



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 06:46 AM UTC

PDB ID : 1DRV / pdb_00001drv
Title : ESCHERICHIA COLI DHPR/ACNADH COMPLEX
Authors : Reddy, S.G.; Scapin, G.; Blanchard, J.S.
Deposited on : 1996-06-28
Resolution : 2.20 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

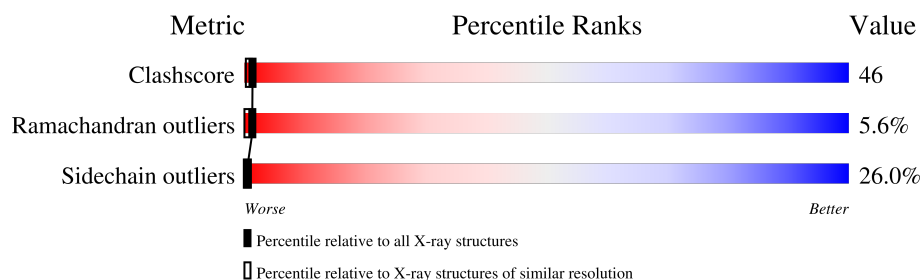
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 Å.

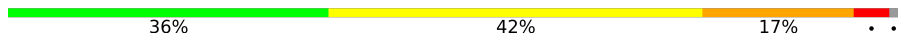
Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)

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	273	

2 Entry composition [i](#)

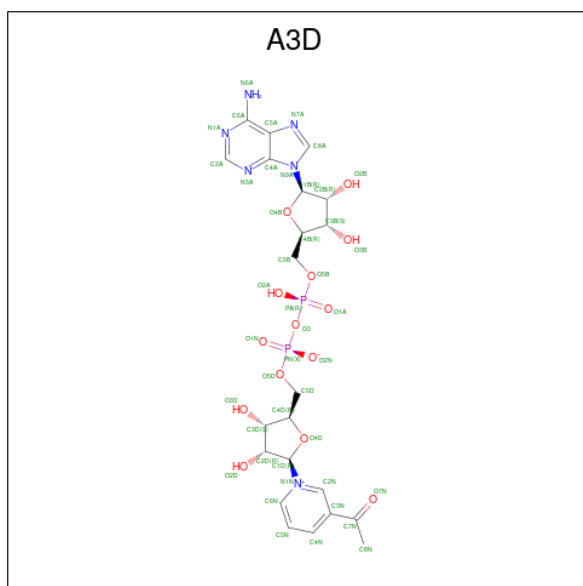
There are 3 unique types of molecules in this entry. The entry contains 2083 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 DIHYDRODIPICOLINATE REDUCTASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	270	1970	1224	356	380	10	0	0	0

- Molecule 2 is 3-ACETILPYRIDINE ADENINE DINUCLEOTIDE (CCD ID: A3D) (formula: $C_{22}H_{28}N_6O_{14}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	A	1	44	22	6	14	2	0	0

- Molecule 3 is water.

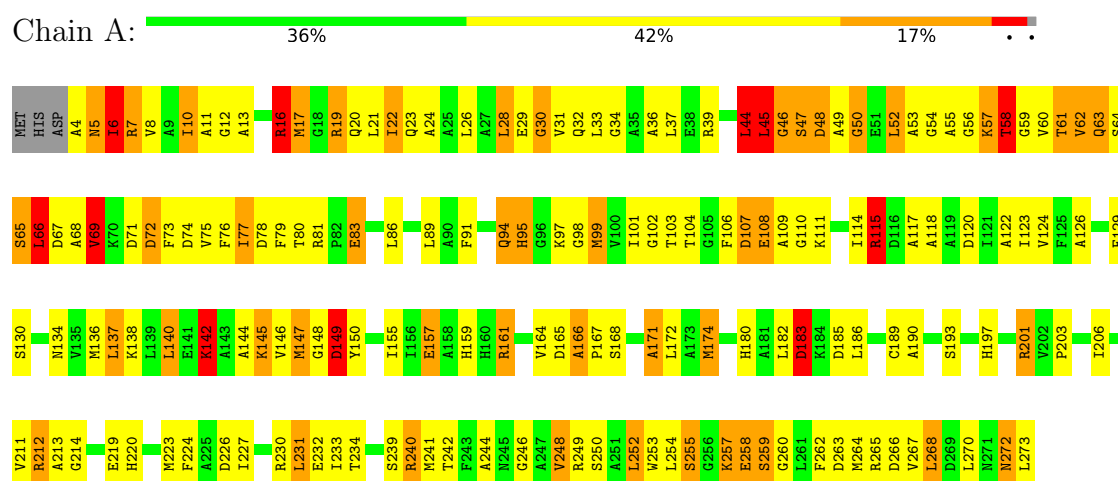
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	69	Total 69 O 69	0	0

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. 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: DIHYDRODIPICOLINATE REDUCTASE



4 Data and refinement statistics

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

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	81.60 Å 84.60 Å 91.90 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.20	Depositor
% Data completeness (in resolution range)	84.0 (20.00-2.20)	Depositor
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
Refinement program	TNT	Depositor
R, R_{free}	0.191 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2083	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

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: A3D

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	1.39	14/1996 (0.7%)	1.79	45/2692 (1.7%)

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	4

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	233	ILE	CA-CB	-9.27	1.42	1.54
1	A	114	ILE	CA-CB	-7.75	1.45	1.54
1	A	99	MET	SD-CE	-7.60	1.60	1.79
1	A	248	VAL	CA-CB	-7.28	1.45	1.54
1	A	206	ILE	CA-CB	-7.02	1.45	1.53
1	A	166	ALA	CA-CB	-6.96	1.45	1.54
1	A	8	VAL	CA-C	-6.44	1.45	1.52
1	A	101	ILE	CA-CB	6.30	1.61	1.53
1	A	10	ILE	CA-CB	5.85	1.60	1.53
1	A	159	HIS	C-O	5.70	1.30	1.23
1	A	172	LEU	CA-C	-5.46	1.45	1.52
1	A	115	ARG	CA-C	-5.35	1.46	1.52
1	A	260	GLY	N-CA	5.34	1.50	1.45
1	A	211	VAL	CA-CB	-5.31	1.47	1.54

All (45) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	149	ASP	O-C-N	12.71	139.49	122.59
1	A	166	ALA	CA-C-N	12.39	135.33	119.84
1	A	166	ALA	C-N-CA	12.39	135.33	119.84
1	A	5	ASN	O-C-N	9.59	135.35	122.59
1	A	142	LYS	N-CA-C	-7.60	102.68	110.97
1	A	107	ASP	N-CA-C	-7.57	101.10	110.41
1	A	144	ALA	N-CA-C	-7.07	103.62	111.82
1	A	11	ALA	N-CA-C	-6.96	101.55	110.53
1	A	171	ALA	N-CA-C	-6.89	103.77	111.28
1	A	120	ASP	N-CA-C	-6.82	105.59	114.04
1	A	174	MET	N-CA-C	-6.81	103.78	111.14
1	A	98	GLY	N-CA-C	-6.76	102.28	113.02
1	A	30	GLY	N-CA-C	-6.70	105.50	114.95
1	A	189	CYS	N-CA-C	6.64	121.05	112.41
1	A	206	ILE	CB-CA-C	-6.58	102.05	111.19
1	A	16	ARG	N-CA-C	6.57	118.99	111.11
1	A	166	ALA	C-N-CD	-6.39	98.78	125.00
1	A	265	ARG	N-CA-C	-6.38	104.41	111.82
1	A	67	ASP	N-CA-C	-6.34	104.42	111.71
1	A	164	VAL	CA-C-N	-6.15	113.72	122.94
1	A	164	VAL	C-N-CA	-6.15	113.72	122.94
1	A	19	ARG	CA-C-N	5.98	129.40	120.31
1	A	19	ARG	C-N-CA	5.98	129.40	120.31
1	A	149	ASP	CA-C-N	5.90	132.28	122.54
1	A	149	ASP	C-N-CA	5.90	132.28	122.54
1	A	104	THR	CB-CA-C	-5.86	99.92	110.36
1	A	213	ALA	N-CA-C	5.84	117.28	108.46
1	A	7	ARG	CA-C-N	-5.83	113.42	122.69
1	A	7	ARG	C-N-CA	-5.83	113.42	122.69
1	A	99	MET	N-CA-C	5.82	118.69	109.50
1	A	201	ARG	NE-CZ-NH1	-5.78	115.72	121.50
1	A	182	LEU	N-CA-C	-5.72	106.15	113.02
1	A	146	VAL	CB-CA-C	-5.71	106.86	111.71
1	A	78	ASP	N-CA-C	5.70	117.99	108.02
1	A	193	SER	N-CA-C	-5.61	100.83	109.52
1	A	37	LEU	CA-CB-CG	-5.59	96.72	116.30
1	A	272	ASN	N-CA-C	-5.55	106.17	113.17
1	A	59	GLY	N-CA-C	-5.50	106.69	115.66
1	A	183	ASP	N-CA-C	5.37	118.81	111.39
1	A	7	ARG	NE-CZ-NH2	5.37	124.03	119.20
1	A	230	ARG	CB-CA-C	-5.36	100.27	109.65
1	A	29	GLU	N-CA-C	5.30	117.39	108.96
1	A	6	ILE	O-C-N	5.24	129.12	122.57

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	126	ALA	N-CA-C	5.20	117.16	108.99
1	A	63	GLN	N-CA-C	5.07	117.34	110.35

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	155	ILE	Mainchain
1	A	16	ARG	Sidechain
1	A	166	ALA	Peptide,Mainchain

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	1970	0	1933	183	0
2	A	44	0	27	5	0
3	A	69	0	0	3	1
All	All	2083	0	1960	183	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 46.

All (183) 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:4:ALA:HB1	1:A:32:GLN:NE2	1.38	1.38
1:A:4:ALA:CB	1:A:32:GLN:HE21	1.51	1.20
1:A:4:ALA:CB	1:A:32:GLN:NE2	2.08	1.12
1:A:44:LEU:HD13	1:A:45:LEU:H	0.99	1.11
1:A:161:ARG:HH11	1:A:161:ARG:HG2	1.16	1.06
1:A:16:ARG:NE	1:A:240:ARG:HH21	1.59	1.00
1:A:57:LYS:O	1:A:58:THR:OG1	1.80	0.97
1:A:4:ALA:HB1	1:A:32:GLN:HE21	0.78	0.95
1:A:5:ASN:HB2	1:A:30:GLY:O	1.67	0.93

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:44:LEU:HD13	1:A:45:LEU:N	1.84	0.93
1:A:66:LEU:HD23	1:A:91:PHE:CG	2.06	0.91
1:A:161:ARG:HH12	1:A:214:GLY:HA2	1.35	0.90
1:A:44:LEU:CD1	1:A:45:LEU:H	1.83	0.89
1:A:65:SER:OG	1:A:68:ALA:HB3	1.72	0.89
1:A:73:PHE:CE1	1:A:97:LYS:HG3	2.07	0.88
1:A:24:ALA:HB2	1:A:241:MET:CE	2.06	0.85
1:A:46:GLY:CA	1:A:61:THR:HG22	2.07	0.84
1:A:44:LEU:HD22	1:A:62:VAL:HG13	1.59	0.84
1:A:16:ARG:HE	1:A:240:ARG:HH21	1.26	0.81
1:A:148:GLY:O	1:A:149:ASP:HB2	1.78	0.81
1:A:17:MET:O	1:A:21:LEU:HD12	1.81	0.79
1:A:24:ALA:HB2	1:A:241:MET:HE3	1.63	0.78
1:A:254:LEU:C	1:A:257:LYS:HG2	2.10	0.77
1:A:161:ARG:HH11	1:A:161:ARG:CG	1.98	0.77
1:A:66:LEU:HD23	1:A:91:PHE:CD2	2.20	0.77
1:A:65:SER:CB	1:A:68:ALA:HB3	2.13	0.76
1:A:48:ASP:O	1:A:52:LEU:HB2	1.85	0.76
1:A:16:ARG:HG3	1:A:240:ARG:NH2	2.02	0.75
1:A:81:ARG:HG2	1:A:81:ARG:HH11	1.52	0.75
1:A:46:GLY:O	1:A:61:THR:HA	1.87	0.74
1:A:65:SER:OG	1:A:66:LEU:N	2.15	0.74
1:A:264:MET:O	1:A:268:LEU:HB2	1.87	0.74
1:A:180:HIS:O	1:A:183:ASP:N	2.22	0.73
1:A:16:ARG:NE	1:A:240:ARG:NH2	2.37	0.73
1:A:252:LEU:O	1:A:252:LEU:HD23	1.90	0.72
1:A:254:LEU:O	1:A:257:LYS:HG2	1.89	0.71
1:A:72:ASP:N	1:A:72:ASP:OD1	2.22	0.71
1:A:223:MET:HE2	1:A:232:GLU:OE2	1.90	0.70
1:A:244:ALA:O	1:A:248:VAL:HG23	1.91	0.70
1:A:16:ARG:CD	1:A:240:ARG:HH21	2.05	0.70
1:A:44:LEU:CD2	1:A:62:VAL:HG13	2.23	0.69
1:A:118:ALA:HA	1:A:123:ILE:HD12	1.75	0.69
1:A:138:LYS:HA	3:A:454:HOH:O	1.93	0.69
1:A:157:GLU:O	1:A:157:GLU:HG3	1.92	0.68
1:A:122:ALA:HB2	1:A:258:GLU:O	1.93	0.68
1:A:44:LEU:O	1:A:45:LEU:HB2	1.93	0.68
1:A:46:GLY:C	1:A:61:THR:HA	2.19	0.68
1:A:36:ALA:O	1:A:62:VAL:HA	1.94	0.67
1:A:4:ALA:HB3	1:A:32:GLN:NE2	2.06	0.67
1:A:111:LYS:HB2	1:A:115:ARG:HH21	1.58	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:46:GLY:HA2	1:A:61:THR:HG22	1.76	0.66
1:A:86:LEU:HD23	1:A:89:LEU:HD23	1.77	0.66
1:A:107:ASP:OD2	1:A:109:ALA:HB3	1.95	0.66
1:A:263:ASP:O	1:A:267:VAL:HG23	1.96	0.65
1:A:21:LEU:HD22	1:A:77:ILE:HG21	1.77	0.65
1:A:157:GLU:HA	1:A:219:GLU:O	1.97	0.65
1:A:268:LEU:HB3	1:A:270:LEU:HG	1.80	0.64
1:A:161:ARG:NH1	1:A:214:GLY:HA2	2.10	0.64
1:A:22:ILE:HG21	1:A:49:ALA:CA	2.28	0.63
1:A:201:ARG:O	1:A:203:PRO:HD3	1.98	0.63
1:A:44:LEU:HD22	1:A:62:VAL:CG1	2.27	0.63
1:A:24:ALA:O	1:A:28:LEU:HB2	1.99	0.63
1:A:142:LYS:NZ	1:A:273:LEU:O	2.32	0.62
1:A:50:GLY:HA2	1:A:56:GLY:HA2	1.82	0.62
1:A:5:ASN:OD1	1:A:30:GLY:HA2	2.00	0.61
1:A:4:ALA:HB1	1:A:32:GLN:CD	2.20	0.61
1:A:58:THR:C	1:A:60:VAL:H	2.08	0.61
1:A:161:ARG:HG2	1:A:161:ARG:NH1	1.95	0.61
1:A:48:ASP:N	1:A:60:VAL:O	2.34	0.60
1:A:226:ASP:OD1	1:A:227:ILE:N	2.29	0.60
1:A:46:GLY:HA3	1:A:61:THR:HG22	1.84	0.60
1:A:80:THR:O	1:A:81:ARG:NH1	2.31	0.59
1:A:157:GLU:OE1	1:A:212:ARG:NH1	2.26	0.59
1:A:83:GLU:H	1:A:83:GLU:CD	2.10	0.58
1:A:81:ARG:HG2	1:A:81:ARG:NH1	2.18	0.58
1:A:49:ALA:O	1:A:50:GLY:O	2.22	0.57
1:A:66:LEU:HB3	1:A:91:PHE:CE2	2.39	0.57
1:A:124:VAL:HG22	1:A:262:PHE:HB2	1.86	0.57
1:A:6:ILE:O	1:A:6:ILE:CG2	2.52	0.57
1:A:24:ALA:HB2	1:A:241:MET:HE1	1.85	0.57
1:A:148:GLY:O	1:A:149:ASP:CB	2.51	0.57
1:A:44:LEU:CD2	1:A:46:GLY:H	2.16	0.57
1:A:17:MET:HG3	1:A:79:PHE:CD2	2.40	0.56
1:A:6:ILE:O	1:A:6:ILE:HG22	2.05	0.56
1:A:44:LEU:HD22	1:A:46:GLY:H	1.69	0.56
1:A:47:SER:HA	1:A:58:THR:O	2.04	0.56
1:A:50:GLY:CA	1:A:56:GLY:HA2	2.35	0.56
1:A:103:THR:O	1:A:106:PHE:HE1	1.88	0.56
1:A:186:LEU:HD12	1:A:190:ALA:HB2	1.88	0.56
1:A:22:ILE:HG22	1:A:23:GLN:N	2.21	0.55
1:A:44:LEU:HD22	1:A:46:GLY:N	2.20	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:62:VAL:HG22	1:A:62:VAL:O	2.05	0.55
1:A:118:ALA:CA	1:A:123:ILE:HD12	2.35	0.55
1:A:34:GLY:HA3	1:A:72:ASP:O	2.07	0.55
1:A:161:ARG:HH12	1:A:214:GLY:CA	2.12	0.55
1:A:69:VAL:HG12	1:A:73:PHE:CD2	2.41	0.55
1:A:111:LYS:HB2	1:A:115:ARG:NH2	2.23	0.54
1:A:268:LEU:HD23	1:A:270:LEU:HD11	1.90	0.54
1:A:65:SER:O	1:A:66:LEU:HB2	2.08	0.54
1:A:108:GLU:OE1	1:A:108:GLU:O	2.26	0.54
1:A:137:LEU:HA	1:A:140:LEU:HD22	1.90	0.54
1:A:69:VAL:CG1	1:A:73:PHE:HD2	2.21	0.53
1:A:89:LEU:HD11	1:A:117:ALA:CB	2.38	0.53
1:A:263:ASP:O	1:A:266:ASP:HB2	2.09	0.53
1:A:53:ALA:O	1:A:55:ALA:N	2.42	0.52
1:A:66:LEU:HD23	1:A:91:PHE:CD1	2.42	0.52
1:A:252:LEU:O	1:A:255:SER:HB2	2.09	0.52
1:A:16:ARG:HE	1:A:240:ARG:NH2	2.02	0.52
1:A:50:GLY:HA2	1:A:56:GLY:CA	2.40	0.52
1:A:145:LYS:O	1:A:145:LYS:HG3	2.09	0.52
1:A:174:MET:SD	1:A:220:HIS:HB3	2.49	0.52
1:A:16:ARG:CG	1:A:240:ARG:NH2	2.72	0.51
1:A:171:ALA:O	1:A:174:MET:HB2	2.11	0.51
1:A:224:PHE:CD1	1:A:224:PHE:N	2.78	0.51
1:A:80:THR:O	1:A:81:ARG:HG2	2.11	0.50
1:A:201:ARG:HG2	1:A:201:ARG:HH11	1.76	0.50
1:A:12:GLY:HA2	2:A:301:A3D:O3B	2.11	0.50
1:A:157:GLU:OE2	1:A:168:SER:HB3	2.12	0.50
1:A:129:PHE:O	1:A:246:GLY:HA3	2.13	0.49
1:A:17:MET:HB2	2:A:301:A3D:O2N	2.12	0.49
1:A:50:GLY:HA2	1:A:56:GLY:C	2.37	0.49
1:A:66:LEU:CD2	1:A:91:PHE:CG	2.90	0.49
1:A:33:LEU:C	1:A:33:LEU:HD12	2.35	0.48
1:A:58:THR:C	1:A:60:VAL:N	2.72	0.48
1:A:223:MET:HA	1:A:231:LEU:O	2.14	0.47
1:A:24:ALA:CA	1:A:241:MET:HE1	2.45	0.47
1:A:69:VAL:HG12	1:A:73:PHE:CE2	2.49	0.47
1:A:73:PHE:CD1	1:A:73:PHE:C	2.93	0.47
1:A:102:GLY:O	2:A:301:A3D:H1D	2.15	0.47
1:A:45:LEU:O	1:A:46:GLY:C	2.58	0.47
1:A:130:SER:O	1:A:134:ASN:ND2	2.48	0.47
1:A:69:VAL:CG1	1:A:73:PHE:CD2	2.99	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:20:GLN:NE2	1:A:241:MET:SD	2.88	0.46
1:A:80:THR:C	1:A:81:ARG:HG2	2.41	0.46
1:A:13:ALA:O	1:A:52:LEU:HD13	2.15	0.46
1:A:33:LEU:HD12	1:A:33:LEU:HA	1.55	0.46
1:A:44:LEU:HB3	3:A:426:HOH:O	2.16	0.46
1:A:31:VAL:HG13	1:A:252:LEU:HD12	1.98	0.46
1:A:253:TRP:HZ3	1:A:262:PHE:CD2	2.34	0.45
1:A:24:ALA:CB	1:A:241:MET:CE	2.88	0.45
1:A:226:ASP:CG	1:A:227:ILE:H	2.22	0.45
1:A:197:HIS:ND1	1:A:197:HIS:C	2.74	0.45
1:A:161:ARG:NH1	1:A:214:GLY:CA	2.77	0.45
1:A:73:PHE:O	1:A:97:LYS:HE3	2.17	0.44
1:A:118:ALA:CB	1:A:123:ILE:HD12	2.47	0.44
1:A:16:ARG:CD	1:A:240:ARG:NH2	2.78	0.44
1:A:75:VAL:CG1	1:A:76:PHE:N	2.80	0.44
1:A:108:GLU:CD	1:A:115:ARG:HH22	2.24	0.44
1:A:46:GLY:O	1:A:48:ASP:N	2.51	0.44
1:A:220:HIS:O	1:A:234:THR:HA	2.18	0.44
1:A:108:GLU:OE2	1:A:115:ARG:NH2	2.43	0.44
1:A:258:GLU:O	1:A:259:SER:C	2.61	0.44
1:A:254:LEU:HA	1:A:257:LYS:CG	2.47	0.43
1:A:22:ILE:HG21	1:A:53:ALA:HB3	2.01	0.43
1:A:48:ASP:C	1:A:50:GLY:N	2.74	0.43
1:A:118:ALA:HB2	1:A:123:ILE:HD12	1.99	0.43
1:A:157:GLU:OE2	1:A:168:SER:N	2.36	0.43
1:A:33:LEU:O	1:A:60:VAL:HG21	2.19	0.43
1:A:47:SER:O	1:A:48:ASP:O	2.37	0.43
1:A:212:ARG:HD3	1:A:212:ARG:HA	1.73	0.43
1:A:254:LEU:HA	1:A:257:LYS:HG3	2.02	0.42
1:A:50:GLY:C	1:A:56:GLY:HA2	2.44	0.42
1:A:102:GLY:O	2:A:301:A3D:H2N	2.19	0.42
1:A:124:VAL:HG23	1:A:254:LEU:HD21	2.00	0.42
1:A:24:ALA:CB	1:A:241:MET:HE1	2.50	0.42
1:A:12:GLY:CA	2:A:301:A3D:O3B	2.68	0.42
1:A:52:LEU:HD23	1:A:52:LEU:HA	1.69	0.42
1:A:66:LEU:CD2	1:A:91:PHE:CD2	2.98	0.42
1:A:138:LYS:HE2	1:A:270:LEU:O	2.20	0.42
1:A:137:LEU:HD12	1:A:140:LEU:CD2	2.49	0.42
1:A:108:GLU:C	1:A:110:GLY:N	2.75	0.41
1:A:147:MET:HA	1:A:150:TYR:CE2	2.55	0.41
1:A:94:GLN:HB3	1:A:95:HIS:ND1	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:111:LYS:C	1:A:115:ARG:HH21	2.28	0.41
1:A:223:MET:O	1:A:223:MET:HG3	2.20	0.41
1:A:180:HIS:CD2	1:A:180:HIS:C	2.95	0.41
1:A:142:LYS:NZ	1:A:273:LEU:C	2.78	0.41
1:A:150:TYR:HE1	3:A:408:HOH:O	2.02	0.41
1:A:7:ARG:HB2	1:A:74:ASP:OD1	2.20	0.41
1:A:24:ALA:N	1:A:241:MET:HE1	2.36	0.41
1:A:95:HIS:ND1	1:A:95:HIS:N	2.68	0.40
1:A:73:PHE:CZ	1:A:97:LYS:HG3	2.54	0.40
1:A:250:SER:O	1:A:253:TRP:HB3	2.20	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)
3:A:406:HOH:O	3:A:406:HOH:O[3_556]	2.19	0.01

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	268/273 (98%)	219 (82%)	34 (13%)	15 (6%)	1 0

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	44	LEU
1	A	45	LEU
1	A	48	ASP
1	A	65	SER
1	A	66	LEU
1	A	149	ASP

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Mol	Chain	Res	Type
1	A	259	SER
1	A	46	GLY
1	A	47	SER
1	A	50	GLY
1	A	54	GLY
1	A	57	LYS
1	A	58	THR
1	A	167	PRO
1	A	69	VAL

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	192/201 (96%)	142 (74%)	50 (26%)	0 0

All (50) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	ILE
1	A	10	ILE
1	A	17	MET
1	A	19	ARG
1	A	22	ILE
1	A	26	LEU
1	A	28	LEU
1	A	39	ARG
1	A	44	LEU
1	A	45	LEU
1	A	52	LEU
1	A	58	THR
1	A	61	THR
1	A	62	VAL
1	A	63	GLN
1	A	64	SER
1	A	66	LEU

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Mol	Chain	Res	Type
1	A	69	VAL
1	A	71	ASP
1	A	72	ASP
1	A	77	ILE
1	A	83	GLU
1	A	94	GLN
1	A	95	HIS
1	A	99	MET
1	A	108	GLU
1	A	115	ARG
1	A	136	MET
1	A	137	LEU
1	A	140	LEU
1	A	142	LYS
1	A	145	LYS
1	A	147	MET
1	A	157	GLU
1	A	161	ARG
1	A	165	ASP
1	A	183	ASP
1	A	185	ASP
1	A	212	ARG
1	A	231	LEU
1	A	239	SER
1	A	240	ARG
1	A	242	THR
1	A	249	ARG
1	A	252	LEU
1	A	255	SER
1	A	257	LYS
1	A	258	GLU
1	A	268	LEU
1	A	272	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	20	GLN
1	A	23	GLN
1	A	32	GLN
1	A	88	HIS
1	A	94	GLN

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Mol	Chain	Res	Type
1	A	180	HIS
1	A	220	HIS

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 ⓘ

1 ligand is 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	$\# Z > 2$
2	A3D	A	301	-	46,48,48	2.25	6 (13%)	64,73,73	2.24	14 (21%)

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
2	A3D	A	301	-	-	11/30/62/62	0/5/5/5

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	301	A3D	PN-O3	-10.60	1.48	1.59
2	A	301	A3D	PA-O3	-7.09	1.51	1.59
2	A	301	A3D	C3N-C7N	-3.99	1.42	1.49
2	A	301	A3D	C2N-C3N	-3.54	1.33	1.39
2	A	301	A3D	O4D-C1D	3.17	1.45	1.40
2	A	301	A3D	C4N-C3N	-2.22	1.36	1.39

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	301	A3D	O3-PN-O1N	-8.38	85.48	110.70
2	A	301	A3D	C2N-C3N-C4N	-7.14	109.95	118.26
2	A	301	A3D	C5N-C4N-C3N	6.65	126.91	120.36
2	A	301	A3D	O2N-PN-O3	5.35	121.73	107.27
2	A	301	A3D	O3-PA-O1A	4.37	123.85	110.70
2	A	301	A3D	C6N-N1N-C2N	-3.59	118.82	121.88
2	A	301	A3D	C2B-C1B-N9A	-2.98	105.89	113.30
2	A	301	A3D	O3D-C3D-C4D	2.77	119.04	111.08
2	A	301	A3D	C4N-C3N-C7N	2.70	126.96	120.83
2	A	301	A3D	C5N-C6N-N1N	-2.54	116.92	120.38
2	A	301	A3D	C3B-C2B-C1B	-2.36	97.00	101.46
2	A	301	A3D	O2B-C2B-C3B	2.34	119.31	111.82
2	A	301	A3D	O3B-C3B-C4B	-2.28	104.53	111.08
2	A	301	A3D	PN-O5D-C5D	-2.16	109.00	121.35

There are no chirality outliers.

All (11) torsion outliers are listed below:

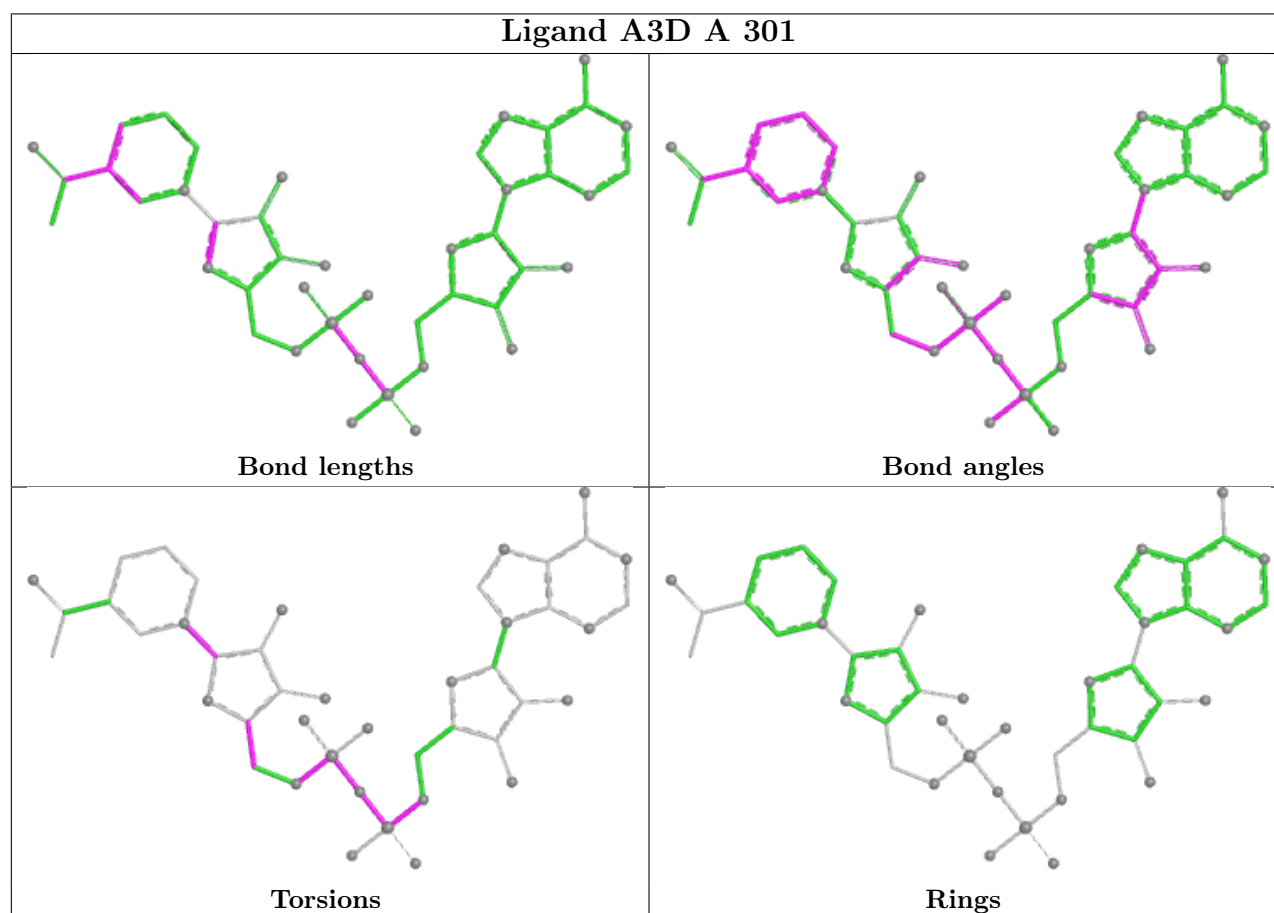
Mol	Chain	Res	Type	Atoms
2	A	301	A3D	C5D-O5D-PN-O3
2	A	301	A3D	O4D-C1D-N1N-C2N
2	A	301	A3D	O4D-C1D-N1N-C6N
2	A	301	A3D	C2D-C1D-N1N-C2N
2	A	301	A3D	C2D-C1D-N1N-C6N
2	A	301	A3D	C3D-C4D-C5D-O5D
2	A	301	A3D	O4D-C4D-C5D-O5D
2	A	301	A3D	PN-O3-PA-O5B
2	A	301	A3D	C5B-O5B-PA-O1A
2	A	301	A3D	PA-O3-PN-O2N
2	A	301	A3D	PA-O3-PN-O1N

There are no ring outliers.

1 monomer is involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	301	A3D	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.



5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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.