



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2014 – 03:15 AM GMT

PDB ID : 2QB4
Title : Crystal Structure Analysis of LeuT complexed with L-leucine, sodium and desipramine
Authors : Singh, S.K.; Yamashita, A.; Gouaux, E.
Deposited on : 2007-06-15
Resolution : 1.90 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

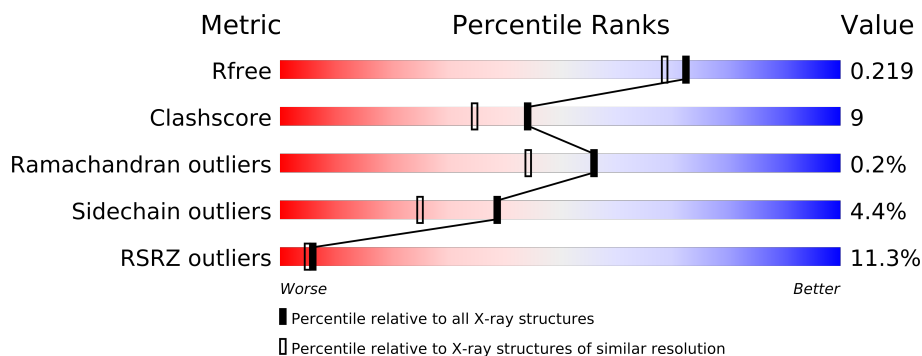
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance


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}	66092	3684 (1.90-1.90)
Clashscore	79885	4465 (1.90-1.90)
Ramachandran outliers	78287	4413 (1.90-1.90)
Sidechain outliers	78261	4414 (1.90-1.90)
RSRZ outliers	66119	3686 (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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	519	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
2	BOG	A	701	-	X
2	BOG	A	702	-	X
2	BOG	A	703	-	X
2	BOG	A	704	-	X
5	DSM	A	802	-	X

2 Entry composition i

There are 6 unique types of molecules in this entry. The entry contains 4341 atoms, of which 0 are hydrogen and 0 are deuterium.

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.

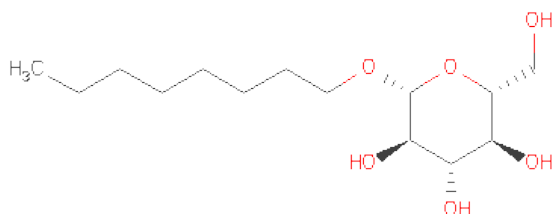
- Molecule 1 is a protein called Transporter.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	510	Total	C	N	O	S	0	1	0
			4049	2741	636	660	12			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	514	GLY	-	CLONING ARTIFACT	UNP O67854
A	515	THR	-	CLONING ARTIFACT	UNP O67854
A	516	LEU	-	CLONING ARTIFACT	UNP O67854
A	517	VAL	-	CLONING ARTIFACT	UNP O67854
A	518	PRO	-	CLONING ARTIFACT	UNP O67854
A	519	ARG	-	CLONING ARTIFACT	UNP O67854

- Molecule 2 is SUGAR (B-OCTYLGLUCOSIDE) (three-letter code: BOG) (formula: C₁₄H₂₈O₆).

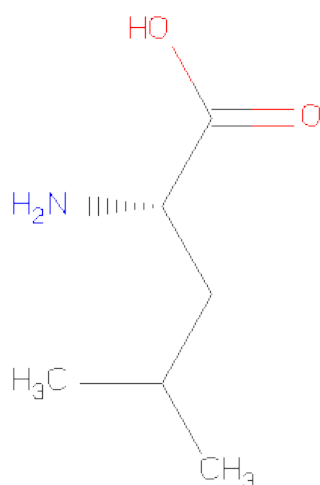


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			20	14	6		
2	A	1	Total	C	O	0	0
			20	14	6		
2	A	1	Total	C	O	0	0
			20	14	6		
2	A	1	Total	C	O	0	0
			20	14	6		
2	A	1	Total	C	O	0	0
			20	14	6		

- Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

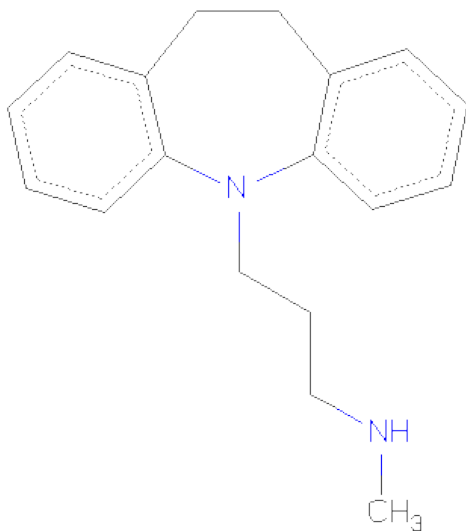
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Na	0	0
			2	2		

- Molecule 4 is LEUCINE (three-letter code: LEU) (formula: C₆H₁₃NO₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			9	6	1	2		

- Molecule 5 is 3-(10,11-DIHYDRO-5H-DIBENZO[B,F]AZEPIN-5-YL)-N-METHYLPROPA N-1-AMINE (three-letter code: DSM) (formula: C₁₈H₂₂N₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	N	0	0
			20	18	2		
5	A	1	Total	C	N	0	0
			20	18	2		

- Molecule 6 is water.

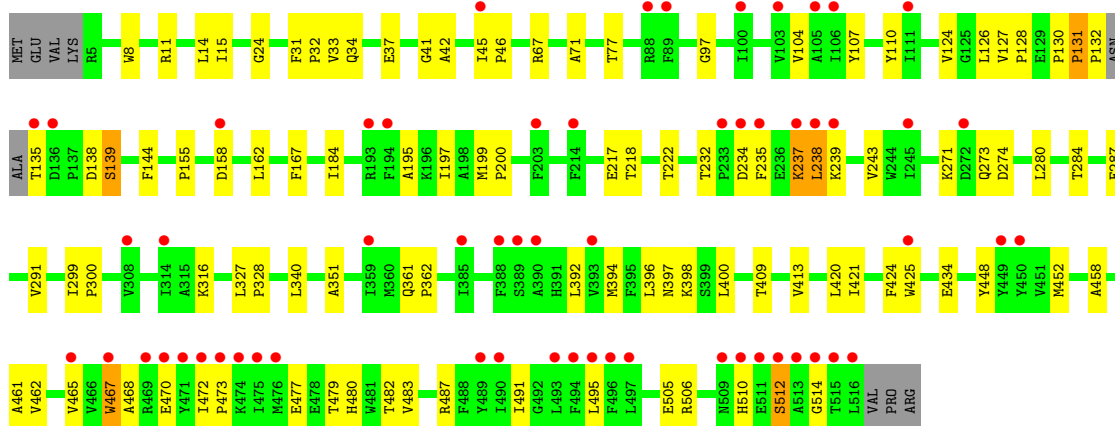
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	141	Total	O	0	0
			141	141		

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: Transporter

Chain A: 



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	88.29Å 86.77Å 81.25Å 90.00° 96.02° 90.00°	Depositor
Resolution (Å)	27.73 – 1.90 47.39 – 1.90	Depositor EDS
% Data completeness (in resolution range)	94.8 (27.73-1.90) 94.9 (47.39-1.90)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.52 (at 1.90Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.201 , 0.217 0.203 , 0.219	Depositor DCC
R_{free} test set	2265 reflections (4.96%)	DCC
Wilson B-factor (Å ²)	33.2	Xtriage
Anisotropy	0.345	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 70.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 48008 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4341	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.90% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NA, DSM, BOG

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	0.40	0/4176	0.59	1/5688 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	131	PRO	C-N-CD	-8.33	102.27	120.60

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4049	0	4127	70	0
2	A	100	0	140	7	0
3	A	2	0	0	0	0
4	A	9	0	10	0	0
5	A	40	0	44	2	0
6	A	141	0	0	3	0
All	All	4341	0	4321	73	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 9.

All (73) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:396:LEU:HB2	2:A:705:BOG:H8'3	1.70	0.73
1:A:237:LYS:HG2	1:A:243:VAL:HG21	1.70	0.71
1:A:45:ILE:HB	1:A:46:PRO:HD3	1.73	0.69
1:A:130:PRO:O	1:A:132:PRO:HD3	1.91	0.69
1:A:299:ILE:HB	1:A:300:PRO:HD3	1.75	0.68
1:A:468:ALA:HA	1:A:472:ILE:HD12	1.75	0.68
1:A:33:VAL:O	1:A:37:GLU:HB2	1.94	0.67
1:A:131:PRO:HD2	1:A:139:SER:O	2.00	0.62
1:A:124:VAL:HG23	1:A:126:LEU:HG	1.83	0.60
2:A:703:BOG:H2'2	2:A:704:BOG:H1'2	1.83	0.60
1:A:128:PRO:O	1:A:130:PRO:HD3	2.02	0.59
1:A:505:GLU:HB2	1:A:506:ARG:NH1	2.19	0.58
1:A:11:ARG:O	1:A:15:ILE:HG12	2.05	0.56
1:A:31:PHE:HB3	1:A:32:PRO:HD3	1.87	0.55
1:A:458:ALA:O	1:A:462:VAL:HG23	2.06	0.55
1:A:155:PRO:HD3	1:A:162:LEU:HD23	1.88	0.54
1:A:506:ARG:HH11	1:A:506:ARG:HG2	1.72	0.54
1:A:34:GLN:HE21	5:A:801:DSM:H3	1.72	0.54
1:A:420:LEU:C	1:A:420:LEU:HD23	2.28	0.54
1:A:77:THR:OG1	1:A:97:GLY:HA3	2.07	0.53
1:A:127:VAL:HB	1:A:128:PRO:HD2	1.89	0.53
1:A:479:THR:HB	1:A:483:VAL:HG11	1.91	0.52
1:A:280:LEU:O	1:A:284:THR:HG23	2.10	0.52
1:A:131:PRO:HG2	1:A:139:SER:OG	2.09	0.52
1:A:45:ILE:CG2	1:A:238:LEU:HD11	2.40	0.51
1:A:409:THR:HG22	1:A:467:TRP:CZ2	2.46	0.51
1:A:506:ARG:NH1	1:A:506:ARG:HG2	2.27	0.50
1:A:71:ALA:CB	1:A:271:LYS:HD2	2.42	0.49
1:A:480:HIS:HD2	1:A:482:THR:OG1	1.96	0.49
1:A:42:ALA:HB2	1:A:234:ASP:HB3	1.92	0.49
1:A:398:LYS:HA	6:A:952:HOH:O	2.13	0.48
1:A:392:LEU:HD13	2:A:704:BOG:H8'1	1.94	0.48
1:A:512:SER:C	1:A:514:GLY:H	2.17	0.48
1:A:110:TYR:CZ	1:A:394:MET:HG2	2.48	0.48
1:A:505:GLU:CB	1:A:506:ARG:NH1	2.77	0.48
1:A:299:ILE:HG12	6:A:928:HOH:O	2.14	0.47
1:A:472:ILE:N	1:A:473:PRO:CD	2.77	0.47
1:A:448:TYR:O	1:A:452:MET:HG3	2.15	0.47
1:A:409:THR:HG22	1:A:467:TRP:CH2	2.50	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:287:GLU:O	1:A:291:VAL:HG22	2.15	0.47
1:A:130:PRO:CG	1:A:144:PHE:HE2	2.28	0.46
1:A:461:ALA:O	1:A:465:VAL:HG23	2.15	0.46
1:A:131:PRO:CG	1:A:139:SER:OG	2.63	0.46
1:A:217:GLU:HG2	1:A:218:THR:N	2.30	0.45
1:A:199:MET:HB2	1:A:200:PRO:HD3	1.98	0.45
1:A:37:GLU:HG3	1:A:316:LYS:HG2	1.98	0.45
1:A:24:GLY:HA2	1:A:351:ALA:HB1	1.99	0.45
1:A:67:ARG:NH1	1:A:273:GLN:O	2.50	0.45
1:A:104:VAL:HA	1:A:107:TYR:CE2	2.52	0.45
1:A:131:PRO:HA	1:A:132:PRO:HD3	1.40	0.45
1:A:130:PRO:HG3	1:A:144:PHE:CE2	2.52	0.44
1:A:420:LEU:O	1:A:420:LEU:HD23	2.17	0.44
1:A:167:PHE:HB2	2:A:702:BOG:H1'1	1.99	0.44
1:A:217:GLU:HG3	1:A:222:THR:HG22	2.00	0.43
1:A:71:ALA:HB2	1:A:271:LYS:HD2	2.00	0.43
1:A:271:LYS:HG3	1:A:271:LYS:O	2.18	0.43
1:A:392:LEU:CD1	2:A:704:BOG:H8'1	2.48	0.43
1:A:505:GLU:HB2	1:A:506:ARG:HH12	1.83	0.43
1:A:468:ALA:CA	1:A:472:ILE:HD12	2.48	0.42
2:A:703:BOG:O6	2:A:704:BOG:H62	2.19	0.42
1:A:327:LEU:N	1:A:328:PRO:HD2	2.34	0.42
1:A:195:ALA:HB2	6:A:1029:HOH:O	2.18	0.42
1:A:400:LEU:C	1:A:400:LEU:HD23	2.40	0.42
2:A:703:BOG:H2	2:A:704:BOG:H5	2.02	0.42
1:A:421:ILE:O	1:A:425:TRP:HB2	2.19	0.42
1:A:413:VAL:HG21	1:A:467:TRP:HZ3	1.85	0.42
1:A:487:ARG:O	1:A:491:ILE:HG13	2.20	0.42
1:A:397:ASN:O	1:A:398:LYS:HB2	2.19	0.41
1:A:197:ILE:HD12	5:A:802:DSM:H10	2.03	0.41
1:A:110:TYR:CE1	1:A:394:MET:HG2	2.56	0.41
1:A:361:GLN:N	1:A:362:PRO:CD	2.84	0.41
1:A:41:GLY:HA3	1:A:232:THR:O	2.20	0.41
1:A:8:TRP:CD1	1:A:14:LEU:HD13	2.55	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	507/519 (98%)	489 (96%)	17 (3%)	1 (0%)	56 44

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	235	PHE

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	410/420 (98%)	392 (96%)	18 (4%)	39 25

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	135	THR
1	A	138	ASP
1	A	139	SER
1	A	158	ASP
1	A	184	ILE
1	A	237	LYS
1	A	238	LEU
1	A	239	LYS
1	A	274	ASP
1	A	340	LEU
1	A	424	PHE
1	A	434	GLU
1	A	467	TRP

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Mol	Chain	Res	Type
1	A	470	GLU
1	A	477	GLU
1	A	495	LEU
1	A	510	HIS
1	A	512	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	34	GLN
1	A	72	GLN
1	A	179	ASN
1	A	333	GLN
1	A	361	GLN
1	A	480	HIS
1	A	509	ASN

5.3.3 RNA ⓘ

There are no RNA chains 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 10 ligands modelled in this entry, 2 are monoatomic - leaving 8 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
4	LEU	A	601	3	8,8,8	0.68	0	10,10,10	0.40	0
2	BOG	A	701	-	20,20,20	1.18	3 (15%)	25,25,25	0.67	0
2	BOG	A	702	-	20,20,20	1.24	3 (15%)	25,25,25	0.90	1 (4%)
2	BOG	A	703	-	20,20,20	1.05	1 (5%)	25,25,25	0.66	0
2	BOG	A	704	-	20,20,20	1.13	2 (10%)	25,25,25	0.75	0
2	BOG	A	705	-	20,20,20	1.13	2 (10%)	25,25,25	0.77	0
5	DSM	A	801	-	22,22,22	1.84	4 (18%)	29,29,29	1.04	1 (3%)
5	DSM	A	802	-	22,22,22	1.90	4 (18%)	29,29,29	0.93	1 (3%)

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
4	LEU	A	601	3	-	0/8/8/8	0/0/0/0
2	BOG	A	701	-	-	0/11/31/31	0/1/1/1
2	BOG	A	702	-	-	0/11/31/31	0/1/1/1
2	BOG	A	703	-	-	0/11/31/31	0/1/1/1
2	BOG	A	704	-	-	0/11/31/31	0/1/1/1
2	BOG	A	705	-	-	0/11/31/31	0/1/1/1
5	DSM	A	801	-	-	0/5/18/18	0/0/3/3
5	DSM	A	802	-	-	0/5/18/18	0/0/3/3

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	802	DSM	C11-C12	6.46	1.49	1.40
5	A	801	DSM	C11-C12	6.18	1.49	1.40
2	A	702	BOG	O5-C1	3.12	1.49	1.41
2	A	705	BOG	O5-C1	2.92	1.49	1.41
2	A	701	BOG	O5-C1	2.91	1.49	1.41
2	A	704	BOG	O5-C1	2.90	1.49	1.41
2	A	703	BOG	O5-C1	2.86	1.49	1.41
2	A	702	BOG	O1-C1	2.58	1.44	1.40
5	A	801	DSM	C6-C12	2.53	1.56	1.51
5	A	802	DSM	C11-N1	2.52	1.47	1.43
2	A	701	BOG	C4-C5	2.41	1.58	1.53
2	A	705	BOG	O1-C1	2.39	1.44	1.40
5	A	802	DSM	C6-C12	2.22	1.56	1.51
2	A	704	BOG	C4-C5	2.22	1.57	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	801	DSM	C11-N1	2.18	1.46	1.43
2	A	701	BOG	O1-C1	2.15	1.44	1.40
2	A	702	BOG	C4-C5	2.14	1.57	1.53
5	A	801	DSM	C1-C14	2.08	1.43	1.39
5	A	802	DSM	C1-C14	2.00	1.43	1.39

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	702	BOG	C1'-O1-C1	2.72	118.86	113.96
5	A	801	DSM	C15-N1-C14	2.51	120.88	117.07
5	A	802	DSM	C12-C11-N1	2.08	124.39	120.12

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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	510/519 (98%)	0.78	59 (11%) 5 5	23, 36, 74, 96	0

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	513	ALA	12.2
1	A	514	GLY	11.9
1	A	515	THR	11.0
1	A	512	SER	10.6
1	A	516	LEU	8.8
1	A	475	ILE	8.5
1	A	510	HIS	8.0
1	A	472	ILE	5.7
1	A	471	TYR	5.4
1	A	511	GLU	5.3
1	A	308	VAL	5.2
1	A	136	ASP	5.0
1	A	238	LEU	4.8
1	A	476	MET	4.6
1	A	235	PHE	4.5
1	A	489	TYR	4.3
1	A	450	TYR	4.1
1	A	203	PHE	3.8
1	A	214	PHE	3.6
1	A	495	LEU	3.6
1	A	469	ARG	3.6
1	A	493	LEU	3.5
1	A	467	TRP	3.4
1	A	158	ASP	3.4
1	A	88	ARG	3.3
1	A	105	ALA	3.2
1	A	496	PHE	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	474	LYS	3.2
1	A	89	PHE	3.1
1	A	359	ILE	3.0
1	A	239	LYS	3.0
1	A	314	ILE	2.9
1	A	103	VAL	2.9
1	A	272	ASP	2.8
1	A	106	ILE	2.8
1	A	509	ASN	2.8
1	A	473	PRO	2.7
1	A	45	ILE	2.7
1	A	393	VAL	2.7
1	A	237	LYS	2.7
1	A	425	TRP	2.6
1	A	385	ILE	2.6
1	A	497	LEU	2.6
1	A	135	THR	2.6
1	A	389	SER	2.6
1	A	490	ILE	2.6
1	A	465	VAL	2.5
1	A	245	ILE	2.4
1	A	449	TYR	2.4
1	A	234	ASP	2.4
1	A	233	PRO	2.4
1	A	193	ARG	2.3
1	A	494	PHE	2.3
1	A	111	ILE	2.3
1	A	388	PHE	2.2
1	A	390	ALA	2.1
1	A	100	ILE	2.1
1	A	194	PHE	2.1
1	A	470	GLU	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	BOG	A	702	20/20	0.48	15.52	82,93,94,94	0
2	BOG	A	704	20/20	0.39	7.38	72,87,90,91	0
5	DSM	A	802	20/20	0.43	4.28	83,85,86,86	0
2	BOG	A	701	20/20	0.27	2.85	71,74,75,76	0
2	BOG	A	703	20/20	0.26	2.82	47,56,71,71	0
2	BOG	A	705	20/20	0.36	1.67	74,87,89,89	0
4	LEU	A	601	9/9	0.16	0.52	23,25,27,28	0
5	DSM	A	801	20/20	0.13	0.04	26,29,36,38	0
3	NA	A	751	1/1	0.08	-1.25	26,26,26,26	0
3	NA	A	752	1/1	0.09	-1.31	28,28,28,28	0

6.5 Other polymers ⓘ

There are no such residues in this entry.