



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 11:39 AM GMT

PDB ID : 3PL2  
Title : Crystal structure of a 5-keto-2-deoxygluconokinase (NCgl0155, Cgl0158) from *Corynebacterium glutamicum* ATCC 13032 KITASATO at 1.89 Å resolution  
Authors : Joint Center for Structural Genomics (JCSG)  
Deposited on : 2010-11-12  
Resolution : 1.89 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](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

---

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

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

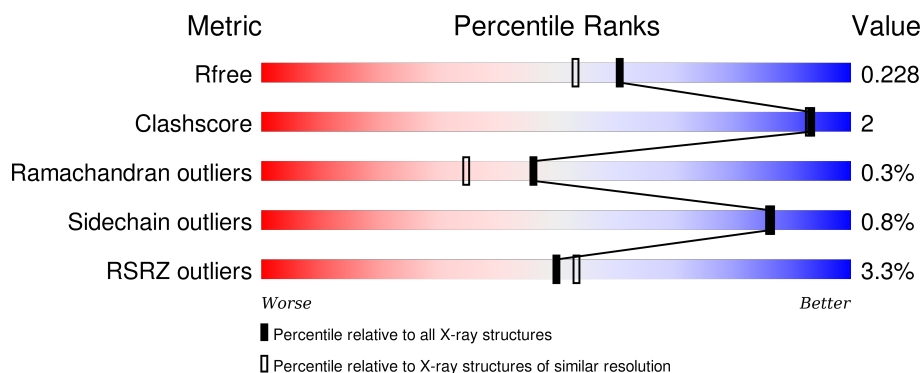
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.89 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))
$R_{free}$	91344	4755 (1.90-1.90)
Clashscore	102246	5398 (1.90-1.90)
Ramachandran outliers	100387	5338 (1.90-1.90)
Sidechain outliers	100360	5339 (1.90-1.90)
RSRZ outliers	91569	4766 (1.90-1.90)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	319	<div> <div>4%</div> <div>91%</div> <div>6% . .</div> </div>
1	B	319	<div> <div>3%</div> <div>91%</div> <div>6% .</div> </div>
1	C	319	<div> <div>2%</div> <div>92%</div> <div>6% .</div> </div>
1	D	319	<div> <div>4%</div> <div>90%</div> <div>7% .</div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10314 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 Sugar kinase, ribokinase family.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	310	Total	C	N	O	S	Se	0	8	0
			2428	1539	409	470	4	6			
1	B	310	Total	C	N	O	S	Se	0	14	0
			2443	1547	408	478	4	6			
1	C	310	Total	C	N	O	S	Se	0	12	0
			2445	1547	413	475	4	6			
1	D	309	Total	C	N	O	S	Se	0	7	0
			2408	1523	407	467	4	7			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	leader sequence	UNP Q8NTZ3
B	0	GLY	-	leader sequence	UNP Q8NTZ3
C	0	GLY	-	leader sequence	UNP Q8NTZ3
D	0	GLY	-	leader sequence	UNP Q8NTZ3

- Molecule 2 is CITRIC ACID (three-letter code: CIT) (formula: C<sub>6</sub>H<sub>8</sub>O<sub>7</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	C	O	0	0
			13	6	7		

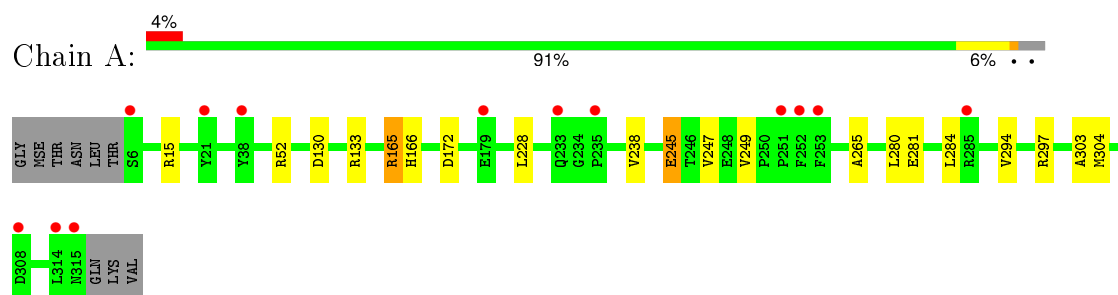
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	128	Total	O	0	0
			128	128		
3	B	164	Total	O	0	0
			164	164		
3	C	162	Total	O	0	0
			162	162		
3	D	123	Total	O	0	0
			123	123		

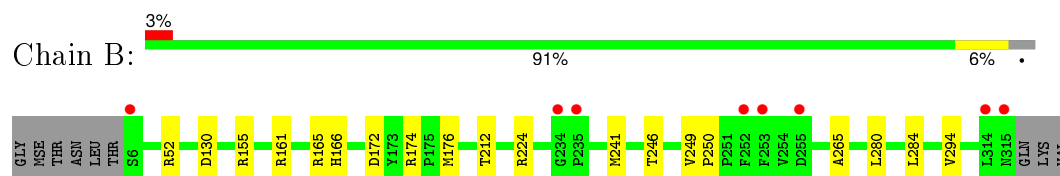
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

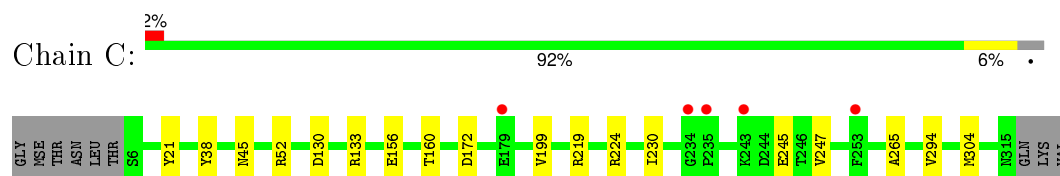
- Molecule 1: Sugar kinase, ribokinase family



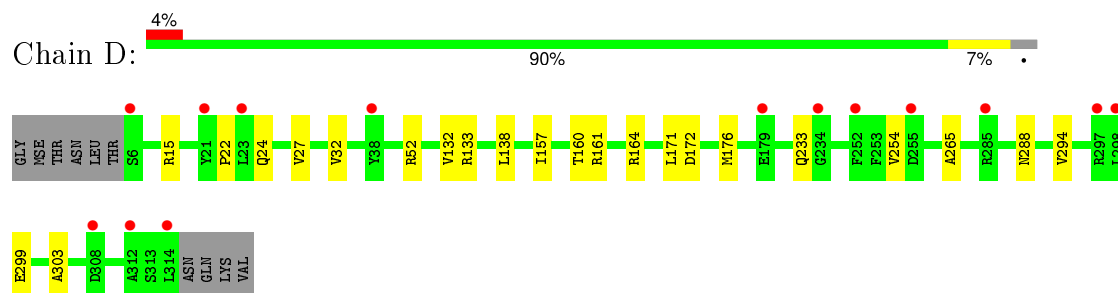
- Molecule 1: Sugar kinase, ribokinase family



- Molecule 1: Sugar kinase, ribokinase family



- Molecule 1: Sugar kinase, ribokinase family



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	45.91Å 79.49Å 81.70Å 101.86° 95.72° 92.11°	Depositor
Resolution (Å)	37.80 – 1.89 37.80 – 1.89	Depositor EDS
% Data completeness (in resolution range)	97.0 (37.80-1.89) 90.2 (37.80-1.89)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.07	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.10 (at 1.89Å)	Xtriage
Refinement program	REFMAC 5.5.0110	Depositor
R, $R_{free}$	0.180 , 0.225 0.186 , 0.228	Depositor DCC
$R_{free}$ test set	4384 reflections (5.29%)	DCC
Wilson B-factor (Å <sup>2</sup> )	25.3	Xtriage
Anisotropy	0.078	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 36.8	EDS
Estimated twinning fraction	0.006 for -h,-l,-k	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 87261 reflections	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	10314	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.42% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: CIT

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	0.74	0/2505	0.80	3/3401 (0.1%)
1	B	0.79	0/2533	0.84	4/3438 (0.1%)
1	C	0.75	0/2534	0.81	5/3442 (0.1%)
1	D	0.74	0/2480	0.81	3/3366 (0.1%)
All	All	0.76	0/10052	0.81	15/13647 (0.1%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	D	0	1

There are no bond length outliers.

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	161	ARG	NE-CZ-NH1	6.60	123.60	120.30
1	C	133	ARG	NE-CZ-NH1	6.14	123.37	120.30
1	A	15	ARG	NE-CZ-NH1	5.77	123.19	120.30
1	C	224	ARG	NE-CZ-NH1	5.75	123.17	120.30
1	B	52	ARG	NE-CZ-NH1	5.49	123.04	120.30
1	C	219	ARG	NE-CZ-NH1	5.48	123.04	120.30
1	D	15	ARG	NE-CZ-NH1	5.44	123.02	120.30
1	B	130	ASP	CB-CG-OD1	5.42	123.17	118.30
1	A	165	ARG	NE-CZ-NH1	5.31	122.95	120.30
1	D	133	ARG	NE-CZ-NH1	5.20	122.90	120.30
1	D	52	ARG	NE-CZ-NH1	5.14	122.87	120.30
1	A	15	ARG	NE-CZ-NH2	-5.13	117.73	120.30

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	52	ARG	NE-CZ-NH1	5.11	122.86	120.30
1	C	224	ARG	NE-CZ-NH2	-5.10	117.75	120.30
1	B	155	ARG	NE-CZ-NH1	5.08	122.84	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	171	LEU	Peptide

## 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	2428	0	2369	10	0
1	B	2443	0	2387	9	0
1	C	2445	0	2386	6	0
1	D	2408	0	2357	12	0
2	B	13	0	5	0	0
3	A	128	0	0	0	0
3	B	164	0	0	0	0
3	C	162	0	0	0	0
3	D	123	0	0	0	0
All	All	10314	0	9504	37	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 2.

All (37) 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:D:27[B]:VAL:HG13	1:D:32:VAL:HG22	1.70	0.73
1:B:280:LEU:O	1:B:284:LEU:HD13	1.96	0.65
1:D:27[A]:VAL:HG23	1:D:32:VAL:HG22	1.83	0.58
1:C:156:GLU:O	1:C:160:THR:HG23	2.06	0.56

*Continued on next page...*



*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:238:VAL:HG22	1:A:284:LEU:HD11	1.89	0.54
1:C:265:ALA:HB2	1:C:294:VAL:HG21	1.90	0.54
1:A:249:VAL:HG23	1:A:284:LEU:HD23	1.89	0.54
1:C:21[B]:TYR:CE1	1:C:38:TYR:CE2	2.95	0.53
1:B:265:ALA:HB2	1:B:294:VAL:HG21	1.91	0.52
1:A:245:GLU:HG3	1:A:247:VAL:HG23	1.92	0.51
1:A:247:VAL:HG11	1:A:281:GLU:HA	1.94	0.49
1:A:265:ALA:HB2	1:A:294:VAL:HG21	1.93	0.49
1:D:27[B]:VAL:HG13	1:D:32:VAL:CG2	2.41	0.49
1:B:176[B]:MSE:H	1:B:176[B]:MSE:HE3	1.80	0.46
1:C:199:VAL:HG22	1:C:230:ILE:HD12	1.98	0.46
1:B:241[B]:MSE:HG2	1:B:246:THR:HG23	1.97	0.45
1:D:27[B]:VAL:CG1	1:D:32:VAL:HG22	2.41	0.45
1:D:299:GLU:HB2	1:D:303:ALA:HB2	2.00	0.44
1:D:265:ALA:HB2	1:D:294:VAL:HG21	1.99	0.44
1:D:132:VAL:CG1	1:D:157:ILE:HG23	2.48	0.44
1:D:22:PRO:HB3	1:D:32:VAL:HG11	2.00	0.43
1:A:228:LEU:HD11	1:A:280:LEU:HD11	1.99	0.43
1:D:176[C]:MSE:HE2	1:D:176[C]:MSE:HB2	1.72	0.43
1:D:161:ARG:NH2	1:D:164:ARG:HD3	2.34	0.42
1:C:245:GLU:HG2	1:C:247:VAL:HG23	2.01	0.42
1:B:249:VAL:HG12	1:B:250:PRO:O	2.20	0.42
1:B:224[A]:ARG:HG3	1:B:224[A]:ARG:NH1	2.35	0.42
1:B:224[A]:ARG:HG3	1:B:224[A]:ARG:HH11	1.85	0.41
1:D:24:GLN:O	1:D:27[B]:VAL:HG12	2.20	0.41
1:B:174[B]:ARG:HE	1:B:176[B]:MSE:SE	2.54	0.41
1:B:165:ARG:HA	1:B:166:HIS:HA	1.92	0.41
1:A:52:ARG:HG3	1:A:304:MSE:HB2	2.02	0.40
1:A:165:ARG:HA	1:A:166:HIS:HA	1.96	0.40
1:A:130:ASP:OD1	1:A:133[A]:ARG:NH2	2.54	0.40
1:D:24:GLN:HB2	1:D:27[B]:VAL:HG11	2.03	0.40
1:C:45:ASN:HB3	1:C:304:MSE:HE3	2.03	0.40
1:A:297[A]:ARG:HD2	1:A:303:ALA:HB1	2.04	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	317/319 (99%)	311 (98%)	5 (2%)	1 (0%)	46	35
1	B	322/319 (101%)	316 (98%)	5 (2%)	1 (0%)	46	35
1	C	321/319 (101%)	317 (99%)	3 (1%)	1 (0%)	46	35
1	D	315/319 (99%)	311 (99%)	3 (1%)	1 (0%)	46	35
All	All	1275/1276 (100%)	1255 (98%)	16 (1%)	4 (0%)	46	35

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	172	ASP
1	B	172	ASP
1	C	172	ASP
1	D	172	ASP

### 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	258/255 (101%)	257 (100%)	1 (0%)	93	94
1	B	263/255 (103%)	262 (100%)	1 (0%)	93	94
1	C	262/255 (103%)	261 (100%)	1 (0%)	93	94
1	D	257/255 (101%)	252 (98%)	5 (2%)	65	59
All	All	1040/1020 (102%)	1032 (99%)	8 (1%)	86	86

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	245	GLU
1	B	212	THR
1	C	130	ASP
1	D	138	LEU
1	D	160	THR
1	D	233	GLN
1	D	254	VAL
1	D	288	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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](#)

1 ligand is 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	CIT	B	400	-	3,12,12	1.02	0	3,17,17	1.00	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	CIT	B	400	-	-	0/6/16/16	0/0/0/0

There are no bond length outliers.

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 ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	306/319 (95%)	0.10	13 (4%) 40 44	25, 37, 64, 84	0
1	B	306/319 (95%)	-0.05	8 (2%) 59 63	20, 30, 49, 61	0
1	C	306/319 (95%)	0.06	5 (1%) 74 78	21, 34, 58, 74	0
1	D	305/319 (95%)	0.09	14 (4%) 36 39	24, 36, 63, 77	0
All	All	1223/1276 (95%)	0.05	40 (3%) 50 53	20, 34, 60, 84	0

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	253	PHE	7.8
1	D	252	PHE	6.8
1	A	253	PHE	5.4
1	A	314	LEU	5.4
1	B	235	PRO	5.4
1	A	315	ASN	5.1
1	C	235	PRO	5.0
1	A	252	PHE	4.6
1	D	234	GLY	4.5
1	B	253	PHE	4.0
1	A	235	PRO	3.6
1	D	38	TYR	3.5
1	D	298	LEU	3.2
1	D	312	ALA	3.2
1	B	234	GLY	3.0
1	A	38	TYR	3.0
1	C	234	GLY	2.9
1	D	314	LEU	2.8
1	A	21[A]	TYR	2.7
1	D	308	ASP	2.7
1	D	297	ARG	2.6

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	B	315	ASN	2.6
1	A	6	SER	2.6
1	B	314	LEU	2.5
1	D	6	SER	2.5
1	D	255	ASP	2.5
1	D	285	ARG	2.4
1	C	243	LYS	2.4
1	D	21	TYR	2.3
1	A	251	PRO	2.3
1	B	255	ASP	2.3
1	B	252	PHE	2.3
1	A	308	ASP	2.3
1	D	179	GLU	2.3
1	A	179	GLU	2.3
1	B	6	SER	2.3
1	C	179	GLU	2.2
1	D	23	LEU	2.1
1	A	285	ARG	2.1
1	A	233	GLN	2.0

## 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	CIT	B	400	13/13	0.88	0.13	1.55	32,38,43,43	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.