



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:19 PM GMT

PDB ID : 3WW3
Title : X-ray structures of Cellulomonas parahominis L-ribose isomerase with no lig-
and
Authors : Terami, Y.; Yoshida, H.; Takata, G.; Kamitori, S.
Deposited on : 2014-06-13
Resolution : 1.90 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 (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

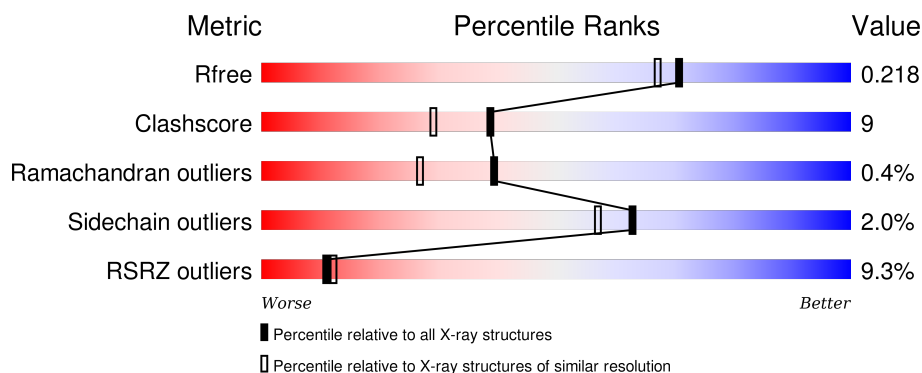
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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 ≥ 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	256	
1	B	256	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4094 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 L-ribose isomerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	240	Total	C	N	O	S	0	0	0
			1866	1184	321	352	9			
1	B	242	Total	C	N	O	S	0	0	0
			1878	1191	323	355	9			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-6	HIS	-	EXPRESSION TAG	UNP L0N3Y0
A	-5	HIS	-	EXPRESSION TAG	UNP L0N3Y0
A	-4	HIS	-	EXPRESSION TAG	UNP L0N3Y0
A	-3	HIS	-	EXPRESSION TAG	UNP L0N3Y0
A	-2	HIS	-	EXPRESSION TAG	UNP L0N3Y0
A	-1	HIS	-	EXPRESSION TAG	UNP L0N3Y0
A	0	GLY	-	EXPRESSION TAG	UNP L0N3Y0
A	1	SER	-	EXPRESSION TAG	UNP L0N3Y0
A	119	LEU	PHE	ENGINEERED MUTATION	UNP L0N3Y0
A	125	PHE	LEU	ENGINEERED MUTATION	UNP L0N3Y0
B	-6	HIS	-	EXPRESSION TAG	UNP L0N3Y0
B	-5	HIS	-	EXPRESSION TAG	UNP L0N3Y0
B	-4	HIS	-	EXPRESSION TAG	UNP L0N3Y0
B	-3	HIS	-	EXPRESSION TAG	UNP L0N3Y0
B	-2	HIS	-	EXPRESSION TAG	UNP L0N3Y0
B	-1	HIS	-	EXPRESSION TAG	UNP L0N3Y0
B	0	GLY	-	EXPRESSION TAG	UNP L0N3Y0
B	1	SER	-	EXPRESSION TAG	UNP L0N3Y0
B	119	LEU	PHE	ENGINEERED MUTATION	UNP L0N3Y0
B	125	PHE	LEU	ENGINEERED MUTATION	UNP L0N3Y0

- Molecule 2 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total 1	Mn 1	0	0
2	A	1	Total 1	Mn 1	0	0

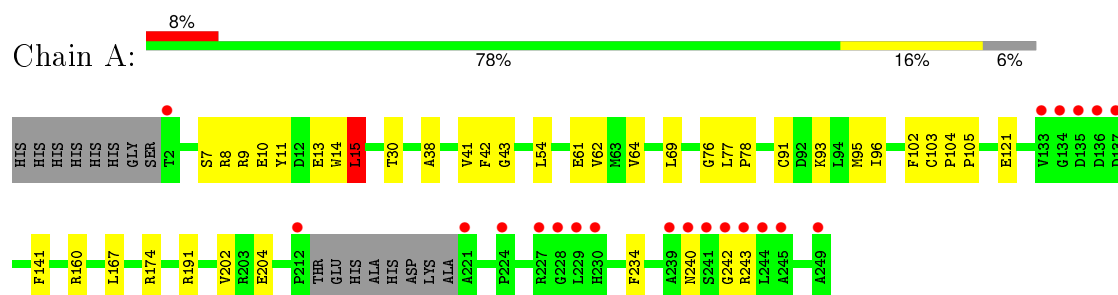
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	181	Total 181	O 181	0	0
3	B	167	Total 167	O 167	0	0

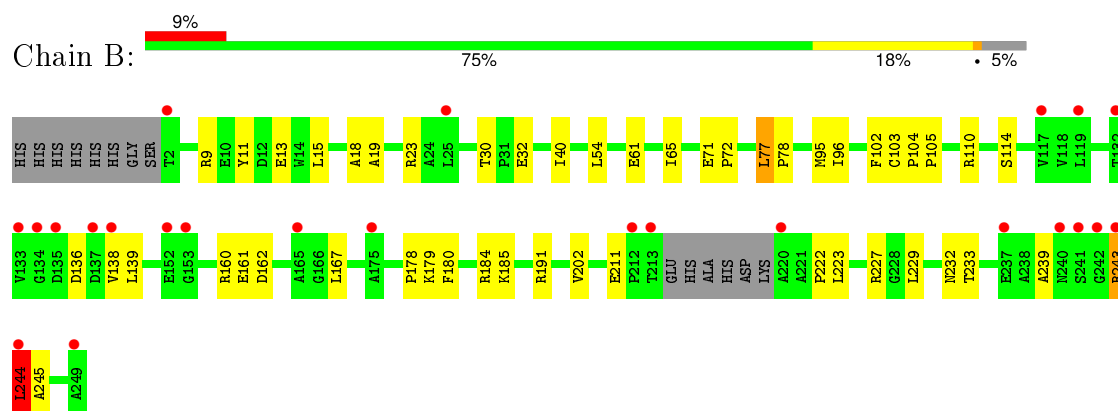
3 Residue-property plots

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: L-ribose isomerase



- Molecule 1: L-ribose isomerase



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	76.52Å 89.07Å 152.58Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.19 – 1.90 46.19 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.6 (46.19-1.90) 99.7 (46.19-1.90)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.74 (at 1.90Å)	Xtriage
Refinement program	CNS 1.3	Depositor
R, R_{free}	0.189 , 0.219 0.189 , 0.218	Depositor DCC
R_{free} test set	4156 reflections (10.06%)	DCC
Wilson B-factor (Å ²)	20.3	Xtriage
Anisotropy	0.017	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 49.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 41301 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4094	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.33% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: MN

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	0.30	0/1923	0.63	2/2624 (0.1%)
1	B	0.30	0/1935	0.60	1/2641 (0.0%)
All	All	0.30	0/3858	0.61	3/5265 (0.1%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	103	CYS	N-CA-C	-7.44	90.91	111.00
1	B	103	CYS	N-CA-C	-7.20	91.56	111.00
1	A	15	LEU	CA-CB-CG	-5.04	103.70	115.30

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	1866	0	1777	31	0
1	B	1878	0	1789	39	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	181	0	0	1	0
3	B	167	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	4094	0	3566	66	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 9.

All (66) 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:64:VAL:CG1	1:A:96:ILE:HB	2.02	0.90
1:A:43:GLY:H	1:A:242:GLY:HA3	1.47	0.80
1:A:64:VAL:HG13	1:A:96:ILE:HB	1.66	0.76
1:B:227:ARG:NH1	1:B:227:ARG:HB2	2.03	0.74
1:B:227:ARG:HH11	1:B:227:ARG:HB2	1.51	0.74
1:A:43:GLY:N	1:A:242:GLY:HA3	2.02	0.73
1:B:239:ALA:O	1:B:243:ARG:HB3	1.92	0.68
1:A:76:GLY:HA2	1:B:138:VAL:HG11	1.76	0.67
1:B:11:TYR:CE1	1:B:15:LEU:HD22	2.34	0.63
1:A:91:CYS:SG	1:A:93:LYS:HE2	2.39	0.63
1:A:15:LEU:HD21	1:A:38:ALA:HB3	1.82	0.62
1:B:77:LEU:HD22	1:B:78:PRO:HD2	1.84	0.60
1:A:14:TRP:CH2	1:A:62:VAL:HG22	2.36	0.60
1:B:19:ALA:O	1:B:23:ARG:HG3	2.02	0.59
1:A:167:LEU:HD12	1:A:167:LEU:N	2.19	0.58
1:A:30:THR:HG21	1:B:178:PRO:HG3	1.85	0.57
1:A:61:GLU:OE1	1:A:160:ARG:NH1	2.38	0.56
1:B:61:GLU:OE1	1:B:160:ARG:NH1	2.38	0.56
1:B:167:LEU:N	1:B:167:LEU:HD12	2.22	0.55
1:B:244:LEU:HD13	1:B:245:ALA:H	1.72	0.54
1:A:234:PHE:HB2	1:A:240:ASN:OD1	2.07	0.53
1:B:233:THR:HG23	3:B:617:HOH:O	2.09	0.53
1:B:227:ARG:HH11	1:B:227:ARG:CB	2.21	0.51
1:B:95:MET:HB2	1:B:202:VAL:HG13	1.93	0.50
1:A:121:GLU:OE2	1:A:174:ARG:HD3	2.12	0.50
1:A:95:MET:HB2	1:A:202:VAL:HG13	1.95	0.49
1:A:77:LEU:HD13	1:A:78:PRO:N	2.28	0.49
1:B:211:GLU:HG2	1:B:232:ASN:HD21	1.78	0.49
1:B:9:ARG:O	1:B:13:GLU:HG3	2.11	0.49
1:A:42:PHE:HA	1:A:242:GLY:O	2.13	0.49
1:B:227:ARG:HG2	3:B:629:HOH:O	2.11	0.49
1:A:96:ILE:HD12	1:A:96:ILE:N	2.29	0.48
1:B:102:PHE:CD2	1:B:191:ARG:HB3	2.49	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:7:SER:OG	1:A:10:GLU:HG3	2.13	0.47
1:A:15:LEU:CD2	1:A:38:ALA:HB3	2.43	0.47
1:A:93:LYS:HG2	1:A:204:GLU:HB3	1.97	0.47
1:B:138:VAL:CG1	1:B:139:LEU:N	2.77	0.47
1:B:243:ARG:HG2	1:B:243:ARG:O	2.15	0.47
1:A:11:TYR:CE2	1:A:15:LEU:HD13	2.49	0.47
1:B:40:ILE:CD1	1:B:65:ILE:HD12	2.45	0.46
1:A:9:ARG:O	1:A:13:GLU:HG3	2.16	0.46
1:B:184:ARG:O	1:B:185:LYS:HB2	2.15	0.46
1:A:69:LEU:HB2	1:B:179:LYS:HG3	1.98	0.45
1:B:136:ASP:HA	3:B:612:HOH:O	2.15	0.45
1:A:102:PHE:CD2	1:A:191:ARG:HB3	2.52	0.44
1:A:167:LEU:CD1	1:A:167:LEU:N	2.80	0.44
1:B:110:ARG:NH2	1:B:229:LEU:HD11	2.33	0.44
1:A:141:PHE:CD1	1:B:222:PRO:HA	2.53	0.44
1:B:114:SER:HA	1:B:180:PHE:O	2.18	0.44
1:B:30:THR:OG1	1:B:32:GLU:HG2	2.18	0.44
1:B:161:GLU:CD	1:B:161:GLU:H	2.22	0.43
1:A:104:PRO:HA	1:A:105:PRO:HD3	1.90	0.43
1:A:11:TYR:CZ	1:A:15:LEU:HD13	2.53	0.43
1:B:77:LEU:CD2	1:B:78:PRO:HD2	2.48	0.43
1:B:40:ILE:HD12	1:B:65:ILE:HD12	2.01	0.43
1:A:41:VAL:HG22	1:A:62:VAL:HG23	2.01	0.42
1:B:243:ARG:C	1:B:243:ARG:HD3	2.40	0.42
1:B:138:VAL:HG12	1:B:139:LEU:O	2.20	0.41
1:B:211:GLU:HG2	1:B:232:ASN:ND2	2.35	0.41
1:B:167:LEU:CD1	1:B:167:LEU:N	2.83	0.41
1:B:18:ALA:HB1	1:B:96:ILE:HD12	2.03	0.41
1:B:71:GLU:HA	1:B:72:PRO:HD3	1.93	0.41
1:B:244:LEU:CD1	1:B:245:ALA:H	2.33	0.41
1:A:77:LEU:HD13	1:A:78:PRO:HD2	2.03	0.40
1:B:104:PRO:HA	1:B:105:PRO:HD3	1.89	0.40
1:A:8:ARG:NH2	3:A:606:HOH:O	2.54	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	236/256 (92%)	229 (97%)	6 (2%)	1 (0%)	39	27
1	B	238/256 (93%)	227 (95%)	10 (4%)	1 (0%)	39	27
All	All	474/512 (93%)	456 (96%)	16 (3%)	2 (0%)	39	27

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	243	ARG
1	B	244	LEU

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	196/209 (94%)	194 (99%)	2 (1%)	82	81
1	B	197/209 (94%)	191 (97%)	6 (3%)	48	38
All	All	393/418 (94%)	385 (98%)	8 (2%)	63	57

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

Mol	Chain	Res	Type
1	A	15	LEU
1	A	54	LEU
1	B	54	LEU
1	B	77	LEU

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Mol	Chain	Res	Type
1	B	162	ASP
1	B	223	LEU
1	B	243	ARG
1	B	244	LEU

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

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

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 ⓘ

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	240/256 (93%)	0.28	21 (8%) 12 14	12, 19, 46, 70	0
1	B	242/256 (94%)	0.52	24 (9%) 9 10	12, 22, 47, 66	0
All	All	482/512 (94%)	0.40	45 (9%) 11 12	12, 20, 47, 70	0

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	242	GLY	8.5
1	B	241	SER	8.5
1	A	244	LEU	7.1
1	A	242	GLY	6.7
1	A	243	ARG	6.7
1	B	243	ARG	6.5
1	A	241	SER	6.4
1	A	136	ASP	6.2
1	B	135	ASP	6.1
1	A	134	GLY	6.0
1	B	134	GLY	5.7
1	B	213	THR	5.1
1	A	230	HIS	4.9
1	A	240	ASN	4.8
1	A	249	ALA	4.1
1	A	227	ARG	4.1
1	A	245	ALA	3.8
1	A	212	PRO	3.7
1	A	135	ASP	3.5
1	A	133	VAL	3.4
1	B	220	ALA	3.4
1	B	249	ALA	3.4
1	B	244	LEU	3.4
1	B	212	PRO	3.2

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Mol	Chain	Res	Type	RSRZ
1	B	153	GLY	3.1
1	B	117	VAL	3.0
1	B	152	GLU	2.8
1	A	2	THR	2.8
1	B	132	THR	2.8
1	A	228	GLY	2.8
1	B	240	ASN	2.7
1	B	133	VAL	2.6
1	A	229	LEU	2.4
1	A	221	ALA	2.4
1	A	224	PRO	2.3
1	B	237	GLU	2.3
1	B	165	ALA	2.3
1	A	239	ALA	2.2
1	B	138	VAL	2.2
1	A	137	ASP	2.1
1	B	175	ALA	2.1
1	B	119	LEU	2.0
1	B	2	THR	2.0
1	B	25	LEU	2.0
1	B	137	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	MN	A	401	1/1	0.99	0.07	-	20,20,20,20	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	MN	B	401	1/1	0.99	0.08	-	21,21,21,21	0

6.5 Other polymers [i](#)

There are no such residues in this entry.