



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:31 AM UTC

PDB ID : 8G18 / pdb_00008g18
Title : Heterodimer of the GluN1b-GluN2B NMDA receptor amino-terminal domains bound to allosteric inhibitor 93-108
Authors : Regan, M.C.; Furukawa, H.
Deposited on : 2023-02-01
Resolution : 2.85 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

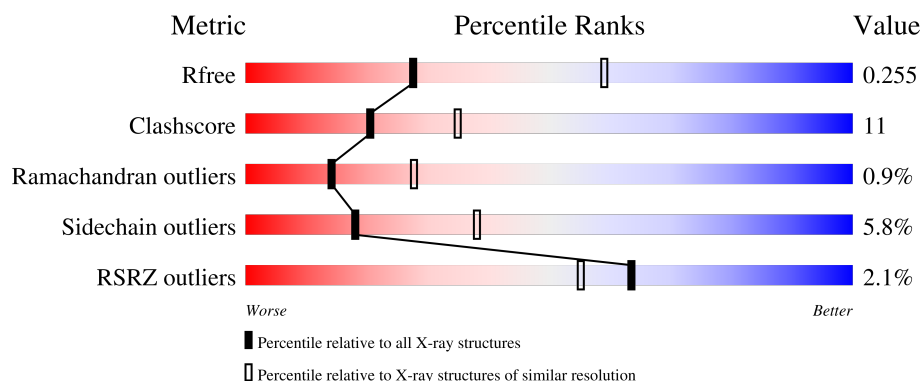
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.85 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	1407 (2.88-2.84)
Clashscore	190562	1446 (2.88-2.84)
Ramachandran outliers	187476	1406 (2.88-2.84)
Sidechain outliers	187428	1407 (2.88-2.84)
RSRZ outliers	180081	1408 (2.88-2.84)

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	385	0.2% 75% 17% 7%
1	C	385	2% 64% 26% 7%
2	B	363	2% 70% 26% •
2	D	363	3% 75% 22% •
3	E	5	40% 60%

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
4	NAG	B	502	X	-	-	-

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 11400 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 Glutamate receptor ionotropic, NMDA 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	359	Total	C	N	O	S	0	0	0
			2765	1763	482	509	11			
1	C	358	Total	C	N	O	S	0	0	0
			2743	1746	477	509	11			

There are 54 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	61	GLN	ASN	conflict	UNP A0A1L8F5J9
A	186	GLY	-	linker	UNP A0A1L8F5J9
A	187	UNK	-	linker	UNP A0A1L8F5J9
A	188	UNK	-	linker	UNP A0A1L8F5J9
A	189	UNK	-	linker	UNP A0A1L8F5J9
A	190	UNK	-	linker	UNP A0A1L8F5J9
A	191	UNK	-	linker	UNP A0A1L8F5J9
A	192	UNK	-	linker	UNP A0A1L8F5J9
A	193	UNK	-	linker	UNP A0A1L8F5J9
A	194	UNK	-	linker	UNP A0A1L8F5J9
A	195	UNK	-	linker	UNP A0A1L8F5J9
A	196	UNK	-	linker	UNP A0A1L8F5J9
A	197	UNK	-	linker	UNP A0A1L8F5J9
A	198	UNK	-	linker	UNP A0A1L8F5J9
A	199	UNK	-	linker	UNP A0A1L8F5J9
A	200	UNK	-	linker	UNP A0A1L8F5J9
A	201	UNK	-	linker	UNP A0A1L8F5J9
A	202	UNK	-	linker	UNP A0A1L8F5J9
A	203	UNK	-	linker	UNP A0A1L8F5J9
A	204	UNK	-	linker	UNP A0A1L8F5J9
A	205	UNK	-	linker	UNP A0A1L8F5J9
A	206	UNK	-	linker	UNP A0A1L8F5J9
A	207	UNK	-	linker	UNP A0A1L8F5J9
A	208	UNK	-	linker	UNP A0A1L8F5J9
A	209	GLY	-	linker	UNP A0A1L8F5J9

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Chain	Residue	Modelled	Actual	Comment	Reference
A	210	PRO	-	linker	UNP A0A1L8F5J9
A	371	GLN	ASN	conflict	UNP A0A1L8F5J9
C	61	GLN	ASN	conflict	UNP A0A1L8F5J9
C	186	GLY	-	linker	UNP A0A1L8F5J9
C	187	UNK	-	linker	UNP A0A1L8F5J9
C	188	UNK	-	linker	UNP A0A1L8F5J9
C	189	UNK	-	linker	UNP A0A1L8F5J9
C	190	UNK	-	linker	UNP A0A1L8F5J9
C	191	UNK	-	linker	UNP A0A1L8F5J9
C	192	UNK	-	linker	UNP A0A1L8F5J9
C	193	UNK	-	linker	UNP A0A1L8F5J9
C	194	UNK	-	linker	UNP A0A1L8F5J9
C	195	UNK	-	linker	UNP A0A1L8F5J9
C	196	UNK	-	linker	UNP A0A1L8F5J9
C	197	UNK	-	linker	UNP A0A1L8F5J9
C	198	UNK	-	linker	UNP A0A1L8F5J9
C	199	UNK	-	linker	UNP A0A1L8F5J9
C	200	UNK	-	linker	UNP A0A1L8F5J9
C	201	UNK	-	linker	UNP A0A1L8F5J9
C	202	UNK	-	linker	UNP A0A1L8F5J9
C	203	UNK	-	linker	UNP A0A1L8F5J9
C	204	UNK	-	linker	UNP A0A1L8F5J9
C	205	UNK	-	linker	UNP A0A1L8F5J9
C	206	UNK	-	linker	UNP A0A1L8F5J9
C	207	UNK	-	linker	UNP A0A1L8F5J9
C	208	UNK	-	linker	UNP A0A1L8F5J9
C	209	GLY	-	linker	UNP A0A1L8F5J9
C	210	PRO	-	linker	UNP A0A1L8F5J9
C	371	GLN	ASN	conflict	UNP A0A1L8F5J9

- Molecule 2 is a protein called Glutamate receptor ionotropic, NMDA 2B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	363	Total	C	N	O	S	0	0	0
			2820	1809	453	542	16			
2	D	363	Total	C	N	O	S	0	0	0
			2757	1771	445	525	16			

There are 2 discrepancies between the modelled and reference sequences:

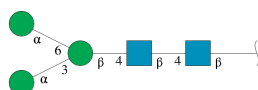
Chain	Residue	Modelled	Actual	Comment	Reference
B	348	ASP	ASN	conflict	UNP Q00960

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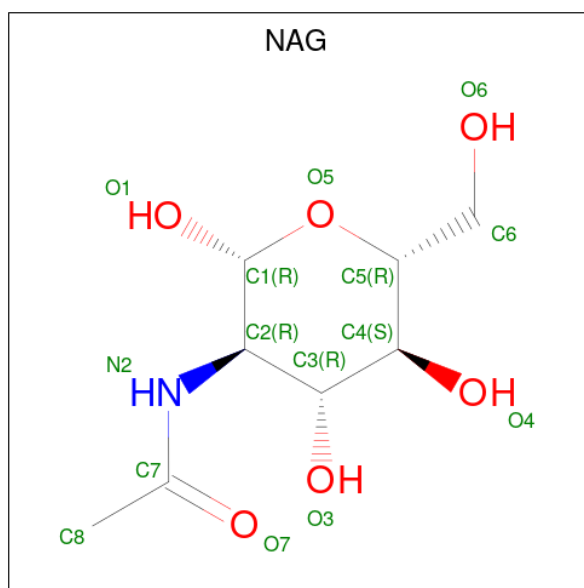
Chain	Residue	Modelled	Actual	Comment	Reference
D	348	ASP	ASN	conflict	UNP Q00960

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



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	E	5	Total	C	N	O	61	0	0
			61	34	2	25			

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆) (labeled as "Ligand of Interest" by depositor).



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

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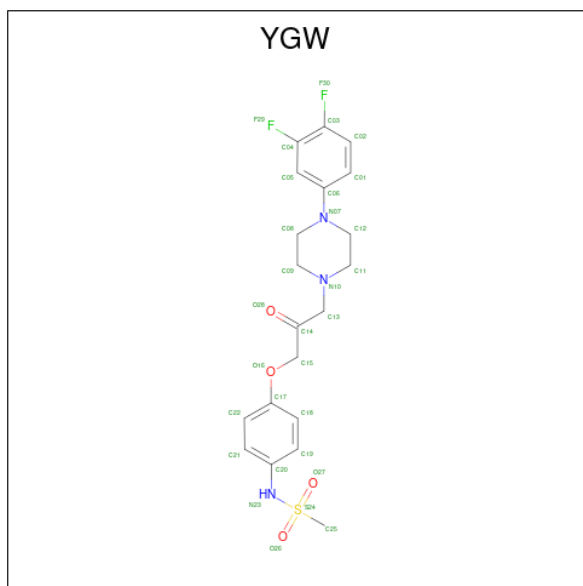
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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	C	1	Total	C	N	O	14	0
			14	8	1	5		
4	D	1	Total	C	N	O	14	0
			14	8	1	5		
4	D	1	Total	C	N	O	14	0
			14	8	1	5		

- Molecule 5 is SODIUM ION (CCD ID: NA) (formula: Na) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Na	1	0
			1	1		
5	C	1	Total	Na	1	0
			1	1		

- Molecule 6 is N-(4-{3-[4-(3,4-difluorophenyl)piperazin-1-yl]-2-oxopropoxy}phenyl)methanesulfonamide (CCD ID: YGW) (formula: C₂₀H₂₃F₂N₃O₄S) (labeled as "Ligand of Interest" by depositor).



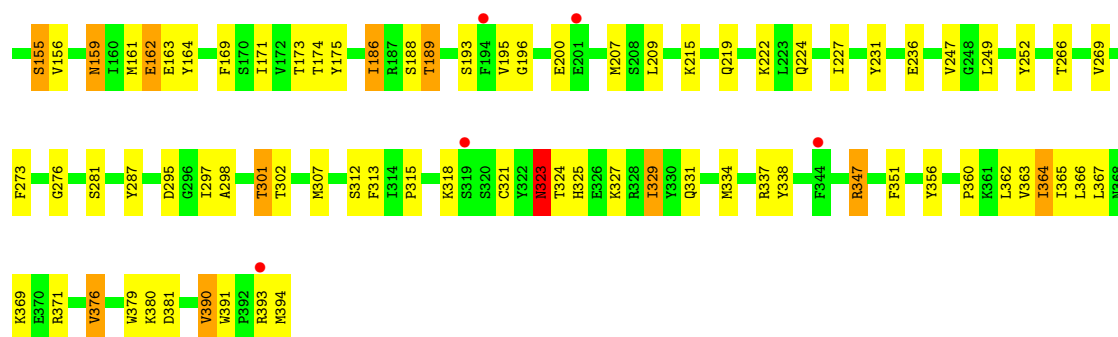
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
6	B	1	Total	C	F	N	O	S	0	0
			30	20	2	3	4	1		
6	D	1	Total	C	F	N	O	S	0	0
			30	20	2	3	4	1		

- Molecule 7 is water.

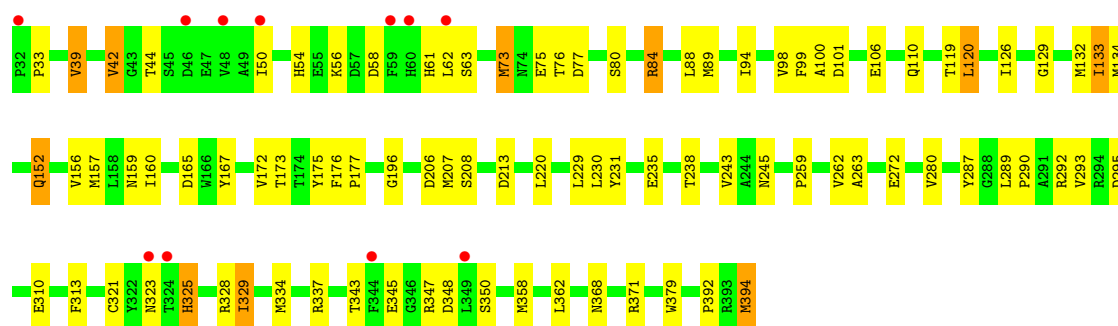
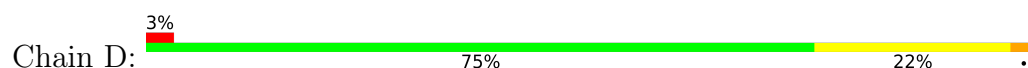
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	36	Total 36	O 36	0	0
7	B	21	Total 21	O 21	0	0
7	C	18	Total 18	O 18	0	0
7	D	19	Total 19	O 19	0	0

- Molecule 1: Glutamate receptor ionotropic, NMDA 1





• Molecule 2: Glutamate receptor ionotropic, NMDA 2B



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



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	269.17Å 60.08Å 145.57Å 90.00° 117.02° 90.00°	Depositor
Resolution (Å)	34.37 – 2.85 34.37 – 2.85	Depositor EDS
% Data completeness (in resolution range)	84.0 (34.37-2.85) 84.0 (34.37-2.85)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.03 (at 2.85Å)	Xtriage
Refinement program	PHENIX 1.20.1-4487	Depositor
R, R_{free}	0.174 , 0.246 (Not available) , 0.255	Depositor DCC
R_{free} test set	2034 reflections (4.15%)	wwPDB-VP
Wilson B-factor (Å ²)	51.2	Xtriage
Anisotropy	0.107	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 73.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	11400	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.48% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MAN, NA, BMA, YGW, NAG

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	0.68	0/2822	1.17	2/3834 (0.1%)
1	C	0.59	0/2799	1.11	3/3805 (0.1%)
2	B	0.62	0/2883	1.13	2/3929 (0.1%)
2	D	0.58	0/2821	1.09	1/3853 (0.0%)
All	All	0.62	0/11325	1.13	8/15421 (0.1%)

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	2

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	106	PRO	CB-CA-C	-7.33	101.44	111.85
1	A	170	ASP	CA-CB-CG	6.89	119.49	112.60
2	D	392	PRO	CB-CA-C	-5.62	103.87	111.12
2	B	104	ASP	CB-CA-C	-5.55	103.43	111.14
1	C	396	ASN	CA-CB-CG	5.37	117.97	112.60
2	B	325	HIS	CA-CB-CG	5.01	118.81	113.80
1	C	230	LEU	CA-C-N	5.01	126.99	120.28
1	C	230	LEU	C-N-CA	5.01	126.99	120.28

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	281	ARG	Sidechain
1	A	398	ARG	Sidechain

5.2 Too-close contacts [i](#)

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	2765	0	2730	37	0
1	C	2743	0	2709	67	3
2	B	2820	0	2683	83	0
2	D	2757	0	2590	64	0
3	E	61	0	52	0	0
4	A	14	0	13	0	0
4	B	28	0	26	0	0
4	C	28	0	26	0	0
4	D	28	0	26	0	0
5	A	1	0	0	0	0
5	C	1	0	0	0	0
6	B	30	0	0	2	0
6	D	30	0	0	3	0
7	A	36	0	0	0	0
7	B	21	0	0	5	3
7	C	18	0	0	3	0
7	D	19	0	0	2	0
All	All	11400	0	10855	245	3

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 11.

All (245) 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)
2:B:61:HIS:HA	7:B:610:HOH:O	1.63	0.97
2:D:343:THR:HG22	2:D:348:ASP:HA	1.55	0.86
2:D:61:HIS:HA	7:D:604:HOH:O	1.77	0.84
1:A:270:VAL:CG1	1:A:274:GLU:HB2	2.08	0.84
2:B:323:ASN:ND2	2:B:323:ASN:O	2.12	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:159:ASN:N	2:B:159:ASN:HD22	1.75	0.82
1:C:270:VAL:CG1	1:C:274:GLU:HB2	2.11	0.81
1:C:270:VAL:HG11	1:C:274:GLU:HB2	1.61	0.81
2:B:163:GLU:O	7:B:601:HOH:O	1.99	0.80
1:C:275:ILE:HD13	1:C:289:GLY:HA3	1.63	0.78
2:D:62:LEU:HA	7:D:602:HOH:O	1.85	0.77
2:B:161:MET:HE3	2:B:227:ILE:HG21	1.66	0.77
1:A:157:LEU:HD23	1:A:158:PHE:CE2	2.21	0.76
2:B:66:PRO:HD3	2:B:301:THR:HG21	1.69	0.75
2:B:173:THR:HG22	2:B:231:TYR:HB3	1.71	0.73
1:C:270:VAL:HG21	1:C:287:ILE:HD11	1.70	0.73
2:D:207:MET:H	6:D:503:YGW:C25	2.01	0.73
2:B:159:ASN:HD22	2:B:159:ASN:H	1.34	0.72
2:D:295:ASP:OD2	2:D:347:ARG:NH1	2.22	0.72
1:A:68:ARG:HG3	1:A:68:ARG:HH11	1.54	0.72
2:B:120:LEU:O	2:B:318:LYS:HE3	1.90	0.72
1:C:227:ALA:HB3	7:C:607:HOH:O	1.91	0.71
2:B:297:ILE:O	2:B:301:THR:OG1	2.08	0.71
2:B:327:LYS:O	2:B:331:GLN:NE2	2.26	0.68
1:C:34:SER:OG	1:C:298:GLU:OE1	2.08	0.66
1:A:285:ASP:OD1	1:A:380:ARG:HG2	1.95	0.66
1:A:270:VAL:HG13	1:A:274:GLU:HB2	1.78	0.65
2:B:351:PHE:HD1	2:B:356:TYR:O	1.80	0.65
2:B:334:MET:HE3	2:B:337:ARG:NH1	2.12	0.64
2:D:280:VAL:HG11	2:D:362:LEU:HD13	1.80	0.64
1:C:224:ASN:ND2	7:C:601:HOH:O	2.26	0.63
2:B:56:LYS:CB	2:B:59:PHE:CB	2.76	0.63
1:C:165:LEU:HD22	1:C:180:LEU:HD23	1.80	0.62
2:D:289:LEU:HB3	2:D:290:PRO:HD3	1.80	0.62
1:C:305:ALA:O	1:C:309:VAL:HG23	1.99	0.62
1:C:378:GLN:HG2	1:C:401:ILE:HG12	1.81	0.62
2:B:175:TYR:HD2	6:B:503:YGW:C25	2.13	0.61
1:A:155:MET:HE3	1:A:163:VAL:HG11	1.82	0.61
1:A:167:VAL:O	1:A:217:GLN:HA	2.00	0.61
2:B:42:VAL:HG11	2:B:76:THR:CG2	2.32	0.60
2:B:161:MET:HE1	2:B:169:PHE:CG	2.36	0.60
1:C:289:GLY:O	1:C:290:LEU:HD23	2.02	0.59
2:D:272:GLU:N	2:D:272:GLU:OE1	2.35	0.59
2:B:287:TYR:OH	2:B:295:ASP:OD2	2.14	0.59
2:B:135:ALA:HB2	2:B:150:ILE:HD11	1.84	0.59
2:B:247:VAL:HG23	2:B:249:LEU:HD12	1.85	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:344:ARG:HG3	1:C:344:ARG:HH11	1.65	0.59
2:D:157:MET:HE3	2:D:229:LEU:HD22	1.86	0.58
2:B:329:ILE:O	2:D:313:PHE:HA	2.04	0.58
1:C:58:ILE:HD11	1:C:314:HIS:ND1	2.18	0.57
1:C:149:LEU:HD23	1:C:183:LEU:HD11	1.86	0.57
2:B:364:ILE:HD13	2:B:364:ILE:N	2.19	0.57
2:D:100:ALA:HB2	2:D:126:ILE:HD11	1.84	0.57
1:C:381:LYS:O	1:C:383:VAL:HG23	2.05	0.57
1:C:42:PHE:HE2	1:C:63:THR:O	1.88	0.56
2:D:175:TYR:HB2	6:D:503:YGW:O27	2.06	0.56
2:B:174:THR:HG21	2:B:236:GLU:OE1	2.05	0.56
1:C:46:VAL:HG21	1:C:62:ALA:HB2	1.87	0.56
2:B:200:GLU:OE1	2:B:224:GLN:HB2	2.06	0.56
1:C:157:LEU:HD23	1:C:158:PHE:CE1	2.41	0.56
2:B:390:VAL:HA	7:B:603:HOH:O	2.05	0.55
1:A:96:PRO:O	1:A:97:ALA:C	2.50	0.55
2:B:159:ASN:OD1	2:B:379:TRP:CZ3	2.59	0.55
2:D:173:THR:HG22	2:D:231:TYR:HB3	1.87	0.55
1:C:138:LEU:HD12	1:C:138:LEU:N	2.22	0.55
2:B:39:VAL:HB	2:B:70:LEU:HD23	1.89	0.55
1:C:72:ILE:HD12	2:D:321:CYS:HB2	1.89	0.55
1:A:68:ARG:HG3	1:A:68:ARG:NH1	2.20	0.55
2:B:318:LYS:HD2	2:B:331:GLN:OE1	2.06	0.55
2:B:159:ASN:H	2:B:159:ASN:ND2	2.04	0.54
2:D:368:ASN:HA	2:D:394:MET:HE2	1.89	0.54
2:D:152:GLN:HE21	2:D:152:GLN:HA	1.71	0.54
1:A:157:LEU:HD23	1:A:158:PHE:CD2	2.42	0.54
1:C:247:ASP:O	1:C:250:THR:HB	2.08	0.54
2:B:164:TYR:CZ	2:B:391:TRP:HZ3	2.26	0.54
2:B:394:MET:HE3	2:B:394:MET:HA	1.90	0.54
2:B:164:TYR:CE1	2:B:391:TRP:CE3	2.95	0.53
2:B:247:VAL:CG2	2:B:249:LEU:HD12	2.38	0.53
1:C:155:MET:HE2	1:C:210:PRO:HB2	1.89	0.53
1:C:304:ASP:OD2	1:C:356:THR:OG1	2.20	0.53
2:B:298:ALA:O	2:B:302:THR:HG23	2.09	0.52
1:C:155:MET:HE2	1:C:210:PRO:CB	2.40	0.52
1:A:250:THR:HG22	1:A:254:LYS:HE3	1.92	0.52
1:A:375:MET:HB3	1:A:382:LEU:HD22	1.92	0.52
2:D:334:MET:HE3	2:D:337:ARG:NH2	2.24	0.52
1:C:387:ILE:O	1:C:387:ILE:HG13	2.11	0.51
2:B:334:MET:HE3	2:B:337:ARG:HH12	1.73	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:307:MET:HE1	2:B:315:PRO:HD3	1.92	0.51
1:A:166:ILE:HD12	1:A:240:ILE:CG2	2.41	0.51
1:C:28:ASN:ND2	1:C:85:SER:O	2.35	0.51
1:C:130:ASP:OD1	1:C:132:SER:HB2	2.11	0.51
2:B:122:PRO:HG3	2:B:142:MET:HE2	1.93	0.51
2:B:159:ASN:OD1	2:B:379:TRP:HZ3	1.94	0.50
2:B:252:TYR:CE1	2:B:393:ARG:HD2	2.46	0.50
1:C:261:MET:HA	1:C:266:TYR:CE2	2.47	0.50
2:D:99:PHE:O	2:D:126:ILE:HG12	2.11	0.50
1:C:264:ALA:HA	1:C:403:PRO:O	2.11	0.49
1:C:311:GLN:NE2	1:C:350:LYS:O	2.41	0.49
2:D:280:VAL:CG1	2:D:362:LEU:HD13	2.41	0.49
2:D:39:VAL:HG13	2:D:98:VAL:HB	1.95	0.49
2:D:54:HIS:C	2:D:56:LYS:N	2.70	0.49
1:C:166:ILE:C	1:C:167:VAL:HG23	2.38	0.49
2:B:207:MET:HB2	6:B:503:YGW:O26	2.12	0.49
2:B:69:GLU:CB	2:B:92:ARG:NH2	2.76	0.49
2:B:366:LEU:HB3	2:B:376:VAL:HG21	1.94	0.49
1:C:152:PHE:CZ	1:C:156:ARG:HD2	2.48	0.49
2:B:276:GLY:HA2	2:B:366:LEU:HD11	1.95	0.49
2:D:343:THR:HG22	2:D:348:ASP:CA	2.35	0.49
1:A:241:ILE:HA	1:A:269:LEU:O	2.13	0.48
2:D:77:ASP:OD2	2:D:80:SER:OG	2.26	0.48
2:D:152:GLN:O	2:D:156:VAL:HG23	2.13	0.48
2:B:323:ASN:ND2	2:B:323:ASN:C	2.71	0.48
1:A:149:LEU:HD23	1:A:183:LEU:HD11	1.95	0.48
1:C:184:LEU:HD22	1:C:210:PRO:O	2.14	0.48
2:D:213:ASP:N	2:D:213:ASP:OD1	2.46	0.48
2:B:215:LYS:O	2:B:219:GLN:HG3	2.13	0.48
2:B:122:PRO:HA	2:B:142:MET:O	2.14	0.47
1:C:222:THR:HB	1:C:225:LEU:HD21	1.95	0.47
1:C:153:GLU:OE2	1:C:156:ARG:NH1	2.47	0.47
2:D:110:GLN:OE1	2:D:134:MET:HE2	2.14	0.47
1:A:57:LYS:HG2	1:A:58:ILE:HG23	1.96	0.47
1:A:155:MET:HE3	1:A:163:VAL:CG1	2.43	0.47
2:D:84:ARG:O	2:D:88:LEU:HG	2.14	0.47
2:B:307:MET:CE	2:B:315:PRO:HD3	2.45	0.47
2:B:369:LYS:H	2:B:394:MET:HE2	1.80	0.47
2:D:207:MET:N	6:D:503:YGW:C25	2.72	0.47
1:A:33:LEU:O	1:A:66:THR:HA	2.14	0.47
2:B:164:TYR:CE1	2:B:391:TRP:CZ3	3.03	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:307:MET:HB2	2:B:338:TYR:CD2	2.50	0.47
1:C:92:VAL:CG2	1:C:107:ILE:HG21	2.45	0.47
1:C:171:HIS:HD1	1:C:171:HIS:C	2.22	0.47
1:A:356:THR:OG1	1:A:367:ARG:NH2	2.48	0.47
2:D:167:TYR:HB3	2:D:196:GLY:O	2.15	0.46
1:A:347:MET:HE2	2:B:209:LEU:HD11	1.97	0.46
2:B:110:GLN:OE1	2:B:134:MET:HE2	2.15	0.46
2:D:39:VAL:HG11	2:D:293:VAL:HG13	1.97	0.46
2:B:171:ILE:HD11	2:B:186:ILE:HG21	1.97	0.46
2:B:307:MET:SD	2:B:315:PRO:HD3	2.55	0.46
1:A:39:GLU:HG2	1:A:64:SER:OG	2.16	0.46
1:C:218:PHE:HB3	1:C:228:LEU:CD1	2.46	0.46
1:A:177:GLN:OE1	1:A:217:GLN:NE2	2.48	0.46
2:D:50:ILE:HG21	2:D:293:VAL:HG11	1.97	0.46
2:D:235:GLU:O	2:D:238:THR:OG1	2.33	0.46
2:B:162:GLU:O	2:B:163:GLU:C	2.58	0.45
1:C:124:ARG:O	1:C:143:PRO:HA	2.16	0.45
2:D:325:HIS:ND1	2:D:325:HIS:N	2.60	0.45
1:C:261:MET:O	1:C:266:TYR:CD2	2.69	0.45
2:D:362:LEU:HB2	2:D:379:TRP:HB3	1.99	0.45
2:B:89:MET:HE1	2:B:119:THR:HB	1.99	0.45
2:D:88:LEU:O	2:D:94:ILE:CD1	2.65	0.45
2:B:66:PRO:CD	2:B:301:THR:HG21	2.43	0.45
1:C:310:ALA:O	1:C:311:GLN:C	2.59	0.45
2:B:164:TYR:CE1	2:B:391:TRP:HE3	2.34	0.45
2:B:379:TRP:CD1	2:B:379:TRP:C	2.94	0.45
1:C:37:LYS:HE3	1:C:41:ILE:HD11	1.99	0.45
1:C:105:THR:N	1:C:106:PRO:CD	2.80	0.45
1:A:104:PRO:HA	1:A:107:ILE:HD12	1.98	0.45
2:D:159:ASN:OD1	2:D:379:TRP:CZ3	2.69	0.45
1:A:270:VAL:HG11	1:A:274:GLU:HB2	1.93	0.45
1:C:58:ILE:HD11	1:C:314:HIS:CE1	2.52	0.44
2:D:42:VAL:HG11	2:D:76:THR:HG23	1.98	0.44
2:D:73:MET:HG3	2:D:84:ARG:CD	2.47	0.44
1:C:344:ARG:HG3	1:C:344:ARG:NH1	2.32	0.44
2:B:318:LYS:CE	7:B:608:HOH:O	2.66	0.44
2:D:73:MET:HG3	2:D:84:ARG:HD3	1.98	0.44
1:A:218:PHE:HB3	1:A:228:LEU:CD1	2.47	0.44
1:A:42:PHE:CE2	1:A:62:ALA:HB1	2.52	0.44
1:A:226:THR:HG22	1:A:259:LEU:HD21	2.00	0.44
2:B:318:LYS:HE2	7:B:608:HOH:O	2.16	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:27:VAL:HG13	1:A:88:TYR:CD1	2.52	0.44
1:A:302:ILE:O	1:A:306:VAL:HG23	2.17	0.44
2:B:321:CYS:HA	2:B:324:THR:CG2	2.48	0.44
2:D:347:ARG:CB	2:D:347:ARG:HH21	2.31	0.44
2:D:159:ASN:OD1	2:D:379:TRP:HZ3	2.00	0.44
1:C:69:PRO:HG2	1:C:73:GLN:NE2	2.32	0.44
2:D:156:VAL:HG21	2:D:362:LEU:HD12	2.00	0.44
2:B:334:MET:HE1	2:D:334:MET:HE1	2.00	0.43
2:B:71:VAL:HG21	2:B:88:LEU:HD11	1.99	0.43
2:B:222:LYS:HE2	2:B:222:LYS:HB3	1.76	0.43
1:C:155:MET:HE3	1:C:163:VAL:HG11	2.00	0.43
1:C:259:LEU:N	1:C:259:LEU:HD23	2.34	0.43
2:D:220:LEU:HD12	2:D:243:VAL:HG12	2.01	0.43
2:B:164:TYR:N	2:B:164:TYR:CD2	2.85	0.43
2:B:193:SER:O	2:B:195:VAL:O	2.36	0.43
1:C:124:ARG:NH1	7:C:604:HOH:O	2.52	0.43
1:A:402:TRP:HD1	1:A:406:GLU:CB	2.31	0.43
2:B:269:VAL:HG11	2:B:367:LEU:HD11	2.01	0.43
2:D:89:MET:HE1	2:D:119:THR:HB	2.01	0.43
2:B:186:ILE:O	2:B:189:THR:N	2.48	0.43
1:C:405:GLY:O	1:C:406:GLU:C	2.61	0.43
1:C:165:LEU:HG	1:C:167:VAL:HG23	2.01	0.42
1:C:228:LEU:C	1:C:228:LEU:HD23	2.44	0.42
1:C:244:ALA:O	1:C:274:GLU:HG3	2.18	0.42
1:C:234:GLU:O	1:C:235:LEU:C	2.62	0.42
2:B:380:LYS:O	2:B:381:ASP:C	2.61	0.42
2:B:360:PRO:HG2	2:B:362:LEU:HD21	2.02	0.42
2:D:33:PRO:HD2	2:D:63:SER:O	2.20	0.42
2:D:287:TYR:CE2	2:D:292:ARG:HG2	2.54	0.42
2:D:88:LEU:O	2:D:94:ILE:HD13	2.19	0.42
1:A:293:ILE:O	1:A:294:ASN:HB2	2.20	0.42
2:D:119:THR:O	2:D:120:LEU:C	2.63	0.42
2:D:172:VAL:O	2:D:230:LEU:HA	2.20	0.42
1:C:23:ASP:N	1:C:24:PRO:CD	2.83	0.42
1:C:293:ILE:O	1:C:294:ASN:HB2	2.19	0.42
2:D:230:LEU:O	2:D:259:PRO:HD3	2.19	0.42
2:B:273:PHE:O	2:B:371:ARG:NH1	2.47	0.42
2:B:313:PHE:HA	2:D:329:ILE:O	2.18	0.42
1:C:308:VAL:HG21	1:C:359:ILE:HG21	2.01	0.42
2:D:101:ASP:OD1	2:D:129:GLY:N	2.50	0.42
2:B:155:SER:O	2:B:156:VAL:C	2.62	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:159:ASN:O	2:D:160:ILE:C	2.62	0.42
1:A:269:LEU:C	1:A:270:VAL:HG23	2.45	0.41
2:B:77:ASP:HB2	2:B:78:PRO:HD2	2.01	0.41
2:B:195:VAL:HG12	2:B:196:GLY:N	2.35	0.41
1:C:171:HIS:C	1:C:171:HIS:ND1	2.78	0.41
2:D:54:HIS:C	2:D:56:LYS:H	2.28	0.41
1:A:400:ILE:HG21	1:A:402:TRP:CZ2	2.55	0.41
2:B:376:VAL:HG12	2:B:376:VAL:O	2.20	0.41
2:D:106:GLU:HB3	2:D:132:MET:SD	2.60	0.41
1:A:156:ARG:O	1:A:157:LEU:C	2.64	0.41
1:C:67:HIS:HE1	1:C:93:SER:OG	2.03	0.41
2:D:176:PHE:O	2:D:177:PRO:C	2.62	0.41
2:B:110:GLN:HA	2:B:134:MET:HE3	2.01	0.41
2:D:165:ASP:HA	2:D:167:TYR:CZ	2.54	0.41
2:B:363:VAL:HG12	2:B:365:ILE:HG23	2.02	0.41
2:B:39:VAL:HB	2:B:70:LEU:CD2	2.51	0.41
1:C:155:MET:HE3	1:C:163:VAL:CG1	2.51	0.41
1:A:338:THR:O	1:A:338:THR:HG22	2.20	0.41
2:D:132:MET:O	2:D:133:ILE:C	2.63	0.41
1:A:339:GLY:N	1:A:340:PRO:CD	2.83	0.41
2:B:73:MET:HE1	2:B:81:ILE:HG12	2.03	0.41
1:A:226:THR:CG2	1:A:259:LEU:HD21	2.51	0.41
2:B:347:ARG:NH1	2:B:347:ARG:HB3	2.35	0.41
1:C:103:THR:HB	1:C:104:PRO:HD2	2.03	0.41
1:C:118:VAL:C	1:C:119:ILE:HD13	2.46	0.41
1:C:331:GLY:HA2	2:D:75:GLU:OE1	2.20	0.41
2:D:50:ILE:CG2	2:D:293:VAL:HG11	2.51	0.41
1:C:96:PRO:HA	1:C:97:ALA:HA	1.84	0.41
2:D:262:VAL:O	2:D:263:ALA:C	2.64	0.41
1:C:269:LEU:C	1:C:270:VAL:HG23	2.46	0.40
1:C:63:THR:O	1:C:64:SER:HB3	2.20	0.40
2:D:206:ASP:OD1	2:D:208:SER:OG	2.35	0.40
1:C:226:THR:HG23	1:C:255:SER:HB3	2.04	0.40
2:B:156:VAL:HG21	2:B:362:LEU:HD12	2.03	0.40
2:D:371:ARG:HA	2:D:371:ARG:HD2	1.97	0.40

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:380:ARG:NE	7:B:601:HOH:O[4_7410]	2.08	0.12

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:380:ARG:NH2	7:B:601:HOH:O[4_7410]	2.18	0.02
1:C:380:ARG:CB	7:B:603:HOH:O[4_7410]	2.19	0.01

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	353/385 (92%)	329 (93%)	22 (6%)	2 (1%)	21	38
1	C	352/385 (91%)	319 (91%)	31 (9%)	2 (1%)	21	38
2	B	361/363 (99%)	312 (86%)	44 (12%)	5 (1%)	9	19
2	D	361/363 (99%)	317 (88%)	40 (11%)	4 (1%)	11	24
All	All	1427/1496 (95%)	1277 (90%)	137 (10%)	13 (1%)	14	28

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	46	ASP
2	D	44	THR
2	D	323	ASN
1	A	405	GLY
1	A	406	GLU
1	C	405	GLY
2	B	120	LEU
2	B	323	ASN
2	B	186	ILE
2	D	120	LEU
2	D	345	GLU
2	B	376	VAL
1	C	167	VAL

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	289/309 (94%)	275 (95%)	14 (5%)	23	46
1	C	290/309 (94%)	273 (94%)	17 (6%)	18	37
2	B	300/326 (92%)	278 (93%)	22 (7%)	13	27
2	D	286/326 (88%)	271 (95%)	15 (5%)	21	43
All	All	1165/1270 (92%)	1097 (94%)	68 (6%)	18	38

All (68) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	65	VAL
1	A	68	ARG
1	A	119	ILE
1	A	125	MET
1	A	135	LEU
1	A	226	THR
1	A	238	ARG
1	A	285	ASP
1	A	347	MET
1	A	358	ARG
1	A	363	GLU
1	A	383	VAL
1	A	391	SER
1	A	401	ILE
2	B	44	THR
2	B	45	SER
2	B	61	HIS
2	B	64	VAL
2	B	68	VAL
2	B	73	MET
2	B	94	ILE
2	B	133	ILE
2	B	155	SER
2	B	159	ASN

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Mol	Chain	Res	Type
2	B	162	GLU
2	B	188	SER
2	B	189	THR
2	B	266	THR
2	B	281	SER
2	B	301	THR
2	B	312	SER
2	B	323	ASN
2	B	329	ILE
2	B	347	ARG
2	B	364	ILE
2	B	390	VAL
1	C	68	ARG
1	C	102	LEU
1	C	131	LYS
1	C	135	LEU
1	C	136	SER
1	C	153	GLU
1	C	214	LYS
1	C	226	THR
1	C	258	MET
1	C	259	LEU
1	C	318	GLU
1	C	363	GLU
1	C	373	SER
1	C	380	ARG
1	C	387	ILE
1	C	391	SER
1	C	396	ASN
2	D	39	VAL
2	D	42	VAL
2	D	58	ASP
2	D	73	MET
2	D	84	ARG
2	D	133	ILE
2	D	152	GLN
2	D	245	ASN
2	D	310	GLU
2	D	325	HIS
2	D	328	ARG
2	D	329	ILE
2	D	350	SER

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Mol	Chain	Res	Type
2	D	358	MET
2	D	394	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	53	HIS
1	A	67	HIS
1	A	73	GLN
1	A	161	ASN
1	A	177	GLN
1	A	217	GLN
1	A	321	ASN
1	A	371	GLN
2	B	60	HIS
2	B	159	ASN
2	B	192	ASN
2	B	245	ASN
2	B	357	GLN
2	B	385	GLN
1	C	67	HIS
1	C	73	GLN
1	C	101	HIS
1	C	146	HIS
1	C	301	HIS
1	C	314	HIS
1	C	395	GLN
2	D	152	GLN
2	D	159	ASN
2	D	217	GLN
2	D	331	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

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	E	1	3	14,14,15	0.48	0	17,19,21	0.85	1 (5%)
3	NAG	E	2	3	14,14,15	0.32	0	17,19,21	1.11	2 (11%)
3	BMA	E	3	3	11,11,12	0.32	0	15,15,17	0.78	0
3	MAN	E	4	3	11,11,12	0.50	0	15,15,17	1.76	3 (20%)
3	MAN	E	5	3	11,11,12	0.49	0	15,15,17	0.87	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	NAG	E	1	3	-	0/6/23/26	0/1/1/1
3	NAG	E	2	3	-	0/6/23/26	0/1/1/1
3	BMA	E	3	3	-	0/2/19/22	0/1/1/1
3	MAN	E	4	3	-	0/2/19/22	0/1/1/1
3	MAN	E	5	3	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	4	MAN	C1-C2-C3	5.07	117.03	109.64
3	E	4	MAN	C2-C3-C4	3.15	116.41	110.86
3	E	2	NAG	C1-C2-N2	-2.42	106.62	110.43
3	E	2	NAG	C2-N2-C7	2.30	125.98	122.90
3	E	1	NAG	O4-C4-C3	-2.14	105.33	110.38
3	E	4	MAN	C3-C4-C5	2.12	114.08	110.23

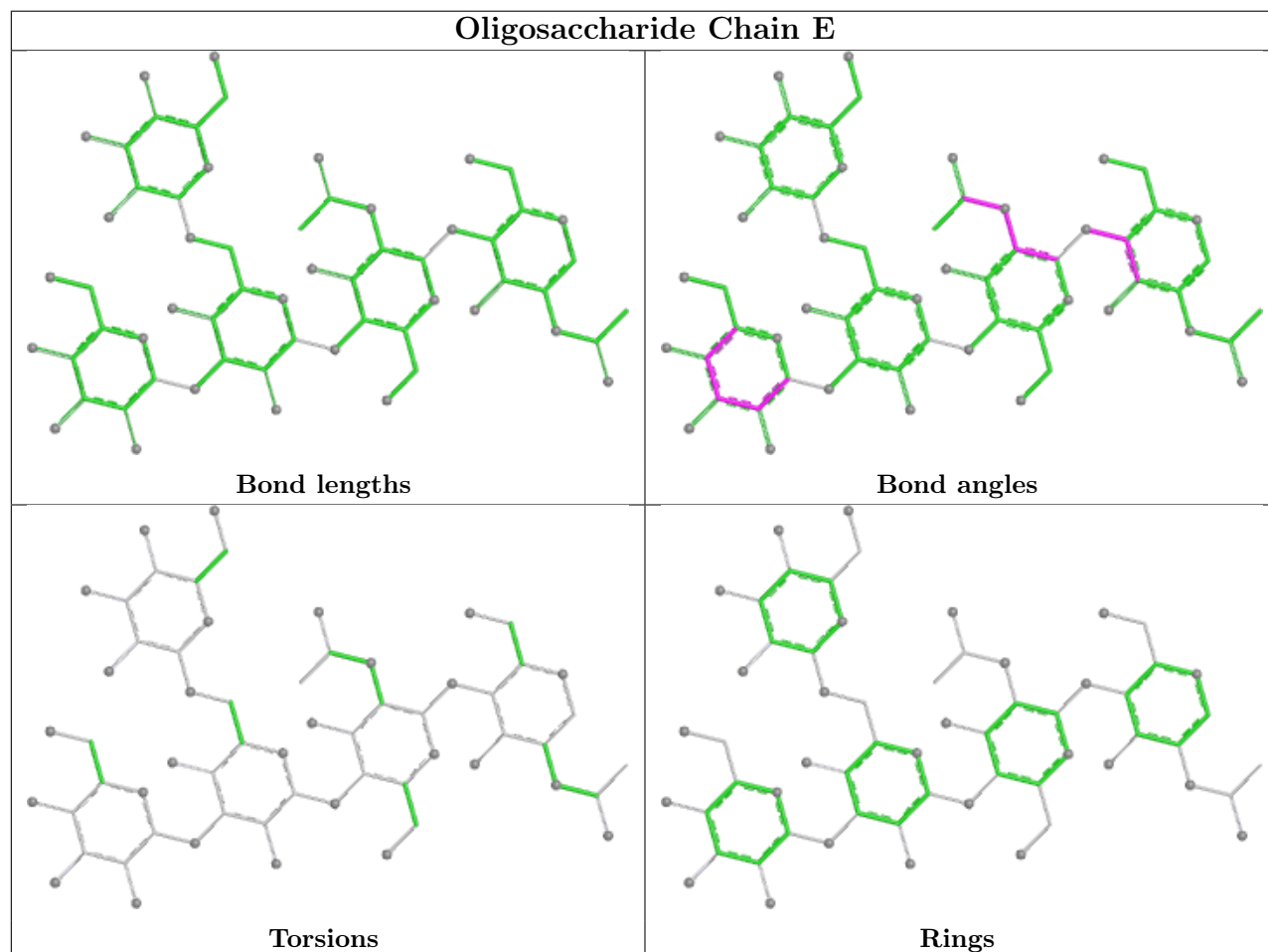
There are no chirality outliers.

There are no torsion outliers.

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 11 ligands modelled in this entry, 2 are monoatomic - leaving 9 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	NAG	C	502	1	14,14,15	0.54	0	17,19,21	1.39	4 (23%)
4	NAG	C	501	1	14,14,15	0.46	0	17,19,21	0.62	0
4	NAG	D	501	-	14,14,15	0.55	0	17,19,21	1.15	1 (5%)
4	NAG	D	502	-	14,14,15	0.29	0	17,19,21	2.41	2 (11%)
4	NAG	B	502	2	14,14,15	0.43	0	17,19,21	2.43	5 (29%)
4	NAG	A	501	1	14,14,15	0.74	1 (7%)	17,19,21	1.53	4 (23%)
6	YGW	D	503	-	32,32,32	3.51	10 (31%)	43,45,45	5.48	14 (32%)
6	YGW	B	503	-	32,32,32	3.44	8 (25%)	43,45,45	5.39	11 (25%)
4	NAG	B	501	2	14,14,15	0.52	0	17,19,21	1.03	1 (5%)

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	NAG	C	502	1	-	4/6/23/26	0/1/1/1
4	NAG	C	501	1	-	0/6/23/26	0/1/1/1
4	NAG	D	501	-	-	3/6/23/26	0/1/1/1
4	NAG	D	502	-	-	3/6/23/26	0/1/1/1
4	NAG	B	502	2	1/1/5/7	4/6/23/26	0/1/1/1
4	NAG	A	501	1	-	2/6/23/26	0/1/1/1
6	YGW	D	503	-	-	10/18/28/28	0/3/3/3
6	YGW	B	503	-	-	4/18/28/28	0/3/3/3
4	NAG	B	501	2	-	0/6/23/26	0/1/1/1

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	503	YGW	O27-S24	14.41	1.67	1.43
6	D	503	YGW	O27-S24	13.88	1.66	1.43
6	D	503	YGW	S24-N23	8.63	1.74	1.63
6	B	503	YGW	C13-N10	-7.88	1.30	1.47
6	B	503	YGW	S24-N23	7.35	1.72	1.63
6	D	503	YGW	C13-N10	-7.04	1.31	1.47
6	D	503	YGW	C13-C14	4.50	1.56	1.51
6	B	503	YGW	C11-N10	-3.44	1.37	1.46
6	D	503	YGW	C11-N10	-3.19	1.38	1.46
6	D	503	YGW	C06-N07	3.01	1.47	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	D	503	YGW	C15-C14	2.75	1.55	1.50
6	D	503	YGW	C09-N10	-2.63	1.39	1.46
6	B	503	YGW	C13-C14	2.59	1.54	1.51
6	B	503	YGW	C25-S24	2.51	1.81	1.75
6	D	503	YGW	C25-S24	2.47	1.81	1.75
6	D	503	YGW	O26-S24	2.47	1.47	1.43
6	B	503	YGW	C09-N10	-2.22	1.40	1.46
4	A	501	NAG	O5-C1	-2.13	1.40	1.43
6	B	503	YGW	C05-C06	-2.01	1.36	1.39

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	B	503	YGW	C25-S24-N23	28.86	139.05	106.56
6	D	503	YGW	C25-S24-N23	24.45	134.08	106.56
6	D	503	YGW	O27-S24-C25	-20.55	75.74	108.26
6	B	503	YGW	O27-S24-C25	-16.12	82.75	108.26
6	D	503	YGW	O27-S24-N23	-9.29	88.54	107.10
6	D	503	YGW	O27-S24-O26	-7.99	108.16	118.87
4	D	502	NAG	C1-O5-C5	7.95	122.84	112.19
4	B	502	NAG	C1-O5-C5	7.63	122.42	112.19
6	B	503	YGW	O27-S24-N23	-7.52	92.08	107.10
6	B	503	YGW	O27-S24-O26	-5.44	111.57	118.87
4	D	502	NAG	O5-C1-C2	4.70	118.57	111.29
6	D	503	YGW	O26-S24-N23	4.08	115.24	107.10
4	D	501	NAG	C1-O5-C5	4.08	117.65	112.19
6	D	503	YGW	C15-O16-C17	4.03	124.85	117.63
4	B	502	NAG	C3-C4-C5	3.44	116.47	110.23
4	B	501	NAG	O5-C1-C2	-3.31	106.18	111.29
4	A	501	NAG	C1-O5-C5	3.11	116.35	112.19
4	B	502	NAG	O5-C5-C4	3.07	118.29	110.83
6	D	503	YGW	C08-N07-C06	3.01	126.31	118.11
6	D	503	YGW	C13-N10-C09	2.92	115.69	111.14
6	B	503	YGW	C01-C06-N07	-2.90	117.37	121.39
6	B	503	YGW	C12-N07-C08	2.86	118.00	111.57
4	C	502	NAG	O3-C3-C2	-2.75	103.69	109.40
6	D	503	YGW	C09-C08-N07	2.67	116.40	110.78
6	D	503	YGW	C22-C21-C20	-2.62	117.29	120.30
6	B	503	YGW	O26-S24-N23	2.60	112.29	107.10
4	C	502	NAG	O6-C6-C5	-2.55	102.65	111.33
4	B	502	NAG	C2-N2-C7	2.52	126.28	122.90
6	D	503	YGW	C05-C06-N07	-2.51	118.62	121.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	501	NAG	C3-C4-C5	2.47	114.72	110.23
6	B	503	YGW	C15-O16-C17	-2.45	113.25	117.63
6	B	503	YGW	O28-C14-C15	-2.43	116.57	120.77
4	C	502	NAG	C3-C4-C5	2.39	114.56	110.23
6	D	503	YGW	C21-C20-C19	2.22	121.98	119.04
4	A	501	NAG	O3-C3-C4	-2.18	105.24	110.38
4	B	502	NAG	C8-C7-N2	2.18	119.73	116.12
6	B	503	YGW	C08-N07-C06	2.15	123.96	118.11
6	B	503	YGW	C11-C12-N07	2.14	115.29	110.78
6	D	503	YGW	C19-C18-C17	-2.13	117.30	119.73
6	D	503	YGW	C19-C20-N23	-2.11	115.58	120.02
4	C	502	NAG	O5-C1-C2	-2.05	108.13	111.29
4	A	501	NAG	C6-C5-C4	-2.00	108.10	113.02

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
4	B	502	NAG	C1

All (30) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	B	503	YGW	C20-N23-S24-O26
6	D	503	YGW	N10-C13-C14-C15
6	D	503	YGW	N10-C13-C14-O28
6	D	503	YGW	C20-N23-S24-O27
4	B	502	NAG	O5-C5-C6-O6
4	D	501	NAG	C4-C5-C6-O6
4	B	502	NAG	C4-C5-C6-O6
4	A	501	NAG	C4-C5-C6-O6
4	D	501	NAG	O5-C5-C6-O6
4	B	502	NAG	C8-C7-N2-C2
4	B	502	NAG	O7-C7-N2-C2
4	C	502	NAG	C8-C7-N2-C2
4	C	502	NAG	O7-C7-N2-C2
4	D	502	NAG	C8-C7-N2-C2
4	D	502	NAG	O7-C7-N2-C2
4	A	501	NAG	O5-C5-C6-O6
6	D	503	YGW	C20-N23-S24-C25
6	D	503	YGW	C22-C17-O16-C15
6	D	503	YGW	C18-C17-O16-C15
6	D	503	YGW	C20-N23-S24-O26

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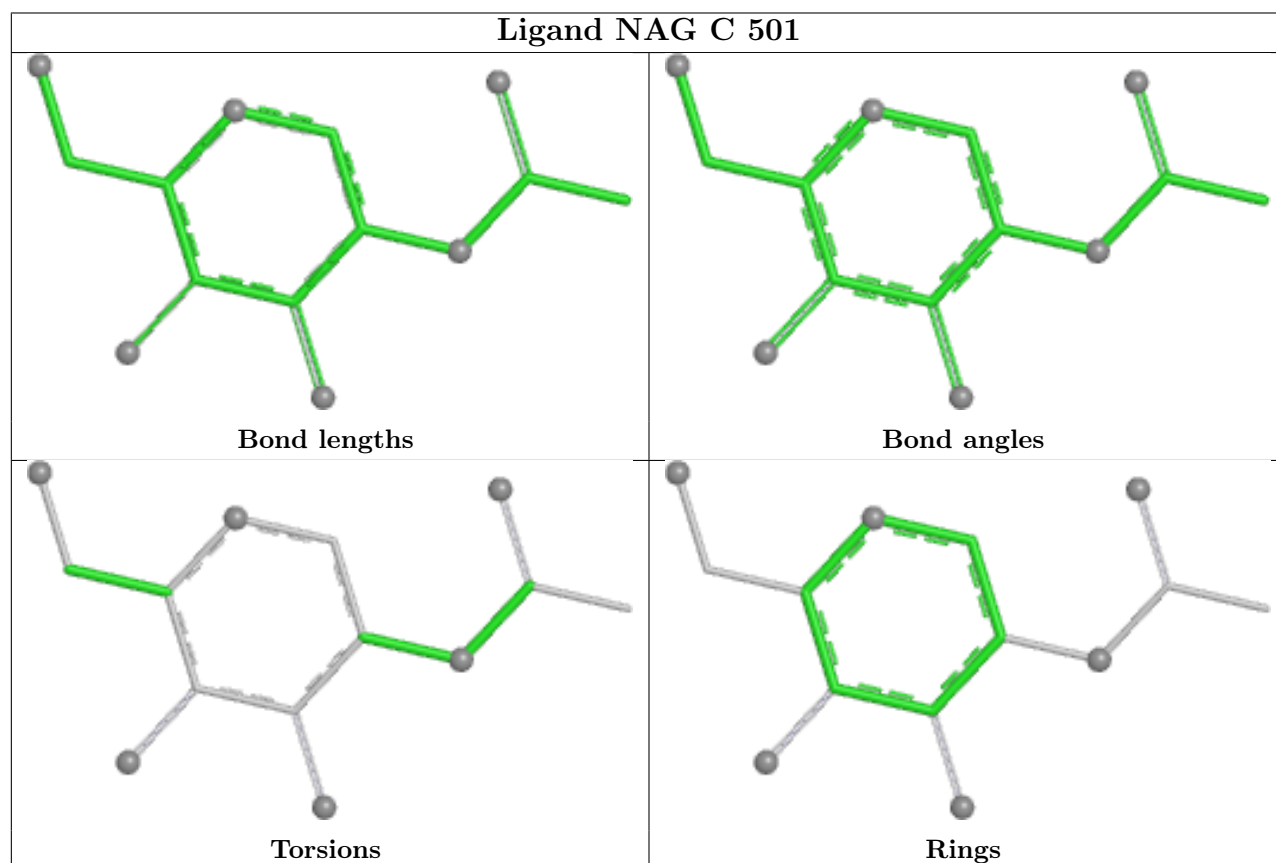
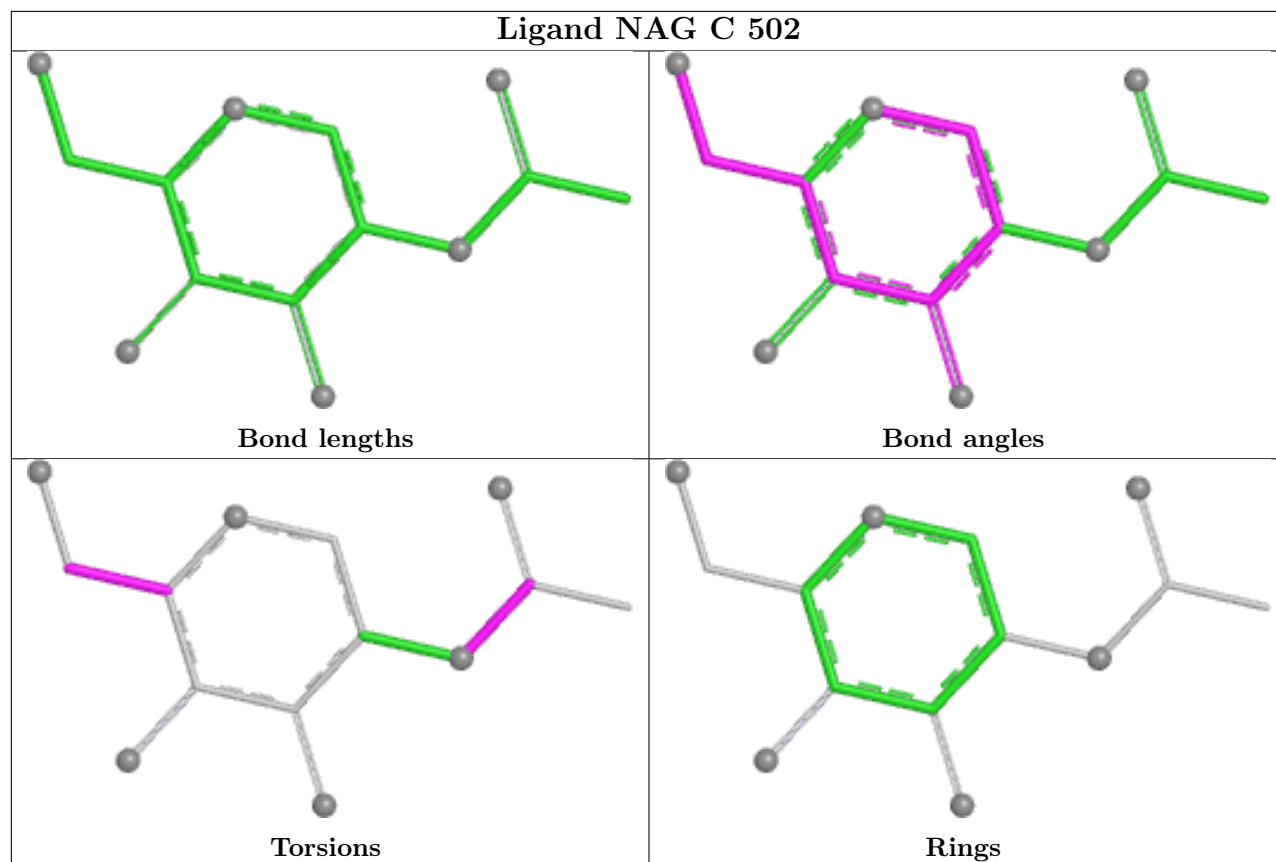
Mol	Chain	Res	Type	Atoms
4	C	502	NAG	C4-C5-C6-O6
6	B	503	YGW	C20-N23-S24-O27
4	D	502	NAG	C4-C5-C6-O6
6	D	503	YGW	O28-C14-C15-O16
4	C	502	NAG	O5-C5-C6-O6
4	D	501	NAG	C3-C2-N2-C7
6	B	503	YGW	O28-C14-C15-O16
6	B	503	YGW	N10-C13-C14-O28
6	D	503	YGW	C14-C13-N10-C09
6	D	503	YGW	C14-C15-O16-C17

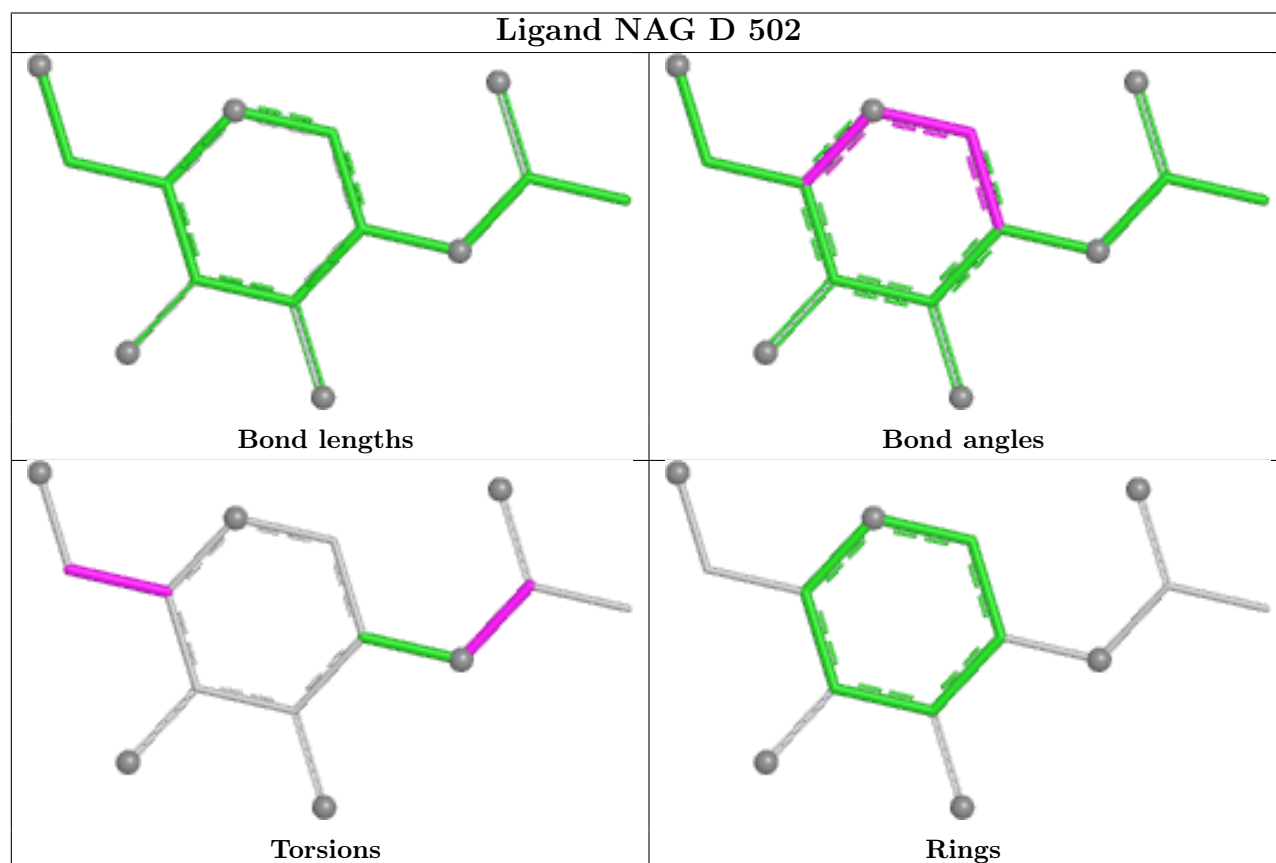
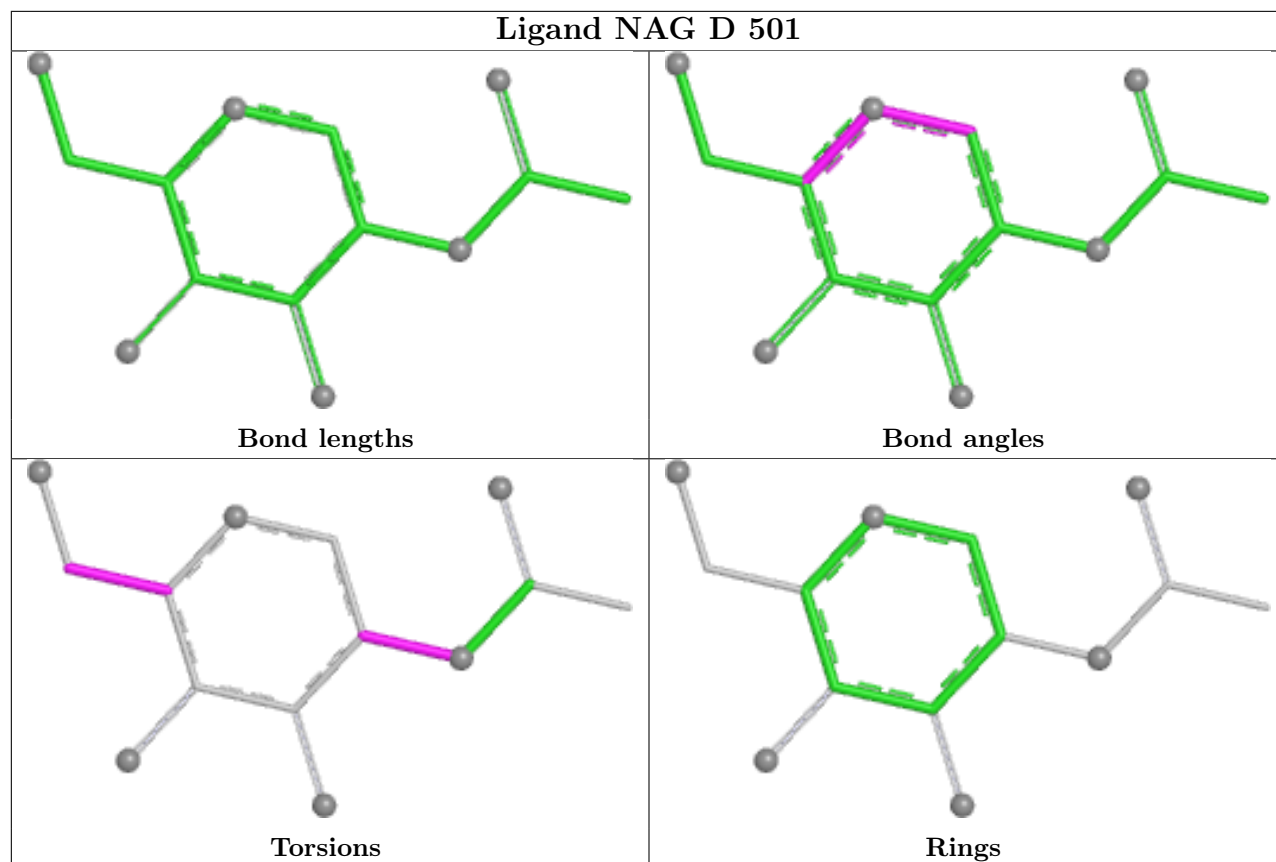
There are no ring outliers.

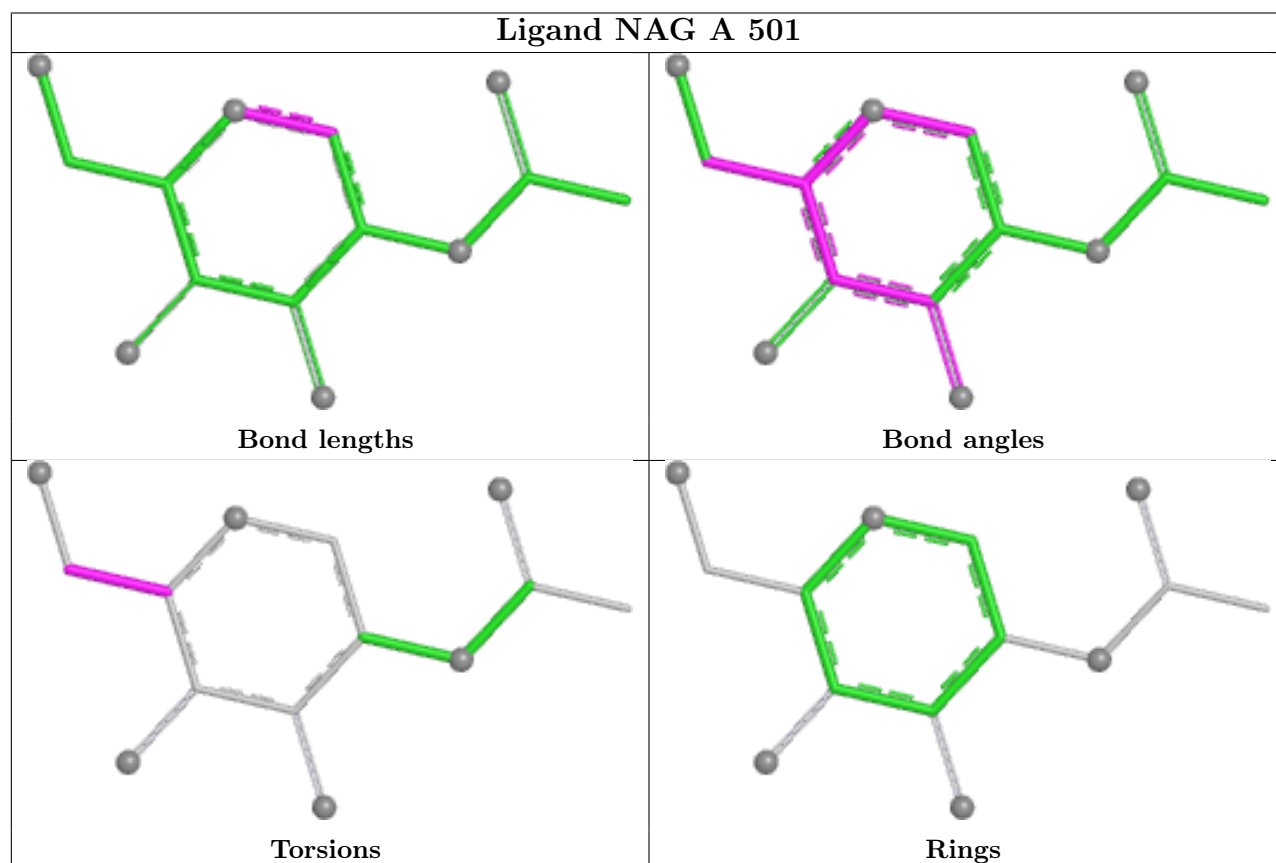
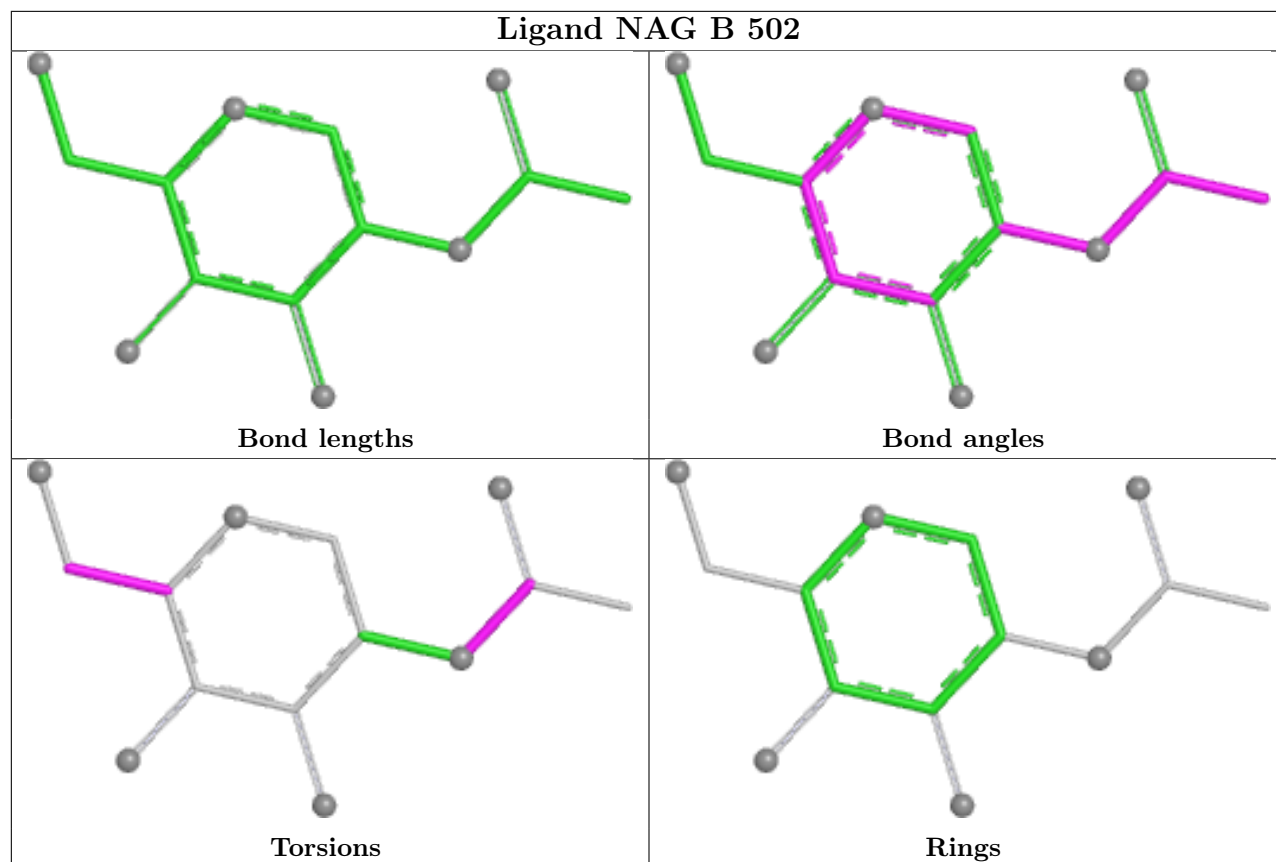
2 monomers are involved in 5 short contacts:

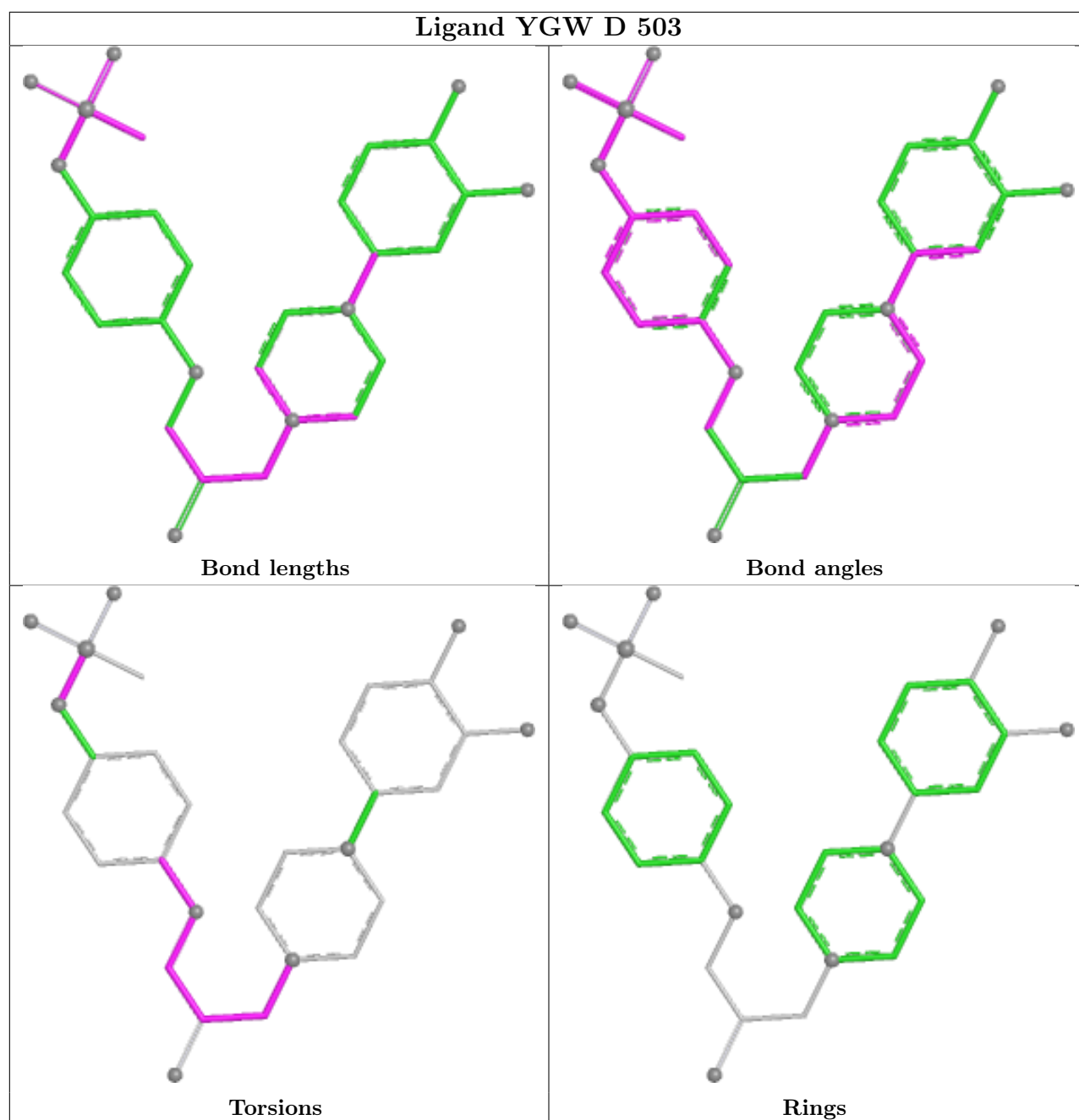
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	D	503	YGW	3	0
6	B	503	YGW	2	0

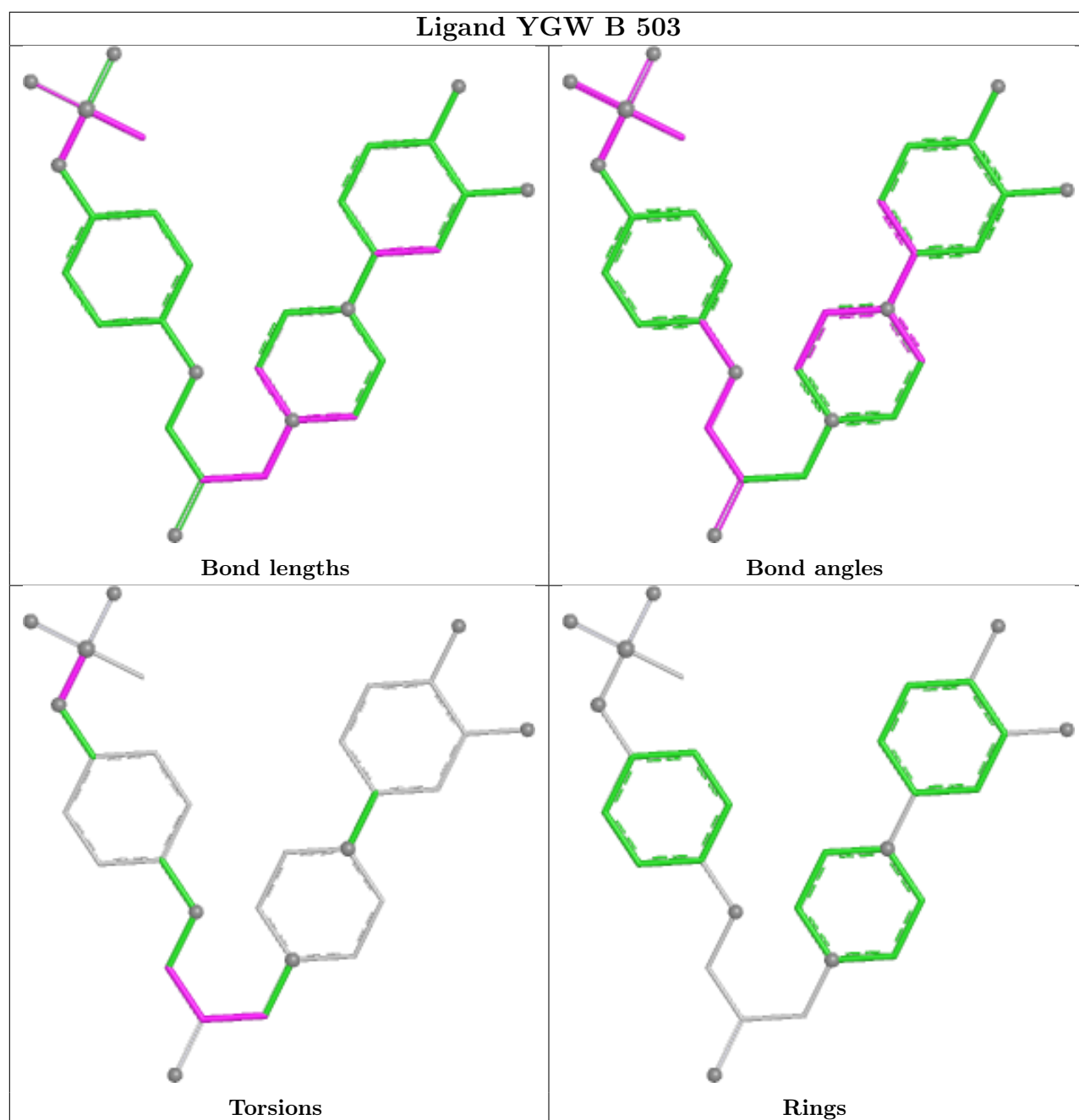
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

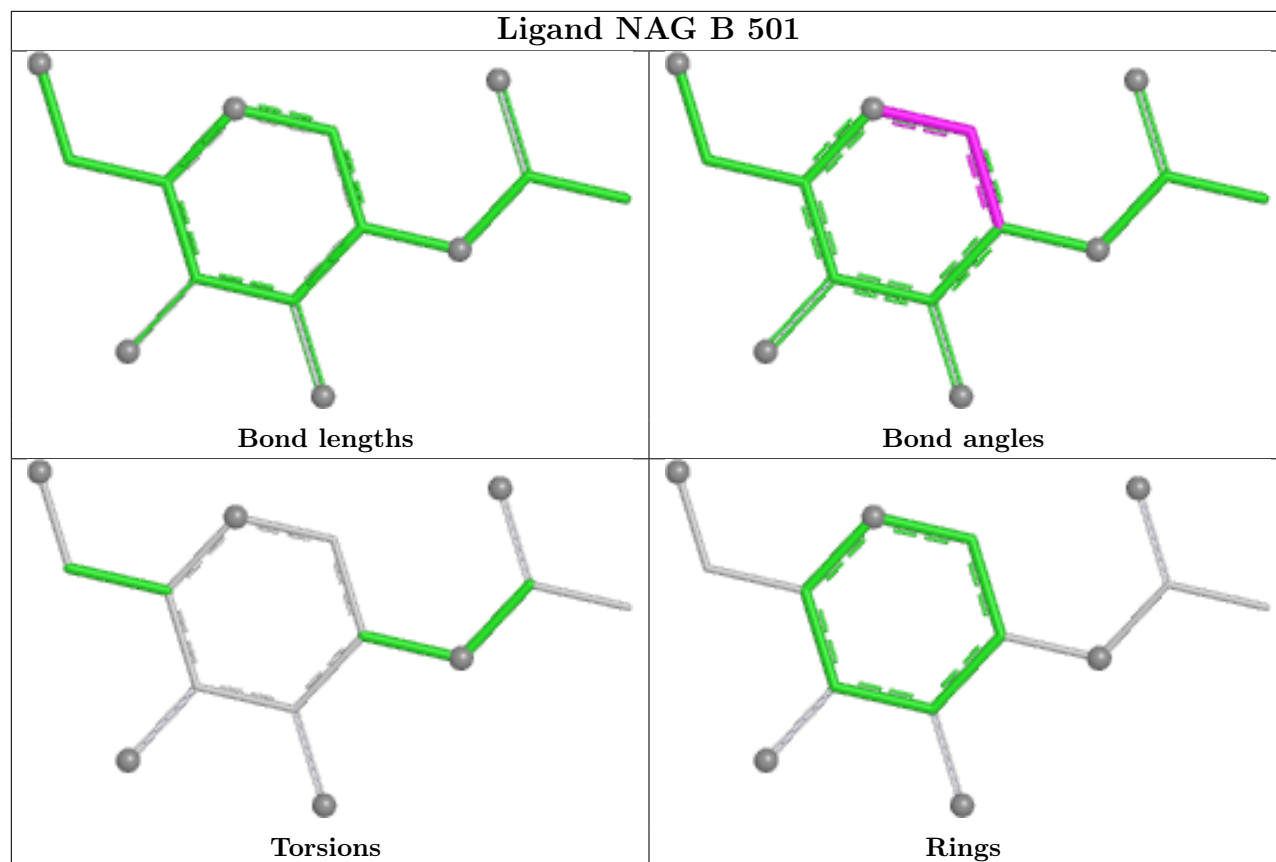












5.7 Other polymers [i](#)

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	359/385 (93%)	-0.45	3 (0%) 82 78	16, 37, 82, 112	0
1	C	358/385 (92%)	0.00	7 (1%) 65 57	31, 62, 102, 124	0
2	B	363/363 (100%)	-0.02	9 (2%) 58 49	17, 53, 99, 134	0
2	D	363/363 (100%)	0.02	11 (3%) 52 44	32, 60, 110, 147	0
All	All	1443/1496 (96%)	-0.11	30 (2%) 63 55	16, 54, 101, 147	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	407	THR	3.8
2	B	319	SER	3.4
1	C	97	ALA	3.3
2	D	344	PHE	3.1
2	B	56	LYS	3.1
1	A	97	ALA	2.9
2	D	48	VAL	2.9
2	D	324	THR	2.7
2	B	201	GLU	2.6
2	B	194	PHE	2.6
2	D	323	ASN	2.6
2	D	59	PHE	2.5
2	D	50	ILE	2.5
2	D	60	HIS	2.5
2	B	55	GLU	2.5
1	C	101	HIS	2.4
1	C	265	GLY	2.4
1	A	96	PRO	2.4
1	C	102	LEU	2.3
2	B	344	PHE	2.3
2	D	349	LEU	2.3

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Mol	Chain	Res	Type	RSRZ
2	B	57	ASP	2.3
2	B	48	VAL	2.2
2	D	32	PRO	2.2
2	D	62	LEU	2.2
2	D	46	ASP	2.2
2	B	393	ARG	2.1
1	C	400	ILE	2.0
1	C	235	LEU	2.0
1	C	238	ARG	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

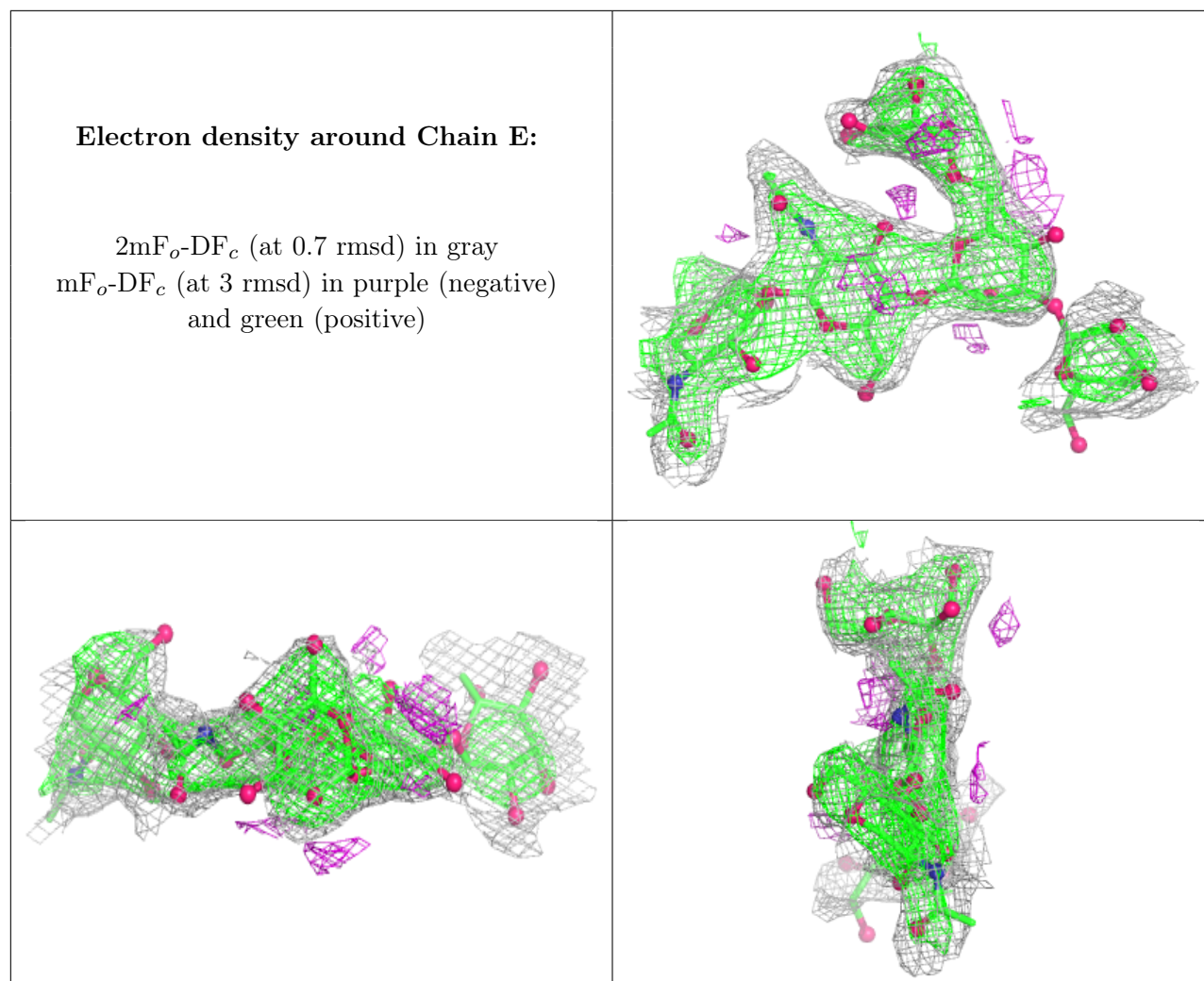
There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	NAG	E	1	14/15	-	-	20,20,20,20	14
3	NAG	E	2	14/15	-	-	20,20,20,20	14
3	BMA	E	3	11/12	-	-	20,20,20,20	11
3	MAN	E	4	11/12	-	-	20,20,20,20	11
3	MAN	E	5	11/12	-	-	20,20,20,20	11

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



6.4 Ligands [i](#)

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

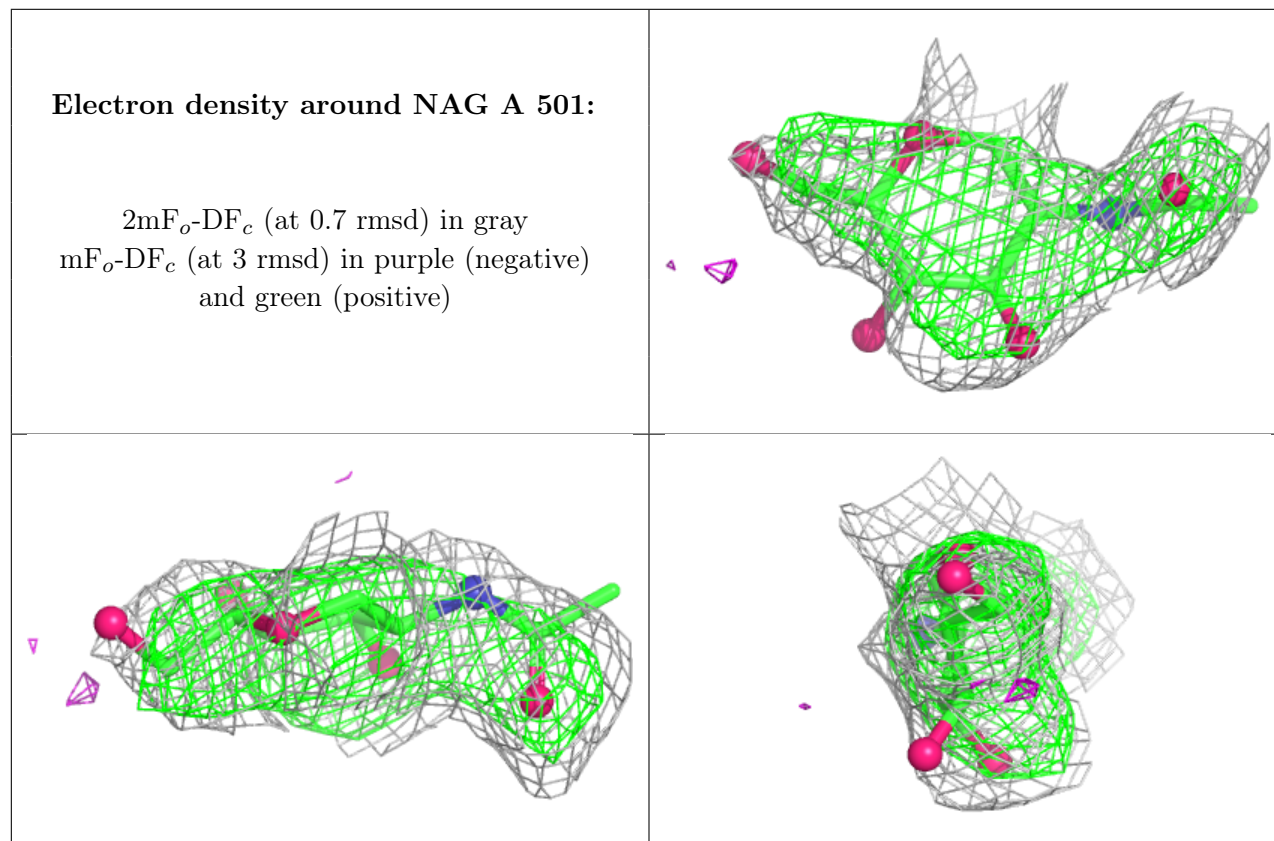
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	NAG	A	501	14/15	-	-	20,20,20,20	14
4	NAG	B	501	14/15	-	-	20,20,20,20	14
4	NAG	B	502	14/15	-	-	20,20,20,20	14
4	NAG	C	501	14/15	-	-	20,20,20,20	14
4	NAG	C	502	14/15	-	-	20,20,20,20	14
4	NAG	D	501	14/15	-	-	20,20,20,20	14
4	NAG	D	502	14/15	-	-	20,20,20,20	14
5	NA	A	502	1/1	-	-	44,44,44,44	1
5	NA	C	503	1/1	-	-	44,44,44,44	1

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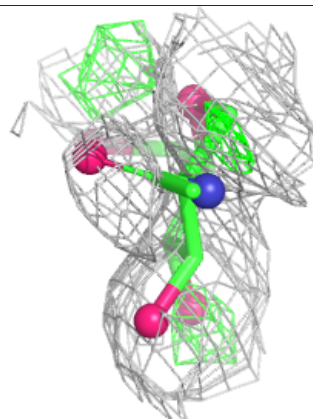
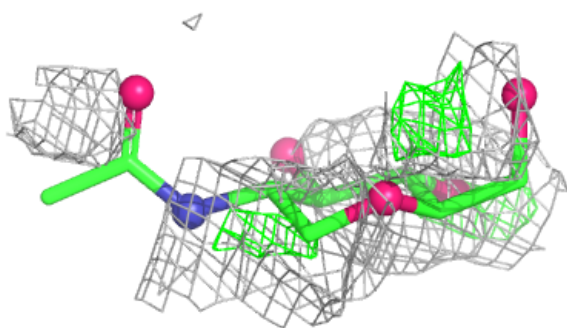
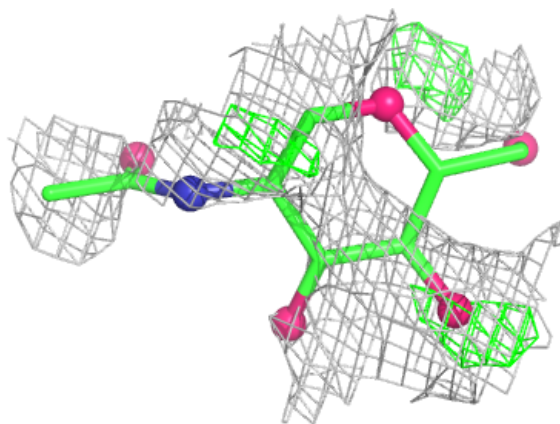
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
6	YGW	B	503	30/30	0.97	0.09	36,42,58,61	0
6	YGW	D	503	30/30	0.97	0.09	47,64,73,91	0

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.

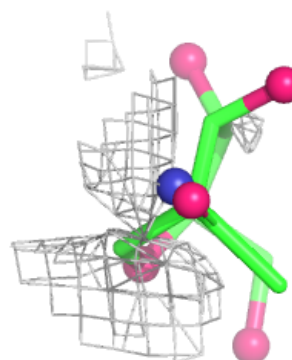
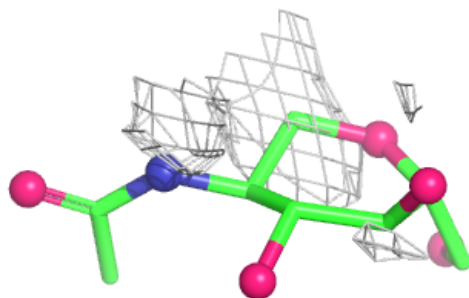
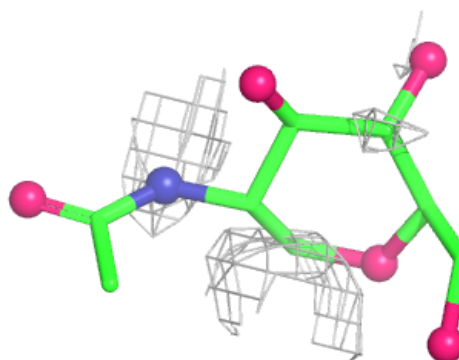


Electron density around NAG B 501:

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

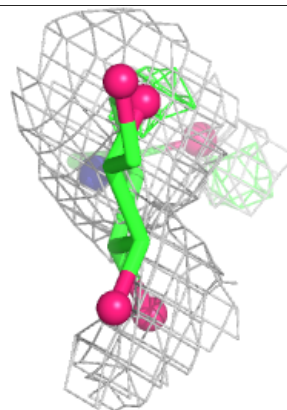
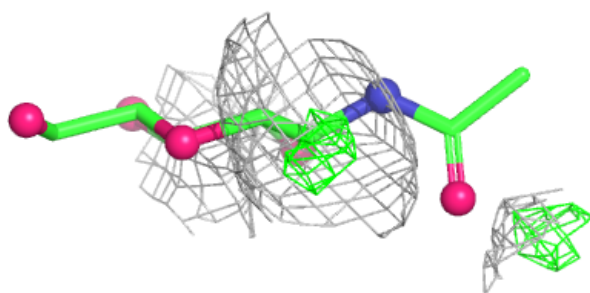
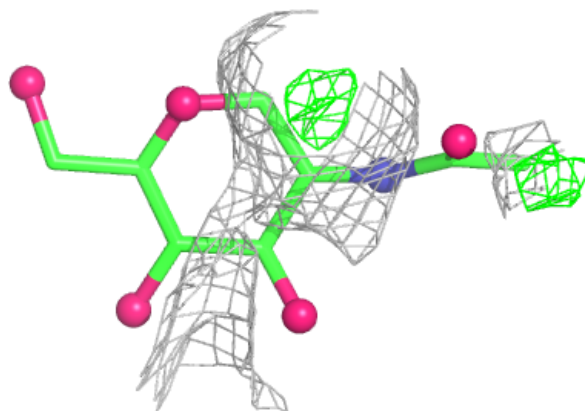
**Electron density around NAG B 502:**

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

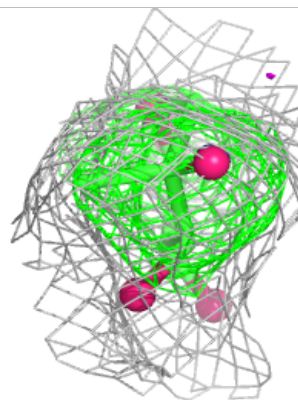
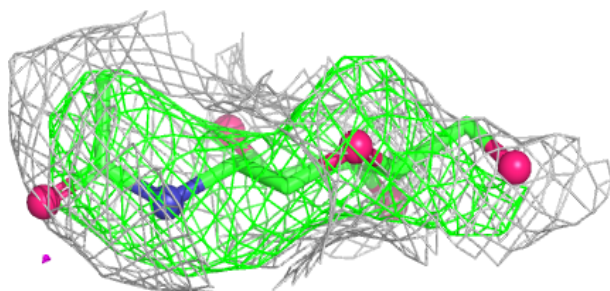
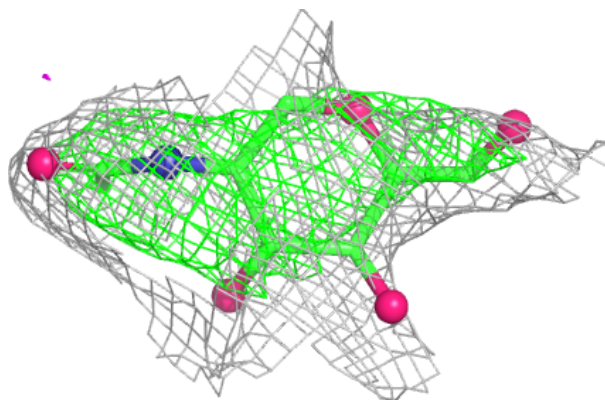


Electron density around NAG C 501:

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

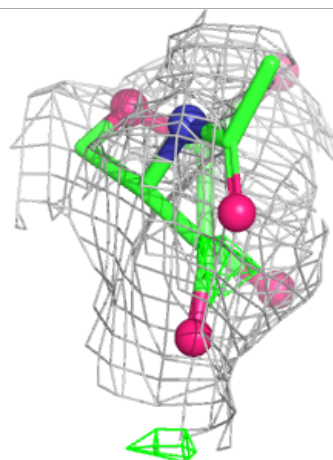
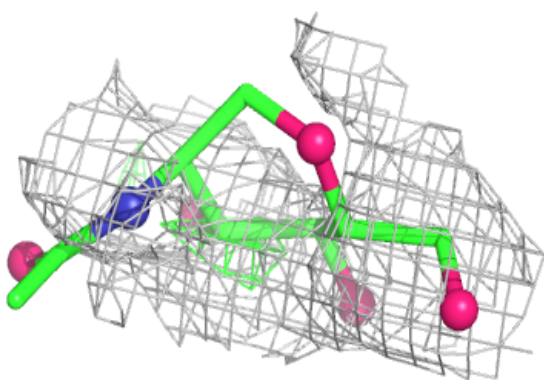
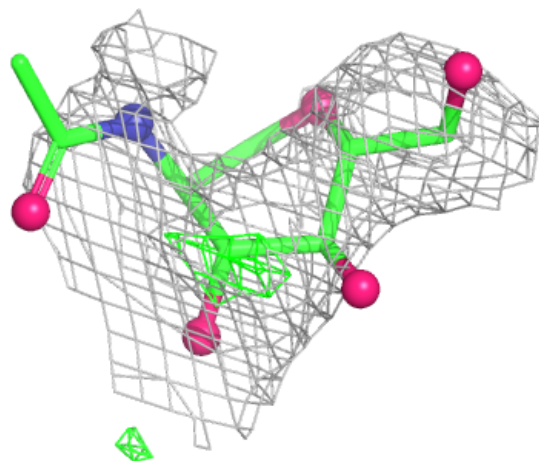
**Electron density around NAG C 502:**

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



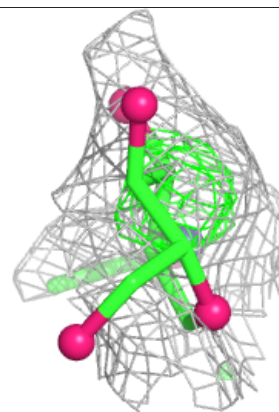
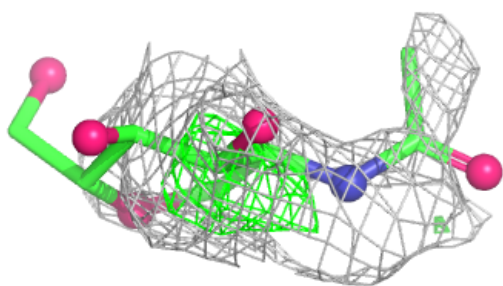
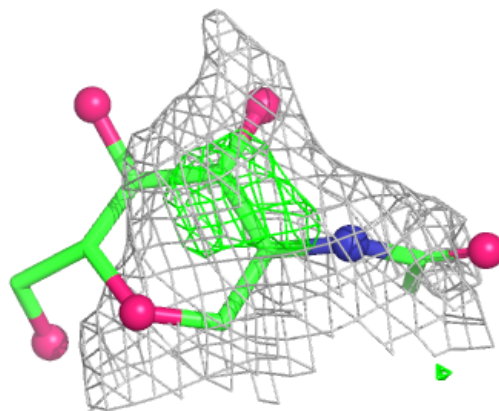
Electron density around NAG D 501:

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



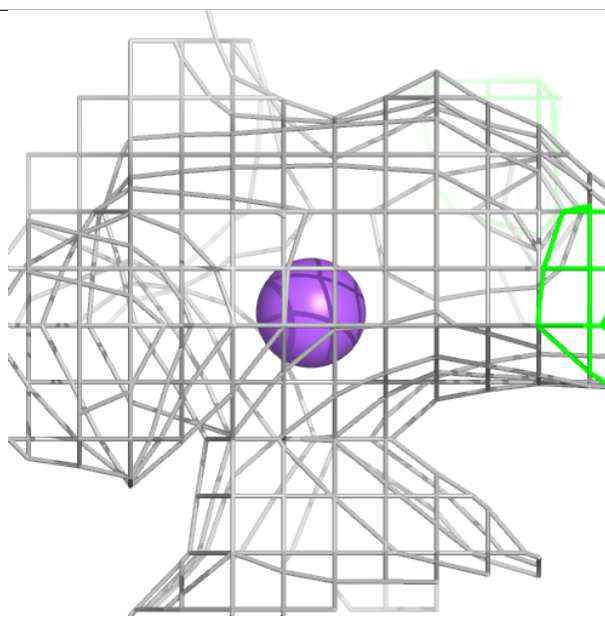
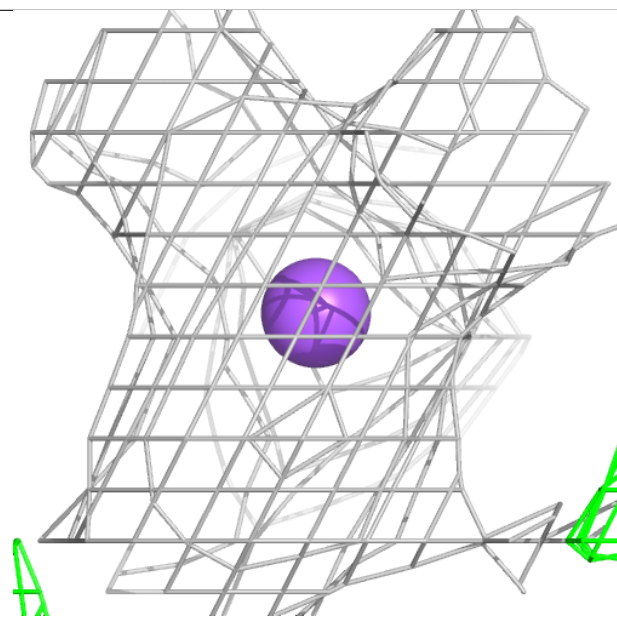
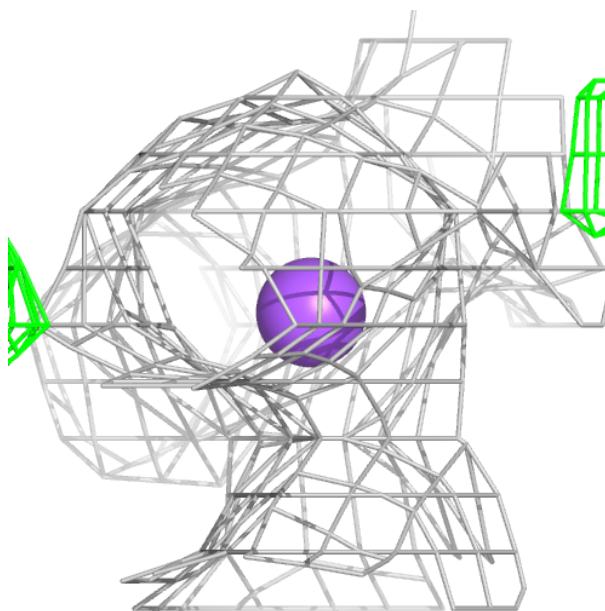
Electron density around NAG D 502:

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



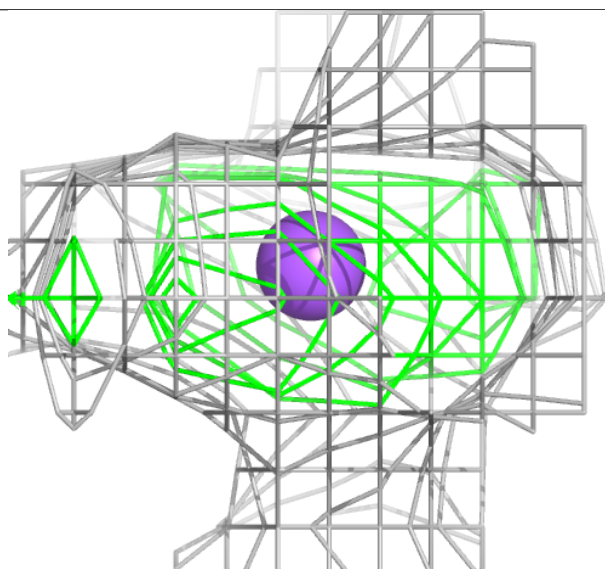
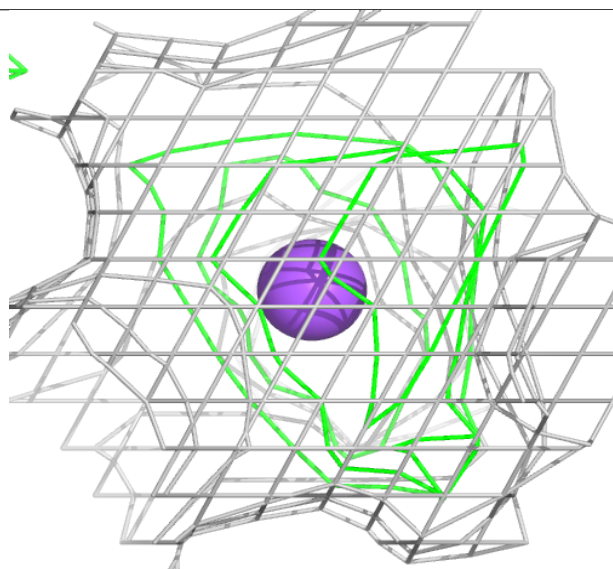
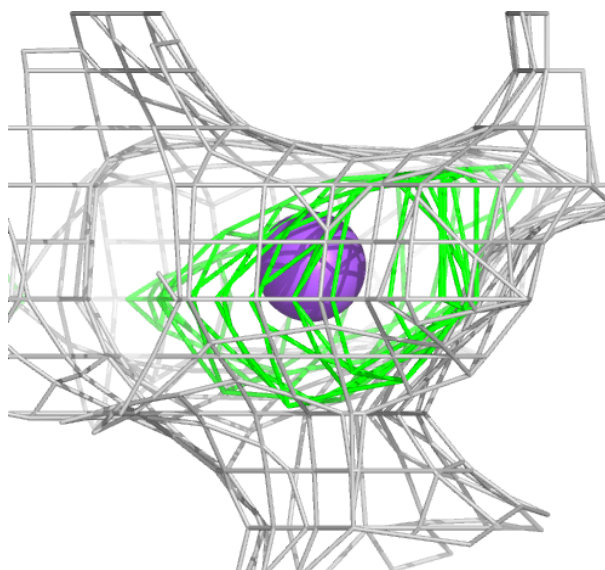
Electron density around NA A 502:

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



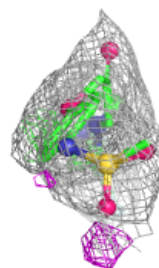
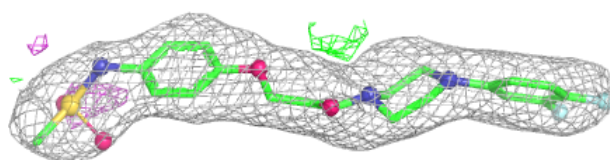
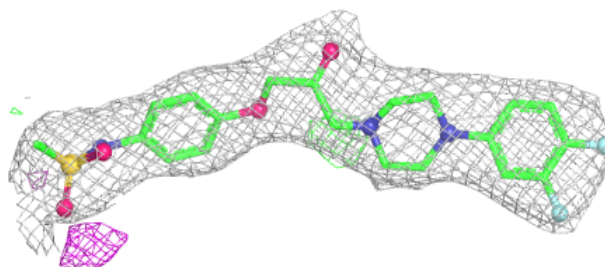
Electron density around NA C 503:

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

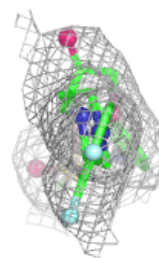
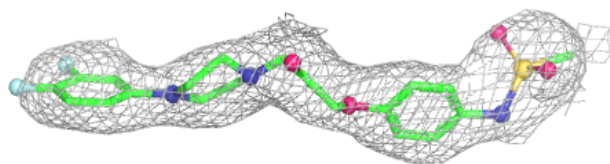
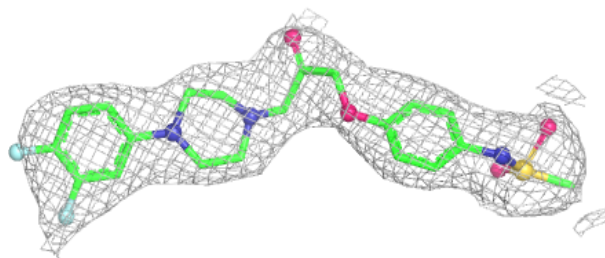


Electron density around YGW B 503:

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

**Electron density around YGW D 503:**

$2mF_o-DF_c$ (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.