2.01	0.92
1:C:303:PHE:O	1:C:308:LYS:NZ	2.05	0.89
1:A:215:CYS:O	1:A:223[A]:ARG:NH2	2.14	0.80

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	479/475 (101%)	463 (97%)	16 (3%)	0	100	100
1	B	474/475 (100%)	458 (97%)	15 (3%)	1 (0%)	43	66
1	C	474/475 (100%)	462 (98%)	12 (2%)	0	100	100
All	All	1427/1425 (100%)	1383 (97%)	43 (3%)	1 (0%)	48	70

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	52	LEU

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	420/416 (101%)	417 (99%)	3 (1%)	76	89
1	B	383/416 (92%)	380 (99%)	3 (1%)	73	88
1	C	389/416 (94%)	386 (99%)	3 (1%)	73	88
All	All	1192/1248 (96%)	1183 (99%)	9 (1%)	73	88

5 of 9 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	115	VAL
1	C	198	VAL
1	B	77	LEU
1	B	198	VAL
1	B	359	VAL

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	444	HIS
1	B	364	GLN
1	C	460	HIS

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](#)

Of 12 ligands modelled in this entry, 1 is monoatomic - leaving 11 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	$\# Z > 2$
2	EDO	A	603	-	3,3,3	0.42	0	2,2,2	0.40	0
2	EDO	A	602	-	3,3,3	0.43	0	2,2,2	0.29	0
2	EDO	A	604	-	3,3,3	0.33	0	2,2,2	0.53	0
2	EDO	A	605	-	3,3,3	0.39	0	2,2,2	0.54	0
2	EDO	A	601	-	3,3,3	0.44	0	2,2,2	0.39	0
2	EDO	B	502	-	3,3,3	0.39	0	2,2,2	0.44	0
2	EDO	A	606	-	3,3,3	0.42	0	2,2,2	0.46	0
2	EDO	B	501	-	3,3,3	0.41	0	2,2,2	0.38	0
2	EDO	A	607	-	3,3,3	0.42	0	2,2,2	0.34	0
2	EDO	B	503	-	3,3,3	0.43	0	2,2,2	0.34	0
2	EDO	A	608	-	3,3,3	0.53	0	2,2,2	0.13	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
2	EDO	A	603	-	-	1/1/1/1	-
2	EDO	A	602	-	-	0/1/1/1	-
2	EDO	A	604	-	-	0/1/1/1	-
2	EDO	A	605	-	-	1/1/1/1	-
2	EDO	A	601	-	-	1/1/1/1	-
2	EDO	B	502	-	-	1/1/1/1	-
2	EDO	A	606	-	-	1/1/1/1	-
2	EDO	B	501	-	-	1/1/1/1	-
2	EDO	A	607	-	-	0/1/1/1	-
2	EDO	B	503	-	-	0/1/1/1	-
2	EDO	A	608	-	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

5 of 7 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	603	EDO	O1-C1-C2-O2
2	A	606	EDO	O1-C1-C2-O2
2	A	608	EDO	O1-C1-C2-O2
2	B	502	EDO	O1-C1-C2-O2
2	A	605	EDO	O1-C1-C2-O2

There are no ring outliers.

4 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	604	EDO	1	0
2	B	502	EDO	2	0
2	B	501	EDO	1	0
2	B	503	EDO	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	1
1	C	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	43:GLU	C	49:ASN	N	11.56
1	C	43:GLU	C	49:ASN	N	3.27

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	475/475 (100%)	0.05	15 (3%)	50	44	24, 56, 102, 147	8 (1%)
1	B	474/475 (99%)	0.74	64 (13%)	7	5	35, 92, 150, 225	3 (0%)
1	C	475/475 (100%)	0.65	32 (6%)	24	19	52, 96, 133, 162	4 (0%)
All	All	1424/1425 (99%)	0.48	111 (7%)	19	15	24, 83, 134, 225	15 (1%)

The worst 5 of 111 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	107	PHE	7.6
1	B	113	HIS	7.4
1	B	111	ASP	7.4
1	C	49	ASN	6.7
1	B	116	THR	6.6

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($\text{\AA}^2$)	Q<0.9
2	EDO	B	503	4/4	0.52	0.67	151,181,194,196	0
2	EDO	B	501	4/4	0.72	0.36	122,146,151,154	0
2	EDO	A	601	4/4	0.77	0.21	70,84,88,94	0
2	EDO	A	608	4/4	0.80	0.24	38,59,71,71	0
2	EDO	A	606	4/4	0.86	0.19	66,79,87,89	0
2	EDO	B	502	4/4	0.88	0.40	130,156,167,174	0
2	EDO	A	603	4/4	0.88	0.16	57,68,72,74	0
2	EDO	A	604	4/4	0.89	0.20	32,46,69,83	0
3	BR	C	501	1/1	0.89	0.30	112,112,112,112	0
2	EDO	A	602	4/4	0.90	0.20	82,98,106,106	0
2	EDO	A	605	4/4	0.93	0.16	34,51,61,73	0
2	EDO	A	607	4/4	0.96	0.19	84,101,102,103	0

6.5 Other polymers [i](#)

There are no such residues in this entry.