



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 10, 2021 – 10:16 AM JST

PDB ID : 7DBK
Title : Crystal structure of human LDHB in complex with NADH
Authors : Sogabe, S.; Miwa, M.
Deposited on : 2020-10-20
Resolution : 1.80 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

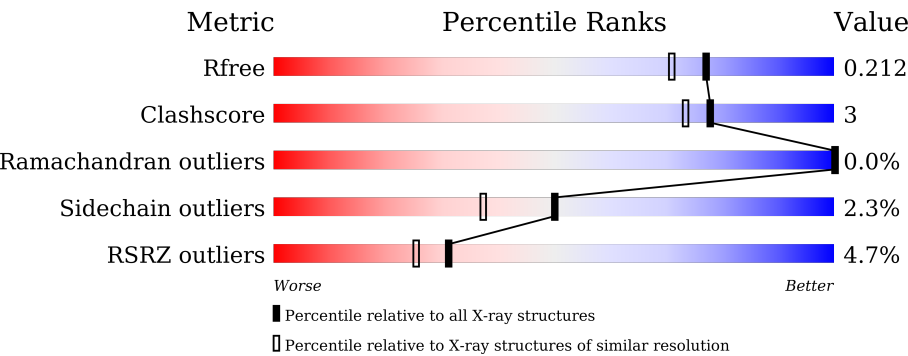
MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.23.2
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.23.2

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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.



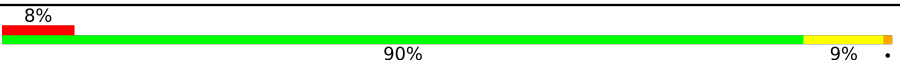

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (1.80-1.80)

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

Mol	Chain	Length	Quality of chain
1	A	333	<div><div>2%</div><div>96%</div><div>.</div></div>
1	B	333	<div><div>3%</div><div>92%</div><div>7%</div><div>.</div></div>
1	C	333	<div><div>7%</div><div>89%</div><div>11%</div></div>
1	D	333	<div><div>6%</div><div>92%</div><div>7%</div><div>.</div></div>
1	E	333	<div><div>%</div><div>92%</div><div>8%</div></div>
1	F	333	<div><div>5%</div><div>89%</div><div>11%</div></div>

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Mol	Chain	Length	Quality of chain
1	G	333	
1	H	333	

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
3	GOL	D	501	-	-	X	-
3	GOL	D	502	-	-	X	-
3	GOL	E	404	-	-	-	X
3	GOL	E	405	-	-	-	X

2 Entry composition

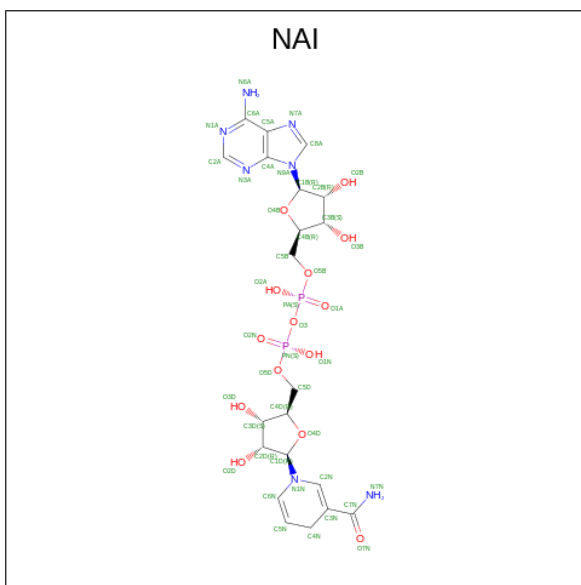
There are 4 unique types of molecules in this entry. The entry contains 21949 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 L-lactate dehydrogenase B chain.

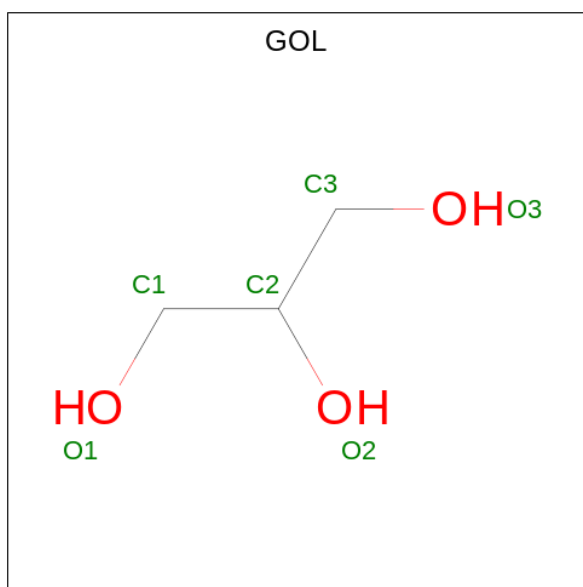
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	333	Total	C	N	O	S	0	0	0
			2561	1628	431	488	14			
1	B	333	Total	C	N	O	S	0	0	0
			2561	1628	431	488	14			
1	C	333	Total	C	N	O	S	0	0	0
			2561	1628	431	488	14			
1	D	333	Total	C	N	O	S	0	0	0
			2561	1628	431	488	14			
1	E	333	Total	C	N	O	S	0	0	0
			2561	1628	431	488	14			
1	F	333	Total	C	N	O	S	0	0	0
			2561	1628	431	488	14			
1	G	333	Total	C	N	O	S	0	0	0
			2561	1628	431	488	14			
1	H	333	Total	C	N	O	S	0	0	0
			2561	1628	431	488	14			

- Molecule 2 is 1,4-DIHYDRONICOTINAMIDE ADENINE DINUCLEOTIDE (three-letter code: NAI) (formula: C₂₁H₂₉N₇O₁₄P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 44	C 21	N 7	O 14	P 2	0	0
2	B	1	Total 44	C 21	N 7	O 14	P 2	0	0
2	C	1	Total 44	C 21	N 7	O 14	P 2	0	0
2	D	1	Total 44	C 21	N 7	O 14	P 2	0	0
2	E	1	Total 44	C 21	N 7	O 14	P 2	0	0
2	F	1	Total 44	C 21	N 7	O 14	P 2	0	0
2	G	1	Total 44	C 21	N 7	O 14	P 2	0	0
2	H	1	Total 44	C 21	N 7	O 14	P 2	0	0

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: $\text{C}_3\text{H}_8\text{O}_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	C	1	Total	C	O	0	0
			6	3	3		
3	D	1	Total	C	O	0	0
			6	3	3		
3	D	1	Total	C	O	0	0
			6	3	3		
3	D	1	Total	C	O	0	0
			6	3	3		
3	D	1	Total	C	O	0	0
			6	3	3		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	E	1	Total	C	O	0	0
			6	3	3		
3	E	1	Total	C	O	0	0
			6	3	3		
3	E	1	Total	C	O	0	0
			6	3	3		
3	E	1	Total	C	O	0	0
			6	3	3		
3	F	1	Total	C	O	0	0
			6	3	3		
3	G	1	Total	C	O	0	0
			6	3	3		
3	H	1	Total	C	O	0	0
			6	3	3		
3	H	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	166	Total	O	0	0
			166	166		
4	B	144	Total	O	0	0
			144	144		
4	C	105	Total	O	0	0
			105	105		
4	D	107	Total	O	0	0
			107	107		
4	E	149	Total	O	0	0
			149	149		
4	F	119	Total	O	0	0
			119	119		
4	G	76	Total	O	0	0
			76	76		
4	H	111	Total	O	0	0
			111	111		

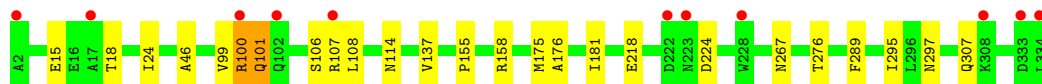
3 Residue-property plots [i](#)

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.

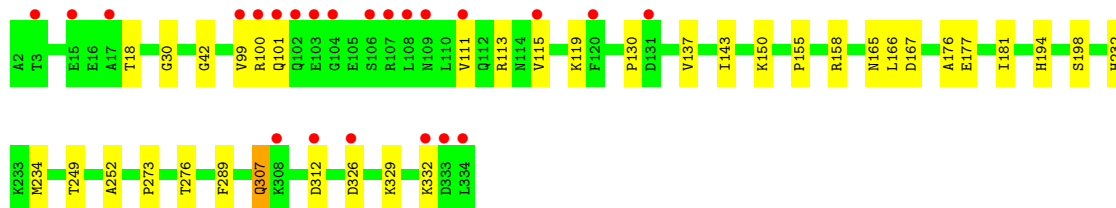
- Molecule 1: L-lactate dehydrogenase B chain



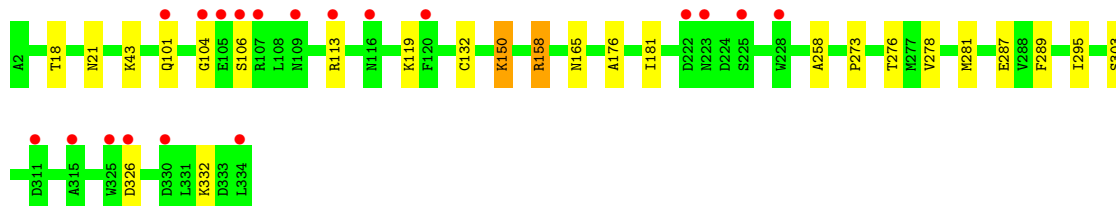
- Molecule 1: L-lactate dehydrogenase B chain



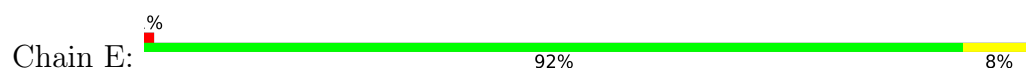
- Molecule 1: L-lactate dehydrogenase B chain



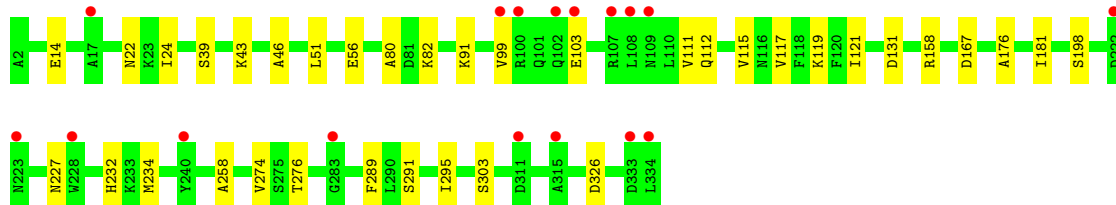
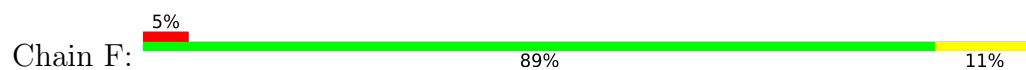
- Molecule 1: L-lactate dehydrogenase B chain



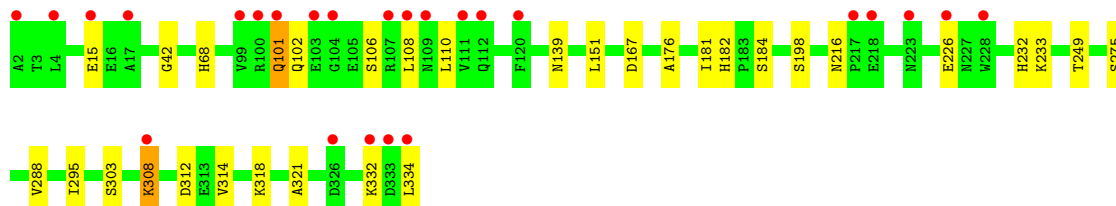
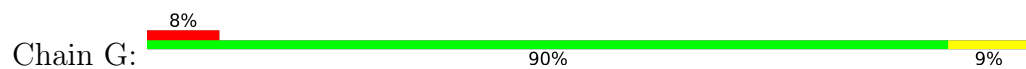
- Molecule 1: L-lactate dehydrogenase B chain



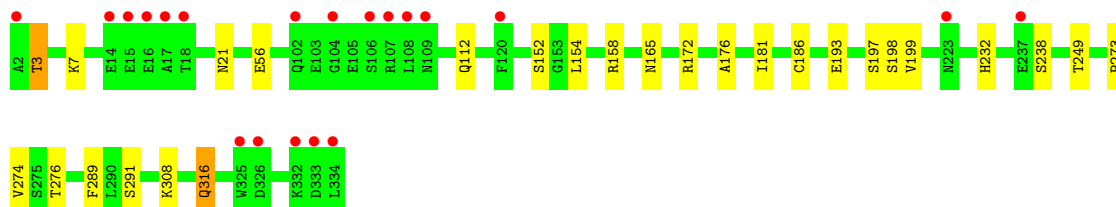
- Molecule 1: L-lactate dehydrogenase B chain



- Molecule 1: L-lactate dehydrogenase B chain



- Molecule 1: L-lactate dehydrogenase B chain



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	232.48Å 84.17Å 156.21Å 90.00° 120.76° 90.00°	Depositor
Resolution (Å)	49.99 – 1.80 49.94 – 1.80	Depositor EDS
% Data completeness (in resolution range)	99.5 (49.99-1.80) 99.5 (49.94-1.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.05 (at 1.81Å)	Xtriage
Refinement program	REFMAC 5.8.0266	Depositor
R, R_{free}	0.170 , 0.203 0.180 , 0.212	Depositor DCC
R_{free} test set	11844 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	23.2	Xtriage
Anisotropy	0.193	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 46.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	21949	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.39% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: NAI, GOL

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	0.79	0/2602	0.82	3/3526 (0.1%)
1	B	0.76	0/2602	0.81	0/3526
1	C	0.77	0/2602	0.80	0/3526
1	D	0.78	0/2602	0.84	3/3526 (0.1%)
1	E	0.77	1/2602 (0.0%)	0.79	1/3526 (0.0%)
1	F	0.76	0/2602	0.83	2/3526 (0.1%)
1	G	0.74	0/2602	0.80	0/3526
1	H	0.76	1/2602 (0.0%)	0.81	2/3526 (0.1%)
All	All	0.77	2/20816 (0.0%)	0.81	11/28208 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	H	56	GLU	CD-OE2	-5.23	1.19	1.25
1	E	262	GLU	CD-OE2	-5.11	1.20	1.25

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	158	ARG	NE-CZ-NH2	-9.22	115.69	120.30
1	F	158	ARG	NE-CZ-NH2	-8.36	116.12	120.30
1	D	158	ARG	NE-CZ-NH1	8.33	124.46	120.30
1	F	158	ARG	NE-CZ-NH1	7.68	124.14	120.30
1	A	158	ARG	NE-CZ-NH1	6.42	123.51	120.30
1	H	158	ARG	NE-CZ-NH1	6.37	123.49	120.30
1	H	158	ARG	NE-CZ-NH2	-6.10	117.25	120.30
1	E	158	ARG	NE-CZ-NH1	-5.43	117.58	120.30
1	D	18	THR	CA-CB-OG1	-5.42	97.63	109.00
1	A	158	ARG	NE-CZ-NH2	-5.09	117.75	120.30
1	A	100	ARG	NE-CZ-NH1	5.03	122.82	120.30

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	2561	0	2630	10	0
1	B	2561	0	2630	14	0
1	C	2561	0	2630	20	0
1	D	2561	0	2630	13	0
1	E	2561	0	2630	20	0
1	F	2561	0	2630	16	0
1	G	2561	0	2630	14	0
1	H	2561	0	2630	18	0
2	A	44	0	27	0	0
2	B	44	0	27	1	0
2	C	44	0	27	2	0
2	D	44	0	27	0	0
2	E	44	0	27	0	0
2	F	44	0	27	0	0
2	G	44	0	27	1	0
2	H	44	0	27	0	0
3	A	18	0	20	0	0
3	B	36	0	45	1	0
3	C	6	0	8	0	0
3	D	24	0	26	6	0
3	E	24	0	30	4	0
3	F	6	0	8	0	0
3	G	6	0	5	1	0
3	H	12	0	12	1	0
4	A	166	0	0	1	0
4	B	144	0	0	0	0
4	C	105	0	0	4	0
4	D	107	0	0	0	0
4	E	149	0	0	1	0
4	F	119	0	0	2	0
4	G	76	0	0	0	0
4	H	111	0	0	2	0
All	All	21949	0	21410	124	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 3.

All (124) 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)
3:D:501:GOL:O2	3:D:502:GOL:O2	1.74	1.04
1:A:21:ASN:HD21	1:D:21:ASN:HD21	1.15	0.89
1:C:42:GLY:HA3	3:D:502:GOL:H2	1.51	0.89
3:D:501:GOL:HO2	3:D:502:GOL:HO2	1.14	0.82
1:B:100:ARG:HH11	1:B:100:ARG:HG2	1.43	0.81
1:D:150:LYS:NZ	1:D:287:GLU:OE1	2.16	0.78
1:A:21:ASN:HD21	1:D:21:ASN:ND2	1.82	0.77
1:G:42:GLY:HA3	3:G:402:GOL:H2	1.71	0.71
1:E:99:VAL:H	1:E:114:ASN:HD21	1.41	0.69
1:H:165:ASN:HD22	1:H:273:PRO:HD2	1.57	0.68
1:A:21:ASN:ND2	1:D:21:ASN:HD21	1.90	0.67
1:G:295:ILE:HD12	1:G:303:SER:HB2	1.78	0.66
3:D:501:GOL:O2	3:D:502:GOL:C2	2.36	0.66
1:B:99:VAL:H	1:B:114:ASN:HD21	1.43	0.65
1:E:233:LYS:HE3	1:E:237:GLU:OE2	1.98	0.63
1:B:100:ARG:HG2	1:B:100:ARG:NH1	2.14	0.62
1:E:139:ASN:HD22	1:E:141:VAL:H	1.47	0.62
1:D:278:VAL:HA	1:D:281:MET:HE3	1.82	0.62
1:E:176:ALA:HB1	1:E:181:ILE:O	2.01	0.60
1:B:295:ILE:HD12	1:B:295:ILE:N	2.17	0.59
1:C:130:PRO:O	1:C:158:ARG:NH2	2.32	0.59
1:H:193:GLU:OE2	1:H:197:SER:OG	2.20	0.59
1:C:198:SER:OG	1:C:232:HIS:HE1	1.86	0.57
1:G:198:SER:OG	1:G:232:HIS:HE1	1.88	0.57
1:H:181:ILE:HG13	1:H:186:CYS:SG	2.44	0.57
1:G:216:ASN:ND2	1:G:226:GLU:OE1	2.35	0.56
1:H:21:ASN:ND2	4:H:503:HOH:O	2.39	0.55
1:A:295:ILE:HD12	1:A:303:SER:HB2	1.88	0.55
1:E:206:ASN:HB3	3:E:403:GOL:H32	1.88	0.55
1:G:176:ALA:HB1	1:G:181:ILE:O	2.06	0.55
1:E:198:SER:OG	1:E:232:HIS:HE1	1.90	0.54
1:C:234:MET:HG2	4:C:598:HOH:O	2.07	0.54
1:C:42:GLY:HA3	3:D:502:GOL:C2	2.31	0.53
1:G:182:HIS:HD2	1:G:184:SER:H	1.55	0.53
1:F:232:HIS:HD2	4:F:615:HOH:O	1.92	0.53
1:H:172:ARG:HH12	3:H:403:GOL:H31	1.74	0.53
1:D:132:CYS:O	1:D:158:ARG:HD2	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:501:GOL:HO2	3:D:502:GOL:C2	2.15	0.52
1:G:182:HIS:CD2	1:G:184:SER:H	2.27	0.52
1:E:175:MET:SD	1:E:186:CYS:HB3	2.49	0.52
1:E:215:LEU:HD23	1:G:308:LYS:HE3	1.92	0.51
1:F:295:ILE:HD12	1:F:303:SER:HB2	1.93	0.51
1:D:176:ALA:HB1	1:D:181:ILE:O	2.11	0.51
1:E:165:ASN:HD22	1:E:273:PRO:HD2	1.76	0.51
1:F:56:GLU:HG3	1:F:82:LYS:HG3	1.91	0.51
1:D:278:VAL:HA	1:D:281:MET:CE	2.41	0.50
1:C:176:ALA:HB1	1:C:181:ILE:O	2.11	0.50
1:E:206:ASN:HB3	3:E:403:GOL:C3	2.41	0.50
1:E:171:PHE:CE2	1:E:175:MET:HE2	2.46	0.50
1:E:181:ILE:HD11	1:E:186:CYS:SG	2.51	0.50
1:F:22:ASN:HA	1:F:91:LYS:HG3	1.93	0.50
1:G:288:VAL:HG11	1:G:321:ALA:HA	1.95	0.49
1:A:198:SER:OG	1:A:232:HIS:HE1	1.94	0.49
1:H:198:SER:OG	1:H:232:HIS:HE1	1.95	0.49
1:C:111:VAL:HG11	1:C:143:ILE:HG21	1.95	0.49
1:C:155:PRO:HG2	1:C:158:ARG:HG3	1.95	0.49
1:F:176:ALA:HB1	1:F:181:ILE:O	2.12	0.49
1:H:199:VAL:CG2	1:H:316:GLN:HG2	2.43	0.49
1:H:232:HIS:HD2	4:H:611:HOH:O	1.96	0.48
1:H:176:ALA:HB1	1:H:181:ILE:O	2.14	0.48
1:A:165:ASN:HD22	1:A:273:PRO:HD2	1.78	0.47
1:C:115:VAL:O	1:C:119:LYS:HB2	2.14	0.47
1:A:210:VAL:HG22	1:C:307:GLN:HG3	1.96	0.47
1:F:234:MET:HG2	4:F:605:HOH:O	2.13	0.47
1:B:18:THR:HG21	4:C:596:HOH:O	2.14	0.47
1:C:326:ASP:HA	1:C:329:LYS:HE3	1.96	0.47
1:D:101:GLN:HE21	1:D:104:GLY:HA2	1.80	0.47
1:B:295:ILE:N	1:B:295:ILE:CD1	2.78	0.47
1:F:111:VAL:O	1:F:115:VAL:HG23	2.14	0.47
1:A:232:HIS:HD2	4:A:660:HOH:O	1.97	0.47
1:E:330:ASP:OD2	4:E:501:HOH:O	2.20	0.46
1:H:165:ASN:HA	1:H:273:PRO:HG2	1.97	0.46
1:C:158:ARG:HD2	4:C:584:HOH:O	2.14	0.46
1:F:198:SER:OG	1:F:232:HIS:HE1	1.99	0.45
1:B:137:VAL:O	2:B:401:NAI:H2N	2.16	0.45
1:C:150:LYS:HA	1:C:150:LYS:HD2	1.77	0.45
1:B:218:GLU:O	1:B:224:ASP:HB2	2.17	0.45
1:A:42:GLY:HA3	3:B:405:GOL:H2	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:165:ASN:HD22	1:C:273:PRO:HD2	1.82	0.45
1:C:137:VAL:O	2:C:401:NAI:H2N	2.17	0.45
1:H:3:THR:O	1:H:7:LYS:HG3	2.17	0.44
1:G:139:ASN:HB2	2:G:401:NAI:O2D	2.17	0.44
1:H:181:ILE:CG1	1:H:186:CYS:SG	3.06	0.44
1:F:276:THR:O	1:F:289:PHE:HA	2.17	0.43
1:G:151:LEU:HD21	1:G:334:LEU:HA	2.00	0.43
1:B:24:ILE:HD12	1:B:46:ALA:HB2	2.01	0.43
1:C:276:THR:O	1:C:289:PHE:HA	2.18	0.43
1:H:181:ILE:HD11	1:H:186:CYS:SG	2.58	0.43
1:B:276:THR:O	1:B:289:PHE:HA	2.18	0.43
1:E:114:ASN:HD22	1:E:114:ASN:HA	1.53	0.43
1:A:21:ASN:O	1:A:91:LYS:CE	2.66	0.43
1:B:101:GLN:H	1:B:101:GLN:HG2	1.70	0.43
3:E:404:GOL:O2	3:E:405:GOL:O2	2.13	0.43
1:G:68:HIS:CE1	1:H:238:SER:HB2	2.53	0.43
1:H:199:VAL:HG22	1:H:316:GLN:HG2	2.00	0.43
1:E:139:ASN:HD22	1:E:141:VAL:N	2.15	0.43
1:F:115:VAL:HG12	1:F:119:LYS:HE3	2.01	0.43
1:D:165:ASN:HA	1:D:273:PRO:HG2	2.01	0.42
1:G:101:GLN:HA	1:G:110:LEU:HD13	2.01	0.42
1:D:43:LYS:HD3	1:D:258:ALA:HB1	2.01	0.42
1:F:274:VAL:O	1:F:291:SER:HA	2.20	0.42
1:B:176:ALA:HB1	1:B:181:ILE:O	2.20	0.41
1:F:117:VAL:HG12	1:F:121:ILE:HD12	2.01	0.41
1:B:267:ASN:OD1	1:B:297:ASN:HB2	2.20	0.41
1:H:274:VAL:O	1:H:291:SER:HA	2.21	0.41
1:D:295:ILE:HD12	1:D:303:SER:HB2	2.02	0.41
1:H:152:SER:HB2	1:H:154:LEU:HG	2.02	0.41
1:E:154:LEU:HD22	1:E:158:ARG:NH1	2.35	0.41
1:C:167:ASP:OD1	1:C:194:HIS:ND1	2.49	0.41
1:F:43:LYS:HD3	1:F:258:ALA:HB1	2.02	0.41
1:B:155:PRO:HG2	1:B:158:ARG:HD3	2.02	0.41
1:C:232:HIS:HD2	4:C:604:HOH:O	2.04	0.41
1:D:276:THR:O	1:D:289:PHE:HA	2.21	0.41
1:E:51:LEU:O	1:E:80:ALA:HA	2.21	0.41
1:E:171:PHE:CE2	1:E:175:MET:CE	3.04	0.41
3:E:404:GOL:H32	1:F:39:SER:HA	2.03	0.41
1:F:24:ILE:HD12	1:F:46:ALA:HB2	2.02	0.41
1:E:218:GLU:O	1:E:224:ASP:HB2	2.21	0.41
1:G:314:VAL:HG12	1:G:318:LYS:HE3	2.02	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:166:LEU:HD11	1:C:252:ALA:HB1	2.03	0.40
1:E:139:ASN:HA	1:E:141:VAL:N	2.36	0.40
1:F:51:LEU:O	1:F:80:ALA:HA	2.21	0.40
1:H:276:THR:O	1:H:289:PHE:HA	2.22	0.40
1:C:30:GLY:HA3	2:C:401:NAI:O5B	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	331/333 (99%)	324 (98%)	7 (2%)	0	100	100
1	B	331/333 (99%)	325 (98%)	6 (2%)	0	100	100
1	C	331/333 (99%)	323 (98%)	8 (2%)	0	100	100
1	D	331/333 (99%)	319 (96%)	11 (3%)	1 (0%)	41	27
1	E	331/333 (99%)	325 (98%)	6 (2%)	0	100	100
1	F	331/333 (99%)	325 (98%)	6 (2%)	0	100	100
1	G	331/333 (99%)	320 (97%)	11 (3%)	0	100	100
1	H	331/333 (99%)	320 (97%)	11 (3%)	0	100	100
All	All	2648/2664 (99%)	2581 (98%)	66 (2%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	106	SER

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	289/289 (100%)	287 (99%)	2 (1%)	84	81
1	B	289/289 (100%)	281 (97%)	8 (3%)	43	30
1	C	289/289 (100%)	279 (96%)	10 (4%)	36	21
1	D	289/289 (100%)	284 (98%)	5 (2%)	60	51
1	E	289/289 (100%)	286 (99%)	3 (1%)	76	71
1	F	289/289 (100%)	281 (97%)	8 (3%)	43	30
1	G	289/289 (100%)	277 (96%)	12 (4%)	30	15
1	H	289/289 (100%)	284 (98%)	5 (2%)	60	51
All	All	2312/2312 (100%)	2259 (98%)	53 (2%)	50	37

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	227	ASN
1	A	230	GLU
1	B	15	GLU
1	B	100	ARG
1	B	101	GLN
1	B	106	SER
1	B	107	ARG
1	B	108	LEU
1	B	175	MET
1	B	307	GLN
1	C	18	THR
1	C	99	VAL
1	C	100	ARG
1	C	101	GLN
1	C	113	ARG
1	C	177	GLU
1	C	249	THR
1	C	307	GLN
1	C	312	ASP

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Mol	Chain	Res	Type
1	C	332	LYS
1	D	113	ARG
1	D	119	LYS
1	D	150	LYS
1	D	326	ASP
1	D	332	LYS
1	E	167	ASP
1	E	229	LYS
1	E	249	THR
1	F	14	GLU
1	F	99	VAL
1	F	103	GLU
1	F	112	GLN
1	F	131	ASP
1	F	167	ASP
1	F	227	ASN
1	F	326	ASP
1	G	15	GLU
1	G	101	GLN
1	G	102	GLN
1	G	106	SER
1	G	108	LEU
1	G	167	ASP
1	G	233	LYS
1	G	249	THR
1	G	275	SER
1	G	308	LYS
1	G	312	ASP
1	G	332	LYS
1	H	3	THR
1	H	112	GLN
1	H	249	THR
1	H	308	LYS
1	H	316	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (37) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	21	ASN
1	A	165	ASN
1	A	232	HIS
1	B	114	ASN

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Mol	Chain	Res	Type
1	B	165	ASN
1	B	232	HIS
1	C	116	ASN
1	C	165	ASN
1	C	232	HIS
1	C	328	GLN
1	D	101	GLN
1	D	165	ASN
1	D	232	HIS
1	D	316	GLN
1	D	328	GLN
1	E	114	ASN
1	E	139	ASN
1	E	165	ASN
1	E	232	HIS
1	F	165	ASN
1	F	232	HIS
1	F	316	GLN
1	G	102	GLN
1	G	109	ASN
1	G	112	GLN
1	G	165	ASN
1	G	182	HIS
1	G	232	HIS
1	G	316	GLN
1	H	101	GLN
1	H	109	ASN
1	H	112	GLN
1	H	116	ASN
1	H	124	GLN
1	H	165	ASN
1	H	232	HIS
1	H	316	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

30 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	GOL	E	405	-	5,5,5	0.11	0	5,5,5	0.58	0
2	NAI	C	401	-	42,48,48	0.57	0	47,73,73	0.82	1 (2%)
3	GOL	A	402	-	5,5,5	0.10	0	5,5,5	0.28	0
3	GOL	B	403	-	5,5,5	0.09	0	5,5,5	0.34	0
2	NAI	A	401	-	42,48,48	0.59	0	47,73,73	0.83	1 (2%)
3	GOL	D	501	-	5,5,5	0.08	0	5,5,5	0.35	0
3	GOL	E	404	-	5,5,5	0.11	0	5,5,5	0.27	0
3	GOL	B	404	-	5,5,5	0.15	0	5,5,5	0.35	0
3	GOL	B	405	-	5,5,5	0.21	0	5,5,5	0.49	0
2	NAI	D	503	-	42,48,48	0.61	0	47,73,73	0.85	2 (4%)
2	NAI	B	401	-	42,48,48	0.58	0	47,73,73	0.88	1 (2%)
3	GOL	F	402	-	5,5,5	0.16	0	5,5,5	0.37	0
3	GOL	E	403	-	5,5,5	0.13	0	5,5,5	0.34	0
3	GOL	C	402	-	5,5,5	0.13	0	5,5,5	0.48	0
3	GOL	H	402	-	5,5,5	0.13	0	5,5,5	0.45	0
3	GOL	H	403	-	5,5,5	0.09	0	5,5,5	0.30	0
3	GOL	A	404	-	5,5,5	0.12	0	5,5,5	0.35	0
2	NAI	H	401	-	42,48,48	0.61	0	47,73,73	0.89	3 (6%)
2	NAI	E	401	-	42,48,48	0.57	0	47,73,73	0.79	0
3	GOL	B	402	-	5,5,5	0.20	0	5,5,5	0.43	0
3	GOL	D	505	-	5,5,5	0.16	0	5,5,5	0.33	0
3	GOL	B	406	-	5,5,5	0.07	0	5,5,5	0.28	0
2	NAI	F	401	-	42,48,48	0.56	0	47,73,73	0.80	1 (2%)
3	GOL	G	402	-	5,5,5	0.13	0	5,5,5	0.39	0
3	GOL	B	407	-	5,5,5	0.12	0	5,5,5	0.32	0
3	GOL	E	402	-	5,5,5	0.13	0	5,5,5	0.30	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAI	G	401	-	42,48,48	0.62	0	47,73,73	0.75	1 (2%)
3	GOL	A	403	-	5,5,5	0.09	0	5,5,5	0.42	0
3	GOL	D	504	-	5,5,5	0.12	0	5,5,5	0.44	0
3	GOL	D	502	-	5,5,5	0.11	0	5,5,5	0.40	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	GOL	E	405	-	-	2/4/4/4	-
2	NAI	C	401	-	-	5/25/72/72	0/5/5/5
3	GOL	A	402	-	-	4/4/4/4	-
3	GOL	B	403	-	-	4/4/4/4	-
2	NAI	A	401	-	-	3/25/72/72	0/5/5/5
3	GOL	D	501	-	-	2/4/4/4	-
3	GOL	E	404	-	-	3/4/4/4	-
3	GOL	B	404	-	-	4/4/4/4	-
3	GOL	B	405	-	-	2/4/4/4	-
2	NAI	D	503	-	-	3/25/72/72	0/5/5/5
2	NAI	B	401	-	-	5/25/72/72	0/5/5/5
3	GOL	F	402	-	-	3/4/4/4	-
3	GOL	E	403	-	-	2/4/4/4	-
3	GOL	C	402	-	-	4/4/4/4	-
3	GOL	H	402	-	-	2/4/4/4	-
3	GOL	H	403	-	-	3/4/4/4	-
3	GOL	A	404	-	-	4/4/4/4	-
2	NAI	H	401	-	-	4/25/72/72	0/5/5/5
2	NAI	E	401	-	-	3/25/72/72	0/5/5/5
3	GOL	B	402	-	-	4/4/4/4	-
3	GOL	D	505	-	-	2/4/4/4	-
3	GOL	B	406	-	-	0/4/4/4	-
2	NAI	F	401	-	-	3/25/72/72	0/5/5/5
3	GOL	G	402	-	-	0/4/4/4	-
3	GOL	B	407	-	-	1/4/4/4	-
3	GOL	E	402	-	-	0/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAI	G	401	-	-	4/25/72/72	0/5/5/5
3	GOL	A	403	-	-	0/4/4/4	-
3	GOL	D	504	-	-	4/4/4/4	-
3	GOL	D	502	-	-	2/4/4/4	-

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	NAI	C5A-C6A-N6A	2.48	124.12	120.35
2	B	401	NAI	C5A-C6A-N6A	2.46	124.09	120.35
2	H	401	NAI	O4D-C1D-C2D	-2.41	101.38	106.64
2	H	401	NAI	O4B-C1B-C2B	-2.32	103.54	106.93
2	H	401	NAI	C5A-C6A-N6A	2.30	123.85	120.35
2	C	401	NAI	C5A-C6A-N6A	2.28	123.82	120.35
2	D	503	NAI	O4B-C1B-C2B	-2.07	103.91	106.93
2	G	401	NAI	C5A-C6A-N6A	2.04	123.45	120.35
2	F	401	NAI	C5A-C6A-N6A	2.02	123.43	120.35
2	D	503	NAI	C5A-C6A-N6A	2.01	123.40	120.35

There are no chirality outliers.

All (82) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	404	GOL	O1-C1-C2-C3
3	A	404	GOL	C1-C2-C3-O3
3	B	402	GOL	O1-C1-C2-C3
3	B	402	GOL	C1-C2-C3-O3
3	B	403	GOL	O1-C1-C2-O2
3	B	403	GOL	O1-C1-C2-C3
3	B	403	GOL	C1-C2-C3-O3
3	B	404	GOL	O1-C1-C2-O2
3	B	404	GOL	O1-C1-C2-C3
3	B	405	GOL	C1-C2-C3-O3
3	C	402	GOL	O1-C1-C2-C3
3	D	501	GOL	O1-C1-C2-C3
3	D	502	GOL	O1-C1-C2-C3
3	D	504	GOL	O1-C1-C2-C3
3	E	404	GOL	C1-C2-C3-O3
3	E	405	GOL	C1-C2-C3-O3

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Mol	Chain	Res	Type	Atoms
3	F	402	GOL	O1-C1-C2-C3
3	H	403	GOL	C1-C2-C3-O3
3	C	402	GOL	O1-C1-C2-O2
3	D	501	GOL	O1-C1-C2-O2
3	D	504	GOL	O1-C1-C2-O2
3	A	402	GOL	O1-C1-C2-C3
3	A	402	GOL	C1-C2-C3-O3
3	B	404	GOL	C1-C2-C3-O3
3	B	407	GOL	C1-C2-C3-O3
3	C	402	GOL	C1-C2-C3-O3
3	E	403	GOL	C1-C2-C3-O3
3	A	404	GOL	O1-C1-C2-O2
3	B	402	GOL	O2-C2-C3-O3
3	B	403	GOL	O2-C2-C3-O3
3	B	404	GOL	O2-C2-C3-O3
3	B	405	GOL	O2-C2-C3-O3
3	D	502	GOL	O1-C1-C2-O2
3	F	402	GOL	O1-C1-C2-O2
3	A	402	GOL	O2-C2-C3-O3
3	B	402	GOL	O1-C1-C2-O2
3	E	404	GOL	O2-C2-C3-O3
3	E	405	GOL	O2-C2-C3-O3
3	A	402	GOL	O1-C1-C2-O2
3	E	403	GOL	O2-C2-C3-O3
3	E	404	GOL	O1-C1-C2-O2
3	H	402	GOL	O2-C2-C3-O3
2	B	401	NAI	C2D-C1D-N1N-C2N
3	C	402	GOL	O2-C2-C3-O3
3	D	504	GOL	O2-C2-C3-O3
3	D	505	GOL	O2-C2-C3-O3
3	H	402	GOL	C1-C2-C3-O3
2	B	401	NAI	C2D-C1D-N1N-C6N
3	H	403	GOL	O1-C1-C2-O2
3	H	403	GOL	O2-C2-C3-O3
2	C	401	NAI	O4D-C1D-N1N-C2N
2	B	401	NAI	O4D-C1D-N1N-C2N
2	G	401	NAI	O4D-C1D-N1N-C2N
2	C	401	NAI	C2D-C1D-N1N-C2N
2	A	401	NAI	O4D-C1D-N1N-C2N
2	D	503	NAI	O4D-C1D-N1N-C2N
2	E	401	NAI	O4D-C1D-N1N-C2N
2	F	401	NAI	O4D-C1D-N1N-C2N

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Mol	Chain	Res	Type	Atoms
2	H	401	NAI	O4D-C1D-N1N-C2N
2	A	401	NAI	C2D-C1D-N1N-C2N
2	D	503	NAI	C2D-C1D-N1N-C2N
2	F	401	NAI	C2D-C1D-N1N-C2N
2	G	401	NAI	C2D-C1D-N1N-C2N
2	H	401	NAI	C2D-C1D-N1N-C2N
2	E	401	NAI	C2D-C1D-N1N-C2N
2	H	401	NAI	O4B-C4B-C5B-O5B
2	B	401	NAI	O4D-C1D-N1N-C6N
3	D	505	GOL	C1-C2-C3-O3
3	F	402	GOL	C1-C2-C3-O3
3	A	404	GOL	O2-C2-C3-O3
2	F	401	NAI	O4B-C4B-C5B-O5B
2	G	401	NAI	O4B-C4B-C5B-O5B
2	C	401	NAI	C2D-C1D-N1N-C6N
2	B	401	NAI	O4B-C4B-C5B-O5B
2	E	401	NAI	O4B-C4B-C5B-O5B
3	D	504	GOL	C1-C2-C3-O3
2	A	401	NAI	O4B-C4B-C5B-O5B
2	C	401	NAI	O4B-C4B-C5B-O5B
2	D	503	NAI	O4B-C4B-C5B-O5B
2	C	401	NAI	O4D-C1D-N1N-C6N
2	G	401	NAI	C2D-C1D-N1N-C6N
2	H	401	NAI	C2D-C1D-N1N-C6N

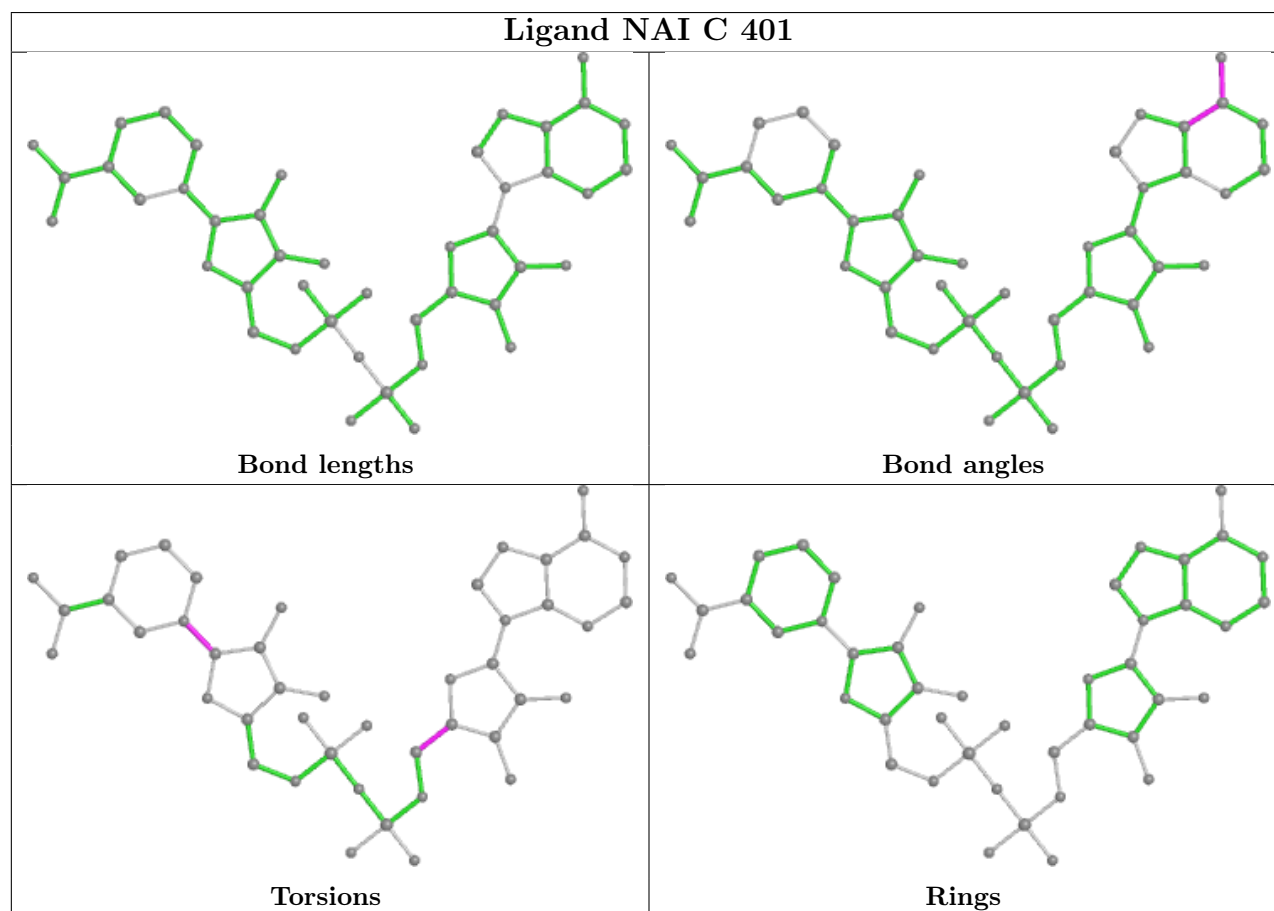
There are no ring outliers.

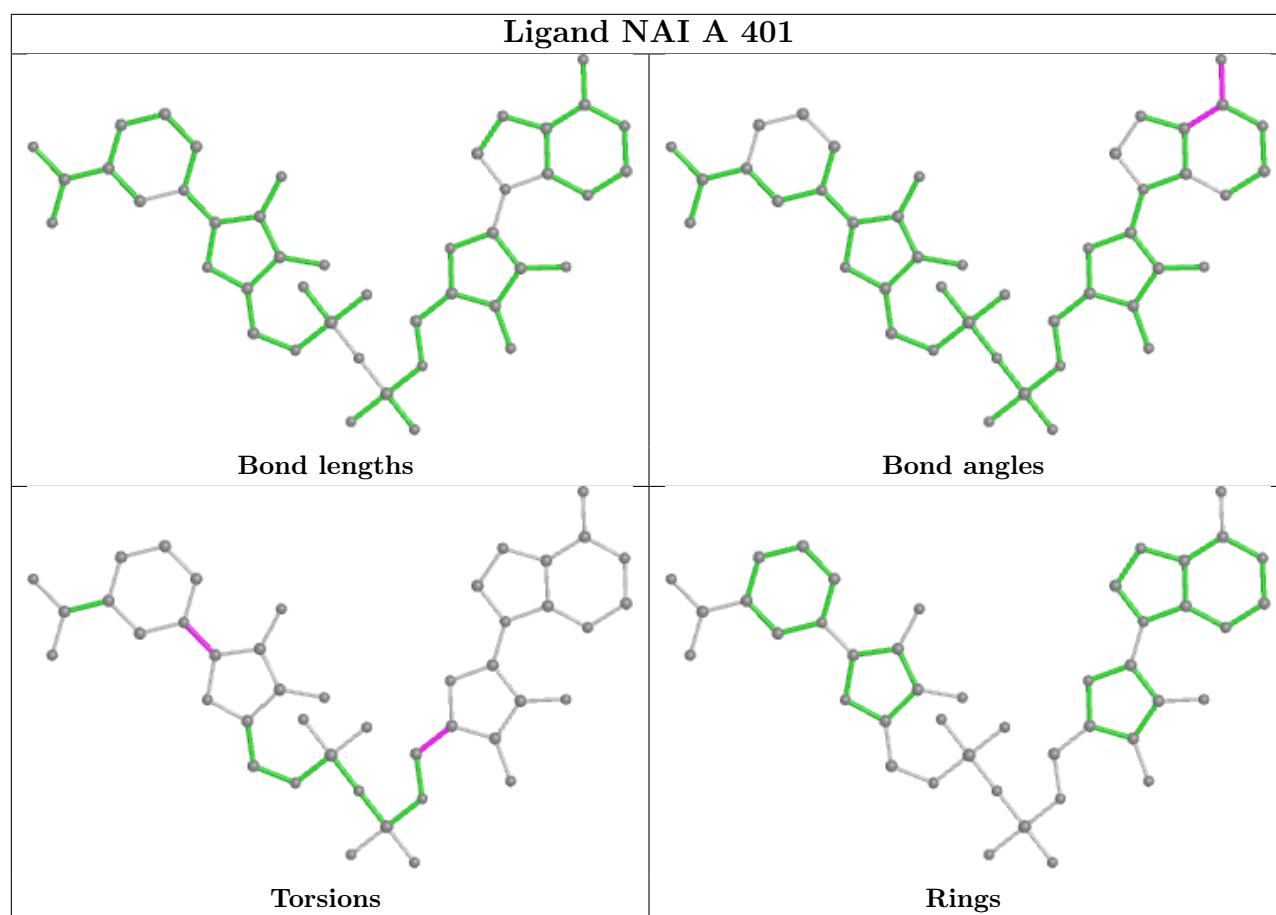
11 monomers are involved in 17 short contacts:

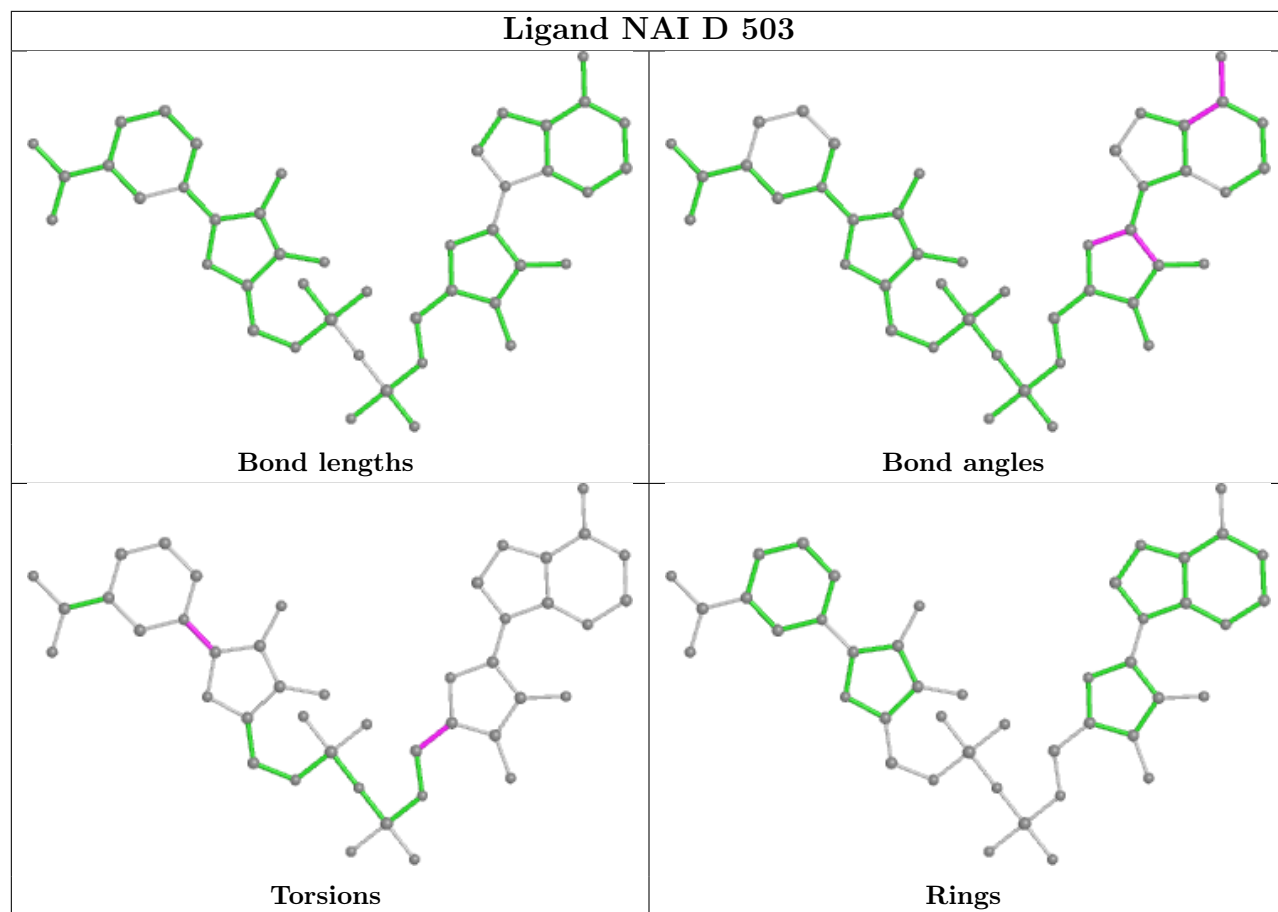
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	E	405	GOL	1	0
2	C	401	NAI	2	0
3	D	501	GOL	4	0
3	E	404	GOL	2	0
3	B	405	GOL	1	0
2	B	401	NAI	1	0
3	E	403	GOL	2	0
3	H	403	GOL	1	0
3	G	402	GOL	1	0
2	G	401	NAI	1	0
3	D	502	GOL	6	0

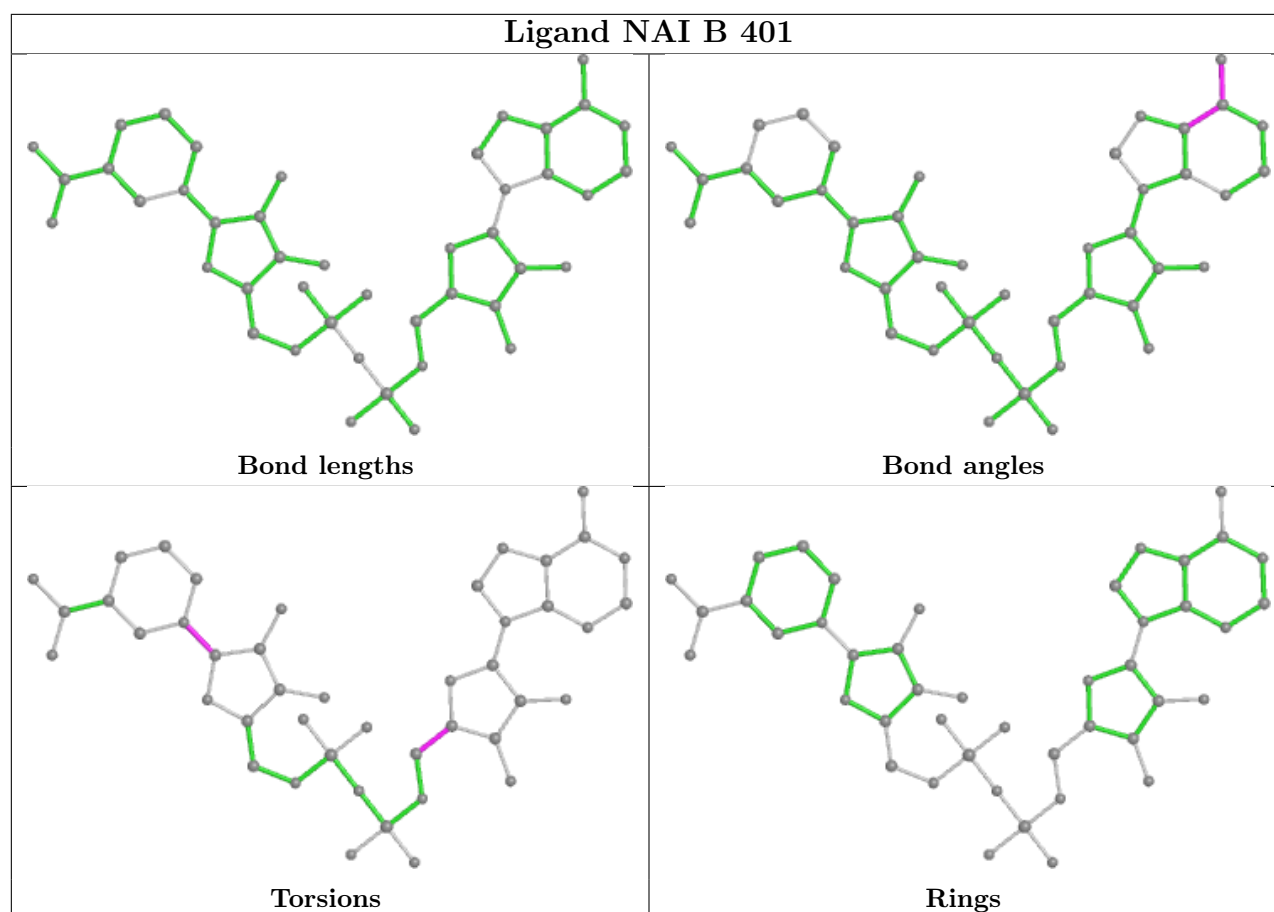
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths,

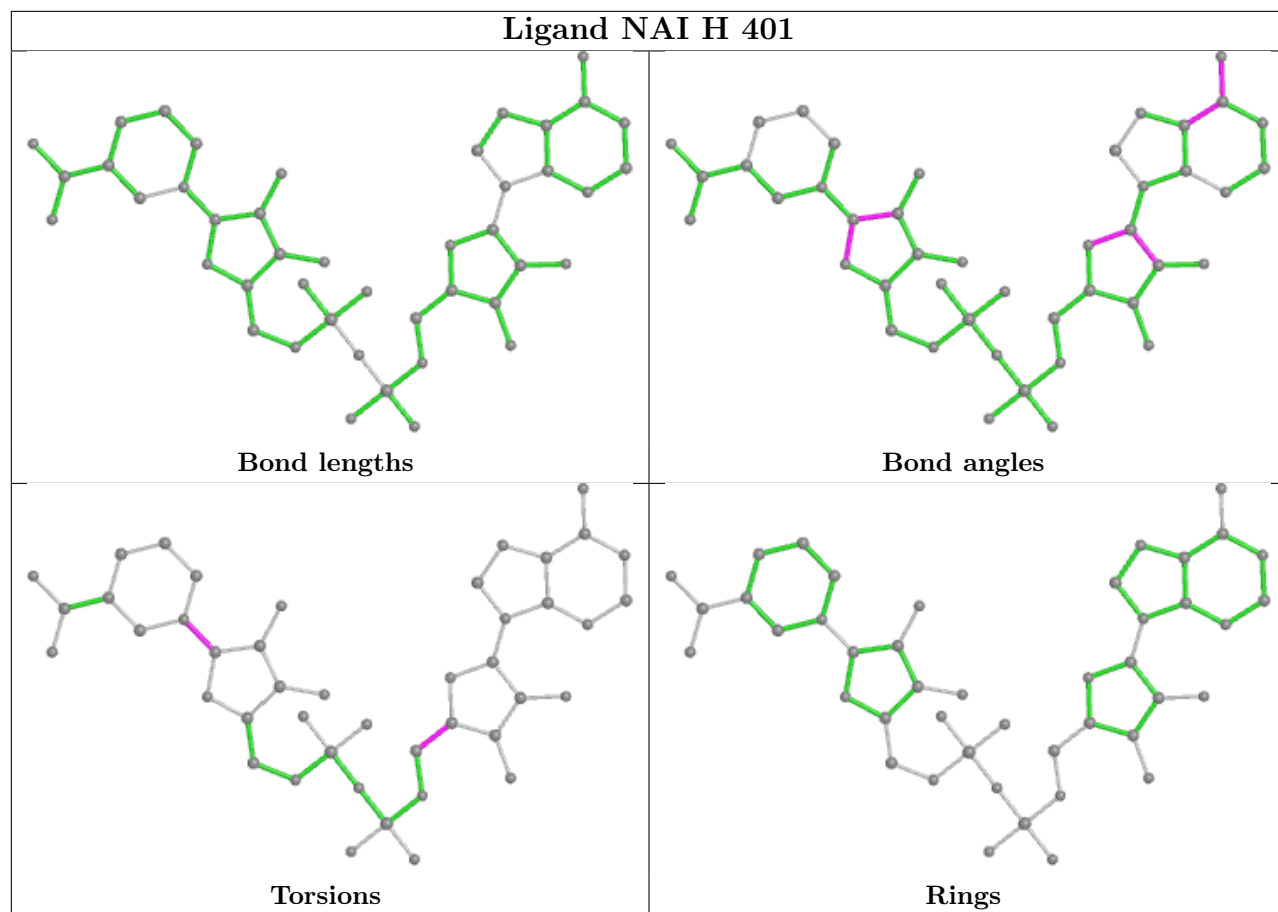
bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within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

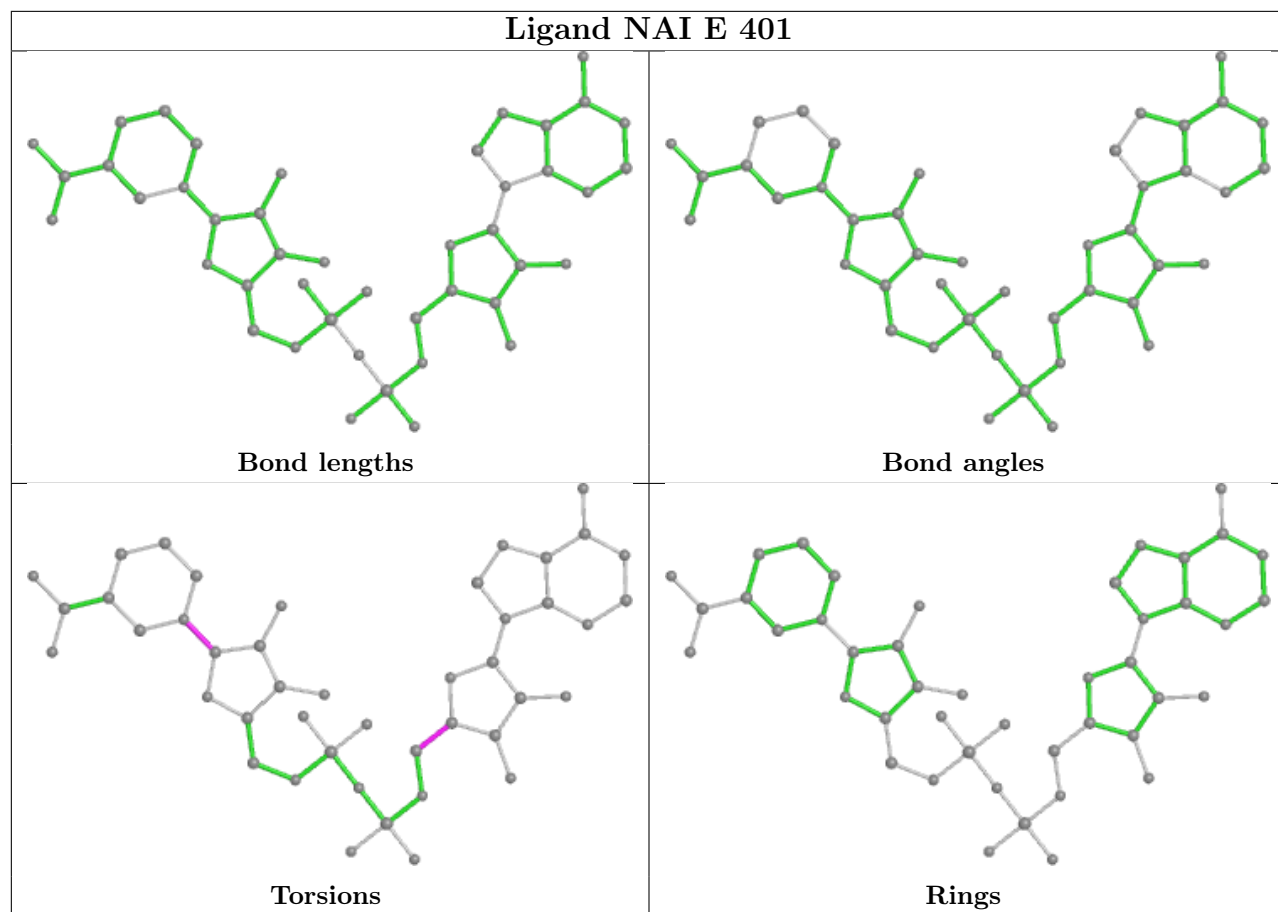


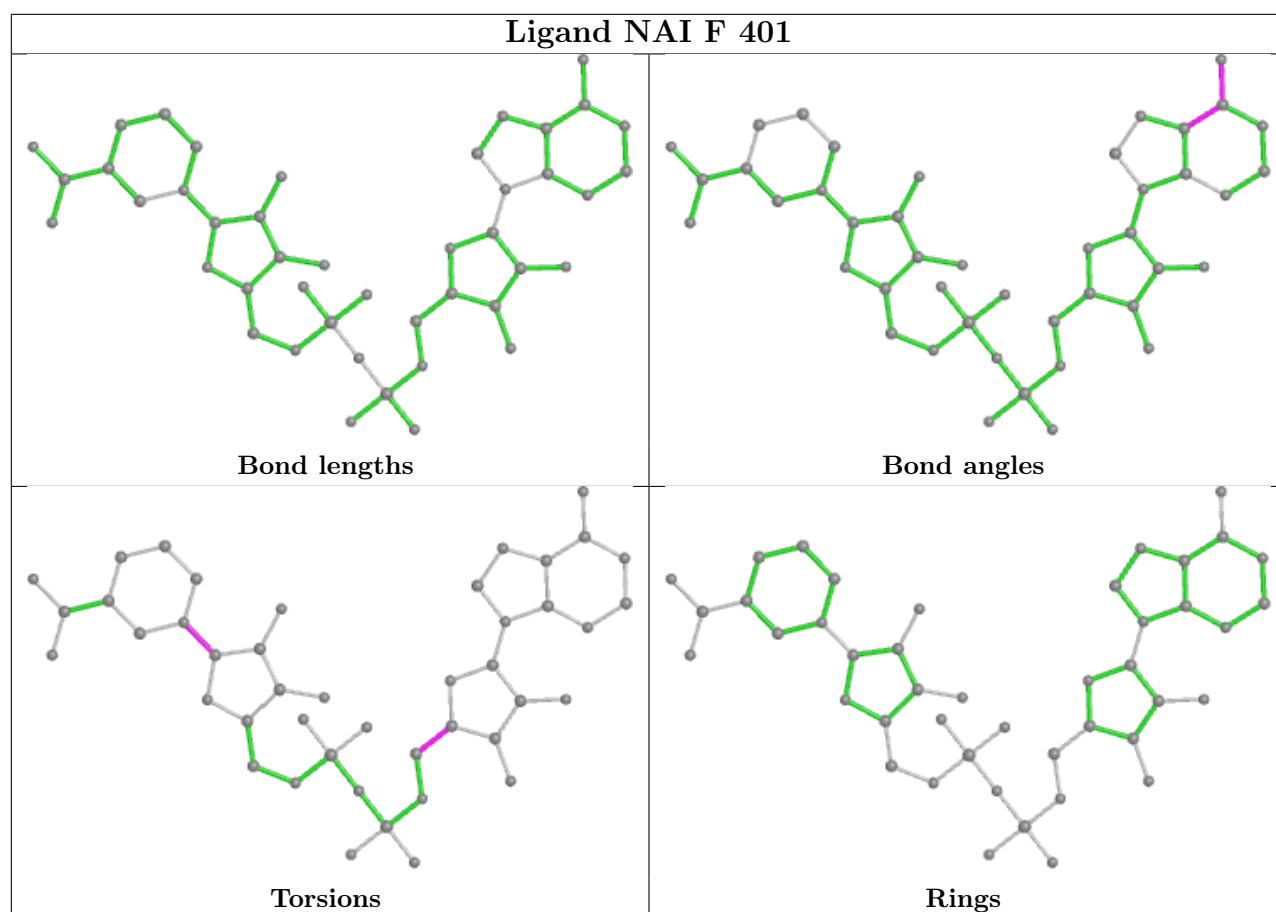


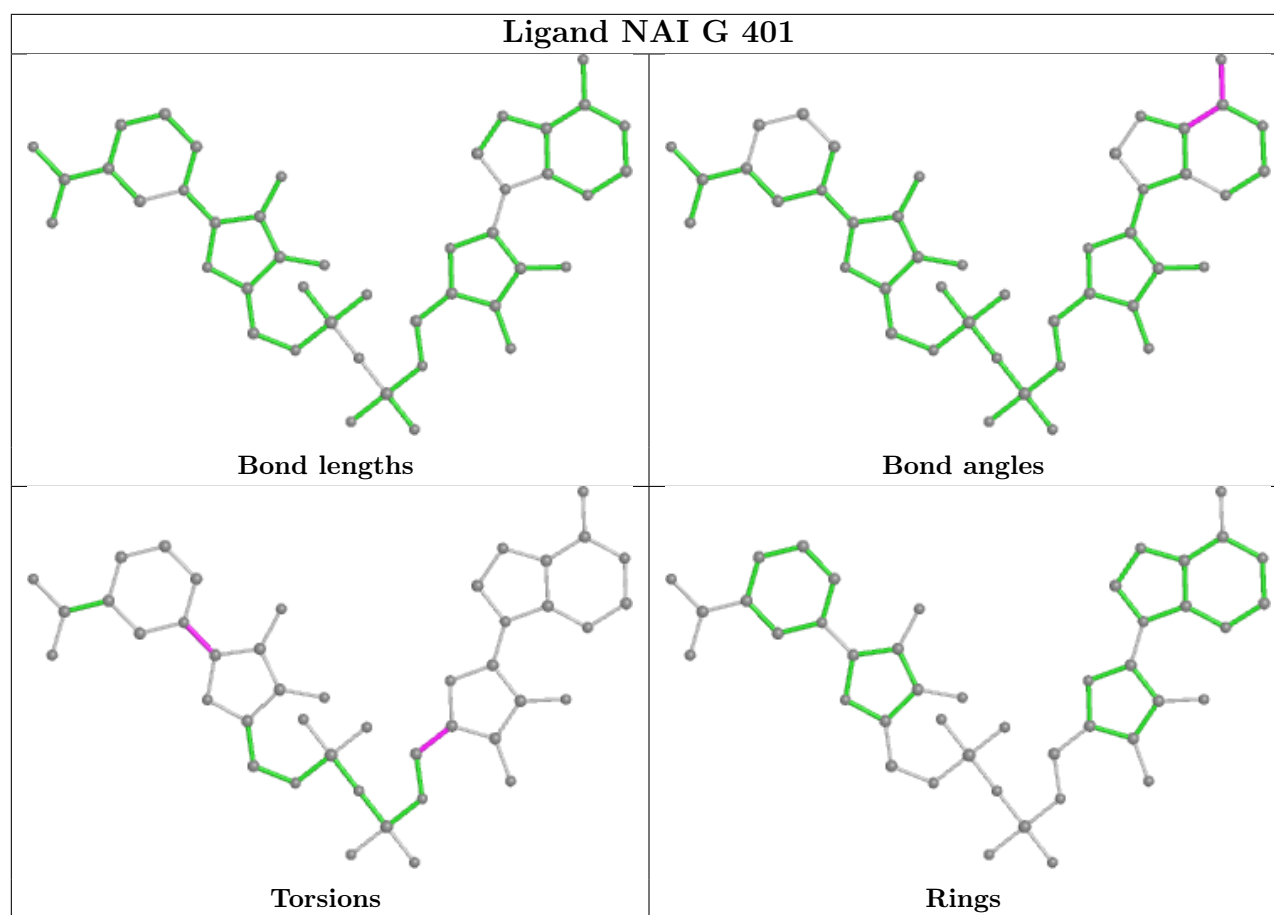












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ > 2	OWAB(Å ²)	Q < 0.9
1	A	333/333 (100%)	-0.30	6 (1%) 68 64	13, 22, 45, 66	0
1	B	333/333 (100%)	-0.29	11 (3%) 46 40	14, 23, 52, 71	0
1	C	333/333 (100%)	0.04	23 (6%) 16 13	17, 29, 56, 88	0
1	D	333/333 (100%)	0.12	19 (5%) 23 19	18, 30, 57, 86	0
1	E	333/333 (100%)	-0.31	4 (1%) 79 76	16, 25, 43, 69	0
1	F	333/333 (100%)	0.03	17 (5%) 28 22	17, 28, 55, 72	0
1	G	333/333 (100%)	0.16	25 (7%) 14 11	21, 35, 67, 96	0
1	H	333/333 (100%)	0.08	20 (6%) 21 17	17, 30, 54, 68	0
All	All	2664/2664 (100%)	-0.06	125 (4%) 31 25	13, 28, 56, 96	0

All (125) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	104	GLY	8.1
1	H	17	ALA	7.6
1	G	101	GLN	7.6
1	D	106	SER	7.5
1	D	105	GLU	6.5
1	H	120	PHE	5.8
1	C	106	SER	5.8
1	F	107	ARG	5.5
1	D	120	PHE	5.3
1	C	107	ARG	5.3
1	C	109	ASN	5.0
1	B	107	ARG	5.0
1	C	103	GLU	4.7
1	C	100	ARG	4.7
1	G	107	ARG	4.7
1	D	222	ASP	4.6

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Mol	Chain	Res	Type	RSRZ
1	G	108	LEU	4.6
1	G	334	LEU	4.6
1	D	107	ARG	4.6
1	A	223	ASN	4.4
1	H	102	GLN	4.4
1	C	101	GLN	4.3
1	F	102	GLN	4.3
1	G	2	ALA	4.2
1	G	111	VAL	4.2
1	D	113	ARG	4.2
1	D	116	ASN	4.2
1	G	100	ARG	4.1
1	A	17	ALA	4.1
1	H	15	GLU	4.0
1	F	222	ASP	3.9
1	D	104	GLY	3.8
1	F	223	ASN	3.8
1	A	15	GLU	3.8
1	G	223	ASN	3.7
1	G	103	GLU	3.7
1	D	101	GLN	3.7
1	D	334	LEU	3.6
1	G	17	ALA	3.6
1	C	99	VAL	3.6
1	G	104	GLY	3.6
1	E	334	LEU	3.5
1	G	120	PHE	3.5
1	C	333	ASP	3.5
1	F	333	ASP	3.4
1	B	222	ASP	3.4
1	F	334	LEU	3.4
1	H	333	ASP	3.3
1	F	103	GLU	3.2
1	H	2	ALA	3.2
1	G	218	GLU	3.1
1	D	330	ASP	3.1
1	A	308	LYS	3.1
1	G	112	GLN	3.1
1	E	225	SER	3.0
1	C	334	LEU	3.0
1	G	15	GLU	3.0
1	H	107	ARG	3.0

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Mol	Chain	Res	Type	RSRZ
1	H	106	SER	3.0
1	F	108	LEU	2.9
1	G	332	LYS	2.9
1	B	100	ARG	2.9
1	B	102	GLN	2.9
1	D	325	TRP	2.8
1	H	326	ASP	2.8
1	D	315	ALA	2.8
1	D	311	ASP	2.8
1	H	109	ASN	2.8
1	B	228	TRP	2.8
1	A	222	ASP	2.8
1	F	100	ARG	2.8
1	B	17	ALA	2.7
1	F	240	TYR	2.7
1	H	14	GLU	2.7
1	H	104	GLY	2.7
1	B	223	ASN	2.6
1	G	4	LEU	2.6
1	C	17	ALA	2.6
1	F	17	ALA	2.6
1	G	326	ASP	2.6
1	F	315	ALA	2.6
1	D	228	TRP	2.6
1	D	109	ASN	2.6
1	C	111	VAL	2.6
1	H	325	TRP	2.5
1	C	102	GLN	2.5
1	F	99	VAL	2.5
1	B	334	LEU	2.5
1	G	99	VAL	2.5
1	C	3	THR	2.5
1	B	308	LYS	2.4
1	G	333	ASP	2.4
1	F	109	ASN	2.4
1	G	217	PRO	2.4
1	H	108	LEU	2.4
1	G	228	TRP	2.4
1	G	226	GLU	2.4
1	C	308	LYS	2.4
1	F	283	GLY	2.4
1	H	223	ASN	2.4

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Mol	Chain	Res	Type	RSRZ
1	E	223	ASN	2.3
1	F	228	TRP	2.3
1	H	18	THR	2.3
1	B	2	ALA	2.3
1	C	332	LYS	2.3
1	D	225	SER	2.3
1	C	312	ASP	2.2
1	C	108	LEU	2.2
1	H	237	GLU	2.2
1	D	223	ASN	2.2
1	D	326	ASP	2.2
1	C	326	ASP	2.2
1	F	311	ASP	2.2
1	C	15	GLU	2.1
1	E	222	ASP	2.1
1	G	308	LYS	2.1
1	H	334	LEU	2.1
1	H	16	GLU	2.1
1	B	333	ASP	2.1
1	G	109	ASN	2.1
1	C	120	PHE	2.1
1	C	115	VAL	2.1
1	C	131	ASP	2.1
1	A	18	THR	2.0
1	H	332	LYS	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 monosaccharides in this entry.

6.4 Ligands ⓘ

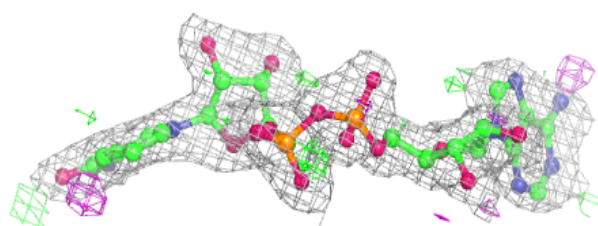
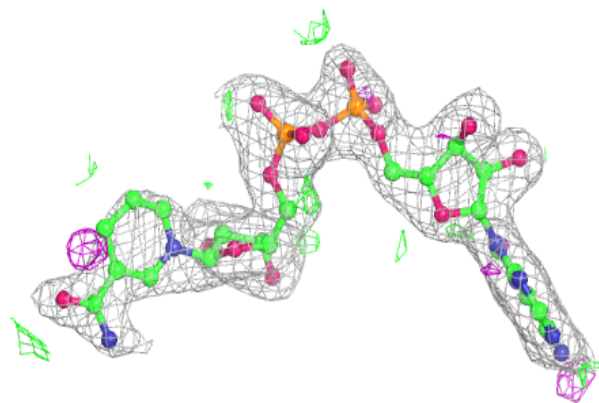
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	GOL	E	404	6/6	0.75	0.49	22,22,24,27	6
3	GOL	A	402	6/6	0.76	0.17	41,45,46,51	0
3	GOL	E	403	6/6	0.79	0.12	44,46,46,48	0
3	GOL	H	403	6/6	0.79	0.20	48,52,55,55	0
3	GOL	E	405	6/6	0.80	0.63	23,24,25,26	6
3	GOL	H	402	6/6	0.81	0.39	30,31,31,31	6
3	GOL	B	405	6/6	0.81	0.27	18,21,21,22	6
3	GOL	A	403	6/6	0.82	0.28	22,25,28,28	6
3	GOL	B	404	6/6	0.83	0.14	37,43,44,46	0
3	GOL	A	404	6/6	0.83	0.14	34,39,40,46	0
3	GOL	G	402	6/6	0.83	0.32	25,27,27,27	6
3	GOL	D	504	6/6	0.83	0.11	40,40,41,41	0
3	GOL	B	402	6/6	0.83	0.12	39,41,43,45	0
3	GOL	D	502	6/6	0.84	0.37	27,30,32,33	6
3	GOL	F	402	6/6	0.85	0.17	38,46,49,53	0
3	GOL	E	402	6/6	0.87	0.14	29,33,35,39	0
3	GOL	D	505	6/6	0.87	0.14	35,40,42,43	0
3	GOL	B	403	6/6	0.89	0.17	32,34,39,39	0
3	GOL	B	406	6/6	0.90	0.22	38,42,44,44	0
3	GOL	B	407	6/6	0.91	0.13	39,47,48,53	0
3	GOL	D	501	6/6	0.93	0.32	23,24,24,25	6
2	NAI	G	401	44/44	0.93	0.11	29,39,46,47	0
3	GOL	C	402	6/6	0.93	0.16	27,37,38,40	0
2	NAI	C	401	44/44	0.95	0.10	26,33,39,42	0
2	NAI	D	503	44/44	0.96	0.07	27,30,35,39	0
2	NAI	H	401	44/44	0.96	0.07	27,30,33,35	0
2	NAI	F	401	44/44	0.96	0.09	24,30,34,37	0
2	NAI	E	401	44/44	0.97	0.07	16,20,25,27	0
2	NAI	B	401	44/44	0.97	0.07	20,25,29,30	0
2	NAI	A	401	44/44	0.98	0.08	13,18,22,25	0

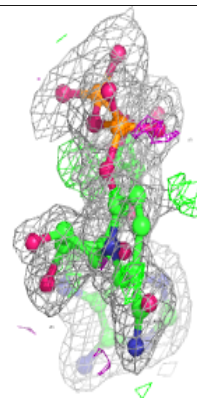
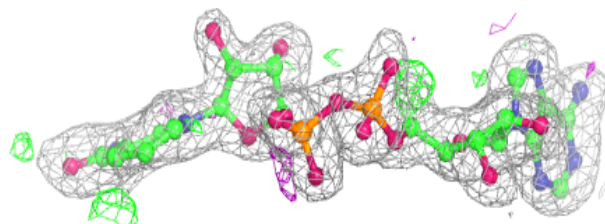
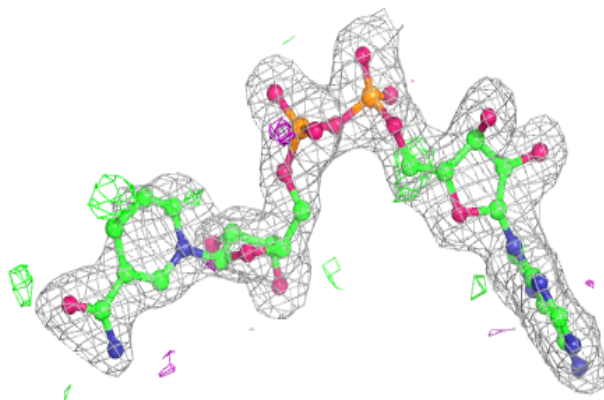
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around NAI G 401:

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

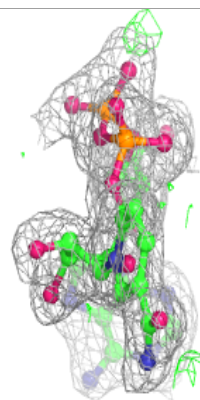
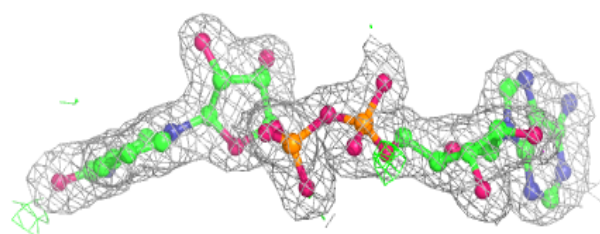
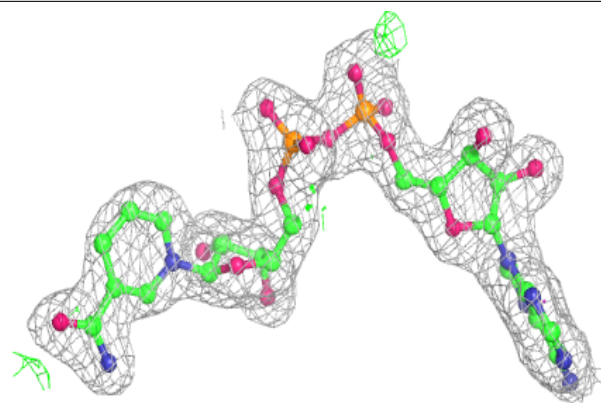
**Electron density around NAI C 401:**

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

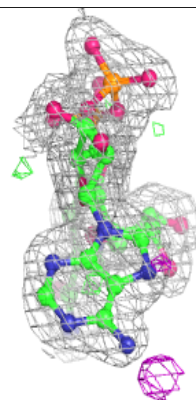
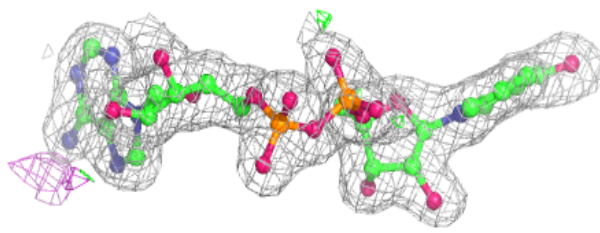
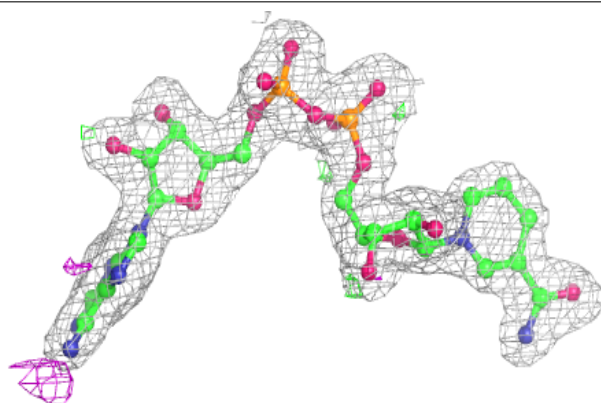


Electron density around NAI D 503:

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

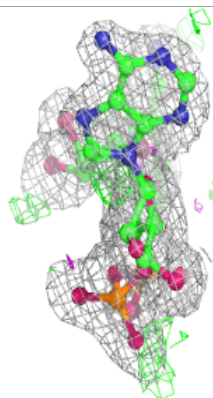
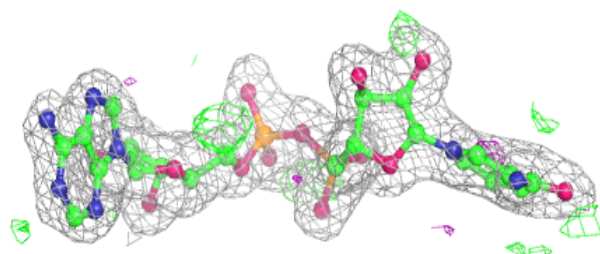
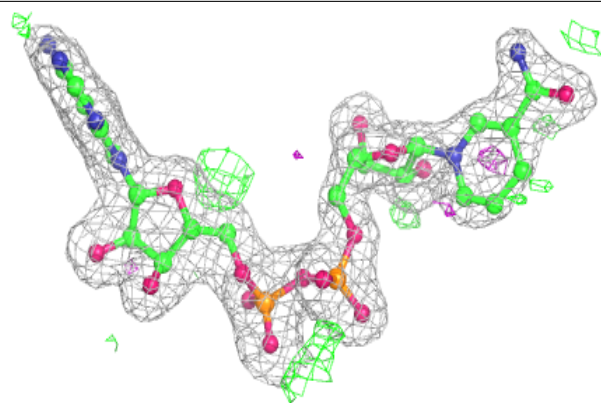
**Electron density around NAI H 401:**

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

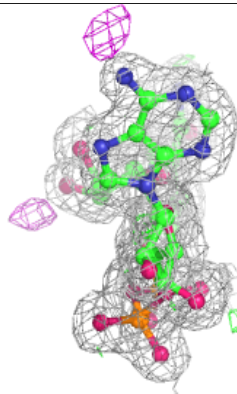
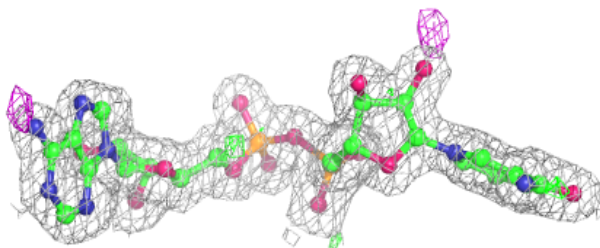
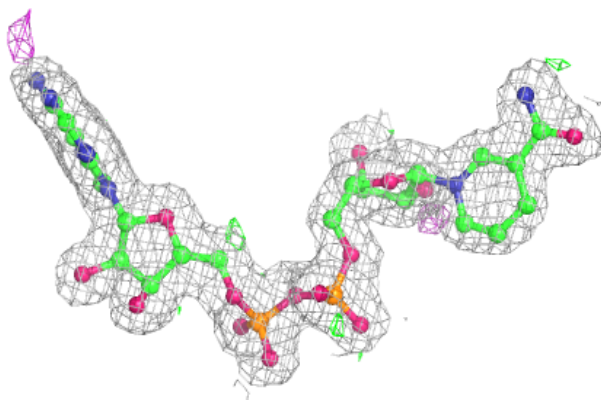


Electron density around NAI F 401:

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

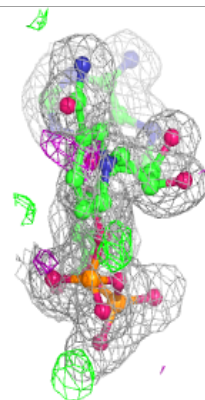
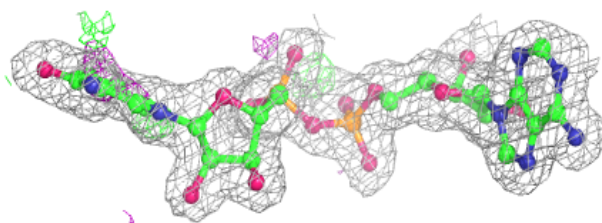
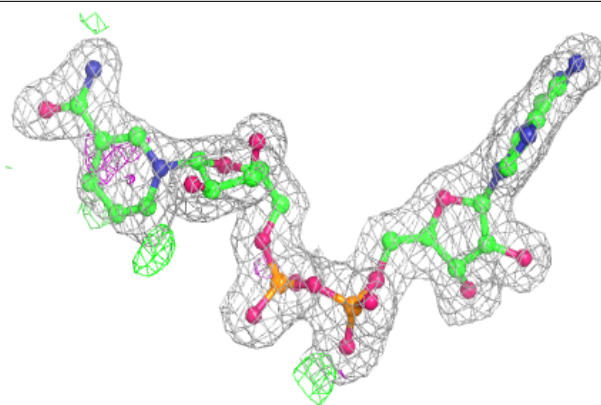
**Electron density around NAI E 401:**

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

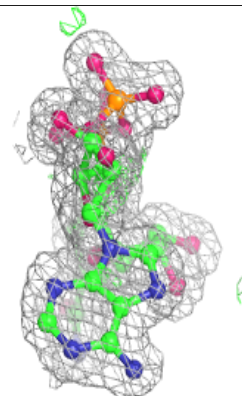
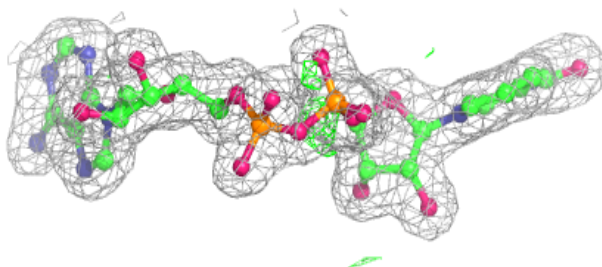
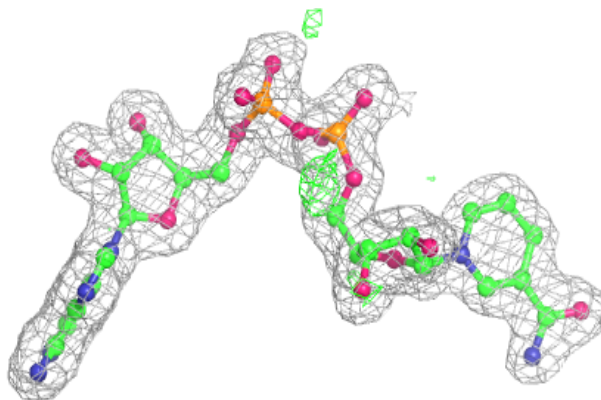


Electron density around NAI B 401:

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

**Electron density around NAI A 401:**

$2mF_o-DF_c$ (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.