



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 14, 2017 – 09:43 PM EDT

PDB ID : 5HGV
Title : Structure of an O-GlcNAc transferase point mutant, D554N in complex with peptide
Authors : Janetzko, J.; Lazarus, M.B.; Walker, S.
Deposited on : unknown
Resolution : 2.05 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 : rb-20029824
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) : rb-20029824

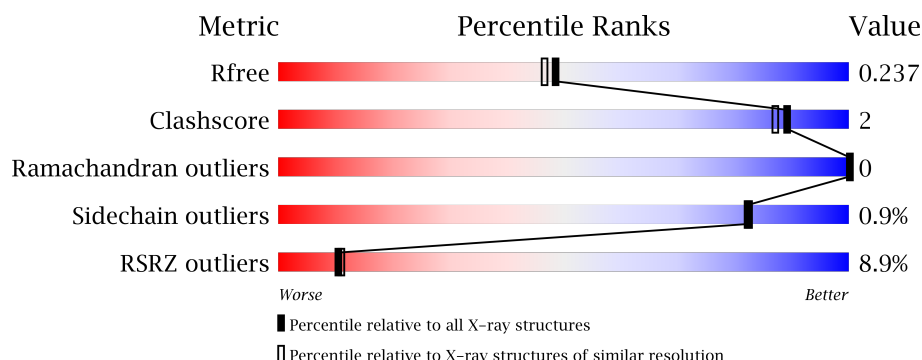
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.05 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	1316 (2.04-2.04)
Clashscore	112137	1394 (2.04-2.04)
Ramachandran outliers	110173	1383 (2.04-2.04)
Sidechain outliers	110143	1383 (2.04-2.04)
RSRZ outliers	101464	1319 (2.04-2.04)

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	719	<div> <div>8%</div> <div>90%</div> <div>7%</div> <div>.</div> </div>
1	C	719	<div> <div>9%</div> <div>88%</div> <div>6%</div> <div>6%</div> </div>
2	B	14	<div> <div>79%</div> <div>21%</div> </div>
2	D	14	<div> <div>86%</div> <div>14%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	SO4	A	1102	-	-	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 11958 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	967	1018	37			
1	C	674	Total	C	N	O	S	0	7	0
			5367	3417	935	977	38			

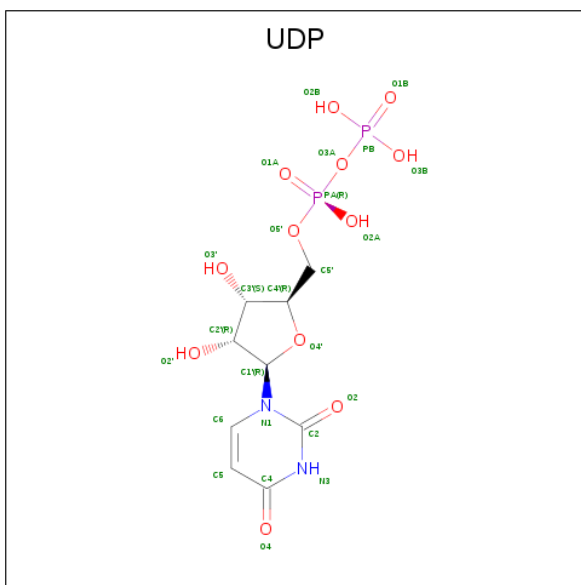
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	554	ASN	ASP	engineered mutation	UNP O15294
C	554	ASN	ASP	engineered mutation	UNP O15294

- Molecule 2 is a protein called TYR-PRO-GLY-GLY-SER-THR-PRO-VAL-SER-SER-ALA-ASN-MET-MET.

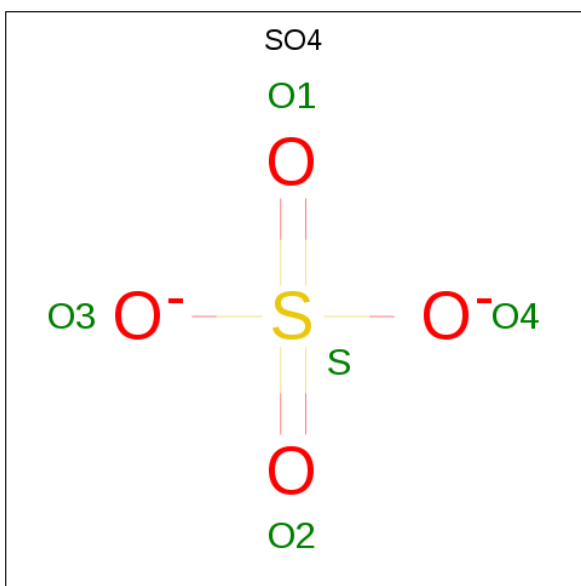
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	14	Total	C	N	O	S	0	0	0
			95	58	15	20	2			
2	D	14	Total	C	N	O	S	0	0	0
			95	58	15	20	2			

- Molecule 3 is URIDINE-5'-DIPHOSPHATE (three-letter code: UDP) (formula: C₉H₁₄N₂O₁₂P₂).



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

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



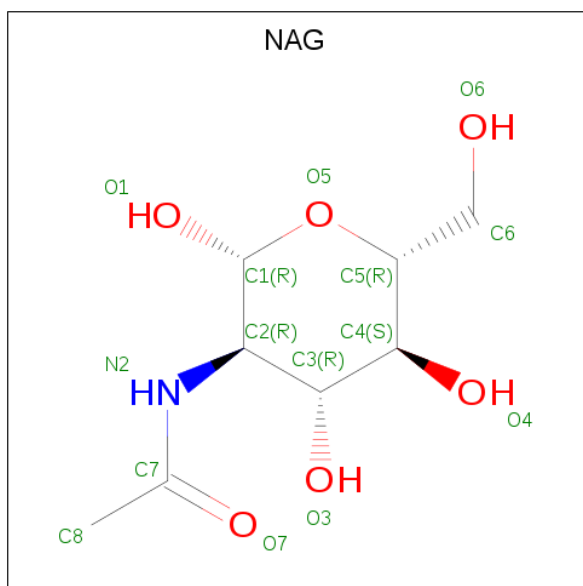
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total 5	O 4	S 1	0	0
4	B	1	Total 5	O 4	S 1	0	0

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

- Molecule 5 is 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		

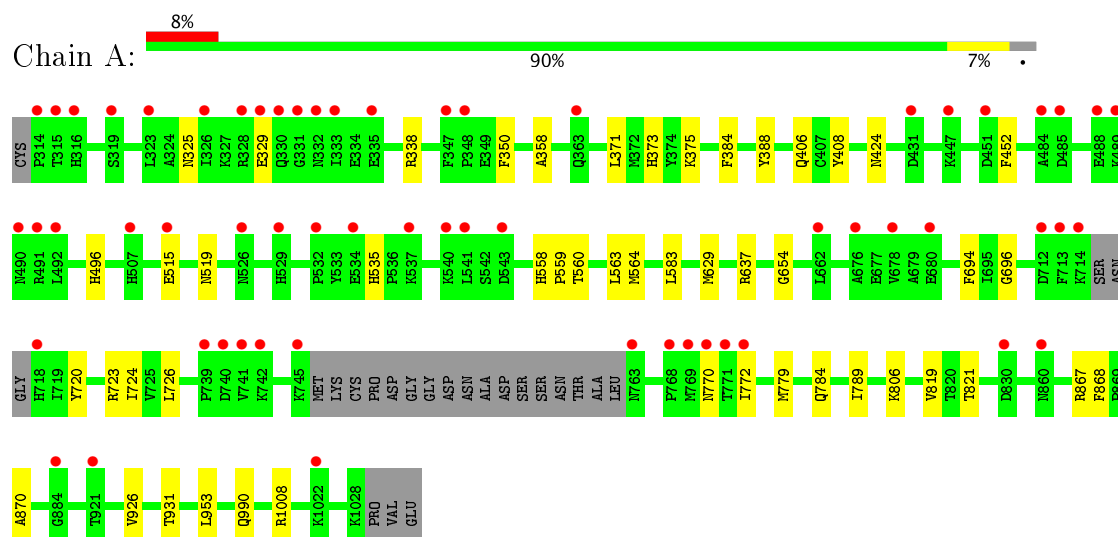
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	370	Total	O	0	0
			370	370		
6	B	14	Total	O	0	0
			14	14		
6	C	376	Total	O	0	0
			376	376		
6	D	12	Total	O	0	0
			12	12		

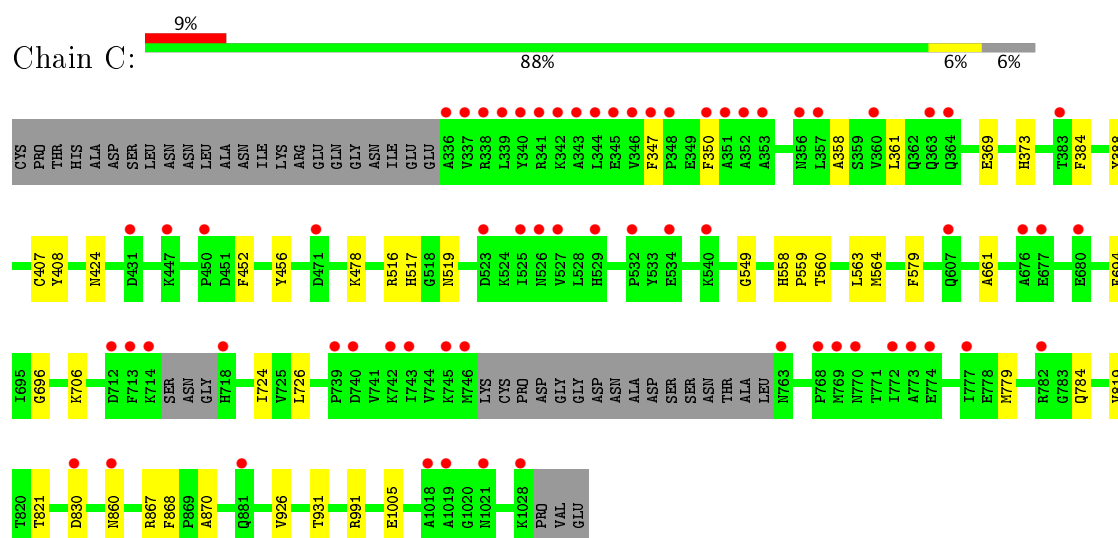
3 Residue-property plots [i](#)

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




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- Molecule 2: TYR-PRO-GLY-GLY-SER-THR-PRO-VAL-SER-SER-ALA-ASN-MET-ME

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Chain B:  79% 21%



- Molecule 2: TYR-PRO-GLY-GLY-SER-THR-PRO-VAL-SER-SER-ALA-ASN-MET-MET

Chain D:  86% 14%



4 Data and refinement statistics

Property	Value	Source
Space group	I 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	98.35Å 137.19Å 153.07Å 90.00° 103.05° 90.00°	Depositor
Resolution (Å)	40.92 – 2.05 40.92 – 2.05	Depositor EDS
% Data completeness (in resolution range)	95.0 (40.92-2.05) 95.0 (40.92-2.05)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.28 (at 2.05Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.217 , 0.237 0.216 , 0.237	Depositor DCC
R_{free} test set	5894 reflections (5.02%)	DCC
Wilson B-factor (Å ²)	23.5	Xtriage
Anisotropy	0.087	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 55.6	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	0.93	EDS
Total number of atoms	11958	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 19.63% 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	0.28	0/5688	0.44	0/7713
1	C	0.28	0/5515	0.45	0/7478
2	B	0.35	0/97	0.53	0/131
2	D	0.36	0/97	0.54	0/131
All	All	0.28	0/11397	0.45	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	5522	26	0
1	C	5367	0	5365	20	0
2	B	95	0	87	2	0
2	D	95	0	87	1	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

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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	370	0	0	7	0
6	B	14	0	0	0	0
6	C	376	0	0	3	0
6	D	12	0	0	0	0
All	All	11958	0	11109	48	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 (48) 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:C:661:ALA:O	6:C:1202:HOH:O	2.05	0.75
2:D:13:TYR:N	2:D:18:THR:HG1	1.93	0.66
1:C:517:HIS:NE2	6:C:1201:HOH:O	1.97	0.64
1:C:516:ARG:NH2	1:C:519:ASN:OD1	2.31	0.62
2:B:13:TYR:N	2:B:18:THR:HG1	1.96	0.62
1:A:338:ARG:O	6:A:1202:HOH:O	2.16	0.61
1:A:1008:ARG:NH2	6:A:1210:HOH:O	2.36	0.58
1:C:358:ALA:HB2	1:C:373:HIS:HB2	1.87	0.56
1:A:371:LEU:HG	1:A:375:LYS:HE3	1.88	0.55
1:C:361:LEU:HD13	1:C:369:GLU:HB3	1.89	0.55
1:C:1005:GLU:HG3	6:C:1474:HOH:O	2.07	0.55
1:A:770:ASN:ND2	6:A:1211:HOH:O	2.38	0.52
1:A:358:ALA:HB2	1:A:373:HIS:HB2	1.92	0.52
1:A:953:LEU:O	1:A:990[B]:GLN:NE2	2.44	0.50
1:C:867:ARG:HB3	1:C:870:ALA:HA	1.92	0.50
1:A:779:MET:HG3	1:A:784:GLN:HB2	1.94	0.49
1:A:806:LYS:NZ	6:A:1205:HOH:O	2.26	0.49
1:C:926:VAL:HG22	1:C:931:THR:HB	1.95	0.48
1:A:926:VAL:HG22	1:A:931:THR:HB	1.96	0.48
1:A:325:ASN:O	1:A:329:GLU:HG2	2.15	0.46
1:A:408:TYR:CZ	1:A:424:ASN:HB3	2.51	0.46
1:A:867:ARG:HB3	1:A:870:ALA:HA	1.97	0.45
1:C:724:ILE:HG23	1:C:821:THR:HG22	1.99	0.45
1:A:535:HIS:ND1	6:A:1209:HOH:O	2.36	0.44
1:A:558:HIS:CG	1:A:559:PRO:HD2	2.52	0.44
1:C:456:TYR:CZ	1:C:478:LYS:HD3	2.52	0.44
1:C:706:LYS:O	1:C:991:ARG:NH2	2.51	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:830:ASP:O	1:C:860:ASN:ND2	2.50	0.44
1:C:408:TYR:CZ	1:C:424:ASN:HB3	2.53	0.44
1:C:726[A]:LEU:CD2	1:C:819:VAL:HG22	2.48	0.43
1:A:406:GLN:OE1	6:A:1204:HOH:O	2.21	0.43
1:A:563:LEU:HA	1:A:696:GLY:HA2	2.00	0.43
1:C:560:THR:O	1:C:564:MET:HG3	2.19	0.42
1:C:558:HIS:CG	1:C:559:PRO:HD2	2.55	0.42
1:C:779:MET:HG3	1:C:784:GLN:HB2	2.01	0.42
1:A:724:ILE:HG23	1:A:821:THR:HG22	2.01	0.42
1:A:560:THR:O	1:A:564:MET:HG3	2.20	0.42
1:A:496:HIS:CE1	2:B:23:ALA:HB1	2.55	0.42
1:A:720:TYR:HB3	1:A:723:ARG:HD3	2.03	0.41
1:C:549:GLY:HA2	1:C:579:PHE:O	2.20	0.41
1:A:388:TYR:OH	6:A:1201:HOH:O	2.10	0.41
1:A:515[A]:GLU:OE2	1:A:519:ASN:ND2	2.53	0.41
1:A:583:LEU:HD22	1:A:637:ARG:HD3	2.03	0.41
1:C:388:TYR:O	1:C:407:CYS:HB3	2.21	0.40
1:A:629:MET:O	1:A:654:GLY:HA3	2.22	0.40
1:A:726:LEU:CD2	1:A:819:VAL:HG22	2.52	0.40
1:A:772:ILE:HG23	1:A:789:ILE:HD13	2.02	0.40
1:C:563:LEU:HA	1:C:696:GLY:HA2	2.03	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	698/719 (97%)	687 (98%)	11 (2%)	0	100	100
1	C	675/719 (94%)	664 (98%)	11 (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

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	1397/1466 (95%)	1375 (98%)	22 (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	606/616 (98%)	601 (99%)	5 (1%)	85	85
1	C	586/616 (95%)	580 (99%)	6 (1%)	80	79
2	B	11/11 (100%)	11 (100%)	0	100	100
2	D	11/11 (100%)	11 (100%)	0	100	100
All	All	1214/1254 (97%)	1203 (99%)	11 (1%)	82	82

All (11) 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	868	PHE
1	C	347	PHE
1	C	350	PHE
1	C	384	PHE
1	C	452	PHE
1	C	694	PHE
1	C	868	PHE

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	406	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	1.07	1 (4%)	22,40,40	1.48	1 (4%)
4	SO4	A	1102	-	4,4,4	0.19	0	6,6,6	0.16	0
5	NAG	B	101	2	14,14,15	0.83	1 (7%)	15,19,21	0.68	1 (6%)
4	SO4	B	102	-	4,4,4	0.13	0	6,6,6	0.18	0
3	UDP	C	1101	-	21,26,26	1.06	1 (4%)	22,40,40	1.48	1 (4%)
5	NAG	D	101	2	14,14,15	0.99	1 (7%)	15,19,21	0.67	1 (6%)
4	SO4	D	102	-	4,4,4	0.14	0	6,6,6	0.21	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

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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 (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	101	NAG	C1-C2	2.80	1.56	1.52
3	A	1101	UDP	C4-N3	2.91	1.38	1.33
3	C	1101	UDP	C4-N3	3.12	1.38	1.33
5	D	101	NAG	C1-C2	3.25	1.56	1.52

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	D	101	NAG	C1-O5-C5	2.04	114.97	112.17
5	B	101	NAG	C1-O5-C5	2.16	115.14	112.17
3	A	1101	UDP	C4-N3-C2	6.31	119.55	114.13
3	C	1101	UDP	C4-N3-C2	6.34	119.58	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	695/719 (96%)	0.47	60 (8%) 11 12	14, 28, 59, 87	0
1	C	674/719 (93%)	0.56	65 (9%) 9 9	14, 28, 60, 87	0
2	B	14/14 (100%)	0.36	0 100 100	17, 25, 54, 54	0
2	D	14/14 (100%)	0.21	0 100 100	16, 26, 51, 52	0
All	All	1397/1466 (95%)	0.51	125 (8%) 10 11	14, 28, 59, 87	0

All (125) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	769	MET	9.2
1	C	746	MET	8.4
1	C	718	HIS	7.8
1	C	347	PHE	7.7
1	C	769	MET	7.5
1	A	745	LYS	7.4
1	C	338	ARG	6.7
1	A	718	HIS	6.6
1	A	740	ASP	6.2
1	A	714	LYS	6.1
1	C	745	LYS	6.1
1	C	337	VAL	5.6
1	C	339	LEU	5.6
1	C	768	PRO	5.5
1	C	344	LEU	5.5
1	A	326	ILE	5.3
1	C	348	PRO	5.3
1	A	713	PHE	5.2
1	C	713	PHE	5.2
1	A	739	PRO	5.1
1	C	740	ASP	5.1

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Mol	Chain	Res	Type	RSRZ
1	C	346	VAL	5.1
1	C	526	ASN	5.1
1	A	491	ARG	5.0
1	C	739	PRO	4.8
1	C	336	ALA	4.8
1	C	342	LYS	4.7
1	C	350	PHE	4.7
1	C	340	TYR	4.5
1	A	526	ASN	4.5
1	A	772	ILE	4.4
1	C	1028	LYS	4.2
1	A	488	GLU	4.2
1	A	314	PRO	4.2
1	A	768	PRO	4.0
1	C	860	ASN	3.9
1	C	529	HIS	3.9
1	C	772	ILE	3.8
1	A	763	ASN	3.7
1	C	714	LYS	3.7
1	A	860	ASN	3.7
1	A	319[A]	SER	3.6
1	C	680	GLU	3.6
1	C	712	ASP	3.6
1	C	770	ASN	3.5
1	C	742	LYS	3.5
1	A	333	ILE	3.4
1	C	527	VAL	3.4
1	A	680	GLU	3.4
1	C	360	VAL	3.4
1	A	447	LYS	3.3
1	A	529	HIS	3.3
1	A	330	GLN	3.3
1	A	662	LEU	3.3
1	C	363	GLN	3.2
1	C	532	PRO	3.2
1	C	351	ALA	3.2
1	A	335	GLU	3.1
1	C	830	ASP	3.1
1	C	1018	ALA	3.1
1	A	532	PRO	3.1
1	C	345	GLU	3.1
1	C	763	ASN	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	347	PHE	3.1
1	C	352	ALA	3.1
1	A	678	VAL	3.1
1	A	328	ARG	3.0
1	A	490	ASN	3.0
1	C	677	GLU	2.9
1	C	356	ASN	2.9
1	A	315	THR	2.9
1	A	712	ASP	2.9
1	A	830	ASP	2.9
1	A	676	ALA	2.9
1	A	742	LYS	2.9
1	C	343	ALA	2.9
1	C	773	ALA	2.8
1	A	1022	LYS	2.8
1	C	743	ILE	2.8
1	A	770	ASN	2.8
1	A	484	ALA	2.8
1	C	676	ALA	2.8
1	A	540	LYS	2.7
1	A	541	LEU	2.7
1	A	489	LYS	2.7
1	C	383	THR	2.7
1	C	782	ARG	2.6
1	A	537	LYS	2.6
1	C	357	LEU	2.6
1	C	353	ALA	2.5
1	C	774	GLU	2.5
1	C	450	PRO	2.5
1	A	332	ASN	2.4
1	A	431	ASP	2.4
1	A	348	PRO	2.4
1	A	884	GLY	2.4
1	C	1021	ASN	2.4
1	C	431	ASP	2.4
1	A	534	GLU	2.4
1	C	341	ARG	2.4
1	C	364	GLN	2.4
1	C	534	GLU	2.3
1	A	331	GLY	2.3
1	C	881	GLN	2.3
1	A	316	HIS	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	515[A]	GLU	2.3
1	A	492	LEU	2.3
1	C	777	ILE	2.3
1	A	771	THR	2.2
1	A	921	THR	2.2
1	A	543	ASP	2.2
1	A	741	VAL	2.2
1	A	329	GLU	2.2
1	A	507	HIS	2.2
1	A	363	GLN	2.2
1	C	1019	ALA	2.2
1	C	525	ILE	2.2
1	C	523	ASP	2.2
1	C	540	LYS	2.1
1	A	485	ASP	2.1
1	A	323	LEU	2.1
1	C	607	GLN	2.1
1	C	447	LYS	2.1
1	A	451	ASP	2.1
1	C	471	ASP	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	A	1102	5/5	0.81	0.32	8.41	78,78,79,79	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	SO4	D	102	5/5	0.97	0.14	-0.02	39,40,42,42	0
5	NAG	D	101	14/15	0.96	0.15	-0.12	13,15,17,18	0
5	NAG	B	101	14/15	0.94	0.16	-0.19	15,17,18,19	0
4	SO4	B	102	5/5	0.98	0.10	-0.87	40,41,42,43	0
3	UDP	C	1101	25/25	0.98	0.13	-0.96	11,15,17,18	0
3	UDP	A	1101	25/25	0.98	0.13	-1.15	10,14,17,19	0

6.5 Other polymers [i](#)

There are no such residues in this entry.