



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 03:41 AM UTC

PDB ID : 3W83 / pdb_00003w83
Title : Structure of Trypanosoma cruzi dihydroorotate dehydrogenase in complex with MII-6-097
Authors : Inaoka, D.K.; Iida, M.; Tabuchi, T.; Lee, N.; Hashimoto, S.; Matsuoka, S.; Kuranaga, T.; Shiba, T.; Sakamoto, K.; Suzuki, S.; Balogun, E.O.; Nara, T.; Aoki, T.; Inoue, M.; Honma, T.; Tanaka, A.; Harada, S.; Kita, K.
Deposited on : 2013-03-12
Resolution : 2.80 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 types of validation reports are described at

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

The following versions of software and data (see [references](#) ①) 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

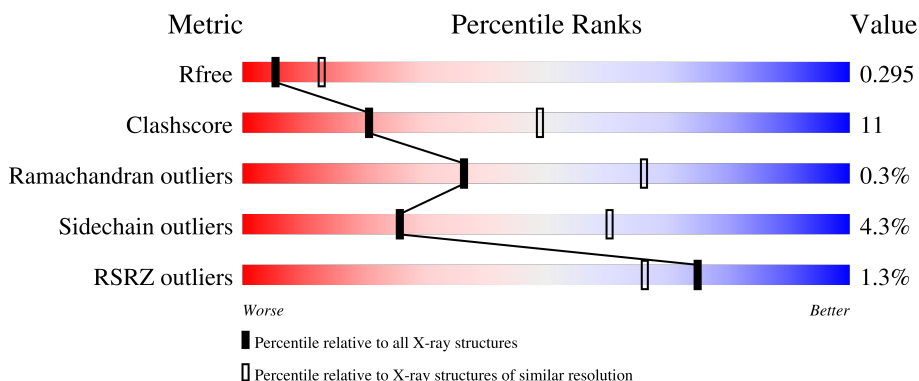
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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}	180053	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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.

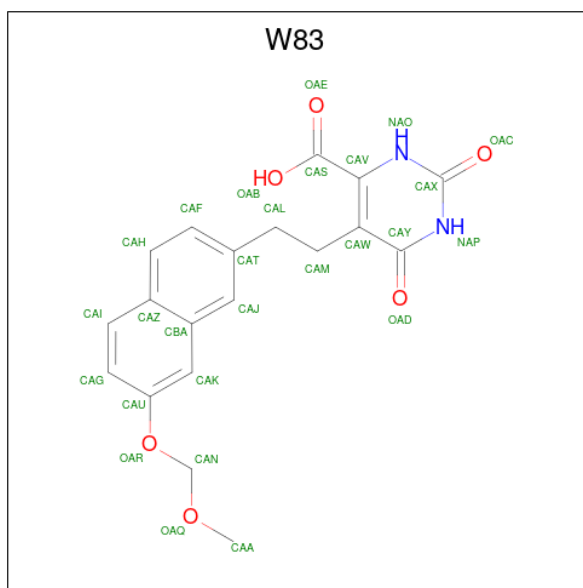
Mol	Chain	Length	Quality of chain
1	A	313	 76% 22% •
1	B	313	 3% 76% 23% •

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 Dihydroorotate dehydrogenase (fumarate).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	313	Total 2388	C 1530	N 396	O 444	S 18	0	0	0
1	B	313	Total 2388	C 1530	N 396	O 444	S 18	0	0	0

- Molecule 2 is 5-{2-[7-(methoxymethoxy)naphthalen-2-yl]ethyl}-2,6-dioxo-1,2,3,6-tetrahydropyrimidine-4-carboxylic acid (CCD ID: W83) (formula: $C_{19}H_{18}N_2O_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total 27	C 19	N 2	O 6	0	0
2	B	1	Total 27	C 19	N 2	O 6	0	0

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



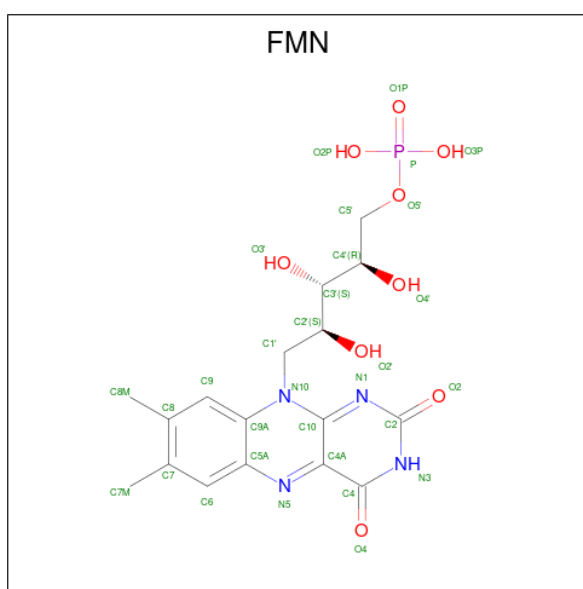
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total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

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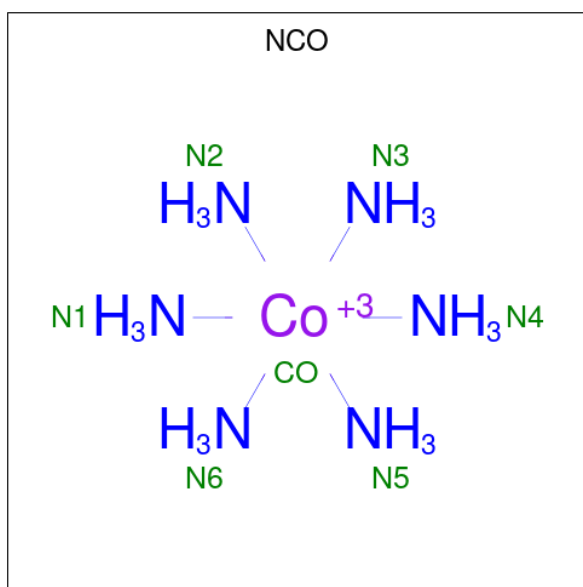
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is FLAVIN MONONUCLEOTIDE (CCD ID: FMN) (formula: $C_{17}H_{21}N_4O_9P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			31	17	4	9	1		
4	B	1	Total	C	N	O	P	0	0
			31	17	4	9	1		

- Molecule 5 is COBALT HEXAMMINE(III) (CCD ID: NCO) (formula: $CoH_{18}N_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	Co	N	0	0
			7	1	6		

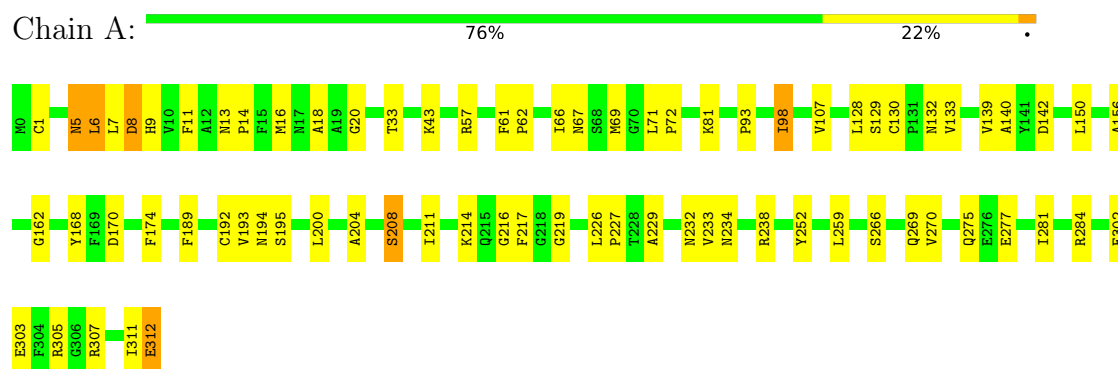
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	170	Total	O	0	0
			170	170		
6	B	152	Total	O	0	0
			152	152		

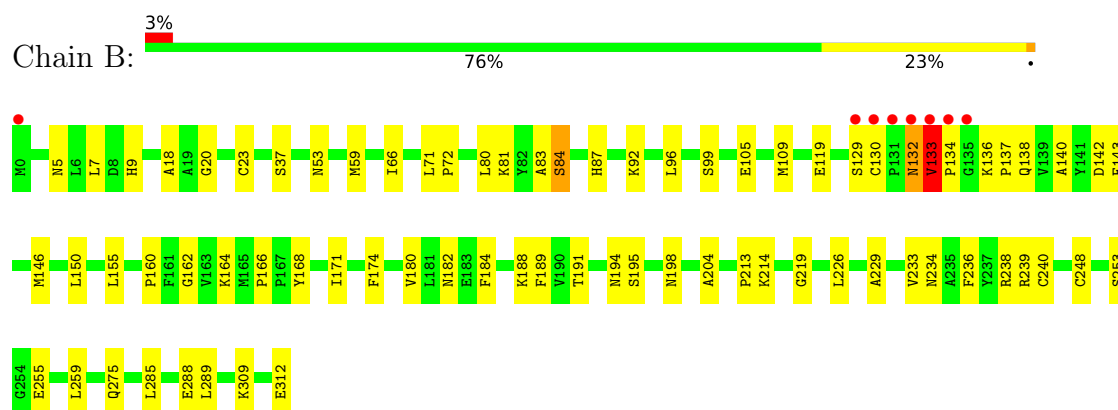
3 Residue-property plots

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

- Molecule 1: Dihydroorotate dehydrogenase (fumarate)



- Molecule 1: Dihydroorotate dehydrogenase (fumarate)



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	68.16Å 71.72Å 129.92Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	34.08 – 2.80 34.08 – 2.80	Depositor EDS
% Data completeness (in resolution range)	97.2 (34.08-2.80) 97.3 (34.08-2.80)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.64 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.194 , 0.297 0.191 , 0.295	Depositor DCC
R_{free} test set	791 reflections (4.83%)	wwPDB-VP
Wilson B-factor (Å ²)	25.8	Xtriage
Anisotropy	0.228	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 30.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.000 for k,h,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5329	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 38.75 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 3.5161e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: W83, GOL, FMN, NCO

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	0.78	0/2440	1.02	3/3305 (0.1%)
1	B	0.79	0/2440	1.00	2/3305 (0.1%)
All	All	0.78	0/4880	1.01	5/6610 (0.1%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	129	SER	N-CA-C	-5.58	105.12	112.72
1	A	1	CYS	N-CA-C	5.45	117.08	108.96
1	B	226	LEU	CA-C-N	-5.28	113.10	118.85
1	B	226	LEU	C-N-CA	-5.28	113.10	118.85
1	A	150	LEU	N-CA-C	5.27	117.11	111.36

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	2388	0	2391	53	0
1	B	2388	0	2391	59	1
2	A	27	0	17	2	0
2	B	27	0	17	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	54	0	72	0	0
3	B	54	0	72	2	0
4	A	31	0	19	4	0
4	B	31	0	19	5	0
5	A	7	0	0	2	1
6	A	170	0	0	5	0
6	B	152	0	0	3	0
All	All	5329	0	4998	111	1

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 11.

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:A:7:LEU:O	1:A:9:HIS:HD2	1.52	0.93
1:B:132:ASN:HB2	1:B:133:VAL:HG13	1.55	0.87
1:A:11:PHE:CE2	1:A:93:PRO:HG2	2.18	0.78
1:B:133:VAL:HG21	1:B:136:LYS:HB2	1.65	0.77
1:B:162:GLY:HA3	1:B:189:PHE:CZ	2.21	0.75
1:B:162:GLY:HA3	1:B:189:PHE:CE1	2.25	0.72
1:B:132:ASN:CB	1:B:133:VAL:HG13	2.20	0.71
1:B:37:SER:OG	1:B:92:LYS:NZ	2.22	0.70
1:B:133:VAL:CG2	1:B:136:LYS:HB2	2.23	0.69
1:A:5:ASN:HD22	1:A:5:ASN:N	1.92	0.67
1:B:160:PRO:HB3	1:B:188:LYS:HG3	1.77	0.67
1:B:132:ASN:HB2	1:B:133:VAL:CG1	2.27	0.64
1:A:7:LEU:O	1:A:9:HIS:CD2	2.44	0.61
1:B:132:ASN:HB2	1:B:133:VAL:HG22	1.82	0.60
1:B:133:VAL:HB	1:B:134:PRO:O	2.01	0.60
1:B:236:PHE:O	1:B:240:CYS:HB2	2.02	0.59
1:B:20:GLY:HA3	4:B:411:FMN:N5	2.17	0.59
1:A:277:GLU:OE1	1:A:281:ILE:HD12	2.03	0.58
1:B:132:ASN:ND2	1:B:137:PRO:O	2.36	0.58
1:B:234:ASN:HD21	1:B:238:ARG:HE	1.52	0.57
1:A:229:ALA:O	1:A:233:VAL:HG23	2.05	0.56
1:A:275:GLN:O	5:A:412:NCO:N3	2.38	0.56
1:B:20:GLY:HA3	4:B:411:FMN:C5A	2.35	0.56
1:A:259:LEU:HD11	6:A:587:HOH:O	2.05	0.56
1:B:18:ALA:HB1	4:B:411:FMN:O2'	2.06	0.54
1:A:234:ASN:HD21	1:A:238:ARG:HE	1.55	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:140:ALA:HB3	1:B:166:PRO:HG3	1.90	0.54
1:B:168:TYR:HB2	1:B:174:PHE:CE1	2.43	0.54
1:A:18:ALA:HB1	4:A:411:FMN:O2'	2.08	0.54
1:A:71:LEU:N	1:A:72:PRO:CD	2.70	0.54
1:A:130:CYS:O	1:A:139:VAL:HG23	2.08	0.54
1:A:226:LEU:HB3	1:A:227:PRO:HD3	1.89	0.53
1:B:80:LEU:O	1:B:84:SER:HB2	2.09	0.53
1:A:132:ASN:HB2	1:A:139:VAL:HG22	1.90	0.53
1:A:259:LEU:HD21	1:B:204:ALA:HB2	1.91	0.52
1:B:87:HIS:HE1	1:B:92:LYS:O	1.93	0.52
1:B:133:VAL:HG23	1:B:136:LYS:H	1.75	0.52
1:B:182:ASN:OD1	6:B:522:HOH:O	2.19	0.52
1:A:11:PHE:CD2	1:A:93:PRO:HG2	2.45	0.51
1:B:184:PHE:HB3	6:B:502:HOH:O	2.10	0.51
1:A:69:MET:O	1:A:72:PRO:HD3	2.11	0.50
1:B:146:MET:HE3	1:B:180:VAL:HB	1.94	0.50
1:B:198:ASN:OD1	1:B:214:LYS:NZ	2.44	0.49
1:A:168:TYR:HB2	1:A:174:PHE:CE1	2.47	0.49
5:A:412:NCO:N4	6:A:636:HOH:O	2.35	0.49
1:A:71:LEU:HB2	2:A:401:W83:H8	1.95	0.49
1:B:143:PHE:HD1	1:B:146:MET:HE2	1.76	0.49
1:B:194:ASN:HA	4:B:411:FMN:O3'	2.13	0.49
1:A:20:GLY:HA3	4:A:411:FMN:N5	2.27	0.48
1:B:146:MET:HG2	1:B:180:VAL:HG11	1.95	0.48
1:A:20:GLY:HA2	1:A:43:LYS:HD2	1.95	0.48
1:A:13:ASN:HB2	1:A:14:PRO:HD2	1.95	0.48
1:B:71:LEU:N	1:B:72:PRO:CD	2.77	0.48
1:B:7:LEU:O	1:B:9:HIS:HD2	1.97	0.48
1:B:59:MET:HE3	6:B:576:HOH:O	2.14	0.47
1:B:162:GLY:HA3	1:B:189:PHE:CE2	2.49	0.47
1:A:204:ALA:HB2	1:B:259:LEU:HD21	1.95	0.47
1:B:53:ASN:HB3	1:B:213:PRO:HG3	1.97	0.47
1:A:232:ASN:ND2	6:A:523:HOH:O	2.47	0.47
1:B:234:ASN:ND2	1:B:238:ARG:HE	2.13	0.46
1:A:192:CYS:HB3	1:A:193:VAL:HG13	1.96	0.46
1:B:99:SER:OG	1:B:130:CYS:HB3	2.16	0.46
1:B:171:ILE:HD12	3:B:403:GOL:H31	1.98	0.46
1:B:229:ALA:O	1:B:233:VAL:HG23	2.16	0.46
1:A:234:ASN:HD21	1:A:238:ARG:NE	2.15	0.45
1:A:252:TYR:O	1:A:284:ARG:NH2	2.49	0.45
1:B:132:ASN:HB2	1:B:133:VAL:CG2	2.45	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:128:LEU:C	1:A:130:CYS:H	2.22	0.45
1:B:132:ASN:HB2	1:B:133:VAL:CB	2.46	0.45
1:A:6:LEU:C	1:A:8:ASP:N	2.72	0.45
1:A:71:LEU:HD13	2:A:401:W83:H7	1.99	0.45
1:B:162:GLY:CA	1:B:189:PHE:CE1	2.99	0.45
1:B:248:CYS:SG	4:B:411:FMN:H3'	2.57	0.45
1:A:302:GLU:HG3	1:A:305:ARG:HH12	1.82	0.45
1:B:239:ARG:HH22	3:B:403:GOL:H11	1.82	0.45
1:B:66:ILE:HA	1:B:219:GLY:O	2.17	0.44
1:B:255:GLU:O	1:B:259:LEU:HG	2.17	0.44
1:A:208:SER:HB2	1:B:309:LYS:HB2	1.99	0.44
1:A:16:MET:O	1:A:269:GLN:HA	2.17	0.44
1:A:162:GLY:HA3	1:A:189:PHE:CZ	2.51	0.44
1:A:6:LEU:O	1:A:7:LEU:C	2.60	0.44
1:B:105:GLU:HG2	1:B:109:MET:HE2	2.00	0.44
1:B:164:LYS:HD2	1:B:191:THR:HB	2.00	0.44
1:B:83:ALA:HB2	1:B:96:LEU:HD12	2.00	0.44
1:A:132:ASN:O	1:A:133:VAL:C	2.62	0.43
1:A:71:LEU:N	1:A:72:PRO:HD2	2.33	0.43
1:A:57:ARG:O	1:A:67:ASN:HA	2.17	0.43
1:A:162:GLY:HA3	1:A:189:PHE:CE2	2.54	0.43
1:B:7:LEU:O	1:B:9:HIS:CD2	2.72	0.42
1:A:140:ALA:HB1	1:A:168:TYR:OH	2.19	0.42
1:A:170:ASP:OD1	1:B:137:PRO:HA	2.19	0.42
1:A:200:LEU:HD13	1:A:217:PHE:CE1	2.55	0.42
4:A:411:FMN:H9	4:A:411:FMN:O4'	2.20	0.42
4:A:411:FMN:HM73	4:A:411:FMN:HM81	1.83	0.42
1:A:61:PHE:O	1:A:62:PRO:C	2.60	0.42
1:B:288:GLU:O	1:B:289:LEU:C	2.62	0.42
1:A:66:ILE:HA	1:A:219:GLY:O	2.20	0.42
1:A:98:ILE:O	1:A:98:ILE:HG13	2.20	0.42
1:A:311:ILE:O	1:A:312:GLU:C	2.62	0.42
1:A:211:ILE:HG13	1:A:216:GLY:HA2	2.01	0.41
1:A:142:ASP:HA	6:A:563:HOH:O	2.19	0.41
1:A:194:ASN:OD1	1:A:195:SER:N	2.48	0.41
1:A:211:ILE:O	1:A:216:GLY:HA2	2.21	0.41
1:B:194:ASN:OD1	1:B:195:SER:N	2.51	0.41
1:A:107:VAL:HG13	1:A:156:ALA:HB2	2.03	0.41
1:A:132:ASN:HB3	6:A:599:HOH:O	2.20	0.41
1:B:146:MET:O	1:B:150:LEU:HG	2.21	0.40
1:B:253:SER:C	1:B:285:LEU:HD23	2.46	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:71:LEU:HB2	2:B:401:W83:H8	2.03	0.40
1:B:140:ALA:C	1:B:142:ASP:H	2.30	0.40
1:A:214:LYS:HE2	1:A:214:LYS:HB3	1.88	0.40

All (1) 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)
1:B:275:GLN:O	5:A:412:NCO:N1[3_545]	2.16	0.04

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	311/313 (99%)	297 (96%)	14 (4%)	0	100	100
1	B	311/313 (99%)	295 (95%)	14 (4%)	2 (1%)	21	51
All	All	622/626 (99%)	592 (95%)	28 (4%)	2 (0%)	36	66

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	23	CYS
1	B	133	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	258/258 (100%)	246 (95%)	12 (5%)	23	57
1	B	258/258 (100%)	248 (96%)	10 (4%)	28	64
All	All	516/516 (100%)	494 (96%)	22 (4%)	26	60

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	ASN
1	A	6	LEU
1	A	8	ASP
1	A	33	THR
1	A	81	LYS
1	A	98	ILE
1	A	208	SER
1	A	266	SER
1	A	270	VAL
1	A	303	GLU
1	A	307	ARG
1	A	312	GLU
1	B	5	ASN
1	B	81	LYS
1	B	84	SER
1	B	119	GLU
1	B	129	SER
1	B	132	ASN
1	B	133	VAL
1	B	138	GLN
1	B	155	LEU
1	B	312	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	5	ASN
1	A	9	HIS
1	A	87	HIS
1	A	151	GLN
1	A	234	ASN
1	A	275	GLN

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Mol	Chain	Res	Type
1	B	9	HIS
1	B	87	HIS
1	B	138	GLN
1	B	234	ASN
1	B	275	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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

23 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	406	-	5,5,5	0.39	0	5,5,5	0.16	0
3	GOL	B	405	-	5,5,5	0.32	0	5,5,5	0.28	0
3	GOL	B	410	-	5,5,5	0.30	0	5,5,5	0.45	0
3	GOL	A	410	-	5,5,5	0.41	0	5,5,5	0.20	0
2	W83	B	401	-	29,29,29	2.04	9 (31%)	38,40,40	2.58	14 (36%)
3	GOL	B	407	-	5,5,5	0.31	0	5,5,5	0.25	0
2	W83	A	401	-	29,29,29	2.16	4 (13%)	38,40,40	2.59	11 (28%)
3	GOL	B	408	-	5,5,5	0.43	0	5,5,5	0.61	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	GOL	B	404	-	5,5,5	0.31	0	5,5,5	0.67	0
3	GOL	A	404	-	5,5,5	0.36	0	5,5,5	0.49	0
3	GOL	A	408	-	5,5,5	0.47	0	5,5,5	0.47	0
3	GOL	A	409	-	5,5,5	0.42	0	5,5,5	0.24	0
3	GOL	A	403	-	5,5,5	0.35	0	5,5,5	0.12	0
3	GOL	A	407	-	5,5,5	0.45	0	5,5,5	0.18	0
3	GOL	B	403	-	5,5,5	0.22	0	5,5,5	0.40	0
3	GOL	A	405	-	5,5,5	0.39	0	5,5,5	0.72	0
3	GOL	A	406	-	5,5,5	0.28	0	5,5,5	0.44	0
4	FMN	A	411	-	33,33,33	0.88	1 (3%)	48,50,50	1.44	11 (22%)
5	NCO	A	412	-	6,6,6	3.38	6 (100%)	-		
4	FMN	B	411	-	33,33,33	0.88	1 (3%)	48,50,50	1.17	6 (12%)
3	GOL	A	402	-	5,5,5	0.39	0	5,5,5	0.43	0
3	GOL	B	402	-	5,5,5	0.39	0	5,5,5	0.29	0
3	GOL	B	409	-	5,5,5	0.39	0	5,5,5	0.82	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. '2' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GOL	B	406	-	-	2/4/4/4	-
3	GOL	B	405	-	-	0/4/4/4	-
3	GOL	B	410	-	-	0/4/4/4	-
3	GOL	A	410	-	-	0/4/4/4	-
2	W83	B	401	-	-	7/13/13/13	0/3/3/3
3	GOL	B	407	-	-	1/4/4/4	-
2	W83	A	401	-	-	8/13/13/13	0/3/3/3
3	GOL	B	408	-	-	1/4/4/4	-
3	GOL	B	404	-	-	1/4/4/4	-
3	GOL	A	404	-	-	2/4/4/4	-
3	GOL	A	408	-	-	2/4/4/4	-
3	GOL	A	409	-	-	0/4/4/4	-
3	GOL	A	403	-	-	2/4/4/4	-
3	GOL	A	407	-	-	4/4/4/4	-
3	GOL	B	403	-	-	2/4/4/4	-
3	GOL	A	405	-	-	2/4/4/4	-
3	GOL	A	406	-	-	0/4/4/4	-
4	FMN	A	411	-	-	7/18/18/18	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	FMN	B	411	-	-	5/18/18/18	0/3/3/3
3	GOL	A	402	-	-	3/4/4/4	-
3	GOL	B	402	-	-	2/4/4/4	-
3	GOL	B	409	-	-	3/4/4/4	-

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	W83	OAD-CAY	6.47	1.35	1.23
2	A	401	W83	OAC-CAX	5.64	1.35	1.23
2	B	401	W83	OAD-CAY	5.46	1.33	1.23
2	A	401	W83	CAY-CAW	-4.32	1.37	1.44
2	B	401	W83	OAC-CAX	4.31	1.32	1.23
5	A	412	NCO	CO-N2	3.92	2.09	1.96
2	B	401	W83	CAY-CAW	-3.68	1.38	1.44
5	A	412	NCO	CO-N6	3.60	2.08	1.96
5	A	412	NCO	CO-N3	3.57	2.08	1.96
5	A	412	NCO	CO-N5	3.50	2.07	1.96
2	B	401	W83	CAV-CAW	3.08	1.43	1.37
5	A	412	NCO	CO-N4	3.01	2.06	1.96
4	B	411	FMN	C4A-N5	2.97	1.37	1.30
4	A	411	FMN	C4A-N5	2.80	1.36	1.30
2	B	401	W83	CAV-CAS	2.51	1.52	1.48
5	A	412	NCO	CO-N1	2.48	2.04	1.96
2	B	401	W83	CAL-CAT	-2.38	1.45	1.51
2	A	401	W83	CAL-CAT	-2.25	1.45	1.51
2	B	401	W83	CAY-NAP	-2.25	1.34	1.38
2	B	401	W83	CAI-CAG	2.07	1.41	1.36
2	B	401	W83	CAH-CAF	2.06	1.41	1.36

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	W83	CAM-CAL-CAT	7.91	140.51	112.70
2	A	401	W83	CAW-CAY-NAP	7.47	121.37	115.70
2	B	401	W83	CAW-CAY-NAP	6.92	120.96	115.70
2	A	401	W83	CAM-CAL-CAT	6.41	135.22	112.70
2	A	401	W83	OAD-CAY-CAW	-5.74	117.48	124.56
2	B	401	W83	CAY-NAP-CAX	-5.00	119.48	126.37
2	A	401	W83	CAM-CAW-CAV	4.69	128.90	119.56
2	A	401	W83	CAY-NAP-CAX	-4.64	119.98	126.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	W83	CAL-CAT-CAJ	-4.37	112.01	121.04
2	B	401	W83	NAP-CAX-NAO	4.14	122.25	115.74
2	B	401	W83	CAM-CAW-CAV	3.89	127.31	119.56
2	B	401	W83	CAL-CAT-CAJ	-3.85	113.09	121.04
2	B	401	W83	OAD-CAY-CAW	-3.73	119.97	124.56
4	A	411	FMN	C4-N3-C2	-3.61	119.24	125.64
4	B	411	FMN	C4A-C10-N10	3.28	121.18	116.48
2	A	401	W83	OAB-CAS-CAV	3.19	121.79	116.73
2	A	401	W83	CAM-CAW-CAY	-3.15	112.42	116.64
4	A	411	FMN	C4A-C10-N10	3.00	120.78	116.48
4	A	411	FMN	O4'-C4'-C3'	2.99	116.25	109.25
4	B	411	FMN	O2'-C2'-C3'	2.82	115.84	109.25
4	A	411	FMN	O4-C4-C4A	-2.75	119.27	126.53
2	B	401	W83	CAS-CAV-CAW	2.69	127.95	123.88
4	A	411	FMN	O2-C2-N1	-2.62	117.45	121.80
4	A	411	FMN	C5'-C4'-C3'	-2.61	107.30	112.22
2	B	401	W83	CAK-CBA-CAJ	-2.56	116.92	122.14
2	B	401	W83	CAM-CAW-CAY	-2.50	113.29	116.64
2	B	401	W83	OAE-CAS-CAV	2.45	124.46	120.20
4	B	411	FMN	C10-C4A-N5	-2.44	119.82	124.81
4	B	411	FMN	O4'-C4'-C3'	2.39	114.84	109.25
4	A	411	FMN	C4A-C4-N3	2.38	119.31	113.25
4	A	411	FMN	C10-C4A-N5	-2.33	120.05	124.81
2	A	401	W83	NAP-CAX-NAO	2.29	119.34	115.74
2	B	401	W83	OAR-CAU-CAK	-2.29	116.63	124.08
2	A	401	W83	OAE-CAS-CAV	-2.25	116.28	120.20
4	A	411	FMN	C4A-C10-N1	-2.16	119.30	124.59
4	B	411	FMN	C4-C4A-C10	2.15	120.63	116.93
2	A	401	W83	CAK-CBA-CAJ	-2.15	117.75	122.14
4	B	411	FMN	O3P-P-O5'	2.13	112.22	106.67
4	A	411	FMN	C9A-C5A-N5	-2.10	120.23	122.45
2	B	401	W83	CAJ-CBA-CAZ	2.08	121.84	118.96
2	B	401	W83	OAC-CAX-NAP	-2.01	118.28	121.86
4	A	411	FMN	C4-C4A-C10	2.01	120.37	116.93

There are no chirality outliers.

All (54) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401	W83	OAB-CAS-CAV-NAO
2	A	401	W83	CAL-CAM-CAW-CAY
2	B	401	W83	OAB-CAS-CAV-NAO

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Mol	Chain	Res	Type	Atoms
2	B	401	W83	CAL-CAM-CAW-CAY
3	A	402	GOL	O1-C1-C2-C3
3	A	403	GOL	O1-C1-C2-C3
3	A	405	GOL	C1-C2-C3-O3
3	A	405	GOL	O2-C2-C3-O3
3	A	407	GOL	O1-C1-C2-C3
3	A	407	GOL	C1-C2-C3-O3
3	A	408	GOL	O1-C1-C2-O2
3	A	408	GOL	O1-C1-C2-C3
3	B	403	GOL	O1-C1-C2-C3
3	B	409	GOL	C1-C2-C3-O3
4	A	411	FMN	C5'-O5'-P-O2P
4	A	411	FMN	C5'-O5'-P-O3P
2	B	401	W83	OAR-CAN-OAQ-CAA
4	B	411	FMN	C2'-C3'-C4'-O4'
4	A	411	FMN	O3'-C3'-C4'-C5'
2	A	401	W83	OAR-CAN-OAQ-CAA
4	A	411	FMN	O3'-C3'-C4'-O4'
4	B	411	FMN	O3'-C3'-C4'-O4'
3	A	403	GOL	O1-C1-C2-O2
4	A	411	FMN	C2'-C3'-C4'-C5'
4	A	411	FMN	C2'-C3'-C4'-O4'
3	A	404	GOL	O1-C1-C2-C3
3	B	402	GOL	C1-C2-C3-O3
3	B	406	GOL	C1-C2-C3-O3
3	A	404	GOL	O1-C1-C2-O2
3	A	407	GOL	O1-C1-C2-O2
3	A	407	GOL	O2-C2-C3-O3
3	B	402	GOL	O2-C2-C3-O3
3	B	403	GOL	O1-C1-C2-O2
4	B	411	FMN	O3'-C3'-C4'-C5'
3	B	409	GOL	O2-C2-C3-O3
4	B	411	FMN	C2'-C3'-C4'-C5'
3	A	402	GOL	O1-C1-C2-O2
2	B	401	W83	OAQ-CAN-OAR-CAU
3	B	409	GOL	O1-C1-C2-O2
2	A	401	W83	OAQ-CAN-OAR-CAU
3	A	402	GOL	O2-C2-C3-O3
3	B	408	GOL	O1-C1-C2-O2
2	A	401	W83	OAE-CAS-CAV-CAW
4	B	411	FMN	C4'-C5'-O5'-P
4	A	411	FMN	C5'-O5'-P-O1P

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Mol	Chain	Res	Type	Atoms
3	B	406	GOL	O2-C2-C3-O3
2	A	401	W83	CAM-CAL-CAT-CAJ
2	A	401	W83	CAM-CAL-CAT-CAF
2	B	401	W83	CAM-CAL-CAT-CAJ
3	B	407	GOL	O1-C1-C2-C3
3	B	404	GOL	C1-C2-C3-O3
2	B	401	W83	CAG-CAU-OAR-CAN
2	A	401	W83	OAE-CAS-CAV-NAO
2	B	401	W83	OAE-CAS-CAV-NAO

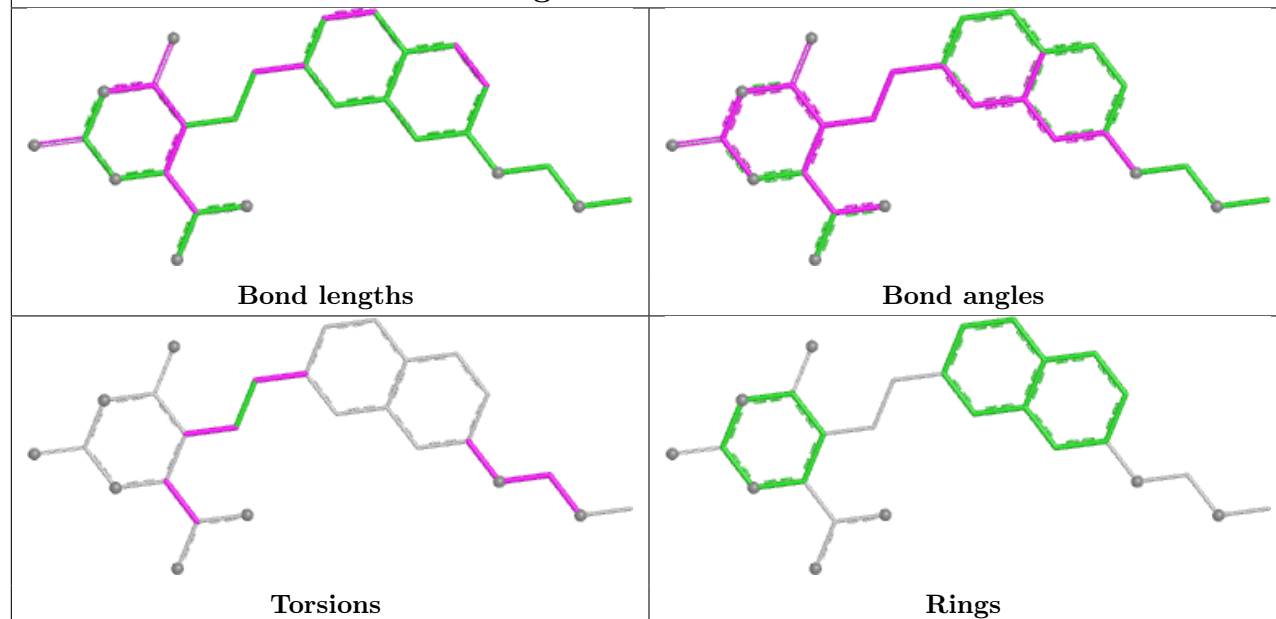
There are no ring outliers.

6 monomers are involved in 17 short contacts:

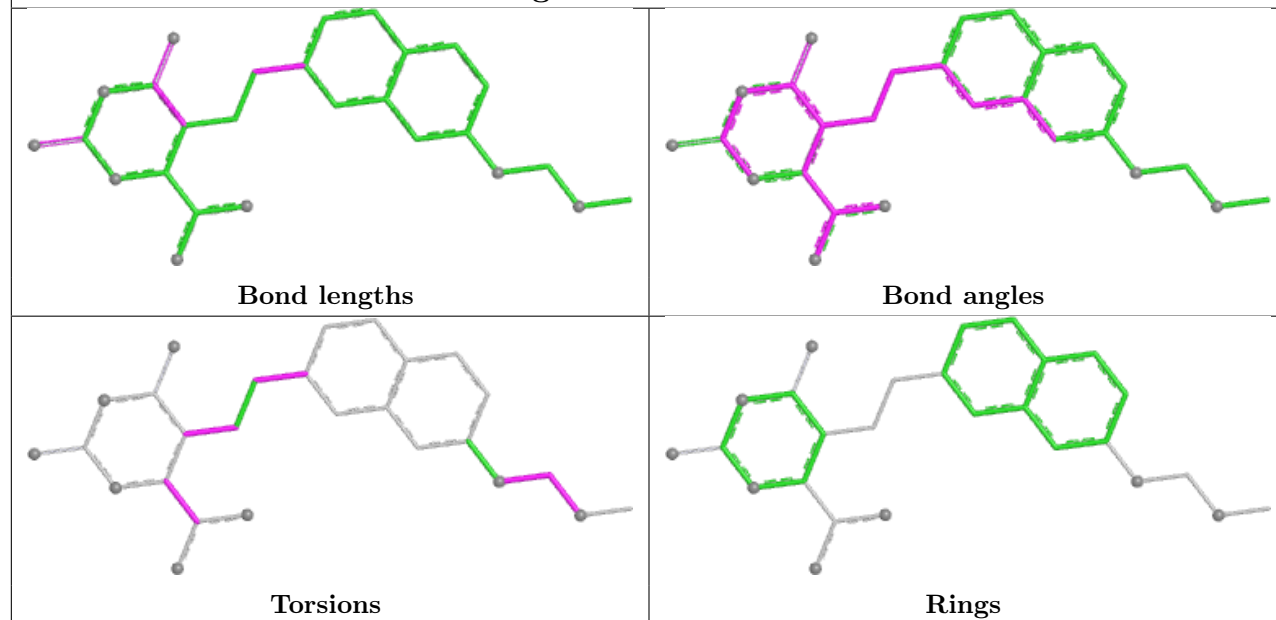
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	401	W83	1	0
2	A	401	W83	2	0
3	B	403	GOL	2	0
4	A	411	FMN	4	0
5	A	412	NCO	2	1
4	B	411	FMN	5	0

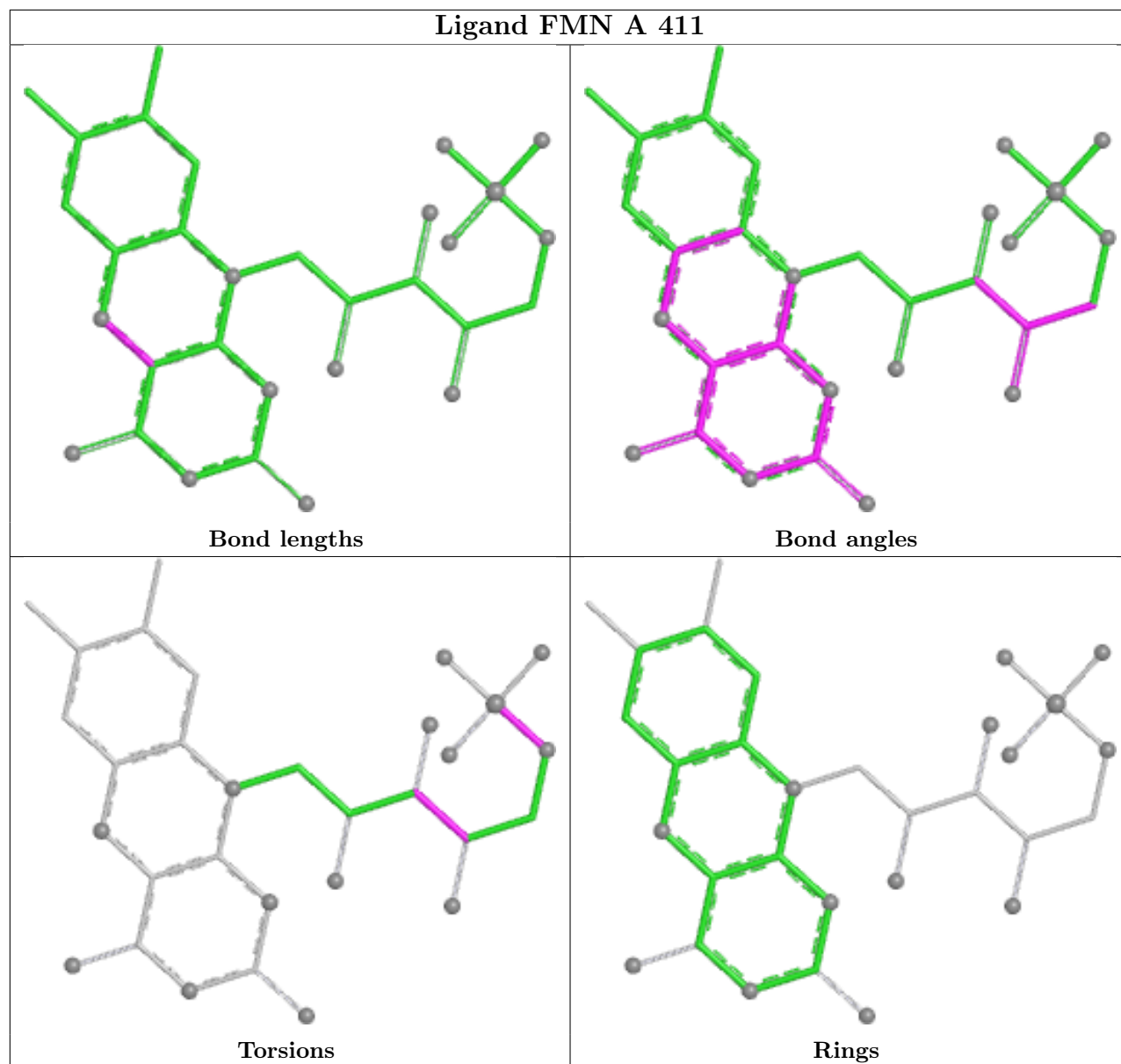
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

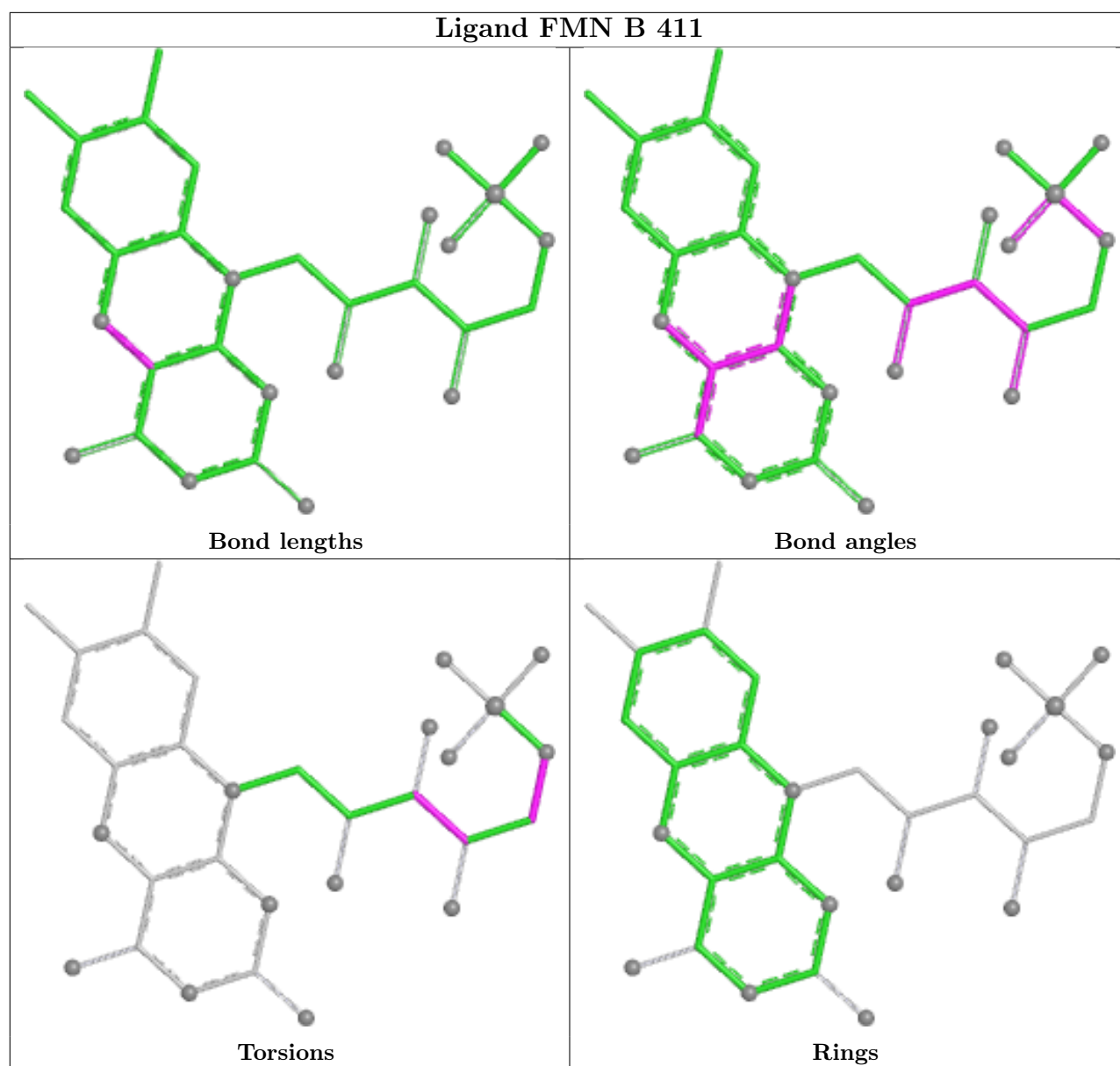
Ligand W83 B 401



Ligand W83 A 401







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	A	313/313 (100%)	-0.14	0 100 100	8, 19, 36, 48	0
1	B	313/313 (100%)	-0.04	8 (2%) 57 47	8, 19, 40, 69	0
All	All	626/626 (100%)	-0.09	8 (1%) 75 66	8, 19, 38, 69	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	132	ASN	3.7
1	B	135	GLY	3.5
1	B	133	VAL	2.8
1	B	134	PRO	2.7
1	B	0	MET	2.6
1	B	130	CYS	2.4
1	B	131	PRO	2.2
1	B	129	SER	2.2

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,

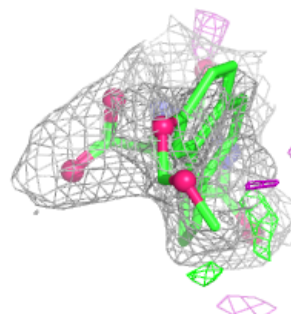
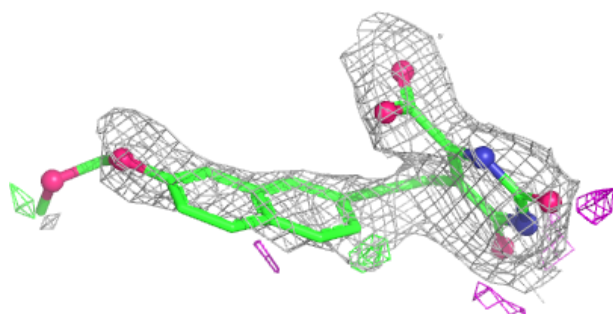
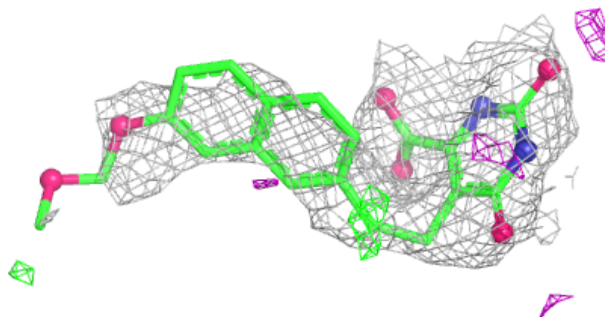
median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	GOL	A	406	6/6	0.70	0.18	36,39,39,40	0
3	GOL	B	406	6/6	0.72	0.21	49,54,55,55	0
3	GOL	B	404	6/6	0.76	0.28	58,59,59,59	0
3	GOL	A	405	6/6	0.76	0.27	44,46,48,48	0
3	GOL	B	409	6/6	0.79	0.13	30,32,32,33	0
3	GOL	A	410	6/6	0.80	0.17	67,68,69,70	0
3	GOL	B	407	6/6	0.83	0.11	50,50,50,51	0
2	W83	B	401	27/27	0.83	0.15	23,38,58,61	0
3	GOL	B	405	6/6	0.85	0.16	70,71,72,72	0
3	GOL	B	402	6/6	0.86	0.10	34,35,35,36	0
3	GOL	A	407	6/6	0.86	0.15	42,44,45,46	0
3	GOL	A	404	6/6	0.88	0.10	39,42,44,45	0
2	W83	A	401	27/27	0.88	0.13	19,30,52,57	0
3	GOL	B	403	6/6	0.88	0.12	36,37,39,41	0
3	GOL	A	409	6/6	0.88	0.10	46,47,48,48	0
3	GOL	B	408	6/6	0.91	0.17	40,41,42,42	0
3	GOL	B	410	6/6	0.91	0.11	47,48,49,49	0
3	GOL	A	408	6/6	0.93	0.11	38,40,41,43	0
3	GOL	A	402	6/6	0.94	0.14	36,37,38,40	0
3	GOL	A	403	6/6	0.94	0.10	21,23,25,25	0
4	FMN	B	411	31/31	0.94	0.09	11,14,15,15	0
4	FMN	A	411	31/31	0.96	0.08	10,13,15,15	0
5	NCO	A	412	7/7	0.96	0.12	10,11,12,12	0

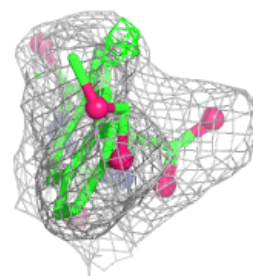
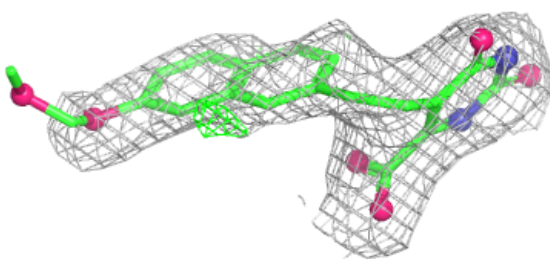
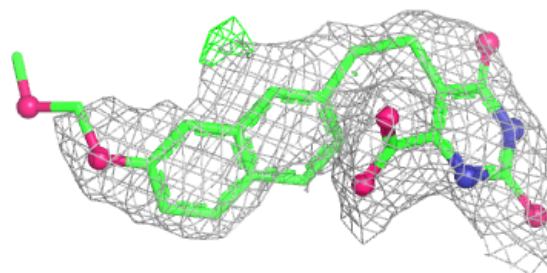
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around W83 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)

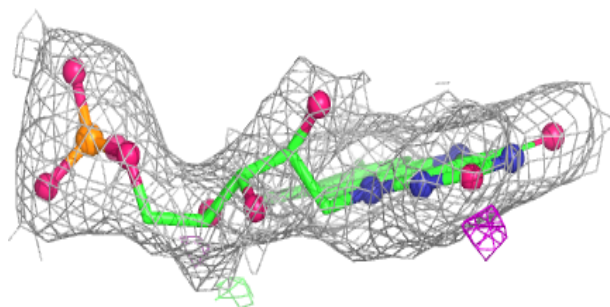
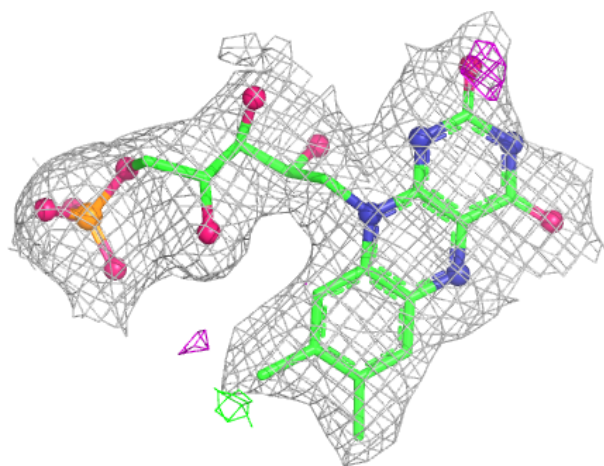
**Electron density around W83 A 401:**

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



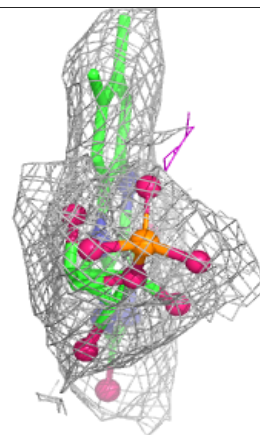
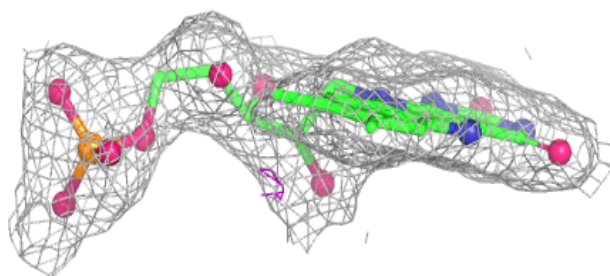
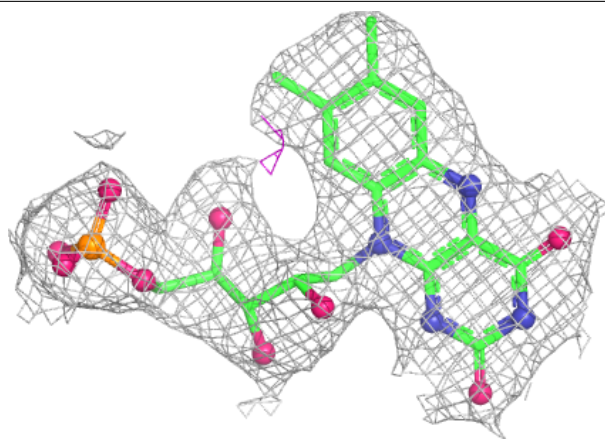
Electron density around FMN B 411:

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



Electron density around FMN A 411:

$2mF_o-DF_c$ (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.