



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:24 PM UTC

PDB ID : 4PSG / pdb_00004psg
Title : Crystal Structure of C.elegans Thymidylate Synthase in complex with an inhibitor N(4)OHdCMP
Authors : Wilk, P.; Jarmula, A.; Rode, W.
Deposited on : 2014-03-07
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 ⓘ](#)) 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

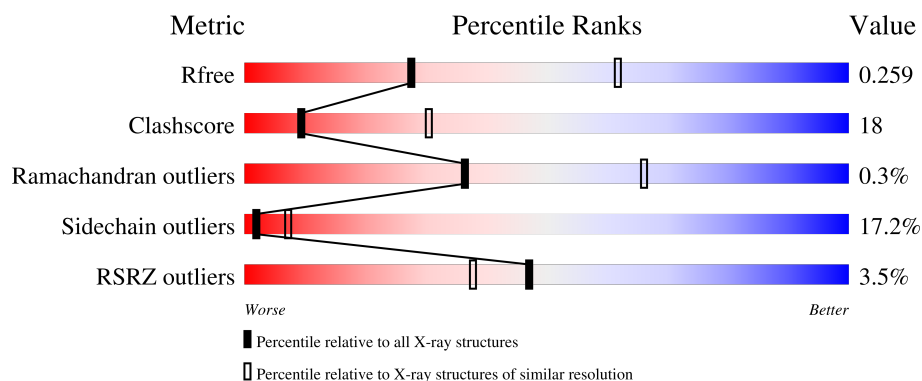
1 Overall quality at a glance

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	315	
1	B	315	

2 Entry composition [i](#)

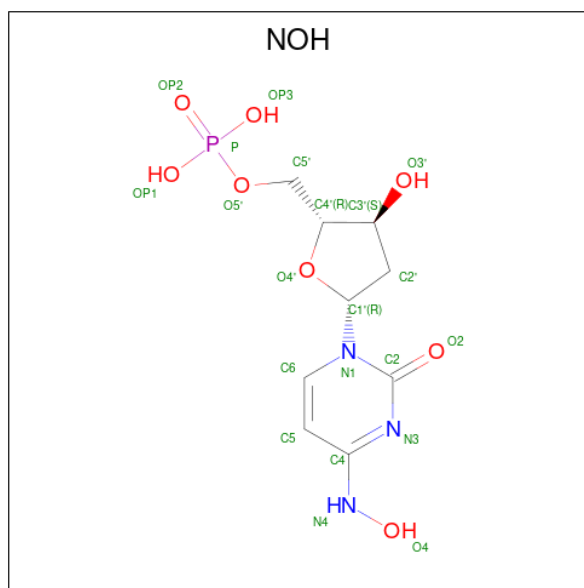
There are 3 unique types of molecules in this entry. The entry contains 4755 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Thymidylate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	289	Total	C	N	O	S	0	2	0
			2327	1475	404	432	16			
1	B	289	Total	C	N	O	S	0	1	0
			2324	1473	404	431	16			

- Molecule 2 is 2'-deoxy-N-hydroxycytidine 5'-(dihydrogen phosphate) (CCD ID: NOH) (formula: C₉H₁₄N₃O₈P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			21	9	3	8	1		
2	B	1	Total	C	N	O	P	0	0
			21	9	3	8	1		

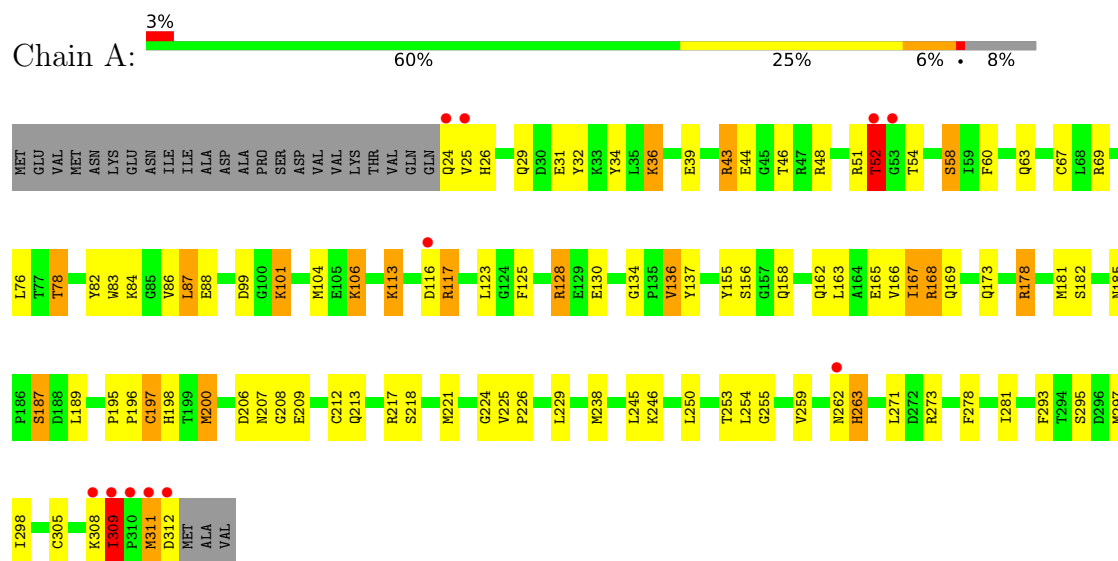
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	30	Total 30	O 30	0	0
3	B	32	Total 32	O 32	0	0

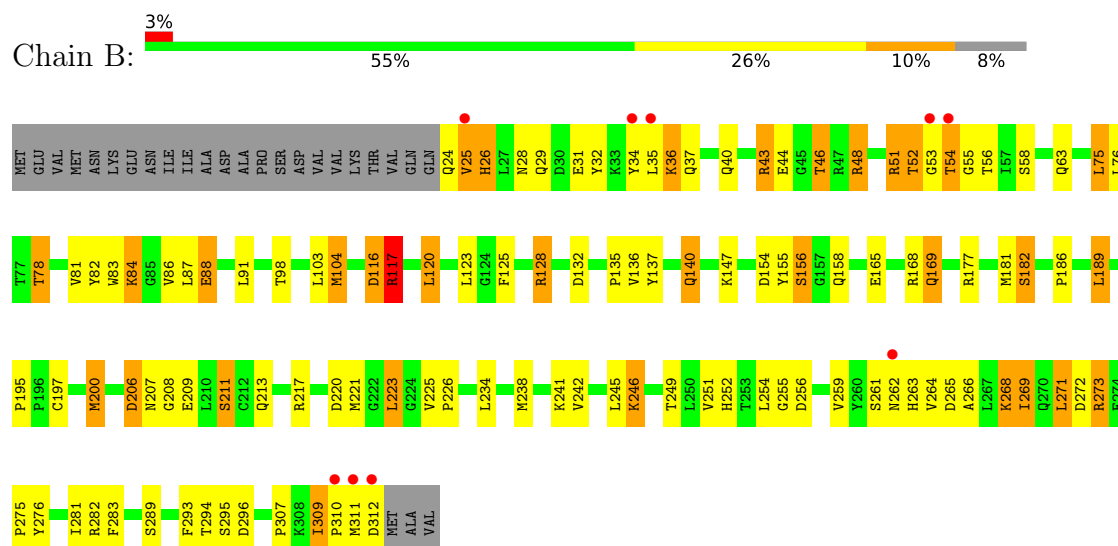
3 Residue-property plots [i](#)

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: Thymidylate synthase



• Molecule 1: Thymidylate synthase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	59.41Å 95.89Å 82.18Å 90.00° 103.81° 90.00°	Depositor
Resolution (Å)	25.11 – 2.80 25.11 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.9 (25.11-2.80) 99.8 (25.11-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.24	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.53 (at 2.80Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.231 , 0.267 0.225 , 0.259	Depositor DCC
R_{free} test set	1130 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å ²)	20.3	Xtriage
Anisotropy	0.184	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 31.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	4755	wwPDB-VP
Average B, all atoms (Å ²)	16.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.60% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

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

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	0.67	0/2387	1.12	11/3226 (0.3%)
1	B	0.74	1/2381 (0.0%)	1.08	7/3218 (0.2%)
All	All	0.71	1/4768 (0.0%)	1.10	18/6444 (0.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	220	ASP	C-O	-5.88	1.16	1.24

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	224	GLY	CA-C-N	-7.80	115.33	120.24
1	A	224	GLY	C-N-CA	-7.80	115.33	120.24
1	B	26	HIS	N-CA-C	-6.20	98.80	108.90
1	A	52	THR	N-CA-C	-6.01	105.93	113.20
1	A	25	VAL	N-CA-C	5.93	112.55	106.21
1	B	117	ARG	N-CA-C	5.80	119.61	112.54
1	B	137	TYR	N-CA-C	5.76	120.87	111.37
1	B	48	ARG	N-CA-C	-5.73	97.12	107.60
1	A	309	ILE	CA-C-N	5.47	124.66	118.97
1	A	309	ILE	C-N-CA	5.47	124.66	118.97
1	B	195	PRO	O-C-N	5.38	123.68	121.15
1	A	278	PHE	CA-C-N	5.28	126.03	120.11
1	A	278	PHE	C-N-CA	5.28	126.03	120.11
1	B	276	TYR	N-CA-C	-5.20	102.76	110.46
1	A	137	TYR	N-CA-C	5.13	119.84	111.37
1	A	197	CYS	N-CA-C	-5.12	106.97	113.43
1	B	32	TYR	N-CA-C	-5.04	105.91	111.71
1	A	58	SER	N-CA-C	5.04	117.66	109.24

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	2327	0	2288	70	0
1	B	2324	0	2283	104	0
2	A	21	0	11	0	0
2	B	21	0	12	0	0
3	A	30	0	0	0	0
3	B	32	0	0	0	0
All	All	4755	0	4594	167	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 18.

All (167) 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:B:26:HIS:HB2	1:B:29:GLN:NE2	1.61	1.13
1:B:51:ARG:HG2	1:B:51:ARG:HH11	0.91	1.05
1:B:51:ARG:HH11	1:B:51:ARG:CG	1.69	1.05
1:A:178[B]:ARG:HH11	1:A:178[B]:ARG:CG	1.74	1.01
1:A:82:TYR:O	1:A:86:VAL:HG23	1.63	0.99
1:A:178[B]:ARG:HH11	1:A:178[B]:ARG:HG2	0.80	0.97
1:B:182:SER:HB2	1:B:200:MET:HB3	1.47	0.94
1:B:206:ASP:O	1:B:207:ASN:HB2	1.65	0.94
1:A:178[B]:ARG:HG2	1:A:178[B]:ARG:NH1	1.60	0.93
1:B:120:LEU:N	1:B:120:LEU:HD23	1.79	0.93
1:B:51:ARG:HG2	1:B:51:ARG:NH1	1.67	0.92
1:B:211:SER:HB3	1:B:249:THR:O	1.73	0.88
1:B:36:LYS:N	1:B:36:LYS:HD2	1.88	0.88
1:B:26:HIS:HB2	1:B:29:GLN:HE22	1.38	0.86
1:B:58:SER:OG	1:B:256:ASP:OD1	1.92	0.85
1:B:26:HIS:HB2	1:B:29:GLN:HE21	1.41	0.85
1:B:117:ARG:O	1:B:117:ARG:HG3	1.75	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:34:TYR:CE2	1:B:76:LEU:HD11	2.13	0.83
1:B:54:THR:O	1:B:54:THR:HG23	1.79	0.81
1:B:165:GLU:O	1:B:169:GLN:HG3	1.81	0.81
1:A:206:ASP:OD2	1:B:46:THR:HG21	1.81	0.81
1:B:262:ASN:OD1	1:B:263:HIS:CE1	2.34	0.81
1:A:255:GLY:HA2	1:B:213:GLN:HE22	1.46	0.80
1:A:165:GLU:O	1:A:169:GLN:HG3	1.83	0.79
1:B:269:ILE:O	1:B:272:ASP:HB2	1.82	0.78
1:B:225:VAL:N	1:B:226:PRO:HD2	1.98	0.78
1:B:78:THR:O	1:B:273:ARG:NH1	2.17	0.78
1:A:34:TYR:OH	1:A:221:MET:O	2.02	0.76
1:B:54:THR:O	1:B:54:THR:CG2	2.33	0.76
1:A:84:LYS:NZ	1:A:106:LYS:O	2.18	0.75
1:A:262:ASN:C	1:A:263:HIS:HD2	1.94	0.75
1:A:262:ASN:O	1:A:263:HIS:HD2	1.70	0.74
1:B:261:SER:O	1:B:264:VAL:HG23	1.87	0.74
1:A:206:ASP:O	1:A:207:ASN:CB	2.36	0.74
1:B:34:TYR:CD2	1:B:76:LEU:HD11	2.24	0.72
1:B:34:TYR:OH	1:B:226:PRO:HD3	1.89	0.72
1:B:311:MET:O	1:B:312:ASP:HB2	1.89	0.71
1:A:311:MET:O	1:A:312:ASP:CB	2.40	0.70
1:A:311:MET:O	1:A:312:ASP:CG	2.35	0.69
1:B:206:ASP:OD1	1:B:206:ASP:C	2.37	0.68
1:B:225:VAL:H	1:B:226:PRO:HD2	1.58	0.67
1:A:262:ASN:C	1:A:263:HIS:CD2	2.72	0.67
1:B:165:GLU:CD	1:B:168:ARG:HH11	2.02	0.67
1:A:60:PHE:CD1	1:A:255:GLY:O	2.48	0.66
1:A:196:PRO:O	1:A:217:ARG:NE	2.27	0.66
1:A:155:TYR:O	1:A:158:GLN:HB2	1.95	0.66
1:A:168:ARG:NH1	1:A:173:GLN:OE1	2.29	0.65
1:B:123:LEU:HD13	1:B:125:PHE:CE1	2.31	0.65
1:B:116:ASP:C	1:B:116:ASP:OD2	2.40	0.64
1:B:225:VAL:N	1:B:226:PRO:CD	2.60	0.64
1:B:165:GLU:OE2	1:B:168:ARG:NH1	2.30	0.64
1:B:117:ARG:NH1	1:B:128:ARG:O	2.20	0.64
1:B:186:PRO:HA	1:B:189:LEU:HD12	1.79	0.64
1:A:69:ARG:HG3	1:A:69:ARG:HH21	1.64	0.62
1:A:206:ASP:OD2	1:B:46:THR:CG2	2.48	0.61
1:B:46:THR:HG22	1:B:46:THR:O	2.00	0.61
1:B:208:GLY:C	1:B:245:LEU:HD22	2.27	0.60
1:B:262:ASN:OD1	1:B:263:HIS:ND1	2.34	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:218:SER:OG	1:B:177:ARG:HD2	2.02	0.60
1:A:99:ASP:OD1	1:A:101:LYS:HB2	2.01	0.60
1:A:163:LEU:O	1:A:167:ILE:HG12	2.02	0.59
1:B:128:ARG:HG2	1:B:132:ASP:HB3	1.82	0.59
1:B:213:GLN:HG3	1:B:251:VAL:HB	1.83	0.59
1:B:217:ARG:HG3	1:B:217:ARG:HH11	1.68	0.58
1:A:206:ASP:O	1:A:207:ASN:HB2	2.02	0.58
1:A:225:VAL:N	1:A:226:PRO:HD2	2.17	0.58
1:A:63:GLN:HG3	1:A:253:THR:OG1	2.02	0.58
1:A:162:GLN:NE2	1:A:182:SER:O	2.32	0.58
1:A:123:LEU:HD13	1:A:125:PHE:CE1	2.38	0.57
1:A:101:LYS:NZ	1:A:130:GLU:O	2.35	0.57
1:A:263:HIS:CD2	1:A:263:HIS:N	2.72	0.57
1:B:225:VAL:HG13	1:B:252:HIS:CE1	2.40	0.56
1:B:182:SER:HB2	1:B:200:MET:CB	2.28	0.56
1:B:238:MET:O	1:B:242:VAL:HG23	2.06	0.56
1:A:238:MET:HG2	1:A:293:PHE:CE1	2.42	0.55
1:B:78:THR:HG22	1:B:273:ARG:HB3	1.87	0.55
1:A:113:LYS:H	1:A:113:LYS:HE2	1.71	0.55
1:B:209:GLU:HA	1:B:246:LYS:O	2.07	0.55
1:A:52:THR:HB	1:A:54:THR:HG23	1.89	0.54
1:B:311:MET:O	1:B:312:ASP:CB	2.54	0.54
1:A:69:ARG:HG3	1:A:69:ARG:NH2	2.23	0.53
1:B:43:ARG:CG	1:B:43:ARG:O	2.55	0.53
1:A:117:ARG:NH1	1:A:128:ARG:O	2.33	0.52
1:B:165:GLU:HA	1:B:165:GLU:OE1	2.08	0.52
1:B:265:ASP:O	1:B:269:ILE:HD12	2.10	0.52
1:A:255:GLY:HA2	1:B:213:GLN:NE2	2.20	0.52
1:B:225:VAL:HG13	1:B:252:HIS:HE1	1.73	0.52
1:A:26:HIS:HB2	1:A:29:GLN:NE2	2.25	0.51
1:A:209:GLU:HA	1:A:246:LYS:O	2.10	0.51
1:B:238:MET:HE3	1:B:293:PHE:CG	2.45	0.51
1:B:211:SER:CB	1:B:249:THR:O	2.53	0.51
1:A:123:LEU:HD13	1:A:125:PHE:HE1	1.74	0.51
1:B:26:HIS:CB	1:B:29:GLN:NE2	2.54	0.51
1:B:34:TYR:OH	1:B:221:MET:O	2.19	0.50
1:B:206:ASP:OD1	1:B:207:ASN:HB2	2.12	0.50
1:A:238:MET:HE3	1:A:293:PHE:CG	2.47	0.50
1:A:213:GLN:HE22	1:B:255:GLY:HA2	1.77	0.49
1:A:31:GLU:HG3	1:A:76:LEU:HD22	1.93	0.49
1:B:75:LEU:HD23	1:B:81:VAL:HB	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:84:LYS:O	1:B:88:GLU:HB2	2.12	0.49
1:B:123:LEU:HD13	1:B:125:PHE:HE1	1.74	0.49
1:A:206:ASP:CG	1:B:46:THR:HG21	2.38	0.49
1:B:51:ARG:CG	1:B:51:ARG:NH1	2.41	0.49
1:B:182:SER:CB	1:B:200:MET:HB3	2.32	0.49
1:B:273:ARG:NH2	1:B:307:PRO:O	2.39	0.49
1:B:309:ILE:HA	1:B:310:PRO:HD3	1.65	0.49
1:B:53:GLY:O	1:B:262:ASN:ND2	2.46	0.49
1:B:217:ARG:HG3	1:B:217:ARG:NH1	2.28	0.48
1:B:266:ALA:HB1	1:B:309:ILE:HG21	1.95	0.48
1:A:178[B]:ARG:CG	1:A:178[B]:ARG:NH1	2.45	0.48
1:B:154:ASP:OD1	1:B:156:SER:OG	2.32	0.48
1:B:31:GLU:HG3	1:B:76:LEU:HD22	1.96	0.47
1:B:309:ILE:HG22	1:B:311:MET:H	1.78	0.47
1:B:52:THR:O	1:B:52:THR:OG1	2.30	0.47
1:B:241:LYS:HD2	1:B:283:PHE:CD2	2.50	0.47
1:A:185:ASN:OD1	1:A:187:SER:HB2	2.14	0.47
1:B:34:TYR:CE2	1:B:76:LEU:CD1	2.94	0.47
1:A:262:ASN:O	1:A:263:HIS:CD2	2.59	0.47
1:B:181:MET:HE2	1:B:181:MET:HB3	1.83	0.47
1:B:28:ASN:HB3	1:B:275:PRO:HB2	1.97	0.46
1:B:55:GLY:O	1:B:261:SER:N	2.39	0.46
1:A:311:MET:O	1:A:312:ASP:HB2	2.15	0.45
1:B:40:GLN:O	1:B:44:GLU:HB2	2.17	0.45
1:B:268:LYS:HE2	1:B:268:LYS:HB3	1.66	0.45
1:B:269:ILE:O	1:B:272:ASP:N	2.44	0.45
1:A:43:ARG:NH1	1:A:44:GLU:OE1	2.49	0.45
1:B:206:ASP:O	1:B:206:ASP:OD1	2.33	0.45
1:B:309:ILE:HG23	1:B:310:PRO:HD2	1.99	0.45
1:A:297:MET:HE2	1:A:297:MET:HB2	1.88	0.45
1:A:212:CYS:HB3	1:A:250:LEU:HD12	1.99	0.45
1:B:26:HIS:CB	1:B:29:GLN:HE22	2.18	0.45
1:B:206:ASP:O	1:B:207:ASN:CB	2.44	0.45
1:B:294:THR:OG1	1:B:296:ASP:OD2	2.33	0.45
1:A:311:MET:O	1:A:312:ASP:OD1	2.35	0.44
1:A:200:MET:HB2	1:A:200:MET:HE2	1.60	0.44
1:B:83:TRP:CH2	1:B:234:LEU:HD22	2.52	0.44
1:B:165:GLU:O	1:B:169:GLN:CG	2.60	0.44
1:A:36:LYS:HD3	1:A:36:LYS:HA	1.71	0.44
1:B:34:TYR:CD2	1:B:76:LEU:CD1	2.98	0.43
1:A:87:LEU:HD21	1:A:298:ILE:HD12	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:82:TYR:O	1:B:86:VAL:HG23	2.19	0.43
1:A:83:TRP:O	1:A:87:LEU:HB2	2.19	0.43
1:B:25:VAL:O	1:B:25:VAL:HG12	2.18	0.43
1:B:104:MET:HE2	1:B:104:MET:HB2	1.88	0.43
1:A:208:GLY:C	1:A:245:LEU:HD22	2.44	0.43
1:B:36:LYS:N	1:B:36:LYS:CD	2.70	0.43
1:A:125:PHE:O	1:A:128:ARG:HB2	2.19	0.43
1:B:208:GLY:O	1:B:245:LEU:HD22	2.19	0.42
1:A:197:CYS:HA	1:A:217:ARG:HG2	2.02	0.42
1:B:46:THR:CG2	1:B:46:THR:O	2.60	0.42
1:B:88:GLU:OE1	1:B:88:GLU:HA	2.19	0.42
1:A:34:TYR:OH	1:A:226:PRO:HD3	2.20	0.42
1:A:134:GLY:O	1:A:136:VAL:N	2.48	0.42
1:B:135:PRO:O	1:B:140:GLN:NE2	2.53	0.42
1:B:223:LEU:HA	1:B:223:LEU:HD23	1.72	0.42
1:A:67:CYS:SG	1:A:69:ARG:NH2	2.93	0.41
1:A:31:GLU:O	1:A:32:TYR:C	2.63	0.41
1:B:155:TYR:O	1:B:158:GLN:HB2	2.20	0.41
1:A:185:ASN:O	1:A:189:LEU:HG	2.21	0.41
1:A:78:THR:O	1:A:273:ARG:NH1	2.52	0.41
1:B:35:LEU:O	1:B:271:LEU:HD11	2.21	0.41
1:A:178[A]:ARG:HE	1:A:178[A]:ARG:HB3	1.47	0.40
1:A:195:PRO:HA	1:A:196:PRO:HD3	1.97	0.40
1:A:309:ILE:O	1:A:309:ILE:HD13	2.21	0.40
1:B:51:ARG:HH11	1:B:51:ARG:CB	2.28	0.40

There are no symmetry-related clashes.

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	289/315 (92%)	280 (97%)	8 (3%)	1 (0%)	36 66

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	288/315 (91%)	276 (96%)	11 (4%)	1 (0%)	36	66
All	All	577/630 (92%)	556 (96%)	19 (3%)	2 (0%)	36	66

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	136	VAL
1	A	136	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	255/276 (92%)	215 (84%)	40 (16%)	2	9
1	B	254/276 (92%)	206 (81%)	48 (19%)	1	5
All	All	509/552 (92%)	421 (83%)	88 (17%)	2	7

All (88) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	GLN
1	A	36	LYS
1	A	39	GLU
1	A	43	ARG
1	A	46	THR
1	A	48	ARG
1	A	51	ARG
1	A	52	THR
1	A	58	SER
1	A	78	THR
1	A	87	LEU
1	A	88	GLU
1	A	101	LYS
1	A	104	MET

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Mol	Chain	Res	Type
1	A	106	LYS
1	A	113	LYS
1	A	116	ASP
1	A	117	ARG
1	A	128	ARG
1	A	156	SER
1	A	166	VAL
1	A	167	ILE
1	A	168	ARG
1	A	178[A]	ARG
1	A	178[B]	ARG
1	A	181	MET
1	A	187	SER
1	A	198	HIS
1	A	200	MET
1	A	229	LEU
1	A	254	LEU
1	A	259	VAL
1	A	263	HIS
1	A	271	LEU
1	A	281	ILE
1	A	295	SER
1	A	305	CYS
1	A	308	LYS
1	A	309	ILE
1	A	311	MET
1	B	24	GLN
1	B	25	VAL
1	B	36	LYS
1	B	37	GLN
1	B	43	ARG
1	B	46	THR
1	B	48	ARG
1	B	51	ARG
1	B	52	THR
1	B	54	THR
1	B	56	THR
1	B	63	GLN
1	B	75	LEU
1	B	78	THR
1	B	84	LYS
1	B	87	LEU

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Mol	Chain	Res	Type
1	B	88	GLU
1	B	91	LEU
1	B	98	THR
1	B	103	LEU
1	B	104	MET
1	B	116	ASP
1	B	117	ARG
1	B	120	LEU
1	B	128	ARG
1	B	140	GLN
1	B	147	LYS
1	B	156	SER
1	B	169	GLN
1	B	182	SER
1	B	189	LEU
1	B	197	CYS
1	B	200	MET
1	B	206	ASP
1	B	211	SER
1	B	223	LEU
1	B	246	LYS
1	B	254	LEU
1	B	259	VAL
1	B	268	LYS
1	B	269	ILE
1	B	271	LEU
1	B	273	ARG
1	B	281	ILE
1	B	282	ARG
1	B	289	SER
1	B	295	SER
1	B	309	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	24	GLN
1	A	26	HIS
1	A	29	GLN
1	A	70	ASN
1	A	114	ASN
1	A	191	GLN

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Mol	Chain	Res	Type
1	A	213	GLN
1	A	263	HIS
1	B	29	GLN
1	B	70	ASN
1	B	169	GLN
1	B	191	GLN
1	B	207	ASN
1	B	213	GLN
1	B	216	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](#)

2 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	$\# Z > 2$
2	NOH	A	401	-	22,22,22	2.27	4 (18%)	28,32,32	1.65	5 (17%)
2	NOH	B	401	-	22,22,22	1.37	3 (13%)	28,32,32	1.17	1 (3%)

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	NOH	A	401	-	-	4/10/24/24	0/2/2/2
2	NOH	B	401	-	-	2/10/24/24	0/2/2/2

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	NOH	O4-N4	-8.45	1.19	1.40
2	B	401	NOH	C2-N1	-3.01	1.33	1.40
2	A	401	NOH	C4-N4	-2.64	1.29	1.35
2	A	401	NOH	C2-N1	-2.54	1.34	1.40
2	B	401	NOH	O4-N4	-2.20	1.34	1.40
2	B	401	NOH	P-OP3	2.08	1.62	1.54
2	A	401	NOH	P-OP3	2.04	1.62	1.54

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	NOH	O4'-C1'-N1	-4.33	100.18	107.86
2	A	401	NOH	O5'-P-OP2	3.77	116.64	106.44
2	A	401	NOH	O2-C2-N3	-3.10	117.45	122.33
2	B	401	NOH	O5'-P-OP2	2.82	114.06	106.44
2	A	401	NOH	C1'-N1-C2	2.53	122.17	117.83
2	A	401	NOH	C2'-C1'-N1	-2.08	108.61	113.81

There are no chirality outliers.

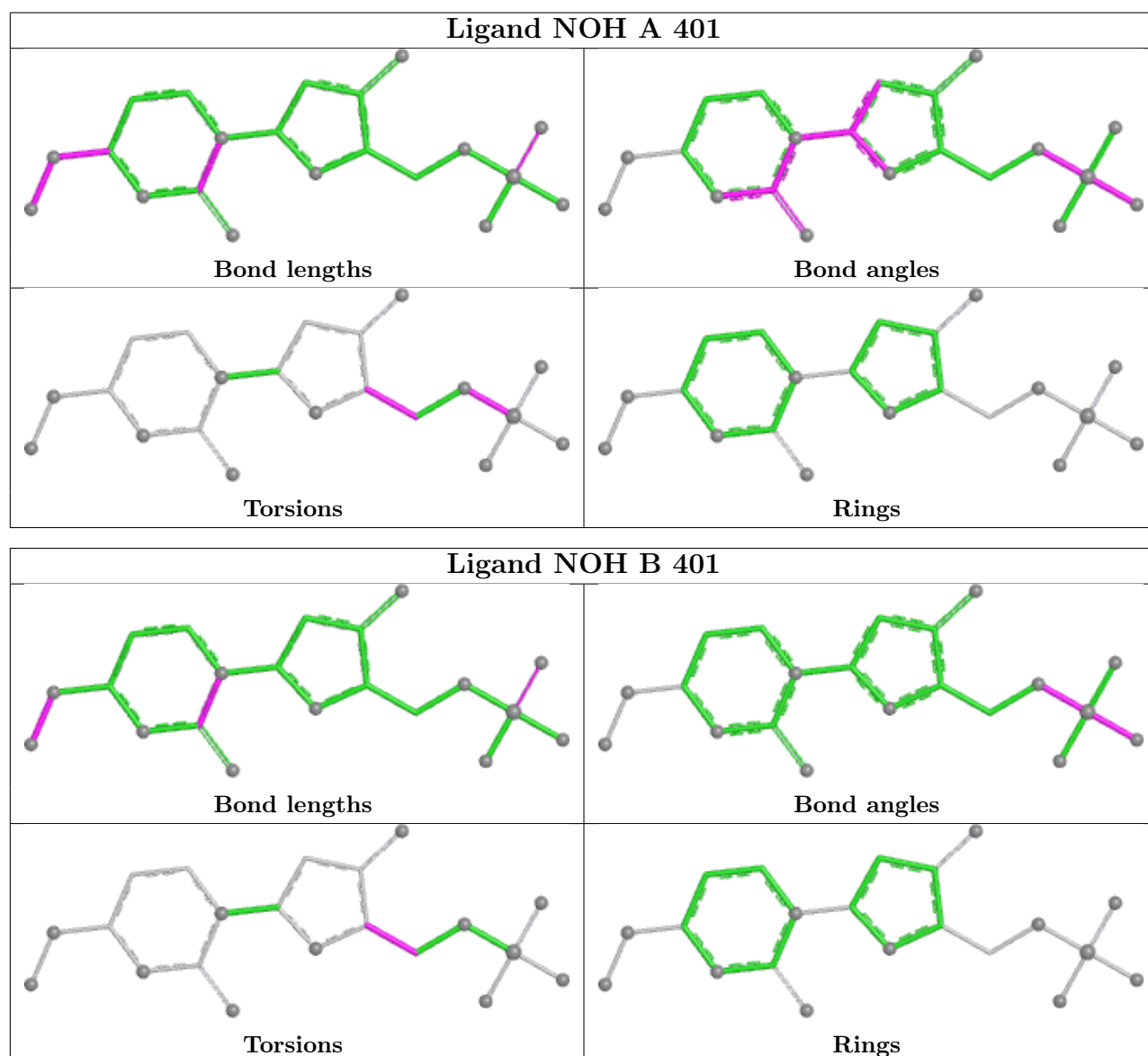
All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401	NOH	C5'-O5'-P-OP3
2	B	401	NOH	O4'-C4'-C5'-O5'
2	A	401	NOH	C3'-C4'-C5'-O5'
2	B	401	NOH	C3'-C4'-C5'-O5'
2	A	401	NOH	O4'-C4'-C5'-O5'
2	A	401	NOH	C5'-O5'-P-OP1

There are no ring outliers.

No monomer is involved in short contacts.

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



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	289/315 (91%)	0.08	11 (3%) 44 35	5, 14, 30, 66	2 (0%)
1	B	289/315 (91%)	0.07	9 (3%) 51 41	6, 14, 31, 70	1 (0%)
All	All	578/630 (91%)	0.08	20 (3%) 47 38	5, 14, 31, 70	3 (0%)

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	25	VAL	4.6
1	A	309	ILE	4.5
1	B	312	ASP	4.5
1	A	311	MET	4.0
1	A	310	PRO	3.9
1	A	24	GLN	3.8
1	B	311	MET	3.7
1	B	54	THR	3.3
1	A	262	ASN	3.2
1	A	116	ASP	3.1
1	B	262	ASN	2.9
1	B	53	GLY	2.7
1	A	312	ASP	2.6
1	B	35	LEU	2.5
1	B	310	PRO	2.3
1	A	308	LYS	2.3
1	B	25	VAL	2.3
1	B	34	TYR	2.2
1	A	53	GLY	2.1
1	A	52	THR	2.0

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

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

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

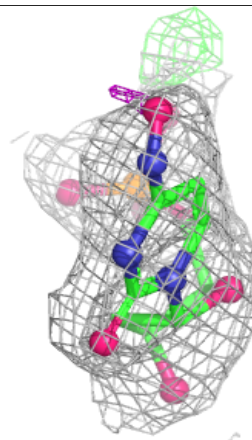
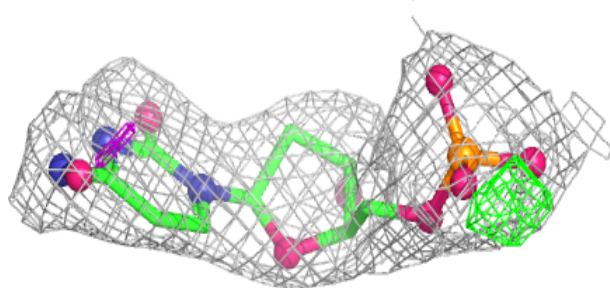
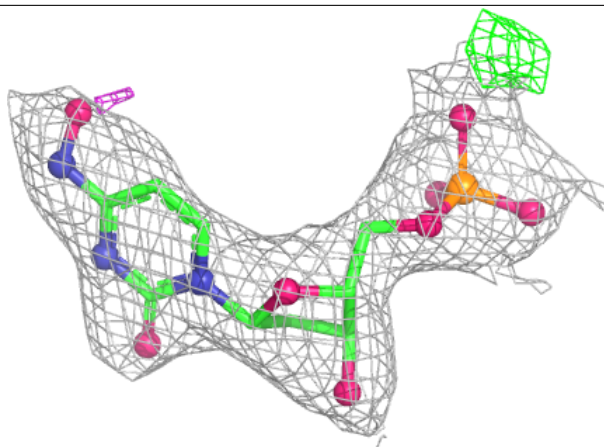
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

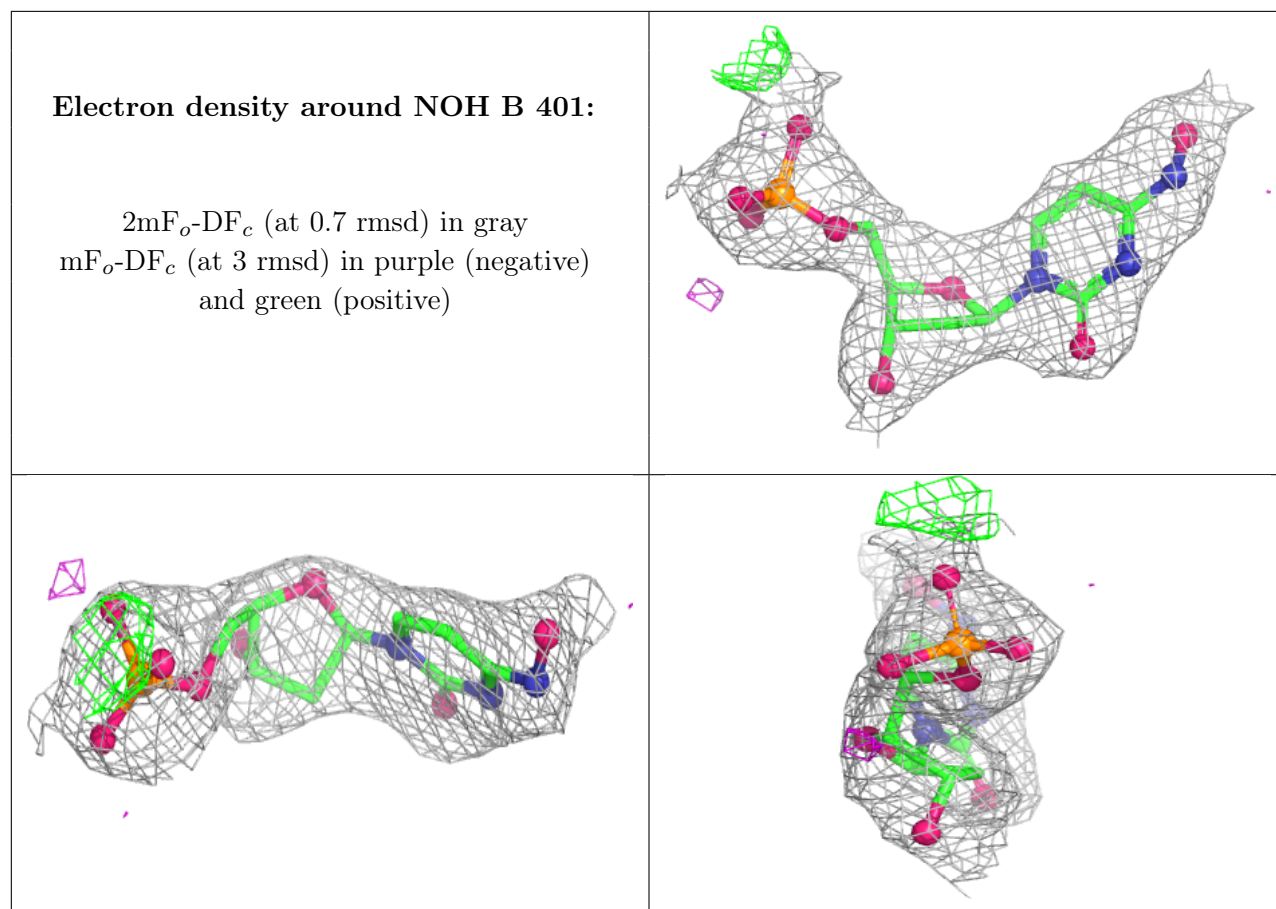
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	NOH	A	401	21/21	0.93	0.10	22,23,24,24	0
2	NOH	B	401	21/21	0.94	0.09	17,18,18,18	0

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 NOH A 401:

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





6.5 Other polymers ⓘ

There are no such residues in this entry.