



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:09 PM GMT

PDB ID : 3WAK
Title : Crystal structure of the Archaeoglobus fulgidus oligosaccharyltransferase (O29867_ARCFU) in the apo form
Authors : Matsumoto, S.; Shimada, A.; Kohda, D.
Deposited on : 2013-05-03
Resolution : 3.41 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

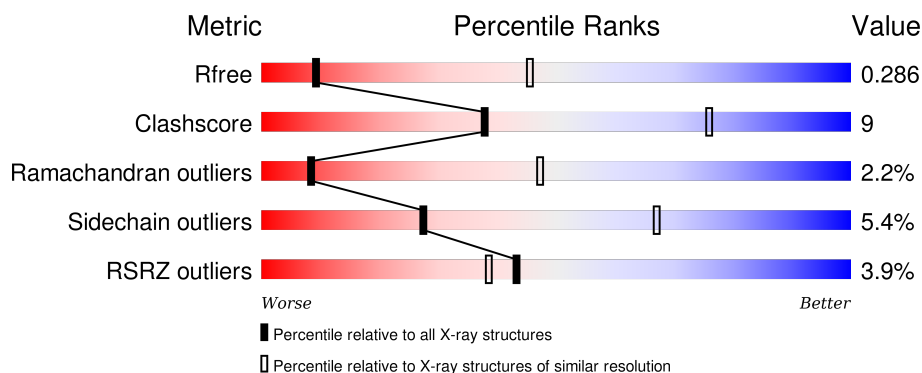
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.41 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	1049 (3.52-3.32)
Clashscore	102246	1032 (3.50-3.34)
Ramachandran outliers	100387	1002 (3.50-3.34)
Sidechain outliers	100360	1003 (3.50-3.34)
RSRZ outliers	91569	1054 (3.52-3.32)

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	875	<div> <div>4%</div> <div>72%</div> <div>24%</div> <div>..</div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6955 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 Transmembrane oligosaccharyl transferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	864	Total	C	N	O	S	0	0	0
			6954	4619	1108	1211	16			

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	869	GLU	-	EXPRESSION TAG	UNP O29867
A	870	LEU	-	EXPRESSION TAG	UNP O29867
A	871	ALA	-	EXPRESSION TAG	UNP O29867
A	872	LEU	-	EXPRESSION TAG	UNP O29867
A	873	VAL	-	EXPRESSION TAG	UNP O29867
A	874	PRO	-	EXPRESSION TAG	UNP O29867
A	875	ARG	-	EXPRESSION TAG	UNP O29867

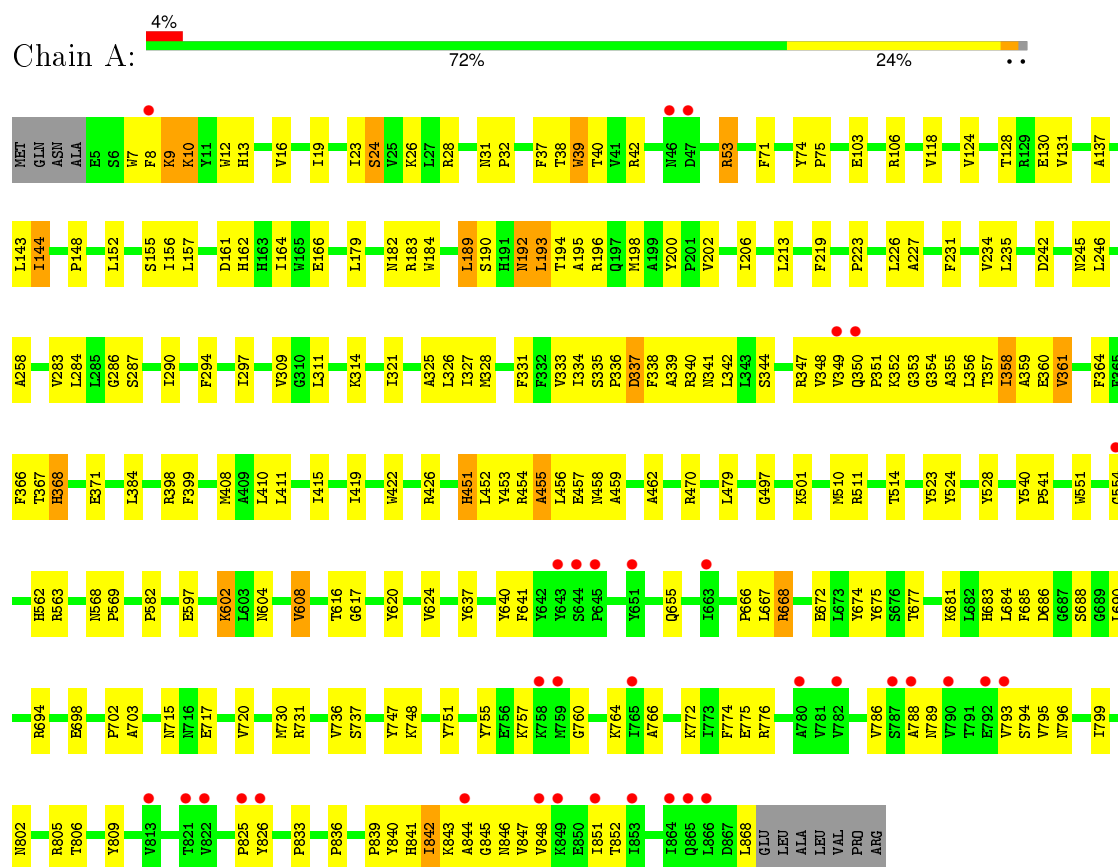
- Molecule 2 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

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

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: Transmembrane oligosaccharyl transferase



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	123.36Å 123.36Å 182.50Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.90 – 3.41 49.90 – 3.41	Depositor EDS
% Data completeness (in resolution range)	98.8 (49.90-3.41) 99.0 (49.90-3.41)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.12 (at 3.40Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, R_{free}	0.219 , 0.272 0.227 , 0.286	Depositor DCC
R_{free} test set	1000 reflections (5.38%)	DCC
Wilson B-factor (Å ²)	134.7	Xtriage
Anisotropy	0.517	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 73.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 19596 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6955	wwPDB-VP
Average B, all atoms (Å ²)	108.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.72% 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

5.1 Standard geometry

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	0.24	0/7182	0.42	0/9803

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

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	6954	0	6852	127	0
2	A	1	0	0	0	0
All	All	6955	0	6852	127	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 (127) 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:124:VAL:HG21	1:A:144:ILE:HG13	1.65	0.79
1:A:131:VAL:HG21	1:A:410:LEU:HD22	1.66	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:352:LYS:HD3	1:A:356:LEU:HG	1.68	0.75
1:A:357:THR:HA	1:A:358:ILE:HG22	1.74	0.70
1:A:192:ASN:N	1:A:192:ASN:OD1	2.24	0.70
1:A:843:LYS:NZ	1:A:845:GLY:O	2.25	0.69
1:A:451:HIS:O	1:A:453:TYR:N	2.26	0.68
1:A:24:SER:HB3	1:A:118:VAL:HG22	1.74	0.68
1:A:354:GLY:HA2	1:A:355:ALA:HB3	1.76	0.67
1:A:358:ILE:HG21	1:A:361:VAL:HG12	1.76	0.67
1:A:764:LYS:NZ	1:A:806:THR:O	2.29	0.66
1:A:26:LYS:O	1:A:31:ASN:ND2	2.27	0.65
1:A:458:ASN:HA	1:A:462:ALA:H	1.63	0.64
1:A:455:ALA:O	1:A:457:GLU:N	2.34	0.61
1:A:684:LEU:HD11	1:A:748:LYS:HG3	1.81	0.61
1:A:183:ARG:NH1	1:A:200:TYR:OH	2.35	0.60
1:A:326:LEU:O	1:A:328:MET:N	2.35	0.60
1:A:455:ALA:HB1	1:A:459:ALA:HB2	1.83	0.60
1:A:690:LEU:O	1:A:809:TYR:OH	2.15	0.59
1:A:352:LYS:HD2	1:A:361:VAL:HG21	1.83	0.59
1:A:28:ARG:HE	1:A:152:LEU:HD12	1.66	0.58
1:A:356:LEU:HD12	1:A:358:ILE:HB	1.86	0.58
1:A:551:TRP:HE3	1:A:569:PRO:HA	1.68	0.57
1:A:802:ASN:HB3	1:A:833:PRO:HB2	1.87	0.57
1:A:336:PRO:HA	1:A:337:ASP:C	2.25	0.57
1:A:334:ILE:HG23	1:A:335:SER:HB3	1.85	0.56
1:A:356:LEU:HA	1:A:357:THR:C	2.25	0.56
1:A:698:GLU:OE1	1:A:805:ARG:NH2	2.31	0.56
1:A:674:TYR:O	1:A:677:THR:OG1	2.18	0.56
1:A:748:LYS:NZ	1:A:766:ALA:O	2.34	0.55
1:A:608:VAL:HG13	1:A:774:PHE:HB2	1.89	0.55
1:A:842:ILE:HD11	1:A:851:ILE:HG12	1.89	0.55
1:A:7:TRP:HA	1:A:9:LYS:H	1.72	0.54
1:A:511:ARG:O	1:A:563:ARG:NH2	2.41	0.54
1:A:193:LEU:HD22	1:A:198:MET:HG3	1.91	0.53
1:A:336:PRO:HA	1:A:338:PHE:N	2.24	0.53
1:A:640:TYR:HE1	1:A:655:GLN:HE22	1.57	0.53
1:A:161:ASP:OD1	1:A:162:HIS:ND1	2.42	0.53
1:A:717:GLU:HA	1:A:720:VAL:HG12	1.91	0.53
1:A:637:TYR:HB3	1:A:668:ARG:HG2	1.89	0.53
1:A:193:LEU:HB2	1:A:198:MET:HB2	1.90	0.52
1:A:562:HIS:O	1:A:563:ARG:NH1	2.37	0.52
1:A:683:HIS:O	1:A:772:LYS:NZ	2.41	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:53:ARG:NH1	1:A:524:TYR:O	2.35	0.52
1:A:42:ARG:NH1	1:A:497:GLY:O	2.43	0.52
1:A:336:PRO:HB3	1:A:339:ALA:HB3	1.91	0.52
1:A:242:ASP:OD2	1:A:245:ASN:ND2	2.43	0.52
1:A:795:VAL:HB	1:A:842:ILE:HG23	1.92	0.51
1:A:28:ARG:HA	1:A:156:ILE:HD12	1.93	0.51
1:A:840:TYR:HB2	1:A:851:ILE:HG13	1.93	0.51
1:A:193:LEU:HD12	1:A:194:THR:H	1.76	0.51
1:A:350:GLN:HB3	1:A:351:PRO:HA	1.93	0.51
1:A:616:THR:OG1	1:A:617:GLY:N	2.44	0.51
1:A:731:ARG:HB3	1:A:736:VAL:HG23	1.92	0.50
1:A:455:ALA:O	1:A:458:ASN:N	2.45	0.50
1:A:161:ASP:OD1	1:A:162:HIS:N	2.41	0.49
1:A:143:LEU:HD22	1:A:384:LEU:HD21	1.95	0.49
1:A:53:ARG:NH2	1:A:528:TYR:OH	2.46	0.49
1:A:258:ALA:HB2	1:A:284:LEU:HB3	1.94	0.49
1:A:793:VAL:HG22	1:A:844:ALA:HB2	1.95	0.48
1:A:360:GLU:HB3	1:A:426:ARG:HB2	1.96	0.48
1:A:355:ALA:H	1:A:624:VAL:HG12	1.78	0.48
1:A:454:ARG:HA	1:A:455:ALA:O	2.13	0.48
1:A:321:ILE:O	1:A:325:ALA:N	2.46	0.48
1:A:219:PHE:HD1	1:A:283:VAL:HG11	1.79	0.48
1:A:788:ALA:HA	1:A:789:ASN:HA	1.44	0.48
1:A:666:PRO:HG2	1:A:737:SER:HA	1.96	0.47
1:A:686:ASP:OD1	1:A:805:ARG:NH2	2.37	0.47
1:A:826:TYR:O	1:A:840:TYR:OH	2.28	0.47
1:A:294:PHE:HA	1:A:297:ILE:HD12	1.96	0.46
1:A:367:THR:OG1	1:A:368:HIS:N	2.47	0.46
1:A:348:VAL:HG12	1:A:350:GLN:HG3	1.96	0.46
1:A:231:PHE:HA	1:A:234:VAL:HG22	1.96	0.46
1:A:582:PRO:HG3	1:A:602:LYS:HG2	1.98	0.46
1:A:194:THR:O	1:A:196:ARG:N	2.48	0.46
1:A:685:PHE:HB3	1:A:688:SER:HB3	1.97	0.46
1:A:411:LEU:O	1:A:415:ILE:HG12	2.16	0.45
1:A:184:TRP:CZ2	1:A:193:LEU:HB3	2.52	0.45
1:A:234:VAL:HG23	1:A:235:LEU:HG	1.98	0.45
1:A:399:PHE:CE2	1:A:408:MET:HB2	2.51	0.45
1:A:39:TRP:CE3	1:A:40:THR:HG22	2.52	0.45
1:A:451:HIS:CE1	1:A:470:ARG:HD2	2.52	0.44
1:A:309:VAL:HG23	1:A:311:LEU:HG	1.99	0.44
1:A:162:HIS:O	1:A:166:GLU:HG3	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:675:TYR:CE1	1:A:751:TYR:HB2	2.53	0.44
1:A:551:TRP:CE3	1:A:569:PRO:HA	2.51	0.43
1:A:19:ILE:O	1:A:23:ILE:HG13	2.18	0.43
1:A:540:TYR:HA	1:A:541:PRO:HD3	1.80	0.43
1:A:455:ALA:HA	1:A:458:ASN:OD1	2.18	0.43
1:A:640:TYR:CZ	1:A:668:ARG:HD3	2.52	0.43
1:A:226:LEU:HD11	1:A:287:SER:HB3	2.00	0.43
1:A:667:LEU:HD21	1:A:747:TYR:HE1	1.84	0.43
1:A:9:LYS:HA	1:A:10:LYS:HA	1.76	0.43
1:A:451:HIS:HB3	1:A:454:ARG:O	2.19	0.43
1:A:103:GLU:HG2	1:A:106:ARG:HH12	1.83	0.43
1:A:356:LEU:HD12	1:A:358:ILE:HD13	2.01	0.43
1:A:366:PHE:HE2	1:A:422:TRP:CZ2	2.37	0.43
1:A:189:LEU:HA	1:A:190:SER:HA	1.71	0.43
1:A:358:ILE:HD12	1:A:358:ILE:HA	1.92	0.42
1:A:358:ILE:HG13	1:A:359:ALA:H	1.84	0.42
1:A:130:GLU:OE1	1:A:182:ASN:ND2	2.45	0.42
1:A:341:ASN:O	1:A:344:SER:OG	2.23	0.42
1:A:799:ILE:HA	1:A:836:PRO:HA	2.00	0.42
1:A:74:TYR:HA	1:A:75:PRO:HA	1.74	0.42
1:A:223:PRO:O	1:A:227:ALA:N	2.48	0.42
1:A:597:GLU:OE1	1:A:776:ARG:NH1	2.44	0.42
1:A:39:TRP:CG	1:A:501:LYS:HE2	2.54	0.42
1:A:355:ALA:O	1:A:356:LEU:HD13	2.19	0.42
1:A:751:TYR:OH	1:A:760:GLY:O	2.38	0.42
1:A:523:TYR:O	1:A:528:TYR:OH	2.30	0.42
1:A:128:THR:HB	1:A:137:ALA:HA	2.01	0.42
1:A:510:MET:HG3	1:A:514:THR:HG21	2.02	0.41
1:A:554:GLY:HA3	1:A:568:ASN:HA	2.02	0.41
1:A:12:TRP:O	1:A:16:VAL:HG12	2.20	0.41
1:A:694:ARG:NH2	1:A:775:GLU:OE1	2.50	0.41
1:A:202:VAL:O	1:A:206:ILE:HG13	2.20	0.41
1:A:620:TYR:O	1:A:624:VAL:HG23	2.20	0.41
1:A:32:PRO:HG3	1:A:156:ILE:HG13	2.02	0.41
1:A:10:LYS:HB2	1:A:13:HIS:NE2	2.35	0.41
1:A:514:THR:OG1	1:A:563:ARG:NE	2.47	0.41
1:A:681:LYS:HA	1:A:685:PHE:HD2	1.87	0.40
1:A:415:ILE:O	1:A:419:ILE:HG12	2.22	0.40
1:A:155:SER:HB3	1:A:164:ILE:HD13	2.03	0.40
1:A:672:GLU:HG3	1:A:755:TYR:HB3	2.03	0.40
1:A:37:PHE:HE2	1:A:157:LEU:HD21	1.86	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:286:GLY:O	1:A:290:ILE:HG13	2.21	0.40
1:A:839:PRO:O	1:A:841:HIS:ND1	2.53	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	862/875 (98%)	775 (90%)	68 (8%)	19 (2%)	8 47

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	358	ILE
1	A	452	LEU
1	A	456	LEU
1	A	455	ALA
1	A	195	ALA
1	A	327	ILE
1	A	337	ASP
1	A	9	LYS
1	A	347	ARG
1	A	353	GLY
1	A	451	HIS
1	A	703	ALA
1	A	148	PRO
1	A	757	LYS
1	A	825	PRO
1	A	702	PRO
1	A	786	VAL
1	A	349	VAL
1	A	333	VAL

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	718/727 (99%)	679 (95%)	39 (5%)	27 67

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	PHE
1	A	10	LYS
1	A	24	SER
1	A	38	THR
1	A	39	TRP
1	A	53	ARG
1	A	71	PHE
1	A	144	ILE
1	A	179	LEU
1	A	189	LEU
1	A	192	ASN
1	A	193	LEU
1	A	213	LEU
1	A	246	LEU
1	A	314	LYS
1	A	331	PHE
1	A	340	ARG
1	A	342	LEU
1	A	361	VAL
1	A	364	PHE
1	A	368	HIS
1	A	371	GLU
1	A	398	ARG
1	A	479	LEU
1	A	602	LYS
1	A	604	ASN
1	A	608	VAL
1	A	641	PHE
1	A	668	ARG
1	A	715	ASN

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Mol	Chain	Res	Type
1	A	730	MET
1	A	794	SER
1	A	796	ASN
1	A	842	ILE
1	A	846	ASN
1	A	847	VAL
1	A	848	VAL
1	A	852	THR
1	A	868	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	245	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 1 ligands modelled in this entry, 1 is 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 [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 ⓘ

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	864/875 (98%)	0.15	34 (3%) 43 38	50, 102, 169, 204	0

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	866	LEU	4.0
1	A	759	MET	4.0
1	A	826	TYR	3.9
1	A	758	LYS	3.8
1	A	8	PHE	3.6
1	A	47	ASP	3.6
1	A	349	VAL	3.5
1	A	350	GLN	3.5
1	A	822	VAL	3.5
1	A	864	ILE	3.4
1	A	865	GLN	3.4
1	A	792	GLU	3.0
1	A	790	VAL	2.9
1	A	782	VAL	2.9
1	A	787	SER	2.9
1	A	851	ILE	2.9
1	A	46	ASN	2.7
1	A	788	ALA	2.6
1	A	780	ALA	2.5
1	A	813	VAL	2.4
1	A	765	ILE	2.4
1	A	844	ALA	2.4
1	A	643	TYR	2.3
1	A	853	ILE	2.3
1	A	645	PRO	2.3
1	A	644	SER	2.3
1	A	651	TYR	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	825	PRO	2.2
1	A	554	GLY	2.2
1	A	849	LYS	2.2
1	A	821	THR	2.2
1	A	793	VAL	2.1
1	A	848	VAL	2.1
1	A	663	ILE	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(\AA^2)	Q<0.9
2	MN	A	1200	1/1	0.98	0.49	-	77,77,77,77	0

6.5 Other polymers [i](#)

There are no such residues in this entry.