



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 6, 2020 – 08:24 AM BST

PDB ID : 1MJT
Title : CRYSTAL STRUCTURE OF SANOS, A BACTERIAL NITRIC OXIDE
SYNTHASE OXYGENASE PROTEIN, IN COMPLEX WITH NAD⁺ AND
SEITU
Authors : Bird, L.E.; Ren, J.; Stammers, D.K.
Deposited on : 2002-08-28
Resolution : 2.40 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.13.1
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.13.1

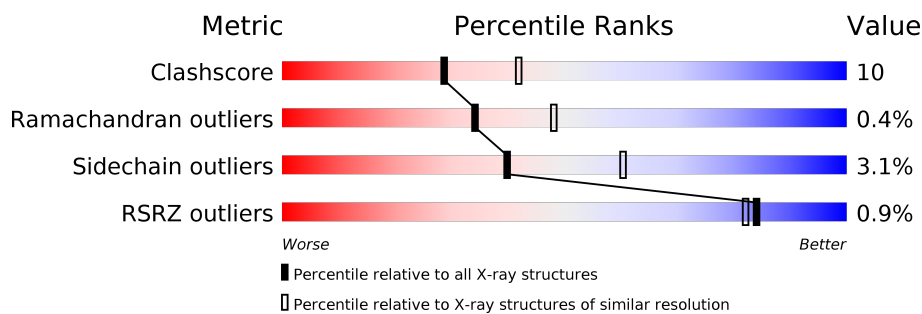
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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 on 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	347	<div> <div style="width: 100%; height: 10px; background-color: red;"></div> <div style="display: flex; justify-content: space-between; align-items: center;"> % <div style="width: 81%; height: 10px; background-color: green;"></div> <div style="width: 17%; height: 10px; background-color: orange;"></div> • </div> <div style="display: flex; justify-content: space-between; align-items: center;"> 81% 17% </div> </div>
1	B	347	<div> <div style="width: 100%; height: 10px; background-color: red;"></div> <div style="display: flex; justify-content: space-between; align-items: center;"> % <div style="width: 76%; height: 10px; background-color: green;"></div> <div style="width: 22%; height: 10px; background-color: orange;"></div> • </div> <div style="display: flex; justify-content: space-between; align-items: center;"> 76% 22% </div> </div>
2	C	2	<div> <div style="width: 100%; height: 10px; background-color: yellow;"></div> <div style="display: flex; justify-content: space-between; align-items: center;"> 100% </div> </div>
2	D	2	<div> <div style="width: 100%; height: 10px; background-color: yellow;"></div> <div style="display: flex; justify-content: space-between; align-items: center;"> 100% </div> </div>

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 6302 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.

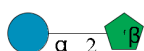
- Molecule 1 is a protein called NITRIC-OXIDE SYNTHASE HOMOLOG.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	346	Total	C	N	O	S	0	0	0
			2857	1843	479	528	7			
1	B	347	Total	C	N	O	S	0	0	0
			2867	1849	482	529	7			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	HIS	-	expression tag	UNP P0A004
A	1	HIS	-	expression tag	UNP P0A004
B	0	HIS	-	expression tag	UNP P0A004
B	1	HIS	-	expression tag	UNP P0A004

- Molecule 2 is an oligosaccharide called beta-D-fructofuranose-(2-1)-alpha-D-glucopyranose.



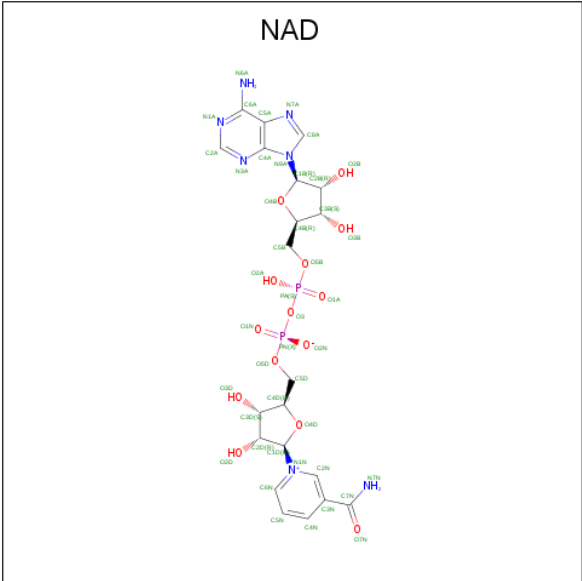
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	C	2	Total	C	O	0	0	0
			23	12	11			
2	D	2	Total	C	O	0	0	0
			23	12	11			

- Molecule 3 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: C₃₄H₃₂FeN₄O₄).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
3	B	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

- Molecule 4 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: C₂₁H₂₇N₇O₁₄P₂).



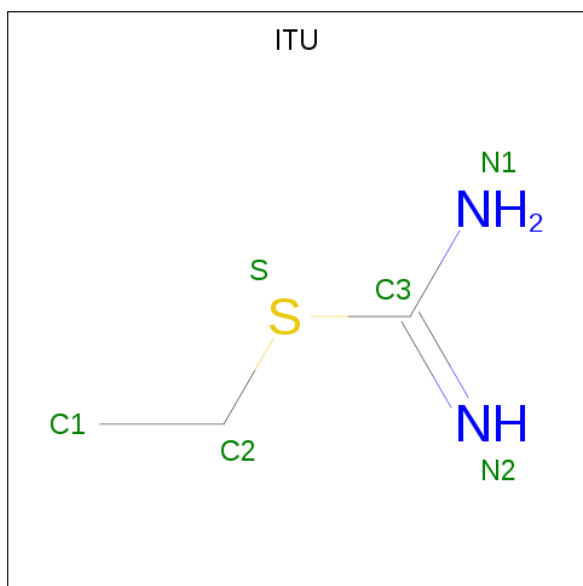
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			26	11	2	11	2		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			26	11	2	11	2		

- Molecule 5 is ETHYLISOTHIOUREA (three-letter code: ITU) (formula: $C_3H_8N_2S$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	S	0	0
			6	3	2	1		
5	B	1	Total	C	N	S	0	0
			6	3	2	1		

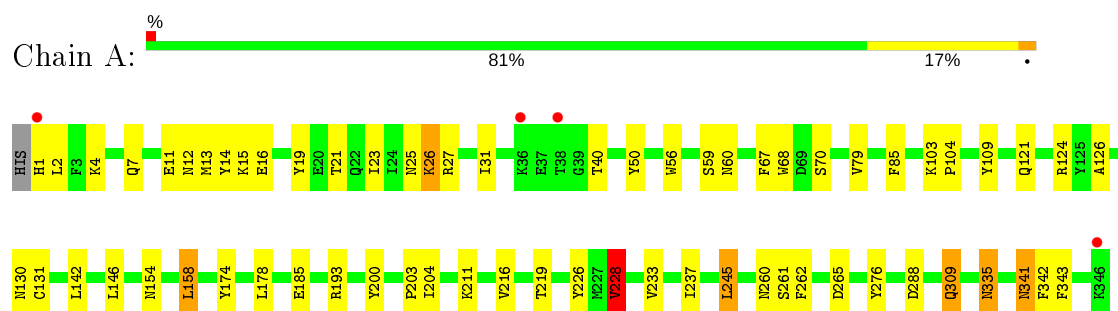
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	191	Total	O	0	0
			191	191		
6	B	191	Total	O	0	0
			191	191		

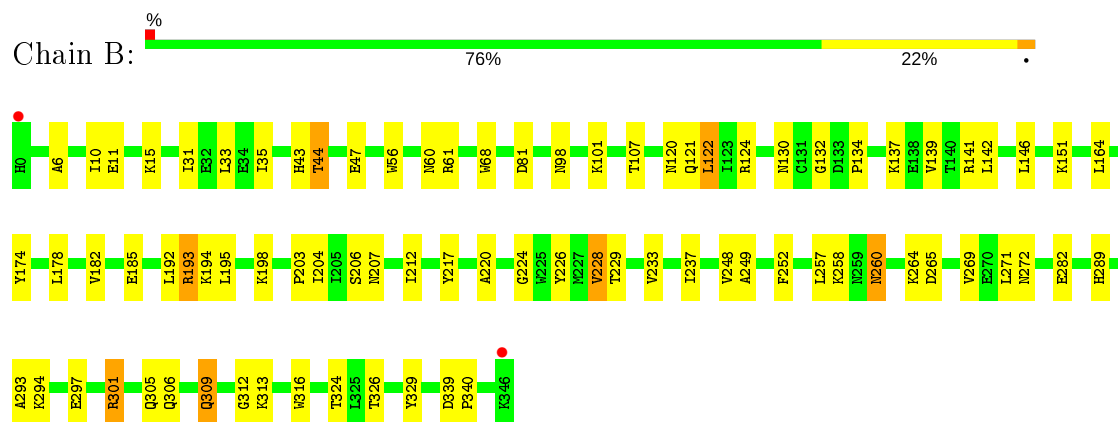
3 Residue-property plots

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

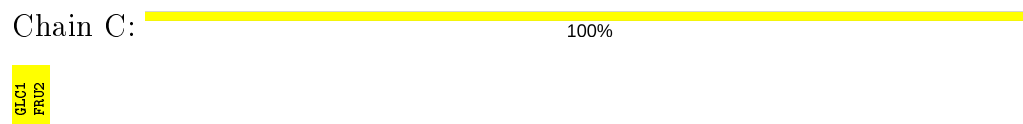
• Molecule 1: NITRIC-OXIDE SYNTHASE HOMOLOG



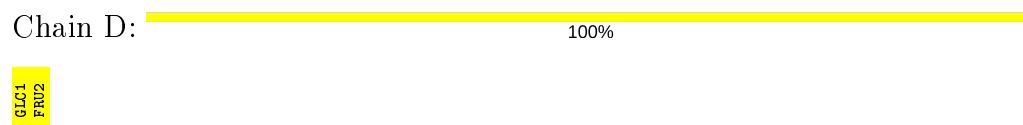
• Molecule 1: NITRIC-OXIDE SYNTHASE HOMOLOG



• Molecule 2: beta-D-fructofuranose-(2-1)-alpha-D-glucopyranose



• Molecule 2: beta-D-fructofuranose-(2-1)-alpha-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	65.78Å 115.14Å 126.02Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.34 – 2.40 29.34 – 2.40	Depositor EDS
% Data completeness (in resolution range)	95.7 (29.34-2.40) 95.8 (29.34-2.40)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.42 (at 2.39Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.182 , 0.241 0.175 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	33.9	Xtriage
Anisotropy	0.308	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 59.1	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	0.96	EDS
Total number of atoms	6302	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.92% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: HEM, GLC, ITU, FRU, NAD

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	0.37	0/2938	0.58	0/3986
1	B	0.38	0/2949	0.59	1/4001 (0.0%)
All	All	0.38	0/5887	0.58	1/7987 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	224	GLY	N-CA-C	-5.03	100.54	113.10

There are no chirality outliers.

There are no planarity outliers.

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	2857	0	2780	49	0
1	B	2867	0	2787	66	0
2	C	23	0	21	0	0
2	D	23	0	21	0	0
3	A	43	0	30	0	0
3	B	43	0	30	1	0
4	A	52	0	28	0	0
5	A	6	0	7	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	B	6	0	7	0	0
6	A	191	0	0	1	0
6	B	191	0	0	5	0
All	All	6302	0	5711	113	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 10.

All (113) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:341:ASN:HD22	1:A:342:PHE:H	1.08	1.00
1:A:60:ASN:HD22	1:A:335:ASN:HD22	1.25	0.85
1:A:13:MET:HG3	1:A:50:TYR:CD1	2.19	0.77
1:A:341:ASN:ND2	1:A:342:PHE:H	1.80	0.77
1:A:19:TYR:HB3	1:A:23:ILE:HD11	1.67	0.74
1:A:341:ASN:HD22	1:A:342:PHE:N	1.84	0.74
1:B:120:ASN:HA	1:B:207:ASN:HD22	1.54	0.73
1:A:60:ASN:HD22	1:A:335:ASN:ND2	1.88	0.72
1:A:1:HIS:HA	1:A:4:LYS:HD3	1.71	0.71
1:A:228:VAL:HG22	1:A:288:ASP:HA	1.72	0.70
1:B:44:THR:HG23	6:B:1098:HOH:O	1.92	0.70
1:B:258:LYS:HE2	1:B:260:ASN:ND2	2.08	0.69
1:B:258:LYS:HE2	1:B:260:ASN:HD21	1.58	0.68
1:B:44:THR:HG22	1:B:47:GLU:H	1.59	0.68
1:B:44:THR:HG21	6:B:1101:HOH:O	1.94	0.68
1:B:121:GLN:HB2	1:B:124:ARG:HG2	1.76	0.67
1:B:56:TRP:CE3	1:B:68:TRP:HA	2.30	0.67
1:B:141:ARG:HG2	1:B:141:ARG:HH11	1.60	0.66
1:B:151:LYS:HD3	6:B:1347:HOH:O	1.95	0.65
1:B:301:ARG:HB3	1:B:301:ARG:HH11	1.63	0.63
1:A:26:LYS:HE2	1:A:26:LYS:HA	1.82	0.62
1:B:43:HIS:HD2	1:B:217:TYR:OH	1.83	0.62
1:B:182:VAL:HG13	1:B:282:GLU:HG3	1.81	0.61
1:A:228:VAL:HG11	1:A:276:TYR:HB2	1.83	0.59
1:B:174:TYR:HB3	1:B:178:LEU:HD23	1.85	0.58
1:A:185:GLU:HB3	1:A:193:ARG:HD3	1.85	0.57
1:A:60:ASN:ND2	1:A:335:ASN:HD22	1.99	0.56
1:B:260:ASN:C	1:B:260:ASN:HD22	2.08	0.56
1:B:120:ASN:CA	1:B:207:ASN:HD22	2.19	0.55
1:B:339:ASP:OD1	1:B:340:PRO:HA	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:11:GLU:O	1:A:15:LYS:HD3	2.07	0.54
1:B:122:LEU:HD13	1:B:206:SER:HB3	1.90	0.54
1:B:141:ARG:NH1	1:B:141:ARG:HG2	2.21	0.54
1:B:137:LYS:O	1:B:141:ARG:HG3	2.08	0.53
1:B:6:ALA:O	1:B:10:ILE:HG12	2.08	0.53
1:A:126:ALA:HB2	1:A:158:LEU:HD13	1.89	0.53
1:B:132:GLY:O	1:B:134:PRO:HD3	2.09	0.52
1:A:12:ASN:O	1:A:16:GLU:HG3	2.09	0.52
1:B:44:THR:HB	1:B:47:GLU:CD	2.30	0.52
1:A:79:VAL:HG11	1:A:85:PHE:HA	1.91	0.51
1:B:192:LEU:HD12	1:B:195:LEU:HD12	1.92	0.51
1:A:245:LEU:HB2	6:A:1005:HOH:O	2.10	0.51
1:A:131:CYS:HB2	1:A:154:ASN:OD1	2.10	0.51
1:B:130:ASN:HB2	6:B:1322:HOH:O	2.10	0.51
1:A:174:TYR:HB3	1:A:178:LEU:HD23	1.92	0.51
1:A:228:VAL:CG2	1:A:288:ASP:HA	2.42	0.50
1:B:249:ALA:HB1	1:B:257:LEU:HD21	1.93	0.50
1:B:122:LEU:HD22	1:B:204:ILE:HG21	1.92	0.49
1:A:7:GLN:HG2	1:A:11:GLU:OE2	2.13	0.49
1:B:11:GLU:O	1:B:15:LYS:HB2	2.11	0.49
1:A:121:GLN:HB2	1:A:124:ARG:HD3	1.94	0.49
1:A:211:LYS:HG2	1:A:216:VAL:HG22	1.93	0.49
1:B:301:ARG:CB	1:B:301:ARG:HH11	2.24	0.49
1:B:289:HIS:CD2	1:B:289:HIS:H	2.31	0.48
1:A:121:GLN:HB2	1:A:124:ARG:HG2	1.95	0.48
1:B:203:PRO:HB3	1:B:226:TYR:CZ	2.49	0.48
1:B:233:VAL:HG23	1:B:272:ASN:HD21	1.78	0.48
1:B:194:LYS:HD2	1:B:194:LYS:N	2.28	0.48
1:A:21:THR:HG22	1:A:25:ASN:HD21	1.78	0.47
1:A:121:GLN:HB2	1:A:124:ARG:CG	2.44	0.47
1:A:203:PRO:HB3	1:A:226:TYR:CZ	2.50	0.47
1:A:265:ASP:OD2	1:B:289:HIS:HE1	1.97	0.47
1:A:19:TYR:CB	1:A:23:ILE:HD11	2.40	0.46
1:A:2:LEU:HD21	1:A:40:THR:HA	1.96	0.46
1:A:341:ASN:HB3	1:A:343:PHE:CE1	2.50	0.46
1:B:228:VAL:HG13	1:B:229:THR:N	2.31	0.46
1:B:324:THR:HG22	1:B:329:TYR:CE1	2.50	0.46
1:B:121:GLN:HB2	1:B:124:ARG:CG	2.46	0.45
1:B:107:THR:O	1:B:220:ALA:HA	2.17	0.45
1:B:339:ASP:CG	1:B:340:PRO:HA	2.37	0.45
1:B:61:ARG:NH1	3:B:501:HEM:HAD1	2.31	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:185:GLU:HB3	1:B:193:ARG:CD	2.47	0.45
1:B:212:ILE:HA	1:B:340:PRO:HB3	1.99	0.45
1:B:31:ILE:O	1:B:35:ILE:HG13	2.17	0.45
1:B:248:VAL:HG12	1:B:252:PHE:CE1	2.52	0.44
1:B:249:ALA:CB	1:B:257:LEU:HD21	2.48	0.44
1:A:142:LEU:O	1:A:146:LEU:HG	2.18	0.44
1:A:233:VAL:O	1:A:237:ILE:HB	2.17	0.44
1:A:56:TRP:CE3	1:A:68:TRP:HA	2.52	0.43
1:B:142:LEU:O	1:B:146:LEU:HG	2.18	0.43
1:B:257:LEU:O	1:B:264:LYS:HE3	2.18	0.43
1:B:15:LYS:HB2	1:B:15:LYS:HE2	1.89	0.43
1:A:309:GLN:HB2	1:A:309:GLN:HE21	1.52	0.43
1:B:265:ASP:O	1:B:269:VAL:HG23	2.18	0.43
1:B:233:VAL:O	1:B:237:ILE:HB	2.19	0.43
1:B:294:LYS:HD2	1:B:297:GLU:OE2	2.19	0.43
1:B:289:HIS:CD2	1:B:289:HIS:N	2.86	0.43
1:A:26:LYS:HE2	1:A:26:LYS:CA	2.48	0.43
1:A:261:SER:O	1:A:262:PHE:HB2	2.19	0.42
1:A:14:TYR:CD1	1:A:14:TYR:N	2.87	0.42
1:B:193:ARG:HD2	1:B:193:ARG:HA	1.91	0.42
1:A:265:ASP:OD2	1:B:289:HIS:CE1	2.71	0.42
1:A:27:ARG:O	1:A:31:ILE:HG13	2.19	0.42
1:B:248:VAL:CG1	1:B:252:PHE:CE1	3.02	0.42
1:B:98:ASN:ND2	1:B:101:LYS:HB2	2.35	0.42
1:B:260:ASN:ND2	1:B:260:ASN:C	2.73	0.42
1:B:309:GLN:HB2	1:B:309:GLN:HE21	1.52	0.42
1:B:60:ASN:HB3	1:B:312:GLY:HA2	2.02	0.42
1:B:237:ILE:HD11	1:B:271:LEU:CD2	2.50	0.42
1:B:293:ALA:HB1	1:B:326:THR:HB	2.02	0.42
1:A:130:ASN:O	1:A:131:CYS:HB3	2.20	0.42
1:A:121:GLN:HA	1:A:204:ILE:O	2.20	0.42
1:B:313:LYS:HD3	1:B:316:TRP:CD1	2.55	0.42
1:A:211:LYS:HD2	1:A:343:PHE:CE1	2.55	0.41
1:B:258:LYS:CE	1:B:260:ASN:HD21	2.29	0.41
1:A:121:GLN:NE2	1:A:200:TYR:OH	2.53	0.41
1:B:182:VAL:O	1:B:198:LYS:HA	2.20	0.41
1:A:109:TYR:CD2	1:A:219:THR:HG22	2.56	0.41
1:B:43:HIS:CD2	1:B:217:TYR:OH	2.69	0.41
1:A:67:PHE:O	1:A:70:SER:HB3	2.19	0.41
1:A:26:LYS:HA	1:A:26:LYS:CE	2.49	0.41
1:A:103:LYS:HA	1:A:104:PRO:HD3	1.94	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:139:VAL:HG23	6:B:1020:HOH:O	2.20	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	344/347 (99%)	325 (94%)	17 (5%)	2 (1%)	25	36
1	B	345/347 (99%)	329 (95%)	15 (4%)	1 (0%)	41	55
All	All	689/694 (99%)	654 (95%)	32 (5%)	3 (0%)	34	48

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	59	SER
1	A	228	VAL
1	B	228	VAL

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	307/308 (100%)	299 (97%)	8 (3%)	46	66
1	B	308/308 (100%)	297 (96%)	11 (4%)	35	54

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	615/616 (100%)	596 (97%)	19 (3%)	40 60

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	26	LYS
1	A	158	LEU
1	A	228	VAL
1	A	245	LEU
1	A	260	ASN
1	A	309	GLN
1	A	335	ASN
1	A	341	ASN
1	B	33	LEU
1	B	44	THR
1	B	81	ASP
1	B	122	LEU
1	B	164	LEU
1	B	193	ARG
1	B	260	ASN
1	B	301	ARG
1	B	305	GLN
1	B	306	GLN
1	B	309	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	25	ASN
1	A	29	HIS
1	A	95	GLN
1	A	121	GLN
1	A	144	ASN
1	A	166	ASN
1	A	235	ASN
1	A	260	ASN
1	A	305	GLN
1	A	309	GLN
1	A	335	ASN
1	A	341	ASN
1	B	12	ASN

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Mol	Chain	Res	Type
1	B	25	ASN
1	B	43	HIS
1	B	95	GLN
1	B	98	ASN
1	B	121	GLN
1	B	144	ASN
1	B	207	ASN
1	B	260	ASN
1	B	289	HIS
1	B	306	GLN
1	B	335	ASN

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 ⓘ

4 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	GLC	C	1	2	11,11,12	1.49	2 (18%)	15,15,17	1.07	2 (13%)
2	FRU	C	2	2	11,12,12	1.53	1 (9%)	10,18,18	0.83	0
2	GLC	D	1	2	11,11,12	1.32	1 (9%)	15,15,17	1.20	2 (13%)
2	FRU	D	2	2	11,12,12	1.34	1 (9%)	10,18,18	0.74	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	GLC	C	1	2	-	2/2/19/22	0/1/1/1
2	FRU	C	2	2	-	0/5/24/24	0/1/1/1
2	GLC	D	1	2	-	2/2/19/22	0/1/1/1
2	FRU	D	2	2	-	2/5/24/24	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	2	FRU	O2-C2	4.66	1.48	1.40
2	D	2	FRU	O2-C2	4.08	1.47	1.40
2	C	1	GLC	O5-C1	2.91	1.48	1.43
2	D	1	GLC	O5-C1	2.90	1.48	1.43
2	C	1	GLC	C2-C3	2.37	1.56	1.52

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	1	GLC	C1-O5-C5	3.34	116.71	112.19
2	C	1	GLC	C1-O5-C5	2.75	115.91	112.19
2	D	1	GLC	O5-C1-C2	-2.27	107.26	110.77
2	C	1	GLC	O5-C1-C2	-2.15	107.45	110.77

There are no chirality outliers.

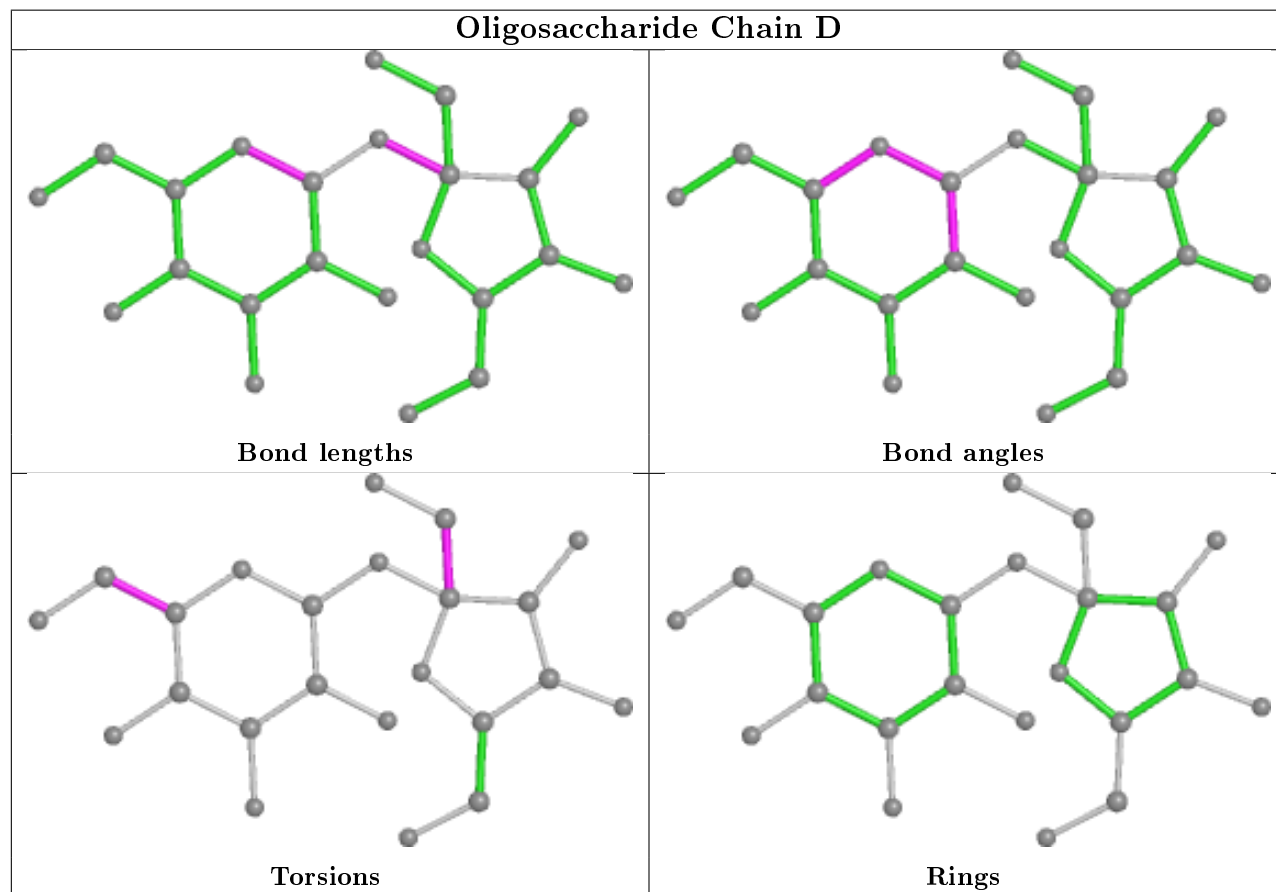
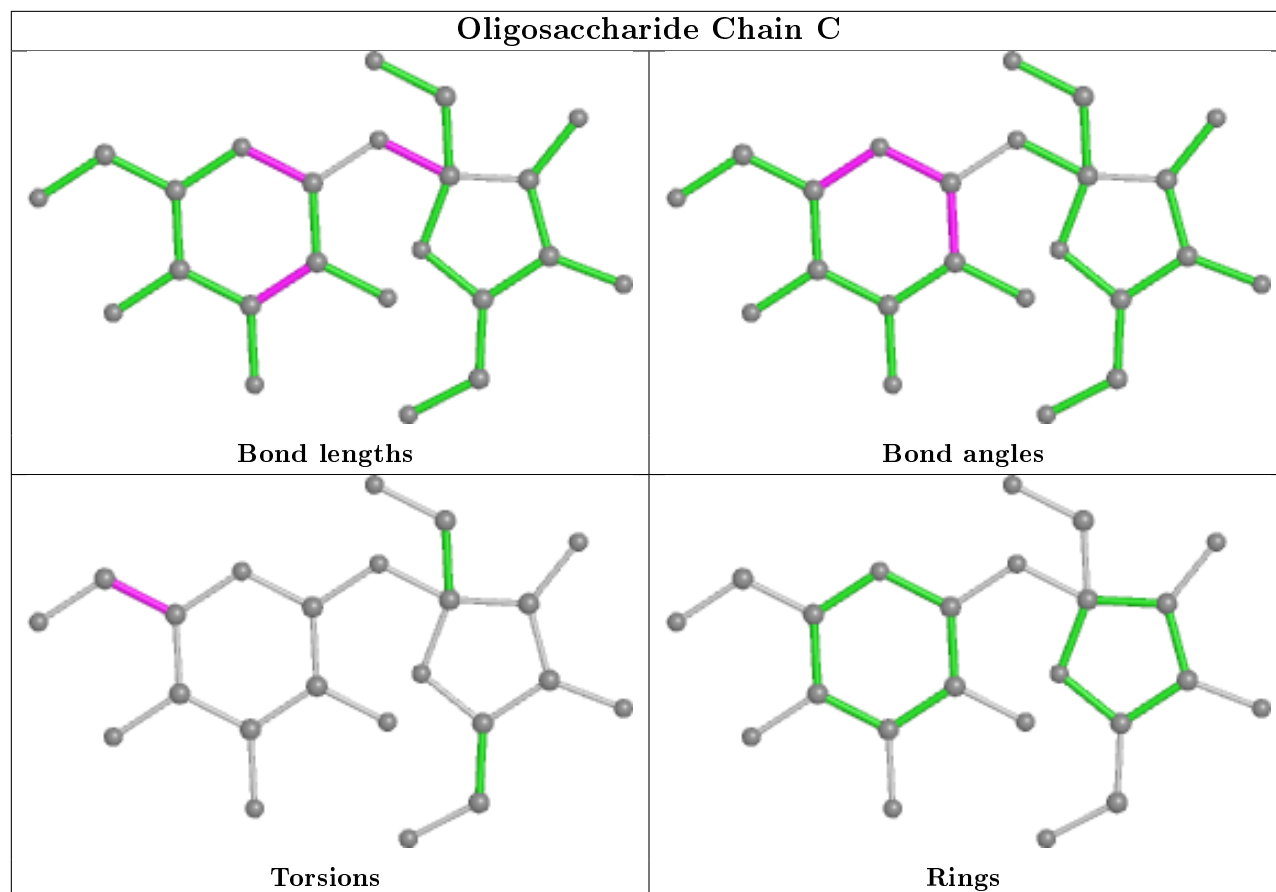
All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	2	FRU	O1-C1-C2-O2
2	D	1	GLC	O5-C5-C6-O6
2	C	1	GLC	C4-C5-C6-O6
2	D	1	GLC	C4-C5-C6-O6
2	C	1	GLC	O5-C5-C6-O6
2	D	2	FRU	O1-C1-C2-O5

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

6 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
5	ITU	B	801	-	4,5,5	0.79	0	4,5,5	2.95	1 (25%)
3	HEM	A	500	1	27,50,50	1.86	10 (37%)	17,82,82	1.14	1 (5%)
4	NAD	A	701	-	24,27,48	2.60	9 (37%)	34,41,73	1.98	10 (29%)
5	ITU	A	800	-	4,5,5	0.84	0	4,5,5	3.09	1 (25%)
3	HEM	B	501	1	27,50,50	1.86	7 (25%)	17,82,82	0.95	1 (5%)
4	NAD	A	700	-	24,27,48	2.68	9 (37%)	34,41,73	1.96	8 (23%)

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
5	ITU	B	801	-	-	1/3/3/3	-
3	HEM	A	500	1	-	0/6/54/54	-
4	NAD	A	701	-	-	4/20/36/62	0/2/2/5
5	ITU	A	800	-	-	1/3/3/3	-
3	HEM	B	501	1	-	0/6/54/54	-
4	NAD	A	700	-	-	5/20/36/62	0/2/2/5

All (35) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	700	NAD	C2N-N1N	6.48	1.42	1.35
4	A	701	NAD	C2N-N1N	6.40	1.42	1.35
4	A	700	NAD	O4D-C1D	5.06	1.48	1.41
4	A	701	NAD	C6N-N1N	4.96	1.47	1.35
4	A	700	NAD	C6N-N1N	4.95	1.47	1.35
4	A	700	NAD	C4N-C3N	4.65	1.47	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	701	NAD	C4N-C3N	4.55	1.47	1.39
4	A	701	NAD	O4D-C1D	4.43	1.47	1.41
4	A	701	NAD	C3N-C7N	4.21	1.56	1.50
4	A	700	NAD	C3N-C7N	4.07	1.56	1.50
3	A	500	HEM	C3B-CAB	-3.83	1.40	1.47
3	A	500	HEM	CBB-CAB	3.57	1.53	1.29
3	B	501	HEM	CBC-CAC	3.56	1.52	1.29
3	B	501	HEM	CBB-CAB	3.56	1.52	1.29
3	A	500	HEM	CBC-CAC	3.39	1.51	1.29
3	B	501	HEM	C3B-CAB	-3.27	1.41	1.47
4	A	700	NAD	PA-O5B	3.25	1.67	1.54
4	A	701	NAD	PA-O5B	3.20	1.67	1.54
3	B	501	HEM	C3C-CAC	-3.16	1.41	1.47
3	A	500	HEM	C3C-C2C	-3.04	1.36	1.40
3	A	500	HEM	C3C-CAC	-2.98	1.41	1.47
4	A	701	NAD	C5N-C4N	2.95	1.45	1.38
4	A	700	NAD	C5N-C4N	2.92	1.45	1.38
3	B	501	HEM	C1A-NA	2.89	1.42	1.36
3	B	501	HEM	C4B-NB	2.77	1.41	1.36
3	A	500	HEM	C3B-C2B	-2.68	1.36	1.40
4	A	701	NAD	C2N-C3N	2.59	1.43	1.39
4	A	700	NAD	C2N-C3N	2.55	1.43	1.39
3	B	501	HEM	C3C-C2C	-2.47	1.36	1.40
3	A	500	HEM	C1C-C2C	2.27	1.47	1.42
3	A	500	HEM	C4B-NB	2.24	1.40	1.36
3	A	500	HEM	C1D-ND	2.19	1.40	1.36
4	A	700	NAD	C7N-N7N	2.18	1.37	1.33
4	A	701	NAD	C7N-N7N	2.03	1.36	1.33
3	A	500	HEM	C4D-C3D	2.02	1.47	1.42

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	800	ITU	C2-S-C3	6.18	109.67	103.20
5	B	801	ITU	C2-S-C3	5.88	109.36	103.20
4	A	701	NAD	C6N-C5N-C4N	4.86	126.50	119.44
4	A	700	NAD	C6N-C5N-C4N	4.86	126.50	119.44
4	A	701	NAD	C5N-C4N-C3N	-4.81	114.66	120.34
4	A	700	NAD	C5N-C4N-C3N	-4.72	114.76	120.34
4	A	701	NAD	C2N-C3N-C4N	3.83	122.60	118.26
4	A	700	NAD	C2N-C3N-C4N	3.72	122.48	118.26
4	A	700	NAD	C5N-C6N-N1N	-3.67	115.14	120.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	701	NAD	C5N-C6N-N1N	-3.62	115.20	120.40
4	A	701	NAD	C2N-N1N-C1D	-3.27	111.86	119.14
4	A	700	NAD	C2N-N1N-C1D	-3.26	111.87	119.14
4	A	701	NAD	C4N-C3N-C7N	-2.74	113.69	121.04
4	A	700	NAD	C4N-C3N-C7N	-2.72	113.76	121.04
3	A	500	HEM	C4C-C3C-C2C	-2.65	105.05	106.90
4	A	700	NAD	C3N-C2N-N1N	-2.51	117.97	120.43
4	A	701	NAD	C3N-C2N-N1N	-2.49	117.99	120.43
4	A	701	NAD	C3D-C2D-C1D	2.10	104.14	100.98
3	B	501	HEM	C4A-C3A-C2A	2.08	108.44	107.00
4	A	700	NAD	O2N-PN-O1N	2.07	122.50	112.24
4	A	701	NAD	O2A-PA-O1A	2.03	118.63	110.68
4	A	701	NAD	O2N-PN-O1N	2.00	122.15	112.24

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	701	NAD	PA-O3-PN-O5D
4	A	701	NAD	O4D-C4D-C5D-O5D
4	A	701	NAD	C3D-C4D-C5D-O5D
4	A	700	NAD	C5D-O5D-PN-O3
4	A	700	NAD	C5D-O5D-PN-O1N
4	A	700	NAD	C3D-C4D-C5D-O5D
5	B	801	ITU	C1-C2-S-C3
5	A	800	ITU	C1-C2-S-C3
4	A	700	NAD	O4D-C4D-C5D-O5D
4	A	700	NAD	PA-O3-PN-O5D
4	A	701	NAD	C5D-O5D-PN-O2N

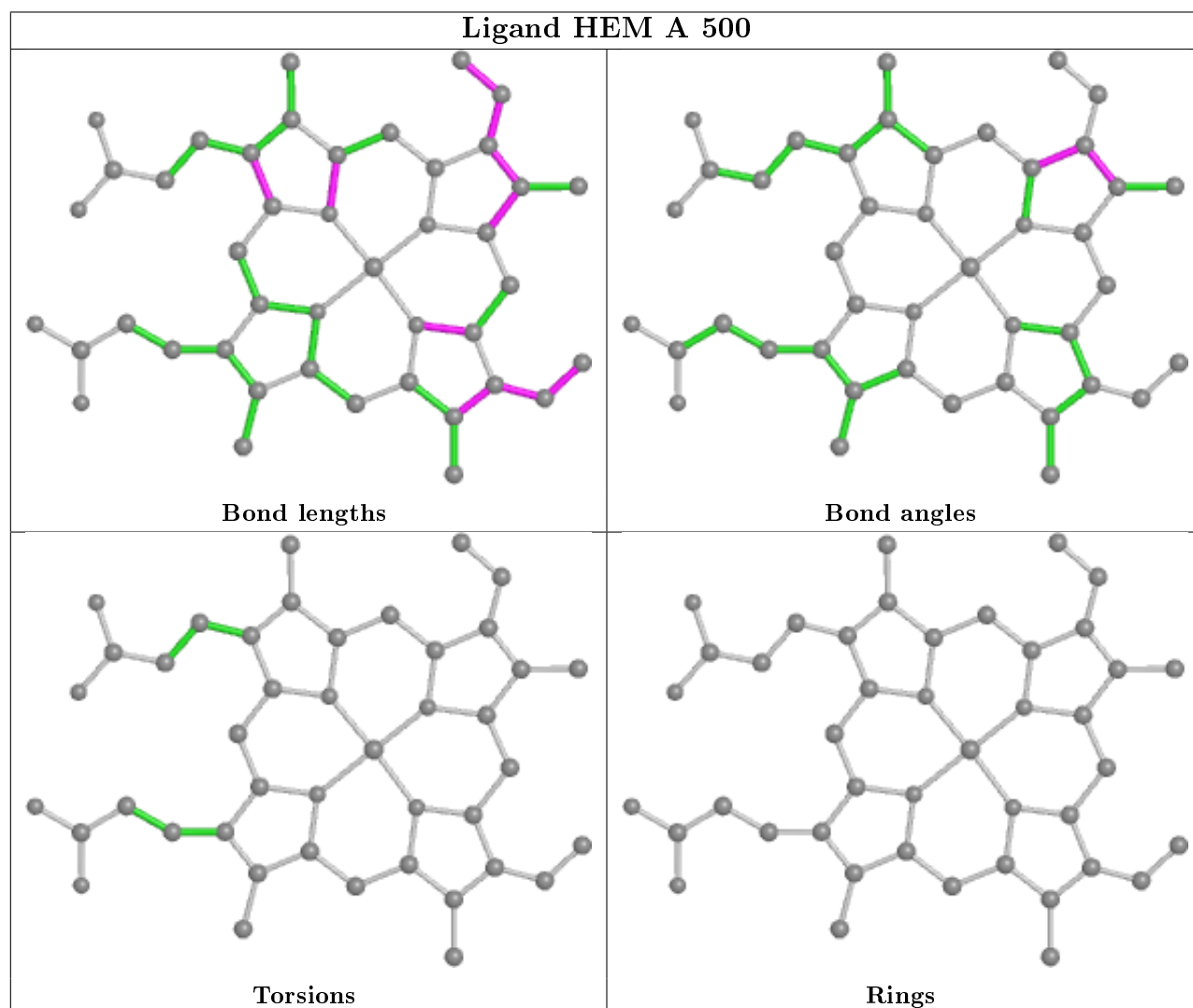
There are no ring outliers.

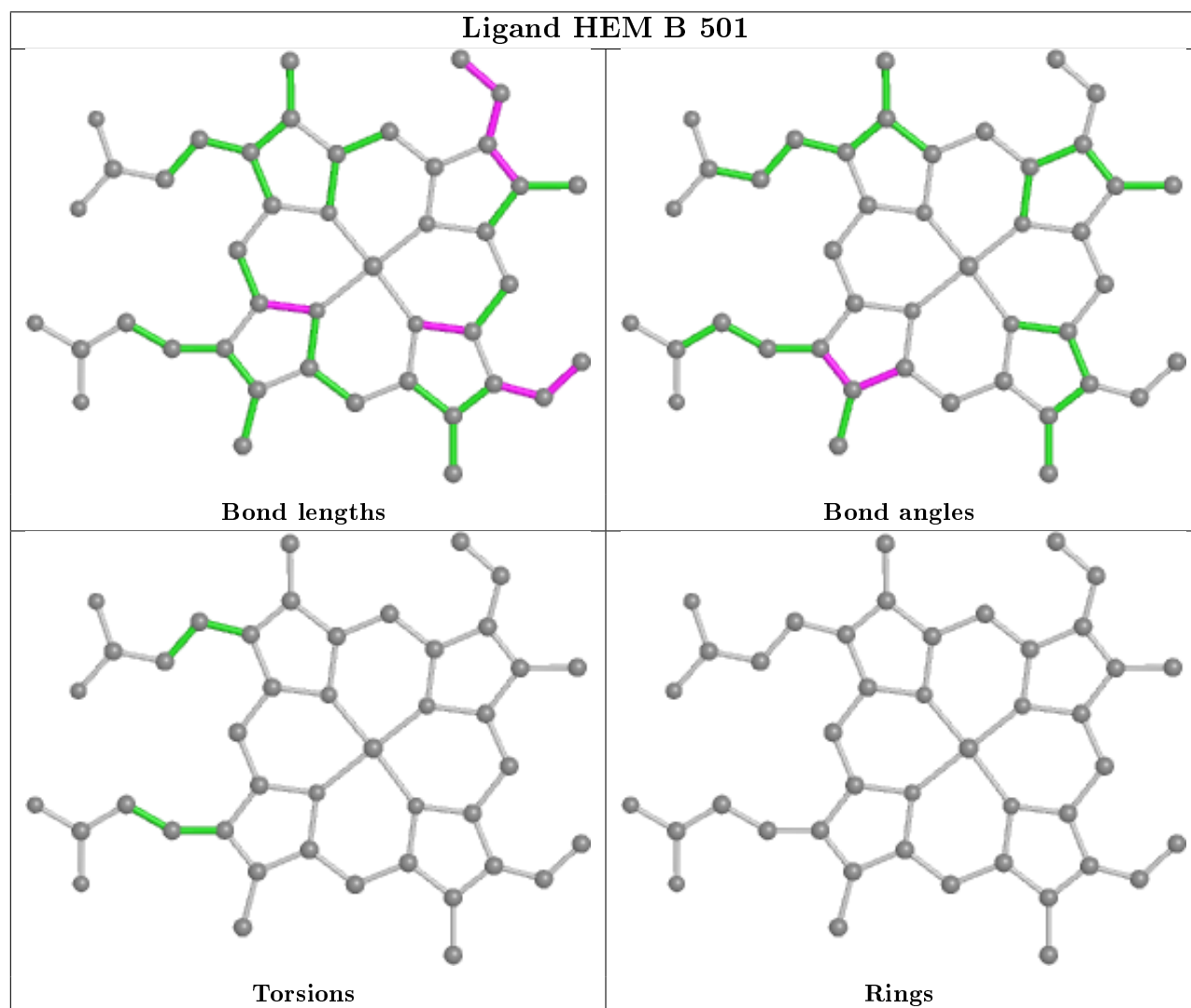
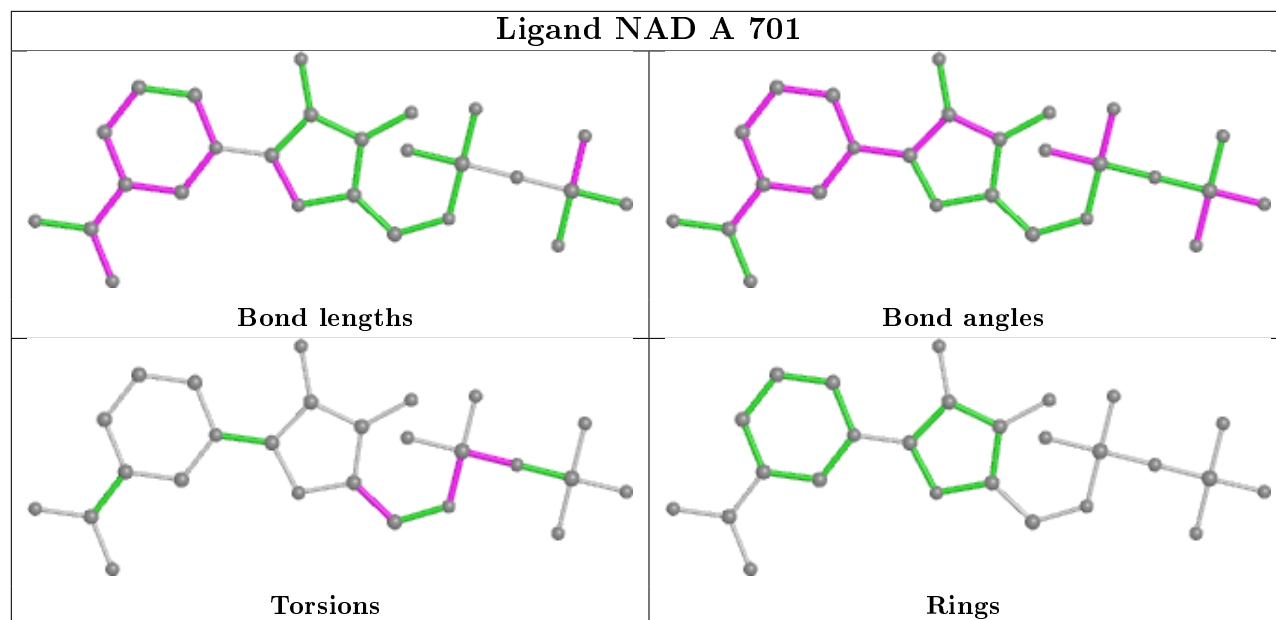
1 monomer is involved in 1 short contact:

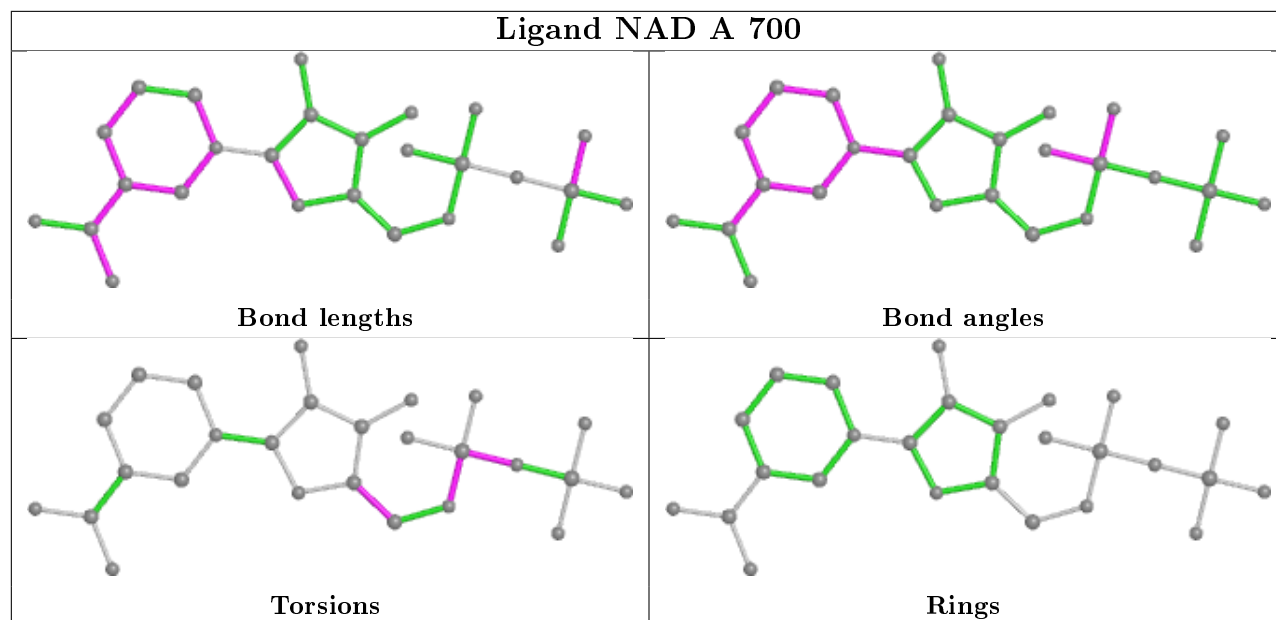
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	501	HEM	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be

highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.







5.7 Other polymers [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 [i](#)

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

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	346/347 (99%)	-0.54	4 (1%) 79 77	20, 33, 72, 90	0
1	B	347/347 (100%)	-0.68	2 (0%) 89 88	16, 30, 56, 85	0
All	All	693/694 (99%)	-0.61	6 (0%) 84 82	16, 32, 66, 90	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	346	LYS	3.5
1	A	346	LYS	2.8
1	B	0	HIS	2.8
1	A	1	HIS	2.5
1	A	38	THR	2.1
1	A	36	LYS	2.1

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

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

6.3 Carbohydrates [i](#)

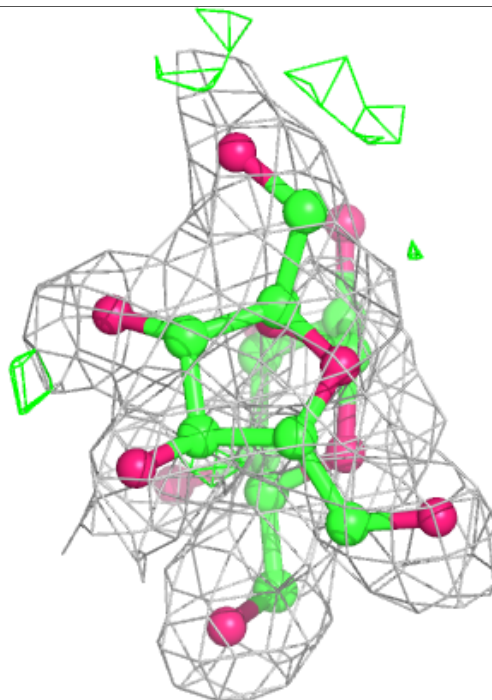
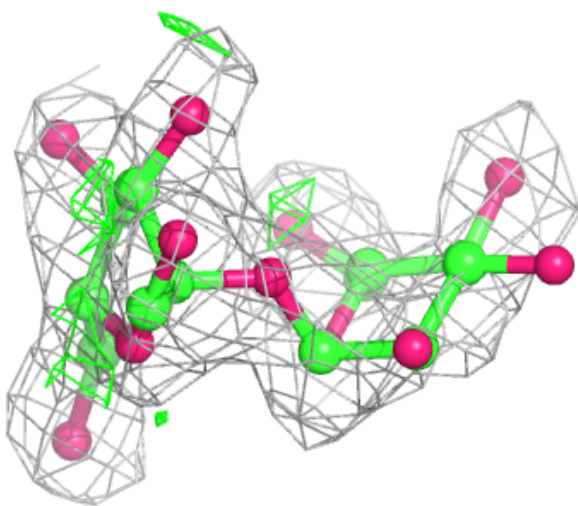
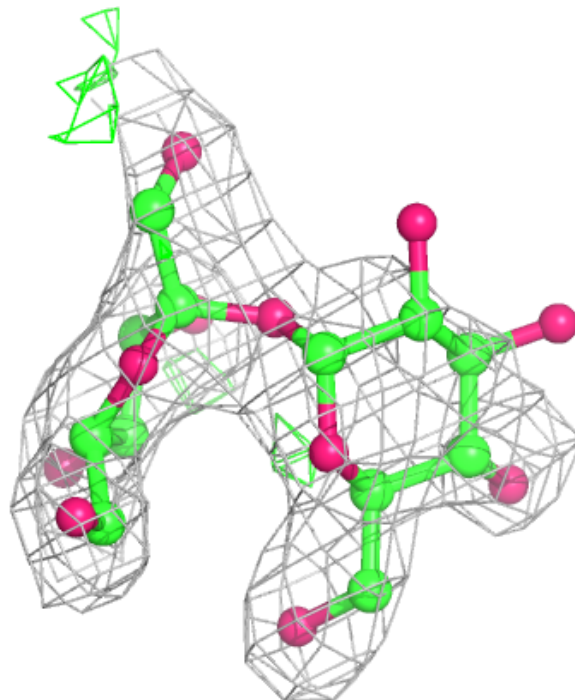
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists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

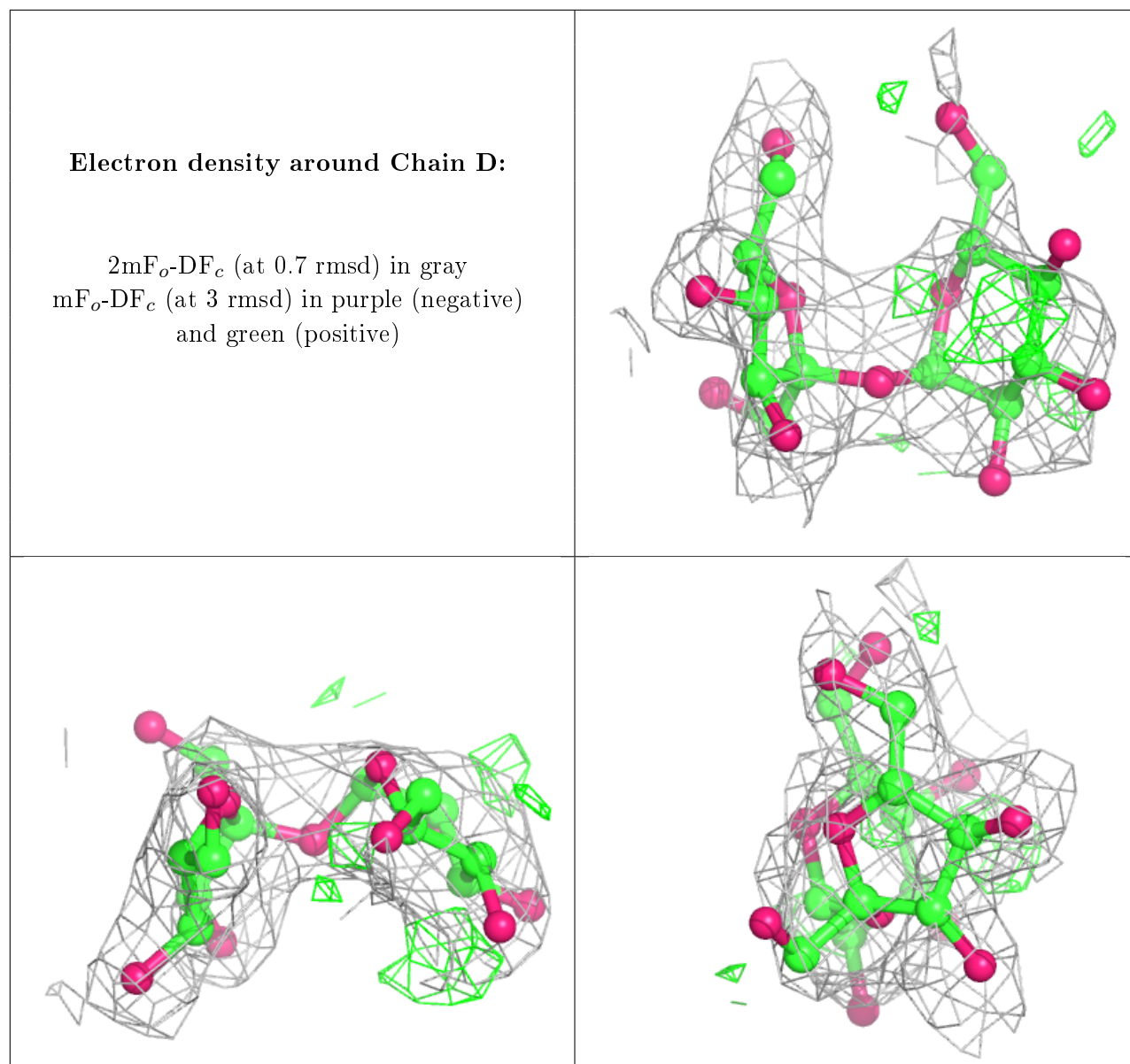
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	GLC	C	1	11/12	0.79	0.37	17,50,65,66	11
2	GLC	D	1	11/12	0.84	0.33	55,63,73,78	11
2	FRU	C	2	12/12	0.88	0.25	22,42,48,49	12
2	FRU	D	2	12/12	0.91	0.22	15,34,47,54	12

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 C:

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





6.4 Ligands [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.

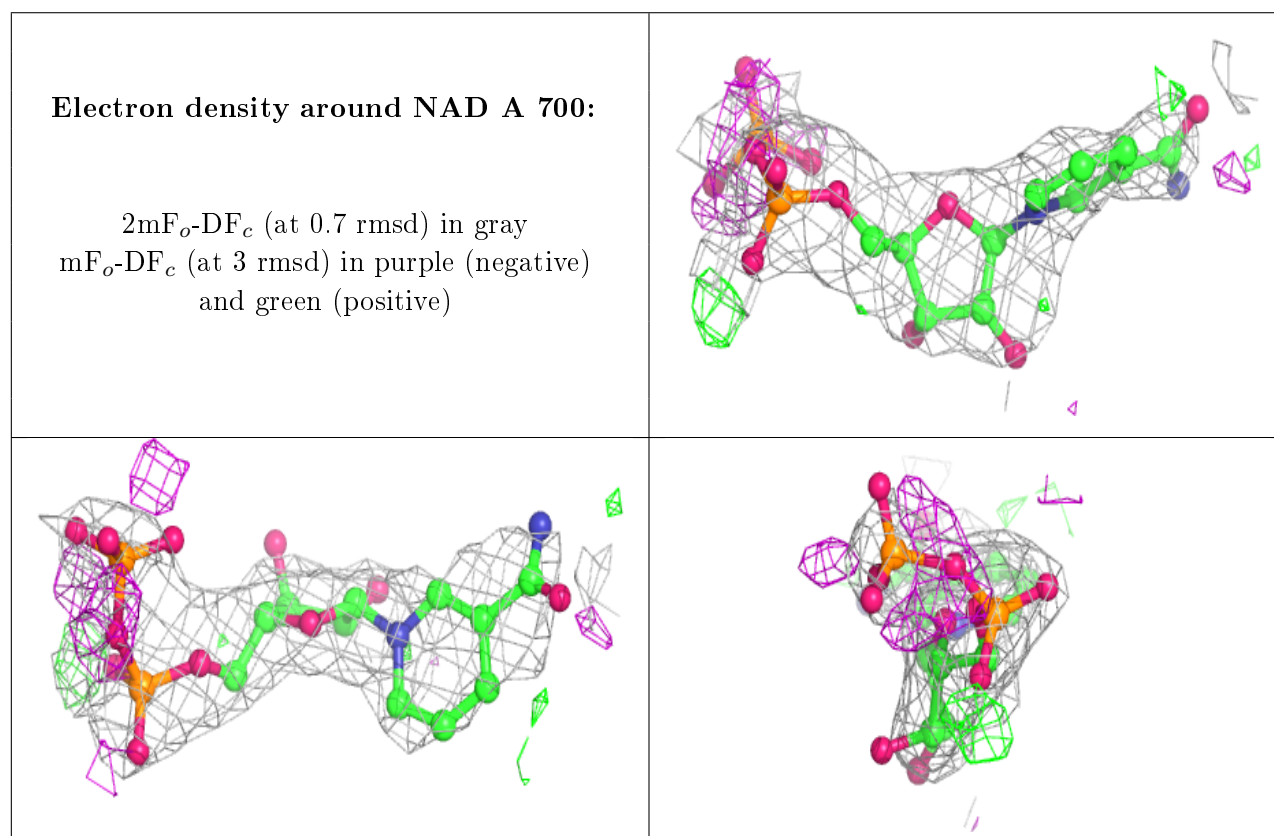
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	NAD	A	700	26/44	0.84	0.23	41,77,100,100	0
4	NAD	A	701	26/44	0.88	0.22	41,71,100,100	0
5	ITU	A	800	6/6	0.98	0.19	32,36,38,49	0
3	HEM	A	500	43/43	0.99	0.13	14,27,46,57	0
3	HEM	B	501	43/43	0.99	0.13	9,22,42,53	0

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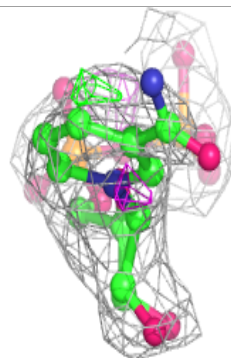
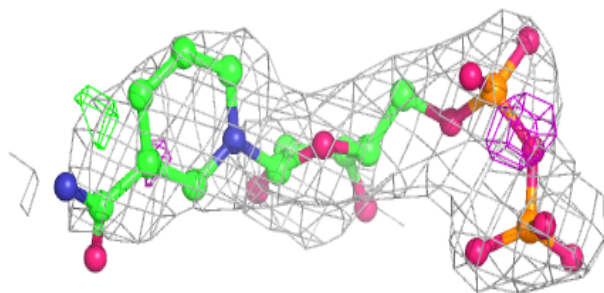
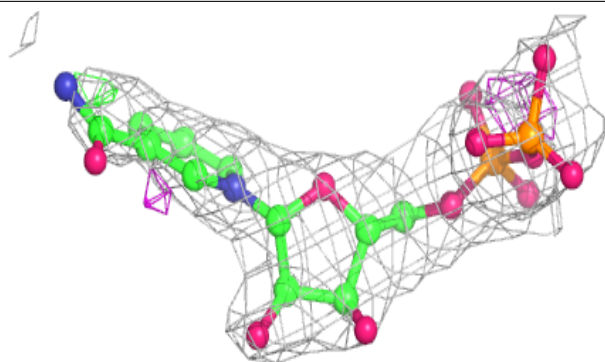
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
5	ITU	B	801	6/6	0.99	0.15	19,23,27,32	0

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.

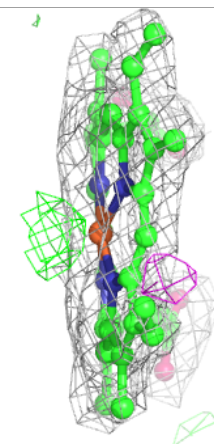
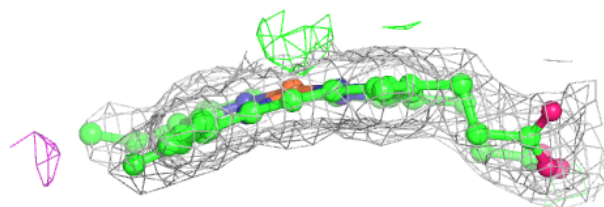
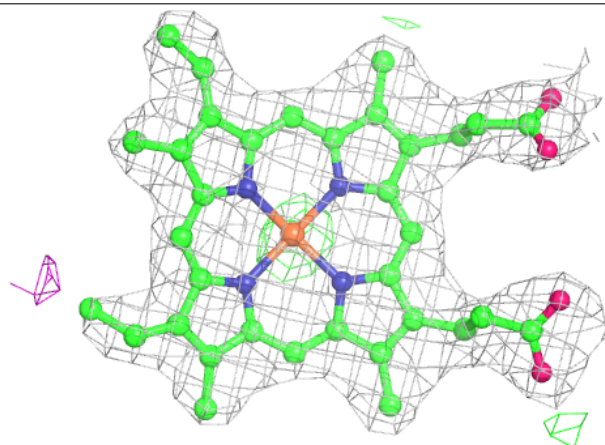


Electron density around NAD A 701:

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

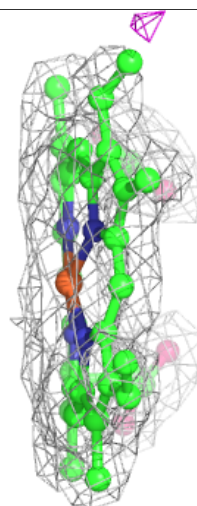
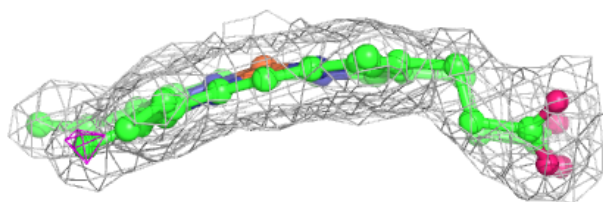
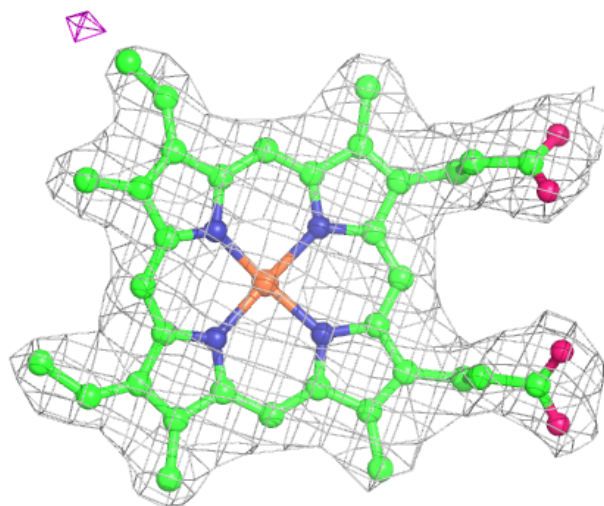
**Electron density around HEM A 500:**

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



Electron density around HEM 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)



6.5 Other polymers [i](#)

There are no such residues in this entry.