



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 05:00 PM UTC

PDB ID : 4H71 / pdb_00004h71
Title : Human Plk1-PBD in complex with Poloxime ((E)-4-(hydroxyimino)-2-isopropyl-5-methylcyclohexa-2,5-dienone)
Authors : Yin, Z.; Rehse, P.H.
Deposited on : 2012-09-19
Resolution : 1.93 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

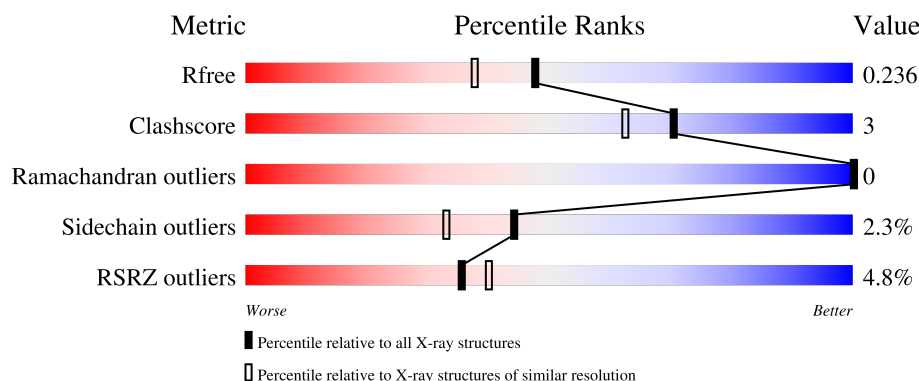
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

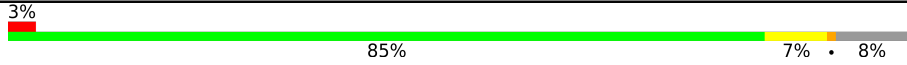
The reported resolution of this entry is 1.93 Å.

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	1452 (1.94-1.94)
Clashscore	190562	1494 (1.94-1.94)
Ramachandran outliers	187476	1479 (1.94-1.94)
Sidechain outliers	187428	1479 (1.94-1.94)
RSRZ outliers	180081	1453 (1.94-1.94)

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	240	
1	B	240	

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
3	PXE	A	702	-	-	X	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 3838 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.

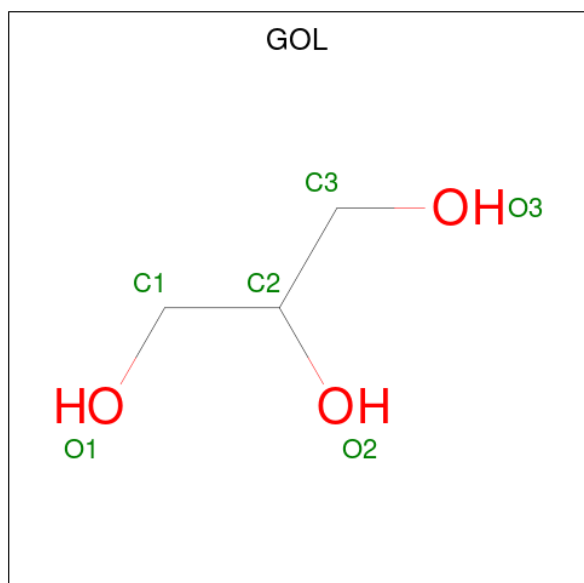
- Molecule 1 is a protein called Serine/threonine-protein kinase PLK1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	216	Total	C	N	O	S	0	0	0
			1722	1096	294	321	11			
1	A	222	Total	C	N	O	S	0	0	0
			1779	1131	306	331	11			

There are 6 discrepancies between the modelled and reference sequences:

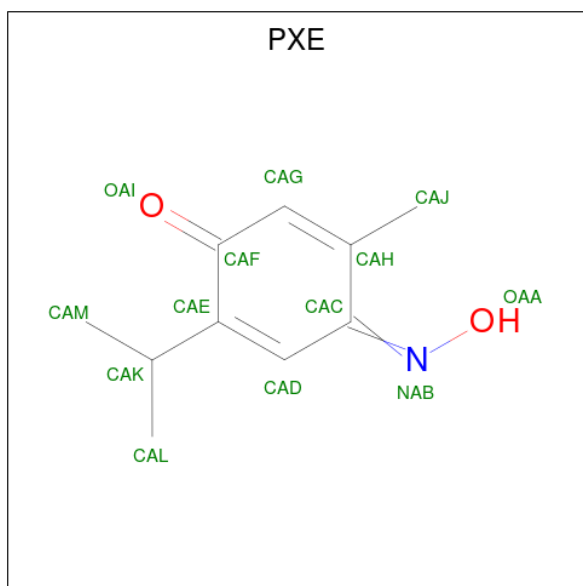
Chain	Residue	Modelled	Actual	Comment	Reference
B	364	SER	-	expression tag	UNP P53350
B	365	ASN	-	expression tag	UNP P53350
B	366	ALA	-	expression tag	UNP P53350
A	364	SER	-	expression tag	UNP P53350
A	365	ASN	-	expression tag	UNP P53350
A	366	ALA	-	expression tag	UNP P53350

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 3 is 2-methyl-5-(1-methylethyl)cyclohexa-2,5-diene-1,4-dione 1-oxime (CCD ID: PXE) (formula: $C_{10}H_{13}NO_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			13	10	1	2		

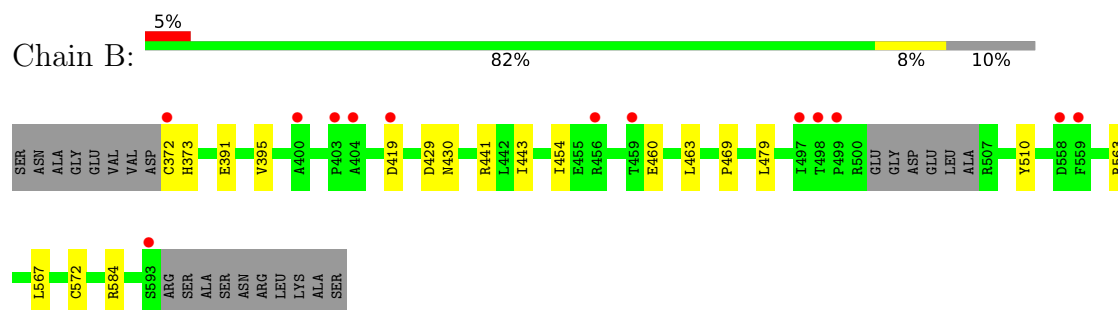
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	136	Total	O	0	0
			136	136		
4	A	182	Total	O	0	0
			182	182		

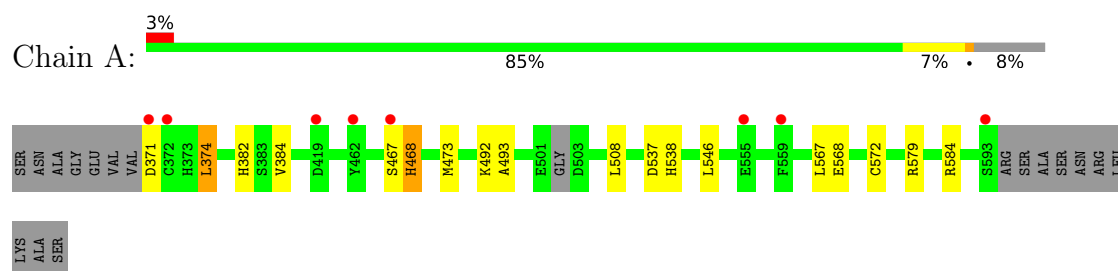
3 Residue-property plots [i](#)

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: Serine/threonine-protein kinase PLK1



- Molecule 1: Serine/threonine-protein kinase PLK1



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	33.26Å 102.75Å 68.49Å 90.00° 93.85° 90.00°	Depositor
Resolution (Å)	33.18 – 1.93 33.18 – 1.93	Depositor EDS
% Data completeness (in resolution range)	92.5 (33.18-1.93) 92.4 (33.18-1.93)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	8.01 (at 1.94Å)	Xtriage
Refinement program	REFMAC 5.5.0110	Depositor
R, R_{free}	0.188 , 0.237 0.188 , 0.236	Depositor DCC
R_{free} test set	1735 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	17.6	Xtriage
Anisotropy	0.149	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 43.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3838	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.55% 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: PXE, 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	0.65	0/1816	0.78	0/2458
1	B	0.62	0/1759	0.79	0/2379
All	All	0.64	0/3575	0.79	0/4837

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 [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	1779	0	1745	11	0
1	B	1722	0	1687	9	0
2	A	6	0	8	0	0
3	A	13	0	13	7	0
4	A	182	0	0	5	0
4	B	136	0	0	0	0
All	All	3838	0	3453	24	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.

All (24) 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)
3:A:702:PXE:OAI	3:A:702:PXE:HAMB	1.72	0.90
3:A:702:PXE:HAJ	4:A:826:HOH:O	1.76	0.86
3:A:702:PXE:HAJB	4:A:802:HOH:O	1.89	0.72
1:A:382:HIS:HD2	1:A:584:ARG:HH22	1.38	0.72
3:A:702:PXE:CAJ	4:A:802:HOH:O	2.43	0.68
1:A:493:ALA:HA	3:A:702:PXE:HALB	1.78	0.66
1:B:454:ILE:HG12	1:B:460:GLU:HG2	1.81	0.63
1:A:538:HIS:ND1	3:A:702:PXE:HAG	2.14	0.63
1:B:443:ILE:HD11	1:B:510:TYR:HB3	1.85	0.56
1:B:479:LEU:C	1:B:479:LEU:HD23	2.30	0.56
1:A:371:ASP:HB2	1:A:374:LEU:HB2	1.87	0.56
1:B:372:CYS:SG	1:B:373:HIS:N	2.76	0.55
1:A:567:LEU:HD22	1:A:572:CYS:HB3	1.90	0.54
1:A:537:ASP:OD2	1:A:579:ARG:NH2	2.49	0.45
1:A:538:HIS:HD2	4:A:873:HOH:O	2.00	0.45
1:B:429:ASP:O	1:B:430:ASN:HB2	2.16	0.44
1:A:473:MET:HE2	4:A:906:HOH:O	2.17	0.44
1:A:384:VAL:HA	1:A:568:GLU:HG2	2.00	0.44
1:B:567:LEU:HD22	1:B:572:CYS:HB3	2.00	0.43
1:A:468:HIS:ND1	1:A:468:HIS:C	2.77	0.42
1:B:584:ARG:C	1:B:584:ARG:HD3	2.45	0.41
1:B:441:ARG:HD3	1:B:510:TYR:CD1	2.56	0.41
1:A:492:LYS:C	3:A:702:PXE:CAL	2.92	0.41
1:B:463:LEU:CD1	1:B:469:PRO:HG3	2.51	0.41

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	218/240 (91%)	213 (98%)	5 (2%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	212/240 (88%)	209 (99%)	3 (1%)	0	100	100
All	All	430/480 (90%)	422 (98%)	8 (2%)	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	195/215 (91%)	190 (97%)	5 (3%)	40	28
1	B	189/215 (88%)	185 (98%)	4 (2%)	47	37
All	All	384/430 (89%)	375 (98%)	9 (2%)	44	33

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	391	GLU
1	B	395	VAL
1	B	419	ASP
1	B	563	ARG
1	A	374	LEU
1	A	467	SER
1	A	468	HIS
1	A	508	LEU
1	A	546	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	373	HIS
1	B	426	GLN
1	B	430	ASN
1	B	470	ASN

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Mol	Chain	Res	Type
1	B	524	HIS
1	A	382	HIS
1	A	385	ASN
1	A	452	GLN
1	A	489	HIS
1	A	524	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](#)

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	$\# Z > 2$
3	PXE	A	702	-	13,13,13	2.75	6 (46%)	17,18,18	2.88	8 (47%)
2	GOL	A	701	-	5,5,5	0.46	0	5,5,5	0.39	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	PXE	A	702	-	-	1/6/22/22	0/1/1/1
2	GOL	A	701	-	-	0/4/4/4	-

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	702	PXE	CAK-CAE	-6.12	1.43	1.52
3	A	702	PXE	CAG-CAF	-4.31	1.33	1.44
3	A	702	PXE	CAC-NAB	4.30	1.37	1.30
3	A	702	PXE	OAI-CAF	-2.27	1.18	1.24
3	A	702	PXE	CAD-CAC	-2.19	1.38	1.42
3	A	702	PXE	CAE-CAF	-2.12	1.44	1.48

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	702	PXE	CAM-CAK-CAE	-7.48	97.12	111.70
3	A	702	PXE	CAJ-CAH-CAC	4.40	125.83	119.24
3	A	702	PXE	CAJ-CAH-CAG	-3.87	113.15	121.96
3	A	702	PXE	CAK-CAE-CAD	3.52	128.32	123.32
3	A	702	PXE	CAK-CAE-CAF	-2.78	114.44	117.30
3	A	702	PXE	OAA-NAB-CAC	2.54	118.67	113.46
3	A	702	PXE	CAC-CAD-CAE	2.24	121.18	118.27
3	A	702	PXE	CAG-CAF-CAE	2.13	122.39	118.29

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	702	PXE	CAF-CAE-CAK-CAM

There are no ring outliers.

1 monomer is involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	702	PXE	7	0

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	222/240 (92%)	0.01	8 (3%) 46 52	7, 16, 27, 37	0
1	B	216/240 (90%)	0.43	13 (6%) 27 30	10, 20, 35, 41	0
All	All	438/480 (91%)	0.22	21 (4%) 35 40	7, 18, 33, 41	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	372	CYS	3.9
1	B	559	PHE	3.8
1	B	403	PRO	3.5
1	B	404	ALA	3.0
1	A	371	ASP	2.9
1	B	459	THR	2.8
1	B	400	ALA	2.7
1	A	467	SER	2.7
1	B	558	ASP	2.7
1	A	372	CYS	2.7
1	B	498	THR	2.4
1	B	593	SER	2.4
1	B	497	ILE	2.3
1	A	462	TYR	2.3
1	B	419	ASP	2.2
1	B	456	ARG	2.2
1	A	419	ASP	2.1
1	A	555	GLU	2.1
1	A	593	SER	2.0
1	A	559	PHE	2.0
1	B	499	PRO	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($\text{\AA}^2$)	Q<0.9
3	PXE	A	702	13/13	0.74	0.22	41,42,44,45	0
2	GOL	A	701	6/6	0.96	0.07	15,21,24,27	0

6.5 Other polymers [i](#)

There are no such residues in this entry.