



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 09:14 pm GMT

PDB ID : 3RHZ
Title : Structure and functional analysis of a new subfamily of glycosyltransferases required for glycosylation of serine-rich streptococcal adhesions
Authors : Zhu, F.; Li, J.; Wu, H.
Deposited on : 2011-04-12
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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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) : recalc28949

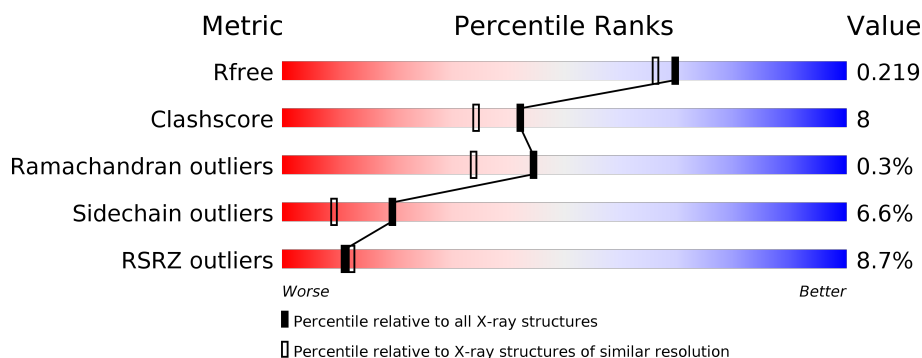
1 Overall quality at a glance

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}	100719	5047 (1.90-1.90)
Clashscore	112137	5731 (1.90-1.90)
Ramachandran outliers	110173	5669 (1.90-1.90)
Sidechain outliers	110143	5670 (1.90-1.90)
RSRZ outliers	101464	5100 (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	339	<div> <div>2%</div> <div> <div></div> <div>81%</div> <div>14%</div> <div>• •</div> </div> </div>
1	B	339	<div> <div>15%</div> <div> <div></div> <div>77%</div> <div>16%</div> <div>• •</div> </div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5727 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.

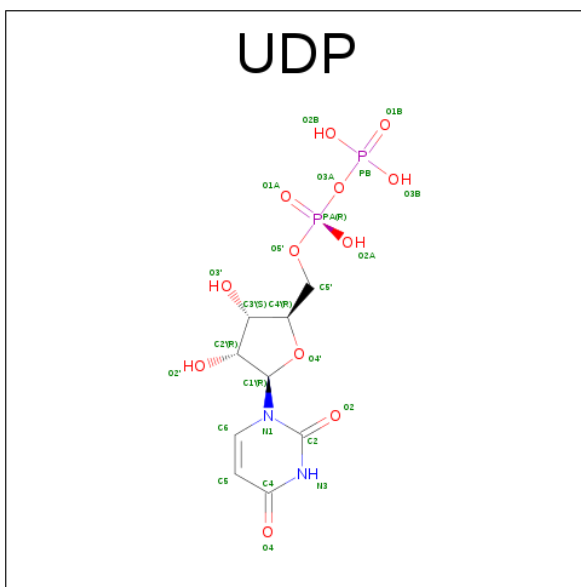
- Molecule 1 is a protein called Nucleotide sugar synthetase-like protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	330	Total	C	N	O	S	0	0	0
			2676	1722	445	492	17			
1	B	326	Total	C	N	O	S	0	0	0
			2651	1706	441	487	17			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-9	SER	-	EXPRESSION TAG	UNP B5A7L9
A	-8	GLU	-	EXPRESSION TAG	UNP B5A7L9
A	-7	PHE	-	EXPRESSION TAG	UNP B5A7L9
A	-6	GLU	-	EXPRESSION TAG	UNP B5A7L9
A	-5	LEU	-	EXPRESSION TAG	UNP B5A7L9
A	-4	ARG	-	EXPRESSION TAG	UNP B5A7L9
A	-3	ARG	-	EXPRESSION TAG	UNP B5A7L9
A	-2	GLN	-	EXPRESSION TAG	UNP B5A7L9
A	-1	ALA	-	EXPRESSION TAG	UNP B5A7L9
A	0	CYS	-	EXPRESSION TAG	UNP B5A7L9
B	-9	SER	-	EXPRESSION TAG	UNP B5A7L9
B	-8	GLU	-	EXPRESSION TAG	UNP B5A7L9
B	-7	PHE	-	EXPRESSION TAG	UNP B5A7L9
B	-6	GLU	-	EXPRESSION TAG	UNP B5A7L9
B	-5	LEU	-	EXPRESSION TAG	UNP B5A7L9
B	-4	ARG	-	EXPRESSION TAG	UNP B5A7L9
B	-3	ARG	-	EXPRESSION TAG	UNP B5A7L9
B	-2	GLN	-	EXPRESSION TAG	UNP B5A7L9
B	-1	ALA	-	EXPRESSION TAG	UNP B5A7L9
B	0	CYS	-	EXPRESSION TAG	UNP B5A7L9

- Molecule 2 is URIDINE-5'-DIPHOSPHATE (three-letter code: UDP) (formula: C₉H₁₄N₂O₁₂P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			25	9	2	12	2		
2	B	1	Total	C	N	O	P	0	0
			25	9	2	12	2		

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

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

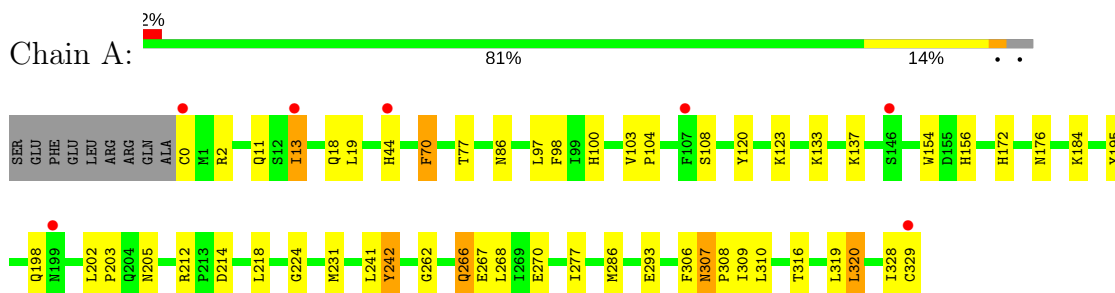
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	212	Total	O	0	0
			212	212		
4	B	137	Total	O	0	0
			137	137		

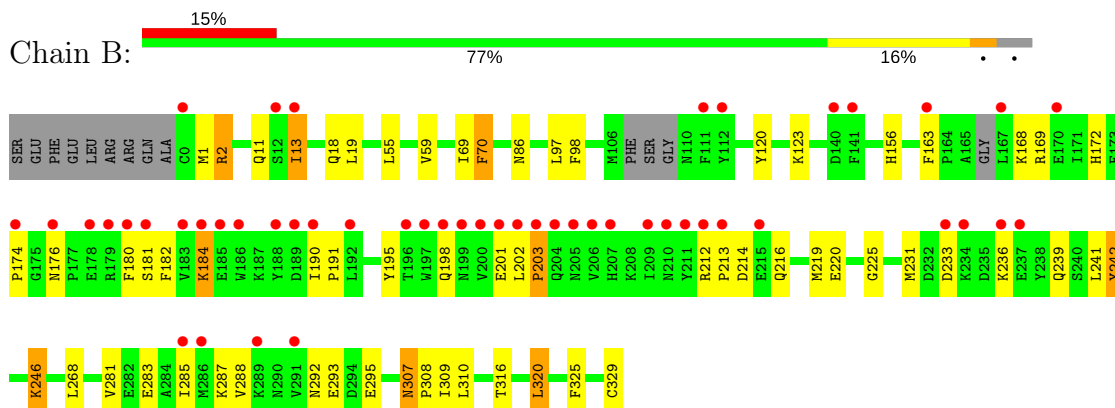
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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: Nucleotide sugar synthetase-like protein



- Molecule 1: Nucleotide sugar synthetase-like protein



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	77.18Å 193.86Å 96.25Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.29 – 1.90 29.29 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.8 (29.29-1.90) 99.8 (29.29-1.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.36 (at 1.89Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.1_351)	Depositor
R, R_{free}	0.182 , 0.223 0.180 , 0.219	Depositor DCC
R_{free} test set	2901 reflections (5.07%)	DCC
Wilson B-factor (Å ²)	23.3	Xtriage
Anisotropy	0.379	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 50.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5727	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.28% 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: UDP, 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	0.39	1/2737 (0.0%)	0.56	0/3706
1	B	0.48	1/2709 (0.0%)	0.57	2/3666 (0.1%)
All	All	0.44	2/5446 (0.0%)	0.56	2/7372 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	329	CYS	CB-SG	-18.23	1.51	1.82
1	A	0	CYS	C-N	-5.61	1.21	1.34

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	2	ARG	N-CA-C	-8.56	87.88	111.00
1	B	1	MET	CB-CA-C	-5.49	99.42	110.40

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	2676	0	2666	41	0
1	B	2651	0	2644	48	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	25	0	11	0	0
2	B	25	0	11	2	0
3	B	1	0	0	0	0
4	A	212	0	0	5	0
4	B	137	0	0	1	0
All	All	5727	0	5332	85	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 8.

All (85) 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:B:212:ARG:HB2	1:B:213:PRO:HD2	1.57	0.86
1:A:11:GLN:H	1:A:18:GLN:HE22	1.25	0.84
1:A:156:HIS:HE1	1:A:214:ASP:OD2	1.67	0.78
1:A:328:ILE:O	1:A:329:CYS:SG	2.43	0.77
1:B:156:HIS:HE1	1:B:214:ASP:OD2	1.69	0.75
1:A:172:HIS:HD2	1:A:195:TYR:OH	1.71	0.74
1:B:283:GLU:O	1:B:287:LYS:HG2	1.89	0.73
1:A:262:GLY:H	1:A:266:GLN:HE22	1.36	0.71
1:B:11:GLN:H	1:B:18:GLN:HE22	1.37	0.70
1:A:44:HIS:HD2	4:A:385:HOH:O	1.77	0.67
1:B:180:PHE:HB3	1:B:182:PHE:CE1	2.31	0.65
1:B:201:GLU:HG3	4:B:377:HOH:O	1.97	0.64
1:A:316:THR:HG22	1:A:320:LEU:HD22	1.79	0.64
1:A:309:ILE:HD11	1:B:325:PHE:CD1	2.33	0.63
1:A:262:GLY:H	1:A:266:GLN:NE2	1.99	0.60
1:B:292:ASN:OD1	1:B:295:GLU:HG3	2.02	0.60
1:A:11:GLN:H	1:A:18:GLN:NE2	1.96	0.59
1:B:213:PRO:HD2	1:B:216:GLN:NE2	2.18	0.59
1:A:86:ASN:HD21	1:A:123:LYS:HZ2	1.50	0.59
1:B:11:GLN:H	1:B:18:GLN:NE2	2.01	0.59
1:A:44:HIS:CD2	4:A:385:HOH:O	2.54	0.58
1:B:202:LEU:HB3	1:B:203:PRO:HD2	1.85	0.58
1:B:233:ASP:HA	1:B:236:LYS:HG3	1.85	0.57
1:A:172:HIS:CD2	1:A:195:TYR:OH	2.56	0.57
1:A:86:ASN:HD21	1:A:123:LYS:NZ	2.03	0.57
1:B:86:ASN:HD21	1:B:123:LYS:NZ	2.02	0.56
1:A:328:ILE:HG21	1:B:309:ILE:HG13	1.88	0.56
1:B:172:HIS:HD2	1:B:195:TYR:OH	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:316:THR:HG22	1:B:320:LEU:HD22	1.87	0.56
1:A:328:ILE:C	1:A:329:CYS:SG	2.85	0.55
1:A:13:ILE:O	1:A:13:ILE:HD13	2.08	0.54
1:A:270:GLU:HG3	1:A:277:ILE:HD11	1.87	0.54
1:B:168:LYS:O	1:B:225:GLY:HA2	2.06	0.54
1:A:270:GLU:HG3	1:A:277:ILE:CD1	2.37	0.54
1:A:286:MET:HG3	4:A:390:HOH:O	2.07	0.54
1:B:163:PHE:HE1	1:B:219:MET:HB3	1.74	0.53
1:A:328:ILE:O	1:A:329:CYS:CB	2.57	0.53
1:B:212:ARG:HB2	1:B:216:GLN:NE2	2.24	0.53
1:B:176:ASN:OD1	1:B:198:GLN:HG3	2.08	0.53
1:B:213:PRO:HG2	1:B:216:GLN:HE21	1.73	0.53
1:A:176:ASN:HB2	4:A:429:HOH:O	2.10	0.52
1:A:307:ASN:HD22	1:A:308:PRO:HD3	1.74	0.52
1:A:133:LYS:HD2	4:A:396:HOH:O	2.10	0.51
1:A:86:ASN:ND2	1:A:123:LYS:HZ2	2.10	0.50
1:A:309:ILE:HD11	1:B:325:PHE:CE1	2.45	0.50
1:B:246:LYS:HE2	2:B:7517:UDP:O3B	2.12	0.50
1:A:176:ASN:OD1	1:A:198:GLN:HG3	2.11	0.50
1:B:86:ASN:HD21	1:B:123:LYS:HZ2	1.59	0.49
1:A:11:GLN:HB2	1:A:18:GLN:NE2	2.27	0.49
1:A:86:ASN:ND2	1:A:123:LYS:NZ	2.60	0.49
1:A:307:ASN:HD22	1:A:307:ASN:C	2.15	0.49
1:A:328:ILE:CG2	1:B:309:ILE:HG13	2.44	0.47
1:B:55:LEU:O	1:B:59:VAL:HG22	2.14	0.47
1:A:267:GLU:OE2	1:A:268:LEU:HD13	2.14	0.47
1:B:307:ASN:HD22	1:B:307:ASN:C	2.17	0.47
1:B:307:ASN:HD22	1:B:308:PRO:HD3	1.80	0.46
1:B:216:GLN:O	1:B:220:GLU:HG3	2.15	0.46
1:A:319:LEU:HD12	1:A:320:LEU:N	2.31	0.46
1:B:212:ARG:HB2	1:B:213:PRO:CD	2.37	0.46
1:B:13:ILE:O	1:B:13:ILE:HD13	2.16	0.46
1:A:172:HIS:HE1	1:A:224:GLY:O	1.99	0.45
1:B:11:GLN:HB2	1:B:18:GLN:NE2	2.31	0.45
1:B:169:ARG:NH1	1:B:288:VAL:O	2.49	0.45
1:B:174:PRO:O	2:B:7517:UDP:H5'1	2.17	0.45
1:A:307:ASN:HD22	1:A:308:PRO:CD	2.30	0.44
1:B:180:PHE:HB3	1:B:182:PHE:HE1	1.81	0.44
1:B:156:HIS:CE1	1:B:214:ASP:OD2	2.60	0.44
1:B:181:SER:HA	1:B:184:LYS:HE2	1.99	0.43
1:B:213:PRO:CG	1:B:216:GLN:HE21	2.30	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:69:ILE:HD13	1:B:320:LEU:HD12	2.00	0.43
1:B:242:TYR:C	1:B:242:TYR:CD1	2.91	0.42
1:B:86:ASN:ND2	1:B:123:LYS:NZ	2.66	0.42
1:B:213:PRO:HD2	1:B:216:GLN:HE21	1.83	0.42
1:A:306:PHE:O	1:A:309:ILE:HG22	2.20	0.42
1:B:202:LEU:HB3	1:B:203:PRO:CD	2.49	0.42
1:B:316:THR:O	1:B:320:LEU:HB2	2.20	0.42
1:A:242:TYR:CD1	1:A:242:TYR:C	2.92	0.42
1:B:190:ILE:HA	1:B:191:PRO:HD3	1.89	0.41
1:B:184:LYS:H	1:B:184:LYS:HD3	1.86	0.41
1:A:100:HIS:CE1	1:A:154:TRP:HH2	2.39	0.41
1:B:281:VAL:O	1:B:285:ILE:HG12	2.21	0.40
1:A:103:VAL:HB	1:A:104:PRO:HD3	2.04	0.40
1:A:202:LEU:HB3	1:A:203:PRO:HD2	2.02	0.40
1:A:70:PHE:C	1:A:70:PHE:CD1	2.94	0.40
1:B:70:PHE:CD1	1:B:70:PHE:C	2.95	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	328/339 (97%)	325 (99%)	2 (1%)	1 (0%)	44	34
1	B	320/339 (94%)	314 (98%)	5 (2%)	1 (0%)	44	34
All	All	648/678 (96%)	639 (99%)	7 (1%)	2 (0%)	44	34

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	203	PRO
1	A	77	THR

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	297/305 (97%)	276 (93%)	21 (7%)	17	7
1	B	295/305 (97%)	277 (94%)	18 (6%)	22	11
All	All	592/610 (97%)	553 (93%)	39 (7%)	19	9

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

Mol	Chain	Res	Type
1	A	2	ARG
1	A	13	ILE
1	A	19	LEU
1	A	70	PHE
1	A	97	LEU
1	A	98	PHE
1	A	108	SER
1	A	120	TYR
1	A	137	LYS
1	A	184	LYS
1	A	205	ASN
1	A	212	ARG
1	A	218	LEU
1	A	231	MET
1	A	241	LEU
1	A	242	TYR
1	A	266	GLN
1	A	293	GLU
1	A	307	ASN
1	A	310	LEU
1	A	320	LEU
1	B	2	ARG
1	B	13	ILE
1	B	19	LEU
1	B	70	PHE
1	B	97	LEU
1	B	98	PHE

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Mol	Chain	Res	Type
1	B	120	TYR
1	B	184	LYS
1	B	231	MET
1	B	239	GLN
1	B	241	LEU
1	B	242	TYR
1	B	246	LYS
1	B	268	LEU
1	B	293	GLU
1	B	307	ASN
1	B	310	LEU
1	B	320	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (31) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	9	ASN
1	A	18	GLN
1	A	22	ASN
1	A	44	HIS
1	A	86	ASN
1	A	110	ASN
1	A	132	GLN
1	A	151	GLN
1	A	156	HIS
1	A	172	HIS
1	A	210	ASN
1	A	216	GLN
1	A	265	ASN
1	A	266	GLN
1	A	307	ASN
1	A	326	GLN
1	B	9	ASN
1	B	18	GLN
1	B	86	ASN
1	B	110	ASN
1	B	151	GLN
1	B	156	HIS
1	B	172	HIS
1	B	207	HIS
1	B	210	ASN
1	B	216	GLN

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Mol	Chain	Res	Type
1	B	239	GLN
1	B	265	ASN
1	B	266	GLN
1	B	307	ASN
1	B	326	GLN

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 3 ligands modelled in this entry, 1 is monoatomic - leaving 2 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	$\# Z > 2$
2	UDP	A	7517	-	21,26,26	0.99	1 (4%)	22,40,40	1.75	2 (9%)
2	UDP	B	7517	-	21,26,26	1.10	1 (4%)	22,40,40	1.64	2 (9%)

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
2	UDP	A	7517	-	-	0/12/32/32	0/2/2/2
2	UDP	B	7517	-	-	0/12/32/32	0/2/2/2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	7517	UDP	C4-N3	2.72	1.38	1.33
2	B	7517	UDP	C4-N3	2.94	1.38	1.33

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	7517	UDP	O4'-C1'-N1	2.34	112.76	108.08
2	A	7517	UDP	O4'-C1'-N1	2.98	114.05	108.08
2	B	7517	UDP	C4-N3-C2	6.79	119.96	114.13
2	A	7517	UDP	C4-N3-C2	6.94	120.10	114.13

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	7517	UDP	2	0

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	330/339 (97%)	-0.02	7 (2%) 64 67	12, 24, 47, 80	0
1	B	326/339 (96%)	0.69	50 (15%) 2 2	11, 33, 97, 150	0
All	All	656/678 (96%)	0.33	57 (8%) 11 12	11, 27, 77, 150	0

All (57) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	206	VAL	11.2
1	B	197	TRP	9.8
1	B	199	ASN	9.5
1	B	188	TYR	7.4
1	B	202	LEU	7.1
1	B	200	VAL	6.9
1	B	190	ILE	5.8
1	B	186	TRP	5.7
1	B	211	TYR	5.5
1	A	329	CYS	5.3
1	B	178	GLU	5.2
1	B	189	ASP	4.8
1	B	181	SER	4.6
1	B	210	ASN	4.5
1	B	192	LEU	4.5
1	B	198	GLN	4.5
1	B	286	MET	4.4
1	B	285	ILE	4.0
1	B	183	VAL	4.0
1	B	0	CYS	4.0
1	A	107	PHE	3.9
1	B	196	THR	3.6
1	B	233	ASP	3.5
1	B	201	GLU	3.5

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Mol	Chain	Res	Type	RSRZ
1	B	184	LYS	3.5
1	B	204	GLN	3.5
1	B	176	ASN	3.4
1	B	180	PHE	3.3
1	B	289	LYS	3.3
1	B	213	PRO	3.2
1	A	13	ILE	3.2
1	B	163	PHE	3.2
1	B	205	ASN	3.2
1	B	207	HIS	3.0
1	B	141	PHE	3.0
1	B	234	LYS	2.9
1	B	111	PHE	2.8
1	B	209	ILE	2.8
1	B	237	GLU	2.6
1	B	13	ILE	2.6
1	B	170	GLU	2.6
1	B	203	PRO	2.5
1	A	0	CYS	2.5
1	B	291	VAL	2.4
1	A	199	ASN	2.4
1	B	140	ASP	2.3
1	B	212	ARG	2.3
1	B	236	LYS	2.3
1	B	174	PRO	2.3
1	B	179	ARG	2.1
1	B	112	TYR	2.1
1	B	12	SER	2.1
1	A	44	HIS	2.1
1	A	146	SER	2.1
1	B	185	GLU	2.1
1	B	167	LEU	2.0
1	B	215	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 [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	UDP	B	7517	25/25	0.91	0.14	-0.85	38,58,81,83	0
2	UDP	A	7517	25/25	0.98	0.09	-1.03	15,20,26,27	0
3	CL	B	330	1/1	0.82	0.11	-	78,78,78,78	0

6.5 Other polymers [i](#)

There are no such residues in this entry.