



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 01:33 AM UTC

PDB ID : 1JX6 / pdb_00001jx6
Title : CRYSTAL STRUCTURE OF LUXP FROM VIBRIO HARVEYI COM-
PLEXED WITH AUTOINDUCER-2
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Deposited on : 2001-09-05
Resolution : 1.50 Å(reported)

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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

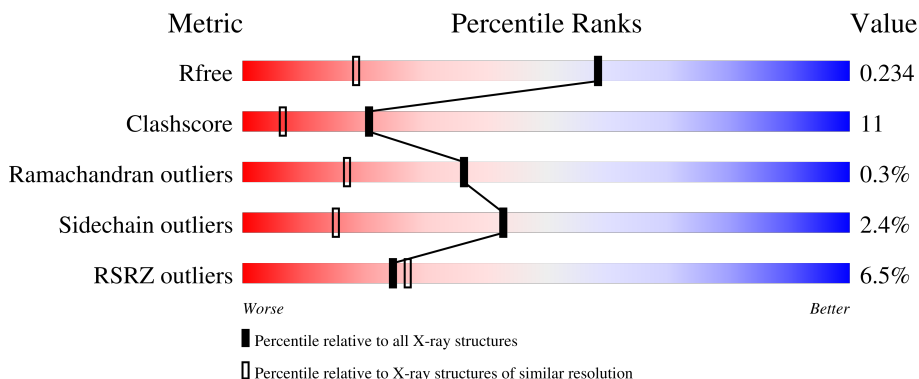
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.50 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	4037 (1.50-1.50)
Clashscore	190562	4235 (1.50-1.50)
Ramachandran outliers	187476	4153 (1.50-1.50)
Sidechain outliers	187428	4150 (1.50-1.50)
RSRZ outliers	180081	4039 (1.50-1.50)

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	342	6% 75% 22% ..

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 3073 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.

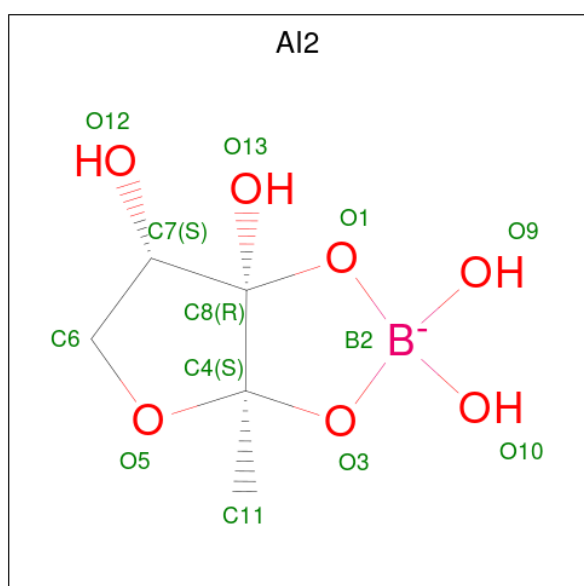
- Molecule 1 is a protein called LUXP PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	338	Total	C	N	O	S	0	0	0
			2727	1737	460	524	6			

- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Ca	0	0
			1	1		

- Molecule 3 is 3A-METHYL-5,6-DIHYDRO-FURO[2,3-D][1,3,2]DIOXABOROLE-2,2,6,6A-TETRAOL (CCD ID: AI2) (formula: C₅H₁₀BO₇).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	B	C	O	0	0
			13	1	5	7		

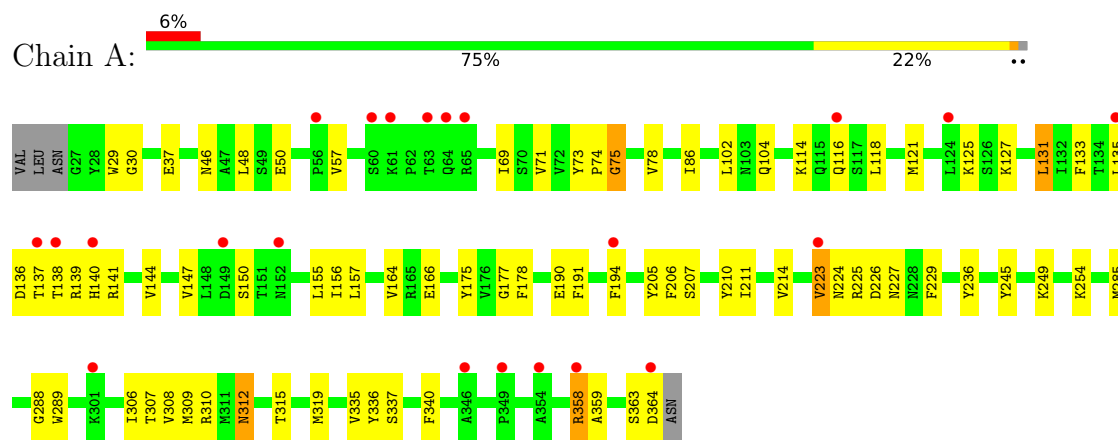
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	332	Total 332	O 332	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: LUXP PROTEIN



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	44.27Å 77.47Å 52.01Å 90.00° 96.37° 90.00°	Depositor
Resolution (Å)	24.58 – 1.50 24.58 – 1.50	Depositor EDS
% Data completeness (in resolution range)	92.3 (24.58-1.50) 96.5 (24.58-1.50)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	0.04	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.35 (at 1.45Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.214 , 0.239 0.210 , 0.234	Depositor DCC
R_{free} test set	5031 reflections (8.17%)	wwPDB-VP
Wilson B-factor (Å ²)	18.5	Xtriage
Anisotropy	0.463	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 40.7	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	0.96	EDS
Total number of atoms	3073	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.54% 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: AI2, CA

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	0.37	1/2792 (0.0%)	0.89	11/3785 (0.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	309	MET	SD-CE	-5.14	1.66	1.79

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	206	PHE	N-CA-C	-7.47	98.70	109.31
1	A	30	GLY	N-CA-C	-7.07	103.49	112.54
1	A	288	GLY	N-CA-C	-6.41	101.48	112.64
1	A	75	GLY	N-CA-C	6.38	121.69	114.67
1	A	207	SER	N-CA-C	-6.21	102.52	110.53
1	A	312	ASN	N-CA-C	6.16	123.91	110.80
1	A	359	ALA	N-CA-C	5.88	117.77	111.36
1	A	78	VAL	N-CA-C	-5.67	104.99	110.72
1	A	74	PRO	N-CA-C	-5.38	104.25	113.12
1	A	175	TYR	N-CA-C	-5.37	99.77	108.52
1	A	57	VAL	N-CA-C	-5.11	103.09	108.15

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	2727	0	2648	60	0
2	A	1	0	0	0	0
3	A	13	0	10	0	0
4	A	332	0	0	4	0
All	All	3073	0	2658	60	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 11.

All (60) 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:223:VAL:HG13	1:A:229:PHE:HB2	1.60	0.81
1:A:137:THR:HG22	1:A:138:THR:H	1.54	0.72
1:A:191:PHE:CZ	1:A:285:MET:HE2	2.29	0.68
1:A:137:THR:HG22	1:A:138:THR:N	2.10	0.66
1:A:139:ARG:HB2	1:A:139:ARG:NH1	2.11	0.66
1:A:223:VAL:CG1	1:A:229:PHE:HB2	2.26	0.65
1:A:135:LEU:CD1	1:A:157:LEU:HD13	2.27	0.64
1:A:225:ARG:HD2	1:A:226:ASP:OD1	1.97	0.64
1:A:71:VAL:HG23	1:A:102:LEU:HD21	1.80	0.64
1:A:178:PHE:CD2	1:A:310:ARG:HB2	2.33	0.63
1:A:363:SER:O	1:A:364:ASP:HB2	1.98	0.62
1:A:135:LEU:HD13	1:A:157:LEU:HD13	1.82	0.61
1:A:139:ARG:HB2	1:A:139:ARG:HH11	1.68	0.59
1:A:131:LEU:HD22	1:A:133:PHE:CD1	2.41	0.55
1:A:135:LEU:CD1	1:A:157:LEU:HD22	2.37	0.55
1:A:156:ILE:HG21	1:A:319:MET:HE3	1.89	0.54
1:A:223:VAL:HG12	1:A:224:ASN:HD22	1.72	0.54
1:A:147:VAL:HG11	1:A:155:LEU:HD21	1.90	0.54
1:A:190:GLU:HG3	1:A:194:PHE:CZ	2.42	0.54
1:A:191:PHE:CE2	1:A:285:MET:HE2	2.43	0.53
1:A:102:LEU:HD13	1:A:102:LEU:C	2.34	0.53
1:A:210:TYR:O	1:A:214:VAL:HG22	2.09	0.53
1:A:136:ASP:OD1	1:A:211:ILE:HG12	2.09	0.52
1:A:121:MET:O	1:A:125:LYS:HG3	2.09	0.52
1:A:135:LEU:HD12	1:A:157:LEU:HD22	1.91	0.51
1:A:137:THR:CG2	1:A:138:THR:H	2.25	0.49
1:A:223:VAL:HG13	1:A:229:PHE:CB	2.36	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:178:PHE:CE2	1:A:310:ARG:HB2	2.48	0.48
1:A:166:GLU:HG2	4:A:468:HOH:O	2.13	0.48
1:A:315:THR:O	1:A:319:MET:HG3	2.13	0.48
1:A:223:VAL:HG12	1:A:224:ASN:ND2	2.30	0.47
1:A:125:LYS:O	1:A:127:LYS:HD2	2.15	0.47
1:A:205:TYR:HB2	1:A:236:TYR:HA	1.97	0.47
1:A:29:TRP:CH2	1:A:37:GLU:HG2	2.50	0.46
1:A:306:ILE:HG12	1:A:307:THR:N	2.31	0.46
1:A:194:PHE:CD1	1:A:285:MET:HE1	2.51	0.46
1:A:131:LEU:HD22	1:A:133:PHE:HD1	1.80	0.46
1:A:358:ARG:NH1	4:A:549:HOH:O	2.47	0.46
1:A:178:PHE:HB3	1:A:340:PHE:CZ	2.52	0.45
1:A:46:ASN:O	1:A:50:GLU:HG3	2.16	0.45
1:A:308:VAL:O	1:A:308:VAL:HG13	2.15	0.45
1:A:69:ILE:CG2	1:A:102:LEU:HD23	2.47	0.44
1:A:137:THR:CG2	1:A:138:THR:N	2.79	0.44
1:A:177:GLY:O	1:A:337:SER:HA	2.18	0.44
1:A:69:ILE:HB	1:A:102:LEU:HD23	2.00	0.43
1:A:116:GLN:OE1	1:A:140:HIS:HE1	2.02	0.42
1:A:223:VAL:HG13	1:A:229:PHE:CD1	2.55	0.42
1:A:194:PHE:HD1	1:A:285:MET:HE1	1.85	0.42
1:A:335:VAL:HG22	1:A:336:TYR:N	2.34	0.41
1:A:73:TYR:C	1:A:75:GLY:H	2.28	0.41
1:A:133:PHE:HE1	1:A:155:LEU:HD22	1.84	0.41
1:A:86:ILE:HG21	1:A:104:GLN:NE2	2.36	0.41
1:A:224:ASN:HB3	4:A:603:HOH:O	2.21	0.41
1:A:245:TYR:CZ	1:A:249:LYS:HD2	2.55	0.41
1:A:141:ARG:HD3	4:A:681:HOH:O	2.20	0.41
1:A:254:LYS:N	1:A:254:LYS:HD2	2.36	0.41
1:A:140:HIS:O	1:A:144:VAL:HG23	2.21	0.40
1:A:164:VAL:HG12	1:A:166:GLU:HG2	2.02	0.40
1:A:116:GLN:CD	1:A:140:HIS:HE1	2.29	0.40
1:A:114:LYS:O	1:A:118:LEU:HG	2.22	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	336/342 (98%)	327 (97%)	8 (2%)	1 (0%)	36	17

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	312	ASN

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	293/297 (99%)	286 (98%)	7 (2%)	43	15

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	48	LEU
1	A	131	LEU
1	A	150	SER
1	A	223	VAL
1	A	227	ASN
1	A	289	TRP
1	A	358	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	64	GLN
1	A	101	GLN
1	A	104	GLN
1	A	110	ASN
1	A	140	HIS
1	A	221	HIS
1	A	222	GLN
1	A	224	ASN
1	A	312	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 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	$\# Z > 2$
3	AI2	A	400	-	13,14,14	2.70	8 (61%)	11,25,25	2.02	4 (36%)

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	AI2	A	400	-	-	-	0/2/2/2

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	400	AI2	O3-C4	5.84	1.47	1.39
3	A	400	AI2	O1-C8	-4.46	1.33	1.39
3	A	400	AI2	B2-O3	2.92	1.51	1.45
3	A	400	AI2	O13-C8	2.69	1.44	1.39
3	A	400	AI2	O5-C4	2.50	1.46	1.42
3	A	400	AI2	O12-C7	-2.48	1.39	1.43
3	A	400	AI2	B2-O1	-2.14	1.41	1.45
3	A	400	AI2	C8-C7	2.05	1.56	1.53

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	400	AI2	O5-C4-C11	3.34	113.16	109.26
3	A	400	AI2	O13-C8-O1	3.34	116.10	109.67
3	A	400	AI2	O3-C4-C11	-2.72	103.62	107.05
3	A	400	AI2	O12-C7-C8	2.28	118.47	110.94

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 [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	338/342 (98%)	0.54	22 (6%)	25 27	12, 22, 38, 64	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	223	VAL	3.5
1	A	194	PHE	3.1
1	A	64	GLN	3.0
1	A	349	PRO	2.8
1	A	138	THR	2.8
1	A	65	ARG	2.7
1	A	358	ARG	2.7
1	A	135	LEU	2.5
1	A	364	ASP	2.5
1	A	137	THR	2.5
1	A	354	ALA	2.4
1	A	61	LYS	2.4
1	A	301	LYS	2.4
1	A	63	THR	2.4
1	A	149	ASP	2.3
1	A	116	GLN	2.2
1	A	152	ASN	2.2
1	A	140	HIS	2.2
1	A	60	SER	2.1
1	A	56	PRO	2.1
1	A	346	ALA	2.0
1	A	124	LEU	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
2	CA	A	401	1/1	0.89	0.18	50,50,50,50	0
3	AI2	A	400	13/13	0.98	0.04	9,11,13,14	0

6.5 Other polymers [i](#)

There are no such residues in this entry.