



Full wwPDB X-ray Structure Validation Report

Feb 26, 2014 – 04:28 PM GMT

PDB ID : 4IRQ
Title : Crystal structure of catalytic domain of human beta1,4galactosyltransferase7
in closed conformation in complex with manganese and UDP
Authors : Tsutsui, Y.; Ramakrishnan, B.; Qasba, P.K.
Deposited on : 2013-01-15
Resolution : 2.30 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

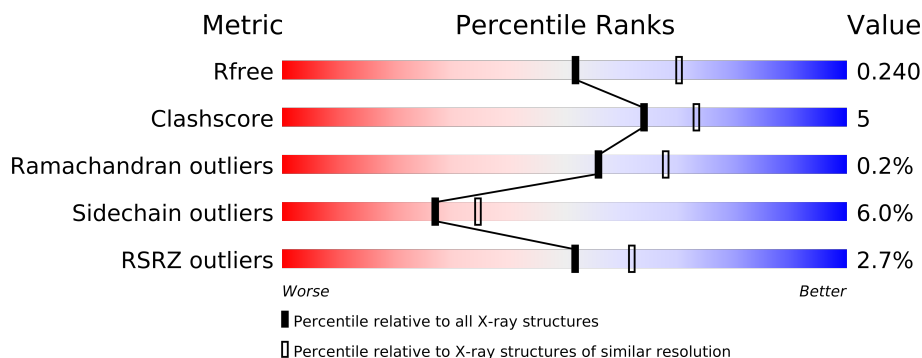
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.30 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}	66092	2929 (2.30-2.30)
Clashscore	79885	3679 (2.30-2.30)
Ramachandran outliers	78287	3642 (2.30-2.30)
Sidechain outliers	78261	3641 (2.30-2.30)
RSRZ outliers	66119	2930 (2.30-2.30)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	251	
1	B	251	
1	C	251	
1	D	251	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
2	MN	B	401	-	X
4	TRS	D	403	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8542 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Beta-1,4-galactosyltransferase7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	246	Total	C	N	O	S	0	1	0
			2037	1304	371	354	8			
1	B	240	Total	C	N	O	S	0	3	0
			1999	1282	366	343	8			
1	C	249	Total	C	N	O	S	0	1	0
			2066	1322	378	358	8			
1	D	243	Total	C	N	O	S	0	2	0
			2015	1289	371	347	8			

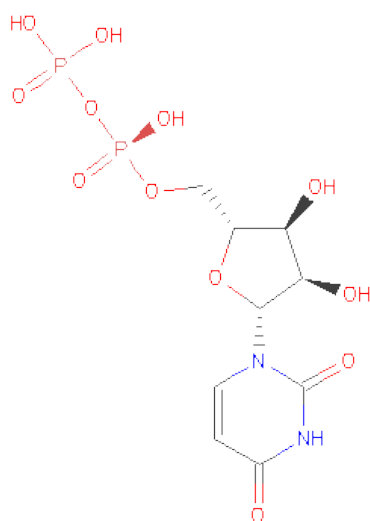
There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	77	GLY	-	EXPRESSION TAG	UNP Q9UBV7
A	78	SER	-	EXPRESSION TAG	UNP Q9UBV7
A	79	ASP	-	EXPRESSION TAG	UNP Q9UBV7
A	80	ILE	-	EXPRESSION TAG	UNP Q9UBV7
B	77	GLY	-	EXPRESSION TAG	UNP Q9UBV7
B	78	SER	-	EXPRESSION TAG	UNP Q9UBV7
B	79	ASP	-	EXPRESSION TAG	UNP Q9UBV7
B	80	ILE	-	EXPRESSION TAG	UNP Q9UBV7
C	77	GLY	-	EXPRESSION TAG	UNP Q9UBV7
C	78	SER	-	EXPRESSION TAG	UNP Q9UBV7
C	79	ASP	-	EXPRESSION TAG	UNP Q9UBV7
C	80	ILE	-	EXPRESSION TAG	UNP Q9UBV7
D	77	GLY	-	EXPRESSION TAG	UNP Q9UBV7
D	78	SER	-	EXPRESSION TAG	UNP Q9UBV7
D	79	ASP	-	EXPRESSION TAG	UNP Q9UBV7
D	80	ILE	-	EXPRESSION TAG	UNP Q9UBV7

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

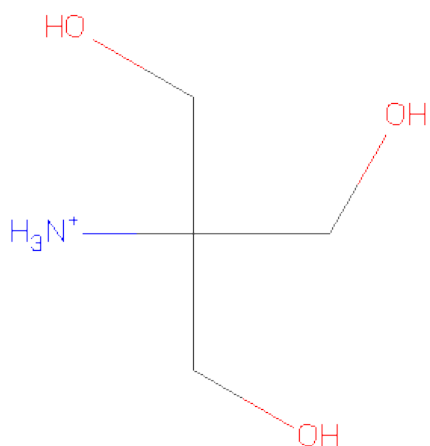
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Mn	0	0
			1	1		
2	A	1	Total	Mn	0	0
			1	1		
2	D	1	Total	Mn	0	0
			1	1		
2	C	1	Total	Mn	0	0
			1	1		

- 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	B	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		
3	D	1	Total	C	N	O	P	0	0
			25	9	2	12	2		

- Molecule 4 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (three-letter code: TRS) (formula: $C_4H_{12}NO_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			8	4	1	3		
4	B	1	Total	C	N	O	0	0
			8	4	1	3		
4	C	1	Total	C	N	O	0	0
			8	4	1	3		
4	D	1	Total	C	N	O	0	0
			8	4	1	3		

- Molecule 5 is water.

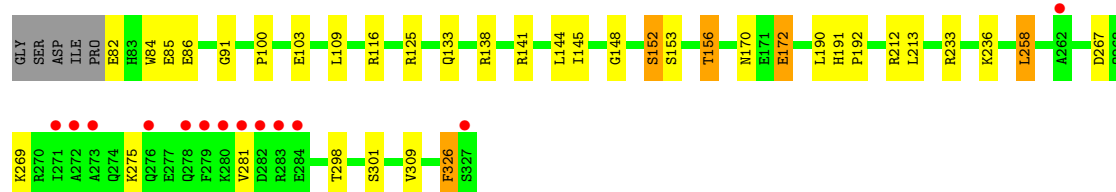
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	85	Total	O	0	0
			85	85		
5	B	66	Total	O	0	0
			66	66		
5	C	64	Total	O	0	0
			64	64		
5	D	74	Total	O	0	0
			74	74		

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.

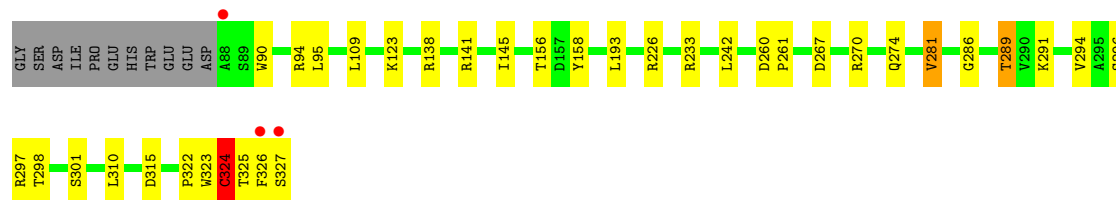
• Molecule 1: Beta-1,4-galactosyltransferase7

Chain A: 



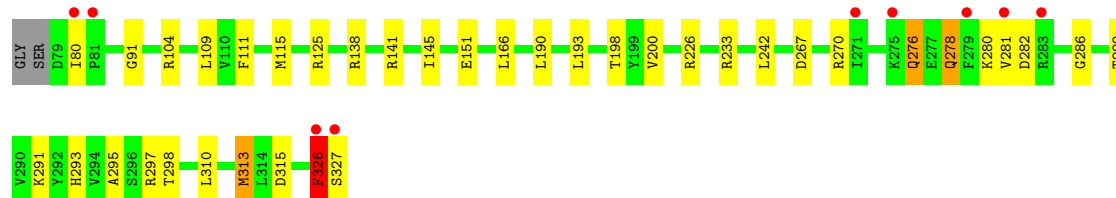
• Molecule 1: Beta-1,4-galactosyltransferase7

Chain B: 



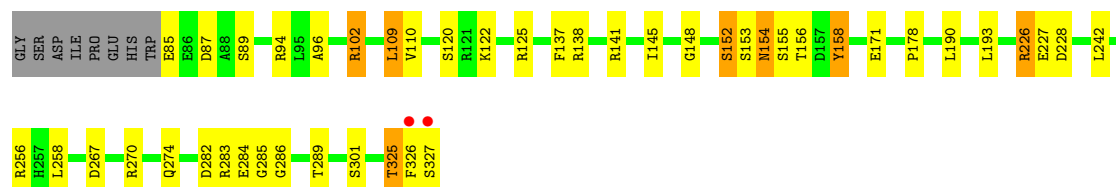
• Molecule 1: Beta-1,4-galactosyltransferase7

Chain C: 



• Molecule 1: Beta-1,4-galactosyltransferase7

Chain D: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	67.12Å 121.53Å 97.14Å 90.00° 95.12° 90.00°	Depositor
Resolution (Å)	37.85 – 2.30 37.85 – 2.30	Depositor EDS
% Data completeness (in resolution range)	96.6 (37.85-2.30) 96.3 (37.85-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.25 (at 2.29Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.1069)	Depositor
R, R_{free}	0.185 , 0.237 0.188 , 0.240	Depositor DCC
R_{free} test set	3406 reflections (5.11%)	DCC
Wilson B-factor (Å ²)	36.1	Xtriage
Anisotropy	0.566	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 35.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Outliers	0 of 66738 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8542	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.29% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: TRS, MN, UDP

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	0.41	0/2102	0.59	0/2844
1	B	0.43	1/2070 (0.0%)	0.58	1/2798 (0.0%)
1	C	0.40	0/2129	0.59	1/2881 (0.0%)
1	D	0.41	0/2080	0.59	0/2811
All	All	0.41	1/8381 (0.0%)	0.59	2/11334 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	324	CYS	CB-SG	-8.78	1.67	1.82

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	326	PHE	CB-CA-C	5.77	121.94	110.40
1	B	324	CYS	CB-CA-C	-5.13	100.14	110.40

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2037	0	1977	22	0
1	B	1999	0	1967	16	0
1	C	2066	0	2007	22	0
1	D	2015	0	1976	28	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	25	0	11	1	0
3	B	25	0	11	0	0
3	C	25	0	11	0	0
3	D	25	0	11	0	0
4	A	8	0	12	1	0
4	B	8	0	12	0	0
4	C	8	0	12	0	0
4	D	8	0	12	3	0
5	A	85	0	0	4	0
5	B	66	0	0	1	0
5	C	64	0	0	2	0
5	D	74	0	0	1	0
All	All	8542	0	8019	88	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 5.

All (88) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:286:GLY:H	1:D:289:THR:HG22	1.45	0.80
1:C:151:GLU:OE1	1:C:297:ARG:NH2	2.19	0.76
1:A:153:SER:O	1:A:156:THR:HG23	1.90	0.72
4:D:403:TRS:O2	5:D:544:HOH:O	2.08	0.71
1:A:258:LEU:HD23	1:A:258:LEU:H	1.56	0.71
1:C:286:GLY:H	1:C:289:THR:HB	1.58	0.68
1:D:227:GLU:OE2	4:D:403:TRS:H21	1.93	0.67
1:D:138:ARG:HD3	1:D:327:SER:HB3	1.75	0.67
1:D:228:ASP:OD2	4:D:403:TRS:H22	1.95	0.66
1:A:298:THR:HG21	1:D:109:LEU:HG	1.78	0.65
1:C:226:ARG:HH21	1:C:233:ARG:HH22	1.45	0.63
1:D:110:VAL:HG11	1:D:256[B]:ARG:NH1	2.15	0.62
1:D:148:GLY:O	1:D:152:SER:OG	2.17	0.61
1:C:226:ARG:HH22	1:C:233:ARG:HH12	1.48	0.60
1:D:94:ARG:HD2	1:D:125:ARG:HB3	1.84	0.59
1:D:96:ALA:HB2	1:D:156:THR:HG21	1.86	0.57

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:138:ARG:NH2	1:D:274:GLN:OE1	2.38	0.56
1:D:256[B]:ARG:NH2	1:D:258:LEU:HD22	2.21	0.56
1:C:138:ARG:HD2	1:C:267:ASP:OD2	2.07	0.55
1:B:286:GLY:H	1:B:289:THR:HB	1.70	0.55
1:C:104[B]:ARG:NH1	5:C:533:HOH:O	2.39	0.55
1:B:138:ARG:HD3	1:B:327:SER:HB3	1.88	0.55
1:A:85:GLU:OE1	1:A:116:ARG:NH1	2.40	0.55
1:B:94:ARG:HG2	1:B:156:THR:HG22	1.88	0.54
1:D:102:ARG:NH2	1:D:137:PHE:O	2.32	0.54
1:D:138:ARG:HD2	1:D:267:ASP:OD2	2.08	0.54
1:A:301:SER:HB3	1:D:301:SER:HB3	1.89	0.54
1:A:138:ARG:HD2	1:A:267:ASP:OD2	2.08	0.53
1:C:91:GLY:O	1:C:125:ARG:HD2	2.09	0.53
1:A:298:THR:HG22	5:A:526:HOH:O	2.09	0.53
1:B:138:ARG:HD2	1:B:267:ASP:OD2	2.10	0.52
1:C:295:ALA:HB2	1:C:313:MET:HE2	1.93	0.51
1:C:291:LYS:HB2	1:C:315:ASP:HB3	1.94	0.50
1:D:228:ASP:N	1:D:228:ASP:OD1	2.43	0.49
1:A:148:GLY:O	1:A:152:SER:OG	2.23	0.49
1:B:141:ARG:NH2	5:B:537:HOH:O	2.31	0.49
1:D:154:ASN:N	1:D:154:ASN:OD1	2.45	0.49
3:A:402:UDP:O3A	4:A:403:TRS:O3	2.31	0.49
1:A:91:GLY:O	1:A:125:ARG:HD2	2.14	0.48
1:B:138:ARG:NH2	1:B:274:GLN:OE1	2.46	0.48
1:A:212:ARG:NH2	5:A:583:HOH:O	2.39	0.47
1:A:141:ARG:O	1:A:145:ILE:HG12	2.14	0.47
1:B:95:LEU:HB2	1:B:158:TYR:CZ	2.49	0.47
1:A:258:LEU:HD23	1:A:258:LEU:N	2.27	0.47
1:D:325:THR:HG22	1:D:326:PHE:H	1.79	0.47
1:C:226:ARG:NH2	1:C:233:ARG:HH12	2.12	0.47
1:D:158:TYR:CD2	1:D:178:PRO:HD3	2.49	0.47
1:C:270:ARG:HA	1:C:270:ARG:HD3	1.68	0.46
1:B:324:CYS:HB3	1:B:325:THR:HG23	1.95	0.46
1:C:138:ARG:NH1	1:C:267:ASP:OD1	2.48	0.46
1:D:85:GLU:HG2	1:D:120:SER:HB3	1.98	0.46
1:B:297[B]:ARG:HG2	1:B:310:LEU:CD2	2.47	0.45
1:C:297:ARG:HG2	1:C:310:LEU:CD2	2.46	0.45
1:B:141:ARG:O	1:B:145:ILE:HG12	2.17	0.45
1:C:141:ARG:O	1:C:145:ILE:HG12	2.17	0.45
1:A:298:THR:HG23	1:A:309:VAL:HB	1.98	0.44
1:A:103:GLU:OE1	5:A:552:HOH:O	2.21	0.44
1:C:276:GLN:HE21	1:C:276:GLN:HB2	1.48	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:233:ARG:CZ	1:B:281:VAL:HG23	2.47	0.44
1:D:122:LYS:NZ	1:D:171:GLU:O	2.51	0.43
1:B:322:PRO:HD2	1:B:323:TRP:CZ3	2.52	0.43
1:B:291:LYS:HB3	1:B:315:ASP:HB3	1.99	0.43
1:C:233:ARG:HE	1:C:281:VAL:HG22	1.83	0.43
1:C:226:ARG:NH2	1:C:233:ARG:HH22	2.15	0.43
1:A:191:HIS:CG	1:A:192:PRO:HD2	2.54	0.43
1:D:110:VAL:HG11	1:D:256[B]:ARG:HH11	1.82	0.43
1:D:284:GLU:HG3	1:D:285:GLY:H	1.84	0.43
1:A:170:ASN:OD1	1:A:172:GLU:HB2	2.19	0.43
1:D:286:GLY:H	1:D:289:THR:CG2	2.22	0.42
1:C:111:PHE:CZ	1:C:115:MET:HG3	2.55	0.42
1:D:153:SER:C	1:D:155:SER:N	2.72	0.42
1:C:280:LYS:HB3	1:C:280:LYS:HE2	1.70	0.42
1:A:84:TRP:CH2	1:A:156:THR:HG22	2.54	0.42
1:D:87:ASP:HB3	1:D:89:SER:H	1.84	0.42
1:D:141:ARG:O	1:D:145:ILE:HG12	2.20	0.42
1:B:90:TRP:CE2	1:B:123:LYS:HG2	2.54	0.42
1:A:269:LYS:HZ3	1:A:326:PHE:HZ	1.67	0.42
1:A:100:PRO:HD3	1:A:145:ILE:HD11	2.03	0.41
1:D:226:ARG:HD3	1:D:226:ARG:HA	1.75	0.41
1:D:256[B]:ARG:NH2	1:D:258:LEU:HB2	2.35	0.41
1:A:133:GLN:HB2	1:A:144:LEU:HD11	2.02	0.41
1:C:138:ARG:CD	1:C:327:SER:HB3	2.50	0.41
1:A:233:ARG:HG2	5:A:563:HOH:O	2.21	0.41
1:C:282:ASP:HA	5:C:548:HOH:O	2.21	0.40
1:A:86:GLU:CB	1:A:125:ARG:HG2	2.52	0.40
1:B:233:ARG:NE	1:B:281:VAL:HG23	2.37	0.40
1:B:260:ASP:HA	1:B:261:PRO:HD3	1.88	0.40
1:C:270:ARG:NH2	1:C:278:GLN:OE1	2.45	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	245/251 (98%)	233 (95%)	12 (5%)	0	100	100
1	B	242/251 (96%)	228 (94%)	13 (5%)	1 (0%)	43	52
1	C	248/251 (99%)	235 (95%)	12 (5%)	1 (0%)	43	52
1	D	243/251 (97%)	229 (94%)	14 (6%)	0	100	100
All	All	978/1004 (97%)	925 (95%)	51 (5%)	2 (0%)	56	68

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	326	PHE
1	C	326	PHE

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	215/218 (99%)	203 (94%)	12 (6%)	30	38
1	B	212/218 (97%)	200 (94%)	12 (6%)	29	37
1	C	218/218 (100%)	204 (94%)	14 (6%)	25	31
1	D	213/218 (98%)	200 (94%)	13 (6%)	26	34
All	All	858/872 (98%)	807 (94%)	51 (6%)	27	35

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	82	GLU
1	A	109	LEU
1	A	152	SER
1	A	156	THR
1	A	172	GLU
1	A	190	LEU
1	A	213	LEU
1	A	236	LYS
1	A	258	LEU
1	A	275	LYS

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Mol	Chain	Res	Type
1	A	281	VAL
1	A	326	PHE
1	B	109	LEU
1	B	193	LEU
1	B	226	ARG
1	B	242	LEU
1	B	270	ARG
1	B	281	VAL
1	B	289	THR
1	B	294	VAL
1	B	296	SER
1	B	298	THR
1	B	301	SER
1	B	324	CYS
1	C	80	ILE
1	C	109	LEU
1	C	166	LEU
1	C	190	LEU
1	C	193	LEU
1	C	198	THR
1	C	200	VAL
1	C	242	LEU
1	C	276	GLN
1	C	278	GLN
1	C	293	HIS
1	C	298	THR
1	C	313	MET
1	C	326	PHE
1	D	102	ARG
1	D	109	LEU
1	D	152	SER
1	D	154	ASN
1	D	158	TYR
1	D	190	LEU
1	D	193	LEU
1	D	226	ARG
1	D	242	LEU
1	D	270	ARG
1	D	282	ASP
1	D	283	ARG
1	D	325	THR

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

sidechains are listed below:

Mol	Chain	Res	Type
1	C	276	GLN

5.3.3 RNA ⓘ

There are no RNA chains 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 12 ligands modelled in this entry, 4 are monoatomic - leaving 8 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	$\# Z > 2$
3	UDP	A	402	2	26,26,26	1.02	1 (3%)	36,40,40	1.42	5 (13%)
4	TRS	A	403	-	7,7,7	1.08	1 (14%)	9,9,9	1.95	3 (33%)
3	UDP	B	402	2	26,26,26	1.02	1 (3%)	36,40,40	1.50	4 (11%)
4	TRS	B	403	-	7,7,7	0.87	1 (14%)	9,9,9	0.56	0
3	UDP	C	402	2	26,26,26	1.08	1 (3%)	36,40,40	1.40	3 (8%)
4	TRS	C	403	-	7,7,7	0.94	1 (14%)	9,9,9	0.75	0
3	UDP	D	402	2	26,26,26	1.06	1 (3%)	36,40,40	1.31	2 (5%)
4	TRS	D	403	-	7,7,7	1.03	1 (14%)	9,9,9	1.46	2 (22%)

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	402	2	-	0/14/32/32	0/2/2/2
4	TRS	A	403	-	-	0/9/9/9	0/0/0/0
3	UDP	B	402	2	-	0/14/32/32	0/2/2/2
4	TRS	B	403	-	-	0/9/9/9	0/0/0/0
3	UDP	C	402	2	-	0/14/32/32	0/2/2/2
4	TRS	C	403	-	-	0/9/9/9	0/0/0/0
3	UDP	D	402	2	-	0/14/32/32	0/2/2/2
4	TRS	D	403	-	-	0/9/9/9	0/0/0/0

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	402	UDP	C6-C5	2.89	1.40	1.36
3	C	402	UDP	C6-C5	2.86	1.40	1.36
4	A	403	TRS	C-N	-2.64	1.46	1.50
3	B	402	UDP	C6-C5	2.63	1.40	1.36
3	D	402	UDP	C6-C5	2.57	1.40	1.36
4	D	403	TRS	C-N	-2.48	1.47	1.50
4	C	403	TRS	C-N	-2.40	1.47	1.50
4	B	403	TRS	C-N	-2.26	1.47	1.50

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	402	UDP	N3-C2-N1	6.62	121.50	115.97
3	D	402	UDP	N3-C2-N1	6.35	121.27	115.97
3	A	402	UDP	N3-C2-N1	6.06	121.03	115.97
3	C	402	UDP	N3-C2-N1	5.54	120.60	115.97
3	C	402	UDP	C2-N1-C1'	3.68	120.51	118.21
4	A	403	TRS	C2-C-C1	-3.24	105.39	110.70
4	A	403	TRS	C3-C-C2	3.10	115.79	110.70
3	A	402	UDP	C2-N1-C1'	2.53	119.79	118.21
4	A	403	TRS	O3-C3-C	-2.50	106.08	111.55
4	D	403	TRS	C1-C-N	2.44	114.94	108.30
3	A	402	UDP	PA-O3A-PB	-2.34	124.81	131.68
3	D	402	UDP	C2-N1-C1'	2.10	119.53	118.21
3	C	402	UDP	C5-C4-N3	2.07	120.61	116.70
3	A	402	UDP	C5-C4-N3	2.07	120.60	116.70
3	A	402	UDP	O4'-C1'-N1	2.05	112.39	108.06

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	403	TRS	C3-C-C1	-2.04	107.35	110.70
3	B	402	UDP	PA-O3A-PB	-2.04	125.72	131.68
3	B	402	UDP	C2-N1-C1'	-2.02	116.94	118.21
3	B	402	UDP	C6-N1-C1'	2.02	124.33	119.33

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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, 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	246/251 (98%)	-0.23	13 (5%) 25 35	30, 41, 83, 99	0
1	B	240/251 (95%)	-0.24	3 (1%) 74 82	31, 45, 68, 80	0
1	C	249/251 (99%)	-0.17	9 (3%) 41 51	32, 46, 76, 98	0
1	D	243/251 (96%)	-0.20	2 (0%) 83 90	33, 45, 69, 85	0
All	All	978/1004 (97%)	-0.21	27 (2%) 52 60	30, 45, 72, 99	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	272	ALA	6.4
1	C	80	ILE	5.7
1	A	281	VAL	5.5
1	A	279	PHE	5.3
1	B	327	SER	5.2
1	D	327	SER	5.0
1	B	88	ALA	5.0
1	C	279	PHE	4.9
1	C	81	PRO	4.4
1	A	271	ILE	3.8
1	A	327	SER	3.6
1	A	282	ASP	3.6
1	B	326	PHE	3.3
1	D	326	PHE	3.3
1	C	327	SER	3.2
1	C	326	PHE	3.1
1	A	273	ALA	3.0
1	C	283	ARG	2.9
1	C	271	ILE	2.9
1	A	283	ARG	2.8
1	A	280	LYS	2.6

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Mol	Chain	Res	Type	RSRZ
1	C	281	VAL	2.6
1	A	276	GLN	2.6
1	A	262	ALA	2.3
1	A	284	GLU	2.2
1	A	278	GLN	2.1
1	C	275	LYS	2.1

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 ⓘ

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
4	TRS	D	403	8/8	0.20	2.34	31,34,37,49	0
2	MN	B	401	1/1	0.13	2.20	41,41,41,41	0
4	TRS	B	403	8/8	0.18	1.39	39,40,43,43	0
4	TRS	A	403	8/8	0.14	1.19	24,33,40,42	0
4	TRS	C	403	8/8	0.13	0.75	32,39,43,44	0
3	UDP	D	402	25/25	0.14	0.20	27,34,39,41	0
3	UDP	B	402	25/25	0.14	0.17	30,37,43,45	0
3	UDP	C	402	25/25	0.12	0.08	31,36,39,42	0
2	MN	C	401	1/1	0.10	-0.35	36,36,36,36	0
3	UDP	A	402	25/25	0.10	-0.44	31,34,40,41	0
2	MN	A	401	1/1	0.07	-1.33	38,38,38,38	0
2	MN	D	401	1/1	0.09	-1.55	33,33,33,33	0

6.5 Other polymers ⓘ

There are no such residues in this entry.