



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 9, 2020 – 04:03 PM BST

PDB ID : 3WSW  
Title : Crystal structure of minor L-lactate dehydrogenase from *Enterococcus mundtii* in the ligands-bound form  
Authors : Matoba, Y.; Sugiyama, M.  
Deposited on : 2014-03-27  
Resolution : 2.30 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](mailto: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.

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

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

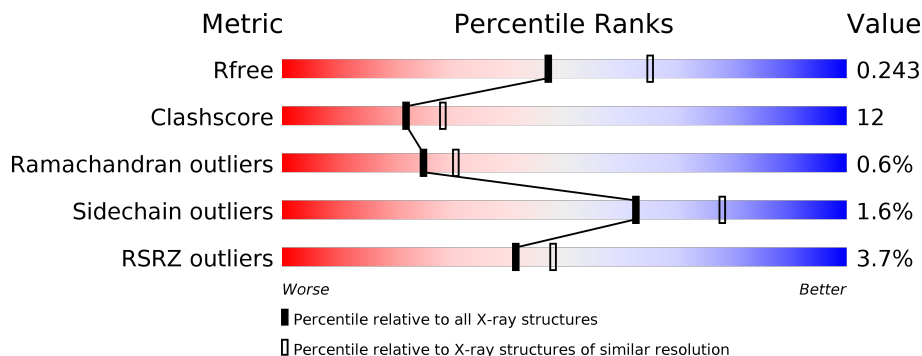
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.30 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	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	322	 3% 73% 25%
1	B	322	 5% 71% 25%
1	C	322	 2% 75% 22%
1	D	322	 4% 73% 22%

## 2 Entry composition [i](#)

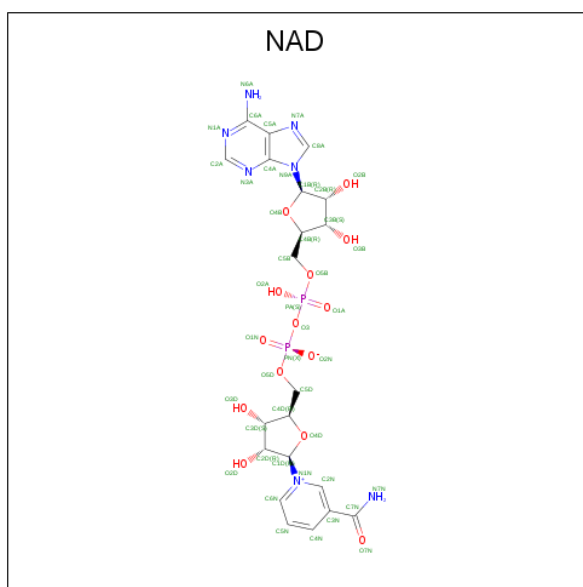
There are 5 unique types of molecules in this entry. The entry contains 10436 atoms, of which 0 are hydrogens and 0 are deuteriums.

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.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	316	Total 2431	C 1532	N 412	O 480	S 7	0	0	0
1	B	313	Total 2402	C 1515	N 405	O 475	S 7	0	0	0
1	C	314	Total 2411	C 1520	N 406	O 478	S 7	0	0	0
1	D	314	Total 2411	C 1520	N 406	O 478	S 7	0	0	0

- Molecule 2 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: C<sub>21</sub>H<sub>27</sub>N<sub>7</sub>O<sub>14</sub>P<sub>2</sub>).



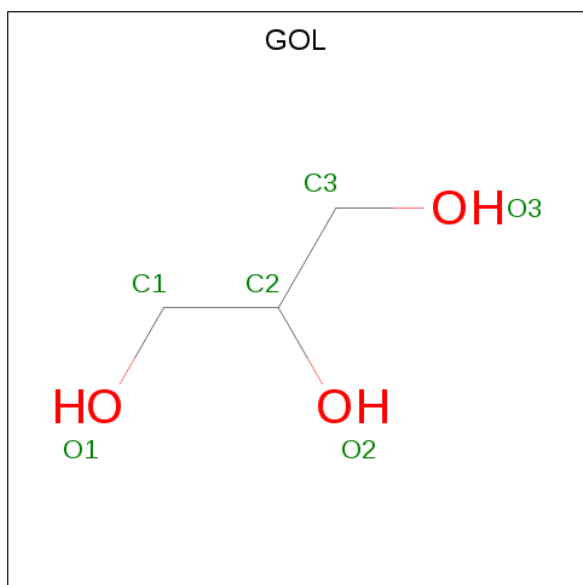
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
			Total	C	N	O			P
2	A	1	Total 44	C 21	N 7	O 14	P 2	0	0
2	B	1	Total 44	C 21	N 7	O 14	P 2	0	0

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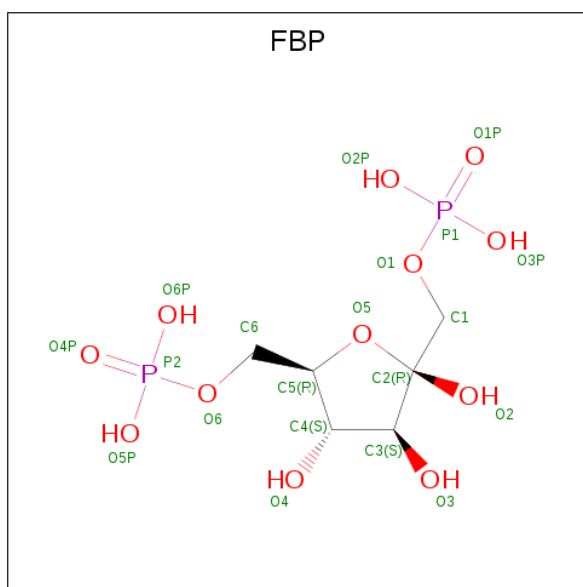
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
2	C	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	D	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula:  $C_3H_8O_3$ ).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	
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		

- Molecule 4 is 1,6-di-O-phosphono-beta-D-fructofuranose (three-letter code: FBP) (formula:  $C_6H_{14}O_{12}P_2$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	
			Total	C	O			P
4	B	1	40	12	24	4	0	1
4	D	1	40	12	24	4	0	1

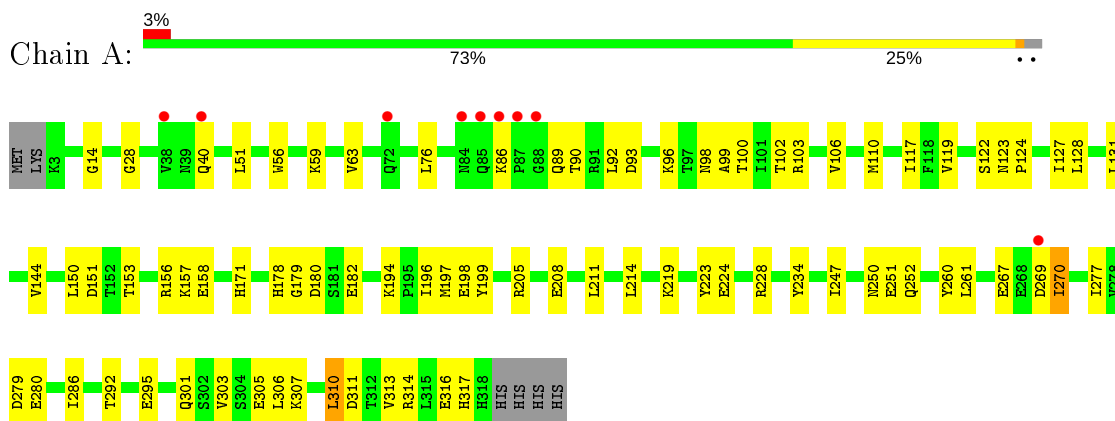
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	A	127	127	127	0	0
5	B	108	108	108	0	0
5	C	114	114	114	0	0
5	D	158	158	158	0	0

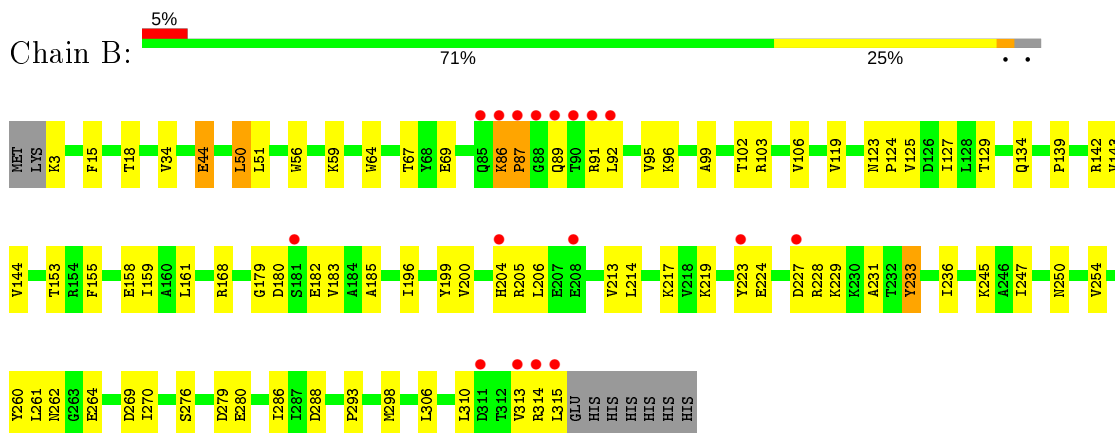
### 3 Residue-property plots [i](#)

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.

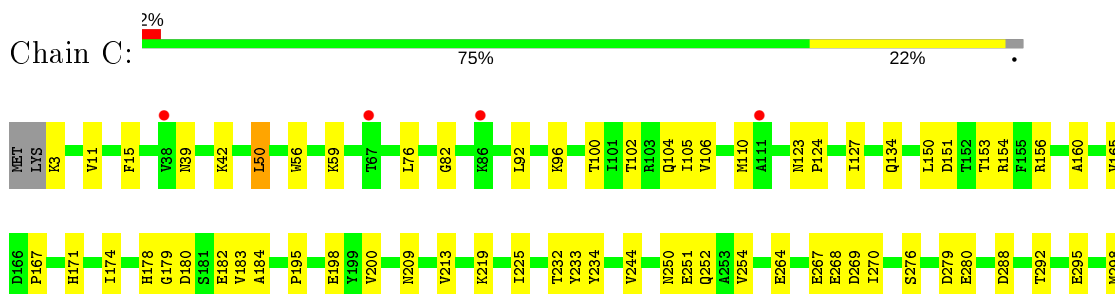
- Molecule 1: L-lactate dehydrogenase

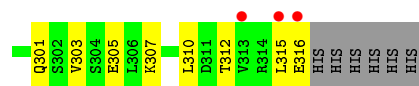


- Molecule 1: L-lactate dehydrogenase

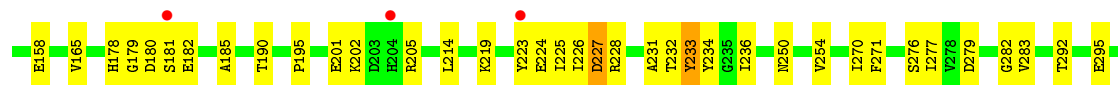
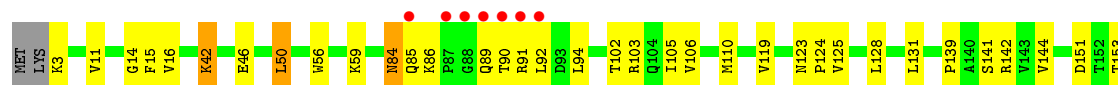
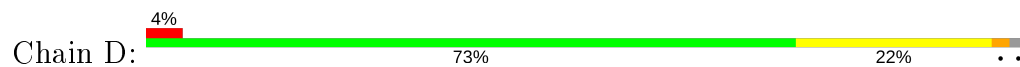


- Molecule 1: L-lactate dehydrogenase





- Molecule 1: L-lactate dehydrogenase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	67.77Å 123.54Å 85.79Å 90.00° 107.75° 90.00°	Depositor
Resolution (Å)	29.67 – 2.30 38.53 – 2.30	Depositor EDS
% Data completeness (in resolution range)	93.1 (29.67-2.30) 95.4 (38.53-2.30)	Depositor EDS
$R_{merge}$	0.11	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.30 (at 2.29Å)	Xtrriage
Refinement program	CNS 1.2	Depositor
R, $R_{free}$	0.195 , 0.246 0.191 , 0.243	Depositor DCC
$R_{free}$ test set	2882 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	23.0	Xtrriage
Anisotropy	1.327	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.39 , 59.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	10436	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	33.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.13% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: GOL, FBP, NAD

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
1	A	0.32	0/2469	0.58	0/3353
1	B	0.33	0/2438	0.58	0/3311
1	C	0.32	0/2447	0.57	0/3323
1	D	0.34	0/2447	0.58	0/3323
All	All	0.33	0/9801	0.58	0/13310

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2431	0	2435	56	0
1	B	2402	0	2415	74	0
1	C	2411	0	2421	50	0
1	D	2411	0	2421	64	0
2	A	44	0	26	2	0
2	B	44	0	26	2	0
2	C	44	0	26	2	0
2	D	44	0	26	3	0
3	A	6	0	8	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	12	0	16	3	0
4	B	40	0	20	1	0
4	D	40	0	20	1	0
5	A	127	0	0	1	0
5	B	108	0	0	4	0
5	C	114	0	0	4	0
5	D	158	0	0	4	0
All	All	10436	0	9860	233	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 (233) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:90:THR:HG22	1:A:92:LEU:H	1.19	1.06
1:D:180:ASP:HA	1:D:219:LYS:HD3	1.52	0.92
1:A:180:ASP:HA	1:A:219:LYS:HD2	1.58	0.85
1:D:84:ASN:HD22	1:D:85:GLN:H	1.23	0.85
1:B:86:LYS:HB3	1:B:87:PRO:HD2	1.57	0.85
1:C:250:ASN:OD1	1:C:280:GLU:HG2	1.79	0.83
1:A:98:ASN:O	1:A:102:THR:HG23	1.80	0.82
1:B:86:LYS:HB2	1:B:89:GLN:HG3	1.62	0.80
1:D:84:ASN:HD22	1:D:85:GLN:N	1.81	0.78
1:A:76:LEU:HD23	1:A:117:ILE:HB	1.69	0.73
1:D:85:GLN:HG2	1:D:94:LEU:HD22	1.70	0.73
1:B:86:LYS:HB2	1:B:89:GLN:CG	2.18	0.73
1:C:100:THR:O	1:C:104:GLN:HG3	1.91	0.71
1:B:86:LYS:CB	1:B:87:PRO:HD2	2.22	0.70
1:B:86:LYS:HB3	1:B:87:PRO:CD	2.23	0.69
1:B:179:GLY:O	1:B:182:GLU:HG2	1.93	0.69
1:B:205:ARG:HG2	1:B:205:ARG:HH11	1.58	0.69
1:B:92:LEU:H	1:B:92:LEU:HD12	1.58	0.68
1:A:76:LEU:CD2	1:A:117:ILE:HB	2.25	0.67
1:A:86:LYS:O	1:A:89:GLN:HB3	1.95	0.66
1:D:90:THR:HG22	1:D:92:LEU:H	1.61	0.65
1:C:106:VAL:O	1:C:110:MET:HG2	1.96	0.65
1:C:59:LYS:HD3	5:C:516:HOH:O	1.95	0.65
1:A:310:LEU:HD22	1:A:314:ARG:HB3	1.79	0.65
1:C:269:ASP:O	1:C:270:ILE:HG23	1.96	0.65
1:D:158:GLU:HB3	1:D:214:LEU:HD22	1.79	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:311:ASP:HA	1:A:314:ARG:HD2	1.78	0.65
1:D:179:GLY:O	1:D:182:GLU:HG2	1.97	0.65
1:A:267:GLU:HG3	1:A:303:VAL:HG21	1.79	0.64
1:A:179:GLY:O	1:A:182:GLU:HG2	1.97	0.64
1:B:250:ASN:OD1	1:B:279:ASP:HB2	1.96	0.64
1:D:205:ARG:HG2	1:D:205:ARG:HH11	1.63	0.64
1:B:270:ILE:HD11	1:B:306:LEU:HD22	1.80	0.64
1:A:123:ASN:HB2	2:A:401:NAD:O2D	1.97	0.64
1:D:270:ILE:HD11	1:D:306:LEU:HD22	1.80	0.63
1:C:102:THR:O	1:C:106:VAL:HG23	1.99	0.62
1:D:85:GLN:CG	1:D:94:LEU:HD22	2.30	0.62
1:D:181:SER:O	1:D:298:MET:HE2	1.99	0.62
1:D:102:THR:O	1:D:106:VAL:HG23	2.00	0.61
1:D:250:ASN:OD1	1:D:279:ASP:HB2	2.00	0.61
1:A:301:GLN:O	1:A:305:GLU:HG3	2.02	0.60
1:D:180:ASP:HA	1:D:219:LYS:CD	2.30	0.60
1:A:76:LEU:HD21	1:A:247:ILE:HG21	1.84	0.59
1:D:201:GLU:O	1:D:202:LYS:HG3	2.03	0.59
1:C:150:LEU:O	1:C:153:THR:HG22	2.02	0.59
1:C:156:ARG:NH1	5:C:521:HOH:O	2.36	0.58
1:B:293:PRO:HG2	5:B:519:HOH:O	2.04	0.58
1:D:310:LEU:HG	1:D:314:ARG:HG3	1.85	0.58
1:B:250:ASN:HB2	1:B:280:GLU:HB3	1.85	0.58
1:B:158:GLU:HB3	1:B:214:LEU:HD22	1.86	0.58
1:A:156:ARG:NH1	5:A:550:HOH:O	2.36	0.57
1:B:224:GLU:HG3	1:B:228:ARG:NH1	2.19	0.57
1:B:89:GLN:HA	1:B:89:GLN:HE21	1.69	0.57
1:B:92:LEU:O	1:B:96:LYS:HG3	2.04	0.57
1:B:67:THR:OG1	1:B:69:GLU:HG3	2.05	0.57
1:B:224:GLU:HG3	1:B:228:ARG:HH12	1.69	0.57
1:A:131:LEU:HD21	1:A:314:ARG:HA	1.87	0.57
1:C:179:GLY:O	1:C:182:GLU:HG2	2.04	0.57
1:A:194:LYS:NZ	1:A:198:GLU:OE2	2.37	0.56
1:D:190:THR:HG22	1:D:195:PRO:HA	1.87	0.56
1:D:254:VAL:HG22	1:D:277:ILE:HD13	1.87	0.56
1:B:264:GLU:HA	1:B:288:ASP:OD2	2.05	0.56
1:C:180:ASP:HA	1:C:219:LYS:HD2	1.86	0.56
1:A:28:GLY:HA3	1:A:59:LYS:HD3	1.87	0.56
1:B:139:PRO:HD2	1:B:142:ARG:HD3	1.87	0.56
1:B:86:LYS:CB	1:B:87:PRO:CD	2.84	0.55
1:C:270:ILE:O	1:C:270:ILE:HD12	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:226:ILE:HG12	1:D:231:ALA:HA	1.88	0.55
3:B:404:GOL:H31	1:C:167:PRO:HD3	1.87	0.55
1:C:124:PRO:HG2	1:C:127:ILE:HB	1.88	0.55
3:B:404:GOL:H32	1:C:165:VAL:O	2.07	0.55
1:B:196:ILE:O	1:B:199:TYR:HB2	2.08	0.54
1:D:224:GLU:HG3	1:D:228:ARG:HH12	1.71	0.54
1:A:124:PRO:HG2	1:A:127:ILE:HB	1.90	0.54
1:A:234:TYR:CD2	1:B:50:LEU:HD13	2.43	0.54
1:A:99:ALA:O	1:A:103:ARG:HB2	2.08	0.54
1:D:139:PRO:HD2	1:D:142:ARG:HG3	1.89	0.54
1:C:134:GLN:OE1	1:C:315:LEU:HG	2.07	0.54
1:A:250:ASN:OD1	1:A:279:ASP:HB2	2.08	0.54
1:A:96:LYS:O	1:A:100:THR:HG23	2.08	0.53
1:D:11:VAL:HG11	1:D:105:ILE:HD13	1.89	0.53
1:A:150:LEU:O	1:A:153:THR:HG22	2.08	0.53
1:B:59:LYS:HG3	5:B:511:HOH:O	2.08	0.53
1:B:102:THR:O	1:B:106:VAL:HG23	2.09	0.53
1:B:261:LEU:HD23	1:B:286:ILE:HD13	1.91	0.53
1:C:151:ASP:OD2	1:C:178:HIS:ND1	2.38	0.53
1:A:197:MET:HE1	1:A:211:LEU:HD11	1.92	0.52
1:D:106:VAL:O	1:D:110:MET:HG2	2.08	0.52
1:B:310:LEU:HD12	1:B:314:ARG:HD2	1.90	0.52
1:A:90:THR:HG22	1:A:92:LEU:N	2.04	0.52
1:C:303:VAL:O	1:C:307:LYS:HG3	2.09	0.52
1:B:200:VAL:O	1:B:204:HIS:HA	2.09	0.52
1:C:267:GLU:O	1:C:268:GLU:HG3	2.09	0.52
1:A:122:SER:O	1:A:128:LEU:HD12	2.10	0.51
1:B:124:PRO:HG2	1:B:127:ILE:HB	1.91	0.51
1:B:247:ILE:O	1:B:280:GLU:HA	2.09	0.51
1:D:14:GLY:HA3	2:D:402:NAD:O5B	2.10	0.51
1:B:205:ARG:NH1	1:B:205:ARG:HG2	2.23	0.51
1:B:123:ASN:HA	1:B:125:VAL:N	2.24	0.51
1:C:156:ARG:HD3	5:C:521:HOH:O	2.08	0.51
1:C:183:VAL:HG22	1:C:298:MET:HG3	1.92	0.51
1:A:260:TYR:CZ	1:A:269:ASP:HA	2.46	0.51
1:C:11:VAL:HG11	1:C:105:ILE:HD13	1.93	0.51
1:B:223:TYR:O	1:B:227:ASP:HB2	2.11	0.51
1:A:251:GLU:O	1:A:252:GLN:HB2	2.11	0.51
1:B:15:PHE:HB3	1:B:233:TYR:CD2	2.46	0.50
1:C:250:ASN:OD1	1:C:279:ASP:HB2	2.11	0.50
1:C:50:LEU:HD13	1:D:234:TYR:CD2	2.46	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:157:LYS:HD2	1:B:51:LEU:HD23	1.94	0.50
1:D:225:ILE:CG2	1:D:232:THR:HG22	2.41	0.50
1:B:224:GLU:O	1:B:228:ARG:NH1	2.44	0.49
1:D:223:TYR:O	1:D:227:ASP:HB2	2.12	0.49
1:A:153:THR:HB	1:B:56:TRP:NE1	2.28	0.49
1:C:209:ASN:O	1:C:213:VAL:HG23	2.12	0.49
1:B:99:ALA:HB1	1:B:313:VAL:HG11	1.95	0.49
1:B:245:LYS:HE3	5:B:572:HOH:O	2.13	0.49
1:A:250:ASN:OD1	1:A:280:GLU:HG2	2.13	0.48
1:D:86:LYS:HG2	1:D:89:GLN:HG2	1.94	0.48
1:B:180:ASP:HA	1:B:219:LYS:HD3	1.95	0.48
1:A:228:ARG:HD3	1:B:44:GLU:HB3	1.94	0.48
1:D:233:TYR:O	1:D:236:ILE:HG22	2.13	0.48
1:D:179:GLY:C	1:D:181:SER:H	2.16	0.48
1:C:251:GLU:O	1:C:252:GLN:HB2	2.14	0.48
1:C:292:THR:OG1	1:C:295:GLU:HG3	2.14	0.48
1:B:129:THR:HG23	1:B:143:VAL:HG12	1.95	0.48
1:B:99:ALA:O	1:B:103:ARG:HG3	2.14	0.47
1:D:124:PRO:O	1:D:128:LEU:HG	2.14	0.47
1:D:119:VAL:HA	1:D:144:VAL:O	2.14	0.47
1:C:76:LEU:HD12	1:C:76:LEU:C	2.35	0.47
1:D:224:GLU:HG2	5:D:545:HOH:O	2.13	0.47
1:A:119:VAL:HA	1:A:144:VAL:O	2.13	0.47
1:C:254:VAL:HA	1:C:276:SER:O	2.14	0.47
1:B:180:ASP:HA	1:B:219:LYS:CD	2.44	0.47
1:B:99:ALA:CB	1:B:313:VAL:HG11	2.45	0.47
1:B:233:TYR:HA	1:B:236:ILE:HG22	1.95	0.47
1:D:185:ALA:HB3	5:D:504:HOH:O	2.15	0.47
1:B:180:ASP:HA	1:B:219:LYS:CE	2.45	0.47
1:D:42:LYS:HE3	1:D:46:GLU:OE2	2.15	0.47
1:B:161:LEU:HD23	3:B:403:GOL:H31	1.97	0.46
1:A:158:GLU:HB3	1:A:214:LEU:HD22	1.97	0.46
1:D:205:ARG:HG2	1:D:205:ARG:NH1	2.30	0.46
1:D:90:THR:HG22	1:D:91:ARG:N	2.30	0.46
1:B:280:GLU:OE1	1:D:3:LYS:HA	2.15	0.46
1:B:183:VAL:HG22	1:B:298:MET:HG3	1.98	0.46
1:C:225:ILE:CG2	1:C:232:THR:HG22	2.45	0.46
1:D:15:PHE:HB3	1:D:233:TYR:CD2	2.51	0.46
1:C:3:LYS:N	5:C:562:HOH:O	2.49	0.45
1:C:39:ASN:CG	1:C:42:LYS:HB2	2.36	0.45
1:D:84:ASN:ND2	1:D:85:GLN:N	2.56	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:34:VAL:HG22	1:B:64:TRP:CE2	2.51	0.45
1:A:292:THR:OG1	1:A:295:GLU:HG3	2.16	0.45
1:A:270:ILE:CG2	1:A:307:LYS:HE2	2.47	0.45
1:D:103:ARG:HD2	1:D:315:LEU:HD21	1.98	0.45
1:B:254:VAL:HA	1:B:276:SER:O	2.16	0.45
1:A:14:GLY:HA3	2:A:401:NAD:O5B	2.16	0.45
1:C:267:GLU:C	1:C:268:GLU:HG3	2.37	0.45
1:A:224:GLU:HG3	1:A:228:ARG:HH12	1.81	0.45
1:B:228:ARG:HG3	1:B:228:ARG:HH11	1.82	0.45
1:C:150:LEU:HD21	1:C:154:ARG:NH2	2.32	0.45
1:A:224:GLU:HG3	1:A:228:ARG:NH1	2.32	0.45
1:A:28:GLY:O	1:A:59:LYS:HD3	2.16	0.45
1:B:180:ASP:HA	1:B:219:LYS:HE2	1.99	0.45
1:C:92:LEU:HG	1:C:96:LYS:HE3	1.99	0.45
1:D:16:VAL:CG2	2:D:402:NAD:H51N	2.48	0.44
1:A:171:HIS:HE1	4:D:401[A]:FBP:O5	2.00	0.44
1:B:123:ASN:OD1	1:B:124:PRO:HA	2.17	0.44
1:A:234:TYR:OH	1:B:18:THR:HB	2.18	0.44
1:A:196:ILE:O	1:A:199:TYR:HB2	2.17	0.44
1:B:310:LEU:HD13	1:B:314:ARG:HB2	2.00	0.44
1:C:160:ALA:HB1	1:C:165:VAL:O	2.17	0.44
1:B:200:VAL:HA	1:B:206:LEU:O	2.16	0.44
1:D:151:ASP:OD2	1:D:178:HIS:ND1	2.48	0.44
1:B:313:VAL:HG13	5:B:604:HOH:O	2.17	0.44
1:B:89:GLN:HA	1:B:89:GLN:NE2	2.32	0.44
1:C:195:PRO:HG2	1:C:198:GLU:HB2	1.99	0.44
1:D:310:LEU:HG	1:D:314:ARG:CG	2.48	0.44
1:D:141:SER:O	1:D:282:GLY:HA3	2.17	0.44
1:A:270:ILE:HG23	1:A:307:LYS:HE2	1.99	0.44
1:A:51:LEU:HB2	1:A:63:VAL:HB	2.00	0.44
4:B:401[A]:FBP:O5	1:C:171:HIS:HE1	2.01	0.44
1:A:316:GLU:HB3	1:A:317:HIS:ND1	2.33	0.43
1:A:28:GLY:CA	1:A:59:LYS:HD3	2.49	0.43
1:B:3:LYS:N	1:B:3:LYS:HD2	2.33	0.43
1:C:76:LEU:HD21	1:C:244:VAL:HG13	2.00	0.43
1:C:153:THR:HB	1:D:56:TRP:NE1	2.33	0.43
1:D:59:LYS:HG3	5:D:530:HOH:O	2.17	0.43
1:C:264:GLU:HA	1:C:288:ASP:OD2	2.18	0.43
1:D:131:LEU:HD11	1:D:313:VAL:CG1	2.48	0.43
1:B:313:VAL:HG12	1:B:313:VAL:O	2.18	0.43
1:C:312:THR:HG22	1:C:312:THR:O	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:194:LYS:HD2	1:A:199:TYR:CZ	2.54	0.42
1:A:223:TYR:C	1:A:223:TYR:CD1	2.92	0.42
1:B:15:PHE:HB3	1:B:233:TYR:CG	2.54	0.42
1:B:119:VAL:HA	1:B:144:VAL:O	2.19	0.42
1:C:92:LEU:HD21	1:C:312:THR:HG21	2.01	0.42
1:C:56:TRP:NE1	1:D:153:THR:HB	2.34	0.42
1:D:254:VAL:HA	1:D:276:SER:O	2.20	0.42
1:B:123:ASN:HB2	2:B:402:NAD:H1D	2.01	0.42
1:B:155:PHE:O	1:B:159:ILE:HG13	2.19	0.42
1:A:277:ILE:HD13	1:D:165:VAL:HG12	2.02	0.42
1:B:306:LEU:HD23	1:B:306:LEU:O	2.19	0.42
1:D:46:GLU:OE1	1:D:46:GLU:HA	2.20	0.42
1:B:134:GLN:OE1	1:B:315:LEU:HG	2.20	0.42
1:C:234:TYR:CD2	1:D:50:LEU:HD13	2.54	0.42
1:C:315:LEU:HD12	1:C:316:GLU:O	2.19	0.42
1:D:123:ASN:HA	1:D:125:VAL:N	2.35	0.42
1:D:224:GLU:HG3	1:D:228:ARG:NH1	2.35	0.42
1:A:151:ASP:OD2	1:A:178:HIS:ND1	2.48	0.41
1:B:229:LYS:CD	1:B:231:ALA:O	2.68	0.41
1:C:301:GLN:O	1:C:305:GLU:HG3	2.20	0.41
2:B:402:NAD:H6N	2:B:402:NAD:H2D	1.66	0.41
1:B:91:ARG:O	1:B:95:VAL:HG23	2.21	0.41
1:D:292:THR:OG1	1:D:295:GLU:HG3	2.20	0.41
1:A:89:GLN:HG3	1:A:93:ASP:CB	2.50	0.41
1:A:106:VAL:O	1:A:110:MET:HG2	2.21	0.41
1:C:123:ASN:HB2	2:C:401:NAD:O2D	2.20	0.41
1:B:213:VAL:HG12	1:B:217:LYS:HE2	2.02	0.41
1:B:260:TYR:CZ	1:B:269:ASP:HA	2.54	0.41
1:B:86:LYS:HB2	1:B:89:GLN:HG2	2.00	0.41
1:C:174:ILE:HD13	1:C:184:ALA:HA	2.03	0.41
1:D:219:LYS:HE3	1:D:219:LYS:HB2	1.94	0.41
1:C:82:GLY:O	2:C:401:NAD:H8A	2.20	0.41
1:D:11:VAL:CG1	1:D:105:ILE:HD13	2.51	0.41
1:A:261:LEU:HD23	1:A:286:ILE:HD13	2.02	0.41
1:D:46:GLU:O	1:D:50:LEU:HB2	2.21	0.41
1:C:15:PHE:HB3	1:C:233:TYR:CD2	2.56	0.40
1:D:144:VAL:HG22	1:D:283:VAL:HG23	2.03	0.40
1:D:271:PHE:HB2	5:D:516:HOH:O	2.21	0.40
1:D:86:LYS:HG2	1:D:89:GLN:CD	2.41	0.40
1:D:123:ASN:ND2	2:D:402:NAD:C2N	2.85	0.40
1:A:311:ASP:HA	1:A:314:ARG:CD	2.48	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:56:TRP:NE1	1:B:153:THR:HB	2.36	0.40
1:D:131:LEU:HD11	1:D:313:VAL:HG12	2.03	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	314/322 (98%)	299 (95%)	14 (4%)	1 (0%)	41 50
1	B	311/322 (97%)	293 (94%)	14 (4%)	4 (1%)	12 12
1	C	312/322 (97%)	296 (95%)	15 (5%)	1 (0%)	41 50
1	D	312/322 (97%)	299 (96%)	12 (4%)	1 (0%)	41 50
All	All	1249/1288 (97%)	1187 (95%)	55 (4%)	7 (1%)	25 31

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	86	LYS
1	B	87	PRO
1	B	185	ALA
1	C	200	VAL
1	D	233	TYR
1	B	233	TYR
1	A	313	VAL

### 5.3.2 Protein sidechains [i](#)

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



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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	265/271 (98%)	259 (98%)	6 (2%)	50	67
1	B	262/271 (97%)	258 (98%)	4 (2%)	65	79
1	C	263/271 (97%)	261 (99%)	2 (1%)	81	91
1	D	263/271 (97%)	258 (98%)	5 (2%)	57	73
All	All	1053/1084 (97%)	1036 (98%)	17 (2%)	62	78

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	40	GLN
1	A	205	ARG
1	A	208	GLU
1	A	270	ILE
1	A	306	LEU
1	A	310	LEU
1	B	44	GLU
1	B	50	LEU
1	B	168	ARG
1	B	262	ASN
1	C	50	LEU
1	C	310	LEU
1	D	42	LYS
1	D	50	LEU
1	D	84	ASN
1	D	227	ASP
1	D	306	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	84	ASN
1	A	104	GLN
1	A	252	GLN
1	A	294	GLN
1	B	84	ASN
1	B	89	GLN
1	B	104	GLN
1	B	262	ASN

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Mol	Chain	Res	Type
1	B	294	GLN
1	B	301	GLN
1	C	40	GLN
1	C	84	ASN
1	C	252	GLN
1	C	294	GLN
1	D	72	GLN
1	D	84	ASN
1	D	209	ASN
1	D	294	GLN

### 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 monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

11 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	B	403	-	5,5,5	0.38	0	5,5,5	0.22	0
2	NAD	D	402	-	42,48,48	1.20	5 (11%)	50,73,73	1.52	8 (16%)
4	FBP	B	401[A]	-	18,20,20	0.85	1 (5%)	23,32,32	0.97	2 (8%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	NAD	C	401	-	42,48,48	1.16	5 (11%)	50,73,73	1.51	7 (14%)
4	FBP	B	401[B]	-	18,20,20	0.89	1 (5%)	23,32,32	0.97	1 (4%)
3	GOL	B	404	-	5,5,5	0.40	0	5,5,5	0.32	0
3	GOL	A	402	-	5,5,5	0.51	0	5,5,5	0.24	0
4	FBP	D	401[A]	-	18,20,20	0.88	1 (5%)	23,32,32	0.98	1 (4%)
2	NAD	B	402	-	42,48,48	1.29	5 (11%)	50,73,73	1.54	9 (18%)
2	NAD	A	401	-	42,48,48	1.17	5 (11%)	50,73,73	1.52	8 (16%)
4	FBP	D	401[B]	-	18,20,20	0.91	1 (5%)	23,32,32	0.99	2 (8%)

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	B	403	-	-	2/4/4/4	-
2	NAD	D	402	-	-	9/26/62/62	0/5/5/5
4	FBP	B	401[A]	-	-	4/13/32/32	0/1/1/1
2	NAD	C	401	-	-	12/26/62/62	0/5/5/5
4	FBP	B	401[B]	-	-	11/13/32/32	0/1/1/1
3	GOL	B	404	-	-	2/4/4/4	-
3	GOL	A	402	-	-	0/4/4/4	-
4	FBP	D	401[A]	-	-	4/13/32/32	0/1/1/1
2	NAD	B	402	-	-	10/26/62/62	0/5/5/5
2	NAD	A	401	-	-	11/26/62/62	0/5/5/5
4	FBP	D	401[B]	-	-	8/13/32/32	0/1/1/1

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	402	NAD	O4D-C1D	4.44	1.47	1.41
2	D	402	NAD	O4D-C1D	3.23	1.45	1.41
2	D	402	NAD	O4B-C1B	3.00	1.45	1.41
2	A	401	NAD	O4B-C1B	2.97	1.45	1.41
2	C	401	NAD	O4D-C1D	2.87	1.45	1.41
2	B	402	NAD	O4B-C1B	2.83	1.45	1.41
2	A	401	NAD	O4D-C1D	2.83	1.45	1.41
2	C	401	NAD	O4B-C1B	2.79	1.45	1.41
2	A	401	NAD	C4N-C3N	2.76	1.44	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	402	NAD	C4N-C3N	2.58	1.43	1.39
2	C	401	NAD	C4N-C3N	2.54	1.43	1.39
2	B	402	NAD	C4N-C3N	2.51	1.43	1.39
2	A	401	NAD	C8A-N7A	-2.42	1.30	1.34
2	B	402	NAD	C2N-C3N	2.40	1.42	1.39
2	C	401	NAD	C8A-N7A	-2.23	1.30	1.34
2	D	402	NAD	C2N-C3N	2.23	1.42	1.39
4	D	401[B]	FBP	O1-C1	2.13	1.50	1.43
4	B	401[A]	FBP	O1-C1	2.12	1.50	1.43
4	D	401[A]	FBP	O1-C1	2.10	1.50	1.43
2	B	402	NAD	C8A-N7A	-2.09	1.31	1.34
2	C	401	NAD	C2N-C3N	2.09	1.42	1.39
4	B	401[B]	FBP	O1-C1	2.05	1.50	1.43
2	A	401	NAD	C2N-C3N	2.02	1.42	1.39
2	D	402	NAD	C8A-N7A	-2.02	1.31	1.34

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	402	NAD	PN-O3-PA	-6.49	110.54	132.83
2	B	402	NAD	PN-O3-PA	-6.45	110.69	132.83
2	C	401	NAD	PN-O3-PA	-6.35	111.02	132.83
2	A	401	NAD	PN-O3-PA	-6.06	112.02	132.83
2	A	401	NAD	N3A-C2A-N1A	-4.67	121.37	128.68
2	B	402	NAD	N3A-C2A-N1A	-4.65	121.42	128.68
2	C	401	NAD	N3A-C2A-N1A	-4.63	121.44	128.68
2	D	402	NAD	N3A-C2A-N1A	-4.62	121.46	128.68
2	B	402	NAD	C3D-C2D-C1D	2.92	105.37	100.98
2	A	401	NAD	O4B-C1B-C2B	-2.74	102.92	106.93
2	B	402	NAD	O4B-C1B-C2B	-2.55	103.20	106.93
2	D	402	NAD	O4B-C1B-C2B	-2.54	103.22	106.93
2	A	401	NAD	O7N-C7N-N7N	-2.49	119.04	122.58
2	D	402	NAD	O7N-C7N-N7N	-2.37	119.20	122.58
2	C	401	NAD	C4A-C5A-N7A	-2.35	106.95	109.40
2	D	402	NAD	PN-O5D-C5D	-2.34	107.95	121.68
2	A	401	NAD	C4A-C5A-N7A	-2.34	106.96	109.40
2	B	402	NAD	C3N-C7N-N7N	2.32	120.53	117.75
2	C	401	NAD	PA-O5B-C5B	-2.31	108.11	121.68
4	D	401[B]	FBP	P1-O1-C1	-2.31	111.94	118.30
4	D	401[A]	FBP	P1-O1-C1	-2.28	112.02	118.30
2	B	402	NAD	O7N-C7N-N7N	-2.28	119.34	122.58
2	A	401	NAD	PN-O5D-C5D	-2.26	108.44	121.68

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	NAD	C3N-C7N-N7N	2.24	120.44	117.75
2	C	401	NAD	O7N-C7N-N7N	-2.21	119.43	122.58
2	B	402	NAD	PA-O5B-C5B	-2.19	108.84	121.68
4	B	401[B]	FBP	P1-O1-C1	-2.18	112.30	118.30
2	A	401	NAD	PA-O5B-C5B	-2.16	109.04	121.68
2	D	402	NAD	C4A-C5A-N7A	-2.14	107.17	109.40
2	D	402	NAD	C3N-C7N-N7N	2.13	120.31	117.75
2	D	402	NAD	PA-O5B-C5B	-2.12	109.25	121.68
4	D	401[B]	FBP	O6-P2-O4P	2.12	112.42	106.47
2	C	401	NAD	PN-O5D-C5D	-2.12	109.26	121.68
2	B	402	NAD	PN-O5D-C5D	-2.09	109.40	121.68
2	C	401	NAD	C3N-C7N-N7N	2.04	120.19	117.75
4	B	401[A]	FBP	P1-O1-C1	-2.03	112.71	118.30
2	B	402	NAD	C4A-C5A-N7A	-2.02	107.30	109.40
4	B	401[A]	FBP	P2-O6-C6	-2.01	112.76	118.30

There are no chirality outliers.

All (73) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	403	GOL	O1-C1-C2-C3
2	D	402	NAD	O4B-C4B-C5B-O5B
2	D	402	NAD	O4D-C1D-N1N-C2N
2	D	402	NAD	O4D-C1D-N1N-C6N
2	D	402	NAD	C2D-C1D-N1N-C2N
2	D	402	NAD	C2D-C1D-N1N-C6N
4	B	401[A]	FBP	C6-O6-P2-O6P
2	C	401	NAD	C5B-O5B-PA-O1A
2	C	401	NAD	C5D-O5D-PN-O3
2	C	401	NAD	O4D-C1D-N1N-C6N
4	B	401[B]	FBP	C1-O1-P1-O3P
4	B	401[B]	FBP	O1-C1-C2-O2
4	B	401[B]	FBP	O1-C1-C2-C3
4	B	401[B]	FBP	O1-C1-C2-O5
4	B	401[B]	FBP	C4-C5-C6-O6
4	B	401[B]	FBP	C6-O6-P2-O5P
4	B	401[B]	FBP	C6-O6-P2-O6P
4	D	401[A]	FBP	C4-C5-C6-O6
4	D	401[A]	FBP	C6-O6-P2-O6P
2	B	402	NAD	O4B-C4B-C5B-O5B
2	B	402	NAD	C5D-O5D-PN-O3
2	B	402	NAD	C5D-O5D-PN-O1N

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Mol	Chain	Res	Type	Atoms
2	B	402	NAD	C5D-O5D-PN-O2N
2	B	402	NAD	O4D-C1D-N1N-C2N
2	B	402	NAD	C2D-C1D-N1N-C2N
2	A	401	NAD	PN-O3-PA-O5B
2	A	401	NAD	O4D-C1D-N1N-C2N
2	A	401	NAD	O4D-C1D-N1N-C6N
4	D	401[B]	FBP	O1-C1-C2-O2
4	D	401[B]	FBP	O1-C1-C2-C3
4	D	401[B]	FBP	O1-C1-C2-O5
4	D	401[B]	FBP	C4-C5-C6-O6
4	D	401[B]	FBP	O5-C5-C6-O6
4	D	401[B]	FBP	C6-O6-P2-O4P
4	D	401[B]	FBP	C6-O6-P2-O5P
4	D	401[B]	FBP	C6-O6-P2-O6P
2	D	402	NAD	C3B-C4B-C5B-O5B
2	D	402	NAD	O4D-C4D-C5D-O5D
2	C	401	NAD	O4B-C4B-C5B-O5B
2	C	401	NAD	C3B-C4B-C5B-O5B
2	A	401	NAD	O4B-C4B-C5B-O5B
2	A	401	NAD	O4D-C4D-C5D-O5D
4	B	401[A]	FBP	C4-C5-C6-O6
4	B	401[B]	FBP	O5-C5-C6-O6
2	B	402	NAD	C3B-C4B-C5B-O5B
2	C	401	NAD	C3D-C4D-C5D-O5D
3	B	404	GOL	O1-C1-C2-C3
3	B	404	GOL	O1-C1-C2-O2
2	C	401	NAD	O4D-C4D-C5D-O5D
4	D	401[A]	FBP	O5-C5-C6-O6
2	A	401	NAD	C3B-C4B-C5B-O5B
4	B	401[B]	FBP	C6-O6-P2-O4P
2	D	402	NAD	C3D-C4D-C5D-O5D
2	D	402	NAD	PN-O3-PA-O5B
2	C	401	NAD	PN-O3-PA-O5B
2	C	401	NAD	PA-O3-PN-O5D
2	B	402	NAD	PN-O3-PA-O5B
4	B	401[A]	FBP	O5-C5-C6-O6
2	A	401	NAD	C3D-C4D-C5D-O5D
4	B	401[B]	FBP	C1-O1-P1-O2P
3	B	403	GOL	O1-C1-C2-O2
2	C	401	NAD	C5D-O5D-PN-O1N
2	C	401	NAD	C5D-O5D-PN-O2N
4	B	401[B]	FBP	C1-O1-P1-O1P

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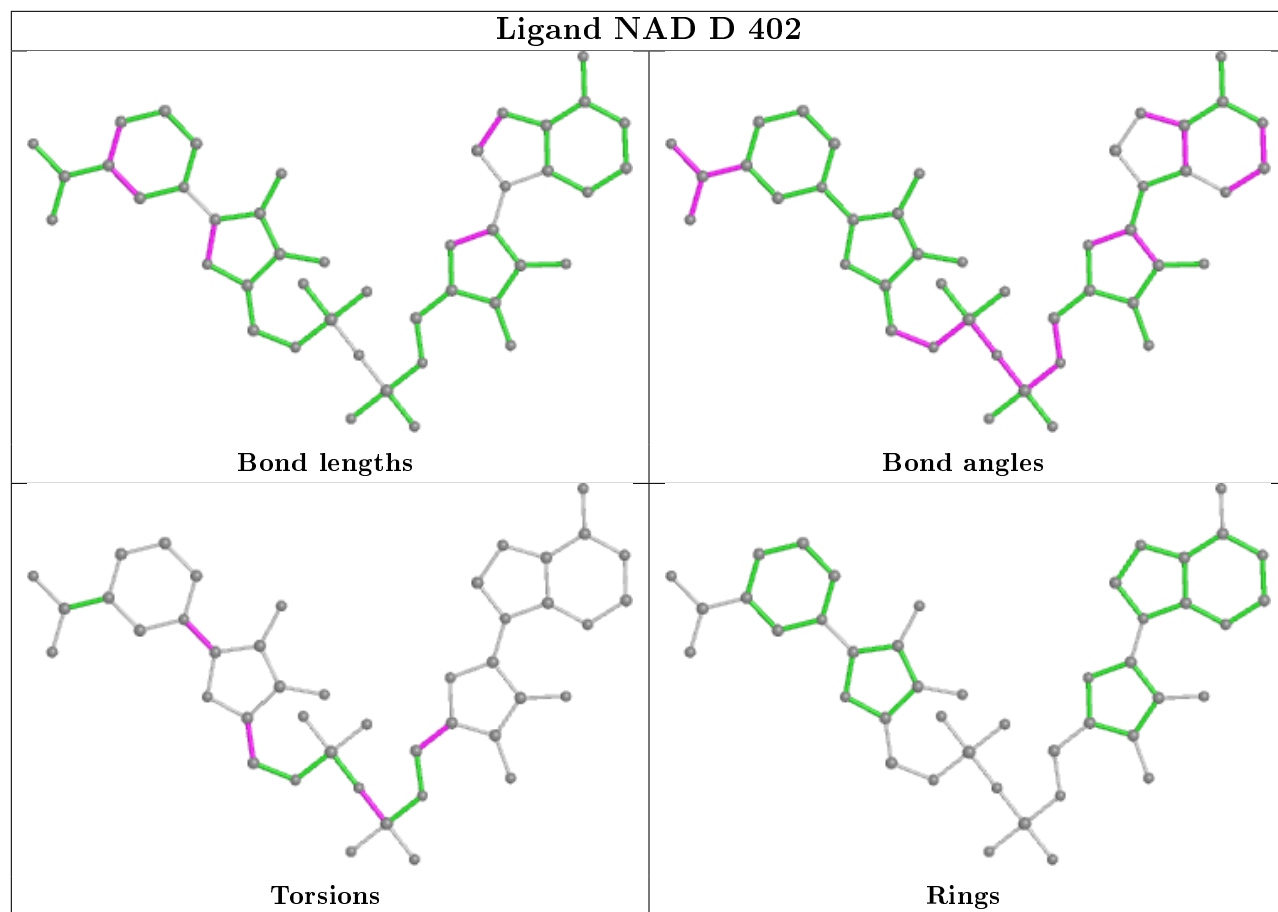
Mol	Chain	Res	Type	Atoms
2	A	401	NAD	PA-O3-PN-O2N
4	B	401[A]	FBP	O1-C1-C2-C3
4	D	401[A]	FBP	O1-C1-C2-C3
2	C	401	NAD	C5B-O5B-PA-O3
2	B	402	NAD	C2D-C1D-N1N-C6N
2	A	401	NAD	C2D-C1D-N1N-C2N
2	A	401	NAD	C2D-C1D-N1N-C6N
2	B	402	NAD	O4D-C4D-C5D-O5D
2	A	401	NAD	PA-O3-PN-O1N

There are no ring outliers.

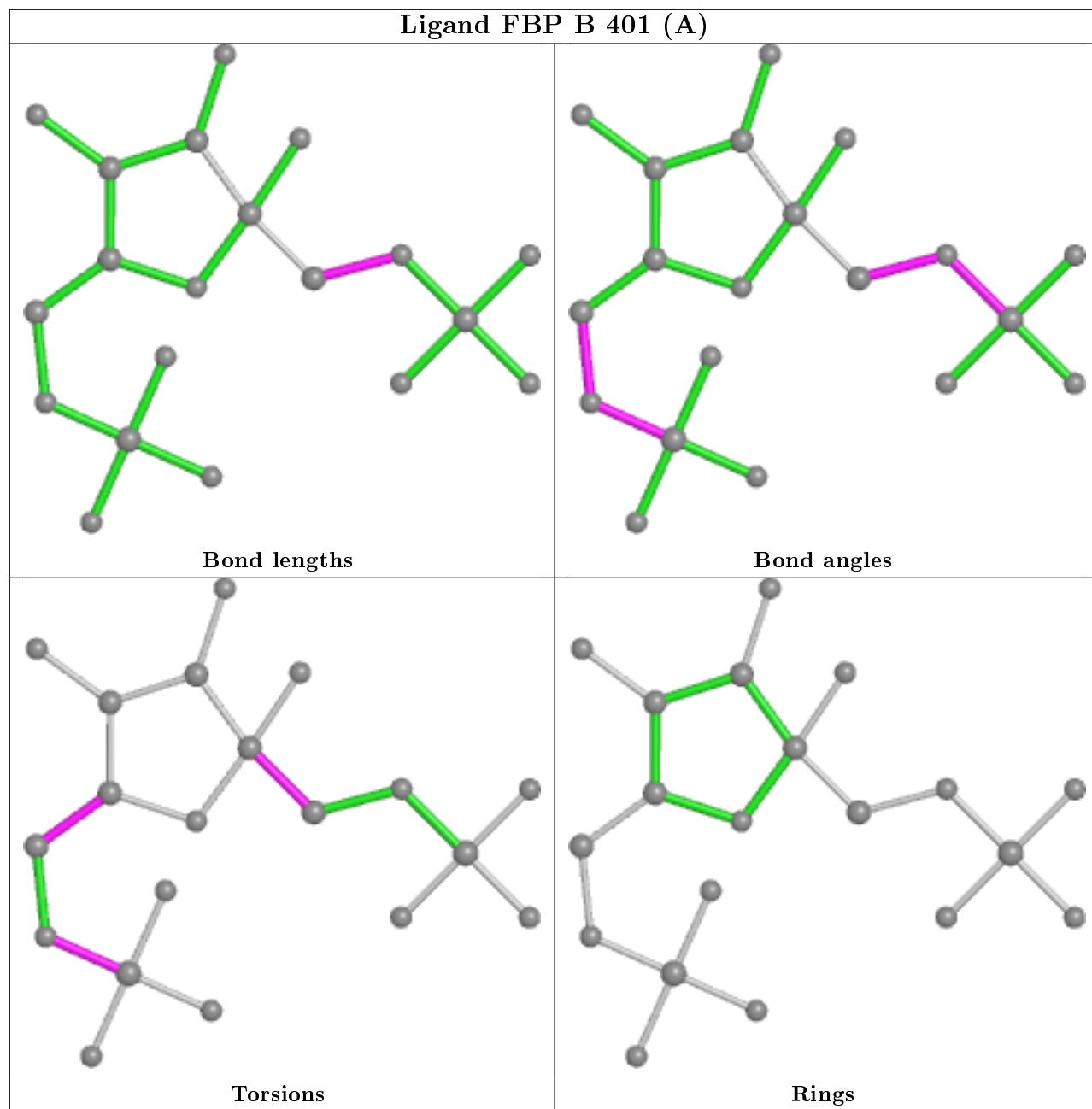
8 monomers are involved in 14 short contacts:

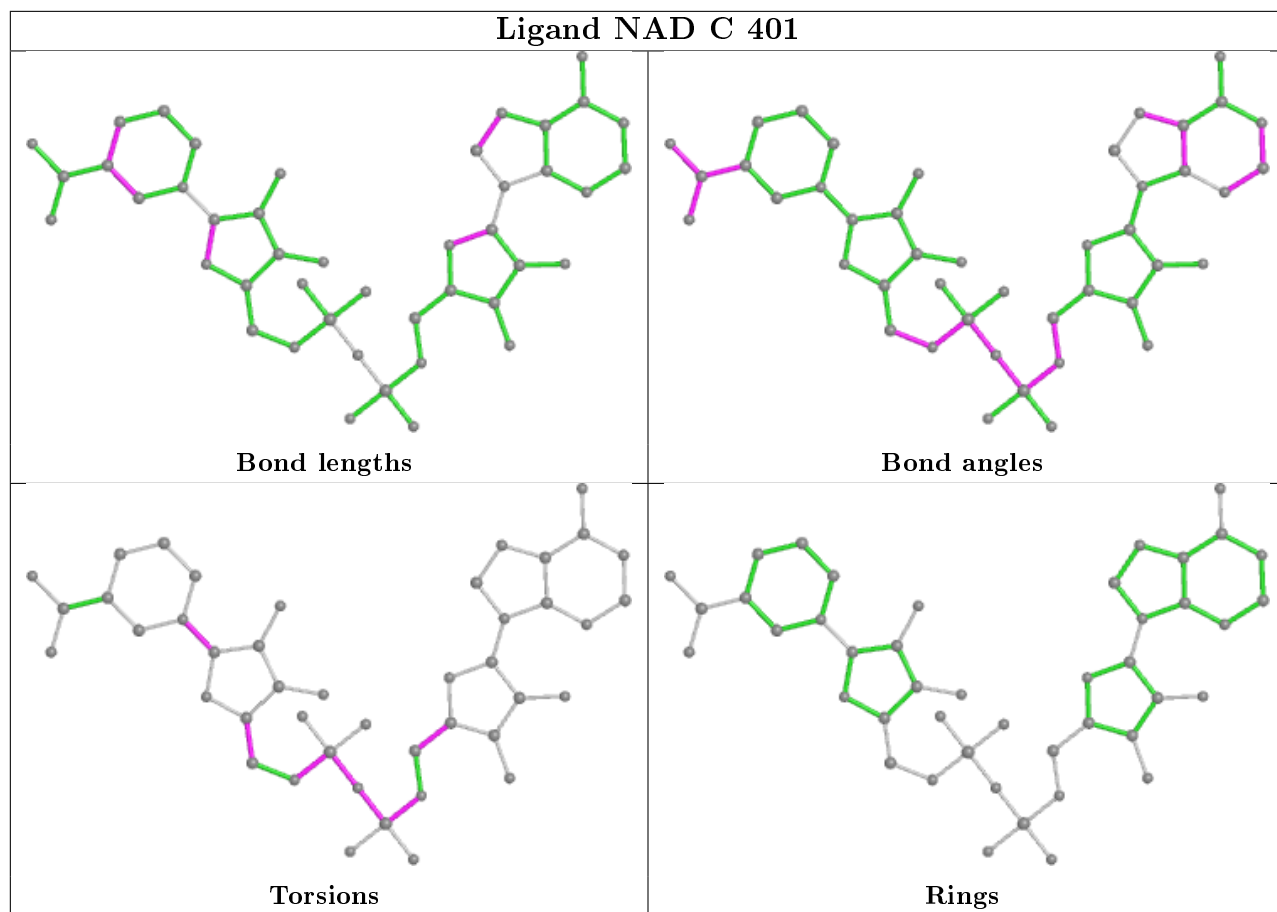
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	403	GOL	1	0
2	D	402	NAD	3	0
4	B	401[A]	FBP	1	0
2	C	401	NAD	2	0
3	B	404	GOL	2	0
4	D	401[A]	FBP	1	0
2	B	402	NAD	2	0
2	A	401	NAD	2	0

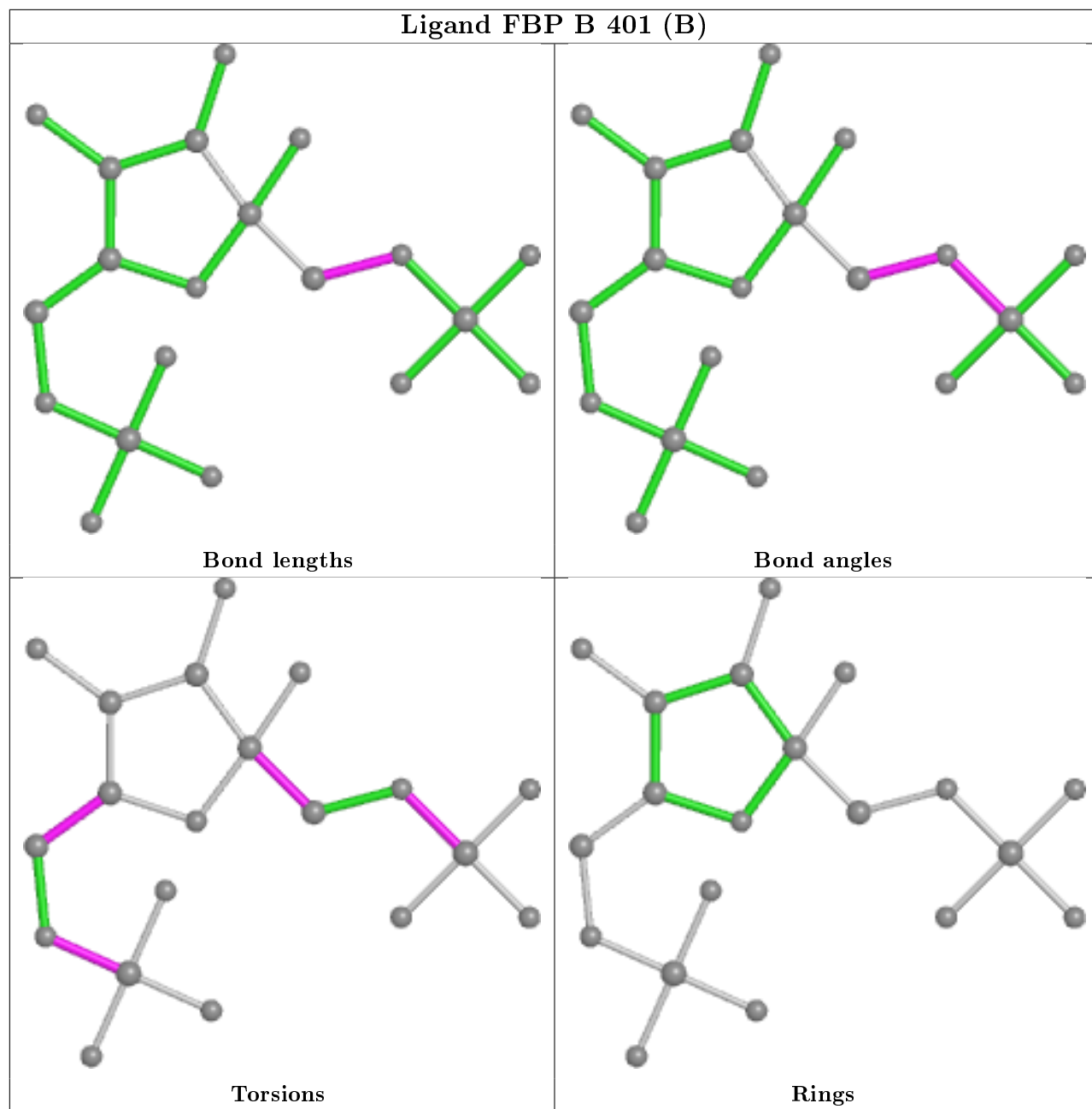
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

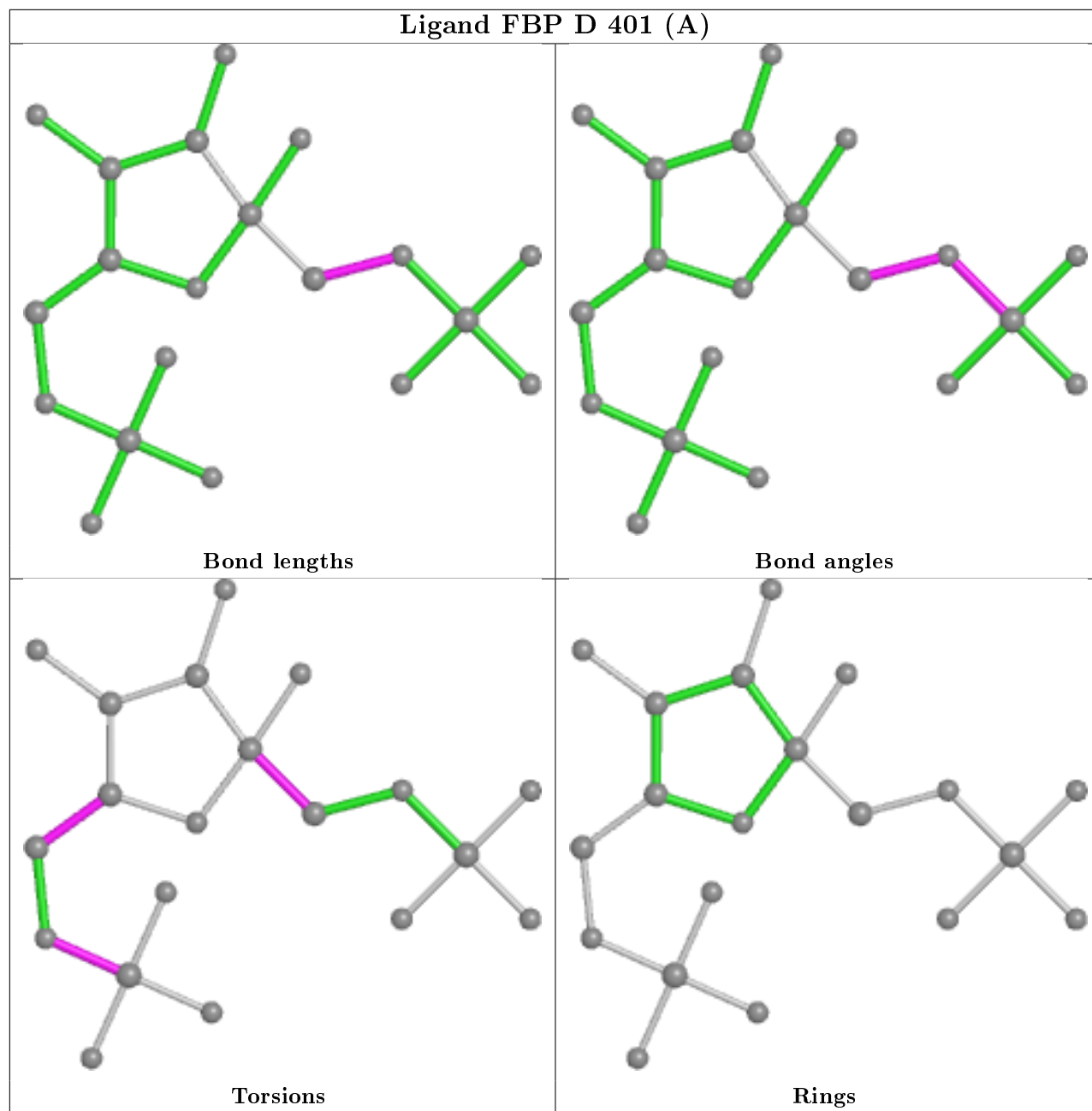


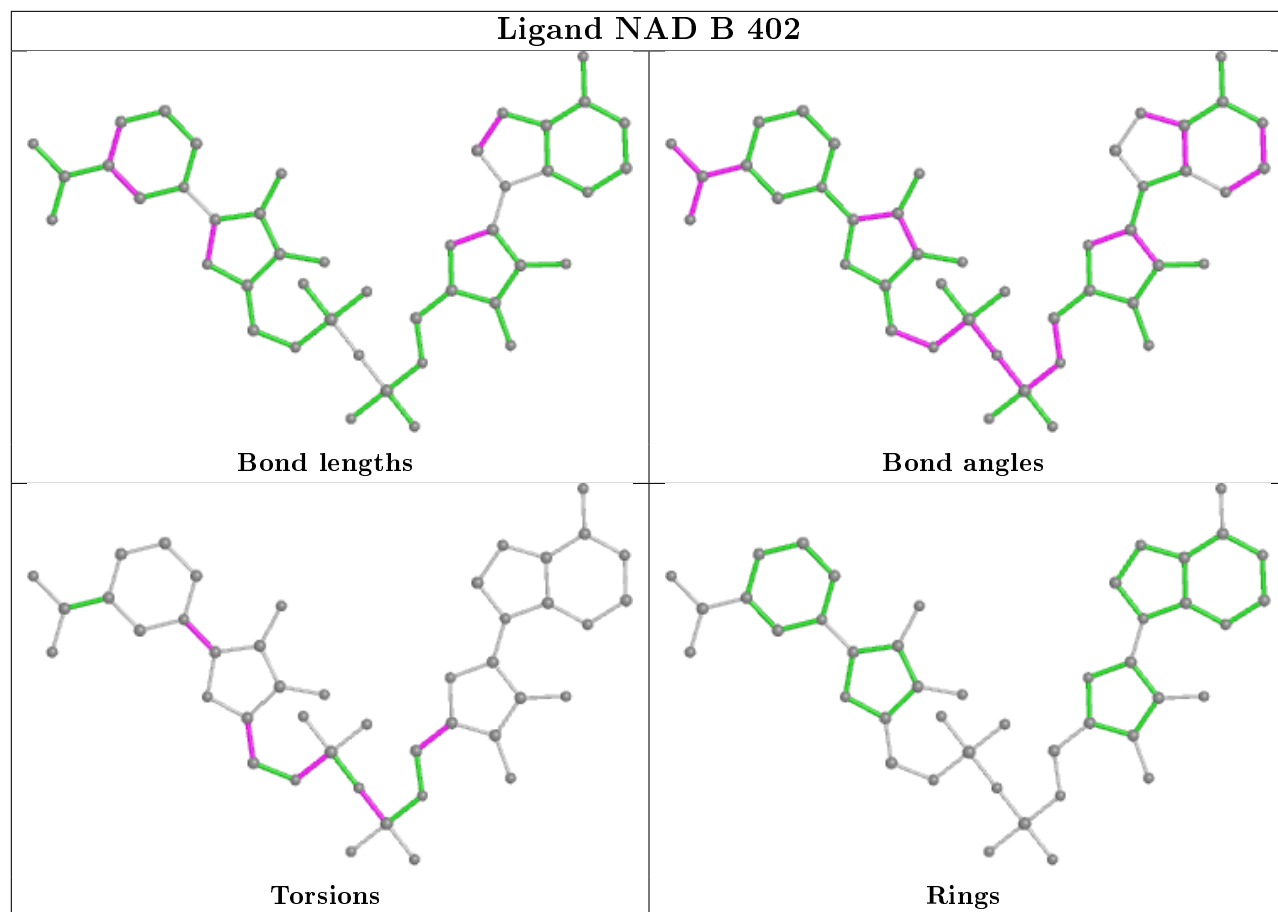


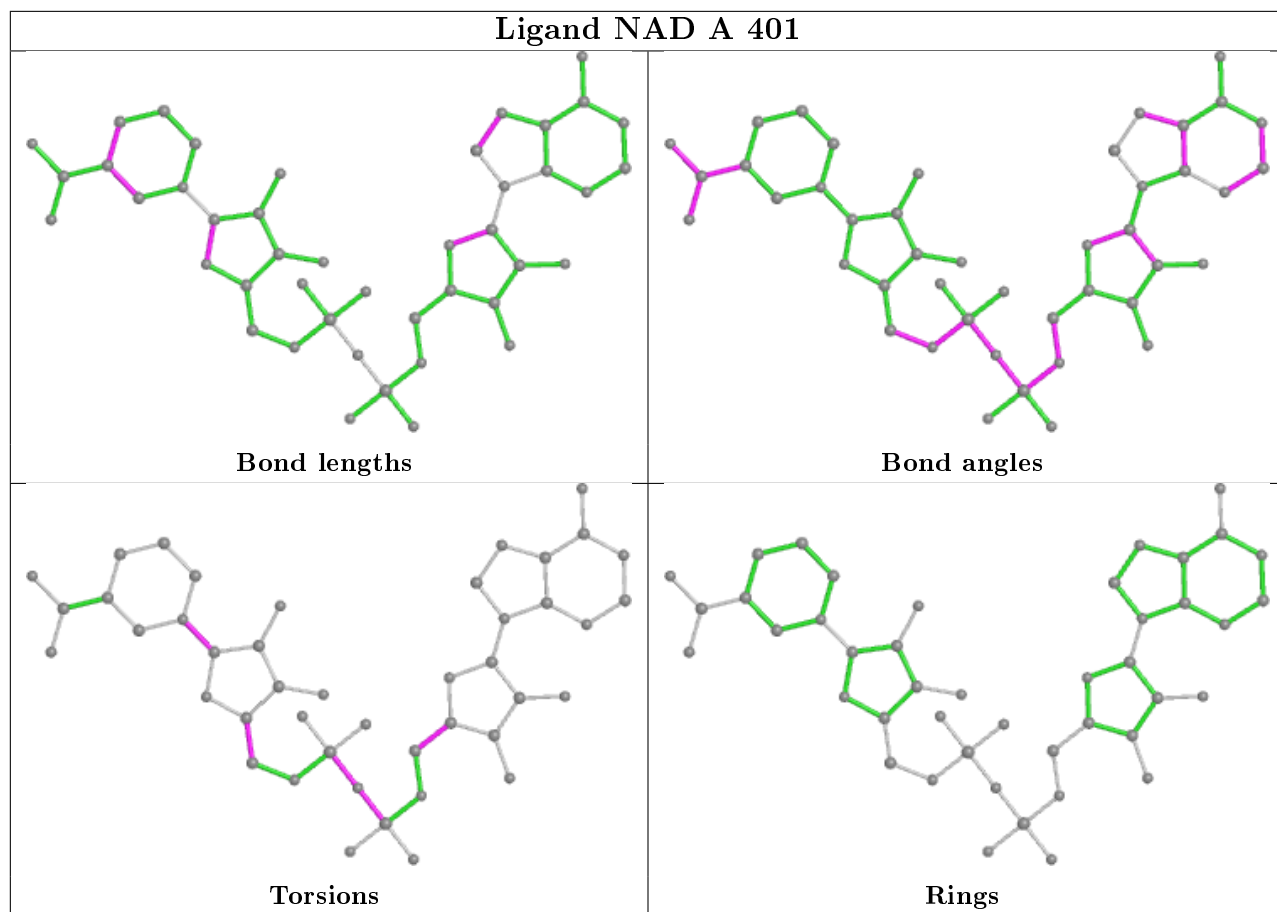


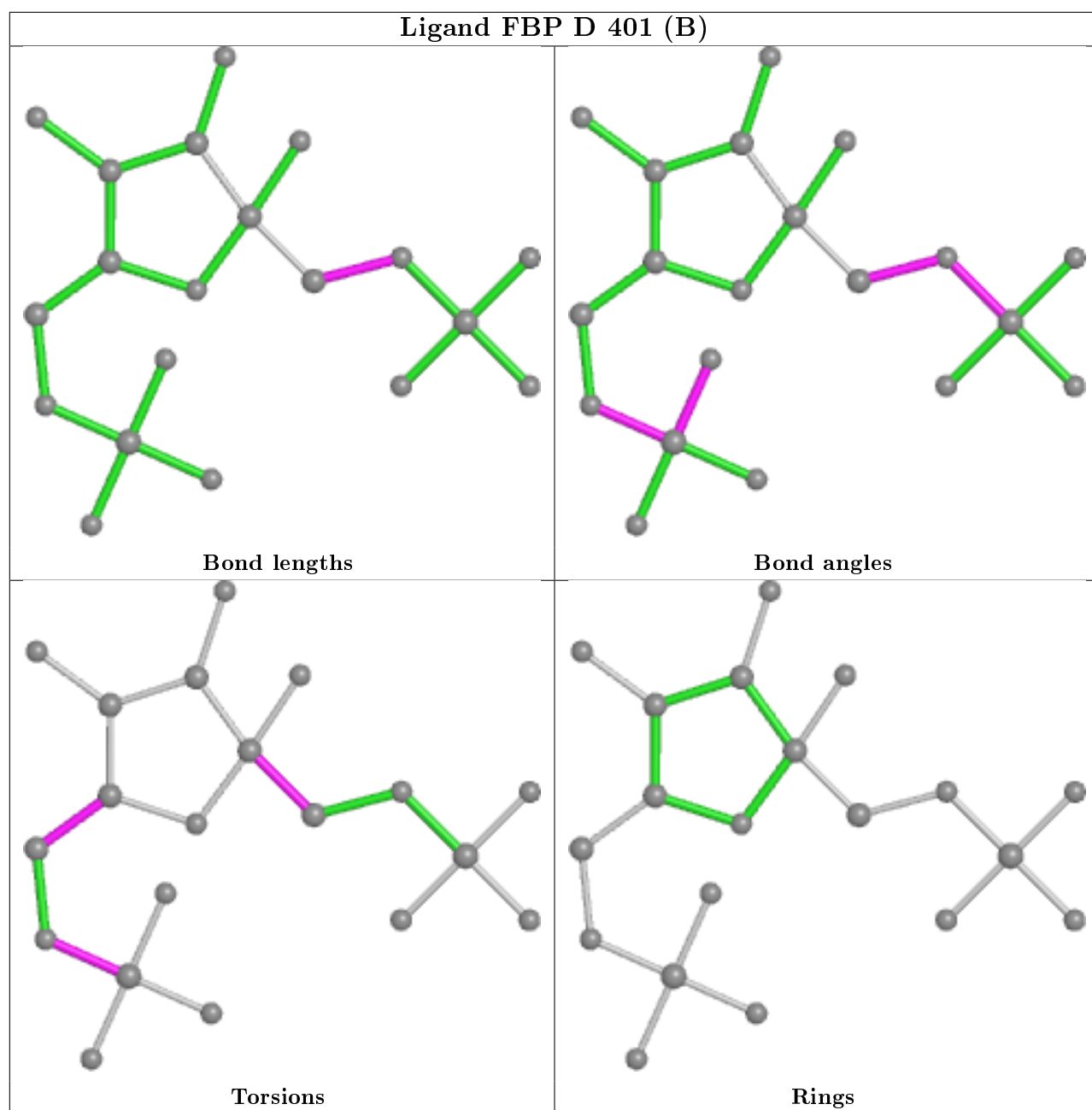












## 5.7 Other polymers [i](#)

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	316/322 (98%)	0.03	9 (2%) 53 60	17, 29, 49, 69	0
1	B	313/322 (97%)	0.29	17 (5%) 25 32	18, 31, 58, 87	0
1	C	314/322 (97%)	-0.06	7 (2%) 62 69	14, 30, 48, 82	0
1	D	314/322 (97%)	0.12	14 (4%) 33 40	15, 28, 56, 84	0
All	All	1257/1288 (97%)	0.09	47 (3%) 41 48	14, 29, 55, 87	0

All (47) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	315	LEU	7.0
1	B	315	LEU	6.9
1	B	88	GLY	6.2
1	B	313	VAL	6.1
1	D	316	GLU	5.7
1	D	87	PRO	5.6
1	B	87	PRO	4.6
1	B	86	LYS	4.3
1	C	316	GLU	4.1
1	B	314	ARG	3.7
1	D	313	VAL	3.6
1	B	227	ASP	3.6
1	D	204	HIS	3.4
1	D	88	GLY	3.3
1	D	89	GLN	3.3
1	D	85	GLN	3.2
1	D	91	ARG	3.2
1	A	85	GLN	3.2
1	B	89	GLN	3.0
1	D	223	TYR	2.9
1	B	85	GLN	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	38	VAL	2.9
1	B	311	ASP	2.8
1	A	40	GLN	2.8
1	A	87	PRO	2.7
1	B	92	LEU	2.7
1	A	86	LYS	2.7
1	B	204	HIS	2.7
1	C	315	LEU	2.6
1	D	92	LEU	2.6
1	B	223	TYR	2.5
1	D	90	THR	2.5
1	B	181	SER	2.5
1	A	84	ASN	2.5
1	B	208	GLU	2.4
1	C	313	VAL	2.4
1	C	86	LYS	2.2
1	C	111	ALA	2.2
1	A	269	ASP	2.2
1	B	91	ARG	2.2
1	C	38	VAL	2.1
1	C	67	THR	2.1
1	A	72	GLN	2.1
1	D	311	ASP	2.1
1	A	88	GLY	2.0
1	D	181	SER	2.0
1	B	90	THR	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 [\(i\)](#)

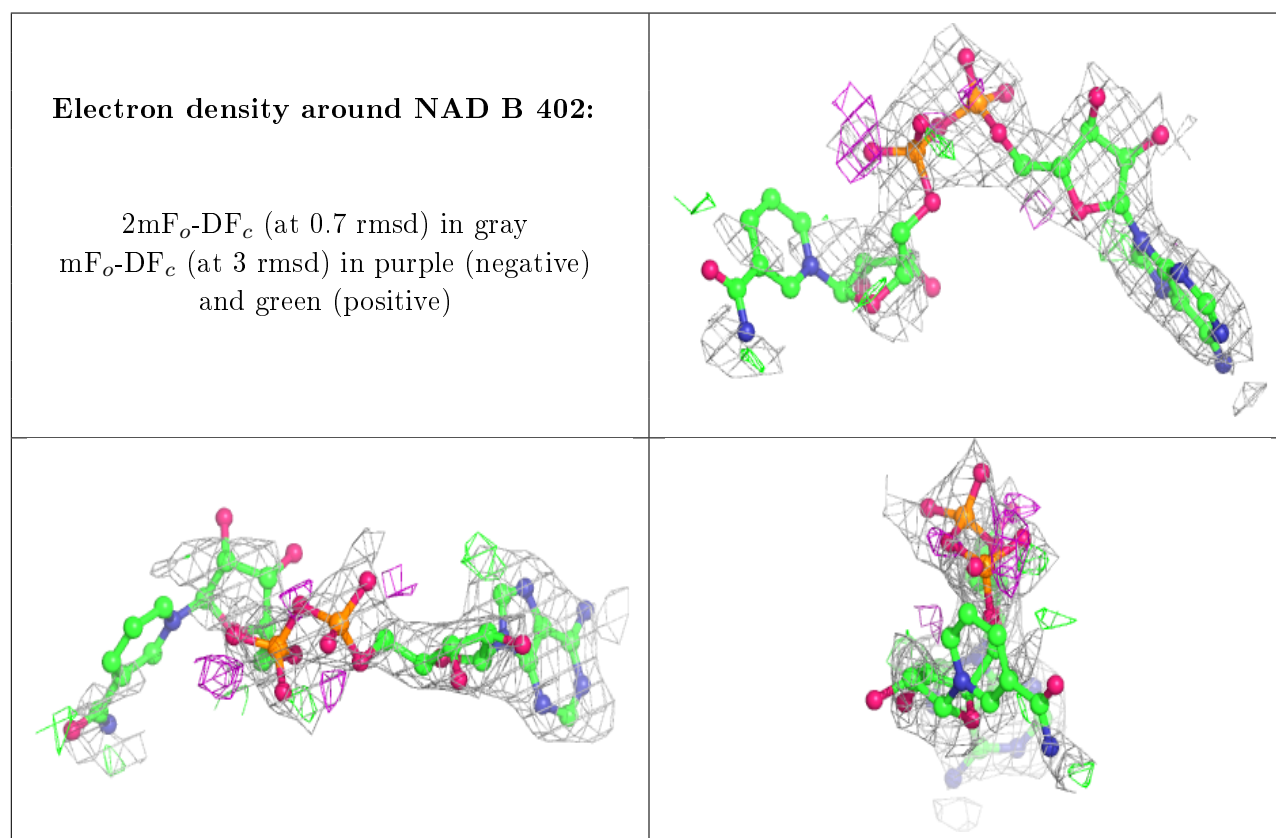
There are no monosaccharides in this entry.

## 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, 95<sup>th</sup> 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.

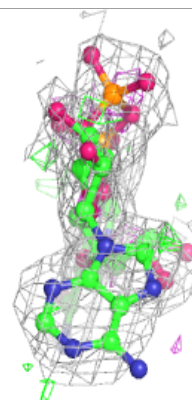
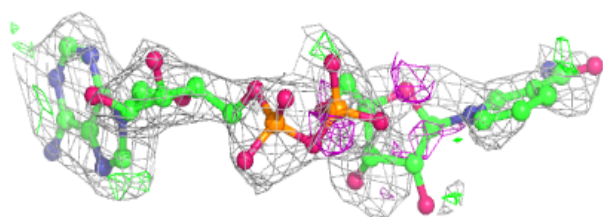
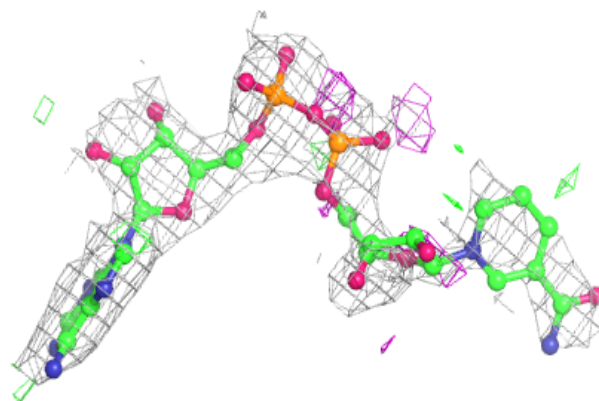
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	NAD	B	402	44/44	0.79	0.27	68,96,111,113	0
2	NAD	D	402	44/44	0.80	0.24	62,83,100,101	0
3	GOL	B	403	6/6	0.84	0.18	48,50,50,51	0
3	GOL	A	402	6/6	0.85	0.18	42,48,49,49	0
2	NAD	C	401	44/44	0.85	0.21	52,65,84,84	0
2	NAD	A	401	44/44	0.86	0.20	52,63,83,85	0
3	GOL	B	404	6/6	0.93	0.15	41,42,43,45	0
4	FBP	B	401[A]	20/20	0.97	0.17	11,17,21,21	20
4	FBP	B	401[B]	20/20	0.97	0.17	26,31,35,36	20
4	FBP	D	401[A]	20/20	0.98	0.14	12,17,19,22	20
4	FBP	D	401[B]	20/20	0.98	0.14	14,20,23,23	20

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

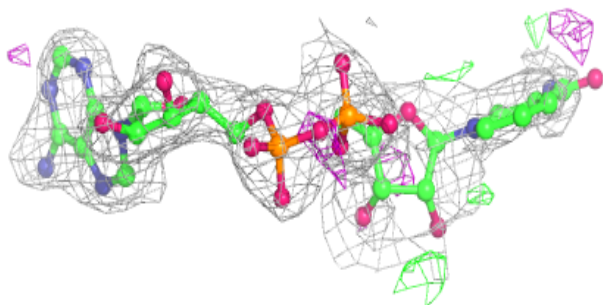
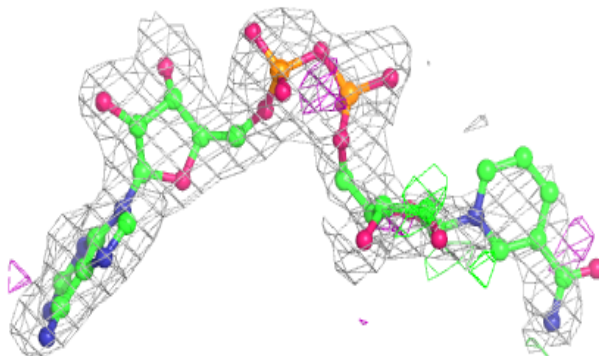


**Electron density around NAD D 402:**

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

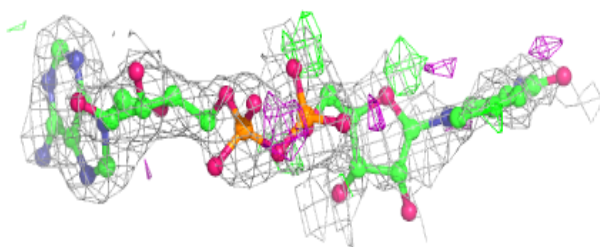
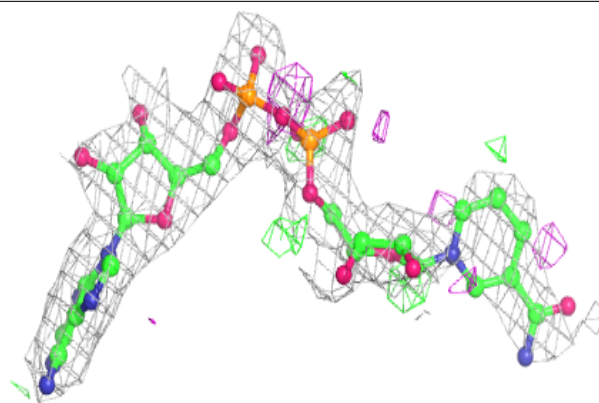
**Electron density around NAD C 401:**

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

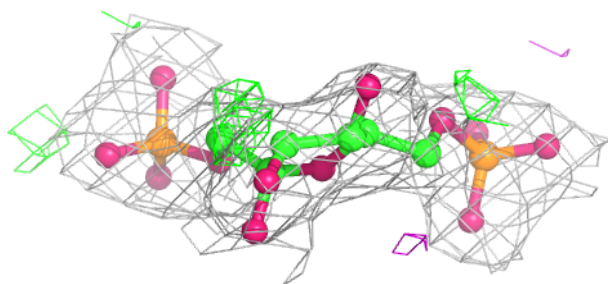
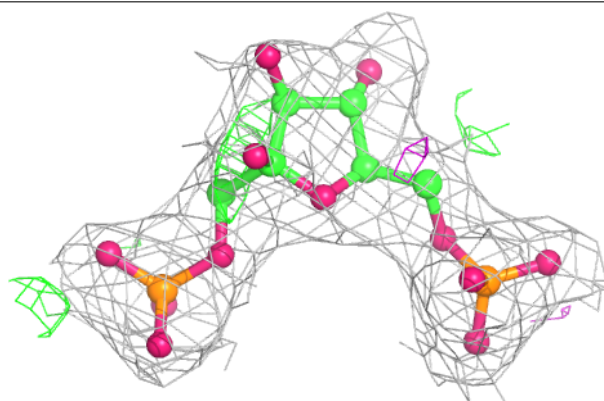


**Electron density around NAD A 401:**

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

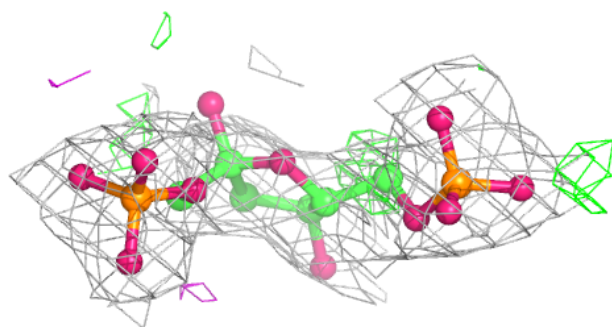
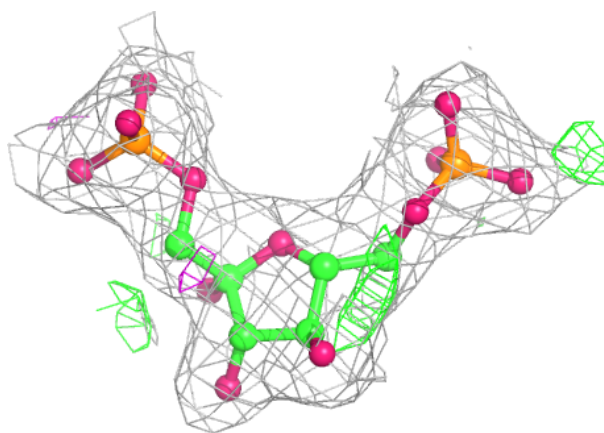
**Electron density around FBP B 401 (A):**

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

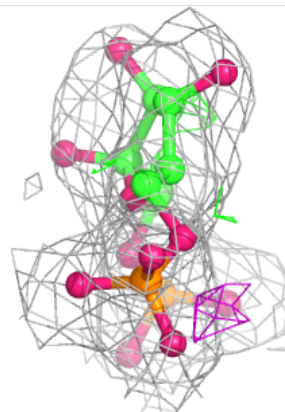
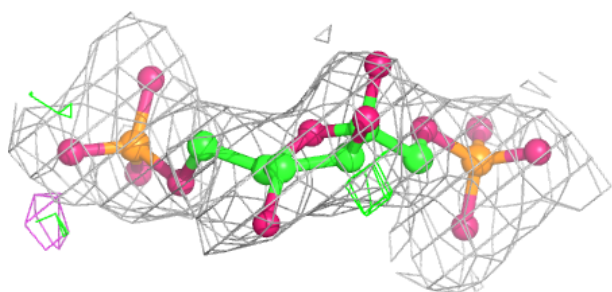
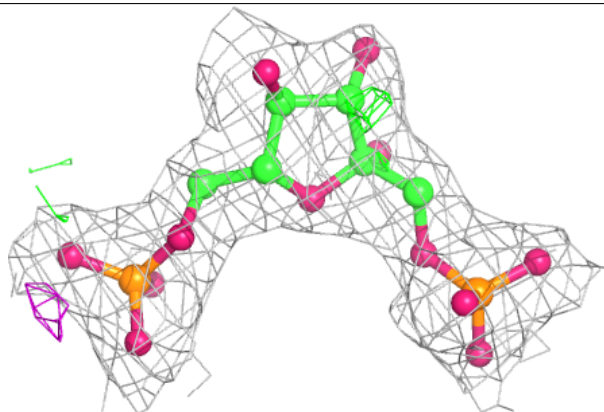


**Electron density around FBP B 401 (B):**

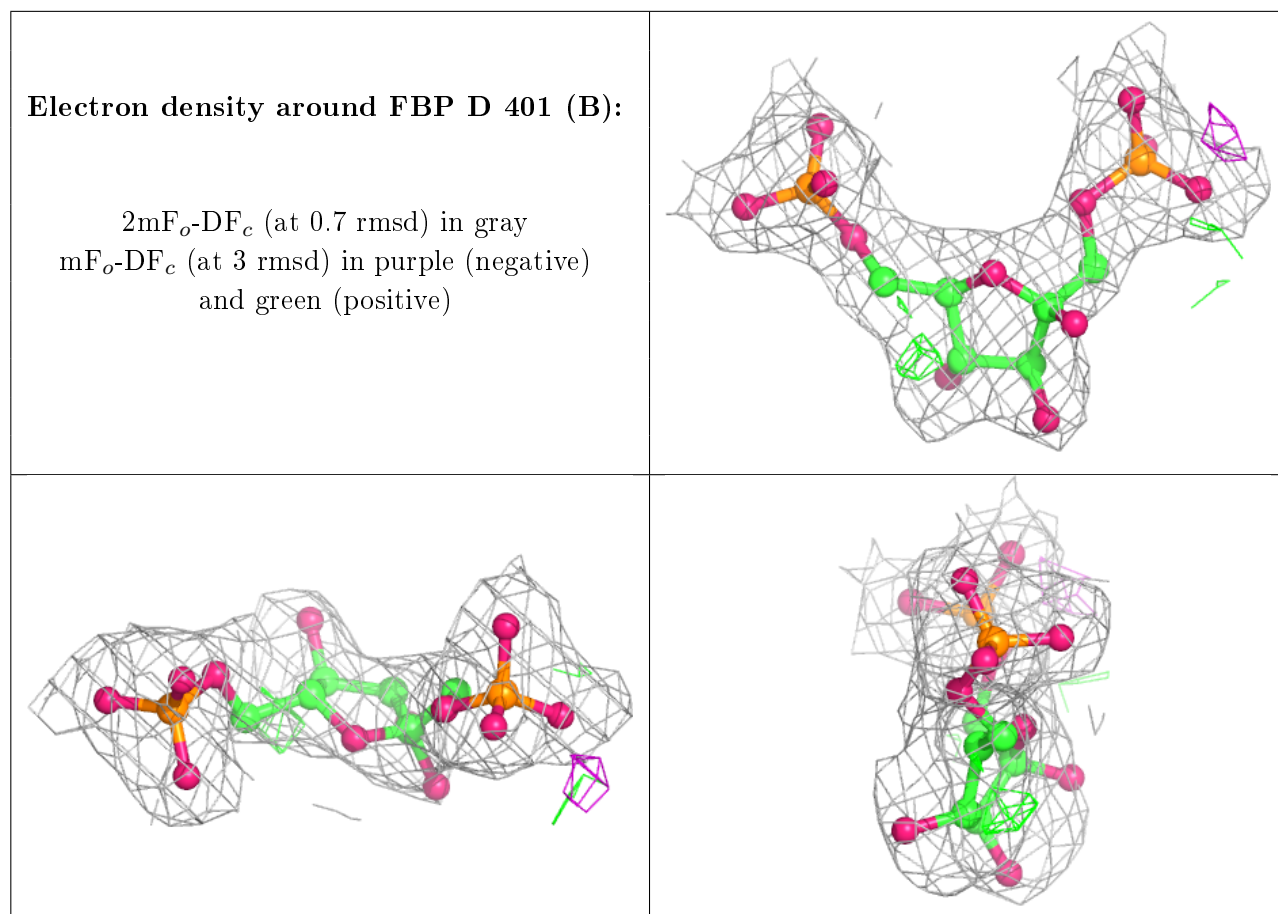
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around FBP D 401 (A):**

$2mF_o-DF_c$  (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.