wwPDB X-ray Structure Validation Summary Report (i)

Jun 4, 2025 – 07:08 pm BST

PDB ID	:	9GQQ / pdb_00009gqq
Title	:	influenza neuraminidase N1/19 hybrid
Authors	:	Bowden, T.A.; Paesen, G.C.; Rijal, P.
Deposited on	:	2024-09-09
Resolution	:	1.78 Å(reported)

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

|		
MolProbity	:	4-5-2 with Phenix2.0rc1
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	2.0rc1
EDS	:	3.0
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.43.1

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 1.78 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	Percentile Ranks	Value
Rfree		0.214
Clashscore		4
Ramachandran outliers		0.3%
Sidechain outliers		1.2%
RSRZ outliers		5.4%

Worse   *Better*

▮ Percentile relative to all X-ray structures

▯ Percentile relative to X-ray structures of similar resolution
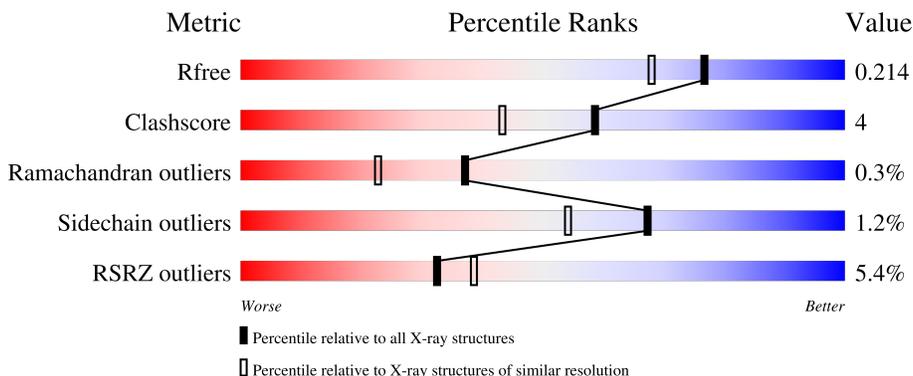
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	164625	1191 (1.78-1.78)
Clashscore	180529	1282 (1.78-1.78)
Ramachandran outliers	177936	1270 (1.78-1.78)
Sidechain outliers	177891	1270 (1.78-1.78)
RSRZ outliers	164620	1191 (1.78-1.78)

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 <=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	511	4% / 70% / 8% / 22%
1	B	511	5% / 69% / 9% / 22%
2	C	2	50% / 50%

2 Entry composition ⓘ

There are 5 unique types of molecules in this entry. The entry contains 6754 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 Neuraminidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	401	Total	C	N	O	S	0	0	0
			3082	1933	527	597	25			
1	B	401	Total	C	N	O	S	0	0	0
			3082	1933	527	597	25			

There are 282 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-41	MET	-	initiating methionine	UNP A0A024CWJ7
A	-40	ASN	-	expression tag	UNP A0A024CWJ7
A	-39	THR	-	expression tag	UNP A0A024CWJ7
A	-38	GLN	-	expression tag	UNP A0A024CWJ7
A	-37	ILE	-	expression tag	UNP A0A024CWJ7
A	-36	LEU	-	expression tag	UNP A0A024CWJ7
A	-35	VAL	-	expression tag	UNP A0A024CWJ7
A	-34	PHE	-	expression tag	UNP A0A024CWJ7
A	-33	ALA	-	expression tag	UNP A0A024CWJ7
A	-32	LEU	-	expression tag	UNP A0A024CWJ7
A	-31	ILE	-	expression tag	UNP A0A024CWJ7
A	-30	ALA	-	expression tag	UNP A0A024CWJ7
A	-29	ILE	-	expression tag	UNP A0A024CWJ7
A	-28	ILE	-	expression tag	UNP A0A024CWJ7
A	-27	PRO	-	expression tag	UNP A0A024CWJ7
A	-26	THR	-	expression tag	UNP A0A024CWJ7
A	-25	ASN	-	expression tag	UNP A0A024CWJ7
A	-24	ALA	-	expression tag	UNP A0A024CWJ7
A	-23	ASP	-	expression tag	UNP A0A024CWJ7
A	-22	LYS	-	expression tag	UNP A0A024CWJ7
A	-21	ILE	-	expression tag	UNP A0A024CWJ7
A	-20	SER	-	expression tag	UNP A0A024CWJ7
A	-19	ALA	-	expression tag	UNP A0A024CWJ7
A	-18	TRP	-	expression tag	UNP A0A024CWJ7
A	-17	SER	-	expression tag	UNP A0A024CWJ7

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-16	HIS	-	expression tag	UNP A0A024CWJ7
A	-15	PRO	-	expression tag	UNP A0A024CWJ7
A	-14	GLN	-	expression tag	UNP A0A024CWJ7
A	-13	PHE	-	expression tag	UNP A0A024CWJ7
A	-12	GLU	-	expression tag	UNP A0A024CWJ7
A	-11	LYS	-	expression tag	UNP A0A024CWJ7
A	-10	GLY	-	expression tag	UNP A0A024CWJ7
A	-9	GLY	-	expression tag	UNP A0A024CWJ7
A	-8	GLY	-	expression tag	UNP A0A024CWJ7
A	-7	GLY	-	expression tag	UNP A0A024CWJ7
A	-6	SER	-	expression tag	UNP A0A024CWJ7
A	-5	HIS	-	expression tag	UNP A0A024CWJ7
A	-4	HIS	-	expression tag	UNP A0A024CWJ7
A	-3	HIS	-	expression tag	UNP A0A024CWJ7
A	-2	HIS	-	expression tag	UNP A0A024CWJ7
A	-1	HIS	-	expression tag	UNP A0A024CWJ7
A	0	HIS	-	expression tag	UNP A0A024CWJ7
A	1	SER	-	expression tag	UNP A0A024CWJ7
A	2	SER	-	expression tag	UNP A0A024CWJ7
A	3	SER	-	expression tag	UNP A0A024CWJ7
A	4	GLY	-	expression tag	UNP A0A024CWJ7
A	5	SER	-	expression tag	UNP A0A024CWJ7
A	6	GLY	-	expression tag	UNP A0A024CWJ7
A	7	ALA	-	expression tag	UNP A0A024CWJ7
A	8	HIS	-	expression tag	UNP A0A024CWJ7
A	9	ILE	-	expression tag	UNP A0A024CWJ7
A	10	VAL	-	expression tag	UNP A0A024CWJ7
A	11	MET	-	expression tag	UNP A0A024CWJ7
A	12	VAL	-	expression tag	UNP A0A024CWJ7
A	13	ASP	-	expression tag	UNP A0A024CWJ7
A	14	ALA	-	expression tag	UNP A0A024CWJ7
A	15	TYR	-	expression tag	UNP A0A024CWJ7
A	16	LYS	-	expression tag	UNP A0A024CWJ7
A	17	PRO	-	expression tag	UNP A0A024CWJ7
A	18	THR	-	expression tag	UNP A0A024CWJ7
A	19	LYS	-	expression tag	UNP A0A024CWJ7
A	20	GLY	-	expression tag	UNP A0A024CWJ7
A	21	GLY	-	expression tag	UNP A0A024CWJ7
A	22	SER	-	expression tag	UNP A0A024CWJ7
A	23	GLY	-	expression tag	UNP A0A024CWJ7
A	24	GLY	-	expression tag	UNP A0A024CWJ7
A	25	GLY	-	expression tag	UNP A0A024CWJ7

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Chain	Residue	Modelled	Actual	Comment	Reference
A	26	GLY	-	expression tag	UNP A0A024CWJ7
A	27	SER	-	expression tag	UNP A0A024CWJ7
A	28	ILE	-	expression tag	UNP A0A024CWJ7
A	29	ILE	-	expression tag	UNP A0A024CWJ7
A	30	ASN	-	expression tag	UNP A0A024CWJ7
A	31	GLU	-	expression tag	UNP A0A024CWJ7
A	32	THR	-	expression tag	UNP A0A024CWJ7
A	33	ALA	-	expression tag	UNP A0A024CWJ7
A	34	ASP	-	expression tag	UNP A0A024CWJ7
A	35	ASP	-	expression tag	UNP A0A024CWJ7
A	36	ILE	-	expression tag	UNP A0A024CWJ7
A	37	VAL	-	expression tag	UNP A0A024CWJ7
A	38	TYR	-	expression tag	UNP A0A024CWJ7
A	39	ARG	-	expression tag	UNP A0A024CWJ7
A	40	LEU	-	expression tag	UNP A0A024CWJ7
A	41	THR	-	expression tag	UNP A0A024CWJ7
A	42	VAL	-	expression tag	UNP A0A024CWJ7
A	43	ILE	-	expression tag	UNP A0A024CWJ7
A	44	ILE	-	expression tag	UNP A0A024CWJ7
A	45	ASP	-	expression tag	UNP A0A024CWJ7
A	46	ASP	-	expression tag	UNP A0A024CWJ7
A	47	ARG	-	expression tag	UNP A0A024CWJ7
A	48	TYR	-	expression tag	UNP A0A024CWJ7
A	49	GLU	-	expression tag	UNP A0A024CWJ7
A	50	SER	-	expression tag	UNP A0A024CWJ7
A	51	LEU	-	expression tag	UNP A0A024CWJ7
A	52	LYS	-	expression tag	UNP A0A024CWJ7
A	53	ASN	-	expression tag	UNP A0A024CWJ7
A	54	LEU	-	expression tag	UNP A0A024CWJ7
A	55	ILE	-	expression tag	UNP A0A024CWJ7
A	56	THR	-	expression tag	UNP A0A024CWJ7
A	57	LEU	-	expression tag	UNP A0A024CWJ7
A	58	ARG	-	expression tag	UNP A0A024CWJ7
A	59	ALA	-	expression tag	UNP A0A024CWJ7
A	60	ASP	-	expression tag	UNP A0A024CWJ7
A	61	ARG	-	expression tag	UNP A0A024CWJ7
A	62	LEU	-	expression tag	UNP A0A024CWJ7
A	63	GLU	-	expression tag	UNP A0A024CWJ7
A	64	MET	-	expression tag	UNP A0A024CWJ7
A	65	ILE	-	expression tag	UNP A0A024CWJ7
A	66	ILE	-	expression tag	UNP A0A024CWJ7
A	67	ASN	-	expression tag	UNP A0A024CWJ7

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Chain	Residue	Modelled	Actual	Comment	Reference
A	68	ASP	-	expression tag	UNP A0A024CWJ7
A	69	ASN	-	expression tag	UNP A0A024CWJ7
A	70	VAL	-	expression tag	UNP A0A024CWJ7
A	71	SER	-	expression tag	UNP A0A024CWJ7
A	72	THR	-	expression tag	UNP A0A024CWJ7
A	73	ILE	-	expression tag	UNP A0A024CWJ7
A	74	LEU	-	expression tag	UNP A0A024CWJ7
A	75	ALA	-	expression tag	UNP A0A024CWJ7
A	76	GLY	-	expression tag	UNP A0A024CWJ7
A	77	GLY	-	expression tag	UNP A0A024CWJ7
A	78	SER	-	expression tag	UNP A0A024CWJ7
A	79	GLY	-	expression tag	UNP A0A024CWJ7
A	80	THR	-	expression tag	UNP A0A024CWJ7
A	81	GLY	-	expression tag	UNP A0A024CWJ7
A	149	ILE	VAL	conflict	UNP A0A024CWJ7
A	206	VAL	LEU	conflict	UNP A0A024CWJ7
A	222	LYS	ASN	conflict	UNP A0A024CWJ7
A	248	ASP	ASN	conflict	UNP A0A024CWJ7
A	269	MET	LEU	conflict	UNP A0A024CWJ7
A	270	LYS	ASN	conflict	UNP A0A024CWJ7
A	287	ASP	GLU	conflict	UNP A0A024CWJ7
A	314	MET	ILE	conflict	UNP A0A024CWJ7
A	331	LYS	GLY	conflict	UNP A0A024CWJ7
A	338	VAL	MET	conflict	UNP A0A024CWJ7
A	344	ASN	TYR	conflict	UNP A0A024CWJ7
A	365	ILE	THR	conflict	UNP A0A024CWJ7
A	369	LYS	SER	conflict	UNP A0A024CWJ7
A	395	GLY	ALA	conflict	UNP A0A024CWJ7
A	397	ASN	THR	conflict	UNP A0A024CWJ7
A	398	GLU	ASP	conflict	UNP A0A024CWJ7
A	412	MET	LEU	conflict	UNP A0A024CWJ7
A	432	GLU	LYS	conflict	UNP A0A024CWJ7
B	-41	MET	-	initiating methionine	UNP A0A024CWJ7
B	-40	ASN	-	expression tag	UNP A0A024CWJ7
B	-39	THR	-	expression tag	UNP A0A024CWJ7
B	-38	GLN	-	expression tag	UNP A0A024CWJ7
B	-37	ILE	-	expression tag	UNP A0A024CWJ7
B	-36	LEU	-	expression tag	UNP A0A024CWJ7
B	-35	VAL	-	expression tag	UNP A0A024CWJ7
B	-34	PHE	-	expression tag	UNP A0A024CWJ7
B	-33	ALA	-	expression tag	UNP A0A024CWJ7
B	-32	LEU	-	expression tag	UNP A0A024CWJ7

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-31	ILE	-	expression tag	UNP A0A024CWJ7
B	-30	ALA	-	expression tag	UNP A0A024CWJ7
B	-29	ILE	-	expression tag	UNP A0A024CWJ7
B	-28	ILE	-	expression tag	UNP A0A024CWJ7
B	-27	PRO	-	expression tag	UNP A0A024CWJ7
B	-26	THR	-	expression tag	UNP A0A024CWJ7
B	-25	ASN	-	expression tag	UNP A0A024CWJ7
B	-24	ALA	-	expression tag	UNP A0A024CWJ7
B	-23	ASP	-	expression tag	UNP A0A024CWJ7
B	-22	LYS	-	expression tag	UNP A0A024CWJ7
B	-21	ILE	-	expression tag	UNP A0A024CWJ7
B	-20	SER	-	expression tag	UNP A0A024CWJ7
B	-19	ALA	-	expression tag	UNP A0A024CWJ7
B	-18	TRP	-	expression tag	UNP A0A024CWJ7
B	-17	SER	-	expression tag	UNP A0A024CWJ7
B	-16	HIS	-	expression tag	UNP A0A024CWJ7
B	-15	PRO	-	expression tag	UNP A0A024CWJ7
B	-14	GLN	-	expression tag	UNP A0A024CWJ7
B	-13	PHE	-	expression tag	UNP A0A024CWJ7
B	-12	GLU	-	expression tag	UNP A0A024CWJ7
B	-11	LYS	-	expression tag	UNP A0A024CWJ7
B	-10	GLY	-	expression tag	UNP A0A024CWJ7
B	-9	GLY	-	expression tag	UNP A0A024CWJ7
B	-8	GLY	-	expression tag	UNP A0A024CWJ7
B	-7	GLY	-	expression tag	UNP A0A024CWJ7
B	-6	SER	-	expression tag	UNP A0A024CWJ7
B	-5	HIS	-	expression tag	UNP A0A024CWJ7
B	-4	HIS	-	expression tag	UNP A0A024CWJ7
B	-3	HIS	-	expression tag	UNP A0A024CWJ7
B	-2	HIS	-	expression tag	UNP A0A024CWJ7
B	-1	HIS	-	expression tag	UNP A0A024CWJ7
B	0	HIS	-	expression tag	UNP A0A024CWJ7
B	1	SER	-	expression tag	UNP A0A024CWJ7
B	2	SER	-	expression tag	UNP A0A024CWJ7
B	3	SER	-	expression tag	UNP A0A024CWJ7
B	4	GLY	-	expression tag	UNP A0A024CWJ7
B	5	SER	-	expression tag	UNP A0A024CWJ7
B	6	GLY	-	expression tag	UNP A0A024CWJ7
B	7	ALA	-	expression tag	UNP A0A024CWJ7
B	8	HIS	-	expression tag	UNP A0A024CWJ7
B	9	ILE	-	expression tag	UNP A0A024CWJ7
B	10	VAL	-	expression tag	UNP A0A024CWJ7

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Chain	Residue	Modelled	Actual	Comment	Reference
B	11	MET	-	expression tag	UNP A0A024CWJ7
B	12	VAL	-	expression tag	UNP A0A024CWJ7
B	13	ASP	-	expression tag	UNP A0A024CWJ7
B	14	ALA	-	expression tag	UNP A0A024CWJ7
B	15	TYR	-	expression tag	UNP A0A024CWJ7
B	16	LYS	-	expression tag	UNP A0A024CWJ7
B	17	PRO	-	expression tag	UNP A0A024CWJ7
B	18	THR	-	expression tag	UNP A0A024CWJ7
B	19	LYS	-	expression tag	UNP A0A024CWJ7
B	20	GLY	-	expression tag	UNP A0A024CWJ7
B	21	GLY	-	expression tag	UNP A0A024CWJ7
B	22	SER	-	expression tag	UNP A0A024CWJ7
B	23	GLY	-	expression tag	UNP A0A024CWJ7
B	24	GLY	-	expression tag	UNP A0A024CWJ7
B	25	GLY	-	expression tag	UNP A0A024CWJ7
B	26	GLY	-	expression tag	UNP A0A024CWJ7
B	27	SER	-	expression tag	UNP A0A024CWJ7
B	28	ILE	-	expression tag	UNP A0A024CWJ7
B	29	ILE	-	expression tag	UNP A0A024CWJ7
B	30	ASN	-	expression tag	UNP A0A024CWJ7
B	31	GLU	-	expression tag	UNP A0A024CWJ7
B	32	THR	-	expression tag	UNP A0A024CWJ7
B	33	ALA	-	expression tag	UNP A0A024CWJ7
B	34	ASP	-	expression tag	UNP A0A024CWJ7
B	35	ASP	-	expression tag	UNP A0A024CWJ7
B	36	ILE	-	expression tag	UNP A0A024CWJ7
B	37	VAL	-	expression tag	UNP A0A024CWJ7
B	38	TYR	-	expression tag	UNP A0A024CWJ7
B	39	ARG	-	expression tag	UNP A0A024CWJ7
B	40	LEU	-	expression tag	UNP A0A024CWJ7
B	41	THR	-	expression tag	UNP A0A024CWJ7
B	42	VAL	-	expression tag	UNP A0A024CWJ7
B	43	ILE	-	expression tag	UNP A0A024CWJ7
B	44	ILE	-	expression tag	UNP A0A024CWJ7
B	45	ASP	-	expression tag	UNP A0A024CWJ7
B	46	ASP	-	expression tag	UNP A0A024CWJ7
B	47	ARG	-	expression tag	UNP A0A024CWJ7
B	48	TYR	-	expression tag	UNP A0A024CWJ7
B	49	GLU	-	expression tag	UNP A0A024CWJ7
B	50	SER	-	expression tag	UNP A0A024CWJ7
B	51	LEU	-	expression tag	UNP A0A024CWJ7
B	52	LYS	-	expression tag	UNP A0A024CWJ7

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Chain	Residue	Modelled	Actual	Comment	Reference
B	53	ASN	-	expression tag	UNP A0A024CWJ7
B	54	LEU	-	expression tag	UNP A0A024CWJ7
B	55	ILE	-	expression tag	UNP A0A024CWJ7
B	56	THR	-	expression tag	UNP A0A024CWJ7
B	57	LEU	-	expression tag	UNP A0A024CWJ7
B	58	ARG	-	expression tag	UNP A0A024CWJ7
B	59	ALA	-	expression tag	UNP A0A024CWJ7
B	60	ASP	-	expression tag	UNP A0A024CWJ7
B	61	ARG	-	expression tag	UNP A0A024CWJ7
B	62	LEU	-	expression tag	UNP A0A024CWJ7
B	63	GLU	-	expression tag	UNP A0A024CWJ7
B	64	MET	-	expression tag	UNP A0A024CWJ7
B	65	ILE	-	expression tag	UNP A0A024CWJ7
B	66	ILE	-	expression tag	UNP A0A024CWJ7
B	67	ASN	-	expression tag	UNP A0A024CWJ7
B	68	ASP	-	expression tag	UNP A0A024CWJ7
B	69	ASN	-	expression tag	UNP A0A024CWJ7
B	70	VAL	-	expression tag	UNP A0A024CWJ7
B	71	SER	-	expression tag	UNP A0A024CWJ7
B	72	THR	-	expression tag	UNP A0A024CWJ7
B	73	ILE	-	expression tag	UNP A0A024CWJ7
B	74	LEU	-	expression tag	UNP A0A024CWJ7
B	75	ALA	-	expression tag	UNP A0A024CWJ7
B	76	GLY	-	expression tag	UNP A0A024CWJ7
B	77	GLY	-	expression tag	UNP A0A024CWJ7
B	78	SER	-	expression tag	UNP A0A024CWJ7
B	79	GLY	-	expression tag	UNP A0A024CWJ7
B	80	THR	-	expression tag	UNP A0A024CWJ7
B	81	GLY	-	expression tag	UNP A0A024CWJ7
B	149	ILE	VAL	conflict	UNP A0A024CWJ7
B	206	VAL	LEU	conflict	UNP A0A024CWJ7
B	222	LYS	ASN	conflict	UNP A0A024CWJ7
B	248	ASP	ASN	conflict	UNP A0A024CWJ7
B	269	MET	LEU	conflict	UNP A0A024CWJ7
B	270	LYS	ASN	conflict	UNP A0A024CWJ7
B	287	ASP	GLU	conflict	UNP A0A024CWJ7
B	314	MET	ILE	conflict	UNP A0A024CWJ7
B	331	LYS	GLY	conflict	UNP A0A024CWJ7
B	338	VAL	MET	conflict	UNP A0A024CWJ7
B	344	ASN	TYR	conflict	UNP A0A024CWJ7
B	365	ILE	THR	conflict	UNP A0A024CWJ7
B	369	LYS	SER	conflict	UNP A0A024CWJ7

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Chain	Residue	Modelled	Actual	Comment	Reference
B	395	GLY	ALA	conflict	UNP A0A024CWJ7
B	397	ASN	THR	conflict	UNP A0A024CWJ7
B	398	GLU	ASP	conflict	UNP A0A024CWJ7
B	412	MET	LEU	conflict	UNP A0A024CWJ7
B	432	GLU	LYS	conflict	UNP A0A024CWJ7

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

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

- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Ca		
3	A	2	2	2	0	0
			Total	Ca		
3	B	2	2	2	0	0

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$) (labeled as "Ligand of Interest" by depositor).
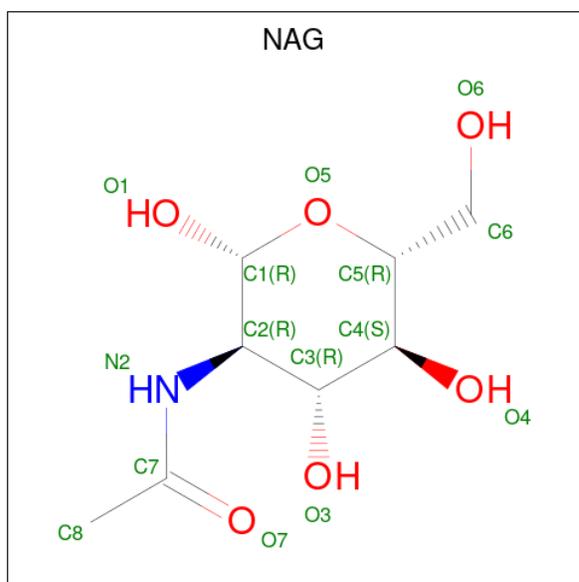
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			14	8	1	5		
4	A	1	Total	C	N	O	0	0
			14	8	1	5		
4	A	1	Total	C	N	O	0	0
			14	8	1	5		
4	B	1	Total	C	N	O	0	0
			14	8	1	5		
4	B	1	Total	C	N	O	0	0
			14	8	1	5		
4	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	236	Total	O	0	0
			236	236		
5	B	238	Total	O	0	0
			238	238		

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: Neuraminidase

Chain A:

• Molecule 1: Neuraminidase

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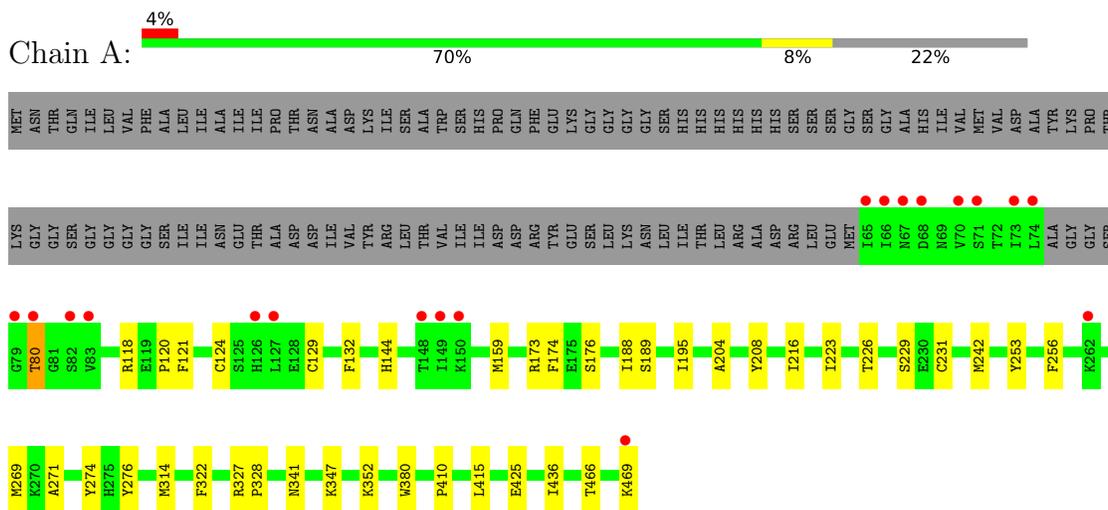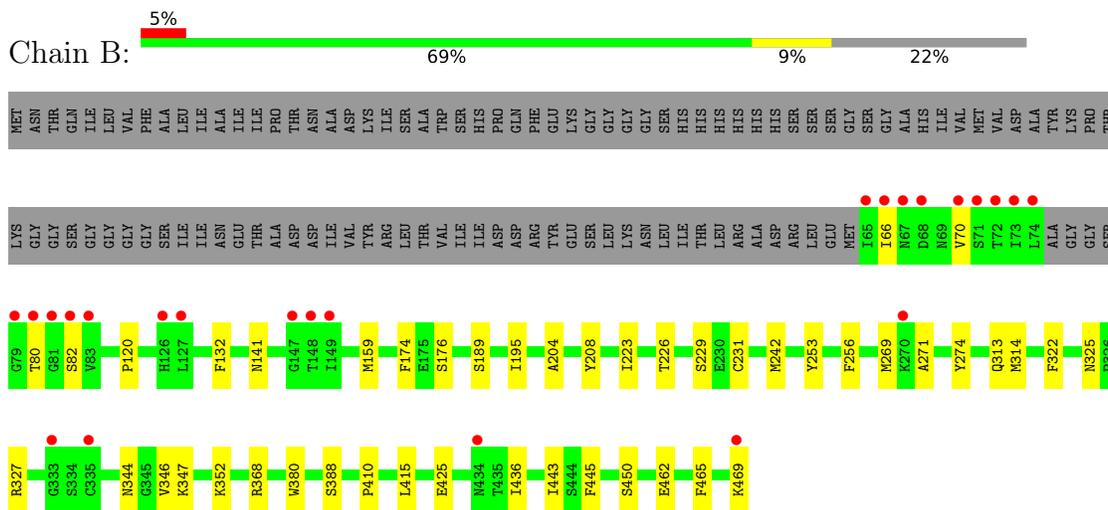
• Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain C:

NAG1 NAG2

4 Data and refinement statistics ⓘ

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α, β, γ	115.41Å 234.37Å 115.65Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	64.69 – 1.78 64.69 – 1.78	Depositor EDS
% Data completeness (in resolution range)	38.9 (64.69-1.78) 36.3 (64.69-1.78)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	0.51 (at 1.77Å)	Xtriage
Refinement program	PHENIX 1.21.1_5286	Depositor
R, R_{free}	0.189 , 0.215 0.189 , 0.214	Depositor DCC
R_{free} test set	56538 reflections (96.56%)	wwPDB-VP
Wilson B-factor (Å²)	11.9	Xtriage
Anisotropy	4.107	Xtriage
Bulk solvent k_{sol}(e/Å³), B_{sol}(Å²)	0.37 , 39.6	EDS
L-test for twinning[2]	$< |L| > = 0.47, < L^2 > = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.87	EDS
Total number of atoms	6754	wwPDB-VP
Average B, all atoms (Å²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.03% of the height of the origin peak. No significant pseudotranslation is detected.*

---

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: NAG, CA

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	0.17	0/3163	0.42	0/4291
1	B	0.17	0/3163	0.42	0/4291
All	All	0.17	0/6326	0.42	0/8582

There are no bond length outliers.

There are no bond angle outliers.

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	3082	0	2910	25	0
1	B	3082	0	2911	28	0
2	C	28	0	25	0	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
4	A	42	0	39	1	0
4	B	42	0	39	1	0
5	A	236	0	0	0	0
5	B	238	0	0	3	0
All	All	6754	0	5924	52	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 4.

The worst 5 of 52 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:159:MET:HG2	1:A:174:PHE:HA	1.71	0.73
1:B:346:VAL:HG12	1:B:368:ARG:HE	1.55	0.70
1:B:159:MET:HG2	1:B:174:PHE:HA	1.82	0.60
1:B:195:ILE:HG12	1:B:204:ALA:HB2	1.82	0.60
1:A:229:SER:HB3	1:A:347:LYS:HE2	1.84	0.59

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	397/511 (78%)	382 (96%)	14 (4%)	1 (0%)	37	23
1	B	397/511 (78%)	382 (96%)	14 (4%)	1 (0%)	37	23
All	All	794/1022 (78%)	764 (96%)	28 (4%)	2 (0%)	37	23

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	223	ILE
1	B	223	ILE

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	343/430 (80%)	340 (99%)	3 (1%)	75	65
1	B	343/430 (80%)	338 (98%)	5 (2%)	60	44
All	All	686/860 (80%)	678 (99%)	8 (1%)	67	54

5 of 8 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	388	SER
1	B	231	CYS
1	B	82	SER
1	B	80	THR
1	B	176	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 8 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	341	ASN
1	B	329	ASN
1	B	221	ASN
1	B	67	ASN
1	B	309	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 ⓘ

2 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	C	1	1,2	14,14,15	0.65	0	17,19,21	0.96	0
2	NAG	C	2	2	14,14,15	0.70	0	17,19,21	1.36	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
2	NAG	C	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	C	2	2	-	1/6/23/26	0/1/1/1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	C	2	NAG	C2-N2-C7	3.54	127.95	122.90

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	1	NAG	C8-C7-N2-C2
2	C	1	NAG	O7-C7-N2-C2
2	C	2	NAG	C1-C2-N2-C7

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.

Oligosaccharide Chain C

Bond lengths

Bond angles

Torsions

Rings
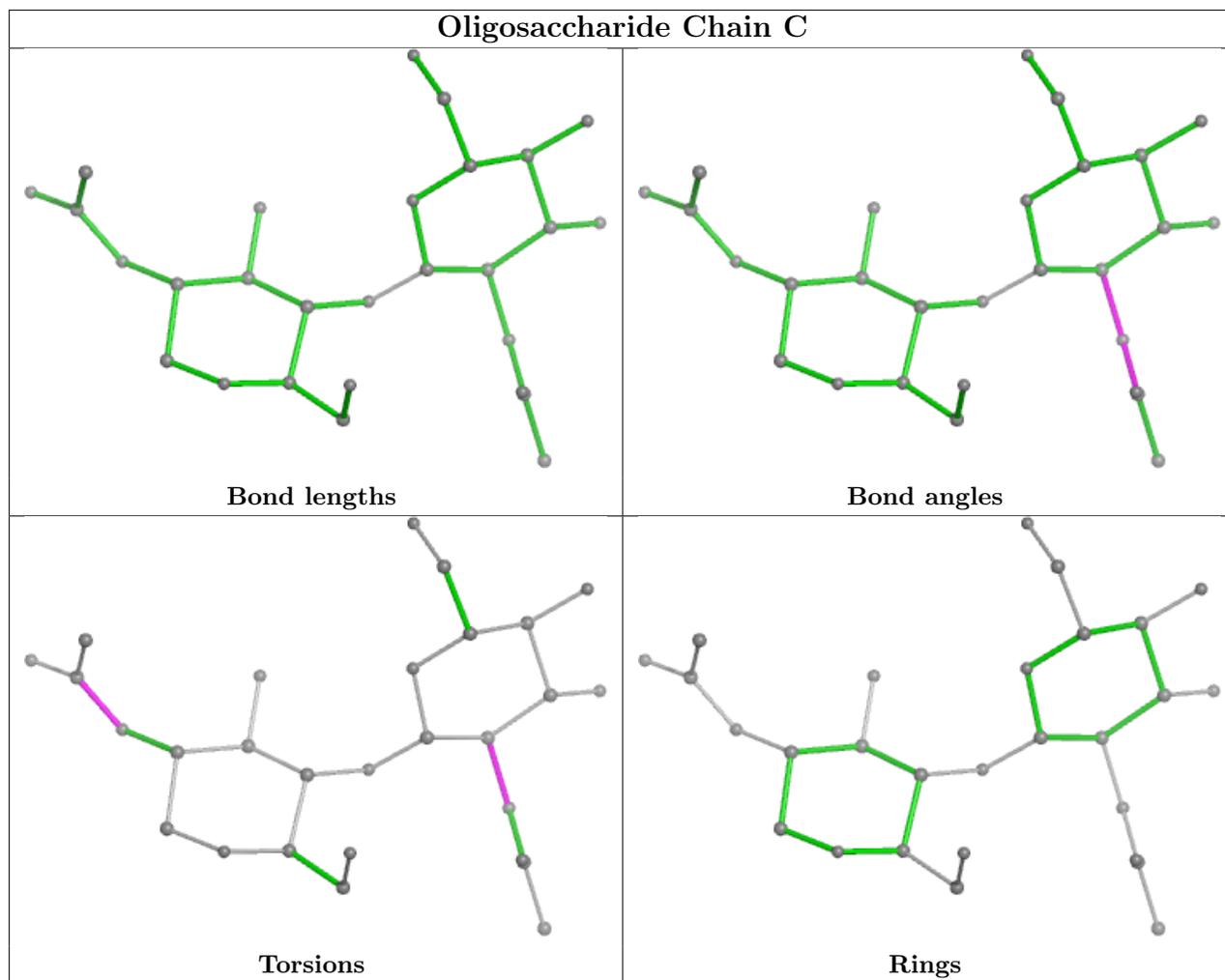
5.6 Ligand geometry ⓘ

Of 10 ligands modelled in this entry, 4 are monoatomic - leaving 6 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	B	504	1	14,14,15	0.74	0	17,19,21	0.96	1 (5%)
4	NAG	A	505	1	14,14,15	0.70	0	17,19,21	1.85	3 (17%)
4	NAG	A	503	1	14,14,15	0.72	0	17,19,21	0.89	0
4	NAG	B	503	1	14,14,15	0.74	0	17,19,21	1.25	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	#|Z| > 2	Counts	RMSZ	#|Z| > 2
4	NAG	B	505	1	14,14,15	0.74	0	17,19,21	1.00	1 (5%)
4	NAG	A	504	1	14,14,15	0.67	0	17,19,21	1.17	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	B	504	1	-	0/6/23/26	0/1/1/1
4	NAG	A	505	1	-	1/6/23/26	0/1/1/1
4	NAG	A	503	1	-	3/6/23/26	0/1/1/1
4	NAG	B	503	1	-	1/6/23/26	0/1/1/1
4	NAG	B	505	1	-	0/6/23/26	0/1/1/1
4	NAG	A	504	1	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

The worst 5 of 7 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	A	505	NAG	C2-N2-C7	5.78	131.14	122.90
4	A	504	NAG	C1-O5-C5	3.44	116.85	112.19
4	B	503	NAG	C2-N2-C7	3.04	127.24	122.90
4	A	505	NAG	O7-C7-N2	2.42	126.39	121.95
4	B	505	NAG	C1-O5-C5	2.21	115.18	112.19

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	505	NAG	C1-C2-N2-C7
4	A	503	NAG	C8-C7-N2-C2
4	A	503	NAG	O7-C7-N2-C2
4	B	503	NAG	C3-C2-N2-C7
4	A	503	NAG	O5-C5-C6-O6

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	505	NAG	1	0
4	A	504	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 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 then 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 NAG B 504

Bond lengths

Bond angles

Torsions

Rings

Ligand NAG A 505

Bond lengths

Bond angles

Torsions

Rings

Ligand NAG A 503

Bond lengths

Bond angles

Torsions

Rings

Ligand NAG B 503

Bond lengths

Bond angles

Torsions

Rings

Ligand NAG B 505

Bond lengths

Bond angles

Torsions

Rings

Ligand NAG A 504

Bond lengths

Bond angles

Torsions

Rings
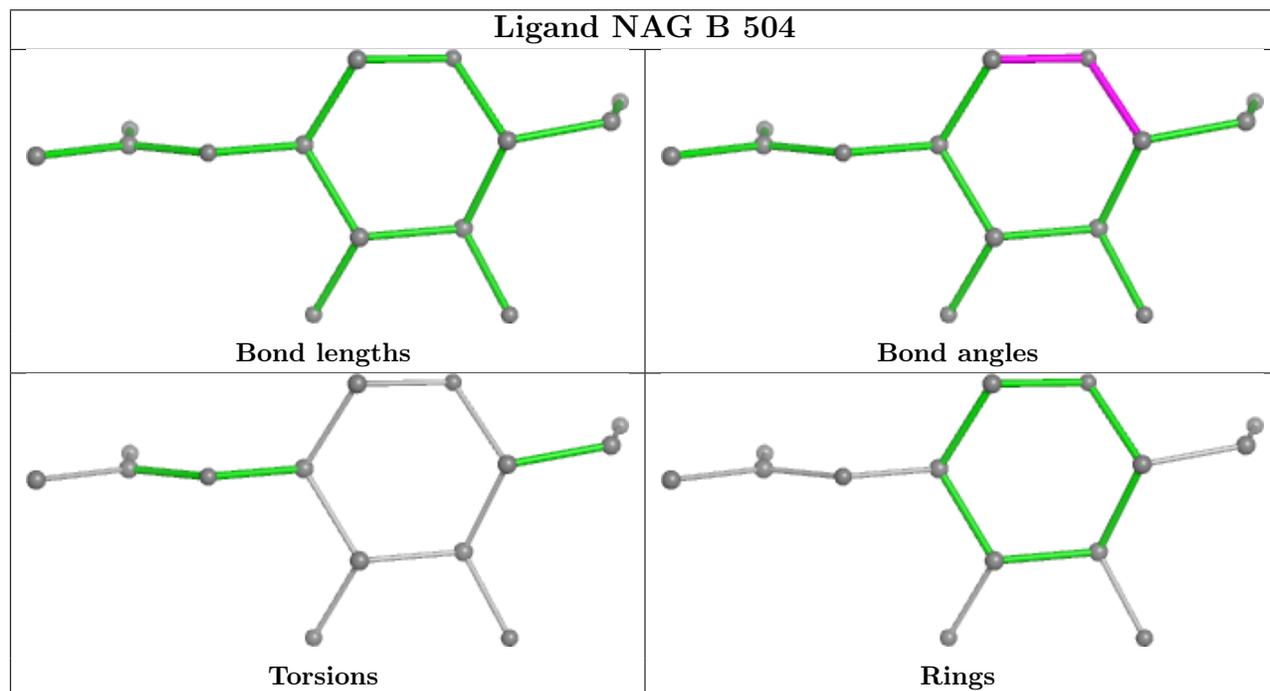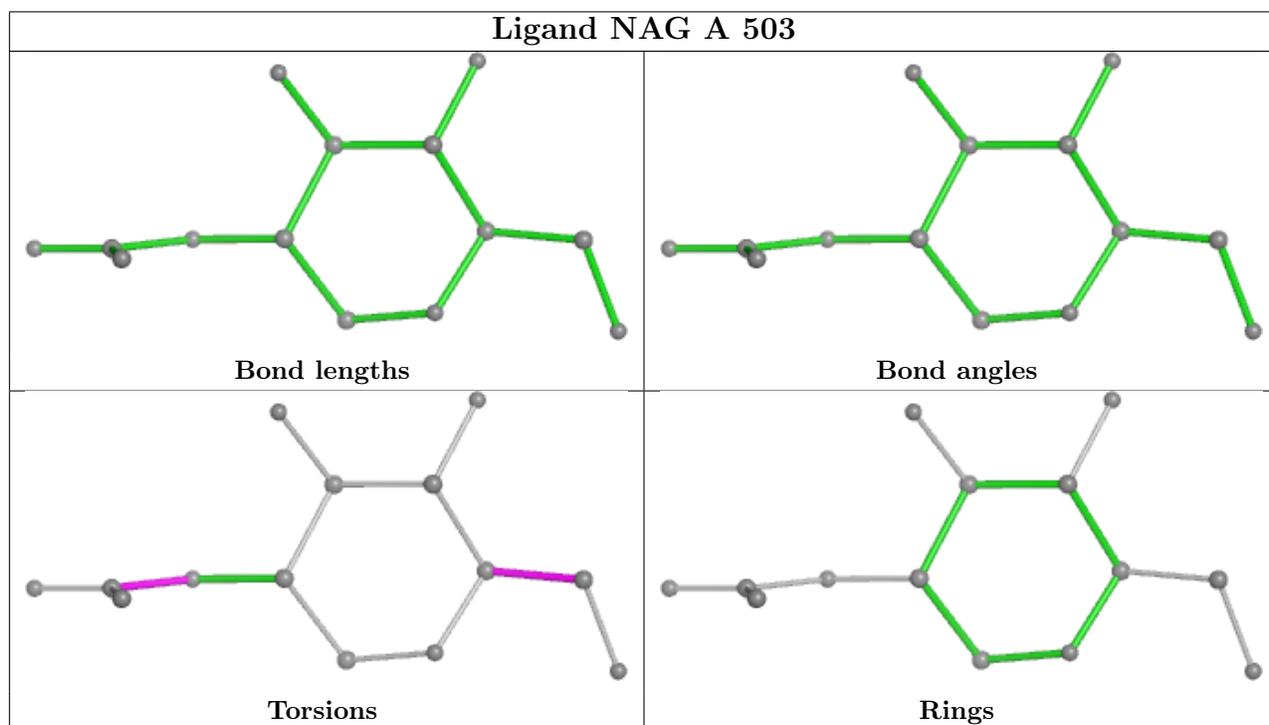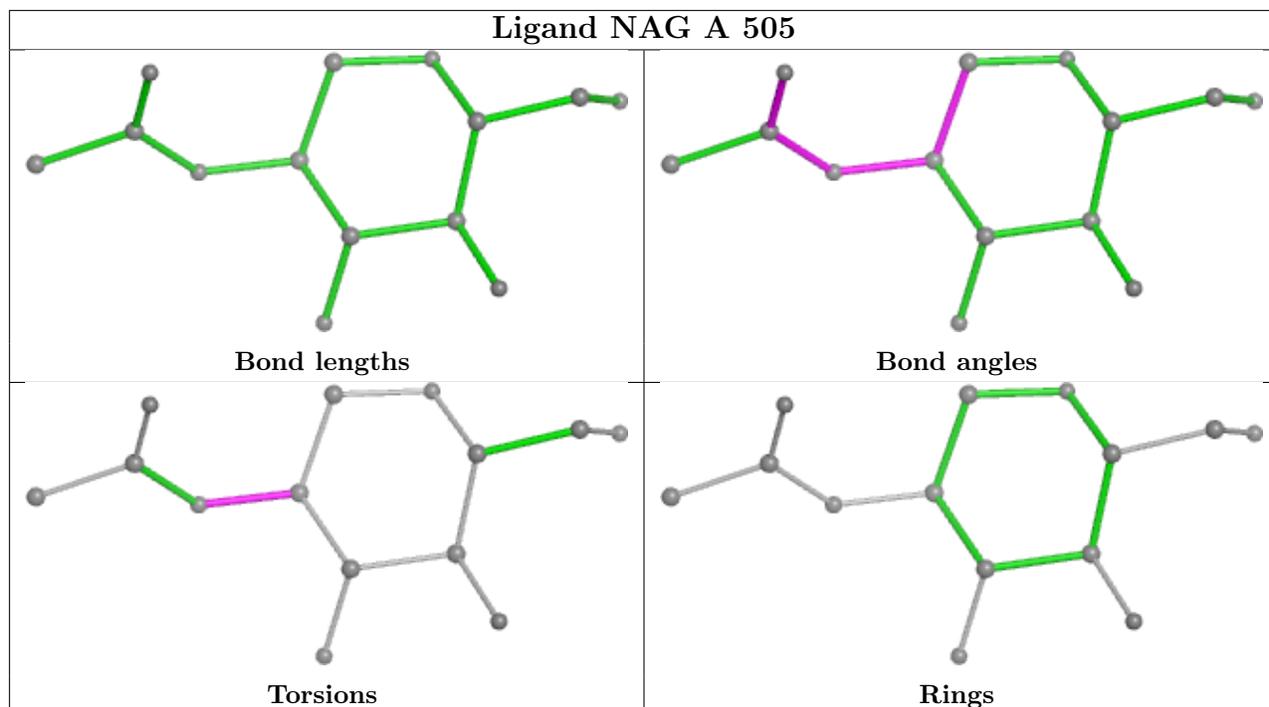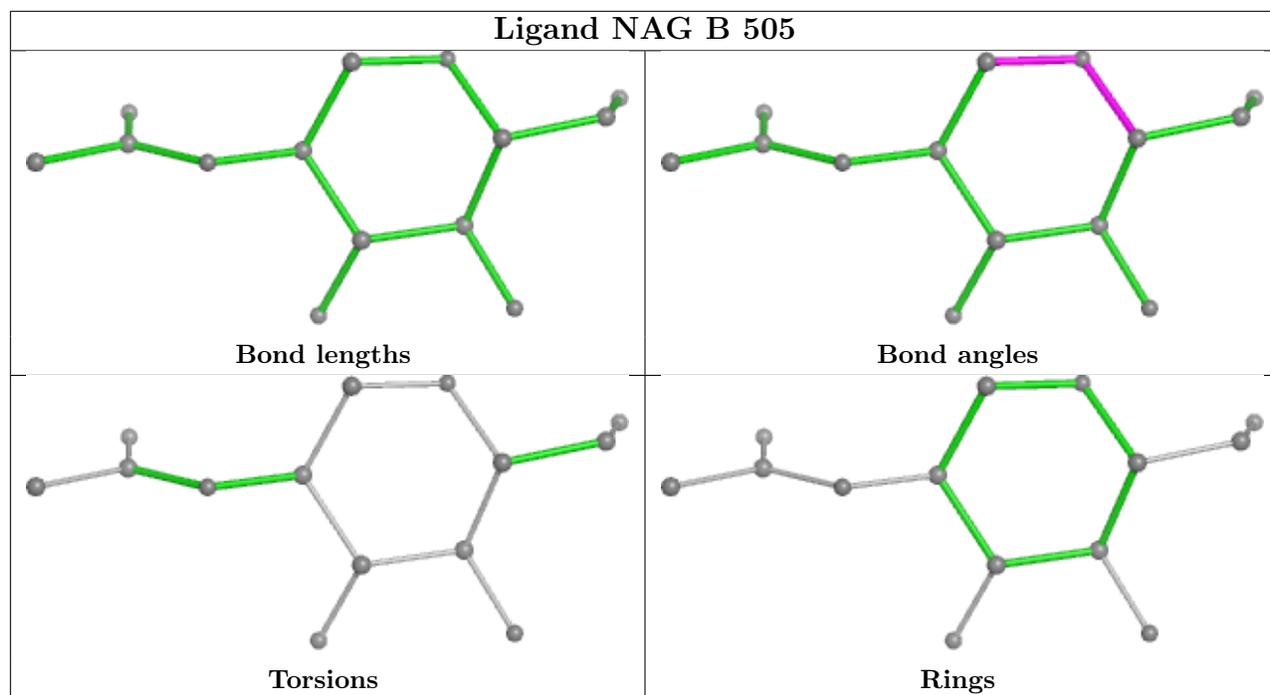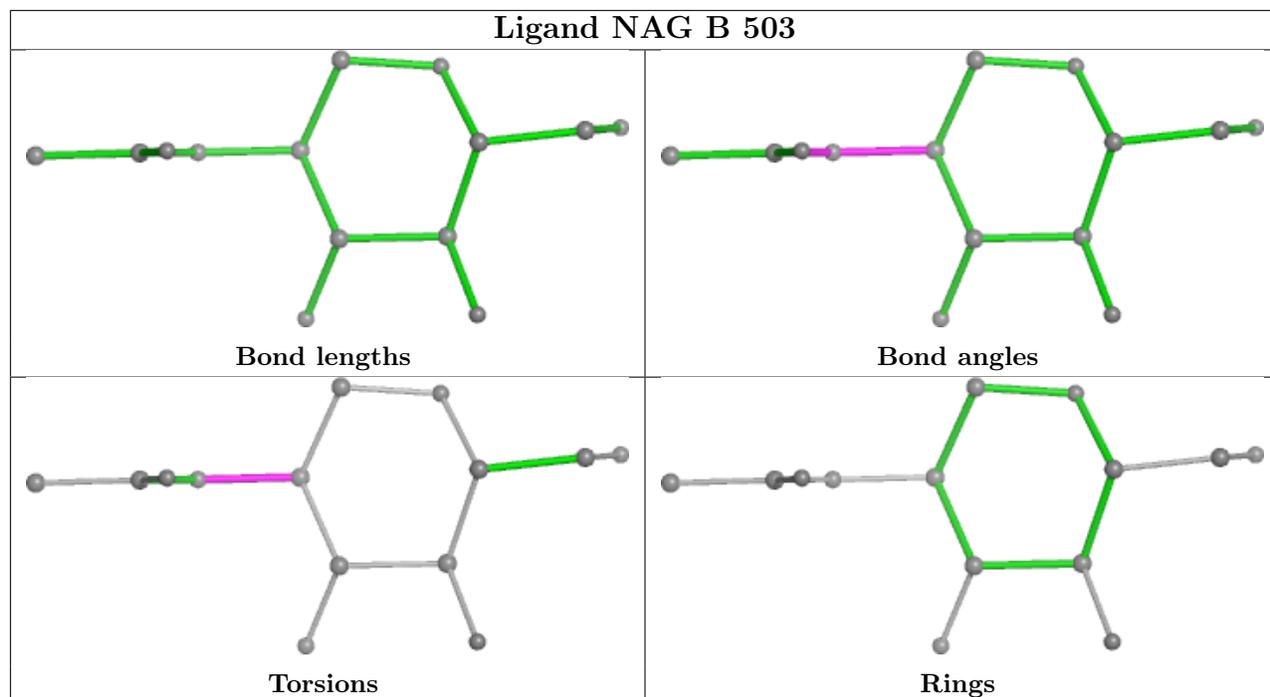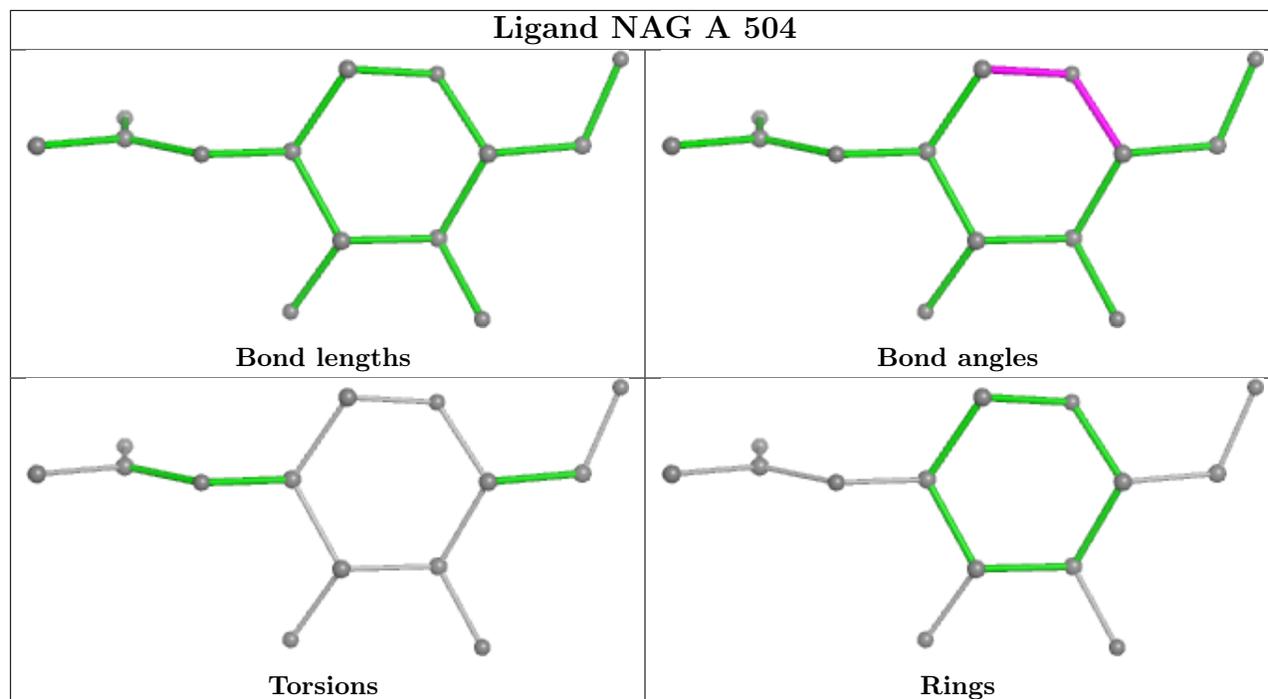
5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	401/511 (78%)	0.37	19 (4%)	37	43	6, 15, 39, 69	0
1	B	401/511 (78%)	0.36	24 (5%)	29	34	6, 15, 40, 73	0
All	All	802/1022 (78%)	0.36	43 (5%)	32	38	6, 15, 40, 73	0

The worst 5 of 43 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	74	LEU	5.5
1	B	79	GLY	4.5
1	A	79	GLY	4.5
1	A	67	ASN	4.4
1	B	65	ILE	4.4

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	NAG	C	2	14/15	0.68	0.21	34,65,84,86	0
2	NAG	C	1	14/15	0.92	0.10	18,36,52,64	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.
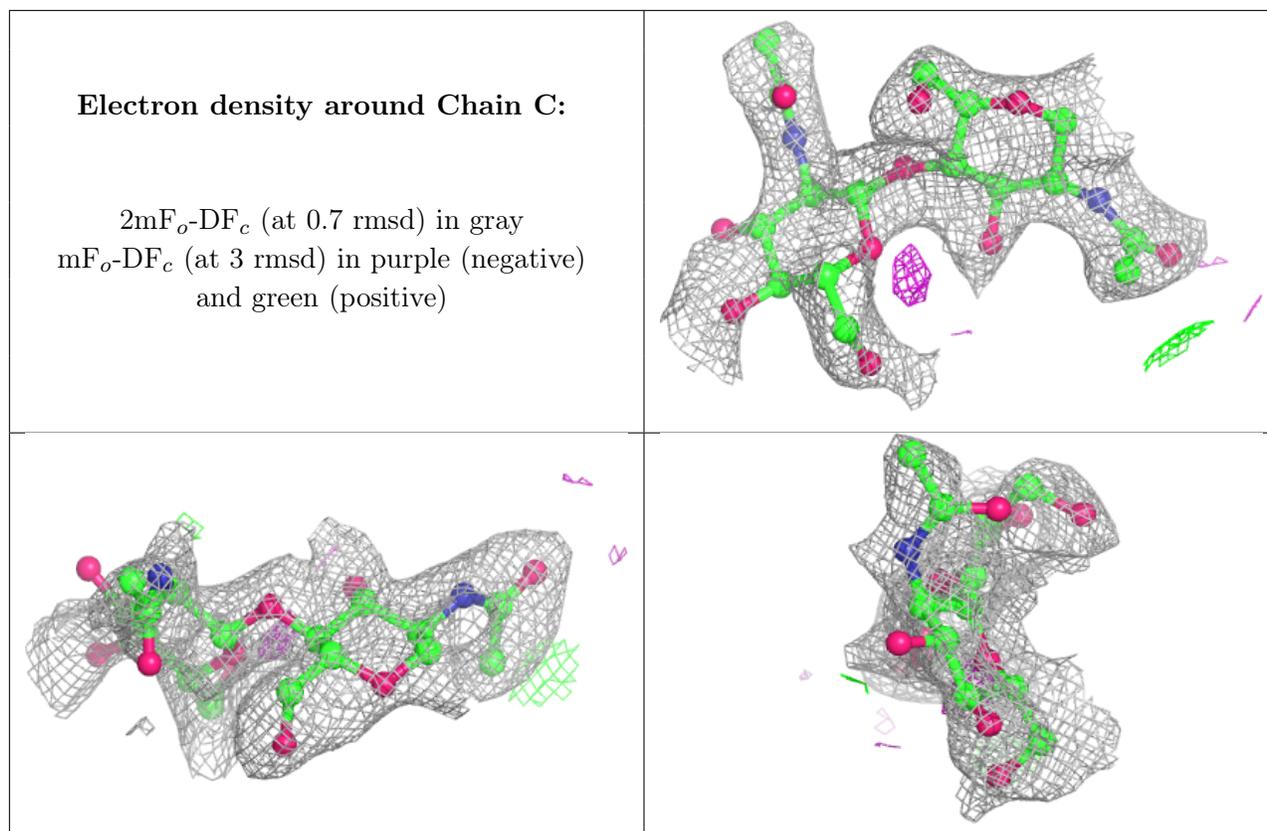
Electron density around Chain C:

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

6.4 Ligands ⓘ

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	NAG	A	503	14/15	0.63	0.21	40,67,72,72	0
4	NAG	A	505	14/15	0.67	0.23	47,71,82,87	0
4	NAG	B	503	14/15	0.67	0.23	62,77,88,89	0
4	NAG	B	505	14/15	0.72	0.20	56,70,78,82	0
4	NAG	A	504	14/15	0.77	0.21	54,69,87,97	0
4	NAG	B	504	14/15	0.92	0.10	19,29,46,47	0
3	CA	B	502	1/1	0.94	0.05	28,28,28,28	0
3	CA	A	502	1/1	0.96	0.04	22,22,22,22	0
3	CA	A	501	1/1	0.99	0.04	14,14,14,14	0
3	CA	B	501	1/1	1.00	0.02	14,14,14,14	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 A 503:

2mF$_o$-DF$_c$ (at 0.7 rmsd) in gray
mF$_o$-DF$_c$ (at 3 rmsd) in purple (negative)
and green (positive)
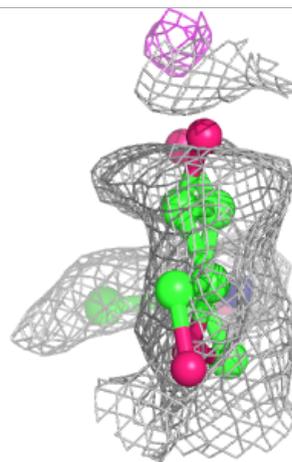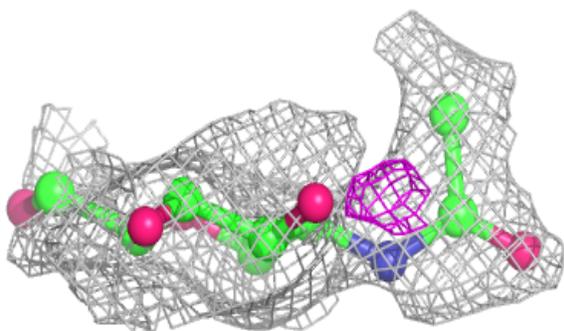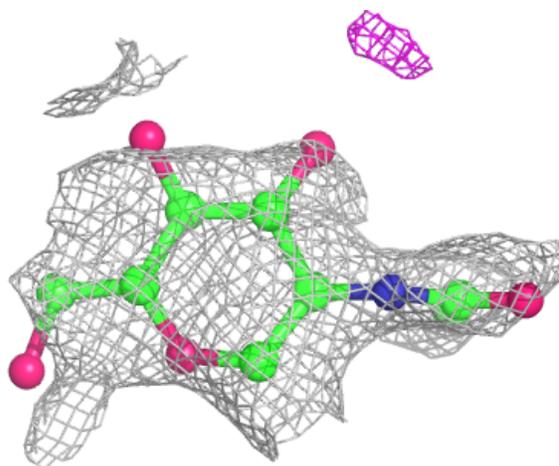
Electron density around NAG A 505:

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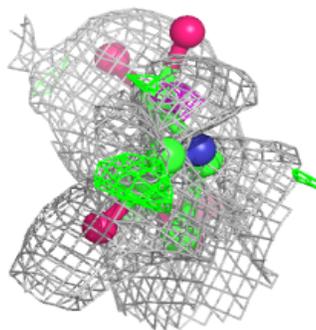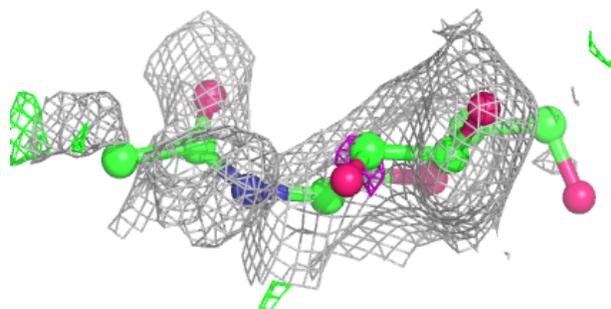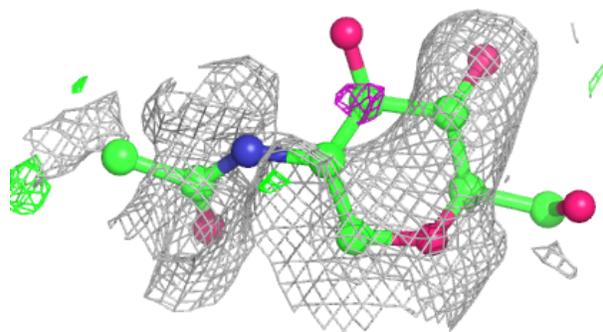


Electron density around NAG B 503:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)
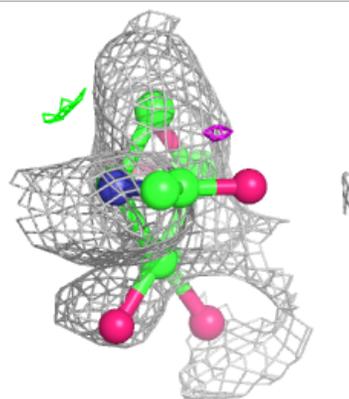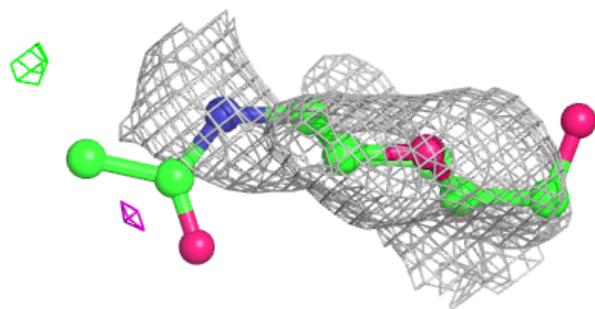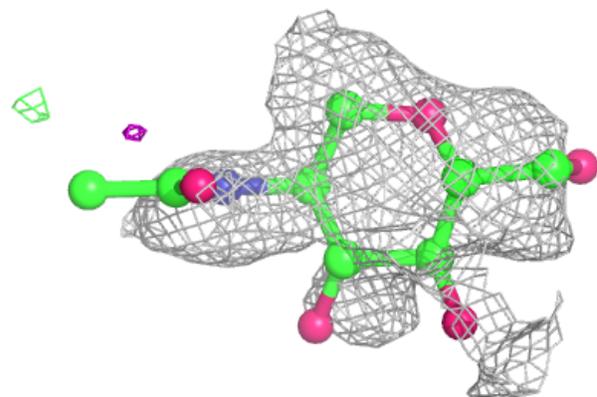
Electron density around NAG B 505:

$2mF_o\text{-}DF_c$ (at 0.7 rmsd) in gray
$mF_o\text{-}DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)
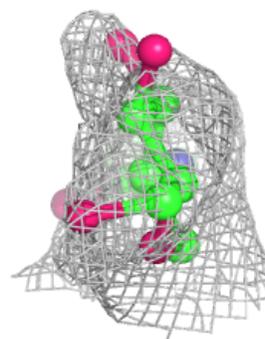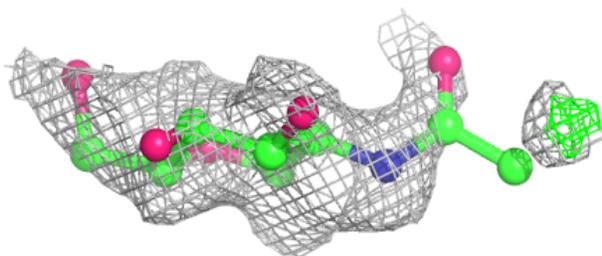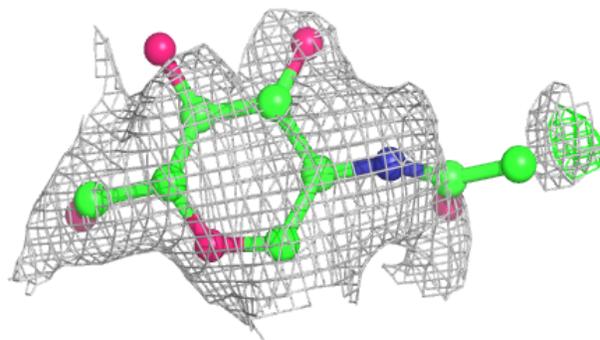
Electron density around NAG A 504:

$2mF_o\text{-}DF_c$ (at 0.7 rmsd) in gray
$mF_o\text{-}DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)
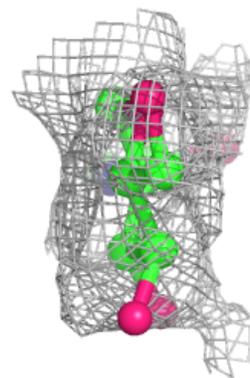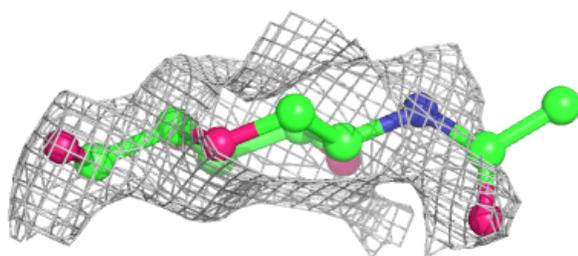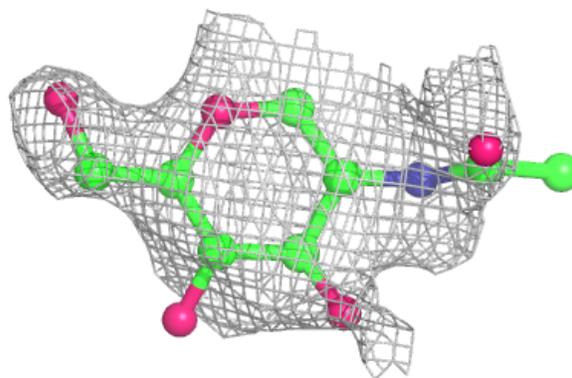
Electron density around NAG B 504:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)
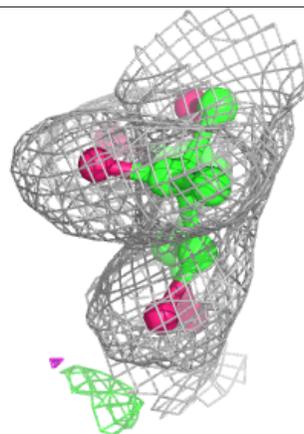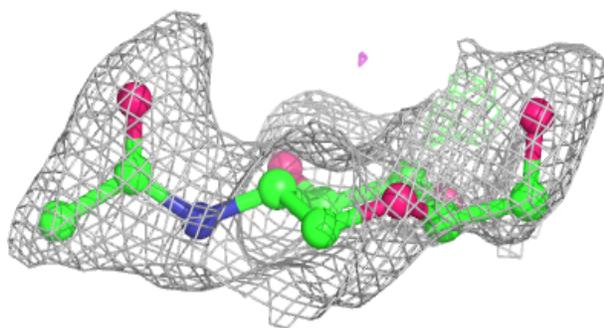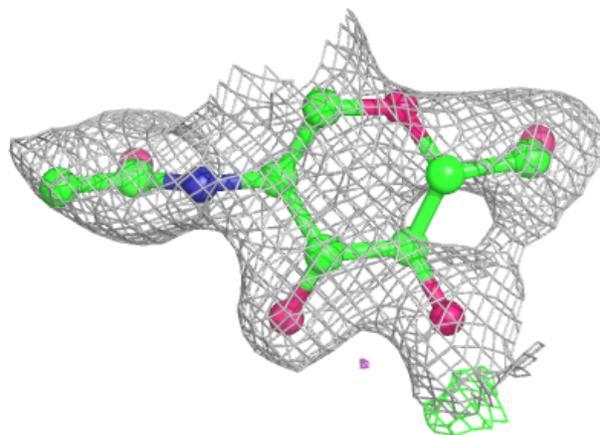
Electron density around CA B 502:

$2mF_o\text{-}DF_c$ (at 0.7 rmsd) in gray
$mF_o\text{-}DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)
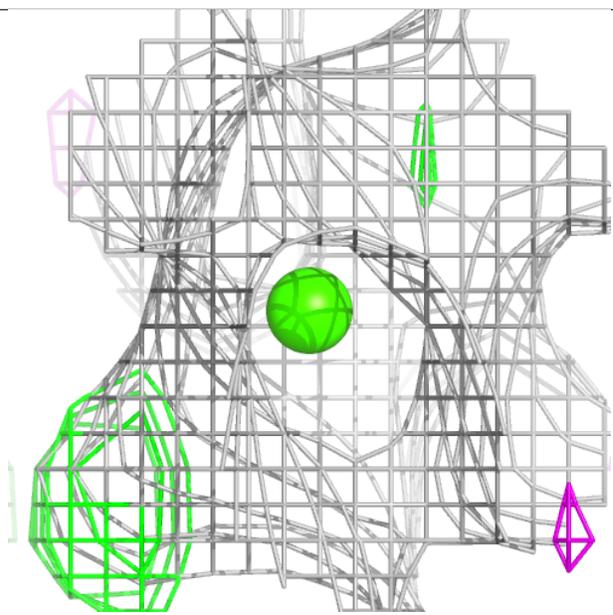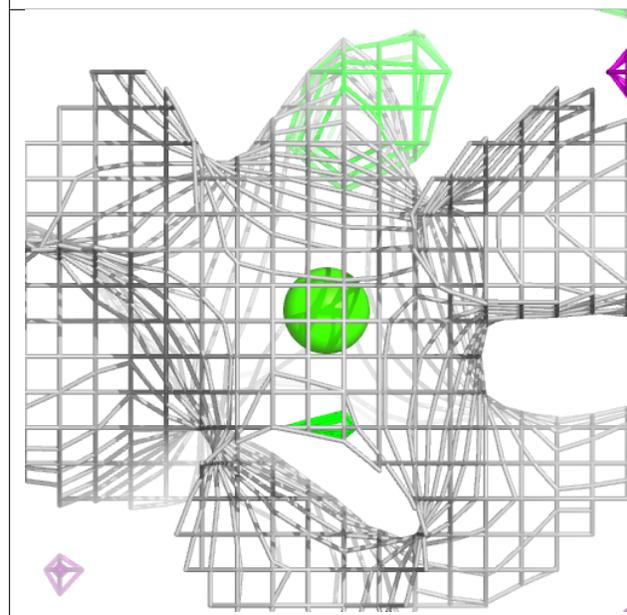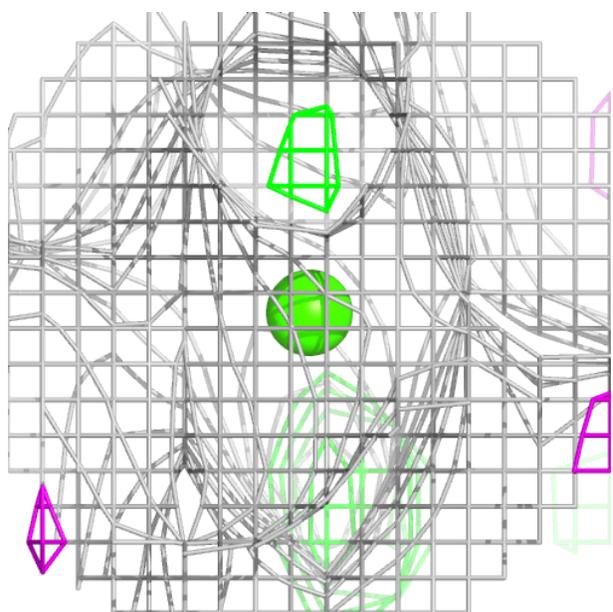
Electron density around CA A 502:

$2mF_o\text{-}DF_c$ (at 0.7 rmsd) in gray
$mF_o\text{-}DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)
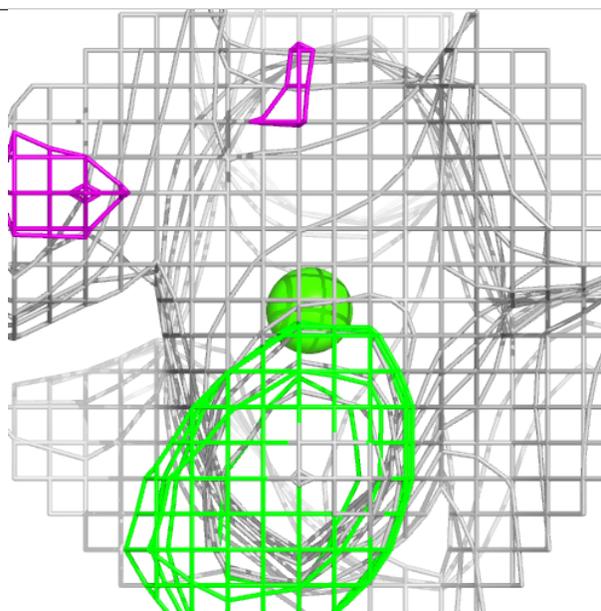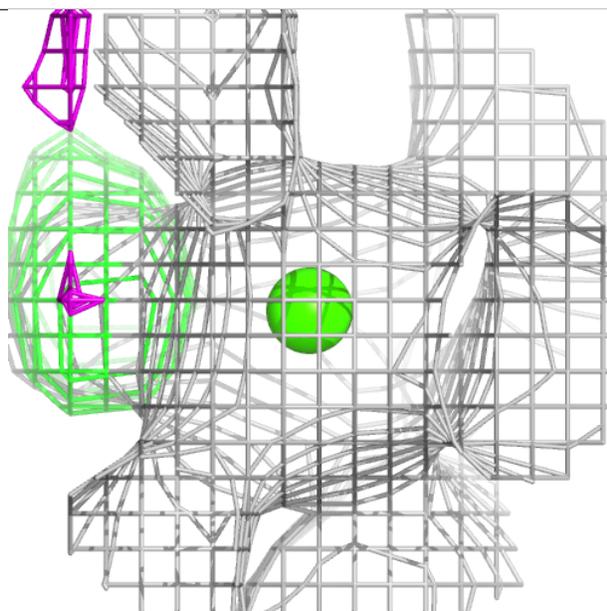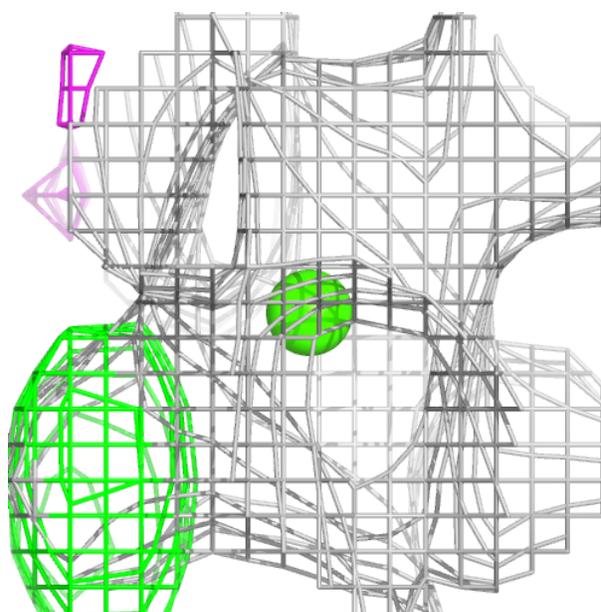
Electron density around CA A 501:

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

Electron density around CA 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 ⓘ

There are no such residues in this entry.