



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 09:39 AM UTC

PDB ID : 7P9Y / pdb_00007p9y
Title : N-acetylglucosamine kinase from Plesiomonas shigelloides complexed with alp
ha-N-acetylglucosamine
Authors : Roy, S.; Isupov, M.N.; Harmer, N.J.; Ames, J.R.
Deposited on : 2021-07-28
Resolution : 1.94 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

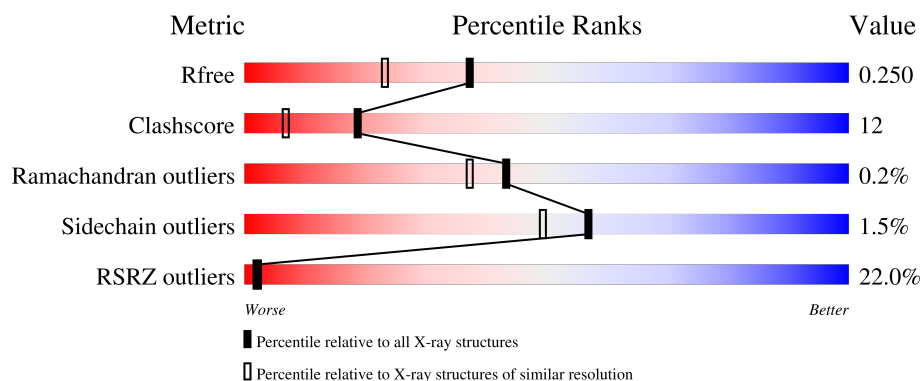
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.94 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	1452 (1.94-1.94)
Clashscore	190562	1494 (1.94-1.94)
Ramachandran outliers	187476	1479 (1.94-1.94)
Sidechain outliers	187428	1479 (1.94-1.94)
RSRZ outliers	180081	1453 (1.94-1.94)

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	AAA	417	18% 58% 15% 27%
1	BBB	417	15% 60% 13% 27%

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
2	MPD	AAA	601	-	-	X	-
3	EDO	AAA	602	-	-	X	-
3	EDO	BBB	504	-	-	X	-
9	PEG	BBB	501	-	-	X	-

2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 5095 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 Ubiquitin-like protein SMT3,N-acetyl-D-glucosamine kinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	AAA	304	Total	C	N	O	S	0	3	0
			2335	1491	405	429	10			
1	BBB	304	Total	C	N	O	S	0	6	0
			2344	1498	405	431	10			

There are 40 discrepancies between the modelled and reference sequences:

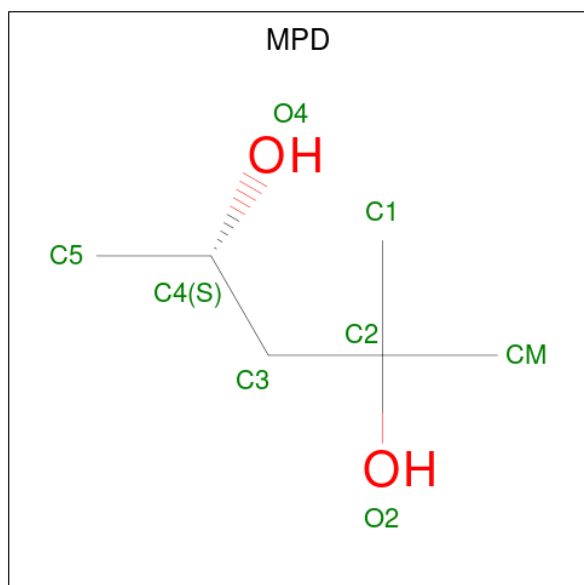
Chain	Residue	Modelled	Actual	Comment	Reference
AAA	-114	MET	-	initiating methionine	UNP Q12306
AAA	-113	ALA	-	expression tag	UNP Q12306
AAA	-112	HIS	-	expression tag	UNP Q12306
AAA	-111	HIS	-	expression tag	UNP Q12306
AAA	-110	HIS	-	expression tag	UNP Q12306
AAA	-109	HIS	-	expression tag	UNP Q12306
AAA	-108	HIS	-	expression tag	UNP Q12306
AAA	-107	HIS	-	expression tag	UNP Q12306
AAA	-106	GLY	-	expression tag	UNP Q12306
AAA	-10	SER	-	linker	UNP Q12306
AAA	-9	SER	-	linker	UNP Q12306
AAA	-8	GLY	-	linker	UNP Q12306
AAA	-7	LEU	-	linker	UNP Q12306
AAA	-6	GLU	-	linker	UNP Q12306
AAA	-5	VAL	-	linker	UNP Q12306
AAA	-4	LEU	-	linker	UNP Q12306
AAA	-3	PHE	-	linker	UNP Q12306
AAA	-2	GLN	-	linker	UNP Q12306
AAA	-1	GLY	-	linker	UNP Q12306
AAA	0	THR	-	linker	UNP Q12306
BBB	-114	MET	-	initiating methionine	UNP Q12306
BBB	-113	ALA	-	expression tag	UNP Q12306
BBB	-112	HIS	-	expression tag	UNP Q12306
BBB	-111	HIS	-	expression tag	UNP Q12306
BBB	-110	HIS	-	expression tag	UNP Q12306

Continued on next page...

Continued from previous page...

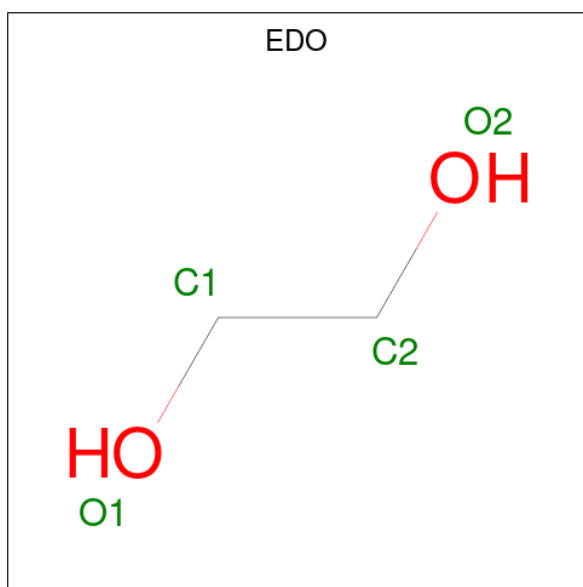
Chain	Residue	Modelled	Actual	Comment	Reference
BBB	-109	HIS	-	expression tag	UNP Q12306
BBB	-108	HIS	-	expression tag	UNP Q12306
BBB	-107	HIS	-	expression tag	UNP Q12306
BBB	-106	GLY	-	expression tag	UNP Q12306
BBB	-10	SER	-	linker	UNP Q12306
BBB	-9	SER	-	linker	UNP Q12306
BBB	-8	GLY	-	linker	UNP Q12306
BBB	-7	LEU	-	linker	UNP Q12306
BBB	-6	GLU	-	linker	UNP Q12306
BBB	-5	VAL	-	linker	UNP Q12306
BBB	-4	LEU	-	linker	UNP Q12306
BBB	-3	PHE	-	linker	UNP Q12306
BBB	-2	GLN	-	linker	UNP Q12306
BBB	-1	GLY	-	linker	UNP Q12306
BBB	0	THR	-	linker	UNP Q12306

- Molecule 2 is (4S)-2-METHYL-2,4-PENTANEDIOL (CCD ID: MPD) (formula: $C_6H_{14}O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	AAA	1	Total	C	O	0	0
			8	6	2		
2	AAA	1	Total	C	O	0	0
			8	6	2		
2	BBB	1	Total	C	O	0	0
			8	6	2		

- Molecule 3 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



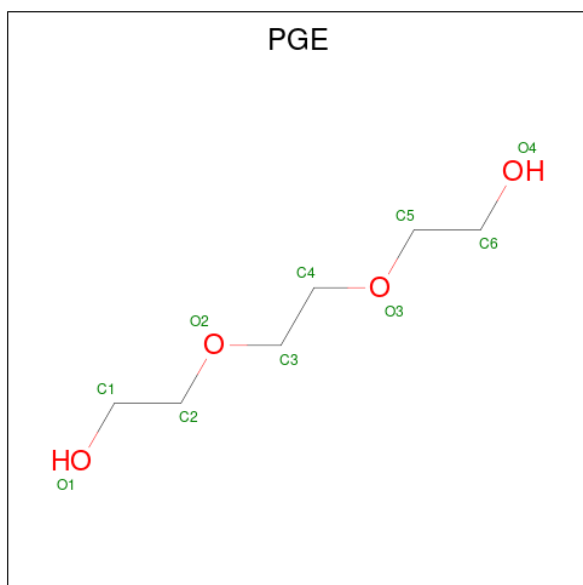
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	AAA	1	Total	C	O	0	0
			4	2	2		
3	AAA	1	Total	C	O	0	0
			4	2	2		
3	AAA	1	Total	C	O	0	0
			4	2	2		
3	AAA	1	Total	C	O	0	0
			4	2	2		
3	AAA	1	Total	C	O	0	0
			4	2	2		
3	AAA	1	Total	C	O	0	0
			4	2	2		
3	AAA	1	Total	C	O	0	0
			4	2	2		
3	BBB	1	Total	C	O	0	0
			4	2	2		
3	BBB	1	Total	C	O	0	0
			4	2	2		
3	BBB	1	Total	C	O	0	0
			4	2	2		
3	BBB	1	Total	C	O	0	0
			4	2	2		

Continued on next page...

Continued from previous page...

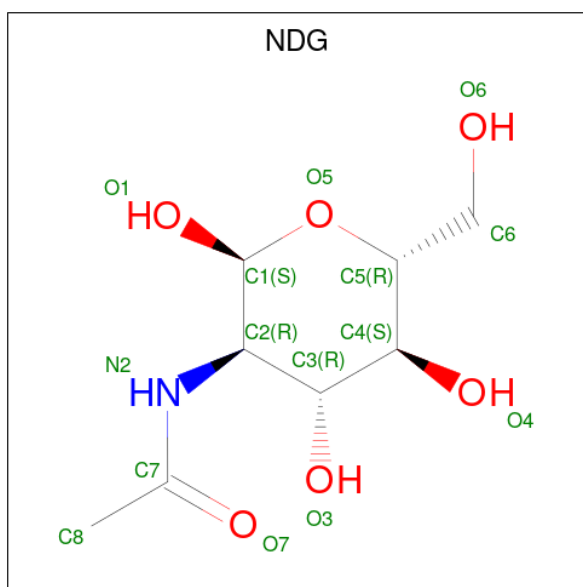
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	BBB	1	Total	C	O	0	0
			4	2	2		
3	BBB	1	Total	C	O	0	0
			4	2	2		

- Molecule 4 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: $C_6H_{14}O_4$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	AAA	1	Total	C	O	0	0
			10	6	4		

- Molecule 5 is 2-acetamido-2-deoxy-alpha-D-glucopyranose (CCD ID: NDG) (formula: $C_8H_{15}NO_6$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	AAA	1	Total	C	N	O	0	0
			15	8	1	6		
5	BBB	1	Total	C	N	O	0	0
			15	8	1	6		

- Molecule 6 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	AAA	2	Total	Zn	0	0
			2	2		
6	BBB	2	Total	Zn	0	0
			2	2		

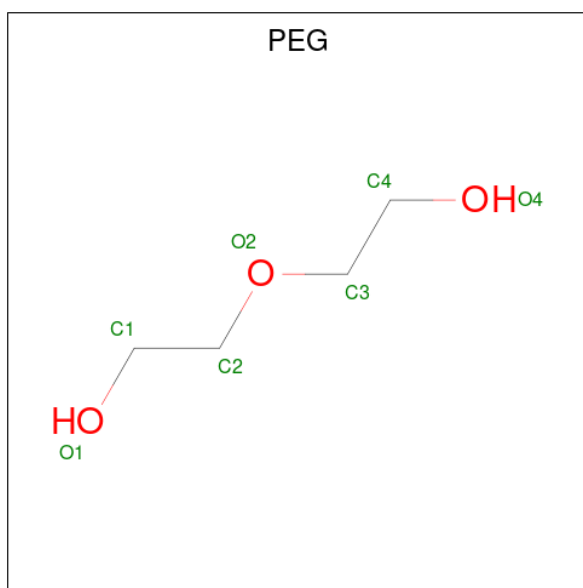
- Molecule 7 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	AAA	1	Total	K	0	0
			1	1		
7	BBB	1	Total	K	0	0
			1	1		

- Molecule 8 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	AAA	2	Total	Cl	0	0
			2	2		

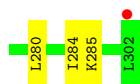
- Molecule 9 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	BBB	1	Total	C	O	0	0
			7	4	3		

- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	AAA	141	Total	O	0	0
			141	141		
10	BBB	132	Total	O	0	0
			132	132		



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	115.48Å 115.48Å 119.66Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	46.14 – 1.94 46.14 – 1.94	Depositor EDS
% Data completeness (in resolution range)	100.0 (46.14-1.94) 99.9 (46.14-1.94)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.27 (at 1.94Å)	Xtriage
Refinement program	REFMAC 5.8.0267, BUSTER	Depositor
R, R_{free}	0.209 , 0.250 0.209 , 0.250	Depositor DCC
R_{free} test set	3517 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	40.0	Xtriage
Anisotropy	0.118	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 48.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.024 for -h,-k,l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5095	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.90% 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: PGE, ZN, EDO, MPD, NDG, CL, K, PEG

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	AAA	0.63	0/2393	1.13	4/3240 (0.1%)
1	BBB	0.64	0/2406	1.11	3/3257 (0.1%)
All	All	0.64	0/4799	1.12	7/6497 (0.1%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AAA	157	HIS	CA-CB-CG	-6.65	107.15	113.80
1	AAA	246	ASP	O-C-N	-5.89	118.82	121.53
1	BBB	246	ASP	O-C-N	-5.38	119.06	121.53
1	AAA	10	THR	CB-CA-C	5.34	118.37	109.56
1	BBB	10	THR	CB-CA-C	5.29	118.29	109.56
1	AAA	163	ASP	CA-CB-CG	5.04	117.64	112.60
1	BBB	157	HIS	CA-CB-CG	-5.01	108.79	113.80

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	AAA	2335	0	2334	62	0
1	BBB	2344	0	2359	51	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	AAA	16	0	28	8	0
2	BBB	8	0	14	2	0
3	AAA	36	0	54	8	0
3	BBB	28	0	42	13	0
4	AAA	10	0	14	2	0
5	AAA	15	0	12	0	0
5	BBB	15	0	12	0	0
6	AAA	2	0	0	0	0
6	BBB	2	0	0	0	0
7	AAA	1	0	0	0	0
7	BBB	1	0	0	0	0
8	AAA	2	0	0	0	0
9	BBB	7	0	10	10	0
10	AAA	141	0	0	2	0
10	BBB	132	0	0	1	0
All	All	5095	0	4879	115	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 12.

All (115) 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:BBB:276:HIS:H	3:BBB:504:EDO:H11	1.08	1.16
1:BBB:141:ASN:HD22	9:BBB:501:PEG:H41	1.20	1.04
1:BBB:140:PHE:HB2	9:BBB:501:PEG:H12	1.41	1.00
1:AAA:141:ASN:OD1	9:BBB:501:PEG:H22	1.66	0.95
1:BBB:280:LEU:H	3:BBB:505:EDO:H11	1.31	0.94
1:AAA:255:GLY:HA3	2:AAA:601:MPD:H52	1.48	0.93
1:AAA:57:LYS:HB2	1:AAA:98:ARG:HH22	1.36	0.91
1:AAA:39:PHE:HE2	1:AAA:80:ILE:HG21	1.34	0.90
1:AAA:57:LYS:HB2	1:AAA:98:ARG:NH2	1.88	0.89
1:AAA:68:ALA:HB3	3:AAA:602:EDO:H21	1.63	0.80
1:BBB:143:LYS:HB3	9:BBB:501:PEG:H21	1.64	0.79
1:BBB:276:HIS:N	3:BBB:504:EDO:H11	1.93	0.79
1:BBB:111[B]:LEU:HD11	1:BBB:144[B]:VAL:HG21	1.63	0.79
1:AAA:57:LYS:CB	1:AAA:98:ARG:HH22	2.03	0.72
1:AAA:260:GLU:H	4:AAA:605:PGE:H42	1.57	0.70
1:BBB:111[B]:LEU:CD1	1:BBB:144[B]:VAL:HG21	2.22	0.70
1:BBB:68:ALA:HB3	3:BBB:502:EDO:H22	1.73	0.69
1:BBB:285:LYS:H	3:BBB:509:EDO:H12	1.58	0.68

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BBB:141:ASN:HD22	9:BBB:501:PEG:C4	2.02	0.67
1:AAA:30:ALA:O	1:AAA:32:PRO:HD3	1.94	0.67
1:BBB:30:ALA:O	1:BBB:32:PRO:HD3	1.95	0.66
1:BBB:103:GLU:HA	3:BBB:502:EDO:H11	1.78	0.66
1:AAA:33:THR:HA	1:AAA:82:ALA:HB2	1.79	0.63
1:AAA:40:LEU:HD23	10:AAA:816:HOH:O	1.97	0.63
1:BBB:111[B]:LEU:HG	1:BBB:144[B]:VAL:CG2	2.30	0.62
1:BBB:255:GLY:HA3	2:BBB:503:MPD:H4	1.80	0.62
1:AAA:39:PHE:CE2	1:AAA:80:ILE:HG21	2.26	0.62
2:AAA:601:MPD:H53	2:AAA:601:MPD:H11	1.81	0.60
1:AAA:43:ILE:O	1:AAA:47:VAL:HG23	2.01	0.60
1:AAA:255:GLY:CA	2:AAA:601:MPD:H52	2.27	0.60
1:BBB:111[B]:LEU:HD11	1:BBB:144[B]:VAL:CG2	2.30	0.59
1:BBB:36:TYR:HB2	1:BBB:82:ALA:O	2.02	0.59
1:BBB:280:LEU:N	3:BBB:505:EDO:H11	2.10	0.59
1:AAA:39:PHE:CE2	1:AAA:83:ALA:HB2	2.38	0.58
1:BBB:141:ASN:ND2	9:BBB:501:PEG:H41	2.05	0.58
1:AAA:51:ASP:HA	1:AAA:56:VAL:O	2.05	0.57
1:AAA:57:LYS:C	1:AAA:98:ARG:HH22	2.13	0.57
1:BBB:98:ARG:HB2	1:BBB:99:PRO:HD2	1.86	0.57
1:BBB:43:ILE:O	1:BBB:47:VAL:HG23	2.04	0.57
1:AAA:36:TYR:HB2	1:AAA:82:ALA:O	2.04	0.56
1:AAA:89:GLN:O	1:AAA:93:GLU:HG3	2.06	0.56
1:BBB:89:GLN:O	1:BBB:93:GLU:HG3	2.05	0.55
1:BBB:140:PHE:HB2	9:BBB:501:PEG:C1	2.25	0.55
1:AAA:255:GLY:HA3	2:AAA:601:MPD:C5	2.31	0.54
1:AAA:57:LYS:CB	1:AAA:98:ARG:NH2	2.64	0.54
1:BBB:33:THR:HA	1:BBB:82:ALA:HB2	1.88	0.54
1:BBB:128[A]:LEU:CD1	1:BBB:155:ILE:HD11	2.37	0.54
1:AAA:115[A]:TRP:HA	1:AAA:115[A]:TRP:CE3	2.43	0.54
1:AAA:86:HIS:HB3	10:AAA:719:HOH:O	2.08	0.53
1:AAA:128:LEU:CD1	1:AAA:155:ILE:HD11	2.38	0.53
1:BBB:128[A]:LEU:HD23	1:BBB:252:LEU:HD22	1.90	0.53
1:BBB:111[B]:LEU:CG	1:BBB:144[B]:VAL:CG2	2.87	0.53
1:AAA:140:PHE:HB2	9:BBB:501:PEG:H32	1.91	0.52
1:AAA:68:ALA:CB	3:AAA:602:EDO:H21	2.38	0.52
1:BBB:111[B]:LEU:CD1	1:BBB:144[B]:VAL:CG2	2.87	0.52
1:AAA:57:LYS:C	1:AAA:98:ARG:NH2	2.68	0.52
1:AAA:75:LEU:HG	1:AAA:77:THR:HG22	1.92	0.52
1:AAA:98:ARG:HB2	1:AAA:99:PRO:HD2	1.91	0.51
1:AAA:39:PHE:HE2	1:AAA:80:ILE:CG2	2.16	0.50

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:AAA:601:MPD:H53	2:AAA:601:MPD:C1	2.40	0.50
1:BBB:280:LEU:H	3:BBB:505:EDO:C1	2.13	0.50
1:AAA:180:CYS:O	1:AAA:181:LYS:HB2	2.12	0.50
1:AAA:193:ARG:HH11	1:AAA:197:GLN:HE22	1.60	0.50
1:AAA:230:PHE:HD2	1:AAA:231:MET:HE2	1.77	0.49
1:AAA:57:LYS:CA	1:AAA:98:ARG:HH22	2.24	0.49
1:AAA:140:PHE:CB	9:BBB:501:PEG:H32	2.42	0.49
1:AAA:248:HIS:HA	1:BBB:148:ARG:HB2	1.95	0.49
1:AAA:2:TYR:CE2	1:AAA:302:LEU:HD12	2.47	0.49
1:AAA:68:ALA:H	3:AAA:602:EDO:C2	2.26	0.49
1:AAA:128:LEU:HD23	1:AAA:252:LEU:HD22	1.94	0.49
1:BBB:150:ASN:ND2	3:BBB:502:EDO:H21	2.28	0.48
1:BBB:284:ILE:HG22	3:BBB:509:EDO:H22	1.95	0.48
1:AAA:57:LYS:HB2	1:AAA:98:ARG:CZ	2.43	0.48
1:BBB:111[B]:LEU:HG	1:BBB:144[B]:VAL:HG22	1.94	0.48
1:AAA:68:ALA:H	3:AAA:602:EDO:H21	1.79	0.48
1:BBB:94:GLU:C	1:BBB:96:LEU:H	2.21	0.48
1:AAA:77:THR:HG23	1:AAA:77:THR:O	2.14	0.48
1:AAA:132:THR:H	3:AAA:604:EDO:H12	1.80	0.47
1:AAA:211:GLU:OE2	3:AAA:609:EDO:H11	2.15	0.47
1:BBB:276:HIS:HB3	3:BBB:504:EDO:H22	1.97	0.46
1:AAA:131:GLY:HA2	2:AAA:601:MPD:H4	1.97	0.46
1:BBB:51:ASP:HA	1:BBB:56:VAL:O	2.16	0.46
1:BBB:180:CYS:O	1:BBB:181:LYS:HB2	2.16	0.45
1:AAA:47:VAL:O	1:AAA:51:ASP:CG	2.59	0.45
1:AAA:128:LEU:HD23	1:AAA:252:LEU:CD2	2.47	0.45
1:BBB:193:ARG:HH11	1:BBB:197:GLN:HE22	1.65	0.45
1:BBB:47:VAL:O	1:BBB:51:ASP:CG	2.60	0.44
1:AAA:11:LYS:HG2	1:AAA:30:ALA:HA	2.00	0.44
1:AAA:7:ILE:O	1:AAA:65:PRO:HD3	2.18	0.43
1:AAA:228:GLU:CD	1:AAA:269:ARG:HH22	2.25	0.43
1:BBB:20:ASP:O	1:BBB:21:LEU:HB2	2.18	0.43
1:AAA:180:CYS:HB2	3:BBB:504:EDO:H12	2.00	0.43
1:BBB:111[B]:LEU:HD21	1:BBB:144[B]:VAL:HG23	2.00	0.43
1:BBB:123:PRO:HB2	1:BBB:248:HIS:CE1	2.54	0.43
1:BBB:220:GLU:HG3	3:BBB:507:EDO:H22	2.00	0.43
1:AAA:192:GLY:HA3	2:AAA:601:MPD:H31	2.01	0.43
1:AAA:20:ASP:O	1:AAA:21:LEU:HB2	2.19	0.42
1:BBB:7:ILE:O	1:BBB:65:PRO:HD3	2.19	0.42
2:BBB:503:MPD:H53	10:BBB:701:HOH:O	2.19	0.42
1:AAA:123:PRO:HB2	1:AAA:248:HIS:CE1	2.54	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AAA:200:ASP:O	3:AAA:608:EDO:H11	2.20	0.42
4:AAA:605:PGE:H1	4:AAA:605:PGE:H52	2.01	0.42
1:AAA:21:LEU:HD12	1:AAA:115[B]:TRP:CZ3	2.55	0.42
1:AAA:207:LEU:CD2	3:AAA:609:EDO:H22	2.50	0.42
1:BBB:77:THR:HG23	1:BBB:77:THR:O	2.20	0.42
1:BBB:44:VAL:HG11	1:BBB:95:ARG:NH2	2.35	0.42
1:BBB:90:ARG:HD3	1:BBB:90:ARG:HA	1.80	0.41
1:AAA:140:PHE:HB3	9:BBB:501:PEG:H11	2.01	0.41
1:AAA:151:ILE:O	1:AAA:151:ILE:CG2	2.68	0.41
1:AAA:209:ALA:HB3	2:AAA:601:MPD:CM	2.50	0.41
1:BBB:128[A]:LEU:HD23	1:BBB:252:LEU:CD2	2.49	0.41
1:AAA:75:LEU:HD23	1:AAA:83:ALA:HB1	2.03	0.41
1:AAA:148:ARG:HB2	1:BBB:248:HIS:HA	2.03	0.41
1:BBB:126:LEU:HG	1:BBB:128[B]:LEU:HG	2.03	0.40
1:BBB:97:GLN:HG3	1:BBB:97:GLN:O	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	AAA	305/417 (73%)	291 (95%)	14 (5%)	0	100	100
1	BBB	308/417 (74%)	292 (95%)	15 (5%)	1 (0%)	36	30
All	All	613/834 (74%)	583 (95%)	29 (5%)	1 (0%)	43	37

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	BBB	96	LEU

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	AAA	241/340 (71%)	236 (98%)	5 (2%)	47	37
1	BBB	244/340 (72%)	242 (99%)	2 (1%)	73	71
All	All	485/680 (71%)	478 (99%)	7 (1%)	57	52

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	AAA	64	ILE
1	AAA	76	LEU
1	AAA	116	ASP
1	AAA	163	ASP
1	AAA	166	LYS
1	BBB	76	LEU
1	BBB	116	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 31 ligands modelled in this entry, 8 are monoatomic - leaving 23 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	$\# Z > 2$
3	EDO	AAA	607	-	3,3,3	0.17	0	2,2,2	0.36	0
2	MPD	AAA	611	-	7,7,7	0.12	0	9,10,10	0.48	0
2	MPD	AAA	601	-	7,7,7	0.44	0	9,10,10	1.16	0
3	EDO	BBB	508	-	3,3,3	0.17	0	2,2,2	0.18	0
3	EDO	AAA	610	-	3,3,3	0.30	0	2,2,2	0.57	0
9	PEG	BBB	501	-	6,6,6	0.57	0	5,5,5	0.27	0
3	EDO	AAA	606	-	3,3,3	0.06	0	2,2,2	0.21	0
3	EDO	BBB	504	-	3,3,3	0.28	0	2,2,2	0.19	0
2	MPD	BBB	503	-	7,7,7	0.45	0	9,10,10	1.24	1 (11%)
3	EDO	BBB	509	-	3,3,3	0.16	0	2,2,2	0.38	0
3	EDO	BBB	506	-	3,3,3	0.12	0	2,2,2	0.26	0
3	EDO	BBB	507	-	3,3,3	0.15	0	2,2,2	0.29	0
3	EDO	AAA	608	-	3,3,3	0.13	0	2,2,2	0.51	0
3	EDO	AAA	604	-	3,3,3	0.31	0	2,2,2	0.69	0
3	EDO	BBB	505	-	3,3,3	0.19	0	2,2,2	0.40	0
3	EDO	AAA	609	-	3,3,3	0.24	0	2,2,2	0.68	0
3	EDO	AAA	603	-	3,3,3	0.09	0	2,2,2	0.61	0
5	NDG	BBB	510	-	15,15,15	0.38	0	21,21,21	1.82	2 (9%)
3	EDO	BBB	502	-	3,3,3	0.44	0	2,2,2	0.26	0
4	PGE	AAA	605	-	9,9,9	0.35	0	8,8,8	0.27	0
3	EDO	AAA	602	-	3,3,3	0.08	0	2,2,2	0.45	0
3	EDO	AAA	616	-	3,3,3	0.39	0	2,2,2	0.10	0
5	NDG	AAA	612	-	15,15,15	0.37	0	21,21,21	2.37	2 (9%)

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	EDO	AAA	607	-	-	1/1/1/1	-

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	MPD	AAA	611	-	-	2/5/5/5	-
2	MPD	AAA	601	-	-	2/5/5/5	-
3	EDO	BBB	508	-	-	1/1/1/1	-
3	EDO	AAA	610	-	-	0/1/1/1	-
9	PEG	BBB	501	-	-	3/4/4/4	-
3	EDO	AAA	606	-	-	0/1/1/1	-
3	EDO	BBB	504	-	-	1/1/1/1	-
2	MPD	BBB	503	-	-	5/5/5/5	-
3	EDO	BBB	509	-	-	0/1/1/1	-
3	EDO	BBB	506	-	-	1/1/1/1	-
3	EDO	BBB	507	-	-	1/1/1/1	-
3	EDO	AAA	608	-	-	0/1/1/1	-
3	EDO	AAA	604	-	-	1/1/1/1	-
3	EDO	BBB	505	-	-	0/1/1/1	-
3	EDO	AAA	609	-	-	1/1/1/1	-
3	EDO	AAA	603	-	-	1/1/1/1	-
5	NDG	BBB	510	-	-	2/6/26/26	0/1/1/1
3	EDO	BBB	502	-	-	0/1/1/1	-
4	PGE	AAA	605	-	-	2/7/7/7	-
3	EDO	AAA	602	-	-	0/1/1/1	-
3	EDO	AAA	616	-	-	1/1/1/1	-
5	NDG	AAA	612	-	-	1/6/26/26	0/1/1/1

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	AAA	612	NDG	C1-C2-N2	7.87	119.84	110.73
5	AAA	612	NDG	O5-C1-C2	6.98	116.53	109.52
5	BBB	510	NDG	C1-C2-N2	6.27	117.99	110.73
5	BBB	510	NDG	O5-C1-C2	4.26	113.80	109.52
2	BBB	503	MPD	C5-C4-C3	3.28	126.90	111.67

There are no chirality outliers.

All (26) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	BBB	503	MPD	C1-C2-C3-C4
2	BBB	503	MPD	C2-C3-C4-O4
2	BBB	503	MPD	C2-C3-C4-C5

Continued on next page...

Continued from previous page...

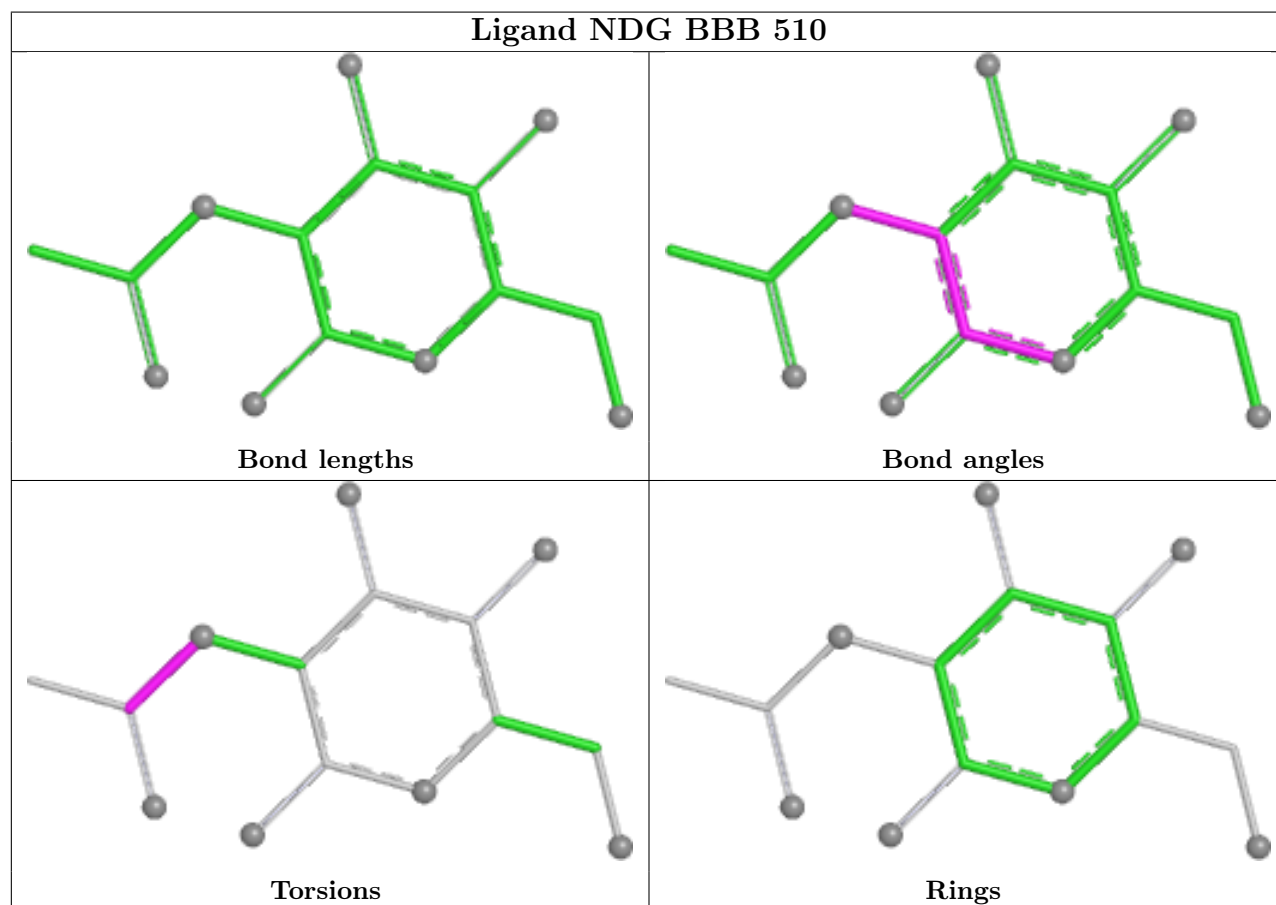
Mol	Chain	Res	Type	Atoms
5	BBB	510	NDG	C8-C7-N2-C2
5	BBB	510	NDG	O7-C7-N2-C2
3	AAA	604	EDO	O1-C1-C2-O2
3	BBB	508	EDO	O1-C1-C2-O2
9	BBB	501	PEG	O2-C3-C4-O4
3	AAA	607	EDO	O1-C1-C2-O2
3	BBB	507	EDO	O1-C1-C2-O2
3	BBB	506	EDO	O1-C1-C2-O2
4	AAA	605	PGE	C4-C3-O2-C2
2	BBB	503	MPD	O2-C2-C3-C4
2	BBB	503	MPD	CM-C2-C3-C4
9	BBB	501	PEG	O1-C1-C2-O2
3	AAA	616	EDO	O1-C1-C2-O2
2	AAA	601	MPD	C2-C3-C4-C5
3	AAA	603	EDO	O1-C1-C2-O2
3	AAA	609	EDO	O1-C1-C2-O2
3	BBB	504	EDO	O1-C1-C2-O2
4	AAA	605	PGE	O2-C3-C4-O3
5	AAA	612	NDG	O7-C7-N2-C2
2	AAA	601	MPD	O2-C2-C3-C4
2	AAA	611	MPD	O2-C2-C3-C4
2	AAA	611	MPD	C1-C2-C3-C4
9	BBB	501	PEG	C1-C2-O2-C3

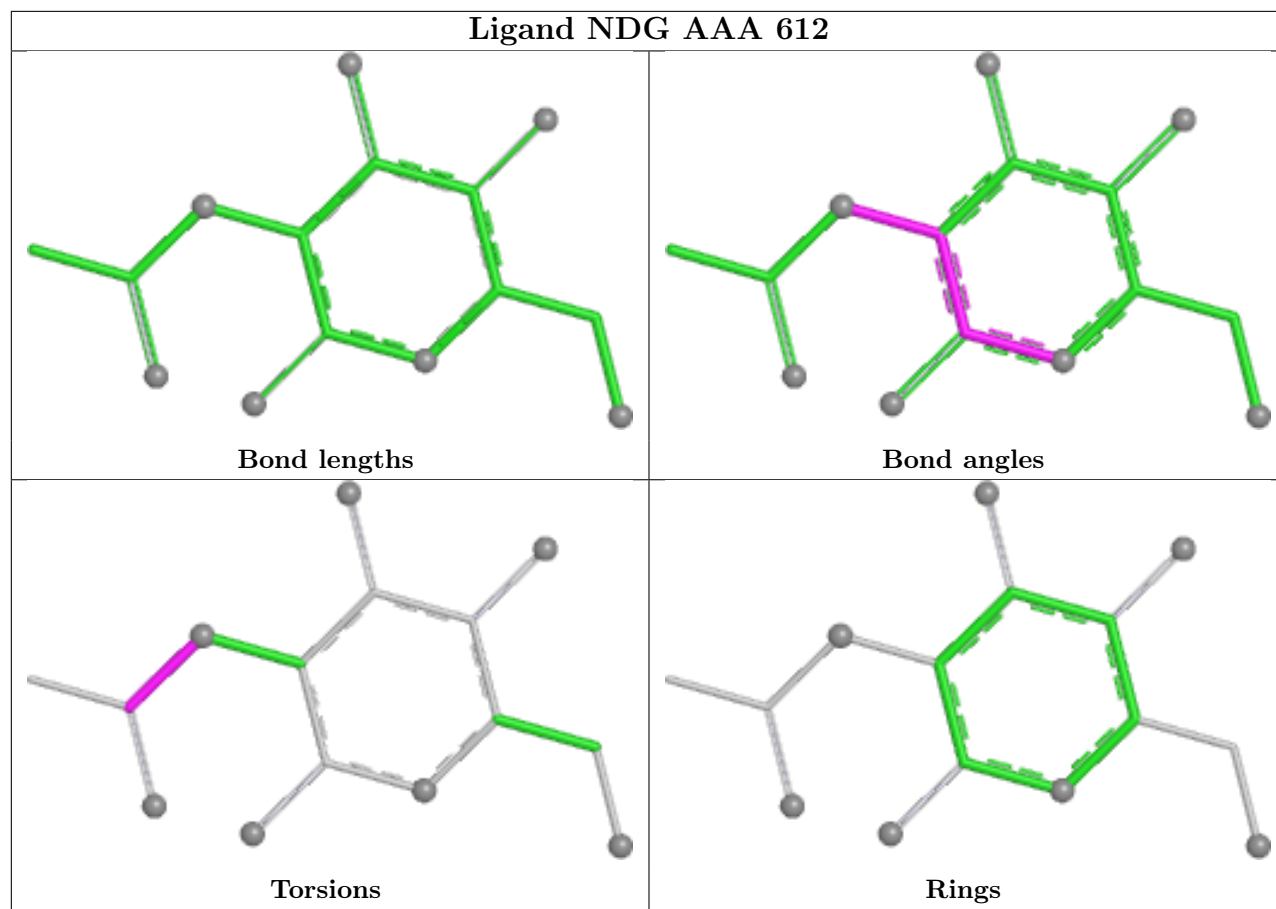
There are no ring outliers.

13 monomers are involved in 43 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	AAA	601	MPD	8	0
9	BBB	501	PEG	10	0
3	BBB	504	EDO	4	0
2	BBB	503	MPD	2	0
3	BBB	509	EDO	2	0
3	BBB	507	EDO	1	0
3	AAA	608	EDO	1	0
3	AAA	604	EDO	1	0
3	BBB	505	EDO	3	0
3	AAA	609	EDO	2	0
3	BBB	502	EDO	3	0
4	AAA	605	PGE	2	0
3	AAA	602	EDO	4	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å²)	Q<0.9
1	AAA	304/417 (72%)	1.14	73 (24%)	2	1	25, 45, 114, 150	3 (0%)
1	BBB	304/417 (72%)	0.98	61 (20%)	3	3	19, 44, 93, 140	6 (1%)
All	All	608/834 (72%)	1.06	134 (22%)	2	2	19, 45, 106, 150	9 (1%)

All (134) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	AAA	43	ILE	8.2
1	BBB	40	LEU	7.6
1	AAA	40	LEU	7.5
1	AAA	58	GLY	7.0
1	AAA	96	LEU	6.8
1	AAA	39	PHE	6.8
1	BBB	44	VAL	6.6
1	AAA	88	LEU	6.6
1	AAA	44	VAL	6.3
1	AAA	92	LEU	6.0
1	AAA	115[A]	TRP	6.0
1	AAA	56	VAL	5.9
1	AAA	75	LEU	5.8
1	AAA	46	LEU	5.7
1	AAA	29	VAL	5.5
1	AAA	12	ILE	5.5
1	AAA	32	PRO	5.4
1	AAA	42	ALA	5.2
1	BBB	36	TYR	5.1
1	BBB	56	VAL	5.1
1	BBB	32	PRO	5.1
1	BBB	54	PHE	5.0
1	AAA	76	LEU	4.9
1	BBB	33	THR	4.7

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	BBB	81	PRO	4.7
1	BBB	96	LEU	4.7
1	BBB	39	PHE	4.6
1	AAA	21	LEU	4.5
1	AAA	30	ALA	4.5
1	AAA	80	ILE	4.4
1	AAA	73	GLY	4.4
1	BBB	12	ILE	4.4
1	AAA	67	ILE	4.3
1	AAA	24	VAL	4.2
1	BBB	38	ALA	4.2
1	BBB	100	VAL	4.1
1	BBB	52	ALA	4.1
1	BBB	29	VAL	4.0
1	AAA	83	ALA	4.0
1	AAA	60	VAL	3.9
1	BBB	92	LEU	3.9
1	BBB	47	VAL	3.8
1	BBB	84	MET	3.8
1	AAA	36	TYR	3.8
1	AAA	3	TYR	3.7
1	AAA	57	LYS	3.7
1	BBB	41	ASP	3.7
1	BBB	75	LEU	3.6
1	BBB	55	GLY	3.6
1	AAA	33	THR	3.6
1	AAA	54	PHE	3.6
1	AAA	66	GLY	3.6
1	AAA	5	PHE	3.6
1	AAA	45	THR	3.5
1	AAA	99	PRO	3.5
1	AAA	64	ILE	3.5
1	AAA	47	VAL	3.5
1	AAA	52	ALA	3.4
1	AAA	65	PRO	3.4
1	BBB	80	ILE	3.4
1	BBB	76	LEU	3.4
1	BBB	98	ARG	3.4
1	BBB	46	LEU	3.3
1	BBB	31	THR	3.3
1	BBB	3	TYR	3.3
1	AAA	98	ARG	3.3

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	BBB	58	GLY	3.3
1	AAA	37	ALA	3.3
1	BBB	88	LEU	3.3
1	AAA	1	MET	3.2
1	AAA	14	PHE	3.2
1	AAA	87	THR	3.2
1	AAA	38	ALA	3.1
1	BBB	10	THR	3.1
1	BBB	0	THR	3.1
1	BBB	9	GLY	3.1
1	AAA	6	ASP	3.1
1	BBB	45	THR	3.1
1	AAA	50	ALA	3.1
1	BBB	83	ALA	3.1
1	AAA	102	ILE	3.1
1	BBB	-1	GLY	3.1
1	AAA	7	ILE	3.0
1	BBB	7	ILE	3.0
1	AAA	100	VAL	3.0
1	AAA	59	THR	3.0
1	BBB	2	TYR	3.0
1	BBB	30	ALA	3.0
1	AAA	55	GLY	2.9
1	AAA	89	GLN	2.9
1	AAA	11	LYS	2.9
1	AAA	74	LYS	2.9
1	BBB	77	THR	2.9
1	AAA	78	SER	2.9
1	BBB	90	ARG	2.8
1	BBB	57	LYS	2.8
1	AAA	302	LEU	2.8
1	BBB	43	ILE	2.8
1	AAA	41	ASP	2.8
1	BBB	118	ASP	2.8
1	AAA	68	ALA	2.7
1	BBB	60	VAL	2.7
1	AAA	94	GLU	2.7
1	BBB	115	TRP	2.7
1	BBB	1	MET	2.7
1	AAA	81	PRO	2.7
1	AAA	31	THR	2.6
1	AAA	-1	GLY	2.6

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	BBB	82	ALA	2.6
1	BBB	50	ALA	2.5
1	AAA	72	THR	2.5
1	BBB	37	ALA	2.5
1	AAA	2	TYR	2.5
1	BBB	42	ALA	2.5
1	AAA	93	GLU	2.4
1	BBB	8	GLY	2.4
1	BBB	86	HIS	2.4
1	BBB	21	LEU	2.4
1	AAA	79	ASN	2.4
1	BBB	73	GLY	2.4
1	AAA	27	GLU	2.3
1	AAA	151	ILE	2.3
1	BBB	302	LEU	2.3
1	AAA	48	ASN	2.3
1	BBB	95	ARG	2.2
1	AAA	62	ILE	2.2
1	AAA	84	MET	2.2
1	BBB	11	LYS	2.2
1	AAA	61	GLY	2.2
1	BBB	65	PRO	2.2
1	BBB	28	ARG	2.1
1	BBB	111[A]	LEU	2.1
1	AAA	4	GLY	2.0
1	BBB	85	GLY	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](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

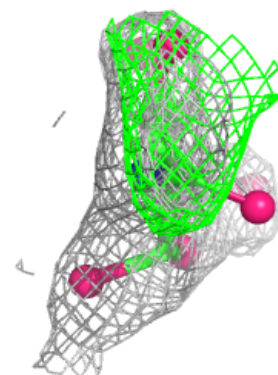
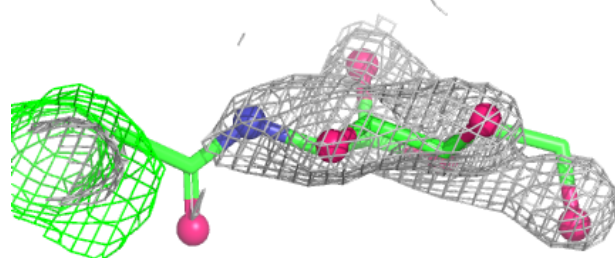
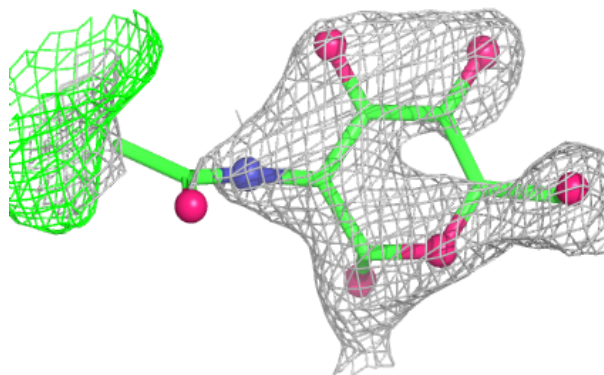
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
3	EDO	AAA	603	4/4	0.69	0.20	56,70,77,87	0
3	EDO	BBB	507	4/4	0.71	0.24	59,62,63,85	4
3	EDO	BBB	508	4/4	0.73	0.19	75,96,97,102	0
3	EDO	AAA	608	4/4	0.75	0.17	64,76,78,86	0
3	EDO	AAA	607	4/4	0.76	0.18	58,70,89,100	0
3	EDO	AAA	610	4/4	0.76	0.20	63,63,65,78	0
9	PEG	BBB	501	7/7	0.78	0.16	32,43,51,57	7
3	EDO	BBB	506	4/4	0.79	0.17	68,83,89,91	0
3	EDO	BBB	509	4/4	0.80	0.17	48,70,70,80	0
3	EDO	AAA	604	4/4	0.80	0.21	56,64,69,81	0
4	PGE	AAA	605	10/10	0.81	0.24	39,65,81,89	10
5	NDG	AAA	612	15/15	0.81	0.24	57,75,100,107	15
2	MPD	AAA	611	8/8	0.81	0.28	67,76,93,96	8
7	K	AAA	615	1/1	0.82	0.29	70,70,70,70	0
5	NDG	BBB	510	15/15	0.82	0.21	49,63,71,86	15
3	EDO	AAA	602	4/4	0.83	0.15	56,60,62,77	0
3	EDO	AAA	609	4/4	0.86	0.17	44,53,62,82	0
2	MPD	BBB	503	8/8	0.86	0.25	36,71,83,88	0
3	EDO	BBB	504	4/4	0.86	0.14	56,58,65,79	0
3	EDO	AAA	606	4/4	0.87	0.12	68,69,77,83	0
2	MPD	AAA	601	8/8	0.88	0.21	37,64,73,78	0
3	EDO	BBB	505	4/4	0.90	0.12	42,54,62,65	0
3	EDO	AAA	616	4/4	0.91	0.25	42,51,58,63	0
3	EDO	BBB	502	4/4	0.93	0.17	43,51,53,53	0
7	K	BBB	513	1/1	0.94	0.22	60,60,60,60	0
8	CL	AAA	617	1/1	0.96	0.07	72,72,72,72	0
6	ZN	BBB	512	1/1	0.97	0.12	56,56,56,56	0
8	CL	AAA	618	1/1	0.98	0.04	41,41,41,41	0
6	ZN	AAA	613	1/1	0.99	0.03	44,44,44,44	0
6	ZN	AAA	614	1/1	0.99	0.04	44,44,44,44	0
6	ZN	BBB	511	1/1	1.00	0.03	38,38,38,38	0

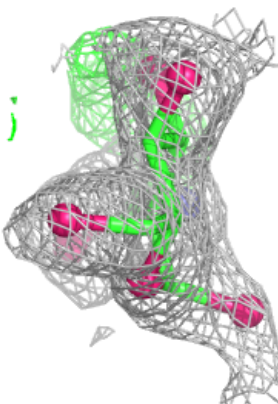
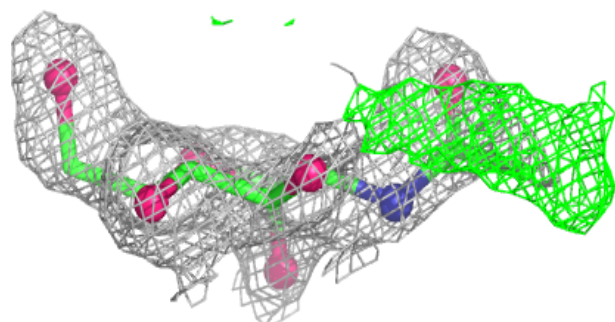
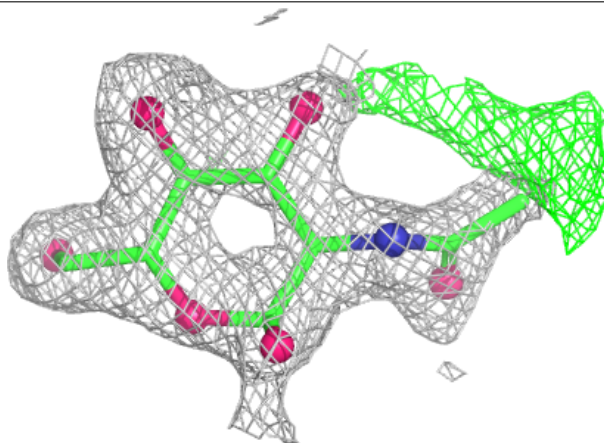
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around NDG AAA 612:

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

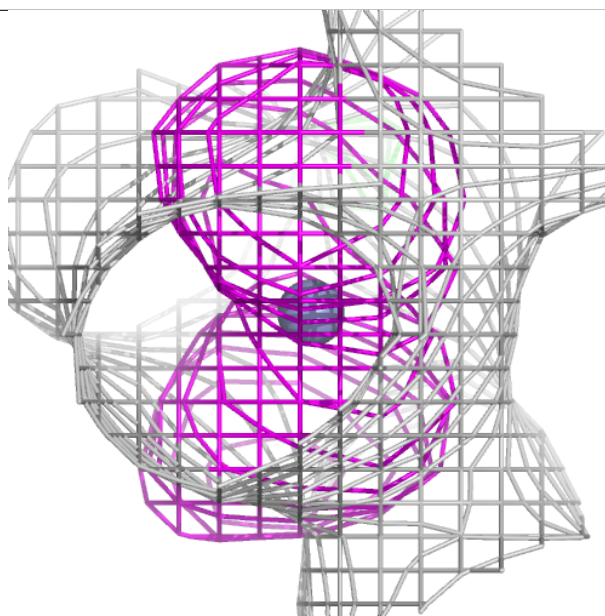
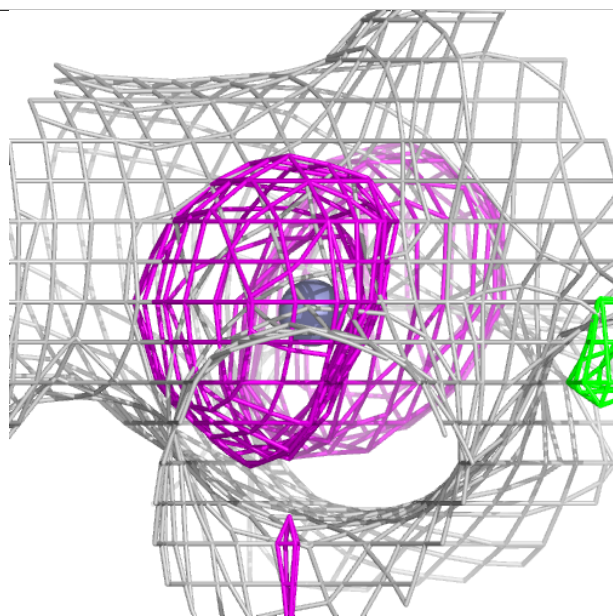
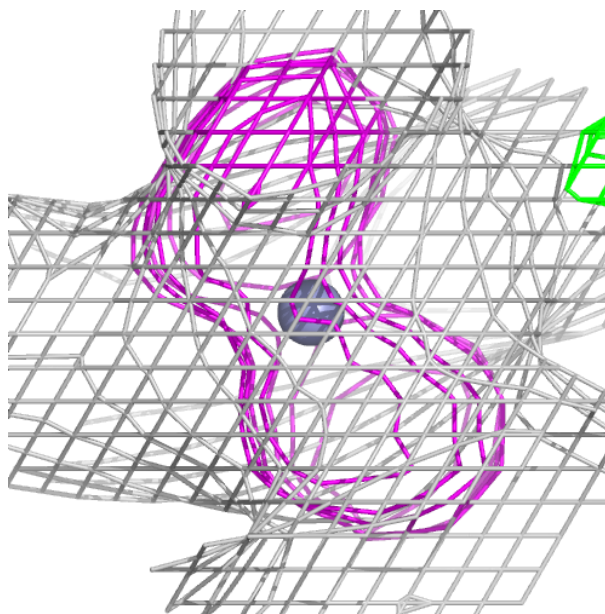
**Electron density around NDG BBB 510:**

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



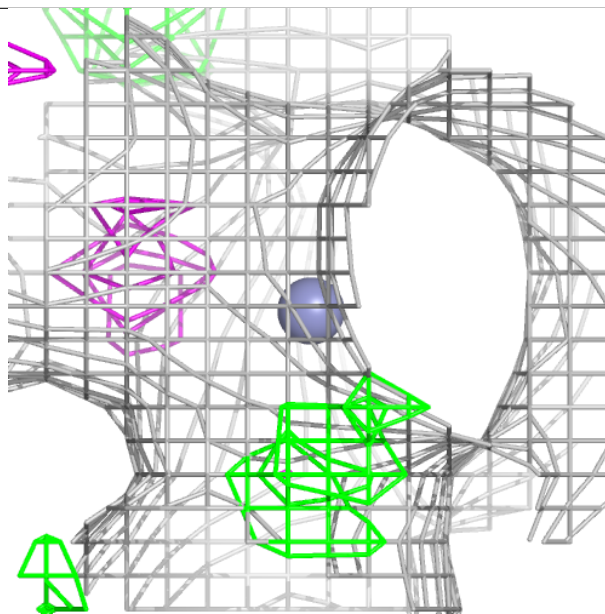
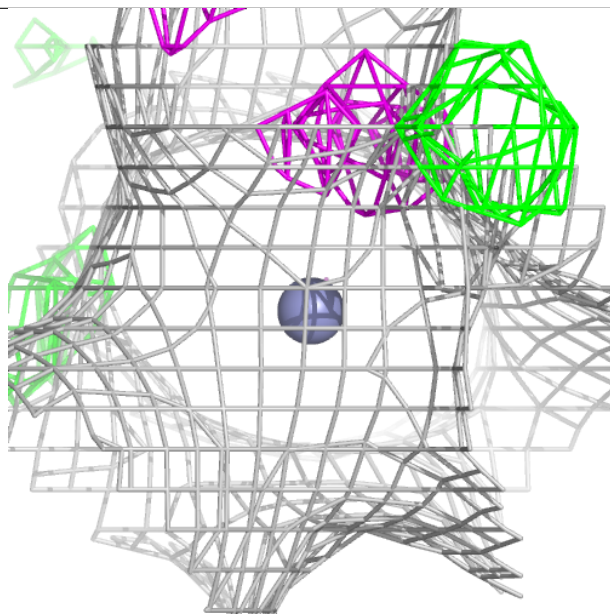
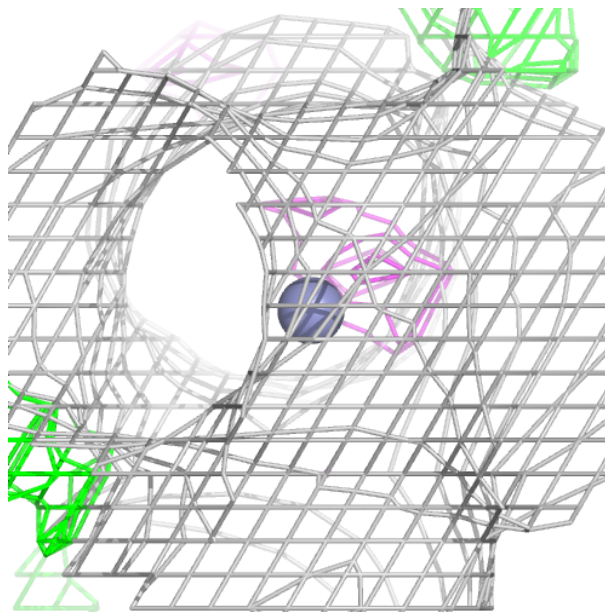
Electron density around ZN BBB 512:

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



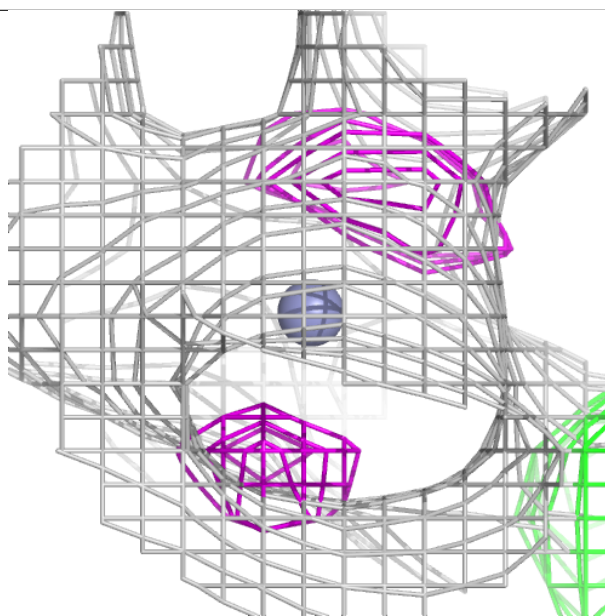
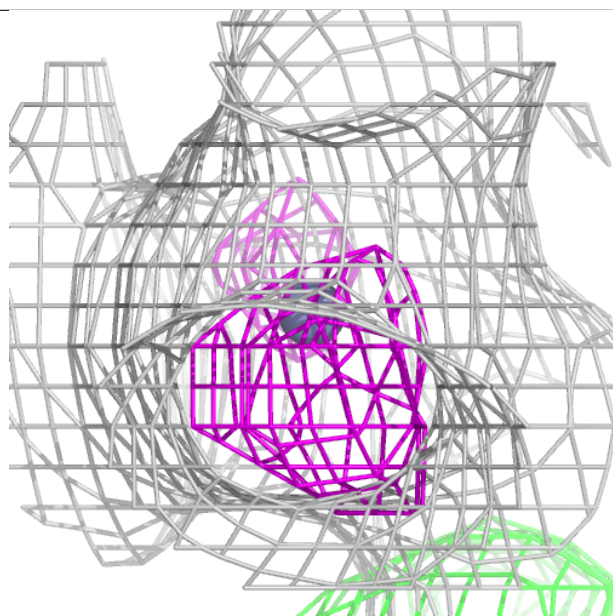
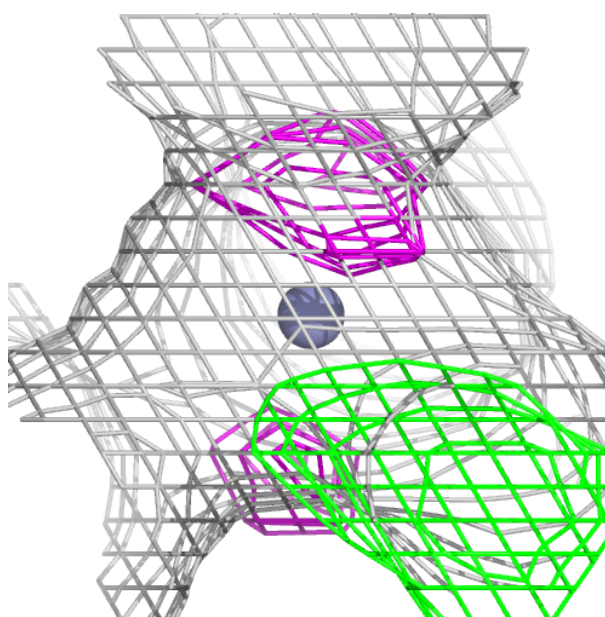
Electron density around ZN AAA 613:

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



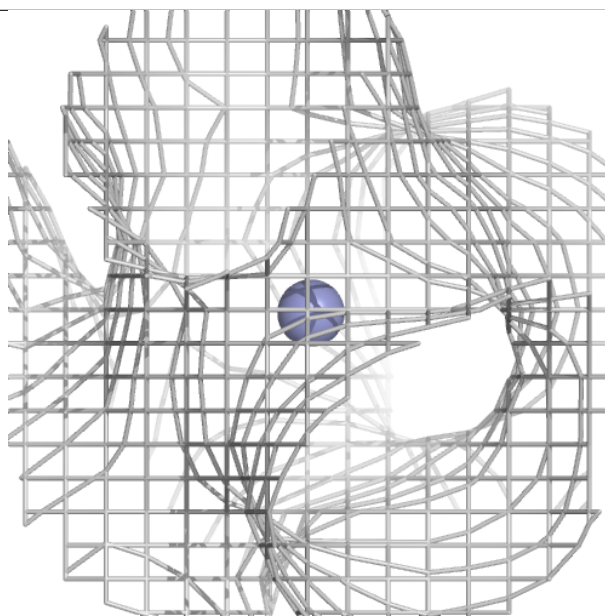
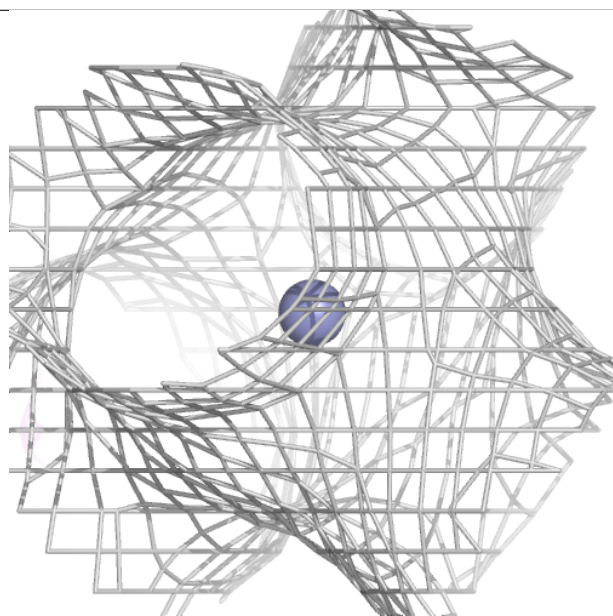
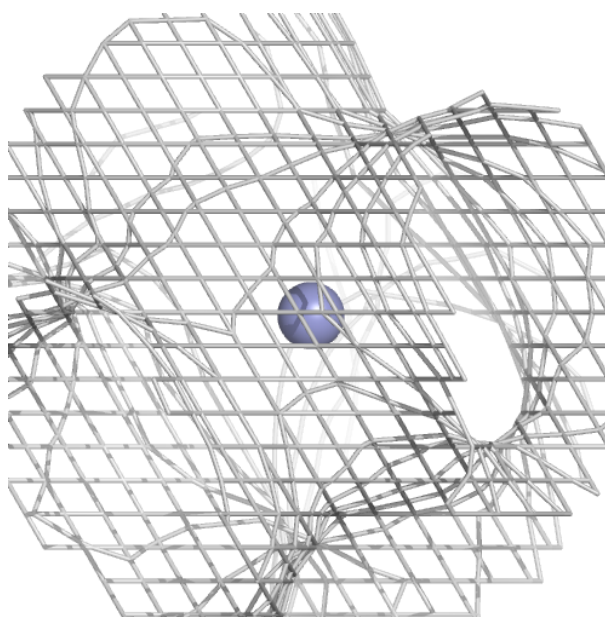
Electron density around ZN AAA 614:

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



Electron density around ZN BBB 511:

$2mF_o-DF_c$ (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.