Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 02:01 PM GMT

PDB ID	:	3VU1
Title	:	Crystal structure of the C-terminal globular domain of oligosaccharyltransferase (PhAglB-L, O74088_PYRHO) from Pyrococcus horikoshii
Authors	:	Nyirenda, J.; Matsumoto, S.; Saitoh, T.; Maita, N.; Noda, N.N.; Inagaki, F.; Kohda, D.
Deposited on	:	2012-06-13
Resolution	:	2.70 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

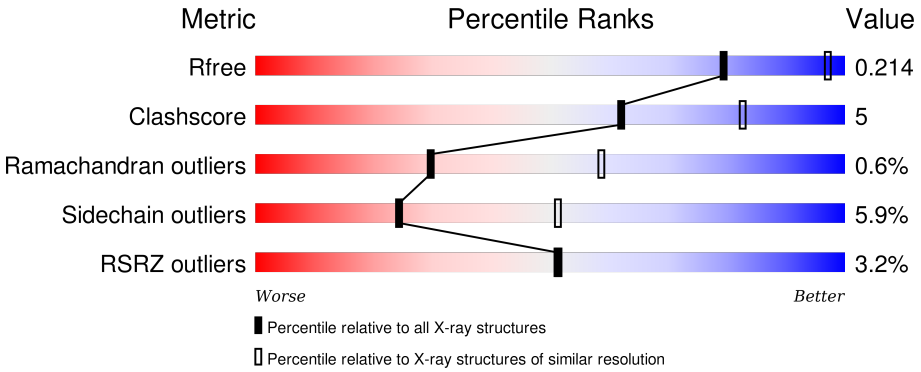
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	2103 (2.70-2.70)
Clashscore	102246	2422 (2.70-2.70)
Ramachandran outliers	100387	2382 (2.70-2.70)
Sidechain outliers	100360	2382 (2.70-2.70)
RSRZ outliers	91569	2107 (2.70-2.70)

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 <=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	506	2% 80% 14%
1	B	506	4% 80% 15%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7936 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 Putative uncharacterized protein PH0242.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	488	Total	C	N	O	S	0	1	0
			3942	2553	671	711	7			
1	B	490	Total	C	N	O	S	0	0	0
			3910	2528	659	716	7			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	481	MET	-	EXPRESSION TAG	UNP O74088
A	977	HIS	-	EXPRESSION TAG	UNP O74088
A	978	HIS	-	EXPRESSION TAG	UNP O74088
A	979	HIS	-	EXPRESSION TAG	UNP O74088
A	980	HIS	-	EXPRESSION TAG	UNP O74088
A	981	HIS	-	EXPRESSION TAG	UNP O74088
A	982	HIS	-	EXPRESSION TAG	UNP O74088
A	983	HIS	-	EXPRESSION TAG	UNP O74088
A	984	HIS	-	EXPRESSION TAG	UNP O74088
A	985	HIS	-	EXPRESSION TAG	UNP O74088
A	986	HIS	-	EXPRESSION TAG	UNP O74088
B	481	MET	-	EXPRESSION TAG	UNP O74088
B	977	HIS	-	EXPRESSION TAG	UNP O74088
B	978	HIS	-	EXPRESSION TAG	UNP O74088
B	979	HIS	-	EXPRESSION TAG	UNP O74088
B	980	HIS	-	EXPRESSION TAG	UNP O74088
B	981	HIS	-	EXPRESSION TAG	UNP O74088
B	982	HIS	-	EXPRESSION TAG	UNP O74088
B	983	HIS	-	EXPRESSION TAG	UNP O74088
B	984	HIS	-	EXPRESSION TAG	UNP O74088
B	985	HIS	-	EXPRESSION TAG	UNP O74088
B	986	HIS	-	EXPRESSION TAG	UNP O74088

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

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

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total Cl 1 1	0	0
3	A	1	Total Cl 1 1	0	0

- Molecule 4 is water.

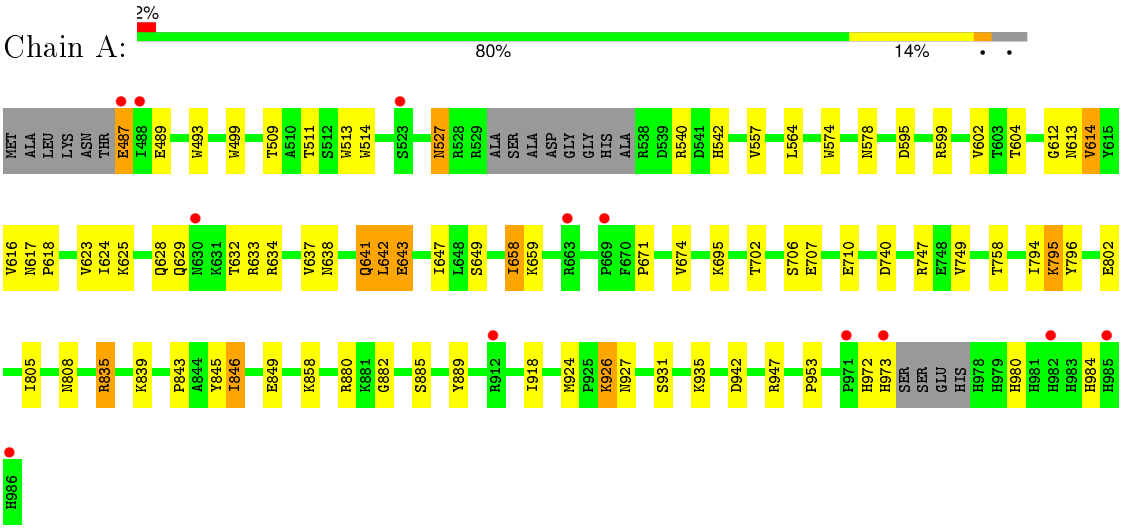
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	46	Total O 46 46	0	0
4	B	34	Total O 34 34	0	0

3 Residue-property plots (i)

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.

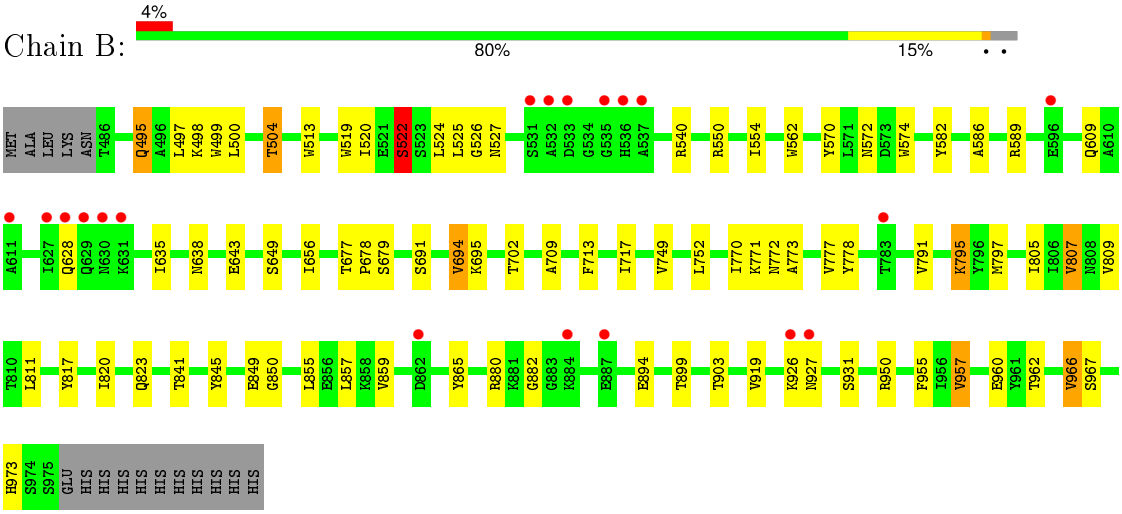
- Molecule 1: Putative uncharacterized protein PH0242

Chain A:



- Molecule 1: Putative uncharacterized protein PH0242

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	83.47Å 94.84Å 186.34Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.70 29.70 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.3 (30.00-2.70) 99.4 (29.70-2.70)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.16 (at 2.68Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.171 , 0.215 0.171 , 0.214	Depositor DCC
R_{free} test set	2081 reflections (5.33%)	DCC
Wilson B-factor (Å^2)	52.0	Xtriage
Anisotropy	0.129	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 42.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Outliers	0 of 41141 reflections	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	7936	wwPDB-VP
Average B, all atoms (Å^2)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.58% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: CA, CL

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	0.66	5/4045 (0.1%)	0.74	1/5477 (0.0%)
1	B	0.66	4/4004 (0.1%)	0.73	0/5425
All	All	0.66	9/8049 (0.1%)	0.74	1/10902 (0.0%)

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	493	TRP	CD2-CE2	5.85	1.48	1.41
1	A	513	TRP	CD2-CE2	5.82	1.48	1.41
1	B	513	TRP	CD2-CE2	5.80	1.48	1.41
1	A	574	TRP	CD2-CE2	5.73	1.48	1.41
1	B	499	TRP	CD2-CE2	5.38	1.47	1.41
1	A	514	TRP	CD2-CE2	5.22	1.47	1.41
1	B	574	TRP	CD2-CE2	5.12	1.47	1.41
1	B	519	TRP	CD2-CE2	5.12	1.47	1.41
1	A	499	TRP	CD2-CE2	5.07	1.47	1.41

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	835	ARG	NE-CZ-NH2	-5.15	117.72	120.30

There are no chirality outliers.

There are no planarity outliers.

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	3942	0	3932	40	0
1	B	3910	0	3914	46	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	46	0	0	0	0
4	B	34	0	0	2	0
All	All	7936	0	7846	84	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 5.

All (84) 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:972:HIS:HB2	1:A:973:HIS:CB	1.72	1.19
1:B:865:TYR:HB3	1:B:957:VAL:HG21	1.52	0.92
1:A:509:THR:HG23	1:A:527:ASN:HB2	1.53	0.90
1:A:846:ILE:HG12	1:A:953:PRO:HD3	1.64	0.78
1:B:497:LEU:HD21	1:B:520:ILE:HD11	1.67	0.75
1:B:771:LYS:H	1:B:823:GLN:HE22	1.36	0.72
1:B:850:GLY:HA2	1:B:926:LYS:HD3	1.69	0.72
1:A:927:ASN:O	1:A:931:SER:HB3	1.91	0.71
1:A:972:HIS:CB	1:A:973:HIS:CB	2.62	0.70
1:B:677:THR:HG22	1:B:679:SER:H	1.58	0.69
1:B:540:ARG:HG2	1:B:562:TRP:CZ2	2.32	0.65
1:B:500:LEU:O	1:B:504:THR:HB	1.97	0.65
1:A:578:ASN:HD21	1:A:604:THR:H	1.45	0.64
1:B:772:ASN:OD1	1:B:795:LYS:HE2	1.98	0.63
1:A:641:GLN:H	1:A:641:GLN:NE2	1.96	0.63
1:A:758:THR:OG1	1:A:808:ASN:ND2	2.32	0.63
1:B:771:LYS:H	1:B:823:GLN:NE2	1.96	0.63
1:B:927:ASN:O	1:B:931:SER:HB2	1.99	0.62
1:A:641:GLN:H	1:A:641:GLN:HE21	1.46	0.62
1:A:795:LYS:HD2	1:A:796:TYR:H	1.63	0.62
1:B:497:LEU:HD21	1:B:520:ILE:CD1	2.30	0.62
1:A:629:GLN:O	1:A:632:THR:HG22	2.00	0.61
1:B:522:SER:O	4:B:1134:HOH:O	2.16	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:773:ALA:HB2	1:B:823:GLN:HE21	1.67	0.60
1:A:595:ASP:HA	1:A:889:TYR:CD2	2.38	0.59
1:A:927:ASN:O	1:A:931:SER:CB	2.51	0.59
1:A:706:SER:O	1:A:707:GLU:HB2	2.02	0.59
1:A:880:ARG:HD2	1:A:882:GLY:O	2.04	0.58
1:A:740:ASP:HB2	1:A:749:VAL:HG11	1.86	0.57
1:B:752:LEU:HD12	1:B:966:VAL:HG21	1.86	0.57
1:B:791:VAL:O	1:B:807:VAL:HG22	2.04	0.57
1:A:695:LYS:HE3	1:A:702:THR:O	2.04	0.56
1:A:509:THR:CG2	1:A:527:ASN:HB2	2.32	0.55
1:B:849:GLU:OE2	1:B:927:ASN:HA	2.07	0.54
1:A:634:ARG:HA	1:A:658:ILE:HD11	1.89	0.54
1:B:865:TYR:CB	1:B:957:VAL:HG21	2.30	0.54
1:B:865:TYR:HB3	1:B:957:VAL:CG2	2.31	0.53
1:B:880:ARG:HD2	1:B:882:GLY:O	2.09	0.52
1:B:497:LEU:HD11	1:B:525:LEU:HD12	1.91	0.52
1:A:487:GLU:N	1:B:498:LYS:HZ3	2.07	0.52
1:B:695:LYS:HE3	1:B:702:THR:O	2.10	0.52
1:B:540:ARG:HD3	1:B:950:ARG:NH2	2.25	0.51
1:B:495:GLN:CD	1:B:495:GLN:H	2.15	0.50
1:B:582:TYR:OH	1:B:589:ARG:HD2	2.11	0.50
1:B:778:TYR:OH	1:B:960:GLU:OE2	2.26	0.50
1:B:927:ASN:O	1:B:931:SER:CB	2.59	0.49
1:B:635:ILE:HG13	1:B:649:SER:CB	2.42	0.49
1:A:642:LEU:H	1:A:642:LEU:HD23	1.78	0.49
1:A:511:THR:HG22	1:A:564:LEU:HD11	1.94	0.49
1:B:677:THR:HG23	1:B:678:PRO:HD2	1.95	0.48
1:B:850:GLY:CA	1:B:926:LYS:HD3	2.42	0.48
1:A:638:ASN:ND2	1:A:643:GLU:HB2	2.29	0.48
1:B:691:SER:OG	1:B:694:VAL:HG13	2.13	0.48
1:A:846:ILE:HD12	1:A:846:ILE:HA	1.63	0.47
1:A:647:ILE:HD11	1:A:671:PRO:HA	1.97	0.47
1:A:794:ILE:HG12	1:A:805:ILE:HD11	1.96	0.47
1:A:843:PRO:HA	1:A:846:ILE:CD1	2.45	0.46
1:B:550:ARG:HD3	1:B:586:ALA:HB2	1.98	0.46
1:A:835:ARG:HA	1:A:839:LYS:O	2.16	0.46
1:A:849:GLU:HA	1:A:924:MET:HB3	1.99	0.45
1:B:859:VAL:HG21	1:B:955:PHE:CE1	2.52	0.45
1:B:797:MET:HE1	1:B:805:ILE:HD11	1.99	0.45
1:B:865:TYR:CG	1:B:957:VAL:HG21	2.52	0.44
1:A:624:ILE:HG12	1:A:637:VAL:HG22	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:614:VAL:HG22	1:A:625:LYS:HB3	1.99	0.44
1:B:855:LEU:O	1:B:919:VAL:HA	2.18	0.44
1:B:770:ILE:HA	1:B:823:GLN:HE22	1.81	0.44
1:A:599:ARG:HG2	1:B:841:THR:HG22	1.99	0.44
1:B:709:ALA:O	1:B:713:PHE:CD2	2.71	0.44
1:A:617:ASN:HA	1:A:618:PRO:HD2	1.85	0.44
1:A:527:ASN:ND2	1:A:540:ARG:HH12	2.15	0.44
1:B:638:ASN:OD1	1:B:643:GLU:HG3	2.18	0.43
1:B:811:LEU:HD13	1:B:817:TYR:CE2	2.53	0.43
1:A:649:SER:HA	1:A:674:VAL:O	2.19	0.43
1:B:570:TYR:CZ	1:B:572:ASN:HB2	2.53	0.43
1:B:526:GLY:N	4:B:1134:HOH:O	2.50	0.43
1:A:926:LYS:O	1:A:927:ASN:HB2	2.19	0.42
1:B:797:MET:CE	1:B:805:ILE:HD11	2.49	0.42
1:B:777:VAL:HG11	1:B:809:VAL:HG21	2.01	0.42
1:A:616:VAL:HG12	1:A:623:VAL:HG22	2.02	0.42
1:A:527:ASN:HD21	1:A:540:ARG:HH12	1.67	0.41
1:A:542:HIS:NE2	1:A:947:ARG:HD2	2.35	0.41
1:B:540:ARG:HG2	1:B:562:TRP:CH2	2.56	0.41
1:A:794:ILE:HG12	1:A:805:ILE:CD1	2.51	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	483/506 (96%)	460 (95%)	21 (4%)	2 (0%)	39	69
1	B	488/506 (96%)	460 (94%)	24 (5%)	4 (1%)	24	51
All	All	971/1012 (96%)	920 (95%)	45 (5%)	6 (1%)	30	59

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	612	GLY
1	B	524	LEU
1	B	522	SER
1	B	527	ASN
1	A	613	ASN
1	B	973	HIS

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	418/432 (97%)	391 (94%)	27 (6%)	21 46
1	B	414/432 (96%)	392 (95%)	22 (5%)	28 57
All	All	832/864 (96%)	783 (94%)	49 (6%)	24 51

All (49) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	487	GLU
1	A	489	GLU
1	A	527	ASN
1	A	557	VAL
1	A	602	VAL
1	A	614	VAL
1	A	628	GLN
1	A	633	ARG
1	A	641	GLN
1	A	642	LEU
1	A	643	GLU
1	A	658	ILE
1	A	659	LYS
1	A	710	GLU
1	A	747	ARG
1	A	795	LYS
1	A	802	GLU
1	A	845	TYR

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Mol	Chain	Res	Type
1	A	846	ILE
1	A	858	LYS
1	A	885	SER
1	A	918	ILE
1	A	926	LYS
1	A	935	LYS
1	A	942	ASP
1	A	980	HIS
1	A	984	HIS
1	B	495	GLN
1	B	504	THR
1	B	522	SER
1	B	554	ILE
1	B	609	GLN
1	B	628	GLN
1	B	656	ILE
1	B	694	VAL
1	B	717	ILE
1	B	749	VAL
1	B	795	LYS
1	B	807	VAL
1	B	820	ILE
1	B	845	TYR
1	B	857	LEU
1	B	894	GLU
1	B	899	THR
1	B	903	THR
1	B	957	VAL
1	B	962	THR
1	B	966	VAL
1	B	967	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	578	ASN
1	A	628	GLN
1	A	638	ASN
1	A	641	GLN
1	A	808	ASN
1	A	815	GLN
1	B	503	ASN

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Mol	Chain	Res	Type
1	B	715	ASN
1	B	781	ASN
1	B	808	ASN
1	B	815	GLN
1	B	823	GLN

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 4 ligands modelled in this entry, 4 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, 95^{th} 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	488/506 (96%)	-0.12	12 (2%)	61	61	26, 44, 82, 131	0
1	B	490/506 (96%)	-0.04	19 (3%)	43	43	26, 46, 85, 129	0
All	All	978/1012 (96%)	-0.08	31 (3%)	51	51	26, 45, 85, 131	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	532	ALA	5.5
1	B	531	SER	5.0
1	A	487	GLU	4.6
1	B	535	GLY	4.5
1	A	986	HIS	4.3
1	A	630	ASN	4.0
1	B	533	ASP	3.9
1	B	537	ALA	3.8
1	B	629	GLN	3.3
1	A	523	SER	3.3
1	B	630	ASN	3.1
1	B	926	LYS	3.1
1	B	884	LYS	2.8
1	A	663	ARG	2.7
1	B	596	GLU	2.7
1	B	536	HIS	2.7
1	B	887	GLU	2.6
1	B	927	ASN	2.6
1	A	488	ILE	2.5
1	A	985	HIS	2.5
1	A	973	HIS	2.5
1	A	982	HIS	2.4
1	B	631	LYS	2.4
1	B	783	THR	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	611	ALA	2.4
1	A	669	PRO	2.3
1	B	862	ASP	2.3
1	B	628	GLN	2.3
1	A	971	PRO	2.2
1	A	912	ARG	2.2
1	B	627	ILE	2.1

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, 95^{th} 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
3	CL	A	1002	1/1	0.98	0.18	-0.13	41,41,41,41	0
2	CA	A	1001	1/1	0.98	0.14	-0.73	49,49,49,49	0
2	CA	B	1001	1/1	0.98	0.11	-2.64	65,65,65,65	0
3	CL	B	1002	1/1	0.94	0.12	-2.95	53,53,53,53	0

6.5 Other polymers (i)

There are no such residues in this entry.