



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 12:24 PM UTC

PDB ID : 3W3D / pdb_00003w3d
Title : Crystal structure of smooth muscle G actin DNase I complex
Authors : Sakabe, N.; Sakabe, K.; Sasaki, K.; Kondo, H.; Shimomur, M.
Deposited on : 2012-12-20
Resolution : 1.80 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

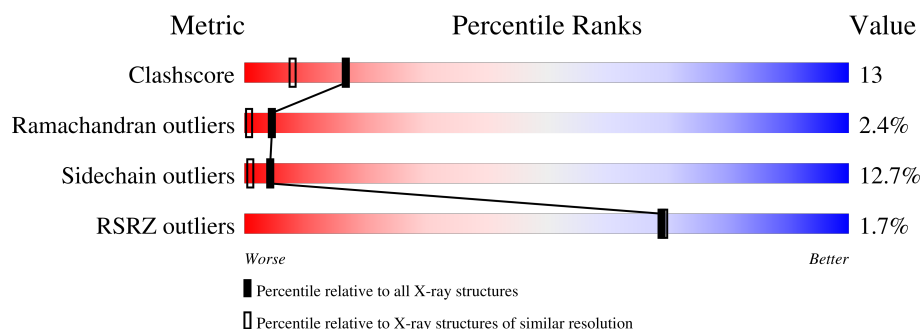
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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))
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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	374	3% 53% 33% 12% .
2	B	260	57% 32% 8% .
3	C	7	100%

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 5472 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 Actin, gamma-enteric smooth muscle.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	374	Total	C	N	O	S	0	0	0
			2924	1849	492	562	21			

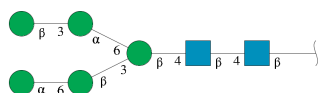
There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	359	GLN	PRO	conflict	UNP P63270

- Molecule 2 is a protein called Deoxyribonuclease-1.

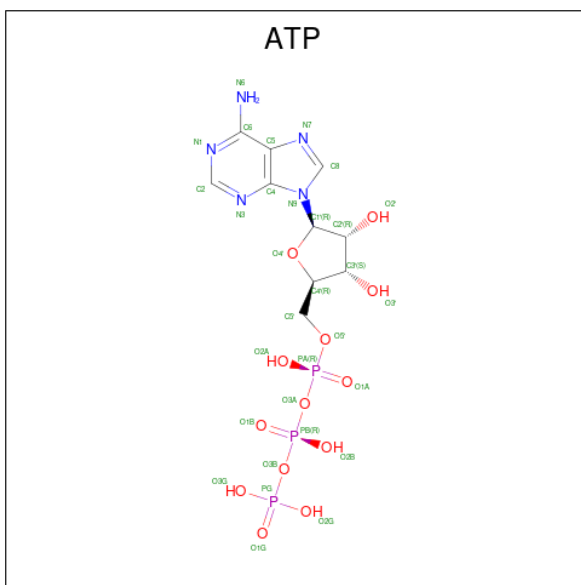
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	260	Total	C	N	O	S	0	0	0
			2049	1298	341	402	8			

- Molecule 3 is an oligosaccharide called alpha-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-3)-[beta-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	7	Total	C	N	O	0	0	0
			83	46	2	35			

- Molecule 4 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: C₁₀H₁₆N₅O₁₃P₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			31	10	5	13	3		

- Molecule 5 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total Ca 1 1	0	0
5	B	1	Total Ca 1 1	0	0

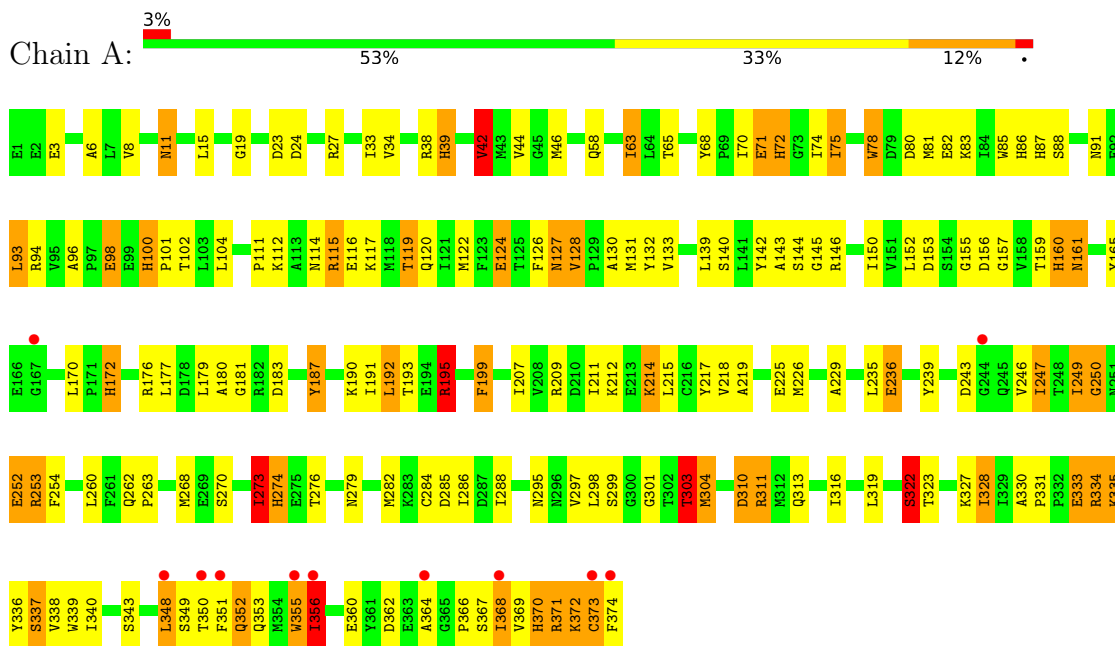
- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	195	Total O 195 195	0	0
6	B	188	Total O 188 188	0	0

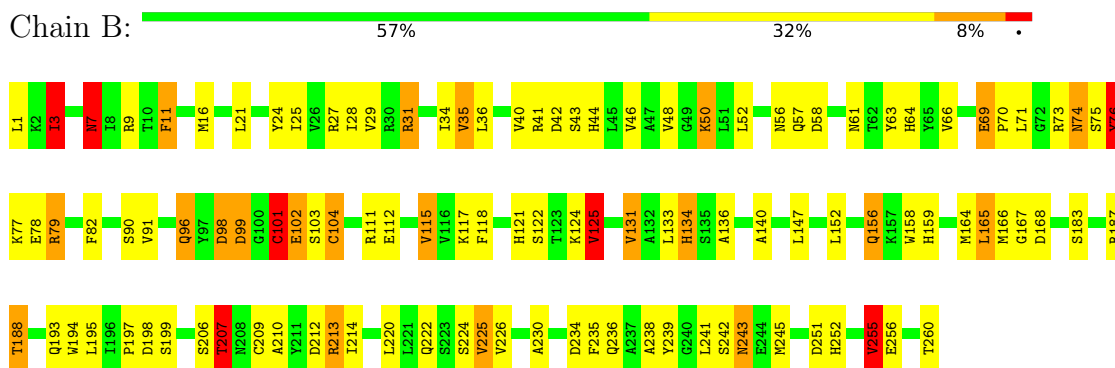
3 Residue-property plots

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: Actin, gamma-enteric smooth muscle



- Molecule 2: Deoxyribonuclease-1



- Molecule 3: alpha-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-3)-[beta-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain C:  100%

MA1
MA2
MA3
MA4
MA5
MA6
MA7

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	42.00Å 225.30Å 77.40Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 1.80 10.00 – 1.80	Depositor EDS
% Data completeness (in resolution range)	(Not available) (10.00-1.80) 77.0 (10.00-1.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$	-	Xtriage
Refinement program	X-PLOR	Depositor
R, R_{free}	0.196 , (Not available) 0.203 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	16.0	Xtriage
Anisotropy	0.364	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.07 , 35.5	EDS
L-test for twinning ¹	$\langle L \rangle = 0.30$, $\langle L^2 \rangle = 0.14$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5472	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.98% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Theoretical 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: CA, NAG, HIC, MAN, ATP, BMA

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	1.09	11/2974 (0.4%)	1.93	82/4025 (2.0%)
2	B	1.22	12/2095 (0.6%)	2.06	65/2853 (2.3%)
All	All	1.15	23/5069 (0.5%)	1.99	147/6878 (2.1%)

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	44	HIS	CD2-NE2	-9.14	1.27	1.37
1	A	160	HIS	CD2-NE2	-8.28	1.28	1.37
1	A	172	HIS	CD2-NE2	-7.22	1.29	1.37
2	B	134	HIS	CD2-NE2	-7.22	1.29	1.37
1	A	356	ILE	CA-CB	7.13	1.63	1.54
1	A	87	HIS	CD2-NE2	-6.89	1.30	1.37
1	A	39	HIS	CD2-NE2	-6.70	1.30	1.37
2	B	252	HIS	CD2-NE2	-6.68	1.30	1.37
2	B	252	HIS	CG-ND1	-6.25	1.31	1.38
2	B	115	VAL	CA-CB	6.25	1.63	1.54
1	A	100	HIS	CD2-NE2	-6.03	1.31	1.37
2	B	121	HIS	CD2-NE2	-6.00	1.31	1.37
2	B	159	HIS	CD2-NE2	-5.89	1.31	1.37
1	A	370	HIS	CD2-NE2	-5.88	1.31	1.37
1	A	274	HIS	CD2-NE2	-5.76	1.31	1.37
2	B	64	HIS	CD2-NE2	-5.66	1.31	1.37
1	A	39	HIS	CG-ND1	-5.56	1.32	1.38
1	A	87	HIS	CG-ND1	-5.49	1.32	1.38
1	A	86	HIS	CD2-NE2	-5.43	1.31	1.37
2	B	82	PHE	CA-CB	-5.36	1.45	1.53
2	B	44	HIS	CG-ND1	-5.29	1.32	1.38
2	B	225	VAL	CA-CB	5.17	1.60	1.54
2	B	158	TRP	CG-CD2	-5.02	1.34	1.43

All (147) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	368	ILE	N-CA-C	15.39	124.88	110.42
1	A	126	PHE	CA-CB-CG	-11.53	102.27	113.80
2	B	255	VAL	N-CA-CB	-11.35	97.16	111.46
2	B	7	ASN	OD1-CG-ND2	-10.72	111.88	122.60
1	A	364	ALA	N-CA-C	10.44	133.04	110.80
2	B	104	CYS	N-CA-C	9.14	124.23	112.34
1	A	63	ILE	N-CA-CB	-8.92	100.34	111.94
2	B	251	ASP	CA-CB-CG	8.91	121.52	112.60
1	A	362	ASP	CA-CB-CG	8.75	121.35	112.60
1	A	311	ARG	CA-CB-CG	-8.71	96.68	114.10
1	A	370	HIS	N-CA-C	8.69	121.53	110.24
2	B	255	VAL	CB-CA-C	8.67	123.62	110.96
1	A	372	LYS	N-CA-C	8.55	129.02	110.80
2	B	226	VAL	N-CA-C	-8.31	101.48	108.63
2	B	31	ARG	NE-CZ-NH2	8.31	126.68	119.20
2	B	222	GLN	OE1-CD-NE2	-8.10	114.50	122.60
1	A	371	ARG	N-CA-C	8.02	127.88	110.80
2	B	125	VAL	N-CA-C	-7.96	99.32	109.58
1	A	322	SER	N-CA-C	7.79	127.40	110.80
1	A	46	MET	N-CA-C	7.78	121.78	109.50
1	A	11	ASN	OD1-CG-ND2	-7.72	114.88	122.60
2	B	79	ARG	NE-CZ-NH2	7.65	126.08	119.20
1	A	65	THR	N-CA-C	-7.64	94.47	107.99
1	A	279	ASN	OD1-CG-ND2	-7.53	115.07	122.60
1	A	338	VAL	N-CA-C	-7.52	103.35	110.42
2	B	255	VAL	N-CA-C	-7.36	97.13	107.80
1	A	334	ARG	N-CA-C	7.24	121.56	112.87
2	B	76	TYR	N-CA-C	-7.12	99.95	110.48
2	B	136	ALA	N-CA-C	-7.03	99.47	109.24
1	A	310	ASP	CA-CB-CG	-7.02	105.58	112.60
2	B	188	THR	N-CA-CB	-6.88	99.95	110.20
1	A	274	HIS	CB-CG-CD2	-6.78	122.39	131.20
2	B	44	HIS	CB-CG-CD2	-6.75	122.43	131.20
1	A	337	SER	N-CA-C	6.72	119.46	111.33
2	B	42	ASP	CA-CB-CG	6.67	119.27	112.60
2	B	99	ASP	CA-CB-CG	6.64	119.24	112.60
2	B	212	ASP	CA-CB-CG	-6.63	105.97	112.60
2	B	238	ALA	N-CA-C	6.62	118.50	111.28
1	A	331	PRO	N-CA-C	6.53	116.74	110.47
1	A	199	PHE	N-CA-C	-6.51	94.87	107.44
1	A	44	VAL	O-C-N	-6.50	115.52	121.89
1	A	253	ARG	NE-CZ-NH2	6.46	125.02	119.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	44	VAL	CA-C-O	-6.44	114.35	121.05
1	A	262	GLN	OE1-CD-NE2	-6.39	116.21	122.60
1	A	86	HIS	CB-CG-CD2	-6.38	122.90	131.20
1	A	243	ASP	CA-CB-CG	6.38	118.98	112.60
2	B	9	ARG	NE-CZ-NH1	-6.36	115.14	121.50
2	B	235	PHE	CA-CB-CG	6.34	120.14	113.80
1	A	274	HIS	CB-CG-ND1	6.33	132.20	122.70
2	B	103	SER	CA-CB-OG	6.32	123.75	111.10
1	A	44	VAL	N-CA-C	6.26	117.75	110.62
1	A	311	ARG	NE-CZ-NH1	-6.25	115.25	121.50
2	B	98	ASP	N-CA-C	-6.20	100.11	108.86
1	A	313	GLN	OE1-CD-NE2	-6.19	116.41	122.60
1	A	15	LEU	N-CA-C	6.19	119.49	109.40
1	A	78	TRP	CG-CD2-CE3	6.17	140.07	133.90
1	A	331	PRO	O-C-N	-6.14	118.49	121.31
2	B	96	GLN	N-CA-C	-6.12	98.92	108.90
2	B	7	ASN	N-CA-C	-6.12	97.73	108.20
1	A	68	TYR	CB-CA-C	6.08	118.24	110.34
2	B	222	GLN	CB-CG-CD	6.08	122.93	112.60
1	A	38	ARG	NE-CZ-NH1	-6.06	115.44	121.50
1	A	371	ARG	NE-CZ-NH2	6.05	124.65	119.20
1	A	273	ILE	N-CA-CB	-6.05	101.25	111.23
2	B	140	ALA	N-CA-C	6.04	118.35	111.11
2	B	121	HIS	CB-CG-CD2	-6.03	123.37	131.20
1	A	311	ARG	NE-CZ-NH2	5.99	124.59	119.20
2	B	58	ASP	CA-CB-CG	5.92	118.52	112.60
2	B	198	ASP	CA-CB-CG	5.89	118.49	112.60
1	A	68	TYR	CA-CB-CG	5.89	124.50	113.90
1	A	23	ASP	CA-CB-CG	5.88	118.48	112.60
1	A	370	HIS	CB-CG-CD2	-5.87	123.57	131.20
2	B	156	GLN	OE1-CD-NE2	-5.84	116.76	122.60
1	A	75	ILE	CB-CG1-CD1	5.83	126.05	113.80
2	B	213	ARG	NE-CZ-NH1	-5.82	115.68	121.50
2	B	34	ILE	O-C-N	-5.80	117.15	123.18
1	A	161	ASN	CA-CB-CG	-5.77	106.83	112.60
1	A	187	TYR	N-CA-C	-5.75	105.09	111.36
2	B	73	ARG	NE-CZ-NH2	5.75	124.38	119.20
1	A	311	ARG	CA-C-O	-5.74	114.34	120.42
2	B	222	GLN	CG-CD-NE2	5.72	124.98	116.40
2	B	193	GLN	OE1-CD-NE2	-5.69	116.91	122.60
2	B	168	ASP	CA-CB-CG	5.69	118.29	112.60
1	A	115	ARG	CA-CB-CG	-5.64	102.82	114.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	125	VAL	N-CA-CB	5.64	118.93	110.13
2	B	31	ARG	NE-CZ-NH1	-5.63	115.87	121.50
1	A	226	MET	N-CA-CB	-5.62	101.57	109.94
1	A	160	HIS	CB-CG-CD2	-5.62	123.90	131.20
1	A	193	THR	CA-CB-OG1	-5.61	101.19	109.60
2	B	165	LEU	O-C-N	-5.56	116.73	123.29
1	A	295	ASN	OD1-CG-ND2	-5.51	117.09	122.60
1	A	218	VAL	CA-C-O	5.50	126.42	120.53
1	A	303	THR	N-CA-CB	-5.50	101.52	110.42
1	A	114	ASN	CB-CA-C	-5.49	101.67	110.79
2	B	61	ASN	CB-CG-ND2	5.49	124.64	116.40
2	B	134	HIS	CB-CG-ND1	5.49	130.94	122.70
2	B	234	ASP	CA-CB-CG	5.46	118.06	112.60
1	A	371	ARG	N-CA-CB	-5.45	101.28	110.49
1	A	128	VAL	CA-C-N	5.45	124.63	118.97
1	A	128	VAL	C-N-CA	5.45	124.63	118.97
2	B	117	LYS	CA-CB-CG	-5.44	103.22	114.10
1	A	246	VAL	N-CA-CB	-5.43	105.88	112.07
2	B	11	PHE	CA-CB-CG	-5.39	108.41	113.80
1	A	262	GLN	CG-CD-NE2	5.37	124.46	116.40
1	A	252	GLU	CB-CG-CD	5.37	121.72	112.60
2	B	134	HIS	CB-CG-CD2	-5.35	124.24	131.20
1	A	39	HIS	CB-CG-CD2	-5.34	124.25	131.20
1	A	214	LYS	N-CA-C	5.34	119.78	112.68
1	A	115	ARG	N-CA-C	-5.34	105.46	111.28
1	A	333	GLU	CA-CB-CG	5.33	124.76	114.10
2	B	124	LYS	CA-CB-CG	-5.31	103.48	114.10
2	B	27	ARG	N-CA-CB	-5.30	102.06	110.28
1	A	27	ARG	N-CA-C	-5.30	106.65	113.01
2	B	242	SER	N-CA-C	-5.30	103.69	110.53
2	B	9	ARG	CA-C-O	-5.29	114.73	120.81
2	B	207	THR	N-CA-CB	-5.27	102.44	110.45
2	B	46	VAL	O-C-N	-5.24	116.78	121.87
1	A	328	ILE	N-CA-C	-5.24	100.35	107.99
1	A	42	VAL	N-CA-CB	-5.23	103.36	111.05
1	A	226	MET	CA-CB-CG	5.23	124.56	114.10
1	A	195	ARG	CB-CG-CD	-5.22	99.30	111.30
2	B	188	THR	CA-CB-OG1	-5.21	101.79	109.60
1	A	119	THR	CA-C-N	5.19	127.19	120.44
1	A	119	THR	C-N-CA	5.19	127.19	120.44
1	A	139	LEU	CA-C-O	-5.19	115.05	120.55
1	A	353	GLN	OE1-CD-NE2	-5.18	117.42	122.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	99	ASP	N-CA-C	-5.17	101.92	109.63
2	B	11	PHE	N-CA-C	-5.14	102.31	109.96
1	A	78	TRP	CE2-CD2-CG	-5.14	101.04	107.20
1	A	63	ILE	CA-CB-CG1	-5.13	101.67	110.40
2	B	48	VAL	N-CA-C	-5.12	105.55	110.72
2	B	111	ARG	CD-NE-CZ	-5.11	117.25	124.40
1	A	330	ALA	CA-C-N	5.10	123.38	119.66
1	A	330	ALA	C-N-CA	5.10	123.38	119.66
2	B	73	ARG	CB-CG-CD	-5.10	99.58	111.30
2	B	195	LEU	CA-C-O	-5.10	114.58	120.24
1	A	339	TRP	CE2-CD2-CG	-5.08	101.10	107.20
2	B	64	HIS	N-CA-C	-5.07	102.69	110.14
2	B	3	ILE	O-C-N	-5.06	117.71	123.18
1	A	303	THR	N-CA-C	-5.05	107.09	113.20
1	A	145	GLY	N-CA-C	-5.04	106.19	114.76
2	B	102	GLU	N-CA-C	5.04	116.86	111.36
1	A	270	SER	O-C-N	-5.04	118.25	123.29
2	B	102	GLU	CA-CB-CG	5.03	124.16	114.10
1	A	78	TRP	CG-CD1-NE1	-5.02	103.67	110.20
2	B	96	GLN	CG-CD-NE2	5.01	123.92	116.40
2	B	3	ILE	CB-CG1-CD1	-5.00	103.30	113.80

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	2924	0	2887	85	0
2	B	2049	0	1981	43	0
3	C	83	0	70	0	0
4	A	31	0	12	4	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
6	A	195	0	0	14	0
6	B	188	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	5472	0	4950	128	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 13.

All (128) 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:119:THR:HG23	1:A:131:MET:HE1	1.59	0.84
2:B:99:ASP:HB3	2:B:104:CYS:HB3	1.59	0.83
2:B:207:THR:HG23	2:B:209:CYS:SG	2.19	0.82
2:B:56:ASN:HD21	2:B:63:TYR:H	1.34	0.74
1:A:219:ALA:HB1	1:A:225:GLU:HG3	1.73	0.70
1:A:239:TYR:HB3	1:A:247:ILE:HG22	1.75	0.69
1:A:191:ILE:HG21	1:A:252:GLU:HG3	1.75	0.67
2:B:74:ASN:HA	2:B:77:LYS:NZ	2.12	0.65
1:A:144:SER:HB2	1:A:146:ARG:HG2	1.78	0.64
1:A:152:LEU:HD11	1:A:273:ILE:HD12	1.79	0.64
2:B:11:PHE:HD2	2:B:40:VAL:HG13	1.64	0.63
2:B:239:TYR:HB2	2:B:241:LEU:HD13	1.82	0.61
1:A:34:VAL:HG21	1:A:80:ASP:HB3	1.82	0.60
1:A:122:MET:HB3	1:A:128:VAL:HG21	1.83	0.60
1:A:187:TYR:O	1:A:190:LYS:HB3	2.03	0.59
1:A:334:ARG:HA	1:A:337:SER:OG	2.02	0.59
2:B:3:ILE:HD11	2:B:164:MET:SD	2.44	0.58
1:A:219:ALA:CB	1:A:225:GLU:HG3	2.34	0.57
1:A:298:LEU:HB3	1:A:303:THR:HG23	1.85	0.57
1:A:195:ARG:NH1	1:A:249:ILE:HA	2.21	0.56
1:A:282:MET:SD	6:A:531:HOH:O	2.58	0.56
1:A:8:VAL:HG21	1:A:343:SER:HA	1.88	0.56
1:A:301:GLY:HA3	4:A:401:ATP:O4'	2.06	0.56
1:A:352:GLN:HA	1:A:355:TRP:HE3	1.71	0.55
1:A:153:ASP:HA	1:A:299:SER:O	2.06	0.55
1:A:161:ASN:HD22	1:A:276:THR:HG22	1.70	0.55
2:B:41:ARG:NE	2:B:76:TYR:HE1	2.05	0.55
2:B:70:PRO:HD3	2:B:79:ARG:NH2	2.22	0.55
1:A:301:GLY:O	1:A:304:MET:HB2	2.07	0.55
1:A:72:HIC:HD2	1:A:157:GLY:O	2.07	0.54
2:B:131:VAL:HG13	2:B:165:LEU:HG	1.88	0.54
1:A:209:ARG:O	1:A:212:LYS:HB3	2.07	0.54
2:B:24:TYR:CE2	2:B:245:MET:HE2	2.43	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:101:CYS:O	2:B:104:CYS:HB2	2.08	0.53
2:B:11:PHE:HE1	2:B:16:MET:HE3	1.71	0.53
1:A:316:ILE:HD12	1:A:328:ILE:HD11	1.90	0.53
2:B:21:LEU:HD23	2:B:245:MET:HE1	1.90	0.53
2:B:74:ASN:HA	2:B:77:LYS:HZ1	1.73	0.53
1:A:217:TYR:O	1:A:254:PHE:HA	2.09	0.53
1:A:98:GLU:HA	1:A:127:ASN:O	2.09	0.52
1:A:88:SER:O	1:A:93:LEU:HB2	2.10	0.52
2:B:56:ASN:ND2	2:B:63:TYR:H	2.04	0.52
1:A:229:ALA:HB2	1:A:235:LEU:HD12	1.91	0.52
2:B:11:PHE:CE1	2:B:16:MET:HE3	2.44	0.52
2:B:224:SER:O	2:B:260:THR:HG22	2.10	0.52
1:A:192:LEU:HD21	1:A:211:ILE:HD11	1.92	0.52
1:A:152:LEU:HD21	1:A:273:ILE:HG23	1.92	0.51
1:A:249:ILE:HB	1:A:252:GLU:HB2	1.92	0.51
1:A:42:VAL:HG12	2:B:66:VAL:HG13	1.92	0.51
2:B:3:ILE:HD12	2:B:166:MET:HE1	1.92	0.51
2:B:133:LEU:O	2:B:167:GLY:HA3	2.11	0.51
1:A:122:MET:HB3	1:A:128:VAL:CG2	2.41	0.51
1:A:102:THR:O	1:A:131:MET:HA	2.11	0.50
1:A:24:ASP:HB2	6:A:516:HOH:O	2.11	0.50
1:A:33:ILE:HD11	1:A:58:GLN:HB2	1.94	0.49
1:A:160:HIS:CE1	1:A:176:ARG:HG3	2.47	0.49
2:B:28:ILE:O	2:B:31:ARG:HG3	2.12	0.49
1:A:112:LYS:HB3	6:A:659:HOH:O	2.13	0.49
1:A:142:TYR:HA	1:A:146:ARG:O	2.12	0.49
2:B:210:ALA:O	2:B:213:ARG:HD3	2.13	0.49
2:B:7:ASN:C	2:B:7:ASN:HD22	2.21	0.49
1:A:159:THR:HG21	1:A:273:ILE:HD11	1.93	0.49
1:A:170:LEU:HD12	1:A:284:CYS:SG	2.53	0.48
1:A:70:ILE:HD11	1:A:81:MET:HE1	1.95	0.48
1:A:71:GLU:HB3	6:A:576:HOH:O	2.13	0.48
2:B:11:PHE:HB3	2:B:40:VAL:HA	1.96	0.48
2:B:78:GLU:OE2	2:B:134:HIS:HD2	1.96	0.47
2:B:236:GLN:NE2	2:B:243:ASN:HA	2.29	0.47
1:A:120:GLN:O	1:A:124:GLU:HB2	2.14	0.47
2:B:50:LYS:HE3	2:B:50:LYS:HA	1.96	0.47
1:A:239:TYR:HB3	1:A:247:ILE:CG2	2.44	0.47
1:A:249:ILE:H	1:A:249:ILE:HD13	1.78	0.47
1:A:263:PRO:HB3	1:A:268:MET:HB3	1.97	0.47
2:B:96:GLN:HG2	6:B:532:HOH:O	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:90:SER:O	2:B:118:PHE:HA	2.15	0.47
2:B:1:LEU:N	6:B:568:HOH:O	2.48	0.46
1:A:39:HIS:HA	6:A:672:HOH:O	2.16	0.46
1:A:19:GLY:HA3	6:A:626:HOH:O	2.16	0.46
2:B:197:PRO:HA	6:B:513:HOH:O	2.15	0.46
1:A:192:LEU:HB3	1:A:199:PHE:HE1	1.81	0.46
1:A:215:LEU:O	1:A:253:ARG:HD2	2.14	0.46
1:A:336:TYR:O	1:A:340:ILE:HG13	2.16	0.45
1:A:183:ASP:HB3	6:A:588:HOH:O	2.16	0.45
1:A:170:LEU:HB3	1:A:373:CYS:HB3	1.97	0.45
1:A:11:ASN:ND2	1:A:85:TRP:HE1	2.14	0.45
1:A:239:TYR:HA	6:A:618:HOH:O	2.16	0.45
1:A:152:LEU:CD2	1:A:273:ILE:HG23	2.47	0.45
1:A:115:ARG:HD3	1:A:368:ILE:HG12	1.99	0.45
1:A:101:PRO:HA	1:A:130:ALA:O	2.18	0.44
1:A:304:MET:HB3	4:A:401:ATP:C6	2.53	0.44
2:B:11:PHE:CD2	2:B:40:VAL:HG13	2.50	0.44
2:B:70:PRO:HB3	2:B:77:LYS:HB3	2.00	0.44
2:B:29:VAL:HG13	2:B:35:VAL:HG11	1.99	0.44
2:B:125:VAL:HG13	2:B:220:LEU:HB3	2.00	0.44
1:A:3:GLU:HB3	6:A:597:HOH:O	2.17	0.44
1:A:165:TYR:HB3	1:A:170:LEU:HD21	2.00	0.44
1:A:369:VAL:HG23	1:A:371:ARG:NH1	2.33	0.44
2:B:214:ILE:HG13	2:B:255:VAL:HG13	1.98	0.44
1:A:239:TYR:HE2	6:A:633:HOH:O	2.00	0.43
1:A:349:SER:HA	1:A:352:GLN:OE1	2.19	0.43
1:A:297:VAL:HG23	6:A:584:HOH:O	2.17	0.43
1:A:348:LEU:HB3	1:A:350:THR:OG1	2.19	0.43
1:A:236:GLU:HA	1:A:250:GLY:HA2	2.00	0.43
1:A:368:ILE:HG21	6:A:646:HOH:O	2.17	0.43
2:B:69:GLU:O	2:B:71:LEU:HG	2.19	0.42
1:A:78:TRP:CD2	1:A:117:LYS:HG2	2.55	0.42
1:A:91:ASN:O	1:A:94:ARG:HD3	2.19	0.42
2:B:16:MET:CE	2:B:25:ILE:HD12	2.49	0.42
1:A:132:TYR:CG	1:A:351:PHE:HZ	2.38	0.42
1:A:335:LYS:HE3	1:A:335:LYS:HB2	1.86	0.42
2:B:147:LEU:HD12	2:B:147:LEU:HA	1.89	0.42
1:A:303:THR:HG22	1:A:334:ARG:HE	1.85	0.42
2:B:230:ALA:HA	2:B:256:GLU:O	2.20	0.41
1:A:131:MET:O	1:A:356:ILE:HG23	2.20	0.41
1:A:155:GLY:O	1:A:180:ALA:HB1	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:3:ILE:HD11	2:B:166:MET:SD	2.60	0.41
1:A:6:ALA:HA	1:A:101:PRO:HD2	2.02	0.41
1:A:115:ARG:HD2	1:A:368:ILE:HD11	2.02	0.41
1:A:153:ASP:O	1:A:159:THR:HA	2.20	0.41
1:A:96:ALA:O	1:A:100:HIS:HD2	2.04	0.41
1:A:156:ASP:HB2	4:A:401:ATP:H5'1	2.03	0.41
1:A:104:LEU:O	1:A:133:VAL:HA	2.20	0.41
1:A:172:HIS:HB2	1:A:373:CYS:O	2.21	0.41
1:A:140:SER:O	1:A:143:ALA:HB3	2.20	0.40
1:A:214:LYS:HE2	6:A:633:HOH:O	2.21	0.40
4:A:401:ATP:H1'	6:A:506:HOH:O	2.21	0.40
1:A:285:ASP:O	1:A:288:ILE:HG12	2.21	0.40
2:B:187:ARG:HD3	2:B:194:TRP:CE2	2.56	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	371/374 (99%)	339 (91%)	22 (6%)	10 (3%)	4	0
2	B	258/260 (99%)	244 (95%)	9 (4%)	5 (2%)	6	1
All	All	629/634 (99%)	583 (93%)	31 (5%)	15 (2%)	4	1

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	322	SER
1	A	367	SER
1	A	373	CYS
1	A	286	ILE
1	A	372	LYS

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Mol	Chain	Res	Type
2	B	57	GLN
1	A	195	ARG
2	B	43	SER
1	A	273	ILE
2	B	75	SER
2	B	74	ASN
2	B	101	CYS
1	A	250	GLY
1	A	366	PRO
1	A	181	GLY

5.3.2 Protein sidechains [i](#)

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	316/316 (100%)	274 (87%)	42 (13%)	4	1
2	B	229/229 (100%)	202 (88%)	27 (12%)	5	1
All	All	545/545 (100%)	476 (87%)	69 (13%)	4	1

All (69) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	42	VAL
1	A	63	ILE
1	A	71	GLU
1	A	74	ILE
1	A	75	ILE
1	A	82	GLU
1	A	83	LYS
1	A	93	LEU
1	A	98	GLU
1	A	111	PRO
1	A	116	GLU
1	A	124	GLU
1	A	127	ASN

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Mol	Chain	Res	Type
1	A	150	ILE
1	A	177	LEU
1	A	179	LEU
1	A	192	LEU
1	A	195	ARG
1	A	207	ILE
1	A	236	GLU
1	A	247	ILE
1	A	249	ILE
1	A	260	LEU
1	A	273	ILE
1	A	274	HIS
1	A	303	THR
1	A	304	MET
1	A	310	ASP
1	A	311	ARG
1	A	319	LEU
1	A	322	SER
1	A	323	THR
1	A	327	LYS
1	A	333	GLU
1	A	335	LYS
1	A	348	LEU
1	A	352	GLN
1	A	355	TRP
1	A	356	ILE
1	A	360	GLU
1	A	370	HIS
1	A	374	PHE
2	B	3	ILE
2	B	7	ASN
2	B	35	VAL
2	B	36	LEU
2	B	50	LYS
2	B	52	LEU
2	B	69	GLU
2	B	76	TYR
2	B	91	VAL
2	B	98	ASP
2	B	101	CYS
2	B	102	GLU
2	B	112	GLU

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Mol	Chain	Res	Type
2	B	115	VAL
2	B	122	SER
2	B	125	VAL
2	B	131	VAL
2	B	152	LEU
2	B	156	GLN
2	B	183	SER
2	B	188	THR
2	B	199	SER
2	B	206	SER
2	B	207	THR
2	B	225	VAL
2	B	243	ASN
2	B	255	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	11	ASN
1	A	40	GLN
1	A	160	HIS
1	A	161	ASN
1	A	313	GLN
2	B	7	ASN
2	B	44	HIS
2	B	56	ASN
2	B	134	HIS
2	B	170	ASN
2	B	193	GLN
2	B	236	GLN
2	B	243	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

1 non-standard protein/DNA/RNA residue is 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$
1	HIC	A	72	1	10,11,12	1.63	1 (10%)	9,14,16	13.27	6 (66%)

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
1	HIC	A	72	1	-	1/5/6/8	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	72	HIC	CD2-NE2	-3.81	1.31	1.37

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	72	HIC	CD2-NE2-CE1	29.36	118.31	106.71
1	A	72	HIC	NE2-CE1-ND1	-26.28	102.60	112.66
1	A	72	HIC	CZ-NE2-CE1	-3.51	118.42	126.71
1	A	72	HIC	CE1-ND1-CG	2.94	110.52	105.80
1	A	72	HIC	CA-CB-CG	2.59	120.15	113.77
1	A	72	HIC	CZ-NE2-CD2	-2.11	123.27	126.03

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	72	HIC	CA-CB-CG-ND1

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	72	HIC	1	0

5.5 Carbohydrates [i](#)

7 monosaccharides are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	NAG	C	1	3,2	14,14,15	1.23	1 (7%)	17,19,21	1.70	4 (23%)
3	NAG	C	2	3	14,14,15	1.37	2 (14%)	17,19,21	1.98	5 (29%)
3	BMA	C	3	3	11,11,12	1.92	3 (27%)	15,15,17	2.09	3 (20%)
3	BMA	C	4	3	11,11,12	1.60	1 (9%)	15,15,17	1.25	2 (13%)
3	MAN	C	5	3	11,11,12	1.28	2 (18%)	15,15,17	2.21	2 (13%)
3	MAN	C	6	3	11,11,12	1.76	4 (36%)	15,15,17	2.11	1 (6%)
3	BMA	C	7	3	11,11,12	1.29	2 (18%)	15,15,17	1.00	1 (6%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	C	1	3,2	-	3/6/23/26	0/1/1/1
3	NAG	C	2	3	-	3/6/23/26	0/1/1/1
3	BMA	C	3	3	-	2/2/19/22	0/1/1/1
3	BMA	C	4	3	-	0/2/19/22	0/1/1/1
3	MAN	C	5	3	-	0/2/19/22	0/1/1/1
3	MAN	C	6	3	-	0/2/19/22	1/1/1/1
3	BMA	C	7	3	-	0/2/19/22	0/1/1/1

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	3	BMA	C2-C3	4.27	1.59	1.52
3	C	4	BMA	C2-C3	4.14	1.58	1.52
3	C	6	MAN	C4-C5	3.43	1.60	1.53
3	C	3	BMA	O3-C3	2.89	1.50	1.43
3	C	2	NAG	O5-C5	2.85	1.49	1.43
3	C	1	NAG	C4-C5	2.80	1.59	1.53
3	C	5	MAN	C4-C5	2.43	1.58	1.53
3	C	7	BMA	C1-C2	2.36	1.57	1.52
3	C	6	MAN	C4-C3	2.32	1.58	1.52
3	C	3	BMA	C4-C3	2.21	1.58	1.52
3	C	7	BMA	C2-C3	2.19	1.55	1.52
3	C	6	MAN	O5-C5	2.12	1.47	1.43
3	C	2	NAG	C1-C2	2.08	1.55	1.52
3	C	5	MAN	C1-C2	2.03	1.57	1.52
3	C	6	MAN	O3-C3	2.03	1.48	1.43

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	6	MAN	C1-O5-C5	7.34	122.02	112.19
3	C	5	MAN	C1-O5-C5	6.55	120.97	112.19
3	C	3	BMA	C1-O5-C5	6.49	120.89	112.19
3	C	2	NAG	C1-O5-C5	4.95	118.82	112.19
3	C	1	NAG	C1-C2-N2	3.86	116.51	110.43
3	C	1	NAG	C1-O5-C5	3.72	117.17	112.19
3	C	5	MAN	C3-C4-C5	3.50	116.58	110.23
3	C	2	NAG	C3-C4-C5	-3.50	103.89	110.23
3	C	2	NAG	C4-C3-C2	-3.30	106.19	111.02
3	C	1	NAG	C2-N2-C7	3.04	126.98	122.90
3	C	3	BMA	C3-C4-C5	2.84	115.39	110.23
3	C	4	BMA	C1-O5-C5	2.65	115.73	112.19
3	C	1	NAG	O5-C1-C2	-2.61	107.25	111.29
3	C	7	BMA	C1-O5-C5	2.58	115.64	112.19
3	C	3	BMA	C1-C2-C3	2.45	113.21	109.64
3	C	4	BMA	O3-C3-C2	2.35	114.84	110.05
3	C	2	NAG	C1-C2-N2	2.22	113.94	110.43
3	C	2	NAG	O4-C4-C5	2.11	114.52	109.32

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	1	NAG	C1-C2-N2-C7

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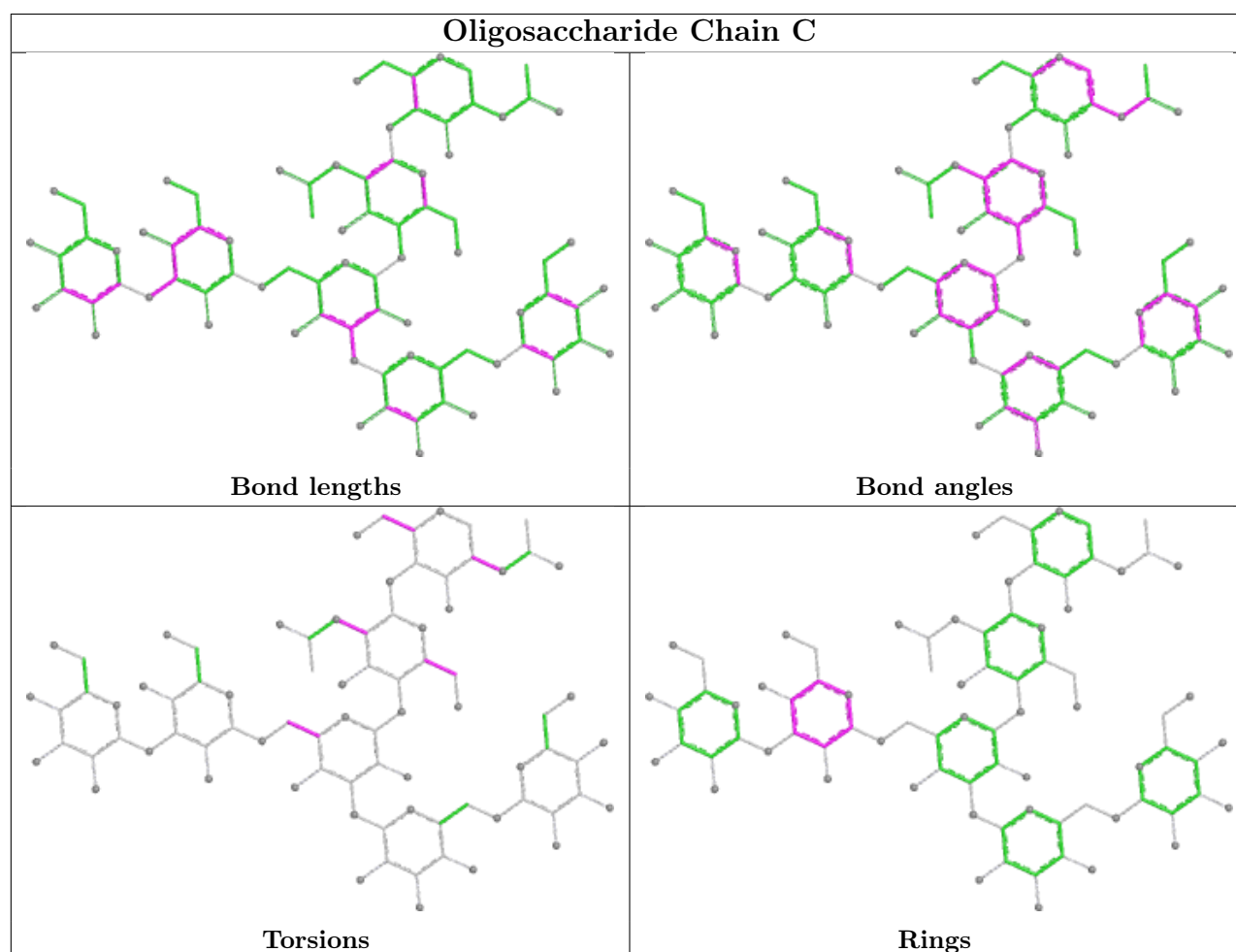
Mol	Chain	Res	Type	Atoms
3	C	2	NAG	C1-C2-N2-C7
3	C	3	BMA	C4-C5-C6-O6
3	C	3	BMA	O5-C5-C6-O6
3	C	1	NAG	C4-C5-C6-O6
3	C	2	NAG	C4-C5-C6-O6
3	C	1	NAG	O5-C5-C6-O6
3	C	2	NAG	O5-C5-C6-O6

All (1) ring outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	6	MAN	C1-C2-C3-C4-C5-O5

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.



5.6 Ligand geometry

Of 3 ligands modelled in this entry, 2 are monoatomic - leaving 1 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	ATP	A	401	5	32,33,33	1.33	2 (6%)	48,52,52	1.18	4 (8%)

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	ATP	A	401	5	-	3/22/38/38	0/3/3/3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	401	ATP	PB-O3B	4.74	1.64	1.59
4	A	401	ATP	PG-O3G	-2.38	1.46	1.54

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	401	ATP	O3B-PB-O1B	-3.49	100.21	110.70
4	A	401	ATP	O4'-C1'-N9	2.81	113.49	108.09
4	A	401	ATP	C4'-O4'-C1'	-2.23	104.55	109.47
4	A	401	ATP	O4'-C4'-C3'	2.15	109.43	105.15

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	401	ATP	PB-O3B-PG-O2G
4	A	401	ATP	C5'-O5'-PA-O1A

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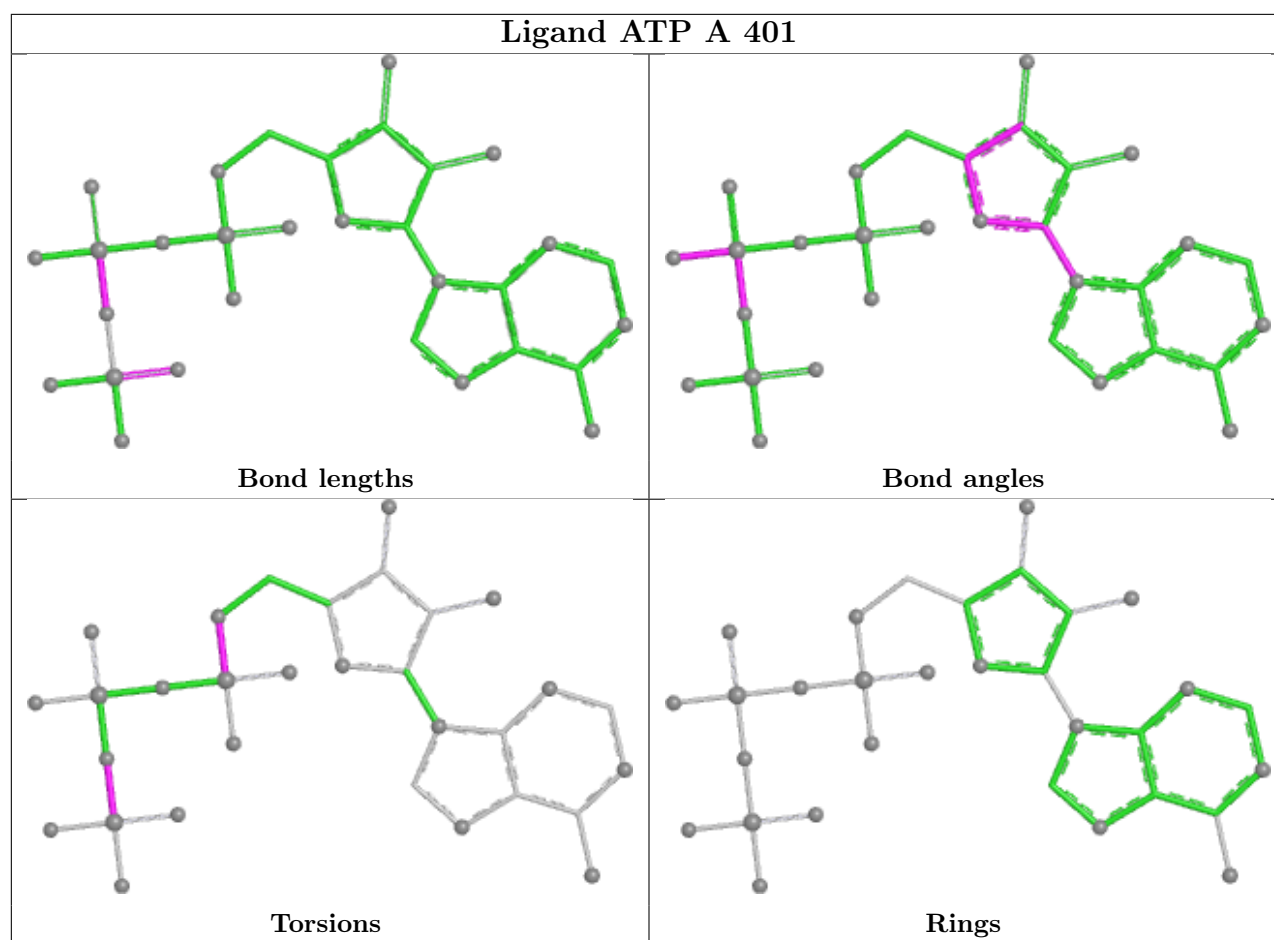
Mol	Chain	Res	Type	Atoms
4	A	401	ATP	PB-O3B-PG-O3G

There are no ring outliers.

1 monomer is involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	401	ATP	4	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	373/374 (99%)	0.02	11 (2%) 53 53	10, 34, 64, 72	0
2	B	260/260 (100%)	-0.72	0 100 100	4, 16, 42, 57	0
All	All	633/634 (99%)	-0.28	11 (1%) 69 69	4, 26, 61, 72	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	374	PHE	4.3
1	A	364	ALA	3.2
1	A	348	LEU	3.1
1	A	368	ILE	2.9
1	A	355	TRP	2.8
1	A	351	PHE	2.7
1	A	244	GLY	2.6
1	A	373	CYS	2.1
1	A	167	GLY	2.1
1	A	356	ILE	2.1
1	A	350	THR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains

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.

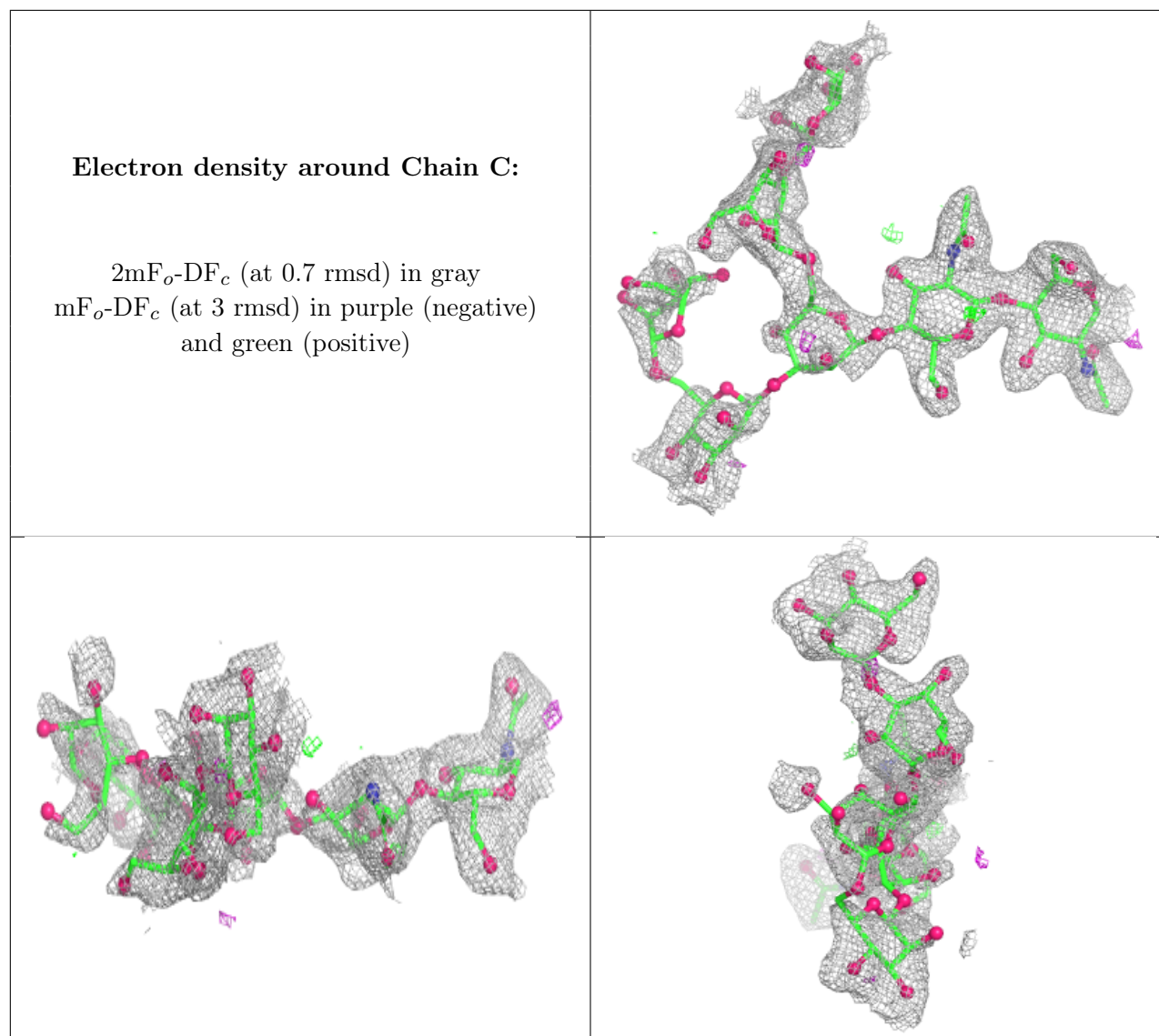
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
1	HIC	A	72	11/12	0.96	0.05	14,18,23,26	0

6.3 Carbohydrates

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	MAN	C	5	11/12	0.62	0.15	65,70,72,73	0
3	MAN	C	6	11/12	0.67	0.11	60,62,63,65	0
3	BMA	C	7	11/12	0.68	0.10	62,64,66,67	0
3	BMA	C	4	11/12	0.70	0.12	66,68,70,71	0
3	BMA	C	3	11/12	0.71	0.12	60,63,65,67	0
3	NAG	C	2	14/15	0.79	0.10	53,58,60,61	0
3	NAG	C	1	14/15	0.88	0.07	36,40,45,49	0

The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.



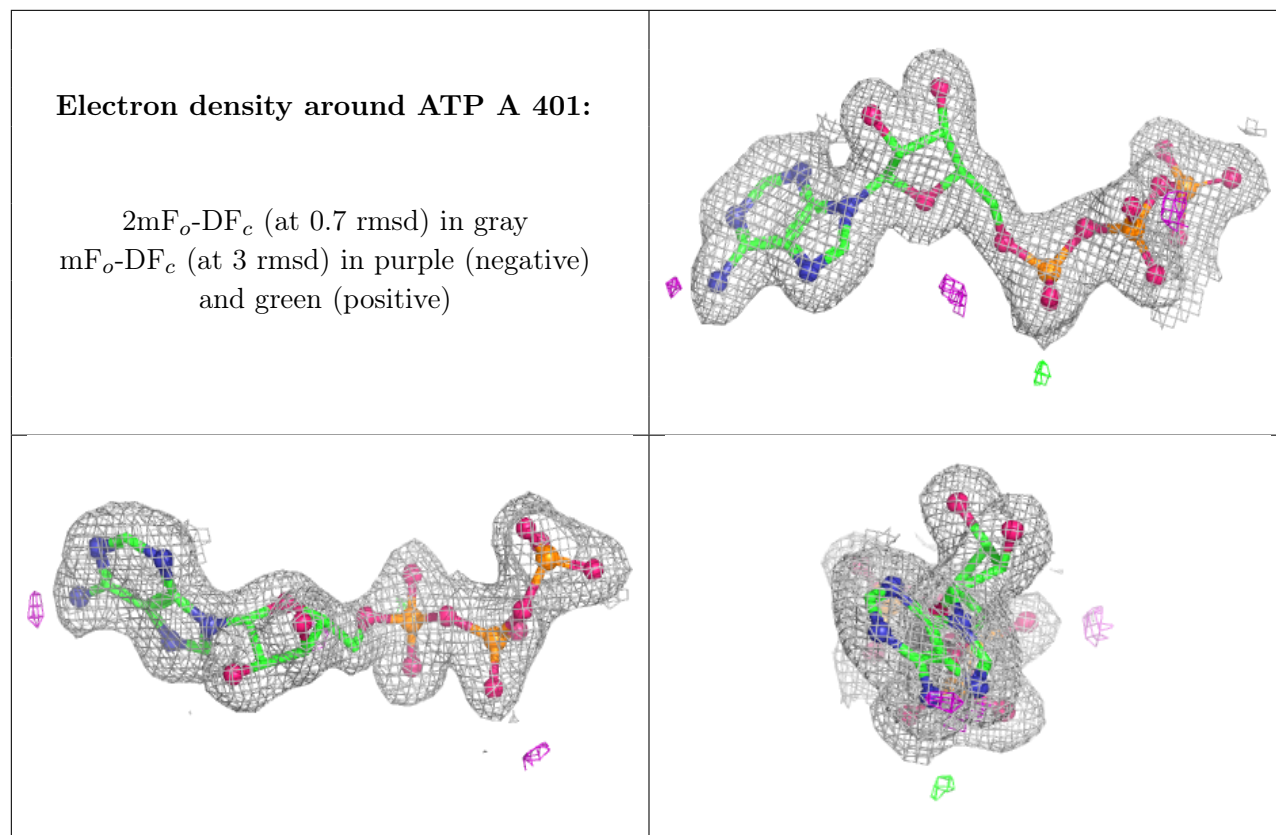
6.4 Ligands [i](#)

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(Å ²)	Q<0.9
5	CA	A	402	1/1	0.98	0.03	29,29,29,29	0
4	ATP	A	401	31/31	0.99	0.04	11,19,26,28	0
5	CA	B	308	1/1	0.99	0.05	15,15,15,15	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.



6.5 Other polymers [i](#)

There are no such residues in this entry.