



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 12:14 AM UTC

PDB ID : 3PE3 / pdb_00003pe3
Title : Structure of human O-GlcNAc transferase and its complex with a peptide substrate
Authors : Lazarus, M.B.; Nam, Y.; Jiang, J.; Sliz, P.; Walker, S.
Deposited on : 2010-10-25
Resolution : 2.78 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

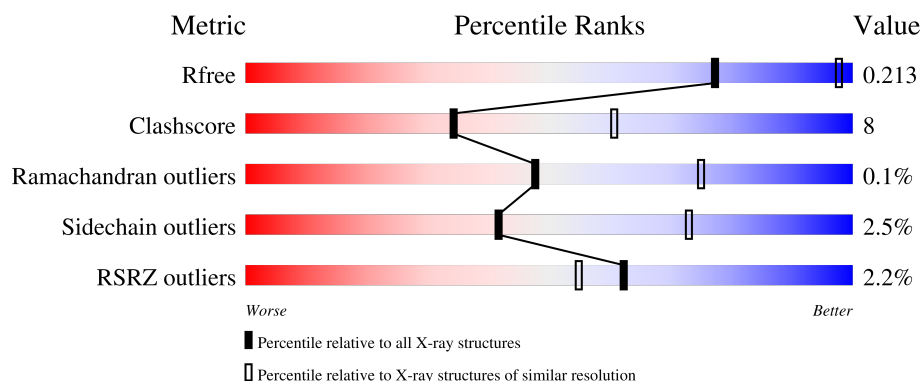
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

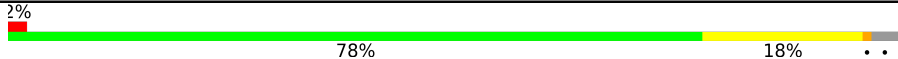
The reported resolution of this entry is 2.78 Å.

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	5248 (2.80-2.76)
Clashscore	190562	5693 (2.80-2.76)
Ramachandran outliers	187476	5590 (2.80-2.76)
Sidechain outliers	187428	5592 (2.80-2.76)
RSRZ outliers	180081	5251 (2.80-2.76)

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	723	
1	B	723	
1	C	723	
1	D	723	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 22501 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.

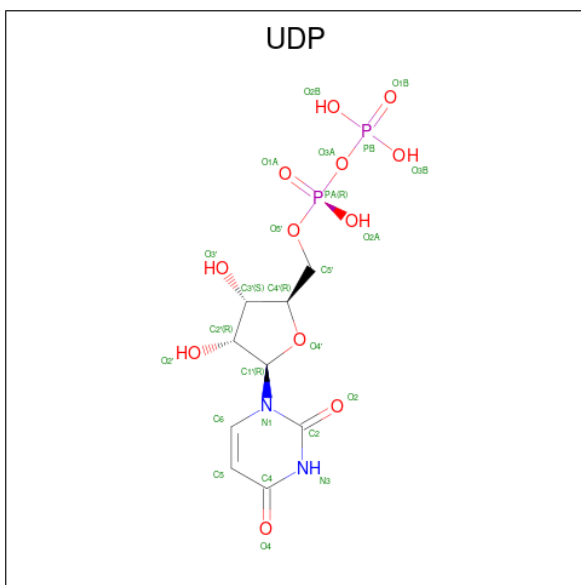
- Molecule 1 is a protein called UDP-N-acetylglucosamine--peptide N-acetylglucosaminyltransferase 110 kDa subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	701	Total	C	N	O	S	0	0	0
			5534	3511	968	1016	39			
1	B	701	Total	C	N	O	S	0	0	0
			5534	3511	968	1016	39			
1	C	701	Total	C	N	O	S	0	0	0
			5534	3511	968	1016	39			
1	D	701	Total	C	N	O	S	0	0	0
			5534	3511	968	1016	39			

There are 16 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
B	309	GLY	-	expression tag	UNP O15294
B	310	PRO	-	expression tag	UNP O15294
B	311	GLY	-	expression tag	UNP O15294
B	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
D	309	GLY	-	expression tag	UNP O15294
D	310	PRO	-	expression tag	UNP O15294
D	311	GLY	-	expression tag	UNP O15294
D	312	SER	-	expression tag	UNP O15294

- Molecule 2 is URIDINE-5'-DIPHOSPHATE (CCD ID: UDP) (formula: C₉H₁₄N₂O₁₂P₂).



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

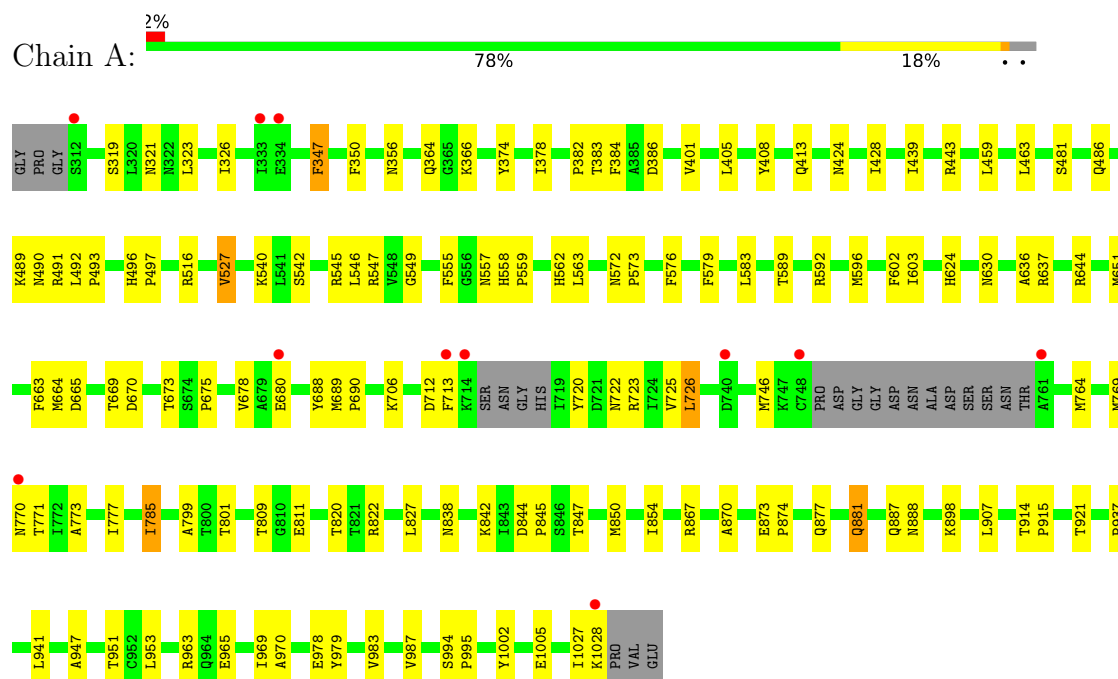
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	83	Total O 83 83	0	0
3	B	70	Total O 70 70	0	0
3	C	35	Total O 35 35	0	0
3	D	77	Total O 77 77	0	0

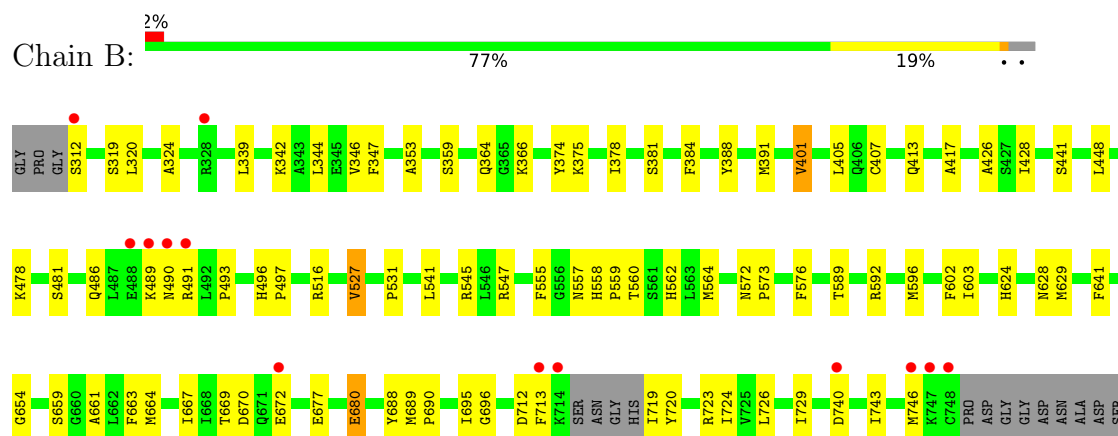
3 Residue-property plots

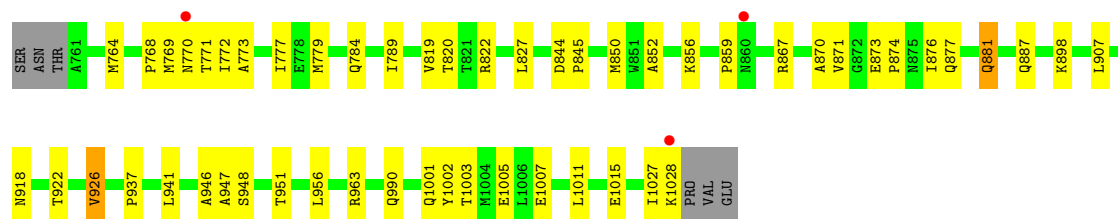
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: UDP-N-acetylglucosamine--peptide N-acetylglucosaminyltransferase 110 kDa subunit

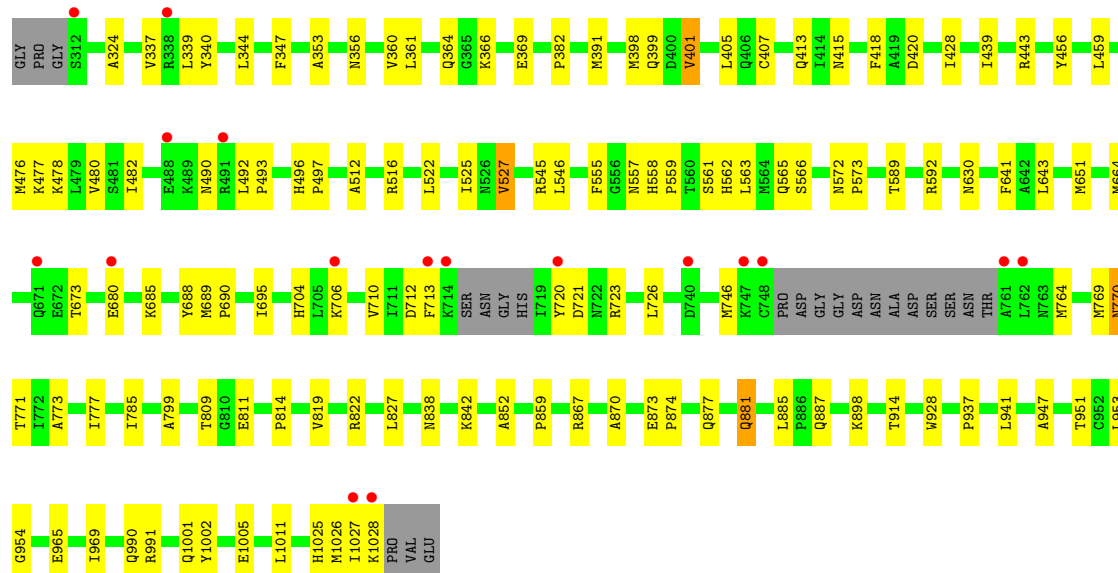
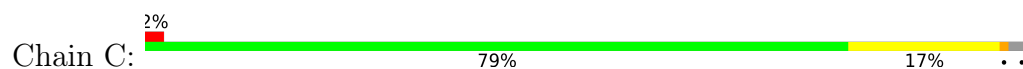


- Molecule 1: UDP-N-acetylglucosamine--peptide N-acetylglucosaminyltransferase 110 kDa subunit

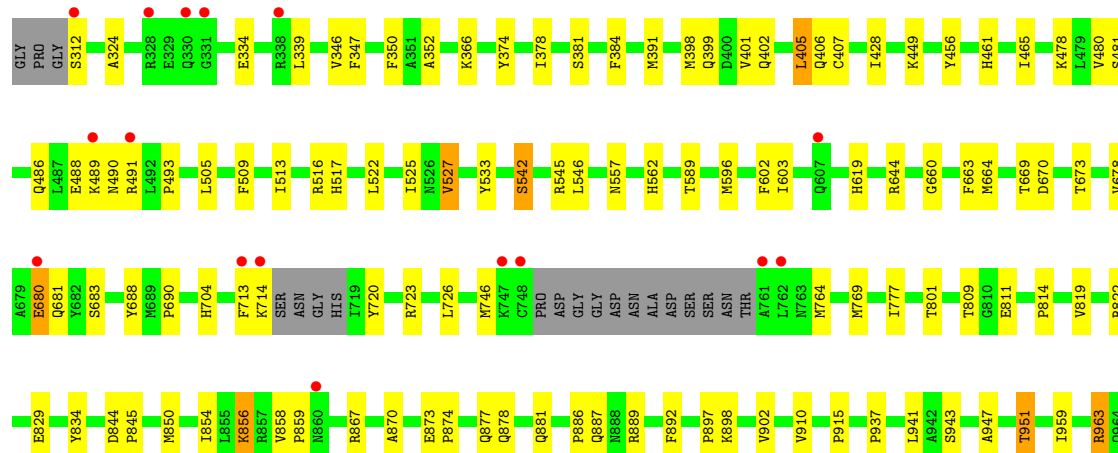
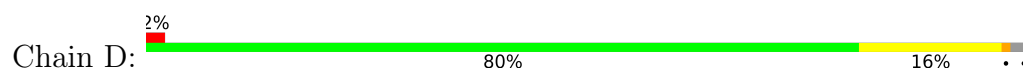




- Molecule 1: UDP-N-acetylglucosamine--peptide N-acetylglucosaminyltransferase 110 kDa subunit



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





4 Data and refinement statistics

Property	Value	Source
Space group	P 3 2 1	Depositor
Cell constants a, b, c, α , β , γ	273.40Å 273.40Å 142.80Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.58 – 2.78 48.58 – 2.78	Depositor EDS
% Data completeness (in resolution range)	97.8 (48.58-2.78) 98.0 (48.58-2.78)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.33 (at 2.77Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.1_357)	Depositor
R, R_{free}	0.185 , 0.218 0.181 , 0.213	Depositor DCC
R_{free} test set	2040 reflections (1.34%)	wwPDB-VP
Wilson B-factor (Å ²)	40.1	Xtriage
Anisotropy	0.511	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 52.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	0.034 for -h,-k,l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	22501	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.52% 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	0.31	0/5661	0.72	2/7676 (0.0%)
1	B	0.31	0/5661	0.73	2/7676 (0.0%)
1	C	0.29	0/5661	0.72	6/7676 (0.1%)
1	D	0.31	0/5661	0.71	4/7676 (0.1%)
All	All	0.31	0/22644	0.72	14/30704 (0.0%)

There are no bond length outliers.

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	347	PHE	CA-C-N	6.73	126.52	119.05
1	C	347	PHE	C-N-CA	6.73	126.52	119.05
1	D	449	LYS	CA-C-N	5.92	126.06	119.32
1	D	449	LYS	C-N-CA	5.92	126.06	119.32
1	B	347	PHE	CA-C-N	5.87	126.01	119.32
1	B	347	PHE	C-N-CA	5.87	126.01	119.32
1	A	347	PHE	CA-C-N	5.37	125.18	119.28
1	A	347	PHE	C-N-CA	5.37	125.18	119.28
1	C	914	THR	CA-C-N	5.32	124.83	119.19
1	C	914	THR	C-N-CA	5.32	124.83	119.19
1	D	347	PHE	CA-C-N	5.09	124.59	119.19
1	D	347	PHE	C-N-CA	5.09	124.59	119.19
1	C	415	ASN	CA-C-N	5.08	124.75	119.56
1	C	415	ASN	C-N-CA	5.08	124.75	119.56

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	5534	0	5512	88	0
1	B	5534	0	5512	95	0
1	C	5534	0	5512	86	0
1	D	5534	0	5512	72	0
2	A	25	0	11	0	0
2	B	25	0	11	0	0
2	C	25	0	11	1	0
2	D	25	0	11	0	0
3	A	83	0	0	2	0
3	B	70	0	0	2	0
3	C	35	0	0	0	0
3	D	77	0	0	1	0
All	All	22501	0	22092	335	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 8.

All (335) 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:547:ARG:H	1:B:624:HIS:HD2	1.11	0.93
1:C:770:ASN:H	1:C:770:ASN:HD22	1.15	0.91
1:A:547:ARG:H	1:A:624:HIS:HD2	1.09	0.90
1:D:850:MET:HE2	1:D:963:ARG:HE	1.37	0.88
1:A:850:MET:HE2	1:A:963:ARG:HE	1.41	0.85
1:B:850:MET:HE2	1:B:963:ARG:HE	1.53	0.74
1:A:321:ASN:HD21	1:A:356:ASN:HD22	1.37	0.72
1:B:547:ARG:H	1:B:624:HIS:CD2	2.01	0.71
1:A:545:ARG:HD3	1:A:573:PRO:O	1.91	0.71
1:C:770:ASN:HD22	1:C:770:ASN:N	1.87	0.71
1:B:719:ILE:N	3:B:196:HOH:O	2.23	0.71
1:A:562:HIS:ND1	1:A:898:LYS:HE3	2.06	0.70
1:D:562:HIS:ND1	1:D:898:LYS:HE3	2.06	0.70
1:A:547:ARG:H	1:A:624:HIS:CD2	2.00	0.70
1:B:629:MET:O	1:B:654:GLY:HA3	1.93	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:873:GLU:HB3	1:B:874:PRO:HD3	1.74	0.69
1:A:937:PRO:HG3	1:A:947:ALA:HB2	1.74	0.69
1:B:672:GLU:OE1	1:B:948:SER:HA	1.93	0.69
1:D:726:LEU:HD22	1:D:819:VAL:HG22	1.76	0.68
1:C:937:PRO:HG3	1:C:947:ALA:HB2	1.74	0.68
1:B:364:GLN:HB3	1:B:527:VAL:HG21	1.75	0.68
1:A:540:LYS:NZ	3:A:219:HOH:O	2.26	0.67
1:A:364:GLN:HB3	1:A:527:VAL:HG21	1.77	0.67
1:C:746:MET:HE3	1:C:764:MET:HB2	1.76	0.67
1:D:902:VAL:O	3:D:141:HOH:O	2.12	0.67
1:C:562:HIS:ND1	1:C:898:LYS:HE3	2.10	0.66
1:C:391:MET:HE3	1:C:407:CYS:SG	2.36	0.66
1:B:366:LYS:HE3	1:B:527:VAL:HG11	1.78	0.66
1:C:773:ALA:O	1:C:777:ILE:HG12	1.96	0.66
1:A:405:LEU:HD13	1:A:428:ILE:HG21	1.79	0.65
1:C:685:LYS:HD2	1:C:1025:HIS:CD2	2.32	0.65
1:B:867:ARG:HB3	1:B:870:ALA:HA	1.78	0.65
1:B:779:MET:HG3	1:B:784:GLN:HB2	1.79	0.65
1:A:583:LEU:HD22	1:A:637:ARG:HD3	1.79	0.65
1:D:947:ALA:O	1:D:951:THR:HG23	1.98	0.64
1:B:557:ASN:HB2	1:B:589:THR:HG21	1.80	0.64
1:A:773:ALA:O	1:A:777:ILE:HG12	1.96	0.64
1:D:690:PRO:HG2	1:D:1005:GLU:CD	2.23	0.64
1:C:852:ALA:HA	1:C:885:LEU:HD11	1.79	0.63
1:C:867:ARG:HB3	1:C:870:ALA:HA	1.80	0.63
1:B:951:THR:HG22	1:B:956:LEU:CD2	2.29	0.63
1:B:726:LEU:HD22	1:B:819:VAL:HG22	1.81	0.62
1:C:324:ALA:HB2	1:C:339:LEU:HB2	1.81	0.62
1:C:490:ASN:HA	1:C:516:ARG:HH22	1.64	0.62
1:C:641:PHE:CD1	1:C:664:MET:HE1	2.34	0.62
1:C:877:GLN:O	1:C:881:GLN:HG2	2.00	0.61
1:C:873:GLU:HB3	1:C:874:PRO:HD3	1.83	0.61
1:B:720:TYR:CB	1:B:723:ARG:HG2	2.31	0.61
1:C:726:LEU:HD22	1:C:819:VAL:HG22	1.82	0.60
1:A:877:GLN:O	1:A:881:GLN:HG2	2.00	0.60
1:D:873:GLU:HB3	1:D:874:PRO:HD3	1.84	0.60
1:A:490:ASN:HA	1:A:516:ARG:HH21	1.66	0.60
1:C:478:LYS:O	1:C:482:ILE:HG13	2.03	0.59
1:A:822:ARG:HB3	1:A:827:LEU:HB2	1.85	0.59
1:C:490:ASN:HA	1:C:516:ARG:NH2	2.18	0.58
1:B:545:ARG:HD3	1:B:573:PRO:O	2.04	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:877:GLN:O	1:B:881:GLN:HG2	2.04	0.57
1:D:480:VAL:HG22	1:D:505:LEU:HD23	1.86	0.57
1:D:850:MET:HE1	1:D:915:PRO:HG2	1.87	0.57
1:B:366:LYS:HE3	1:B:527:VAL:CG1	2.35	0.56
1:D:873:GLU:HG3	1:D:892:PHE:CD2	2.41	0.56
1:B:772:ILE:HG23	1:B:789:ILE:HD13	1.86	0.56
1:B:531:PRO:O	3:B:181:HOH:O	2.18	0.55
1:C:476:MET:O	1:C:480:VAL:HG23	2.06	0.55
1:A:413:GLN:HE21	1:C:413:GLN:HG3	1.72	0.55
1:C:947:ALA:O	1:C:951:THR:HG23	2.06	0.55
1:B:937:PRO:HG3	1:B:947:ALA:HB2	1.89	0.55
1:C:366:LYS:HE3	1:C:527:VAL:HG11	1.87	0.55
1:D:486:GLN:OE1	1:D:493:PRO:HA	2.07	0.55
1:D:720:TYR:CB	1:D:723:ARG:HG2	2.37	0.55
1:D:461:HIS:O	1:D:465:ILE:HG13	2.07	0.54
1:C:704:HIS:CD2	1:C:814:PRO:HG2	2.43	0.54
1:D:312:SER:HA	1:D:346:VAL:O	2.06	0.54
1:A:408:TYR:CZ	1:A:424:ASN:HB3	2.42	0.54
1:A:557:ASN:HB2	1:A:589:THR:HG21	1.89	0.54
1:B:720:TYR:HB3	1:B:723:ARG:HG2	1.89	0.54
1:D:867:ARG:HB3	1:D:870:ALA:HA	1.88	0.54
1:A:558:HIS:CG	1:A:559:PRO:HD2	2.43	0.54
1:C:405:LEU:HD13	1:C:428:ILE:HG21	1.90	0.53
1:A:770:ASN:CG	1:A:771:THR:H	2.15	0.53
1:A:366:LYS:HE3	1:A:527:VAL:HG11	1.90	0.53
1:B:562:HIS:ND1	1:B:898:LYS:HE3	2.23	0.53
1:D:509:PHE:O	1:D:513:ILE:HG13	2.08	0.53
1:A:603:ILE:N	1:A:603:ILE:HD12	2.23	0.53
1:A:688:TYR:CE2	1:A:1027:ILE:HD12	2.44	0.53
1:C:838:ASN:HB3	1:C:842:LYS:HD2	1.90	0.53
1:C:492:LEU:HD12	1:C:493:PRO:HD2	1.90	0.52
1:A:563:LEU:HD21	1:A:921:THR:HG23	1.91	0.52
1:C:690:PRO:HG2	1:C:1005:GLU:CD	2.35	0.52
1:A:770:ASN:H	1:A:773:ALA:HB3	1.74	0.52
1:A:850:MET:HE2	1:A:963:ARG:NE	2.19	0.52
1:D:660:GLY:HA2	1:D:683:SER:HB3	1.91	0.52
1:A:630:ASN:HB2	3:A:77:HOH:O	2.09	0.52
1:C:770:ASN:H	1:C:770:ASN:ND2	1.92	0.52
1:C:1001:GLN:O	1:C:1005:GLU:HG2	2.09	0.52
1:B:381:SER:O	1:B:384:PHE:HB2	2.10	0.52
1:B:486:GLN:OE1	1:B:493:PRO:HA	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:401:VAL:HG13	1:C:382:PRO:HG2	1.91	0.51
1:C:954:GLY:HA3	1:C:990:GLN:HE22	1.76	0.51
1:A:492:LEU:HD12	1:A:493:PRO:HD2	1.91	0.51
1:A:690:PRO:HG2	1:A:1005:GLU:CD	2.34	0.51
1:A:720:TYR:CD2	1:A:723:ARG:HG3	2.45	0.51
1:A:844:ASP:HB2	1:A:845:PRO:HD2	1.92	0.51
1:C:726:LEU:CD2	1:C:819:VAL:HG22	2.40	0.51
1:D:562:HIS:HD1	1:D:898:LYS:HE3	1.74	0.51
1:D:596:MET:HG2	1:D:602:PHE:CD2	2.46	0.51
1:B:820:THR:HG22	1:B:907:LEU:HD21	1.92	0.51
1:C:525:ILE:HB	1:C:643:LEU:HD21	1.92	0.51
1:C:898:LYS:HE2	2:C:1201:UDP:O2'	2.10	0.51
1:A:867:ARG:HB3	1:A:870:ALA:HA	1.93	0.51
1:B:695:ILE:HG13	1:B:696:GLY:N	2.25	0.51
1:D:723:ARG:HD2	1:D:822:ARG:HD2	1.93	0.51
1:D:663:PHE:CD1	1:D:664:MET:HE2	2.46	0.51
1:D:850:MET:HE2	1:D:963:ARG:NE	2.18	0.51
1:A:382:PRO:HG2	1:C:401:VAL:HG22	1.93	0.51
1:C:673:THR:CG2	1:C:941:LEU:HG	2.41	0.51
1:A:663:PHE:CD1	1:A:664:MET:HE2	2.46	0.50
1:B:558:HIS:CG	1:B:559:PRO:HD2	2.45	0.50
1:A:384:PHE:CE2	1:A:386:ASP:HB3	2.46	0.50
1:B:1003:THR:O	1:B:1007:GLU:HG3	2.12	0.50
1:D:374:TYR:O	1:D:378:ILE:HG12	2.12	0.50
1:B:374:TYR:O	1:B:378:ILE:HG12	2.12	0.50
1:B:770:ASN:CG	1:B:771:THR:H	2.19	0.50
1:B:723:ARG:HD2	1:B:822:ARG:HD2	1.94	0.50
1:D:1028:LYS:NZ	1:D:1028:LYS:HB3	2.27	0.50
1:C:785:ILE:HD12	1:C:799:ALA:CB	2.41	0.50
1:D:350:PHE:CE1	1:D:352:ALA:HB3	2.46	0.50
1:D:381:SER:O	1:D:384:PHE:HB2	2.12	0.49
1:A:675:PRO:HG2	1:A:678:VAL:HG22	1.94	0.49
1:B:555:PHE:O	1:B:592:ARG:HD3	2.12	0.49
1:C:720:TYR:CB	1:C:723:ARG:HG2	2.42	0.49
1:B:773:ALA:O	1:B:777:ILE:HG12	2.11	0.49
1:D:1001:GLN:O	1:D:1005:GLU:HG2	2.12	0.49
1:A:366:LYS:HE3	1:A:527:VAL:CG1	2.42	0.49
1:A:722:ASN:OD1	1:A:723:ARG:NH1	2.45	0.49
1:D:809:THR:OG1	1:D:811:GLU:HG3	2.13	0.48
1:B:661:ALA:HB1	1:B:663:PHE:CE2	2.48	0.48
1:C:688:TYR:CE2	1:C:1027:ILE:HD12	2.48	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:603:ILE:N	1:D:603:ILE:HD12	2.27	0.48
1:D:663:PHE:HD1	1:D:664:MET:HE2	1.77	0.48
1:D:713:PHE:N	1:D:713:PHE:CD1	2.81	0.48
1:B:541:LEU:HD12	1:B:541:LEU:H	1.79	0.48
1:D:456:TYR:CZ	1:D:478:LYS:HD3	2.49	0.48
1:B:312:SER:HA	1:B:346:VAL:O	2.13	0.48
1:C:723:ARG:HD2	1:C:822:ARG:HD2	1.94	0.48
1:A:941:LEU:C	1:A:941:LEU:HD23	2.38	0.48
1:B:712:ASP:O	1:B:769:MET:HE2	2.14	0.48
1:A:854:ILE:HD13	1:A:970:ALA:HB3	1.95	0.48
1:C:809:THR:OG1	1:C:811:GLU:HG3	2.14	0.48
1:C:713:PHE:HB3	1:C:769:MET:HE3	1.96	0.48
1:B:667:ILE:HG23	1:B:667:ILE:O	2.14	0.47
1:C:563:LEU:HD22	1:C:695:ILE:O	2.13	0.47
1:C:688:TYR:CD2	1:C:1027:ILE:HG23	2.49	0.47
1:B:690:PRO:HG2	1:B:1005:GLU:CD	2.39	0.47
1:C:557:ASN:HB2	1:C:589:THR:HG21	1.96	0.47
1:B:489:LYS:HD3	1:B:491:ARG:NH2	2.29	0.47
1:C:522:LEU:HA	1:C:525:ILE:HG12	1.96	0.47
1:D:673:THR:CG2	1:D:941:LEU:HG	2.44	0.47
1:A:873:GLU:HB3	1:A:874:PRO:HD3	1.95	0.47
1:B:344:LEU:HD21	1:B:353:ALA:HB3	1.96	0.47
1:B:746:MET:HE3	1:B:764:MET:HB2	1.97	0.47
1:C:689:MET:HE1	1:C:1005:GLU:HB2	1.97	0.47
1:C:712:ASP:O	1:C:769:MET:HE2	2.14	0.47
1:A:947:ALA:O	1:A:951:THR:HG23	2.14	0.47
1:C:713:PHE:HB3	1:C:769:MET:CE	2.44	0.47
1:A:347:PHE:O	1:A:350:PHE:HB2	2.15	0.47
1:A:374:TYR:O	1:A:378:ILE:HG12	2.15	0.47
1:D:324:ALA:HB2	1:D:339:LEU:HB2	1.96	0.47
1:C:689:MET:HE2	1:C:1002:TYR:CE1	2.50	0.47
1:C:558:HIS:CG	1:C:559:PRO:HD2	2.50	0.46
1:D:856:LYS:HZ2	1:D:856:LYS:HB2	1.80	0.46
1:A:555:PHE:O	1:A:592:ARG:HD3	2.15	0.46
1:A:651:MET:HG2	1:A:664:MET:HG2	1.98	0.46
1:B:426:ALA:HB2	1:B:441:SER:HB2	1.98	0.46
1:D:941:LEU:C	1:D:941:LEU:HD23	2.40	0.46
1:A:489:LYS:HD3	1:A:491:ARG:NH2	2.31	0.46
1:A:809:THR:OG1	1:A:811:GLU:HG3	2.16	0.46
1:B:490:ASN:HA	1:B:516:ARG:HH21	1.81	0.46
1:C:953:LEU:O	1:C:990:GLN:NE2	2.48	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:628:ASN:HB2	1:B:641:PHE:CE1	2.51	0.46
1:A:770:ASN:CG	1:A:771:THR:N	2.74	0.46
1:A:965:GLU:O	1:A:969:ILE:HG13	2.16	0.46
1:C:546:LEU:HD22	1:C:1011:LEU:HD23	1.98	0.46
1:D:877:GLN:O	1:D:881:GLN:HG2	2.15	0.46
1:B:941:LEU:HD23	1:B:941:LEU:C	2.41	0.46
1:C:965:GLU:O	1:C:969:ILE:HG13	2.15	0.46
1:D:680:GLU:H	1:D:680:GLU:HG3	1.51	0.46
1:A:706:LYS:HE2	1:A:706:LYS:HB3	1.75	0.45
1:C:704:HIS:NE2	1:C:814:PRO:HG2	2.30	0.45
1:A:720:TYR:HB2	1:A:723:ARG:HG2	1.99	0.45
1:D:596:MET:HG2	1:D:602:PHE:CG	2.51	0.45
1:A:630:ASN:HB3	1:A:636:ALA:HB2	1.99	0.45
1:A:725:VAL:C	1:A:726:LEU:HD23	2.41	0.45
1:C:398:MET:O	1:C:399:GLN:HB2	2.17	0.45
1:C:418:PHE:CZ	1:C:420:ASP:HB2	2.51	0.45
1:D:490:ASN:HA	1:D:516:ARG:HH21	1.82	0.45
1:D:723:ARG:NE	1:D:829:GLU:O	2.41	0.45
1:A:785:ILE:HD12	1:A:799:ALA:CB	2.47	0.45
1:C:344:LEU:HD21	1:C:353:ALA:HB3	1.97	0.45
1:C:822:ARG:HB3	1:C:827:LEU:HB2	1.98	0.45
1:A:669:THR:OG1	1:A:670:ASP:N	2.50	0.45
1:A:953:LEU:HD21	1:A:987:VAL:HG13	1.98	0.45
1:D:720:TYR:CD2	1:D:723:ARG:HG3	2.52	0.45
1:D:713:PHE:HB3	1:D:769:MET:CE	2.47	0.45
1:B:405:LEU:HD13	1:B:428:ILE:HG21	1.99	0.45
1:B:918:ASN:HB3	1:B:946:ALA:HB2	1.99	0.45
1:C:477:LYS:HB2	1:C:477:LYS:HE3	1.72	0.45
1:D:366:LYS:HE3	1:D:527:VAL:HG21	1.98	0.45
1:D:746:MET:CE	1:D:764:MET:HB2	2.47	0.45
1:B:688:TYR:CE2	1:B:1027:ILE:HD12	2.52	0.44
1:C:439:ILE:O	1:C:443:ARG:HG3	2.16	0.44
1:C:555:PHE:O	1:C:592:ARG:HD3	2.17	0.44
1:D:669:THR:OG1	1:D:670:ASP:N	2.50	0.44
1:A:486:GLN:OE1	1:A:493:PRO:HA	2.16	0.44
1:B:391:MET:HE3	1:B:407:CYS:SG	2.57	0.44
1:B:603:ILE:N	1:B:603:ILE:HD12	2.32	0.44
1:C:439:ILE:HG23	1:C:459:LEU:HD11	1.99	0.44
1:B:713:PHE:CD1	1:B:713:PHE:N	2.85	0.44
1:C:512:ALA:O	1:C:516:ARG:HG2	2.18	0.44
1:B:388:TYR:O	1:B:407:CYS:HB3	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:417:ALA:HA	1:B:448:LEU:HD22	1.99	0.44
1:A:439:ILE:O	1:A:443:ARG:HG3	2.17	0.44
1:D:489:LYS:HD2	1:D:491:ARG:HH12	1.83	0.44
1:D:546:LEU:HD22	1:D:1011:LEU:HD23	1.99	0.44
1:B:572:ASN:HA	1:B:573:PRO:HD3	1.89	0.44
1:B:496:HIS:CG	1:B:497:PRO:HD2	2.52	0.44
1:A:688:TYR:CD2	1:A:1027:ILE:HD12	2.53	0.44
1:C:337:VAL:HG22	1:C:360:VAL:HG21	1.98	0.44
1:B:689:MET:HE2	1:B:1002:TYR:CE1	2.52	0.43
1:B:951:THR:HG22	1:B:956:LEU:HD23	2.00	0.43
1:C:324:ALA:HB2	1:C:339:LEU:CB	2.45	0.43
1:D:391:MET:HE3	1:D:407:CYS:SG	2.58	0.43
1:A:673:THR:HG22	1:A:941:LEU:HG	1.99	0.43
1:A:979:TYR:O	1:A:983:VAL:HG23	2.18	0.43
1:D:405:LEU:HD12	1:D:405:LEU:HA	1.78	0.43
1:D:951:THR:HG22	1:D:959:ILE:HD11	2.01	0.43
1:B:663:PHE:CD1	1:B:664:MET:HE2	2.54	0.43
1:B:770:ASN:ND2	1:B:771:THR:H	2.16	0.43
1:B:576:PHE:CE2	1:B:1007:GLU:HB3	2.53	0.43
1:B:720:TYR:CG	1:B:723:ARG:HG2	2.53	0.43
1:C:364:GLN:HB3	1:C:527:VAL:HG21	2.00	0.43
1:A:689:MET:HE1	1:A:1005:GLU:HB2	2.01	0.43
1:C:720:TYR:CD2	1:C:723:ARG:HG3	2.54	0.43
1:C:785:ILE:HD12	1:C:799:ALA:HB1	2.01	0.43
1:A:838:ASN:HB3	1:A:842:LYS:HD2	1.99	0.43
1:A:914:THR:HA	1:A:915:PRO:HD3	1.81	0.43
1:A:994:SER:HB2	1:A:995:PRO:HD2	1.99	0.43
1:B:659:SER:C	1:B:661:ALA:H	2.26	0.43
1:C:456:TYR:CZ	1:C:478:LYS:HD3	2.54	0.43
1:A:847:THR:HA	1:A:850:MET:HE3	2.01	0.43
1:B:688:TYR:CD2	1:B:1027:ILE:HD12	2.54	0.43
1:D:542:SER:HB2	1:D:545:ARG:O	2.18	0.43
1:A:746:MET:HE2	1:A:764:MET:HB2	2.00	0.43
1:B:669:THR:OG1	1:B:670:ASP:N	2.51	0.43
1:B:596:MET:HG2	1:B:602:PHE:CD2	2.53	0.42
1:B:596:MET:HG2	1:B:602:PHE:CG	2.54	0.42
1:C:572:ASN:HA	1:C:573:PRO:HD3	1.86	0.42
1:D:398:MET:O	1:D:399:GLN:HB2	2.19	0.42
1:D:897:PRO:O	1:D:898:LYS:C	2.62	0.42
1:B:951:THR:HG22	1:B:956:LEU:HD21	2.00	0.42
1:C:641:PHE:HB3	1:C:664:MET:CE	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:689:MET:HE2	1:A:1002:TYR:CE1	2.54	0.42
1:C:770:ASN:N	1:C:770:ASN:ND2	2.59	0.42
1:A:854:ILE:HD13	1:A:970:ALA:CB	2.48	0.42
1:A:459:LEU:O	1:A:463:LEU:HG	2.20	0.42
1:A:644:ARG:HH22	1:A:665:ASP:CG	2.27	0.42
1:A:673:THR:CG2	1:A:941:LEU:HG	2.49	0.42
1:A:720:TYR:CB	1:A:723:ARG:HG2	2.48	0.42
1:B:822:ARG:HB3	1:B:827:LEU:HB2	2.02	0.42
1:B:547:ARG:N	1:B:624:HIS:HD2	1.94	0.42
1:B:413:GLN:HG3	1:C:413:GLN:HE21	1.84	0.42
1:B:844:ASP:HB2	1:B:845:PRO:HD2	2.02	0.42
1:D:937:PRO:HA	1:D:943:SER:O	2.19	0.42
1:A:820:THR:HG22	1:A:907:LEU:HD21	2.02	0.42
1:D:886:PRO:HD2	1:D:889:ARG:HG3	2.02	0.42
1:B:740:ASP:O	1:B:768:PRO:HG3	2.20	0.42
1:A:323:LEU:HD23	1:A:326:ILE:HD12	2.01	0.41
1:B:560:THR:O	1:B:564:MET:HG3	2.20	0.41
1:C:651:MET:HE3	1:C:651:MET:HB3	2.00	0.41
1:D:858:VAL:HA	1:D:859:PRO:HD2	1.88	0.41
1:A:572:ASN:HA	1:A:573:PRO:HD3	1.89	0.41
1:C:928:TRP:O	1:C:991:ARG:NH1	2.50	0.41
1:A:663:PHE:HD1	1:A:664:MET:HE2	1.84	0.41
1:D:522:LEU:HA	1:D:525:ILE:HG12	2.02	0.41
1:D:678:VAL:HB	1:D:681:GLN:NE2	2.35	0.41
1:A:547:ARG:HG2	1:A:579:PHE:HE2	1.85	0.41
1:A:592:ARG:O	1:A:596:MET:HG3	2.20	0.41
1:A:602:PHE:C	1:A:603:ILE:HD12	2.45	0.41
1:C:706:LYS:HE3	1:C:706:LYS:HB2	1.76	0.41
1:D:965:GLU:O	1:D:969:ILE:HG13	2.20	0.41
1:D:704:HIS:NE2	1:D:814:PRO:HG2	2.36	0.41
1:D:714:LYS:HE2	1:D:714:LYS:HB2	1.89	0.41
1:A:712:ASP:O	1:A:769:MET:HE2	2.21	0.41
1:A:713:PHE:CD1	1:A:713:PHE:N	2.89	0.41
1:B:641:PHE:HB3	1:B:664:MET:HE1	2.02	0.41
1:B:852:ALA:O	1:B:856:LYS:HG3	2.20	0.41
1:B:937:PRO:HG3	1:B:947:ALA:CB	2.50	0.41
1:B:320:LEU:HD13	1:B:342:LYS:HB3	2.02	0.41
1:B:324:ALA:HB2	1:B:339:LEU:HB2	2.01	0.41
1:B:1001:GLN:O	1:B:1005:GLU:HG2	2.21	0.41
1:A:384:PHE:HE2	1:A:386:ASP:HB3	1.85	0.41
1:A:496:HIS:CG	1:A:497:PRO:HD2	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:405:LEU:HD12	1:B:405:LEU:HA	1.93	0.41
1:B:413:GLN:HG3	1:C:413:GLN:NE2	2.36	0.41
1:B:413:GLN:CD	1:C:413:GLN:NE2	2.78	0.41
1:B:672:GLU:CD	1:B:948:SER:HA	2.46	0.41
1:C:561:SER:O	1:C:565:GLN:HB3	2.20	0.41
1:A:746:MET:CE	1:A:764:MET:HB2	2.51	0.41
1:B:871:VAL:O	1:B:874:PRO:HD2	2.21	0.41
1:C:340:TYR:CZ	1:C:356:ASN:HB3	2.56	0.41
1:C:361:LEU:HD13	1:C:369:GLU:HG2	2.03	0.41
1:C:545:ARG:HD3	1:C:573:PRO:O	2.20	0.41
1:C:710:VAL:HG11	1:C:721:ASP:HA	2.02	0.41
1:D:513:ILE:O	1:D:517:HIS:HD2	2.04	0.41
1:D:533:TYR:HB3	1:D:619:HIS:CD2	2.56	0.41
1:D:644:ARG:HH11	1:D:663:PHE:HA	1.86	0.41
1:D:834:TYR:HA	1:D:910:VAL:O	2.20	0.41
1:A:546:LEU:HD23	1:A:576:PHE:CE1	2.56	0.41
1:D:688:TYR:CE2	1:D:1027:ILE:HD12	2.55	0.41
1:D:405:LEU:HD13	1:D:428:ILE:HG21	2.03	0.40
1:D:844:ASP:HB2	1:D:845:PRO:HD2	2.03	0.40
1:B:680:GLU:H	1:B:680:GLU:HG3	1.50	0.40
1:D:746:MET:HE2	1:D:764:MET:HB2	2.03	0.40
1:B:724:ILE:HD13	1:B:777:ILE:HD11	2.03	0.40
1:B:726:LEU:CD2	1:B:819:VAL:HG22	2.51	0.40
1:B:922:THR:O	1:B:926:VAL:HG22	2.22	0.40
1:D:854:ILE:HD13	1:D:970:ALA:HB3	2.03	0.40
1:B:729:ILE:HD13	1:B:729:ILE:HA	1.87	0.40
1:B:1011:LEU:O	1:B:1015:GLU:HG2	2.22	0.40
1:C:366:LYS:HE3	1:C:527:VAL:CG1	2.52	0.40
1:C:496:HIS:CG	1:C:497:PRO:HD2	2.57	0.40
1:D:557:ASN:HB2	1:D:589:THR:HG21	2.03	0.40
1:A:549:GLY:HA2	1:A:579:PHE:O	2.21	0.40
1:B:478:LYS:HA	1:B:481:SER:HB3	2.03	0.40
1:B:576:PHE:CZ	1:B:1007:GLU:HB3	2.57	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	695/723 (96%)	673 (97%)	22 (3%)	0	100	100
1	B	695/723 (96%)	677 (97%)	17 (2%)	1 (0%)	48	75
1	C	695/723 (96%)	669 (96%)	25 (4%)	1 (0%)	48	75
1	D	695/723 (96%)	671 (96%)	24 (4%)	0	100	100
All	All	2780/2892 (96%)	2690 (97%)	88 (3%)	2 (0%)	48	75

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	859	PRO
1	B	859	PRO

5.3.2 Protein sidechains [i](#)

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	602/618 (97%)	587 (98%)	15 (2%)	42	73
1	B	602/618 (97%)	588 (98%)	14 (2%)	44	75
1	C	602/618 (97%)	591 (98%)	11 (2%)	51	79
1	D	602/618 (97%)	583 (97%)	19 (3%)	34	67
All	All	2408/2472 (97%)	2349 (98%)	59 (2%)	42	73

All (59) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	319	SER
1	A	383	THR
1	A	401	VAL
1	A	481	SER
1	A	527	VAL
1	A	542	SER
1	A	680	GLU
1	A	726	LEU
1	A	785	ILE
1	A	801	THR
1	A	881	GLN
1	A	887	GLN
1	A	888	ASN
1	A	978	GLU
1	A	1028	LYS
1	B	319	SER
1	B	359	SER
1	B	375	LYS
1	B	401	VAL
1	B	527	VAL
1	B	677	GLU
1	B	680	GLU
1	B	743	ILE
1	B	876	ILE
1	B	881	GLN
1	B	887	GLN
1	B	926	VAL
1	B	990	GLN
1	B	1028	LYS
1	C	401	VAL
1	C	527	VAL
1	C	566	SER
1	C	630	ASN
1	C	680	GLU
1	C	770	ASN
1	C	771	THR
1	C	881	GLN
1	C	887	GLN
1	C	1026	MET
1	C	1028	LYS
1	D	334	GLU
1	D	401	VAL
1	D	402	GLN

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Mol	Chain	Res	Type
1	D	405	LEU
1	D	406	GLN
1	D	481	SER
1	D	488	GLU
1	D	527	VAL
1	D	542	SER
1	D	680	GLU
1	D	777	ILE
1	D	801	THR
1	D	856	LYS
1	D	878	GLN
1	D	887	GLN
1	D	951	THR
1	D	963	ARG
1	D	978	GLU
1	D	1028	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (57) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	356	ASN
1	A	363	GLN
1	A	368	GLN
1	A	413	GLN
1	A	434	ASN
1	A	517	HIS
1	A	529	HIS
1	A	620	GLN
1	A	624	HIS
1	A	681	GLN
1	A	727	ASN
1	A	770	ASN
1	A	860	ASN
1	A	888	ASN
1	A	906	GLN
1	A	990	GLN
1	B	321	ASN
1	B	362	GLN
1	B	402	GLN
1	B	413	GLN
1	B	434	ASN
1	B	517	HIS

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Mol	Chain	Res	Type
1	B	529	HIS
1	B	620	GLN
1	B	624	HIS
1	B	691	HIS
1	B	763	ASN
1	B	770	ASN
1	B	784	GLN
1	B	881	GLN
1	B	1012	GLN
1	C	321	ASN
1	C	356	ASN
1	C	402	GLN
1	C	413	GLN
1	C	517	HIS
1	C	620	GLN
1	C	681	GLN
1	C	727	ASN
1	C	763	ASN
1	C	770	ASN
1	C	784	GLN
1	C	881	GLN
1	C	906	GLN
1	C	990	GLN
1	C	1016	HIS
1	C	1025	HIS
1	D	316	HIS
1	D	321	ASN
1	D	356	ASN
1	D	363	GLN
1	D	517	HIS
1	D	620	GLN
1	D	681	GLN
1	D	784	GLN
1	D	881	GLN
1	D	1025	HIS

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 ⓘ

4 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	A	1201	-	25,26,26	1.05	1 (4%)	38,40,40	1.54	5 (13%)
2	UDP	C	1201	-	25,26,26	1.02	1 (4%)	38,40,40	1.66	6 (15%)
2	UDP	D	1201	-	25,26,26	1.02	1 (4%)	38,40,40	1.50	4 (10%)
2	UDP	B	1201	-	25,26,26	1.01	1 (4%)	38,40,40	1.60	6 (15%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	UDP	A	1201	-	-	3/16/32/32	0/2/2/2
2	UDP	C	1201	-	-	5/16/32/32	0/2/2/2
2	UDP	D	1201	-	-	5/16/32/32	0/2/2/2
2	UDP	B	1201	-	-	5/16/32/32	0/2/2/2

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	1201	UDP	PA-O3A	2.10	1.61	1.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	1201	UDP	C6-C5	2.10	1.39	1.35
2	B	1201	UDP	C6-C5	2.06	1.39	1.35
2	A	1201	UDP	PA-O3A	2.03	1.61	1.59

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1201	UDP	C4-N3-C2	-5.55	119.72	126.61
2	C	1201	UDP	C4-N3-C2	-5.51	119.77	126.61
2	A	1201	UDP	C4-N3-C2	-5.46	119.84	126.61
2	D	1201	UDP	C4-N3-C2	-5.27	120.07	126.61
2	C	1201	UDP	N3-C2-N1	4.35	120.56	114.89
2	B	1201	UDP	N3-C2-N1	4.30	120.48	114.89
2	D	1201	UDP	N3-C2-N1	4.15	120.30	114.89
2	A	1201	UDP	C5-C4-N3	3.89	120.25	114.80
2	A	1201	UDP	N3-C2-N1	3.73	119.75	114.89
2	B	1201	UDP	C5-C4-N3	3.59	119.83	114.80
2	D	1201	UDP	C5-C4-N3	3.56	119.79	114.80
2	C	1201	UDP	C5-C4-N3	3.32	119.45	114.80
2	B	1201	UDP	O4-C4-C5	-2.74	120.44	125.16
2	C	1201	UDP	O4-C4-C5	-2.62	120.64	125.16
2	A	1201	UDP	O4-C4-C5	-2.60	120.68	125.16
2	B	1201	UDP	O2-C2-N1	-2.50	119.54	122.80
2	C	1201	UDP	O4'-C1'-N1	2.45	113.92	108.36
2	B	1201	UDP	O2A-PA-O3A	2.24	113.33	107.27
2	D	1201	UDP	O4-C4-C5	-2.23	121.31	125.16
2	C	1201	UDP	C5'-C4'-C3'	-2.18	107.38	115.21
2	A	1201	UDP	O2-C2-N1	-2.06	120.12	122.80

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	1201	UDP	O4'-C4'-C5'-O5'
2	B	1201	UDP	C5'-O5'-PA-O1A
2	C	1201	UDP	C5'-O5'-PA-O1A
2	C	1201	UDP	C5'-O5'-PA-O2A
2	D	1201	UDP	C3'-C4'-C5'-O5'
2	D	1201	UDP	O4'-C4'-C5'-O5'
2	B	1201	UDP	C3'-C4'-C5'-O5'
2	A	1201	UDP	O4'-C4'-C5'-O5'
2	C	1201	UDP	C3'-C4'-C5'-O5'

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Mol	Chain	Res	Type	Atoms
2	A	1201	UDP	C3'-C4'-C5'-O5'
2	C	1201	UDP	O4'-C4'-C5'-O5'
2	A	1201	UDP	C5'-O5'-PA-O1A
2	B	1201	UDP	C5'-O5'-PA-O2A
2	B	1201	UDP	C5'-O5'-PA-O3A
2	C	1201	UDP	C5'-O5'-PA-O3A
2	D	1201	UDP	C5'-O5'-PA-O1A
2	D	1201	UDP	C5'-O5'-PA-O2A
2	D	1201	UDP	C5'-O5'-PA-O3A

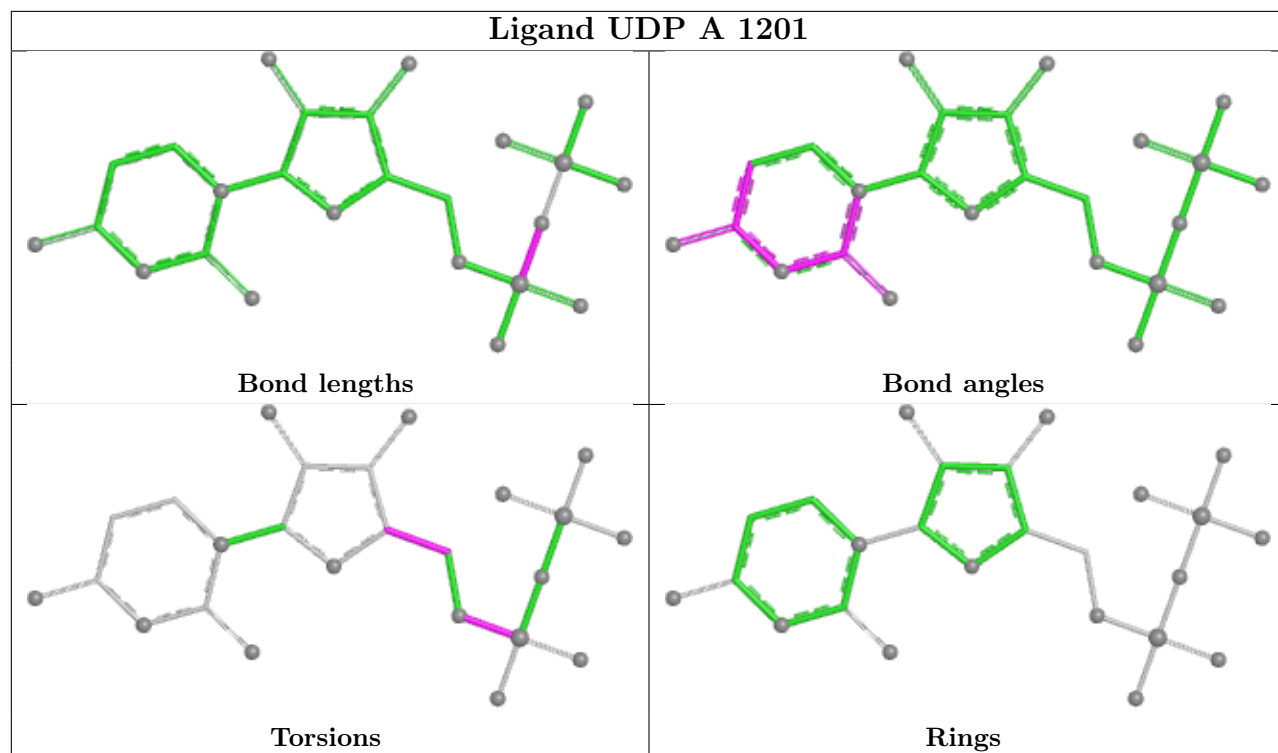
There are no ring outliers.

1 monomer is involved in 1 short contact:

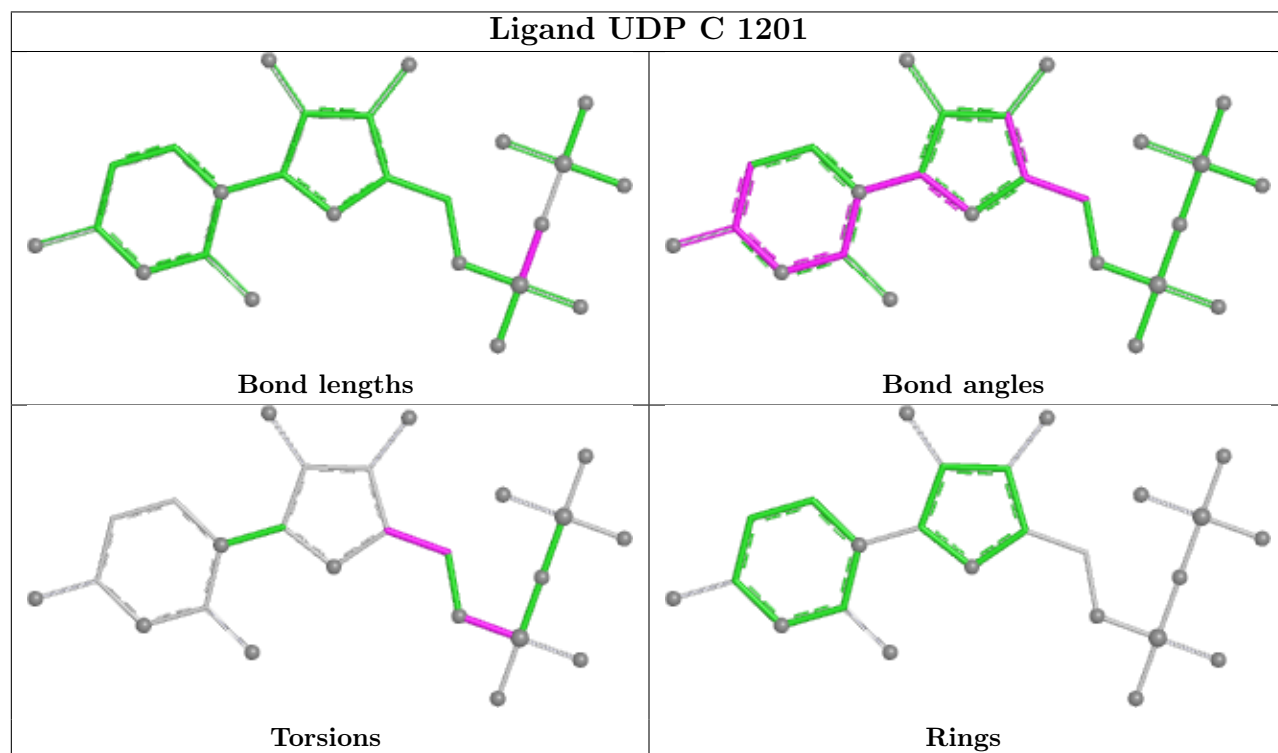
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	1201	UDP	1	0

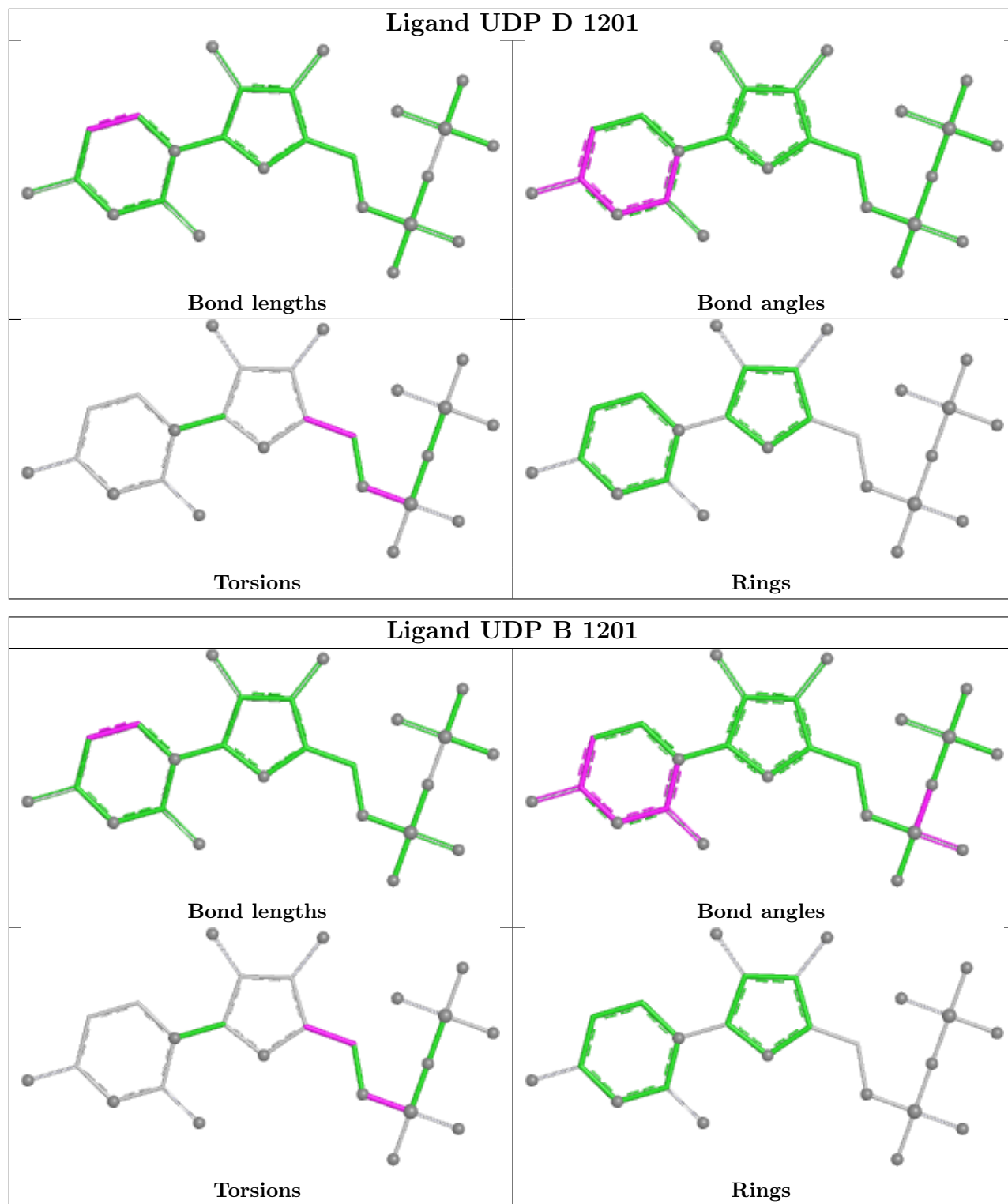
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

Ligand UDP A 1201



Ligand UDP C 1201





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	701/723 (96%)	-0.41	11 (1%) 70 63	27, 42, 73, 113	0
1	B	701/723 (96%)	-0.34	16 (2%) 61 53	22, 45, 79, 120	0
1	C	701/723 (96%)	-0.09	17 (2%) 59 52	27, 56, 87, 123	0
1	D	701/723 (96%)	-0.37	18 (2%) 57 49	24, 42, 74, 119	0
All	All	2804/2892 (96%)	-0.30	62 (2%) 62 55	22, 46, 80, 123	0

All (62) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	714	LYS	6.3
1	C	714	LYS	6.2
1	D	761	ALA	5.4
1	A	714	LYS	5.1
1	B	714	LYS	4.9
1	B	312	SER	4.4
1	A	1028	LYS	4.4
1	C	761	ALA	4.2
1	B	1028	LYS	4.1
1	A	770	ASN	4.1
1	B	748	CYS	4.0
1	A	713	PHE	4.0
1	C	713	PHE	3.9
1	A	312	SER	3.9
1	D	713	PHE	3.8
1	C	747	LYS	3.6
1	D	338	ARG	3.6
1	D	491	ARG	3.6
1	B	672	GLU	3.5
1	C	488	GLU	3.4
1	C	312	SER	3.4

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Mol	Chain	Res	Type	RSRZ
1	C	1028	LYS	3.3
1	C	680	GLU	3.3
1	C	762	LEU	3.3
1	D	748	CYS	3.3
1	B	747	LYS	3.2
1	D	330	GLN	3.2
1	C	740	ASP	3.2
1	A	761	ALA	3.1
1	C	1027	ILE	3.1
1	D	860	ASN	3.1
1	B	488	GLU	3.1
1	C	491	ARG	3.1
1	D	747	LYS	3.1
1	D	1028	LYS	3.1
1	D	680	GLU	3.0
1	A	740	ASP	3.0
1	B	770	ASN	2.9
1	C	748	CYS	2.9
1	B	491	ARG	2.8
1	B	860	ASN	2.8
1	B	740	ASP	2.8
1	C	671	GLN	2.7
1	D	1027	ILE	2.5
1	B	489	LYS	2.4
1	B	713	PHE	2.4
1	B	490	ASN	2.4
1	A	334	GLU	2.4
1	D	762	LEU	2.3
1	A	333	ILE	2.3
1	C	338	ARG	2.3
1	C	706	LYS	2.2
1	D	328	ARG	2.2
1	D	331	GLY	2.2
1	B	328	ARG	2.2
1	D	607	GLN	2.1
1	A	680	GLU	2.1
1	D	489	LYS	2.1
1	A	748	CYS	2.1
1	B	746	MET	2.0
1	C	720	TYR	2.0
1	D	312	SER	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](#)

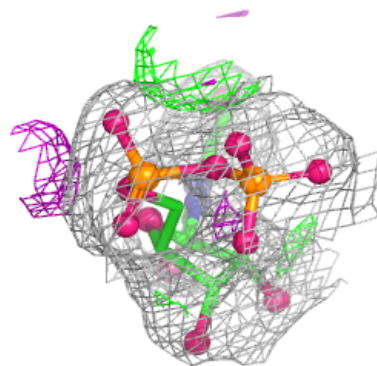
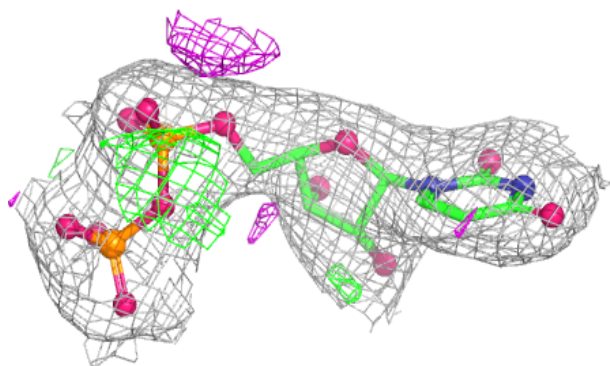
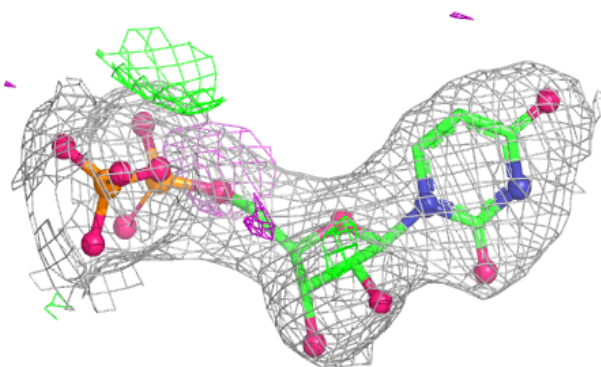
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	C	1201	25/25	0.97	0.07	29,47,59,62	0
2	UDP	B	1201	25/25	0.98	0.05	26,36,42,45	0
2	UDP	A	1201	25/25	0.98	0.05	24,33,38,49	0
2	UDP	D	1201	25/25	0.98	0.06	24,31,36,39	0

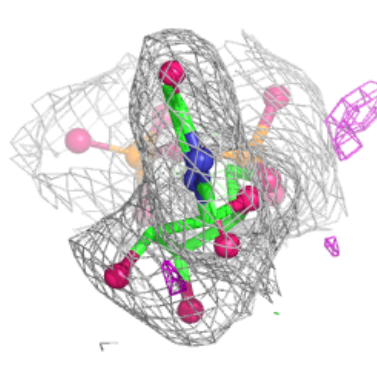
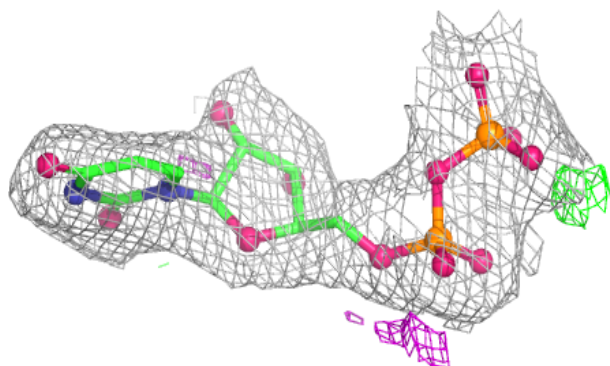
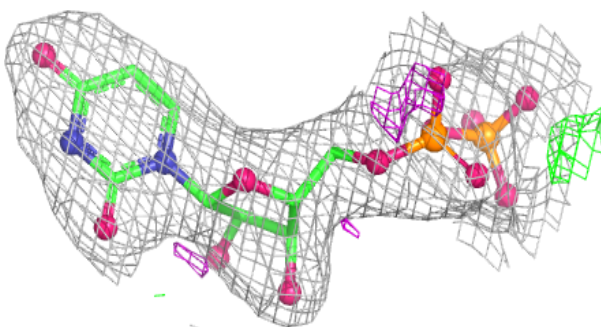
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around UDP C 1201:

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

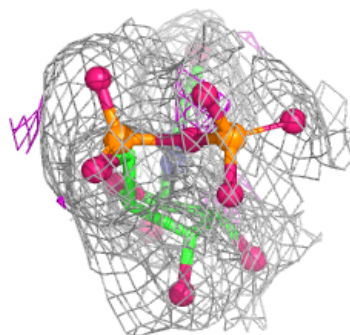
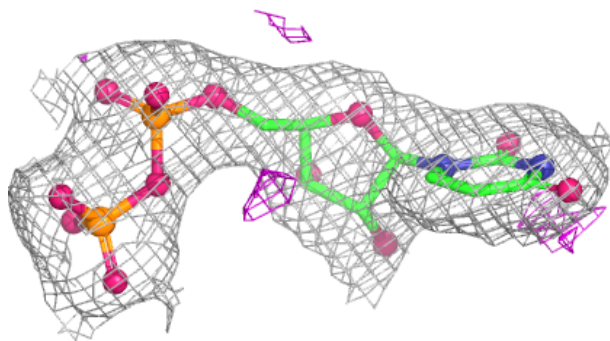
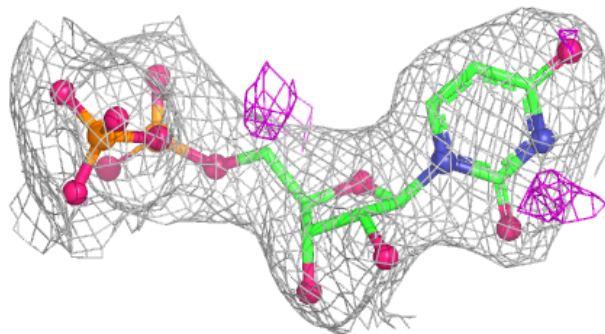
**Electron density around UDP B 1201:**

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

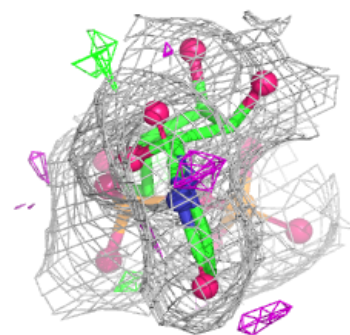
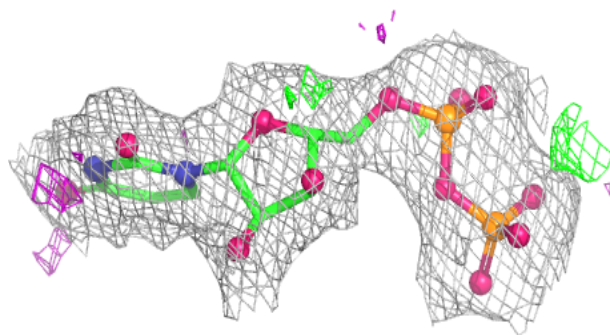
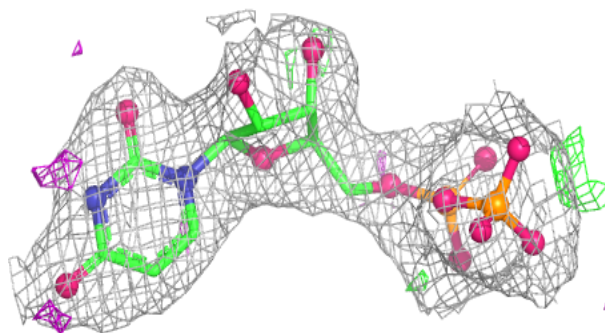


Electron density around UDP A 1201:

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

**Electron density around UDP D 1201:**

$2mF_o-DF_c$ (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.