



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 07:28 PM UTC

PDB ID : 5MQL / pdb_00005mql
Title : Crystal structure of dCK mutant C3S in complex with masitinib and UDP
Authors : Rebuffet, E.; Hammam, K.; Saez-Ayala, M.; Gros, L.; Lopez, S.; Hajem, B.;
Humbert, M.; Baudalet, E.; Audebert, S.; Betzi, S.; Lugari, A.; Combes, S.;
Pez, D.; Letard, S.; Mansfield, C.; Moussy, A.; de Sepulveda, P.; Morelli, X.;
Dubreuil, P.
Deposited on : 2016-12-20
Resolution : 3.25 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](#) ①) 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

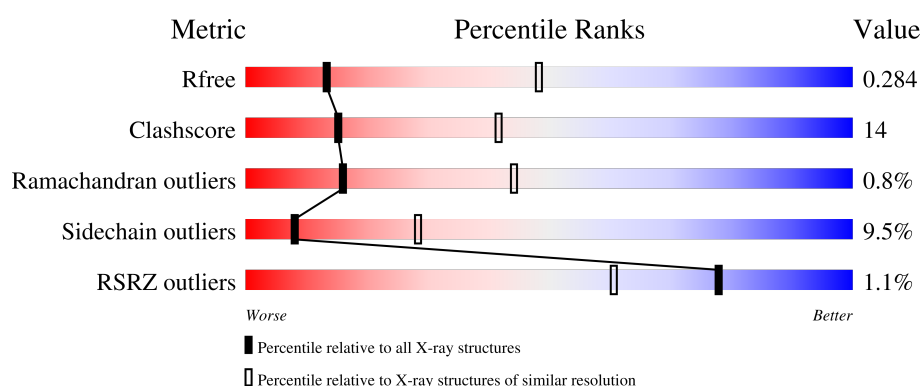
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.25 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	1605 (3.30-3.22)
Clashscore	190562	1660 (3.30-3.22)
Ramachandran outliers	187476	1630 (3.30-3.22)
Sidechain outliers	187428	1629 (3.30-3.22)
RSRZ outliers	180081	1605 (3.30-3.22)

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	289	% 57% 19% • 20%
1	B	289	% 55% 20% • 21%
1	C	289	53% 22% • 21%
1	D	289	% 54% 19% 5% 22%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 7527 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	231	Total	C	N	O	S	0	0	0
			1858	1192	305	353	8			
1	B	227	Total	C	N	O	S	0	0	0
			1806	1167	299	333	7			
1	C	228	Total	C	N	O	S	0	0	0
			1853	1190	308	348	7			
1	D	225	Total	C	N	O	S	0	0	0
			1815	1169	294	345	7			

There are 128 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-28	MET	-	initiating methionine	UNP P27707
A	-27	SER	-	expression tag	UNP P27707
A	-26	TYR	-	expression tag	UNP P27707
A	-25	TYR	-	expression tag	UNP P27707
A	-24	HIS	-	expression tag	UNP P27707
A	-23	HIS	-	expression tag	UNP P27707
A	-22	HIS	-	expression tag	UNP P27707
A	-21	HIS	-	expression tag	UNP P27707
A	-20	HIS	-	expression tag	UNP P27707
A	-19	HIS	-	expression tag	UNP P27707
A	-18	LEU	-	expression tag	UNP P27707
A	-17	GLU	-	expression tag	UNP P27707
A	-16	SER	-	expression tag	UNP P27707
A	-15	THR	-	expression tag	UNP P27707
A	-14	SER	-	expression tag	UNP P27707
A	-13	LEU	-	expression tag	UNP P27707
A	-12	TYR	-	expression tag	UNP P27707
A	-11	LYS	-	expression tag	UNP P27707
A	-10	LYS	-	expression tag	UNP P27707
A	-9	ALA	-	expression tag	UNP P27707
A	-8	GLY	-	expression tag	UNP P27707

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	LEU	-	expression tag	UNP P27707
A	-6	GLU	-	expression tag	UNP P27707
A	-5	ASN	-	expression tag	UNP P27707
A	-4	LEU	-	expression tag	UNP P27707
A	-3	TYR	-	expression tag	UNP P27707
A	-2	PHE	-	expression tag	UNP P27707
A	-1	GLN	-	expression tag	UNP P27707
A	0	GLY	-	expression tag	UNP P27707
A	9	SER	CYS	engineered mutation	UNP P27707
A	45	SER	CYS	engineered mutation	UNP P27707
A	59	SER	CYS	engineered mutation	UNP P27707
B	-28	MET	-	initiating methionine	UNP P27707
B	-27	SER	-	expression tag	UNP P27707
B	-26	TYR	-	expression tag	UNP P27707
B	-25	TYR	-	expression tag	UNP P27707
B	-24	HIS	-	expression tag	UNP P27707
B	-23	HIS	-	expression tag	UNP P27707
B	-22	HIS	-	expression tag	UNP P27707
B	-21	HIS	-	expression tag	UNP P27707
B	-20	HIS	-	expression tag	UNP P27707
B	-19	HIS	-	expression tag	UNP P27707
B	-18	LEU	-	expression tag	UNP P27707
B	-17	GLU	-	expression tag	UNP P27707
B	-16	SER	-	expression tag	UNP P27707
B	-15	THR	-	expression tag	UNP P27707
B	-14	SER	-	expression tag	UNP P27707
B	-13	LEU	-	expression tag	UNP P27707
B	-12	TYR	-	expression tag	UNP P27707
B	-11	LYS	-	expression tag	UNP P27707
B	-10	LYS	-	expression tag	UNP P27707
B	-9	ALA	-	expression tag	UNP P27707
B	-8	GLY	-	expression tag	UNP P27707
B	-7	LEU	-	expression tag	UNP P27707
B	-6	GLU	-	expression tag	UNP P27707
B	-5	ASN	-	expression tag	UNP P27707
B	-4	LEU	-	expression tag	UNP P27707
B	-3	TYR	-	expression tag	UNP P27707
B	-2	PHE	-	expression tag	UNP P27707
B	-1	GLN	-	expression tag	UNP P27707
B	0	GLY	-	expression tag	UNP P27707
B	9	SER	CYS	engineered mutation	UNP P27707
B	45	SER	CYS	engineered mutation	UNP P27707

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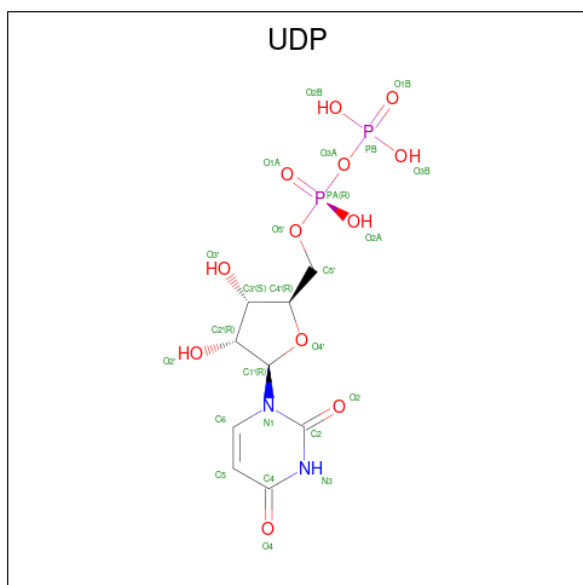
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B	59	SER	CYS	engineered mutation	UNP P27707
C	-28	MET	-	initiating methionine	UNP P27707
C	-27	SER	-	expression tag	UNP P27707
C	-26	TYR	-	expression tag	UNP P27707
C	-25	TYR	-	expression tag	UNP P27707
C	-24	HIS	-	expression tag	UNP P27707
C	-23	HIS	-	expression tag	UNP P27707
C	-22	HIS	-	expression tag	UNP P27707
C	-21	HIS	-	expression tag	UNP P27707
C	-20	HIS	-	expression tag	UNP P27707
C	-19	HIS	-	expression tag	UNP P27707
C	-18	LEU	-	expression tag	UNP P27707
C	-17	GLU	-	expression tag	UNP P27707
C	-16	SER	-	expression tag	UNP P27707
C	-15	THR	-	expression tag	UNP P27707
C	-14	SER	-	expression tag	UNP P27707
C	-13	LEU	-	expression tag	UNP P27707
C	-12	TYR	-	expression tag	UNP P27707
C	-11	LYS	-	expression tag	UNP P27707
C	-10	LYS	-	expression tag	UNP P27707
C	-9	ALA	-	expression tag	UNP P27707
C	-8	GLY	-	expression tag	UNP P27707
C	-7	LEU	-	expression tag	UNP P27707
C	-6	GLU	-	expression tag	UNP P27707
C	-5	ASN	-	expression tag	UNP P27707
C	-4	LEU	-	expression tag	UNP P27707
C	-3	TYR	-	expression tag	UNP P27707
C	-2	PHE	-	expression tag	UNP P27707
C	-1	GLN	-	expression tag	UNP P27707
C	0	GLY	-	expression tag	UNP P27707
C	9	SER	CYS	engineered mutation	UNP P27707
C	45	SER	CYS	engineered mutation	UNP P27707
C	59	SER	CYS	engineered mutation	UNP P27707
D	-28	MET	-	initiating methionine	UNP P27707
D	-27	SER	-	expression tag	UNP P27707
D	-26	TYR	-	expression tag	UNP P27707
D	-25	TYR	-	expression tag	UNP P27707
D	-24	HIS	-	expression tag	UNP P27707
D	-23	HIS	-	expression tag	UNP P27707
D	-22	HIS	-	expression tag	UNP P27707
D	-21	HIS	-	expression tag	UNP P27707
D	-20	HIS	-	expression tag	UNP P27707

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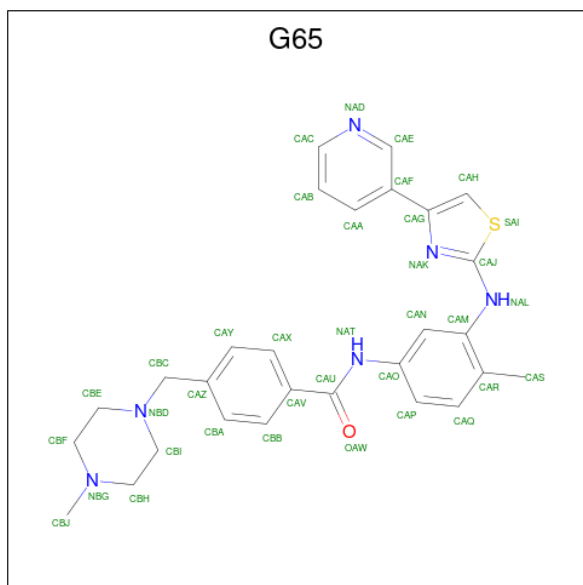
Chain	Residue	Modelled	Actual	Comment	Reference
D	-19	HIS	-	expression tag	UNP P27707
D	-18	LEU	-	expression tag	UNP P27707
D	-17	GLU	-	expression tag	UNP P27707
D	-16	SER	-	expression tag	UNP P27707
D	-15	THR	-	expression tag	UNP P27707
D	-14	SER	-	expression tag	UNP P27707
D	-13	LEU	-	expression tag	UNP P27707
D	-12	TYR	-	expression tag	UNP P27707
D	-11	LYS	-	expression tag	UNP P27707
D	-10	LYS	-	expression tag	UNP P27707
D	-9	ALA	-	expression tag	UNP P27707
D	-8	GLY	-	expression tag	UNP P27707
D	-7	LEU	-	expression tag	UNP P27707
D	-6	GLU	-	expression tag	UNP P27707
D	-5	ASN	-	expression tag	UNP P27707
D	-4	LEU	-	expression tag	UNP P27707
D	-3	TYR	-	expression tag	UNP P27707
D	-2	PHE	-	expression tag	UNP P27707
D	-1	GLN	-	expression tag	UNP P27707
D	0	GLY	-	expression tag	UNP P27707
D	9	SER	CYS	engineered mutation	UNP P27707
D	45	SER	CYS	engineered mutation	UNP P27707
D	59	SER	CYS	engineered mutation	UNP P27707

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



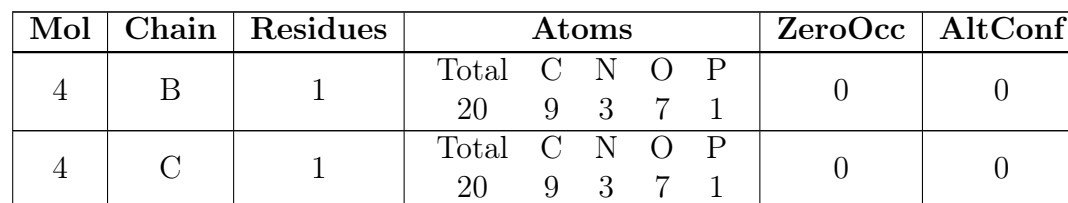
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			25	9	2	12	2		
2	B	1	Total	C	N	O	P	0	0
			25	9	2	12	2		
2	C	1	Total	C	N	O	P	0	0
			25	9	2	12	2		
2	D	1	Total	C	N	O	P	0	0
			25	9	2	12	2		

- Molecule 3 is Masitinib (CCD ID: G65) (formula: $C_{28}H_{30}N_6OS$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			36	28	6	1	1		

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



- | Mol | Chain | Residues | Atoms | ZeroOcc | AltConf |
|-----|-------|----------|-----------------|---------|---------|
| 5 | C | 1 | Total Mg
1 1 | 0 | 0 |

- | Mol | Chain | Residues | Atoms | ZeroOcc | AltConf |
|-----|-------|----------|----------------|---------|---------|
| 6 | A | 5 | Total O
5 5 | 0 | 0 |
| 6 | B | 3 | Total O
3 3 | 0 | 0 |
| 6 | C | 5 | Total O
5 5 | 0 | 0 |
| 6 | D | 5 | Total O
5 5 | 0 | 0 |

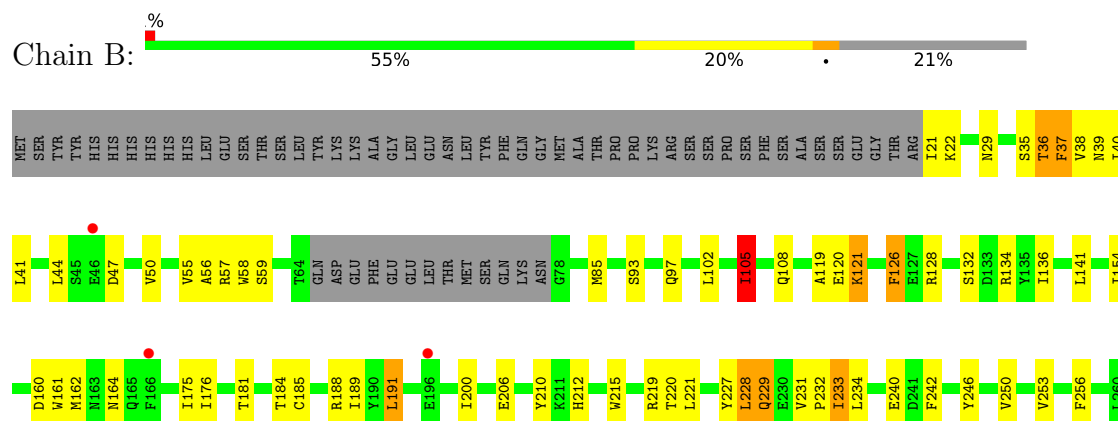
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

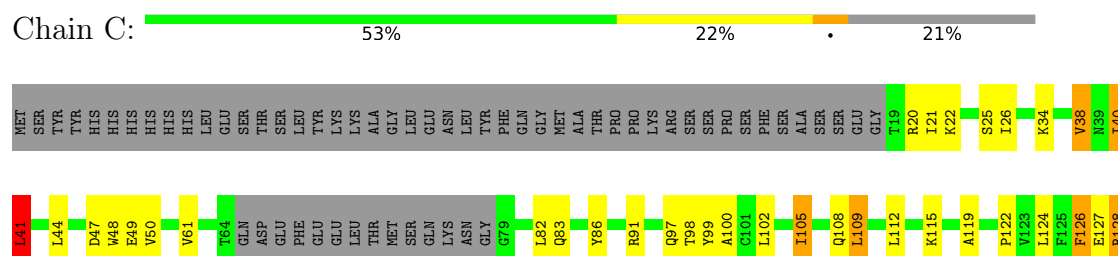
• Molecule 1: Deoxycytidine kinase

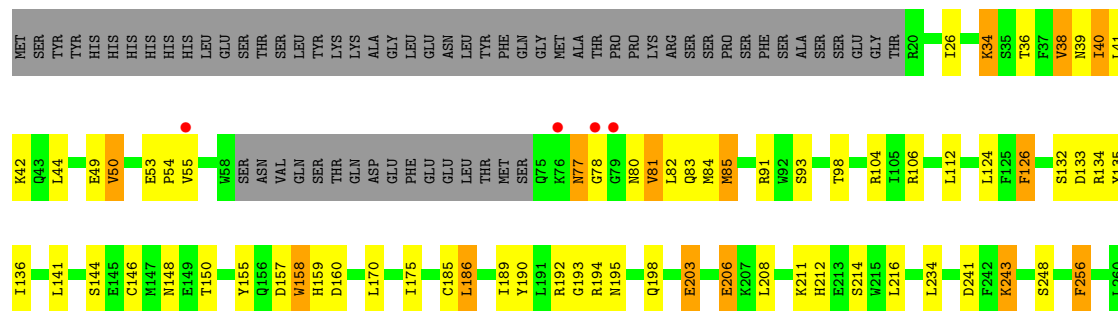


• Molecule 1: Deoxycytidine kinase



• 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	88.33Å 88.33Å 342.53Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.80 – 3.25 47.80 – 3.25	Depositor EDS
% Data completeness (in resolution range)	99.9 (47.80-3.25) 99.9 (47.80-3.25)	Depositor EDS
R_{merge}	0.27	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.55 (at 3.25Å)	Xtriage
Refinement program	BUSTER 2.10.3	Depositor
R, R_{free}	0.215 , 0.277 0.225 , 0.284	Depositor DCC
R_{free} test set	1120 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	100.5	Xtriage
Anisotropy	0.443	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 100.0	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	0.93	EDS
Total number of atoms	7527	wwPDB-VP
Average B, all atoms (Å ²)	97.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.12% 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: DCM, G65, UDP, MG

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	0.93	0/1904	1.43	16/2590 (0.6%)
1	B	0.90	1/1852 (0.1%)	1.39	10/2521 (0.4%)
1	C	0.96	0/1899	1.41	14/2581 (0.5%)
1	D	0.96	0/1861	1.40	13/2532 (0.5%)
All	All	0.94	1/7516 (0.0%)	1.41	53/10224 (0.5%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	121	LYS	CA-C	6.69	1.58	1.53

All (53) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	126	PHE	CA-CB-CG	10.05	123.85	113.80
1	A	216	LEU	N-CA-C	8.38	120.49	111.36
1	C	196	GLU	CA-C-N	8.36	133.02	120.31
1	C	196	GLU	C-N-CA	8.36	133.02	120.31
1	A	135	TYR	N-CA-C	7.87	119.94	111.36
1	A	42	LYS	CA-C-N	7.40	130.94	120.28
1	A	42	LYS	C-N-CA	7.40	130.94	120.28
1	C	126	PHE	CA-CB-CG	7.15	120.95	113.80
1	B	57	ARG	N-CA-C	-6.93	103.82	112.90
1	A	166	PHE	CA-CB-CG	-6.92	106.88	113.80
1	D	170	LEU	N-CA-C	-6.75	103.12	112.30
1	C	41	LEU	N-CA-C	6.68	118.64	111.36
1	B	37	PHE	CA-CB-CG	6.68	120.48	113.80
1	A	125	PHE	CA-CB-CG	6.66	120.46	113.80
1	D	133	ASP	CA-CB-CG	6.65	119.25	112.60
1	A	51	VAL	CB-CA-C	6.54	117.29	111.08

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	42	LYS	N-CA-C	-6.37	104.25	111.07
1	C	83	GLN	CA-C-N	6.17	128.84	120.38
1	C	83	GLN	C-N-CA	6.17	128.84	120.38
1	A	126	PHE	CA-CB-CG	5.99	119.78	113.80
1	D	158	TRP	CA-C-N	5.91	128.20	120.28
1	D	158	TRP	C-N-CA	5.91	128.20	120.28
1	B	126	PHE	CA-CB-CG	5.89	119.69	113.80
1	C	40	ILE	CB-CA-C	-5.89	104.43	111.97
1	D	157	ASP	CA-C-N	5.86	128.39	120.65
1	D	157	ASP	C-N-CA	5.86	128.39	120.65
1	C	191	LEU	CA-C-N	5.86	128.13	120.28
1	C	191	LEU	C-N-CA	5.86	128.13	120.28
1	C	100	ALA	N-CA-C	-5.73	104.67	111.03
1	A	20	ARG	CA-C-N	5.66	129.66	121.18
1	A	20	ARG	C-N-CA	5.66	129.66	121.18
1	B	58	TRP	N-CA-C	-5.60	104.80	111.69
1	C	166	PHE	CA-CB-CG	5.55	119.35	113.80
1	A	225	PHE	CA-C-N	5.53	128.01	120.54
1	A	225	PHE	C-N-CA	5.53	128.01	120.54
1	B	36	THR	CA-C-N	5.37	127.91	120.29
1	B	36	THR	C-N-CA	5.37	127.91	120.29
1	A	168	GLN	CA-C-N	5.29	127.69	120.54
1	A	168	GLN	C-N-CA	5.29	127.69	120.54
1	B	191	LEU	CA-C-N	5.29	127.37	120.28
1	B	191	LEU	C-N-CA	5.29	127.37	120.28
1	D	256	PHE	CA-CB-CG	5.25	119.06	113.80
1	C	130	VAL	CA-C-N	5.25	128.29	120.31
1	C	130	VAL	C-N-CA	5.25	128.29	120.31
1	B	105	ILE	CA-C-N	5.22	127.54	120.65
1	B	105	ILE	C-N-CA	5.22	127.54	120.65
1	C	127	GLU	N-CA-C	-5.09	100.87	108.96
1	D	195	ASN	CA-C-N	5.04	127.53	120.28
1	D	195	ASN	C-N-CA	5.04	127.53	120.28
1	A	19	THR	CA-C-N	5.01	127.00	120.28
1	A	19	THR	C-N-CA	5.01	127.00	120.28
1	D	54	PRO	CA-C-N	5.01	129.22	120.86
1	D	54	PRO	C-N-CA	5.01	129.22	120.86

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	1858	0	1734	42	0
1	B	1806	0	1690	62	0
1	C	1853	0	1750	58	0
1	D	1815	0	1688	43	0
2	A	25	0	11	0	0
2	B	25	0	11	0	0
2	C	25	0	11	1	0
2	D	25	0	11	1	0
3	A	36	0	0	3	0
4	B	20	0	12	1	0
4	C	20	0	12	1	0
5	C	1	0	0	0	0
6	A	5	0	0	0	0
6	B	3	0	0	0	0
6	C	5	0	0	0	0
6	D	5	0	0	0	0
All	All	7527	0	6930	196	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 14.

All (196) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:219:ARG:NH1	1:B:233:ILE:HD12	1.32	1.38
1:B:219:ARG:HH11	1:B:233:ILE:CD1	1.40	1.34
1:B:219:ARG:NH1	1:B:233:ILE:CD1	1.91	1.30
1:B:47:ASP:OD1	1:B:119:ALA:HB1	1.36	1.25
1:C:41:LEU:HD11	1:C:48:TRP:CZ3	1.80	1.15
1:D:85:MET:CE	1:D:93:SER:HA	1.85	1.06
1:D:85:MET:HE3	1:D:93:SER:HA	1.07	1.05
1:B:47:ASP:OD1	1:B:119:ALA:CB	2.05	1.04
1:B:35:SER:O	1:B:38:VAL:HG12	1.56	1.02
1:C:41:LEU:HD11	1:C:48:TRP:HZ3	1.12	0.98
1:A:49:GLU:HB2	1:A:116:LEU:HD11	1.46	0.97

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:21:ILE:HG22	1:B:21:ILE:O	1.70	0.90
1:C:41:LEU:CD1	1:C:48:TRP:HZ3	1.89	0.85
1:B:219:ARG:HH12	1:B:233:ILE:HD12	1.37	0.85
1:C:41:LEU:CD1	1:C:48:TRP:CZ3	2.60	0.85
1:B:219:ARG:NH1	1:B:233:ILE:HD13	1.92	0.84
1:C:38:VAL:HG11	1:C:50:VAL:HG11	1.62	0.81
1:C:38:VAL:CG1	1:C:50:VAL:HG11	2.10	0.80
1:D:85:MET:HE3	1:D:93:SER:CA	2.01	0.80
1:C:41:LEU:HD12	1:C:41:LEU:O	1.81	0.79
1:B:219:ARG:HH11	1:B:233:ILE:HD12	1.04	0.78
1:B:38:VAL:HG22	1:B:50:VAL:HG21	1.65	0.78
1:D:194:ARG:O	1:D:198:GLN:HG2	1.84	0.78
1:A:97:GLN:HE22	3:A:302:G65:CAE	1.98	0.76
1:B:35:SER:C	1:B:38:VAL:HG12	2.12	0.75
1:A:203:GLU:N	1:A:203:GLU:OE2	2.21	0.74
1:A:40:ILE:HD13	1:A:246:TYR:CG	2.24	0.72
1:B:105:ILE:HD13	1:B:162:MET:HE3	1.72	0.71
1:A:202:LEU:HD12	1:A:202:LEU:O	1.90	0.71
1:C:197:GLU:O	1:C:200:ILE:HG13	1.91	0.71
1:D:85:MET:CE	1:D:93:SER:CA	2.65	0.70
1:B:35:SER:HA	1:B:38:VAL:CG1	2.21	0.69
1:B:108:GLN:NE2	1:B:126:PHE:HB3	2.06	0.69
1:B:97:GLN:HE22	4:B:302:DCM:HN41	1.41	0.69
1:C:61:VAL:HG11	1:D:150:THR:HG23	1.74	0.69
1:A:185:CYS:O	1:A:189:ILE:HG13	1.93	0.69
1:B:35:SER:HA	1:B:38:VAL:HG12	1.75	0.68
1:C:175:ILE:HG21	1:C:216:LEU:HD22	1.76	0.67
1:D:134:ARG:NH1	1:D:160:ASP:OD1	2.28	0.67
1:A:232:PRO:HG2	1:A:260:LEU:CD2	2.24	0.66
1:B:219:ARG:HH11	1:B:233:ILE:HD11	1.52	0.66
1:A:173:ASP:O	1:A:260:LEU:HD22	1.95	0.66
1:C:44:LEU:HD23	1:C:44:LEU:N	2.10	0.66
1:B:38:VAL:HG22	1:B:50:VAL:CG2	2.26	0.65
1:A:232:PRO:HG2	1:A:260:LEU:HD21	1.77	0.65
1:C:41:LEU:HD21	1:C:253:VAL:HG11	1.79	0.65
1:D:93:SER:OG	1:D:141:LEU:HD13	1.97	0.64
1:B:21:ILE:O	1:B:21:ILE:CG2	2.44	0.63
1:A:40:ILE:HD13	1:A:246:TYR:CD2	2.34	0.62
1:D:80:ASN:O	1:D:83:GLN:N	2.34	0.61
1:B:35:SER:CA	1:B:38:VAL:HG12	2.32	0.59
1:D:193:GLY:O	1:D:194:ARG:C	2.45	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:38:VAL:HG11	1:C:50:VAL:CG1	2.34	0.58
1:C:238:VAL:HG23	1:C:238:VAL:O	2.01	0.58
1:C:136:ILE:HD11	1:C:215:TRP:HE3	1.68	0.57
1:C:161:TRP:CD1	1:D:106:ARG:HD3	2.39	0.57
1:B:108:GLN:NE2	1:B:126:PHE:CG	2.73	0.57
1:B:38:VAL:HG13	1:B:39:ASN:N	2.18	0.57
1:C:38:VAL:HG13	1:C:50:VAL:HG21	1.87	0.57
1:C:161:TRP:CD1	1:D:106:ARG:CD	2.88	0.57
1:A:116:LEU:O	1:A:117:LYS:C	2.47	0.56
1:B:121:LYS:HG3	1:B:121:LYS:O	2.06	0.56
1:B:38:VAL:CG1	1:B:39:ASN:N	2.69	0.56
1:C:246:TYR:CD1	1:C:246:TYR:N	2.73	0.56
1:C:41:LEU:HD12	1:C:41:LEU:C	2.31	0.55
1:B:227:TYR:CD1	1:B:227:TYR:C	2.85	0.55
1:D:26:ILE:HG13	1:D:38:VAL:HG23	1.88	0.54
1:A:109:LEU:HA	1:A:112:LEU:HD12	1.90	0.53
1:B:47:ASP:OD1	1:B:119:ALA:HB3	2.04	0.53
1:C:105:ILE:O	1:C:109:LEU:HB2	2.09	0.52
1:C:41:LEU:CD1	1:C:48:TRP:CE3	2.92	0.52
1:C:189:ILE:HD12	1:C:205:LEU:HD11	1.91	0.52
1:A:177:TYR:HB2	1:A:216:LEU:HD13	1.92	0.52
1:B:40:ILE:HD13	1:B:246:TYR:CD2	2.44	0.52
1:C:38:VAL:CG1	1:C:50:VAL:CG1	2.87	0.52
1:D:136:ILE:HG21	1:D:212:HIS:CE1	2.45	0.52
1:C:97:GLN:HE22	4:C:303:DCM:HN41	1.59	0.51
1:B:232:PRO:HB2	1:B:256:PHE:HE1	1.76	0.51
1:D:77:ASN:ND2	1:D:77:ASN:H	2.08	0.51
1:C:41:LEU:HD11	1:C:48:TRP:CE3	2.43	0.51
1:A:116:LEU:HD23	1:A:116:LEU:N	2.26	0.50
1:C:21:ILE:HG23	1:C:122:PRO:HB2	1.93	0.50
1:C:105:ILE:HD13	1:C:162:MET:HE3	1.93	0.50
1:C:119:ALA:HB3	1:C:122:PRO:HB3	1.94	0.50
1:A:81:VAL:HG12	1:A:96:PHE:HD1	1.76	0.49
1:A:136:ILE:HD13	1:A:212:HIS:CE1	2.47	0.49
1:C:242:PHE:CD1	2:C:302:UDP:O4	2.65	0.49
1:A:116:LEU:O	1:A:119:ALA:HB3	2.12	0.49
1:D:81:VAL:CG1	1:D:82:LEU:N	2.73	0.49
1:C:161:TRP:CD1	1:D:106:ARG:HD2	2.48	0.49
1:A:115:LYS:O	1:A:116:LEU:C	2.56	0.49
1:B:40:ILE:HD12	1:B:41:LEU:HD23	1.94	0.48
1:B:108:GLN:HE22	1:B:126:PHE:HB3	1.77	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:134:ARG:NH2	1:D:135:TYR:OH	2.45	0.48
1:D:77:ASN:ND2	1:D:77:ASN:N	2.60	0.48
1:A:44:LEU:N	1:A:44:LEU:HD23	2.28	0.48
1:B:128:ARG:HD3	1:B:132:SER:OG	2.14	0.48
1:D:36:THR:O	1:D:40:ILE:HG13	2.13	0.48
1:B:234:LEU:HD13	1:B:256:PHE:HB2	1.95	0.48
1:B:215:TRP:HB2	1:B:221:LEU:HD22	1.96	0.47
1:B:250:VAL:HA	1:B:253:VAL:HG12	1.95	0.47
1:D:132:SER:HA	1:D:136:ILE:HD13	1.95	0.47
1:B:47:ASP:CG	1:B:120:GLU:O	2.58	0.47
1:A:81:VAL:HG12	1:A:96:PHE:CD1	2.50	0.47
1:C:49:GLU:HB3	1:C:124:LEU:HD12	1.97	0.47
1:A:116:LEU:O	1:A:119:ALA:N	2.48	0.47
1:D:80:ASN:O	1:D:81:VAL:C	2.55	0.47
1:D:234:LEU:HD13	1:D:256:PHE:HB2	1.97	0.47
1:A:81:VAL:HG11	1:A:96:PHE:HA	1.96	0.47
1:B:21:ILE:O	1:B:22:LYS:C	2.57	0.47
1:B:56:ALA:HA	1:B:59:SER:HB2	1.96	0.47
1:C:173:ASP:O	1:C:260:LEU:HD22	2.15	0.47
1:A:242:PHE:HB2	1:A:249:LEU:HD21	1.96	0.46
1:D:26:ILE:HG22	1:D:34:LYS:HG2	1.98	0.46
1:C:200:ILE:HG23	1:C:204:TYR:HD2	1.80	0.46
1:D:134:ARG:HD3	1:D:159:HIS:ND1	2.31	0.46
1:B:188:ARG:HA	1:B:191:LEU:HD12	1.98	0.45
1:C:38:VAL:HG13	1:C:50:VAL:CG2	2.47	0.45
1:A:106:ARG:HG3	1:B:161:TRP:NE1	2.32	0.45
1:C:181:THR:HG23	1:C:239:ASN:HD21	1.81	0.45
1:C:136:ILE:HG21	1:C:212:HIS:CE1	2.52	0.45
1:A:35:SER:HA	1:A:38:VAL:HG22	1.98	0.45
1:A:176:ILE:HG12	1:A:234:LEU:HD23	1.99	0.45
1:C:91:ARG:CD	1:D:91:ARG:HD3	2.47	0.45
1:C:132:SER:O	1:C:136:ILE:HB	2.17	0.45
1:C:134:ARG:HB2	1:C:159:HIS:CD2	2.52	0.44
1:D:203:GLU:HA	1:D:206:GLU:HG2	1.99	0.44
1:D:185:CYS:O	1:D:189:ILE:HG12	2.17	0.44
1:A:161:TRP:CZ2	1:A:165:GLN:HG3	2.52	0.44
1:B:108:GLN:NE2	1:B:126:PHE:CB	2.78	0.44
1:A:202:LEU:HD12	1:A:202:LEU:C	2.38	0.44
1:C:136:ILE:HD11	1:C:215:TRP:CE3	2.51	0.44
1:C:161:TRP:NE1	1:D:106:ARG:HD3	2.31	0.44
1:C:161:TRP:O	1:C:165:GLN:HG2	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:200:ILE:CG2	1:C:204:TYR:HD2	2.31	0.44
1:A:196:GLU:H	1:A:196:GLU:CD	2.26	0.44
1:B:93:SER:HB3	1:B:141:LEU:HD13	1.99	0.44
1:C:34:LYS:HE3	1:C:128:ARG:CZ	2.48	0.44
1:C:200:ILE:HG23	1:C:204:TYR:CD2	2.53	0.43
1:B:134:ARG:NH1	1:B:160:ASP:OD1	2.44	0.43
1:D:38:VAL:HG13	1:D:50:VAL:HG11	2.00	0.43
1:A:135:TYR:CE2	1:A:223:THR:HB	2.53	0.43
1:B:175:ILE:O	1:B:233:ILE:HA	2.17	0.43
1:B:227:TYR:O	1:B:228:LEU:C	2.60	0.43
1:A:97:GLN:NE2	3:A:302:G65:CAE	2.73	0.43
1:B:29:ASN:HB3	1:B:185:CYS:SG	2.59	0.43
1:C:47:ASP:O	1:C:122:PRO:HA	2.18	0.43
1:B:227:TYR:HE1	1:B:231:VAL:HG21	1.84	0.43
1:B:227:TYR:O	1:B:229:GLN:N	2.52	0.43
1:C:234:LEU:HD13	1:C:256:PHE:HB2	2.01	0.43
1:D:190:TYR:HA	1:D:198:GLN:HE22	1.83	0.43
1:B:35:SER:O	1:B:38:VAL:CG1	2.47	0.43
1:C:82:LEU:HD11	1:C:86:TYR:HE2	1.84	0.43
1:C:61:VAL:HG23	1:C:99:TYR:CE2	2.54	0.42
1:D:175:ILE:HG21	1:D:216:LEU:HD22	2.01	0.42
1:A:40:ILE:CD1	1:A:246:TYR:CD2	3.01	0.42
1:B:35:SER:HA	1:B:38:VAL:HG11	1.96	0.42
1:B:36:THR:O	1:B:40:ILE:HG23	2.19	0.42
1:B:206:GLU:HG2	1:B:210:TYR:CE2	2.54	0.42
1:C:124:LEU:HD23	1:C:126:PHE:HE1	1.84	0.42
1:A:30:ILE:O	1:A:185:CYS:HB3	2.20	0.42
1:B:176:ILE:HG12	1:B:234:LEU:HD23	2.01	0.42
1:D:77:ASN:HD22	1:D:78:GLY:H	1.67	0.42
1:D:81:VAL:HA	1:D:84:MET:HB2	2.00	0.42
1:D:81:VAL:O	1:D:82:LEU:C	2.60	0.42
1:A:75:GLN:O	1:A:76:LYS:C	2.63	0.42
1:A:81:VAL:CG1	1:A:96:PHE:HB2	2.49	0.42
1:D:192:ARG:HD3	2:D:500:UDP:H5'1	2.01	0.42
1:A:106:ARG:HG3	1:B:161:TRP:CD1	2.55	0.42
1:A:81:VAL:CG1	1:A:96:PHE:HD1	2.33	0.42
1:D:124:LEU:HD23	1:D:126:PHE:HE1	1.84	0.42
1:A:195:ASN:HA	1:A:198:GLN:HG2	2.02	0.42
1:A:232:PRO:HB2	1:A:256:PHE:HE1	1.84	0.42
1:C:25:SER:HB2	1:C:172:LEU:HD22	2.02	0.42
1:D:40:ILE:O	1:D:44:LEU:HG	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:215:TRP:O	1:A:219:ARG:HA	2.19	0.41
1:B:227:TYR:CD1	1:B:228:LEU:N	2.89	0.41
1:A:81:VAL:HG23	1:B:154:ILE:HD11	2.02	0.41
1:B:108:GLN:HE21	1:B:126:PHE:HB3	1.82	0.41
1:C:242:PHE:CD1	1:C:242:PHE:N	2.86	0.41
1:B:242:PHE:O	1:B:246:TYR:HB3	2.21	0.41
1:C:231:VAL:HA	1:C:232:PRO:HD3	1.88	0.41
1:B:38:VAL:HG21	1:B:50:VAL:CG1	2.51	0.41
1:B:136:ILE:HD13	1:B:212:HIS:CE1	2.56	0.41
1:C:177:TYR:HB2	1:C:216:LEU:HD13	2.03	0.41
1:C:98:THR:HG22	1:C:155:TYR:HD1	1.86	0.41
1:D:53:GLU:HG3	1:D:104:ARG:HH22	1.86	0.41
1:D:155:TYR:O	1:D:158:TRP:HB3	2.21	0.41
1:A:55:VAL:HG21	3:A:302:G65:CAA	2.50	0.40
1:B:37:PHE:O	1:B:40:ILE:HG13	2.21	0.40
1:B:219:ARG:CZ	1:B:233:ILE:HD13	2.49	0.40
1:D:85:MET:CE	1:D:93:SER:CB	2.99	0.40
1:D:186:LEU:HD22	1:D:190:TYR:CZ	2.56	0.40
1:C:20:ARG:O	1:C:20:ARG:HG2	2.21	0.40
1:C:22:LYS:HA	1:C:173:ASP:OD2	2.21	0.40
1:D:85:MET:HE1	1:D:93:SER:OG	2.22	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	227/289 (78%)	202 (89%)	24 (11%)	1 (0%)	30	59
1	B	223/289 (77%)	200 (90%)	21 (9%)	2 (1%)	14	42
1	C	224/289 (78%)	209 (93%)	13 (6%)	2 (1%)	14	42
1	D	221/289 (76%)	188 (85%)	31 (14%)	2 (1%)	14	42

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	895/1156 (77%)	799 (89%)	89 (10%)	7 (1%)	16	44

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	128	ARG
1	C	115	LYS
1	A	120	GLU
1	B	228	LEU
1	D	241	ASP
1	D	243	LYS
1	B	55	VAL

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	195/265 (74%)	178 (91%)	17 (9%)	9	31
1	B	185/265 (70%)	172 (93%)	13 (7%)	14	40
1	C	196/265 (74%)	177 (90%)	19 (10%)	8	27
1	D	189/265 (71%)	165 (87%)	24 (13%)	4	18
All	All	765/1060 (72%)	692 (90%)	73 (10%)	8	28

All (73) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	THR
1	A	36	THR
1	A	40	ILE
1	A	44	LEU
1	A	50	VAL
1	A	51	VAL
1	A	55	VAL
1	A	73	MET

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Mol	Chain	Res	Type
1	A	103	SER
1	A	111	SER
1	A	130	VAL
1	A	148	ASN
1	A	171	GLU
1	A	189	ILE
1	A	207	LYS
1	A	249	LEU
1	A	251	GLU
1	B	44	LEU
1	B	85	MET
1	B	102	LEU
1	B	105	ILE
1	B	164	ASN
1	B	181	THR
1	B	184	THR
1	B	189	ILE
1	B	200	ILE
1	B	220	THR
1	B	229	GLN
1	B	233	ILE
1	B	240	GLU
1	C	26	ILE
1	C	38	VAL
1	C	40	ILE
1	C	41	LEU
1	C	102	LEU
1	C	105	ILE
1	C	108	GLN
1	C	109	LEU
1	C	112	LEU
1	C	183	GLU
1	C	197	GLU
1	C	200	ILE
1	C	203	GLU
1	C	214	SER
1	C	224	ASN
1	C	231	VAL
1	C	238	VAL
1	C	257	LEU
1	C	260	LEU
1	D	34	LYS

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Mol	Chain	Res	Type
1	D	38	VAL
1	D	39	ASN
1	D	40	ILE
1	D	41	LEU
1	D	49	GLU
1	D	50	VAL
1	D	55	VAL
1	D	77	ASN
1	D	81	VAL
1	D	85	MET
1	D	98	THR
1	D	112	LEU
1	D	144	SER
1	D	146	CYS
1	D	148	ASN
1	D	186	LEU
1	D	203	GLU
1	D	206	GLU
1	D	208	LEU
1	D	211	LYS
1	D	214	SER
1	D	243	LYS
1	D	248	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	39	ASN
1	A	97	GLN
1	A	113	ASN
1	A	187	HIS
1	A	212	HIS
1	B	80	ASN
1	B	97	GLN
1	B	108	GLN
1	B	140	ASN
1	B	163	ASN
1	B	212	HIS
1	C	97	GLN
1	C	108	GLN
1	C	156	GLN
1	D	77	ASN

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Mol	Chain	Res	Type
1	D	97	GLN
1	D	156	GLN
1	D	218	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](#)

Of 8 ligands modelled in this entry, 1 is monoatomic - leaving 7 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	$\# Z > 2$
2	UDP	C	302	-	25,26,26	0.39	0	38,40,40	0.44	0
2	UDP	D	500	-	25,26,26	0.44	0	38,40,40	0.73	1 (2%)
4	DCM	C	303	5	21,21,21	0.63	0	30,31,31	0.43	0
2	UDP	B	301	-	25,26,26	0.42	0	38,40,40	0.59	0
2	UDP	A	301	-	25,26,26	0.54	0	38,40,40	0.59	0
3	G65	A	302	-	40,40,40	2.43	10 (25%)	55,55,55	2.16	13 (23%)
4	DCM	B	302	-	21,21,21	0.50	0	30,31,31	0.51	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	C	302	-	-	6/16/32/32	0/2/2/2
2	UDP	D	500	-	-	0/16/32/32	0/2/2/2
4	DCM	C	303	5	-	4/10/22/22	0/2/2/2
2	UDP	B	301	-	-	0/16/32/32	0/2/2/2
2	UDP	A	301	-	-	0/16/32/32	0/2/2/2
3	G65	A	302	-	-	9/20/30/30	0/5/5/5
4	DCM	B	302	-	-	4/10/22/22	0/2/2/2

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	302	G65	CAF-CAG	-7.77	1.36	1.47
3	A	302	G65	CAS-CAR	-6.35	1.39	1.51
3	A	302	G65	CBC-CAZ	-6.32	1.40	1.51
3	A	302	G65	CAV-CAU	-4.97	1.39	1.50
3	A	302	G65	CAO-NAT	-3.65	1.34	1.41
3	A	302	G65	CAE-NAD	3.53	1.41	1.34
3	A	302	G65	CAJ-SAI	-3.06	1.69	1.74
3	A	302	G65	CAM-NAL	-3.05	1.35	1.41
3	A	302	G65	CAC-NAD	2.42	1.40	1.33
3	A	302	G65	CAJ-NAK	2.32	1.34	1.31

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	302	G65	CAH-SAI-CAJ	7.40	94.15	88.40
3	A	302	G65	SAI-CAJ-NAK	-6.87	108.47	115.69
3	A	302	G65	CAJ-NAK-CAG	6.20	114.97	109.96
3	A	302	G65	CBH-NBG-CBF	3.88	115.71	109.54
3	A	302	G65	CAS-CAR-CAM	3.60	125.31	121.23
2	D	500	UDP	O2B-PB-O3A	2.90	114.35	104.64
3	A	302	G65	SAI-CAJ-NAL	2.69	129.22	121.37
3	A	302	G65	CAF-CAG-CAH	-2.54	123.53	126.54
3	A	302	G65	CBJ-NBG-CBH	-2.51	105.86	110.63
3	A	302	G65	CAG-CAH-SAI	-2.47	108.17	111.16
3	A	302	G65	CAR-CAM-NAL	2.40	122.66	118.82
3	A	302	G65	CAS-CAR-CAQ	-2.33	115.79	120.28
3	A	302	G65	CAO-NAT-CAU	-2.26	120.64	126.61
3	A	302	G65	CAX-CAV-CBB	2.13	121.27	118.57

There are no chirality outliers.

All (23) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	302	UDP	C5'-O5'-PA-O1A
2	C	302	UDP	C5'-O5'-PA-O3A
2	C	302	UDP	PA-O3A-PB-O3B
4	B	302	DCM	C5'-O5'-P-O2P
4	B	302	DCM	C5'-O5'-P-O3P
4	C	303	DCM	C5'-O5'-P-O1P
4	C	303	DCM	C5'-O5'-P-O3P
3	A	302	G65	CAZ-CBC-NBD-CBE
3	A	302	G65	CAZ-CBC-NBD-CBI
3	A	302	G65	NAT-CAU-CAV-CAX
3	A	302	G65	OAW-CAU-CAV-CAX
3	A	302	G65	CAN-CAO-NAT-CAU
3	A	302	G65	NAT-CAU-CAV-CBB
3	A	302	G65	CAP-CAO-NAT-CAU
3	A	302	G65	OAW-CAU-CAV-CBB
4	B	302	DCM	C5'-O5'-P-O1P
2	C	302	UDP	PA-O3A-PB-O2B
2	C	302	UDP	C5'-O5'-PA-O2A
3	A	302	G65	CAN-CAM-NAL-CAJ
4	C	303	DCM	O4'-C4'-C5'-O5'
4	B	302	DCM	C3'-C4'-C5'-O5'
4	C	303	DCM	C5'-O5'-P-O2P
2	C	302	UDP	C4'-C5'-O5'-PA

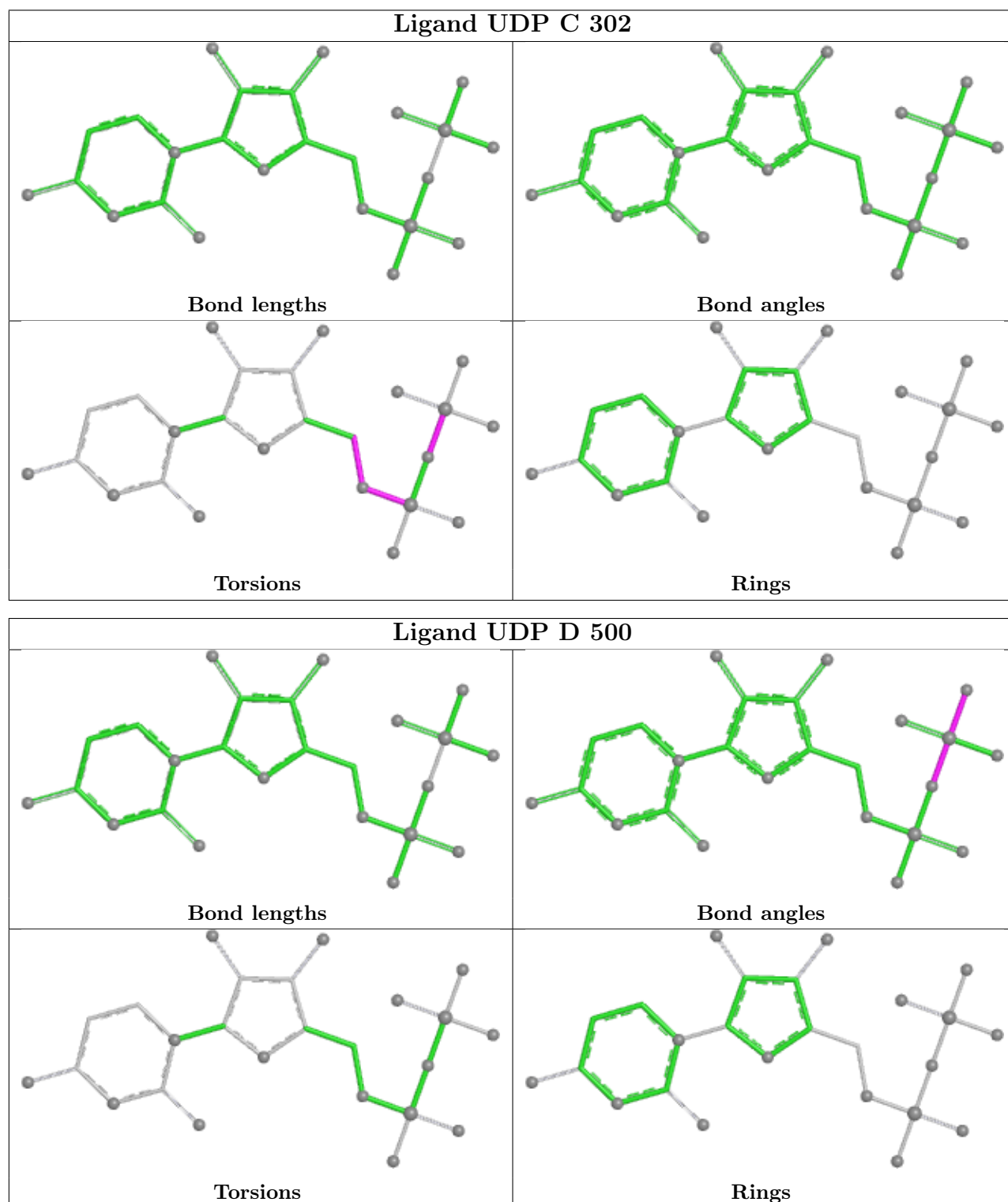
There are no ring outliers.

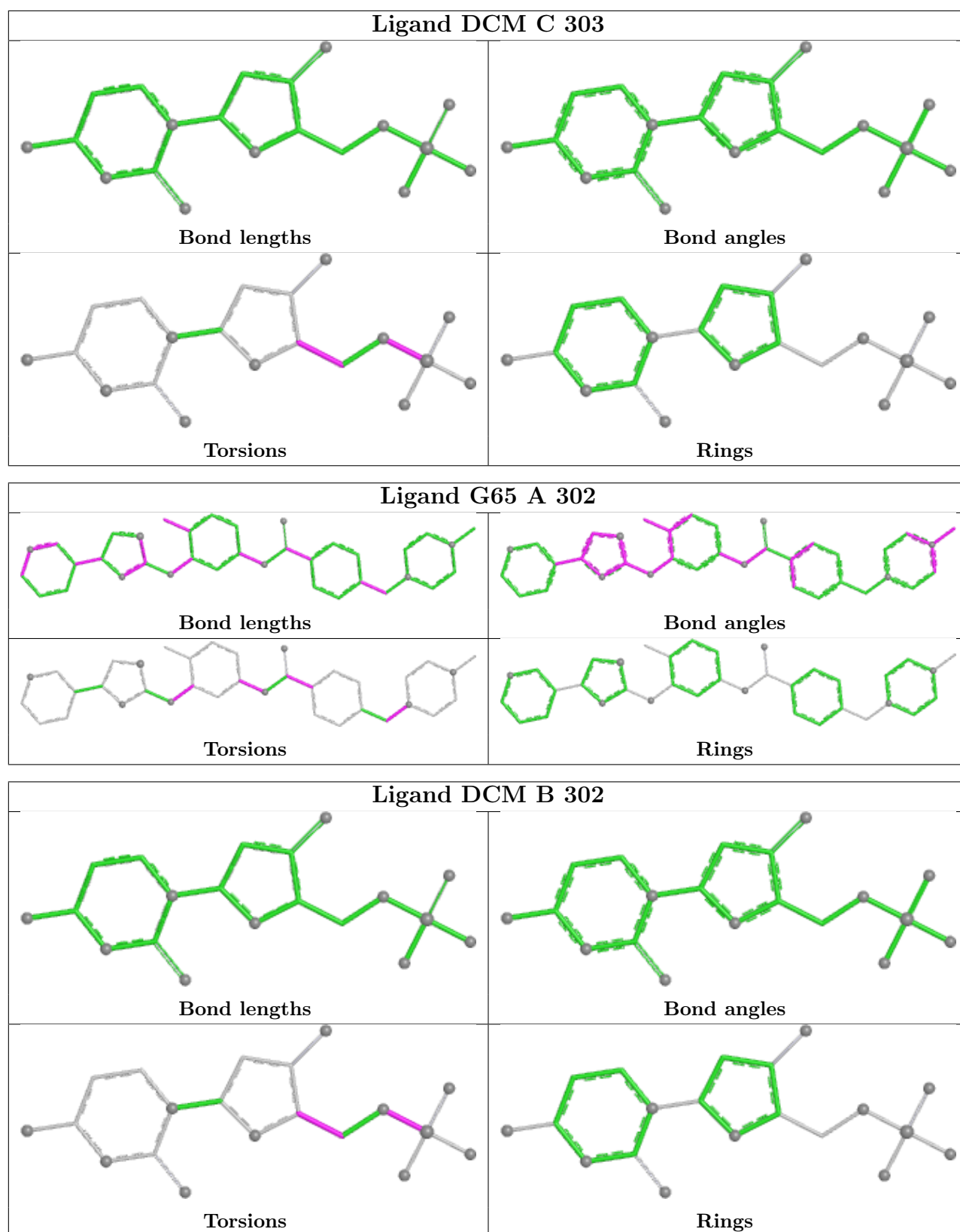
5 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	302	UDP	1	0
2	D	500	UDP	1	0
4	C	303	DCM	1	0
3	A	302	G65	3	0
4	B	302	DCM	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 ⓘ

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	231/289 (79%)	0.07	3 (1%) 75 56	75, 99, 128, 153	0
1	B	227/289 (78%)	0.07	3 (1%) 75 56	70, 96, 134, 145	0
1	C	228/289 (78%)	-0.05	0 100 100	67, 88, 125, 170	0
1	D	225/289 (77%)	0.06	4 (1%) 67 48	72, 95, 127, 141	0
All	All	911/1156 (78%)	0.04	10 (1%) 78 60	67, 95, 129, 170	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	78	GLY	5.1
1	D	55	VAL	5.0
1	B	166	PHE	2.5
1	B	196	GLU	2.4
1	D	76	LYS	2.3
1	D	79	GLY	2.3
1	A	73	MET	2.2
1	B	46	GLU	2.1
1	A	166	PHE	2.1
1	A	200	ILE	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 ⓘ

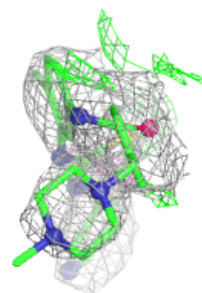
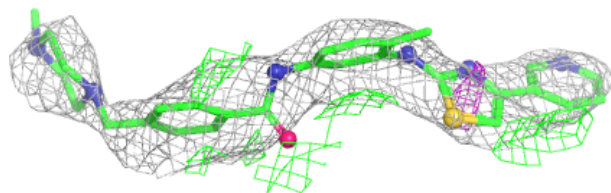
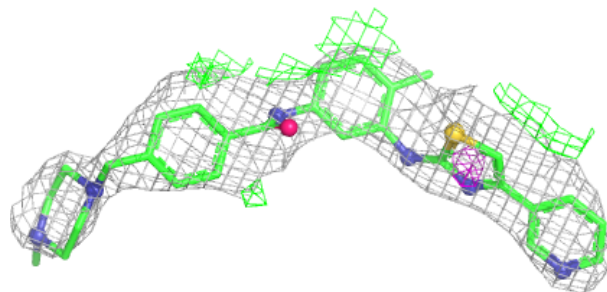
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	G65	A	302	36/36	0.83	0.20	122,133,142,143	0
5	MG	C	301	1/1	0.90	0.10	67,67,67,67	0
4	DCM	B	302	20/20	0.92	0.13	105,114,127,129	0
2	UDP	A	301	25/25	0.94	0.10	113,119,124,125	0
2	UDP	B	301	25/25	0.95	0.08	106,109,111,111	0
2	UDP	C	302	25/25	0.95	0.08	91,95,105,107	0
2	UDP	D	500	25/25	0.95	0.09	98,100,104,106	0
4	DCM	C	303	20/20	0.96	0.12	80,89,108,109	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.

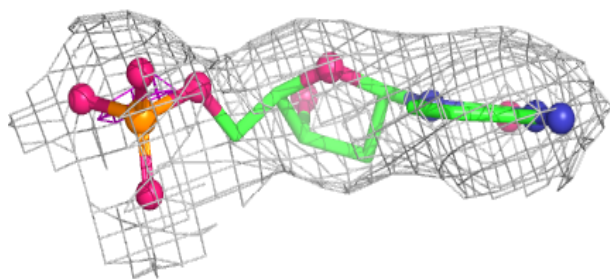
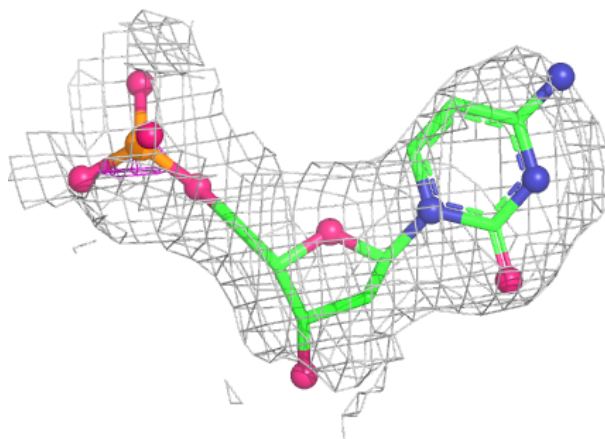
Electron density around G65 A 302:

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



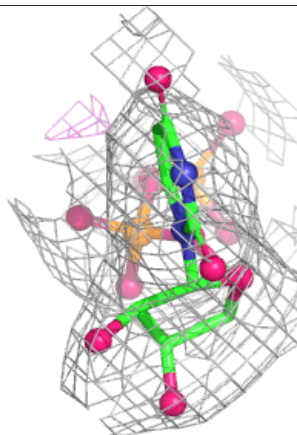
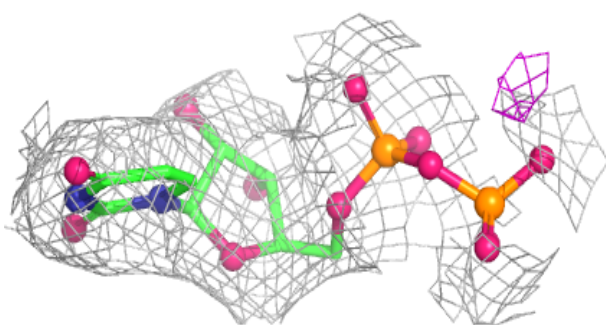
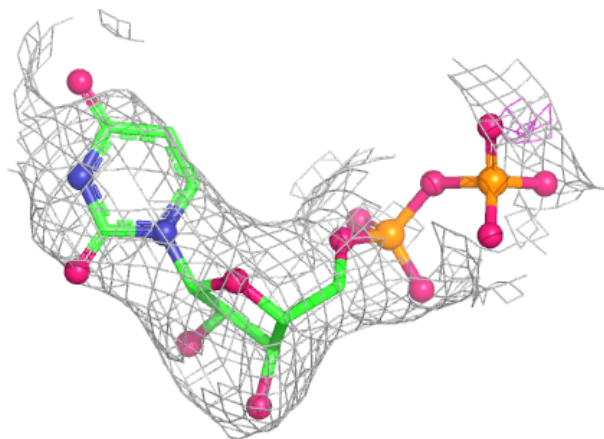
Electron density around DCM B 302:

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



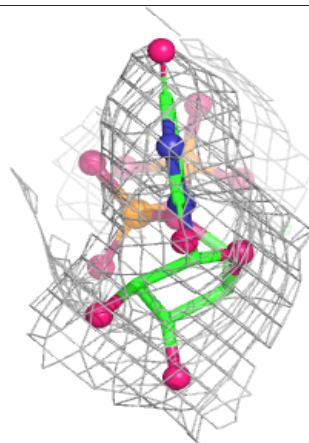
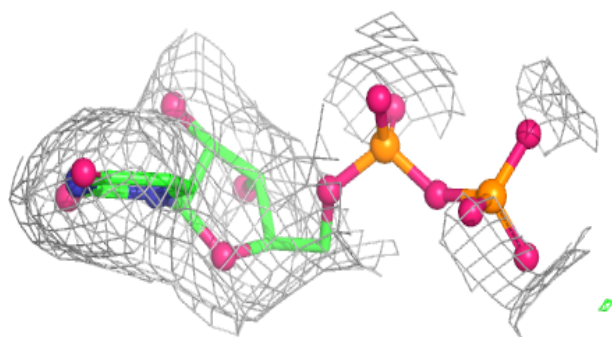
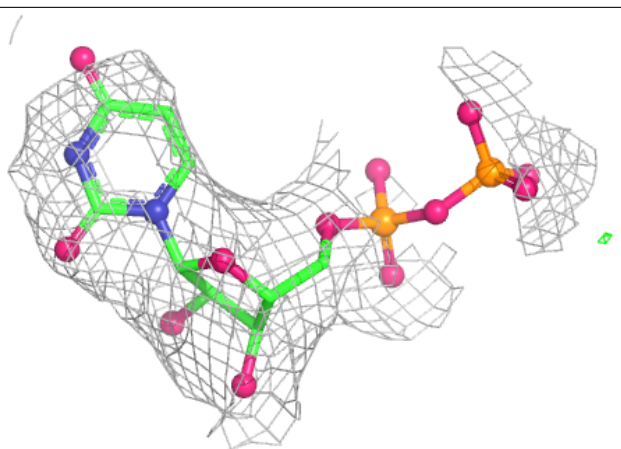
Electron density around UDP C 302:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

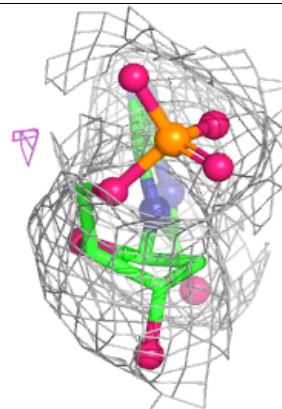
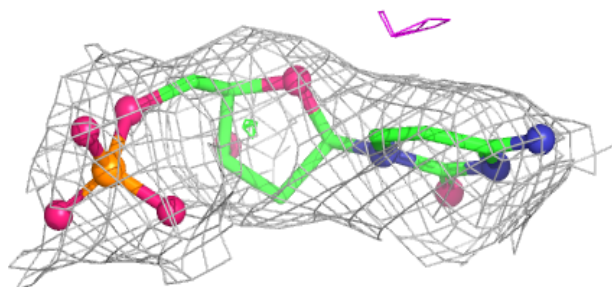
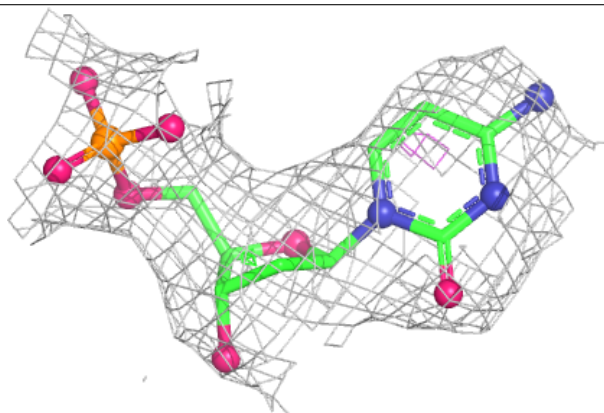


Electron density around UDP D 500:

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

**Electron density around DCM C 303:**

$2mF_o-DF_c$ (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.