



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 15, 2017 – 12:46 am GMT

PDB ID : 4GYW
Title : Crystal structure of human O-GlcNAc Transferase in complex with UDP and a glycopeptide
Authors : Lazarus, M.B.; Jiang, J.; Gloster, T.M.; Zandberg, W.F.; Vocadlo, D.J.; Walker, S.
Deposited on : 2012-09-05
Resolution : 1.70 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 ⓘ 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

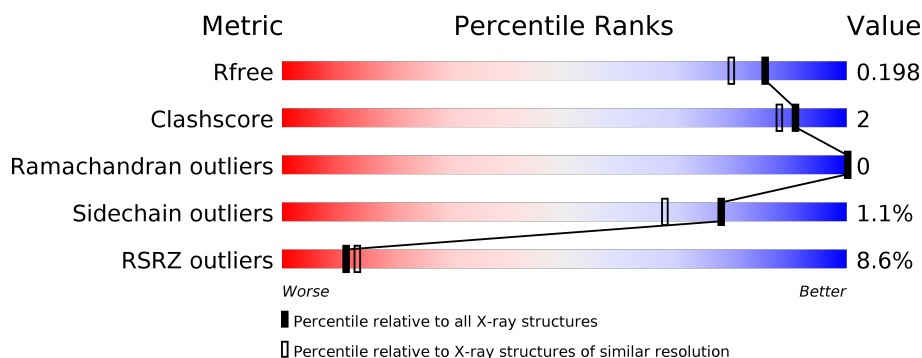
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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}	100719	3453 (1.70-1.70)
Clashscore	112137	3876 (1.70-1.70)
Ramachandran outliers	110173	3815 (1.70-1.70)
Sidechain outliers	110143	3815 (1.70-1.70)
RSRZ outliers	101464	3491 (1.70-1.70)

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 $\leq 5\%$. 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	723	<div> <div>8%</div> <div>91%</div> <div>5%</div> </div>
1	C	723	<div> <div>8%</div> <div>87%</div> <div>6%</div> <div>7%</div> </div>
2	B	14	<div> <div>14%</div> <div>79%</div> <div>21%</div> </div>
2	D	14	<div> <div>14%</div> <div>71%</div> <div>29%</div> </div>

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 12156 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 UDP-N-acetylglucosamine--peptide N-acetylglucosaminyltransferase 110 kDa subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	695	Total	C	N	O	S	0	8	0
			5536	3514	966	1019	37			
1	C	674	Total	C	N	O	S	0	7	0
			5367	3417	934	978	38			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	309	GLY	-	EXPRESSION TAG	UNP O15294
A	310	PRO	-	EXPRESSION TAG	UNP O15294
A	311	GLY	-	EXPRESSION TAG	UNP O15294
A	312	SER	-	EXPRESSION TAG	UNP O15294
C	309	GLY	-	EXPRESSION TAG	UNP O15294
C	310	PRO	-	EXPRESSION TAG	UNP O15294
C	311	GLY	-	EXPRESSION TAG	UNP O15294
C	312	SER	-	EXPRESSION TAG	UNP O15294

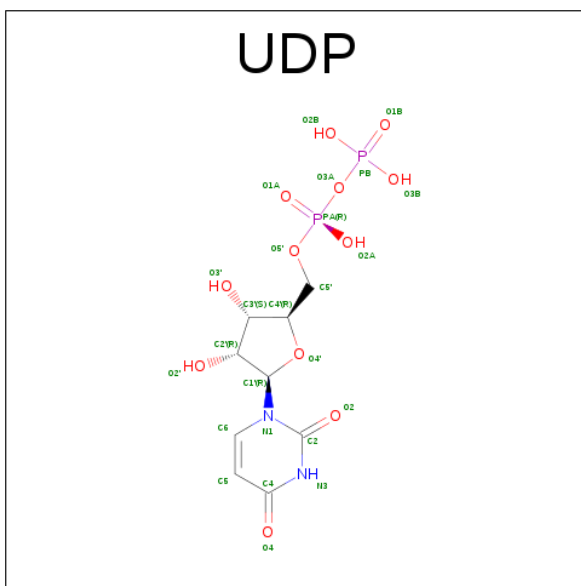
- Molecule 2 is a protein called Casein kinase II subunit alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	14	Total	C	N	O	S	0	0	0
			95	58	15	20	2			
2	D	14	Total	C	N	O	S	0	0	0
			95	58	15	20	2			

There are 2 discrepancies between the modelled and reference sequences:

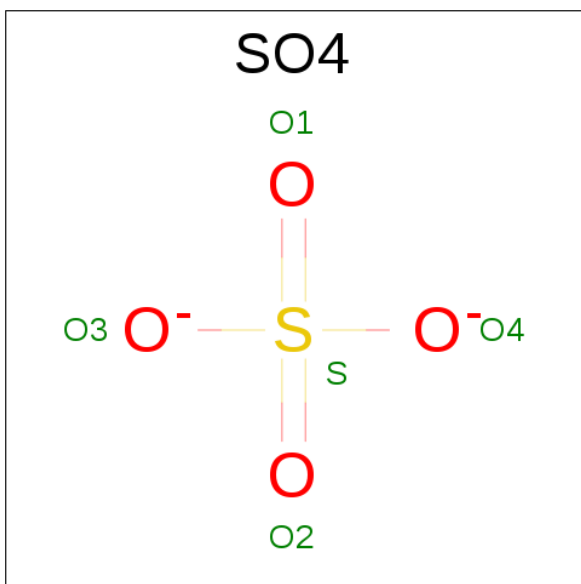
Chain	Residue	Modelled	Actual	Comment	Reference
B	13	TYR	-	EXPRESSION TAG	UNP P68400
D	13	TYR	-	EXPRESSION TAG	UNP P68400

- Molecule 3 is URIDINE-5'-DIPHOSPHATE (three-letter code: UDP) (formula: $C_9H_{14}N_2O_{12}P_2$).



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

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



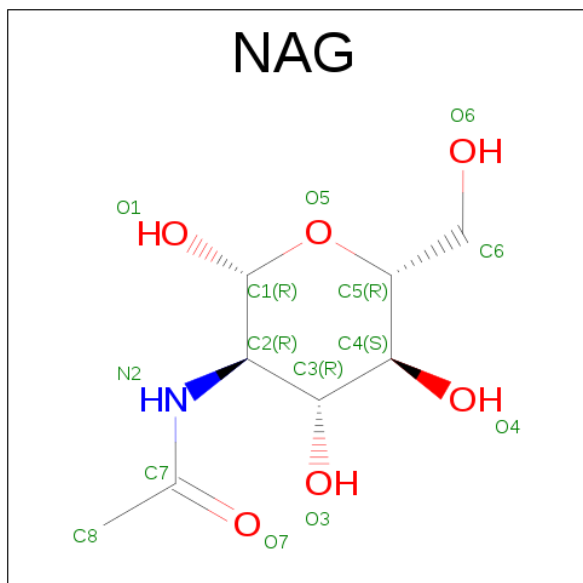
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 5 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	B	1	Total	C	N	O	0	0
			14	8	1	5		
5	D	1	Total	C	N	O	0	0
			14	8	1	5		

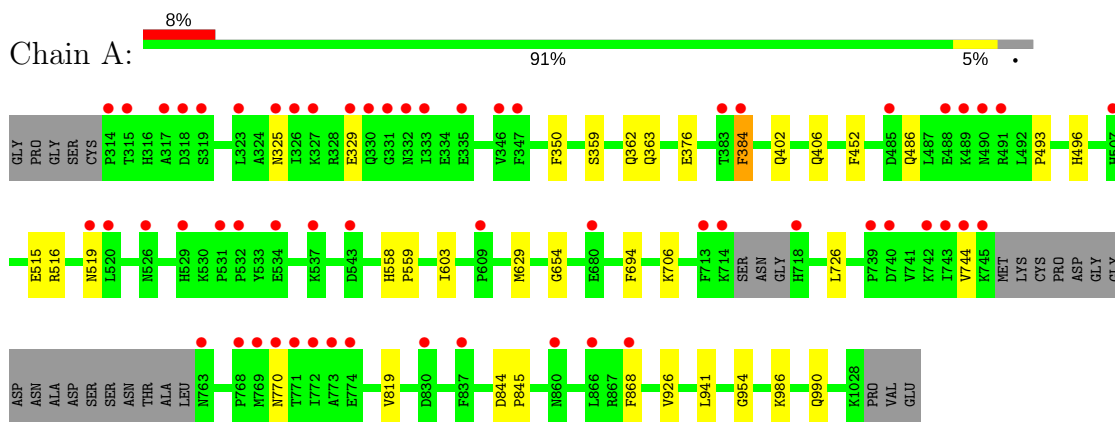
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	487	Total	O	0	0
			487	487		
6	B	19	Total	O	0	0
			19	19		
6	C	450	Total	O	0	0
			450	450		
6	D	14	Total	O	0	0
			14	14		

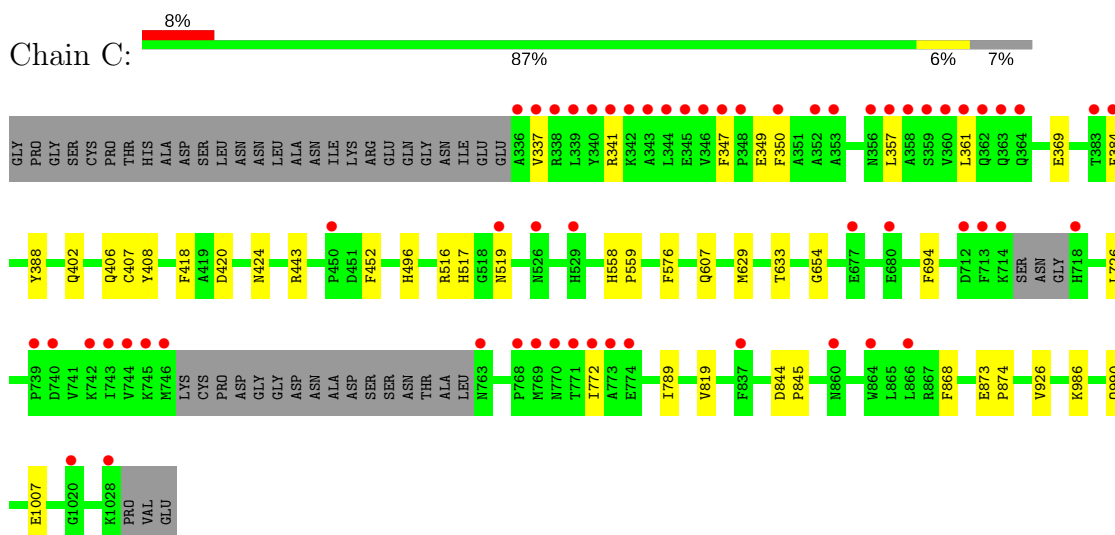
3 Residue-property plots [i](#)

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: UDP-N-acetylglucosamine--peptide N-acetylglucosaminyltransferase 110 kDa subunit



- Molecule 1: UDP-N-acetylglucosamine--peptide N-acetylglucosaminyltransferase 110 kDa subunit

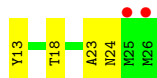


- Molecule 2: Casein kinase II subunit alpha





- Molecule 2: Casein kinase II subunit alpha



4 Data and refinement statistics

Property	Value	Source
Space group	I 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	99.00Å 137.77Å 153.57Å 90.00° 102.90° 90.00°	Depositor
Resolution (Å)	43.90 – 1.70 43.90 – 1.70	Depositor EDS
% Data completeness (in resolution range)	94.2 (43.90-1.70) 94.2 (43.90-1.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.24 (at 1.70Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7_650)	Depositor
R, R_{free}	0.187 , 0.205 0.179 , 0.198	Depositor DCC
R_{free} test set	10313 reflections (4.99%)	DCC
Wilson B-factor (Å ²)	18.9	Xtriage
Anisotropy	0.149	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 52.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	12156	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 3.53% 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, NAG, SO4

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	0/5688	0.48	0/7713
1	C	0.30	0/5515	0.48	0/7478
2	B	0.32	0/97	0.45	0/131
2	D	0.29	0/97	0.42	0/131
All	All	0.30	0/11397	0.48	0/15453

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 [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	5536	0	5520	20	0
1	C	5367	0	5363	22	0
2	B	95	0	87	2	0
2	D	95	0	87	3	0
3	A	25	0	11	0	0
3	C	25	0	11	0	0
4	A	5	0	0	0	0
4	B	5	0	0	0	0
4	D	5	0	0	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	B	14	0	13	0	0
5	D	14	0	13	0	0
6	A	487	0	0	5	0
6	B	19	0	0	0	0
6	C	450	0	0	3	0
6	D	14	0	0	0	0
All	All	12156	0	11105	44	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 2.

All (44) 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)
2:D:13:TYR:N	2:D:18:THR:HG1	1.92	0.68
1:A:726:LEU:CD2	1:A:819:VAL:HG22	2.26	0.66
2:B:13:TYR:N	2:B:18:THR:HG1	1.97	0.62
1:A:706:LYS:HE3	6:A:1654:HOH:O	2.04	0.58
1:A:986:LYS:HG3	1:A:990[B]:GLN:OE1	2.04	0.56
1:C:443:ARG:NH2	6:C:1621:HOH:O	2.36	0.56
1:C:726[A]:LEU:CD2	1:C:819:VAL:HG22	2.37	0.55
1:C:607:GLN:NE2	6:C:1582:HOH:O	2.39	0.54
1:C:986:LYS:HG3	1:C:990[B]:GLN:OE1	2.09	0.53
1:A:558:HIS:CG	1:A:559:PRO:HD2	2.46	0.50
1:A:376:GLU:HG2	6:A:1424:HOH:O	2.12	0.50
1:A:402:GLN:O	1:A:406:GLN:HG2	2.12	0.49
1:C:772:ILE:HG23	1:C:789:ILE:HD13	1.94	0.48
1:A:325:ASN:O	1:A:329:GLU:HG2	2.14	0.47
1:C:337:VAL:O	1:C:341:ARG:HG3	2.14	0.47
1:A:770:ASN:ND2	6:A:1682:HOH:O	2.47	0.47
1:A:496:HIS:CE1	2:B:23:ALA:HB1	2.51	0.46
1:A:954:GLY:O	1:A:986:LYS:HE2	2.16	0.46
1:C:516:ARG:NH2	1:C:519:ASN:OD1	2.48	0.46
1:A:362:GLN:HG2	6:A:1639:HOH:O	2.15	0.46
1:C:402:GLN:O	1:C:406:GLN:HG2	2.17	0.45
1:A:359:SER:O	1:A:363:GLN:HG3	2.17	0.44
1:C:517:HIS:NE2	6:C:1569:HOH:O	2.31	0.44
1:C:347:PHE:CE1	1:C:349:GLU:HB2	2.53	0.44
1:A:516:ARG:NH2	6:A:1680:HOH:O	2.51	0.43
1:C:496:HIS:CE1	2:D:23:ALA:HB1	2.53	0.43
1:C:558:HIS:CG	1:C:559:PRO:HD2	2.55	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:408:TYR:CZ	1:C:424:ASN:HB3	2.55	0.42
1:C:629:MET:O	1:C:654:GLY:HA3	2.20	0.42
1:A:486:GLN:OE1	1:A:493:PRO:HA	2.19	0.42
1:A:515[A]:GLU:OE2	1:A:519:ASN:ND2	2.51	0.42
1:C:576:PHE:CZ	1:C:1007:GLU:HB3	2.55	0.41
1:C:361:LEU:HD13	1:C:369:GLU:HB3	2.02	0.41
1:C:844:ASP:HB2	1:C:845:PRO:HD2	2.01	0.41
1:A:844:ASP:HB2	1:A:845:PRO:HD2	2.01	0.41
1:A:603:ILE:N	1:A:603:ILE:HD12	2.35	0.41
1:C:418:PHE:CE1	1:C:420:ASP:HB2	2.55	0.41
1:C:873:GLU:N	1:C:874:PRO:HD2	2.36	0.41
1:A:941:LEU:HD23	1:A:941:LEU:C	2.41	0.41
1:A:629:MET:O	1:A:654:GLY:HA3	2.21	0.40
1:C:388:TYR:O	1:C:407:CYS:HB3	2.21	0.40
1:C:633:THR:O	2:D:24:ASN:HB3	2.21	0.40
1:C:357:LEU:O	1:C:361:LEU:HG	2.21	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	698/723 (96%)	686 (98%)	12 (2%)	0	100	100
1	C	675/723 (93%)	663 (98%)	12 (2%)	0	100	100
2	B	12/14 (86%)	12 (100%)	0	0	100	100
2	D	12/14 (86%)	12 (100%)	0	0	100	100
All	All	1397/1474 (95%)	1373 (98%)	24 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	606/618 (98%)	599 (99%)	7 (1%)	75	64
1	C	586/618 (95%)	580 (99%)	6 (1%)	80	71
2	B	11/11 (100%)	11 (100%)	0	100	100
2	D	11/11 (100%)	11 (100%)	0	100	100
All	All	1214/1258 (96%)	1201 (99%)	13 (1%)	78	68

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	350	PHE
1	A	384	PHE
1	A	452	PHE
1	A	694	PHE
1	A	744	VAL
1	A	868	PHE
1	A	926	VAL
1	C	350	PHE
1	C	384	PHE
1	C	452	PHE
1	C	694	PHE
1	C	868	PHE
1	C	926	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	607	GLN
1	A	763	ASN
1	C	607	GLN
1	C	784	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 ⓘ

7 ligands are modelled in this entry.

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
3	UDP	A	1101	-	21,26,26	1.13	1 (4%)	22,40,40	1.86	1 (4%)
4	SO4	A	1102	-	4,4,4	0.15	0	6,6,6	0.07	0
5	NAG	B	101	2	14,14,15	0.51	0	15,19,21	0.86	1 (6%)
4	SO4	B	102	-	4,4,4	0.15	0	6,6,6	0.08	0
3	UDP	C	1101	-	21,26,26	1.06	1 (4%)	22,40,40	1.72	1 (4%)
5	NAG	D	101	2	14,14,15	0.41	0	15,19,21	0.82	0
4	SO4	D	102	-	4,4,4	0.18	0	6,6,6	0.11	0

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
3	UDP	A	1101	-	-	0/12/32/32	0/2/2/2
4	SO4	A	1102	-	-	0/0/0/0	0/0/0/0

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	B	101	2	-	0/6/23/26	0/1/1/1
4	SO4	B	102	-	-	0/0/0/0	0/0/0/0
3	UDP	C	1101	-	-	0/12/32/32	0/2/2/2
5	NAG	D	101	2	-	0/6/23/26	0/1/1/1
4	SO4	D	102	-	-	0/0/0/0	0/0/0/0

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	1101	UDP	C4-N3	2.82	1.38	1.33
3	A	1101	UDP	C4-N3	3.04	1.38	1.33

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	101	NAG	C1-O5-C5	2.10	115.06	112.17
3	C	1101	UDP	C4-N3-C2	7.53	120.60	114.13
3	A	1101	UDP	C4-N3-C2	7.97	120.97	114.13

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	695/723 (96%)	0.47	58 (8%) 12 14	9, 20, 44, 91	0
1	C	674/723 (93%)	0.60	58 (8%) 11 13	10, 19, 47, 119	0
2	B	14/14 (100%)	0.47	2 (14%) 3 3	14, 19, 57, 64	0
2	D	14/14 (100%)	0.46	2 (14%) 3 3	13, 20, 53, 59	0
All	All	1397/1474 (94%)	0.53	120 (8%) 11 13	9, 19, 46, 119	0

All (120) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	337	VAL	15.8
1	C	339	LEU	15.6
1	C	346	VAL	12.2
1	C	344	LEU	11.7
1	C	347	PHE	11.5
1	C	348	PRO	9.5
1	C	338	ARG	9.5
1	C	340	TYR	9.3
1	C	342	LYS	9.0
1	C	336	ALA	8.8
1	A	718	HIS	8.8
1	C	350	PHE	8.2
1	A	772	ILE	7.7
1	C	718	HIS	7.3
1	A	714	LYS	6.9
1	A	769	MET	6.5
1	A	314	PRO	6.5
1	C	360	VAL	6.4
1	A	713	PHE	5.8
1	C	1028	LYS	5.7
1	C	713	PHE	5.6

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	315	THR	5.1
1	C	343	ALA	5.1
1	C	746	MET	5.1
1	C	769	MET	4.7
1	A	347	PHE	4.6
1	A	491	ARG	4.5
1	C	714	LYS	4.5
1	C	745	LYS	4.4
1	A	740	ASP	4.4
1	A	745	LYS	4.4
1	C	345	GLU	4.2
1	C	356	ASN	3.9
1	C	740	ASP	3.7
1	C	361	LEU	3.7
1	C	771	THR	3.7
1	A	323	LEU	3.5
1	C	743	ILE	3.5
1	A	331	GLY	3.5
1	A	739	PRO	3.5
1	A	768	PRO	3.4
1	C	341	ARG	3.4
1	C	772	ILE	3.4
1	C	352	ALA	3.3
1	C	742	LYS	3.3
1	A	490	ASN	3.2
1	C	770	ASN	3.2
1	A	332	ASN	3.2
1	A	771	THR	3.2
1	A	326	ILE	3.2
1	A	488	GLU	3.2
1	A	860	ASN	3.2
2	D	26	MET	3.0
1	A	534	GLU	3.0
1	C	529	HIS	3.0
1	A	680	GLU	3.0
1	A	335	GLU	3.0
1	C	860	ASN	2.9
1	C	739	PRO	2.9
1	C	763	ASN	2.9
1	A	384	PHE	2.9
1	A	770	ASN	2.9
1	A	742	LYS	2.8

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	532	PRO	2.8
1	C	357	LEU	2.8
1	C	712	ASP	2.8
1	A	383	THR	2.7
1	C	363	GLN	2.7
2	B	26	MET	2.7
1	C	383	THR	2.6
1	A	507	HIS	2.6
1	A	537	LYS	2.6
1	A	333	ILE	2.6
1	A	520	LEU	2.6
1	C	866	LEU	2.6
1	A	743	ILE	2.6
1	A	866	LEU	2.6
1	A	329	GLU	2.5
1	C	359	SER	2.5
2	B	25	MET	2.5
1	C	768	PRO	2.5
1	A	744	VAL	2.5
1	A	519	ASN	2.5
1	A	346	VAL	2.5
1	A	526	ASN	2.4
1	A	489	LYS	2.4
1	C	1020	GLY	2.4
1	C	837	PHE	2.4
1	A	830	ASP	2.4
1	A	774	GLU	2.4
1	A	609	PRO	2.3
2	D	25	MET	2.3
1	C	680	GLU	2.3
1	C	353	ALA	2.3
1	C	450	PRO	2.3
1	A	763	ASN	2.3
1	C	526	ASN	2.3
1	A	330	GLN	2.3
1	A	837	PHE	2.3
1	A	319[A]	SER	2.3
1	A	773	ALA	2.2
1	A	543	ASP	2.2
1	C	744	VAL	2.2
1	A	325	ASN	2.2
1	C	519	ASN	2.2

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	485	ASP	2.1
1	C	774	GLU	2.1
1	A	327	LYS	2.1
1	A	868	PHE	2.1
1	A	529	HIS	2.1
1	C	677	GLU	2.1
1	C	362	GLN	2.1
1	C	364	GLN	2.1
1	C	384	PHE	2.1
1	A	318	ASP	2.1
1	A	531	PRO	2.0
1	A	317	ALA	2.0
1	C	358	ALA	2.0
1	C	773	ALA	2.0
1	C	864	TRP	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
4	SO4	D	102	5/5	0.95	0.14	0.85	43,45,48,50	0
4	SO4	B	102	5/5	0.97	0.14	0.70	43,43,47,48	0
5	NAG	B	101	14/15	0.94	0.12	-0.11	12,14,18,18	0
5	NAG	D	101	14/15	0.95	0.12	-0.18	11,14,17,19	0
3	UDP	A	1101	25/25	0.98	0.12	-0.99	8,10,13,15	0
3	UDP	C	1101	25/25	0.98	0.11	-1.09	9,11,13,15	0
4	SO4	A	1102	5/5	0.91	0.16	-	57,57,59,60	0

6.5 Other polymers [i](#)

There are no such residues in this entry.