



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 19, 2026 – 03:39 AM UTC

PDB ID : 6O2Z / pdb_00006o2z
Title : Crystal structure of IDH1 R132H mutant in complex with compound 32
Authors : Toms, A.V.; Lin, J.
Deposited on : 2019-02-25
Resolution : 2.50 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

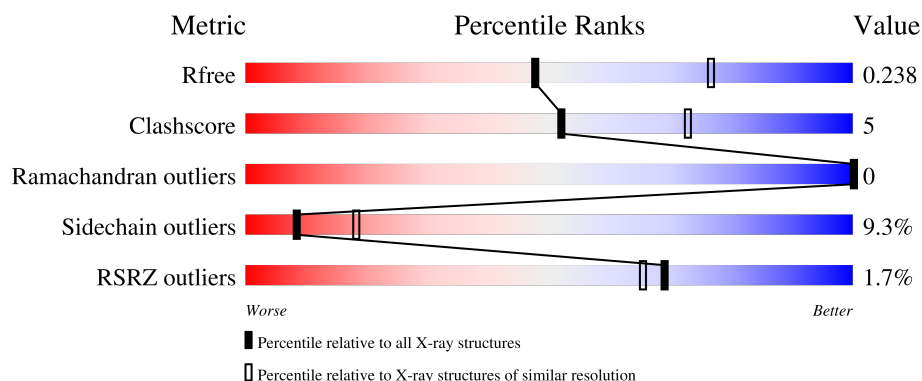
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

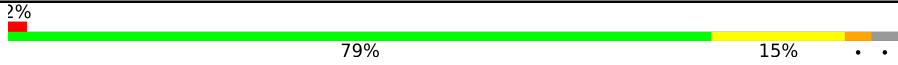
The reported resolution of this entry is 2.50 Å.

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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	425	
1	B	425	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6774 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.

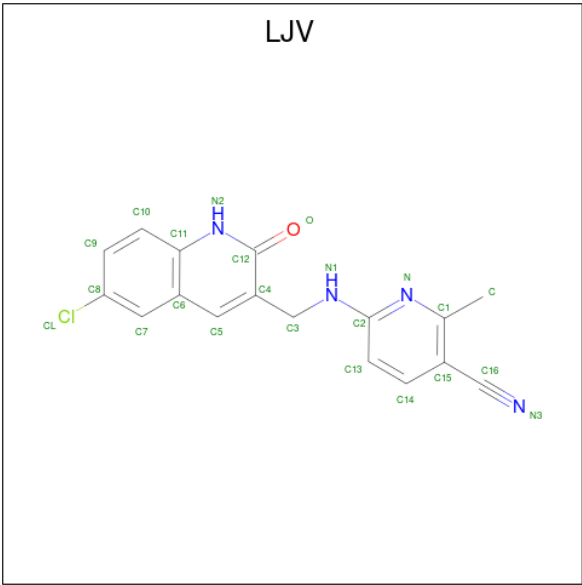
- Molecule 1 is a protein called Isocitrate dehydrogenase [NADP] cytoplasmic.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	411	Total	C	N	O	S	0	0	0
			3243	2060	549	616	18			
1	B	411	Total	C	N	O	S	0	0	0
			3237	2057	546	616	18			

There are 24 discrepancies between the modelled and reference sequences:

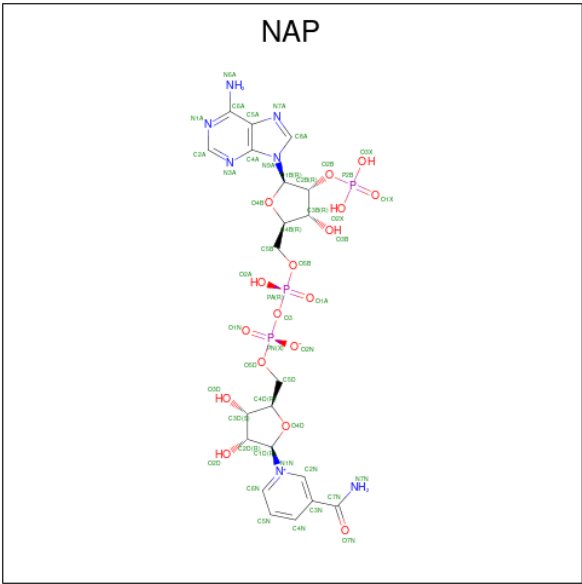
Chain	Residue	Modelled	Actual	Comment	Reference
A	132	HIS	ARG	engineered mutation	UNP O75874
A	415	SER	-	expression tag	UNP O75874
A	416	LEU	-	expression tag	UNP O75874
A	417	GLU	-	expression tag	UNP O75874
A	418	HIS	-	expression tag	UNP O75874
A	419	HIS	-	expression tag	UNP O75874
A	420	HIS	-	expression tag	UNP O75874
A	421	HIS	-	expression tag	UNP O75874
A	422	HIS	-	expression tag	UNP O75874
A	423	HIS	-	expression tag	UNP O75874
A	424	HIS	-	expression tag	UNP O75874
A	425	HIS	-	expression tag	UNP O75874
B	132	HIS	ARG	engineered mutation	UNP O75874
B	415	SER	-	expression tag	UNP O75874
B	416	LEU	-	expression tag	UNP O75874
B	417	GLU	-	expression tag	UNP O75874
B	418	HIS	-	expression tag	UNP O75874
B	419	HIS	-	expression tag	UNP O75874
B	420	HIS	-	expression tag	UNP O75874
B	421	HIS	-	expression tag	UNP O75874
B	422	HIS	-	expression tag	UNP O75874
B	423	HIS	-	expression tag	UNP O75874
B	424	HIS	-	expression tag	UNP O75874
B	425	HIS	-	expression tag	UNP O75874

- Molecule 2 is 6-{{[(6-chloro-2-oxo-1,2-dihydroquinolin-3-yl)methyl]amino}-2-methylpyridine-3-carbonitrile (CCD ID: LJV) (formula: C₁₇H₁₃ClN₄O) (labeled as "Ligand of Interest" by depositor).



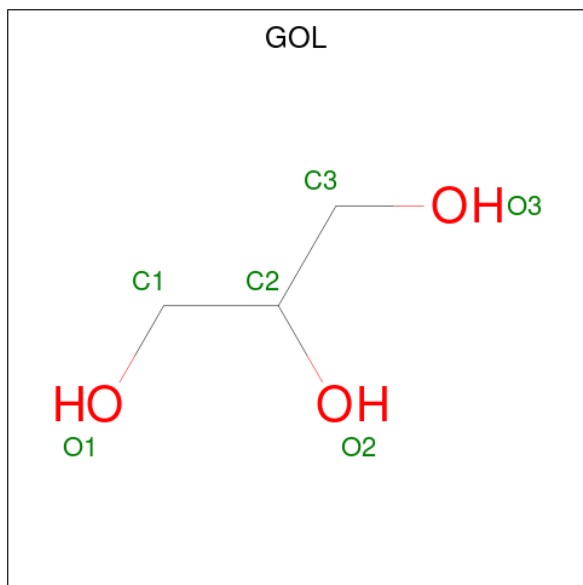
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	Cl	N	O		
2	A	1	23	17	1	4	1	0	0
2	B	1	23	17	1	4	1	0	0

- Molecule 3 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (CCD ID: NAP) (formula: C₂₁H₂₈N₇O₁₇P₃).



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

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

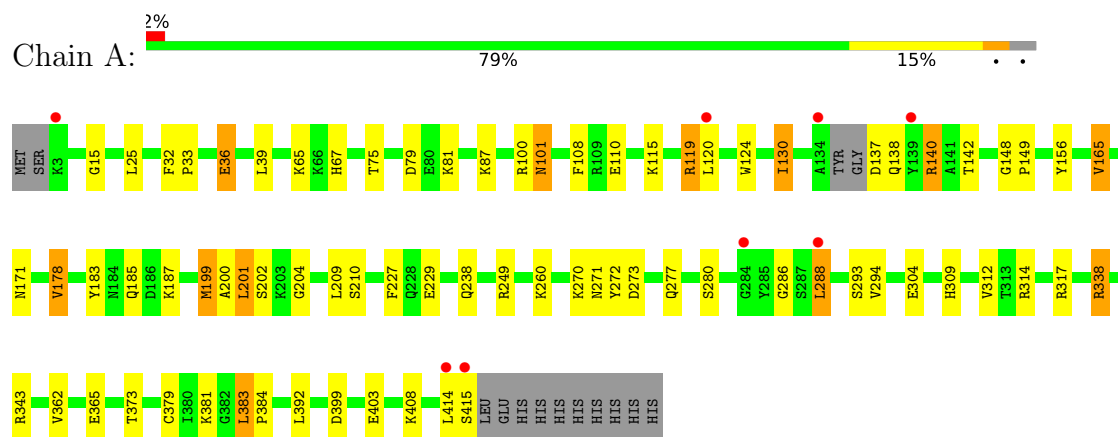
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	77	Total	O	0	0
			77	77		
5	B	63	Total	O	0	0
			63	63		

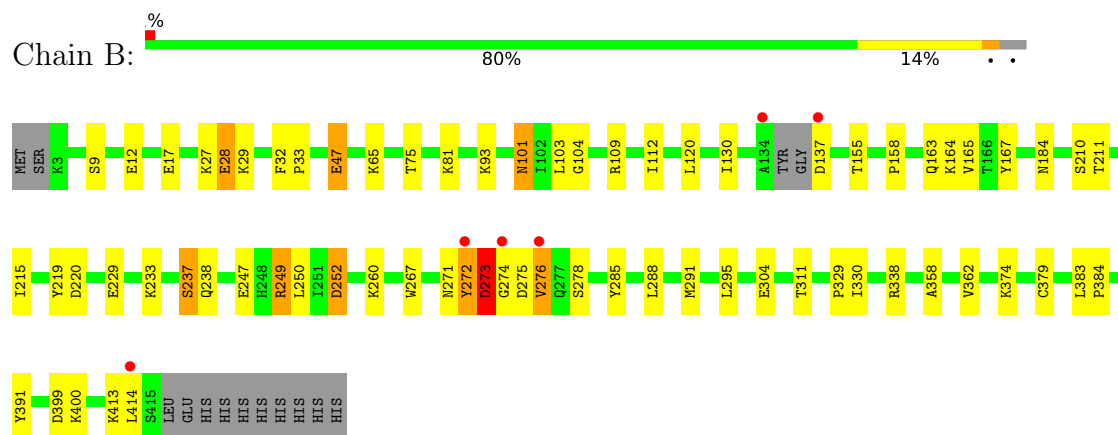
3 Residue-property plots [i](#)

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: 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	82.34Å 82.34Å 303.74Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.25 – 2.50 46.25 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.5 (46.25-2.50) 99.7 (46.25-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.44 (at 2.48Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.172 , 0.234 0.178 , 0.238	Depositor DCC
R_{free} test set	1894 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	44.4	Xtriage
Anisotropy	0.037	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 28.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6774	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.98% 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: LJV, GOL, NAP

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	1.00	0/3310	1.45	4/4466 (0.1%)
1	B	0.99	0/3304	1.41	2/4459 (0.0%)
All	All	1.00	0/6614	1.43	6/8925 (0.1%)

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	B	0	1

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	273	ASP	CA-CB-CG	5.81	118.41	112.60
1	B	101	ASN	CB-CA-C	5.72	121.65	110.67
1	A	399	ASP	CA-CB-CG	5.33	117.93	112.60
1	A	15	GLY	CA-C-O	-5.24	117.01	121.78
1	A	373	THR	CA-C-N	5.20	127.25	120.28
1	A	373	THR	C-N-CA	5.20	127.25	120.28

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	273	ASP	Peptide

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	3243	0	3192	37	0
1	B	3237	0	3181	36	0
2	A	23	0	0	2	0
2	B	23	0	0	0	0
3	A	48	0	25	1	0
3	B	48	0	25	1	0
4	A	6	0	8	0	0
4	B	6	0	8	0	0
5	A	77	0	0	2	0
5	B	63	0	0	3	0
All	All	6774	0	6439	66	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 5.

All (66) 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:288:LEU:HB3	1:A:309:HIS:HB3	1.56	0.86
1:B:29:LYS:NZ	1:B:399:ASP:OD1	2.13	0.81
1:A:286:GLY:HA2	1:A:379:CYS:SG	2.33	0.68
1:B:137:ASP:OD2	5:B:601:HOH:O	2.12	0.67
1:A:119:ARG:HD2	1:A:124:TRP:O	1.94	0.66
1:B:75:THR:O	3:B:501:NAP:H2N	2.03	0.59
1:B:247:GLU:OE2	1:B:249:ARG:NH2	2.31	0.59
1:B:12:GLU:OE2	1:B:27:LYS:NZ	2.33	0.57
1:A:178:VAL:HG13	1:B:219:TYR:CE2	2.40	0.57
1:A:119:ARG:CD	1:A:124:TRP:O	2.53	0.57
1:A:140:ARG:HB3	1:A:183:TYR:CE1	2.43	0.54
1:A:178:VAL:HG13	1:B:219:TYR:CD2	2.42	0.54
1:B:47:GLU:H	1:B:47:GLU:CD	2.18	0.52
1:A:101:ASN:HD22	1:A:101:ASN:C	2.18	0.52
1:A:119:ARG:NH2	5:A:605:HOH:O	2.42	0.51
1:A:210:SER:HA	1:A:249:ARG:O	2.09	0.51
1:B:210:SER:HA	1:B:249:ARG:O	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:211:THR:HB	1:B:220:ASP:HB3	1.93	0.50
1:A:156:TYR:HB3	1:A:165:VAL:HG22	1.94	0.49
1:B:109:ARG:HD3	1:B:276:VAL:HB	1.94	0.49
1:B:130:ILE:HD13	1:B:267:TRP:HB3	1.94	0.49
1:A:142:THR:HG21	1:B:167:TYR:HB3	1.96	0.48
1:A:36:GLU:OE2	1:A:36:GLU:HA	2.14	0.48
1:A:178:VAL:CG2	1:B:219:TYR:HA	2.44	0.48
1:A:277:GLN:HB3	1:B:252:ASP:HB3	1.97	0.47
1:A:138:GLN:HG3	1:B:215:ILE:HG21	1.98	0.46
1:A:383:LEU:N	1:A:384:PRO:CD	2.79	0.46
1:B:17:GLU:HB2	1:B:311:THR:HB	1.98	0.46
1:B:272:TYR:CD1	1:B:272:TYR:N	2.84	0.46
1:B:288:LEU:HD11	1:B:379:CYS:SG	2.56	0.46
1:B:158:PRO:CG	1:B:163:GLN:HG2	2.47	0.45
1:B:28:GLU:HG2	1:B:29:LYS:HG3	1.99	0.45
1:B:383:LEU:N	1:B:384:PRO:CD	2.80	0.45
1:A:101:ASN:C	1:A:101:ASN:ND2	2.74	0.45
1:A:130:ILE:HG13	2:A:500:LJV:C4	2.46	0.44
1:B:237:SER:HB2	5:B:657:HOH:O	2.17	0.44
1:B:238:GLN:HG3	5:B:657:HOH:O	2.18	0.43
1:A:271:ASN:ND2	1:A:273:ASP:O	2.51	0.43
1:A:272:TYR:HB3	1:B:272:TYR:CE1	2.53	0.43
1:A:67:HIS:O	1:A:343:ARG:NH2	2.51	0.43
1:B:272:TYR:C	1:B:274:GLY:H	2.27	0.43
1:B:32:PHE:N	1:B:33:PRO:CD	2.83	0.42
1:A:185:GLN:NE2	5:A:612:HOH:O	2.51	0.42
1:B:291:MET:HE3	1:B:291:MET:HB2	1.87	0.42
1:B:374:LYS:HG3	1:B:391:TYR:CE2	2.54	0.42
1:A:79:ASP:OD1	1:A:79:ASP:C	2.62	0.42
1:A:365:GLU:OE1	1:A:408:LYS:NZ	2.52	0.42
1:A:75:THR:O	3:A:501:NAP:H2N	2.20	0.42
1:B:104:GLY:O	1:B:295:LEU:HD11	2.20	0.41
1:B:229:GLU:O	1:B:233:LYS:HG3	2.20	0.41
1:A:209:LEU:HD22	1:A:227:PHE:CG	2.54	0.41
1:A:110:GLU:CD	1:A:338:ARG:HH12	2.28	0.41
1:A:201:LEU:HD12	1:A:201:LEU:HA	1.94	0.41
1:B:158:PRO:HG2	1:B:163:GLN:HG2	2.02	0.41
1:A:32:PHE:N	1:A:33:PRO:CD	2.84	0.41
1:A:187:LYS:HA	1:A:187:LYS:HD2	1.85	0.41
1:A:199:MET:CE	1:A:294:VAL:HG21	2.51	0.41
1:A:200:ALA:O	1:A:204:GLY:N	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:358:ALA:O	1:B:362:VAL:HG23	2.21	0.41
1:A:130:ILE:HG13	2:A:500:LJV:C12	2.51	0.41
1:B:112:ILE:HD13	1:B:330:ILE:HG22	2.03	0.41
1:A:108:PHE:O	1:A:293:SER:HA	2.22	0.40
1:B:65:LYS:HG3	1:B:103:LEU:CD2	2.51	0.40
1:A:148:GLY:O	1:A:149:PRO:C	2.64	0.40
1:A:362:VAL:HG23	1:A:408:LYS:HD2	2.04	0.40
1:B:155:THR:HA	1:B:165:VAL:O	2.21	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	407/425 (96%)	391 (96%)	16 (4%)	0	100	100
1	B	407/425 (96%)	382 (94%)	25 (6%)	0	100	100
All	All	814/850 (96%)	773 (95%)	41 (5%)	0	100	100

There are no Ramachandran outliers to report.

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	345/361 (96%)	308 (89%)	37 (11%)	6	13

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	344/361 (95%)	317 (92%)	27 (8%)	11	24
All	All	689/722 (95%)	625 (91%)	64 (9%)	8	18

All (64) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	LEU
1	A	36	GLU
1	A	39	LEU
1	A	65	LYS
1	A	81	LYS
1	A	87	LYS
1	A	100	ARG
1	A	101	ASN
1	A	115	LYS
1	A	119	ARG
1	A	120	LEU
1	A	130	ILE
1	A	137	ASP
1	A	140	ARG
1	A	165	VAL
1	A	171	ASN
1	A	178	VAL
1	A	199	MET
1	A	201	LEU
1	A	202	SER
1	A	229	GLU
1	A	238	GLN
1	A	260	LYS
1	A	270	LYS
1	A	280	SER
1	A	288	LEU
1	A	304	GLU
1	A	312	VAL
1	A	314	ARG
1	A	317	ARG
1	A	338	ARG
1	A	381	LYS
1	A	383	LEU
1	A	392	LEU
1	A	403	GLU
1	A	414	LEU

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Mol	Chain	Res	Type
1	A	415	SER
1	B	9	SER
1	B	28	GLU
1	B	47	GLU
1	B	81	LYS
1	B	93	LYS
1	B	101	ASN
1	B	120	LEU
1	B	164	LYS
1	B	184	ASN
1	B	237	SER
1	B	249	ARG
1	B	250	LEU
1	B	252	ASP
1	B	260	LYS
1	B	271	ASN
1	B	272	TYR
1	B	273	ASP
1	B	275	ASP
1	B	276	VAL
1	B	278	SER
1	B	285	TYR
1	B	304	GLU
1	B	329	PRO
1	B	338	ARG
1	B	400	LYS
1	B	413	LYS
1	B	414	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	53	ASN
1	A	68	ASN
1	A	101	ASN
1	A	138	GLN
1	A	228	GLN
1	A	320	GLN
1	B	53	ASN
1	B	68	ASN
1	B	170	HIS
1	B	198	GLN

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Mol	Chain	Res	Type
1	B	228	GLN
1	B	242	GLN
1	B	271	ASN

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](#)

6 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	NAP	B	501	-	50,52,52	1.56	5 (10%)	71,80,80	1.16	3 (4%)
4	GOL	A	502	-	5,5,5	0.09	0	5,5,5	0.29	0
2	LJV	A	500	-	24,25,25	2.41	5 (20%)	29,35,35	1.63	5 (17%)
3	NAP	A	501	-	50,52,52	1.60	6 (12%)	71,80,80	1.21	3 (4%)
4	GOL	B	502	-	5,5,5	0.13	0	5,5,5	0.31	0
2	LJV	B	500	-	24,25,25	2.37	3 (12%)	29,35,35	1.44	6 (20%)

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	NAP	B	501	-	-	9/35/67/67	0/5/5/5
4	GOL	A	502	-	-	2/4/4/4	-
2	LJV	A	500	-	-	2/7/7/7	0/3/3/3
3	NAP	A	501	-	-	9/35/67/67	0/5/5/5
4	GOL	B	502	-	-	2/4/4/4	-
2	LJV	B	500	-	-	1/7/7/7	0/3/3/3

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	500	LJV	C15-C16	-9.72	1.30	1.44
2	A	500	LJV	C15-C16	-9.26	1.31	1.44
3	B	501	NAP	C4N-C3N	7.37	1.50	1.39
3	A	501	NAP	C4N-C3N	6.97	1.50	1.39
3	A	501	NAP	C5N-C4N	6.17	1.49	1.38
3	B	501	NAP	C5N-C4N	5.78	1.48	1.38
2	A	500	LJV	C5-C4	3.72	1.40	1.34
2	B	500	LJV	C5-C4	3.21	1.39	1.34
2	A	500	LJV	C11-N2	-2.96	1.34	1.39
2	A	500	LJV	C12-C4	-2.95	1.38	1.45
3	A	501	NAP	C6N-C5N	-2.93	1.32	1.38
3	B	501	NAP	C2N-N1N	2.69	1.37	1.35
2	B	500	LJV	C11-N2	-2.67	1.35	1.39
3	B	501	NAP	C6N-C5N	-2.65	1.33	1.38
2	A	500	LJV	C7-C8	2.61	1.42	1.38
3	A	501	NAP	C2N-N1N	2.46	1.37	1.35
3	A	501	NAP	P2B-O2B	2.41	1.63	1.59
3	A	501	NAP	C6N-N1N	2.20	1.40	1.35
3	B	501	NAP	C6N-N1N	2.15	1.40	1.35

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	501	NAP	C5N-C4N-C3N	-6.84	113.64	120.36
3	A	501	NAP	C5N-C4N-C3N	-6.48	113.99	120.36
2	A	500	LJV	C2-N-C1	4.57	124.50	118.16
3	B	501	NAP	C6N-N1N-C2N	-3.74	118.69	121.88
2	A	500	LJV	O-C12-C4	-3.54	118.75	124.71
3	A	501	NAP	O2B-P2B-O1X	-3.44	97.08	109.33
2	A	500	LJV	C4-C12-N2	3.29	120.48	115.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	500	LJV	C4-C12-N2	3.03	120.07	115.30
2	A	500	LJV	C13-C2-N	-3.00	118.12	123.15
2	B	500	LJV	C15-C16-N3	-2.78	173.33	177.90
2	B	500	LJV	C11-N2-C12	-2.71	121.02	124.82
3	A	501	NAP	C6N-N1N-C2N	-2.69	119.59	121.88
2	B	500	LJV	C6-C5-C4	-2.62	118.14	121.90
2	A	500	LJV	C15-C1-N	-2.46	119.48	121.60
3	B	501	NAP	P2B-O2B-C2B	-2.25	117.43	123.43
2	B	500	LJV	C6-C11-N2	2.17	120.50	118.66
2	B	500	LJV	O-C12-C4	-2.14	121.12	124.71

There are no chirality outliers.

All (25) torsion outliers are listed below:

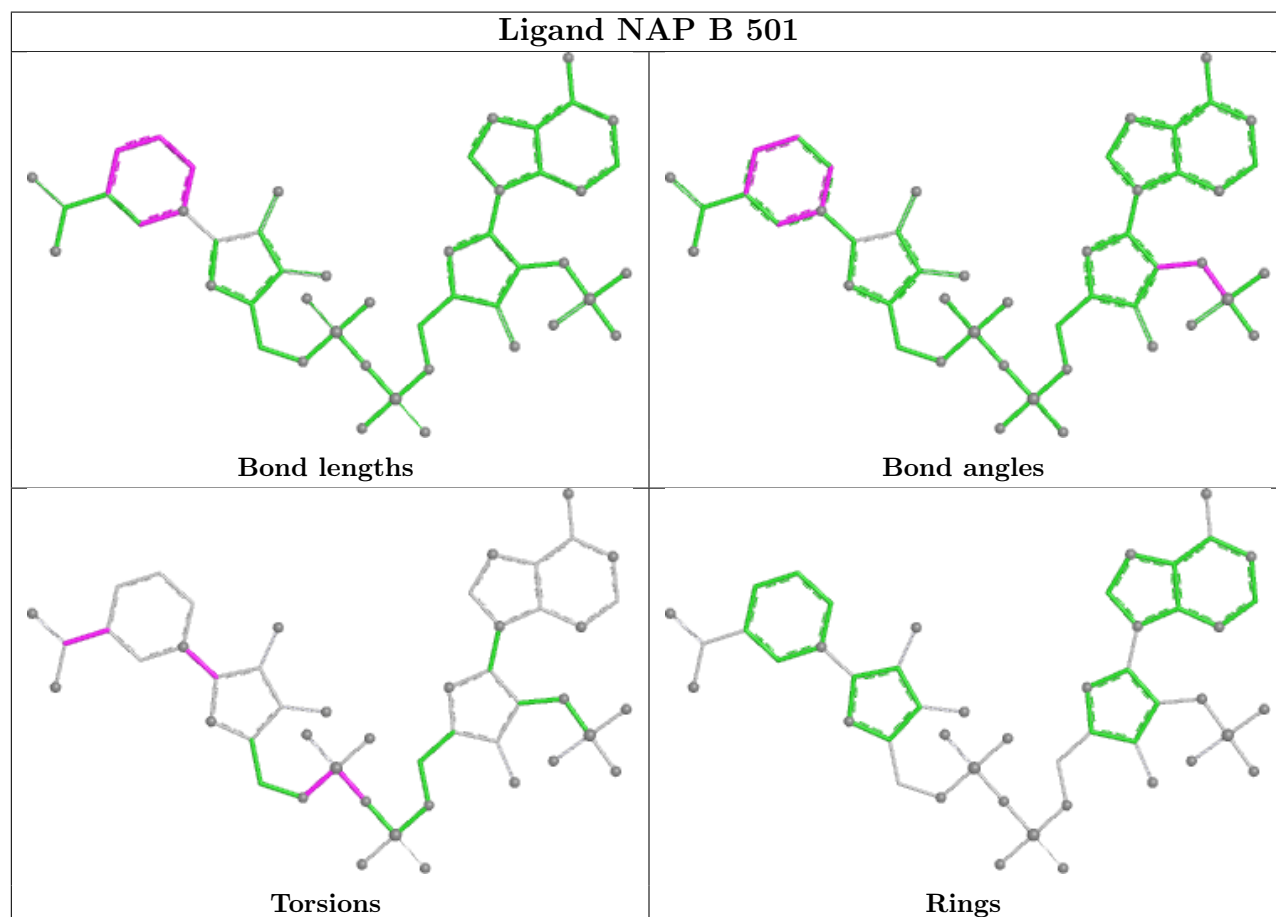
Mol	Chain	Res	Type	Atoms
2	A	500	LJV	N1-C3-C4-C5
2	A	500	LJV	C4-C3-N1-C2
2	B	500	LJV	C4-C3-N1-C2
3	A	501	NAP	C5D-O5D-PN-O3
3	A	501	NAP	C5D-O5D-PN-O2N
3	A	501	NAP	O4D-C1D-N1N-C2N
3	A	501	NAP	O4D-C1D-N1N-C6N
3	A	501	NAP	C2D-C1D-N1N-C2N
3	A	501	NAP	C2D-C1D-N1N-C6N
3	B	501	NAP	C5D-O5D-PN-O3
3	B	501	NAP	C5D-O5D-PN-O1N
3	B	501	NAP	C5D-O5D-PN-O2N
3	B	501	NAP	O4D-C1D-N1N-C6N
4	A	502	GOL	C1-C2-C3-O3
4	B	502	GOL	O1-C1-C2-C3
4	A	502	GOL	O2-C2-C3-O3
4	B	502	GOL	O1-C1-C2-O2
3	A	501	NAP	O4B-C4B-C5B-O5B
3	A	501	NAP	C3B-C4B-C5B-O5B
3	B	501	NAP	PA-O3-PN-O1N
3	A	501	NAP	C5D-O5D-PN-O1N
3	B	501	NAP	PA-O3-PN-O2N
3	B	501	NAP	O4D-C1D-N1N-C2N
3	B	501	NAP	C4N-C3N-C7N-N7N
3	B	501	NAP	C4N-C3N-C7N-O7N

There are no ring outliers.

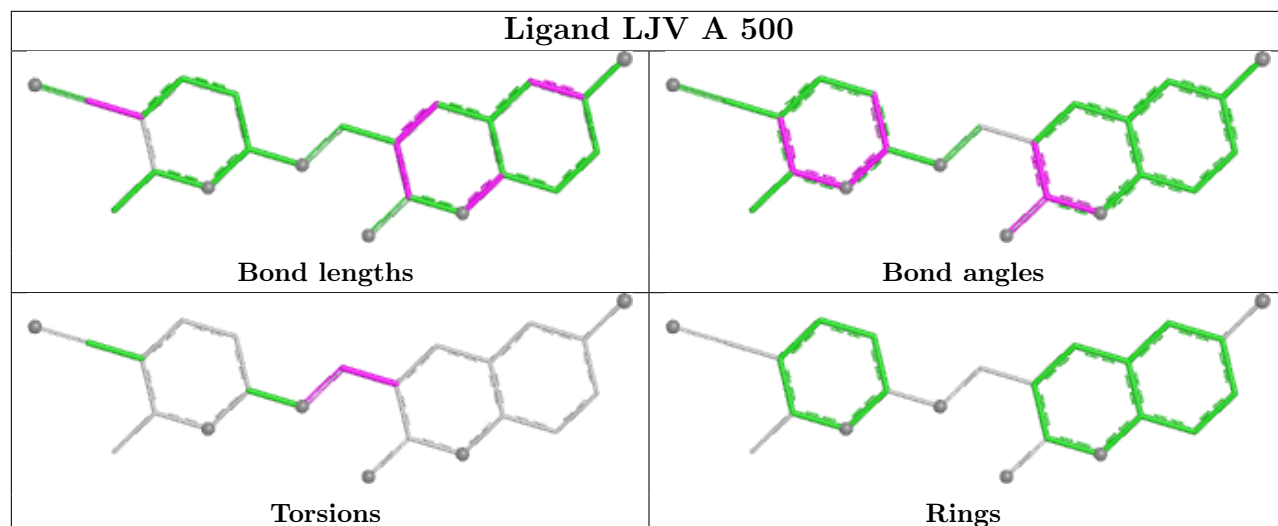
3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	501	NAP	1	0
2	A	500	LJV	2	0
3	A	501	NAP	1	0

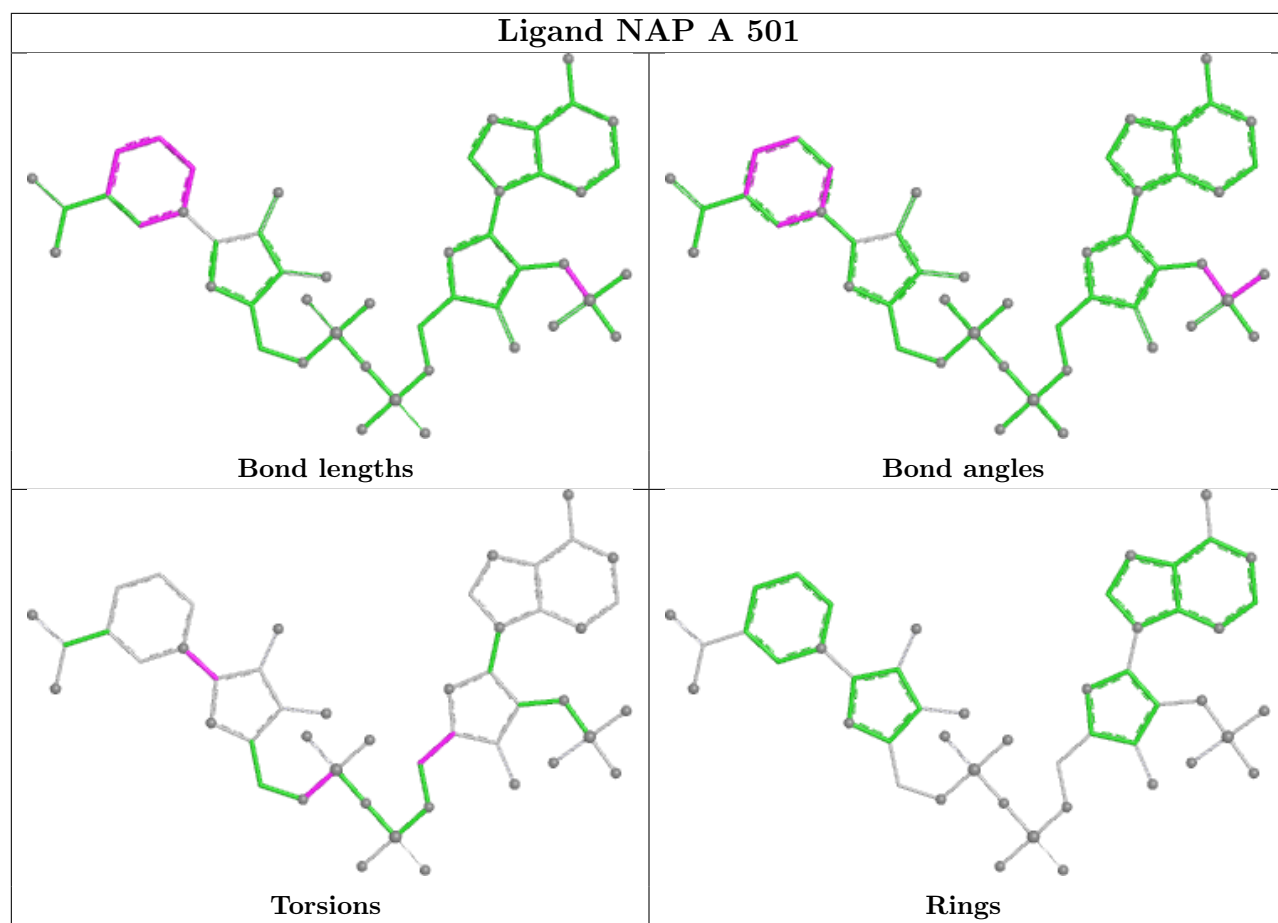
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

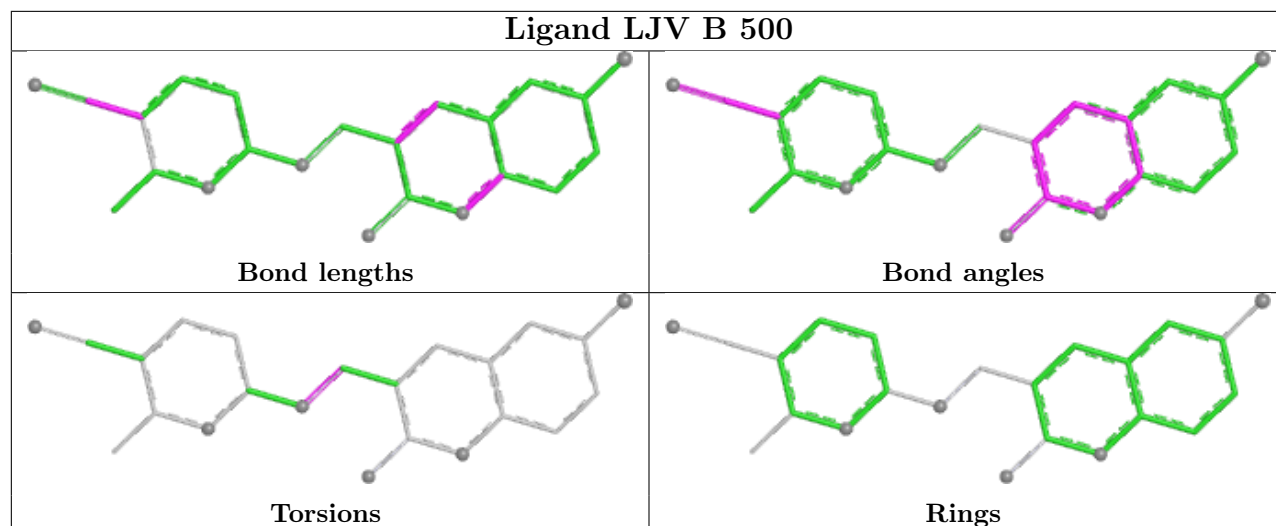


Ligand LJV A 500



Ligand NAP A 501





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	411/425 (96%)	-0.16	8 (1%) 66 62	27, 43, 72, 106	0
1	B	411/425 (96%)	-0.18	6 (1%) 72 68	29, 45, 72, 130	0
All	All	822/850 (96%)	-0.17	14 (1%) 69 65	27, 44, 72, 130	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	134	ALA	7.3
1	B	414	LEU	4.7
1	B	272	TYR	3.4
1	A	414	LEU	3.3
1	B	134	ALA	2.5
1	A	415	SER	2.5
1	A	284	GLY	2.5
1	B	276	VAL	2.4
1	A	120	LEU	2.4
1	A	3	LYS	2.2
1	B	137	ASP	2.1
1	A	288	LEU	2.1
1	A	139	TYR	2.0
1	B	274	GLY	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](#)

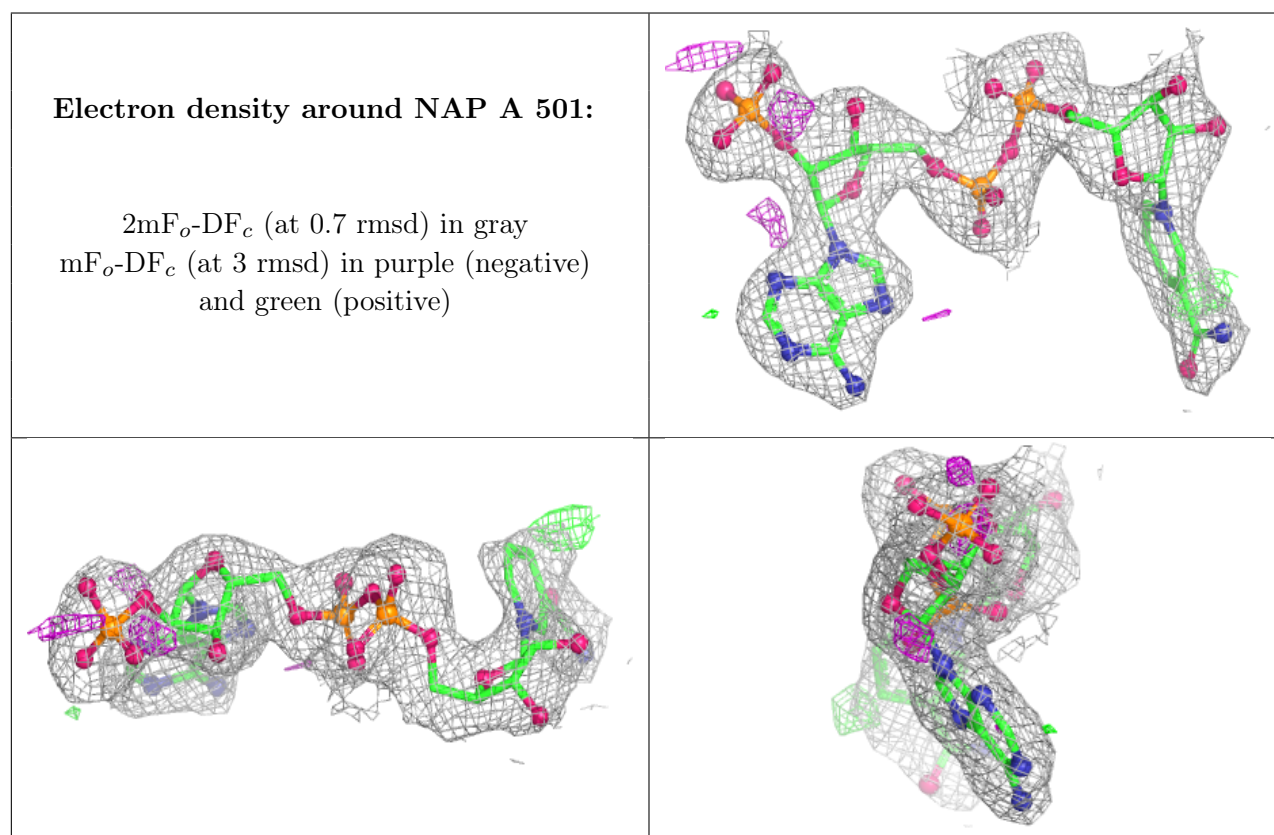
There are no oligosaccharides in this entry.

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.

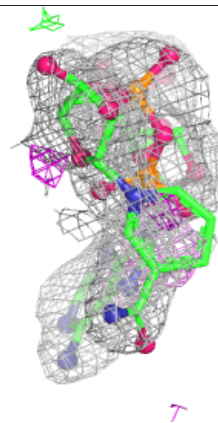
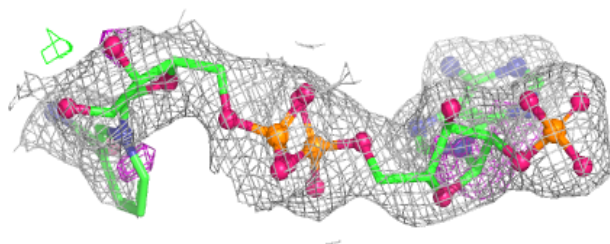
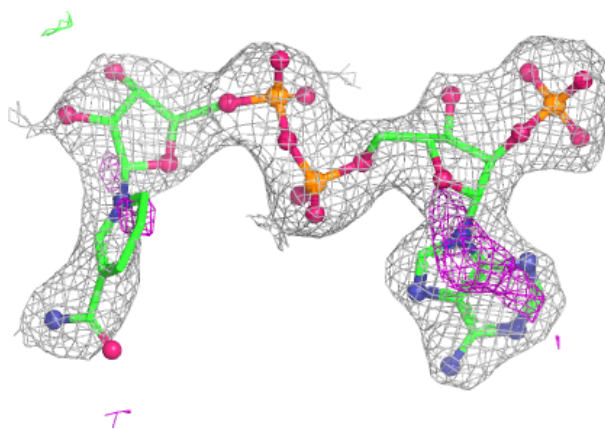
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	GOL	A	502	6/6	0.89	0.14	57,66,71,73	0
4	GOL	B	502	6/6	0.91	0.13	56,67,77,78	0
3	NAP	A	501	48/48	0.92	0.10	41,58,76,80	0
3	NAP	B	501	48/48	0.92	0.10	37,62,87,99	0
2	LJV	A	500	23/23	0.96	0.07	34,37,40,44	0
2	LJV	B	500	23/23	0.97	0.06	31,36,37,41	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.



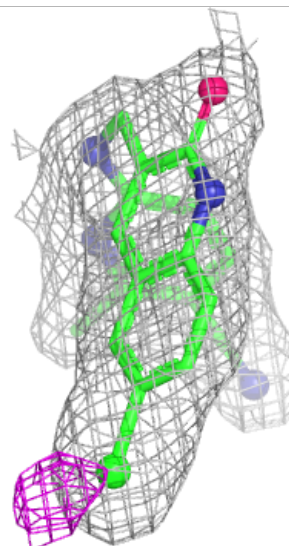
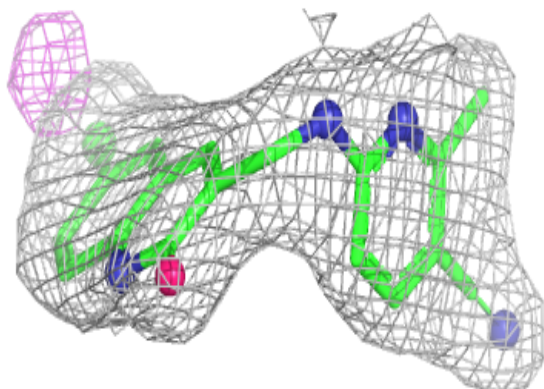
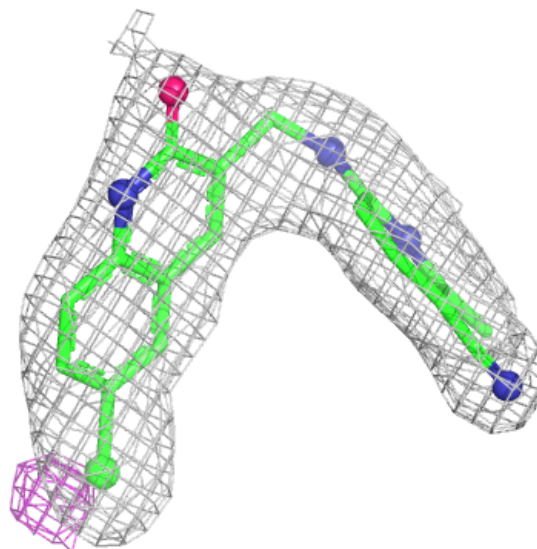
Electron density around NAP B 501:

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



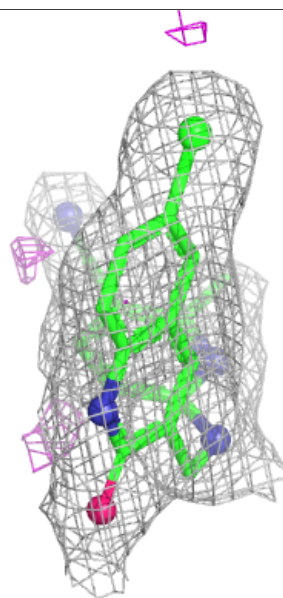
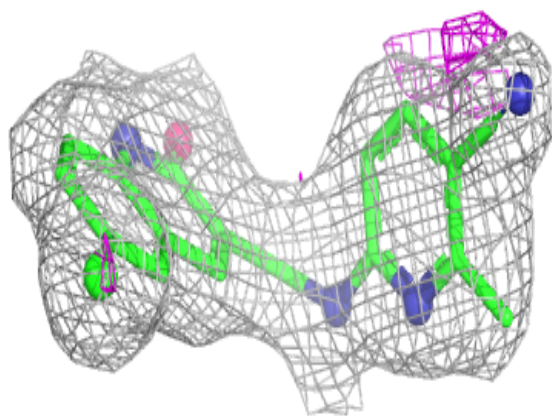
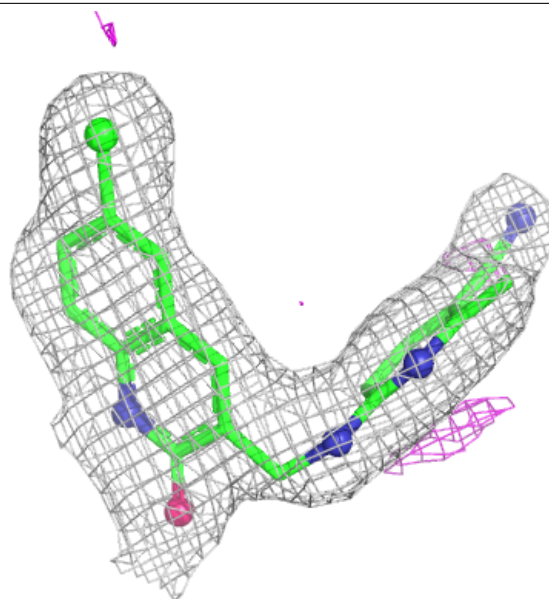
Electron density around LJV A 500:

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



Electron density around LJV B 500:

$2mF_o-DF_c$ (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.