



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:21 PM UTC

PDB ID : 4I3L / pdb_00004i3l
Title : Crystal structure of a metabolic reductase with 6-benzyl-1-hydroxy-4-methylpyridin-2(1H)-one
Authors : Zheng, B.; Yao, Y.; Liu, Z.; Deng, L.; Anglin, J.L.; Jiang, H.; Prasad, B.V.V.; Song, Y.
Deposited on : 2012-11-26
Resolution : 3.29 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

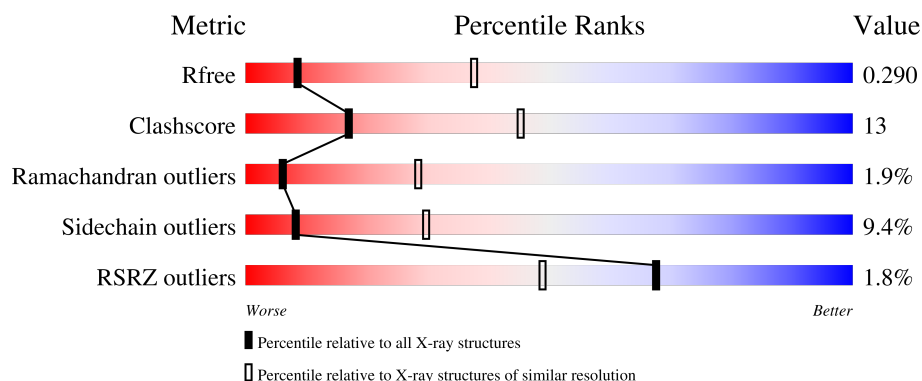
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.29 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	1169 (3.32-3.28)
Clashscore	190562	1209 (3.32-3.28)
Ramachandran outliers	187476	1188 (3.32-3.28)
Sidechain outliers	187428	1187 (3.32-3.28)
RSRZ outliers	180081	1169 (3.32-3.28)

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	414	60% 27% • 11%
1	B	414	2% 54% 28% 5% 14%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 5888 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.

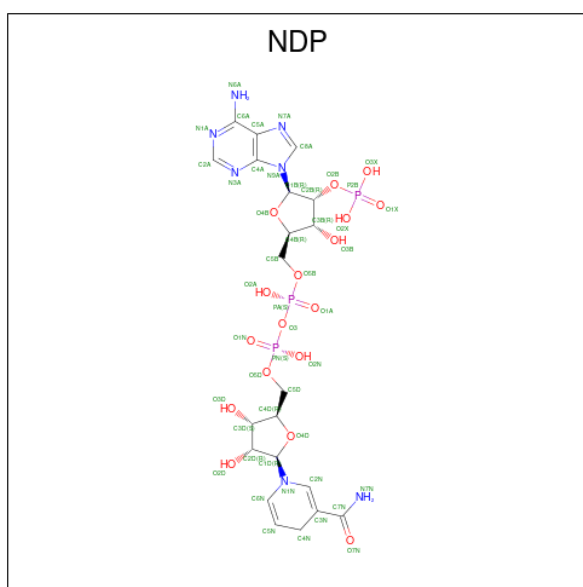
- Molecule 1 is a protein called Isocitrate dehydrogenase [NADP] cytoplasmic.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	367	Total	C	N	O	S	0	0	0
			2922	1863	493	549	17			
1	B	358	Total	C	N	O	S	0	0	0
			2844	1814	478	535	17			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	132	HIS	ARG	engineered mutation	UNP O75874
B	132	HIS	ARG	engineered mutation	UNP O75874

- Molecule 2 is NADPH DIHYDRO-NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (CCD ID: NDP) (formula: $C_{21}H_{30}N_7O_{17}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			48	21	7	17	3		

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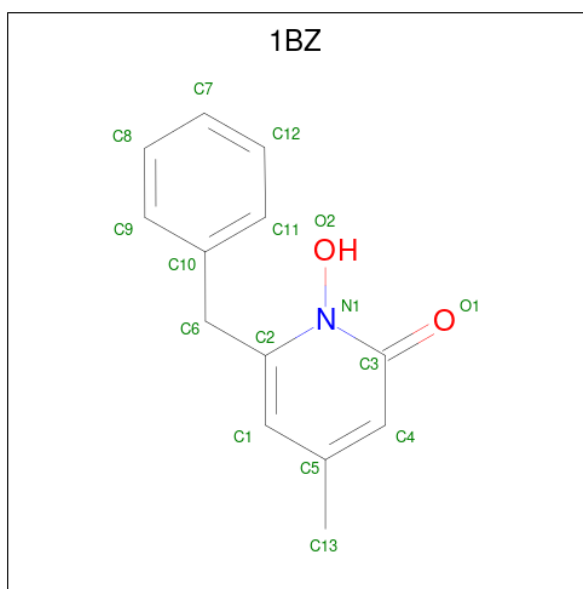
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	N	O	P	0	0
			48	21	7	17	3		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is 6-benzyl-1-hydroxy-4-methylpyridin-2(1H)-one (CCD ID: 1BZ) (formula: C₁₃H₁₃NO₂).

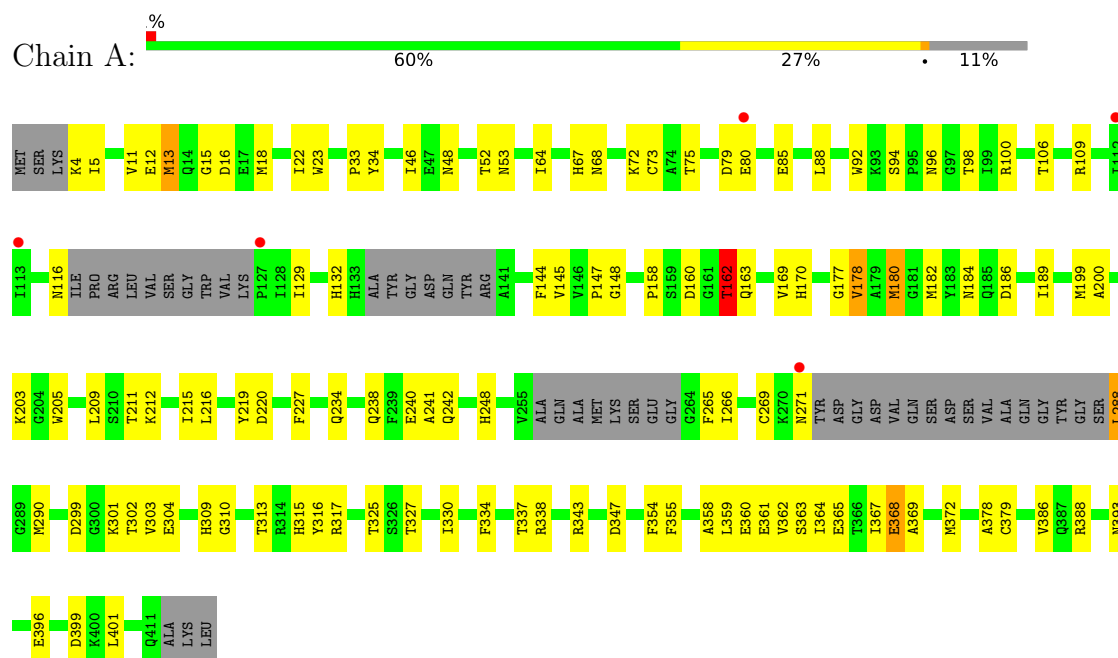


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
4	A	1	16	13	1	2	0	0

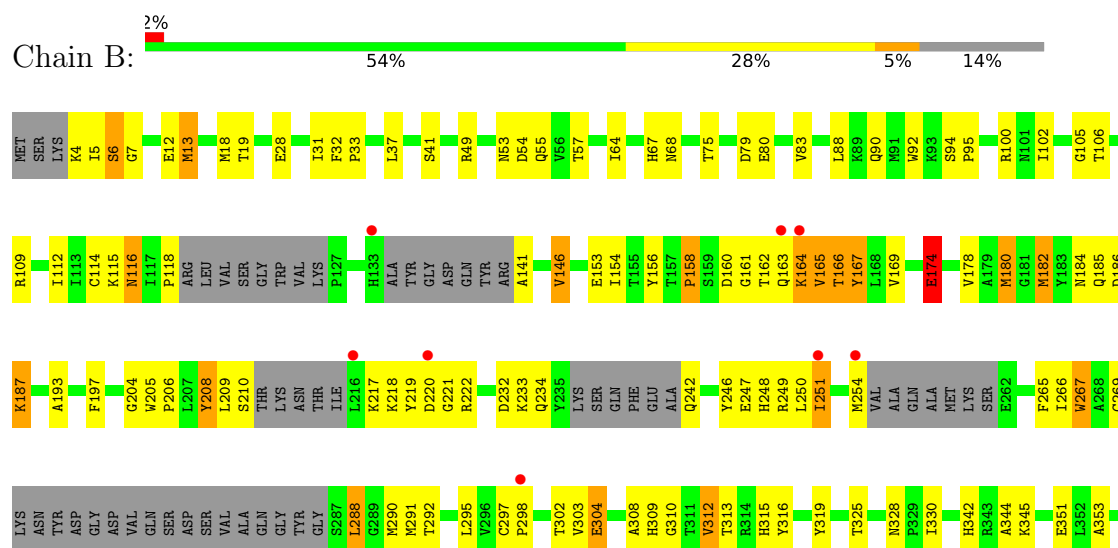
3 Residue-property plots

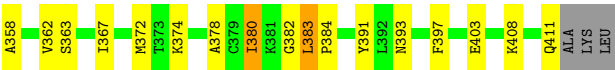
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Isocitrate dehydrogenase [NADP] cytoplasmic



- Molecule 1: Isocitrate dehydrogenase [NADP] cytoplasmic





4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	80.85Å 80.85Å 304.84Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.73 – 3.29 45.73 – 3.29	Depositor EDS
% Data completeness (in resolution range)	95.8 (45.73-3.29) 95.8 (45.73-3.29)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.43 (at 3.32Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.220 , 0.293 0.220 , 0.290	Depositor DCC
R_{free} test set	816 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	52.3	Xtriage
Anisotropy	0.271	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 32.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	5888	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.30% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, 1BZ, NDP

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	0.76	0/2980	1.00	6/4012 (0.1%)
1	B	0.78	0/2900	1.01	6/3903 (0.2%)
All	All	0.77	0/5880	1.01	12/7915 (0.2%)

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	372	MET	N-CA-C	6.97	118.78	108.60
1	B	163	GLN	N-CA-C	6.80	118.13	108.74
1	A	148	GLY	CA-C-N	6.46	126.83	119.92
1	A	148	GLY	C-N-CA	6.46	126.83	119.92
1	B	180	MET	N-CA-C	6.06	118.07	108.79
1	A	94	SER	CA-C-N	5.88	125.09	118.97
1	A	94	SER	C-N-CA	5.88	125.09	118.97
1	B	94	SER	CA-C-N	5.82	125.51	119.05
1	B	94	SER	C-N-CA	5.82	125.51	119.05
1	B	146	VAL	CB-CA-C	5.37	116.95	110.78
1	B	267	TRP	N-CA-C	5.30	116.08	107.23
1	A	16	ASP	N-CA-C	5.19	117.56	109.52

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen

atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2922	0	2899	74	0
1	B	2844	0	2811	76	0
2	A	48	0	26	5	0
2	B	48	0	26	1	0
3	A	10	0	0	0	0
4	A	16	0	13	1	0
All	All	5888	0	5775	148	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 13.

All (148) 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:13:MET:HE1	1:B:64:ILE:CD1	1.73	1.18
1:B:13:MET:CE	1:B:64:ILE:HD11	1.74	1.16
1:A:200:ALA:HA	1:A:266:ILE:CD1	2.10	0.81
1:B:206:PRO:HB2	1:B:208:TYR:HE1	1.45	0.81
1:B:83:VAL:HA	1:B:88:LEU:HD12	1.63	0.79
1:A:361:GLU:O	1:A:365:GLU:HG2	1.87	0.73
1:A:325:THR:O	1:A:393:ASN:HB2	1.89	0.72
1:B:18:MET:HB2	1:B:316:TYR:HB2	1.72	0.71
1:B:7:GLY:HA3	1:B:37:LEU:HD23	1.74	0.69
1:A:288:LEU:C	1:A:290:MET:H	2.01	0.68
1:B:291:MET:HB3	1:B:308:ALA:HB3	1.76	0.67
1:A:75:THR:H	2:A:501:NDP:H71N	1.40	0.67
1:B:303:VAL:HB	1:B:342:HIS:CD2	2.31	0.65
1:A:367:ILE:C	1:A:369:ALA:H	2.05	0.64
1:B:167:TYR:CD1	1:B:167:TYR:N	2.65	0.64
1:A:13:MET:HE1	1:A:64:ILE:HD11	1.81	0.63
1:A:378:ALA:HA	1:A:386:VAL:HG21	1.81	0.62
1:B:13:MET:HE1	1:B:64:ILE:HD11	0.81	0.62
1:B:297:CYS:HB3	1:B:298:PRO:HD2	1.81	0.62
1:B:164:LYS:O	1:B:166:THR:HG22	2.00	0.62
1:A:330:ILE:HD12	1:A:363:SER:HB3	1.81	0.61
1:B:164:LYS:O	1:B:165:VAL:C	2.43	0.61
1:A:200:ALA:HA	1:A:266:ILE:HD12	1.82	0.61
1:B:313:THR:OG1	2:B:501:NDP:H51N	2.01	0.60
1:A:334:PHE:O	1:A:338:ARG:HG2	2.01	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:167:TYR:H	1:B:167:TYR:HD1	1.48	0.60
1:A:288:LEU:HD22	1:A:309:HIS:HB3	1.83	0.60
1:A:53:ASN:HA	1:A:92:TRP:CH2	2.38	0.58
1:B:219:TYR:C	1:B:221:GLY:N	2.60	0.58
1:B:325:THR:O	1:B:393:ASN:HB2	2.04	0.58
1:A:96:ASN:O	1:A:100:ARG:HB2	2.04	0.58
1:B:164:LYS:O	1:B:164:LYS:CG	2.51	0.58
1:A:5:ILE:HD12	1:A:34:TYR:O	2.03	0.58
1:A:162:THR:HG22	1:A:163:GLN:H	1.70	0.57
1:A:158:PRO:HG2	1:A:162:THR:HB	1.86	0.56
1:A:177:GLY:O	1:A:178:VAL:HG13	2.05	0.56
1:B:403:GLU:OE1	1:B:403:GLU:HA	2.06	0.56
1:A:23:TRP:CD2	1:A:73:CYS:HB2	2.42	0.55
1:A:378:ALA:HA	1:A:386:VAL:CG2	2.36	0.55
1:B:164:LYS:O	1:B:164:LYS:HG3	2.06	0.55
1:B:112:ILE:HD11	1:B:292:THR:OG1	2.06	0.55
1:A:367:ILE:C	1:A:369:ALA:N	2.65	0.55
1:B:105:GLY:HA3	1:B:297:CYS:HA	1.89	0.54
1:B:351:GLU:H	1:B:351:GLU:CD	2.16	0.54
1:A:337:THR:HG23	1:A:359:LEU:HD23	1.90	0.54
1:A:358:ALA:O	1:A:362:VAL:HG23	2.07	0.54
1:A:75:THR:O	2:A:501:NDP:H2N	2.08	0.54
1:A:211:THR:OG1	1:A:248:HIS:HE1	1.90	0.54
1:B:309:HIS:HD2	1:B:310:GLY:O	1.91	0.54
1:A:367:ILE:O	1:A:369:ALA:N	2.41	0.53
1:A:378:ALA:CA	1:A:386:VAL:HG21	2.37	0.53
1:A:240:GLU:C	1:A:242:GLN:H	2.16	0.53
1:B:5:ILE:HG22	1:B:6:SER:N	2.24	0.53
1:A:205:TRP:HB3	1:A:265:PHE:HA	1.91	0.53
1:A:313:THR:OG1	2:A:501:NDP:H51N	2.09	0.53
1:A:5:ILE:HD12	1:A:5:ILE:H	1.74	0.53
1:A:211:THR:HB	1:A:220:ASP:HB3	1.90	0.53
1:B:219:TYR:C	1:B:221:GLY:H	2.18	0.52
1:B:67:HIS:O	1:B:68:ASN:HB2	2.10	0.52
1:B:248:HIS:O	1:B:249:ARG:HD2	2.10	0.52
2:A:501:NDP:H42N	4:A:504:1BZ:H5	1.92	0.52
1:A:169:VAL:O	1:A:170:HIS:HB2	2.11	0.51
1:B:304:GLU:O	1:B:304:GLU:HG3	2.10	0.51
1:B:141:ALA:HB2	1:B:182:MET:HG3	1.91	0.51
1:B:154:ILE:HG22	1:B:169:VAL:HG22	1.91	0.51
1:B:372:MET:HA	1:B:397:PHE:CE1	2.47	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:48:ASN:O	1:A:52:THR:HG23	2.12	0.50
1:A:288:LEU:C	1:A:290:MET:N	2.65	0.49
1:B:247:GLU:OE2	1:B:249:ARG:NH1	2.44	0.49
1:B:267:TRP:CD1	1:B:267:TRP:C	2.90	0.49
1:A:309:HIS:HD2	1:A:310:GLY:O	1.95	0.49
1:A:337:THR:CG2	1:A:359:LEU:HD23	2.43	0.49
1:A:18:MET:HB2	1:A:316:TYR:HB2	1.96	0.48
1:B:358:ALA:O	1:B:362:VAL:HG23	2.13	0.48
1:A:72:LYS:NZ	2:A:501:NDP:H72N	2.12	0.48
1:B:383:LEU:N	1:B:384:PRO:HD2	2.29	0.48
1:B:68:ASN:HA	1:B:302:THR:HG23	1.96	0.48
1:A:129:ILE:HG21	1:A:199:MET:HE3	1.95	0.47
1:A:158:PRO:HB2	1:A:160:ASP:OD1	2.14	0.47
1:B:100:ARG:HG2	1:B:295:LEU:HD22	1.96	0.47
1:A:15:GLY:HA2	1:A:75:THR:HG22	1.96	0.47
1:A:13:MET:HG3	1:A:72:LYS:HA	1.97	0.47
1:A:299:ASP:C	1:A:301:LYS:H	2.23	0.47
1:B:116:ASN:O	1:B:380:ILE:CD1	2.63	0.46
1:A:354:PHE:CD2	1:A:354:PHE:C	2.92	0.46
1:B:49:ARG:NH1	1:B:95:PRO:HD3	2.30	0.46
1:A:240:GLU:C	1:A:242:GLN:N	2.74	0.46
1:B:112:ILE:HD12	1:B:290:MET:O	2.16	0.46
1:B:344:ALA:HB1	1:B:353:ALA:HB2	1.97	0.46
1:B:374:LYS:HD2	1:B:391:TYR:CZ	2.50	0.46
1:A:186:ASP:HA	1:A:189:ILE:HD12	1.98	0.46
1:A:343:ARG:NE	1:A:347:ASP:OD2	2.49	0.46
1:B:53:ASN:HA	1:B:92:TRP:CH2	2.51	0.46
1:B:174:GLU:H	1:B:174:GLU:CD	2.23	0.46
1:A:363:SER:HA	1:A:401:LEU:HD21	1.98	0.45
1:A:396:GLU:O	1:A:399:ASP:HB2	2.16	0.45
1:B:79:ASP:O	1:B:83:VAL:HG23	2.17	0.45
1:B:53:ASN:O	1:B:55:GLN:N	2.49	0.45
1:B:233:LYS:HG3	1:B:234:GLN:HG3	1.98	0.45
1:A:334:PHE:CE2	1:A:360:GLU:HG2	2.52	0.44
1:B:109:ARG:HD3	1:B:291:MET:HE1	1.98	0.44
1:B:205:TRP:HB3	1:B:265:PHE:HA	1.99	0.44
1:A:67:HIS:O	1:A:68:ASN:HB2	2.18	0.44
1:A:212:LYS:HD3	1:A:215:ILE:HD12	1.99	0.44
1:A:355:PHE:O	1:A:358:ALA:HB3	2.18	0.44
1:B:204:GLY:O	1:B:205:TRP:HD1	2.01	0.44
1:B:160:ASP:O	1:B:162:THR:HG23	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:4:LYS:HE2	1:A:33:PRO:O	2.19	0.43
1:B:193:ALA:O	1:B:197:PHE:CD1	2.72	0.43
1:B:251:ILE:H	1:B:251:ILE:HG13	1.55	0.43
1:A:209:LEU:HD22	1:A:227:PHE:CG	2.53	0.43
1:B:209:LEU:HD12	1:B:210:SER:N	2.34	0.43
1:B:328:ASN:OD1	1:B:330:ILE:HG12	2.19	0.43
1:A:216:LEU:HB3	1:A:219:TYR:HB3	2.00	0.43
1:B:206:PRO:HB2	1:B:208:TYR:CE1	2.37	0.42
1:B:378:ALA:O	1:B:382:GLY:N	2.49	0.42
1:B:205:TRP:HA	1:B:206:PRO:HD2	1.83	0.42
1:B:246:TYR:CZ	1:B:247:GLU:O	2.73	0.42
1:A:116:ASN:OD1	1:A:368:GLU:HA	2.18	0.42
1:B:316:TYR:O	1:B:319:TYR:HB3	2.19	0.42
1:B:4:LYS:HE2	1:B:32:PHE:O	2.19	0.42
1:A:211:THR:OG1	1:A:248:HIS:CE1	2.71	0.42
1:B:184:ASN:HB2	1:B:219:TYR:OH	2.19	0.42
1:A:46:ILE:HG13	1:A:88:LEU:HD21	2.01	0.42
1:A:12:GLU:HG2	1:A:13:MET:N	2.34	0.42
1:A:315:HIS:O	1:A:325:THR:HG21	2.20	0.41
1:A:145:VAL:HG23	1:B:218:LYS:HB3	2.01	0.41
1:A:79:ASP:O	1:A:80:GLU:C	2.63	0.41
1:A:13:MET:HE1	1:A:64:ILE:CD1	2.48	0.41
1:A:11:VAL:HG21	1:A:64:ILE:HA	2.01	0.41
1:A:182:MET:SD	1:B:180:MET:HE3	2.60	0.41
1:A:85:GLU:OE2	1:A:317:ARG:NH2	2.53	0.41
1:A:234:GLN:O	1:A:238:GLN:NE2	2.54	0.41
1:B:154:ILE:HG22	1:B:169:VAL:CG2	2.50	0.41
1:B:288:LEU:C	1:B:290:MET:H	2.28	0.41
1:B:254:MET:H	1:B:254:MET:HG2	1.67	0.41
1:A:144:PHE:CD2	1:A:144:PHE:C	2.99	0.41
1:A:22:ILE:HD11	1:A:327:THR:HB	2.03	0.41
1:A:68:ASN:HA	1:A:302:THR:HG23	2.01	0.41
1:B:312:VAL:HG12	1:B:315:HIS:HB2	2.02	0.41
1:A:180:MET:HG2	1:B:219:TYR:CE1	2.56	0.40
1:B:116:ASN:HD21	1:B:367:ILE:HG22	1.86	0.40
1:B:156:TYR:CE1	1:B:158:PRO:HD3	2.57	0.40
1:B:186:ASP:O	1:B:187:LYS:C	2.63	0.40
1:B:193:ALA:O	1:B:197:PHE:HD1	2.04	0.40
1:B:31:ILE:C	1:B:33:PRO:HD2	2.47	0.40
1:A:129:ILE:HD11	1:A:203:LYS:HE3	2.03	0.40
1:B:209:LEU:HD12	1:B:210:SER:H	1.87	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	357/414 (86%)	323 (90%)	29 (8%)	5 (1%)	9	33
1	B	344/414 (83%)	298 (87%)	38 (11%)	8 (2%)	5	25
All	All	701/828 (85%)	621 (89%)	67 (10%)	13 (2%)	6	28

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	54	ASP
1	B	220	ASP
1	A	132	HIS
1	B	161	GLY
1	B	217	LYS
1	A	241	ALA
1	B	158	PRO
1	B	174	GLU
1	B	380	ILE
1	A	368	GLU
1	B	251	ILE
1	A	147	PRO
1	A	162	THR

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	314/350 (90%)	298 (95%)	16 (5%)	21	50
1	B	305/350 (87%)	263 (86%)	42 (14%)	3	15
All	All	619/700 (88%)	561 (91%)	58 (9%)	8	30

All (58) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	MET
1	A	98	THR
1	A	106	THR
1	A	109	ARG
1	A	162	THR
1	A	178	VAL
1	A	180	MET
1	A	184	ASN
1	A	269	CYS
1	A	271	ASN
1	A	288	LEU
1	A	303	VAL
1	A	304	GLU
1	A	364	ILE
1	A	379	CYS
1	A	388	ARG
1	B	6	SER
1	B	12	GLU
1	B	13	MET
1	B	19	THR
1	B	28	GLU
1	B	41	SER
1	B	57	THR
1	B	75	THR
1	B	80	GLU
1	B	90	GLN
1	B	102	ILE
1	B	106	THR
1	B	114	CYS
1	B	115	LYS
1	B	116	ASN
1	B	118	PRO
1	B	146	VAL
1	B	153	GLU
1	B	164	LYS

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Mol	Chain	Res	Type
1	B	165	VAL
1	B	166	THR
1	B	167	TYR
1	B	174	GLU
1	B	178	VAL
1	B	182	MET
1	B	185	GLN
1	B	187	LYS
1	B	208	TYR
1	B	222	ARG
1	B	232	ASP
1	B	242	GLN
1	B	250	LEU
1	B	266	ILE
1	B	269	CYS
1	B	288	LEU
1	B	304	GLU
1	B	312	VAL
1	B	345	LYS
1	B	363	SER
1	B	383	LEU
1	B	408	LYS
1	B	411	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	14	GLN
1	A	101	ASN
1	A	234	GLN
1	A	238	GLN
1	A	242	GLN
1	A	248	HIS
1	A	309	HIS
1	A	348	ASN
1	B	116	ASN
1	B	132	HIS
1	B	171	ASN
1	B	309	HIS
1	B	348	ASN
1	B	387	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 ⓘ

5 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	NDP	A	501	-	51,52,52	1.39	7 (13%)	71,80,80	1.60	11 (15%)
2	NDP	B	501	-	51,52,52	1.24	5 (9%)	71,80,80	1.53	10 (14%)
3	SO4	A	502	-	4,4,4	0.45	0	6,6,6	0.13	0
3	SO4	A	503	-	4,4,4	0.48	0	6,6,6	0.09	0
4	1BZ	A	504	-	16,17,17	1.76	1 (6%)	21,23,23	4.68	8 (38%)

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
4	1BZ	A	504	-	-	2/4/4/4	0/2/2/2
2	NDP	A	501	-	-	9/34/77/77	0/5/5/5
2	NDP	B	501	-	-	15/34/77/77	0/5/5/5

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	504	1BZ	O2-N1	-6.12	1.31	1.38
2	A	501	NDP	C5A-C4A	5.26	1.48	1.39
2	B	501	NDP	C5A-C4A	4.74	1.47	1.39
2	A	501	NDP	C5A-C6A	3.20	1.49	1.41
2	A	501	NDP	PN-O3	3.17	1.62	1.59
2	B	501	NDP	C5A-N7A	-2.83	1.33	1.39
2	B	501	NDP	C6N-C5N	2.69	1.41	1.33
2	A	501	NDP	C8A-N7A	2.38	1.36	1.31
2	A	501	NDP	C6N-C5N	2.29	1.40	1.33
2	A	501	NDP	PA-O3	2.27	1.61	1.59
2	A	501	NDP	C5A-N7A	-2.27	1.34	1.39
2	B	501	NDP	C5A-C6A	2.19	1.47	1.41
2	B	501	NDP	C8A-N7A	2.12	1.35	1.31

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	504	1BZ	C4-C3-N1	14.67	120.03	114.32
4	A	504	1BZ	C3-N1-C2	-12.54	119.58	126.30
2	A	501	NDP	C5A-C4A-N3A	-6.46	117.82	126.72
2	B	501	NDP	C5A-C4A-N3A	-5.76	118.79	126.72
4	A	504	1BZ	O2-N1-C3	5.30	120.55	116.75
2	A	501	NDP	N3A-C4A-N9A	4.64	135.06	127.17
2	B	501	NDP	N3A-C4A-N9A	4.63	135.04	127.17
4	A	504	1BZ	C1-C2-N1	4.46	119.70	117.11
2	A	501	NDP	C4A-C5A-N7A	-3.90	106.13	110.58
2	A	501	NDP	C2A-N3A-C4A	3.78	121.07	111.83
4	A	504	1BZ	C3-C4-C5	-3.49	120.30	122.49
2	B	501	NDP	C1D-N1N-C2N	-3.46	115.43	121.14
2	B	501	NDP	C4A-C5A-N7A	-3.32	106.78	110.58
4	A	504	1BZ	C10-C6-C2	-3.24	107.01	113.79
2	B	501	NDP	C2A-N3A-C4A	3.12	119.46	111.83
2	A	501	NDP	N3A-C2A-N1A	-2.80	124.34	128.58
2	B	501	NDP	C5A-N7A-C8A	2.74	107.76	103.45
2	A	501	NDP	O2B-P2B-O1X	-2.70	99.70	109.33
2	B	501	NDP	N3A-C2A-N1A	-2.58	124.68	128.58
4	A	504	1BZ	O1-C3-C4	-2.55	119.31	125.61
2	A	501	NDP	C5A-N7A-C8A	2.51	107.39	103.45
2	B	501	NDP	C4A-N9A-C8A	2.41	108.27	105.74
2	A	501	NDP	C3D-C2D-C1D	2.23	105.69	101.46
4	A	504	1BZ	O2-N1-C2	2.22	119.87	117.33
2	B	501	NDP	O4D-C1D-N1N	2.19	112.26	108.08

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	NDP	O4D-C1D-N1N	2.19	112.26	108.08
2	A	501	NDP	C6A-C5A-N7A	2.10	136.15	132.09
2	B	501	NDP	N9A-C8A-N7A	-2.05	111.02	113.94
2	A	501	NDP	O4B-C1B-N9A	2.03	111.99	108.09

There are no chirality outliers.

All (26) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	NDP	C5D-O5D-PN-O3
2	B	501	NDP	C5D-O5D-PN-O3
2	B	501	NDP	O4D-C4D-C5D-O5D
2	B	501	NDP	C2N-C3N-C7N-N7N
2	A	501	NDP	O4D-C4D-C5D-O5D
2	B	501	NDP	C3D-C4D-C5D-O5D
2	B	501	NDP	C1B-C2B-O2B-P2B
2	B	501	NDP	C3B-C2B-O2B-P2B
2	A	501	NDP	C3D-C4D-C5D-O5D
4	A	504	1BZ	N1-C2-C6-C10
2	A	501	NDP	C2B-O2B-P2B-O1X
2	A	501	NDP	C5D-O5D-PN-O1N
2	B	501	NDP	C5B-O5B-PA-O1A
2	B	501	NDP	C5B-O5B-PA-O2A
2	B	501	NDP	C5B-O5B-PA-O3
2	B	501	NDP	C5D-O5D-PN-O1N
4	A	504	1BZ	C1-C2-C6-C10
2	B	501	NDP	O4D-C1D-N1N-C2N
2	A	501	NDP	O4D-C1D-N1N-C2N
2	A	501	NDP	C2N-C3N-C7N-N7N
2	B	501	NDP	O4B-C4B-C5B-O5B
2	B	501	NDP	C2D-C1D-N1N-C2N
2	A	501	NDP	C2D-C1D-N1N-C2N
2	B	501	NDP	O4D-C1D-N1N-C6N
2	B	501	NDP	C2B-O2B-P2B-O1X
2	A	501	NDP	PN-O3-PA-O2A

There are no ring outliers.

3 monomers are involved in 6 short contacts:

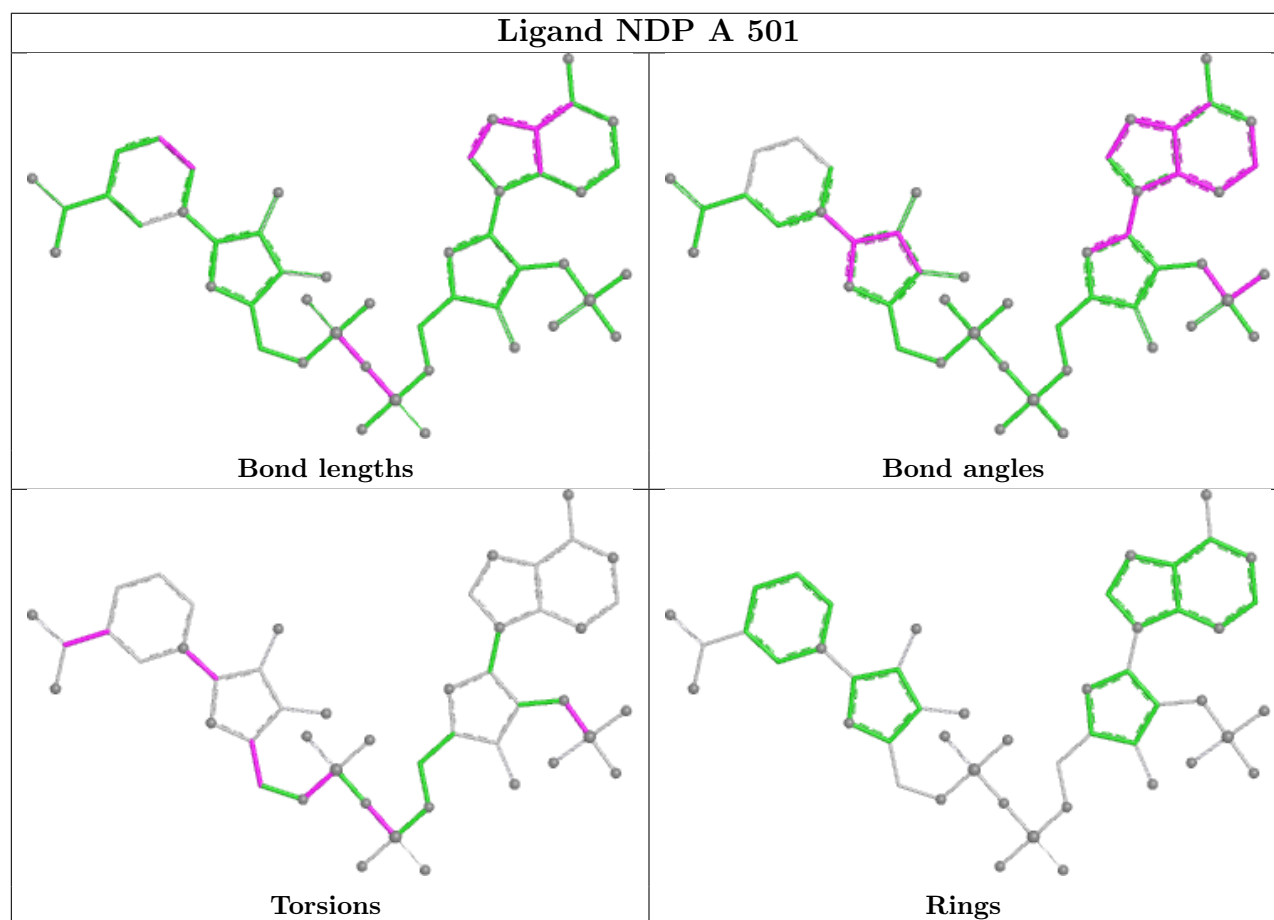
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	NDP	5	0

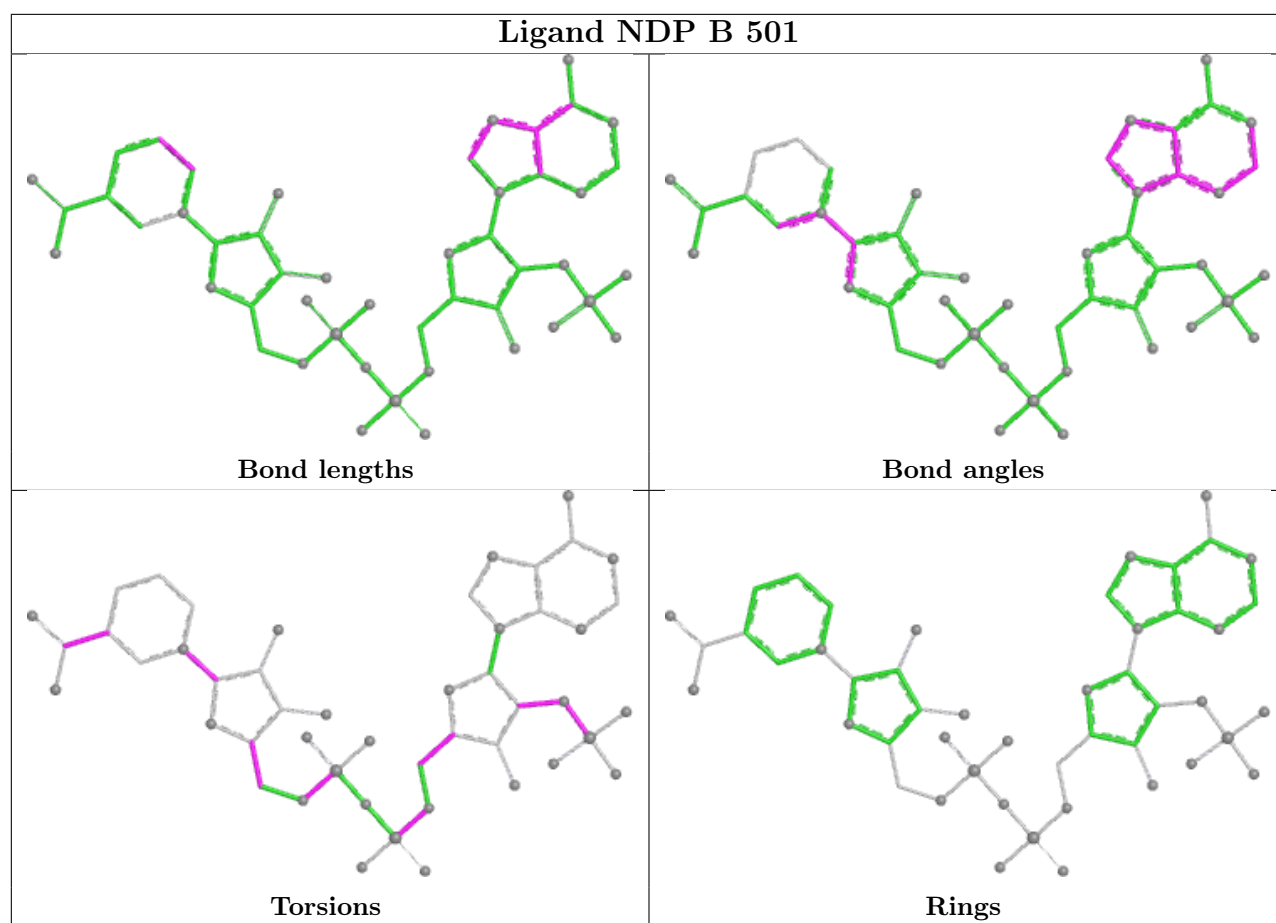
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	501	NDP	1	0
4	A	504	1BZ	1	0

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 [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	367/414 (88%)	-0.17	5 (1%) 73 55	26, 45, 76, 100	0
1	B	358/414 (86%)	-0.05	8 (2%) 62 43	23, 44, 81, 115	0
All	All	725/828 (87%)	-0.11	13 (1%) 67 49	23, 45, 78, 115	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	163	GLN	4.1
1	B	164	LYS	3.8
1	A	127	PRO	3.2
1	B	251	ILE	2.7
1	B	298	PRO	2.5
1	B	216	LEU	2.5
1	A	80	GLU	2.5
1	B	254	MET	2.4
1	A	112	ILE	2.3
1	B	133	HIS	2.3
1	B	220	ASP	2.2
1	A	113	ILE	2.1
1	A	271	ASN	2.0

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

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

6.3 Carbohydrates [i](#)

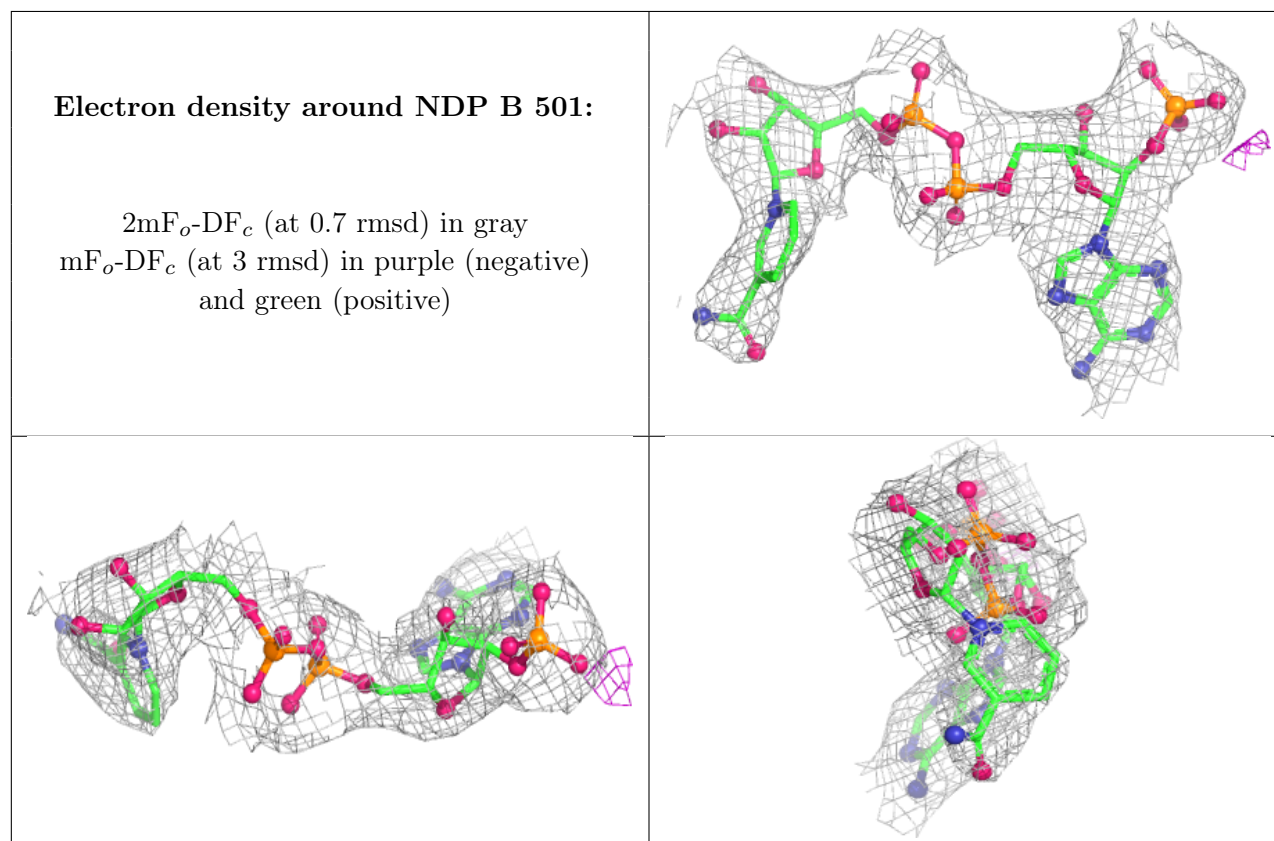
There are no oligosaccharides in this entry.

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.

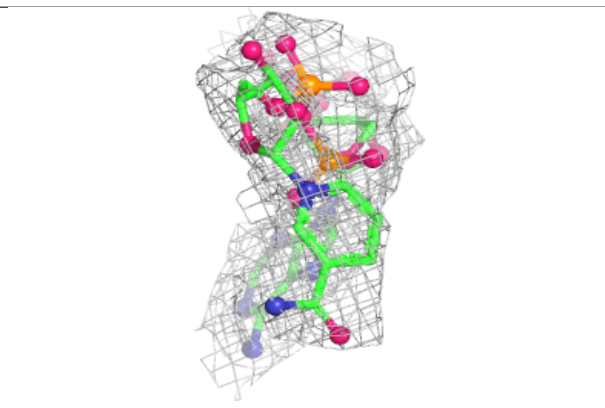
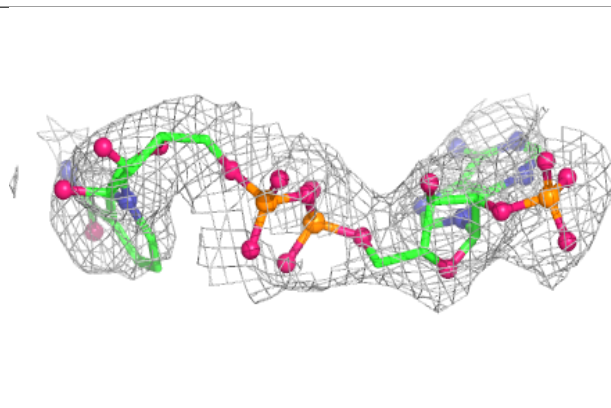
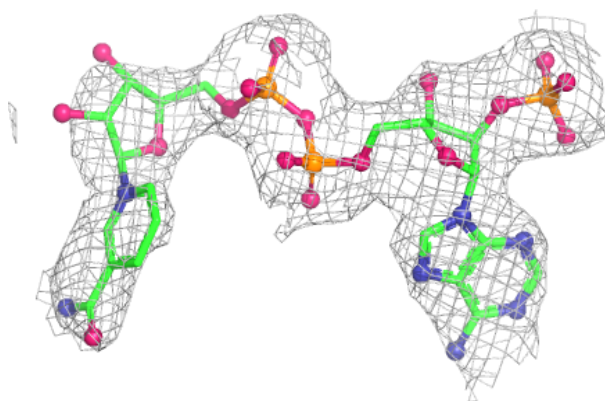
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	1BZ	A	504	16/16	0.81	0.21	88,94,100,101	0
3	SO4	A	502	5/5	0.82	0.12	89,92,94,96	0
3	SO4	A	503	5/5	0.86	0.09	83,83,86,88	0
2	NDP	B	501	48/48	0.94	0.09	32,45,69,70	0
2	NDP	A	501	48/48	0.94	0.08	49,58,79,82	0

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 NDP A 501:

$2mF_o-DF_c$ (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.