



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 07:29 am GMT

PDB ID : 1C7Q
Title : THE CRYSTAL STRUCTURE OF PHOSPHOGLUCOSE ISOMERASE/AUTOCRINE MOTILITY FACTOR/NEUROLEUKIN COMPLEXED WITH ITS CARBOHYDRATE PHOSPHATE INHIBITORS AND ITS SUBSTRATE RECOGNITION MECHANISM
Authors : Chou, C.-C.; Meng, M.; Sun, Y.-J.; Hsiao, C.-D.
Deposited on : 2000-03-02
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

<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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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) : recalc28949

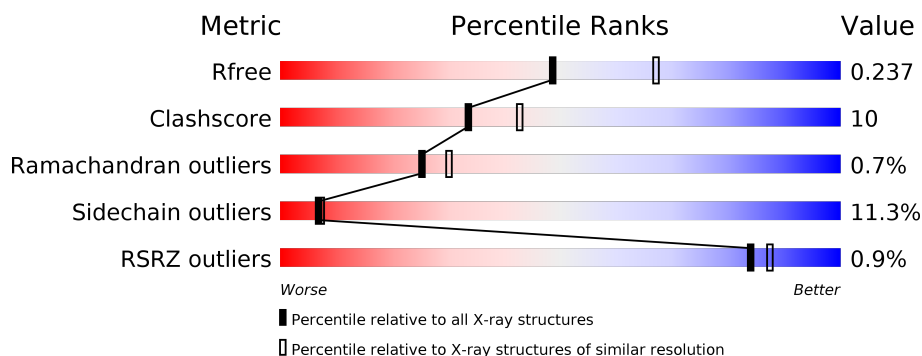
1 Overall quality at a glance

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}	100719	4130 (2.30-2.30)
Clashscore	112137	4751 (2.30-2.30)
Ramachandran outliers	110173	4705 (2.30-2.30)
Sidechain outliers	110143	4704 (2.30-2.30)
RSRZ outliers	101464	4156 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 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	445	

2 Entry composition [i](#)

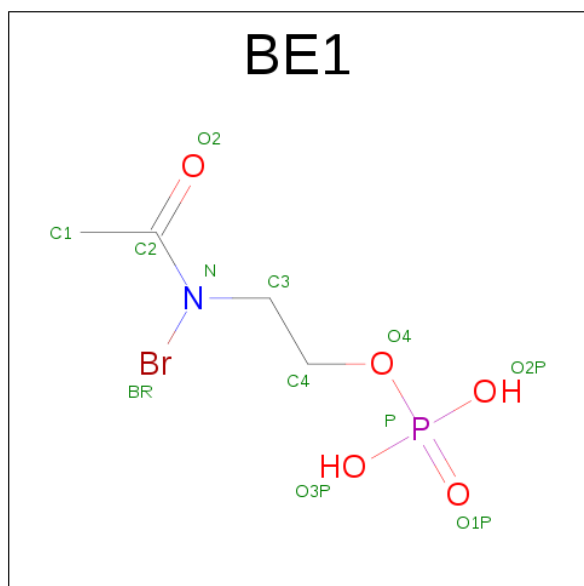
There are 3 unique types of molecules in this entry. The entry contains 3659 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 PHOSPHOGLUCOSE ISOMERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	442	Total	C	N	O	S	0	0	0
			3514	2243	594	665	12			

- Molecule 2 is N-BROMOACETYL-AMINOETHYL PHOSPHATE (three-letter code: BE1) (formula: C₄H₉BrNO₅P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			11	4	1	5	1		

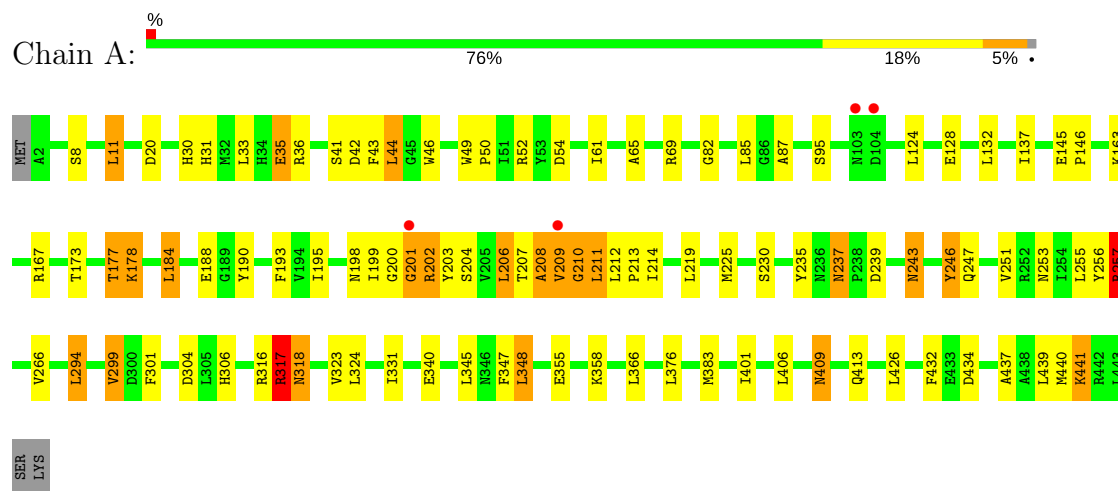
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	134	Total	O	0	0
			134	134		

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 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: PHOSPHOGLUCOSE ISOMERASE



4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	74.94Å 93.64Å 171.99Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 2.30 8.00 – 2.30	Depositor EDS
% Data completeness (in resolution range)	95.8 (8.00-2.30) 95.8 (8.00-2.30)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.21 (at 2.30Å)	Xtriage
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.172 , 0.242 0.173 , 0.237	Depositor DCC
R_{free} test set	2528 reflections (9.92%)	DCC
Wilson B-factor (Å ²)	23.0	Xtriage
Anisotropy	0.124	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 63.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3659	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.44% 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: BE1

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	0.50	1/3592 (0.0%)	0.63	0/4853

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	A	0	13

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	306	HIS	C-N	5.60	1.47	1.34

There are no bond angle outliers.

There are no chirality outliers.

All (13) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	190	TYR	Sidechain
1	A	202	ARG	Mainchain,Peptide
1	A	208	ALA	Mainchain,Peptide
1	A	210	GLY	Mainchain
1	A	235	TYR	Sidechain
1	A	246	TYR	Sidechain
1	A	257	ARG	Sidechain
1	A	301	PHE	Mainchain,Peptide
1	A	317	ARG	Sidechain

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Group
1	A	43	PHE	Sidechain

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	3514	0	3451	67	0
2	A	11	0	5	1	0
3	A	134	0	0	2	0
All	All	3659	0	3456	68	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 10.

All (68) 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:87:ALA:HB2	1:A:209:VAL:HB	1.30	1.10
1:A:201:GLY:O	1:A:202:ARG:HG2	1.74	0.86
1:A:41:SER:HA	1:A:44:LEU:HD22	1.61	0.82
1:A:204:SER:O	1:A:210:GLY:HA3	1.85	0.75
1:A:256:TYR:HD2	1:A:294:LEU:HD13	1.53	0.72
1:A:202:ARG:HB2	1:A:203:TYR:HD1	1.55	0.70
1:A:46:TRP:HB3	1:A:203:TYR:HD2	1.59	0.67
1:A:50:PRO:HA	1:A:225:MET:HE1	1.76	0.67
1:A:82:GLY:HA3	1:A:202:ARG:HA	1.78	0.66
1:A:8:SER:HA	1:A:11:LEU:HD22	1.78	0.66
1:A:202:ARG:HG3	1:A:203:TYR:CE1	2.32	0.65
1:A:256:TYR:CD2	1:A:294:LEU:HD13	2.36	0.60
1:A:137:ILE:CD1	1:A:209:VAL:HG13	2.32	0.59
1:A:30:HIS:HD2	1:A:246:TYR:OH	1.85	0.59
1:A:247:GLN:O	1:A:251:VAL:HG13	2.02	0.58
1:A:200:GLY:O	1:A:202:ARG:N	2.38	0.57
1:A:317:ARG:HD2	1:A:317:ARG:N	2.20	0.56
1:A:46:TRP:HB3	1:A:203:TYR:CD2	2.42	0.55
1:A:137:ILE:HD12	1:A:209:VAL:HG13	1.90	0.54

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:212:LEU:HB3	1:A:213:PRO:HD3	1.89	0.53
1:A:193:PHE:CG	1:A:207:THR:HG21	2.44	0.53
2:A:500:BE1:O3P	3:A:571:HOH:O	2.19	0.52
1:A:243:ASN:C	1:A:243:ASN:HD22	2.12	0.52
1:A:347:PHE:CE1	1:A:348:LEU:HD13	2.45	0.52
1:A:195:ILE:HG23	1:A:206:LEU:HD23	1.92	0.51
1:A:177:THR:OG1	1:A:178:LYS:HE3	2.11	0.51
1:A:163:LYS:O	1:A:167:ARG:HG3	2.10	0.50
1:A:437:ALA:O	1:A:441:LYS:HD2	2.11	0.50
1:A:87:ALA:HB2	1:A:209:VAL:CB	2.22	0.49
1:A:210:GLY:HA2	1:A:214:ILE:CD1	2.42	0.49
1:A:137:ILE:HD13	1:A:209:VAL:H	1.78	0.48
1:A:409:ASN:C	1:A:409:ASN:HD22	2.17	0.48
1:A:299:VAL:HG22	1:A:304:ASP:CB	2.44	0.48
1:A:199:ILE:CD1	1:A:206:LEU:HD11	2.44	0.47
1:A:253:ASN:O	1:A:257:ARG:HD2	2.15	0.47
1:A:208:ALA:O	1:A:209:VAL:HG22	2.15	0.47
1:A:318:ASN:HD22	1:A:318:ASN:H	1.62	0.47
1:A:46:TRP:CB	1:A:203:TYR:HD2	2.27	0.46
1:A:49:TRP:CE2	1:A:206:LEU:HD12	2.50	0.46
1:A:61:ILE:HD11	1:A:207:THR:OG1	2.16	0.45
1:A:266:VAL:HG22	1:A:323:VAL:HG13	1.98	0.45
1:A:195:ILE:HG12	1:A:208:ALA:HB2	1.98	0.45
1:A:206:LEU:O	1:A:211:LEU:HB2	2.18	0.44
1:A:432:PHE:N	1:A:432:PHE:CD2	2.85	0.44
1:A:145:GLU:HB2	1:A:146:PRO:HD3	1.99	0.44
1:A:31:HIS:O	1:A:35:GLU:HB2	2.18	0.43
1:A:46:TRP:CG	1:A:203:TYR:HE2	2.36	0.43
1:A:199:ILE:HD13	1:A:206:LEU:HD11	2.01	0.43
1:A:299:VAL:HG22	1:A:304:ASP:HB2	2.00	0.43
1:A:237:ASN:HD21	1:A:239:ASP:HB3	1.83	0.43
1:A:42:ASP:O	1:A:198:ASN:HB2	2.19	0.43
1:A:209:VAL:HG12	1:A:213:PRO:HG2	2.01	0.42
1:A:95:SER:HB3	3:A:576:HOH:O	2.20	0.42
1:A:202:ARG:HG3	1:A:203:TYR:CD1	2.54	0.42
1:A:347:PHE:CD1	1:A:348:LEU:HD13	2.55	0.42
1:A:324:LEU:HA	1:A:324:LEU:HD23	1.88	0.42
1:A:345:LEU:O	1:A:348:LEU:HB2	2.20	0.42
1:A:316:ARG:C	1:A:317:ARG:HD2	2.40	0.41
1:A:246:TYR:CE2	1:A:401:ILE:HG13	2.55	0.41
1:A:210:GLY:HA2	1:A:214:ILE:HD11	2.02	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:137:ILE:HG23	1:A:173:THR:HB	2.02	0.41
1:A:184:LEU:HD22	1:A:188:GLU:HG2	2.02	0.41
1:A:65:ALA:O	1:A:69:ARG:HG3	2.20	0.41
1:A:49:TRP:HB3	1:A:50:PRO:HD3	2.02	0.41
1:A:46:TRP:CD1	1:A:203:TYR:HE2	2.39	0.40
1:A:163:LYS:HE3	1:A:163:LYS:HB3	1.86	0.40
1:A:210:GLY:O	1:A:213:PRO:HD2	2.21	0.40
1:A:409:ASN:C	1:A:409:ASN:ND2	2.74	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	440/445 (99%)	420 (96%)	17 (4%)	3 (1%)	25	30

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	201	GLY
1	A	209	VAL
1	A	54	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	371/374 (99%)	329 (89%)	42 (11%)	7 7

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	LEU
1	A	20	ASP
1	A	33	LEU
1	A	35	GLU
1	A	36	ARG
1	A	44	LEU
1	A	52	ARG
1	A	85	LEU
1	A	124	LEU
1	A	128	GLU
1	A	132	LEU
1	A	177	THR
1	A	178	LYS
1	A	184	LEU
1	A	206	LEU
1	A	211	LEU
1	A	219	LEU
1	A	230	SER
1	A	237	ASN
1	A	243	ASN
1	A	255	LEU
1	A	257	ARG
1	A	294	LEU
1	A	299	VAL
1	A	317	ARG
1	A	318	ASN
1	A	331	ILE
1	A	340	GLU
1	A	348	LEU
1	A	355	GLU
1	A	358	LYS
1	A	366	LEU
1	A	376	LEU
1	A	383	MET
1	A	406	LEU
1	A	409	ASN
1	A	413	GLN
1	A	426	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	434	ASP
1	A	439	LEU
1	A	440	MET
1	A	441	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	30	HIS
1	A	34	HIS
1	A	70	ASN
1	A	135	ASN
1	A	233	HIS
1	A	237	ASN
1	A	243	ASN
1	A	247	GLN
1	A	281	GLN
1	A	313	GLN
1	A	318	ASN
1	A	357	ASN
1	A	409	ASN
1	A	413	GLN
1	A	422	ASN

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	BE1	A	500	1	10,10,11	0.79	0	12,13,15	2.75	8 (66%)

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	BE1	A	500	1	-	0/8/8/11	0/0/0/0

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	500	BE1	O2-C2-C1	-5.82	111.47	122.06
2	A	500	BE1	O3P-P-O2P	-3.30	94.28	107.61
2	A	500	BE1	O2-C2-N	-2.66	116.80	121.89
2	A	500	BE1	C1-C2-N	-2.52	111.50	116.03
2	A	500	BE1	O2P-P-O4	-2.10	101.14	106.73
2	A	500	BE1	O2P-P-O1P	2.50	120.30	110.50
2	A	500	BE1	P-O4-C4	2.58	125.40	118.30
2	A	500	BE1	O3P-P-O4	3.15	115.13	106.73

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	500	BE1	1	0

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	442/445 (99%)	-0.74	4 (0%) 84 87	5, 18, 41, 61	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	201	GLY	5.7
1	A	209	VAL	2.5
1	A	103	ASN	2.2
1	A	104	ASP	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, 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	LLDF	B-factors(Å ²)	Q<0.9
2	BE1	A	500	11/12	0.89	0.16	0.74	24,38,42,43	0

6.5 Other polymers [i](#)

There are no such residues in this entry.