



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 17, 2026 – 10:28 PM UTC

PDB ID : 1DYI / pdb_00001dyi
Title : ISOMORPHOUS CRYSTAL STRUCTURES OF ESCHERICHIA COLI DIHYDROFOLATE REDUCTASE COMPLEXED WITH FOLATE, 5-DEAZAFOLATE AND 5,10-DIDEAZATETRAHYDROFOLATE: MECHANISTIC IMPLICATIONS
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Deposited on : 1994-10-26
Resolution : 1.90 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

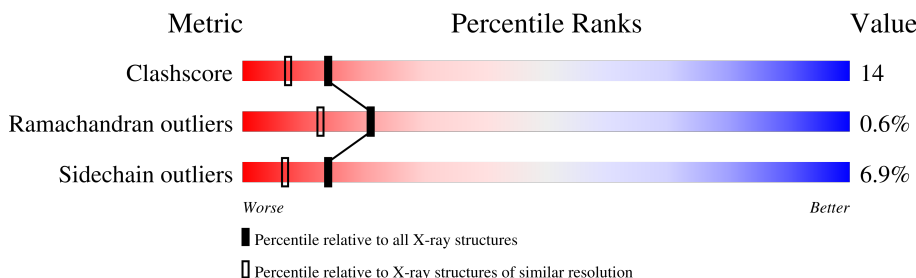
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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))
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	159	
1	B	159	

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	FOL	A	161	-	X	-	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 2994 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 DIHYDROFOLATE REDUCTASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	159	Total	C	N	O	S	0	0	0
			1242	791	210	234	7			
1	B	159	Total	C	N	O	S	0	0	0
			1259	802	216	234	7			

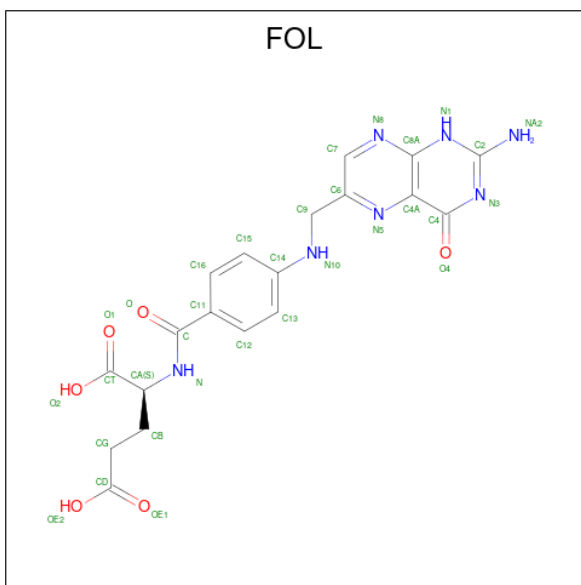
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	37	ASP	ASN	conflict	UNP P0ABQ4
B	37	ASP	ASN	conflict	UNP P0ABQ4

- Molecule 2 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Cl	0	0
			1	1		
2	B	1	Total	Cl	0	0
			1	1		

- Molecule 3 is FOLIC ACID (CCD ID: FOL) (formula: C₁₉H₁₉N₇O₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total 32	C 19	N 7	O 6	0	0
3	B	1	Total 29	C 18	N 7	O 4	0	0

- Molecule 4 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total Ca 1 1	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	235	Total O 235 235	0	0
5	B	194	Total O 194 194	0	0

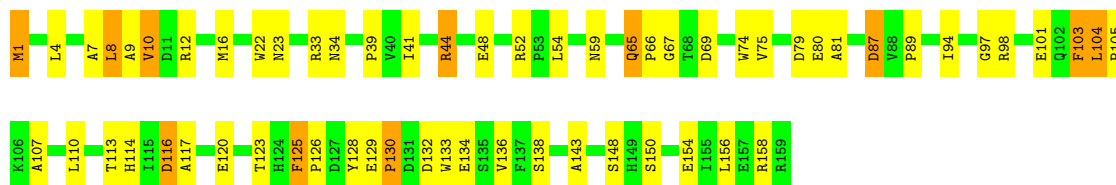
3 Residue-property plots [i](#)

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. 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. 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.

Note EDS was not executed.

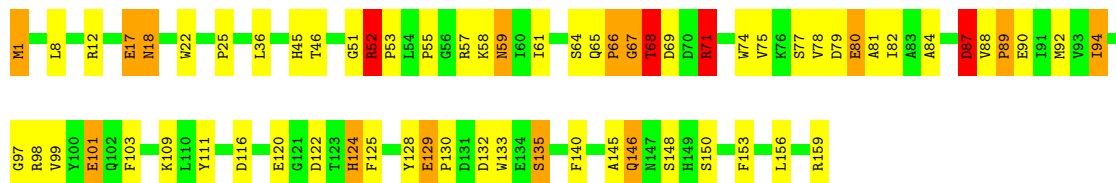
• Molecule 1: DIHYDROFOLATE REDUCTASE

Chain A: 



• Molecule 1: DIHYDROFOLATE REDUCTASE

Chain B: 



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α , β , γ	93.07Å 93.07Å 73.81Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	(Not available) – 1.90	Depositor
% Data completeness (in resolution range)	(Not available) ((Not available)-1.90)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	TNT	Depositor
R, R_{free}	0.137 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2994	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CA, FOL, CL

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	1.24	0/1275	1.91	30/1735 (1.7%)
1	B	1.33	3/1293 (0.2%)	2.05	52/1758 (3.0%)
All	All	1.28	3/2568 (0.1%)	1.98	82/3493 (2.3%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	1	1
All	All	1	2

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	87	ASP	CG-OD2	7.21	1.39	1.25
1	B	135	SER	C-O	-5.51	1.17	1.24
1	B	88	VAL	N-CA	-5.35	1.41	1.46

All (82) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	87	ASP	CA-CB-CG	-12.26	100.34	112.60
1	A	116	ASP	CA-CB-CG	-9.67	102.93	112.60
1	B	94	ILE	N-CA-C	8.94	121.07	111.77
1	B	55	PRO	CB-CA-C	-8.86	99.38	110.98
1	A	94	ILE	N-CA-C	8.30	120.10	112.29
1	B	120	GLU	CA-C-N	-8.27	112.73	122.16

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	120	GLU	C-N-CA	-8.27	112.73	122.16
1	A	87	ASP	CA-CB-CG	-8.07	104.53	112.60
1	B	68	THR	CA-CB-OG1	7.91	121.46	109.60
1	B	153	PHE	CA-CB-CG	-7.72	106.08	113.80
1	B	101	GLU	CB-CA-C	-7.59	98.97	110.88
1	B	103	PHE	CA-CB-CG	7.47	121.27	113.80
1	A	79	ASP	N-CA-CB	7.43	120.84	110.07
1	B	129	GLU	N-CA-CB	7.37	119.89	110.23
1	B	148	SER	N-CA-C	7.31	121.46	112.54
1	A	97	GLY	N-CA-C	7.31	121.48	112.64
1	A	125	PHE	N-CA-C	-7.10	100.81	109.83
1	B	129	GLU	CA-C-O	6.89	124.42	119.32
1	A	130	PRO	N-CA-CB	6.86	110.45	103.25
1	A	79	ASP	CA-C-O	-6.79	113.71	120.70
1	A	54	LEU	N-CA-C	-6.62	101.31	109.65
1	A	103	PHE	CA-CB-CG	-6.56	107.24	113.80
1	B	71	ARG	NE-CZ-NH2	-6.51	113.34	119.20
1	A	69	ASP	N-CA-C	6.42	119.98	109.06
1	B	68	THR	CA-CB-CG2	6.17	120.99	110.50
1	A	120	GLU	CA-C-N	-6.14	110.97	121.17
1	A	120	GLU	C-N-CA	-6.14	110.97	121.17
1	B	17	GLU	N-CA-CB	6.13	119.79	110.28
1	B	109	LYS	N-CA-CB	-6.03	100.31	111.53
1	A	148	SER	N-CA-C	6.01	118.62	111.71
1	B	116	ASP	CA-CB-CG	5.96	118.56	112.60
1	B	101	GLU	CB-CG-CD	-5.95	102.49	112.60
1	B	145	ALA	CA-C-N	-5.94	112.82	122.26
1	B	145	ALA	C-N-CA	-5.94	112.82	122.26
1	A	34	ASN	CA-C-N	-5.92	112.94	122.23
1	A	34	ASN	C-N-CA	-5.92	112.94	122.23
1	B	52	ARG	CD-NE-CZ	5.90	132.66	124.40
1	B	135	SER	CA-C-O	-5.89	113.75	120.58
1	B	80	GLU	N-CA-CB	-5.89	101.47	110.01
1	A	107	ALA	N-CA-C	5.82	118.70	110.50
1	B	87	ASP	CB-CA-C	-5.82	103.21	112.05
1	B	45	HIS	CA-CB-CG	-5.75	108.05	113.80
1	B	140	PHE	CA-CB-CG	-5.71	108.09	113.80
1	B	18	ASN	CA-CB-CG	-5.63	106.97	112.60
1	A	136	VAL	CB-CA-C	-5.59	104.27	111.20
1	B	98	ARG	NE-CZ-NH2	-5.58	114.18	119.20
1	B	12	ARG	CD-NE-CZ	5.56	132.18	124.40
1	B	51	GLY	CA-C-N	-5.54	115.98	123.96

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	51	GLY	C-N-CA	-5.54	115.98	123.96
1	B	92	MET	CA-C-N	-5.54	116.32	123.19
1	B	92	MET	C-N-CA	-5.54	116.32	123.19
1	B	125	PHE	N-CA-C	-5.46	102.90	109.83
1	B	25	PRO	O-C-N	-5.43	115.21	122.22
1	B	122	ASP	N-CA-C	-5.42	105.48	112.68
1	B	57	ARG	CA-C-N	-5.40	115.01	122.30
1	B	57	ARG	C-N-CA	-5.40	115.01	122.30
1	A	39	PRO	N-CA-CB	5.32	108.31	103.15
1	B	146	GLN	CB-CG-CD	-5.29	103.61	112.60
1	B	89	PRO	CA-C-N	-5.27	112.66	121.94
1	B	89	PRO	C-N-CA	-5.27	112.66	121.94
1	A	80	GLU	CB-CG-CD	-5.27	103.65	112.60
1	A	44	ARG	N-CA-C	5.26	117.01	111.28
1	B	66	PRO	CA-C-N	5.24	131.69	121.41
1	B	66	PRO	C-N-CA	5.24	131.69	121.41
1	A	113	THR	CA-CB-OG1	-5.23	101.75	109.60
1	B	124	HIS	N-CA-CB	-5.23	102.45	111.55
1	A	114	HIS	CA-CB-CG	-5.23	108.57	113.80
1	B	8	LEU	CB-CA-C	-5.22	100.16	109.71
1	A	65	GLN	CA-C-N	5.21	125.62	119.99
1	A	65	GLN	C-N-CA	5.21	125.62	119.99
1	A	130	PRO	CB-CA-C	5.21	120.16	111.56
1	A	48	GLU	CB-CA-C	-5.21	102.15	110.79
1	B	55	PRO	N-CA-CB	5.15	108.12	103.33
1	A	143	ALA	N-CA-C	-5.13	103.75	110.53
1	B	150	SER	O-C-N	-5.08	117.06	122.85
1	A	123	THR	N-CA-C	5.08	118.39	110.32
1	B	46	THR	N-CA-C	-5.08	105.66	111.14
1	B	124	HIS	CA-CB-CG	-5.07	108.73	113.80
1	B	111	TYR	N-CA-C	-5.06	99.16	108.02
1	B	79	ASP	N-CA-CB	5.03	117.52	110.12
1	A	16	MET	CG-SD-CE	5.01	111.92	100.90
1	B	120	GLU	O-C-N	-5.01	117.37	123.13

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	B	68	THR	CB

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	23	ASN	Sidechain
1	B	87	ASP	Sidechain

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1242	0	1172	33	1
1	B	1259	0	1215	37	1
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	32	0	15	0	0
3	B	29	0	14	2	0
4	B	1	0	0	1	0
5	A	235	0	0	10	5
5	B	194	0	0	11	8
All	All	2994	0	2416	71	9

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 14.

All (71) 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)
4:B:420:CA:CA	5:B:477:HOH:O	1.24	1.12
1:B:124:HIS:HB3	5:B:597:HOH:O	1.76	0.84
1:B:67:GLY:O	1:B:68:THR:HB	1.82	0.79
1:B:65:GLN:HB3	1:B:66:PRO:HD2	1.64	0.77
1:B:135:SER:HA	1:B:156:LEU:HD23	1.67	0.76
1:A:134:GLU:HG3	5:A:601:HOH:O	1.88	0.73
1:B:69:ASP:OD2	1:B:71:ARG:NH1	2.23	0.71
1:A:1:MET:HE1	5:A:452:HOH:O	1.93	0.68
1:A:10:VAL:HG13	1:A:117:ALA:O	1.93	0.68
1:B:68:THR:HG23	5:B:585:HOH:O	1.95	0.65
1:A:130:PRO:N	5:A:574:HOH:O	2.29	0.65
1:B:65:GLN:HB3	1:B:66:PRO:CD	2.27	0.63
1:B:1:MET:HE3	1:B:90:GLU:HG3	1.80	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:44:ARG:HB3	5:A:573:HOH:O	2.00	0.62
1:A:128:TYR:O	1:A:130:PRO:HD3	2.00	0.62
1:A:67:GLY:HA2	1:A:74:TRP:CE2	2.34	0.61
1:A:67:GLY:HA2	1:A:74:TRP:CD2	2.35	0.61
1:B:146:GLN:HG3	5:B:451:HOH:O	2.02	0.59
1:A:12:ARG:HD2	5:A:546:HOH:O	2.04	0.55
1:B:66:PRO:O	1:B:67:GLY:O	2.25	0.53
1:A:4:LEU:HB2	1:A:110:LEU:HD23	1.89	0.53
1:B:132:ASP:HB3	5:B:614:HOH:O	2.10	0.51
1:B:18:ASN:HB2	5:B:537:HOH:O	2.10	0.51
1:A:125:PHE:CD1	1:A:126:PRO:HD2	2.46	0.51
1:A:98:ARG:NH1	1:A:101:GLU:OE1	2.44	0.50
1:B:77:SER:OG	1:B:80:GLU:HB2	2.12	0.49
1:B:1:MET:N	1:B:89:PRO:O	2.46	0.49
1:A:66:PRO:N	5:A:514:HOH:O	2.45	0.48
1:B:52:ARG:HB2	1:B:53:PRO:HD2	1.96	0.48
1:A:104:LEU:N	1:A:105:PRO:CD	2.78	0.47
1:B:75:VAL:HG21	1:B:81:ALA:HA	1.95	0.47
1:A:9:ALA:HB2	1:A:22:TRP:HZ3	1.78	0.47
1:A:130:PRO:O	1:A:133:TRP:O	2.32	0.47
1:A:103:PHE:C	1:A:105:PRO:HD2	2.40	0.47
1:B:75:VAL:HG21	1:B:81:ALA:CA	2.45	0.47
1:A:129:GLU:O	1:A:132:ASP:HB2	2.15	0.47
1:A:7:ALA:C	1:A:8:LEU:HD23	2.40	0.47
1:A:116:ASP:HB3	5:A:620:HOH:O	2.14	0.47
1:A:44:ARG:HH21	1:A:65:GLN:HB2	1.80	0.46
1:B:52:ARG:HD2	1:B:53:PRO:O	2.16	0.45
1:A:158:ARG:HB3	5:A:621:HOH:O	2.17	0.45
1:B:67:GLY:HA2	1:B:74:TRP:CD2	2.52	0.45
1:B:1:MET:HE3	1:B:90:GLU:CG	2.47	0.44
1:B:78:VAL:O	1:B:82:ILE:HG13	2.16	0.44
1:B:84:ALA:HB2	5:B:548:HOH:O	2.18	0.44
1:A:154:GLU:HG2	1:A:156:LEU:CD1	2.48	0.44
1:A:4:LEU:HB2	1:A:110:LEU:CD2	2.48	0.44
1:B:129:GLU:O	1:B:133:TRP:HD1	2.02	0.43
1:B:64:SER:HB2	1:B:65:GLN:NE2	2.33	0.43
1:B:128:TYR:O	1:B:130:PRO:HD3	2.19	0.43
1:A:154:GLU:HG2	1:A:156:LEU:HD11	1.99	0.43
1:B:97:GLY:O	1:B:101:GLU:HG3	2.19	0.43
1:B:99:VAL:HG23	5:B:541:HOH:O	2.17	0.43
1:A:1:MET:HE2	1:A:89:PRO:O	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:12:ARG:NH2	5:A:597:HOH:O	2.29	0.43
1:B:69:ASP:OD1	1:B:71:ARG:HD2	2.19	0.43
1:B:94:ILE:O	3:B:161:FOL:H7	2.18	0.42
1:B:58:LYS:HA	5:B:604:HOH:O	2.19	0.42
1:B:129:GLU:HA	1:B:130:PRO:HD3	1.80	0.42
1:B:36:LEU:HA	1:B:36:LEU:HD12	1.77	0.42
1:A:75:VAL:HG21	1:A:81:ALA:HB2	2.01	0.41
1:B:52:ARG:NH2	3:B:161:FOL:O	2.54	0.41
1:B:69:ASP:OD1	1:B:71:ARG:CD	2.69	0.41
1:A:8:LEU:HD23	1:A:8:LEU:N	2.35	0.41
1:A:41:ILE:HD13	1:A:81:ALA:HB1	2.02	0.41
1:A:125:PHE:CG	1:A:126:PRO:HD2	2.56	0.40
1:B:22:TRP:HZ2	5:B:515:HOH:O	2.02	0.40
1:A:44:ARG:HB3	5:A:602:HOH:O	2.19	0.40
1:B:59:ASN:HD22	1:B:59:ASN:N	2.19	0.40
1:B:71:ARG:NH2	5:B:576:HOH:O	2.53	0.40
1:A:67:GLY:CA	1:A:74:TRP:CE2	3.04	0.40

All (9) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:B:424:HOH:O	5:B:477:HOH:O[3_564]	0.65	1.55
5:A:622:HOH:O	5:B:612:HOH:O[3_564]	0.92	1.28
5:A:624:HOH:O	5:A:625:HOH:O[5_564]	1.61	0.59
5:A:615:HOH:O	5:B:609:HOH:O[6_654]	1.89	0.31
5:B:424:HOH:O	5:B:460:HOH:O[3_564]	1.98	0.22
5:A:622:HOH:O	5:B:459:HOH:O[2_664]	2.12	0.08
1:B:159:ARG:NH1	5:B:557:HOH:O[2_665]	2.12	0.08
1:A:134:GLU:OE1	5:B:609:HOH:O[6_654]	2.15	0.05
5:A:466:HOH:O	5:B:526:HOH:O[3_564]	2.15	0.05

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	157/159 (99%)	151 (96%)	6 (4%)	0	100	100
1	B	157/159 (99%)	152 (97%)	3 (2%)	2 (1%)	9	3
All	All	314/318 (99%)	303 (96%)	9 (3%)	2 (1%)	21	13

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	67	GLY
1	B	68	THR

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	128/136 (94%)	118 (92%)	10 (8%)	11	5
1	B	133/136 (98%)	125 (94%)	8 (6%)	17	9
All	All	261/272 (96%)	243 (93%)	18 (7%)	14	7

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	8	LEU
1	A	10	VAL
1	A	33	ARG
1	A	52	ARG
1	A	59	ASN
1	A	87	ASP
1	A	104	LEU
1	A	138	SER
1	A	150	SER
1	B	1	MET

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Mol	Chain	Res	Type
1	B	17	GLU
1	B	52	ARG
1	B	59	ASN
1	B	61	ILE
1	B	68	THR
1	B	71	ARG
1	B	87	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	18	ASN
1	A	65	GLN
1	B	45	HIS
1	B	65	GLN
1	B	124	HIS
1	B	146	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 5 ligands modelled in this entry, 3 are monoatomic - leaving 2 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the

expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	FOL	A	161	-	34,34,34	4.70	26 (76%)	43,47,47	2.36	16 (37%)
3	FOL	B	161	-	31,31,34	3.37	18 (58%)	39,43,47	2.65	13 (33%)

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	FOL	A	161	-	-	4/22/22/22	0/3/3/3
3	FOL	B	161	-	-	2/19/19/22	0/3/3/3

All (44) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	161	FOL	C16-C15	9.88	1.54	1.38
3	A	161	FOL	CG-CD	-9.80	1.27	1.50
3	A	161	FOL	C7-N8	-8.61	1.16	1.34
3	A	161	FOL	OE1-CD	7.14	1.45	1.22
3	A	161	FOL	O-C	7.04	1.39	1.23
3	A	161	FOL	C8A-N1	-6.84	1.29	1.38
3	B	161	FOL	C15-C14	6.55	1.50	1.39
3	A	161	FOL	C2-NA2	-6.24	1.19	1.34
3	B	161	FOL	C16-C11	-6.23	1.30	1.39
3	A	161	FOL	C13-C12	5.96	1.48	1.38
3	B	161	FOL	C7-N8	-5.64	1.22	1.34
3	A	161	FOL	C15-C14	-5.08	1.30	1.39
3	B	161	FOL	C16-C15	-5.05	1.30	1.38
3	B	161	FOL	C4A-C8A	-5.02	1.28	1.42
3	B	161	FOL	CA-N	-4.94	1.35	1.45
3	A	161	FOL	C4-N3	-4.93	1.28	1.38
3	A	161	FOL	C9-C6	-4.81	1.40	1.50
3	B	161	FOL	C-N	-4.60	1.23	1.34
3	B	161	FOL	C9-N10	4.54	1.59	1.45
3	A	161	FOL	C13-C14	4.47	1.46	1.39
3	B	161	FOL	O-C	4.46	1.33	1.23
3	A	161	FOL	C4A-C8A	-4.30	1.30	1.42
3	A	161	FOL	C16-C11	-4.19	1.33	1.39
3	A	161	FOL	CA-CT	-4.08	1.42	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	161	FOL	C4-N3	4.06	1.47	1.38
3	B	161	FOL	C8A-N1	3.98	1.43	1.38
3	A	161	FOL	O1-CT	3.80	1.33	1.22
3	A	161	FOL	CB-CG	-3.79	1.41	1.52
3	B	161	FOL	C4A-C4	-3.76	1.40	1.47
3	A	161	FOL	O4-C4	-3.66	1.17	1.23
3	A	161	FOL	CB-CA	3.63	1.61	1.53
3	A	161	FOL	O2-CT	-3.62	1.19	1.30
3	B	161	FOL	C13-C12	3.57	1.44	1.38
3	A	161	FOL	C9-N10	3.16	1.54	1.45
3	B	161	FOL	C14-N10	-3.08	1.30	1.38
3	A	161	FOL	C4A-N5	2.81	1.40	1.34
3	B	161	FOL	C9-C6	2.69	1.56	1.50
3	B	161	FOL	O1-CT	2.46	1.29	1.22
3	B	161	FOL	CG-CB	2.44	1.65	1.51
3	A	161	FOL	C4A-C4	-2.27	1.43	1.47
3	A	161	FOL	C2-N1	-2.24	1.32	1.37
3	A	161	FOL	C8A-N8	2.17	1.38	1.35
3	B	161	FOL	C7-C6	2.11	1.43	1.39
3	A	161	FOL	CA-N	-2.03	1.41	1.45

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	161	FOL	C6-C7-N8	6.28	129.17	123.14
3	B	161	FOL	C15-C16-C11	6.16	127.38	120.80
3	B	161	FOL	NA2-C2-N1	6.13	129.67	116.77
3	B	161	FOL	C12-C13-C14	-5.93	113.47	120.30
3	B	161	FOL	N1-C2-N3	-5.62	113.03	123.32
3	B	161	FOL	C16-C15-C14	-5.05	114.49	120.30
3	A	161	FOL	NA2-C2-N1	5.03	127.36	116.77
3	A	161	FOL	OE1-CD-CG	-4.90	107.55	123.09
3	A	161	FOL	CB-CA-CT	-4.22	100.34	110.35
3	B	161	FOL	C7-C6-N5	-4.08	118.22	120.87
3	A	161	FOL	C13-C12-C11	4.04	125.11	120.80
3	A	161	FOL	OE2-CD-OE1	3.88	133.31	123.33
3	A	161	FOL	C7-C6-N5	-3.43	118.65	120.87
3	B	161	FOL	C13-C14-N10	-3.26	113.59	120.96
3	A	161	FOL	C16-C15-C14	3.25	124.05	120.30
3	A	161	FOL	CB-CG-CD	-3.07	104.31	112.49
3	A	161	FOL	N1-C2-N3	-2.97	117.89	123.32
3	A	161	FOL	C16-C11-C12	-2.93	114.84	118.57

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	161	FOL	C9-N10-C14	2.84	130.98	121.87
3	B	161	FOL	C6-C7-N8	2.66	125.69	123.14
3	A	161	FOL	CG-CB-CA	-2.65	108.28	113.16
3	B	161	FOL	O4-C4-N3	2.57	124.14	120.23
3	B	161	FOL	C15-C14-C13	2.55	122.42	119.04
3	A	161	FOL	NA2-C2-N3	-2.53	114.73	119.67
3	A	161	FOL	C12-C13-C14	-2.53	117.39	120.30
3	B	161	FOL	C8A-C4A-C4	-2.48	117.95	119.56
3	A	161	FOL	C9-C6-C7	2.25	125.41	121.38
3	A	161	FOL	O4-C4-C4A	2.11	124.12	120.42
3	B	161	FOL	O-C-N	-2.05	118.56	122.47

There are no chirality outliers.

All (6) torsion outliers are listed below:

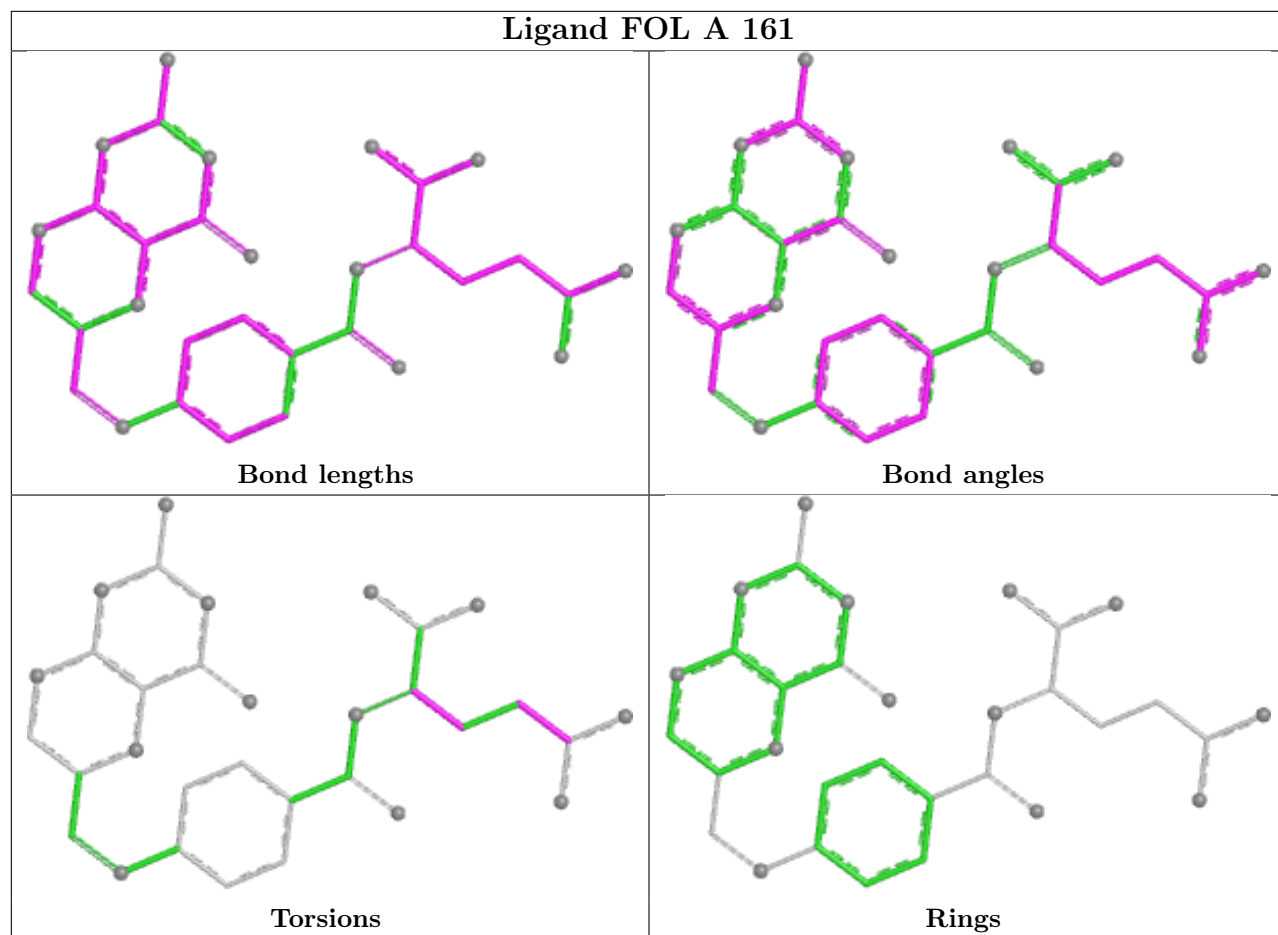
Mol	Chain	Res	Type	Atoms
3	B	161	FOL	CT-CA-CB-CG
3	A	161	FOL	N-CA-CB-CG
3	A	161	FOL	CT-CA-CB-CG
3	B	161	FOL	N-CA-CB-CG
3	A	161	FOL	OE1-CD-CG-CB
3	A	161	FOL	OE2-CD-CG-CB

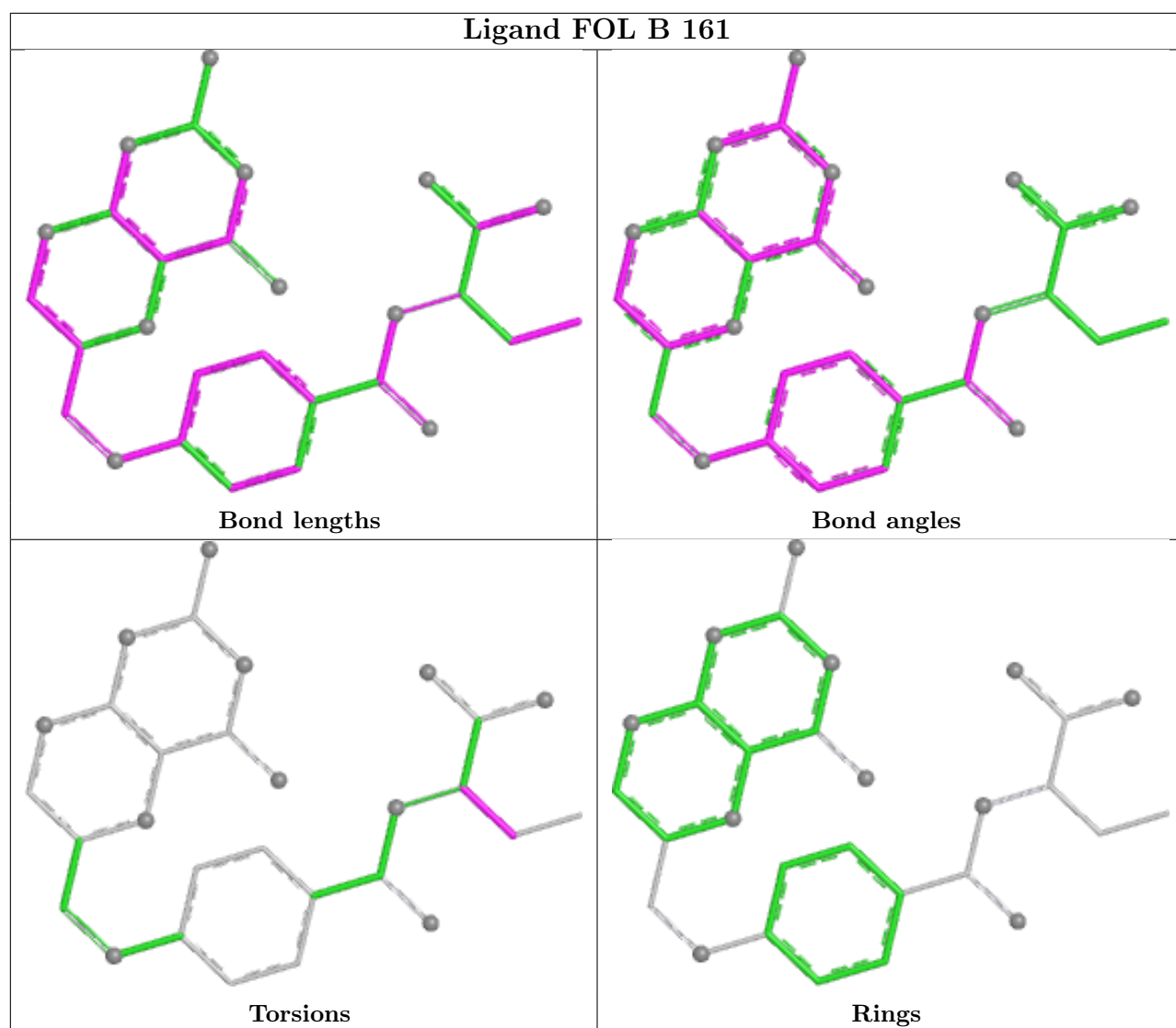
There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	161	FOL	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

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.