



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 02:54 PM UTC

PDB ID : 9RO7 / pdb_00009ro7
Title : Crystal structure of human CD22 Ig domains 1-3 in complex with modified sialoside 7-012
Authors : Ereno-Orbea, J.; Sicard, T.; Julien, J.-P.
Deposited on : 2025-06-20
Resolution : 3.15 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:

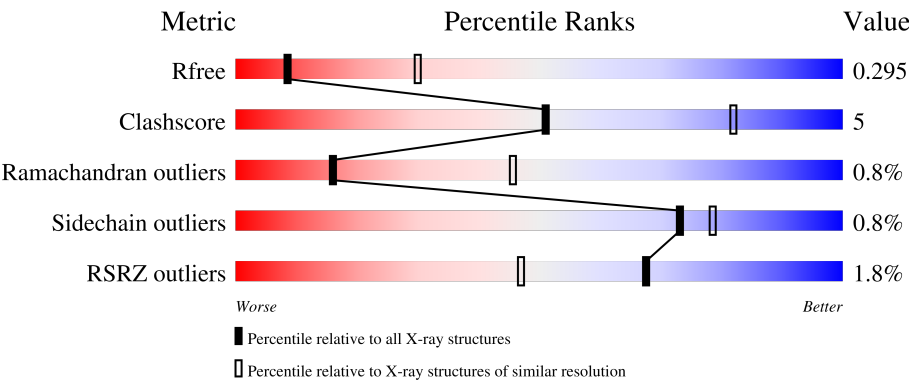
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal 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 i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.15 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.



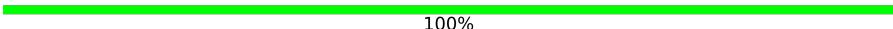
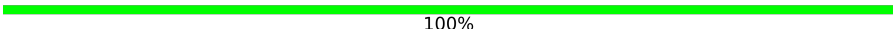
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R _{free}	180053	2361 (3.20-3.12)
Clashscore	190562	2486 (3.20-3.12)
Ramachandran outliers	187476	2405 (3.20-3.12)
Sidechain outliers	187428	2404 (3.20-3.12)
RSRZ outliers	180081	2361 (3.20-3.12)

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	324	% 83%11%5%
1	B	324	% 80%14%6%
1	C	324	3% 74%14%12%
1	D	324	% 77%14%9%
2	E	3	67%33%

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Mol	Chain	Length	Quality of chain
2	H	3	 67% 33%
2	J	3	 100%
2	L	3	 100%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9883 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 B-cell receptor CD22.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	307	Total	C	N	O	S	0	0	0
			2438	1543	413	469	13			
1	B	306	Total	C	N	O	S	0	0	0
			2432	1540	412	467	13			
1	C	286	Total	C	N	O	S	0	0	0
			2277	1444	382	438	13			
1	D	295	Total	C	N	O	S	0	0	0
			2343	1486	395	449	13			

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	17	GLU	-	expression tag	UNP P20273
A	18	THR	-	expression tag	UNP P20273
A	19	GLY	-	expression tag	UNP P20273
A	67	ALA	ASN	engineered mutation	UNP P20273
A	112	ALA	ASN	engineered mutation	UNP P20273
A	135	ALA	ASN	engineered mutation	UNP P20273
A	164	ALA	ASN	engineered mutation	UNP P20273
A	231	ALA	ASN	engineered mutation	UNP P20273
A	331	GLY	-	expression tag	UNP P20273
A	332	GLY	-	expression tag	UNP P20273
A	333	THR	-	expression tag	UNP P20273
A	334	LYS	-	expression tag	UNP P20273
A	335	HIS	-	expression tag	UNP P20273
A	336	HIS	-	expression tag	UNP P20273
A	337	HIS	-	expression tag	UNP P20273
A	338	HIS	-	expression tag	UNP P20273
A	339	HIS	-	expression tag	UNP P20273
A	340	HIS	-	expression tag	UNP P20273
B	17	GLU	-	expression tag	UNP P20273
B	18	THR	-	expression tag	UNP P20273
B	19	GLY	-	expression tag	UNP P20273

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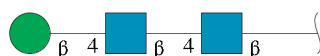
Chain	Residue	Modelled	Actual	Comment	Reference
B	67	ALA	ASN	engineered mutation	UNP P20273
B	112	ALA	ASN	engineered mutation	UNP P20273
B	135	ALA	ASN	engineered mutation	UNP P20273
B	164	ALA	ASN	engineered mutation	UNP P20273
B	231	ALA	ASN	engineered mutation	UNP P20273
B	331	GLY	-	expression tag	UNP P20273
B	332	GLY	-	expression tag	UNP P20273
B	333	THR	-	expression tag	UNP P20273
B	334	LYS	-	expression tag	UNP P20273
B	335	HIS	-	expression tag	UNP P20273
B	336	HIS	-	expression tag	UNP P20273
B	337	HIS	-	expression tag	UNP P20273
B	338	HIS	-	expression tag	UNP P20273
B	339	HIS	-	expression tag	UNP P20273
B	340	HIS	-	expression tag	UNP P20273
C	17	GLU	-	expression tag	UNP P20273
C	18	THR	-	expression tag	UNP P20273
C	19	GLY	-	expression tag	UNP P20273
C	67	ALA	ASN	engineered mutation	UNP P20273
C	112	ALA	ASN	engineered mutation	UNP P20273
C	135	ALA	ASN	engineered mutation	UNP P20273
C	164	ALA	ASN	engineered mutation	UNP P20273
C	231	ALA	ASN	engineered mutation	UNP P20273
C	331	GLY	-	expression tag	UNP P20273
C	332	GLY	-	expression tag	UNP P20273
C	333	THR	-	expression tag	UNP P20273
C	334	LYS	-	expression tag	UNP P20273
C	335	HIS	-	expression tag	UNP P20273
C	336	HIS	-	expression tag	UNP P20273
C	337	HIS	-	expression tag	UNP P20273
C	338	HIS	-	expression tag	UNP P20273
C	339	HIS	-	expression tag	UNP P20273
C	340	HIS	-	expression tag	UNP P20273
D	17	GLU	-	expression tag	UNP P20273
D	18	THR	-	expression tag	UNP P20273
D	19	GLY	-	expression tag	UNP P20273
D	67	ALA	ASN	engineered mutation	UNP P20273
D	112	ALA	ASN	engineered mutation	UNP P20273
D	135	ALA	ASN	engineered mutation	UNP P20273
D	164	ALA	ASN	engineered mutation	UNP P20273
D	231	ALA	ASN	engineered mutation	UNP P20273
D	331	GLY	-	expression tag	UNP P20273

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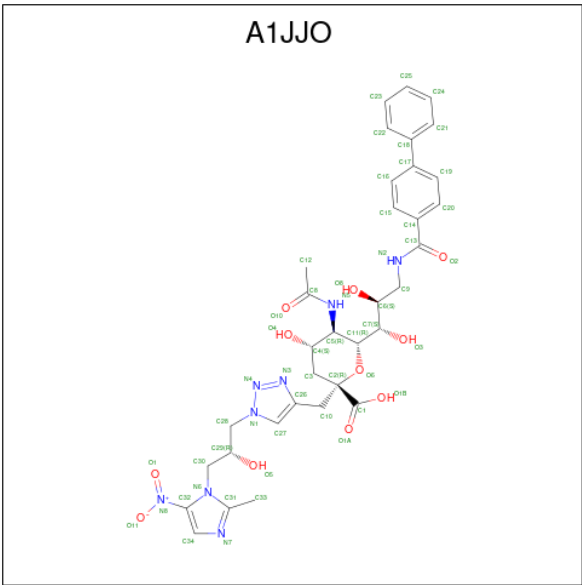
Chain	Residue	Modelled	Actual	Comment	Reference
D	332	GLY	-	expression tag	UNP P20273
D	333	THR	-	expression tag	UNP P20273
D	334	LYS	-	expression tag	UNP P20273
D	335	HIS	-	expression tag	UNP P20273
D	336	HIS	-	expression tag	UNP P20273
D	337	HIS	-	expression tag	UNP P20273
D	338	HIS	-	expression tag	UNP P20273
D	339	HIS	-	expression tag	UNP P20273
D	340	HIS	-	expression tag	UNP P20273

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



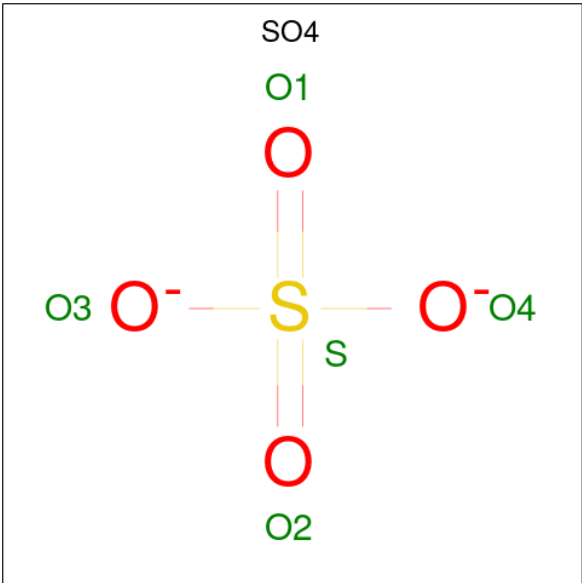
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	E	3	Total	C	N	O	0	0	0
			39	22	2	15			
2	H	3	Total	C	N	O	0	0	0
			39	22	2	15			
2	J	3	Total	C	N	O	0	0	0
			39	22	2	15			
2	L	3	Total	C	N	O	0	0	0
			39	22	2	15			

- Molecule 3 is (2R,4S,5R,6R)-5-acetamido-6-[(1S,2S)-1,2-bis(oxidanyl)-3-[(4-phenylphenyl)carbonylamino]propyl]-2-[[1-[(2R)-3-(2-methyl-5-nitro-imidazol-1-yl)-2-oxidanyl-propyl]-1,2,3-triazol-4-yl]methyl]-4-oxidanyl-oxane-2-carboxylic acid (CCD ID: A1JJO) (formula: C₃₄H₄₀N₈O₁₁) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			53	34	8	11		
3	B	1	Total	C	N	O	0	0
			53	34	8	11		
3	C	1	Total	C	N	O	0	0
			53	34	8	11		
3	D	1	Total	C	N	O	0	0
			53	34	8	11		

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S) (labeled as "Ligand of Interest" by depositor).

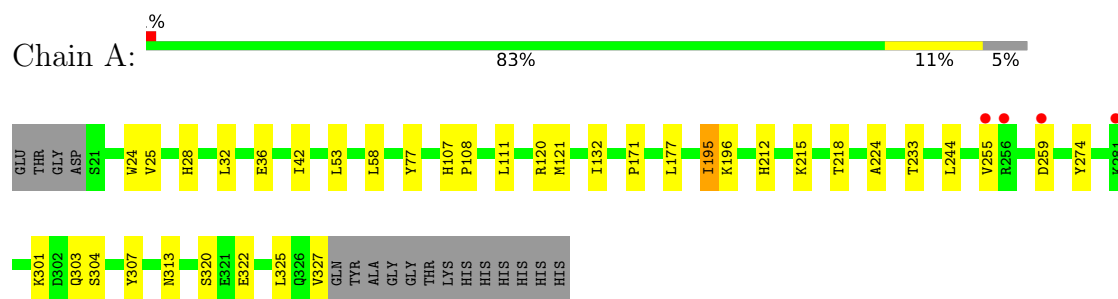


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		

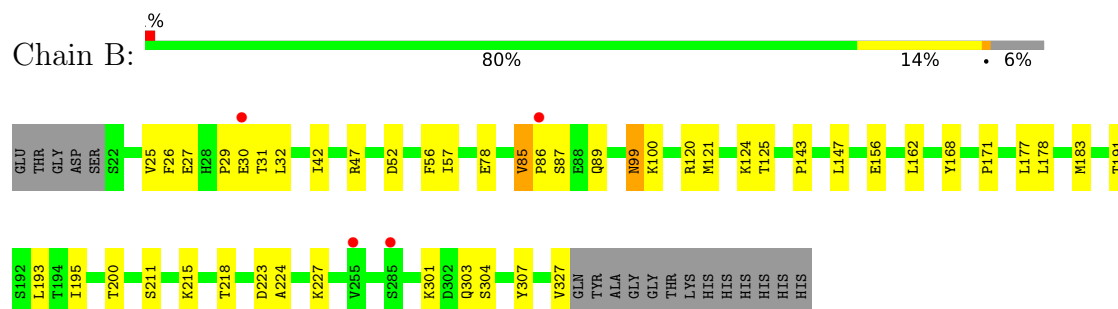
3 Residue-property plots [i](#)

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.

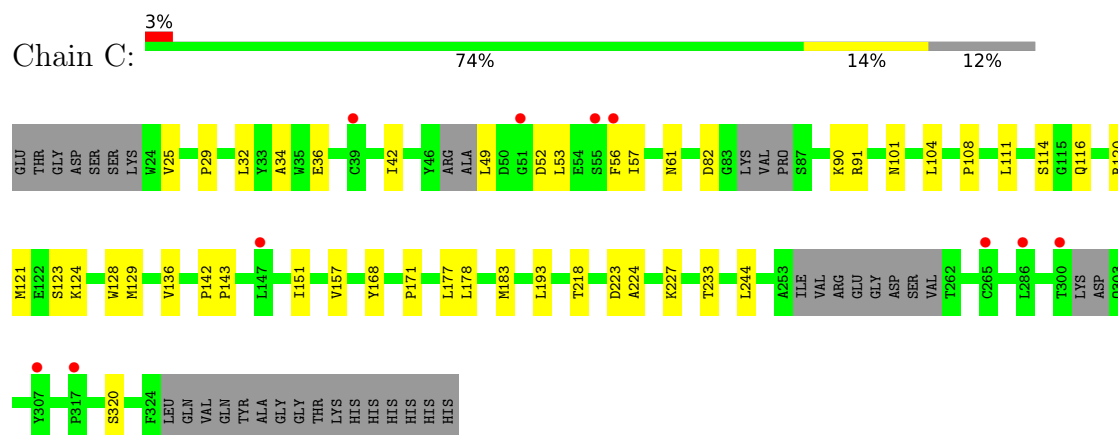
• Molecule 1: B-cell receptor CD22



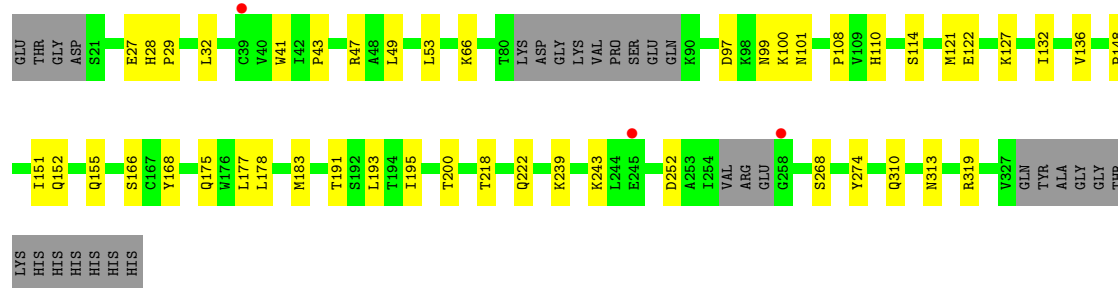
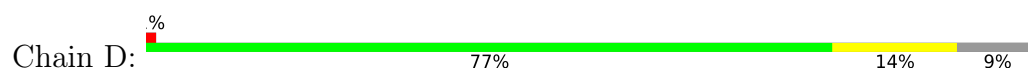
• Molecule 1: B-cell receptor CD22



• Molecule 1: B-cell receptor CD22



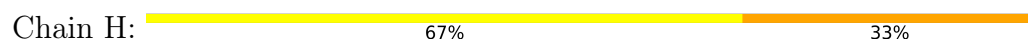
• Molecule 1: B-cell receptor CD22



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



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



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



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



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	102.39Å 113.29Å 138.37Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.42 – 3.15 39.42 – 3.15	Depositor EDS
% Data completeness (in resolution range)	99.8 (39.42-3.15) 99.8 (39.42-3.15)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.79 (at 3.12Å)	Xtriage
Refinement program	PHENIX (1.13_2998: ???)	Depositor
R, R_{free}	0.239 , 0.298 0.239 , 0.295	Depositor DCC
R_{free} test set	1425 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	61.6	Xtriage
Anisotropy	0.493	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 38.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	9883	wwPDB-VP
Average B, all atoms (Å ²)	66.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 30.24 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.3511e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: BMA, NAG, SO4, A1JJO

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	0.21	0/2497	0.39	0/3390
1	B	0.22	0/2491	0.40	0/3382
1	C	0.19	0/2331	0.44	3/3161 (0.1%)
1	D	0.18	0/2399	0.35	0/3256
All	All	0.20	0/9718	0.40	3/13189 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	123	SER	CA-C-O	8.72	123.82	118.33
1	C	123	SER	CB-CA-C	-8.58	103.92	116.53
1	C	123	SER	N-CA-C	6.67	115.16	108.75

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	2438	0	2391	23	0
1	B	2432	0	2386	27	0
1	C	2277	0	2212	25	0
1	D	2343	0	2293	27	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	E	39	0	34	0	0
2	H	39	0	34	2	0
2	J	39	0	34	0	0
2	L	39	0	34	0	0
3	A	53	0	0	3	0
3	B	53	0	0	2	0
3	C	53	0	0	3	0
3	D	53	0	0	0	0
4	A	5	0	0	0	0
4	B	5	0	0	0	0
4	C	10	0	0	0	0
4	D	5	0	0	0	0
All	All	9883	0	9418	100	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 5.

All (100) 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:85:VAL:HB	1:B:86:PRO:HD3	1.70	0.73
1:B:178:LEU:HB2	1:B:183:MET:HE2	1.72	0.70
1:B:177:LEU:HB2	1:B:218:THR:HB	1.74	0.68
1:D:175:GLN:OE1	1:D:222:GLN:NE2	2.27	0.67
1:C:57:ILE:HB	1:C:120:ARG:HB3	1.76	0.67
1:C:52:ASP:H	1:C:124:LYS:HE3	1.62	0.64
1:D:178:LEU:HB2	1:D:183:MET:HE2	1.80	0.63
1:B:171:PRO:HG2	1:B:224:ALA:HB2	1.82	0.61
1:A:53:LEU:HD11	1:A:121:MET:HE3	1.83	0.61
1:C:151:ILE:HG23	1:C:157:VAL:HG11	1.81	0.61
1:D:53:LEU:HD11	1:D:121:MET:HE3	1.83	0.60
1:B:32:LEU:HD22	1:B:42:ILE:HG12	1.83	0.60
1:B:303:GLN:O	1:B:307:TYR:OH	2.20	0.60
1:D:27:GLU:HB3	1:D:195:ILE:HD13	1.82	0.59
1:A:195:ILE:HD12	1:A:196:LYS:HG3	1.84	0.59
1:C:244:LEU:HB2	1:C:320:SER:HB3	1.85	0.59
1:C:53:LEU:HD11	1:C:121:MET:HB3	1.85	0.57
1:B:124:LYS:HG3	1:B:125:THR:HG23	1.87	0.56
1:C:120:ARG:HD3	1:C:129:MET:HE2	1.88	0.56
1:C:178:LEU:HB2	1:C:183:MET:HE2	1.87	0.56
1:B:27:GLU:HB3	1:B:195:ILE:HD13	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:28:HIS:CE1	1:D:132:ILE:HG12	2.41	0.55
1:A:120:ARG:HH12	3:A:401:A1JJO:C1	2.21	0.54
1:A:304:SER:HA	1:A:325:LEU:HB3	1.90	0.54
1:D:243:LYS:O	1:D:268:SER:OG	2.26	0.54
1:C:177:LEU:HB2	1:C:218:THR:HB	1.89	0.54
1:C:171:PRO:HG2	1:C:224:ALA:HB2	1.90	0.54
1:A:304:SER:HB3	1:A:327:VAL:HG13	1.90	0.53
1:D:152:GLN:HG2	1:D:239:LYS:HB2	1.90	0.53
1:B:29:PRO:O	1:B:31:THR:N	2.37	0.53
1:A:36:GLU:OE1	1:A:111:LEU:N	2.32	0.52
1:B:168:TYR:CZ	1:B:193:LEU:HD23	2.44	0.52
2:H:1:NAG:H61	2:H:2:NAG:C7	2.39	0.52
1:B:47:ARG:HD2	1:D:49:LEU:HD11	1.91	0.51
1:B:120:ARG:NH1	3:B:401:A1JJO:O1A	2.33	0.51
2:H:2:NAG:H83	2:H:2:NAG:H3	1.92	0.51
1:B:301:LYS:HA	1:B:327:VAL:HG21	1.91	0.51
1:B:56:PHE:HD1	1:B:121:MET:HG3	1.76	0.50
1:C:142:PRO:HD2	1:D:66:LYS:HG2	1.94	0.50
1:A:212:HIS:HA	1:A:215:LYS:HD2	1.93	0.50
1:C:120:ARG:NH1	3:C:401:A1JJO:O1A	2.45	0.49
1:C:29:PRO:HG2	1:C:32:LEU:HD11	1.94	0.49
3:C:401:A1JJO:O3	3:C:401:A1JJO:N5	2.45	0.49
1:D:108:PRO:HG2	1:D:110:HIS:CE1	2.48	0.49
1:B:120:ARG:HH12	3:B:401:A1JJO:C1	2.24	0.49
1:B:223:ASP:OD1	1:B:227:LYS:N	2.46	0.49
1:A:32:LEU:HD22	1:A:42:ILE:HG12	1.95	0.48
1:B:52:ASP:OD1	1:D:47:ARG:NH2	2.45	0.48
1:C:168:TYR:CZ	1:C:193:LEU:HD23	2.48	0.48
1:D:97:ASP:H	1:D:101:ASN:HB3	1.79	0.48
1:B:304:SER:HB2	1:B:327:VAL:HG23	1.95	0.47
1:B:87:SER:C	1:B:89:GLN:H	2.21	0.47
1:C:82:ASP:OD1	1:C:90:LYS:NZ	2.45	0.47
1:C:114:SER:HB2	1:C:136:VAL:HB	1.97	0.47
1:D:191:THR:HG22	1:D:200:THR:HG23	1.96	0.47
1:B:99:ASN:ND2	1:D:49:LEU:O	2.47	0.47
1:B:26:PHE:CE2	1:B:121:MET:HE1	2.51	0.46
1:D:41:TRP:O	1:D:43:PRO:HD3	2.15	0.46
1:B:47:ARG:HG3	1:D:49:LEU:HD21	1.97	0.46
1:C:120:ARG:HH12	3:C:401:A1JJO:C1	2.29	0.46
1:A:171:PRO:HG2	1:A:224:ALA:HB2	1.97	0.46
1:A:274:TYR:HA	1:A:313:ASN:HB3	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:28:HIS:CE1	1:A:132:ILE:HG12	2.52	0.45
1:C:61:ASN:OD1	1:C:116:GLN:N	2.41	0.45
1:D:29:PRO:HG2	1:D:32:LEU:HD11	1.97	0.45
1:A:120:ARG:NH1	3:A:401:A1JJO:O1A	2.50	0.45
1:C:223:ASP:OD1	1:C:227:LYS:N	2.47	0.45
1:A:255:VAL:HG12	1:A:327:VAL:HA	1.99	0.44
1:C:218:THR:OG1	1:C:233:THR:HG22	2.18	0.44
1:D:274:TYR:HA	1:D:313:ASN:HB3	2.00	0.43
1:A:177:LEU:HB2	1:A:218:THR:HB	1.99	0.43
1:A:218:THR:HG23	1:A:233:THR:HG22	1.99	0.43
1:D:310:GLN:HE21	1:D:319:ARG:HE	1.66	0.43
1:A:255:VAL:HG22	1:A:259:ASP:HB3	2.00	0.43
1:A:107:HIS:HA	1:A:108:PRO:HA	1.82	0.43
1:B:191:THR:HG22	1:B:200:THR:HG23	2.01	0.43
1:B:100:LYS:HG2	1:D:49:LEU:HB3	1.99	0.43
1:C:42:ILE:HB	1:C:104:LEU:HB2	2.00	0.42
1:A:58:LEU:HB3	1:A:77:TYR:HB3	2.01	0.42
1:A:244:LEU:HB2	1:A:320:SER:HB3	2.00	0.42
1:A:24:TRP:HZ3	1:A:121:MET:HE2	1.85	0.42
1:A:301:LYS:HD2	1:A:327:VAL:HG21	2.01	0.42
1:C:56:PHE:CZ	1:C:104:LEU:HD21	2.55	0.42
1:B:211:SER:O	1:B:215:LYS:HE2	2.20	0.41
1:D:148:PRO:HD2	1:D:151:ILE:HD11	2.03	0.41
1:A:307:TYR:O	1:A:322:GLU:HA	2.20	0.41
1:B:57:ILE:HG23	1:B:78:GLU:HG2	2.02	0.41
1:C:91:ARG:HG3	1:C:108:PRO:HD2	2.03	0.41
1:D:122:GLU:HG2	1:D:127:LYS:HG2	2.02	0.41
1:D:177:LEU:HB2	1:D:218:THR:HB	2.01	0.41
1:D:114:SER:HB2	1:D:136:VAL:HB	2.01	0.41
1:D:252:ASP:OD1	1:D:252:ASP:N	2.51	0.41
3:A:401:A1JJO:O1	3:A:401:A1JJO:O5	2.38	0.41
1:C:36:GLU:OE1	1:C:111:LEU:N	2.45	0.41
1:A:195:ILE:H	1:A:195:ILE:HG13	1.61	0.41
1:B:85:VAL:HB	1:B:86:PRO:CD	2.47	0.41
1:D:152:GLN:HB2	1:D:155:GLN:CD	2.46	0.40
1:C:34:ALA:O	1:C:136:VAL:HA	2.20	0.40
1:C:121:MET:HE3	1:C:128:TRP:HD1	1.87	0.40
1:D:168:TYR:CE2	1:D:193:LEU:HD23	2.57	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	305/324 (94%)	276 (90%)	28 (9%)	1 (0%)	36	64
1	B	304/324 (94%)	277 (91%)	23 (8%)	4 (1%)	9	36
1	C	276/324 (85%)	257 (93%)	18 (6%)	1 (0%)	30	59
1	D	289/324 (89%)	270 (93%)	16 (6%)	3 (1%)	12	41
All	All	1174/1296 (91%)	1080 (92%)	85 (7%)	9 (1%)	16	46

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	85	VAL
1	A	303	GLN
1	B	147	LEU
1	D	99	ASN
1	D	166	SER
1	B	99	ASN
1	C	143	PRO
1	D	100	LYS
1	B	143	PRO

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	279/292 (96%)	277 (99%)	2 (1%)	76	80
1	B	278/292 (95%)	274 (99%)	4 (1%)	59	73

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	260/292 (89%)	257 (99%)	3 (1%)	63	74
1	D	268/292 (92%)	268 (100%)	0	100	100
All	All	1085/1168 (93%)	1076 (99%)	9 (1%)	73	78

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	VAL
1	A	195	ILE
1	B	25	VAL
1	B	30	GLU
1	B	156	GLU
1	B	162	LEU
1	C	25	VAL
1	C	49	LEU
1	C	101	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	235	GLN
1	B	60	HIS
1	B	93	GLN
1	B	144	HIS
1	C	99	ASN
1	D	60	HIS
1	D	116	GLN
1	D	303	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates

12 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
2	NAG	E	1	2,1	14,14,15	0.57	0	17,19,21	0.54	0
2	NAG	E	2	2	14,14,15	0.33	0	17,19,21	0.87	1 (5%)
2	BMA	E	3	2	11,11,12	0.89	0	15,15,17	0.79	0
2	NAG	H	1	2,1	14,14,15	0.37	0	17,19,21	0.51	0
2	NAG	H	2	2	14,14,15	0.45	0	17,19,21	1.46	3 (17%)
2	BMA	H	3	2	11,11,12	0.78	0	15,15,17	1.04	1 (6%)
2	NAG	J	1	2,1	14,14,15	0.47	0	17,19,21	0.58	0
2	NAG	J	2	2	14,14,15	0.26	0	17,19,21	0.51	0
2	BMA	J	3	2	11,11,12	0.86	0	15,15,17	0.81	0
2	NAG	L	1	2,1	14,14,15	0.33	0	17,19,21	0.61	0
2	NAG	L	2	2	14,14,15	0.36	0	17,19,21	0.54	0
2	BMA	L	3	2	11,11,12	0.84	0	15,15,17	0.75	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	NAG	E	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	E	2	2	-	4/6/23/26	0/1/1/1
2	BMA	E	3	2	-	0/2/19/22	0/1/1/1
2	NAG	H	1	2,1	-	2/6/23/26	0/1/1/1
2	NAG	H	2	2	-	4/6/23/26	0/1/1/1
2	BMA	H	3	2	-	2/2/19/22	0/1/1/1
2	NAG	J	1	2,1	-	2/6/23/26	0/1/1/1
2	NAG	J	2	2	-	1/6/23/26	0/1/1/1
2	BMA	J	3	2	-	1/2/19/22	0/1/1/1
2	NAG	L	1	2,1	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	L	2	2	-	2/6/23/26	0/1/1/1
2	BMA	L	3	2	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	2	NAG	C2-N2-C7	4.39	128.79	122.90
2	H	3	BMA	C1-O5-C5	2.57	115.63	112.19
2	E	2	NAG	C1-O5-C5	2.47	115.50	112.19
2	H	2	NAG	C1-O5-C5	2.42	115.43	112.19
2	H	2	NAG	C1-C2-N2	2.19	113.89	110.43

There are no chirality outliers.

All (20) torsion outliers are listed below:

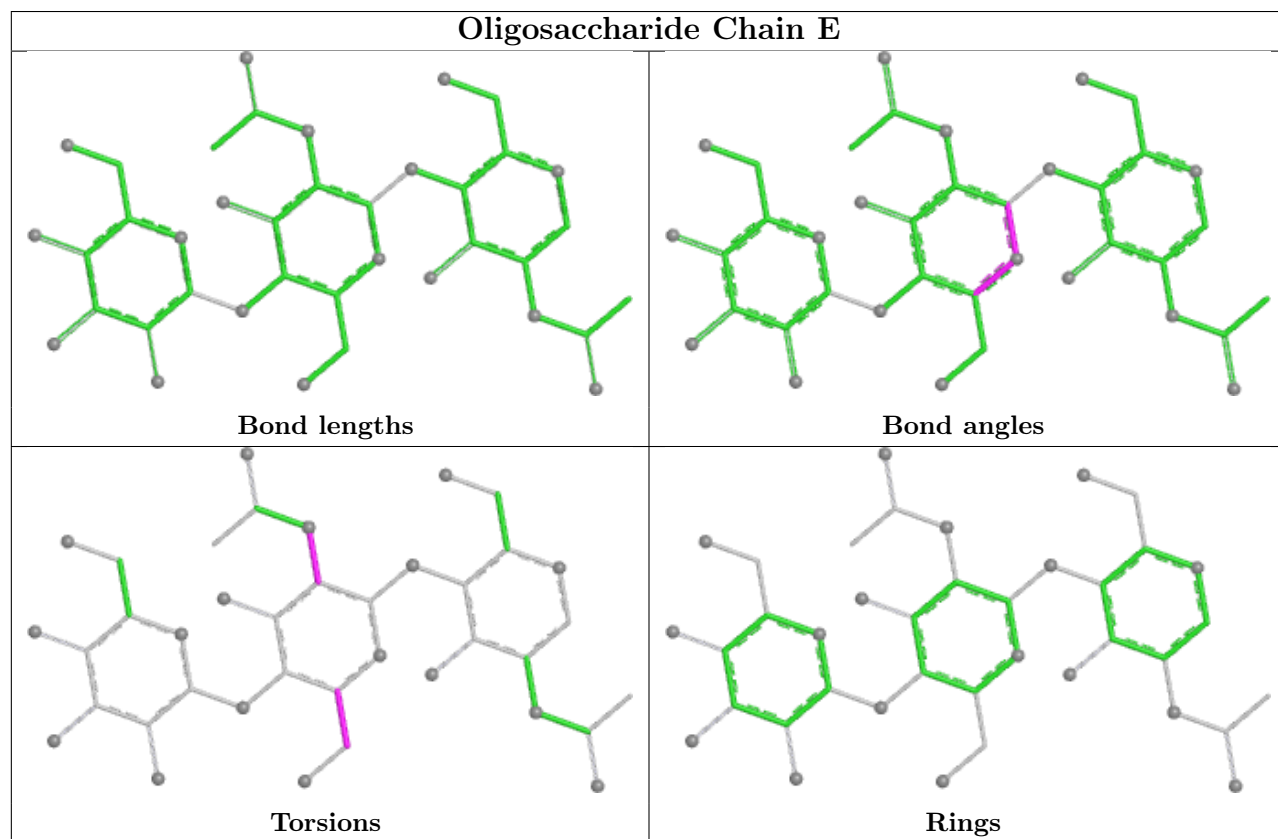
Mol	Chain	Res	Type	Atoms
2	E	2	NAG	O5-C5-C6-O6
2	L	2	NAG	C4-C5-C6-O6
2	L	1	NAG	O5-C5-C6-O6
2	L	2	NAG	O5-C5-C6-O6
2	L	1	NAG	C4-C5-C6-O6
2	E	2	NAG	C4-C5-C6-O6
2	H	1	NAG	C8-C7-N2-C2
2	H	1	NAG	O7-C7-N2-C2
2	H	2	NAG	C8-C7-N2-C2
2	H	2	NAG	O7-C7-N2-C2
2	J	2	NAG	O5-C5-C6-O6
2	J	3	BMA	O5-C5-C6-O6
2	H	3	BMA	C4-C5-C6-O6
2	H	3	BMA	O5-C5-C6-O6
2	J	1	NAG	C4-C5-C6-O6
2	E	2	NAG	C3-C2-N2-C7
2	J	1	NAG	O5-C5-C6-O6
2	E	2	NAG	C1-C2-N2-C7
2	H	2	NAG	C1-C2-N2-C7
2	H	2	NAG	C3-C2-N2-C7

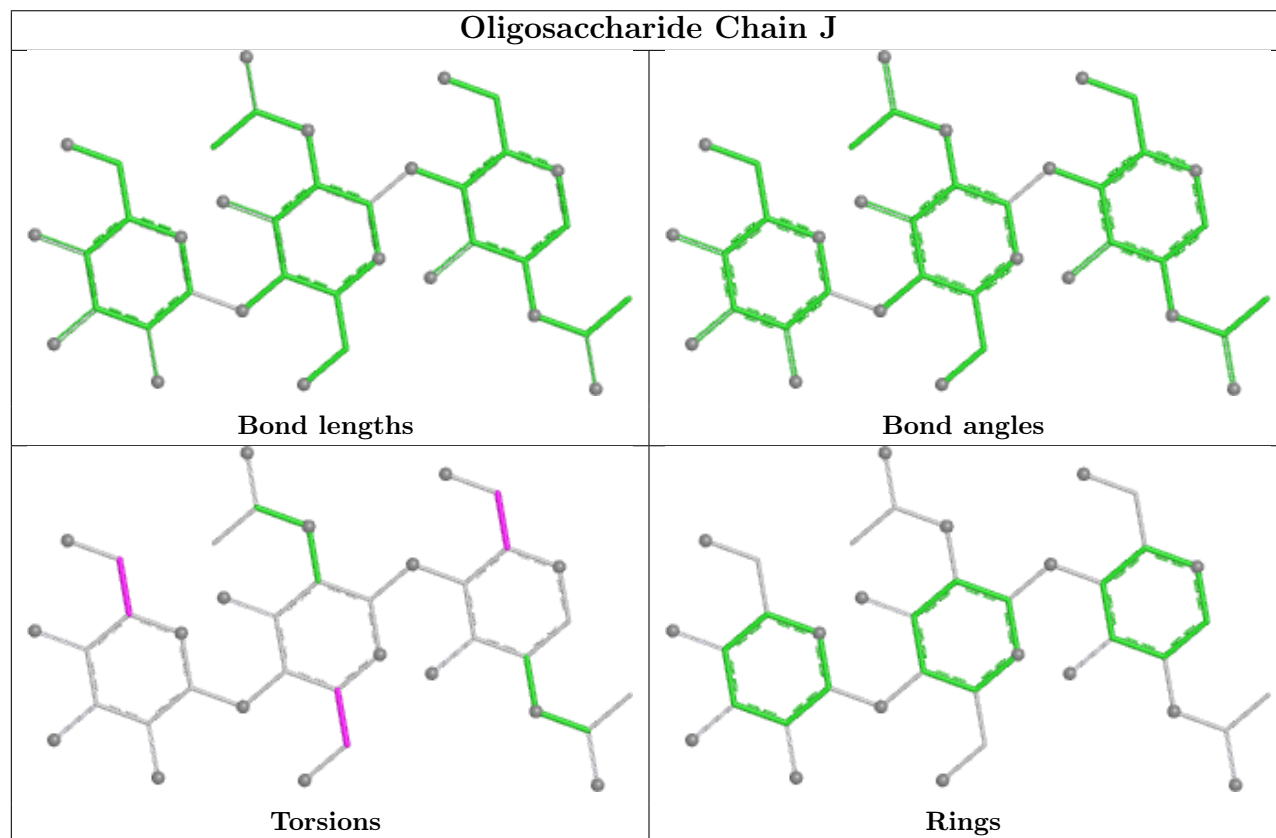
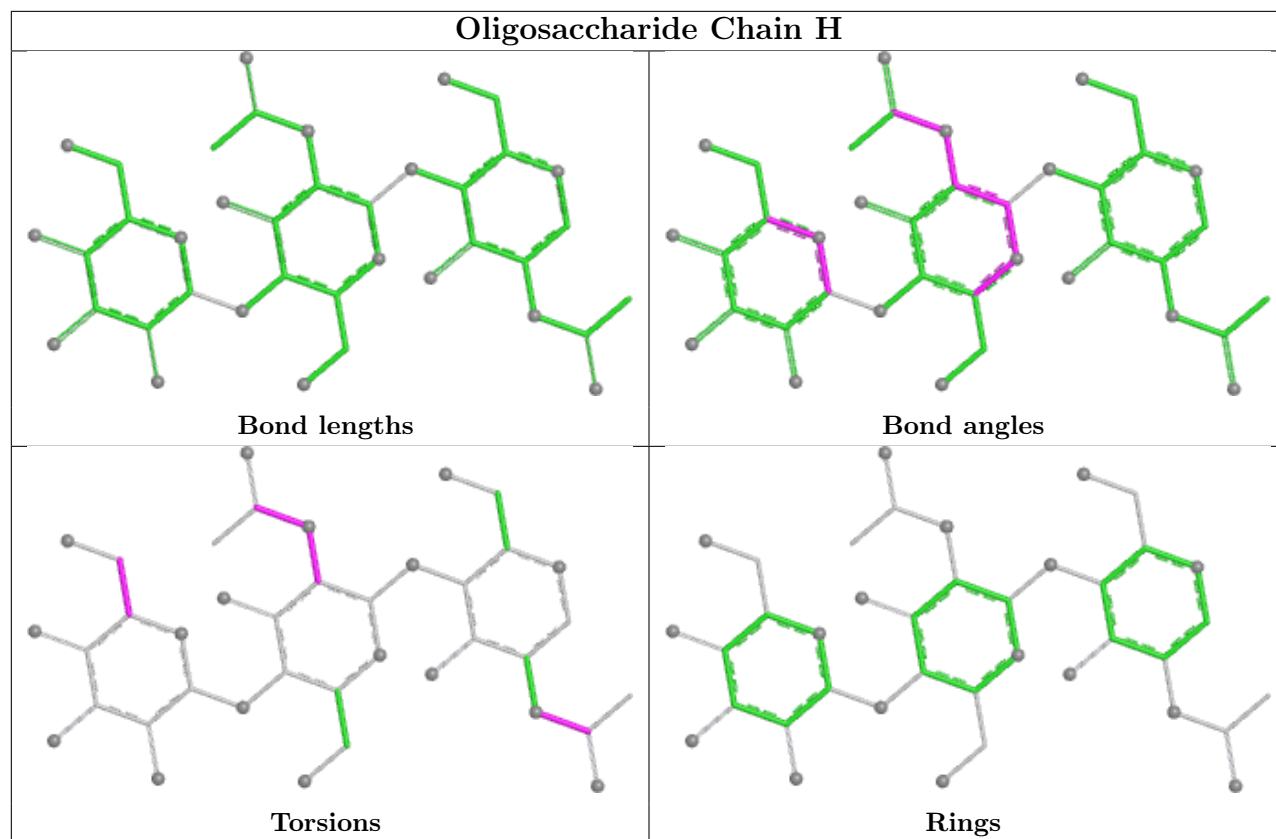
There are no ring outliers.

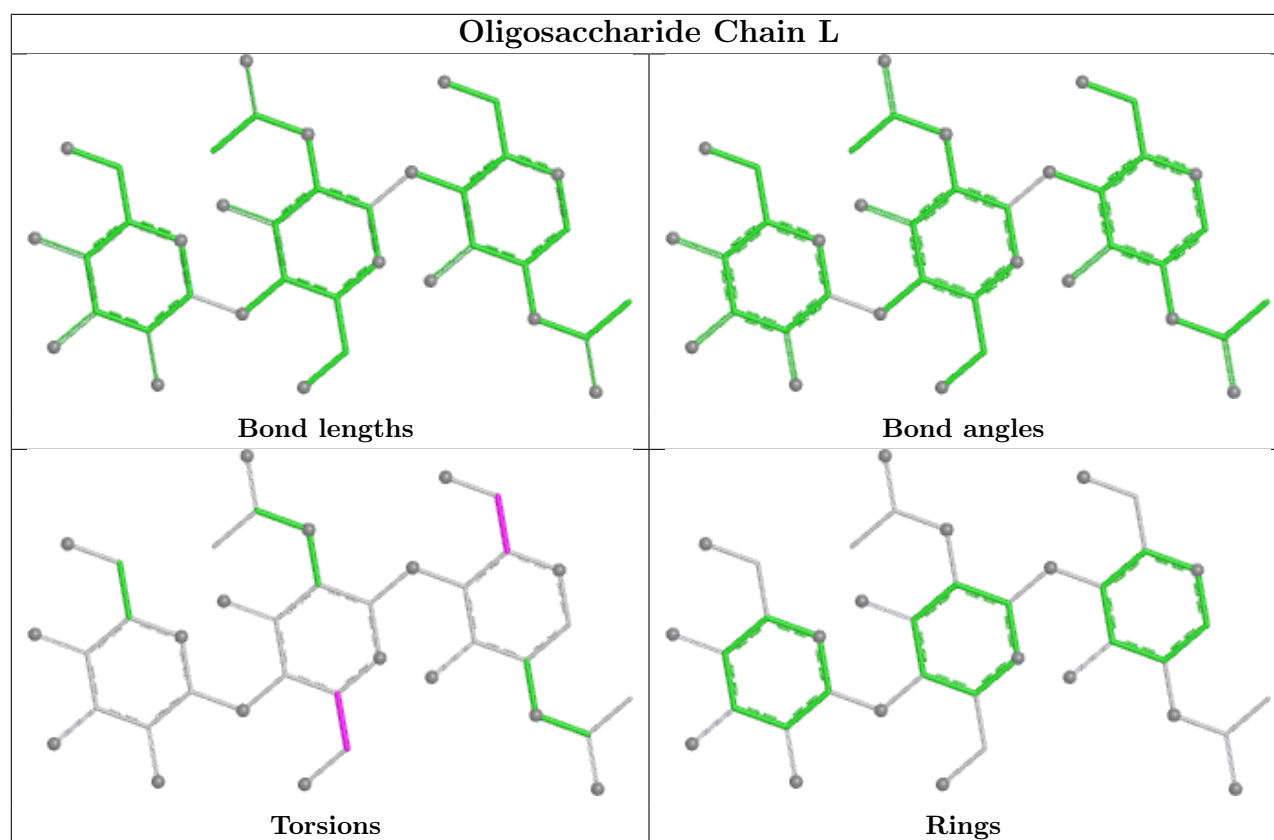
2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	H	2	NAG	2	0
2	H	1	NAG	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 oligosaccharide.







5.6 Ligand geometry [i](#)

9 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
3	A1JJO	D	401	-	55,57,57	0.48	1 (1%)	71,82,82	0.62	2 (2%)
4	SO4	B	402	-	4,4,4	0.25	0	6,6,6	0.12	0
4	SO4	A	402	-	4,4,4	0.25	0	6,6,6	0.18	0
3	A1JJO	A	401	-	55,57,57	0.49	1 (1%)	71,82,82	0.62	1 (1%)
4	SO4	C	403	-	4,4,4	0.23	0	6,6,6	0.31	0
3	A1JJO	B	401	-	55,57,57	0.49	1 (1%)	71,82,82	0.81	2 (2%)
3	A1JJO	C	401	-	55,57,57	0.48	1 (1%)	71,82,82	0.67	2 (2%)
4	SO4	C	402	-	4,4,4	0.23	0	6,6,6	0.16	0
4	SO4	D	402	-	4,4,4	0.25	0	6,6,6	0.13	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
3	A1JJO	C	401	-	-	19/46/66/66	0/5/5/5
3	A1JJO	B	401	-	-	17/46/66/66	0/5/5/5
3	A1JJO	D	401	-	-	21/46/66/66	0/5/5/5
3	A1JJO	A	401	-	-	18/46/66/66	0/5/5/5

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	401	A1JJO	C2-C1	2.13	1.54	1.52
3	B	401	A1JJO	C2-C1	2.10	1.54	1.52
3	C	401	A1JJO	C2-C1	2.08	1.54	1.52
3	D	401	A1JJO	C2-C1	2.07	1.54	1.52

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	401	A1JJO	O6-C11-C5	5.18	114.59	109.84
3	C	401	A1JJO	O6-C11-C5	3.18	112.75	109.84
3	D	401	A1JJO	O6-C11-C5	3.16	112.74	109.84
3	A	401	A1JJO	O6-C11-C5	2.68	112.30	109.84
3	C	401	A1JJO	C10-C2-C1	-2.63	102.83	108.69
3	D	401	A1JJO	O6-C2-C10	2.50	112.65	104.44
3	B	401	A1JJO	O6-C2-C10	2.12	111.40	104.44

There are no chirality outliers.

All (75) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	401	A1JJO	C5-C11-C7-C6
3	A	401	A1JJO	C5-C11-C7-O3
3	A	401	A1JJO	O6-C11-C7-C6
3	A	401	A1JJO	O6-C11-C7-O3
3	A	401	A1JJO	C26-C10-C2-C1
3	A	401	A1JJO	C26-C10-C2-C3
3	A	401	A1JJO	C26-C10-C2-O6
3	A	401	A1JJO	C2-C10-C26-C27
3	A	401	A1JJO	C2-C10-C26-N3

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Mol	Chain	Res	Type	Atoms
3	A	401	A1JJO	C7-C6-C9-N2
3	A	401	A1JJO	O8-C6-C9-N2
3	A	401	A1JJO	C28-C29-C30-N6
3	B	401	A1JJO	C5-C11-C7-C6
3	B	401	A1JJO	C5-C11-C7-O3
3	B	401	A1JJO	O6-C11-C7-C6
3	B	401	A1JJO	O6-C11-C7-O3
3	B	401	A1JJO	C26-C10-C2-C1
3	B	401	A1JJO	C26-C10-C2-C3
3	B	401	A1JJO	C26-C10-C2-O6
3	B	401	A1JJO	C2-C10-C26-C27
3	B	401	A1JJO	C2-C10-C26-N3
3	B	401	A1JJO	C7-C6-C9-N2
3	B	401	A1JJO	O8-C6-C9-N2
3	C	401	A1JJO	C5-C11-C7-C6
3	C	401	A1JJO	C5-C11-C7-O3
3	C	401	A1JJO	C26-C10-C2-C1
3	C	401	A1JJO	C26-C10-C2-C3
3	C	401	A1JJO	C26-C10-C2-O6
3	C	401	A1JJO	C2-C10-C26-C27
3	C	401	A1JJO	C2-C10-C26-N3
3	C	401	A1JJO	C7-C6-C9-N2
3	C	401	A1JJO	O8-C6-C9-N2
3	D	401	A1JJO	C5-C11-C7-C6
3	D	401	A1JJO	C5-C11-C7-O3
3	D	401	A1JJO	O6-C11-C7-C6
3	D	401	A1JJO	O6-C11-C7-O3
3	D	401	A1JJO	C26-C10-C2-C1
3	D	401	A1JJO	C26-C10-C2-C3
3	D	401	A1JJO	C26-C10-C2-O6
3	D	401	A1JJO	C2-C10-C26-C27
3	D	401	A1JJO	C2-C10-C26-N3
3	D	401	A1JJO	C7-C6-C9-N2
3	D	401	A1JJO	O8-C6-C9-N2
3	D	401	A1JJO	C28-C29-C30-N6
3	B	401	A1JJO	C28-C29-C30-N6
3	C	401	A1JJO	C28-C29-C30-N6
3	D	401	A1JJO	O8-C6-C7-C11
3	A	401	A1JJO	C29-C28-N1-N4
3	B	401	A1JJO	C29-C28-N1-N4
3	C	401	A1JJO	C29-C28-N1-N4
3	D	401	A1JJO	C29-C28-N1-N4

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Mol	Chain	Res	Type	Atoms
3	A	401	A1JJO	C29-C28-N1-C27
3	B	401	A1JJO	C29-C28-N1-C27
3	C	401	A1JJO	C29-C28-N1-C27
3	D	401	A1JJO	C29-C28-N1-C27
3	C	401	A1JJO	O6-C11-C7-O3
3	A	401	A1JJO	N1-C28-C29-C30
3	C	401	A1JJO	N1-C28-C29-C30
3	C	401	A1JJO	N1-C28-C29-O5
3	A	401	A1JJO	O1A-C1-C2-C10
3	A	401	A1JJO	O1B-C1-C2-C10
3	B	401	A1JJO	O1A-C1-C2-C10
3	B	401	A1JJO	O1B-C1-C2-C10
3	C	401	A1JJO	O6-C11-C7-C6
3	C	401	A1JJO	O1A-C1-C2-C10
3	C	401	A1JJO	O1B-C1-C2-C10
3	C	401	A1JJO	O1B-C1-C2-C3
3	D	401	A1JJO	O1A-C1-C2-C10
3	D	401	A1JJO	O1B-C1-C2-C10
3	D	401	A1JJO	O5-C29-C30-N6
3	B	401	A1JJO	O8-C6-C7-C11
3	D	401	A1JJO	O8-C6-C7-O3
3	A	401	A1JJO	O1B-C1-C2-C3
3	D	401	A1JJO	O1A-C1-C2-C3
3	D	401	A1JJO	O1B-C1-C2-C3

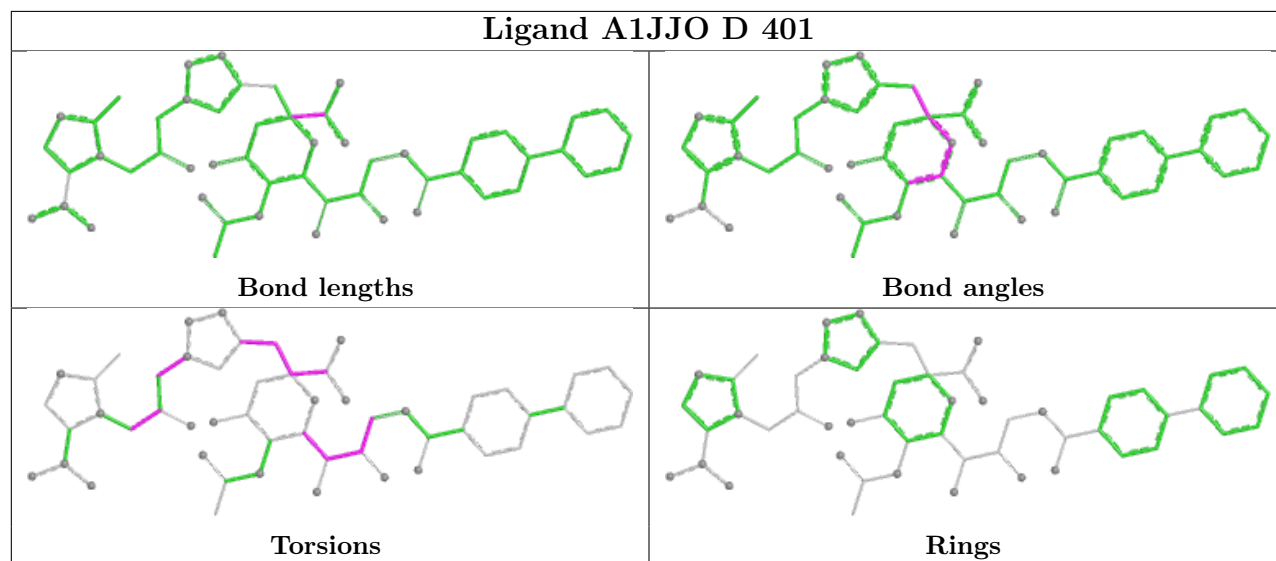
There are no ring outliers.

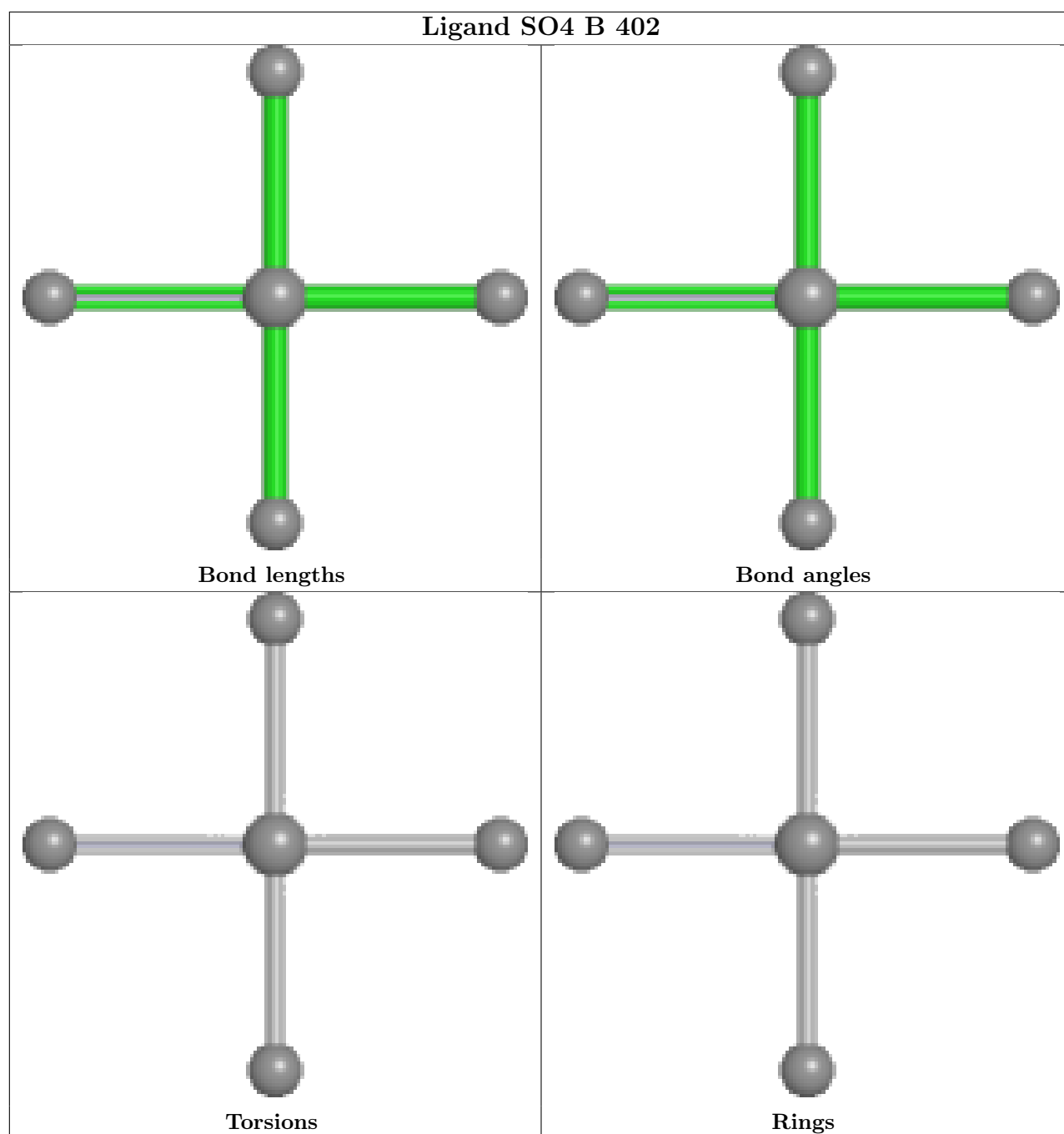
3 monomers are involved in 8 short contacts:

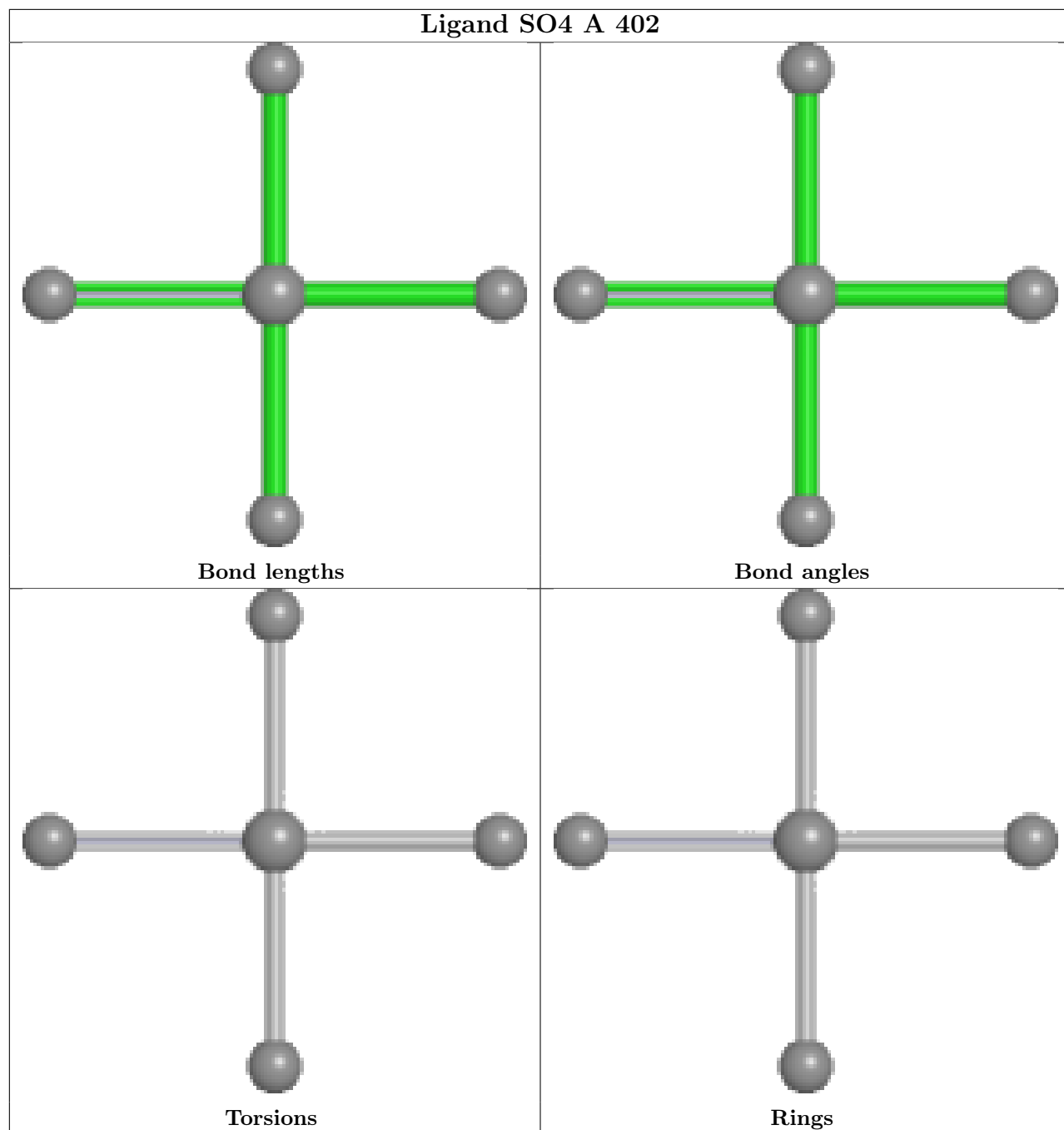
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	401	A1JJO	3	0
3	B	401	A1JJO	2	0
3	C	401	A1JJO	3	0

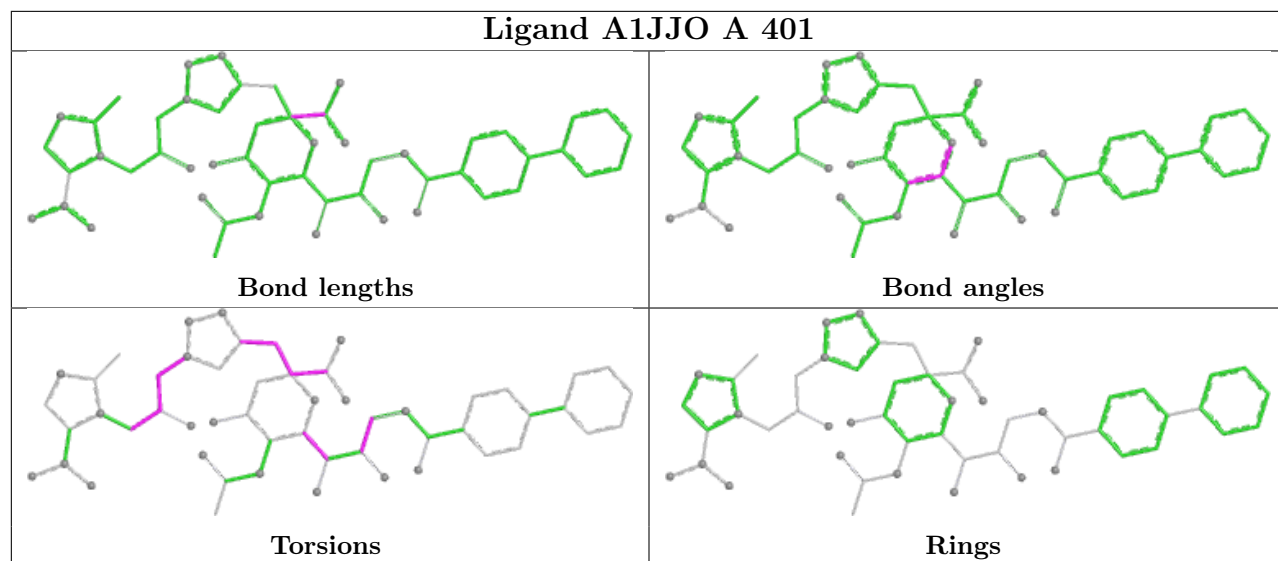
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and

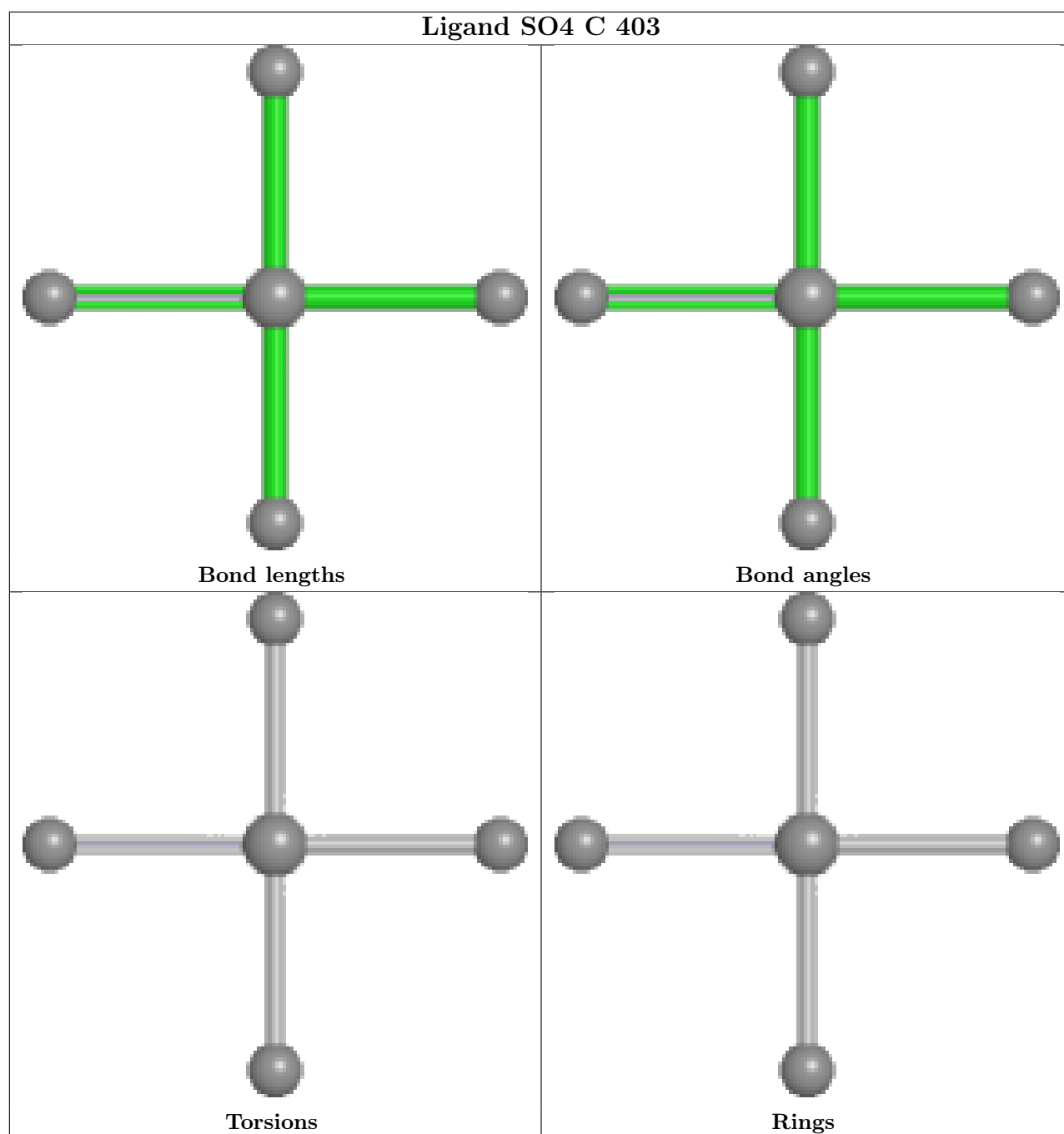
any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



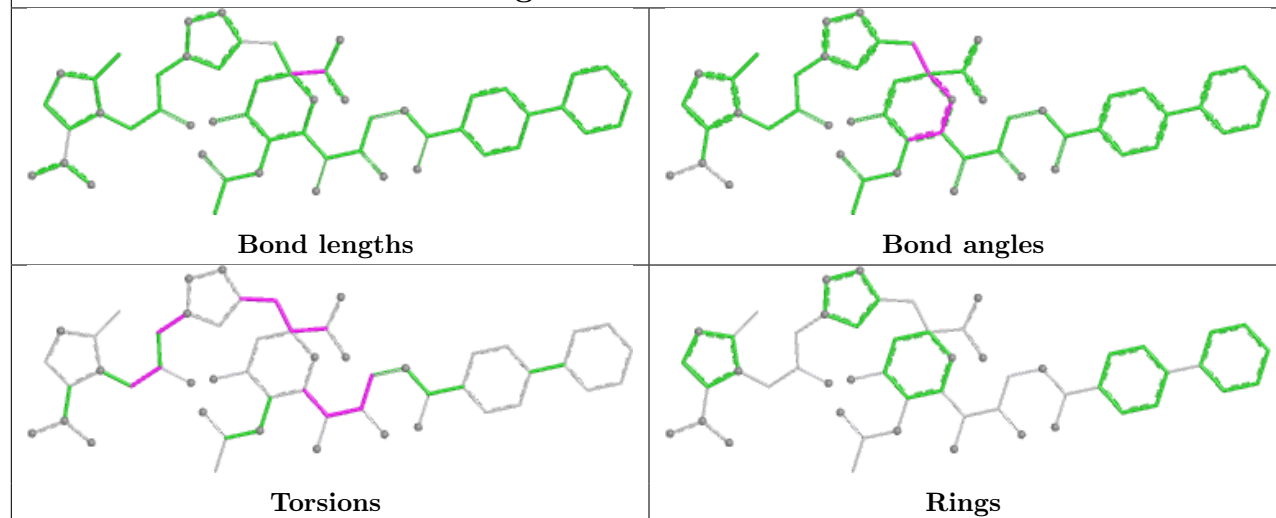




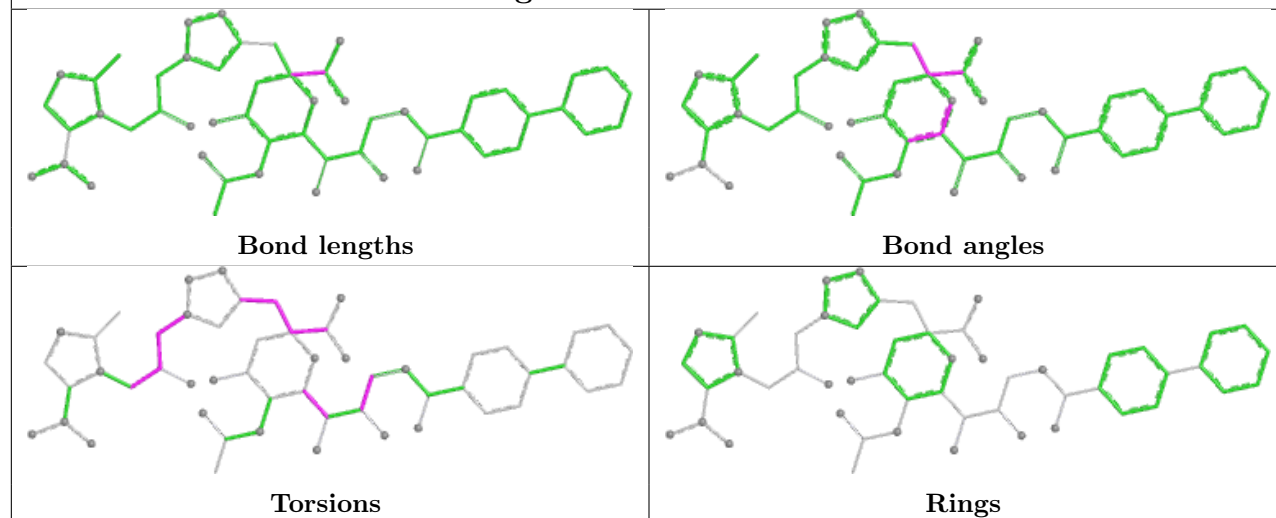


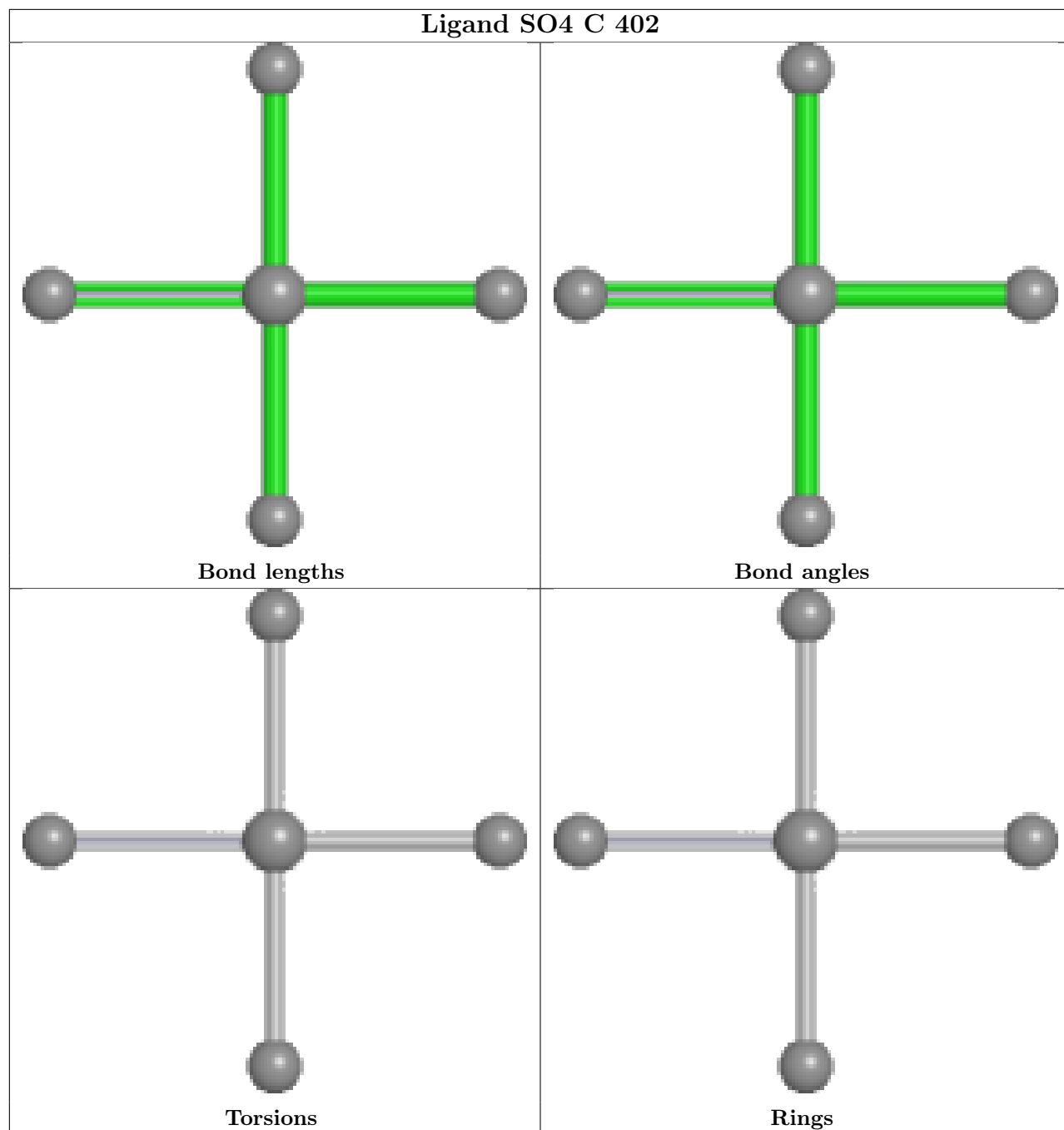


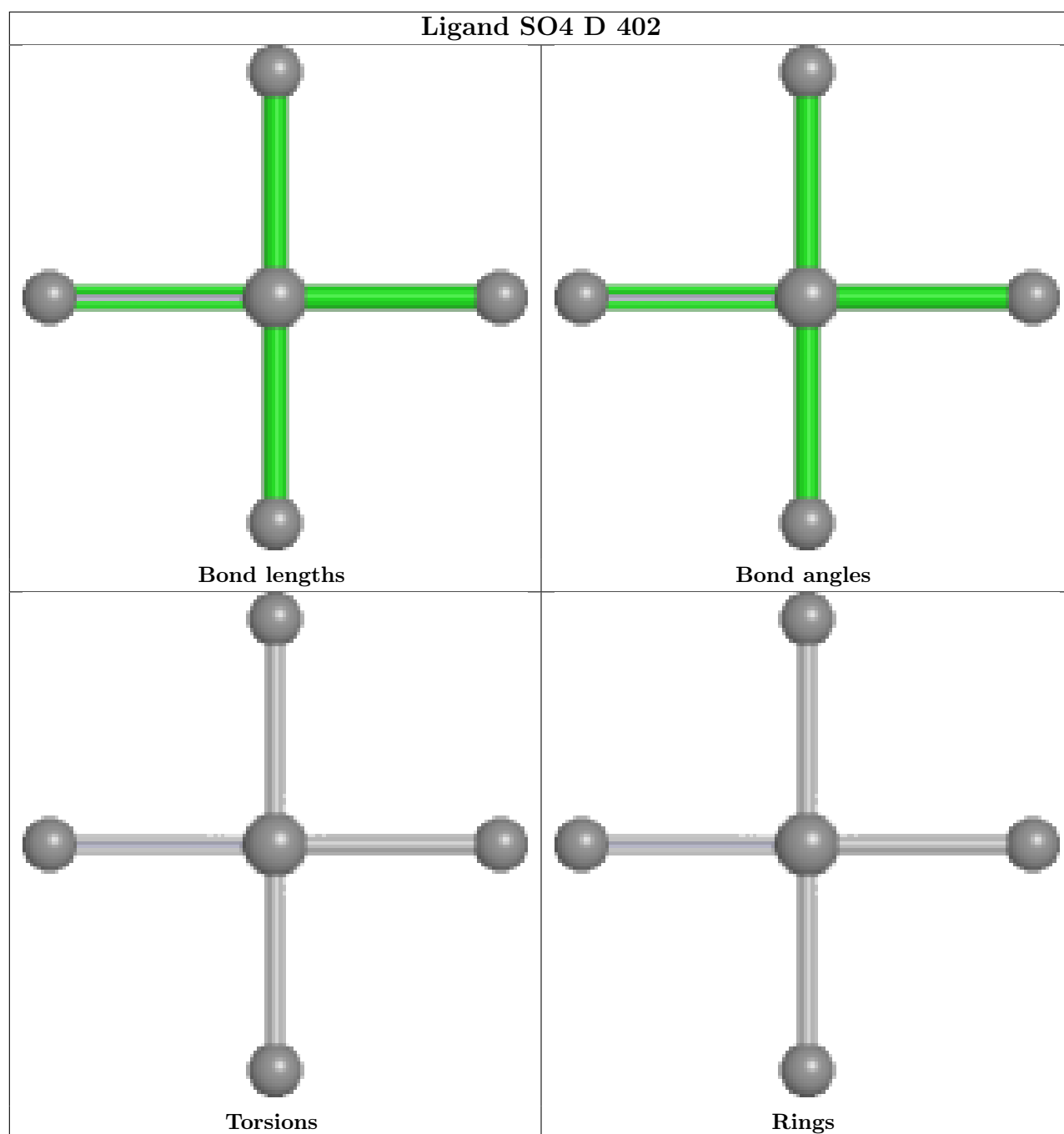
Ligand A1JJO B 401



Ligand A1JJO C 401







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	307/324 (94%)	0.02	4 (1%) 75 55	31, 55, 90, 116	0
1	B	306/324 (94%)	0.08	4 (1%) 75 55	31, 56, 100, 126	0
1	C	286/324 (88%)	0.48	10 (3%) 47 28	44, 74, 128, 149	0
1	D	295/324 (91%)	0.21	3 (1%) 79 62	37, 65, 94, 118	0
All	All	1194/1296 (92%)	0.19	21 (1%) 67 47	31, 62, 113, 149	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	258	GLY	2.8
1	A	281	LYS	2.8
1	C	317	PRO	2.7
1	C	307	TYR	2.7
1	B	86	PRO	2.6
1	A	259	ASP	2.5
1	D	39	CYS	2.4
1	A	256	ARG	2.3
1	B	255	VAL	2.3
1	C	39	CYS	2.3
1	C	55	SER	2.3
1	A	255	VAL	2.2
1	D	245	GLU	2.2
1	C	147	LEU	2.2
1	B	285	SER	2.1
1	C	265	CYS	2.1
1	C	51	GLY	2.1
1	B	30	GLU	2.1
1	C	286	LEU	2.0
1	C	300	THR	2.0
1	C	56	PHE	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

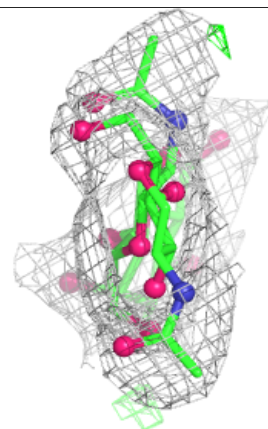
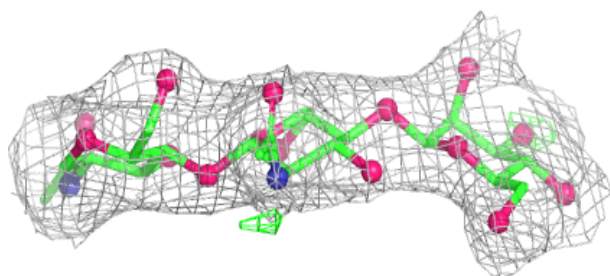
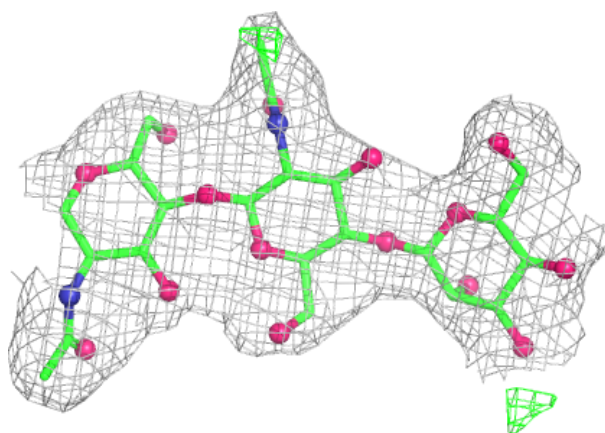
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	BMA	H	3	11/12	0.58	0.13	47,50,63,73	0
2	BMA	E	3	11/12	0.84	0.09	45,53,58,58	0
2	NAG	H	2	14/15	0.85	0.10	45,54,61,62	0
2	NAG	H	1	14/15	0.85	0.10	40,48,53,55	0
2	NAG	E	2	14/15	0.88	0.10	33,48,50,52	0
2	NAG	E	1	14/15	0.94	0.09	31,41,45,47	0
2	NAG	J	1	14/15	-	-	45,64,74,76	0
2	NAG	J	2	14/15	-	-	65,72,77,83	0
2	BMA	J	3	11/12	-	-	63,70,81,82	0
2	NAG	L	1	14/15	-	-	43,61,69,82	0
2	NAG	L	2	14/15	-	-	60,78,80,82	0
2	BMA	L	3	11/12	-	-	59,75,84,85	0

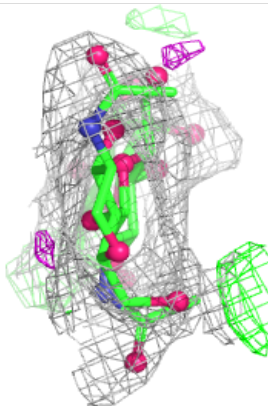
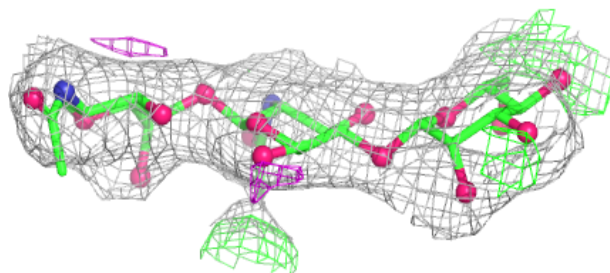
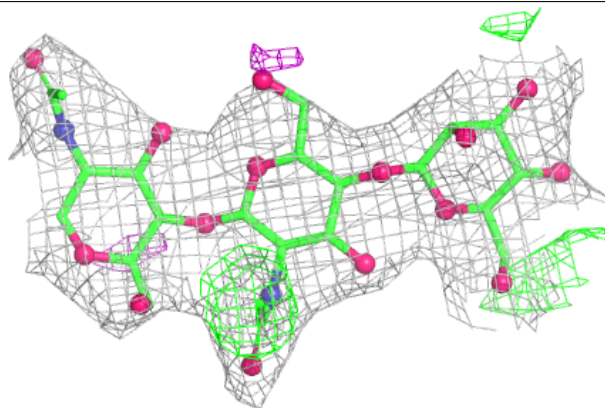
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain E:

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

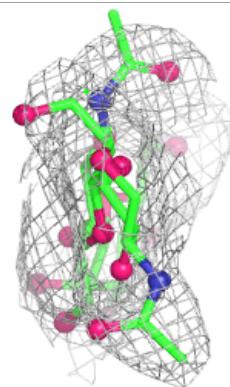
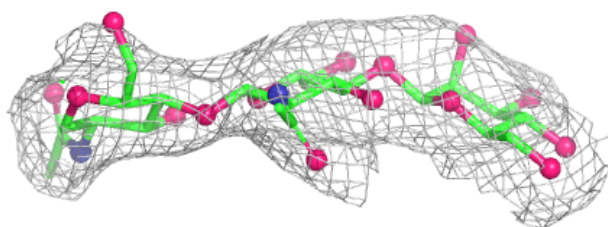
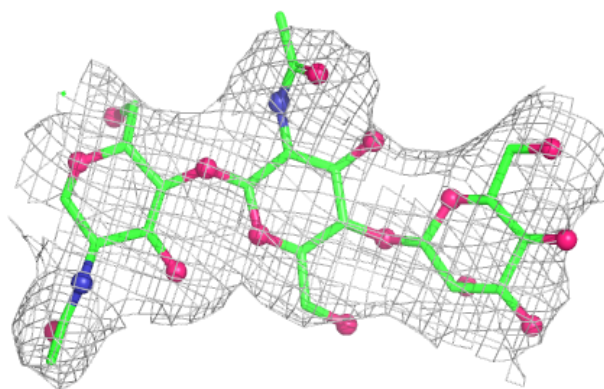
**Electron density around Chain H:**

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

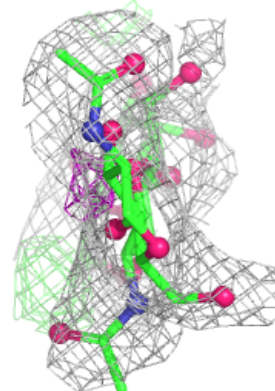
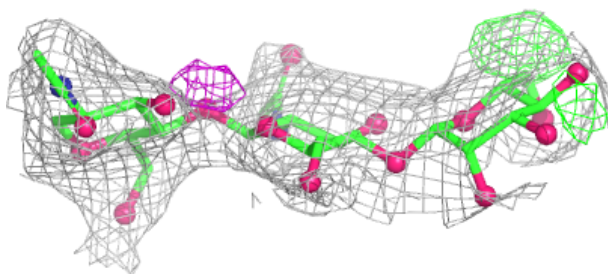
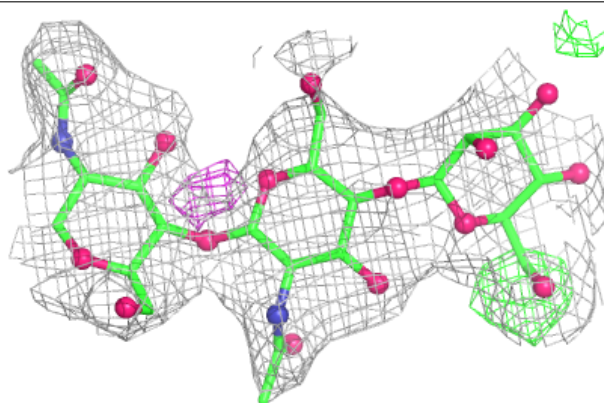


Electron density around Chain J:

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

**Electron density around Chain L:**

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



6.4 Ligands ⓘ

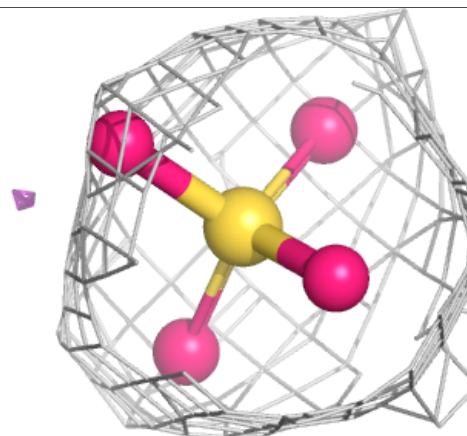
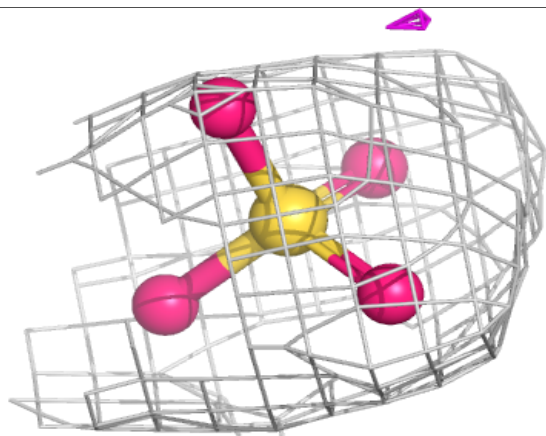
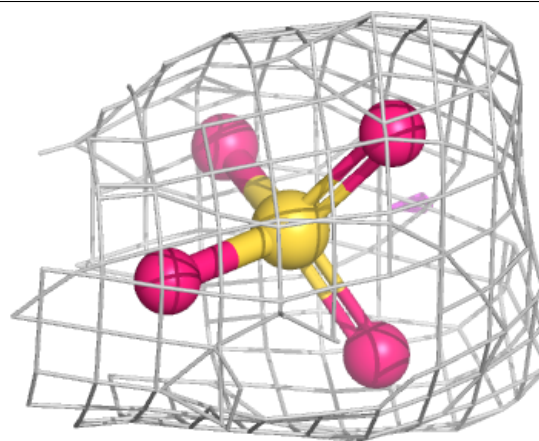
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	SO4	C	403	5/5	0.73	0.14	63,74,100,108	0
4	SO4	C	402	5/5	0.74	0.27	84,102,113,128	0
4	SO4	D	402	5/5	0.78	0.27	90,107,123,130	0
4	SO4	B	402	5/5	0.81	0.18	68,70,94,96	0
3	A1JJO	D	401	53/53	0.84	0.14	46,66,82,86	0
3	A1JJO	C	401	53/53	0.87	0.11	53,64,74,79	0
4	SO4	A	402	5/5	0.88	0.21	74,76,96,103	0
3	A1JJO	A	401	53/53	0.89	0.11	32,47,63,78	0
3	A1JJO	B	401	53/53	0.90	0.11	26,42,56,61	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.

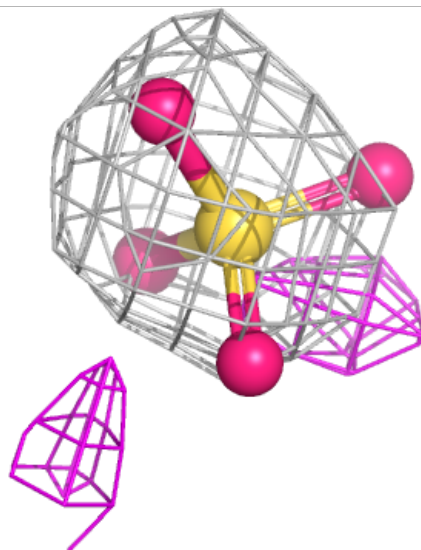
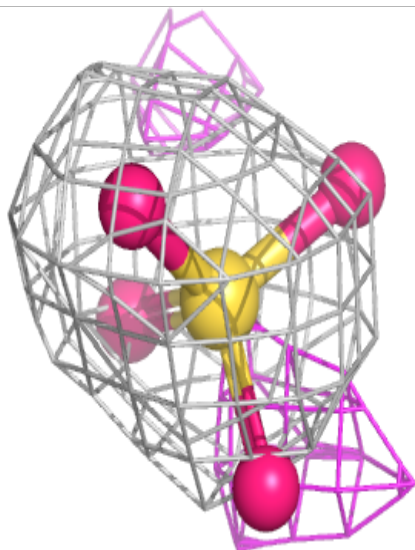
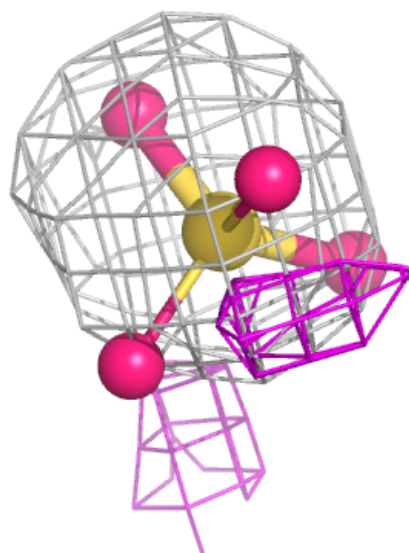
Electron density around SO4 C 403:

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



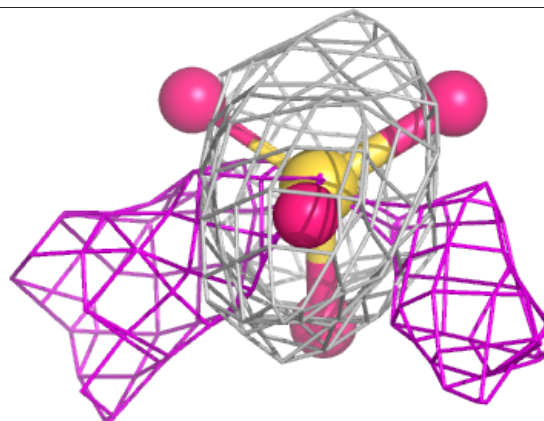
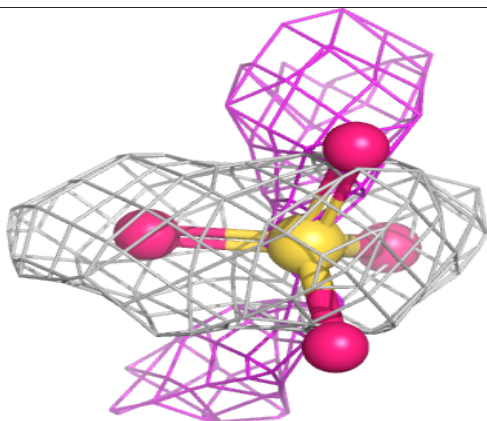
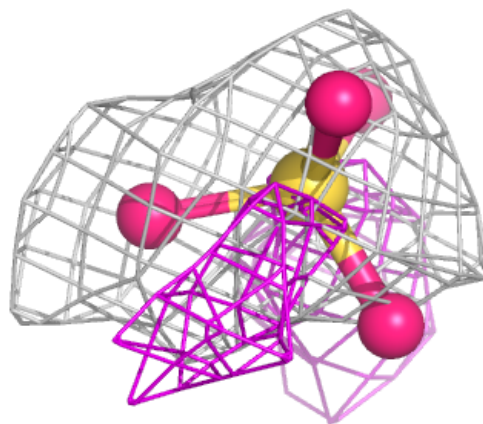
Electron density around SO4 C 402:

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

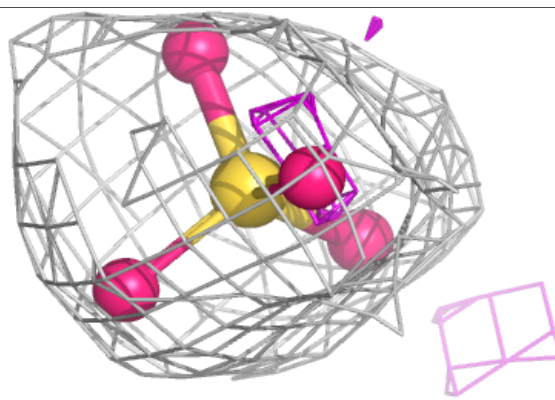
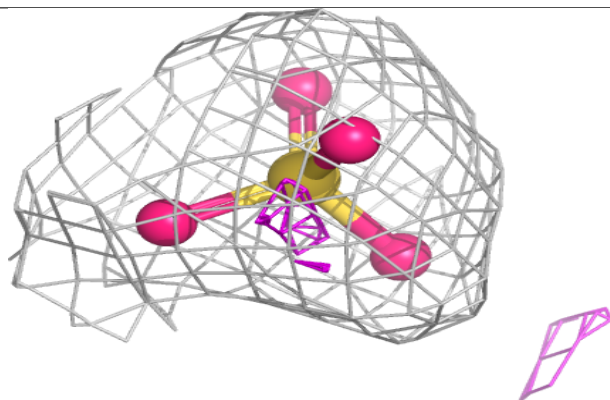
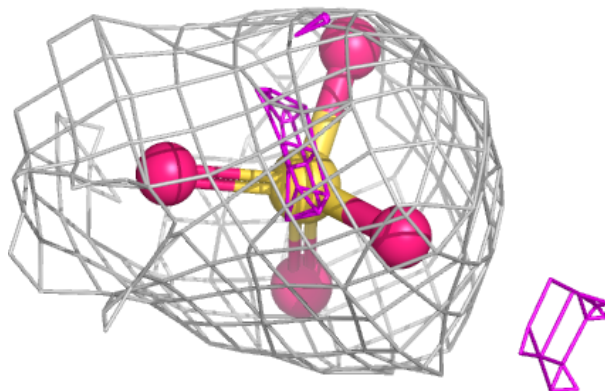


Electron density around SO4 D 402:

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

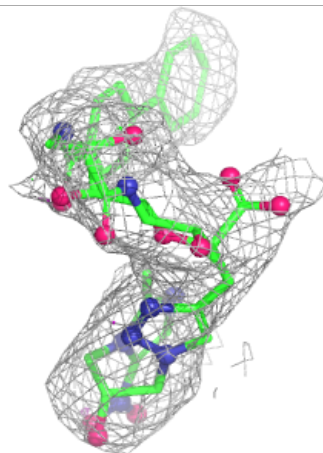
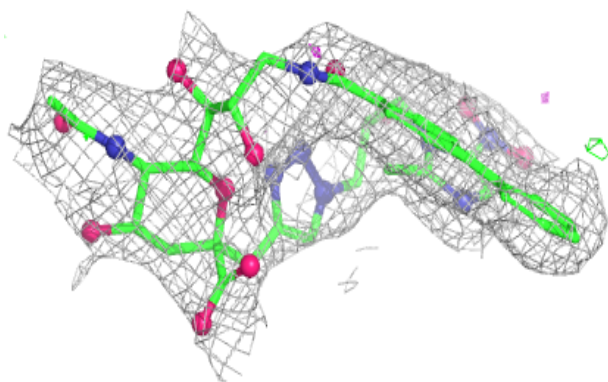
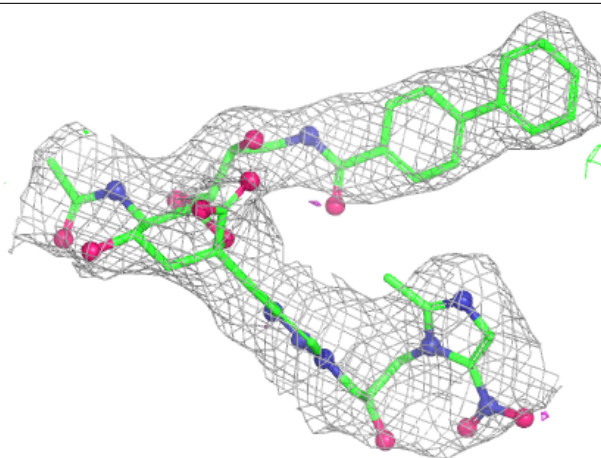
**Electron density around SO4 B 402:**

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



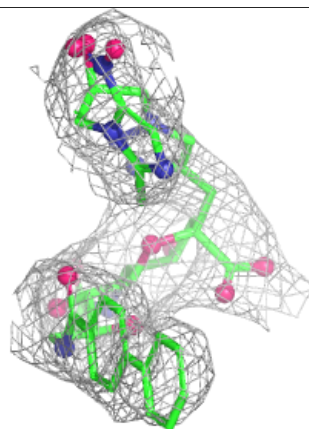
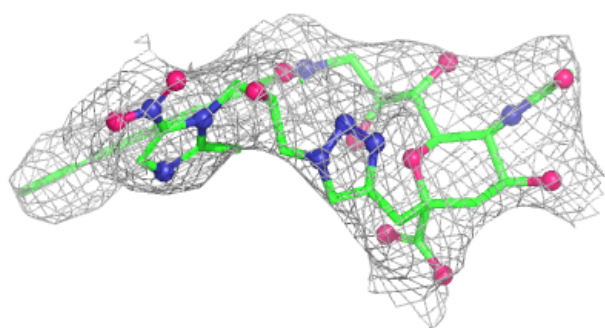
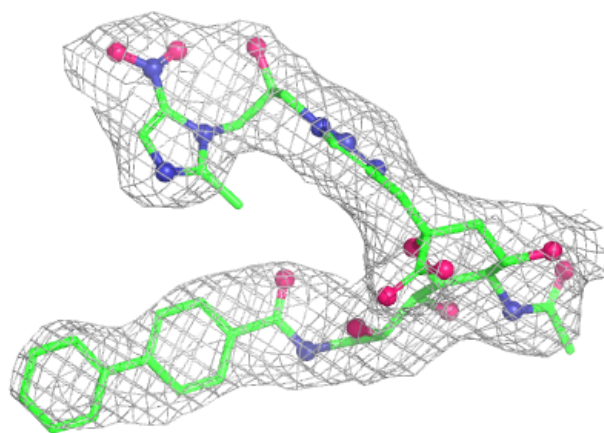
Electron density around A1JJO D 401:

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



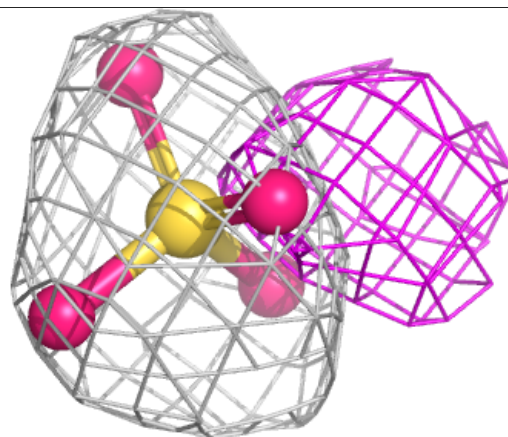
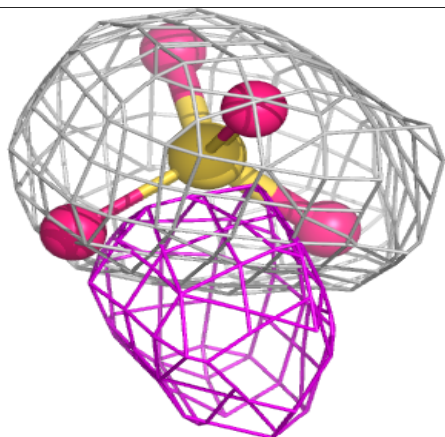
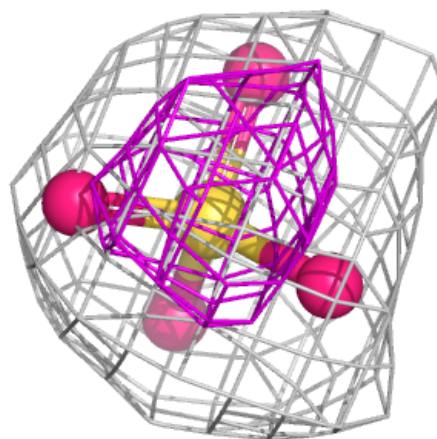
Electron density around A1JJO C 401:

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



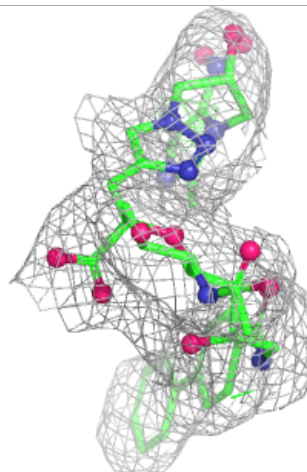
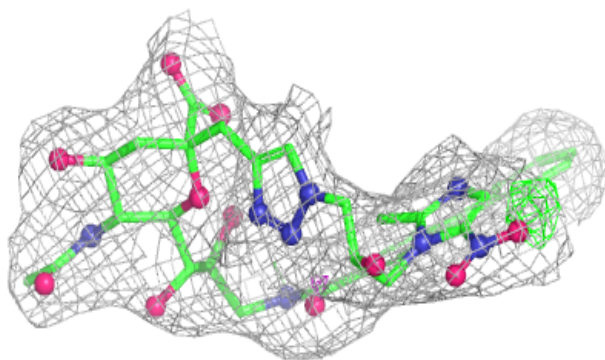
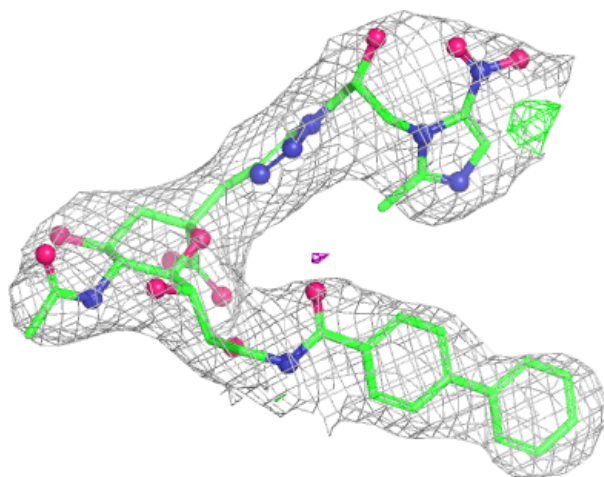
Electron density around SO4 A 402:

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



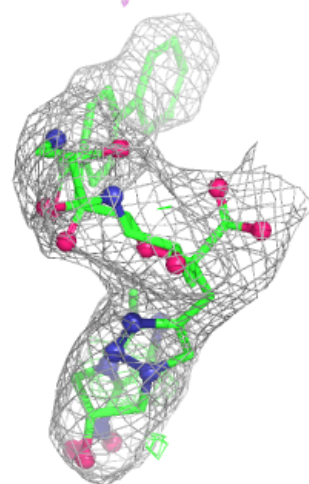
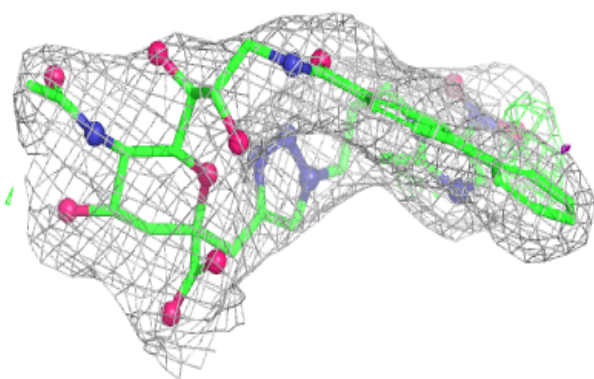
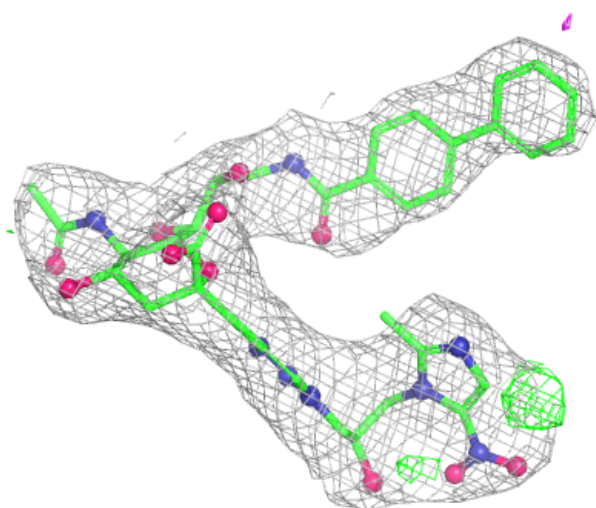
Electron density around A1JJO A 401:

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



Electron density around A1JJO B 401:

$2mF_o-DF_c$ (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.