



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 02:13 AM UTC

PDB ID : 8FE7 / pdb_00008fe7
Title : Crystal structure of human O-GlcNAc transferase (OGT) in complex with an exosite-binding peptide (SMG9) and UDP-GlcNAc
Authors : Xie, J.; Jiang, J.
Deposited on : 2022-12-05
Resolution : 2.98 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

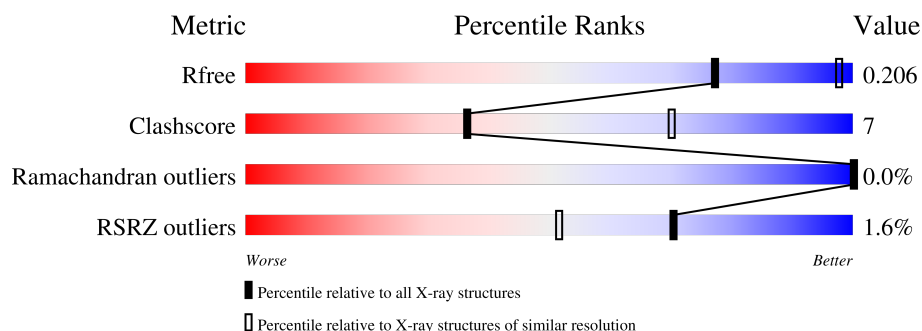
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.98 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	3580 (3.00-2.96)
Clashscore	190562	3904 (3.00-2.96)
Ramachandran outliers	187476	3761 (3.00-2.96)
RSRZ outliers	180081	3579 (3.00-2.96)

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	2% 79% 18% .
1	C	723	% 80% 17% .
1	E	723	% 83% 14% .
1	G	723	2% 79% 18% .
2	B	12	8% 67% 25% 8%
2	D	12	8% 83% 8% 8%

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Mol	Chain	Length	Quality of chain
2	F	12	8% 83% 8% 8%
2	H	12	83% 8% 8%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 22788 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	705	Total	C	N	O	S	0	0	0
			5563	3528	975	1022	38			
1	C	705	Total	C	N	O	S	0	0	0
			5563	3528	975	1022	38			
1	E	705	Total	C	N	O	S	0	0	0
			5563	3528	975	1022	38			
1	G	705	Total	C	N	O	S	0	0	0
			5563	3528	975	1022	38			

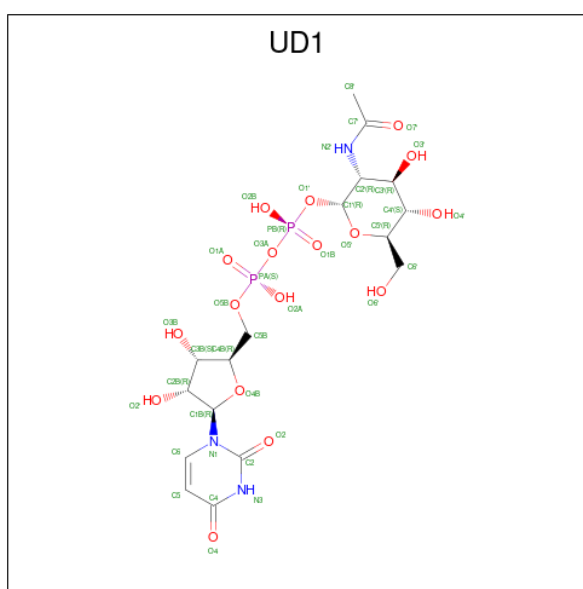
There are 16 discrepancies between the modelled and reference sequences:

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
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
E	309	GLY	-	expression tag	UNP O15294
E	310	PRO	-	expression tag	UNP O15294
E	311	GLY	-	expression tag	UNP O15294
E	312	SER	-	expression tag	UNP O15294
G	309	GLY	-	expression tag	UNP O15294
G	310	PRO	-	expression tag	UNP O15294
G	311	GLY	-	expression tag	UNP O15294
G	312	SER	-	expression tag	UNP O15294

- Molecule 2 is a protein called Nonsense-mediated mRNA decay factor SMG9.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	11	Total	C	N	O	0	0	0
			95	60	18	17			
2	D	11	Total	C	N	O	0	0	0
			95	60	18	17			
2	F	11	Total	C	N	O	0	0	0
			95	60	18	17			
2	H	11	Total	C	N	O	0	0	0
			95	60	18	17			

- Molecule 3 is URIDINE-DIPHOSPHATE-N-ACETYLGLUCOSAMINE (CCD ID: UD1) (formula: $C_{17}H_{27}N_3O_{17}P_2$) (labeled as "Ligand of Interest" by depositor).

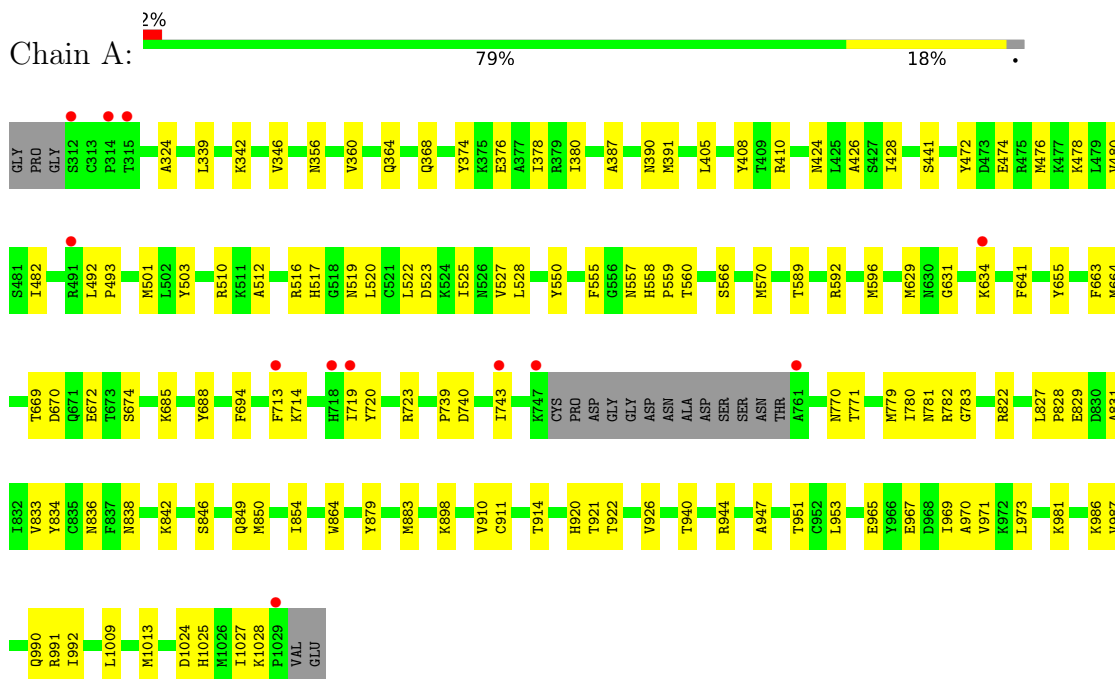


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			39	17	3	17	2		
3	C	1	Total	C	N	O	P	0	0
			39	17	3	17	2		
3	E	1	Total	C	N	O	P	0	0
			39	17	3	17	2		
3	G	1	Total	C	N	O	P	0	0
			39	17	3	17	2		

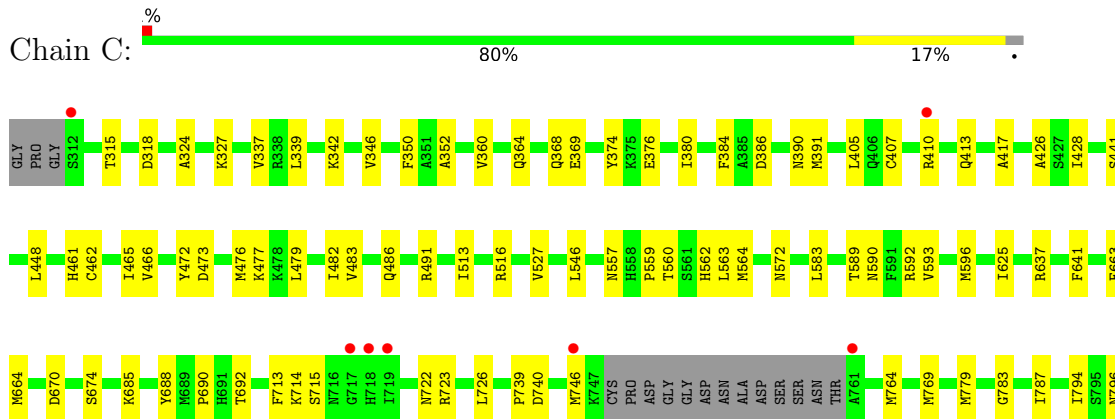
3 Residue-property plots

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

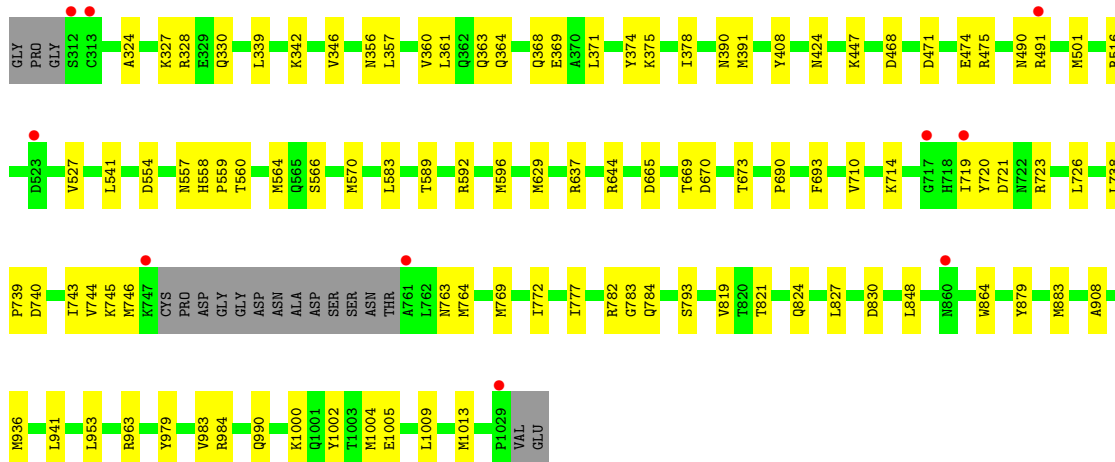
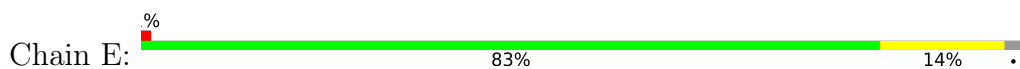


- Molecule 1: UDP-N-acetylglucosamine--peptide N-acetylglucosaminyltransferase 110 kDa subunit

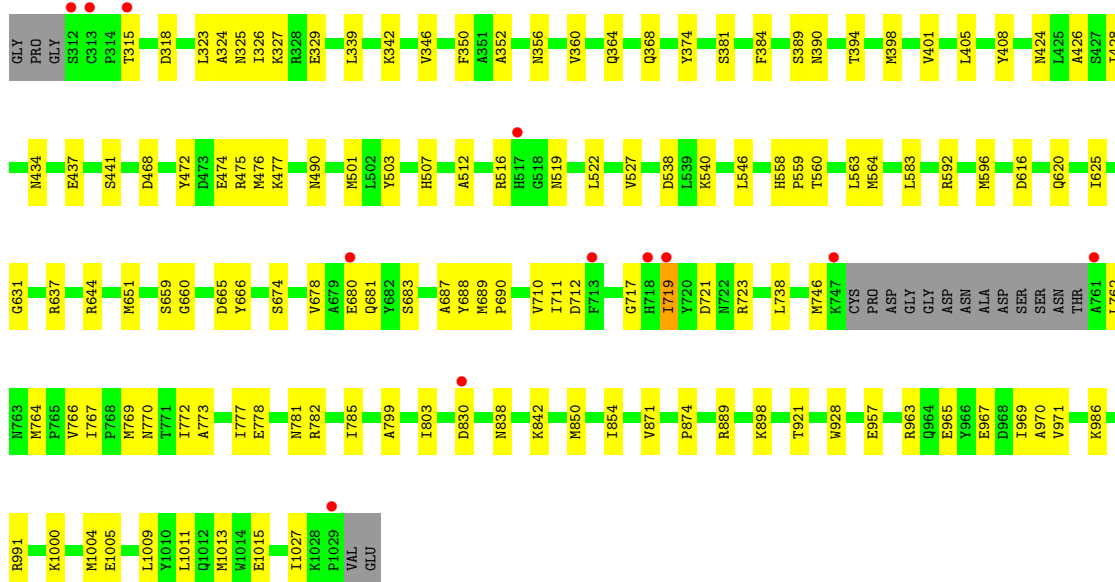
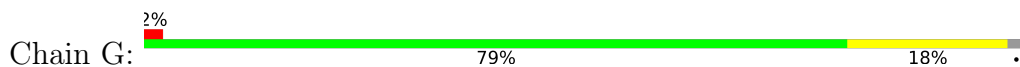




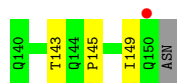
- Molecule 1: UDP-N-acetylglucosamine--peptide N-acetylglucosaminyltransferase 110 kDa subunit



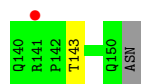
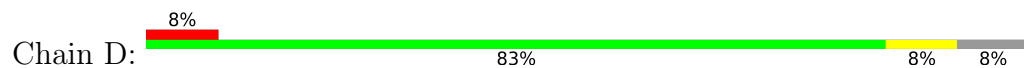
- Molecule 1: UDP-N-acetylglucosamine--peptide N-acetylglucosaminyltransferase 110 kDa subunit



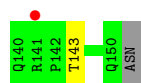
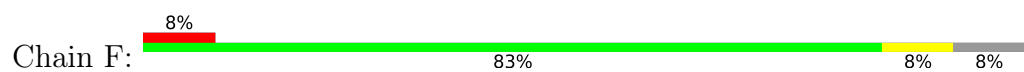
- Molecule 2: Nonsense-mediated mRNA decay factor SMG9



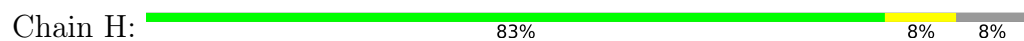
- Molecule 2: Nonsense-mediated mRNA decay factor SMG9



- Molecule 2: Nonsense-mediated mRNA decay factor SMG9



- Molecule 2: Nonsense-mediated mRNA decay factor SMG9



4 Data and refinement statistics

Property	Value	Source
Space group	P 3 2 1	Depositor
Cell constants a, b, c, α , β , γ	273.67Å 273.67Å 143.14Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	118.50 – 2.98 118.50 – 2.98	Depositor EDS
% Data completeness (in resolution range)	99.4 (118.50-2.98) 99.4 (118.50-2.98)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.25 (at 2.96Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.186 , 0.217 (Not available) , 0.206	Depositor DCC
R_{free} test set	1987 reflections (1.58%)	wwPDB-VP
Wilson B-factor (Å ²)	71.9	Xtriage
Anisotropy	0.398	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 46.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.009 for -h,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	22788	wwPDB-VP
Average B, all atoms (Å ²)	74.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.38% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: UD1

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	0.14	0/5693	0.34	0/7722
1	C	0.15	0/5693	0.34	1/7722 (0.0%)
1	E	0.13	0/5693	0.32	0/7722
1	G	0.14	0/5693	0.34	0/7722
2	B	0.15	0/97	0.49	0/132
2	D	0.14	0/97	0.36	0/132
2	F	0.11	0/97	0.30	0/132
2	H	0.10	0/97	0.31	0/132
All	All	0.14	0/23160	0.34	1/31416 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	413	GLN	CA-CB-CG	-5.22	103.66	114.10

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5563	0	5536	90	0
1	C	5563	0	5536	79	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	5563	0	5536	72	0
1	G	5563	0	5536	85	0
2	B	95	0	94	3	0
2	D	95	0	94	1	0
2	F	95	0	94	1	0
2	H	95	0	94	1	0
3	A	39	0	25	5	0
3	C	39	0	25	3	0
3	E	39	0	25	3	0
3	G	39	0	25	5	0
All	All	22788	0	22620	327	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 7.

All (327) 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:A:378:ILE:HD12	1:A:391:MET:HE2	1.55	0.86
1:A:770:ASN:OD1	1:A:771:THR:N	2.17	0.77
1:G:364:GLN:HB3	1:G:527:VAL:HG21	1.68	0.75
1:G:368:GLN:OE1	1:G:368:GLN:N	2.18	0.73
1:C:878:GLN:NE2	1:C:881:GLN:OE1	2.22	0.72
1:E:559:PRO:HB2	3:E:1101:UD1:H6'2	1.72	0.71
1:C:563:LEU:HD21	1:C:921:THR:HG23	1.73	0.70
1:G:315:THR:HG23	1:G:318:ASP:H	1.57	0.70
1:G:616:ASP:O	1:G:620:GLN:HG3	1.92	0.70
1:C:670:ASP:OD2	1:C:692:THR:HA	1.92	0.69
1:G:850:MET:O	1:G:854:ILE:HD12	1.92	0.69
1:E:583:LEU:HD22	1:E:637:ARG:HD3	1.74	0.69
1:G:583:LEU:HD22	1:G:637:ARG:HD3	1.75	0.69
1:G:746:MET:HE3	1:G:762:LEU:HB3	1.74	0.68
1:A:836:ASN:HD21	1:A:914:THR:HG23	1.59	0.68
1:A:476:MET:HE1	1:A:503:TYR:CD1	2.29	0.67
1:A:720:TYR:HD2	1:A:723:ARG:HG3	1.60	0.67
1:C:557:ASN:HB2	1:C:589:THR:HG21	1.76	0.66
1:E:745:LYS:H	1:E:745:LYS:HD2	1.59	0.66
1:A:991:ARG:HG2	1:A:992:ILE:HD12	1.76	0.66
1:G:519:ASN:HA	1:G:522:LEU:HD23	1.78	0.65
1:G:559:PRO:HB2	3:G:1101:UD1:H6'2	1.78	0.65
1:G:592:ARG:O	1:G:596:MET:HG3	1.95	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:559:PRO:HB2	3:C:1101:UD1:H6'2	1.78	0.65
1:A:474:GLU:OE2	1:A:474:GLU:N	2.17	0.65
1:A:592:ARG:HE	1:A:596:MET:HE2	1.61	0.65
1:C:955:CYS:HB3	1:C:958:LEU:HD12	1.79	0.65
1:C:482:ILE:O	1:C:486:GLN:HG3	1.98	0.64
1:E:564:MET:HE2	1:E:629:MET:HE1	1.79	0.63
1:E:740:ASP:OD1	1:E:740:ASP:N	2.32	0.62
1:C:787:ILE:HG13	1:C:794:ILE:HB	1.80	0.62
1:C:590:ASN:HA	1:C:593:VAL:HG12	1.81	0.62
1:E:719:ILE:HG21	1:E:744:VAL:HG21	1.83	0.61
1:G:356:ASN:O	1:G:360:VAL:HG23	2.00	0.61
1:C:740:ASP:OD1	1:C:740:ASP:N	2.34	0.61
1:C:846:SER:O	1:C:850:MET:HG3	2.01	0.61
1:E:324:ALA:HB2	1:E:339:LEU:HB2	1.83	0.61
1:G:512:ALA:O	1:G:516:ARG:HG2	2.01	0.61
1:A:558:HIS:HD2	1:A:560:THR:H	1.50	0.60
1:A:559:PRO:HB2	3:A:1101:UD1:H6'2	1.84	0.60
1:A:410:ARG:HH21	1:C:410:ARG:NH2	1.99	0.60
1:C:690:PRO:HG2	1:C:1005:GLU:HG2	1.83	0.60
1:A:947:ALA:O	1:A:951:THR:HG23	2.02	0.60
1:C:461:HIS:CE1	1:C:465:ILE:HD11	2.37	0.60
1:A:849:GLN:HA	1:A:883:MET:HE1	1.84	0.59
1:C:324:ALA:HB2	1:C:339:LEU:HB2	1.85	0.59
1:G:773:ALA:O	1:G:777:ILE:HD12	2.02	0.59
1:E:979:TYR:O	1:E:983:VAL:HG23	2.03	0.59
1:A:525:ILE:HA	1:A:528:LEU:HD23	1.84	0.59
1:A:965:GLU:O	1:A:969:ILE:HG13	2.03	0.59
1:C:405:LEU:HD13	1:C:428:ILE:HG21	1.84	0.59
1:E:557:ASN:HB2	1:E:589:THR:HG21	1.83	0.59
1:A:476:MET:O	1:A:480:VAL:HG23	2.02	0.58
1:G:738:LEU:HD22	1:G:772:ILE:HD13	1.84	0.58
1:C:376:GLU:O	1:C:380:ILE:HD12	2.04	0.58
1:A:472:TYR:O	1:A:476:MET:HG2	2.04	0.57
1:A:740:ASP:OD1	1:A:740:ASP:N	2.35	0.57
1:A:476:MET:HE1	1:A:503:TYR:HD1	1.69	0.57
1:G:770:ASN:ND2	1:G:772:ILE:H	2.03	0.57
1:C:592:ARG:O	1:C:596:MET:HG3	2.04	0.56
1:E:644:ARG:NH2	1:E:665:ASP:OD2	2.38	0.56
1:A:376:GLU:O	1:A:380:ILE:HD12	2.04	0.56
1:E:471:ASP:HB3	1:E:474:GLU:HG2	1.87	0.56
1:C:713:PHE:HD1	1:C:713:PHE:H	1.53	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:563:LEU:HD21	1:G:921:THR:HG23	1.87	0.56
1:C:723:ARG:NE	1:C:829:GLU:O	2.34	0.56
1:C:714:LYS:HD3	1:C:715:SER:N	2.20	0.56
1:C:337:VAL:HG13	1:C:360:VAL:HG11	1.89	0.55
1:A:854:ILE:HG12	1:A:970:ALA:HB3	1.87	0.55
1:E:491:ARG:N	1:E:516:ARG:HH22	2.04	0.55
1:E:368:GLN:H	1:E:368:GLN:CD	2.12	0.55
1:G:674:SER:O	1:G:688:TYR:OH	2.24	0.55
1:G:678:VAL:O	1:G:681:GLN:HG2	2.06	0.55
1:E:592:ARG:O	1:E:596:MET:HG3	2.07	0.55
1:C:560:THR:O	1:C:564:MET:HG3	2.07	0.55
1:G:546:LEU:HD21	1:G:625:ILE:HD12	1.88	0.55
1:E:558:HIS:CD2	1:E:559:PRO:HD2	2.42	0.55
1:G:342:LYS:O	1:G:346:VAL:HG23	2.07	0.55
1:A:739:PRO:HG2	1:A:740:ASP:OD1	2.07	0.54
1:C:739:PRO:HG2	1:C:740:ASP:OD1	2.07	0.54
1:C:583:LEU:HD22	1:C:637:ARG:HD3	1.88	0.54
1:E:501:MET:HE1	3:E:1101:UD1:H8'1	1.88	0.54
1:A:558:HIS:CD2	1:A:560:THR:H	2.26	0.54
1:A:634:LYS:N	1:A:634:LYS:HD2	2.23	0.54
1:A:364:GLN:HB3	1:A:527:VAL:HG21	1.89	0.54
1:C:783:GLY:O	2:D:143:THR:OG1	2.25	0.54
1:E:936:MET:HE1	1:E:963:ARG:HG3	1.88	0.54
1:G:501:MET:HE1	3:G:1101:UD1:H8'1	1.90	0.54
1:E:374:TYR:CZ	1:E:390:ASN:HB3	2.42	0.54
1:G:631:GLY:HA2	1:G:651:MET:HE1	1.89	0.54
1:A:557:ASN:HB2	1:A:589:THR:HG21	1.91	0.53
1:E:739:PRO:HG2	1:E:740:ASP:OD1	2.08	0.53
1:C:688:TYR:CD2	1:C:1027:ILE:HG23	2.44	0.53
1:E:782:ARG:HD3	1:E:784:GLN:NE2	2.23	0.53
1:G:401:VAL:HG13	1:G:428:ILE:HD12	1.90	0.53
1:E:490:ASN:C	1:E:516:ARG:HH22	2.17	0.53
1:C:364:GLN:HB3	1:C:527:VAL:HG11	1.90	0.53
1:A:922:THR:O	1:A:926:VAL:HG12	2.10	0.52
1:A:879:TYR:O	1:A:883:MET:HG3	2.07	0.52
1:A:986:LYS:O	1:A:990:GLN:HG2	2.09	0.52
1:G:723:ARG:NH2	1:G:830:ASP:HA	2.24	0.52
1:C:670:ASP:OD2	1:C:948:SER:OG	2.28	0.52
1:A:555:PHE:O	1:A:592:ARG:HD3	2.09	0.52
1:A:967:GLU:O	1:A:971:VAL:HG13	2.09	0.52
1:E:560:THR:O	1:E:564:MET:HG3	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:378:ILE:CD1	1:A:391:MET:HE2	2.35	0.52
1:G:778:GLU:OE2	1:G:782:ARG:HD2	2.10	0.52
1:E:327:LYS:HD3	1:E:330:GLN:OE1	2.10	0.51
1:E:491:ARG:C	1:E:516:ARG:HH22	2.19	0.51
1:G:688:TYR:HD2	1:G:1027:ILE:HG23	1.74	0.51
1:A:781:ASN:HD21	2:B:149:ILE:H	1.59	0.51
1:A:342:LYS:O	1:A:346:VAL:HG23	2.10	0.51
1:A:405:LEU:HD13	1:A:428:ILE:HG21	1.91	0.51
1:E:848:LEU:HG	1:E:883:MET:HE1	1.92	0.51
1:A:674:SER:O	1:A:688:TYR:OH	2.25	0.51
1:C:828:PRO:HG2	1:C:831:ALA:HB3	1.93	0.51
1:E:491:ARG:O	1:E:516:ARG:NH2	2.44	0.51
1:G:468:ASP:O	1:G:475:ARG:NH2	2.43	0.51
1:G:434:ASN:HB3	1:G:437:GLU:OE1	2.11	0.50
1:G:560:THR:O	1:G:564:MET:HG3	2.11	0.50
1:G:957:GLU:OE2	1:G:986:LYS:HE2	2.11	0.50
1:G:854:ILE:HG12	1:G:970:ALA:HB3	1.92	0.50
1:G:963:ARG:O	1:G:967:GLU:HG3	2.11	0.50
1:G:476:MET:HE1	1:G:503:TYR:HD1	1.76	0.50
1:E:328:ARG:NH2	1:E:363:GLN:OE1	2.44	0.50
1:E:592:ARG:HG3	1:E:596:MET:HE2	1.93	0.50
1:G:838:ASN:HB3	1:G:842:LYS:HD2	1.93	0.50
1:G:325:ASN:O	1:G:329:GLU:HG2	2.12	0.50
1:C:426:ALA:HB2	1:C:441:SER:HB2	1.93	0.50
1:C:391:MET:HE3	1:C:407:CYS:SG	2.52	0.50
1:G:1000:LYS:O	1:G:1004:MET:HG3	2.11	0.50
1:A:842:LYS:NZ	3:A:1101:UD1:O1B	2.41	0.49
1:C:516:ARG:HG2	1:C:516:ARG:HH11	1.76	0.49
1:E:726:LEU:HG	1:E:819:VAL:HG22	1.94	0.49
1:G:690:PRO:HG2	1:G:1005:GLU:HG2	1.94	0.49
1:E:468:ASP:O	1:E:475:ARG:NH2	2.45	0.49
3:E:1101:UD1:H1'	3:E:1101:UD1:O1A	2.11	0.49
1:A:663:PHE:HD1	1:A:664:MET:HE2	1.76	0.49
1:E:720:TYR:CD2	1:E:723:ARG:HG3	2.48	0.49
1:E:738:LEU:HD11	1:E:772:ILE:HG21	1.95	0.49
1:G:688:TYR:CD2	1:G:1027:ILE:HG23	2.47	0.49
1:C:878:GLN:HE21	1:C:878:GLN:HA	1.77	0.49
1:A:503:TYR:O	1:A:510:ARG:NH1	2.46	0.49
1:E:447:LYS:HB3	1:E:447:LYS:HE2	1.53	0.49
1:G:644:ARG:NH2	1:G:665:ASP:OD1	2.46	0.49
1:A:846:SER:O	1:A:850:MET:HG3	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:315:THR:H	1:G:318:ASP:HB2	1.78	0.48
1:C:878:GLN:NE2	1:C:878:GLN:HA	2.28	0.48
1:E:541:LEU:HD12	1:E:541:LEU:H	1.78	0.48
1:G:538:ASP:OD1	1:G:540:LYS:HG3	2.13	0.48
1:E:566:SER:HB3	1:E:570:MET:HE3	1.96	0.48
1:E:783:GLY:O	2:F:143:THR:OG1	2.31	0.48
1:G:898:LYS:NZ	3:G:1101:UD1:O3B	2.47	0.48
1:A:713:PHE:HE2	1:A:829:GLU:CD	2.21	0.48
1:A:1009:LEU:O	1:A:1013:MET:HG3	2.12	0.48
1:C:350:PHE:CE1	1:C:352:ALA:HB3	2.49	0.48
1:E:723:ARG:HH21	1:E:830:ASP:H	1.61	0.48
1:A:1027:ILE:HG22	1:A:1028:LYS:HD3	1.96	0.48
1:A:921:THR:N	3:A:1101:UD1:O2B	2.46	0.48
1:C:417:ALA:HA	1:C:448:LEU:HD22	1.95	0.48
1:C:852:ALA:HA	1:C:885:LEU:HD11	1.95	0.48
1:A:828:PRO:HB2	1:A:831:ALA:HB3	1.96	0.48
1:E:710:VAL:HG11	1:E:721:ASP:HA	1.96	0.48
1:G:781:ASN:HD21	2:H:149:ILE:H	1.61	0.48
1:G:1011:LEU:O	1:G:1015:GLU:HG2	2.14	0.47
1:G:324:ALA:HB2	1:G:339:LEU:HB2	1.96	0.47
1:G:921:THR:N	3:G:1101:UD1:O2B	2.47	0.47
1:G:323:LEU:HA	1:G:326:ILE:HG12	1.97	0.47
1:G:957:GLU:OE1	1:G:957:GLU:N	2.26	0.47
1:A:374:TYR:CZ	1:A:390:ASN:HB3	2.50	0.47
1:A:378:ILE:HD11	1:A:387:ALA:C	2.39	0.47
1:A:368:GLN:CD	1:A:368:GLN:H	2.22	0.47
1:E:342:LYS:O	1:E:346:VAL:HG12	2.15	0.47
1:E:356:ASN:O	1:E:360:VAL:HG22	2.15	0.47
1:E:371:LEU:HD21	1:E:375:LYS:HE3	1.96	0.47
1:E:769:MET:HE1	1:E:777:ILE:HD12	1.96	0.47
1:E:1009:LEU:HG	1:E:1013:MET:HE3	1.97	0.47
1:G:408:TYR:CZ	1:G:424:ASN:HB3	2.50	0.47
1:G:799:ALA:O	1:G:803:ILE:HD12	2.14	0.47
1:G:721:ASP:HB3	1:G:764:MET:HE1	1.96	0.47
3:G:1101:UD1:H1'	3:G:1101:UD1:O1A	2.15	0.47
1:C:342:LYS:O	1:C:346:VAL:HG13	2.15	0.46
1:E:364:GLN:HB3	1:E:527:VAL:HG21	1.96	0.46
1:E:821:THR:OG1	1:E:824:GLN:HG3	2.15	0.46
1:A:685:LYS:HD2	1:A:1025:HIS:CE1	2.50	0.46
1:A:782:ARG:HH11	1:A:782:ARG:HG2	1.81	0.46
1:A:953:LEU:HD21	1:A:987:VAL:HG13	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:491:ARG:HD2	1:C:491:ARG:HA	1.80	0.46
1:E:378:ILE:HD12	1:E:391:MET:HE2	1.97	0.46
1:E:1000:LYS:O	1:E:1004:MET:HG3	2.15	0.46
1:G:651:MET:HE2	1:G:659:SER:HB3	1.98	0.46
1:G:778:GLU:O	1:G:782:ARG:HG3	2.14	0.46
1:C:572:ASN:OD1	1:E:793:SER:HB3	2.15	0.46
1:E:690:PRO:HG2	1:E:1005:GLU:HG2	1.96	0.46
1:G:326:ILE:HG13	1:G:327:LYS:N	2.30	0.46
1:A:723:ARG:HA	1:A:822:ARG:HG3	1.96	0.46
1:G:967:GLU:O	1:G:971:VAL:HG23	2.15	0.46
1:G:472:TYR:CZ	1:G:476:MET:HE2	2.51	0.46
1:A:910:VAL:HG21	1:A:973:LEU:HB3	1.98	0.46
1:C:472:TYR:CZ	1:C:476:MET:HE3	2.51	0.46
1:C:726:LEU:HD22	1:C:819:VAL:HG22	1.97	0.46
1:G:389:SER:OG	1:G:424:ASN:ND2	2.49	0.46
1:C:473:ASP:O	1:C:477:LYS:HG3	2.16	0.45
1:C:546:LEU:HD21	1:C:625:ILE:HD12	1.97	0.45
1:C:963:ARG:H	1:C:963:ARG:HE	1.63	0.45
1:A:324:ALA:HB2	1:A:339:LEU:HB2	1.98	0.45
1:C:722:ASN:OD1	1:C:723:ARG:NH1	2.39	0.45
1:G:871:VAL:O	1:G:874:PRO:HD2	2.17	0.45
1:E:743:ILE:HG23	1:E:763:ASN:HB2	1.99	0.45
1:A:898:LYS:NZ	3:A:1101:UD1:O3B	2.49	0.45
1:E:673:THR:HG22	1:E:941:LEU:HG	1.99	0.45
1:C:327:LYS:HA	1:C:327:LYS:HD3	1.70	0.45
1:C:746:MET:HE3	1:C:764:MET:HB2	1.98	0.45
1:C:641:PHE:CD1	1:C:664:MET:HE1	2.51	0.45
1:G:712:ASP:OD1	1:G:717:GLY:HA3	2.16	0.45
1:A:478:LYS:O	1:A:482:ILE:HG13	2.16	0.45
1:A:827:LEU:HD13	1:A:833:VAL:HG11	1.97	0.45
1:E:669:THR:OG1	1:E:670:ASP:N	2.49	0.44
1:A:631:GLY:HA3	1:A:655:TYR:HB2	1.98	0.44
1:A:356:ASN:O	1:A:360:VAL:HG23	2.18	0.44
3:A:1101:UD1:H1'	3:A:1101:UD1:O1A	2.17	0.44
1:C:479:LEU:O	1:C:483:VAL:HG23	2.18	0.44
1:G:405:LEU:HG	1:G:428:ILE:HG21	2.00	0.44
1:G:476:MET:HE1	1:G:503:TYR:CD1	2.53	0.44
1:C:779:MET:HE2	1:C:779:MET:HB3	1.88	0.44
1:A:713:PHE:CG	1:A:714:LYS:N	2.85	0.44
1:C:806:LYS:HB3	1:C:812:GLU:HG3	1.99	0.44
1:C:1025:HIS:HB3	1:C:1027:ILE:HD11	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:491:ARG:C	1:E:516:ARG:NH2	2.76	0.44
1:A:669:THR:OG1	1:A:670:ASP:N	2.51	0.44
1:A:854:ILE:HD11	1:A:967:GLU:HA	1.99	0.44
1:G:426:ALA:HB2	1:G:441:SER:CB	2.48	0.44
1:A:408:TYR:CZ	1:A:424:ASN:HB3	2.52	0.44
1:A:834:TYR:O	1:A:864:TRP:N	2.47	0.44
1:E:644:ARG:HH22	1:E:665:ASP:CG	2.26	0.44
1:C:462:CYS:O	1:C:466:VAL:HG22	2.17	0.43
1:G:394:THR:O	1:G:398:MET:HG3	2.18	0.43
1:E:714:LYS:HD3	1:E:714:LYS:HA	1.71	0.43
1:E:721:ASP:OD1	1:E:721:ASP:C	2.61	0.43
1:G:666:TYR:CG	1:G:1013:MET:HE2	2.53	0.43
1:G:889:ARG:HD2	1:G:889:ARG:HA	1.76	0.43
1:G:965:GLU:O	1:G:969:ILE:HG13	2.18	0.43
1:A:940:THR:O	1:A:944:ARG:HG3	2.18	0.43
3:C:1101:UD1:H1'	3:C:1101:UD1:O1A	2.18	0.43
1:A:566:SER:HB3	1:A:570:MET:HE3	2.01	0.43
1:A:782:ARG:HG2	1:A:782:ARG:NH1	2.33	0.43
1:C:324:ALA:HB2	1:C:339:LEU:CB	2.48	0.43
1:E:361:LEU:HD13	1:E:369:GLU:HB3	2.00	0.43
1:G:374:TYR:CZ	1:G:390:ASN:HB3	2.53	0.43
1:A:779:MET:HE2	1:A:779:MET:HB3	1.94	0.43
1:C:879:TYR:O	1:C:883:MET:HG3	2.19	0.43
1:E:738:LEU:HD11	1:E:772:ILE:HG12	2.00	0.43
1:G:680:GLU:H	1:G:680:GLU:HG3	1.66	0.43
1:E:879:TYR:O	1:E:883:MET:HG3	2.19	0.43
1:A:512:ALA:O	1:A:516:ARG:HG2	2.19	0.43
1:A:780:ILE:O	2:B:145:PRO:HG3	2.19	0.43
1:A:981:LYS:HB2	1:A:981:LYS:HE3	1.80	0.43
1:C:663:PHE:HD1	1:C:664:MET:HE2	1.83	0.43
1:E:408:TYR:CZ	1:E:424:ASN:HB3	2.53	0.43
1:G:507:HIS:HB3	1:G:681:GLN:HB3	2.01	0.43
1:G:711:ILE:HG12	1:G:767:ILE:HB	2.01	0.43
1:A:410:ARG:HH21	1:C:410:ARG:HH22	1.66	0.43
1:A:911:CYS:CB	1:A:926:VAL:HG21	2.49	0.43
1:C:562:HIS:ND1	1:C:898:LYS:HE3	2.34	0.43
1:C:921:THR:N	3:C:1101:UD1:O2B	2.52	0.42
1:E:378:ILE:HD12	1:E:391:MET:CE	2.48	0.42
1:G:928:TRP:O	1:G:991:ARG:NH1	2.50	0.42
1:A:550:TYR:HB3	1:A:629:MET:SD	2.59	0.42
1:C:796:ASN:OD1	1:C:824:GLN:NE2	2.51	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:855:LEU:O	1:C:889:ARG:NH2	2.44	0.42
1:C:1009:LEU:HG	1:C:1013:MET:HE3	2.02	0.42
1:G:558:HIS:CG	1:G:559:PRO:HD2	2.54	0.42
1:C:714:LYS:HD3	1:C:715:SER:H	1.85	0.42
1:G:689:MET:HE3	1:G:689:MET:HB3	1.81	0.42
1:A:519:ASN:O	1:A:523:ASP:OD2	2.38	0.42
1:C:713:PHE:HA	1:C:769:MET:HG2	2.01	0.42
1:A:522:LEU:O	1:A:525:ILE:HG13	2.19	0.42
1:C:746:MET:CE	1:C:764:MET:HB2	2.49	0.42
1:E:357:LEU:O	1:E:361:LEU:HG	2.19	0.42
1:E:746:MET:HE2	1:E:764:MET:HE2	2.00	0.42
1:E:953:LEU:HD12	1:E:990:GLN:HB3	2.02	0.42
1:A:672:GLU:O	1:A:944:ARG:NH2	2.50	0.42
1:G:660:GLY:HA2	1:G:683:SER:HB3	2.02	0.42
1:G:785:ILE:HG23	1:G:803:ILE:HD11	2.01	0.42
1:C:483:VAL:HG11	1:C:513:ILE:HD11	2.02	0.42
1:A:838:ASN:HB3	1:A:842:LYS:HD2	2.02	0.42
1:G:770:ASN:HD21	1:G:772:ILE:HB	1.85	0.42
1:A:719:ILE:O	1:A:719:ILE:HD12	2.20	0.41
1:C:822:ARG:NH2	1:C:909:ASP:OD1	2.53	0.41
1:C:1011:LEU:HD23	1:C:1011:LEU:HA	1.90	0.41
1:G:769:MET:HE2	1:G:769:MET:HB2	1.87	0.41
1:E:908:ALA:O	1:E:984:ARG:NH1	2.53	0.41
1:A:492:LEU:HD12	1:A:493:PRO:HD2	2.02	0.41
1:C:685:LYS:HG2	1:C:1024:ASP:HA	2.01	0.41
1:C:910:VAL:HG21	1:C:973:LEU:HB3	2.02	0.41
1:C:368:GLN:HG2	1:C:369:GLU:N	2.36	0.41
1:C:641:PHE:HB3	1:C:664:MET:HE1	2.02	0.41
1:G:360:VAL:O	1:G:364:GLN:HG3	2.21	0.41
1:E:721:ASP:CG	1:E:984:ARG:HH21	2.28	0.41
1:G:719:ILE:HG13	1:G:719:ILE:O	2.21	0.41
1:A:694:PHE:CZ	1:A:920:HIS:HB3	2.56	0.41
1:G:381:SER:HB3	1:G:384:PHE:HB2	2.02	0.41
1:G:490:ASN:HA	1:G:516:ARG:NH2	2.35	0.41
1:C:674:SER:O	1:C:688:TYR:OH	2.31	0.41
1:E:745:LYS:HD2	1:E:745:LYS:N	2.31	0.41
1:A:663:PHE:CD1	1:A:664:MET:HE2	2.55	0.41
1:A:685:LYS:HE3	1:A:1024:ASP:HA	2.02	0.41
1:A:783:GLY:O	2:B:143:THR:OG1	2.38	0.41
1:A:911:CYS:HB2	1:A:926:VAL:HG21	2.02	0.41
1:E:693:PHE:O	1:E:1002:TYR:OH	2.30	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:827:LEU:HD11	1:E:864:TRP:CZ3	2.56	0.41
1:G:474:GLU:OE2	1:G:477:LYS:HE2	2.21	0.41
1:G:350:PHE:CE1	1:G:352:ALA:HB3	2.56	0.41
1:C:315:THR:HG23	1:C:318:ASP:OD1	2.20	0.40
1:A:641:PHE:CG	1:A:664:MET:HE1	2.56	0.40
1:A:743:ILE:N	1:A:743:ILE:HD12	2.36	0.40
1:A:426:ALA:HB2	1:A:441:SER:CB	2.51	0.40
1:E:327:LYS:HD3	1:E:327:LYS:HA	1.79	0.40
1:E:554:ASP:HB3	1:E:558:HIS:HD1	1.86	0.40
1:A:501:MET:HE2	1:A:501:MET:HB3	1.73	0.40
1:C:384:PHE:CE1	1:C:386:ASP:HB3	2.56	0.40
1:E:490:ASN:C	1:E:516:ARG:NH2	2.80	0.40
1:G:687:ALA:HB1	1:G:1009:LEU:HD21	2.03	0.40
1:A:517:HIS:HA	1:A:520:LEU:HD12	2.04	0.40
1:C:374:TYR:CZ	1:C:390:ASN:HB3	2.57	0.40
1:C:871:VAL:O	1:C:874:PRO:HD2	2.21	0.40
1:G:710:VAL:HG23	1:G:766:VAL:HG13	2.03	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	701/723 (97%)	680 (97%)	21 (3%)	0	100	100
1	C	701/723 (97%)	685 (98%)	16 (2%)	0	100	100
1	E	701/723 (97%)	688 (98%)	13 (2%)	0	100	100
1	G	701/723 (97%)	682 (97%)	18 (3%)	1 (0%)	48	78
2	B	9/12 (75%)	9 (100%)	0	0	100	100
2	D	9/12 (75%)	7 (78%)	2 (22%)	0	100	100
2	F	9/12 (75%)	9 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	H	9/12 (75%)	9 (100%)	0	0	100	100
All	All	2840/2940 (97%)	2769 (98%)	70 (2%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	719	ILE

5.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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	$\# Z > 2$
3	UD1	A	1101	-	40,41,41	0.28	0	59,62,62	0.58	1 (1%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	UD1	G	1101	-	40,41,41	0.27	0	59,62,62	0.53	1 (1%)
3	UD1	C	1101	-	40,41,41	0.27	0	59,62,62	0.51	1 (1%)
3	UD1	E	1101	-	40,41,41	0.30	0	59,62,62	0.55	1 (1%)

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	UD1	A	1101	-	-	5/26/63/63	0/3/3/3
3	UD1	G	1101	-	-	3/26/63/63	0/3/3/3
3	UD1	C	1101	-	-	5/26/63/63	0/3/3/3
3	UD1	E	1101	-	-	7/26/63/63	0/3/3/3

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1101	UD1	PB-O1'-C1'	-3.51	106.92	121.21
3	E	1101	UD1	PB-O1'-C1'	-3.15	108.40	121.21
3	C	1101	UD1	PB-O1'-C1'	-2.98	109.06	121.21
3	G	1101	UD1	PB-O1'-C1'	-2.94	109.22	121.21

There are no chirality outliers.

All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	1101	UD1	C3B-C4B-C5B-O5B
3	C	1101	UD1	O4B-C4B-C5B-O5B
3	E	1101	UD1	O4B-C4B-C5B-O5B
3	E	1101	UD1	C3B-C4B-C5B-O5B
3	A	1101	UD1	O4B-C4B-C5B-O5B
3	G	1101	UD1	O4B-C4B-C5B-O5B
3	G	1101	UD1	C3B-C4B-C5B-O5B
3	A	1101	UD1	C3B-C4B-C5B-O5B
3	C	1101	UD1	C4'-C5'-C6'-O6'
3	A	1101	UD1	O5'-C1'-O1'-PB
3	C	1101	UD1	O5'-C1'-O1'-PB

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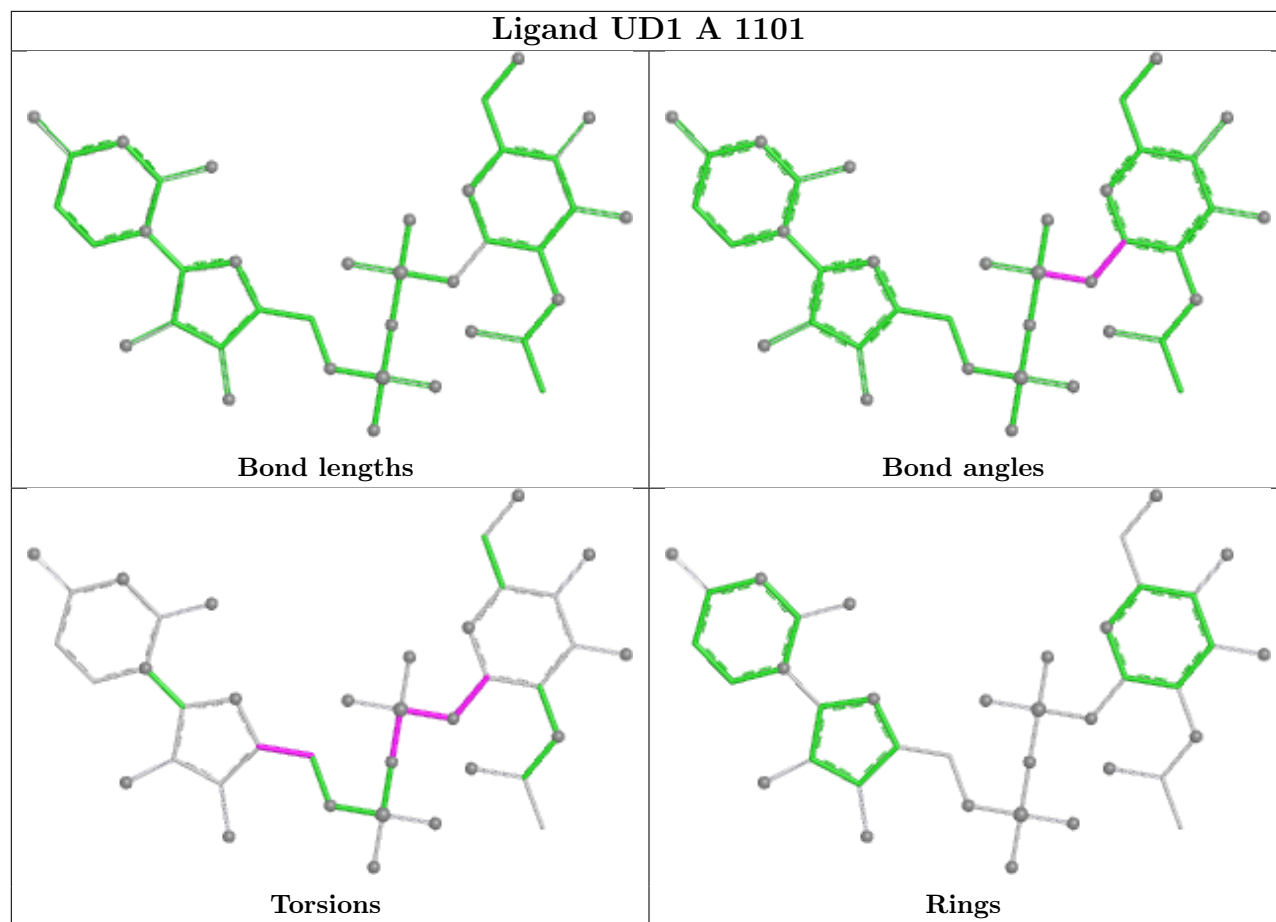
Mol	Chain	Res	Type	Atoms
3	E	1101	UD1	O5'-C1'-O1'-PB
3	G	1101	UD1	O5'-C1'-O1'-PB
3	C	1101	UD1	O5'-C5'-C6'-O6'
3	E	1101	UD1	C4'-C5'-C6'-O6'
3	A	1101	UD1	C1'-O1'-PB-O1B
3	E	1101	UD1	C1'-O1'-PB-O1B
3	E	1101	UD1	O5'-C5'-C6'-O6'
3	A	1101	UD1	PA-O3A-PB-O2B
3	E	1101	UD1	PA-O3A-PB-O2B

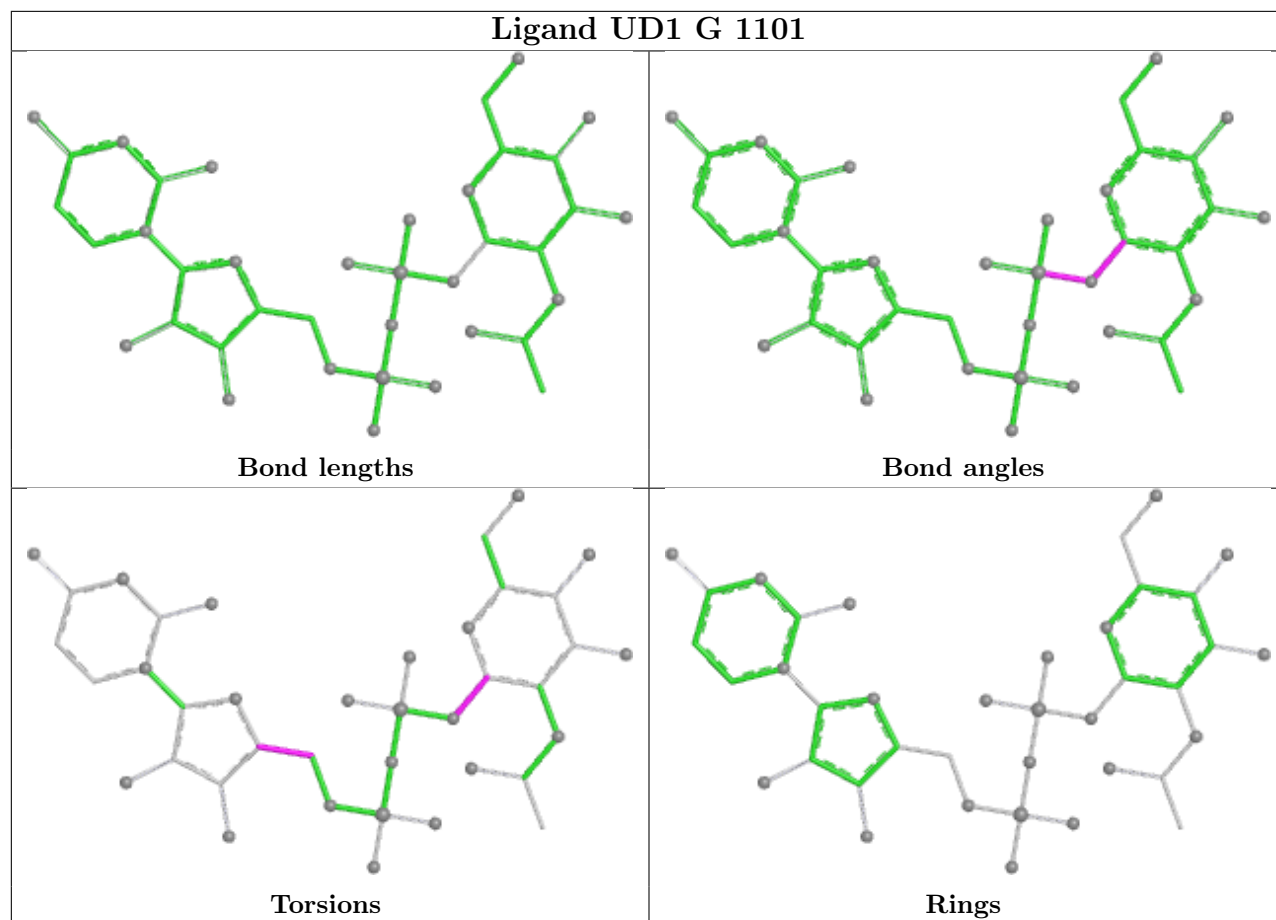
There are no ring outliers.

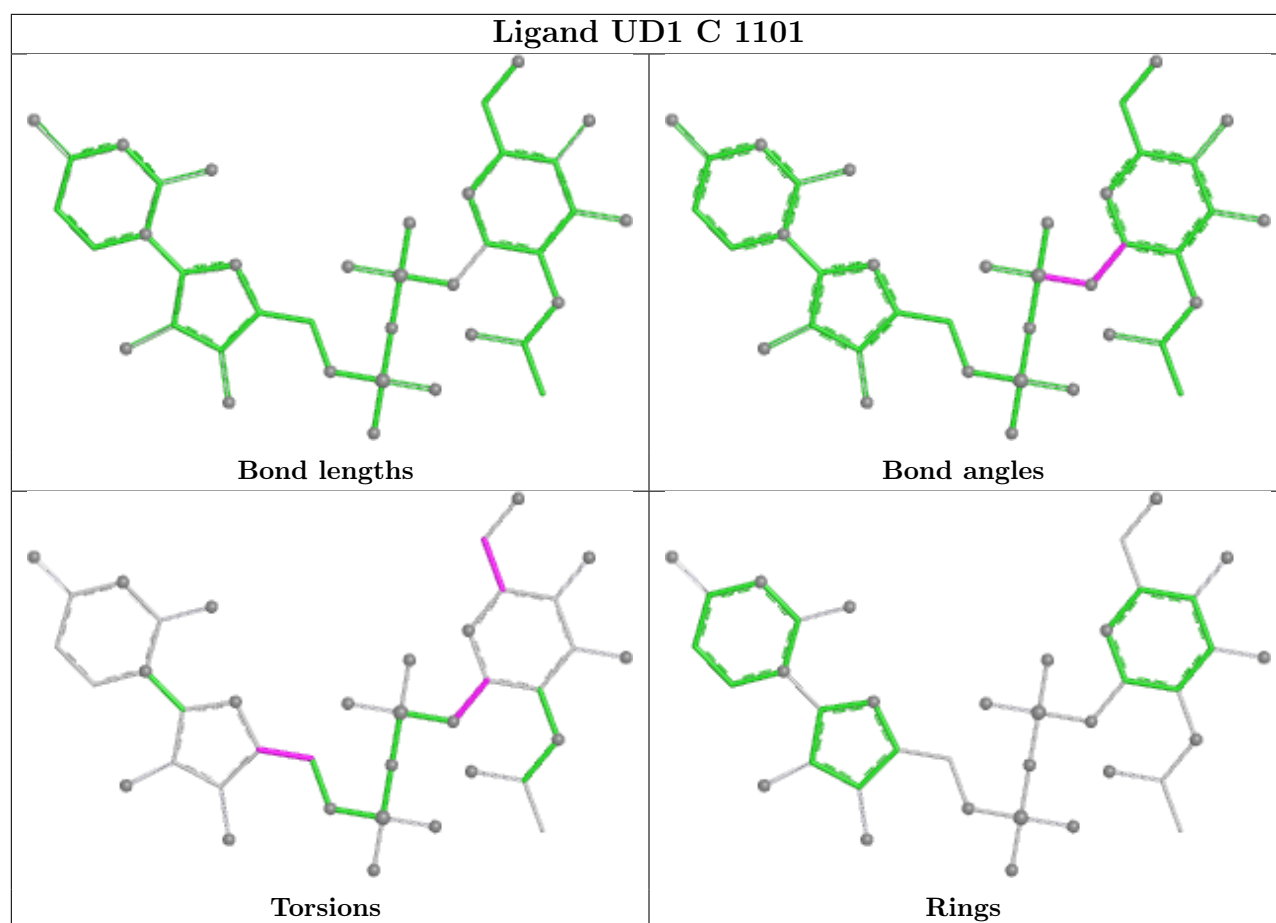
4 monomers are involved in 16 short contacts:

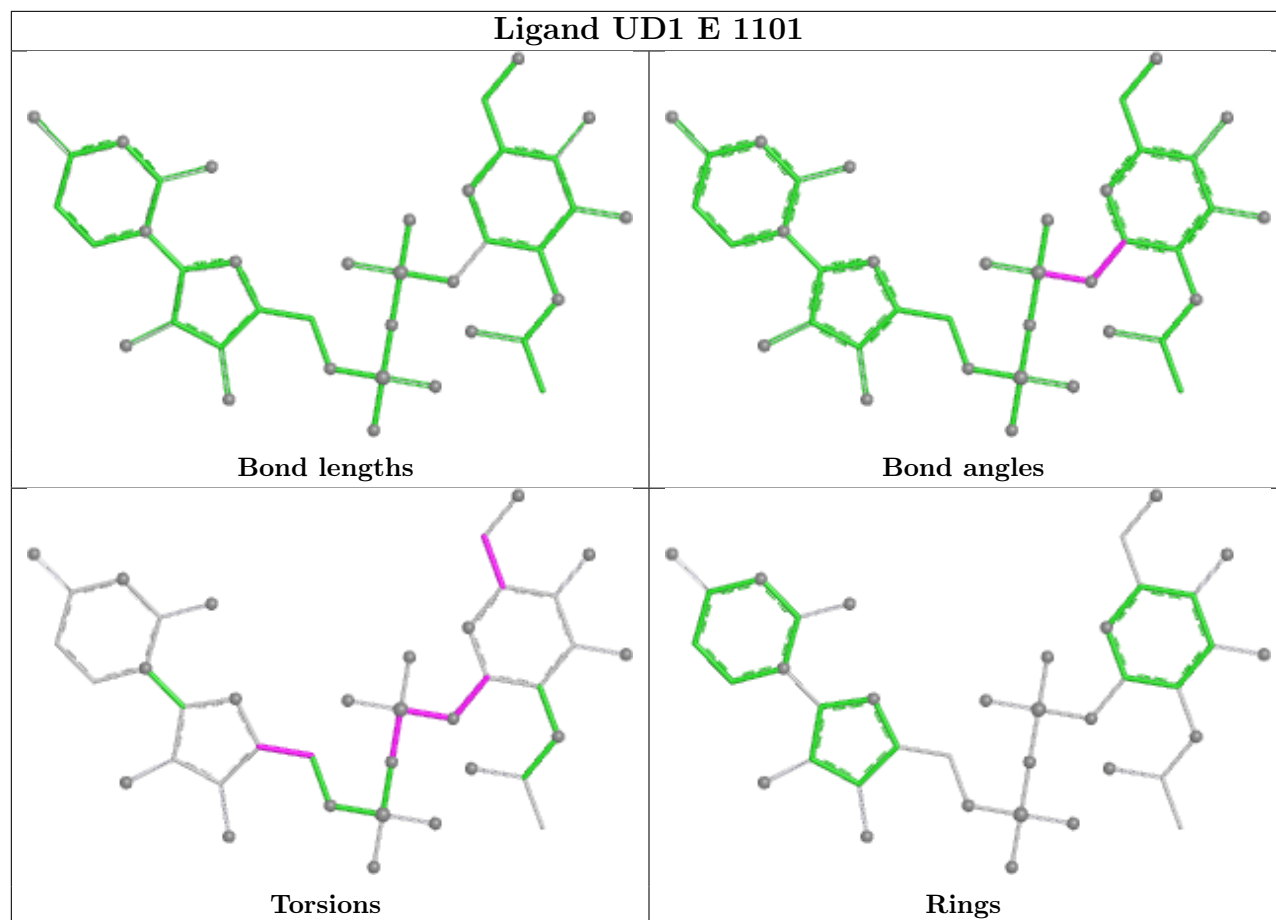
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1101	UD1	5	0
3	G	1101	UD1	5	0
3	C	1101	UD1	3	0
3	E	1101	UD1	3	0

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.









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	705/723 (97%)	-0.02	12 (1%) 69 50	53, 80, 108, 172	0
1	C	705/723 (97%)	-0.17	8 (1%) 78 61	50, 69, 101, 156	0
1	E	705/723 (97%)	-0.18	10 (1%) 73 55	48, 66, 101, 154	0
1	G	705/723 (97%)	-0.21	12 (1%) 69 50	52, 66, 98, 166	0
2	B	11/12 (91%)	0.87	1 (9%) 15 9	78, 84, 122, 137	0
2	D	11/12 (91%)	0.52	1 (9%) 15 9	74, 85, 123, 133	0
2	F	11/12 (91%)	0.47	1 (9%) 15 9	57, 65, 114, 130	0
2	H	11/12 (91%)	0.14	0 100 100	62, 71, 108, 122	0
All	All	2864/2940 (97%)	-0.14	45 (1%) 70 51	48, 70, 105, 172	0

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	719	ILE	5.5
1	C	1029	PRO	5.0
1	A	1029	PRO	4.9
1	E	1029	PRO	4.8
1	G	747	LYS	4.7
1	C	761	ALA	4.3
1	E	761	ALA	4.1
1	G	1029	PRO	3.9
1	A	761	ALA	3.9
1	A	719	ILE	3.7
1	G	312	SER	3.5
1	G	761	ALA	3.5
1	A	718	HIS	3.3
1	G	313	CYS	3.3
1	E	717	GLY	3.3
1	A	312	SER	3.1

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Mol	Chain	Res	Type	RSRZ
1	C	718	HIS	3.0
1	E	312	SER	2.9
1	G	713	PHE	2.9
1	A	747	LYS	2.7
2	B	150	GLN	2.7
1	E	860	ASN	2.6
1	E	719	ILE	2.6
1	G	680	GLU	2.6
1	C	312	SER	2.6
1	C	410	ARG	2.5
1	G	718	HIS	2.5
1	A	491	ARG	2.4
1	G	315	THR	2.4
1	C	746	MET	2.4
1	C	717	GLY	2.4
2	F	141	ARG	2.3
1	E	523	ASP	2.3
2	D	141	ARG	2.3
1	G	719	ILE	2.3
1	A	713	PHE	2.3
1	E	491	ARG	2.3
1	E	747	LYS	2.3
1	A	314	PRO	2.3
1	A	315	THR	2.2
1	E	313	CYS	2.2
1	G	517	HIS	2.1
1	A	743	ILE	2.1
1	G	830	ASP	2.1
1	A	634	LYS	2.0

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 ⓘ

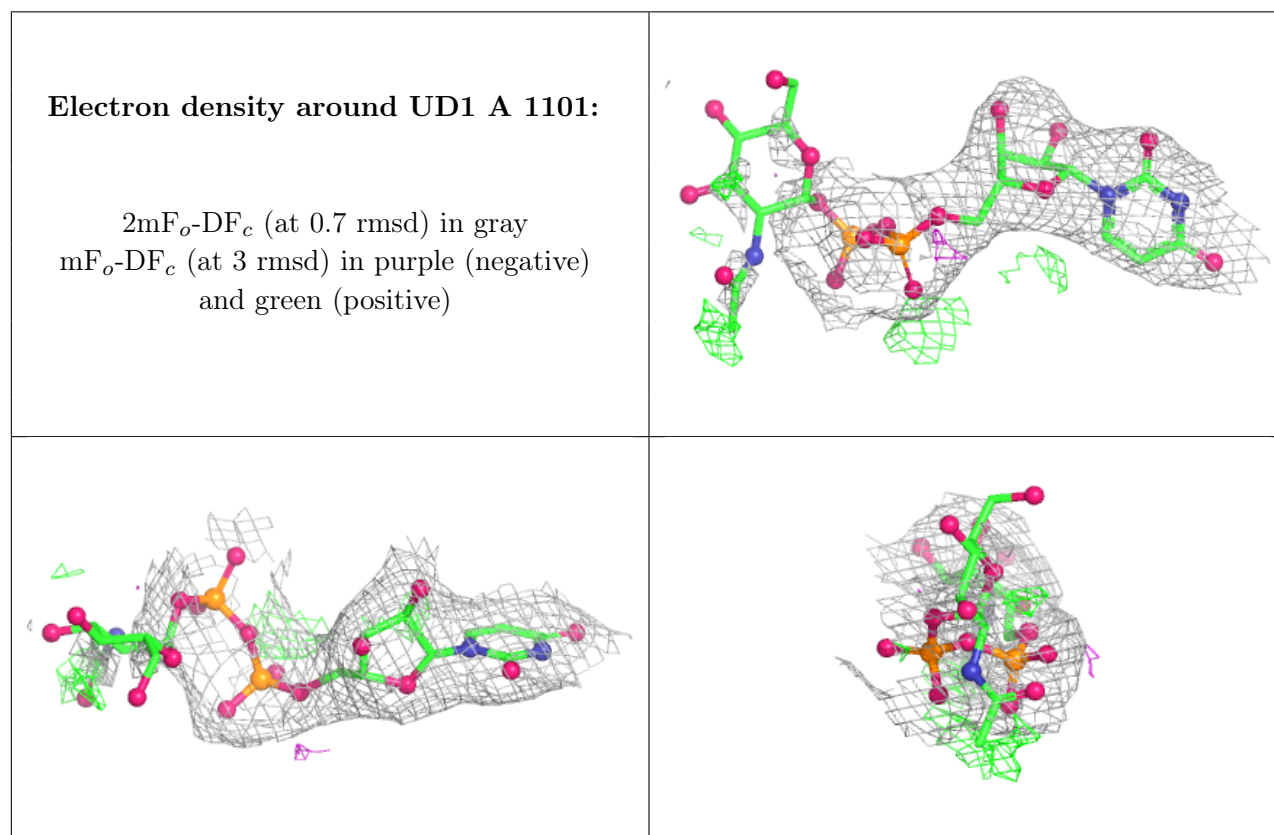
There are no oligosaccharides in this entry.

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.

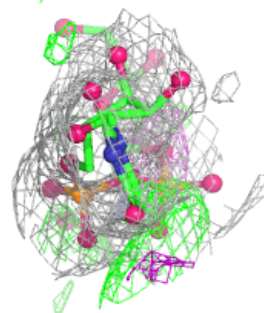
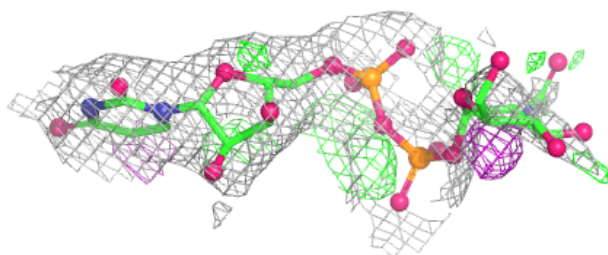
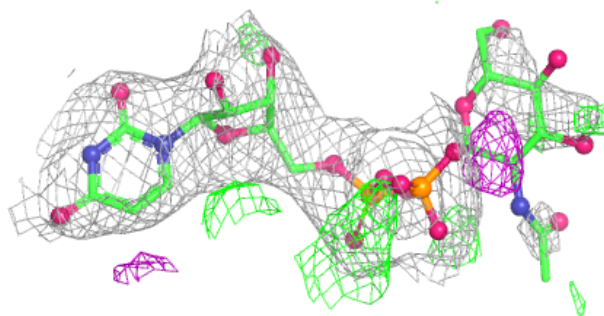
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	UD1	A	1101	39/39	0.96	0.12	61,76,97,98	15
3	UD1	G	1101	39/39	0.96	0.11	47,63,83,87	15
3	UD1	E	1101	39/39	0.97	0.10	46,63,83,86	15
3	UD1	C	1101	39/39	0.97	0.11	51,62,83,84	15

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.

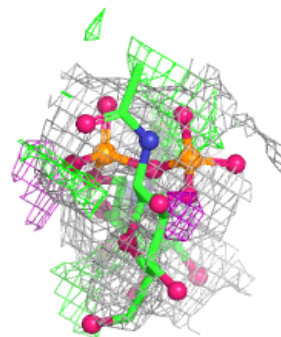
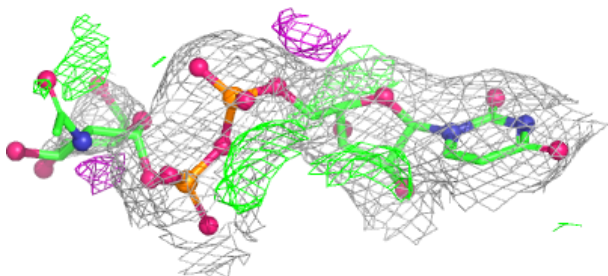
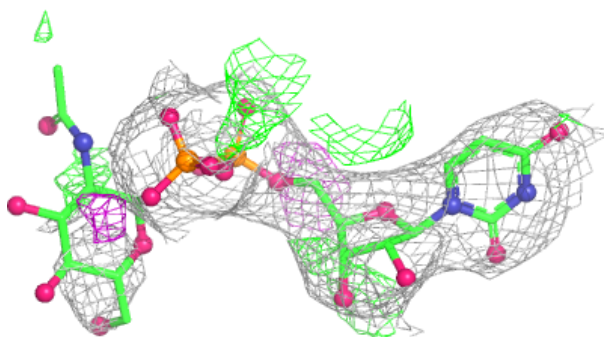


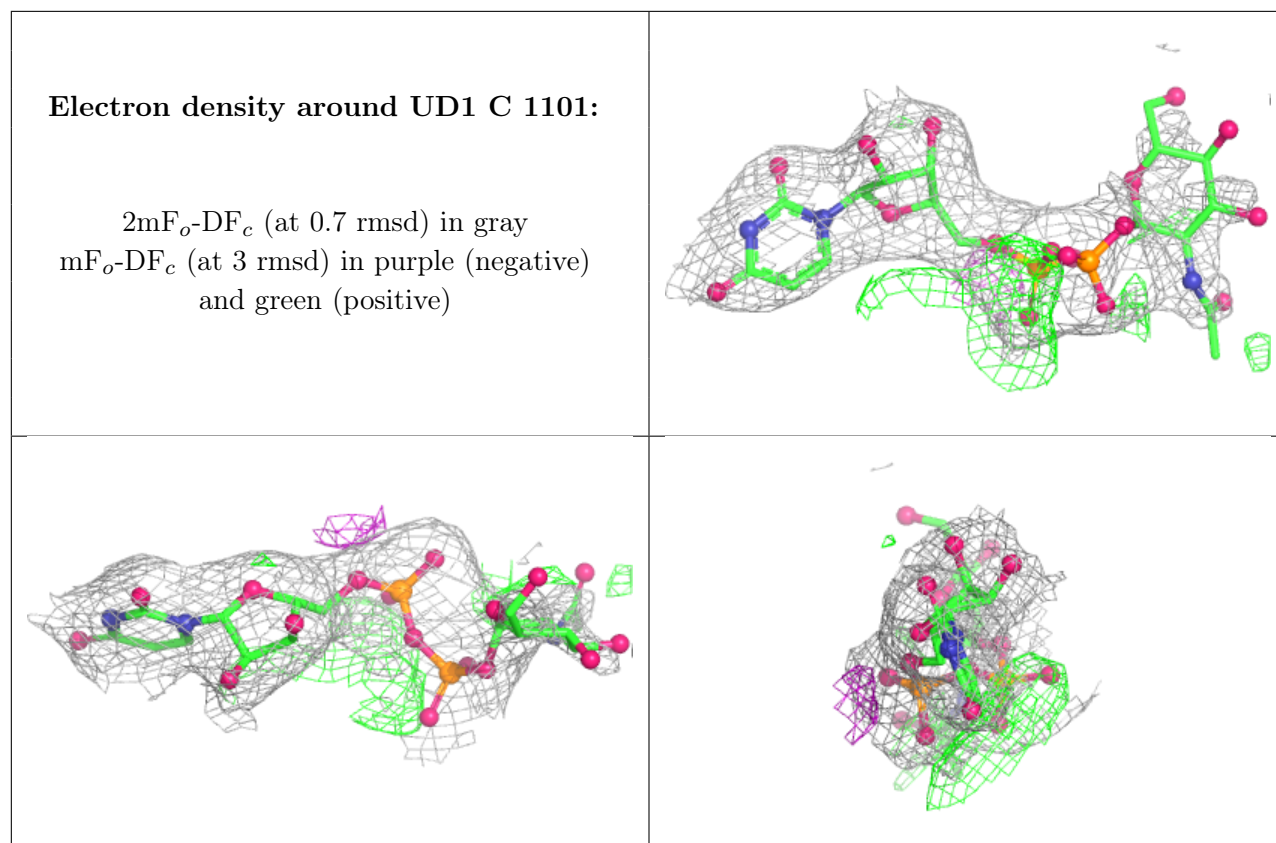
Electron density around UD1 G 1101:

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

**Electron density around UD1 E 1101:**

$2mF_o-DF_c$ (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.