



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 01:35 AM UTC

PDB ID : 2VSN / pdb_00002vsn
Title : Structure and topological arrangement of an O-GlcNAc transferase homolog:
insight into molecular control of intracellular glycosylation
Authors : Martinez-Fleites, C.; Macauley, M.S.; He, Y.; Shen, D.; Vocadlo, D.; Davies,
G.J.
Deposited on : 2008-04-28
Resolution : 2.75 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

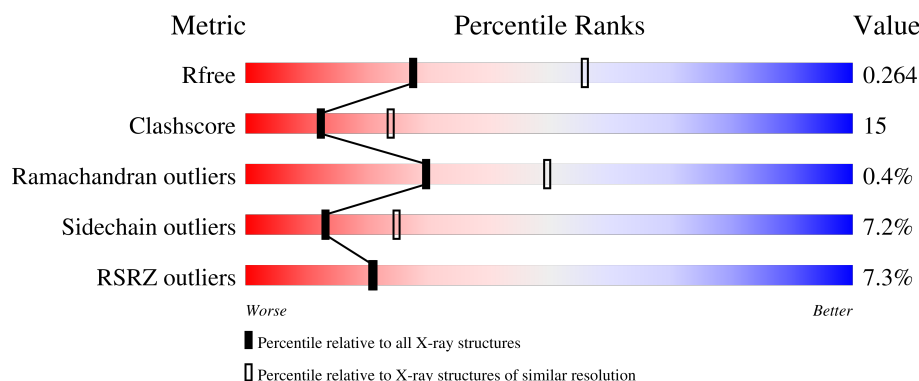
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.75 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}	180053	1009 (2.76-2.76)
Clashscore	190562	1044 (2.76-2.76)
Ramachandran outliers	187476	1024 (2.76-2.76)
Sidechain outliers	187428	1024 (2.76-2.76)
RSRZ outliers	180081	1009 (2.76-2.76)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	568	6% 62% 27% 5% 6%
1	B	568	7% 65% 24% 5% 6%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 8234 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.

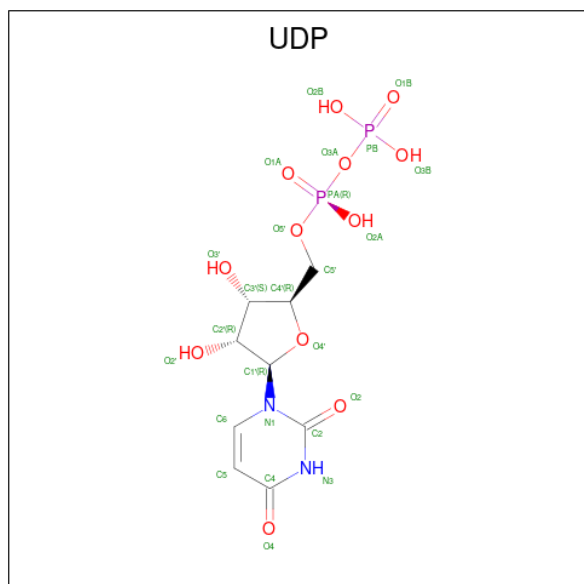
- Molecule 1 is a protein called XCOGT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	534	Total	C	N	O	S	0	0	0
			4055	2563	746	731	15			
1	B	534	Total	C	N	O	S	0	0	0
			4058	2567	749	727	15			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	290	ALA	ARG	engineered mutation	UNP Q4URB5
B	290	ALA	ARG	engineered mutation	UNP Q4URB5

- Molecule 2 is URIDINE-5'-DIPHOSPHATE (CCD ID: UDP) (formula: $C_9H_{14}N_2O_{12}P_2$).



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

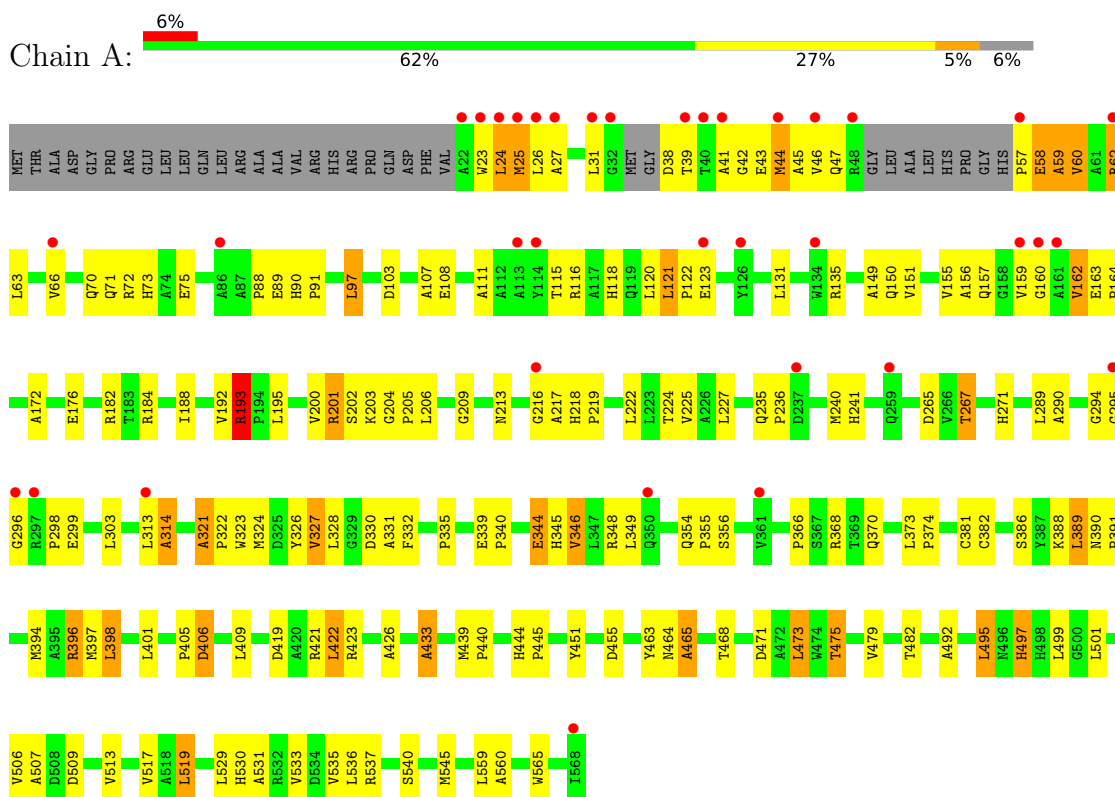
- Molecule 3 is water.

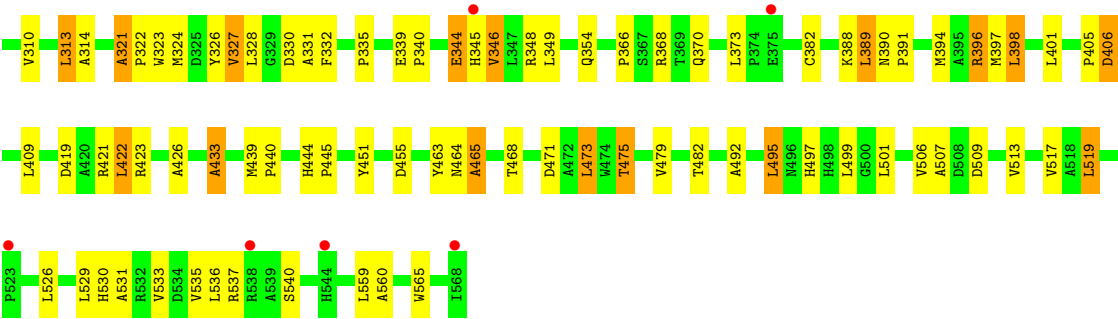
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	41	Total	O	0	0
			41	41		
3	B	30	Total	O	0	0
			30	30		

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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: XCOGT





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	86.69Å 105.69Å 151.04Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	75.52 – 2.75 75.52 – 2.75	Depositor EDS
% Data completeness (in resolution range)	99.7 (75.52-2.75) 99.8 (75.52-2.75)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.73 (at 2.73Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.241 , 0.265 0.241 , 0.264	Depositor DCC
R_{free} test set	1818 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	48.0	Xtriage
Anisotropy	0.044	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 41.6	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	0.91	EDS
Total number of atoms	8234	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.05% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.48	0/4158	1.05	32/5674 (0.6%)
1	B	0.46	0/4161	1.02	25/5676 (0.4%)
All	All	0.47	0/8319	1.03	57/11350 (0.5%)

There are no bond length outliers.

All (57) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	313	LEU	CB-CA-C	-11.78	92.76	111.48
1	A	58	GLU	N-CA-C	-11.75	99.77	112.93
1	A	313	LEU	CB-CA-C	-11.50	95.01	111.80
1	B	201	ARG	N-CA-C	11.06	123.34	111.28
1	A	201	ARG	N-CA-C	11.00	123.27	111.28
1	B	313	LEU	N-CA-C	9.36	124.07	112.47
1	A	314	ALA	N-CA-CB	8.52	124.89	110.49
1	A	313	LEU	N-CA-C	8.41	122.98	111.54
1	B	121	LEU	CA-C-N	8.07	128.15	119.28
1	B	121	LEU	C-N-CA	8.07	128.15	119.28
1	B	465	ALA	N-CA-C	-8.05	96.43	109.96
1	A	121	LEU	CA-C-N	7.66	127.70	119.28
1	A	121	LEU	C-N-CA	7.66	127.70	119.28
1	A	465	ALA	N-CA-C	-7.57	97.24	109.96
1	A	295	GLY	N-CA-C	-7.42	95.61	113.18
1	A	213	ASN	N-CA-C	-7.41	103.37	113.30
1	B	213	ASN	N-CA-C	-7.34	103.47	113.30
1	A	31	LEU	CB-CA-C	6.78	123.61	110.46
1	B	59	ALA	N-CA-C	-6.55	103.64	111.69
1	A	354	GLN	N-CA-C	6.46	118.86	109.04
1	B	354	GLN	N-CA-C	6.44	118.83	109.04
1	A	60	VAL	N-CA-C	-6.37	103.67	110.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	224	THR	N-CA-C	6.27	121.09	113.38
1	B	182	ARG	CB-CA-C	-6.16	100.95	110.81
1	A	182	ARG	CB-CA-C	-6.13	101.00	110.81
1	A	107	ALA	N-CA-C	5.96	117.47	110.97
1	B	471	ASP	N-CA-C	-5.96	104.71	111.14
1	B	224	THR	N-CA-C	5.92	120.66	113.38
1	A	24	LEU	N-CA-C	5.90	118.19	111.11
1	B	507	ALA	N-CA-C	5.88	120.52	113.23
1	A	507	ALA	N-CA-C	5.84	120.48	113.23
1	A	314	ALA	N-CA-C	-5.72	98.61	110.80
1	A	321	ALA	CA-C-N	5.69	125.36	119.56
1	A	321	ALA	C-N-CA	5.69	125.36	119.56
1	A	59	ALA	N-CA-C	-5.68	105.37	112.93
1	B	225	VAL	N-CA-C	5.65	116.29	110.36
1	B	107	ALA	N-CA-C	5.61	117.08	110.97
1	A	346	VAL	N-CA-C	5.54	116.21	108.89
1	B	346	VAL	N-CA-C	5.53	116.18	108.89
1	B	200	VAL	N-CA-C	-5.36	102.67	109.80
1	A	225	VAL	N-CA-C	5.33	115.95	110.36
1	B	32	GLU	N-CA-C	5.32	117.77	111.33
1	A	63	LEU	N-CA-C	-5.30	105.50	111.28
1	A	193	ARG	CA-C-N	5.29	125.29	119.89
1	A	193	ARG	C-N-CA	5.29	125.29	119.89
1	A	386	SER	N-CA-C	5.26	117.70	111.33
1	B	290	ALA	N-CA-C	-5.24	106.02	112.72
1	B	63	LEU	N-CA-C	-5.23	105.58	111.28
1	A	62	ARG	CG-CD-NE	-5.20	100.56	112.00
1	B	62	ARG	CG-CD-NE	-5.19	100.59	112.00
1	B	321	ALA	CA-C-N	5.16	124.95	119.28
1	B	321	ALA	C-N-CA	5.16	124.95	119.28
1	A	200	VAL	N-CA-C	-5.14	102.97	109.80
1	B	192	VAL	N-CA-C	5.12	116.03	108.45
1	B	310	VAL	N-CA-C	5.08	115.79	108.48
1	A	471	ASP	N-CA-C	-5.07	105.67	111.14
1	A	192	VAL	N-CA-C	5.01	115.87	108.45

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	4055	0	3977	128	0
1	B	4058	0	3993	126	0
2	A	25	0	11	0	0
2	B	25	0	11	0	0
3	A	41	0	0	3	0
3	B	30	0	0	1	0
All	All	8234	0	7992	249	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 15.

All (249) 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:202:SER:HA	1:B:565:TRP:HA	1.57	0.87
1:A:202:SER:HA	1:A:565:TRP:HA	1.57	0.87
1:B:206:LEU:HD21	1:B:560:ALA:HB2	1.59	0.82
1:A:332:PHE:CD1	1:A:497:HIS:HD2	1.96	0.82
1:A:206:LEU:HD21	1:A:560:ALA:HB2	1.61	0.81
1:A:38:ASP:HB3	1:A:41:ALA:HB3	1.62	0.81
1:A:332:PHE:HD1	1:A:497:HIS:HD2	1.28	0.79
1:B:23:VAL:HG12	1:B:24:ALA:H	1.47	0.79
1:B:328:LEU:HD11	1:B:349:LEU:HG	1.64	0.78
1:A:328:LEU:HD11	1:A:349:LEU:HG	1.67	0.77
1:B:451:TYR:HB3	1:B:475:THR:HG21	1.67	0.76
1:A:327:VAL:HG13	1:A:344:GLU:HG3	1.66	0.76
1:B:29:ALA:O	1:B:32:GLU:HG2	1.85	0.76
1:B:327:VAL:HG13	1:B:344:GLU:HG3	1.67	0.76
1:B:366:PRO:HB3	1:B:370:GLN:HE21	1.51	0.75
1:A:366:PRO:HB3	1:A:370:GLN:HE21	1.53	0.72
1:B:396:ARG:HD2	1:B:509:ASP:OD2	1.89	0.72
1:A:451:TYR:HB3	1:A:475:THR:HG21	1.71	0.71
1:A:58:GLU:OE2	1:A:90:HIS:NE2	2.21	0.71
1:B:332:PHE:CD1	1:B:497:HIS:HD2	2.09	0.70
1:B:62:ARG:O	1:B:66:VAL:HG23	1.92	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:62:ARG:O	1:A:66:VAL:HG23	1.93	0.69
1:A:159:VAL:HG22	1:A:160:GLY:H	1.58	0.68
1:A:464:ASN:HD22	1:A:492:ALA:H	1.41	0.68
1:B:44:GLY:O	1:B:48:VAL:HG23	1.94	0.67
1:B:451:TYR:CB	1:B:475:THR:HG21	2.23	0.67
1:A:157:GLN:HA	1:B:182:ARG:HD2	1.76	0.67
1:A:451:TYR:CB	1:A:475:THR:HG21	2.25	0.67
1:A:396:ARG:HD2	1:A:509:ASP:OD2	1.93	0.67
1:B:464:ASN:HD22	1:B:492:ALA:H	1.41	0.66
1:B:240:MET:HG3	3:B:2010:HOH:O	1.96	0.66
1:B:332:PHE:HD1	1:B:497:HIS:HD2	1.45	0.65
1:B:58:GLU:O	1:B:58:GLU:HG3	1.95	0.65
1:B:290:ALA:HB1	1:B:296:GLY:HA2	1.77	0.65
1:A:327:VAL:HG13	1:A:344:GLU:CG	2.27	0.64
1:B:159:VAL:HG22	1:B:160:GLY:H	1.61	0.64
1:A:327:VAL:HG22	1:A:344:GLU:HG2	1.80	0.64
1:B:290:ALA:CB	1:B:296:GLY:HA2	2.28	0.64
1:A:58:GLU:O	1:A:62:ARG:CD	2.46	0.63
1:B:465:ALA:HB1	1:B:468:THR:HB	1.80	0.63
1:A:157:GLN:HA	1:B:182:ARG:CD	2.29	0.63
1:A:401:LEU:HD11	1:A:409:LEU:HB2	1.81	0.62
1:B:401:LEU:HD11	1:B:409:LEU:HB2	1.81	0.62
1:B:389:LEU:HG	1:B:394:MET:HE3	1.81	0.62
1:B:327:VAL:HG13	1:B:344:GLU:CG	2.28	0.62
1:A:332:PHE:CD1	1:A:497:HIS:CD2	2.84	0.61
1:B:115:THR:CG2	1:B:135:ARG:HH21	2.13	0.61
1:B:206:LEU:HD21	1:B:560:ALA:CB	2.28	0.61
1:A:43:GLU:HB2	1:A:66:VAL:HG11	1.83	0.61
1:B:382:CYS:SG	1:B:397:MET:HE1	2.40	0.61
1:B:30:ASP:OD1	1:B:62:ARG:NH1	2.33	0.61
1:B:519:LEU:HD21	1:B:529:LEU:HD22	1.81	0.61
1:A:398:LEU:HD21	1:A:426:ALA:HA	1.83	0.61
1:A:519:LEU:HD21	1:A:529:LEU:HD22	1.83	0.60
1:A:374:PRO:HD3	3:A:2028:HOH:O	2.00	0.60
1:A:419:ASP:O	1:A:423:ARG:HG3	2.01	0.60
1:A:465:ALA:HB1	1:A:468:THR:HB	1.82	0.60
1:B:327:VAL:HG22	1:B:344:GLU:HG2	1.83	0.60
1:A:184:ARG:HH21	1:A:188:ILE:HD11	1.67	0.60
1:A:72:ARG:HB3	1:A:75:GLU:OE1	2.02	0.60
1:B:72:ARG:HB3	1:B:75:GLU:OE1	2.02	0.60
1:A:222:LEU:HD21	1:A:444:HIS:CE1	2.36	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:115:THR:CG2	1:A:135:ARG:HH21	2.15	0.60
1:B:184:ARG:HH21	1:B:188:ILE:HD11	1.67	0.59
1:A:58:GLU:O	1:A:62:ARG:HD2	2.02	0.59
1:A:389:LEU:HG	1:A:394:MET:HE3	1.85	0.59
1:B:332:PHE:CD1	1:B:497:HIS:CD2	2.89	0.59
1:B:121:LEU:N	1:B:122:PRO:HD3	2.17	0.59
1:B:398:LEU:HD21	1:B:426:ALA:HA	1.85	0.59
1:B:222:LEU:HD21	1:B:444:HIS:CE1	2.38	0.58
1:A:536:LEU:O	1:A:540:SER:HB3	2.04	0.58
1:A:149:ALA:HB3	1:B:340:PRO:HG2	1.84	0.58
1:A:121:LEU:N	1:A:122:PRO:HD3	2.19	0.58
1:A:382:CYS:SG	1:A:397:MET:HE1	2.43	0.58
1:A:206:LEU:HD21	1:A:560:ALA:CB	2.30	0.58
1:A:155:VAL:HG22	1:A:162:VAL:HG11	1.85	0.57
1:B:405:PRO:O	1:B:406:ASP:HB2	2.05	0.57
1:B:536:LEU:O	1:B:540:SER:HB3	2.05	0.57
1:B:419:ASP:O	1:B:423:ARG:HG3	2.04	0.57
1:A:42:GLY:O	1:A:46:VAL:HG23	2.04	0.56
1:A:501:LEU:HD11	1:A:533:VAL:HG22	1.87	0.56
1:A:405:PRO:O	1:A:406:ASP:HB2	2.04	0.56
1:B:155:VAL:HG22	1:B:162:VAL:HG11	1.88	0.56
1:B:296:GLY:O	1:B:297:ARG:HB2	2.05	0.56
1:B:339:GLU:HG2	1:B:346:VAL:HG21	1.86	0.56
1:B:26:LEU:HD23	1:B:62:ARG:HD2	1.88	0.56
1:B:501:LEU:HD11	1:B:533:VAL:HG22	1.87	0.56
1:A:330:ASP:C	1:A:348:ARG:HD3	2.31	0.55
1:A:58:GLU:O	1:A:62:ARG:NE	2.39	0.55
1:B:366:PRO:HB3	1:B:370:GLN:NE2	2.21	0.55
1:A:339:GLU:HG2	1:A:346:VAL:HG21	1.87	0.55
1:B:444:HIS:HB3	1:B:445:PRO:HD3	1.89	0.55
1:B:330:ASP:C	1:B:348:ARG:HD3	2.31	0.55
1:B:473:LEU:HD13	1:B:479:VAL:HG23	1.89	0.54
1:B:33:LEU:N	1:B:33:LEU:HD12	2.22	0.54
1:B:390:ASN:HB2	1:B:391:PRO:HD2	1.90	0.54
1:B:482:THR:HG22	1:B:506:VAL:HG23	1.90	0.54
1:A:290:ALA:HA	3:A:2005:HOH:O	2.06	0.53
1:A:290:ALA:HB1	1:A:296:GLY:HA2	1.90	0.53
1:A:482:THR:HG22	1:A:506:VAL:HG23	1.91	0.53
1:A:294:GLY:O	1:A:296:GLY:N	2.42	0.53
1:B:388:LYS:HE2	1:B:463:TYR:CD1	2.44	0.52
1:B:88:PRO:O	1:B:89:GLU:HG3	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:473:LEU:HD13	1:A:479:VAL:HG23	1.91	0.52
1:B:390:ASN:HB2	1:B:391:PRO:CD	2.39	0.52
1:B:455:ASP:O	1:B:530:HIS:HE1	1.92	0.52
1:B:332:PHE:HD1	1:B:497:HIS:CD2	2.26	0.52
1:A:88:PRO:O	1:A:89:GLU:HG3	2.10	0.52
1:A:444:HIS:HB3	1:A:445:PRO:HD3	1.90	0.52
1:B:155:VAL:CG2	1:B:162:VAL:HG11	2.40	0.52
1:A:390:ASN:HB2	1:A:391:PRO:HD2	1.92	0.52
1:A:156:ALA:O	1:B:182:ARG:NH1	2.43	0.51
1:A:235:GLN:N	1:A:236:PRO:HD3	2.25	0.51
1:B:235:GLN:N	1:B:236:PRO:HD3	2.25	0.51
1:A:388:LYS:HE2	1:A:463:TYR:CD1	2.45	0.51
1:A:155:VAL:CG2	1:A:162:VAL:HG11	2.40	0.51
1:A:390:ASN:HB2	1:A:391:PRO:CD	2.41	0.51
1:A:323:TRP:CD1	1:A:324:MET:HE2	2.46	0.50
1:A:455:ASP:O	1:A:530:HIS:HE1	1.94	0.50
1:A:57:PRO:C	1:A:59:ALA:H	2.18	0.50
1:A:73:HIS:CE1	1:A:103:ASP:HB3	2.47	0.49
1:B:327:VAL:HG23	1:B:346:VAL:HA	1.94	0.49
1:A:366:PRO:HB3	1:A:370:GLN:NE2	2.22	0.49
1:B:115:THR:HG22	1:B:135:ARG:HH21	1.76	0.49
1:B:201:ARG:HB3	1:B:201:ARG:CZ	2.42	0.49
1:B:513:VAL:O	1:B:517:VAL:HG23	2.11	0.49
1:A:25:MET:HE3	1:A:26:LEU:H	1.77	0.49
1:B:331:ALA:N	1:B:348:ARG:HD3	2.27	0.49
1:A:218:HIS:ND1	1:A:219:PRO:HD2	2.27	0.49
1:B:204:GLY:CA	1:B:565:TRP:HE1	2.25	0.49
1:B:289:LEU:C	1:B:314:ALA:HB2	2.38	0.49
1:A:201:ARG:CZ	1:A:201:ARG:HB3	2.42	0.49
1:B:45:GLU:O	1:B:49:GLN:HG2	2.13	0.49
1:A:193:ARG:NH1	1:A:193:ARG:HG2	2.28	0.48
1:B:218:HIS:ND1	1:B:219:PRO:HD2	2.28	0.48
1:A:513:VAL:O	1:A:517:VAL:HG23	2.13	0.48
1:B:73:HIS:CE1	1:B:103:ASP:HB3	2.49	0.48
1:A:204:GLY:CA	1:A:565:TRP:HE1	2.26	0.48
1:B:423:ARG:HB3	1:B:433:ALA:HB1	1.95	0.48
1:B:193:ARG:NH1	1:B:193:ARG:HG2	2.28	0.48
1:A:327:VAL:HG23	1:A:346:VAL:HA	1.96	0.48
1:B:23:VAL:HG12	1:B:24:ALA:N	2.23	0.47
1:B:46:MET:O	1:B:50:ARG:HG2	2.14	0.47
1:B:58:GLU:O	1:B:58:GLU:CG	2.61	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:97:LEU:HD22	1:A:97:LEU:O	2.14	0.47
1:A:115:THR:HG22	1:A:135:ARG:HH21	1.78	0.47
1:A:423:ARG:HB3	1:A:433:ALA:HB1	1.95	0.47
1:B:193:ARG:HG2	1:B:193:ARG:HH11	1.80	0.47
1:A:193:ARG:HG2	1:A:193:ARG:HH11	1.79	0.47
1:A:265:ASP:OD1	1:A:267:THR:HB	2.15	0.47
1:A:298:PRO:HG2	1:A:299:GLU:OE2	2.15	0.47
1:B:26:LEU:HD12	1:B:26:LEU:H	1.80	0.47
1:A:326:TYR:CD1	1:A:345:HIS:HB3	2.50	0.46
1:A:331:ALA:N	1:A:348:ARG:HD3	2.30	0.46
1:B:23:VAL:CG1	1:B:24:ALA:H	2.19	0.46
1:A:394:MET:HA	1:A:394:MET:HE2	1.97	0.46
1:A:150:GLN:HA	1:B:340:PRO:HB3	1.98	0.46
1:B:323:TRP:CD1	1:B:324:MET:HE2	2.51	0.46
1:A:290:ALA:CB	1:A:296:GLY:HA2	2.46	0.46
1:B:326:TYR:CD1	1:B:345:HIS:HB3	2.51	0.46
1:A:24:LEU:HD23	1:A:62:ARG:CD	2.46	0.45
1:B:313:LEU:O	1:B:313:LEU:HG	2.16	0.45
1:B:509:ASP:O	1:B:513:VAL:HG23	2.17	0.45
1:B:70:GLN:O	1:B:72:ARG:HG3	2.16	0.45
1:A:271:HIS:H	1:A:271:HIS:HD1	1.64	0.45
1:A:451:TYR:HB2	1:A:475:THR:HG21	1.98	0.45
1:A:59:ALA:O	1:A:60:VAL:C	2.59	0.45
1:A:91:PRO:HB3	1:A:120:LEU:HB2	1.98	0.45
1:B:70:GLN:C	1:B:72:ARG:H	2.24	0.45
1:A:70:GLN:O	1:A:72:ARG:HG3	2.17	0.45
1:B:91:PRO:HB3	1:B:120:LEU:HB2	1.99	0.45
1:B:339:GLU:N	1:B:340:PRO:CD	2.80	0.45
1:A:201:ARG:HG3	1:A:204:GLY:H	1.83	0.44
1:A:339:GLU:N	1:A:340:PRO:CD	2.80	0.44
1:B:394:MET:HE2	1:B:394:MET:HA	1.98	0.44
1:A:398:LEU:HD12	1:A:398:LEU:HA	1.84	0.44
1:B:172:ALA:HB1	1:B:176:GLU:HB2	1.98	0.44
1:A:70:GLN:C	1:A:72:ARG:H	2.25	0.44
1:A:327:VAL:HG22	1:A:344:GLU:CG	2.46	0.44
1:A:24:LEU:HD23	1:A:62:ARG:HD2	1.99	0.44
1:B:184:ARG:O	1:B:188:ILE:HG12	2.18	0.44
1:A:57:PRO:HD2	1:A:60:VAL:HG23	2.00	0.44
1:B:265:ASP:OD1	1:B:267:THR:HB	2.18	0.44
1:B:332:PHE:O	1:B:335:PRO:HD3	2.18	0.44
1:B:201:ARG:HD2	1:B:205:PRO:O	2.17	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:201:ARG:HD2	1:A:205:PRO:O	2.18	0.43
1:A:388:LYS:HE2	1:A:463:TYR:CE1	2.54	0.43
1:A:321:ALA:HA	1:A:322:PRO:HD3	1.78	0.43
1:B:388:LYS:HE2	1:B:463:TYR:CE1	2.54	0.43
1:B:33:LEU:HD12	1:B:33:LEU:H	1.83	0.43
1:B:108:GLU:O	1:B:111:ALA:HB3	2.19	0.43
1:A:509:ASP:O	1:A:513:VAL:HG23	2.18	0.43
1:B:57:PRO:HB2	1:B:58:GLU:H	1.62	0.43
1:A:27:ALA:HB2	1:A:45:ALA:HB3	2.00	0.43
1:B:97:LEU:HD22	1:B:97:LEU:O	2.19	0.43
1:A:111:ALA:O	1:A:115:THR:HG23	2.19	0.43
1:A:495:LEU:O	1:A:499:LEU:HB2	2.19	0.42
1:B:111:ALA:O	1:B:115:THR:HG23	2.18	0.42
1:A:218:HIS:CG	1:A:219:PRO:HD2	2.54	0.42
1:A:332:PHE:O	1:A:335:PRO:HD3	2.19	0.42
1:B:217:ALA:H	1:B:253:LEU:HD12	1.84	0.42
1:B:67:ARG:HA	1:B:67:ARG:HD3	1.86	0.42
1:A:163:GLU:HA	1:A:164:PRO:HD3	1.96	0.42
1:B:201:ARG:HG3	1:B:204:GLY:H	1.84	0.42
1:B:439:MET:HA	1:B:440:PRO:HD3	1.90	0.42
1:B:339:GLU:CD	1:B:348:ARG:HH22	2.27	0.42
1:A:439:MET:HA	1:A:440:PRO:HD3	1.90	0.42
1:A:204:GLY:HA2	1:A:205:PRO:C	2.45	0.42
1:B:30:ASP:CG	1:B:62:ARG:NH1	2.78	0.42
1:A:115:THR:HG22	1:A:131:LEU:HD11	2.02	0.41
1:B:526:LEU:HD23	1:B:526:LEU:HA	1.85	0.41
1:A:172:ALA:HB1	1:A:176:GLU:HB2	2.01	0.41
1:A:289:LEU:C	1:A:314:ALA:HB2	2.44	0.41
1:A:184:ARG:O	1:A:188:ILE:HG12	2.21	0.41
1:A:108:GLU:O	1:A:111:ALA:HB3	2.20	0.41
1:A:155:VAL:HG22	1:A:162:VAL:CG1	2.49	0.41
1:B:32:GLU:OE1	1:B:32:GLU:N	2.53	0.41
1:B:108:GLU:CD	1:B:421:ARG:HH22	2.29	0.41
1:B:115:THR:HG22	1:B:131:LEU:HD11	2.03	0.41
1:A:45:ALA:C	1:A:47:GLN:N	2.78	0.41
1:A:116:ARG:C	1:A:118:HIS:N	2.78	0.41
1:A:240:MET:HG3	3:A:2010:HOH:O	2.19	0.41
1:A:323:TRP:NE1	1:A:324:MET:HE2	2.35	0.41
1:B:155:VAL:HG22	1:B:162:VAL:CG1	2.51	0.41
1:B:327:VAL:HG22	1:B:344:GLU:CG	2.50	0.41
1:A:44:MET:HA	1:A:44:MET:HE2	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:355:PRO:CB	1:A:545:MET:HE1	2.50	0.41
1:A:206:LEU:HD12	1:A:206:LEU:HA	1.95	0.41
1:A:531:ALA:O	1:A:535:VAL:HG23	2.21	0.41
1:B:70:GLN:C	1:B:72:ARG:N	2.79	0.41
1:A:108:GLU:CD	1:A:421:ARG:HH22	2.29	0.41
1:A:497:HIS:ND1	1:A:497:HIS:C	2.78	0.41
1:B:389:LEU:HD23	1:B:422:LEU:HD21	2.01	0.41
1:A:216:GLY:O	1:A:217:ALA:HB3	2.21	0.41
1:B:495:LEU:O	1:B:499:LEU:HB2	2.21	0.41
1:B:531:ALA:O	1:B:535:VAL:HG23	2.21	0.41
1:A:389:LEU:HD23	1:A:422:LEU:HD21	2.02	0.40
1:B:218:HIS:CG	1:B:219:PRO:HD2	2.56	0.40
1:A:151:VAL:O	1:A:155:VAL:HG23	2.20	0.40
1:A:209:GLY:HA2	1:A:241:HIS:O	2.22	0.40
1:B:313:LEU:O	1:B:313:LEU:CG	2.70	0.40
1:B:116:ARG:C	1:B:118:HIS:N	2.78	0.40
1:B:225:VAL:HG12	1:B:257:LEU:HG	2.03	0.40
1:A:66:VAL:O	1:A:70:GLN:HG3	2.22	0.40
1:B:64:GLY:HA2	1:B:79:LEU:HD12	2.03	0.40
1:B:451:TYR:HB2	1:B:475:THR:HG21	1.99	0.40
1:B:163:GLU:HA	1:B:164:PRO:HD3	1.97	0.40
1:B:321:ALA:HA	1:B:322:PRO:HD3	1.78	0.40
1:B:499:LEU:HA	1:B:499:LEU:HD12	1.81	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	528/568 (93%)	496 (94%)	30 (6%)	2 (0%)	30	50
1	B	528/568 (93%)	492 (93%)	34 (6%)	2 (0%)	30	50

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	1056/1136 (93%)	988 (94%)	64 (6%)	4 (0%)	30 50

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	433	ALA
1	B	433	ALA
1	A	71	GLN
1	B	71	GLN

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	401/428 (94%)	370 (92%)	31 (8%)	12 22
1	B	401/428 (94%)	374 (93%)	27 (7%)	15 28
All	All	802/856 (94%)	744 (93%)	58 (7%)	13 25

All (58) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	23	TRP
1	A	25	MET
1	A	39	THR
1	A	44	MET
1	A	97	LEU
1	A	123	GLU
1	A	162	VAL
1	A	193	ARG
1	A	195	LEU
1	A	203	LYS
1	A	227	LEU
1	A	267	THR
1	A	303	LEU
1	A	327	VAL

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Mol	Chain	Res	Type
1	A	344	GLU
1	A	356	SER
1	A	368	ARG
1	A	373	LEU
1	A	381	CYS
1	A	389	LEU
1	A	396	ARG
1	A	398	LEU
1	A	406	ASP
1	A	422	LEU
1	A	473	LEU
1	A	475	THR
1	A	495	LEU
1	A	497	HIS
1	A	519	LEU
1	A	537	ARG
1	A	559	LEU
1	B	32	GLU
1	B	45	GLU
1	B	46	MET
1	B	58	GLU
1	B	97	LEU
1	B	123	GLU
1	B	162	VAL
1	B	193	ARG
1	B	195	LEU
1	B	203	LYS
1	B	227	LEU
1	B	267	THR
1	B	327	VAL
1	B	344	GLU
1	B	368	ARG
1	B	373	LEU
1	B	389	LEU
1	B	396	ARG
1	B	398	LEU
1	B	406	ASP
1	B	422	LEU
1	B	473	LEU
1	B	475	THR
1	B	495	LEU
1	B	519	LEU

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Mol	Chain	Res	Type
1	B	537	ARG
1	B	559	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (36) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	47	GLN
1	A	82	GLN
1	A	118	HIS
1	A	119	GLN
1	A	130	GLN
1	A	150	GLN
1	A	157	GLN
1	A	239	GLN
1	A	277	HIS
1	A	311	ASN
1	A	345	HIS
1	A	350	GLN
1	A	370	GLN
1	A	376	GLN
1	A	385	ASN
1	A	392	GLN
1	A	497	HIS
1	A	530	HIS
1	A	557	GLN
1	B	82	GLN
1	B	118	HIS
1	B	119	GLN
1	B	130	GLN
1	B	177	GLN
1	B	239	GLN
1	B	311	ASN
1	B	345	HIS
1	B	350	GLN
1	B	370	GLN
1	B	376	GLN
1	B	385	ASN
1	B	392	GLN
1	B	466	HIS
1	B	497	HIS
1	B	530	HIS
1	B	557	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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

2 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$
2	UDP	B	900	-	25,26,26	1.36	4 (16%)	38,40,40	2.13	10 (26%)
2	UDP	A	900	-	25,26,26	1.47	4 (16%)	38,40,40	1.99	8 (21%)

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	B	900	-	-	8/16/32/32	0/2/2/2
2	UDP	A	900	-	-	5/16/32/32	0/2/2/2

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	900	UDP	PA-O3A	-3.38	1.55	1.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	900	UDP	PB-O3B	-2.92	1.43	1.54
2	B	900	UDP	PA-O3A	-2.78	1.56	1.59
2	B	900	UDP	PB-O3B	-2.69	1.44	1.54
2	B	900	UDP	C5-C4	-2.66	1.37	1.43
2	A	900	UDP	C5-C4	-2.63	1.38	1.43
2	A	900	UDP	PA-O2A	-2.40	1.44	1.55
2	B	900	UDP	PA-O2A	-2.19	1.45	1.55

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	900	UDP	C4-N3-C2	-5.88	119.31	126.61
2	A	900	UDP	C4-N3-C2	-5.18	120.19	126.61
2	B	900	UDP	O4'-C1'-N1	-5.07	96.87	108.36
2	B	900	UDP	N3-C2-N1	4.53	120.79	114.89
2	A	900	UDP	N3-C2-N1	4.37	120.58	114.89
2	A	900	UDP	O2A-PA-O3A	4.34	119.00	107.27
2	B	900	UDP	C5-C4-N3	3.92	120.29	114.80
2	B	900	UDP	O2A-PA-O3A	3.31	116.22	107.27
2	A	900	UDP	C5-C4-N3	3.15	119.21	114.80
2	A	900	UDP	C5'-C4'-C3'	-3.14	103.91	115.21
2	A	900	UDP	O4'-C1'-N1	-2.98	101.60	108.36
2	A	900	UDP	O2-C2-N1	-2.84	119.10	122.80
2	B	900	UDP	O4-C4-C5	-2.80	120.33	125.16
2	A	900	UDP	O4-C4-C5	-2.74	120.43	125.16
2	B	900	UDP	O2-C2-N1	-2.70	119.28	122.80
2	B	900	UDP	O3A-PA-O1A	-2.52	103.12	110.70
2	B	900	UDP	O4'-C1'-C2'	-2.21	101.89	106.62
2	B	900	UDP	C2'-C1'-N1	2.07	119.02	113.25

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	900	UDP	C5'-O5'-PA-O3A
2	B	900	UDP	C5'-O5'-PA-O1A
2	B	900	UDP	C5'-O5'-PA-O2A
2	B	900	UDP	C5'-O5'-PA-O3A
2	B	900	UDP	PA-O3A-PB-O2B
2	B	900	UDP	PA-O3A-PB-O3B
2	A	900	UDP	C3'-C4'-C5'-O5'
2	A	900	UDP	O4'-C4'-C5'-O5'

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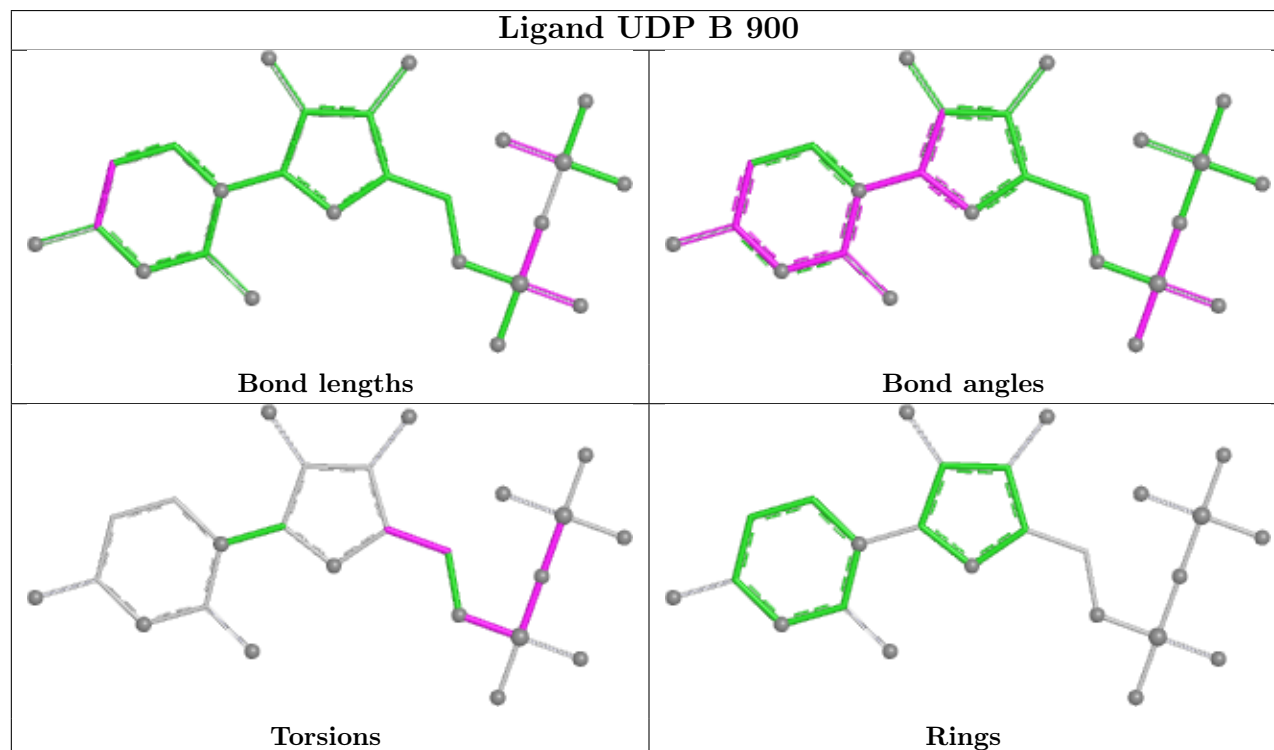
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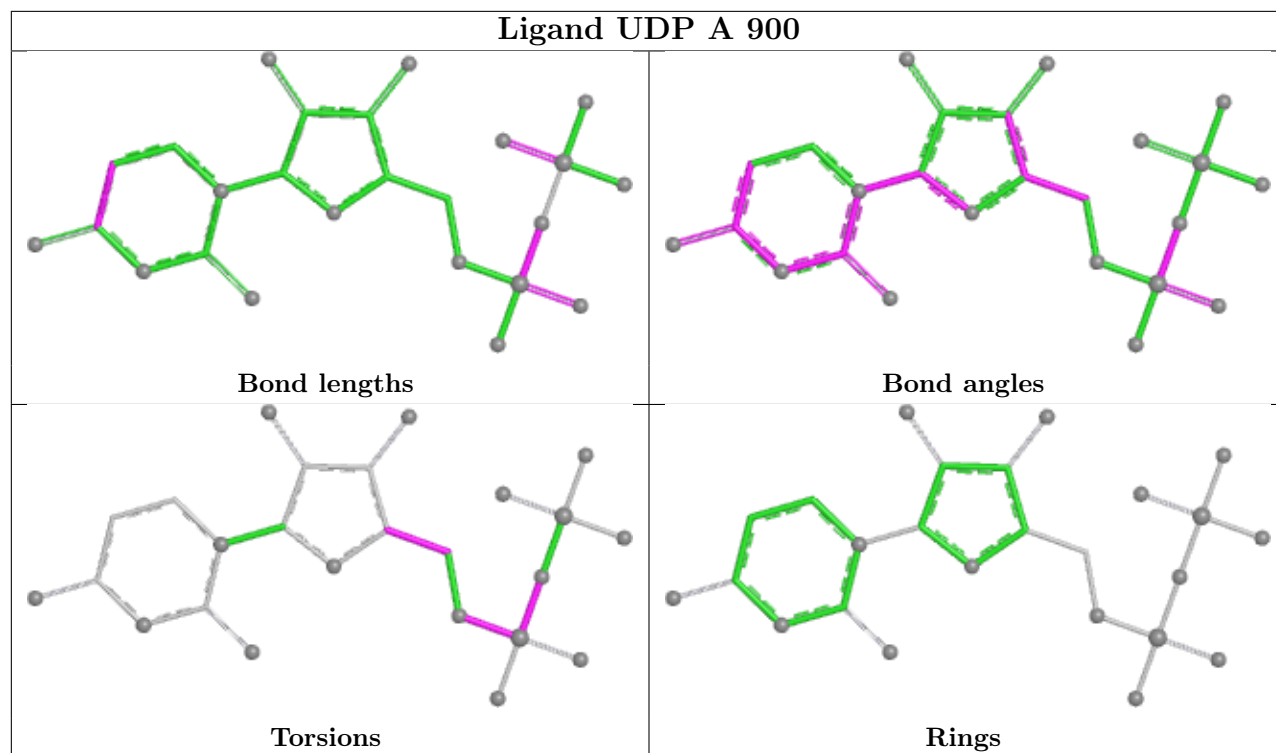
Mol	Chain	Res	Type	Atoms
2	B	900	UDP	O4'-C4'-C5'-O5'
2	B	900	UDP	C3'-C4'-C5'-O5'
2	A	900	UDP	PB-O3A-PA-O2A
2	B	900	UDP	PB-O3A-PA-O1A
2	A	900	UDP	PB-O3A-PA-O1A

There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





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	534/568 (94%)	0.54	36 (6%) 24 23	22, 42, 74, 111	0
1	B	534/568 (94%)	0.52	42 (7%) 18 19	22, 43, 74, 103	0
All	All	1068/1136 (94%)	0.53	78 (7%) 21 21	22, 43, 74, 111	0

All (78) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	23	TRP	6.7
1	A	48	ARG	6.3
1	B	52	LEU	5.5
1	B	23	VAL	5.4
1	B	57	PRO	5.2
1	A	22	ALA	5.2
1	A	57	PRO	5.2
1	A	31	LEU	5.1
1	A	26	LEU	4.6
1	A	297	ARG	4.5
1	B	24	ALA	4.4
1	A	361	VAL	4.1
1	A	24	LEU	4.0
1	B	568	ILE	3.9
1	B	203	LYS	3.7
1	B	87	ALA	3.7
1	B	34	GLY	3.3
1	A	39	THR	3.3
1	A	259	GLN	3.2
1	A	237	ASP	3.1
1	A	32	GLY	3.1
1	A	568	ILE	3.1
1	B	93	ILE	3.1
1	B	43	ALA	3.0

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Mol	Chain	Res	Type	RSRZ
1	B	204	GLY	2.9
1	A	159	VAL	2.9
1	B	46	MET	2.9
1	B	61	ALA	2.9
1	A	25	MET	2.8
1	A	123	GLU	2.8
1	B	59	ALA	2.8
1	A	40	THR	2.7
1	A	161	ALA	2.7
1	A	44	MET	2.6
1	A	126	TYR	2.6
1	B	202	SER	2.6
1	B	32	GLU	2.6
1	A	160	GLY	2.6
1	B	92	GLY	2.6
1	A	216	GLY	2.5
1	A	62	ARG	2.5
1	B	33	LEU	2.5
1	B	523	PRO	2.5
1	A	113	ALA	2.5
1	B	25	TRP	2.5
1	B	28	LEU	2.4
1	B	88	PRO	2.4
1	B	114	TYR	2.4
1	A	66	VAL	2.4
1	B	91	PRO	2.4
1	A	350	GLN	2.4
1	B	60	VAL	2.4
1	A	296	GLY	2.3
1	B	123	GLU	2.3
1	B	296	GLY	2.3
1	A	46	VAL	2.3
1	A	313	LEU	2.3
1	B	26	LEU	2.3
1	B	58	GLU	2.2
1	A	114	TYR	2.2
1	A	134	TRP	2.2
1	A	27	ALA	2.2
1	A	295	GLY	2.2
1	B	159	VAL	2.2
1	B	216	GLY	2.2
1	B	544	HIS	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	71	GLN	2.1
1	B	27	MET	2.1
1	B	271	HIS	2.1
1	B	345	HIS	2.1
1	A	86	ALA	2.1
1	B	137	ARG	2.0
1	A	41	ALA	2.0
1	B	31	ALA	2.0
1	B	375	GLU	2.0
1	B	538	ARG	2.0
1	B	186	GLN	2.0
1	B	295	GLY	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

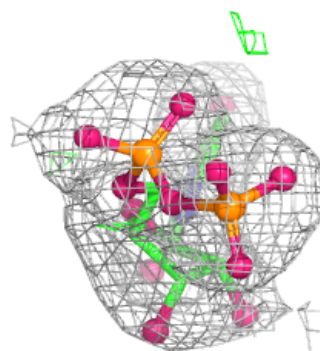
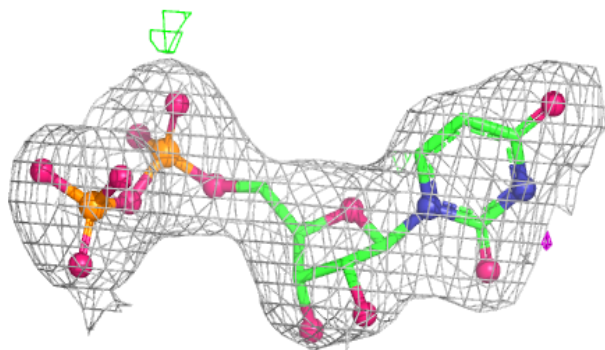
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	UDP	A	900	25/25	0.95	0.09	37,42,47,48	0
2	UDP	B	900	25/25	0.96	0.07	41,46,49,51	0

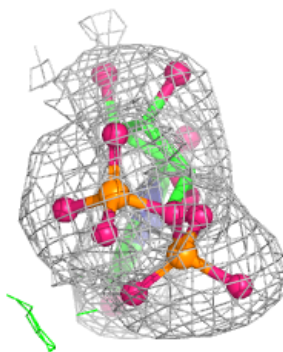
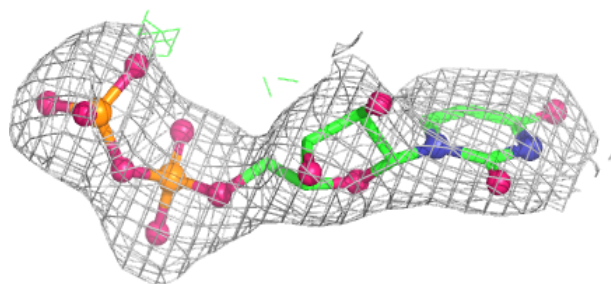
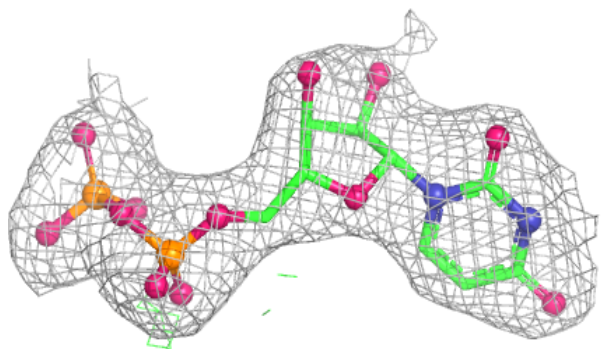
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around UDP A 900:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around UDP B 900:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.