



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 07:00 PM UTC

PDB ID : 3C3O / pdb_00003c3o
Title : ALIX Bro1-domain:CHMIP4A co-crystal structure
Authors : McCullough, J.B.; Fisher, R.D.; Whitby, F.G.; Sundquist, W.I.; Hill, C.P.
Deposited on : 2008-01-28
Resolution : 2.15 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

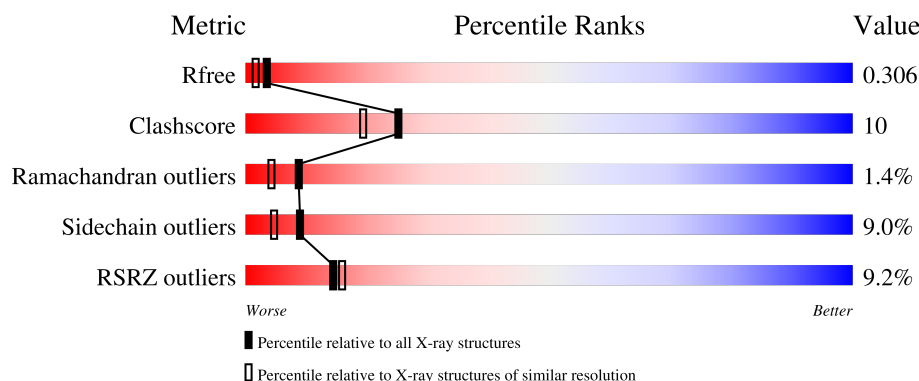
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.15 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	2057 (2.16-2.16)
Clashscore	190562	2159 (2.16-2.16)
Ramachandran outliers	187476	2134 (2.16-2.16)
Sidechain outliers	187428	2133 (2.16-2.16)
RSRZ outliers	180081	2059 (2.16-2.16)

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	380	9% 72% 18% • • 6%
2	B	13	8% 54% 46%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3043 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 Programmed cell death 6-interacting protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	357	Total	C	N	O	S	0	0	0
			2803	1792	471	529	11			

There are 21 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-20	MET	-	expression tag	UNP Q8WUM4
A	-19	HIS	-	expression tag	UNP Q8WUM4
A	-18	HIS	-	expression tag	UNP Q8WUM4
A	-17	HIS	-	expression tag	UNP Q8WUM4
A	-16	HIS	-	expression tag	UNP Q8WUM4
A	-15	HIS	-	expression tag	UNP Q8WUM4
A	-14	HIS	-	expression tag	UNP Q8WUM4
A	-13	HIS	-	expression tag	UNP Q8WUM4
A	-12	HIS	-	expression tag	UNP Q8WUM4
A	-11	HIS	-	expression tag	UNP Q8WUM4
A	-10	HIS	-	expression tag	UNP Q8WUM4
A	-9	SER	-	expression tag	UNP Q8WUM4
A	-8	GLY	-	expression tag	UNP Q8WUM4
A	-7	GLN	-	expression tag	UNP Q8WUM4
A	-6	ASN	-	expression tag	UNP Q8WUM4
A	-5	LEU	-	expression tag	UNP Q8WUM4
A	-4	TYR	-	expression tag	UNP Q8WUM4
A	-3	PHE	-	expression tag	UNP Q8WUM4
A	-2	GLN	-	expression tag	UNP Q8WUM4
A	-1	GLY	-	expression tag	UNP Q8WUM4
A	0	HIS	-	expression tag	UNP Q8WUM4

- Molecule 2 is a protein called Charged multivesicular body protein 4a peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	13	Total	C	N	O	0	0	0
			107	67	16	24			

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	126	Total	O	0	0
			126	126		
4	B	1	Total	O	0	0
			1	1		

- Molecule 1: Programmed cell death 6-interacting protein



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	120.67Å 62.68Å 76.12Å 90.00° 122.13° 90.00°	Depositor
Resolution (Å)	32.22 – 2.15 32.22 – 2.15	Depositor EDS
% Data completeness (in resolution range)	94.9 (32.22-2.15) 94.9 (32.22-2.15)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.49 (at 2.16Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.214 , 0.284 0.233 , 0.306	Depositor DCC
R_{free} test set	1799 reflections (6.83%)	wwPDB-VP
Wilson B-factor (Å ²)	43.5	Xtriage
Anisotropy	0.093	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 40.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3043	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.22% 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: GOL

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	1.13	4/2856 (0.1%)	1.22	12/3854 (0.3%)
2	B	0.72	0/108	1.04	1/144 (0.7%)
All	All	1.12	4/2964 (0.1%)	1.21	13/3998 (0.3%)

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	1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	280	ALA	CA-CB	7.07	1.64	1.53
1	A	242	PHE	CA-C	5.75	1.58	1.52
1	A	148	ILE	CA-CB	5.65	1.62	1.54
1	A	249	HIS	CA-C	5.17	1.59	1.52

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	167	VAL	CB-CA-C	-8.39	101.14	111.88
1	A	86	GLU	N-CA-C	-5.87	105.20	112.72
2	B	221	VAL	N-CA-C	-5.73	105.22	113.07
1	A	64	ALA	N-CA-C	5.68	118.20	111.33
1	A	297	VAL	N-CA-CB	-5.66	104.75	112.28
1	A	41	ARG	N-CA-C	-5.40	107.06	113.97
1	A	332	ILE	N-CA-C	-5.38	102.64	109.58

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	167	VAL	N-CA-CB	5.34	116.44	110.51
1	A	202	LYS	N-CA-C	-5.13	105.58	111.07
1	A	197	GLU	N-CA-C	-5.13	105.77	111.36
1	A	176	THR	CB-CA-C	5.09	119.89	109.76
1	A	153	TYR	N-CA-C	-5.03	105.99	111.82
1	A	231	CYS	CB-CA-C	-5.01	101.13	109.55

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	28	THR	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	2803	0	2828	57	0
2	B	107	0	98	3	0
3	A	6	0	8	1	0
4	A	126	0	0	3	0
4	B	1	0	0	0	0
All	All	3043	0	2934	57	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 (57) 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:19:LYS:HB3	1:A:20:PRO:HD3	1.51	0.92
1:A:163:ILE:O	1:A:167:VAL:HG22	1.75	0.85
1:A:46:LEU:HD22	1:A:78:ILE:HD11	1.70	0.71
1:A:28:THR:O	1:A:29:TYR:HB2	1.92	0.68
1:A:176:THR:HG23	1:A:178:ASP:OD1	1.94	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:29:TYR:H	1:A:30:PRO:HD3	1.60	0.67
1:A:39:TYR:HE2	1:A:358:MET:HB2	1.61	0.65
1:A:89:ILE:HG22	1:A:90:CYS:H	1.62	0.65
1:A:46:LEU:HD13	1:A:74:GLN:CG	2.30	0.61
1:A:112:ALA:O	1:A:113:LEU:HD23	1.99	0.61
1:A:46:LEU:HD13	1:A:74:GLN:HG2	1.83	0.60
1:A:87:ASN:HA	1:A:88:GLN:HB2	1.83	0.59
1:A:84:PHE:O	1:A:87:ASN:HB2	2.05	0.56
1:A:158:GLY:HA3	1:A:341:THR:O	2.06	0.56
1:A:24:PHE:HE2	1:A:83:PRO:HG2	1.70	0.56
1:A:39:TYR:CE2	1:A:358:MET:HB2	2.40	0.56
1:A:306:ARG:NH1	4:A:431:HOH:O	2.39	0.56
1:A:87:ASN:CA	1:A:88:GLN:HB2	2.36	0.55
1:A:46:LEU:CD2	1:A:78:ILE:HD11	2.38	0.54
1:A:177:VAL:HG22	4:A:433:HOH:O	2.08	0.54
1:A:38:GLN:C	1:A:40:CYS:H	2.18	0.51
1:A:337:LEU:CD1	2:B:217:LEU:HD13	2.41	0.50
1:A:89:ILE:HG22	1:A:90:CYS:N	2.27	0.50
1:A:176:THR:HG21	4:A:372:HOH:O	2.11	0.50
1:A:151:LYS:HG2	3:A:360:GOL:O1	2.12	0.49
1:A:158:GLY:HA2	1:A:340:SER:HB3	1.95	0.48
1:A:91:LEU:O	1:A:114:ALA:HA	2.13	0.48
1:A:26:GLN:HA	1:A:36:GLN:OE1	2.14	0.48
1:A:19:LYS:HB3	1:A:20:PRO:CD	2.35	0.47
1:A:46:LEU:CD2	1:A:78:ILE:CD1	2.92	0.47
1:A:78:ILE:HG22	1:A:116:LEU:HD21	1.96	0.47
1:A:173:ARG:CG	1:A:174:GLU:N	2.77	0.47
1:A:113:LEU:HD21	1:A:173:ARG:HD3	1.96	0.47
1:A:147:LYS:CD	2:B:220:TRP:O	2.63	0.46
1:A:10:LYS:NZ	1:A:96:LYS:O	2.47	0.46
1:A:82:PHE:HA	1:A:83:PRO:HD3	1.78	0.46
1:A:50:ARG:HD2	1:A:71:TYR:OH	2.17	0.45
1:A:147:LYS:HD3	2:B:220:TRP:O	2.16	0.45
1:A:38:GLN:C	1:A:40:CYS:N	2.75	0.45
1:A:46:LEU:HD13	1:A:74:GLN:HG3	1.97	0.44
1:A:173:ARG:CG	1:A:174:GLU:H	2.29	0.44
1:A:17:LEU:HA	1:A:91:LEU:HD21	1.99	0.44
1:A:98:ALA:N	1:A:178:ASP:OD2	2.36	0.44
1:A:108:SER:HB2	1:A:110:LYS:HG3	2.00	0.44
1:A:228:PHE:CD1	1:A:228:PHE:C	2.96	0.44
1:A:10:LYS:HG3	1:A:98:ALA:HB2	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:240:GLU:H	1:A:240:GLU:CD	2.23	0.44
1:A:15:VAL:HG12	1:A:17:LEU:HD23	1.99	0.43
1:A:71:TYR:CZ	1:A:75:ILE:HD11	2.53	0.43
1:A:83:PRO:HB2	1:A:88:GLN:HB3	1.99	0.43
1:A:84:PHE:HZ	1:A:91:LEU:HD12	1.84	0.43
1:A:87:ASN:N	1:A:88:GLN:HB2	2.34	0.42
1:A:228:PHE:HE2	1:A:249:HIS:CD2	2.37	0.42
1:A:46:LEU:CD1	1:A:74:GLN:HG2	2.47	0.42
1:A:208:MET:HE2	1:A:212:ILE:HD13	2.02	0.42
1:A:29:TYR:N	1:A:30:PRO:HD3	2.32	0.42
1:A:173:ARG:HG3	1:A:174:GLU:H	1.86	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	355/380 (93%)	330 (93%)	20 (6%)	5 (1%)	9	4
2	B	11/13 (85%)	8 (73%)	3 (27%)	0	100	100
All	All	366/393 (93%)	338 (92%)	23 (6%)	5 (1%)	9	4

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	29	TYR
1	A	88	GLN
1	A	87	ASN
1	A	108	SER
1	A	33	GLY

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	299/320 (93%)	274 (92%)	25 (8%)	10	6
2	B	11/11 (100%)	8 (73%)	3 (27%)	0	0
All	All	310/331 (94%)	282 (91%)	28 (9%)	9	5

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	LEU
1	A	29	TYR
1	A	35	GLU
1	A	36	GLN
1	A	41	ARG
1	A	44	GLU
1	A	60	LYS
1	A	77	SER
1	A	78	ILE
1	A	79	GLU
1	A	81	LYS
1	A	85	SER
1	A	110	LYS
1	A	116	LEU
1	A	146	LEU
1	A	167	VAL
1	A	169	SER
1	A	171	LEU
1	A	184	VAL
1	A	228	PHE
1	A	297	VAL
1	A	298	LYS
1	A	338	VAL
1	A	345	VAL
1	A	357	LYS
2	B	211	GLU
2	B	214	LEU

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Mol	Chain	Res	Type
2	B	216	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	87	ASN
1	A	88	GLN
1	A	138	GLN
1	A	230	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](#)

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	$\# Z > 2$
3	GOL	A	360	-	5,5,5	0.38	0	5,5,5	0.56	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
3	GOL	A	360	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	360	GOL	C1-C2-C3-O3
3	A	360	GOL	O2-C2-C3-O3

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	360	GOL	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

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, 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	357/380 (93%)	0.36	33 (9%) 14 16	15, 27, 68, 84	0
2	B	13/13 (100%)	1.05	1 (7%) 19 22	37, 50, 60, 62	0
All	All	370/393 (94%)	0.38	34 (9%) 14 16	15, 28, 68, 84	0

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	29	TYR	5.4
1	A	25	ILE	5.4
1	A	28	THR	5.2
1	A	30	PRO	5.1
1	A	2	ALA	4.8
1	A	84	PHE	4.6
1	A	85	SER	3.7
1	A	82	PHE	3.4
1	A	37	ALA	3.4
1	A	106	GLY	3.4
1	A	27	GLN	3.4
1	A	21	LEU	3.3
1	A	107	GLY	3.2
1	A	24	PHE	3.2
1	A	83	PRO	3.1
1	A	32	GLY	3.0
1	A	91	LEU	2.9
1	A	105	PHE	2.8
1	A	39	TYR	2.8
1	A	358	MET	2.6
1	A	22	VAL	2.6
1	A	167	VAL	2.6
1	A	61	HIS	2.6
1	A	58	LEU	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	36	GLN	2.5
1	A	40	CYS	2.5
1	A	31	SER	2.4
1	A	108	SER	2.4
2	B	221	VAL	2.4
1	A	88	GLN	2.2
1	A	80	PRO	2.2
1	A	23	LYS	2.1
1	A	89	ILE	2.0
1	A	42	ALA	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 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
3	GOL	A	360	6/6	0.91	0.21	79,82,84,84	0

6.5 Other polymers [i](#)

There are no such residues in this entry.