



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 16, 2025 – 06:32 am GMT

PDB ID : 9GBP
Title : Human Angiotensin-1 converting enzyme N-domain in complex with a diprolyl inhibitor- SG3
Authors : Gregory, K.S.; Cozier, G.E.; Acharya, K.R.
Deposited on : 2024-07-31
Resolution : 2.00 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.4, CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 3.0
buster-report : 1.1.7 (2018)
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4 : 9.0.003 (Gargrove)
Density-Fitness : 1.0.11
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.40

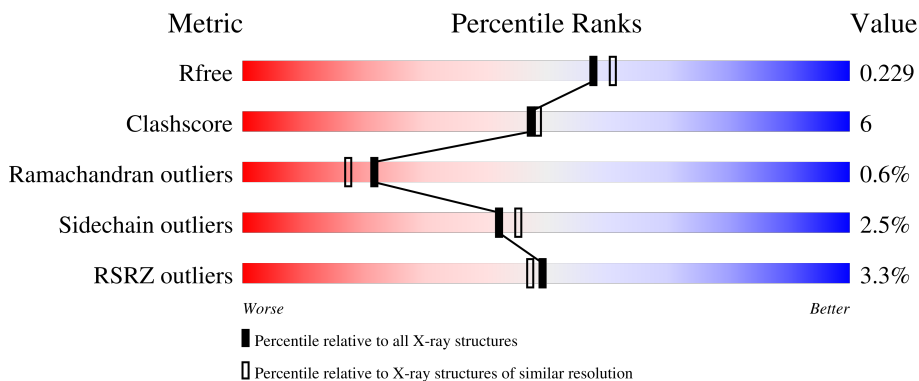
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION



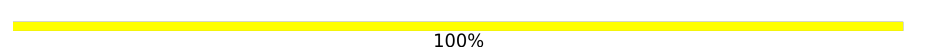
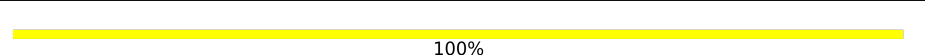
The reported resolution of this entry is 2.00 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	164625	9409 (2.00-2.00)
Clashscore	180529	10737 (2.00-2.00)
Ramachandran outliers	177936	10628 (2.00-2.00)
Sidechain outliers	177891	10627 (2.00-2.00)
RSRZ outliers	164620	9409 (2.00-2.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	628	
1	B	628	
2	E	2	
3	C	3	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
7	PEG	A	704	-	-	X	-
8	EDO	B	709	-	-	X	-

2 Entry composition [i](#)

There are 13 unique types of molecules in this entry. The entry contains 11143 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.

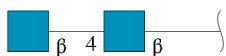
- Molecule 1 is a protein called Angiotensin-converting enzyme, soluble form.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	609	Total	C	N	O	S	0	5	0
			5011	3218	858	915	20			
1	B	604	Total	C	N	O	S	0	3	0
			4948	3175	848	906	19			

There are 16 discrepancies between the modelled and reference sequences:

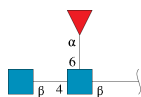
Chain	Residue	Modelled	Actual	Comment	Reference
A	9	GLN	ASN	engineered mutation	UNP P12821
A	25	GLN	ASN	engineered mutation	UNP P12821
A	82	GLN	ASN	engineered mutation	UNP P12821
A	117	GLN	ASN	engineered mutation	UNP P12821
A	131	GLN	ASN	engineered mutation	UNP P12821
A	289	GLN	ASN	engineered mutation	UNP P12821
A	545	ARG	GLN	engineered mutation	UNP P12821
A	576	LEU	PRO	engineered mutation	UNP P12821
B	9	GLN	ASN	engineered mutation	UNP P12821
B	25	GLN	ASN	engineered mutation	UNP P12821
B	82	GLN	ASN	engineered mutation	UNP P12821
B	117	GLN	ASN	engineered mutation	UNP P12821
B	131	GLN	ASN	engineered mutation	UNP P12821
B	289	GLN	ASN	engineered mutation	UNP P12821
B	545	ARG	GLN	engineered mutation	UNP P12821
B	576	LEU	PRO	engineered mutation	UNP P12821

- Molecule 2 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



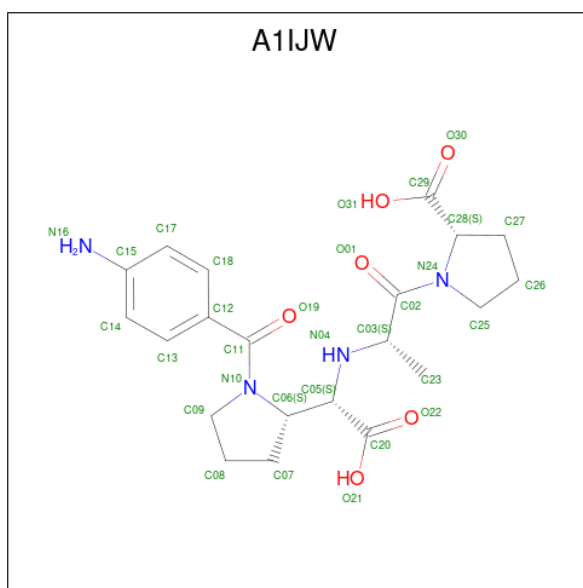
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	E	2	28	16	2	10	0	0	0

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



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
3	C	3	38	22	2	14	0	0	0

- Molecule 4 is (2S)-1-[(2S)-2-[[[(1S)-1-[(2S)-1-(4-aminophenyl)carbonylpyrrolidin-2-yl]-2-oxidanyl-2-oxidanylidene-ethyl]amino]propanoyl]pyrrolidine-2-carboxylic acid (three-letter code: A1IJW) (formula: C₂₁H₂₈N₄O₆) (labeled as "Ligand of Interest" by depositor).



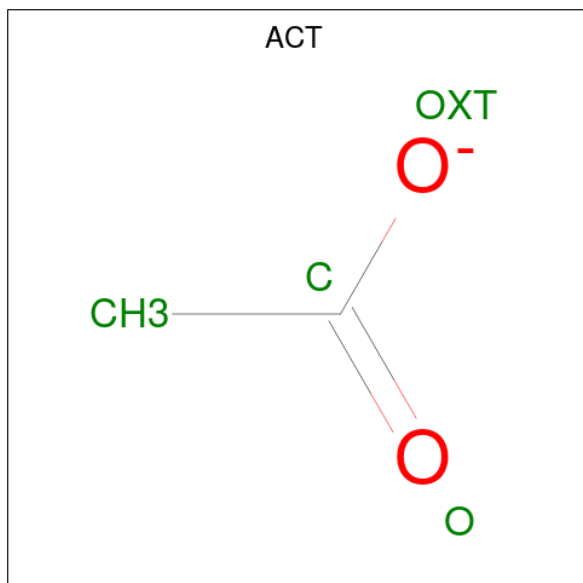
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
4	A	1	31	21	4	6	0	0
4	B	1	31	21	4	6	0	0

- Molecule 5 is TETRAETHYLENE GLYCOL (three-letter code: PG4) (formula: C₈H₁₈O₅).



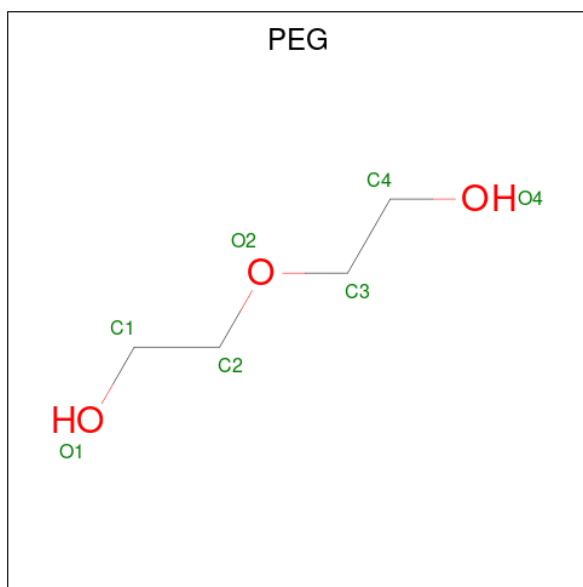
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	1
			26	16	10		
5	B	1	Total	C	O	0	0
			13	8	5		
5	B	1	Total	C	O	0	1
			26	16	10		
5	B	1	Total	C	O	0	0
			13	8	5		

- Molecule 6 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	1	Total C O 4 2 2	0	0
6	B	1	Total C O 4 2 2	0	0

- Molecule 7 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C₄H₁₀O₃).



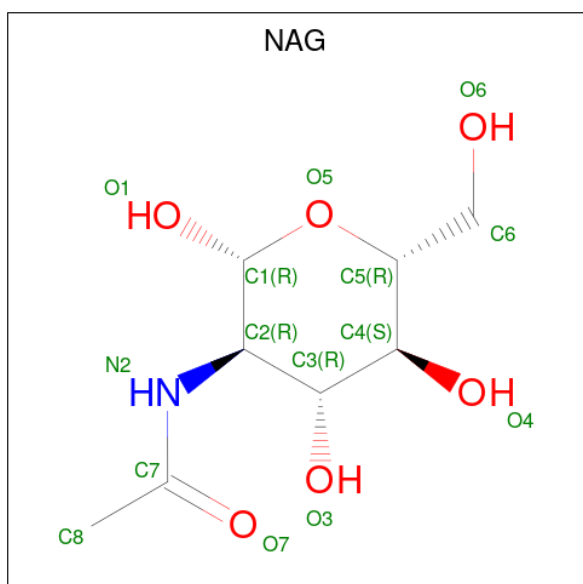
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	1	Total C O 7 4 3	0	0
7	A	1	Total C O 7 4 3	0	0
7	B	1	Total C O 7 4 3	0	0
7	B	1	Total C O 7 4 3	0	0

- Molecule 8 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	1	Total C O 4 2 2	0	0
8	B	1	Total C O 4 2 2	0	0
8	B	1	Total C O 4 2 2	0	0
8	B	1	Total C O 4 2 2	0	0

- Molecule 9 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
9	A	1	Total	C	N	O	0	0
			14	8	1	5		
9	B	1	Total	C	N	O	0	0
			14	8	1	5		
9	B	1	Total	C	N	O	0	0
			14	8	1	5		
9	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 10 is ZINC ION (three-letter code: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	1	Total	Zn	0	0
			1	1		
10	B	1	Total	Zn	0	0
			1	1		

- Molecule 11 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	A	1	Total	Cl	0	0
			1	1		
11	B	1	Total	Cl	0	0
			1	1		

- Molecule 12 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
12	A	1	Total	Mg	0	0
			1	1		
12	B	1	Total	Mg	0	0
			1	1		

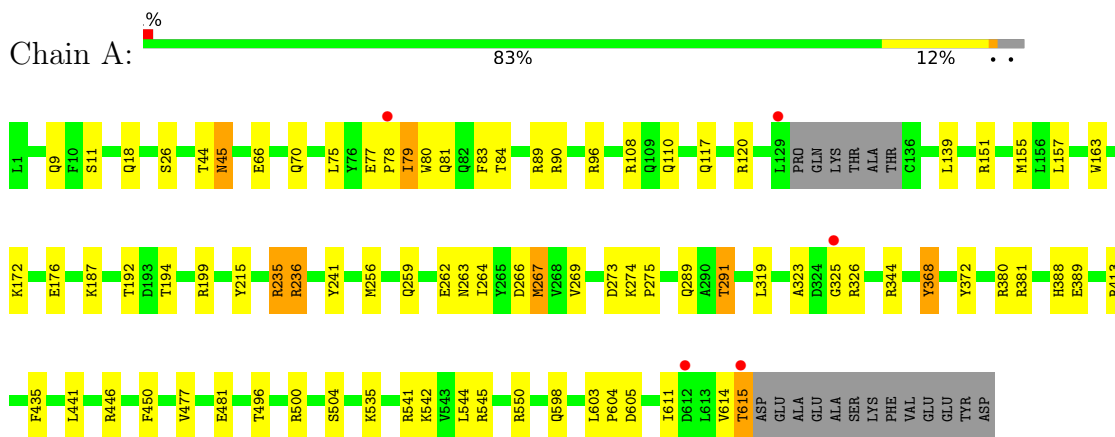
- Molecule 13 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
13	A	519	Total	O	0	0
			519	519		
13	B	345	Total	O	0	0
			345	345		

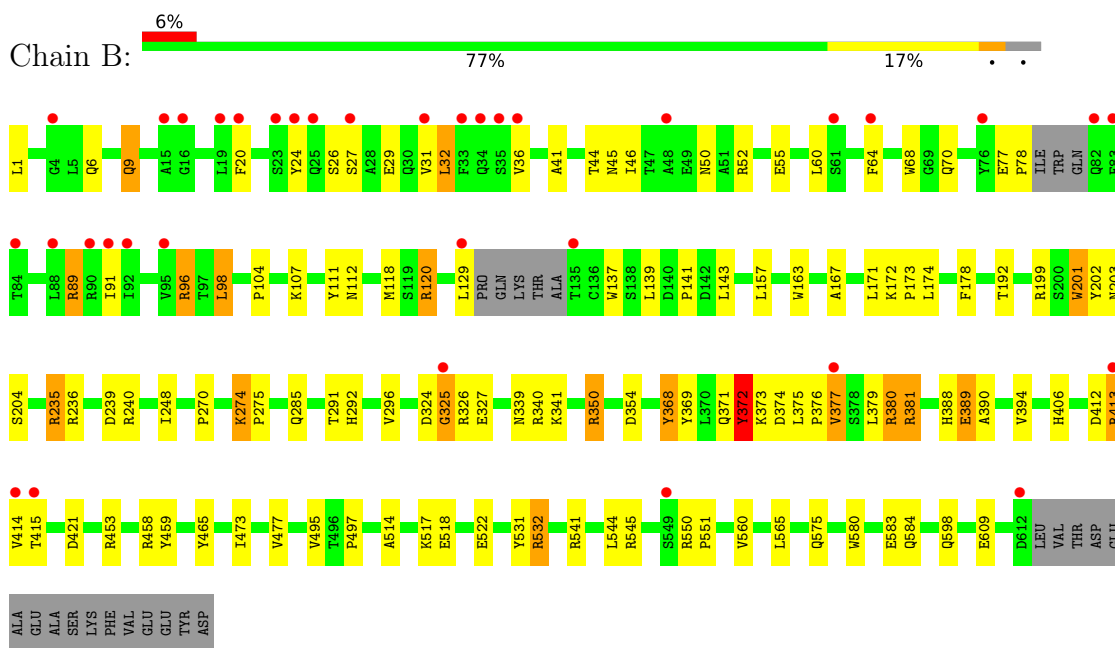
3 Residue-property plots [i](#)

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


- Molecule 1: Angiotensin-converting enzyme, soluble form



- Molecule 1: Angiotensin-converting enzyme, soluble form




- Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain E:  100%

MAG1
MAG2

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

Chain C:  100%

MAG1
MAG2
FUC3

4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	73.09Å 77.97Å 83.31Å 88.27° 64.11° 74.36°	Depositor
Resolution (Å)	74.70 – 2.00 74.70 – 2.00	Depositor EDS
% Data completeness (in resolution range)	98.0 (74.70-2.00) 96.4 (74.70-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.86 (at 2.00Å)	Xtrriage
Refinement program	REFMAC 5.8.0425	Depositor
R, R_{free}	0.183 , 0.230 0.183 , 0.229	Depositor DCC
R_{free} test set	5131 reflections (4.89%)	wwPDB-VP
Wilson B-factor (Å ²)	27.2	Xtrriage
Anisotropy	0.061	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 40.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	11143	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.12% 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 i

5.1 Standard geometry i

Bond lengths and bond angles in the following residue types are not validated in this section: NAG, FUC, ACT, PEG, CL, MG, ZN, PG4, A1IJW, EDO

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	0.71	1/5166 (0.0%)	1.26	37/7035 (0.5%)
1	B	0.66	1/5100 (0.0%)	1.25	34/6943 (0.5%)
All	All	0.69	2/10266 (0.0%)	1.25	71/13978 (0.5%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3
1	B	0	5
All	All	0	8

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	262	GLU	CD-OE1	5.68	1.31	1.25
1	B	389	GLU	CD-OE2	5.43	1.31	1.25

All (71) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	500	ARG	NE-CZ-NH2	-13.63	113.48	120.30
1	A	267	MET	CG-SD-CE	-10.80	82.93	100.20
1	B	44	THR	CA-CB-OG1	-8.84	90.44	109.00
1	A	326	ARG	CB-CA-C	-8.80	92.81	110.40
1	A	263	ASN	CB-CA-C	8.38	127.16	110.40
1	B	350	ARG	NE-CZ-NH2	-8.14	116.23	120.30
1	B	118	MET	CG-SD-CE	8.01	113.02	100.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	413	ARG	NE-CZ-NH1	-7.85	116.38	120.30
1	B	192	THR	CA-CB-OG1	-7.79	92.64	109.00
1	B	235	ARG	NE-CZ-NH1	-7.73	116.43	120.30
1	B	32	LEU	CB-CG-CD2	-7.57	98.12	111.00
1	B	9	GLN	N-CA-CB	7.40	123.92	110.60
1	B	518	GLU	CB-CA-C	7.26	124.92	110.40
1	A	481	GLU	CG-CD-OE2	-7.23	103.83	118.30
1	B	458	ARG	NE-CZ-NH1	7.18	123.89	120.30
1	B	583	GLU	CB-CA-C	-7.13	96.14	110.40
1	B	240	ARG	CD-NE-CZ	6.94	133.32	123.60
1	B	120	ARG	CB-CG-CD	6.79	129.24	111.60
1	B	584	GLN	CB-CA-C	-6.74	96.93	110.40
1	A	18	GLN	N-CA-CB	6.73	122.72	110.60
1	B	532	ARG	NE-CZ-NH2	-6.56	117.02	120.30
1	A	236	ARG	NE-CZ-NH1	-6.52	117.04	120.30
1	A	273	ASP	CB-CA-C	6.45	123.30	110.40
1	B	89	ARG	CA-CB-CG	6.40	127.49	113.40
1	A	155[A]	MET	CG-SD-CE	6.36	110.38	100.20
1	A	155[B]	MET	CG-SD-CE	6.36	110.38	100.20
1	A	615	THR	CA-CB-OG1	-6.29	95.79	109.00
1	A	236	ARG	CD-NE-CZ	6.27	132.38	123.60
1	A	236	ARG	NE-CZ-NH2	6.27	123.43	120.30
1	B	240	ARG	NE-CZ-NH1	6.15	123.38	120.30
1	B	239	ASP	CB-CA-C	-6.09	98.22	110.40
1	A	18	GLN	CB-CA-C	-6.06	98.28	110.40
1	A	381	ARG	CA-CB-CG	6.05	126.71	113.40
1	B	98	LEU	CB-CG-CD2	-6.03	100.75	111.00
1	A	481	GLU	CG-CD-OE1	6.01	130.32	118.30
1	B	174	LEU	CB-CG-CD2	-6.00	100.80	111.00
1	A	215	TYR	CA-CB-CG	-5.99	102.01	113.40
1	A	199	ARG	NE-CZ-NH1	-5.97	117.32	120.30
1	A	291	THR	CA-CB-OG1	-5.92	96.58	109.00
1	A	90	ARG	NE-CZ-NH2	5.89	123.25	120.30
1	B	377	VAL	N-CA-CB	5.85	124.37	111.50
1	B	55	GLU	N-CA-CB	5.83	121.08	110.60
1	A	413	ARG	NE-CZ-NH2	5.82	123.21	120.30
1	B	381	ARG	CB-CA-C	-5.80	98.80	110.40
1	A	603	LEU	CB-CG-CD1	-5.77	101.20	111.00
1	A	256	MET	CG-SD-CE	-5.74	91.01	100.20
1	B	120	ARG	CG-CD-NE	-5.74	99.74	111.80
1	A	90	ARG	CG-CD-NE	-5.63	99.98	111.80
1	B	598	GLN	CB-CA-C	-5.61	99.18	110.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	372	TYR	CA-CB-CG	5.61	124.05	113.40
1	B	354	ASP	CB-CA-C	5.40	121.19	110.40
1	A	192	THR	CA-CB-OG1	-5.39	97.68	109.00
1	B	52	ARG	CD-NE-CZ	5.35	131.09	123.60
1	A	605	ASP	CB-CA-C	-5.33	99.75	110.40
1	A	90	ARG	CD-NE-CZ	5.32	131.05	123.60
1	B	380	ARG	NE-CZ-NH2	-5.30	117.65	120.30
1	A	96	ARG	N-CA-CB	-5.29	101.09	110.60
1	A	550	ARG	NE-CZ-NH2	-5.29	117.66	120.30
1	B	459	TYR	CB-CG-CD2	-5.29	117.83	121.00
1	A	192	THR	N-CA-CB	-5.27	100.28	110.30
1	A	446	ARG	CA-CB-CG	5.25	124.94	113.40
1	B	174	LEU	CB-CG-CD1	5.23	119.89	111.00
1	A	542	LYS	N-CA-CB	5.19	119.94	110.60
1	B	369	TYR	CB-CG-CD2	-5.15	117.91	121.00
1	B	32	LEU	CB-CG-CD1	5.13	119.73	111.00
1	A	235	ARG	CD-NE-CZ	5.13	130.78	123.60
1	B	458	ARG	NE-CZ-NH2	-5.07	117.76	120.30
1	A	319	LEU	CB-CG-CD2	5.07	119.62	111.00
1	A	545	ARG	NE-CZ-NH2	-5.06	117.77	120.30
1	B	199	ARG	CD-NE-CZ	5.06	130.68	123.60
1	A	266	ASP	CB-CA-C	5.03	120.45	110.40

There are no chirality outliers.

All (8) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	235	ARG	Sidechain
1	A	380	ARG	Sidechain
1	A	541	ARG	Sidechain
1	B	120	ARG	Sidechain
1	B	235	ARG	Sidechain
1	B	236	ARG	Sidechain
1	B	326	ARG	Sidechain
1	B	89	ARG	Sidechain

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	5011	0	4789	33	0
1	B	4948	0	4724	65	0
2	E	28	0	25	0	0
3	C	38	0	34	0	0
4	A	31	0	0	0	0
4	B	31	0	0	1	0
5	A	26	0	36	3	0
5	B	52	0	72	10	0
6	A	4	0	3	1	0
6	B	4	0	3	0	0
7	A	14	0	20	4	0
7	B	14	0	20	1	0
8	A	4	0	6	0	0
8	B	12	0	18	5	0
9	A	14	0	13	0	0
9	B	42	0	39	0	0
10	A	1	0	0	0	0
10	B	1	0	0	0	0
11	A	1	0	0	0	0
11	B	1	0	0	0	0
12	A	1	0	0	0	0
12	B	1	0	0	0	0
13	A	519	0	0	9	0
13	B	345	0	0	9	0
All	All	11143	0	9802	111	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 6.

All (111) 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)
5:A:702[B]:PG4:H51	13:A:845:HOH:O	1.63	0.95
5:B:708[B]:PG4:H51	8:B:709:EDO:H22	1.54	0.89
7:A:704:PEG:H12	13:A:1153:HOH:O	1.83	0.79
7:A:704:PEG:C1	13:A:1153:HOH:O	2.32	0.77
1:B:24:TYR:CE1	1:B:64:PHE:HE1	2.06	0.74
1:B:91:ILE:HD11	1:B:376:PRO:CB	2.19	0.73
1:B:465:TYR:CZ	5:B:710:PG4:H42	2.25	0.71
1:A:157:LEU:HD11	1:A:477:VAL:HG13	1.74	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:91:ILE:HD11	1:B:376:PRO:HB2	1.76	0.68
1:A:83:PHE:HB2	1:A:89:ARG:HG2	1.79	0.65
1:B:41:ALA:O	1:B:45:ASN:HB3	1.97	0.65
1:B:31:VAL:HG21	1:B:64:PHE:CD2	2.31	0.64
5:B:708[B]:PG4:H51	8:B:709:EDO:C2	2.28	0.64
1:B:24:TYR:CE1	1:B:64:PHE:CE1	2.87	0.61
1:B:201:TRP:HZ3	1:B:497:PRO:HG2	1.65	0.61
1:B:24:TYR:HE1	1:B:64:PHE:HE1	1.45	0.61
1:A:66:GLU:O	1:A:70:GLN:HG3	2.00	0.61
1:A:264:ILE:HB	1:A:267:MET:HE3	1.83	0.61
4:B:701:A1IJW:C13	5:B:708[A]:PG4:H41	2.33	0.58
1:A:496:THR:OG1	5:A:702[A]:PG4:H71	2.03	0.58
1:B:274:LYS:HB3	1:B:275:PRO:HD2	1.84	0.58
5:B:708[B]:PG4:C5	8:B:709:EDO:H22	2.32	0.58
1:B:324:ASP:O	1:B:325:GLY:C	2.43	0.57
1:A:78:PRO:C	1:A:79:ILE:HG23	2.26	0.56
1:A:110:GLN:NE2	13:A:808:HOH:O	2.39	0.55
1:B:20:PHE:CZ	1:B:68:TRP:HE3	2.23	0.55
1:B:390:ALA:O	1:B:394:VAL:HG23	2.07	0.55
1:B:324:ASP:O	1:B:324:ASP:OD1	2.25	0.54
1:B:1:LEU:HD23	1:B:6:GLN:HG2	1.89	0.54
1:B:406:HIS:NE2	13:B:805:HOH:O	2.33	0.54
1:B:340:ARG:HD3	1:B:373:LYS:O	2.07	0.54
1:B:412:ASP:O	1:B:413:ARG:HB2	2.06	0.54
1:B:201:TRP:CZ3	1:B:497:PRO:HG2	2.43	0.53
1:B:60:LEU:HD23	1:B:60:LEU:O	2.08	0.53
1:A:151:ARG:HD2	1:A:267:MET:CE	2.39	0.53
1:B:550:ARG:HB3	1:B:551:PRO:HD2	1.90	0.52
1:B:292:HIS:NE2	5:B:704:PG4:O2	2.32	0.52
1:B:541:ARG:HD2	1:B:545:ARG:HH11	1.74	0.52
1:B:167:ALA:O	1:B:171:LEU:HD12	2.10	0.51
5:B:708[A]:PG4:H12	13:B:1032:HOH:O	2.09	0.51
1:A:172:LYS:O	1:A:176:GLU:HG3	2.11	0.50
1:A:44:THR:O	1:A:45:ASN:HB2	2.11	0.50
1:A:66:GLU:CD	1:A:108:ARG:HH22	2.16	0.49
1:B:389:GLU:OE2	1:B:389:GLU:HA	2.13	0.49
7:A:704:PEG:H11	13:A:1153:HOH:O	2.08	0.49
1:B:371:GLN:HB3	1:B:545:ARG:HG3	1.94	0.49
1:B:514:ALA:HB2	1:B:560:VAL:HG12	1.95	0.49
1:A:187:LYS:HE2	13:A:1103:HOH:O	2.13	0.49
5:A:702[B]:PG4:H12	5:A:702[B]:PG4:H41	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:29:GLU:HG2	1:B:377:VAL:HG11	1.95	0.48
1:A:80:TRP:CE2	1:A:81:GLN:HG3	2.49	0.48
1:B:96:ARG:O	1:B:98:LEU:HD12	2.13	0.48
1:B:270:PRO:HB3	1:B:580:TRP:CH2	2.48	0.47
1:B:371:GLN:O	1:B:545:ARG:HG2	2.15	0.47
1:B:296:VAL:HG22	7:B:703:PEG:H41	1.96	0.47
1:A:84:THR:HA	13:A:1001:HOH:O	2.14	0.47
1:A:78:PRO:C	1:A:79:ILE:CG2	2.82	0.47
1:A:77:GLU:HA	1:A:77:GLU:OE1	2.15	0.47
1:A:344:ARG:NH2	7:A:704:PEG:H32	2.30	0.46
1:B:514:ALA:CB	1:B:560:VAL:HG12	2.45	0.46
1:A:323:ALA:C	1:A:325:GLY:H	2.18	0.46
1:B:495:VAL:HG12	1:B:495:VAL:O	2.16	0.46
1:B:46:ILE:HB	1:B:327:GLU:HB2	1.97	0.45
1:B:522:GLU:N	13:B:817:HOH:O	2.47	0.45
1:B:172:LYS:HB3	1:B:173:PRO:HD3	1.99	0.45
1:B:202:TYR:O	1:B:203:ASN:C	2.54	0.45
1:A:241:TYR:OH	1:A:604:PRO:HA	2.16	0.45
1:A:139:LEU:HD22	1:A:163:TRP:CZ2	2.52	0.45
1:B:379:LEU:HD23	1:B:379:LEU:HA	1.58	0.45
1:A:289:GLN:NE2	13:A:837:HOH:O	2.50	0.45
1:B:157:LEU:HD11	1:B:477:VAL:HG13	1.99	0.44
1:B:374:ASP:HB2	13:B:803:HOH:O	2.17	0.44
1:B:141:PRO:HB3	1:B:350:ARG:HD3	2.00	0.44
1:B:453:ARG:HD3	13:B:1092:HOH:O	2.18	0.44
1:A:9:GLN:NE2	1:A:75:LEU:HD21	2.32	0.44
1:A:389:GLU:HB2	1:A:504[A]:SER:HB2	1.99	0.44
1:A:598:GLN:HB2	13:B:980:HOH:O	2.16	0.43
6:A:703:ACT:H1	13:A:1000:HOH:O	2.17	0.43
1:B:565:LEU:HD23	1:B:565:LEU:HA	1.81	0.43
1:B:129:LEU:HD21	1:B:137:TRP:CZ2	2.53	0.43
5:B:708[A]:PG4:H21	8:B:709:EDO:O2	2.18	0.43
1:A:259:GLN:O	1:A:435:PHE:HA	2.19	0.43
1:B:291:THR:HG22	13:B:1036:HOH:O	2.18	0.43
1:B:372:TYR:CE1	1:B:380:ARG:HB3	2.52	0.43
1:B:368:TYR:CE1	1:B:544:LEU:HA	2.54	0.43
1:B:375:LEU:HB3	1:B:376:PRO:HD2	2.01	0.43
1:B:541:ARG:HG2	1:B:545:ARG:HD2	2.01	0.42
1:A:274:LYS:HB3	1:A:275:PRO:CD	2.49	0.42
1:A:368:TYR:CE1	1:A:544:LEU:HA	2.54	0.42
1:A:151:ARG:HD2	1:A:267:MET:HE2	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:441:LEU:HD12	1:A:441:LEU:C	2.40	0.42
1:B:248:ILE:O	1:B:473:ILE:HA	2.19	0.42
1:B:139:LEU:HA	1:B:143:LEU:HB2	2.02	0.42
1:A:194:THR:HB	1:A:450:PHE:CD1	2.55	0.41
1:B:111:TYR:HD1	1:B:112:ASN:ND2	2.17	0.41
1:B:104:PRO:HG2	1:B:107:LYS:HD2	2.01	0.41
1:B:531:TYR:O	1:B:532:ARG:HB2	2.20	0.41
1:A:611:ILE:HD13	1:A:611:ILE:HA	1.77	0.41
1:B:204:SER:O	8:B:705:EDO:H21	2.21	0.41
1:A:117[A]:GLN:NE2	1:A:120:ARG:HE	2.19	0.41
1:B:45:ASN:ND2	1:B:50:ASN:HD21	2.18	0.41
1:B:292:HIS:NE2	5:B:704:PG4:H11	2.36	0.41
1:B:32:LEU:O	1:B:36:VAL:HG23	2.21	0.40
1:A:236:ARG:HD2	1:A:267:MET:SD	2.62	0.40
1:B:27:SER:HB3	1:B:64:PHE:HE2	1.87	0.40
1:B:339:ASN:O	1:B:340:ARG:HB2	2.21	0.40
1:B:77:GLU:O	1:B:78:PRO:C	2.60	0.40
1:B:107:LYS:HE2	13:B:1073:HOH:O	2.21	0.40
1:B:374:ASP:CB	13:B:803:HOH:O	2.68	0.40
1:B:139:LEU:HD22	1:B:163:TRP:CZ2	2.56	0.40
5:B:704:PG4:H61	5:B:704:PG4:H81	1.67	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	610/628 (97%)	598 (98%)	10 (2%)	2 (0%)	37	35
1	B	601/628 (96%)	579 (96%)	17 (3%)	5 (1%)	16	12
All	All	1211/1256 (96%)	1177 (97%)	27 (2%)	7 (1%)	22	17

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	413	ARG
1	B	325	GLY
1	B	201	TRP
1	B	274	LYS
1	B	575	GLN
1	A	45	ASN
1	A	79	ILE

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	529/540 (98%)	519 (98%)	10 (2%)	52 57
1	B	522/540 (97%)	506 (97%)	16 (3%)	35 36
All	All	1051/1080 (97%)	1025 (98%)	26 (2%)	42 45

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	SER
1	A	26	SER
1	A	269	VAL
1	A	291	THR
1	A	368	TYR
1	A	372	TYR
1	A	388	HIS
1	A	535	LYS
1	A	614	VAL
1	A	615	THR
1	B	9	GLN
1	B	26	SER
1	B	70	GLN
1	B	96	ARG
1	B	178	PHE
1	B	285	GLN

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Mol	Chain	Res	Type
1	B	341	LYS
1	B	368	TYR
1	B	372	TYR
1	B	381	ARG
1	B	388	HIS
1	B	414	VAL
1	B	415	THR
1	B	421	ASP
1	B	517	LYS
1	B	609	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	18	GLN
1	A	110	GLN
1	A	263	ASN
1	B	25	GLN
1	B	30	GLN
1	B	586	GLN

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](#)

5 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	3,1	14,14,15	0.68	0	17,19,21	1.30	2 (11%)
3	NAG	C	2	3	14,14,15	0.44	0	17,19,21	1.30	3 (17%)
3	FUC	C	3	3	10,10,11	0.63	0	14,14,16	1.18	2 (14%)
2	NAG	E	1	2,1	14,14,15	0.43	0	17,19,21	1.46	2 (11%)
2	NAG	E	2	2	14,14,15	0.37	0	17,19,21	1.18	2 (11%)

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	3,1	-	0/6/23/26	0/1/1/1
3	NAG	C	2	3	-	2/6/23/26	0/1/1/1
3	FUC	C	3	3	-	-	0/1/1/1
2	NAG	E	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	E	2	2	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	1	NAG	C2-N2-C7	-3.73	117.59	122.90
2	E	1	NAG	C1-O5-C5	3.38	116.77	112.19
3	C	2	NAG	C1-C2-N2	3.03	115.66	110.49
3	C	1	NAG	C1-C2-N2	-2.94	105.46	110.49
3	C	3	FUC	O5-C1-C2	-2.84	106.39	110.77
3	C	2	NAG	O3-C3-C2	-2.60	104.08	109.47
2	E	2	NAG	C2-N2-C7	2.42	126.35	122.90
3	C	2	NAG	C1-O5-C5	2.37	115.41	112.19
3	C	3	FUC	O2-C2-C3	2.33	114.80	110.14
3	C	1	NAG	O3-C3-C2	2.24	114.10	109.47
2	E	2	NAG	C1-C2-N2	2.20	114.25	110.49

There are no chirality outliers.

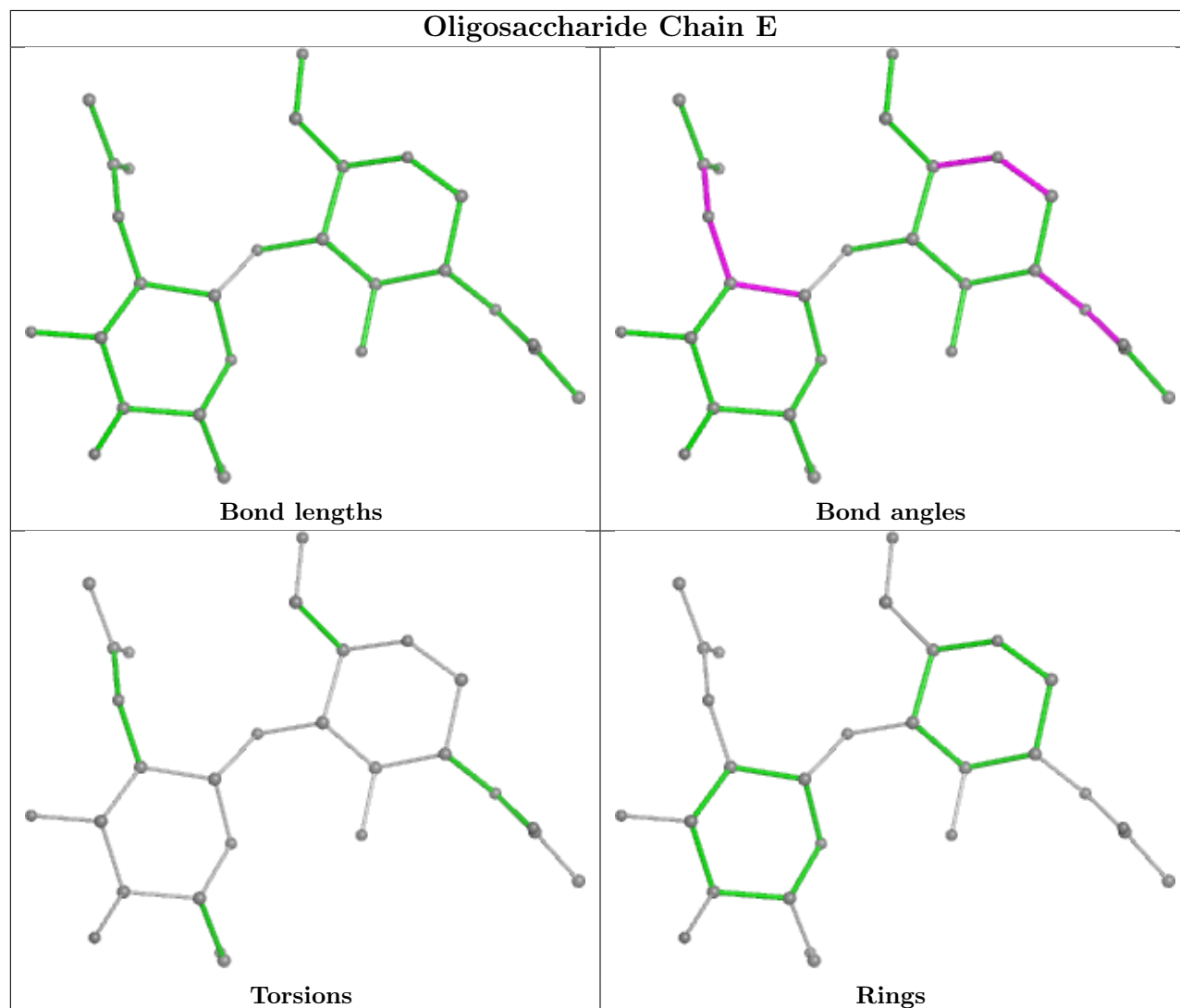
All (2) torsion outliers are listed below:

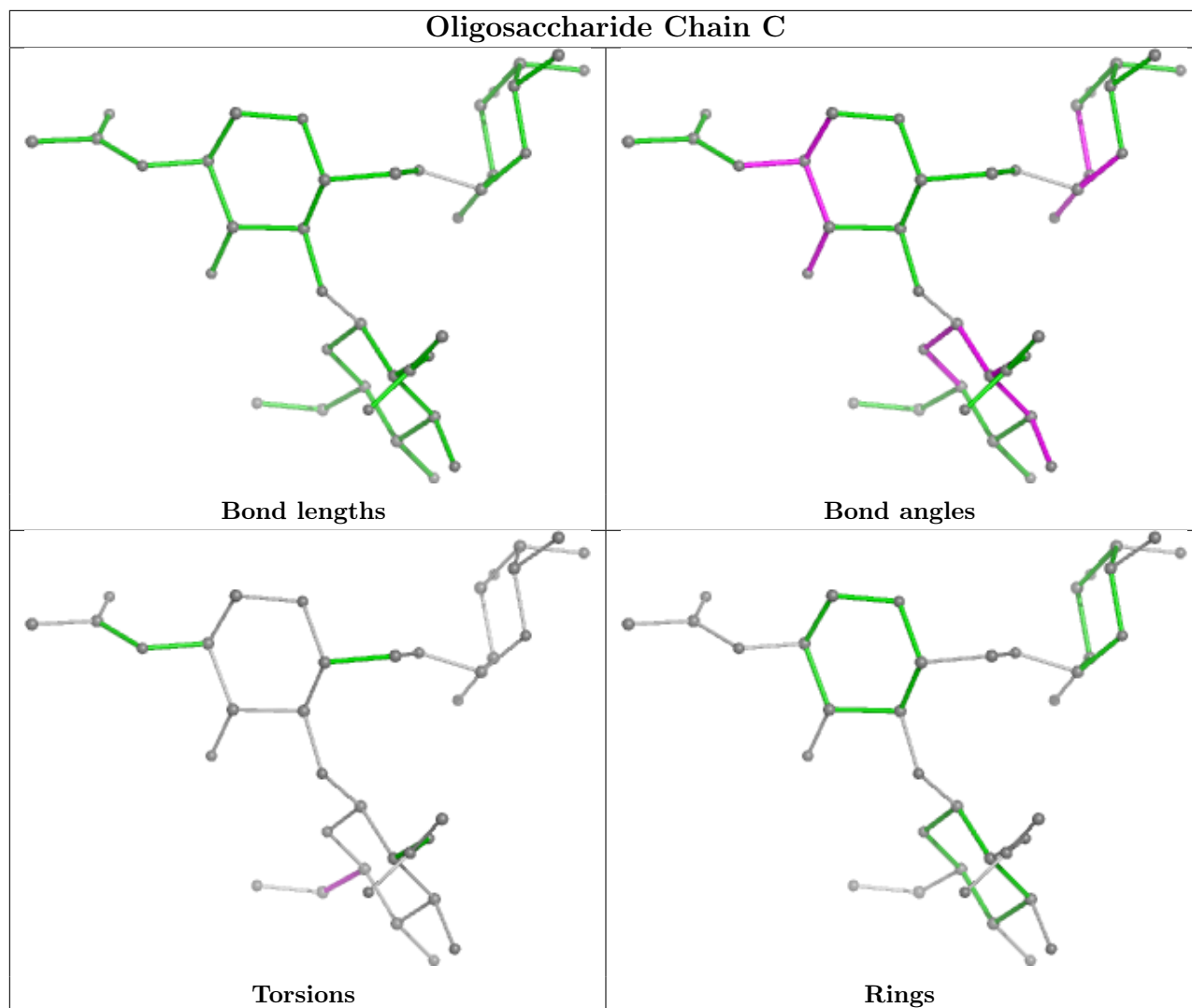
Mol	Chain	Res	Type	Atoms
3	C	2	NAG	O5-C5-C6-O6
3	C	2	NAG	C4-C5-C6-O6

There are no ring outliers.

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 [i](#)

Of 28 ligands modelled in this entry, 6 are monoatomic - leaving 22 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	A1IJW	B	701	10	33,33,33	3.45	14 (42%)	40,47,47	2.36	15 (37%)
6	ACT	A	703	-	3,3,3	1.26	0	3,3,3	0.66	0
9	NAG	B	715	1	14,14,15	0.47	0	17,19,21	1.64	3 (17%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
9	NAG	A	707	1	14,14,15	0.64	0	17,19,21	1.84	2 (11%)
7	PEG	A	706	-	6,6,6	0.28	0	5,5,5	0.24	0
8	EDO	B	706	-	3,3,3	0.18	0	2,2,2	0.26	0
8	EDO	B	709	-	3,3,3	0.06	0	2,2,2	0.37	0
5	PG4	B	708[A]	-	12,12,12	0.15	0	11,11,11	0.30	0
5	PG4	A	702[A]	-	12,12,12	0.31	0	11,11,11	0.21	0
7	PEG	B	702	-	6,6,6	0.28	0	5,5,5	0.40	0
7	PEG	B	703	-	6,6,6	0.30	0	5,5,5	0.19	0
6	ACT	B	707	-	3,3,3	1.02	0	3,3,3	1.20	0
5	PG4	B	708[B]	-	12,12,12	0.18	0	11,11,11	0.34	0
5	PG4	A	702[B]	-	12,12,12	0.23	0	11,11,11	0.37	0
5	PG4	B	704	-	12,12,12	0.44	0	11,11,11	0.31	0
5	PG4	B	710	-	12,12,12	0.43	0	11,11,11	0.35	0
8	EDO	B	705	-	3,3,3	0.70	0	2,2,2	1.11	0
4	A1IJW	A	701	10	33,33,33	2.82	15 (45%)	40,47,47	2.56	17 (42%)
7	PEG	A	704	-	6,6,6	0.46	0	5,5,5	0.38	0
8	EDO	A	705	-	3,3,3	0.71	0	2,2,2	0.94	0
9	NAG	B	714	1	14,14,15	0.61	0	17,19,21	2.46	4 (23%)
9	NAG	B	713	1	14,14,15	0.75	0	17,19,21	3.55	6 (35%)

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	A1IJW	B	701	10	-	3/32/52/52	0/3/3/3
9	NAG	B	715	1	-	2/6/23/26	0/1/1/1
9	NAG	A	707	1	-	3/6/23/26	0/1/1/1
7	PEG	A	706	-	-	3/4/4/4	-
8	EDO	B	706	-	-	1/1/1/1	-
8	EDO	B	709	-	-	1/1/1/1	-
5	PG4	B	708[A]	-	-	5/10/10/10	-
5	PG4	A	702[A]	-	-	3/10/10/10	-
7	PEG	B	702	-	-	3/4/4/4	-
7	PEG	B	703	-	-	4/4/4/4	-
5	PG4	B	708[B]	-	-	8/10/10/10	-
5	PG4	A	702[B]	-	-	7/10/10/10	-
5	PG4	B	704	-	-	5/10/10/10	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	PG4	B	710	-	-	5/10/10/10	-
8	EDO	B	705	-	-	0/1/1/1	-
4	A1IJW	A	701	10	-	5/32/52/52	0/3/3/3
7	PEG	A	704	-	-	2/4/4/4	-
8	EDO	A	705	-	-	0/1/1/1	-
9	NAG	B	714	1	-	4/6/23/26	0/1/1/1
9	NAG	B	713	1	-	2/6/23/26	0/1/1/1

All (29) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	701	A1IJW	C06-N10	-11.84	1.31	1.47
4	A	701	A1IJW	C06-N10	-8.65	1.36	1.47
4	B	701	A1IJW	C09-N10	7.08	1.61	1.47
4	A	701	A1IJW	C02-N24	5.99	1.48	1.34
4	B	701	A1IJW	C11-N10	5.39	1.47	1.34
4	B	701	A1IJW	C05-C06	5.28	1.60	1.54
4	A	701	A1IJW	C25-N24	5.25	1.57	1.47
4	A	701	A1IJW	C09-N10	4.98	1.57	1.47
4	B	701	A1IJW	C27-C28	-4.53	1.42	1.53
4	B	701	A1IJW	C02-N24	4.46	1.44	1.34
4	B	701	A1IJW	C25-N24	4.00	1.55	1.47
4	B	701	A1IJW	C26-C27	3.71	1.67	1.51
4	A	701	A1IJW	C26-C27	3.65	1.67	1.51
4	B	701	A1IJW	C05-N04	3.36	1.50	1.46
4	A	701	A1IJW	C27-C28	-3.06	1.46	1.53
4	A	701	A1IJW	C26-C25	-3.06	1.41	1.51
4	A	701	A1IJW	C28-N24	3.00	1.53	1.47
4	B	701	A1IJW	C28-N24	2.97	1.53	1.47
4	A	701	A1IJW	C07-C06	2.86	1.58	1.53
4	B	701	A1IJW	C07-C06	2.76	1.58	1.53
4	A	701	A1IJW	C11-N10	2.64	1.40	1.34
4	B	701	A1IJW	C26-C25	-2.62	1.42	1.51
4	A	701	A1IJW	C05-C06	2.48	1.57	1.54
4	A	701	A1IJW	C15-N16	2.44	1.46	1.38
4	A	701	A1IJW	C08-C09	-2.20	1.44	1.51
4	B	701	A1IJW	O19-C11	-2.12	1.18	1.22
4	B	701	A1IJW	O01-C02	-2.12	1.18	1.22
4	A	701	A1IJW	C05-N04	2.08	1.48	1.46
4	A	701	A1IJW	C08-C07	2.04	1.60	1.51

All (47) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	B	713	NAG	C2-N2-C7	-11.43	106.62	122.90
9	B	714	NAG	C1-C2-N2	6.81	122.13	110.49
4	B	701	A1IJW	O21-C20-O22	-5.74	111.06	124.09
4	A	701	A1IJW	O21-C20-O22	-5.70	111.16	124.09
9	B	714	NAG	C2-N2-C7	5.69	131.01	122.90
9	B	713	NAG	C1-O5-C5	5.55	119.71	112.19
4	B	701	A1IJW	O19-C11-N10	-5.23	113.12	121.59
9	A	707	NAG	C1-C2-N2	5.07	119.14	110.49
4	A	701	A1IJW	C18-C17-C15	-4.92	114.34	120.67
4	A	701	A1IJW	C14-C13-C12	-4.87	115.11	120.78
4	A	701	A1IJW	O19-C11-N10	-4.43	114.42	121.59
4	A	701	A1IJW	C07-C06-N10	-4.42	98.99	102.64
4	B	701	A1IJW	C18-C17-C15	-4.22	115.25	120.67
4	B	701	A1IJW	O21-C20-C05	4.17	128.93	114.21
4	A	701	A1IJW	C29-C28-N24	-4.15	103.84	112.26
4	A	701	A1IJW	C25-N24-C28	-4.10	105.50	112.00
4	A	701	A1IJW	O21-C20-C05	3.95	128.14	114.21
4	B	701	A1IJW	C08-C09-N10	-3.94	96.33	103.25
9	B	715	NAG	C1-C2-N2	-3.64	104.26	110.49
9	A	707	NAG	C1-O5-C5	3.63	117.11	112.19
4	A	701	A1IJW	C27-C28-N24	3.41	108.09	103.03
9	B	715	NAG	C2-N2-C7	-3.36	118.12	122.90
4	A	701	A1IJW	C08-C07-C06	-3.29	98.62	104.16
4	B	701	A1IJW	C08-C07-C06	-3.28	98.65	104.16
9	B	714	NAG	C1-O5-C5	3.23	116.56	112.19
9	B	713	NAG	C4-C3-C2	3.18	115.67	111.02
4	A	701	A1IJW	O31-C29-O30	-3.17	116.88	124.09
9	B	715	NAG	C4-C3-C2	3.11	115.58	111.02
4	B	701	A1IJW	C07-C08-C09	3.03	113.57	104.98
9	B	714	NAG	O4-C4-C5	2.94	116.61	109.30
4	B	701	A1IJW	C14-C15-N16	-2.85	115.60	120.91
4	A	701	A1IJW	O31-C29-C28	2.83	123.19	113.38
4	B	701	A1IJW	C14-C13-C12	-2.81	117.51	120.78
4	B	701	A1IJW	C29-C28-N24	-2.70	106.79	112.26
9	B	713	NAG	O7-C7-N2	-2.65	117.07	121.95
9	B	713	NAG	O5-C1-C2	2.64	115.46	111.29
9	B	713	NAG	O3-C3-C2	-2.57	104.15	109.47
4	B	701	A1IJW	C17-C18-C12	2.55	123.74	120.78
4	B	701	A1IJW	O19-C11-C12	2.49	125.07	120.23
4	B	701	A1IJW	O01-C02-N24	-2.46	116.99	121.38
4	B	701	A1IJW	C17-C15-C14	2.45	121.95	118.15
4	A	701	A1IJW	C17-C15-C14	2.43	121.93	118.15

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	701	A1IJW	C18-C12-C13	2.38	121.97	118.59
4	A	701	A1IJW	C18-C12-C11	-2.30	114.46	120.29
4	B	701	A1IJW	C23-C03-C02	-2.17	105.80	109.73
4	A	701	A1IJW	C12-C11-N10	2.08	122.90	118.77
4	A	701	A1IJW	C13-C14-C15	2.04	123.30	120.67

There are no chirality outliers.

All (66) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
9	B	714	NAG	C1-C2-N2-C7
9	B	714	NAG	C8-C7-N2-C2
9	B	714	NAG	O7-C7-N2-C2
5	B	704	PG4	C8-C7-O4-C6
5	B	708[A]	PG4	C8-C7-O4-C6
7	A	704	PEG	C4-C3-O2-C2
5	B	710	PG4	O4-C7-C8-O5
9	B	715	NAG	C8-C7-N2-C2
9	B	715	NAG	O7-C7-N2-C2
5	B	704	PG4	O3-C5-C6-O4
5	B	710	PG4	O2-C3-C4-O3
9	A	707	NAG	O5-C5-C6-O6
5	B	710	PG4	O3-C5-C6-O4
5	B	708[B]	PG4	O1-C1-C2-O2
5	B	708[A]	PG4	O2-C3-C4-O3
7	A	706	PEG	O1-C1-C2-O2
5	B	708[A]	PG4	O3-C5-C6-O4
5	A	702[A]	PG4	C5-C6-O4-C7
9	B	713	NAG	O5-C5-C6-O6
9	B	713	NAG	C4-C5-C6-O6
5	A	702[A]	PG4	O1-C1-C2-O2
5	B	704	PG4	O1-C1-C2-O2
7	B	703	PEG	O1-C1-C2-O2
9	B	714	NAG	O5-C5-C6-O6
5	A	702[B]	PG4	O2-C3-C4-O3
5	B	708[A]	PG4	C4-C3-O2-C2
5	B	708[A]	PG4	O1-C1-C2-O2
7	B	702	PEG	O1-C1-C2-O2
7	B	703	PEG	O2-C3-C4-O4
8	B	709	EDO	O1-C1-C2-O2
9	A	707	NAG	C4-C5-C6-O6
5	B	704	PG4	O2-C3-C4-O3

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Mol	Chain	Res	Type	Atoms
5	A	702[B]	PG4	O4-C7-C8-O5
7	A	706	PEG	O2-C3-C4-O4
5	B	708[B]	PG4	O2-C3-C4-O3
5	A	702[A]	PG4	O3-C5-C6-O4
4	A	701	A1IJW	C20-C05-N04-C03
4	B	701	A1IJW	C20-C05-N04-C03
5	B	708[B]	PG4	C6-C5-O3-C4
4	A	701	A1IJW	N04-C05-C06-C07
4	B	701	A1IJW	N04-C05-C06-C07
5	B	708[B]	PG4	C1-C2-O2-C3
7	A	706	PEG	C4-C3-O2-C2
5	B	704	PG4	C5-C6-O4-C7
4	A	701	A1IJW	C20-C05-C06-C07
4	B	701	A1IJW	C06-C05-N04-C03
8	B	706	EDO	O1-C1-C2-O2
5	A	702[B]	PG4	C8-C7-O4-C6
5	A	702[B]	PG4	C1-C2-O2-C3
7	B	703	PEG	C4-C3-O2-C2
7	B	703	PEG	C1-C2-O2-C3
5	A	702[B]	PG4	O1-C1-C2-O2
5	B	708[B]	PG4	O4-C7-C8-O5
9	A	707	NAG	C3-C2-N2-C7
4	A	701	A1IJW	N04-C05-C20-O21
5	A	702[B]	PG4	O3-C5-C6-O4
5	B	708[B]	PG4	C8-C7-O4-C6
5	B	710	PG4	C1-C2-O2-C3
5	A	702[B]	PG4	C5-C6-O4-C7
5	B	710	PG4	C5-C6-O4-C7
5	B	708[B]	PG4	C4-C3-O2-C2
5	B	708[B]	PG4	C5-C6-O4-C7
7	A	704	PEG	O2-C3-C4-O4
7	B	702	PEG	O2-C3-C4-O4
4	A	701	A1IJW	C20-C05-C06-N10
7	B	702	PEG	C1-C2-O2-C3

There are no ring outliers.

12 monomers are involved in 20 short contacts:

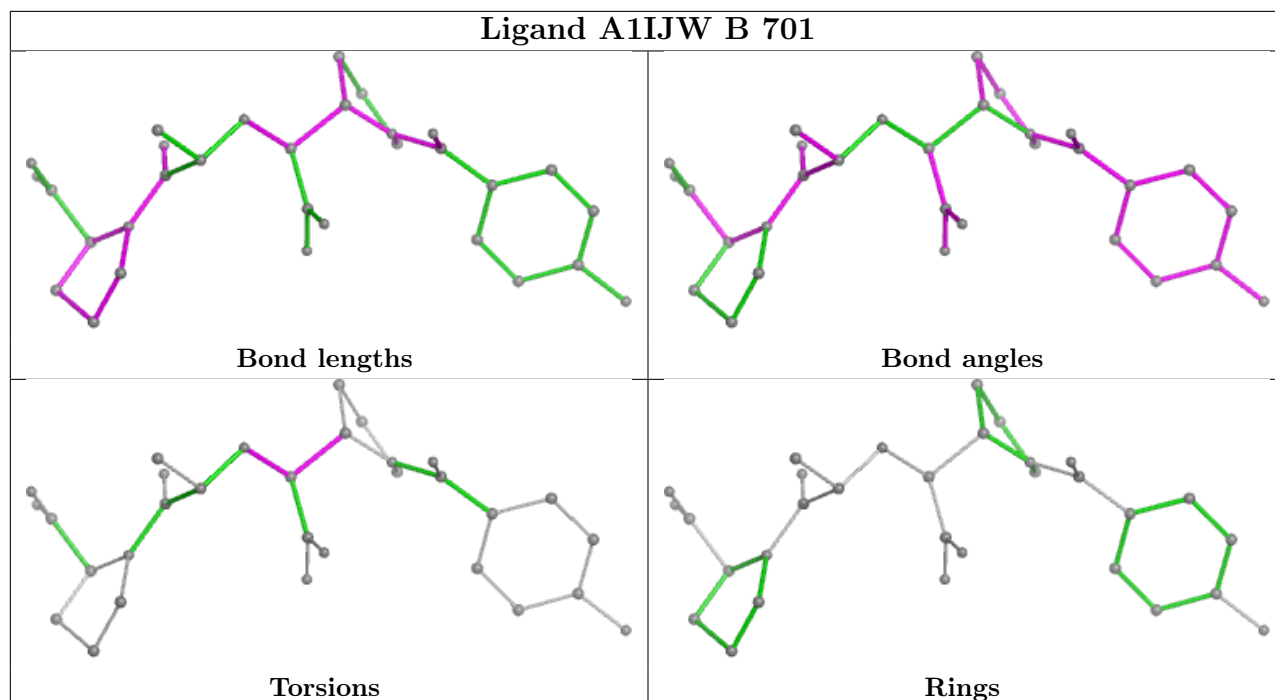
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	701	A1IJW	1	0
6	A	703	ACT	1	0
8	B	709	EDO	4	0

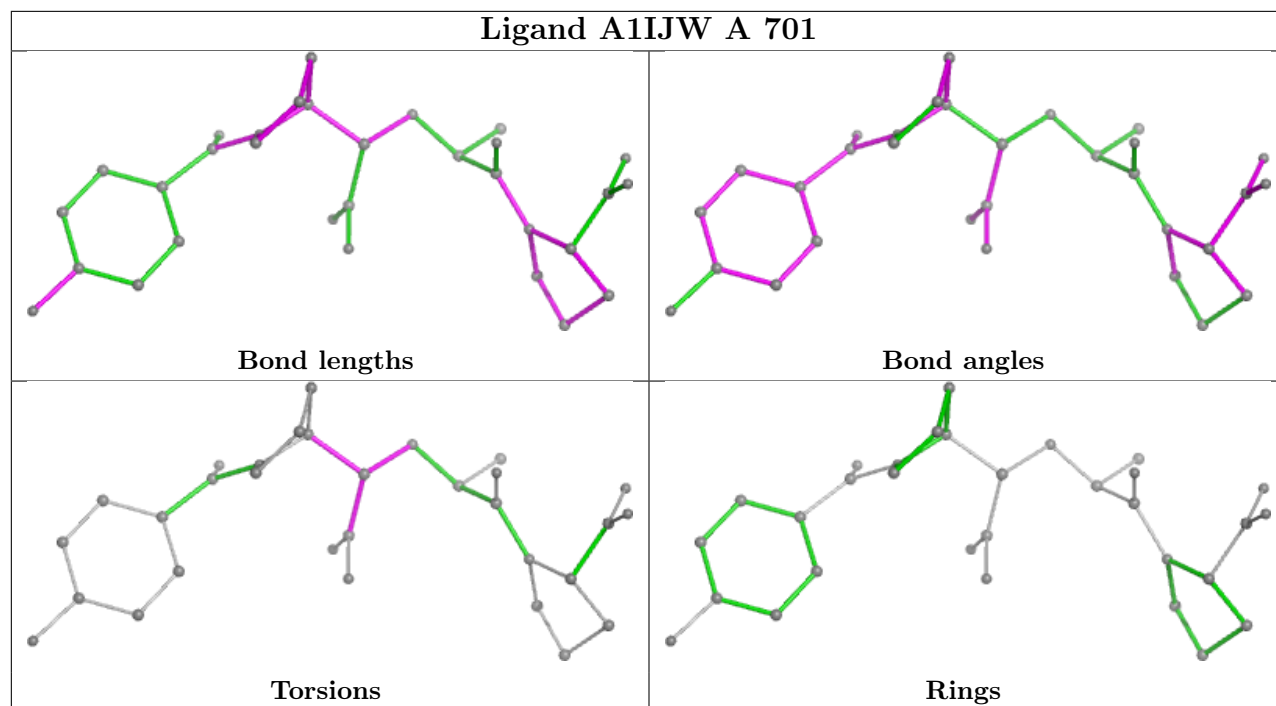
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	708[A]	PG4	3	0
5	A	702[A]	PG4	1	0
7	B	703	PEG	1	0
5	B	708[B]	PG4	3	0
5	A	702[B]	PG4	2	0
5	B	704	PG4	3	0
5	B	710	PG4	1	0
8	B	705	EDO	1	0
7	A	704	PEG	4	0

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	609/628 (96%)	-0.60	5 (0%) 82 82	7, 24, 44, 75	5 (0%)
1	B	604/628 (96%)	0.08	35 (5%) 30 28	9, 33, 62, 99	3 (0%)
All	All	1213/1256 (96%)	-0.26	40 (3%) 49 47	7, 28, 56, 99	8 (0%)

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	83	PHE	6.2
1	A	129	LEU	5.5
1	B	88	LEU	5.1
1	B	414	VAL	5.1
1	B	91	ILE	4.5
1	B	95	VAL	3.6
1	A	615	THR	3.6
1	A	325	GLY	3.3
1	B	24	TYR	3.3
1	B	16	GLY	3.1
1	B	23	SER	3.1
1	B	325	GLY	3.0
1	B	25	GLN	2.9
1	B	82	GLN	2.9
1	B	19	LEU	2.9
1	B	20	PHE	2.9
1	B	64	PHE	2.8
1	B	31	VAL	2.7
1	B	61	SER	2.7
1	B	413	ARG	2.6
1	B	48	ALA	2.5
1	B	129	LEU	2.5
1	B	84	THR	2.5
1	B	76	TYR	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	549	SER	2.4
1	B	35	SER	2.4
1	B	27	SER	2.4
1	B	15	ALA	2.3
1	B	377	VAL	2.3
1	B	612	ASP	2.3
1	B	92	ILE	2.3
1	B	90	ARG	2.3
1	A	78	PRO	2.2
1	B	33	PHE	2.2
1	B	34	GLN	2.2
1	B	135	THR	2.2
1	A	612	ASP	2.1
1	B	415	THR	2.1
1	B	4	GLY	2.1
1	B	36	VAL	2.1

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](#)

SUGAR-RSR INFOmissingINFO

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
8	EDO	B	705	4/4	0.75	0.18	45,45,55,56	0
9	NAG	B	714	14/15	0.76	0.12	59,63,67,67	0
9	NAG	B	715	14/15	0.76	0.17	49,87,103,108	0
8	EDO	B	706	4/4	0.81	0.23	69,71,74,81	0
9	NAG	B	713	14/15	0.82	0.13	30,38,50,53	0
5	PG4	A	702[A]	13/13	0.82	0.19	22,47,52,56	13
5	PG4	A	702[B]	13/13	0.82	0.19	21,26,38,41	13

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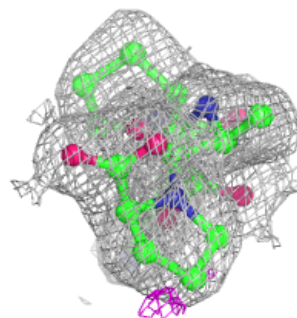
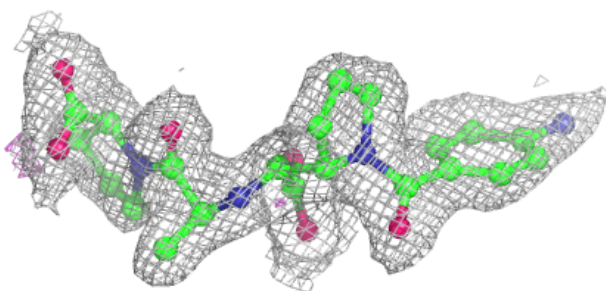
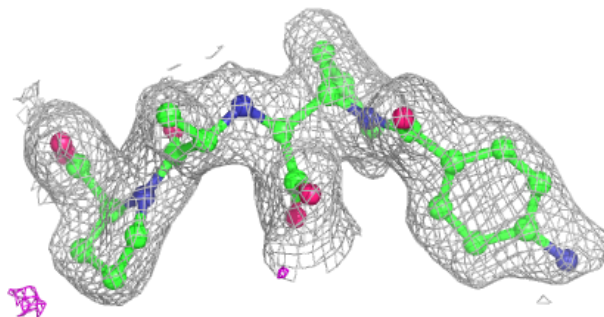
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
8	EDO	A	705	4/4	0.84	0.15	46,47,51,52	0
5	PG4	B	704	13/13	0.86	0.13	44,54,66,68	0
5	PG4	B	710	13/13	0.86	0.15	44,58,65,66	0
5	PG4	B	708[B]	13/13	0.87	0.18	23,39,47,49	13
5	PG4	B	708[A]	13/13	0.87	0.18	18,42,48,56	13
7	PEG	A	704	7/7	0.87	0.13	48,55,57,72	0
9	NAG	A	707	14/15	0.89	0.10	31,35,44,55	0
8	EDO	B	709	4/4	0.89	0.14	46,47,56,60	0
7	PEG	B	702	7/7	0.90	0.15	40,46,50,52	0
7	PEG	A	706	7/7	0.91	0.12	47,51,53,56	0
7	PEG	B	703	7/7	0.91	0.11	38,51,58,65	0
6	ACT	A	703	4/4	0.93	0.12	19,31,35,37	0
6	ACT	B	707	4/4	0.95	0.11	29,37,40,44	0
4	A1IJW	B	701	31/31	0.97	0.06	17,21,36,39	0
12	MG	A	710	1/1	0.97	0.05	25,25,25,25	0
10	ZN	A	708	1/1	0.98	0.03	16,16,16,16	0
11	CL	B	712	1/1	0.98	0.04	29,29,29,29	0
4	A1IJW	A	701	31/31	0.98	0.05	15,17,25,32	0
12	MG	B	716	1/1	0.98	0.03	26,26,26,26	0
10	ZN	B	711	1/1	0.99	0.02	20,20,20,20	0
11	CL	A	709	1/1	1.00	0.05	19,19,19,19	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.

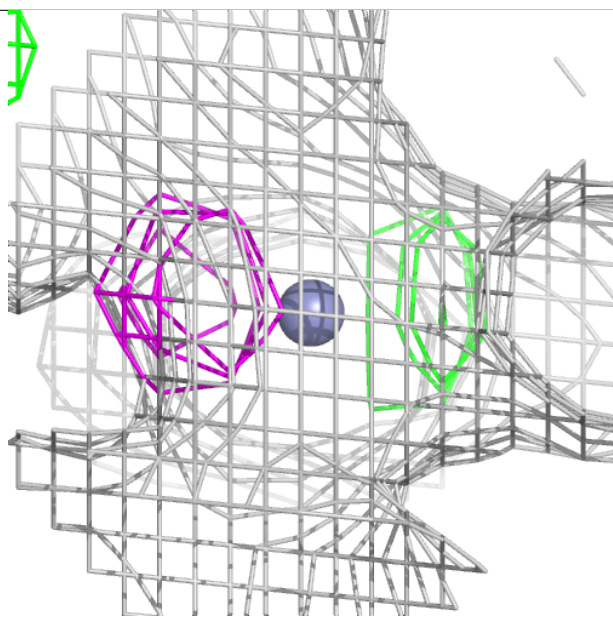
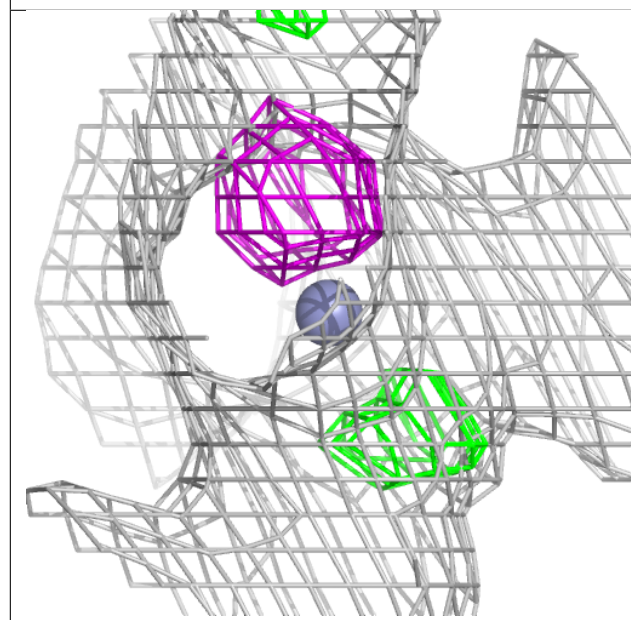
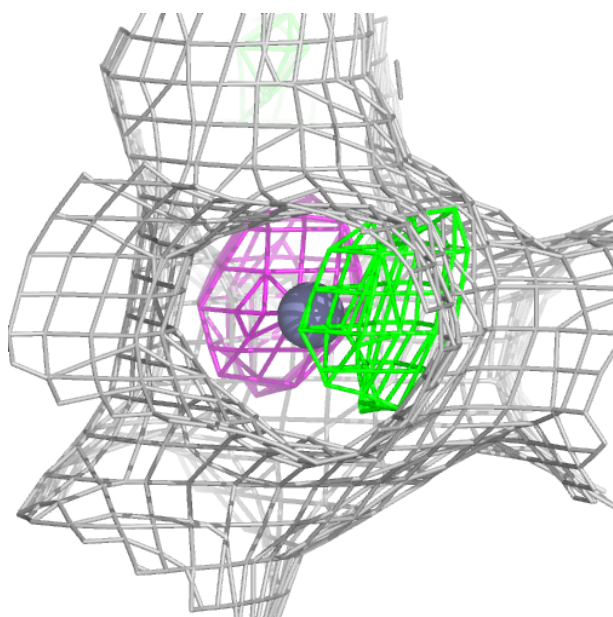
Electron density around A1IJW B 701:

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



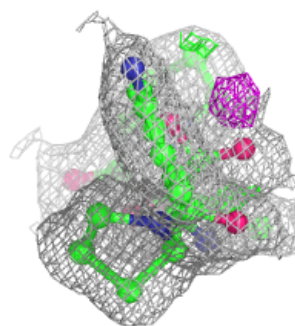
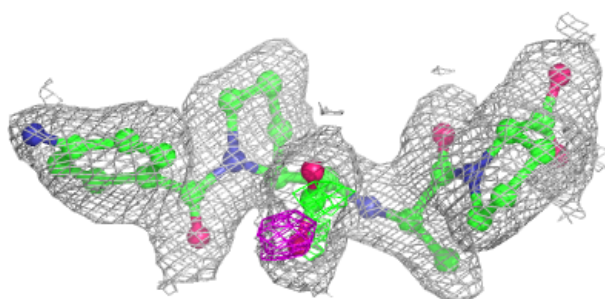
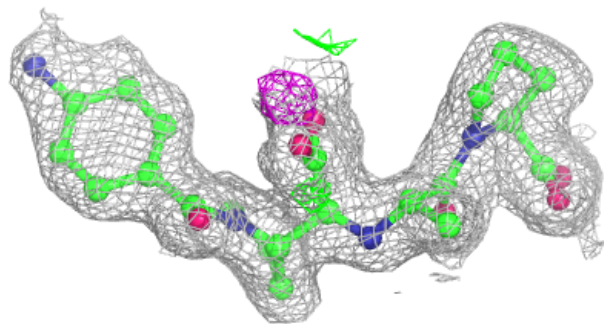
Electron density around ZN A 708:

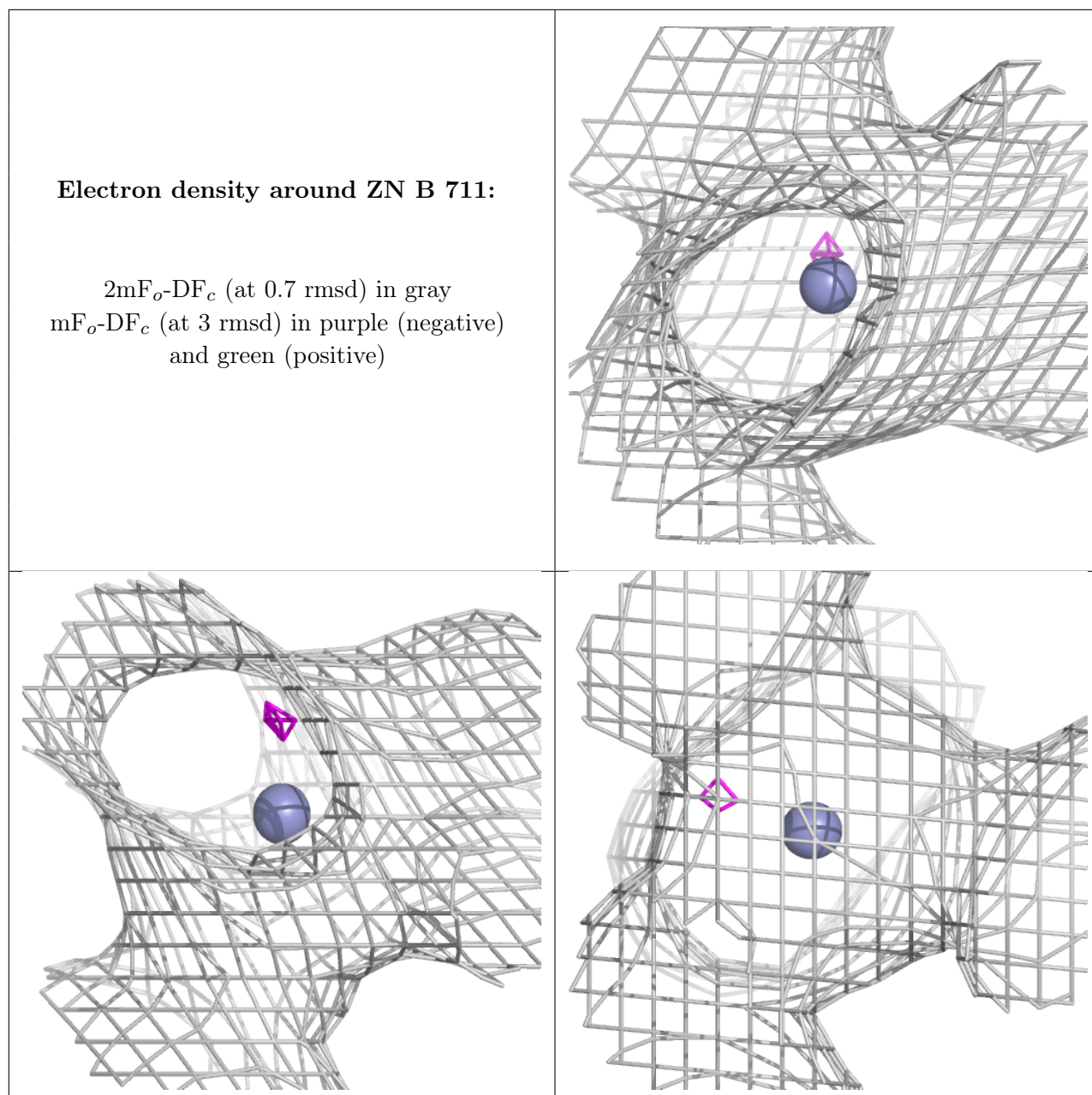
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around A1IJW A 701:

$2mF_o-DF_c$ (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.