



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 08:45 AM UTC

PDB ID : 6T94 / pdb_00006t94
Title : NAD⁺-dependent fungal formate dehydrogenase from *Chaetomium thermophilum*: A complex of N120C mutant protein with the reduced form of the cofactor NADH.
Authors : Isupov, M.N.; Yelmazer, B.; De Rose, S.A.; Littlechild, J.A.
Deposited on : 2019-10-25
Resolution : 1.15 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

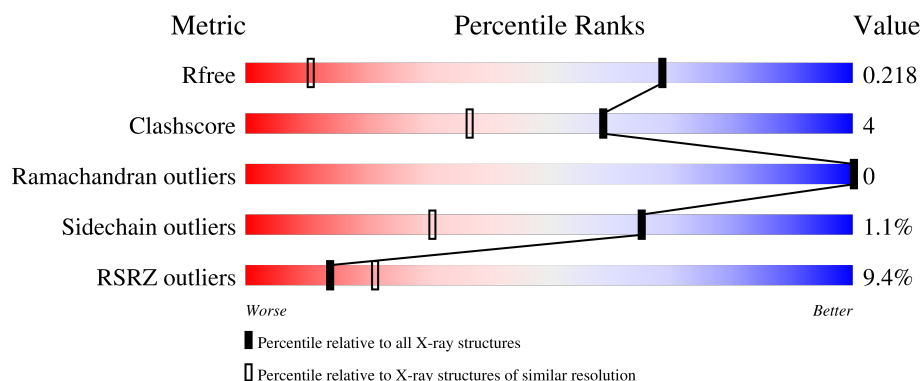
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.15 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1380 (1.16-1.12)
Clashscore	190562	1393 (1.16-1.12)
Ramachandran outliers	187476	1369 (1.16-1.12)
Sidechain outliers	187428	1369 (1.16-1.12)
RSRZ outliers	180081	1379 (1.16-1.12)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	AAA	410	2% 82% 8% 10%
1	BBB	410	9% 83% 8% 9%
1	CCC	410	18% 82% 8% 10%
1	DDD	410	5% 81% 8% 10%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 27255 atoms, of which 12811 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 Formate dehydrogenase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	AAA	367	Total	C	H	N	O	S	119	33	0
			6210	1935	3148	547	566	14			
1	BBB	374	Total	C	H	N	O	S	119	34	0
			6312	1973	3195	549	582	13			
1	CCC	371	Total	C	H	N	O	S	115	31	0
			6259	1951	3177	550	569	12			
1	DDD	367	Total	C	H	N	O	S	120	28	0
			6104	1906	3085	536	563	14			

There are 164 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
AAA	-33	MET	-	initiating methionine	UNP G0SGU4
AAA	-32	ALA	-	expression tag	UNP G0SGU4
AAA	-31	HIS	-	expression tag	UNP G0SGU4
AAA	-30	HIS	-	expression tag	UNP G0SGU4
AAA	-29	HIS	-	expression tag	UNP G0SGU4
AAA	-28	HIS	-	expression tag	UNP G0SGU4
AAA	-27	HIS	-	expression tag	UNP G0SGU4
AAA	-26	HIS	-	expression tag	UNP G0SGU4
AAA	-25	VAL	-	expression tag	UNP G0SGU4
AAA	-24	GLY	-	expression tag	UNP G0SGU4
AAA	-23	THR	-	expression tag	UNP G0SGU4
AAA	-22	GLY	-	expression tag	UNP G0SGU4
AAA	-21	SER	-	expression tag	UNP G0SGU4
AAA	-20	ASN	-	expression tag	UNP G0SGU4
AAA	-19	ASP	-	expression tag	UNP G0SGU4
AAA	-18	ASP	-	expression tag	UNP G0SGU4
AAA	-17	ASP	-	expression tag	UNP G0SGU4
AAA	-16	ASP	-	expression tag	UNP G0SGU4
AAA	-15	LYS	-	expression tag	UNP G0SGU4
AAA	-14	SER	-	expression tag	UNP G0SGU4
AAA	-13	PRO	-	expression tag	UNP G0SGU4

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Chain	Residue	Modelled	Actual	Comment	Reference
AAA	-12	ASP	-	expression tag	UNP G0SGU4
AAA	-11	PRO	-	expression tag	UNP G0SGU4
AAA	-10	ASN	-	expression tag	UNP G0SGU4
AAA	-9	TRP	-	expression tag	UNP G0SGU4
AAA	-8	GLU	-	expression tag	UNP G0SGU4
AAA	-7	LEU	-	expression tag	UNP G0SGU4
AAA	-6	VAL	-	expression tag	UNP G0SGU4
AAA	-5	TYR	-	expression tag	UNP G0SGU4
AAA	-4	THR	-	expression tag	UNP G0SGU4
AAA	-3	ALA	-	expression tag	UNP G0SGU4
AAA	-2	ARG	-	expression tag	UNP G0SGU4
AAA	-1	LEU	-	expression tag	UNP G0SGU4
AAA	0	GLN	-	expression tag	UNP G0SGU4
AAA	120	CYS	ASN	engineered mutation	UNP G0SGU4
AAA	371	HIS	-	expression tag	UNP G0SGU4
AAA	372	HIS	-	expression tag	UNP G0SGU4
AAA	373	HIS	-	expression tag	UNP G0SGU4
AAA	374	HIS	-	expression tag	UNP G0SGU4
AAA	375	HIS	-	expression tag	UNP G0SGU4
AAA	376	HIS	-	expression tag	UNP G0SGU4
BBB	-33	MET	-	initiating methionine	UNP G0SGU4
BBB	-32	ALA	-	expression tag	UNP G0SGU4
BBB	-31	HIS	-	expression tag	UNP G0SGU4
BBB	-30	HIS	-	expression tag	UNP G0SGU4
BBB	-29	HIS	-	expression tag	UNP G0SGU4
BBB	-28	HIS	-	expression tag	UNP G0SGU4
BBB	-27	HIS	-	expression tag	UNP G0SGU4
BBB	-26	HIS	-	expression tag	UNP G0SGU4
BBB	-25	VAL	-	expression tag	UNP G0SGU4
BBB	-24	GLY	-	expression tag	UNP G0SGU4
BBB	-23	THR	-	expression tag	UNP G0SGU4
BBB	-22	GLY	-	expression tag	UNP G0SGU4
BBB	-21	SER	-	expression tag	UNP G0SGU4
BBB	-20	ASN	-	expression tag	UNP G0SGU4
BBB	-19	ASP	-	expression tag	UNP G0SGU4
BBB	-18	ASP	-	expression tag	UNP G0SGU4
BBB	-17	ASP	-	expression tag	UNP G0SGU4
BBB	-16	ASP	-	expression tag	UNP G0SGU4
BBB	-15	LYS	-	expression tag	UNP G0SGU4
BBB	-14	SER	-	expression tag	UNP G0SGU4
BBB	-13	PRO	-	expression tag	UNP G0SGU4
BBB	-12	ASP	-	expression tag	UNP G0SGU4

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Chain	Residue	Modelled	Actual	Comment	Reference
BBB	-11	PRO	-	expression tag	UNP G0SGU4
BBB	-10	ASN	-	expression tag	UNP G0SGU4
BBB	-9	TRP	-	expression tag	UNP G0SGU4
BBB	-8	GLU	-	expression tag	UNP G0SGU4
BBB	-7	LEU	-	expression tag	UNP G0SGU4
BBB	-6	VAL	-	expression tag	UNP G0SGU4
BBB	-5	TYR	-	expression tag	UNP G0SGU4
BBB	-4	THR	-	expression tag	UNP G0SGU4
BBB	-3	ALA	-	expression tag	UNP G0SGU4
BBB	-2	ARG	-	expression tag	UNP G0SGU4
BBB	-1	LEU	-	expression tag	UNP G0SGU4
BBB	0	GLN	-	expression tag	UNP G0SGU4
BBB	120	CYS	ASN	engineered mutation	UNP G0SGU4
BBB	371	HIS	-	expression tag	UNP G0SGU4
BBB	372	HIS	-	expression tag	UNP G0SGU4
BBB	373	HIS	-	expression tag	UNP G0SGU4
BBB	374	HIS	-	expression tag	UNP G0SGU4
BBB	375	HIS	-	expression tag	UNP G0SGU4
BBB	376	HIS	-	expression tag	UNP G0SGU4
CCC	-33	MET	-	initiating methionine	UNP G0SGU4
CCC	-32	ALA	-	expression tag	UNP G0SGU4
CCC	-31	HIS	-	expression tag	UNP G0SGU4
CCC	-30	HIS	-	expression tag	UNP G0SGU4
CCC	-29	HIS	-	expression tag	UNP G0SGU4
CCC	-28	HIS	-	expression tag	UNP G0SGU4
CCC	-27	HIS	-	expression tag	UNP G0SGU4
CCC	-26	HIS	-	expression tag	UNP G0SGU4
CCC	-25	VAL	-	expression tag	UNP G0SGU4
CCC	-24	GLY	-	expression tag	UNP G0SGU4
CCC	-23	THR	-	expression tag	UNP G0SGU4
CCC	-22	GLY	-	expression tag	UNP G0SGU4
CCC	-21	SER	-	expression tag	UNP G0SGU4
CCC	-20	ASN	-	expression tag	UNP G0SGU4
CCC	-19	ASP	-	expression tag	UNP G0SGU4
CCC	-18	ASP	-	expression tag	UNP G0SGU4
CCC	-17	ASP	-	expression tag	UNP G0SGU4
CCC	-16	ASP	-	expression tag	UNP G0SGU4
CCC	-15	LYS	-	expression tag	UNP G0SGU4
CCC	-14	SER	-	expression tag	UNP G0SGU4
CCC	-13	PRO	-	expression tag	UNP G0SGU4
CCC	-12	ASP	-	expression tag	UNP G0SGU4
CCC	-11	PRO	-	expression tag	UNP G0SGU4

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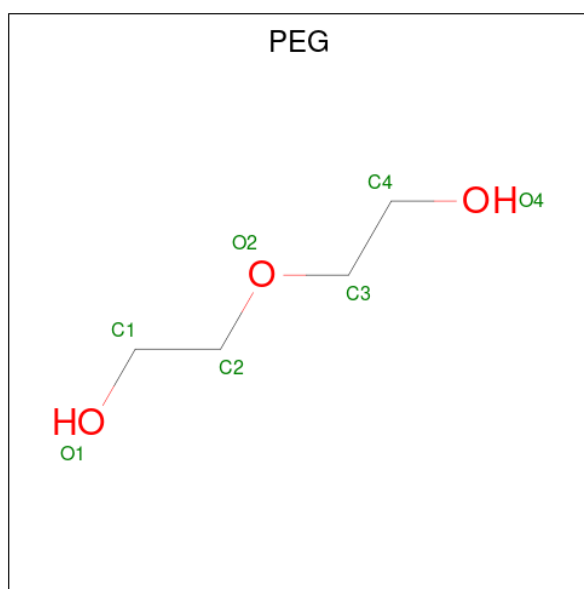
Chain	Residue	Modelled	Actual	Comment	Reference
CCC	-10	ASN	-	expression tag	UNP G0SGU4
CCC	-9	TRP	-	expression tag	UNP G0SGU4
CCC	-8	GLU	-	expression tag	UNP G0SGU4
CCC	-7	LEU	-	expression tag	UNP G0SGU4
CCC	-6	VAL	-	expression tag	UNP G0SGU4
CCC	-5	TYR	-	expression tag	UNP G0SGU4
CCC	-4	THR	-	expression tag	UNP G0SGU4
CCC	-3	ALA	-	expression tag	UNP G0SGU4
CCC	-2	ARG	-	expression tag	UNP G0SGU4
CCC	-1	LEU	-	expression tag	UNP G0SGU4
CCC	0	GLN	-	expression tag	UNP G0SGU4
CCC	120	CYS	ASN	engineered mutation	UNP G0SGU4
CCC	371	HIS	-	expression tag	UNP G0SGU4
CCC	372	HIS	-	expression tag	UNP G0SGU4
CCC	373	HIS	-	expression tag	UNP G0SGU4
CCC	374	HIS	-	expression tag	UNP G0SGU4
CCC	375	HIS	-	expression tag	UNP G0SGU4
CCC	376	HIS	-	expression tag	UNP G0SGU4
DDD	-33	MET	-	initiating methionine	UNP G0SGU4
DDD	-32	ALA	-	expression tag	UNP G0SGU4
DDD	-31	HIS	-	expression tag	UNP G0SGU4
DDD	-30	HIS	-	expression tag	UNP G0SGU4
DDD	-29	HIS	-	expression tag	UNP G0SGU4
DDD	-28	HIS	-	expression tag	UNP G0SGU4
DDD	-27	HIS	-	expression tag	UNP G0SGU4
DDD	-26	HIS	-	expression tag	UNP G0SGU4
DDD	-25	VAL	-	expression tag	UNP G0SGU4
DDD	-24	GLY	-	expression tag	UNP G0SGU4
DDD	-23	THR	-	expression tag	UNP G0SGU4
DDD	-22	GLY	-	expression tag	UNP G0SGU4
DDD	-21	SER	-	expression tag	UNP G0SGU4
DDD	-20	ASN	-	expression tag	UNP G0SGU4
DDD	-19	ASP	-	expression tag	UNP G0SGU4
DDD	-18	ASP	-	expression tag	UNP G0SGU4
DDD	-17	ASP	-	expression tag	UNP G0SGU4
DDD	-16	ASP	-	expression tag	UNP G0SGU4
DDD	-15	LYS	-	expression tag	UNP G0SGU4
DDD	-14	SER	-	expression tag	UNP G0SGU4
DDD	-13	PRO	-	expression tag	UNP G0SGU4
DDD	-12	ASP	-	expression tag	UNP G0SGU4
DDD	-11	PRO	-	expression tag	UNP G0SGU4
DDD	-10	ASN	-	expression tag	UNP G0SGU4

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Chain	Residue	Modelled	Actual	Comment	Reference
DDD	-9	TRP	-	expression tag	UNP G0SGU4
DDD	-8	GLU	-	expression tag	UNP G0SGU4
DDD	-7	LEU	-	expression tag	UNP G0SGU4
DDD	-6	VAL	-	expression tag	UNP G0SGU4
DDD	-5	TYR	-	expression tag	UNP G0SGU4
DDD	-4	THR	-	expression tag	UNP G0SGU4
DDD	-3	ALA	-	expression tag	UNP G0SGU4
DDD	-2	ARG	-	expression tag	UNP G0SGU4
DDD	-1	LEU	-	expression tag	UNP G0SGU4
DDD	0	GLN	-	expression tag	UNP G0SGU4
DDD	120	CYS	ASN	engineered mutation	UNP G0SGU4
DDD	371	HIS	-	expression tag	UNP G0SGU4
DDD	372	HIS	-	expression tag	UNP G0SGU4
DDD	373	HIS	-	expression tag	UNP G0SGU4
DDD	374	HIS	-	expression tag	UNP G0SGU4
DDD	375	HIS	-	expression tag	UNP G0SGU4
DDD	376	HIS	-	expression tag	UNP G0SGU4

- Molecule 2 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



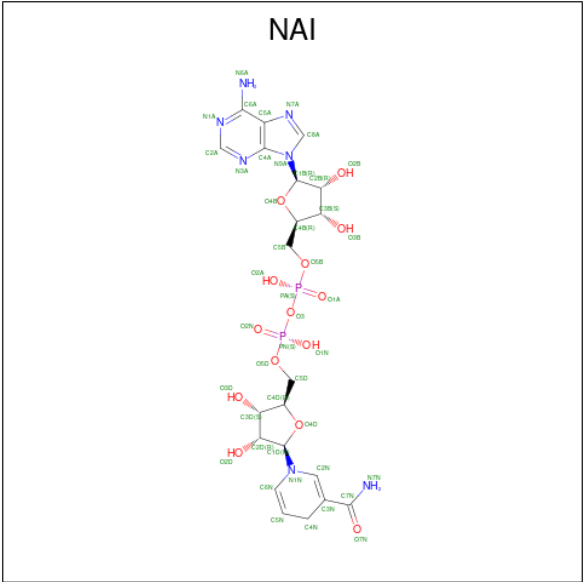
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	AAA	1	Total	C	H	O	1	0
			17	4	10	3		
2	AAA	1	Total	C	H	O	1	0
			17	4	10	3		
2	CCC	1	Total	C	H	O	1	0
			17	4	10	3		

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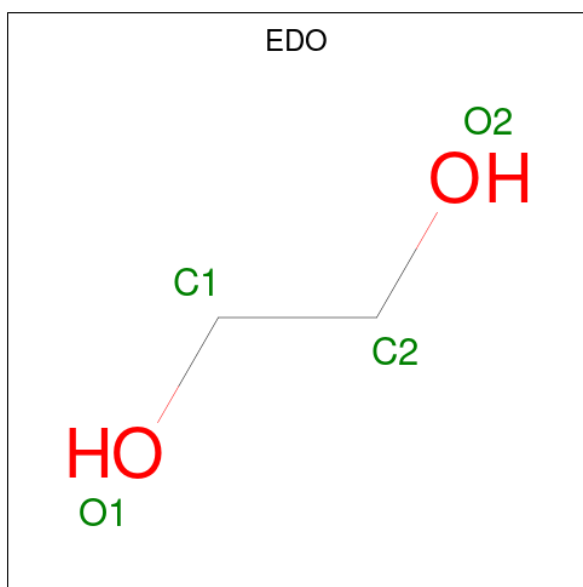
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	H	O		
2	DDD	1	17	4	10	3	1	0
2	DDD	1	17	4	10	3	1	0

- Molecule 3 is 1,4-DIHYDRONICOTINAMIDE ADENINE DINUCLEOTIDE (CCD ID: NAI) (formula: C₂₁H₂₉N₇O₁₄P₂) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
			Total	C	H	N	O	P		
3	AAA	1	71	21	27	7	14	2	4	0
3	BBB	1	71	21	27	7	14	2	4	0
3	CCC	1	71	21	27	7	14	2	4	0
3	DDD	1	71	21	27	7	14	2	4	0

- Molecule 4 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	AAA	1	Total	C	H	O	1	0
			10	2	6	2		
4	AAA	1	Total	C	H	O	1	0
			10	2	6	2		
4	AAA	1	Total	C	H	O	1	0
			10	2	6	2		
4	AAA	1	Total	C	H	O	1	0
			10	2	6	2		
4	BBB	1	Total	C	H	O	1	0
			10	2	6	2		
4	CCC	1	Total	C	H	O	1	0
			10	2	6	2		
4	DDD	1	Total	C	H	O	1	0
			10	2	6	2		
4	DDD	1	Total	C	H	O	1	0
			10	2	6	2		

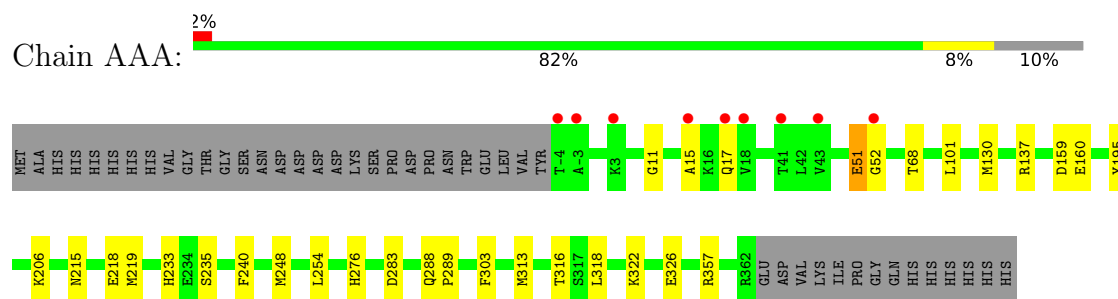
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	AAA	515	Total	O	0	0
			515	515		
5	BBB	454	Total	O	0	0
			454	454		
5	CCC	448	Total	O	0	0
			448	448		
5	DDD	504	Total	O	0	0
			504	504		

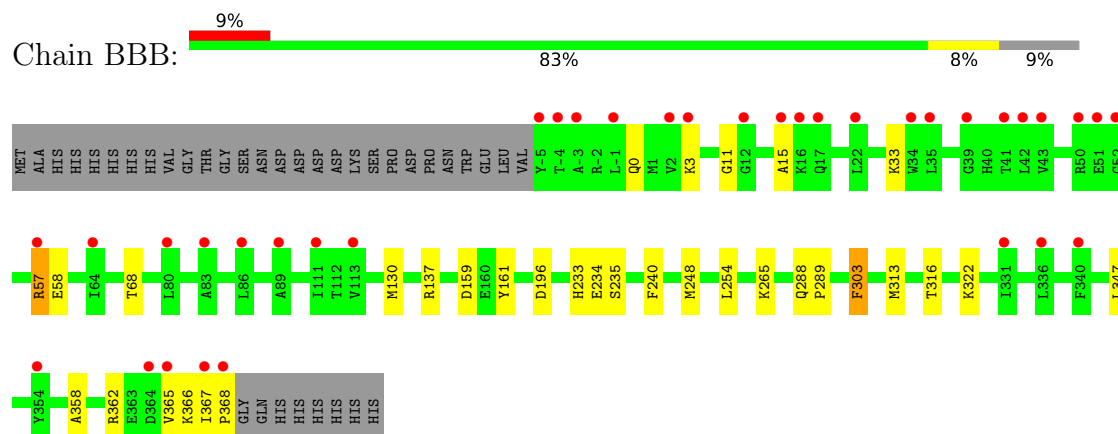
3 Residue-property plots [i](#)

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

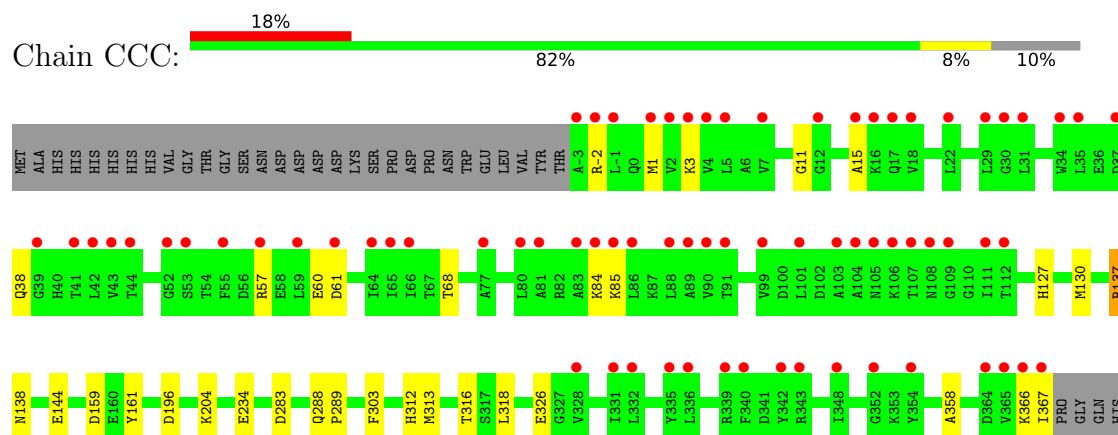
• Molecule 1: Formate dehydrogenase



• Molecule 1: Formate dehydrogenase



• Molecule 1: Formate dehydrogenase



[illegible]

4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	50.73Å 94.75Å 94.60Å 85.96° 89.97° 81.58°	Depositor
Resolution (Å)	68.90 – 1.15 68.90 – 1.15	Depositor EDS
% Data completeness (in resolution range)	86.8 (68.90-1.15) 86.8 (68.90-1.15)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.09 (at 1.15Å)	Xtriage
Refinement program	REFMAC 5.8.0257	Depositor
R, R_{free}	0.196 , 0.218 0.196 , 0.218	Depositor DCC
R_{free} test set	26891 reflections (4.35%)	wwPDB-VP
Wilson B-factor (Å ²)	13.3	Xtriage
Anisotropy	0.530	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 49.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	27255	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.47% 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: NAI, PEG, EDO

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	1.05	2/3198 (0.1%)	1.22	5/4306 (0.1%)
1	BBB	1.04	0/3261	1.20	5/4395 (0.1%)
1	CCC	1.09	3/3218 (0.1%)	1.23	4/4333 (0.1%)
1	DDD	1.13	3/3146 (0.1%)	1.26	5/4241 (0.1%)
All	All	1.08	8/12823 (0.1%)	1.23	19/17275 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	AAA	0	1
1	BBB	0	1
1	CCC	0	1
1	DDD	0	1
All	All	0	4

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	DDD	127	HIS	CE1-NE2	7.93	1.40	1.32
1	CCC	127	HIS	CE1-NE2	6.57	1.39	1.32
1	CCC	144	GLU	C-O	5.61	1.30	1.24
1	CCC	137	ARG	NE-CZ	-5.33	1.27	1.33
1	AAA	276	HIS	CE1-NE2	5.24	1.37	1.32
1	DDD	225	VAL	C-N	-5.07	1.26	1.33
1	AAA	160	GLU	C-O	5.01	1.30	1.24
1	DDD	233	HIS	CE1-NE2	5.00	1.37	1.32

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	DDD	168	VAL	CA-C-N	7.71	127.96	122.33
1	DDD	168	VAL	C-N-CA	7.71	127.96	122.33
1	AAA	68	THR	CB-CA-C	6.01	117.46	109.65
1	AAA	283	ASP	CA-CB-CG	5.84	118.44	112.60
1	DDD	68	THR	CB-CA-C	5.78	117.17	109.65
1	CCC	68	THR	CB-CA-C	5.72	117.09	109.65
1	CCC	138	ASN	OD1-CG-ND2	5.64	128.24	122.60
1	BBB	68	THR	CB-CA-C	5.60	116.94	109.65
1	DDD	130	MET	O-C-N	5.52	127.83	122.09
1	DDD	283	ASP	CA-CB-CG	5.25	117.85	112.60
1	AAA	303[A]	PHE	CA-CB-CG	5.20	119.00	113.80
1	AAA	303[B]	PHE	CA-CB-CG	5.20	119.00	113.80
1	BBB	130	MET	O-C-N	5.18	127.48	122.09
1	AAA	130	MET	O-C-N	5.15	127.44	122.09
1	BBB	196	ASP	CA-CB-CG	5.13	117.73	112.60
1	CCC	196	ASP	CA-CB-CG	5.09	117.69	112.60
1	CCC	130	MET	O-C-N	5.00	127.29	122.09
1	BBB	303[A]	PHE	CA-CB-CG	5.00	118.80	113.80
1	BBB	303[B]	PHE	CA-CB-CG	5.00	118.80	113.80

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	AAA	137	ARG	Sidechain
1	BBB	137	ARG	Sidechain
1	CCC	137	ARG	Sidechain
1	DDD	137	ARG	Sidechain

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	AAA	3062	3148	3163	28	0
1	BBB	3117	3195	3214	23	0
1	CCC	3082	3177	3194	33	0
1	DDD	3019	3085	3099	26	0
2	AAA	14	20	20	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	CCC	7	10	10	0	0
2	DDD	14	20	20	0	0
3	AAA	44	27	27	1	0
3	BBB	44	27	27	2	0
3	CCC	44	27	27	2	0
3	DDD	44	27	27	1	0
4	AAA	16	24	24	5	0
4	BBB	4	6	6	0	0
4	CCC	4	6	6	1	0
4	DDD	8	12	12	0	0
5	AAA	515	0	0	8	2
5	BBB	454	0	0	3	2
5	CCC	448	0	0	6	0
5	DDD	504	0	0	3	0
All	All	14444	12811	12876	111	2

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.

All (111) 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:CCC:3[A]:LYS:NZ	1:CCC:85[A]:LYS:HZ1	1.55	1.03
1:CCC:3[A]:LYS:HZ3	1:CCC:85[A]:LYS:NZ	1.67	0.93
1:AAA:195:TYR:HB2	1:AAA:219[B]:MET:HE2	1.49	0.92
1:CCC:159[A]:ASP:OD1	1:CCC:303[A]:PHE:CE2	2.25	0.89
1:DDD:195:TYR:HB2	1:DDD:219[B]:MET:HE2	1.54	0.88
1:DDD:195:TYR:N	1:DDD:219[B]:MET:HE1	1.89	0.88
1:CCC:-2:ARG:HH21	1:CCC:38:GLN:NE2	1.75	0.85
1:BBB:248[B]:MET:HE1	1:BBB:254:LEU:HD13	1.59	0.85
4:CCC:403:EDO:H22	5:CCC:632:HOH:O	1.75	0.85
1:DDD:195:TYR:HB2	1:DDD:219[B]:MET:CE	2.08	0.84
1:BBB:240:PHE:HE2	1:BBB:248[B]:MET:CE	1.93	0.82
1:CCC:3[A]:LYS:NZ	1:CCC:85[A]:LYS:NZ	2.26	0.81
1:CCC:3[A]:LYS:HZ3	1:CCC:85[A]:LYS:HZ1	0.81	0.81
1:DDD:195:TYR:CB	1:DDD:219[B]:MET:CE	2.63	0.77
1:CCC:159[A]:ASP:OD1	1:CCC:303[A]:PHE:HE2	1.67	0.77
1:DDD:248[B]:MET:HE1	1:DDD:254:LEU:HD13	1.67	0.76
1:CCC:-2:ARG:NH2	1:CCC:38:GLN:NE2	2.34	0.75
1:DDD:240:PHE:HE2	1:DDD:248[B]:MET:HE2	1.50	0.74
1:CCC:85[A]:LYS:HD2	5:CCC:844:HOH:O	1.92	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:DDD:240:PHE:HE2	1:DDD:248[B]:MET:CE	2.06	0.69
1:AAA:240:PHE:HE2	1:AAA:248[B]:MET:HE2	1.58	0.68
1:CCC:159[A]:ASP:OD1	1:CCC:303[A]:PHE:CD2	2.46	0.67
1:AAA:101:LEU:N	4:AAA:507:EDO:H21	2.10	0.67
1:AAA:326[A]:GLU:HG3	5:AAA:630:HOH:O	1.95	0.66
1:AAA:195:TYR:CB	1:AAA:219[B]:MET:HE2	2.24	0.66
1:CCC:159[A]:ASP:OD2	1:CCC:303[A]:PHE:CD2	2.49	0.66
1:AAA:206[B]:LYS:HE2	5:AAA:637:HOH:O	1.96	0.65
1:CCC:159[A]:ASP:CG	1:CCC:303[A]:PHE:CD2	2.75	0.65
1:AAA:195:TYR:HB2	1:AAA:219[B]:MET:CE	2.26	0.64
1:DDD:288:GLN:HA	1:DDD:289:PRO:C	2.23	0.63
1:CCC:288:GLN:HA	1:CCC:289:PRO:C	2.23	0.63
1:BBB:288:GLN:HA	1:BBB:289:PRO:C	2.23	0.63
1:AAA:195:TYR:CB	1:AAA:219[B]:MET:CE	2.77	0.62
1:AAA:206[B]:LYS:CE	5:AAA:637:HOH:O	2.46	0.62
1:DDD:240:PHE:CE2	1:DDD:248[B]:MET:CE	2.83	0.61
1:CCC:-2:ARG:NH2	1:CCC:38:GLN:HE22	1.99	0.60
1:AAA:288:GLN:HA	1:AAA:289:PRO:C	2.25	0.60
1:CCC:159[A]:ASP:CG	1:CCC:303[A]:PHE:HD2	2.09	0.60
1:BBB:240:PHE:CE2	1:BBB:248[B]:MET:CE	2.80	0.58
1:DDD:326[A]:GLU:HG3	5:DDD:619:HOH:O	2.03	0.58
1:AAA:357[B]:ARG:NH2	5:AAA:613:HOH:O	2.36	0.58
1:CCC:159[A]:ASP:OD2	1:CCC:303[A]:PHE:HD2	1.86	0.58
1:DDD:195:TYR:CB	1:DDD:219[B]:MET:HE1	2.33	0.58
1:AAA:248[B]:MET:HE1	1:AAA:254:LEU:HD13	1.86	0.57
1:BBB:234[B]:GLU:HG2	5:BBB:1232:HOH:O	2.05	0.56
1:DDD:240:PHE:CE2	1:DDD:248[B]:MET:HE2	2.36	0.56
1:AAA:101:LEU:H	4:AAA:507:EDO:H21	1.70	0.56
1:AAA:215:ASN:OD1	1:AAA:218[B]:GLU:HG2	2.05	0.56
1:AAA:240:PHE:CE2	1:AAA:248[B]:MET:HE2	2.41	0.55
1:BBB:347:LEU:HD12	1:BBB:368:PRO:CG	2.37	0.55
1:BBB:362:ARG:HB2	1:BBB:365:VAL:HG13	1.88	0.55
1:AAA:51[A]:GLU:OE2	1:AAA:51[A]:GLU:O	2.26	0.53
1:AAA:233:HIS:CE1	1:AAA:235[A]:SER:OG	2.62	0.53
1:AAA:17[A]:GLN:OE1	1:BBB:303[A]:PHE:CE1	2.62	0.53
1:CCC:326[A]:GLU:HG3	5:CCC:612:HOH:O	2.08	0.53
1:DDD:17[B]:GLN:NE2	5:DDD:509:HOH:O	2.43	0.52
1:DDD:194:TYR:C	1:DDD:219[B]:MET:HE1	2.34	0.52
3:AAA:503:NAI:H42N	5:AAA:938:HOH:O	2.10	0.50
1:DDD:195:TYR:HB3	1:DDD:219[B]:MET:CE	2.41	0.50
1:CCC:-2:ARG:HH21	1:CCC:38:GLN:HE22	1.50	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CCC:61:ASP:HB2	1:CCC:85[A]:LYS:HE3	1.95	0.48
1:AAA:51[A]:GLU:HG3	1:AAA:52[A]:GLY:N	2.29	0.48
3:CCC:402:NAI:H42N	5:CCC:789:HOH:O	2.13	0.48
3:BBB:901:NAI:H42N	5:BBB:1306:HOH:O	2.13	0.48
1:AAA:195:TYR:HB3	1:AAA:219[B]:MET:CE	2.44	0.48
1:BBB:265[A]:LYS:NZ	5:BBB:1024:HOH:O	2.47	0.48
1:DDD:84[A]:LYS:NZ	5:DDD:512:HOH:O	2.46	0.47
4:AAA:504:EDO:H22	5:AAA:606:HOH:O	2.15	0.47
1:BBB:240:PHE:HE2	1:BBB:248[B]:MET:HE2	1.77	0.47
1:CCC:61:ASP:HB2	1:CCC:85[A]:LYS:CE	2.44	0.47
1:BBB:57[A]:ARG:HH11	1:BBB:57[A]:ARG:HB3	1.80	0.46
1:BBB:347:LEU:HD12	1:BBB:368:PRO:HG3	1.98	0.46
1:BBB:347:LEU:HD12	1:BBB:368:PRO:HG2	1.97	0.45
1:AAA:313:MET:HA	1:AAA:316:THR:HG22	1.98	0.45
1:CCC:11:GLY:HA3	1:CCC:15:ALA:HB2	1.98	0.45
1:AAA:206[B]:LYS:HG3	5:AAA:783:HOH:O	2.16	0.45
1:BBB:233:HIS:CE1	1:BBB:235[B]:SER:OG	2.70	0.44
1:AAA:248[B]:MET:HE1	1:AAA:254:LEU:HD22	2.00	0.44
1:CCC:313:MET:HA	1:CCC:316:THR:HG22	1.98	0.44
1:AAA:11:GLY:HA3	1:AAA:15:ALA:HB2	2.00	0.44
1:CCC:303[A]:PHE:CE1	1:DDD:17[A]:GLN:OE1	2.70	0.44
1:BBB:240:PHE:CE2	1:BBB:248[B]:MET:HE3	2.52	0.44
1:DDD:10[A]:ASP:OD2	1:DDD:46:SER:OG	2.35	0.44
1:DDD:215:ASN:O	1:DDD:218[B]:GLU:HB3	2.18	0.43
1:DDD:11:GLY:HA3	1:DDD:15:ALA:HB2	2.00	0.43
1:BBB:313:MET:HA	1:BBB:316:THR:HG22	2.01	0.43
4:AAA:505:EDO:O1	1:BBB:322[B]:LYS:NZ	2.50	0.43
1:CCC:204[B]:LYS:HE3	5:CCC:532:HOH:O	2.17	0.43
1:CCC:-2:ARG:NH2	1:CCC:38:GLN:HE21	2.13	0.43
1:AAA:51[A]:GLU:CD	1:AAA:51[A]:GLU:C	2.85	0.43
1:AAA:318:LEU:HD12	1:BBB:161:TYR:CZ	2.53	0.43
1:CCC:161:TYR:CZ	1:DDD:318:LEU:HD12	2.54	0.43
1:BBB:11:GLY:HA3	1:BBB:15:ALA:HB2	2.01	0.42
1:AAA:195:TYR:HB3	1:AAA:219[B]:MET:HE1	2.00	0.42
1:CCC:60:GLU:O	1:CCC:84[B]:LYS:HG3	2.20	0.42
1:BBB:57[B]:ARG:NH2	1:BBB:58:GLU:OE2	2.53	0.42
2:AAA:502:PEG:H32	4:AAA:507:EDO:H22	2.01	0.42
1:AAA:322[B]:LYS:HD2	5:AAA:644:HOH:O	2.20	0.42
1:DDD:120[B]:CYS:HB3	1:DDD:320:ALA:HB1	2.02	0.42
1:BBB:358:ALA:O	3:BBB:901:NAI:H8A	2.19	0.41
1:CCC:318:LEU:HD12	1:DDD:161:TYR:CZ	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:DDD:313:MET:HA	1:DDD:316:THR:HG22	2.02	0.41
1:CCC:303[B]:PHE:CD2	5:CCC:571:HOH:O	2.57	0.41
1:BBB:3[A]:LYS:HE3	1:BBB:3[A]:LYS:HB3	1.85	0.41
1:CCC:3[A]:LYS:HZ1	1:CCC:85[A]:LYS:NZ	2.14	0.41
1:BBB:33[B]:LYS:HD3	1:BBB:33[B]:LYS:HA	1.94	0.41
1:CCC:358:ALA:O	3:CCC:402:NAI:H8A	2.21	0.40
1:CCC:283:ASP:O	1:CCC:312:HIS:HA	2.21	0.40
1:DDD:358:ALA:O	3:DDD:403:NAI:H8A	2.22	0.40
1:CCC:61:ASP:HA	1:CCC:85[A]:LYS:HG3	2.03	0.40
1:DDD:59:LEU:C	1:DDD:59:LEU:HD12	2.47	0.40

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

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AAA:647:HOH:O	5:BBB:1387:HOH:O[1_455]	2.19	0.01
5:AAA:894:HOH:O	5:BBB:1074:HOH:O[1_455]	2.19	0.01

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	398/410 (97%)	388 (98%)	10 (2%)	0	100	100
1	BBB	406/410 (99%)	394 (97%)	12 (3%)	0	100	100
1	CCC	400/410 (98%)	389 (97%)	11 (3%)	0	100	100
1	DDD	393/410 (96%)	382 (97%)	11 (3%)	0	100	100
All	All	1597/1640 (97%)	1553 (97%)	44 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	AAA	333/340 (98%)	329 (99%)	4 (1%)	63	26
1	BBB	341/340 (100%)	334 (98%)	7 (2%)	47	9
1	CCC	335/340 (98%)	328 (98%)	7 (2%)	47	9
1	DDD	328/340 (96%)	326 (99%)	2 (1%)	78	52
All	All	1337/1360 (98%)	1317 (98%)	20 (2%)	65	19

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	AAA	51[A]	GLU
1	AAA	51[B]	GLU
1	AAA	159[A]	ASP
1	AAA	159[B]	ASP
1	BBB	0	GLN
1	BBB	57[A]	ARG
1	BBB	57[B]	ARG
1	BBB	159[A]	ASP
1	BBB	159[B]	ASP
1	BBB	366	LYS
1	BBB	367	ILE
1	CCC	1	MET
1	CCC	57[A]	ARG
1	CCC	57[B]	ARG
1	CCC	234[A]	GLU
1	CCC	234[B]	GLU
1	CCC	366	LYS
1	CCC	367	ILE
1	DDD	159[A]	ASP
1	DDD	159[B]	ASP

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

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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

17 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
3	NAI	BBB	901	-	47,48,48	1.24	4 (8%)	64,73,73	1.99	11 (17%)
2	PEG	DDD	401	-	6,6,6	0.36	0	5,5,5	0.18	0
4	EDO	AAA	504	-	3,3,3	0.13	0	2,2,2	0.45	0
4	EDO	CCC	403	-	3,3,3	0.14	0	2,2,2	0.72	0
4	EDO	DDD	405	-	3,3,3	0.44	0	2,2,2	0.57	0
2	PEG	AAA	501	-	6,6,6	0.21	0	5,5,5	0.13	0
3	NAI	AAA	503	-	47,48,48	1.27	7 (14%)	64,73,73	1.51	14 (21%)
3	NAI	DDD	403	-	47,48,48	1.84	5 (10%)	64,73,73	2.07	16 (25%)
2	PEG	CCC	401	-	6,6,6	0.25	0	5,5,5	0.13	0
4	EDO	AAA	506	-	3,3,3	0.34	0	2,2,2	0.40	0
4	EDO	AAA	507	-	3,3,3	0.21	0	2,2,2	0.52	0
3	NAI	CCC	402	-	47,48,48	1.57	7 (14%)	64,73,73	2.05	17 (26%)
4	EDO	AAA	505	-	3,3,3	0.14	0	2,2,2	0.39	0
4	EDO	DDD	404	-	3,3,3	0.12	0	2,2,2	0.29	0
4	EDO	BBB	902	-	3,3,3	0.25	0	2,2,2	0.34	0
2	PEG	AAA	502	-	6,6,6	0.22	0	5,5,5	0.11	0
2	PEG	DDD	402	-	6,6,6	0.15	0	5,5,5	0.12	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAI	BBB	901	-	-	2/29/72/72	0/5/5/5
2	PEG	DDD	401	-	-	4/4/4/4	-
4	EDO	AAA	504	-	-	1/1/1/1	-
4	EDO	CCC	403	-	-	1/1/1/1	-
4	EDO	DDD	405	-	-	0/1/1/1	-
2	PEG	AAA	501	-	-	1/4/4/4	-
3	NAI	AAA	503	-	-	2/29/72/72	0/5/5/5
3	NAI	DDD	403	-	-	2/29/72/72	0/5/5/5
2	PEG	CCC	401	-	-	1/4/4/4	-
4	EDO	AAA	506	-	-	1/1/1/1	-
4	EDO	AAA	507	-	-	0/1/1/1	-
3	NAI	CCC	402	-	-	1/29/72/72	0/5/5/5
4	EDO	AAA	505	-	-	1/1/1/1	-
4	EDO	DDD	404	-	-	1/1/1/1	-
4	EDO	BBB	902	-	-	1/1/1/1	-
2	PEG	AAA	502	-	-	3/4/4/4	-
2	PEG	DDD	402	-	-	1/4/4/4	-

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	DDD	403	NAI	PN-O3	7.26	1.67	1.59
3	DDD	403	NAI	C5A-C4A	5.69	1.49	1.39
3	CCC	402	NAI	C5A-C4A	4.77	1.47	1.39
3	BBB	901	NAI	C5A-C4A	4.18	1.46	1.39
3	DDD	403	NAI	C2A-N1A	4.08	1.41	1.33
3	CCC	402	NAI	PN-O3	3.84	1.63	1.59
3	AAA	503	NAI	C4A-N9A	-3.62	1.30	1.37
3	CCC	402	NAI	C4N-C5N	-3.48	1.40	1.49
3	AAA	503	NAI	C5A-N7A	-3.24	1.33	1.39
3	DDD	403	NAI	C5A-N7A	-2.88	1.33	1.39
3	CCC	402	NAI	C6N-C5N	2.81	1.41	1.33
3	CCC	402	NAI	C4A-N9A	-2.74	1.32	1.37
3	CCC	402	NAI	PA-O3	-2.64	1.56	1.59
3	BBB	901	NAI	O3D-C3D	2.55	1.49	1.43
3	AAA	503	NAI	C2A-N1A	2.52	1.38	1.33
3	AAA	503	NAI	PN-O3	2.44	1.62	1.59
3	AAA	503	NAI	C6N-C5N	2.26	1.40	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	BBB	901	NAI	C4A-N9A	-2.26	1.33	1.37
3	AAA	503	NAI	C5A-C4A	2.24	1.43	1.39
3	DDD	403	NAI	C4N-C5N	-2.21	1.43	1.49
3	BBB	901	NAI	PN-O3	2.06	1.61	1.59
3	CCC	402	NAI	C4A-N3A	2.04	1.38	1.34
3	AAA	503	NAI	PA-O5B	-2.03	1.51	1.59

All (58) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	CCC	402	NAI	C5A-C4A-N3A	-6.60	117.64	126.72
3	DDD	403	NAI	C4A-N9A-C8A	6.39	112.45	105.74
3	CCC	402	NAI	N3A-C4A-N9A	6.24	137.79	127.17
3	BBB	901	NAI	C4A-C5A-N7A	-5.99	103.74	110.58
3	DDD	403	NAI	N3A-C2A-N1A	-5.79	119.82	128.58
3	BBB	901	NAI	C5A-C4A-N3A	-5.55	119.08	126.72
3	DDD	403	NAI	C6A-C5A-C4A	-5.55	109.60	117.18
3	BBB	901	NAI	N3A-C2A-N1A	-5.34	120.50	128.58
3	BBB	901	NAI	C6A-C5A-N7A	5.02	141.76	132.09
3	AAA	503	NAI	N3A-C2A-N1A	-4.95	121.10	128.58
3	DDD	403	NAI	C5A-C4A-N9A	-4.71	100.68	105.81
3	BBB	901	NAI	C2A-N3A-C4A	4.68	123.25	111.83
3	CCC	402	NAI	C4A-N9A-C8A	4.60	110.57	105.74
3	CCC	402	NAI	C2A-N3A-C4A	4.58	123.00	111.83
3	DDD	403	NAI	C6A-C5A-N7A	4.13	140.04	132.09
3	CCC	402	NAI	N3A-C2A-N1A	-4.08	122.40	128.58
3	CCC	402	NAI	O2A-PA-O3	3.94	117.93	107.27
3	BBB	901	NAI	C5A-N7A-C8A	3.83	109.47	103.45
3	BBB	901	NAI	N3A-C4A-N9A	3.71	133.47	127.17
3	DDD	403	NAI	C4A-N9A-C1B	-3.58	118.27	126.63
3	DDD	403	NAI	N3A-C4A-N9A	3.53	133.17	127.17
3	DDD	403	NAI	N9A-C8A-N7A	-3.42	109.08	113.94
3	AAA	503	NAI	C6A-C5A-N7A	3.32	138.50	132.09
3	BBB	901	NAI	C4A-N9A-C8A	3.25	109.15	105.74
3	AAA	503	NAI	C6N-N1N-C2N	3.02	122.55	119.32
3	DDD	403	NAI	C2A-N3A-C4A	2.93	118.98	111.83
3	DDD	403	NAI	O4B-C1B-C2B	-2.90	100.41	106.62
3	CCC	402	NAI	C4A-N9A-C1B	-2.88	119.89	126.63
3	DDD	403	NAI	O1N-PN-O2N	2.87	125.79	112.44
3	CCC	402	NAI	C6A-C5A-N7A	2.84	137.56	132.09
3	CCC	402	NAI	C1D-N1N-C6N	-2.71	115.04	120.77
3	CCC	402	NAI	C6N-N1N-C2N	2.68	122.19	119.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	BBB	901	NAI	N9A-C8A-N7A	-2.65	110.17	113.94
3	BBB	901	NAI	O1N-PN-O2N	2.61	124.56	112.44
3	AAA	503	NAI	C6A-C5A-C4A	-2.60	113.62	117.18
3	DDD	403	NAI	C5A-C6A-N1A	2.57	124.03	117.51
3	AAA	503	NAI	C3D-C2D-C1D	2.54	106.27	101.46
3	AAA	503	NAI	C4A-N9A-C8A	2.54	108.40	105.74
3	CCC	402	NAI	C5A-C6A-N1A	2.53	123.94	117.51
3	AAA	503	NAI	C4A-C5A-N7A	-2.53	107.69	110.58
3	BBB	901	NAI	O3-PA-O1A	2.52	118.28	110.70
3	CCC	402	NAI	N9A-C8A-N7A	-2.49	110.41	113.94
3	AAA	503	NAI	N9A-C8A-N7A	-2.45	110.47	113.94
3	DDD	403	NAI	C5A-C6A-N6A	-2.41	117.33	123.29
3	AAA	503	NAI	C2A-N3A-C4A	2.41	117.71	111.83
3	CCC	402	NAI	C4A-C5A-N7A	-2.33	107.92	110.58
3	AAA	503	NAI	C2D-C3D-C4D	-2.32	98.13	102.61
3	CCC	402	NAI	C5A-C6A-N6A	-2.28	117.65	123.29
3	AAA	503	NAI	O3-PA-O1A	-2.20	104.09	110.70
3	CCC	402	NAI	O4D-C4D-C3D	-2.20	100.79	105.15
3	CCC	402	NAI	O4B-C1B-N9A	2.19	112.29	108.09
3	AAA	503	NAI	C5A-N7A-C8A	2.17	106.86	103.45
3	DDD	403	NAI	O3B-C3B-C4B	2.16	117.30	111.08
3	CCC	402	NAI	O3-PA-O1A	-2.15	104.23	110.70
3	DDD	403	NAI	O4D-C4D-C3D	-2.11	100.97	105.15
3	AAA	503	NAI	O1N-PN-O2N	2.10	122.23	112.44
3	DDD	403	NAI	C5A-N7A-C8A	2.06	106.69	103.45
3	AAA	503	NAI	O2A-PA-O3	2.06	112.83	107.27

There are no chirality outliers.

All (23) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	DDD	401	PEG	C1-C2-O2-C3
2	CCC	401	PEG	O1-C1-C2-O2
2	DDD	401	PEG	O1-C1-C2-O2
2	DDD	402	PEG	O2-C3-C4-O4
4	AAA	505	EDO	O1-C1-C2-O2
2	AAA	502	PEG	C4-C3-O2-C2
3	DDD	403	NAI	C2B-C1B-N9A-C8A
2	AAA	501	PEG	O1-C1-C2-O2
3	AAA	503	NAI	O4D-C1D-N1N-C2N
3	CCC	402	NAI	O4D-C1D-N1N-C2N
3	DDD	403	NAI	O4D-C1D-N1N-C2N

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Mol	Chain	Res	Type	Atoms
2	DDD	401	PEG	O2-C3-C4-O4
3	BBB	901	NAI	O4D-C1D-N1N-C2N
2	AAA	502	PEG	O2-C3-C4-O4
2	DDD	401	PEG	C4-C3-O2-C2
4	AAA	504	EDO	O1-C1-C2-O2
4	BBB	902	EDO	O1-C1-C2-O2
4	CCC	403	EDO	O1-C1-C2-O2
4	DDD	404	EDO	O1-C1-C2-O2
3	AAA	503	NAI	C2B-C1B-N9A-C8A
4	AAA	506	EDO	O1-C1-C2-O2
2	AAA	502	PEG	O1-C1-C2-O2
3	BBB	901	NAI	C2D-C1D-N1N-C2N

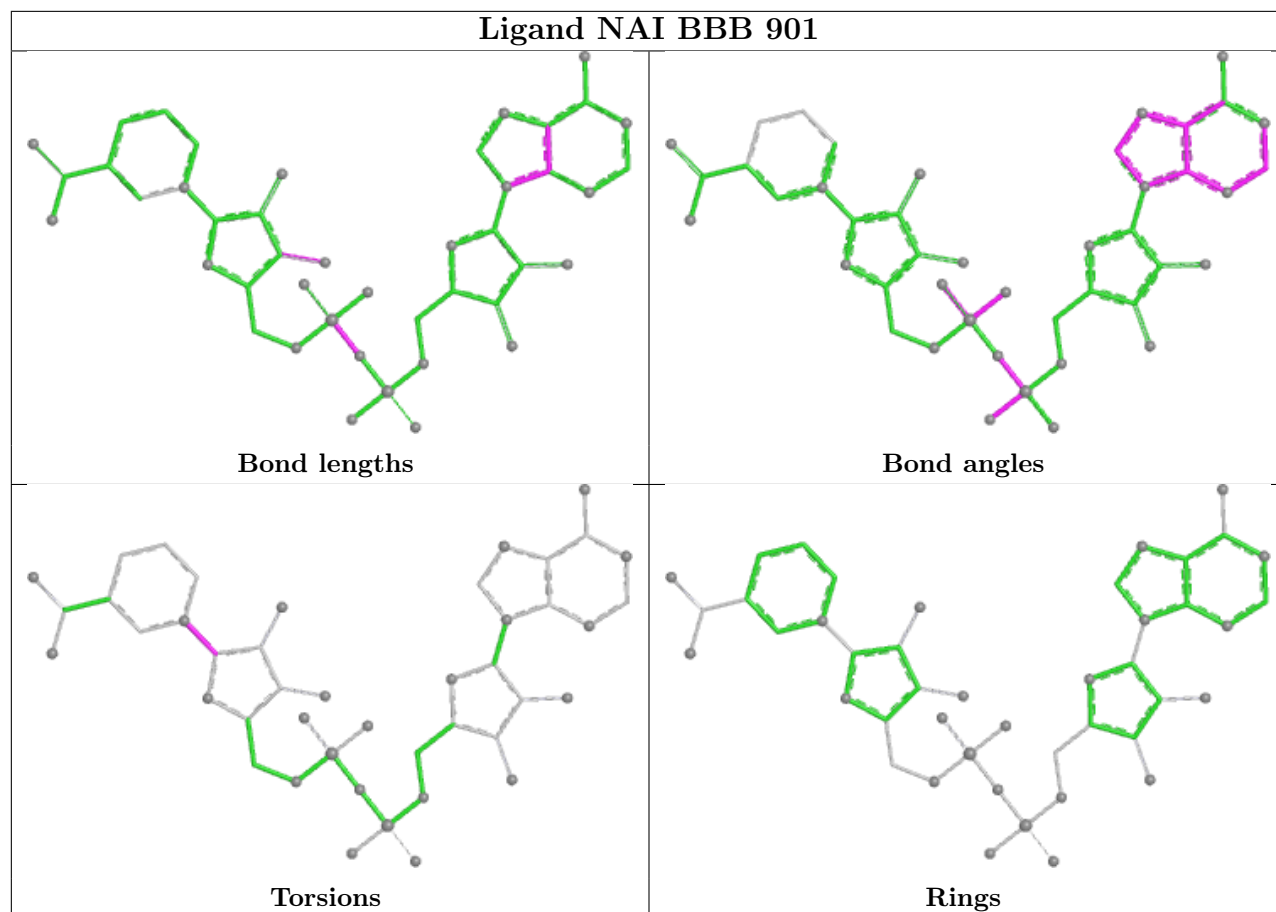
There are no ring outliers.

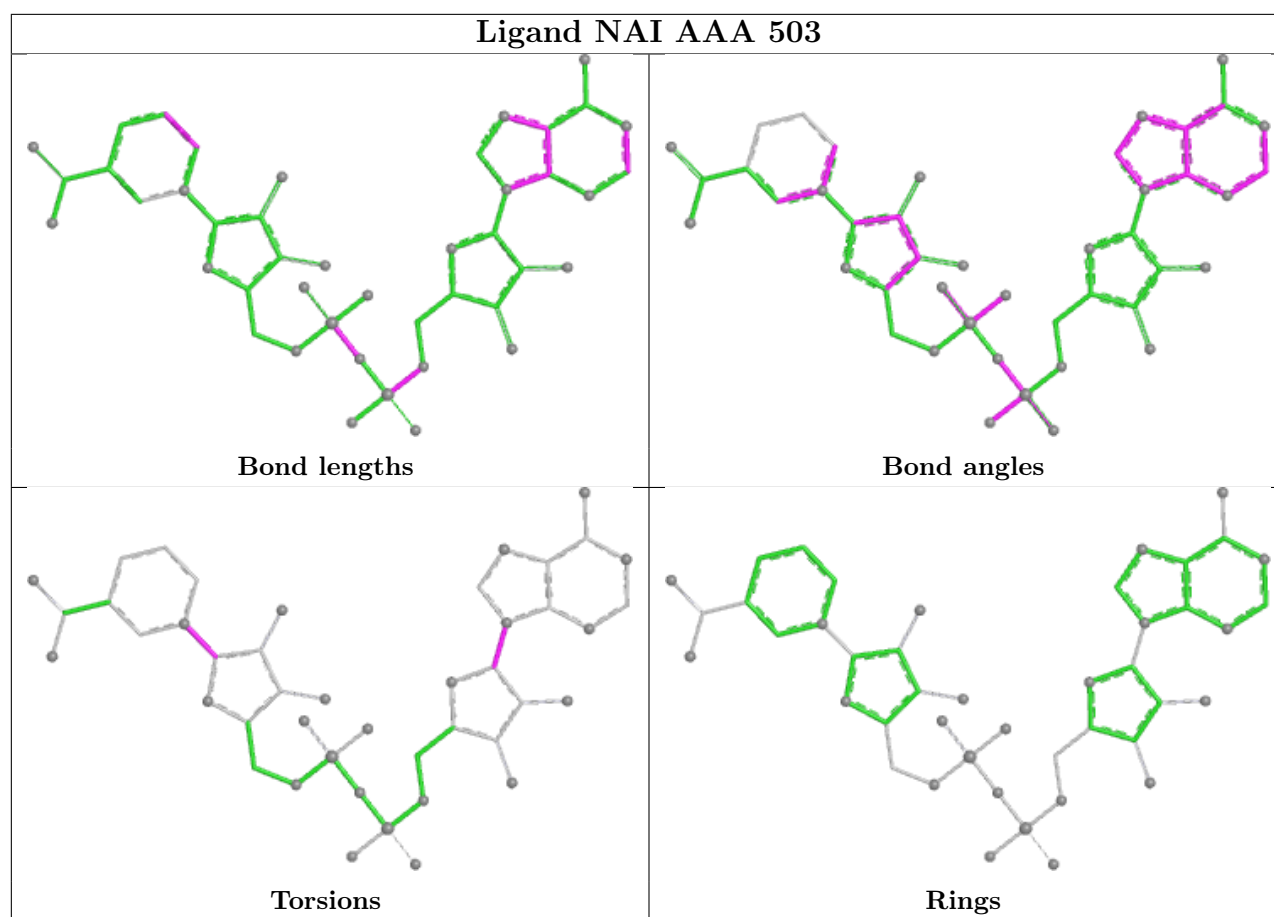
9 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	BBB	901	NAI	2	0
4	AAA	504	EDO	1	0
4	CCC	403	EDO	1	0
3	AAA	503	NAI	1	0
3	DDD	403	NAI	1	0
4	AAA	507	EDO	3	0
3	CCC	402	NAI	2	0
4	AAA	505	EDO	1	0
2	AAA	502	PEG	1	0

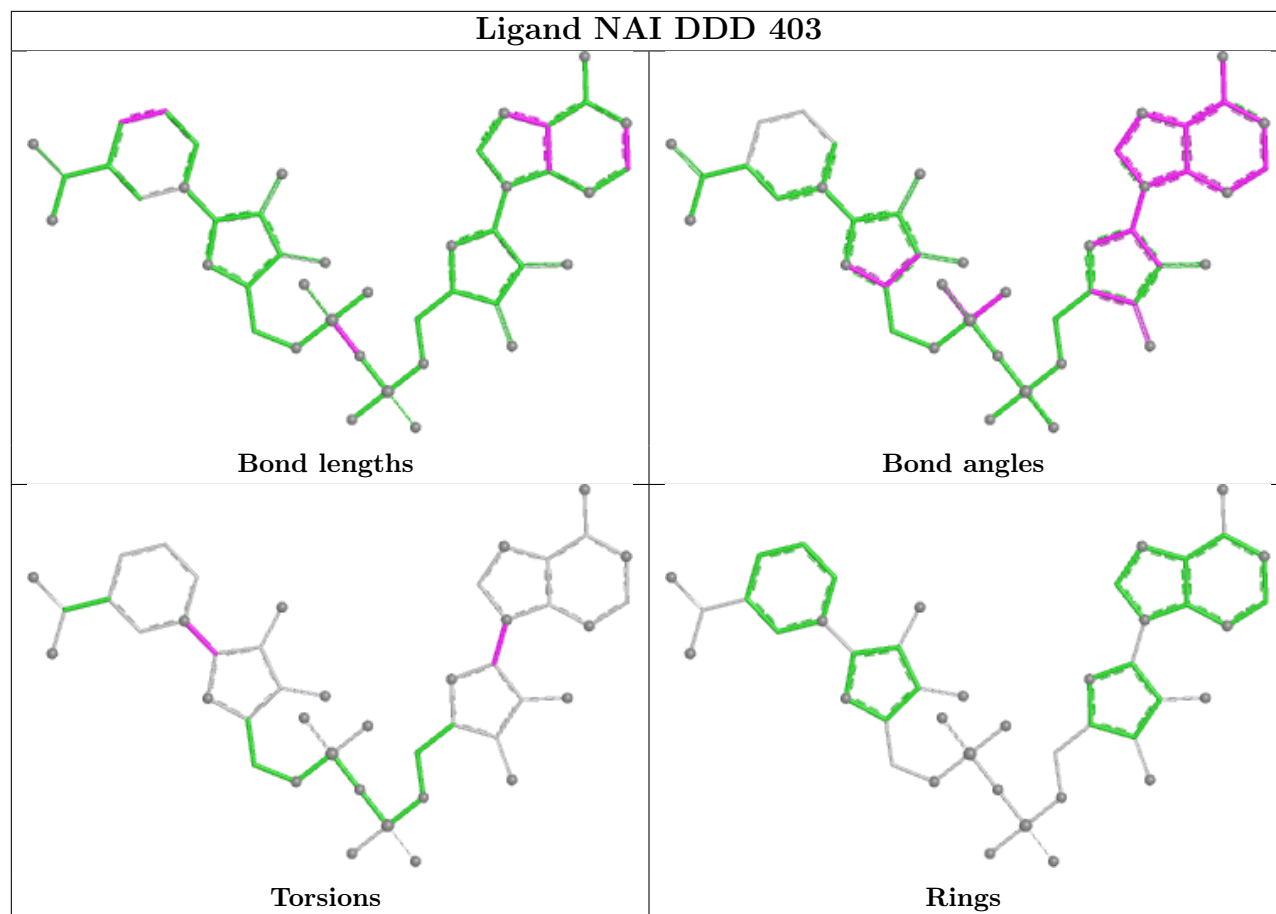
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

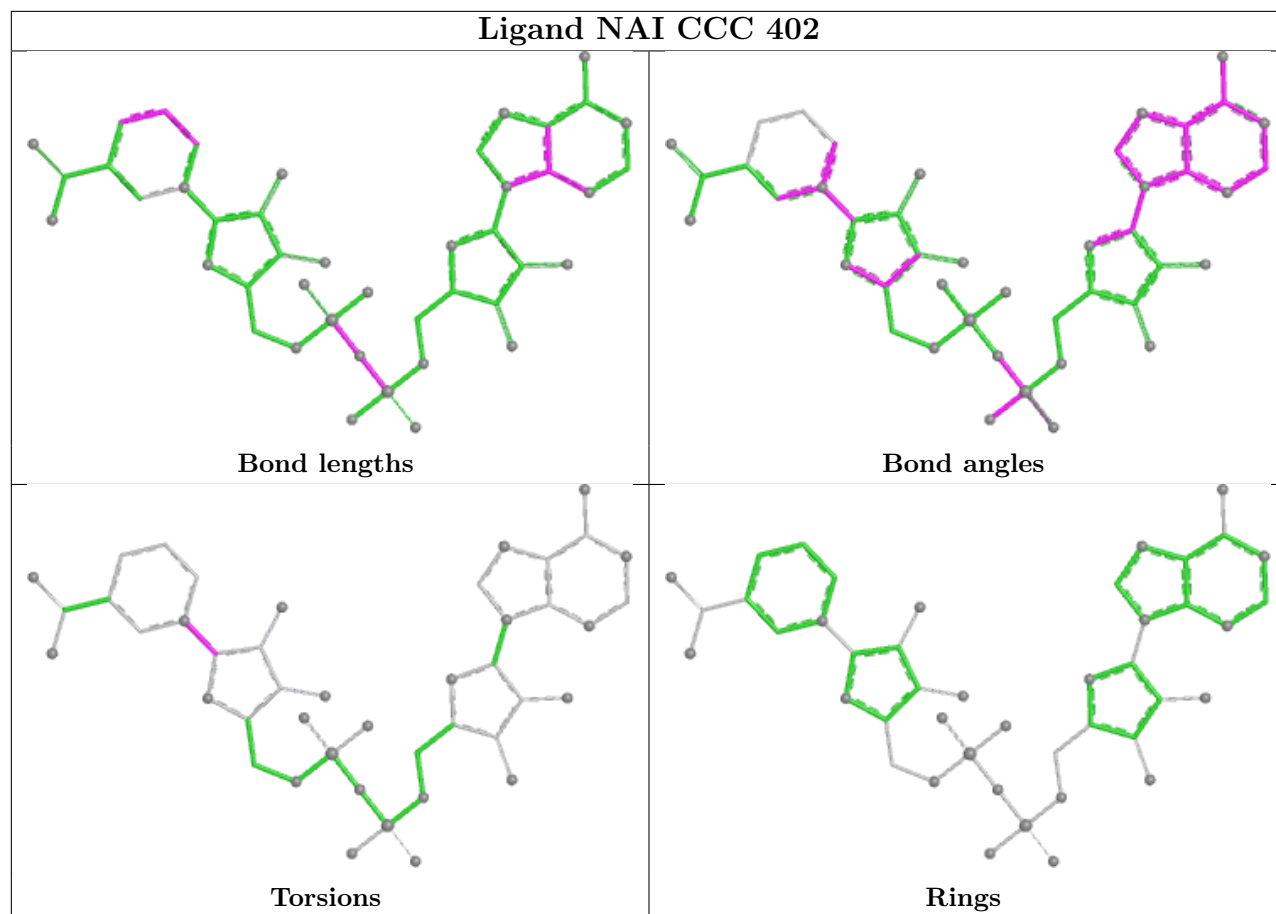
Ligand NAI BBB 901





Ligand NAI DDD 403





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	AAA	367/410 (89%)	0.39	9 (2%) 58 67	6, 16, 30, 83	33 (8%)
1	BBB	374/410 (91%)	0.70	36 (9%) 13 21	7, 18, 44, 113	34 (9%)
1	CCC	371/410 (90%)	0.96	73 (19%) 3 7	7, 18, 40, 129	31 (8%)
1	DDD	367/410 (89%)	0.48	21 (5%) 29 38	6, 15, 32, 73	28 (7%)
All	All	1479/1640 (90%)	0.63	139 (9%) 14 21	6, 17, 39, 129	126 (8%)

All (139) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	CCC	367	ILE	6.6
1	BBB	365	VAL	6.4
1	CCC	-1	LEU	6.3
1	CCC	83	ALA	5.6
1	CCC	365	VAL	5.4
1	BBB	-5	TYR	5.0
1	CCC	-3	ALA	4.9
1	AAA	-4	THR	4.9
1	CCC	336	LEU	4.6
1	CCC	81	ALA	4.5
1	CCC	1	MET	4.5
1	CCC	34	TRP	4.2
1	DDD	-4	THR	4.0
1	CCC	2	VAL	3.9
1	CCC	335	TYR	3.8
1	CCC	328	VAL	3.8
1	CCC	65	ILE	3.8
1	CCC	4	VAL	3.7
1	BBB	368	PRO	3.7
1	BBB	-4	THR	3.6
1	CCC	35	LEU	3.6

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Mol	Chain	Res	Type	RSRZ
1	BBB	86	LEU	3.6
1	BBB	43	VAL	3.5
1	BBB	52[A]	GLY	3.5
1	CCC	57[A]	ARG	3.5
1	DDD	-2	ARG	3.5
1	BBB	367	ILE	3.4
1	CCC	-2	ARG	3.4
1	CCC	366	LYS	3.3
1	CCC	89	ALA	3.3
1	CCC	86	LEU	3.2
1	CCC	3[A]	LYS	3.2
1	AAA	41	THR	3.1
1	CCC	111	ILE	3.1
1	CCC	352	GLY	3.1
1	CCC	103	ALA	3.1
1	DDD	-3	ALA	3.1
1	BBB	2	VAL	3.1
1	AAA	-3	ALA	3.1
1	BBB	89	ALA	3.1
1	CCC	41	THR	3.1
1	CCC	5	LEU	3.0
1	CCC	43	VAL	2.9
1	CCC	31	LEU	2.9
1	DDD	41	THR	2.9
1	CCC	77	ALA	2.9
1	CCC	55	PHE	2.9
1	BBB	336	LEU	2.8
1	BBB	364	ASP	2.8
1	CCC	16	LYS	2.8
1	CCC	104	ALA	2.8
1	CCC	52[A]	GLY	2.8
1	BBB	111	ILE	2.8
1	CCC	107	THR	2.8
1	CCC	364	ASP	2.7
1	BBB	15	ALA	2.7
1	DDD	16	LYS	2.7
1	DDD	57	ARG	2.7
1	CCC	42	LEU	2.7
1	CCC	7	VAL	2.7
1	CCC	105	ASN	2.7
1	DDD	15	ALA	2.7
1	DDD	83	ALA	2.7

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Mol	Chain	Res	Type	RSRZ
1	DDD	52[A]	GLY	2.6
1	CCC	88	LEU	2.6
1	AAA	52[A]	GLY	2.6
1	BBB	35	LEU	2.6
1	DDD	-1	LEU	2.6
1	CCC	340	PHE	2.6
1	BBB	-1	LEU	2.5
1	BBB	57[A]	ARG	2.5
1	CCC	106[A]	LYS	2.5
1	BBB	331	ILE	2.5
1	CCC	64	ILE	2.5
1	BBB	80	LEU	2.5
1	BBB	17	GLN	2.5
1	CCC	343[A]	ARG	2.5
1	CCC	348	ILE	2.4
1	BBB	12	GLY	2.4
1	CCC	85[A]	LYS	2.4
1	CCC	342	TYR	2.4
1	CCC	29	LEU	2.4
1	BBB	-3	ALA	2.4
1	CCC	331	ILE	2.4
1	CCC	18	VAL	2.4
1	BBB	39	GLY	2.3
1	CCC	354	TYR	2.3
1	CCC	84[A]	LYS	2.3
1	CCC	66	ILE	2.3
1	BBB	51[A]	GLU	2.3
1	AAA	18	VAL	2.3
1	AAA	43	VAL	2.3
1	BBB	113	VAL	2.3
1	CCC	12	GLY	2.3
1	CCC	30	GLY	2.3
1	CCC	109	GLY	2.3
1	BBB	354	TYR	2.3
1	DDD	12	GLY	2.3
1	AAA	15	ALA	2.3
1	CCC	101	LEU	2.3
1	CCC	91	THR	2.3
1	BBB	340	PHE	2.2
1	CCC	37	ASP	2.2
1	BBB	50	ARG	2.2
1	BBB	64	ILE	2.2

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Mol	Chain	Res	Type	RSRZ
1	CCC	90	VAL	2.2
1	CCC	108	ASN	2.2
1	AAA	17[A]	GLN	2.2
1	BBB	3[A]	LYS	2.2
1	CCC	15	ALA	2.2
1	CCC	112	THR	2.2
1	BBB	22	LEU	2.2
1	CCC	22	LEU	2.2
1	CCC	59	LEU	2.2
1	CCC	80	LEU	2.2
1	DDD	2	VAL	2.2
1	DDD	43	VAL	2.2
1	CCC	339	ARG	2.2
1	BBB	41	THR	2.1
1	BBB	83	ALA	2.1
1	DDD	51[A]	GLU	2.1
1	DDD	17[A]	GLN	2.1
1	DDD	337	SER	2.1
1	CCC	17	GLN	2.1
1	CCC	44	THR	2.1
1	DDD	62	ALA	2.1
1	DDD	1	MET	2.1
1	CCC	332	LEU	2.1
1	CCC	61	ASP	2.1
1	CCC	53	SER	2.0
1	DDD	75	LEU	2.0
1	AAA	3[A]	LYS	2.0
1	BBB	16	LYS	2.0
1	CCC	99	VAL	2.0
1	DDD	4	VAL	2.0
1	BBB	34	TRP	2.0
1	DDD	84[A]	LYS	2.0
1	CCC	39	GLY	2.0
1	BBB	42	LEU	2.0

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

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

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

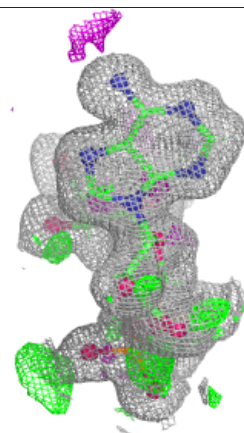
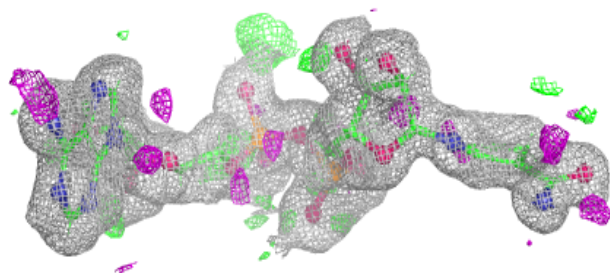
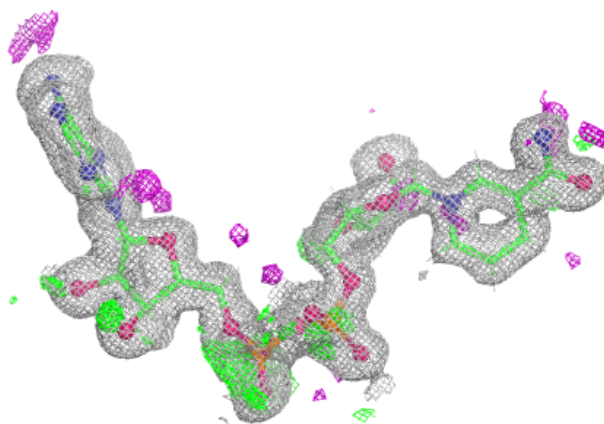
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	EDO	DDD	404	4/4	0.75	0.17	49,53,66,66	1
4	EDO	AAA	505	4/4	0.76	0.17	49,51,60,60	1
4	EDO	AAA	507	4/4	0.82	0.14	31,39,52,52	1
2	PEG	CCC	401	7/7	0.83	0.14	27,31,48,48	1
2	PEG	DDD	401	7/7	0.85	0.14	20,24,47,47	1
4	EDO	AAA	506	4/4	0.85	0.18	38,39,41,41	1
2	PEG	AAA	501	7/7	0.86	0.17	25,27,80,80	1
2	PEG	AAA	502	7/7	0.86	0.13	33,40,70,70	1
4	EDO	DDD	405	4/4	0.87	0.14	30,32,37,37	1
2	PEG	DDD	402	7/7	0.88	0.15	40,59,84,84	1
4	EDO	CCC	403	4/4	0.90	0.14	25,37,55,55	1
4	EDO	BBB	902	4/4	0.92	0.08	25,26,37,37	1
4	EDO	AAA	504	4/4	0.94	0.10	24,33,57,57	1
3	NAI	CCC	402	44/44	0.97	0.06	12,15,18,19	4
3	NAI	BBB	901	44/44	0.98	0.05	13,16,20,20	4
3	NAI	AAA	503	44/44	0.99	0.04	10,13,15,16	4
3	NAI	DDD	403	44/44	0.99	0.04	10,12,14,17	4

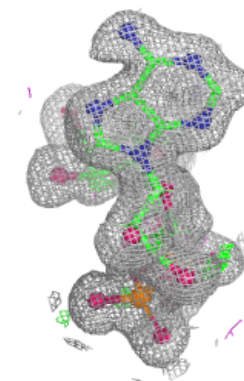
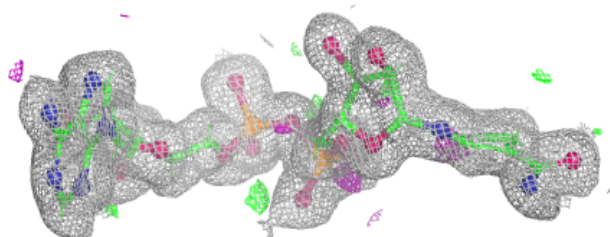
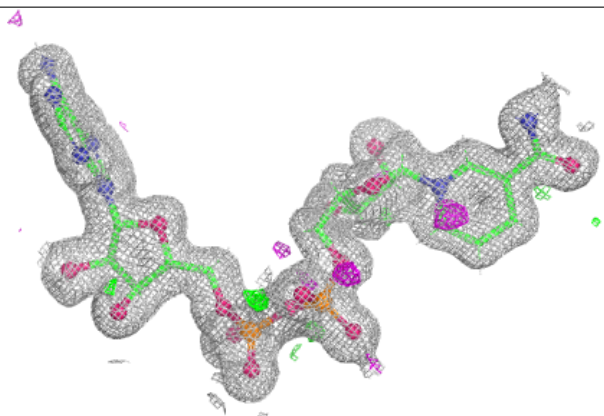
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around NAI CCC 402:

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

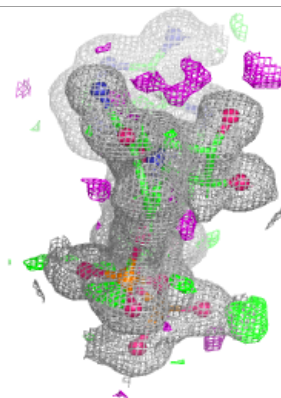
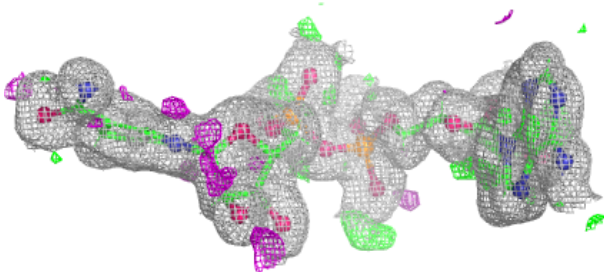
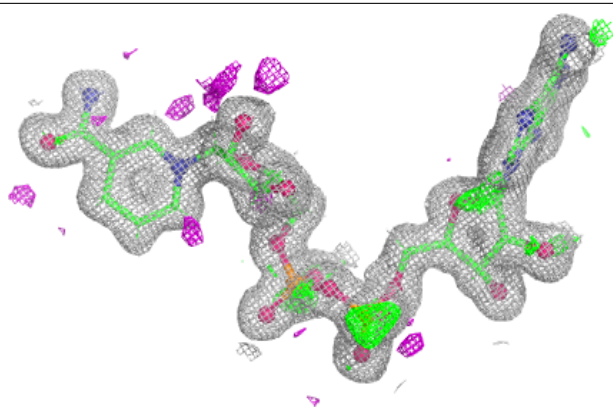
**Electron density around NAI BBB 901:**

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

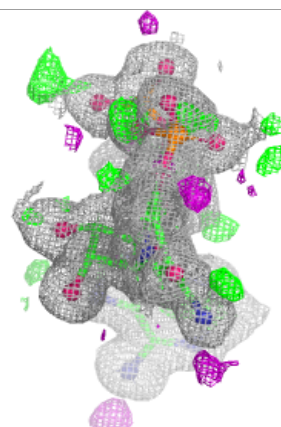
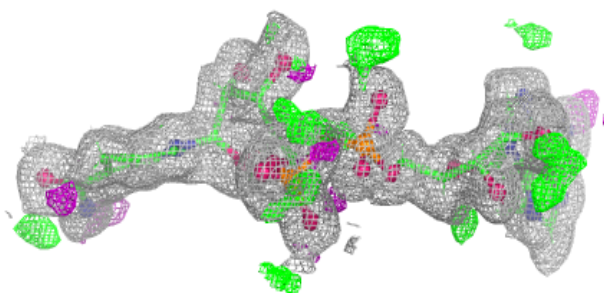
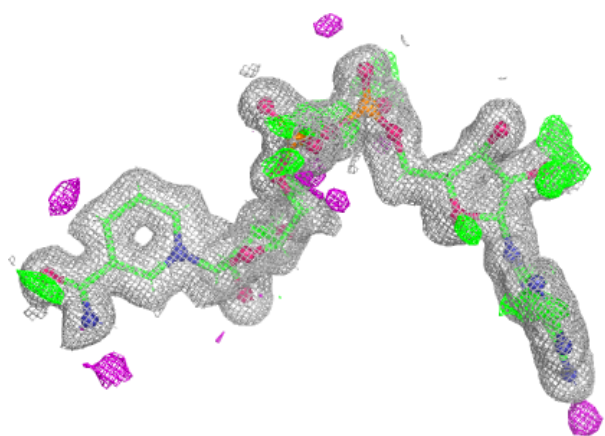


Electron density around NAI AAA 503:

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

**Electron density around NAI DDD 403:**

$2mF_o-DF_c$ (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.