



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:39 PM UTC

PDB ID : 6IJ9 / pdb_00006ij9
Title : Crystal Structure of Arabidopsis thaliana UGT89C1 complexed with UDP
Authors : Zong, G.N.; Wang, X.Q.
Deposited on : 2018-10-09
Resolution : 3.00 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

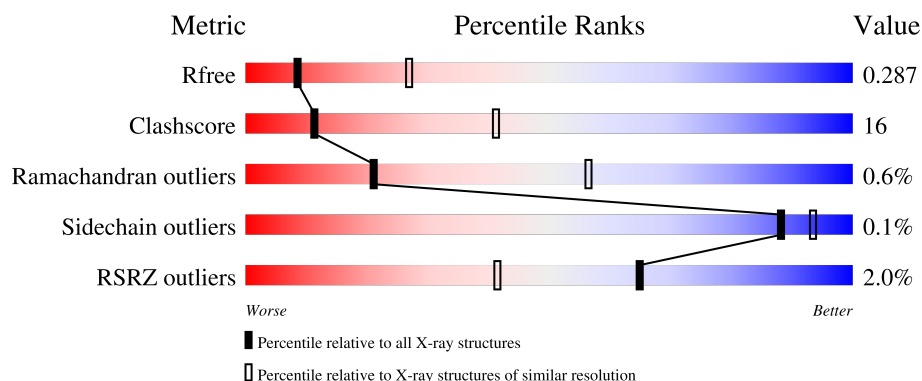
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	435	% 67% 26% • 6%
1	B	435	2% 63% 30% • 6%
1	C	435	% 65% 29% • 5%
1	D	435	3% 52% 37% • 11%

2 Entry composition [i](#)

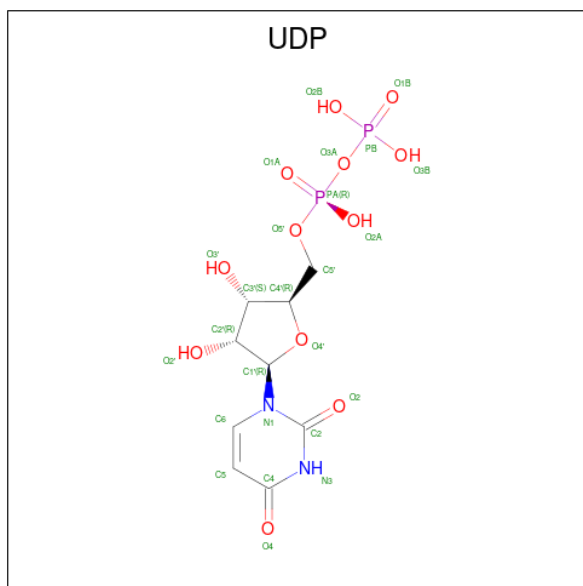
There are 3 unique types of molecules in this entry. The entry contains 12804 atoms, of which 0 are hydrogens and 0 are deuteriums.

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-glycosyltransferase 89C1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	410	Total	C	N	O	S	0	0	0
			3197	2051	549	584	13			
1	B	408	Total	C	N	O	S	0	0	0
			3174	2038	544	580	12			
1	C	413	Total	C	N	O	S	0	0	0
			3232	2074	554	591	13			
1	D	388	Total	C	N	O	S	0	0	0
			3040	1959	518	551	12			

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



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

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

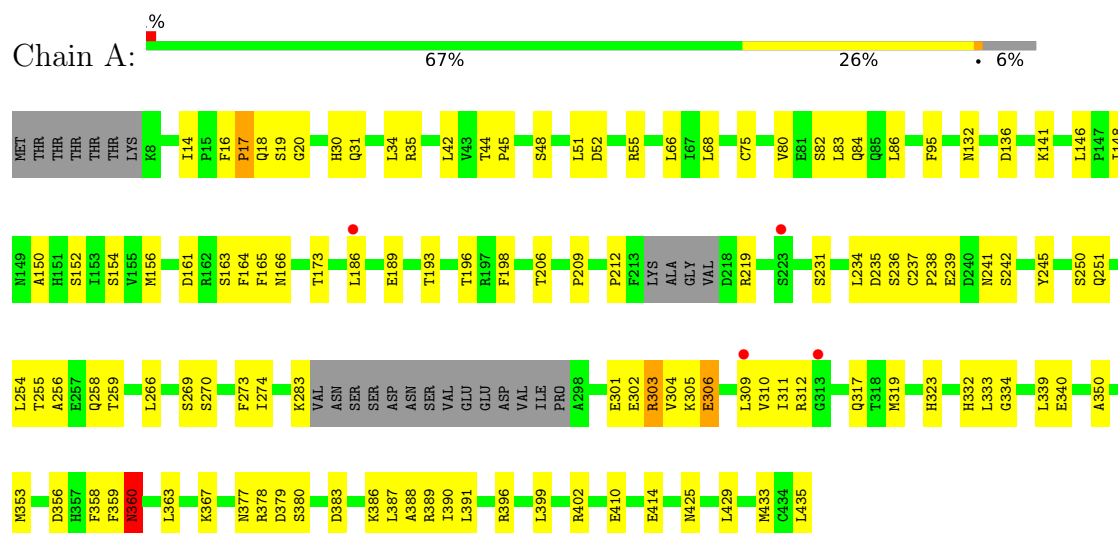
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	14	Total	O	0	0
			14	14		
3	B	18	Total	O	0	0
			18	18		
3	C	14	Total	O	0	0
			14	14		
3	D	15	Total	O	0	0
			15	15		

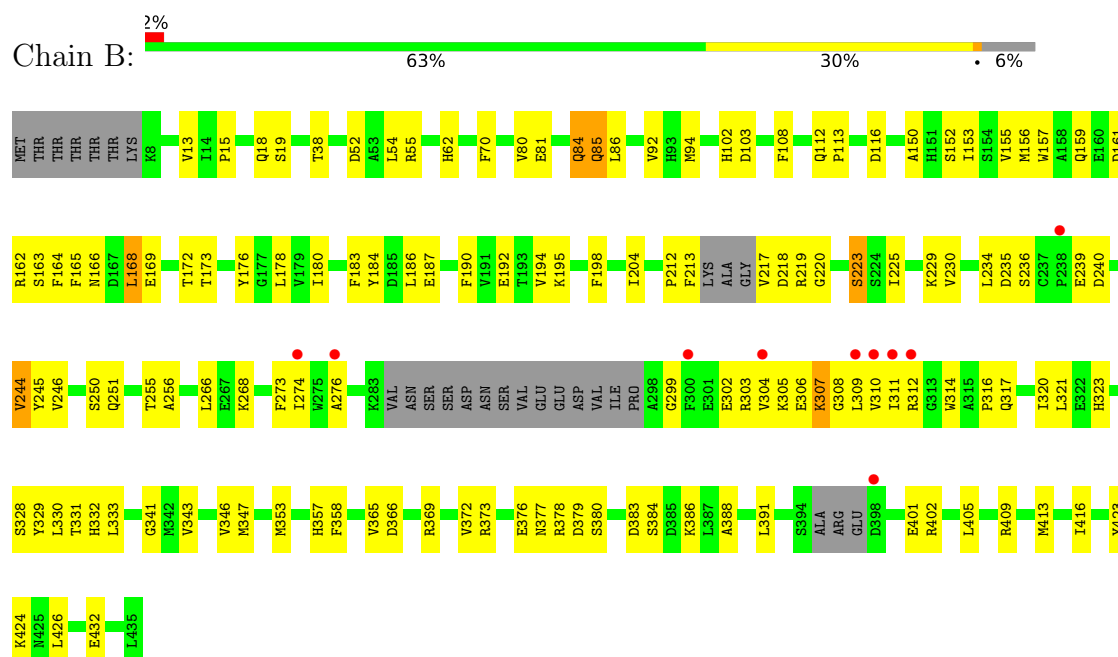
3 Residue-property plots [i](#)

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.

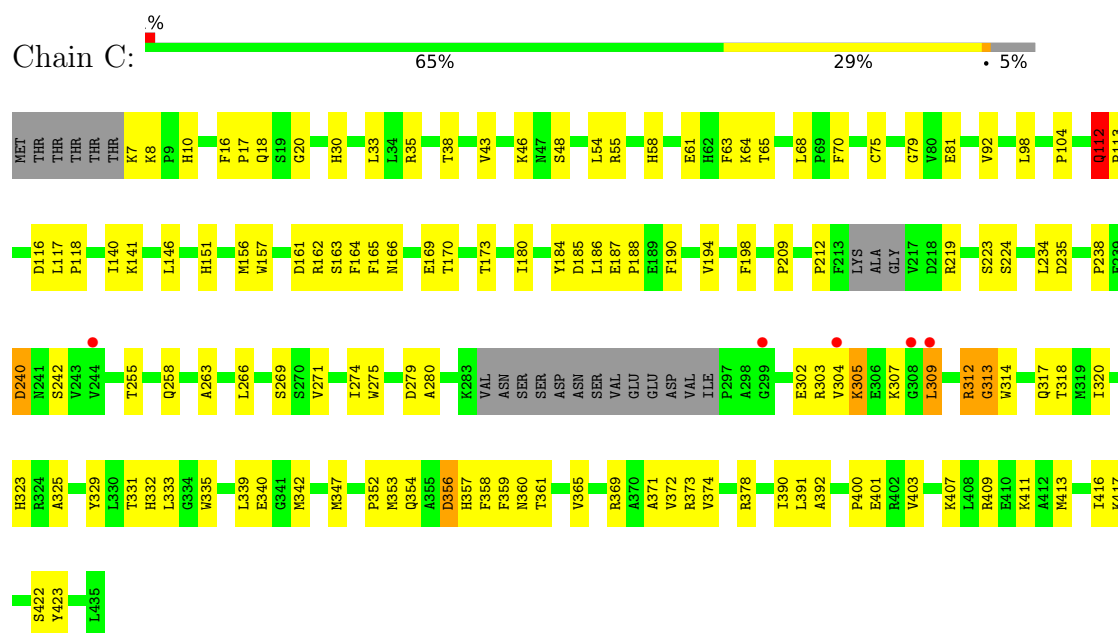
• Molecule 1: UDP-glycosyltransferase 89C1



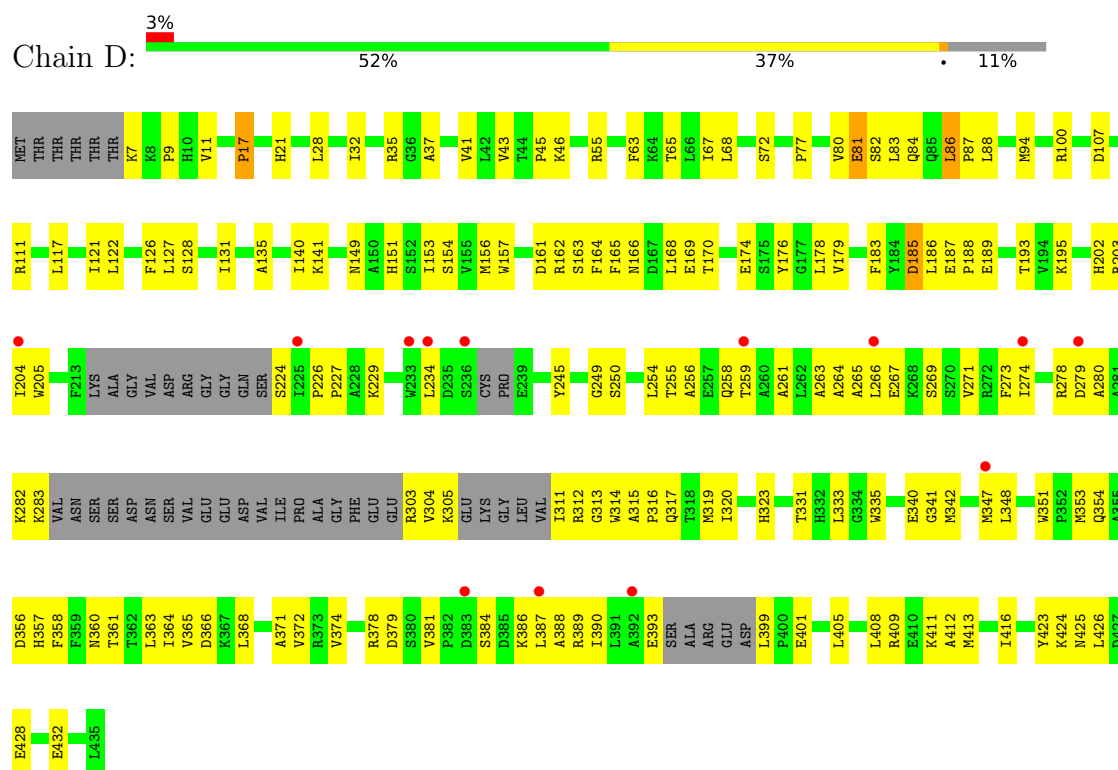
• Molecule 1: UDP-glycosyltransferase 89C1



• Molecule 1: UDP-glycosyltransferase 89C1



• Molecule 1: UDP-glycosyltransferase 89C1



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	92.98Å 84.14Å 129.15Å 90.00° 108.90° 90.00°	Depositor
Resolution (Å)	29.93 – 3.00 29.93 – 3.00	Depositor EDS
% Data completeness (in resolution range)	83.1 (29.93-3.00) 83.1 (29.93-3.00)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.86 (at 3.00Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.217 , 0.274 0.228 , 0.287	Depositor DCC
R_{free} test set	1575 reflections (4.14%)	wwPDB-VP
Wilson B-factor (Å ²)	43.4	Xtriage
Anisotropy	0.314	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 41.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.013 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	12804	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 37.04 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 4.5595e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical 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	0.61	0/3277	0.86	8/4458 (0.2%)
1	B	0.55	0/3253	0.98	7/4426 (0.2%)
1	C	0.64	2/3313 (0.1%)	0.98	9/4504 (0.2%)
1	D	0.57	0/3115	1.06	9/4235 (0.2%)
All	All	0.59	2/12958 (0.0%)	0.97	33/17623 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3
1	B	0	3
1	C	0	2
1	D	0	2
All	All	0	10

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	356	ASP	CB-CG	-6.14	1.36	1.52
1	C	112	GLN	CD-OE1	-5.46	1.13	1.23

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	425	ASN	CB-CG-ND2	-27.80	74.70	116.40
1	B	85	GLN	CG-CD-NE2	-27.64	74.94	116.40
1	D	425	ASN	OD1-CG-ND2	-26.57	96.03	122.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	112	GLN	OE1-CD-NE2	-23.91	98.69	122.60
1	B	85	GLN	OE1-CD-NE2	-23.73	98.87	122.60
1	C	112	GLN	CG-CD-NE2	-23.01	81.89	116.40
1	D	425	ASN	CB-CG-OD1	20.86	162.51	120.80
1	B	85	GLN	CG-CD-OE1	18.97	158.73	120.80
1	C	112	GLN	CG-CD-OE1	15.16	151.11	120.80
1	C	112	GLN	CB-CG-CD	11.77	132.61	112.60
1	A	359	PHE	CA-C-N	-10.85	101.88	122.06
1	A	359	PHE	C-N-CA	-10.85	101.88	122.06
1	C	112	GLN	CA-CB-CG	-10.66	92.78	114.10
1	A	360	ASN	N-CA-CB	9.23	124.95	110.46
1	A	360	ASN	OD1-CG-ND2	-8.55	114.05	122.60
1	D	425	ASN	CA-CB-CG	8.40	121.00	112.60
1	D	425	ASN	N-CA-CB	7.74	122.22	110.30
1	D	424	LYS	CA-C-N	-7.43	108.04	120.68
1	D	424	LYS	C-N-CA	-7.43	108.04	120.68
1	D	425	ASN	N-CA-C	-7.21	103.15	112.23
1	A	360	ASN	N-CA-C	-7.03	104.71	113.28
1	C	356	ASP	CB-CA-C	-6.47	97.08	109.95
1	A	360	ASN	CB-CG-OD1	-6.04	108.72	120.80
1	A	196	THR	CA-C-N	-6.03	110.03	121.54
1	A	196	THR	C-N-CA	-6.03	110.03	121.54
1	B	85	GLN	N-CA-C	5.93	119.48	112.72
1	D	86	LEU	CA-CB-CG	5.74	136.37	116.30
1	B	84	GLN	CA-C-N	5.56	131.69	122.73
1	B	84	GLN	C-N-CA	5.56	131.69	122.73
1	B	168	LEU	CA-CB-CG	5.24	134.65	116.30
1	C	313	GLY	N-CA-C	5.07	128.01	113.30
1	C	312	ARG	CA-C-N	5.01	130.71	121.70
1	C	312	ARG	C-N-CA	5.01	130.71	121.70

There are no chirality outliers.

All (10) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	303	ARG	Peptide
1	A	306	GLU	Peptide
1	A	360	ASN	Sidechain
1	B	307	LYS	Peptide
1	B	308	GLY	Peptide
1	B	85	GLN	Sidechain
1	C	112	GLN	Sidechain

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Mol	Chain	Res	Type	Group
1	C	309	LEU	Peptide
1	D	185	ASP	Peptide
1	D	81	GLU	Peptide

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	3197	0	3195	88	0
1	B	3174	0	3164	99	0
1	C	3232	0	3244	107	0
1	D	3040	0	3051	134	0
2	A	25	0	11	3	0
2	B	25	0	11	1	0
2	C	25	0	11	3	0
2	D	25	0	11	1	0
3	A	14	0	0	0	0
3	B	18	0	0	1	0
3	C	14	0	0	1	0
3	D	15	0	0	2	0
All	All	12804	0	12698	412	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 16.

All (412) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:112:GLN:HE21	1:C:112:GLN:HA	1.20	1.01
1:D:224:SER:N	3:D:1001:HOH:O	1.93	1.01
1:C:112:GLN:HG3	1:C:113:PRO:HD2	1.43	0.96
1:D:312:ARG:NH1	1:D:315:ALA:HA	1.88	0.89
1:B:166:ASN:ND2	1:D:379:ASP:OD2	2.08	0.87
1:B:379:ASP:OD2	1:D:166:ASN:ND2	2.09	0.85
1:C:219:ARG:HH11	1:C:318:THR:HG21	1.40	0.85
1:D:179:VAL:HA	1:D:205:TRP:HB2	1.59	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:227:PRO:HB3	1:D:319:MET:HE1	1.59	0.84
1:C:112:GLN:HE21	1:C:112:GLN:CA	1.91	0.82
1:D:282:LYS:HG2	1:D:283:LYS:H	1.43	0.80
1:D:17:PRO:O	1:D:82:SER:HB2	1.81	0.80
1:A:186:LEU:HD21	1:A:363:LEU:HD21	1.63	0.80
1:D:312:ARG:HH12	1:D:314:TRP:C	1.89	0.80
1:D:186:LEU:HB3	1:D:187:GLU:HG2	1.64	0.79
1:A:379:ASP:OD2	1:C:166:ASN:ND2	2.14	0.78
1:D:274:ILE:HD11	1:D:320:ILE:HD13	1.64	0.77
1:C:219:ARG:NH1	1:C:340:GLU:OE2	2.18	0.77
1:A:80:VAL:HG11	1:A:86:LEU:HD21	1.68	0.76
1:B:38:THR:HG23	1:B:62:HIS:HA	1.68	0.75
1:D:84:GLN:NE2	3:D:1002:HOH:O	2.19	0.75
1:A:301:GLU:HB3	1:A:302:GLU:HB2	1.68	0.74
1:C:17:PRO:HG2	1:C:98:LEU:HD11	1.70	0.73
1:D:46:LYS:HB2	1:D:81:GLU:HG2	1.70	0.73
1:A:353:MET:HG2	1:A:378:ARG:HA	1.71	0.73
1:B:373:ARG:HH21	1:B:376:GLU:HA	1.54	0.73
1:D:269:SER:HB2	1:D:388:ALA:HB1	1.71	0.72
1:A:254:LEU:HB3	1:A:258:GLN:HG3	1.72	0.72
1:B:244:VAL:HG22	1:B:328:SER:HB2	1.70	0.71
1:D:361:THR:O	1:D:365:VAL:HG22	1.90	0.71
1:D:226:PRO:HB2	1:D:229:LYS:HD3	1.72	0.70
1:D:255:THR:HG23	1:D:258:GLN:H	1.55	0.70
1:C:112:GLN:HG3	1:C:113:PRO:CD	2.18	0.70
1:B:369:ARG:HH12	1:C:369:ARG:HD2	1.56	0.70
1:C:331:THR:HG22	1:C:333:LEU:H	1.56	0.70
1:D:203:ARG:NH2	1:D:428:GLU:OE2	2.24	0.70
1:B:388:ALA:HA	1:B:391:LEU:HD12	1.74	0.69
1:C:417:LYS:HG2	1:C:423:TYR:HD2	1.58	0.68
1:D:195:LYS:HB2	1:D:204:ILE:HG23	1.75	0.68
1:C:55:ARG:HH21	1:C:65:THR:HG21	1.57	0.68
1:D:186:LEU:HB2	1:D:368:LEU:HD21	1.75	0.68
1:B:409:ARG:O	1:B:413:MET:HG2	1.93	0.68
1:D:341:GLY:HA3	1:D:348:LEU:HD21	1.75	0.68
1:D:353:MET:HG2	1:D:378:ARG:HA	1.74	0.68
1:B:358:PHE:CE2	1:D:161:ASP:HB2	2.29	0.68
1:B:161:ASP:HB2	1:D:358:PHE:CE2	2.29	0.68
1:A:206:THR:O	1:A:425:ASN:ND2	2.22	0.67
1:D:312:ARG:NH1	1:D:315:ALA:CA	2.57	0.67
1:A:219:ARG:NH2	1:A:340:GLU:OE2	2.23	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:256:ALA:O	1:D:259:THR:HG22	1.95	0.67
1:C:303:ARG:H	1:C:305:LYS:HG3	1.60	0.66
1:C:151:HIS:N	1:C:187:GLU:OE1	2.26	0.66
1:B:383:ASP:HB3	1:B:386:LYS:HG3	1.75	0.66
1:B:369:ARG:NH1	1:C:369:ARG:HD2	2.11	0.66
1:C:92:VAL:HG22	1:C:164:PHE:CE2	2.30	0.66
1:D:282:LYS:HG2	1:D:283:LYS:N	2.11	0.66
1:A:152:SER:O	1:A:156:MET:HG3	1.97	0.65
1:A:82:SER:HB3	1:A:84:GLN:OE1	1.97	0.65
1:B:239:GLU:HB3	1:B:240:ASP:HB3	1.77	0.65
1:B:353:MET:O	1:B:378:ARG:HB2	1.97	0.65
1:B:321:LEU:HD21	1:B:341:GLY:HA2	1.79	0.65
1:A:18:GLN:HG3	1:A:83:LEU:HD23	1.77	0.64
1:C:146:LEU:HB2	1:C:180:ILE:HG22	1.79	0.64
1:A:163:SER:HA	1:A:166:ASN:H	1.62	0.63
1:A:310:VAL:HG13	1:A:312:ARG:HH12	1.62	0.63
1:B:307:LYS:HA	1:B:309:LEU:HB3	1.80	0.63
1:C:173:THR:HG22	1:C:198:PHE:O	1.98	0.63
1:A:14:ILE:HG12	1:A:42:LEU:HB3	1.80	0.63
1:D:365:VAL:HG12	1:D:371:ALA:HB3	1.81	0.63
1:C:30:HIS:HB2	1:C:54:LEU:HD21	1.82	0.62
1:C:374:VAL:HG12	1:C:390:ILE:HD13	1.81	0.62
1:D:72:SER:O	1:D:100:ARG:NH2	2.33	0.62
1:B:54:LEU:HD13	1:B:213:PHE:CE2	2.34	0.62
1:C:353:MET:HG2	1:C:378:ARG:HA	1.82	0.62
1:B:401:GLU:OE2	1:C:369:ARG:NE	2.25	0.62
1:C:112:GLN:HG2	1:C:116:ASP:HB2	1.82	0.62
1:C:356:ASP:HB2	1:C:360:ASN:ND2	2.14	0.61
1:D:186:LEU:HA	1:D:188:PRO:HD3	1.81	0.61
1:B:372:VAL:HG23	1:B:402:ARG:HH21	1.65	0.60
1:D:162:ARG:HD2	1:D:169:GLU:OE1	2.00	0.60
1:C:10:HIS:HB3	1:C:118:PRO:HA	1.84	0.60
1:C:33:LEU:HD13	1:C:58:HIS:CD2	2.36	0.60
1:B:311:ILE:HG22	1:B:312:ARG:H	1.66	0.60
1:B:80:VAL:HG11	1:B:86:LEU:HD21	1.84	0.60
1:B:212:PRO:HG3	1:B:219:ARG:NE	2.16	0.59
1:D:46:LYS:HZ2	1:D:81:GLU:HA	1.68	0.59
1:C:164:PHE:HD1	1:C:164:PHE:H	1.50	0.59
1:B:276:ALA:HA	1:B:312:ARG:HG3	1.85	0.59
1:C:307:LYS:HA	1:C:309:LEU:HD13	1.84	0.59
1:B:230:VAL:HG21	1:B:312:ARG:HH12	1.68	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:223:SER:OG	3:C:1001:HOH:O	2.02	0.59
1:B:163:SER:OG	1:B:164:PHE:HA	2.04	0.58
1:C:400:PRO:O	1:C:403:VAL:HG12	2.03	0.58
1:D:365:VAL:HG23	1:D:366:ASP:OD1	2.04	0.58
1:A:378:ARG:NH1	1:C:163:SER:OG	2.37	0.57
1:A:18:GLN:NE2	1:A:251:GLN:OE1	2.33	0.57
1:D:408:LEU:HD23	1:D:411:LYS:HD3	1.86	0.57
1:C:186:LEU:HD21	1:C:335:TRP:HH2	1.68	0.57
1:D:28:LEU:HD21	1:D:122:LEU:HD21	1.87	0.57
1:D:163:SER:OG	1:D:164:PHE:HA	2.04	0.57
1:D:312:ARG:CZ	1:D:315:ALA:HA	2.35	0.57
1:C:303:ARG:C	1:C:305:LYS:H	2.12	0.57
1:D:312:ARG:NH1	1:D:314:TRP:C	2.63	0.56
1:A:356:ASP:O	1:A:360:ASN:ND2	2.39	0.56
1:B:347:MET:CB	1:B:402:ARG:HH11	2.19	0.56
1:D:149:ASN:HB3	1:D:335:TRP:CE2	2.41	0.56
1:B:184:TYR:OH	1:B:192:GLU:OE2	2.21	0.56
1:C:269:SER:HB2	1:C:392:ALA:HB2	1.87	0.56
1:C:329:TYR:CZ	1:C:331:THR:HG23	2.40	0.56
1:D:363:LEU:HD22	1:D:364:ILE:HD13	1.87	0.56
1:A:304:VAL:HA	1:A:306:GLU:OE1	2.06	0.55
1:A:389:ARG:HG3	1:A:390:ILE:N	2.20	0.55
1:C:170:THR:O	1:C:173:THR:OG1	2.16	0.55
1:C:331:THR:CG2	1:C:333:LEU:H	2.19	0.55
1:D:342:MET:HA	1:D:405:LEU:HD11	1.87	0.55
1:A:306:GLU:HB2	1:A:309:LEU:HG	1.88	0.55
1:B:223:SER:HA	1:B:314:TRP:HD1	1.70	0.55
1:A:340:GLU:OE1	2:A:900:UDP:O2'	2.21	0.55
1:C:372:VAL:HG11	1:C:390:ILE:HD11	1.89	0.55
1:A:161:ASP:HB2	1:C:358:PHE:CD2	2.41	0.55
1:B:299:GLY:O	3:B:1001:HOH:O	2.18	0.55
1:D:83:LEU:HB2	1:D:94:MET:HG3	1.89	0.55
1:D:358:PHE:O	1:D:361:THR:HG22	2.07	0.55
1:B:250:SER:OG	2:B:900:UDP:O3B	2.23	0.54
1:B:343:VAL:HA	1:B:409:ARG:HG3	1.88	0.54
1:C:407:LYS:HE2	1:C:411:LYS:HE3	1.89	0.54
1:B:168:LEU:O	1:B:172:THR:OG1	2.13	0.54
1:B:217:VAL:HG13	1:B:218:ASP:H	1.72	0.54
1:D:165:PHE:O	1:D:169:GLU:HG3	2.07	0.54
1:A:383:ASP:OD2	1:A:386:LYS:HG3	2.08	0.54
1:C:365:VAL:HG13	1:C:371:ALA:O	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:409:ARG:O	1:C:413:MET:HG2	2.07	0.54
1:A:396:ARG:HB3	1:A:399:LEU:HD13	1.90	0.54
1:B:235:ASP:OD2	1:B:323:HIS:NE2	2.39	0.54
1:B:92:VAL:HG22	1:B:164:PHE:CZ	2.43	0.54
1:D:279:ASP:OD1	1:D:280:ALA:N	2.32	0.54
1:A:16:PHE:O	1:A:18:GLN:N	2.37	0.54
1:C:185:ASP:HA	1:C:188:PRO:HG3	1.90	0.53
1:D:245:TYR:OH	1:D:317:GLN:HG2	2.07	0.53
1:A:250:SER:HB2	2:A:900:UDP:O3B	2.08	0.53
1:D:126:PHE:CD1	1:D:168:LEU:HD11	2.44	0.53
1:B:234:LEU:HD21	1:B:274:ILE:HD11	1.91	0.53
1:C:33:LEU:HD11	1:C:63:PHE:HD1	1.74	0.53
1:B:402:ARG:O	1:B:405:LEU:HB3	2.08	0.53
1:C:190:PHE:O	1:C:194:VAL:HG12	2.08	0.53
1:D:205:TRP:HH2	1:D:432:GLU:HG3	1.73	0.53
1:A:234:LEU:HB3	1:A:323:HIS:NE2	2.23	0.53
1:D:409:ARG:HH11	1:D:413:MET:HG3	1.74	0.53
1:C:416:ILE:HA	1:C:422:SER:HB2	1.91	0.53
1:A:66:LEU:HD11	1:A:68:LEU:HD21	1.91	0.52
1:D:121:ILE:HD12	1:D:135:ALA:HB2	1.91	0.52
1:D:412:ALA:O	1:D:416:ILE:HG12	2.09	0.52
1:B:377:ASN:N	1:B:380:SER:OG	2.30	0.52
1:D:164:PHE:HD1	1:D:164:PHE:H	1.56	0.52
1:A:163:SER:HA	1:A:165:PHE:N	2.24	0.52
1:B:195:LYS:HE3	1:B:204:ILE:HG22	1.91	0.52
1:D:265:ALA:HA	1:D:384:SER:OG	2.09	0.52
1:D:141:LYS:HD3	1:D:176:TYR:CZ	2.44	0.52
1:B:152:SER:O	1:B:156:MET:HG3	2.10	0.52
1:D:55:ARG:HG3	1:D:63:PHE:CE2	2.45	0.52
1:D:255:THR:HG22	1:D:258:GLN:HB2	1.92	0.52
1:A:173:THR:HG23	1:A:198:PHE:O	2.11	0.51
1:B:266:LEU:HD22	1:B:273:PHE:CG	2.45	0.51
1:C:279:ASP:OD1	1:C:280:ALA:N	2.40	0.51
1:D:264:ALA:HA	1:D:267:GLU:CD	2.35	0.51
1:B:372:VAL:CG2	1:B:402:ARG:HH21	2.23	0.51
1:C:7:LYS:HD3	1:C:35:ARG:O	2.11	0.51
1:A:239:GLU:O	1:A:242:SER:OG	2.20	0.51
1:B:102:HIS:ND1	1:B:103:ASP:OD1	2.39	0.51
1:B:268:LYS:NZ	1:B:384:SER:HB2	2.26	0.51
1:D:379:ASP:O	1:D:381:VAL:HG23	2.10	0.51
1:A:358:PHE:CE2	1:C:161:ASP:HB2	2.46	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:189:GLU:O	1:D:193:THR:OG1	2.24	0.51
1:D:195:LYS:O	1:D:202:HIS:ND1	2.44	0.51
1:B:332:HIS:HD2	1:B:357:HIS:NE2	2.10	0.50
1:C:163:SER:OG	1:C:164:PHE:HA	2.12	0.50
1:C:43:VAL:HG21	1:C:48:SER:HA	1.93	0.50
1:D:55:ARG:HG3	1:D:63:PHE:HE2	1.76	0.50
1:A:75:CYS:CB	1:C:75:CYS:HG	2.25	0.50
1:B:155:VAL:O	1:B:159:GLN:HG3	2.12	0.50
1:C:302:GLU:HB2	1:C:303:ARG:HG2	1.94	0.50
1:C:416:ILE:O	1:C:423:TYR:HB2	2.12	0.50
1:D:151:HIS:HB3	1:D:363:LEU:HD12	1.94	0.50
1:D:163:SER:HA	1:D:166:ASN:H	1.77	0.49
1:D:249:GLY:O	1:D:278:ARG:HD3	2.12	0.49
1:D:86:LEU:HB2	1:D:87:PRO:CD	2.42	0.49
1:B:153:ILE:HG23	1:B:157:TRP:CE3	2.48	0.49
1:A:377:ASN:H	1:A:380:SER:HB2	1.77	0.49
1:B:161:ASP:HB2	1:D:358:PHE:HE2	1.75	0.49
1:C:10:HIS:HD2	1:C:38:THR:HG22	1.77	0.49
1:D:170:THR:O	1:D:174:GLU:HG2	2.13	0.49
1:C:112:GLN:CA	1:C:112:GLN:NE2	2.67	0.49
1:D:356:ASP:OD2	1:D:360:ASN:ND2	2.46	0.49
1:A:311:ILE:O	1:A:312:ARG:HD3	2.13	0.49
1:C:212:PRO:HG3	1:C:219:ARG:CZ	2.42	0.49
1:D:9:PRO:HD2	1:D:37:ALA:HA	1.94	0.49
1:D:107:ASP:OD1	1:D:111:ARG:NH1	2.45	0.49
1:D:416:ILE:O	1:D:423:TYR:HB2	2.12	0.49
1:A:255:THR:HG22	1:A:256:ALA:N	2.28	0.49
1:C:234:LEU:HD11	1:C:320:ILE:HG23	1.94	0.49
1:C:335:TRP:NE1	1:C:339:LEU:HD11	2.28	0.49
1:D:186:LEU:HB3	1:D:187:GLU:CG	2.40	0.49
1:D:312:ARG:NH2	1:D:316:PRO:HD3	2.28	0.49
1:D:386:LYS:O	1:D:390:ILE:HG13	2.12	0.49
1:A:132:ASN:O	1:A:136:ASP:HB2	2.13	0.48
1:D:43:VAL:HG23	1:D:67:ILE:HG12	1.96	0.48
1:A:209:PRO:HG3	1:A:339:LEU:HD22	1.93	0.48
1:B:176:TYR:OH	1:B:432:GLU:OE1	2.31	0.48
1:D:254:LEU:HB3	1:D:258:GLN:HB3	1.93	0.48
1:D:86:LEU:HB2	1:D:87:PRO:HD2	1.94	0.48
1:D:282:LYS:O	1:D:283:LYS:HB2	2.13	0.48
1:B:52:ASP:OD1	1:B:55:ARG:NH1	2.46	0.48
1:B:165:PHE:HA	1:B:168:LEU:HD13	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:20:GLY:HA3	2:A:900:UDP:O1B	2.13	0.48
1:C:46:LYS:HE3	1:C:79:GLY:O	2.13	0.48
1:B:306:GLU:O	1:B:309:LEU:HD13	2.14	0.48
1:D:255:THR:CG2	1:D:258:GLN:HB2	2.44	0.48
1:A:30:HIS:HE1	1:A:34:LEU:HD21	1.79	0.48
1:D:88:LEU:HD22	1:D:354:GLN:OE1	2.14	0.47
1:A:254:LEU:HB2	1:A:259:THR:HG23	1.95	0.47
1:C:274:ILE:HG23	1:C:312:ARG:HH12	1.79	0.47
1:D:426:LEU:HA	1:D:426:LEU:HD23	1.64	0.47
1:A:173:THR:HG23	1:A:198:PHE:C	2.38	0.47
1:A:164:PHE:HD1	1:A:164:PHE:H	1.63	0.47
1:B:150:ALA:N	1:B:187:GLU:OE2	2.43	0.47
1:A:274:ILE:HG12	1:A:310:VAL:HG12	1.96	0.47
1:D:203:ARG:HD2	1:D:203:ARG:HA	1.69	0.47
1:B:274:ILE:HG23	1:B:310:VAL:HB	1.96	0.47
1:B:347:MET:HA	1:B:402:ARG:HD2	1.97	0.47
1:C:61:GLU:O	1:C:64:LYS:NZ	2.47	0.47
1:A:333:LEU:HD13	1:A:350:ALA:HB2	1.95	0.47
1:A:212:PRO:HB3	1:A:219:ARG:CD	2.45	0.47
1:A:234:LEU:HB3	1:A:323:HIS:CD2	2.50	0.47
1:B:163:SER:HB3	1:D:378:ARG:NH1	2.29	0.47
1:A:245:TYR:OH	1:A:317:GLN:HG2	2.15	0.47
1:B:316:PRO:HD2	1:B:320:ILE:HD11	1.97	0.47
1:B:331:THR:HG22	1:B:333:LEU:H	1.79	0.47
1:C:347:MET:HE1	1:C:401:GLU:HB2	1.96	0.47
1:B:212:PRO:HG3	1:B:219:ARG:HE	1.80	0.46
1:A:410:GLU:O	1:A:414:GLU:HG2	2.15	0.46
1:D:65:THR:HG22	1:D:67:ILE:HG13	1.96	0.46
1:D:11:VAL:HG21	1:D:32:ILE:HD12	1.96	0.46
1:D:258:GLN:HE22	1:D:379:ASP:HA	1.81	0.46
1:A:19:SER:OG	1:A:84:GLN:HG2	2.16	0.46
1:D:304:VAL:HG13	1:D:305:LYS:H	1.80	0.46
1:A:241:ASN:HD21	1:A:402:ARG:HH22	1.64	0.46
1:A:266:LEU:HD22	1:A:273:PHE:CG	2.50	0.46
1:B:18:GLN:NE2	1:B:251:GLN:HG2	2.31	0.46
1:C:16:PHE:O	1:C:18:GLN:N	2.46	0.46
1:D:84:GLN:O	1:D:84:GLN:HG2	2.16	0.46
1:A:274:ILE:HG23	1:A:310:VAL:HG13	1.96	0.46
1:A:383:ASP:HB3	1:A:386:LYS:HB2	1.98	0.46
1:C:240:ASP:O	1:C:325:ALA:HA	2.16	0.46
1:D:409:ARG:NH1	1:D:413:MET:HG3	2.30	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:424:LYS:HD2	1:B:424:LYS:HA	1.64	0.45
1:C:117:LEU:HD23	1:C:117:LEU:HA	1.82	0.45
1:C:186:LEU:HD21	1:C:335:TRP:CH2	2.50	0.45
1:D:311:ILE:O	1:D:312:ARG:HG2	2.15	0.45
1:A:150:ALA:O	1:A:154:SER:OG	2.33	0.45
1:B:255:THR:HG22	1:B:256:ALA:N	2.32	0.45
1:C:255:THR:HG23	1:C:258:GLN:OE1	2.16	0.45
1:C:184:TYR:HD2	1:C:185:ASP:OD1	1.99	0.45
1:C:317:GLN:OE1	2:C:900:UDP:H2'	2.16	0.45
1:D:153:ILE:HG23	1:D:157:TRP:CE3	2.51	0.45
1:D:303:ARG:HA	1:D:304:VAL:HA	1.43	0.45
1:C:312:ARG:HB3	1:C:313:GLY:HA3	1.99	0.45
1:B:183:PHE:CZ	1:B:186:LEU:HB2	2.51	0.45
1:B:190:PHE:O	1:B:194:VAL:HG23	2.16	0.45
1:B:353:MET:HG2	1:B:378:ARG:HA	1.98	0.45
1:D:156:MET:HE3	1:D:156:MET:HB3	1.86	0.45
1:C:162:ARG:HD2	1:C:169:GLU:OE1	2.17	0.45
1:D:185:ASP:HB2	1:D:186:LEU:HD23	1.99	0.45
1:A:273:PHE:O	1:A:309:LEU:HA	2.17	0.44
1:A:332:HIS:CE1	1:A:334:GLY:HA3	2.52	0.44
1:B:245:TYR:OH	1:B:317:GLN:HG2	2.18	0.44
1:A:48:SER:O	1:A:51:LEU:HB2	2.17	0.44
1:A:212:PRO:HB3	1:A:219:ARG:HD3	1.99	0.44
1:B:173:THR:HG23	1:B:198:PHE:O	2.17	0.44
1:C:7:LYS:HA	1:C:8:LYS:HA	1.56	0.44
1:A:311:ILE:HG22	1:A:312:ARG:H	1.83	0.44
1:A:141:LYS:NZ	1:A:435:LEU:O	2.49	0.44
1:B:266:LEU:HB3	1:B:273:PHE:CE1	2.52	0.44
1:B:18:GLN:HE21	1:B:251:GLN:HG2	1.83	0.44
1:C:417:LYS:HG2	1:C:423:TYR:CD2	2.46	0.44
1:D:46:LYS:NZ	1:D:81:GLU:HA	2.33	0.44
1:B:19:SER:OG	1:B:84:GLN:CD	2.61	0.44
1:C:113:PRO:O	1:C:116:ASP:N	2.49	0.44
1:C:342:MET:HE2	1:C:342:MET:HB3	1.95	0.44
1:D:7:LYS:HB3	1:D:35:ARG:O	2.18	0.44
1:C:140:ILE:HG22	1:C:141:LYS:O	2.18	0.44
1:C:223:SER:HA	1:C:314:TRP:O	2.18	0.44
1:C:307:LYS:HA	1:C:309:LEU:HB2	2.00	0.44
1:D:261:ALA:O	1:D:264:ALA:HB3	2.18	0.44
1:A:52:ASP:HA	1:A:55:ARG:NH1	2.33	0.44
1:C:20:GLY:HA2	2:C:900:UDP:H5'1	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:367:LYS:HE2	1:A:367:LYS:HB3	1.87	0.43
1:B:358:PHE:CD2	1:D:161:ASP:HB2	2.52	0.43
1:C:303:ARG:C	1:C:305:LYS:N	2.76	0.43
1:D:387:LEU:HD12	1:D:390:ILE:HD12	2.00	0.43
1:B:70:PHE:CD2	1:B:81:GLU:HB3	2.54	0.43
1:B:108:PHE:O	1:B:112:GLN:HG2	2.18	0.43
1:B:310:VAL:HG12	1:B:311:ILE:H	1.83	0.43
1:A:283:LYS:NZ	1:A:283:LYS:HB3	2.33	0.43
1:D:88:LEU:O	1:D:88:LEU:HG	2.17	0.43
1:A:273:PHE:CZ	1:A:309:LEU:HD23	2.54	0.43
1:C:219:ARG:NH1	1:C:318:THR:HG21	2.21	0.43
1:D:127:LEU:O	1:D:131:ILE:HG13	2.18	0.43
1:C:307:LYS:C	1:C:309:LEU:HB2	2.43	0.43
1:D:21:HIS:CE1	1:D:126:PHE:CE2	3.06	0.43
1:D:128:SER:HA	1:D:131:ILE:HB	2.00	0.43
1:A:16:PHE:CD1	1:A:17:PRO:HD2	2.53	0.43
1:B:268:LYS:HZ2	1:B:384:SER:HB2	1.83	0.43
1:A:387:LEU:O	1:A:391:LEU:HG	2.18	0.43
1:B:113:PRO:O	1:B:116:ASP:N	2.51	0.43
1:C:169:GLU:O	1:C:173:THR:HG23	2.19	0.43
1:C:352:PRO:HA	1:C:357:HIS:HB3	2.01	0.43
1:D:413:MET:SD	1:D:416:ILE:HD11	2.59	0.43
1:B:239:GLU:HA	1:B:240:ASP:HA	1.56	0.43
1:C:224:SER:OG	1:C:313:GLY:HA3	2.18	0.43
1:D:340:GLU:OE2	2:D:900:UDP:O3'	2.26	0.43
1:B:212:PRO:HG3	1:B:219:ARG:CZ	2.49	0.43
1:B:246:VAL:HG13	1:B:330:LEU:HD23	2.00	0.43
1:D:351:TRP:CD1	1:D:351:TRP:C	2.93	0.43
1:A:270:SER:O	1:A:270:SER:OG	2.35	0.42
1:A:332:HIS:CD2	1:A:332:HIS:H	2.36	0.42
1:B:304:VAL:HG11	1:B:309:LEU:HD22	2.01	0.42
1:A:146:LEU:HD13	1:A:148:ILE:HD11	2.01	0.42
1:B:321:LEU:CD2	1:B:341:GLY:HA2	2.49	0.42
1:B:426:LEU:HD23	1:B:426:LEU:HA	1.89	0.42
1:C:235:ASP:OD1	1:C:323:HIS:NE2	2.47	0.42
1:A:31:GLN:OE1	1:A:35:ARG:NH2	2.48	0.42
1:D:45:PRO:HD3	1:D:68:LEU:O	2.19	0.42
1:D:378:ARG:HG3	1:D:379:ASP:CG	2.45	0.42
1:B:307:LYS:O	1:B:307:LYS:HG3	2.19	0.42
1:C:55:ARG:NH2	1:C:65:THR:HG21	2.30	0.42
1:C:332:HIS:HD2	1:C:357:HIS:NE2	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:341:GLY:O	1:B:346:VAL:HG22	2.20	0.42
1:C:240:ASP:O	1:C:242:SER:N	2.49	0.42
1:D:117:LEU:HD23	1:D:117:LEU:HA	1.91	0.42
1:B:373:ARG:HH21	1:B:376:GLU:CA	2.29	0.42
1:C:354:GLN:O	1:C:357:HIS:CG	2.73	0.42
1:A:303:ARG:HB2	1:A:305:LYS:HG2	2.01	0.42
1:D:77:PRO:O	1:D:80:VAL:HG12	2.20	0.42
1:A:16:PHE:C	1:A:18:GLN:H	2.27	0.42
1:A:234:LEU:O	1:A:237:CYS:N	2.42	0.42
1:A:304:VAL:HG23	1:A:304:VAL:O	2.20	0.42
1:B:372:VAL:HG23	1:B:402:ARG:NH2	2.32	0.42
1:D:347:MET:HE1	1:D:401:GLU:HB2	2.02	0.42
1:B:162:ARG:HD2	1:B:169:GLU:OE1	2.20	0.41
1:B:329:TYR:CZ	1:B:331:THR:HG23	2.55	0.41
1:C:98:LEU:HD23	1:C:98:LEU:HA	1.73	0.41
1:C:209:PRO:HD2	1:C:416:ILE:HG23	2.01	0.41
1:B:13:VAL:HG12	1:B:15:PRO:HD3	2.03	0.41
1:B:156:MET:HE2	1:B:156:MET:HB3	1.81	0.41
1:C:266:LEU:HG	1:C:391:LEU:HD21	2.02	0.41
1:D:250:SER:O	1:D:278:ARG:NE	2.53	0.41
1:A:193:THR:HG22	1:D:189:GLU:OE1	2.20	0.41
1:B:178:LEU:HG	1:B:180:ILE:HG23	2.01	0.41
1:B:225:ILE:HG23	1:B:229:LYS:NZ	2.36	0.41
1:B:365:VAL:HG12	1:B:366:ASP:OD1	2.20	0.41
1:D:46:LYS:HD2	1:D:81:GLU:HA	2.02	0.41
1:D:372:VAL:HG23	1:D:374:VAL:HG13	2.03	0.41
1:D:393:GLU:OE2	1:D:399:LEU:HD21	2.20	0.41
1:C:68:LEU:HD13	1:C:104:PRO:HB2	2.03	0.41
1:C:263:ALA:HB2	1:C:275:TRP:CZ3	2.55	0.41
1:D:185:ASP:HB3	1:D:188:PRO:HG3	2.01	0.41
1:A:44:THR:HB	1:A:45:PRO:CD	2.50	0.41
1:C:70:PHE:CD2	1:C:81:GLU:HB3	2.56	0.41
1:D:41:VAL:HG12	1:D:43:VAL:HG13	2.02	0.41
1:D:140:ILE:HD13	1:D:140:ILE:HG21	1.75	0.41
1:A:163:SER:HB3	1:A:166:ASN:HB2	2.02	0.41
1:A:269:SER:HB3	1:A:388:ALA:O	2.21	0.41
1:C:156:MET:HE2	1:C:359:PHE:HE2	1.85	0.41
1:C:157:TRP:CZ2	1:C:169:GLU:HA	2.56	0.41
1:C:340:GLU:OE2	2:C:900:UDP:O3'	2.30	0.41
1:D:354:GLN:O	1:D:357:HIS:CG	2.74	0.41
1:D:384:SER:OG	1:D:384:SER:O	2.14	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:386:LYS:O	1:D:389:ARG:HG2	2.21	0.41
1:A:189:GLU:O	1:A:193:THR:HG23	2.20	0.41
1:A:231:SER:OG	1:A:319:MET:SD	2.79	0.41
1:A:235:ASP:HA	1:A:236:SER:HA	1.55	0.41
1:A:238:PRO:HB2	1:A:239:GLU:HG3	2.02	0.41
1:C:54:LEU:HD23	1:C:54:LEU:HA	1.80	0.41
1:C:302:GLU:HA	1:C:303:ARG:HA	1.46	0.41
1:A:429:LEU:O	1:A:433:MET:HG3	2.20	0.41
1:D:263:ALA:HA	1:D:266:LEU:HD13	2.02	0.41
1:D:331:THR:HG23	1:D:333:LEU:H	1.85	0.41
1:B:255:THR:HG22	1:B:256:ALA:H	1.86	0.41
1:D:46:LYS:H	1:D:46:LYS:HG2	1.40	0.41
1:D:224:SER:HB3	1:D:314:TRP:O	2.21	0.41
1:A:16:PHE:CE1	1:A:95:PHE:HE1	2.39	0.40
1:C:361:THR:HG21	1:C:373:ARG:HD2	2.03	0.40
1:D:178:LEU:O	1:D:205:TRP:HD1	2.04	0.40
1:D:183:PHE:CD2	1:D:185:ASP:O	2.74	0.40
1:A:360:ASN:HD22	1:A:360:ASN:N	2.20	0.40
1:A:377:ASN:H	1:A:380:SER:CB	2.33	0.40
1:B:235:ASP:HA	1:B:236:SER:HA	1.68	0.40
1:C:271:VAL:CG1	1:C:392:ALA:HA	2.50	0.40
1:D:186:LEU:HA	1:D:187:GLU:HA	1.80	0.40
1:D:195:LYS:HB2	1:D:204:ILE:CG2	2.48	0.40
1:C:55:ARG:HH11	1:C:55:ARG:HD2	1.74	0.40
1:C:356:ASP:O	1:C:359:PHE:N	2.54	0.40
1:D:151:HIS:O	1:D:154:SER:OG	2.38	0.40
1:A:358:PHE:C	1:A:360:ASN:N	2.78	0.40
1:B:94:MET:HE2	1:B:94:MET:HB2	1.86	0.40
1:B:220:GLY:HA3	1:B:314:TRP:HZ2	1.86	0.40
1:C:165:PHE:O	1:C:169:GLU:HG3	2.22	0.40
1:D:234:LEU:HD23	1:D:323:HIS:CG	2.55	0.40
1:D:271:VAL:O	1:D:273:PHE:HD1	2.04	0.40
1:B:302:GLU:O	1:B:303:ARG:C	2.65	0.40
1:B:416:ILE:O	1:B:423:TYR:HB2	2.22	0.40
1:D:354:GLN:HE21	1:D:354:GLN:HB3	1.51	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	404/435 (93%)	382 (95%)	21 (5%)	1 (0%)	43	76
1	B	400/435 (92%)	378 (94%)	20 (5%)	2 (0%)	24	60
1	C	407/435 (94%)	380 (93%)	23 (6%)	4 (1%)	12	45
1	D	376/435 (86%)	355 (94%)	19 (5%)	2 (0%)	24	60
All	All	1587/1740 (91%)	1495 (94%)	83 (5%)	9 (1%)	21	56

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	304	VAL
1	B	305	LYS
1	C	240	ASP
1	C	238	PRO
1	C	305	LYS
1	D	313	GLY
1	B	223	SER
1	A	17	PRO
1	D	17	PRO

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	351/377 (93%)	351 (100%)	0	100	100
1	B	348/377 (92%)	347 (100%)	1 (0%)	86	91

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	357/377 (95%)	357 (100%)	0	100	100
1	D	336/377 (89%)	336 (100%)	0	100	100
All	All	1392/1508 (92%)	1391 (100%)	1 (0%)	88	93

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	244	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	360	ASN
1	B	84	GLN
1	B	251	GLN
1	B	377	ASN
1	C	85	GLN
1	C	112	GLN
1	C	360	ASN
1	D	354	GLN
1	D	425	ASN

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	$\# Z > 2$
2	UDP	C	900	-	25,26,26	1.19	1 (4%)	38,40,40	0.73	1 (2%)
2	UDP	A	900	-	25,26,26	0.43	0	38,40,40	0.48	0
2	UDP	B	900	-	25,26,26	0.54	0	38,40,40	0.47	0
2	UDP	D	900	-	25,26,26	0.85	1 (4%)	38,40,40	0.76	1 (2%)

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	C	900	-	-	7/16/32/32	0/2/2/2
2	UDP	A	900	-	-	3/16/32/32	0/2/2/2
2	UDP	B	900	-	-	3/16/32/32	0/2/2/2
2	UDP	D	900	-	-	0/16/32/32	0/2/2/2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	900	UDP	PA-O3A	-5.03	1.54	1.59
2	D	900	UDP	PA-O3A	3.64	1.63	1.59

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	900	UDP	O3B-PB-O3A	2.30	112.36	104.64
2	D	900	UDP	O3B-PB-O3A	2.30	112.33	104.64

There are no chirality outliers.

All (13) torsion outliers are listed below:

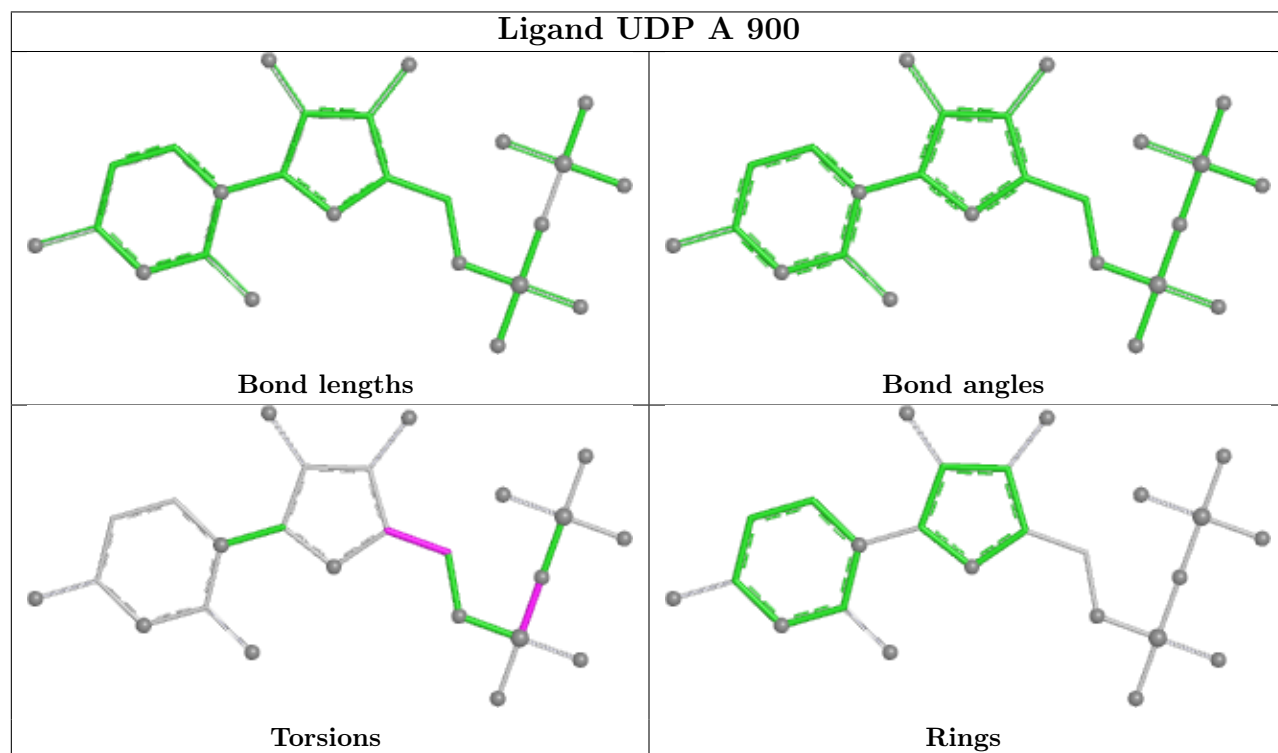
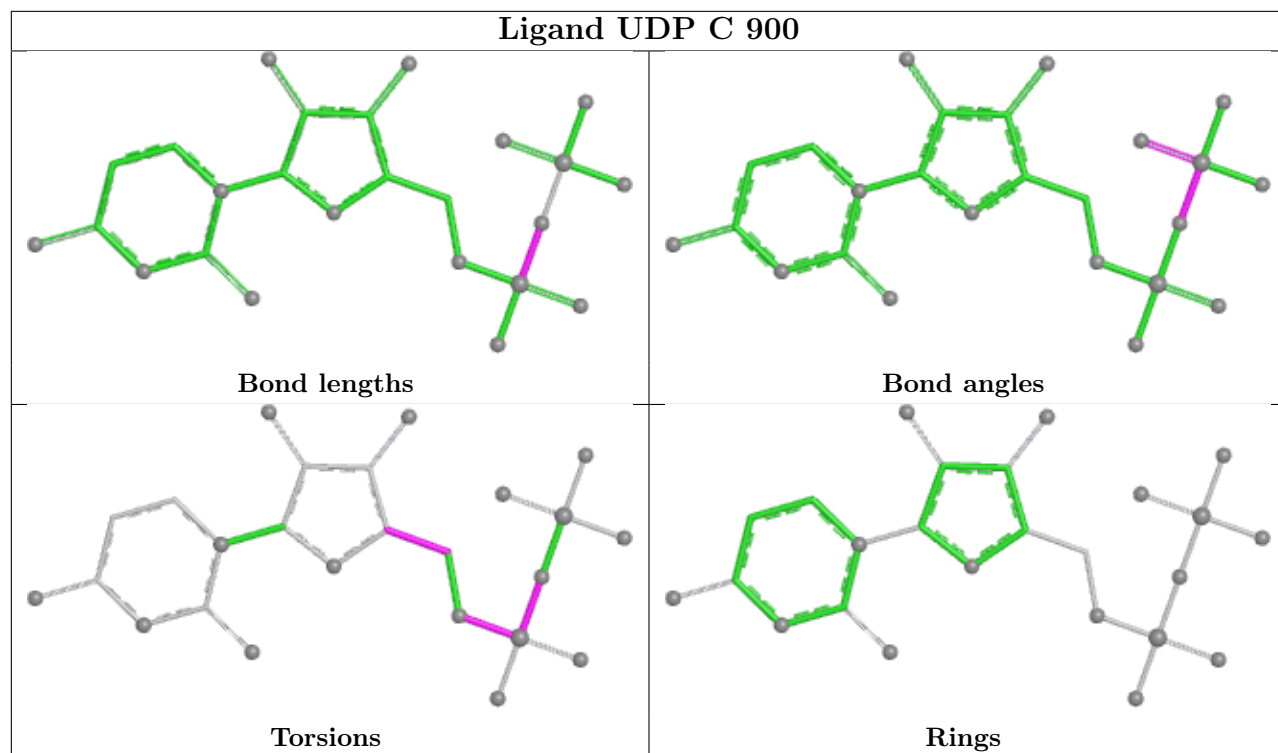
Mol	Chain	Res	Type	Atoms
2	B	900	UDP	O4'-C4'-C5'-O5'
2	C	900	UDP	C3'-C4'-C5'-O5'
2	C	900	UDP	O4'-C4'-C5'-O5'
2	B	900	UDP	C3'-C4'-C5'-O5'
2	A	900	UDP	O4'-C4'-C5'-O5'
2	A	900	UDP	C3'-C4'-C5'-O5'
2	A	900	UDP	PB-O3A-PA-O5'
2	B	900	UDP	PB-O3A-PA-O5'
2	C	900	UDP	C5'-O5'-PA-O1A
2	C	900	UDP	C5'-O5'-PA-O2A
2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

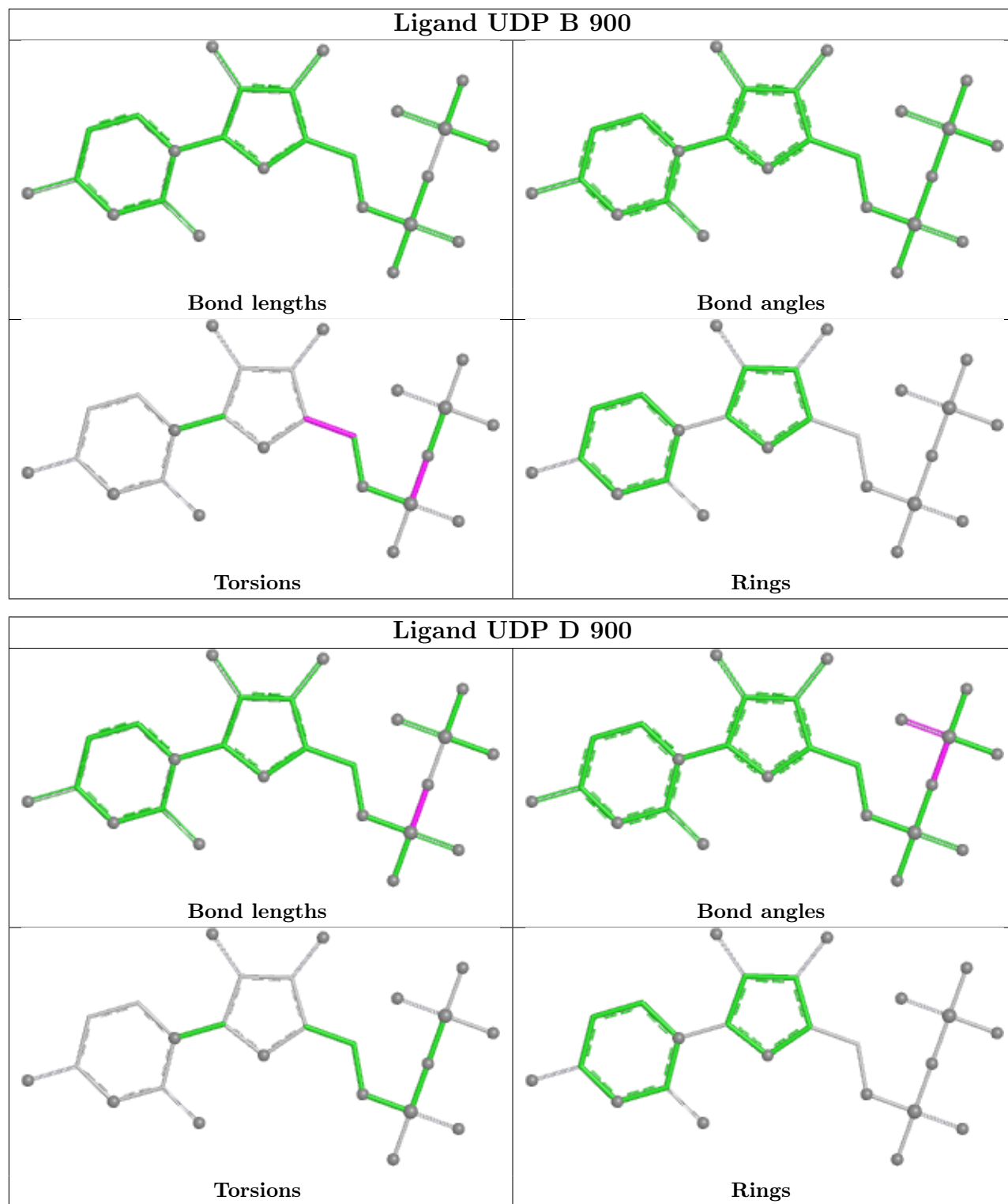
There are no ring outliers.

4 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	900	UDP	3	0
2	A	900	UDP	3	0
2	B	900	UDP	1	0
2	D	900	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.





5.7 Other polymers [i](#)

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	410/435 (94%)	-0.11	4 (0%) 79 59	17, 30, 58, 70	0
1	B	408/435 (93%)	0.06	10 (2%) 58 35	15, 39, 82, 96	0
1	C	413/435 (94%)	-0.05	5 (1%) 76 55	14, 33, 73, 90	0
1	D	388/435 (89%)	0.24	13 (3%) 48 27	21, 42, 87, 102	0
All	All	1619/1740 (93%)	0.03	32 (1%) 65 41	14, 36, 79, 102	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	309	LEU	4.0
1	D	204	ILE	3.5
1	B	311	ILE	3.4
1	B	309	LEU	3.2
1	D	387	LEU	3.1
1	A	223	SER	2.9
1	C	244	VAL	2.9
1	B	398	ASP	2.9
1	B	276	ALA	2.8
1	A	186	LEU	2.7
1	C	304	VAL	2.7
1	D	236	SER	2.6
1	D	225	ILE	2.4
1	D	274	ILE	2.4
1	B	304	VAL	2.4
1	C	309	LEU	2.4
1	B	312	ARG	2.3
1	D	266	LEU	2.3
1	C	299	GLY	2.3
1	D	392	ALA	2.3
1	B	238	PRO	2.3

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Mol	Chain	Res	Type	RSRZ
1	D	233	TRP	2.2
1	C	308	GLY	2.2
1	D	259	THR	2.2
1	B	274	ILE	2.2
1	D	279	ASP	2.1
1	B	310	VAL	2.1
1	D	347	MET	2.1
1	B	300	PHE	2.1
1	A	313	GLY	2.1
1	D	234	LEU	2.0
1	D	383	ASP	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

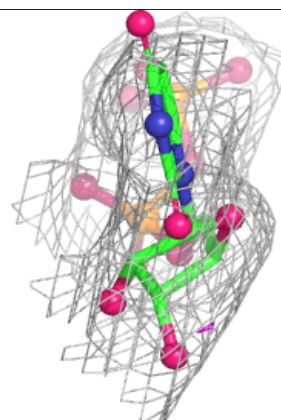
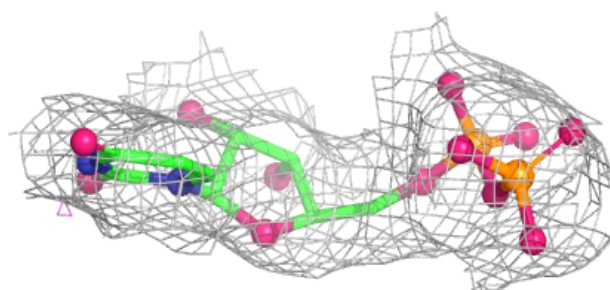
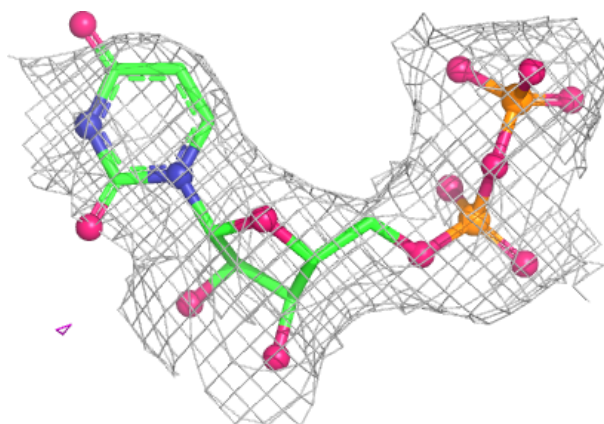
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	D	900	25/25	0.93	0.09	39,51,62,64	0
2	UDP	A	900	25/25	0.94	0.08	18,30,41,43	0
2	UDP	B	900	25/25	0.95	0.09	30,43,53,55	0
2	UDP	C	900	25/25	0.96	0.08	21,34,44,46	0

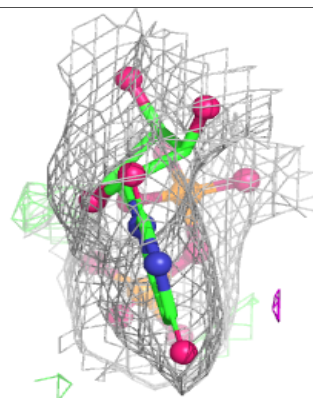
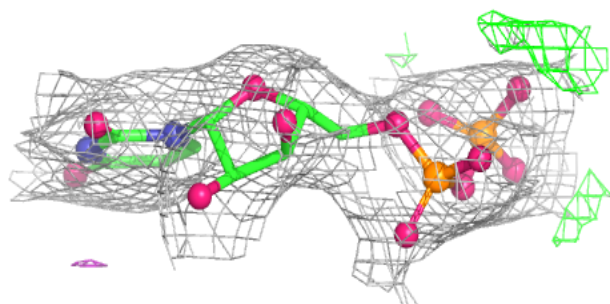
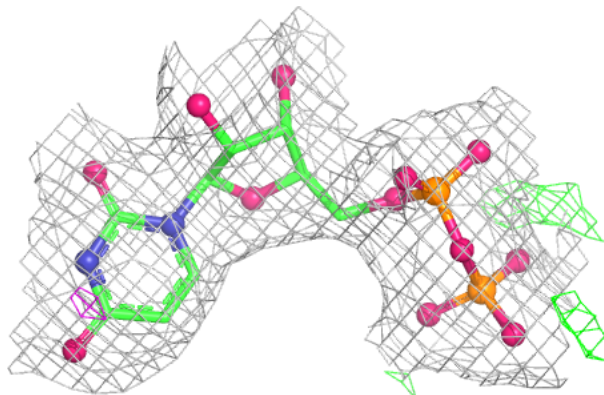
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around UDP D 900:

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

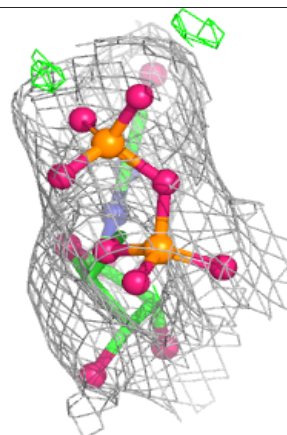
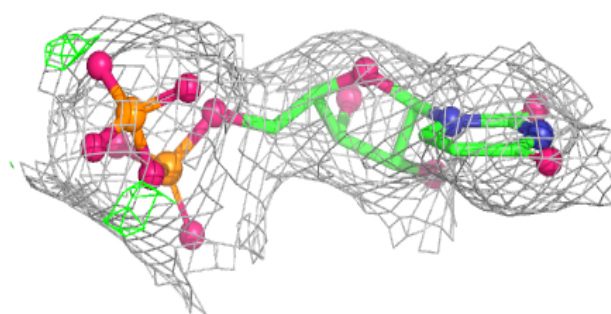
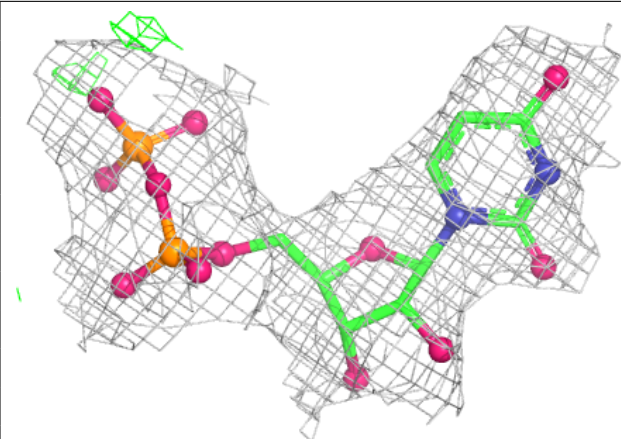
**Electron density around UDP A 900:**

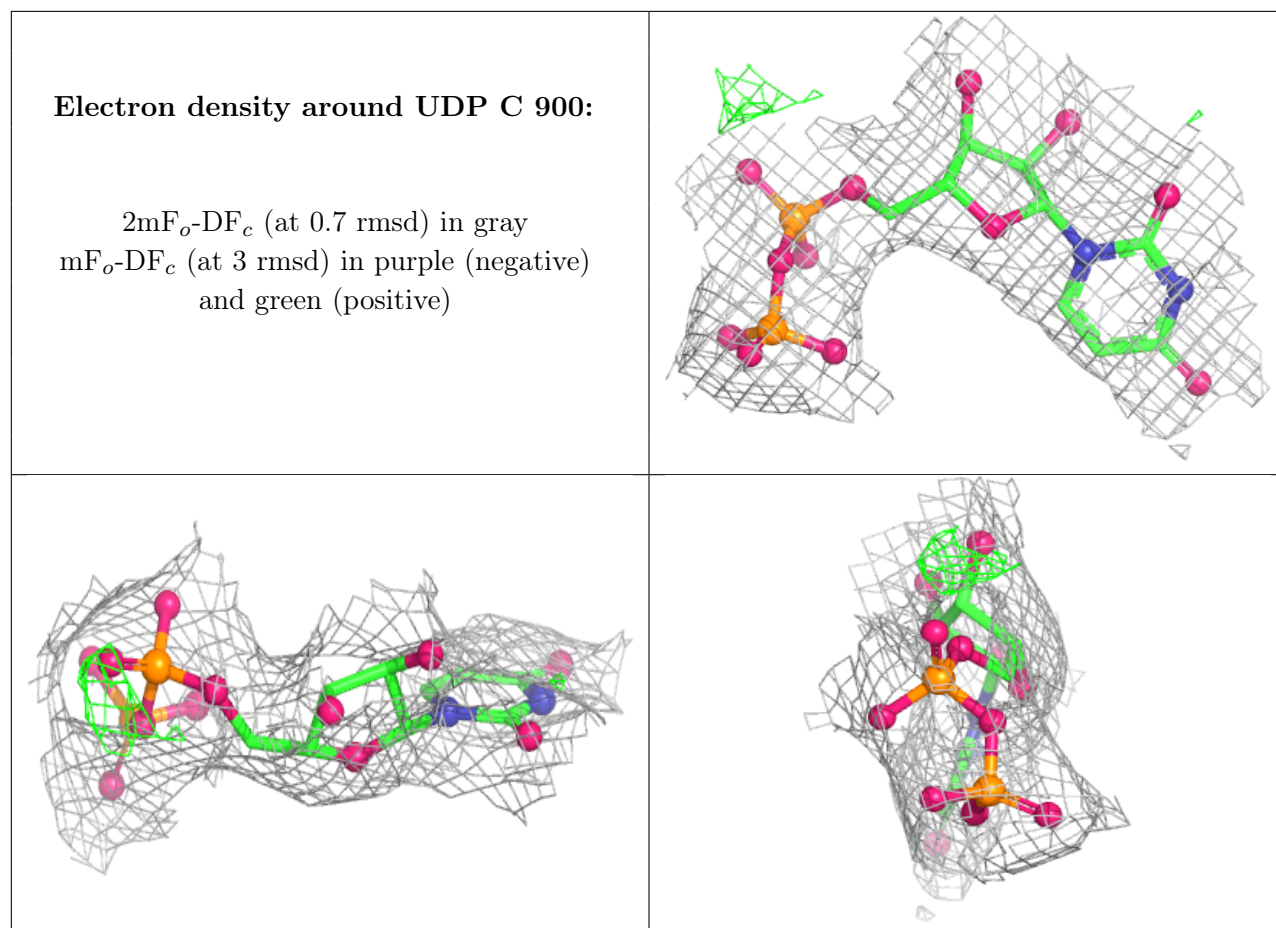
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



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 ⓘ

There are no such residues in this entry.