



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 05:56 PM UTC

PDB ID : 2QRN / pdb_00002qrn
Title : Human Deoxycytidine kinase dCMP, UDP, Mg ion product complex
Authors : Ealick, S.E.; Soriano, E.V.
Deposited on : 2007-07-28
Resolution : 3.40 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

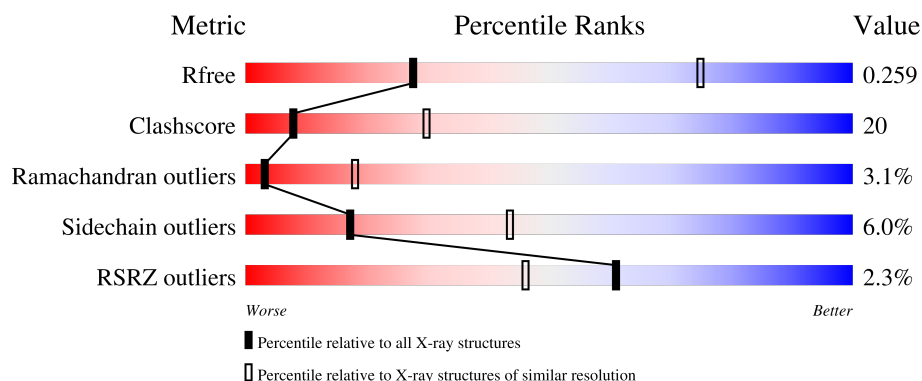
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.40 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	1001 (3.44-3.36)
Clashscore	190562	1022 (3.44-3.36)
Ramachandran outliers	187476	1012 (3.44-3.36)
Sidechain outliers	187428	1012 (3.44-3.36)
RSRZ outliers	180081	1001 (3.44-3.36)

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	280	0% 51% 24% 6% 19%
1	B	280	2% 51% 25% 5% 19%
1	C	280	2% 51% 24% 6% 19%
1	D	280	3% 51% 25% 5% 19%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7808 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 Deoxycytidine kinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	228	Total	C	N	O	S	0	0	0
			1906	1223	317	357	9			
1	B	228	Total	C	N	O	S	0	0	0
			1906	1223	317	357	9			
1	C	228	Total	C	N	O	S	0	0	0
			1906	1223	317	357	9			
1	D	228	Total	C	N	O	S	0	0	0
			1906	1223	317	357	9			

There are 80 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	expression tag	UNP P27707
A	-18	GLY	-	expression tag	UNP P27707
A	-17	SER	-	expression tag	UNP P27707
A	-16	SER	-	expression tag	UNP P27707
A	-15	HIS	-	expression tag	UNP P27707
A	-14	HIS	-	expression tag	UNP P27707
A	-13	HIS	-	expression tag	UNP P27707
A	-12	HIS	-	expression tag	UNP P27707
A	-11	HIS	-	expression tag	UNP P27707
A	-10	HIS	-	expression tag	UNP P27707
A	-9	SER	-	expression tag	UNP P27707
A	-8	SER	-	expression tag	UNP P27707
A	-7	GLY	-	expression tag	UNP P27707
A	-6	LEU	-	expression tag	UNP P27707
A	-5	VAL	-	expression tag	UNP P27707
A	-4	PRO	-	expression tag	UNP P27707
A	-3	ARG	-	expression tag	UNP P27707
A	-2	GLY	-	expression tag	UNP P27707
A	-1	SER	-	expression tag	UNP P27707
A	0	HIS	-	expression tag	UNP P27707
B	-19	MET	-	expression tag	UNP P27707

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-18	GLY	-	expression tag	UNP P27707
B	-17	SER	-	expression tag	UNP P27707
B	-16	SER	-	expression tag	UNP P27707
B	-15	HIS	-	expression tag	UNP P27707
B	-14	HIS	-	expression tag	UNP P27707
B	-13	HIS	-	expression tag	UNP P27707
B	-12	HIS	-	expression tag	UNP P27707
B	-11	HIS	-	expression tag	UNP P27707
B	-10	HIS	-	expression tag	UNP P27707
B	-9	SER	-	expression tag	UNP P27707
B	-8	SER	-	expression tag	UNP P27707
B	-7	GLY	-	expression tag	UNP P27707
B	-6	LEU	-	expression tag	UNP P27707
B	-5	VAL	-	expression tag	UNP P27707
B	-4	PRO	-	expression tag	UNP P27707
B	-3	ARG	-	expression tag	UNP P27707
B	-2	GLY	-	expression tag	UNP P27707
B	-1	SER	-	expression tag	UNP P27707
B	0	HIS	-	expression tag	UNP P27707
C	-19	MET	-	expression tag	UNP P27707
C	-18	GLY	-	expression tag	UNP P27707
C	-17	SER	-	expression tag	UNP P27707
C	-16	SER	-	expression tag	UNP P27707
C	-15	HIS	-	expression tag	UNP P27707
C	-14	HIS	-	expression tag	UNP P27707
C	-13	HIS	-	expression tag	UNP P27707
C	-12	HIS	-	expression tag	UNP P27707
C	-11	HIS	-	expression tag	UNP P27707
C	-10	HIS	-	expression tag	UNP P27707
C	-9	SER	-	expression tag	UNP P27707
C	-8	SER	-	expression tag	UNP P27707
C	-7	GLY	-	expression tag	UNP P27707
C	-6	LEU	-	expression tag	UNP P27707
C	-5	VAL	-	expression tag	UNP P27707
C	-4	PRO	-	expression tag	UNP P27707
C	-3	ARG	-	expression tag	UNP P27707
C	-2	GLY	-	expression tag	UNP P27707
C	-1	SER	-	expression tag	UNP P27707
C	0	HIS	-	expression tag	UNP P27707
D	-19	MET	-	expression tag	UNP P27707
D	-18	GLY	-	expression tag	UNP P27707
D	-17	SER	-	expression tag	UNP P27707

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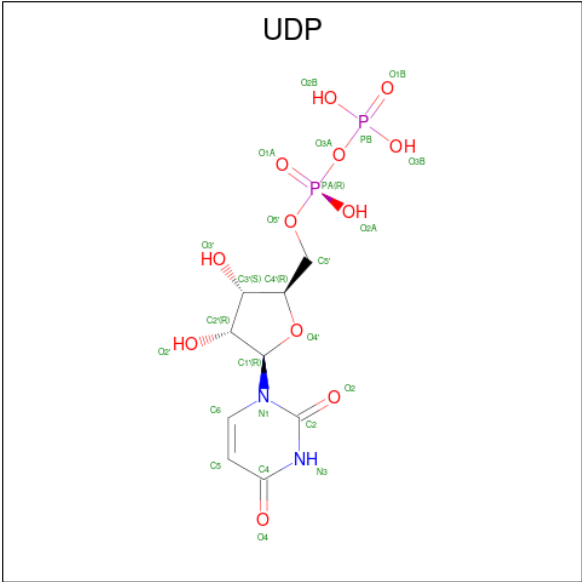
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Chain	Residue	Modelled	Actual	Comment	Reference
D	-16	SER	-	expression tag	UNP P27707
D	-15	HIS	-	expression tag	UNP P27707
D	-14	HIS	-	expression tag	UNP P27707
D	-13	HIS	-	expression tag	UNP P27707
D	-12	HIS	-	expression tag	UNP P27707
D	-11	HIS	-	expression tag	UNP P27707
D	-10	HIS	-	expression tag	UNP P27707
D	-9	SER	-	expression tag	UNP P27707
D	-8	SER	-	expression tag	UNP P27707
D	-7	GLY	-	expression tag	UNP P27707
D	-6	LEU	-	expression tag	UNP P27707
D	-5	VAL	-	expression tag	UNP P27707
D	-4	PRO	-	expression tag	UNP P27707
D	-3	ARG	-	expression tag	UNP P27707
D	-2	GLY	-	expression tag	UNP P27707
D	-1	SER	-	expression tag	UNP P27707
D	0	HIS	-	expression tag	UNP P27707

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

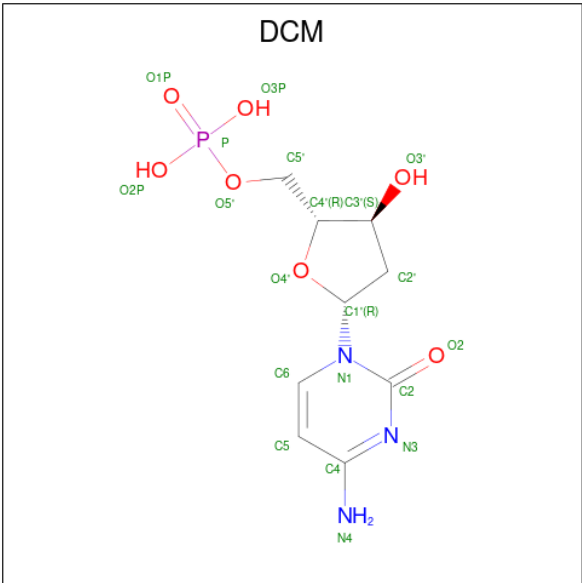
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Mg 1 1	0	0
2	B	1	Total Mg 1 1	0	0
2	C	1	Total Mg 1 1	0	0
2	D	1	Total Mg 1 1	0	0

- Molecule 3 is URIDINE-5'-DIPHOSPHATE (CCD ID: UDP) (formula: C₉H₁₄N₂O₁₂P₂).



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

- Molecule 4 is 2'-DEOXYCYTIDINE-5'-MONOPHOSPHATE (CCD ID: DCM) (formula: $C_9H_{14}N_3O_7P$).

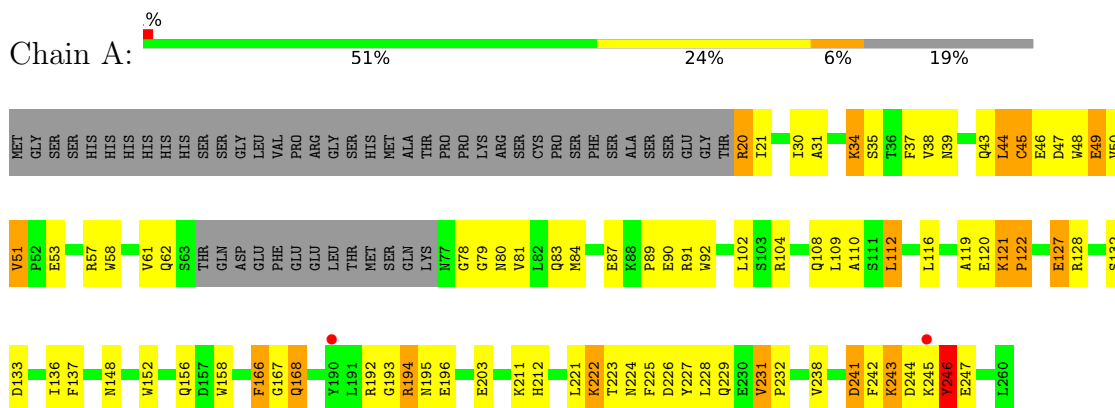


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			20	9	3	7	1		
4	B	1	Total	C	N	O	P	0	0
			20	9	3	7	1		
4	C	1	Total	C	N	O	P	0	0
			20	9	3	7	1		
4	D	1	Total	C	N	O	P	0	0
			20	9	3	7	1		

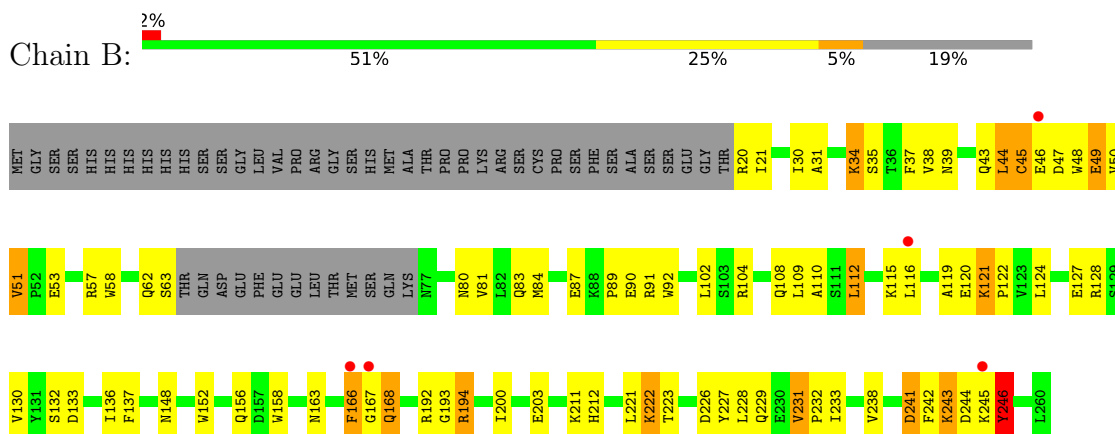
3 Residue-property plots

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.

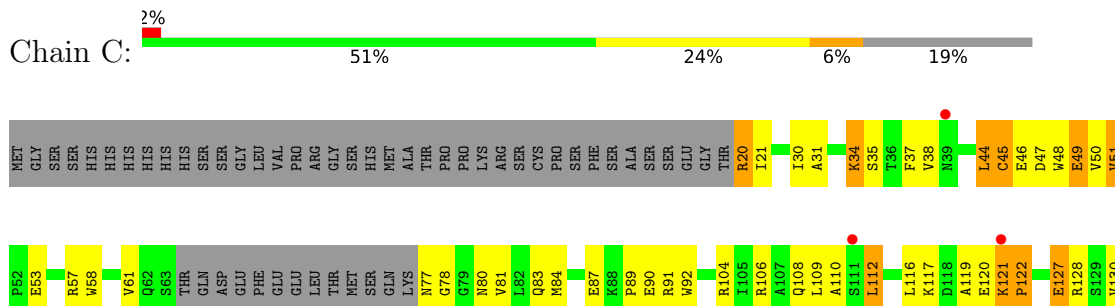
- Molecule 1: Deoxycytidine kinase



- Molecule 1: Deoxycytidine kinase

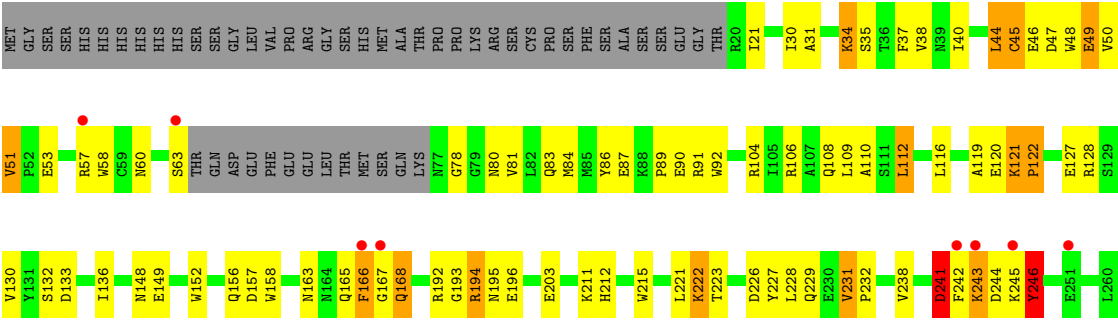


- Molecule 1: Deoxycytidine kinase





• Molecule 1: Deoxycytidine kinase



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	94.56Å 94.56Å 335.33Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.12 – 3.40 48.12 – 3.40	Depositor EDS
% Data completeness (in resolution range)	89.9 (48.12-3.40) 89.8 (48.12-3.40)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	0.13	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.07 (at 3.19Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.217 , 0.269 0.218 , 0.259	Depositor DCC
R_{free} test set	1138 reflections (4.82%)	wwPDB-VP
Wilson B-factor (Å ²)	65.6	Xtriage
Anisotropy	0.403	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 76.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	7808	wwPDB-VP
Average B, all atoms (Å ²)	70.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.97% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MG, DCM, 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	0.50	0/1952	0.97	8/2641 (0.3%)
1	B	0.52	0/1952	0.98	10/2641 (0.4%)
1	C	0.51	0/1952	0.98	9/2641 (0.3%)
1	D	0.53	0/1952	0.97	7/2641 (0.3%)
All	All	0.52	0/7808	0.98	34/10564 (0.3%)

There are no bond length outliers.

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	121	LYS	CA-C-N	7.36	127.28	119.85
1	C	121	LYS	C-N-CA	7.36	127.28	119.85
1	A	121	LYS	CA-C-N	7.06	126.98	119.85
1	A	121	LYS	C-N-CA	7.06	126.98	119.85
1	B	121	LYS	CA-C-N	7.05	126.97	119.85
1	B	121	LYS	C-N-CA	7.05	126.97	119.85
1	D	231	VAL	CA-C-N	6.92	126.88	119.76
1	D	231	VAL	C-N-CA	6.92	126.88	119.76
1	B	231	VAL	CA-C-N	6.91	126.87	119.76
1	B	231	VAL	C-N-CA	6.91	126.87	119.76
1	B	137	PHE	N-CA-C	6.88	118.57	111.14
1	C	61	VAL	N-CA-C	6.86	117.11	107.37
1	C	231	VAL	CA-C-N	6.81	126.78	119.76
1	C	231	VAL	C-N-CA	6.81	126.78	119.76
1	A	231	VAL	CA-C-N	6.50	126.46	119.76
1	A	231	VAL	C-N-CA	6.50	126.46	119.76
1	D	121	LYS	CA-C-N	6.28	126.80	120.14
1	D	121	LYS	C-N-CA	6.28	126.80	120.14
1	C	127	GLU	N-CA-C	-6.10	98.59	108.41
1	A	137	PHE	N-CA-C	6.03	117.72	111.03

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	51	VAL	CA-C-N	6.00	125.92	119.85
1	D	51	VAL	C-N-CA	6.00	125.92	119.85
1	B	51	VAL	CA-C-N	5.89	125.80	119.85
1	B	51	VAL	C-N-CA	5.89	125.80	119.85
1	C	137	PHE	N-CA-C	5.84	117.52	111.03
1	A	51	VAL	CA-C-N	5.68	125.58	119.85
1	A	51	VAL	C-N-CA	5.68	125.58	119.85
1	C	51	VAL	CA-C-N	5.51	125.42	119.85
1	C	51	VAL	C-N-CA	5.51	125.42	119.85
1	D	215	TRP	N-CA-C	5.38	117.96	111.40
1	B	233	ILE	N-CA-C	5.21	116.19	107.18
1	A	127	GLU	N-CA-C	-5.14	98.94	107.99
1	B	200	ILE	CA-C-N	5.04	126.16	120.66
1	B	200	ILE	C-N-CA	5.04	126.16	120.66

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	1906	0	1854	80	0
1	B	1906	0	1855	78	0
1	C	1906	0	1855	87	0
1	D	1906	0	1855	80	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	25	0	11	5	0
3	B	25	0	11	6	0
3	C	25	0	11	5	0
3	D	25	0	11	4	0
4	A	20	0	12	1	0
4	B	20	0	12	1	0
4	C	20	0	12	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	D	20	0	12	1	0
All	All	7808	0	7511	303	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 20.

All (303) 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:222:LYS:HA	1:A:222:LYS:HE2	1.54	0.88
1:D:222:LYS:HA	1:D:222:LYS:HE2	1.56	0.87
1:B:222:LYS:HE2	1:B:222:LYS:HA	1.56	0.84
1:C:222:LYS:HE2	1:C:222:LYS:HA	1.58	0.84
1:D:53:GLU:HG2	1:D:104:ARG:HH12	1.44	0.80
1:B:43:GLN:NE2	1:C:195:ASN:HB2	1.97	0.79
1:D:128:ARG:HG3	1:D:128:ARG:HH11	1.49	0.78
1:C:53:GLU:HG2	1:C:104:ARG:HH12	1.49	0.78
1:C:20:ARG:HD2	1:C:20:ARG:O	1.84	0.77
1:A:53:GLU:HG2	1:A:104:ARG:HH12	1.51	0.76
1:B:53:GLU:HG2	1:B:104:ARG:HH12	1.51	0.75
1:B:152:TRP:O	1:B:156:GLN:HG3	1.87	0.74
1:B:43:GLN:HE22	1:C:195:ASN:HB2	1.53	0.73
1:C:152:TRP:O	1:C:156:GLN:HG3	1.88	0.72
1:A:152:TRP:O	1:A:156:GLN:HG3	1.90	0.72
1:C:241:ASP:HA	3:C:402:UDP:O4	1.91	0.70
1:A:34:LYS:HD3	3:A:402:UDP:O2B	1.92	0.70
1:D:152:TRP:O	1:D:156:GLN:HG3	1.91	0.69
1:B:128:ARG:HG3	1:B:128:ARG:HH11	1.57	0.69
1:C:242:PHE:O	1:C:244:ASP:N	2.27	0.68
1:B:242:PHE:O	1:B:244:ASP:N	2.27	0.67
1:C:77:ASN:HD22	1:C:78:GLY:N	1.93	0.67
1:B:43:GLN:NE2	1:C:195:ASN:CG	2.53	0.67
1:A:242:PHE:O	1:A:244:ASP:N	2.27	0.67
1:D:242:PHE:O	1:D:244:ASP:N	2.29	0.66
1:C:165:GLN:HG3	1:D:166:PHE:HE2	1.61	0.66
1:B:34:LYS:HD3	3:B:402:UDP:O2B	1.95	0.66
1:D:128:ARG:HG3	1:D:128:ARG:NH1	2.09	0.66
1:A:128:ARG:HH11	1:A:128:ARG:HG3	1.59	0.65
1:C:242:PHE:C	1:C:244:ASP:H	2.04	0.65
1:C:128:ARG:HG3	1:C:128:ARG:HH11	1.62	0.64
1:C:166:PHE:HE2	1:D:165:GLN:HG3	1.63	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:222:LYS:HA	1:C:222:LYS:CE	2.28	0.64
1:D:243:LYS:O	1:D:243:LYS:HG2	1.98	0.64
1:C:245:LYS:O	1:C:245:LYS:HG2	1.98	0.64
1:B:43:GLN:NE2	1:C:195:ASN:CB	2.60	0.63
1:D:222:LYS:HA	1:D:222:LYS:CE	2.27	0.63
1:C:243:LYS:HG2	1:C:243:LYS:O	1.97	0.63
1:A:242:PHE:C	1:A:244:ASP:H	2.06	0.63
1:B:245:LYS:O	1:B:245:LYS:HG2	1.99	0.63
1:C:242:PHE:O	1:C:246:TYR:HB3	1.99	0.63
1:C:223:THR:HG21	1:C:228:LEU:HD12	1.80	0.63
1:B:242:PHE:C	1:B:244:ASP:H	2.04	0.62
1:A:222:LYS:HA	1:A:222:LYS:CE	2.26	0.62
1:A:243:LYS:O	1:A:243:LYS:HG2	1.98	0.62
1:D:242:PHE:C	1:D:244:ASP:H	2.07	0.62
1:B:243:LYS:O	1:B:243:LYS:HG2	1.97	0.62
1:A:242:PHE:O	1:A:246:TYR:HB3	2.00	0.61
1:A:128:ARG:HG3	1:A:128:ARG:NH1	2.15	0.61
1:B:242:PHE:O	1:B:246:TYR:HB3	2.00	0.61
1:D:34:LYS:N	3:D:402:UDP:O3B	2.33	0.61
1:A:223:THR:HG21	1:A:228:LEU:HD12	1.83	0.61
1:D:245:LYS:O	1:D:245:LYS:HG2	2.01	0.61
1:A:245:LYS:HG2	1:A:245:LYS:O	1.99	0.61
1:B:128:ARG:HG3	1:B:128:ARG:NH1	2.16	0.60
1:C:238:VAL:HB	1:C:242:PHE:CZ	2.36	0.60
1:D:238:VAL:HB	1:D:242:PHE:CZ	2.37	0.60
1:B:222:LYS:HA	1:B:222:LYS:CE	2.28	0.60
1:C:77:ASN:HD22	1:C:77:ASN:C	2.08	0.59
1:D:223:THR:HG21	1:D:228:LEU:HD12	1.83	0.59
1:B:238:VAL:HB	1:B:242:PHE:CZ	2.37	0.59
1:D:242:PHE:O	1:D:246:TYR:HB3	2.02	0.59
1:D:44:LEU:O	1:D:45:CYS:HB2	2.03	0.59
1:B:120:GLU:HG3	1:B:121:LYS:H	1.68	0.59
1:B:34:LYS:N	3:B:402:UDP:O3B	2.36	0.58
1:B:223:THR:HG21	1:B:228:LEU:HD12	1.85	0.58
1:D:120:GLU:HG3	1:D:121:LYS:H	1.68	0.58
1:C:120:GLU:HG3	1:C:121:LYS:H	1.68	0.58
1:A:120:GLU:HG3	1:A:121:LYS:H	1.68	0.58
1:C:128:ARG:HG3	1:C:128:ARG:NH1	2.18	0.58
1:A:35:SER:N	3:A:402:UDP:O3B	2.37	0.58
1:B:37:PHE:HB2	1:B:242:PHE:HE2	1.67	0.57
1:A:238:VAL:HB	1:A:242:PHE:CZ	2.39	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:44:LEU:O	1:B:45:CYS:HB2	2.02	0.57
1:C:44:LEU:O	1:C:45:CYS:HB2	2.04	0.57
1:B:193:GLY:O	1:B:194:ARG:C	2.48	0.57
1:C:242:PHE:C	1:C:244:ASP:N	2.62	0.56
1:A:44:LEU:O	1:A:45:CYS:HB2	2.05	0.56
1:A:83:GLN:HG2	1:A:87:GLU:OE2	2.06	0.56
1:A:37:PHE:HB2	1:A:242:PHE:HE2	1.70	0.56
1:D:83:GLN:HG2	1:D:87:GLU:OE2	2.07	0.55
1:C:37:PHE:HB2	1:C:242:PHE:HE2	1.71	0.55
1:D:37:PHE:HB2	1:D:242:PHE:HE2	1.72	0.55
1:C:106:ARG:NH1	1:D:157:ASP:OD2	2.39	0.54
1:B:43:GLN:HE22	1:C:195:ASN:CB	2.17	0.54
1:D:80:ASN:O	1:D:84:MET:HG3	2.07	0.54
1:D:242:PHE:C	1:D:244:ASP:N	2.64	0.54
1:D:34:LYS:HD3	3:D:402:UDP:O2B	2.08	0.54
1:C:157:ASP:OD2	1:D:106:ARG:NH1	2.41	0.54
1:D:241:ASP:HA	3:D:402:UDP:O4	2.08	0.54
1:D:31:ALA:HA	3:D:402:UDP:O3A	2.07	0.53
1:A:20:ARG:HD2	1:A:20:ARG:O	2.09	0.53
1:C:109:LEU:HD11	1:C:166:PHE:HB3	1.90	0.53
1:D:193:GLY:O	1:D:194:ARG:C	2.51	0.53
1:D:60:ASN:OD1	1:D:78:GLY:N	2.41	0.53
1:A:242:PHE:C	1:A:244:ASP:N	2.63	0.53
1:B:128:ARG:NH1	4:B:403:DCM:O2P	2.38	0.53
1:C:193:GLY:O	1:C:194:ARG:C	2.52	0.53
1:B:115:LYS:NZ	3:C:402:UDP:H5'2	2.24	0.53
1:D:53:GLU:OE1	1:D:128:ARG:NH1	2.40	0.53
1:D:109:LEU:HD11	1:D:166:PHE:HB3	1.91	0.53
1:A:193:GLY:O	1:A:194:ARG:C	2.52	0.53
1:C:47:ASP:OD2	1:C:119:ALA:HB1	2.09	0.53
1:B:20:ARG:O	1:B:20:ARG:HG2	2.08	0.52
1:A:61:VAL:HG22	1:A:79:GLY:H	1.73	0.52
1:B:47:ASP:OD2	1:B:119:ALA:HB1	2.09	0.52
1:C:83:GLN:HG2	1:C:87:GLU:OE2	2.09	0.52
1:D:53:GLU:HG2	1:D:104:ARG:NH1	2.20	0.52
1:C:166:PHE:HZ	1:D:166:PHE:HZ	1.56	0.52
1:B:83:GLN:HG2	1:B:87:GLU:OE2	2.09	0.52
1:A:109:LEU:HD11	1:A:166:PHE:HB3	1.92	0.52
1:B:90:GLU:OE1	1:B:90:GLU:N	2.39	0.51
1:A:47:ASP:OD2	1:A:119:ALA:HB1	2.11	0.51
1:A:53:GLU:OE1	1:A:128:ARG:NH1	2.41	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:53:GLU:HG2	1:B:104:ARG:NH1	2.24	0.51
1:B:109:LEU:HD11	1:B:166:PHE:HB3	1.91	0.51
1:A:148:ASN:HD21	1:B:92:TRP:HH2	1.58	0.51
1:A:49:GLU:HB2	1:A:116:LEU:HD22	1.93	0.51
1:A:120:GLU:HG3	1:A:121:LYS:N	2.26	0.51
1:A:158:TRP:HD1	1:B:102:LEU:HD22	1.75	0.51
1:C:34:LYS:CD	1:C:34:LYS:H	2.24	0.51
1:C:227:TYR:C	1:C:229:GLN:N	2.68	0.51
1:D:49:GLU:HB2	1:D:116:LEU:HD22	1.93	0.51
1:A:166:PHE:O	1:A:168:GLN:N	2.44	0.51
1:A:238:VAL:HB	1:A:242:PHE:CE1	2.47	0.50
1:B:34:LYS:H	1:B:34:LYS:CD	2.23	0.50
1:B:242:PHE:C	1:B:244:ASP:N	2.63	0.50
1:C:53:GLU:OE1	1:C:128:ARG:NH1	2.42	0.50
1:D:127:GLU:O	1:D:128:ARG:HB2	2.09	0.50
1:C:238:VAL:HB	1:C:242:PHE:CE1	2.46	0.50
1:A:241:ASP:HA	3:A:402:UDP:O4	2.12	0.50
1:B:238:VAL:HB	1:B:242:PHE:CE1	2.47	0.50
1:D:90:GLU:OE1	1:D:90:GLU:N	2.42	0.50
1:B:192:ARG:HG2	1:B:193:GLY:N	2.27	0.50
1:D:34:LYS:CD	1:D:34:LYS:H	2.24	0.50
1:B:53:GLU:OE1	1:B:128:ARG:NH1	2.41	0.50
1:C:166:PHE:O	1:C:168:GLN:N	2.45	0.50
1:A:34:LYS:CD	1:A:34:LYS:H	2.24	0.50
1:B:166:PHE:O	1:B:168:GLN:N	2.45	0.50
1:C:120:GLU:HG3	1:C:121:LYS:N	2.27	0.49
1:D:47:ASP:OD2	1:D:119:ALA:HB1	2.12	0.49
1:D:120:GLU:HG3	1:D:121:LYS:N	2.26	0.49
1:B:49:GLU:HB2	1:B:116:LEU:HD22	1.94	0.49
1:D:166:PHE:O	1:D:168:GLN:N	2.44	0.49
1:C:46:GLU:N	1:C:46:GLU:OE2	2.45	0.49
1:C:148:ASN:HD21	1:D:92:TRP:HH2	1.60	0.49
1:A:90:GLU:OE1	1:A:90:GLU:N	2.41	0.49
1:B:120:GLU:HG3	1:B:121:LYS:N	2.27	0.49
1:C:34:LYS:N	3:C:402:UDP:O3B	2.46	0.49
1:C:132:SER:O	1:C:136:ILE:HB	2.13	0.49
1:D:46:GLU:OE2	1:D:46:GLU:N	2.45	0.49
1:C:49:GLU:HB2	1:C:116:LEU:HD22	1.95	0.49
1:C:203:GLU:CD	1:C:203:GLU:H	2.21	0.49
1:D:53:GLU:OE1	1:D:128:ARG:HG3	2.13	0.49
1:C:80:ASN:O	1:C:84:MET:HG3	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:21:ILE:HG23	1:D:122:PRO:HB2	1.94	0.48
1:C:90:GLU:OE1	1:C:90:GLU:N	2.44	0.48
1:D:238:VAL:HB	1:D:242:PHE:CE1	2.48	0.48
1:A:221:LEU:C	1:A:221:LEU:HD23	2.38	0.48
1:A:227:TYR:C	1:A:229:GLN:N	2.71	0.48
1:A:46:GLU:N	1:A:46:GLU:OE2	2.46	0.48
1:C:35:SER:O	1:C:38:VAL:HG12	2.14	0.48
1:C:47:ASP:O	1:C:48:TRP:HD1	1.97	0.48
1:A:34:LYS:N	3:A:402:UDP:O3B	2.47	0.47
1:C:192:ARG:HG2	1:C:193:GLY:N	2.28	0.47
1:C:158:TRP:HE1	1:D:158:TRP:HE1	1.61	0.47
1:D:132:SER:O	1:D:136:ILE:HB	2.14	0.47
1:A:128:ARG:NH1	4:A:403:DCM:O2P	2.46	0.47
1:B:80:ASN:O	1:B:84:MET:HG3	2.13	0.47
1:B:203:GLU:CD	1:B:203:GLU:H	2.22	0.47
1:C:77:ASN:C	1:C:77:ASN:ND2	2.73	0.47
1:C:166:PHE:HZ	1:D:166:PHE:CZ	2.32	0.47
1:C:166:PHE:CZ	1:D:166:PHE:CZ	3.02	0.47
1:C:92:TRP:HH2	1:D:148:ASN:HD21	1.61	0.47
1:B:46:GLU:OE2	1:B:46:GLU:N	2.46	0.47
1:A:127:GLU:O	1:A:128:ARG:HB2	2.14	0.47
1:B:21:ILE:HG23	1:B:122:PRO:HB2	1.96	0.47
1:C:110:ALA:C	1:C:112:LEU:N	2.72	0.47
1:A:53:GLU:HG2	1:A:104:ARG:NH1	2.24	0.47
1:B:35:SER:O	1:B:38:VAL:HG12	2.14	0.47
1:A:50:VAL:HG12	1:A:51:VAL:N	2.29	0.46
1:A:110:ALA:C	1:A:112:LEU:N	2.73	0.46
1:D:128:ARG:NH1	1:D:128:ARG:CG	2.76	0.46
1:D:47:ASP:O	1:D:48:TRP:HD1	1.98	0.46
1:A:21:ILE:HG23	1:A:122:PRO:HB2	1.96	0.46
1:C:21:ILE:HG23	1:C:122:PRO:HB2	1.98	0.46
1:C:127:GLU:O	1:C:128:ARG:HB2	2.16	0.46
1:D:89:PRO:C	1:D:91:ARG:H	2.24	0.46
1:D:192:ARG:HG2	1:D:193:GLY:N	2.29	0.46
1:A:47:ASP:O	1:A:48:TRP:HD1	1.97	0.46
1:A:61:VAL:O	1:A:62:GLN:OE1	2.34	0.46
1:A:203:GLU:CD	1:A:203:GLU:H	2.23	0.46
1:A:92:TRP:HH2	1:B:148:ASN:HD21	1.64	0.46
1:D:30:ILE:O	1:D:31:ALA:HB3	2.16	0.46
1:A:136:ILE:HG21	1:A:212:HIS:CE1	2.51	0.46
1:B:115:LYS:HZ1	3:C:402:UDP:H5'2	1.80	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:40:ILE:HG21	1:D:246:TYR:CD1	2.51	0.46
1:D:86:TYR:HH	4:D:403:DCM:HO3'	1.63	0.46
1:D:203:GLU:H	1:D:203:GLU:CD	2.24	0.46
1:A:242:PHE:N	3:A:402:UDP:O4	2.49	0.45
1:C:89:PRO:C	1:C:91:ARG:H	2.25	0.45
1:D:227:TYR:C	1:D:229:GLN:N	2.72	0.45
1:A:35:SER:O	1:A:38:VAL:HG12	2.15	0.45
1:A:58:TRP:O	1:A:81:VAL:HG12	2.17	0.45
1:A:192:ARG:HG2	1:A:193:GLY:N	2.30	0.45
1:A:20:ARG:HD2	1:A:20:ARG:C	2.41	0.45
1:A:104:ARG:O	1:A:108:GLN:HG3	2.16	0.45
1:B:47:ASP:O	1:B:48:TRP:HD1	1.98	0.45
1:B:127:GLU:O	1:B:128:ARG:HB2	2.17	0.45
1:C:20:ARG:HD2	1:C:20:ARG:C	2.41	0.45
1:C:136:ILE:HG21	1:C:212:HIS:CE1	2.51	0.45
1:B:62:GLN:O	1:B:63:SER:HB2	2.17	0.45
1:C:50:VAL:HG12	1:C:51:VAL:N	2.30	0.45
1:D:231:VAL:HA	1:D:232:PRO:HD3	1.78	0.45
1:B:89:PRO:C	1:B:91:ARG:H	2.24	0.45
1:A:132:SER:O	1:A:136:ILE:HB	2.16	0.44
1:C:166:PHE:CE2	1:D:165:GLN:HG3	2.48	0.44
1:B:50:VAL:HG12	1:B:51:VAL:N	2.32	0.44
1:D:110:ALA:C	1:D:112:LEU:N	2.75	0.44
1:A:61:VAL:O	1:A:62:GLN:HG2	2.17	0.44
1:C:53:GLU:HG2	1:C:104:ARG:NH1	2.24	0.44
1:A:89:PRO:C	1:A:91:ARG:H	2.25	0.44
1:C:31:ALA:HA	3:C:402:UDP:O3A	2.18	0.44
1:D:136:ILE:CD1	1:D:211:LYS:HB3	2.48	0.44
1:B:53:GLU:OE1	1:B:128:ARG:HG3	2.18	0.44
1:B:110:ALA:C	1:B:112:LEU:N	2.72	0.44
1:B:227:TYR:C	1:B:229:GLN:N	2.71	0.44
1:B:89:PRO:C	1:B:91:ARG:N	2.76	0.44
1:D:35:SER:O	1:D:38:VAL:HG12	2.17	0.44
1:C:227:TYR:C	1:C:229:GLN:H	2.24	0.43
1:D:104:ARG:O	1:D:108:GLN:HG3	2.18	0.43
1:B:231:VAL:HA	1:B:232:PRO:HD3	1.78	0.43
1:D:50:VAL:HG12	1:D:51:VAL:N	2.32	0.43
1:A:91:ARG:HD2	1:A:91:ARG:O	2.19	0.43
1:A:110:ALA:C	1:A:112:LEU:H	2.27	0.43
1:B:110:ALA:C	1:B:112:LEU:H	2.25	0.43
1:B:221:LEU:HD23	1:B:221:LEU:C	2.44	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:148:ASN:O	1:D:149:GLU:C	2.61	0.43
1:D:221:LEU:C	1:D:221:LEU:HD23	2.43	0.43
1:A:223:THR:HB	1:A:224:ASN:H	1.39	0.43
1:A:227:TYR:C	1:A:229:GLN:H	2.26	0.43
1:C:30:ILE:O	1:C:31:ALA:HB3	2.18	0.43
1:C:225:PHE:CD1	1:C:225:PHE:N	2.87	0.43
1:B:31:ALA:HA	3:B:402:UDP:O3A	2.19	0.43
1:B:104:ARG:O	1:B:108:GLN:HG3	2.18	0.43
1:C:89:PRO:C	1:C:91:ARG:N	2.75	0.43
1:C:104:ARG:O	1:C:108:GLN:HG3	2.19	0.43
1:A:34:LYS:O	1:A:35:SER:C	2.62	0.43
1:B:242:PHE:N	3:B:402:UDP:O4	2.52	0.43
1:A:166:PHE:C	1:A:168:GLN:N	2.77	0.42
1:A:89:PRO:C	1:A:91:ARG:N	2.77	0.42
1:D:166:PHE:C	1:D:168:GLN:N	2.77	0.42
1:A:231:VAL:HA	1:A:232:PRO:HD3	1.79	0.42
1:B:136:ILE:CD1	1:B:211:LYS:HB3	2.49	0.42
1:C:110:ALA:C	1:C:112:LEU:H	2.27	0.42
1:C:128:ARG:NH1	4:C:403:DCM:O2P	2.50	0.42
1:D:89:PRO:C	1:D:91:ARG:N	2.75	0.42
1:B:227:TYR:C	1:B:229:GLN:H	2.27	0.42
1:B:51:VAL:HG23	1:B:124:LEU:HD11	2.02	0.42
1:B:58:TRP:O	1:B:81:VAL:HG12	2.19	0.42
1:B:132:SER:O	1:B:136:ILE:HB	2.19	0.42
1:C:166:PHE:CZ	1:D:166:PHE:HZ	2.35	0.42
1:B:136:ILE:HG21	1:B:212:HIS:CE1	2.55	0.42
1:C:53:GLU:OE1	1:C:128:ARG:HG3	2.20	0.42
1:A:194:ARG:C	1:A:196:GLU:H	2.28	0.42
1:A:245:LYS:O	1:A:247:GLU:N	2.53	0.42
1:B:166:PHE:C	1:B:168:GLN:N	2.78	0.42
1:C:130:VAL:HG21	1:C:163:ASN:HD21	1.85	0.42
1:C:224:ASN:ND2	1:C:224:ASN:O	2.53	0.42
1:A:225:PHE:CD1	1:A:225:PHE:N	2.88	0.42
1:B:130:VAL:HG21	1:B:163:ASN:HD21	1.85	0.42
1:A:136:ILE:CD1	1:A:211:LYS:HB3	2.49	0.41
1:C:221:LEU:C	1:C:221:LEU:HD23	2.45	0.41
1:D:58:TRP:O	1:D:81:VAL:HG12	2.19	0.41
1:D:136:ILE:HG21	1:D:212:HIS:CE1	2.54	0.41
1:A:80:ASN:O	1:A:84:MET:HG3	2.20	0.41
1:B:35:SER:N	3:B:402:UDP:O3B	2.51	0.41
1:B:241:ASP:HA	3:B:402:UDP:O4	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:102:LEU:HD22	1:B:158:TRP:HD1	1.85	0.41
1:A:112:LEU:HA	1:A:112:LEU:HD23	1.87	0.41
1:D:130:VAL:HG21	1:D:163:ASN:HD21	1.86	0.41
1:A:50:VAL:CG1	1:A:51:VAL:N	2.84	0.41
1:C:148:ASN:ND2	1:D:92:TRP:HH2	2.19	0.41
1:C:194:ARG:C	1:C:196:GLU:H	2.29	0.41
1:D:194:ARG:C	1:D:196:GLU:H	2.29	0.41
1:A:30:ILE:O	1:A:31:ALA:HB3	2.21	0.41
1:A:39:ASN:O	1:A:43:GLN:HG2	2.21	0.41
1:A:166:PHE:C	1:A:168:GLN:H	2.29	0.41
1:B:30:ILE:O	1:B:31:ALA:HB3	2.20	0.41
1:B:39:ASN:O	1:B:43:GLN:HG2	2.21	0.41
1:B:246:TYR:CD1	1:B:246:TYR:C	2.99	0.41
1:C:34:LYS:O	1:C:35:SER:C	2.63	0.41
1:D:110:ALA:C	1:D:112:LEU:H	2.29	0.41
1:D:227:TYR:C	1:D:229:GLN:H	2.27	0.41
1:C:166:PHE:C	1:C:168:GLN:H	2.29	0.41
1:A:53:GLU:OE1	1:A:128:ARG:HG3	2.21	0.40
1:C:223:THR:HB	1:C:224:ASN:H	1.38	0.40
1:C:231:VAL:HA	1:C:232:PRO:HD3	1.79	0.40
1:D:91:ARG:HD2	1:D:91:ARG:O	2.21	0.40
1:C:117:LYS:HB3	1:C:117:LYS:HE2	1.91	0.40
1:A:246:TYR:CD1	1:A:246:TYR:C	3.00	0.40
1:C:58:TRP:O	1:C:81:VAL:HG12	2.20	0.40
1:D:34:LYS:O	1:D:35:SER:C	2.64	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	224/280 (80%)	184 (82%)	32 (14%)	8 (4%)	2 16

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	224/280 (80%)	185 (83%)	33 (15%)	6 (3%)	4	20
1	C	224/280 (80%)	188 (84%)	30 (13%)	6 (3%)	4	20
1	D	224/280 (80%)	187 (84%)	29 (13%)	8 (4%)	2	16
All	All	896/1120 (80%)	744 (83%)	124 (14%)	28 (3%)	3	18

All (28) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	194	ARG
1	A	243	LYS
1	B	194	ARG
1	B	243	LYS
1	C	194	ARG
1	C	243	LYS
1	D	194	ARG
1	D	243	LYS
1	A	78	GLY
1	A	167	GLY
1	A	226	ASP
1	B	167	GLY
1	B	226	ASP
1	C	167	GLY
1	C	226	ASP
1	D	167	GLY
1	D	226	ASP
1	A	246	TYR
1	B	45	CYS
1	B	246	TYR
1	D	45	CYS
1	D	241	ASP
1	A	45	CYS
1	C	45	CYS
1	C	246	TYR
1	D	246	TYR
1	A	195	ASN
1	D	195	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	210/256 (82%)	197 (94%)	13 (6%)	16	43
1	B	210/256 (82%)	199 (95%)	11 (5%)	21	48
1	C	210/256 (82%)	197 (94%)	13 (6%)	16	43
1	D	210/256 (82%)	197 (94%)	13 (6%)	16	43
All	All	840/1024 (82%)	790 (94%)	50 (6%)	17	43

All (50) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	ARG
1	A	34	LYS
1	A	44	LEU
1	A	49	GLU
1	A	57	ARG
1	A	112	LEU
1	A	122	PRO
1	A	133	ASP
1	A	166	PHE
1	A	168	GLN
1	A	222	LYS
1	A	241	ASP
1	A	246	TYR
1	B	34	LYS
1	B	44	LEU
1	B	49	GLU
1	B	57	ARG
1	B	112	LEU
1	B	133	ASP
1	B	166	PHE
1	B	168	GLN
1	B	222	LYS
1	B	241	ASP
1	B	246	TYR
1	C	20	ARG
1	C	34	LYS
1	C	44	LEU
1	C	49	GLU

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Mol	Chain	Res	Type
1	C	57	ARG
1	C	112	LEU
1	C	122	PRO
1	C	133	ASP
1	C	166	PHE
1	C	168	GLN
1	C	222	LYS
1	C	241	ASP
1	C	246	TYR
1	D	34	LYS
1	D	44	LEU
1	D	49	GLU
1	D	57	ARG
1	D	63	SER
1	D	112	LEU
1	D	122	PRO
1	D	133	ASP
1	D	166	PHE
1	D	168	GLN
1	D	222	LYS
1	D	241	ASP
1	D	246	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (31) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	43	GLN
1	A	77	ASN
1	A	156	GLN
1	A	163	ASN
1	A	187	HIS
1	A	209	HIS
1	A	224	ASN
1	B	43	GLN
1	B	77	ASN
1	B	140	ASN
1	B	156	GLN
1	B	163	ASN
1	B	187	HIS
1	B	209	HIS
1	B	224	ASN
1	C	43	GLN

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Mol	Chain	Res	Type
1	C	62	GLN
1	C	77	ASN
1	C	156	GLN
1	C	163	ASN
1	C	165	GLN
1	C	187	HIS
1	C	209	HIS
1	C	224	ASN
1	D	43	GLN
1	D	156	GLN
1	D	163	ASN
1	D	165	GLN
1	D	187	HIS
1	D	209	HIS
1	D	224	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](#)

Of 12 ligands modelled in this entry, 4 are monoatomic - leaving 8 for Mogul analysis.

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
4	DCM	B	403	-	21,21,21	1.28	4 (19%)	30,31,31	0.91	0
3	UDP	C	402	2	25,26,26	1.33	2 (8%)	38,40,40	1.63	6 (15%)
4	DCM	C	403	-	21,21,21	1.37	4 (19%)	30,31,31	0.91	0
4	DCM	A	403	-	21,21,21	1.35	4 (19%)	30,31,31	0.92	0
3	UDP	A	402	-	25,26,26	1.11	3 (12%)	38,40,40	1.63	6 (15%)
3	UDP	D	402	-	25,26,26	1.64	4 (16%)	38,40,40	1.62	6 (15%)
4	DCM	D	403	-	21,21,21	1.39	4 (19%)	30,31,31	0.90	0
3	UDP	B	402	-	25,26,26	1.23	2 (8%)	38,40,40	1.62	6 (15%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	DCM	B	403	-	-	0/10/22/22	0/2/2/2
3	UDP	C	402	2	-	0/16/32/32	0/2/2/2
4	DCM	C	403	-	-	0/10/22/22	0/2/2/2
4	DCM	A	403	-	-	0/10/22/22	0/2/2/2
3	UDP	A	402	-	-	0/16/32/32	0/2/2/2
3	UDP	D	402	-	-	0/16/32/32	0/2/2/2
4	DCM	D	403	-	-	0/10/22/22	0/2/2/2
3	UDP	B	402	-	-	0/16/32/32	0/2/2/2

All (27) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	402	UDP	PA-O3A	5.38	1.65	1.59
3	C	402	UDP	PA-O3A	4.82	1.64	1.59
4	D	403	DCM	P-O3P	3.15	1.66	1.54
3	B	402	UDP	PA-O3A	3.08	1.62	1.59
4	C	403	DCM	P-O3P	2.88	1.65	1.54
4	A	403	DCM	P-O3P	2.74	1.65	1.54
4	B	403	DCM	P-O3P	2.71	1.64	1.54
3	A	402	UDP	PB-O2B	-2.67	1.44	1.54
3	D	402	UDP	O5'-C5'	-2.66	1.34	1.44
4	B	403	DCM	O4'-C4'	2.51	1.50	1.45
4	C	403	DCM	O4'-C4'	2.51	1.50	1.45
3	A	402	UDP	PA-O3A	2.49	1.62	1.59
3	B	402	UDP	PB-O2B	-2.48	1.45	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	403	DCM	O4'-C4'	2.37	1.50	1.45
4	A	403	DCM	O4'-C4'	2.37	1.50	1.45
4	A	403	DCM	O4'-C1'	2.25	1.47	1.42
4	A	403	DCM	P-O2P	-2.23	1.46	1.54
4	B	403	DCM	P-O2P	-2.22	1.46	1.54
4	D	403	DCM	P-O2P	-2.21	1.46	1.54
4	D	403	DCM	O4'-C1'	2.15	1.47	1.42
4	C	403	DCM	P-O2P	-2.15	1.46	1.54
3	D	402	UDP	C5'-C4'	-2.14	1.45	1.51
4	C	403	DCM	O4'-C1'	2.10	1.47	1.42
3	C	402	UDP	O5'-C5'	-2.08	1.36	1.44
4	B	403	DCM	O4'-C1'	2.08	1.47	1.42
3	A	402	UDP	O4-C4	-2.07	1.20	1.24
3	D	402	UDP	PB-O2B	-2.05	1.47	1.54

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	402	UDP	C4-N3-C2	-5.18	120.19	126.61
3	C	402	UDP	C4-N3-C2	-5.12	120.25	126.61
3	D	402	UDP	C4-N3-C2	-5.11	120.26	126.61
3	B	402	UDP	C4-N3-C2	-5.10	120.28	126.61
3	C	402	UDP	N3-C2-N1	4.58	120.86	114.89
3	D	402	UDP	N3-C2-N1	4.57	120.84	114.89
3	A	402	UDP	N3-C2-N1	4.56	120.83	114.89
3	B	402	UDP	N3-C2-N1	4.51	120.76	114.89
3	A	402	UDP	C5-C4-N3	3.32	119.45	114.80
3	D	402	UDP	C5-C4-N3	3.31	119.44	114.80
3	C	402	UDP	C5-C4-N3	3.31	119.44	114.80
3	B	402	UDP	C5-C4-N3	3.29	119.41	114.80
3	B	402	UDP	O4-C4-C5	-2.84	120.27	125.16
3	A	402	UDP	O4-C4-C5	-2.83	120.28	125.16
3	D	402	UDP	O4-C4-C5	-2.81	120.31	125.16
3	C	402	UDP	O4-C4-C5	-2.81	120.31	125.16
3	B	402	UDP	PA-O5'-C5'	-2.52	106.89	121.35
3	A	402	UDP	PA-O5'-C5'	-2.52	106.93	121.35
3	C	402	UDP	PA-O5'-C5'	-2.49	107.06	121.35
3	D	402	UDP	PA-O5'-C5'	-2.46	107.26	121.35
3	A	402	UDP	O2-C2-N1	-2.43	119.63	122.80
3	B	402	UDP	O2-C2-N1	-2.41	119.66	122.80
3	C	402	UDP	O2-C2-N1	-2.41	119.66	122.80
3	D	402	UDP	O2-C2-N1	-2.39	119.68	122.80

There are no chirality outliers.

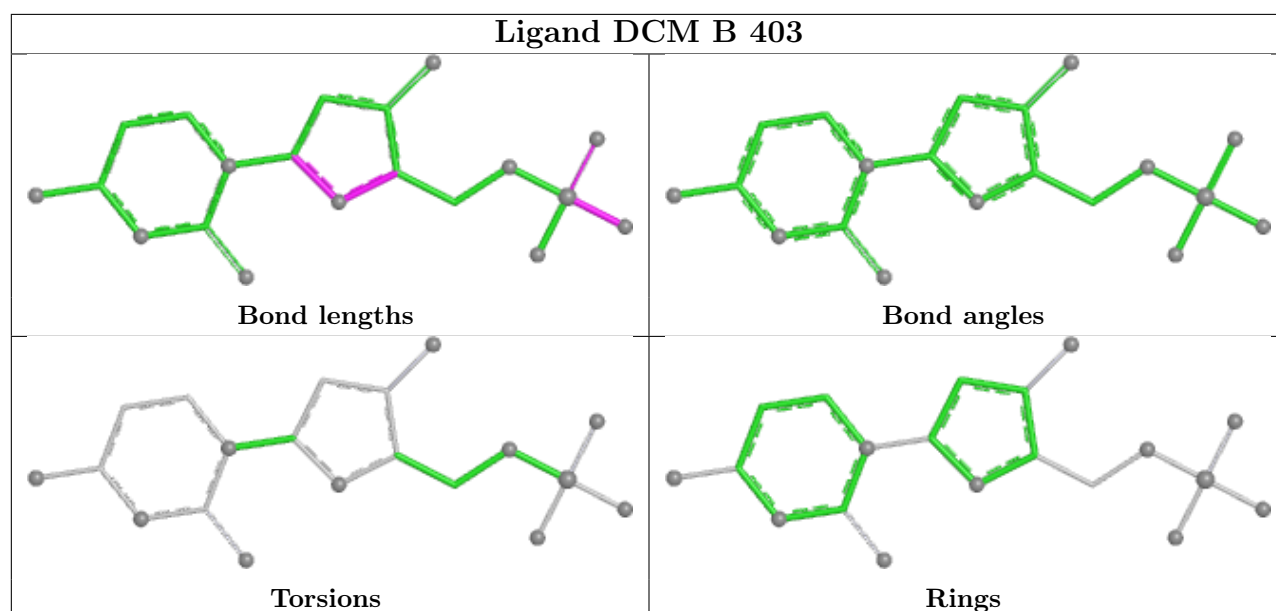
There are no torsion outliers.

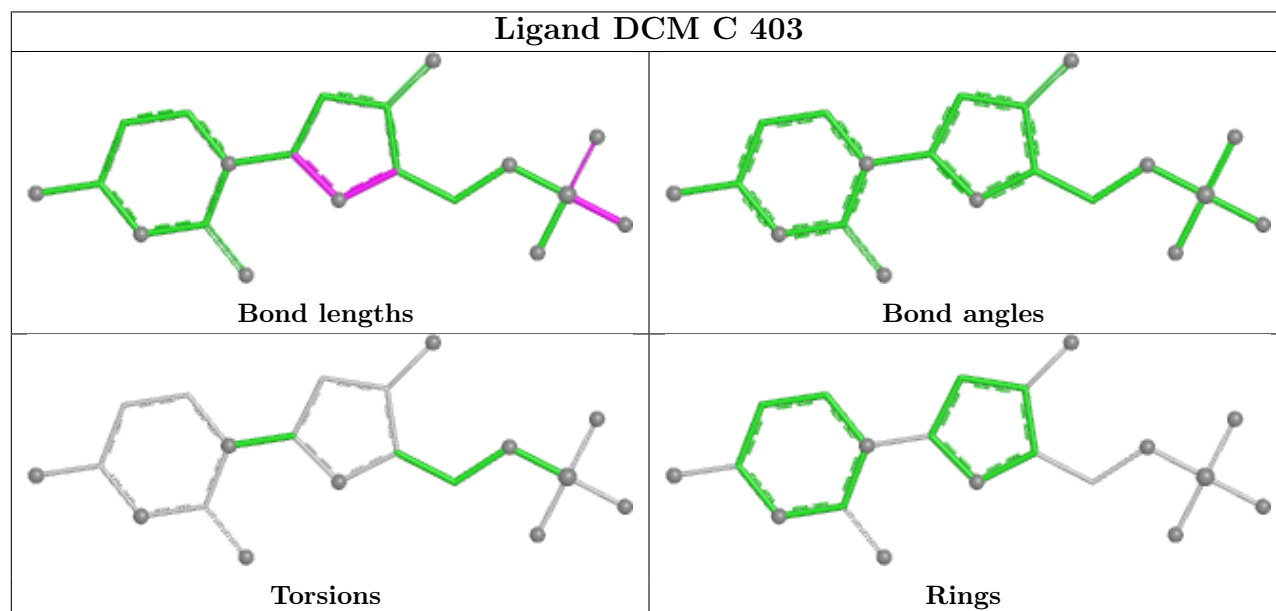
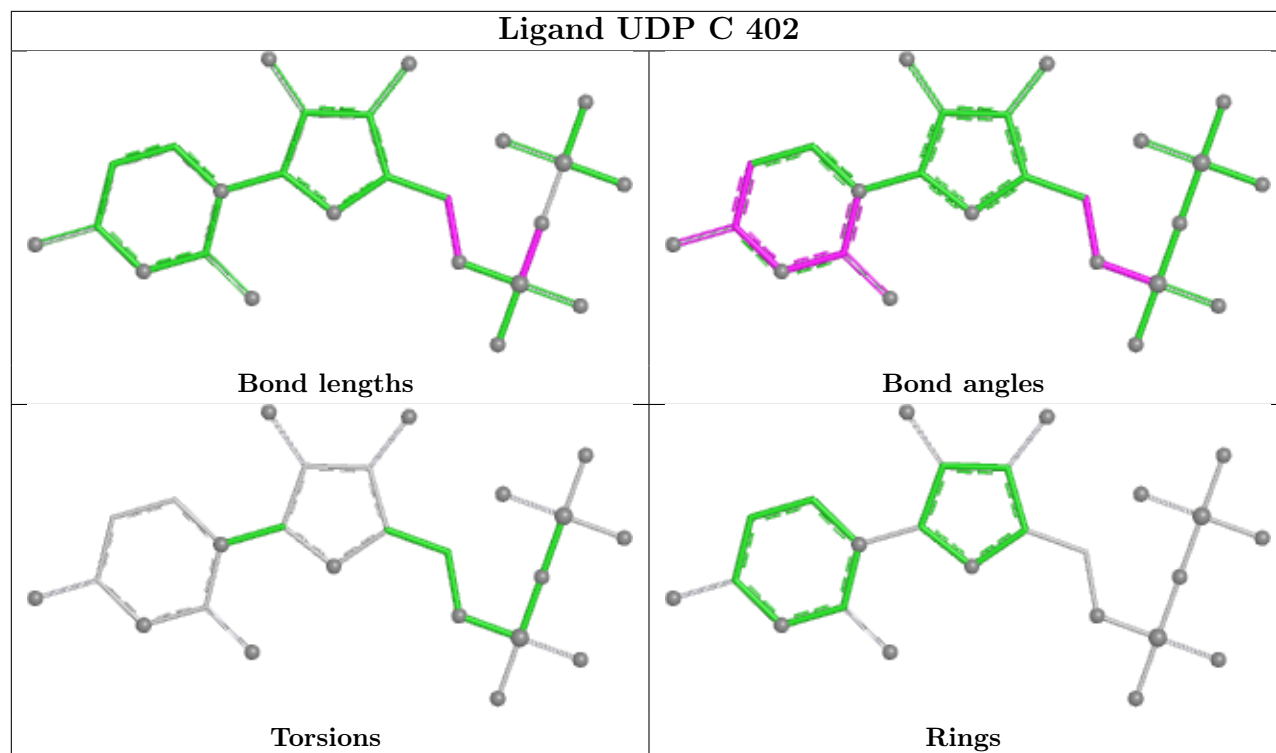
There are no ring outliers.

8 monomers are involved in 24 short contacts:

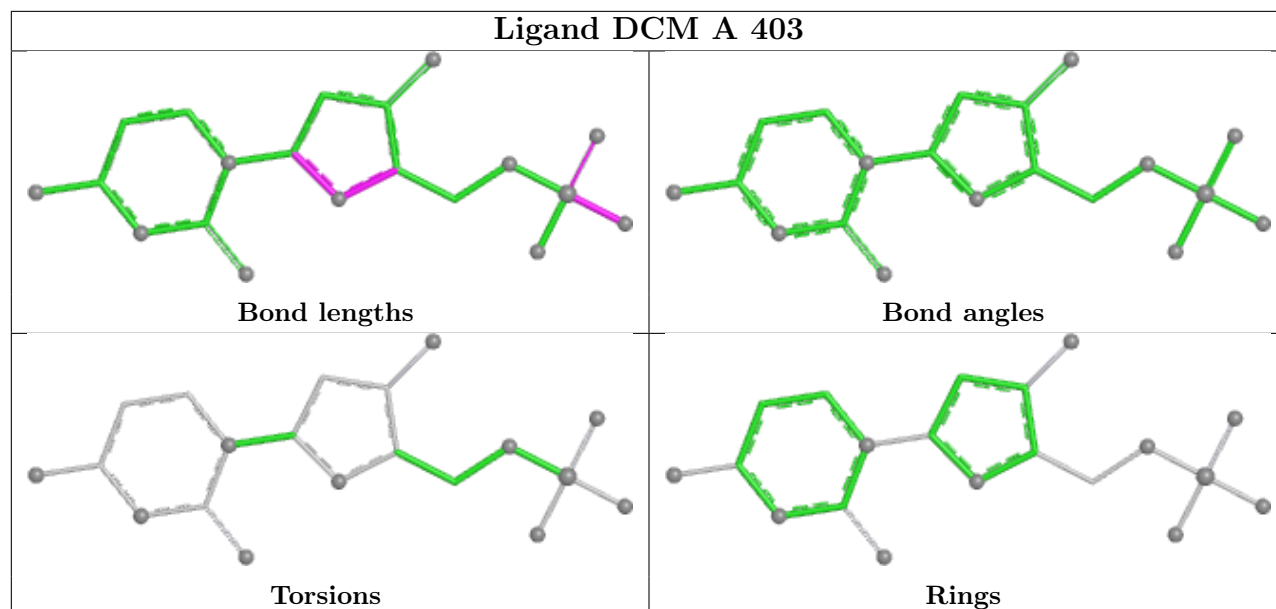
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	403	DCM	1	0
3	C	402	UDP	5	0
4	C	403	DCM	1	0
4	A	403	DCM	1	0
3	A	402	UDP	5	0
3	D	402	UDP	4	0
4	D	403	DCM	1	0
3	B	402	UDP	6	0

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

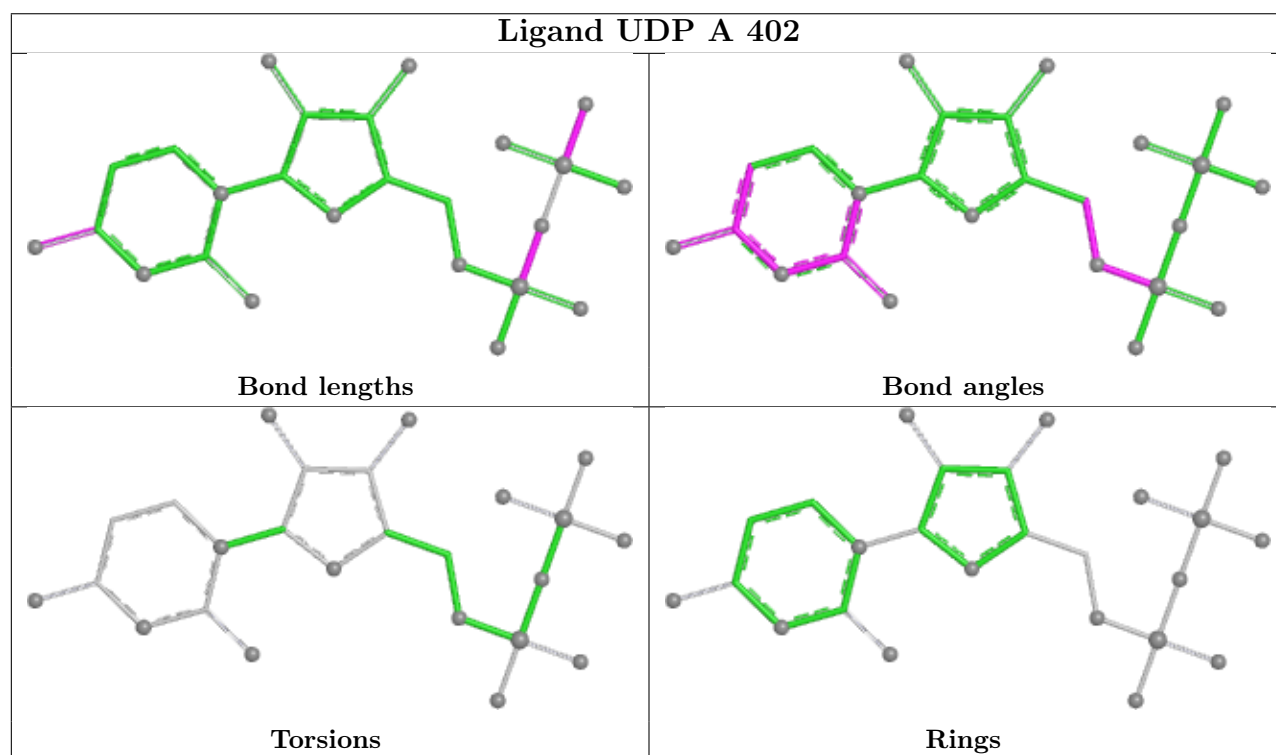


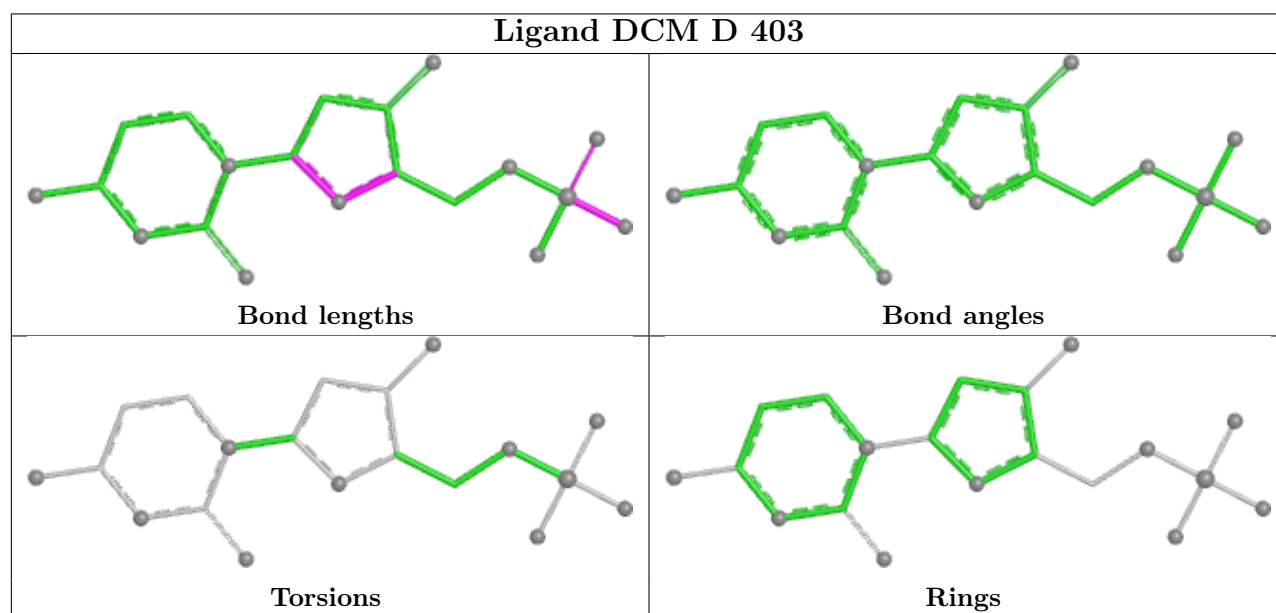
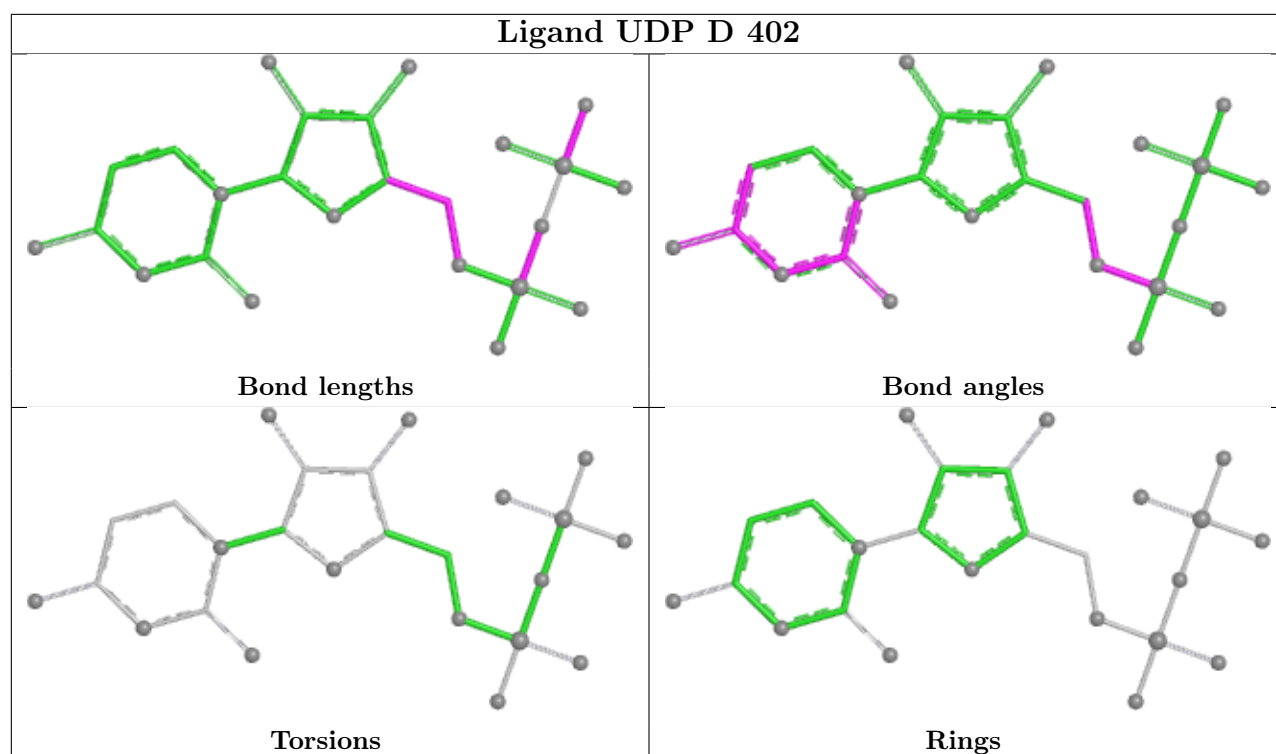


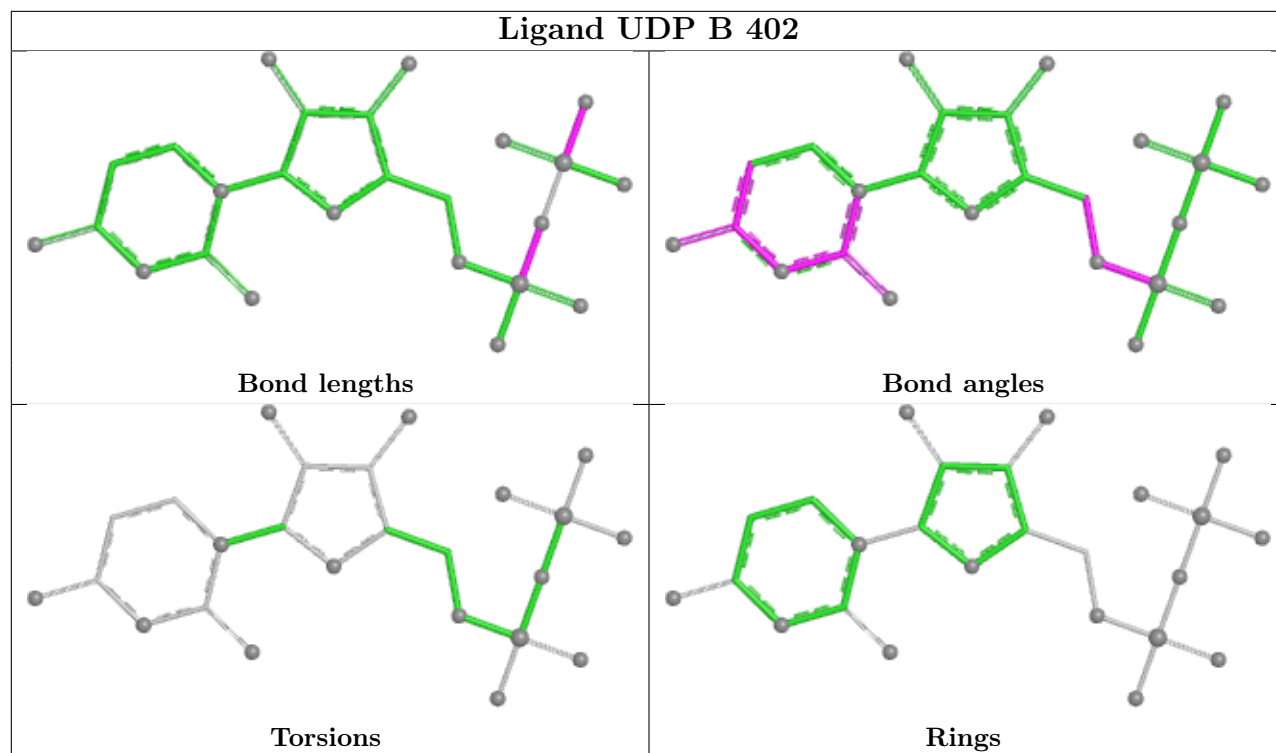
Ligand DCM A 403



Ligand UDP A 402







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	228/280 (81%)	0.06	2 (0%) 81 68	23, 63, 132, 173	0
1	B	228/280 (81%)	0.11	5 (2%) 62 47	26, 63, 135, 190	0
1	C	228/280 (81%)	0.01	6 (2%) 57 42	24, 59, 134, 161	0
1	D	228/280 (81%)	0.19	8 (3%) 47 34	26, 62, 128, 193	0
All	All	912/1120 (81%)	0.09	21 (2%) 61 46	23, 62, 134, 193	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	166	PHE	3.4
1	D	245	LYS	3.1
1	B	166	PHE	3.0
1	A	245	LYS	3.0
1	D	167	GLY	3.0
1	D	242	PHE	2.9
1	B	245	LYS	2.9
1	D	251	GLU	2.8
1	B	167	GLY	2.8
1	C	245	LYS	2.7
1	C	111	SER	2.7
1	C	166	PHE	2.6
1	B	46	GLU	2.6
1	D	63	SER	2.5
1	C	39	ASN	2.5
1	D	243	LYS	2.4
1	C	121	LYS	2.3
1	C	167	GLY	2.3
1	A	190	TYR	2.2
1	B	116	LEU	2.1
1	D	57	ARG	2.1

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 ⓘ

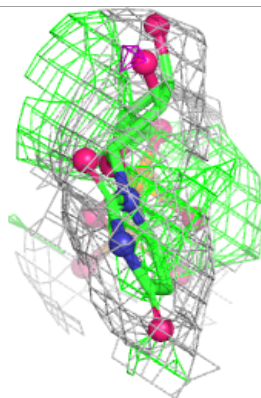
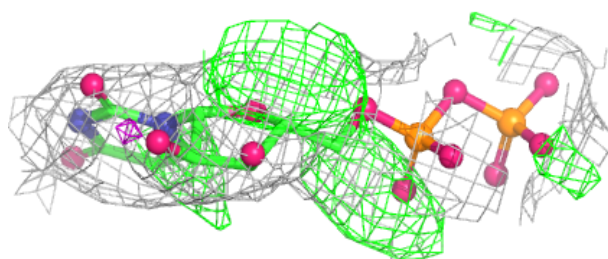
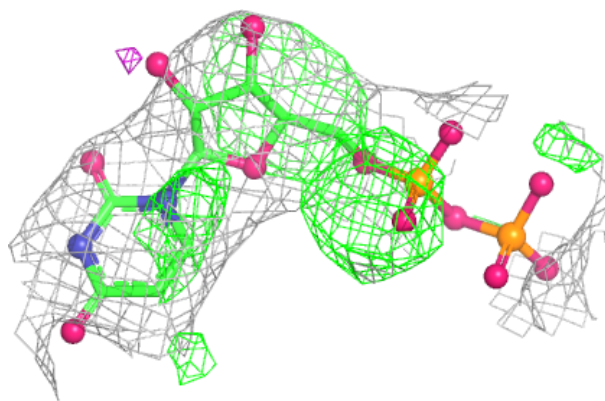
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	UDP	D	402	25/25	0.82	0.17	66,82,87,88	0
3	UDP	C	402	25/25	0.85	0.16	66,82,87,88	0
3	UDP	B	402	25/25	0.87	0.14	66,82,87,88	0
3	UDP	A	402	25/25	0.89	0.15	66,82,87,88	0
2	MG	B	401	1/1	0.91	0.23	56,56,56,56	0
2	MG	C	401	1/1	0.91	0.19	56,56,56,56	0
2	MG	D	401	1/1	0.92	0.18	56,56,56,56	0
4	DCM	C	403	20/20	0.92	0.18	51,80,89,89	0
4	DCM	D	403	20/20	0.92	0.17	51,80,89,89	0
4	DCM	A	403	20/20	0.93	0.17	51,80,89,89	0
4	DCM	B	403	20/20	0.93	0.16	51,80,89,89	0
2	MG	A	401	1/1	0.97	0.14	56,56,56,56	0

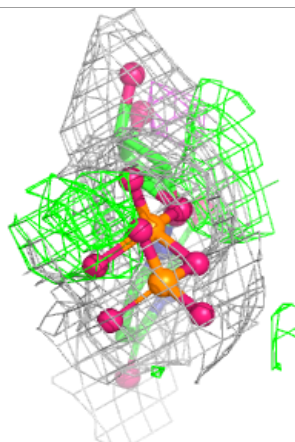
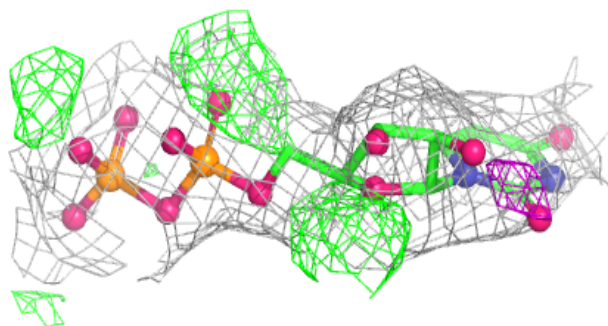
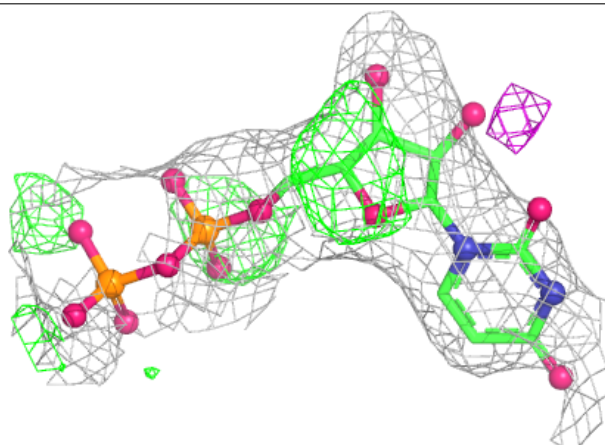
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around UDP D 402:

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

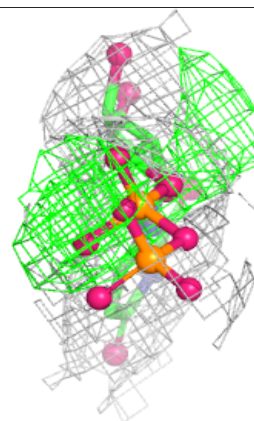
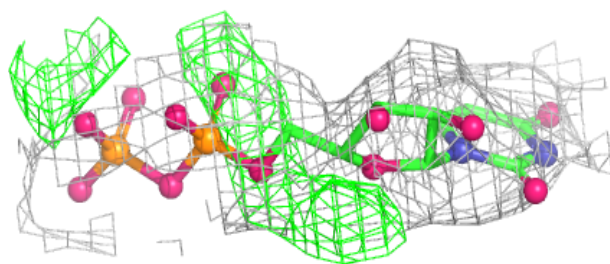
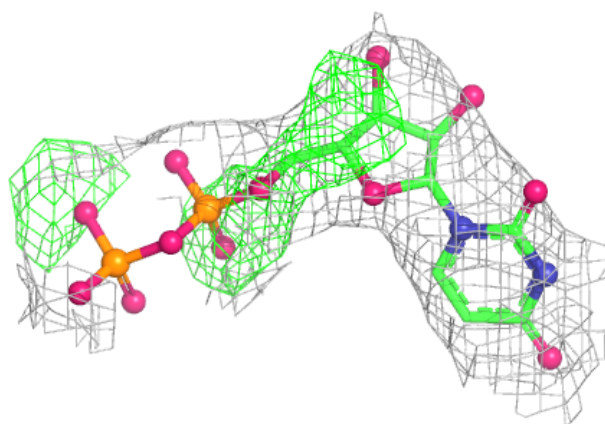
**Electron density around UDP C 402:**

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



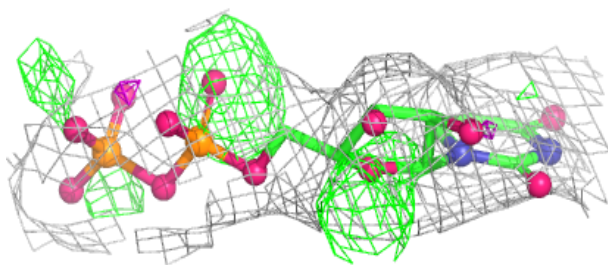
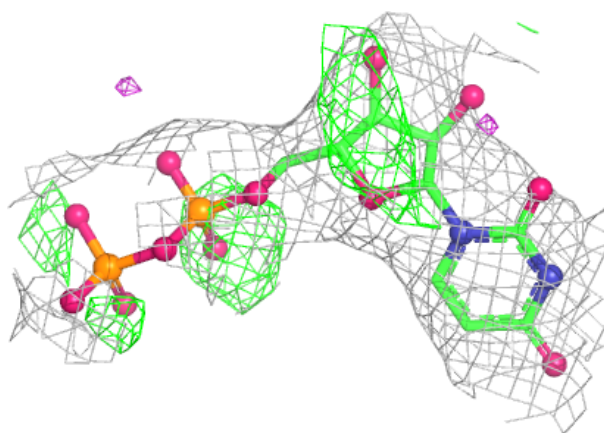
Electron density around UDP B 402:

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



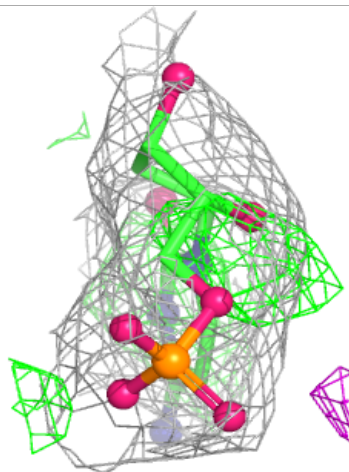
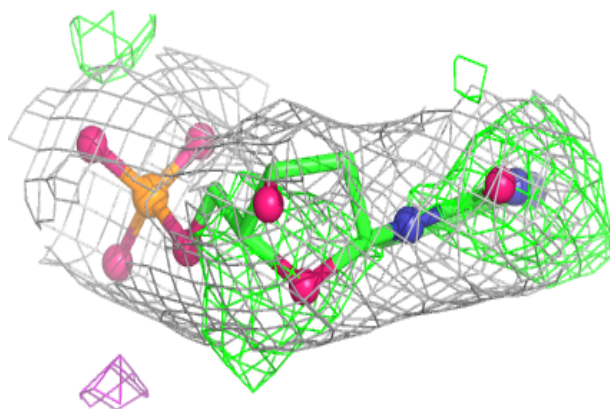
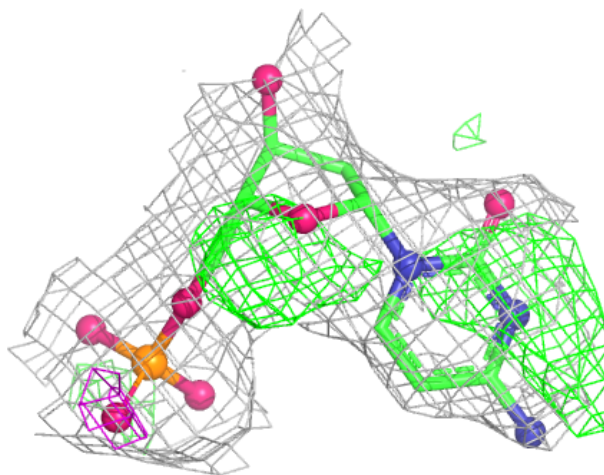
Electron density around UDP A 402:

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



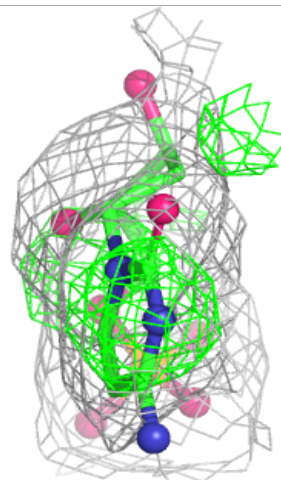
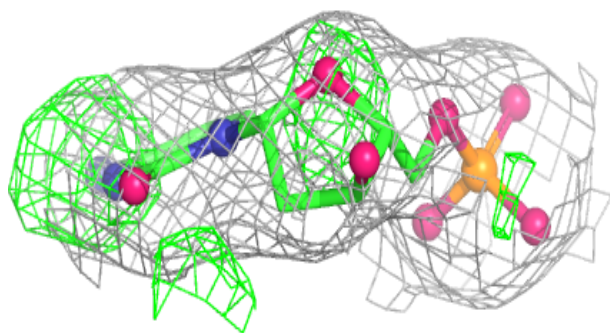
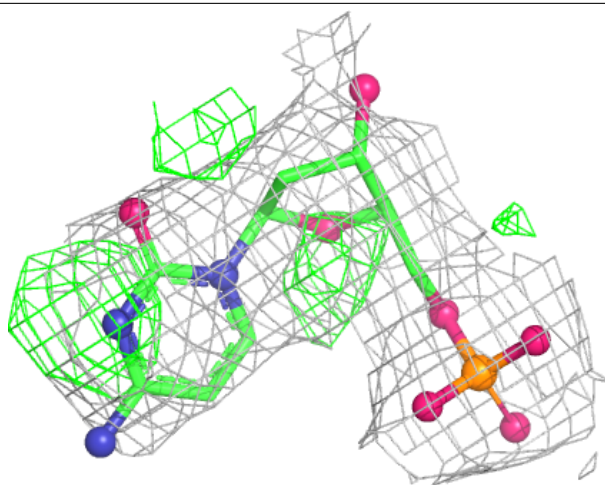
Electron density around DCM C 403:

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



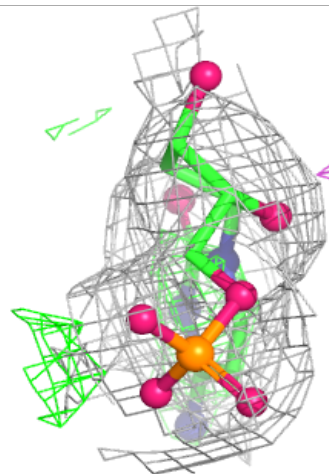
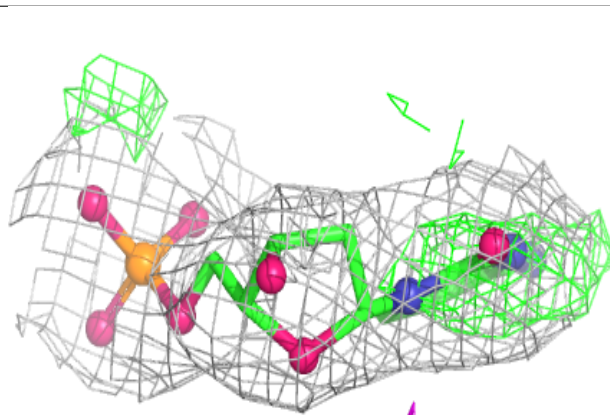
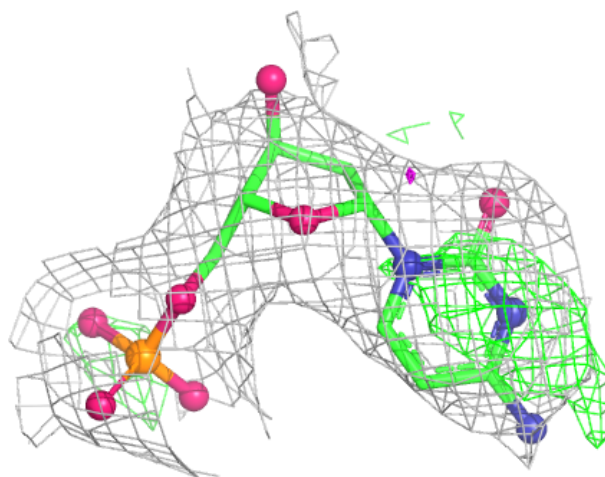
Electron density around DCM D 403:

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



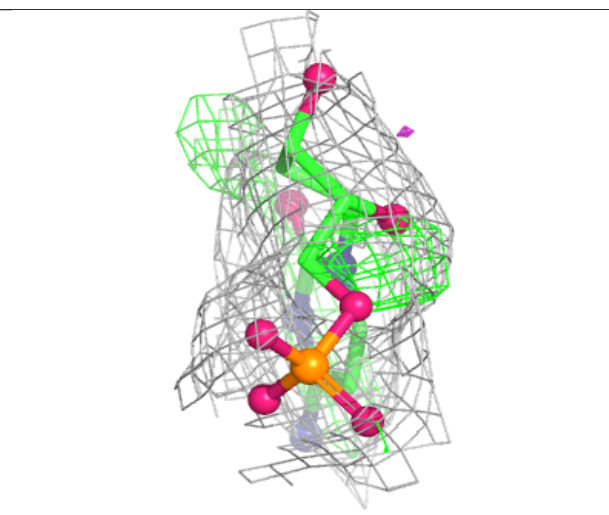
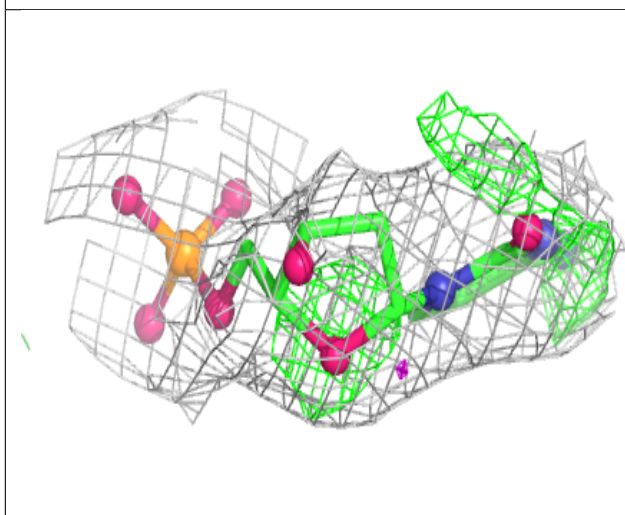
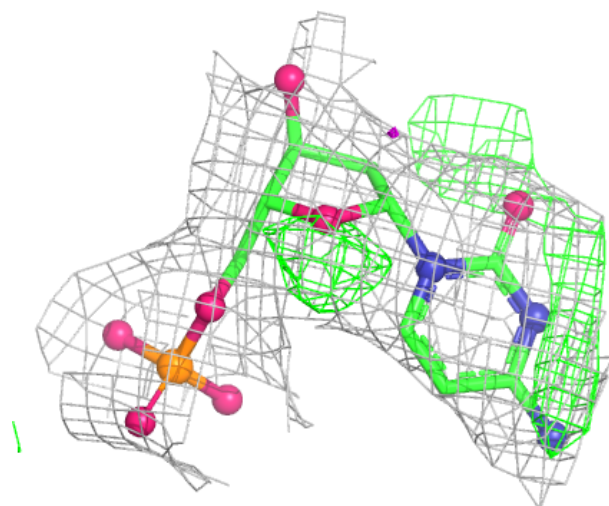
Electron density around DCM A 403:

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



Electron density around DCM B 403:

$2mF_o-DF_c$ (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.