



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 07:58 AM UTC

PDB ID : 4X6L / pdb_00004x6l
Title : Crystal structure of S. aureus TarM in complex with UDP
Authors : Worrall, L.J.; Sobhanifar, S.; Gruninger, R.J.; Strynadka, N.C.
Deposited on : 2014-12-08
Resolution : 3.19 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

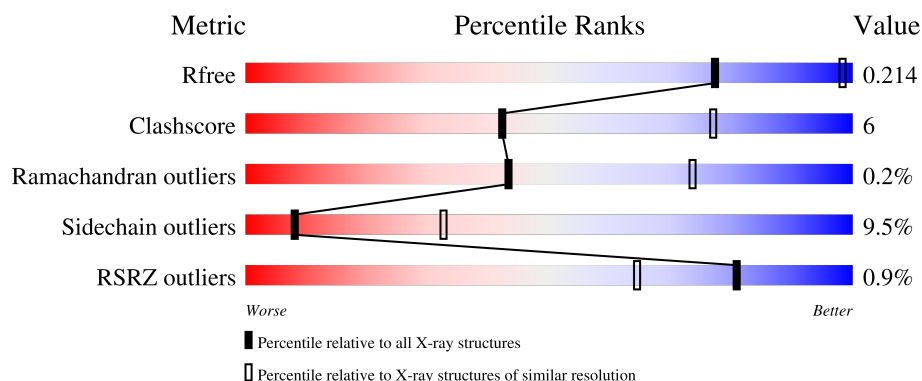
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

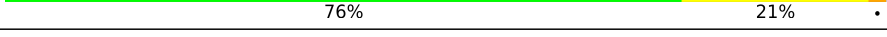
The reported resolution of this entry is 3.19 Å.

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	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

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	493	 76% 20% .
1	B	493	 77% 19% . .
1	C	493	 76% 20% . .
1	D	493	 3% 76% 21% .

2 Entry composition [i](#)

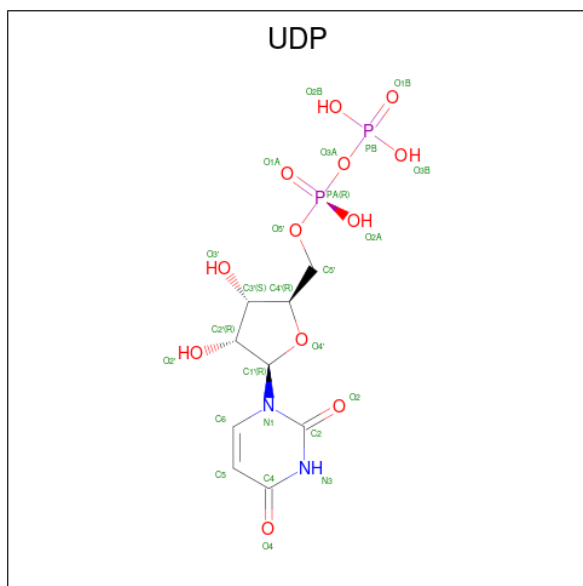
There are 2 unique types of molecules in this entry. The entry contains 16207 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 TarM.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	493	Total	C	N	O	S	0	0	0
			4033	2581	681	752	19			
1	B	493	Total	C	N	O	S	0	0	0
			4033	2581	681	752	19			
1	C	493	Total	C	N	O	S	0	0	0
			4033	2581	681	752	19			
1	D	493	Total	C	N	O	S	0	0	0
			4033	2581	681	752	19			

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



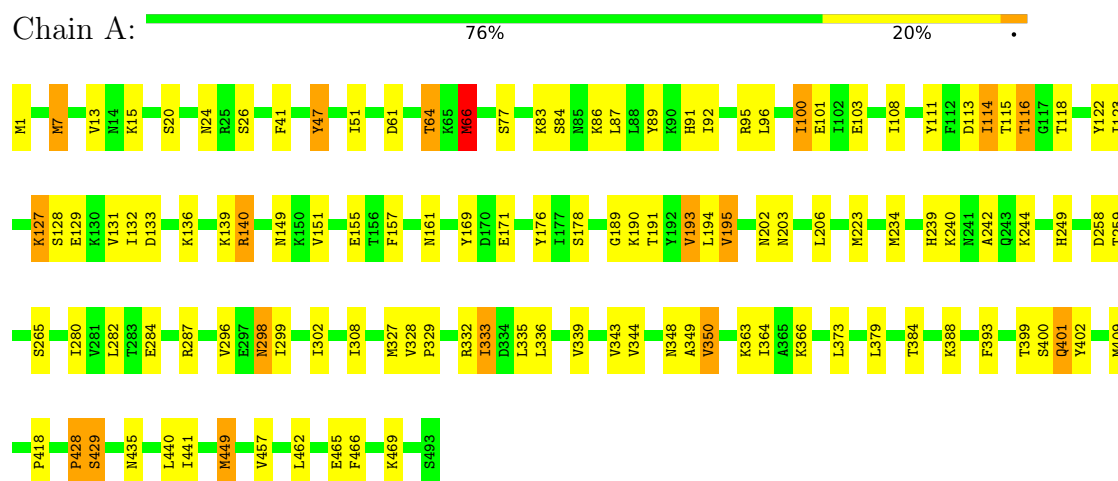
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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	C	1	Total	C	N	O	P	0	0
			25	9	2	12	2		

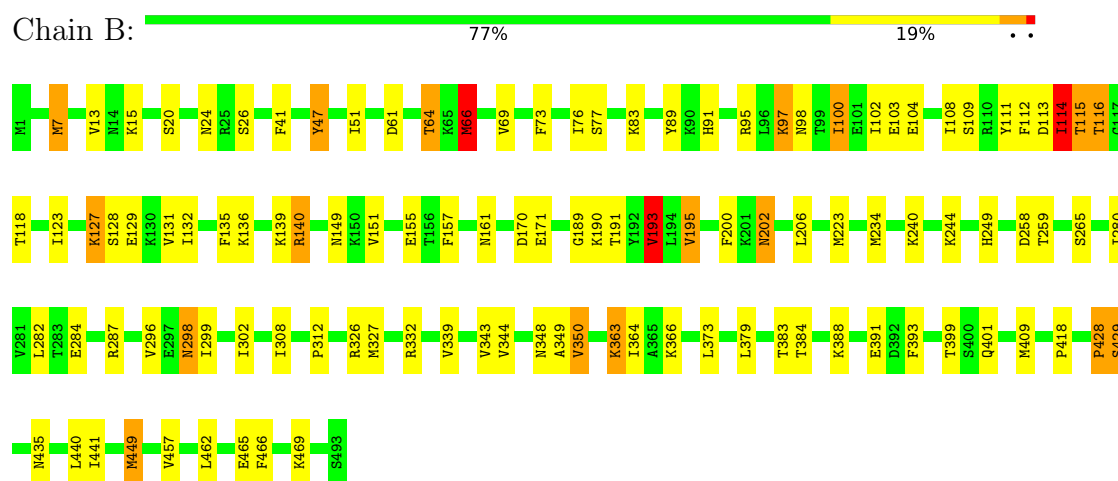
3 Residue-property plots [i](#)

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.

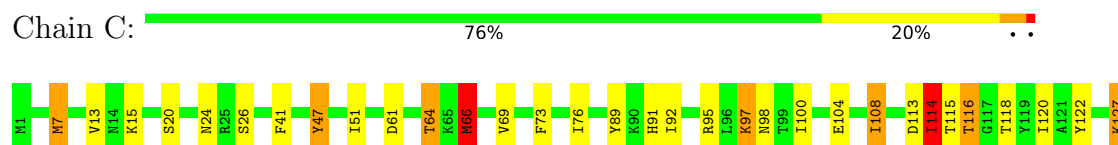
• Molecule 1: TarM

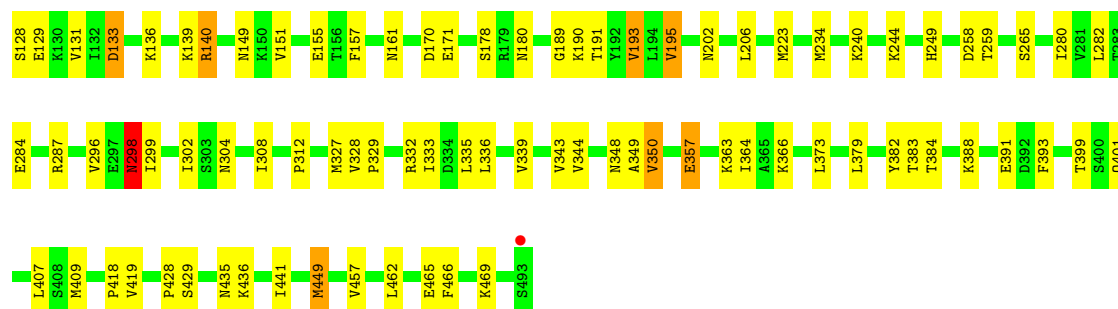


• Molecule 1: TarM

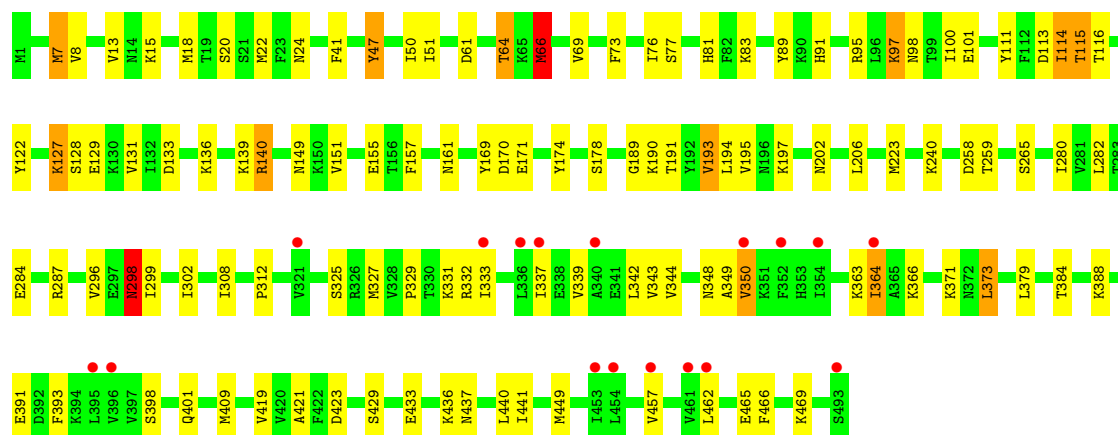
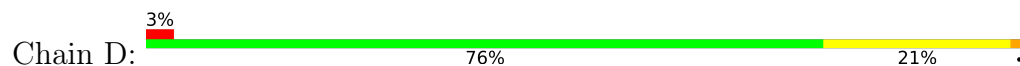


• Molecule 1: TarM





● Molecule 1: TarM



4 Data and refinement statistics

Property	Value	Source
Space group	P 63	Depositor
Cell constants a, b, c, α , β , γ	162.29Å 162.29Å 228.02Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	66.11 – 3.19 66.11 – 3.19	Depositor EDS
% Data completeness (in resolution range)	100.0 (66.11-3.19) 99.9 (66.11-3.19)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.01 (at 3.19Å)	Xtriage
Refinement program	BUSTER-TNT, REFMAC	Depositor
R, R_{free}	0.165 , 0.191 0.188 , 0.214	Depositor DCC
R_{free} test set	2842 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	99.2	Xtriage
Anisotropy	0.187	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 82.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.055 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	16207	wwPDB-VP
Average B, all atoms (Å ²)	120.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.72% 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: 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	0.82	3/4110 (0.1%)	1.39	21/5516 (0.4%)
1	B	0.86	3/4110 (0.1%)	1.38	20/5516 (0.4%)
1	C	0.85	3/4110 (0.1%)	1.42	27/5516 (0.5%)
1	D	0.83	1/4110 (0.0%)	1.41	21/5516 (0.4%)
All	All	0.84	10/16440 (0.1%)	1.40	89/22064 (0.4%)

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	66	MET	SD-CE	-8.35	1.58	1.79
1	C	66	MET	SD-CE	-7.66	1.60	1.79
1	D	66	MET	SD-CE	-7.42	1.61	1.79
1	C	449	MET	SD-CE	-7.38	1.61	1.79
1	A	66	MET	SD-CE	-7.09	1.61	1.79
1	B	449	MET	SD-CE	-6.43	1.63	1.79
1	A	1	MET	SD-CE	5.87	1.94	1.79
1	A	449	MET	SD-CE	-5.63	1.65	1.79
1	C	114	ILE	CA-CB	5.18	1.61	1.54
1	B	114	ILE	CG1-CD1	5.05	1.71	1.51

All (89) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	97	LYS	N-CA-C	7.67	121.57	111.75
1	B	97	LYS	N-CA-C	7.59	120.43	111.71
1	C	15	LYS	N-CA-C	7.15	119.53	110.24
1	B	15	LYS	N-CA-C	7.13	119.50	110.24
1	C	97	LYS	N-CA-C	7.01	120.72	111.75
1	A	15	LYS	N-CA-C	6.91	119.22	110.24
1	B	349	ALA	N-CA-C	6.85	118.75	111.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	15	LYS	N-CA-C	6.82	119.53	110.53
1	D	349	ALA	N-CA-C	6.78	118.67	111.28
1	C	180	ASN	CA-CB-CG	6.61	119.21	112.60
1	C	97	LYS	CA-C-N	-6.61	113.36	122.87
1	C	97	LYS	C-N-CA	-6.61	113.36	122.87
1	C	13	VAL	N-CA-C	6.43	117.43	111.45
1	C	349	ALA	N-CA-C	6.39	118.24	111.28
1	A	349	ALA	N-CA-C	6.30	118.95	111.33
1	B	13	VAL	N-CA-C	6.27	117.28	111.45
1	C	193	VAL	CA-C-N	6.27	128.97	120.38
1	C	193	VAL	C-N-CA	6.27	128.97	120.38
1	A	195	VAL	N-CA-CB	6.23	117.42	110.51
1	C	20	SER	CA-C-N	6.16	128.45	120.44
1	C	20	SER	C-N-CA	6.16	128.45	120.44
1	A	108	ILE	N-CA-C	6.07	117.12	108.93
1	B	97	LYS	CA-C-N	-6.06	114.14	122.87
1	B	97	LYS	C-N-CA	-6.06	114.14	122.87
1	A	20	SER	CA-C-N	6.02	128.26	120.44
1	A	20	SER	C-N-CA	6.02	128.26	120.44
1	C	195	VAL	N-CA-CB	5.92	118.59	110.54
1	B	170	ASP	CA-CB-CG	5.87	118.47	112.60
1	D	342	LEU	CA-C-N	5.80	127.98	120.56
1	D	342	LEU	C-N-CA	5.80	127.98	120.56
1	A	195	VAL	N-CA-C	5.77	116.42	110.36
1	C	348	ASN	CA-C-N	5.72	127.95	120.28
1	C	348	ASN	C-N-CA	5.72	127.95	120.28
1	C	195	VAL	N-CA-C	5.72	116.45	110.62
1	B	116	THR	N-CA-C	5.71	121.12	114.04
1	D	348	ASN	CA-C-N	5.65	127.85	120.28
1	D	348	ASN	C-N-CA	5.65	127.85	120.28
1	D	193	VAL	N-CA-CB	5.63	120.52	111.23
1	B	348	ASN	CA-C-N	5.63	127.82	120.28
1	B	348	ASN	C-N-CA	5.63	127.82	120.28
1	D	329	PRO	CA-C-N	5.59	128.04	120.38
1	D	329	PRO	C-N-CA	5.59	128.04	120.38
1	B	20	SER	CA-C-N	5.59	127.71	120.44
1	B	20	SER	C-N-CA	5.59	127.71	120.44
1	C	284	GLU	CA-C-N	5.58	127.70	120.44
1	C	284	GLU	C-N-CA	5.58	127.70	120.44
1	A	350	VAL	N-CA-C	5.57	116.89	109.37
1	A	348	ASN	CA-C-N	5.55	127.99	120.38
1	A	348	ASN	C-N-CA	5.55	127.99	120.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	97	LYS	CA-C-N	-5.54	114.89	122.87
1	D	97	LYS	C-N-CA	-5.54	114.89	122.87
1	D	20	SER	CA-C-N	5.54	127.64	120.44
1	D	20	SER	C-N-CA	5.54	127.64	120.44
1	B	350	VAL	N-CA-C	5.51	116.81	109.37
1	D	284	GLU	CA-C-N	5.50	127.59	120.44
1	D	284	GLU	C-N-CA	5.50	127.59	120.44
1	C	108	ILE	N-CA-C	5.46	116.30	108.93
1	C	350	VAL	N-CA-C	5.45	116.73	109.37
1	D	282	LEU	N-CA-C	5.43	118.37	111.69
1	D	393	PHE	CA-CB-CG	5.39	119.19	113.80
1	A	193	VAL	N-CA-CB	5.39	117.80	111.66
1	A	284	GLU	CA-C-N	5.39	127.44	120.44
1	A	284	GLU	C-N-CA	5.39	127.44	120.44
1	B	284	GLU	CA-C-N	5.38	127.44	120.44
1	B	284	GLU	C-N-CA	5.38	127.44	120.44
1	A	393	PHE	CA-CB-CG	5.37	119.17	113.80
1	B	393	PHE	CA-CB-CG	5.35	119.15	113.80
1	D	350	VAL	N-CA-C	5.34	116.58	109.37
1	A	428	PRO	CA-C-N	5.34	127.69	120.38
1	A	428	PRO	C-N-CA	5.34	127.69	120.38
1	C	170	ASP	CA-CB-CG	5.31	117.91	112.60
1	C	393	PHE	CA-CB-CG	5.25	119.05	113.80
1	B	428	PRO	CA-C-N	5.20	127.56	120.54
1	B	428	PRO	C-N-CA	5.20	127.56	120.54
1	D	170	ASP	CA-CB-CG	5.15	117.75	112.60
1	C	92	ILE	CA-C-N	5.10	127.07	120.44
1	C	92	ILE	C-N-CA	5.10	127.07	120.44
1	A	13	VAL	N-CA-C	5.09	116.18	111.45
1	C	133	ASP	CA-CB-CG	5.07	117.67	112.60
1	A	400	SER	CA-C-N	5.05	129.15	120.71
1	A	400	SER	C-N-CA	5.05	129.15	120.71
1	C	178	SER	N-CA-C	-5.05	101.69	109.52
1	C	193	VAL	N-CA-CB	5.05	116.96	112.06
1	A	92	ILE	CA-C-N	5.04	126.99	120.44
1	A	92	ILE	C-N-CA	5.04	126.99	120.44
1	D	466	PHE	CA-CB-CG	5.04	118.84	113.80
1	B	326	ARG	N-CA-C	-5.03	102.41	109.96
1	B	193	VAL	N-CA-CB	5.03	117.39	111.66
1	C	193	VAL	CB-CA-C	5.01	118.39	112.68

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	4033	0	4057	48	0
1	B	4033	0	4057	53	0
1	C	4033	0	4057	53	0
1	D	4033	0	4057	49	0
2	A	25	0	11	0	0
2	B	25	0	11	1	0
2	C	25	0	11	1	0
All	All	16207	0	16261	203	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 6.

All (203) 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:100:ILE:HG23	1:A:114:ILE:HG22	1.31	1.10
1:C:100:ILE:HG23	1:C:114:ILE:HG22	1.46	0.95
1:D:100:ILE:HG23	1:D:114:ILE:HG22	1.55	0.87
1:B:339:VAL:HG21	1:B:449:MET:HE2	1.56	0.87
1:D:332:ARG:HG3	1:D:401:GLN:HG2	1.63	0.80
1:D:339:VAL:HG21	1:D:449:MET:HE2	1.64	0.80
1:A:339:VAL:HG21	1:A:449:MET:HE2	1.63	0.79
1:C:47:TYR:CE1	1:C:66:MET:HE1	2.19	0.77
1:B:332:ARG:HH11	1:B:401:GLN:HE22	1.31	0.76
1:D:47:TYR:CE1	1:D:66:MET:HE1	2.21	0.75
1:A:47:TYR:CE1	1:A:66:MET:HE1	2.22	0.74
1:B:47:TYR:CE1	1:B:66:MET:HE1	2.23	0.72
1:C:339:VAL:HG21	1:C:449:MET:HE2	1.70	0.71
1:C:191:THR:HG21	1:C:206:LEU:HD22	1.75	0.68
1:C:113:ASP:HB3	1:C:116:THR:HG22	1.78	0.65
1:C:140:ARG:HD3	1:C:157:PHE:CD2	2.32	0.64
1:D:343:VAL:HG12	1:D:350:VAL:HG11	1.80	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:339:VAL:O	1:C:343:VAL:HG23	1.98	0.63
1:B:249:HIS:HA	1:B:282:LEU:HD11	1.81	0.62
1:C:335:LEU:O	1:C:339:VAL:HG23	2.00	0.61
1:A:332:ARG:HB3	1:A:335:LEU:HD12	1.82	0.60
1:D:77:SER:HB2	1:D:195:VAL:HG11	1.82	0.60
1:C:249:HIS:HA	1:C:282:LEU:HD11	1.83	0.60
1:B:100:ILE:HG23	1:B:114:ILE:HG22	1.83	0.60
1:A:77:SER:HB2	1:A:195:VAL:HG11	1.85	0.59
1:D:333:ILE:HG22	1:D:364:ILE:HG12	1.84	0.58
1:C:357:GLU:HG2	1:C:382:TYR:HB2	1.84	0.58
1:A:7:MET:HE1	1:A:223:MET:HG2	1.86	0.57
1:D:47:TYR:CZ	1:D:66:MET:HE1	2.39	0.57
1:A:249:HIS:HA	1:A:282:LEU:HD11	1.88	0.56
1:D:140:ARG:HD3	1:D:157:PHE:CD1	2.40	0.56
1:D:280:ILE:HG12	1:D:302:ILE:HD12	1.87	0.56
1:A:47:TYR:CZ	1:A:66:MET:HE1	2.39	0.56
1:A:343:VAL:HG12	1:A:350:VAL:HG11	1.87	0.56
1:C:409:MET:HG2	1:C:419:VAL:HG11	1.87	0.56
1:C:122:TYR:HB3	1:C:133:ASP:HB2	1.87	0.56
1:B:280:ILE:HG12	1:B:302:ILE:HD12	1.86	0.56
1:B:191:THR:HG21	1:B:206:LEU:HD22	1.87	0.56
1:C:343:VAL:HG12	1:C:350:VAL:HG11	1.87	0.56
1:A:280:ILE:HG12	1:A:302:ILE:HD12	1.88	0.55
1:B:47:TYR:CZ	1:B:66:MET:HE1	2.42	0.55
1:A:140:ARG:HD3	1:A:157:PHE:CD1	2.41	0.55
1:B:343:VAL:HG12	1:B:350:VAL:HG11	1.88	0.55
1:B:332:ARG:HH11	1:B:401:GLN:NE2	2.01	0.55
1:C:47:TYR:CZ	1:C:66:MET:HE1	2.41	0.55
1:C:280:ILE:HG12	1:C:302:ILE:HD12	1.88	0.55
1:D:98:ASN:HD21	1:D:114:ILE:HG13	1.71	0.55
1:B:98:ASN:HD21	1:B:114:ILE:HG13	1.70	0.55
1:C:113:ASP:CB	1:C:116:THR:HG22	2.37	0.55
1:D:113:ASP:HB3	1:D:116:THR:HG22	1.89	0.55
1:B:140:ARG:HD3	1:B:157:PHE:CD1	2.42	0.55
1:D:61:ASP:HB3	1:D:64:THR:HG22	1.90	0.54
1:A:189:GLY:O	1:A:202:ASN:HA	2.07	0.54
1:A:399:THR:HG23	1:A:449:MET:HE1	1.90	0.54
1:B:98:ASN:ND2	1:B:115:THR:HG22	2.22	0.54
1:B:113:ASP:HB3	1:B:116:THR:HG22	1.90	0.54
1:B:418:PRO:HB3	1:B:466:PHE:HB3	1.89	0.53
1:A:191:THR:HG21	1:A:206:LEU:HD22	1.89	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:333:ILE:HG23	1:A:336:LEU:HD23	1.89	0.53
1:D:7:MET:HE1	1:D:223:MET:HG2	1.90	0.53
1:B:7:MET:HE1	1:B:223:MET:HG2	1.91	0.52
1:B:409:MET:HE1	1:B:428:PRO:HG3	1.90	0.52
1:C:61:ASP:HB3	1:C:64:THR:HG22	1.92	0.52
1:D:100:ILE:CG2	1:D:114:ILE:HG22	2.32	0.52
1:A:61:ASP:HB3	1:A:64:THR:HG22	1.92	0.52
1:C:98:ASN:ND2	1:C:115:THR:HG22	2.25	0.52
1:D:77:SER:CB	1:D:195:VAL:HG11	2.39	0.52
1:C:189:GLY:O	1:C:202:ASN:HA	2.09	0.52
1:C:383:THR:O	2:C:500:UDP:H5	1.92	0.52
1:C:399:THR:HG23	1:C:449:MET:HE1	1.92	0.52
1:B:123:ILE:HG12	1:B:132:ILE:HG12	1.92	0.51
1:B:287:ARG:HG3	1:B:299:ILE:HG22	1.91	0.51
1:D:191:THR:HG21	1:D:206:LEU:HD22	1.92	0.51
1:B:61:ASP:HB3	1:B:64:THR:HG22	1.92	0.51
1:B:399:THR:HG23	1:B:449:MET:HE1	1.91	0.51
1:A:287:ARG:HG3	1:A:299:ILE:HG22	1.92	0.51
1:A:429:SER:HA	1:A:440:LEU:HD21	1.93	0.51
1:C:7:MET:HE1	1:C:223:MET:HG2	1.92	0.51
1:C:116:THR:HG23	1:C:118:THR:OG1	2.11	0.51
1:C:98:ASN:HD21	1:C:114:ILE:HG13	1.76	0.51
1:D:423:ASP:HB2	1:D:440:LEU:HG	1.93	0.50
1:A:418:PRO:HB3	1:A:466:PHE:HB3	1.93	0.50
1:C:332:ARG:HB3	1:C:335:LEU:HD12	1.93	0.50
1:D:98:ASN:ND2	1:D:115:THR:HG22	2.26	0.50
1:D:287:ARG:HG3	1:D:299:ILE:HG22	1.94	0.50
1:D:409:MET:HG2	1:D:419:VAL:HG11	1.94	0.50
1:A:122:TYR:HB3	1:A:133:ASP:HB2	1.94	0.50
1:C:73:PHE:HA	1:C:76:ILE:HD12	1.93	0.50
1:D:178:SER:HB3	1:D:194:LEU:HD11	1.93	0.49
1:C:287:ARG:HG3	1:C:299:ILE:HG22	1.94	0.49
1:C:418:PRO:HB3	1:C:466:PHE:HB3	1.94	0.49
1:B:441:ILE:HG21	1:B:449:MET:HB2	1.95	0.48
1:B:116:THR:HG23	1:B:118:THR:H	1.78	0.48
1:C:409:MET:HE1	1:C:428:PRO:HG3	1.96	0.48
1:B:73:PHE:HA	1:B:76:ILE:HD12	1.96	0.48
1:A:101:GLU:HB2	1:A:111:TYR:CE2	2.49	0.48
1:C:441:ILE:HG21	1:C:449:MET:HB2	1.95	0.48
1:D:73:PHE:HA	1:D:76:ILE:HD12	1.96	0.48
1:B:339:VAL:O	1:B:343:VAL:HG23	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:409:MET:HE1	1:A:428:PRO:HG3	1.97	0.47
1:D:122:TYR:HB3	1:D:133:ASP:HB2	1.97	0.47
1:D:433:GLU:HB2	1:D:437:ASN:HD21	1.80	0.46
1:A:328:VAL:HB	1:A:329:PRO:CD	2.46	0.46
1:A:441:ILE:HG21	1:A:449:MET:HB2	1.98	0.46
1:B:111:TYR:CE2	1:B:123:ILE:HD12	2.50	0.46
1:B:127:LYS:HG2	1:B:128:SER:H	1.80	0.46
1:B:429:SER:HA	1:B:440:LEU:HD21	1.96	0.46
1:B:189:GLY:O	1:B:202:ASN:HA	2.16	0.46
1:B:113:ASP:CB	1:B:116:THR:HG22	2.46	0.46
1:A:113:ASP:HB3	1:A:116:THR:HG22	1.98	0.46
1:A:123:ILE:HG12	1:A:132:ILE:HG12	1.98	0.46
1:A:333:ILE:HD12	1:A:333:ILE:H	1.80	0.46
1:D:127:LYS:HG2	1:D:128:SER:H	1.80	0.46
1:A:401:GLN:HB3	1:A:402:TYR:CD2	2.51	0.45
1:D:77:SER:O	1:D:81:HIS:HD2	1.99	0.45
1:D:441:ILE:HG21	1:D:449:MET:HB2	1.98	0.45
1:C:104:GLU:HB3	1:C:108:ILE:HG22	1.98	0.45
1:D:189:GLY:O	1:D:202:ASN:HA	2.16	0.45
1:A:140:ARG:HD2	1:A:161:ASN:OD1	2.16	0.45
1:B:7:MET:HG2	1:B:41:PHE:CZ	2.52	0.45
1:C:332:ARG:HH11	1:C:401:GLN:NE2	2.15	0.45
1:D:337:ILE:HG22	1:D:371:LYS:HD2	1.97	0.45
1:C:328:VAL:HB	1:C:329:PRO:CD	2.48	0.45
1:B:41:PHE:HA	1:B:69:VAL:HG12	1.99	0.44
1:D:13:VAL:HG13	1:D:50:ILE:HG23	2.00	0.44
1:B:140:ARG:HD2	1:B:161:ASN:OD1	2.18	0.44
1:B:383:THR:O	2:B:500:UDP:H5	1.99	0.44
1:A:7:MET:HG2	1:A:41:PHE:CZ	2.52	0.44
1:A:339:VAL:O	1:A:343:VAL:HG23	2.17	0.44
1:C:127:LYS:HG2	1:C:128:SER:H	1.83	0.44
1:A:91:HIS:O	1:A:95:ARG:HG2	2.18	0.44
1:C:91:HIS:O	1:C:95:ARG:HG2	2.18	0.43
1:C:7:MET:HG2	1:C:41:PHE:CZ	2.53	0.43
1:C:304:ASN:HD22	1:C:407:LEU:HD21	1.83	0.43
1:D:91:HIS:CE1	1:D:95:ARG:HH11	2.36	0.43
1:A:116:THR:HG23	1:A:118:THR:OG1	2.19	0.43
1:B:98:ASN:ND2	1:B:114:ILE:HG13	2.33	0.43
1:D:91:HIS:O	1:D:95:ARG:HG2	2.18	0.43
1:A:91:HIS:CE1	1:A:95:ARG:HH11	2.36	0.43
1:A:127:LYS:HG2	1:A:128:SER:H	1.84	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:91:HIS:CE1	1:B:95:ARG:HH11	2.37	0.43
1:C:333:ILE:HG23	1:C:336:LEU:HD23	2.01	0.43
1:B:193:VAL:CG1	1:B:200:PHE:HE1	2.31	0.43
1:B:312:PRO:HG3	1:B:391:GLU:HG2	2.00	0.43
1:C:140:ARG:HD2	1:C:161:ASN:OD1	2.19	0.43
1:D:398:SER:HB3	1:D:421:ALA:HB2	2.00	0.43
1:D:7:MET:HG2	1:D:41:PHE:CZ	2.54	0.42
1:D:169:TYR:HA	1:D:174:TYR:O	2.18	0.42
1:B:399:THR:CG2	1:B:449:MET:HE1	2.49	0.42
1:C:332:ARG:HH11	1:C:401:GLN:HE22	1.67	0.42
1:D:113:ASP:CB	1:D:116:THR:HG22	2.49	0.42
1:B:234:MET:HG2	1:B:244:LYS:HD3	2.01	0.42
1:C:344:VAL:HG21	1:C:373:LEU:HD22	2.01	0.42
1:B:135:PHE:CE2	1:B:140:ARG:HB2	2.53	0.42
1:D:325:SER:HB2	1:D:331:LYS:HD2	2.00	0.42
1:D:101:GLU:HG3	1:D:111:TYR:CE1	2.55	0.42
1:A:84:SER:HB3	1:A:86:LYS:HE3	2.01	0.42
1:A:327:MET:HE1	1:A:379:LEU:HD22	2.01	0.42
1:C:41:PHE:HA	1:C:69:VAL:HG12	2.00	0.42
1:C:98:ASN:HD21	1:C:115:THR:HG22	1.84	0.42
1:C:91:HIS:CE1	1:C:95:ARG:HH11	2.38	0.42
1:D:41:PHE:HA	1:D:69:VAL:HG12	2.01	0.42
1:B:102:ILE:HD12	1:B:112:PHE:CD1	2.55	0.42
1:C:312:PRO:HG3	1:C:391:GLU:HG2	2.02	0.42
1:D:327:MET:HE1	1:D:379:LEU:HD22	2.01	0.42
1:A:77:SER:CB	1:A:195:VAL:HG11	2.49	0.42
1:B:26:SER:OG	1:B:64:THR:HB	2.20	0.42
1:B:91:HIS:O	1:B:95:ARG:HG2	2.20	0.42
1:C:26:SER:OG	1:C:64:THR:HB	2.20	0.42
1:C:234:MET:HG2	1:C:244:LYS:HD3	2.01	0.42
1:A:169:TYR:C	1:A:176:TYR:HE1	2.28	0.41
1:B:98:ASN:HD21	1:B:115:THR:HG22	1.85	0.41
1:C:113:ASP:HB2	1:C:120:ILE:HD11	2.02	0.41
1:D:312:PRO:HG3	1:D:391:GLU:HG2	2.01	0.41
1:A:87:LEU:HD23	1:A:87:LEU:HA	1.85	0.41
1:B:104:GLU:HB3	1:B:108:ILE:HG22	2.02	0.41
1:A:332:ARG:HG3	1:A:401:GLN:HE21	1.86	0.41
1:A:344:VAL:HG21	1:A:373:LEU:HD22	2.03	0.41
1:B:103:GLU:HG2	1:B:109:SER:OG	2.21	0.41
1:C:339:VAL:HG21	1:C:449:MET:CE	2.44	0.41
1:D:89:TYR:CZ	1:D:151:VAL:HG12	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:298:ASN:HD22	1:D:298:ASN:HA	1.70	0.41
1:A:26:SER:OG	1:A:64:THR:HB	2.20	0.41
1:A:89:TYR:CZ	1:A:151:VAL:HG12	2.55	0.41
1:C:298:ASN:HD22	1:C:298:ASN:HA	1.70	0.41
1:D:140:ARG:HD2	1:D:161:ASN:OD1	2.20	0.41
1:A:178:SER:HB3	1:A:194:LEU:HD11	2.03	0.41
1:B:77:SER:HA	1:B:195:VAL:HG11	2.03	0.41
1:B:327:MET:HE1	1:B:379:LEU:HD22	2.03	0.41
1:B:344:VAL:HG21	1:B:373:LEU:HD22	2.02	0.41
1:C:89:TYR:CZ	1:C:151:VAL:HG12	2.56	0.41
1:D:337:ILE:HD11	1:D:364:ILE:HG23	2.03	0.41
1:A:234:MET:HG2	1:A:244:LYS:HD3	2.03	0.40
1:B:89:TYR:CZ	1:B:151:VAL:HG12	2.57	0.40
1:A:399:THR:CG2	1:A:449:MET:HE1	2.50	0.40
1:B:127:LYS:HG2	1:B:128:SER:N	2.37	0.40
1:C:327:MET:HE1	1:C:379:LEU:HD22	2.02	0.40
1:D:8:VAL:HB	1:D:22:MET:HE1	2.04	0.40
1:C:116:THR:HG23	1:C:118:THR:H	1.86	0.40
1:D:344:VAL:HG21	1:D:373:LEU:HD22	2.02	0.40
1:D:127:LYS:HG2	1:D:128:SER:N	2.36	0.40
1:A:239:HIS:HB3	1:A:242:ALA:HB3	2.04	0.40
1:B:363:LYS:H	1:B:363:LYS:HG2	1.71	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	491/493 (100%)	457 (93%)	33 (7%)	1 (0%)	43	73
1	B	491/493 (100%)	462 (94%)	28 (6%)	1 (0%)	43	73
1	C	491/493 (100%)	461 (94%)	29 (6%)	1 (0%)	43	73

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	491/493 (100%)	456 (93%)	34 (7%)	1 (0%)	43	73
All	All	1964/1972 (100%)	1836 (94%)	124 (6%)	4 (0%)	43	73

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	298	ASN
1	B	298	ASN
1	C	298	ASN
1	D	298	ASN

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	449/449 (100%)	404 (90%)	45 (10%)	7	30
1	B	449/449 (100%)	407 (91%)	42 (9%)	8	33
1	C	449/449 (100%)	408 (91%)	41 (9%)	9	34
1	D	449/449 (100%)	407 (91%)	42 (9%)	8	33
All	All	1796/1796 (100%)	1626 (90%)	170 (10%)	8	32

All (170) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	MET
1	A	24	ASN
1	A	47	TYR
1	A	51	ILE
1	A	64	THR
1	A	66	MET
1	A	83	LYS
1	A	96	LEU
1	A	100	ILE
1	A	103	GLU

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Mol	Chain	Res	Type
1	A	114	ILE
1	A	115	THR
1	A	116	THR
1	A	127	LYS
1	A	129	GLU
1	A	131	VAL
1	A	136	LYS
1	A	139	LYS
1	A	140	ARG
1	A	149	ASN
1	A	155	GLU
1	A	171	GLU
1	A	190	LYS
1	A	193	VAL
1	A	203	ASN
1	A	240	LYS
1	A	258	ASP
1	A	259	THR
1	A	265	SER
1	A	296	VAL
1	A	298	ASN
1	A	308	ILE
1	A	333	ILE
1	A	363	LYS
1	A	364	ILE
1	A	366	LYS
1	A	384	THR
1	A	388	LYS
1	A	401	GLN
1	A	429	SER
1	A	435	ASN
1	A	457	VAL
1	A	462	LEU
1	A	465	GLU
1	A	469	LYS
1	B	7	MET
1	B	24	ASN
1	B	47	TYR
1	B	51	ILE
1	B	64	THR
1	B	66	MET
1	B	83	LYS

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Mol	Chain	Res	Type
1	B	97	LYS
1	B	100	ILE
1	B	114	ILE
1	B	115	THR
1	B	127	LYS
1	B	129	GLU
1	B	131	VAL
1	B	136	LYS
1	B	139	LYS
1	B	140	ARG
1	B	149	ASN
1	B	155	GLU
1	B	171	GLU
1	B	190	LYS
1	B	193	VAL
1	B	195	VAL
1	B	202	ASN
1	B	240	LYS
1	B	258	ASP
1	B	259	THR
1	B	265	SER
1	B	296	VAL
1	B	298	ASN
1	B	308	ILE
1	B	363	LYS
1	B	364	ILE
1	B	366	LYS
1	B	384	THR
1	B	388	LYS
1	B	429	SER
1	B	435	ASN
1	B	457	VAL
1	B	462	LEU
1	B	465	GLU
1	B	469	LYS
1	C	7	MET
1	C	24	ASN
1	C	47	TYR
1	C	51	ILE
1	C	64	THR
1	C	66	MET
1	C	97	LYS

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Mol	Chain	Res	Type
1	C	114	ILE
1	C	116	THR
1	C	127	LYS
1	C	129	GLU
1	C	131	VAL
1	C	136	LYS
1	C	139	LYS
1	C	140	ARG
1	C	149	ASN
1	C	155	GLU
1	C	171	GLU
1	C	190	LYS
1	C	193	VAL
1	C	195	VAL
1	C	240	LYS
1	C	258	ASP
1	C	259	THR
1	C	265	SER
1	C	296	VAL
1	C	298	ASN
1	C	308	ILE
1	C	357	GLU
1	C	363	LYS
1	C	364	ILE
1	C	366	LYS
1	C	384	THR
1	C	388	LYS
1	C	429	SER
1	C	435	ASN
1	C	436	LYS
1	C	457	VAL
1	C	462	LEU
1	C	465	GLU
1	C	469	LYS
1	D	7	MET
1	D	18	MET
1	D	24	ASN
1	D	47	TYR
1	D	51	ILE
1	D	64	THR
1	D	66	MET
1	D	83	LYS

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Mol	Chain	Res	Type
1	D	97	LYS
1	D	114	ILE
1	D	115	THR
1	D	127	LYS
1	D	129	GLU
1	D	131	VAL
1	D	136	LYS
1	D	139	LYS
1	D	140	ARG
1	D	149	ASN
1	D	155	GLU
1	D	171	GLU
1	D	190	LYS
1	D	193	VAL
1	D	197	LYS
1	D	240	LYS
1	D	258	ASP
1	D	259	THR
1	D	265	SER
1	D	296	VAL
1	D	298	ASN
1	D	308	ILE
1	D	363	LYS
1	D	364	ILE
1	D	366	LYS
1	D	373	LEU
1	D	384	THR
1	D	388	LYS
1	D	429	SER
1	D	436	LYS
1	D	457	VAL
1	D	462	LEU
1	D	465	GLU
1	D	469	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (45) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	9	HIS
1	A	91	HIS
1	A	241	ASN
1	A	249	HIS

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Mol	Chain	Res	Type
1	A	292	ASN
1	A	298	ASN
1	A	348	ASN
1	A	372	ASN
1	A	443	ASN
1	A	444	HIS
1	B	81	HIS
1	B	91	HIS
1	B	98	ASN
1	B	203	ASN
1	B	241	ASN
1	B	292	ASN
1	B	298	ASN
1	B	348	ASN
1	B	401	GLN
1	B	443	ASN
1	B	444	HIS
1	C	9	HIS
1	C	81	HIS
1	C	91	HIS
1	C	98	ASN
1	C	203	ASN
1	C	241	ASN
1	C	292	ASN
1	C	298	ASN
1	C	348	ASN
1	C	372	ASN
1	C	401	GLN
1	C	443	ASN
1	D	9	HIS
1	D	14	ASN
1	D	81	HIS
1	D	91	HIS
1	D	98	ASN
1	D	203	ASN
1	D	241	ASN
1	D	255	ASN
1	D	298	ASN
1	D	348	ASN
1	D	437	ASN
1	D	444	HIS

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](#)

3 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	500	-	25,26,26	0.47	0	38,40,40	0.62	0
2	UDP	C	500	-	25,26,26	0.45	0	38,40,40	0.66	0
2	UDP	B	500	-	25,26,26	0.67	2 (8%)	38,40,40	0.66	0

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	500	-	-	5/16/32/32	0/2/2/2
2	UDP	C	500	-	-	5/16/32/32	0/2/2/2
2	UDP	B	500	-	-	4/16/32/32	0/2/2/2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	500	UDP	PB-O3B	2.15	1.62	1.54
2	B	500	UDP	PA-O3A	2.09	1.61	1.59

There are no bond angle outliers.

There are no chirality outliers.

All (14) torsion outliers are listed below:

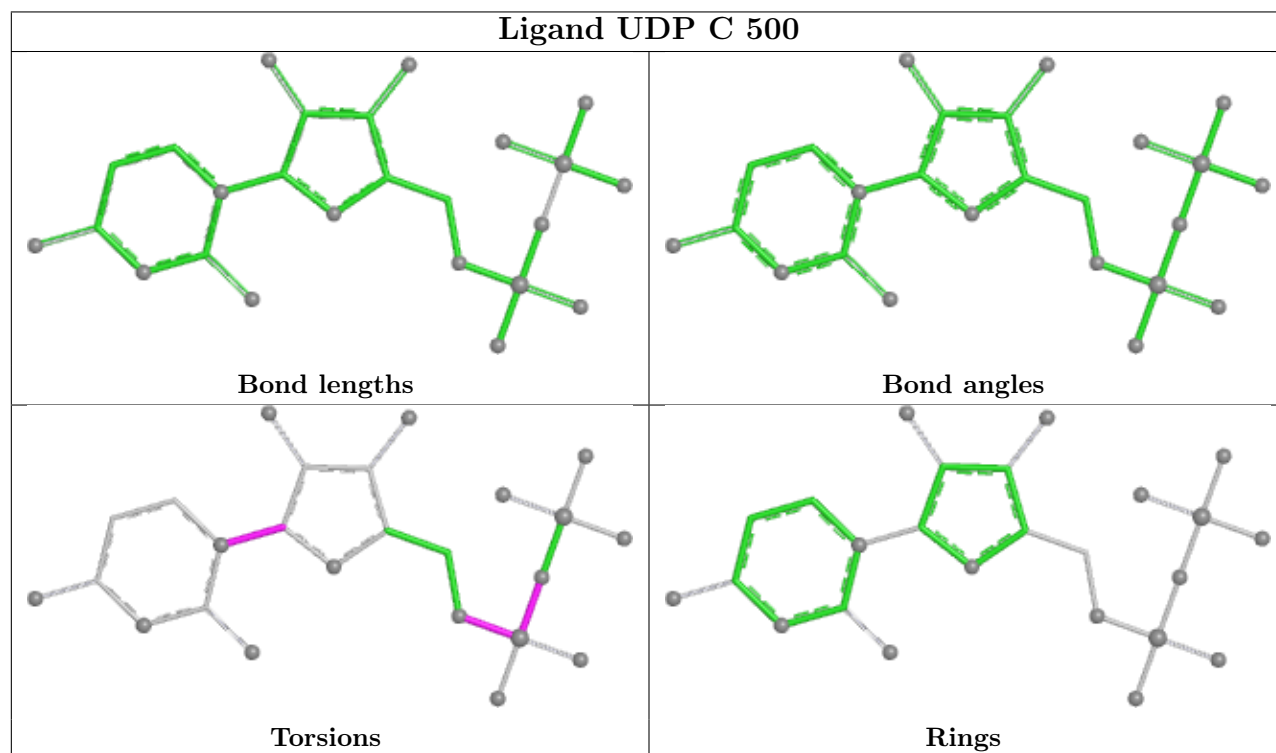
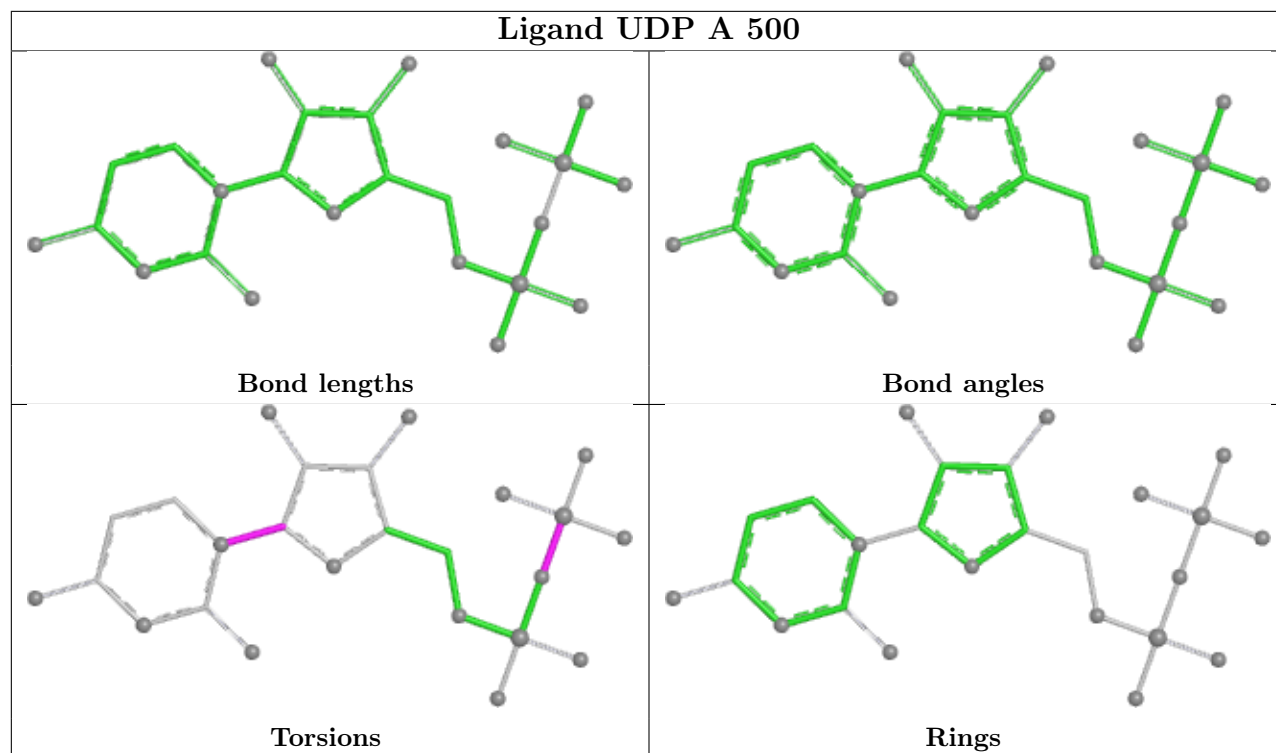
Mol	Chain	Res	Type	Atoms
2	A	500	UDP	O4'-C1'-N1-C2
2	A	500	UDP	O4'-C1'-N1-C6
2	A	500	UDP	PA-O3A-PB-O3B
2	B	500	UDP	O4'-C1'-N1-C2
2	B	500	UDP	O4'-C1'-N1-C6
2	B	500	UDP	PA-O3A-PB-O3B
2	C	500	UDP	O4'-C1'-N1-C2
2	C	500	UDP	O4'-C1'-N1-C6
2	C	500	UDP	C5'-O5'-PA-O1A
2	C	500	UDP	C5'-O5'-PA-O3A
2	C	500	UDP	PB-O3A-PA-O5'
2	A	500	UDP	PA-O3A-PB-O1B
2	B	500	UDP	PA-O3A-PB-O1B
2	A	500	UDP	PA-O3A-PB-O2B

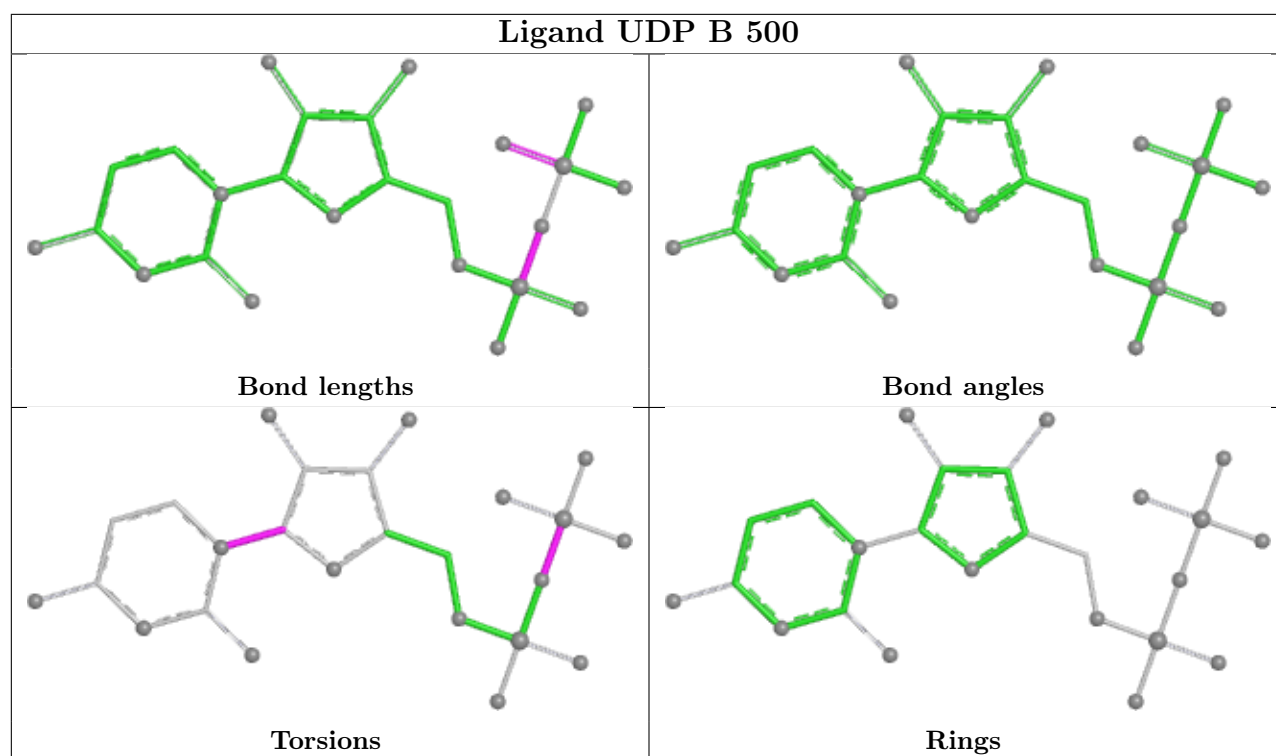
There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	500	UDP	1	0
2	B	500	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 [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	493/493 (100%)	-0.35	0	100	100	77, 112, 147, 163	0
1	B	493/493 (100%)	-0.42	0	100	100	62, 97, 139, 163	0
1	C	493/493 (100%)	-0.37	1 (0%)	91	85	62, 103, 140, 159	0
1	D	493/493 (100%)	0.06	17 (3%)	48	30	67, 148, 269, 277	0
All	All	1972/1972 (100%)	-0.27	18 (0%)	81	64	62, 109, 252, 277	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	453	ILE	6.9
1	D	396	VAL	4.1
1	D	336	LEU	4.0
1	D	337	ILE	3.4
1	D	333	ILE	3.2
1	D	350	VAL	2.9
1	D	457	VAL	2.7
1	D	352	PHE	2.7
1	D	493	SER	2.5
1	D	340	ALA	2.4
1	D	462	LEU	2.4
1	C	493	SER	2.4
1	D	454	LEU	2.3
1	D	321	VAL	2.3
1	D	354	ILE	2.2
1	D	364	ILE	2.2
1	D	395	LEU	2.2
1	D	461	VAL	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](#)

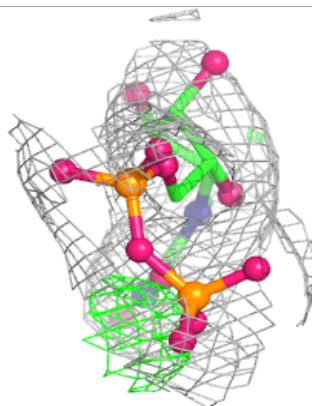
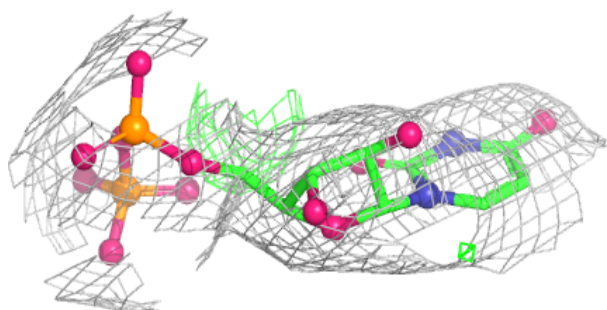
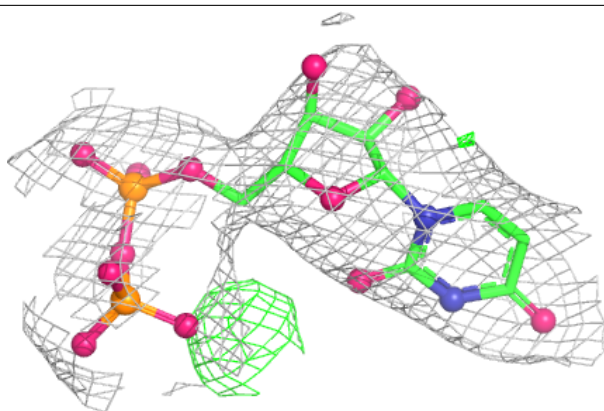
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	500	25/25	0.93	0.09	107,122,129,130	0
2	UDP	B	500	25/25	0.94	0.09	107,118,124,127	0
2	UDP	A	500	25/25	0.96	0.07	117,131,134,140	0

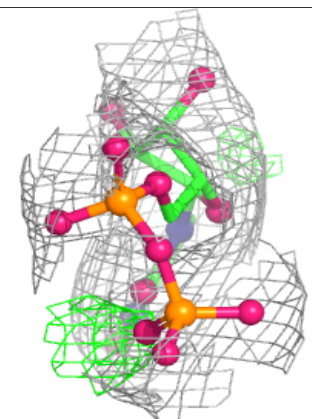
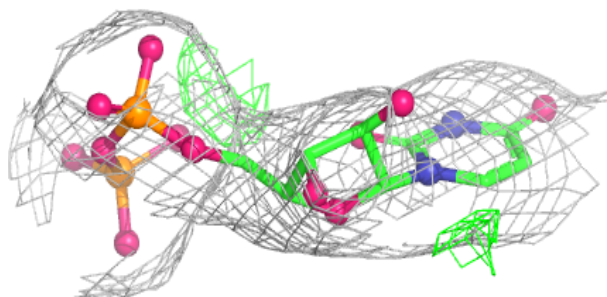
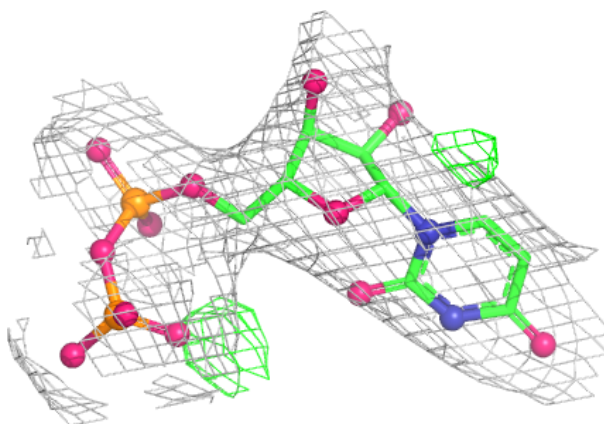
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

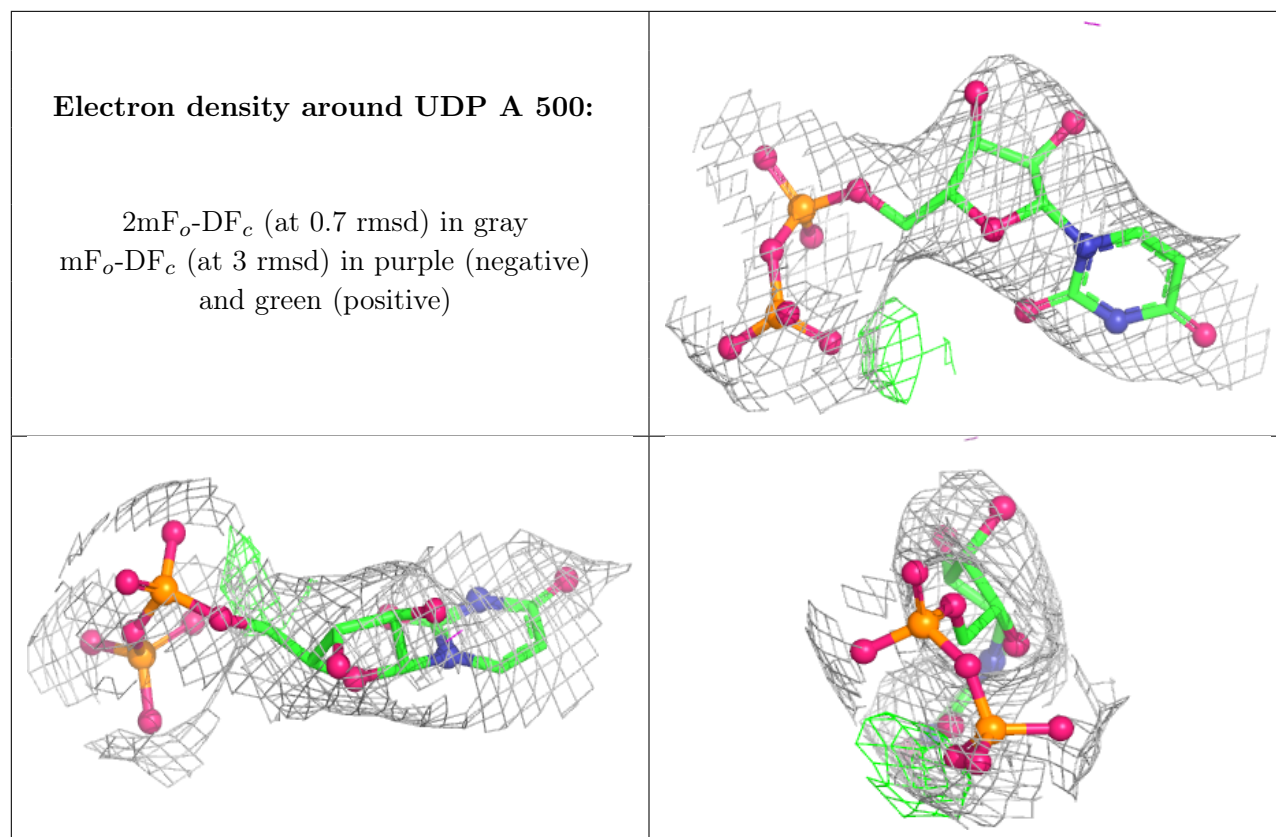
Electron density around UDP C 500:

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

**Electron density around UDP B 500:**

$2mF_o-DF_c$ (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.