



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 12:09 AM UTC

PDB ID : 4KP7 / pdb_00004kp7
Title : Structure of Plasmodium IspC in complex with a beta-thia-isostere derivative of Fosmidomycin
Authors : Kunfermann, A.; Bacher, A.; Groll, M.
Deposited on : 2013-05-13
Resolution : 2.00 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

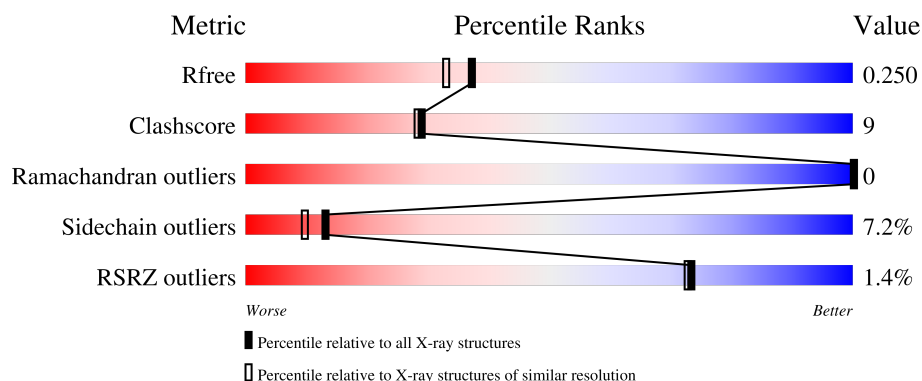
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	427	 80% 13% • •
1	B	427	 78% 14% • 6%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6790 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 1-deoxy-D-xylulose 5-phosphate reductoisomerase, apicoplast.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	409	Total	C	N	O	S	0	0	0
			3268	2099	535	614	20			
1	B	402	Total	C	N	O	S	0	0	0
			3200	2053	520	607	20			

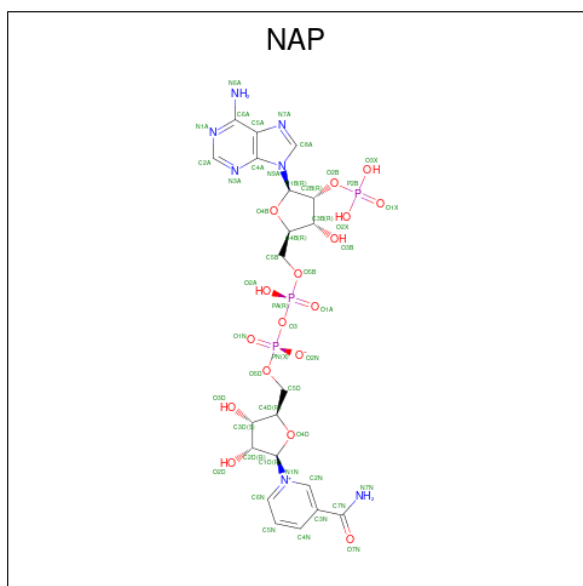
There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	62	MET	-	expression tag	UNP O96693
A	63	ARG	-	expression tag	UNP O96693
A	64	GLY	-	expression tag	UNP O96693
A	65	SER	-	expression tag	UNP O96693
A	66	HIS	-	expression tag	UNP O96693
A	67	HIS	-	expression tag	UNP O96693
A	68	HIS	-	expression tag	UNP O96693
A	69	HIS	-	expression tag	UNP O96693
A	70	HIS	-	expression tag	UNP O96693
A	71	HIS	-	expression tag	UNP O96693
A	72	GLY	-	expression tag	UNP O96693
A	73	SER	-	expression tag	UNP O96693
B	62	MET	-	expression tag	UNP O96693
B	63	ARG	-	expression tag	UNP O96693
B	64	GLY	-	expression tag	UNP O96693
B	65	SER	-	expression tag	UNP O96693
B	66	HIS	-	expression tag	UNP O96693
B	67	HIS	-	expression tag	UNP O96693
B	68	HIS	-	expression tag	UNP O96693
B	69	HIS	-	expression tag	UNP O96693
B	70	HIS	-	expression tag	UNP O96693
B	71	HIS	-	expression tag	UNP O96693
B	72	GLY	-	expression tag	UNP O96693
B	73	SER	-	expression tag	UNP O96693

- Molecule 2 is MANGANESE (III) ION (CCD ID: MN3) (formula: Mn).

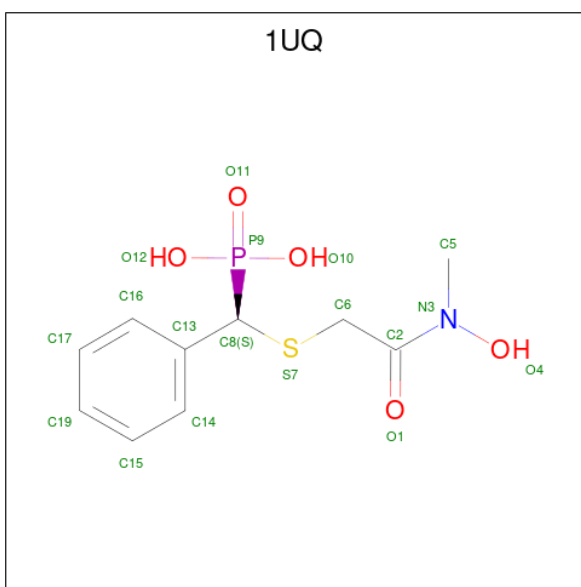
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mn	0	0
			1	1		
2	B	1	Total	Mn	0	0
			1	1		

- Molecule 3 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (CCD ID: NAP) (formula: $C_{21}H_{28}N_7O_{17}P_3$).



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

- Molecule 4 is [(S)-({2-[hydroxy(methyl)amino]-2-oxoethyl}sulfanyl)(phenyl)methyl]phosphonic acid (CCD ID: 1UQ) (formula: $C_{10}H_{14}NO_5PS$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	S	0	0
			18	10	1	5	1	1		
4	B	1	Total	C	N	O	P	S	0	0
			18	10	1	5	1	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	129	Total	O	0	0
			129	129		
5	B	107	Total	O	0	0
			107	107		

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	51.46Å 78.12Å 99.99Å 90.00° 91.53° 90.00°	Depositor
Resolution (Å)	15.00 – 2.00 15.00 – 2.00	Depositor EDS
% Data completeness (in resolution range)	97.1 (15.00-2.00) 96.8 (15.00-2.00)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.03 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.185 , 0.249 0.183 , 0.250	Depositor DCC
R_{free} test set	2594 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	27.1	Xtriage
Anisotropy	0.240	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 47.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.039 for h,-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6790	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.86% 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: 1UQ, NAP, MN3

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	0.73	1/3331 (0.0%)	0.90	2/4498 (0.0%)
1	B	0.70	2/3257 (0.1%)	0.88	0/4397
All	All	0.71	3/6588 (0.0%)	0.89	2/8895 (0.0%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	335	HIS	CG-ND1	-7.75	1.29	1.38
1	B	335	HIS	CG-ND1	-7.58	1.29	1.38
1	B	155	LYS	C-O	-5.03	1.21	1.25

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	155	LYS	CA-C-N	6.94	127.11	120.31
1	A	155	LYS	C-N-CA	6.94	127.11	120.31

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	3268	0	3316	57	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	3200	0	3245	62	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	48	0	25	8	0
4	A	18	0	11	0	0
4	B	18	0	11	0	0
5	A	129	0	0	6	0
5	B	107	0	0	6	0
All	All	6790	0	6608	115	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 9.

All (115) 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:145:LYS:O	1:B:149:LYS:HE2	1.34	1.23
1:B:152:LYS:H	1:B:152:LYS:HE3	0.92	1.05
1:B:145:LYS:O	1:B:149:LYS:CE	2.10	1.00
1:B:152:LYS:HE3	1:B:152:LYS:N	1.76	0.99
1:B:282:LEU:HD22	1:B:437:TYR:CE2	2.01	0.96
1:B:146:GLU:HA	1:B:149:LYS:NZ	1.82	0.94
1:B:122:TYR:HD1	1:B:122:TYR:C	1.75	0.92
1:B:282:LEU:HD22	1:B:437:TYR:HE2	1.31	0.92
1:B:122:TYR:HD1	1:B:122:TYR:O	1.60	0.85
1:B:122:TYR:C	1:B:122:TYR:CD1	2.50	0.84
1:B:197:ASN:HD22	1:B:224:ASN:HB2	1.40	0.84
1:A:122:TYR:CD2	1:A:147:LEU:HB3	2.14	0.83
1:B:146:GLU:HA	1:B:149:LYS:HZ2	1.41	0.83
1:A:185:GLN:HE21	3:A:502:NAP:H62A	1.24	0.82
1:A:122:TYR:HD1	1:A:122:TYR:C	1.87	0.81
1:B:437:TYR:O	1:B:437:TYR:HD1	1.64	0.81
1:A:122:TYR:C	1:A:122:TYR:CD1	2.60	0.80
1:A:255:ASN:ND2	1:A:258:LYS:HE2	1.97	0.80
1:A:122:TYR:HD2	1:A:147:LEU:HB3	1.46	0.77
1:B:437:TYR:HD1	1:B:437:TYR:C	1.95	0.75
1:B:437:TYR:C	1:B:437:TYR:CD1	2.67	0.73
1:A:284:ASN:HB2	5:A:628:HOH:O	1.88	0.72
1:A:359:ASP:HB3	5:A:615:HOH:O	1.91	0.70
1:B:132:TYR:HD2	1:B:168:ILE:HG12	1.56	0.69
1:B:122:TYR:CD2	1:B:147:LEU:HB3	2.28	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:341:HIS:CE1	5:B:634:HOH:O	2.45	0.69
1:A:99:GLU:HA	1:A:99:GLU:OE1	1.92	0.68
1:A:340:ILE:HD12	1:A:391:PHE:HZ	1.59	0.67
1:B:130:PRO:O	1:B:154:TYR:OH	2.08	0.66
1:B:340:ILE:HD12	1:B:391:PHE:HZ	1.62	0.64
1:B:122:TYR:O	1:B:122:TYR:CD1	2.48	0.64
1:A:185:GLN:NE2	3:A:502:NAP:N6A	2.41	0.63
1:B:146:GLU:HA	1:B:149:LYS:CE	2.29	0.62
1:B:113:TYR:HE2	1:B:165:MET:HG3	1.65	0.61
1:A:150:ASN:N	1:A:150:ASN:HD22	1.98	0.61
1:A:217:LYS:HE3	1:A:458:GLU:O	2.01	0.61
1:B:99:GLU:HA	1:B:99:GLU:OE1	2.02	0.59
1:B:217:LYS:HE3	1:B:458:GLU:O	2.03	0.58
1:A:135:ILE:CD1	1:A:140:VAL:HG23	2.35	0.57
1:B:197:ASN:HA	1:B:225:ALA:HB2	1.86	0.57
1:A:355:MET:HE1	1:B:340:ILE:HD13	1.87	0.57
1:A:185:GLN:HG2	3:A:502:NAP:N6A	2.20	0.56
1:A:282:LEU:HD22	1:A:437:TYR:CE2	2.41	0.56
1:B:217:LYS:CE	1:B:458:GLU:O	2.53	0.56
1:A:250:LYS:HE3	5:A:642:HOH:O	2.04	0.56
1:B:255:ASN:ND2	1:B:258:LYS:HE3	2.21	0.56
1:A:132:TYR:HD2	1:A:168:ILE:HG12	1.70	0.56
1:A:185:GLN:NE2	3:A:502:NAP:H62A	1.98	0.56
1:B:122:TYR:CD2	1:B:147:LEU:HD22	2.41	0.55
1:B:152:LYS:O	1:B:152:LYS:HG2	2.05	0.55
1:B:165:MET:SD	1:B:189:SER:HA	2.48	0.54
1:B:135:ILE:CD1	1:B:140:VAL:HG23	2.37	0.54
1:A:253:GLN:HG3	5:A:642:HOH:O	2.06	0.54
1:A:113:TYR:CE1	3:A:502:NAP:H2A	2.44	0.53
1:B:77:PRO:N	5:B:640:HOH:O	2.43	0.52
1:A:122:TYR:CD2	1:A:147:LEU:HD22	2.45	0.52
1:A:213:PHE:HB2	5:A:625:HOH:O	2.10	0.51
1:A:340:ILE:HD13	1:B:355:MET:HE1	1.92	0.51
1:A:355:MET:HE3	1:B:353:SER:OG	2.09	0.51
1:B:131:GLU:C	1:B:132:TYR:HD1	2.18	0.51
1:A:113:TYR:CZ	3:A:502:NAP:H2A	2.45	0.51
1:A:353:SER:OG	1:B:355:MET:HE3	2.11	0.51
1:B:113:TYR:CE2	1:B:165:MET:HG3	2.44	0.51
1:A:122:TYR:HD1	1:A:122:TYR:O	1.92	0.51
1:A:132:TYR:CE1	1:A:157:ILE:HD12	2.47	0.50
1:A:144:LEU:O	1:A:148:VAL:HG23	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:135:ILE:HD11	1:A:140:VAL:HG23	1.93	0.49
1:B:339:ILE:O	1:B:341:HIS:HD2	1.95	0.49
1:A:217:LYS:CE	1:A:458:GLU:O	2.61	0.48
1:A:335:HIS:HE1	1:A:339:ILE:H	1.61	0.48
1:B:132:TYR:CE1	1:B:157:ILE:HD12	2.48	0.48
1:A:150:ASN:N	1:A:150:ASN:ND2	2.60	0.48
1:A:255:ASN:HD22	1:A:258:LYS:HE2	1.74	0.48
1:A:122:TYR:CD1	1:A:122:TYR:O	2.66	0.47
1:B:122:TYR:HD2	1:B:147:LEU:HB3	1.78	0.47
1:B:312:LYS:O	1:B:316:VAL:HG23	2.14	0.47
1:A:185:GLN:HG2	3:A:502:NAP:H61A	1.78	0.47
1:B:135:ILE:HD11	1:B:140:VAL:HG23	1.95	0.47
1:B:339:ILE:O	1:B:341:HIS:CD2	2.67	0.47
1:B:163:GLU:OE1	1:B:163:GLU:N	2.45	0.47
1:A:165:MET:SD	1:A:189:SER:HA	2.55	0.47
1:A:206:GLU:HG3	1:A:420:ASN:HD21	1.80	0.47
1:A:335:HIS:CE1	1:A:339:ILE:H	2.33	0.47
1:B:228:ILE:HG21	1:B:367:SER:HA	1.95	0.47
1:A:135:ILE:HD12	1:A:140:VAL:HG23	1.98	0.45
1:A:261:ASN:HB3	1:A:347:ILE:HG12	1.98	0.45
1:B:206:GLU:HG3	1:B:420:ASN:HD21	1.81	0.45
1:B:359:ASP:HB3	5:B:704:HOH:O	2.17	0.45
1:B:393:LYS:HG2	5:B:690:HOH:O	2.16	0.45
1:B:89:ILE:HD13	5:B:653:HOH:O	2.16	0.45
1:A:77:PRO:HB2	5:A:720:HOH:O	2.17	0.45
1:B:341:HIS:HE1	5:B:634:HOH:O	1.94	0.44
1:A:429:LEU:HB3	1:A:435:ILE:HG12	1.99	0.44
1:A:191:MET:HE3	1:A:214:PHE:HZ	1.83	0.43
1:B:135:ILE:HD12	1:B:140:VAL:HG23	2.01	0.43
1:B:165:MET:HE2	1:B:165:MET:HB3	1.92	0.43
1:B:266:PHE:HB2	1:B:343:CYS:HB2	2.00	0.43
1:A:163:GLU:OE1	1:A:163:GLU:N	2.43	0.43
1:A:197:ASN:HA	1:A:225:ALA:HB2	2.01	0.42
1:B:113:TYR:HA	1:B:134:CYS:O	2.20	0.42
1:B:264:LYS:HB3	1:B:345:GLU:HB3	2.01	0.42
1:A:138:LYS:HE2	1:A:160:CYS:SG	2.60	0.42
1:A:204:ASN:HA	3:A:502:NAP:O3D	2.19	0.42
1:B:152:LYS:O	1:B:153:ASP:HB3	2.19	0.41
1:B:340:ILE:HG12	1:B:355:MET:HE2	2.02	0.41
1:A:266:PHE:HB2	1:A:343:CYS:HB2	2.01	0.41
1:B:122:TYR:CE2	1:B:147:LEU:HD22	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:132:TYR:HE1	1:A:157:ILE:HD12	1.84	0.41
1:A:131:GLU:C	1:A:132:TYR:HD1	2.28	0.41
1:A:254:ASP:O	1:A:255:ASN:HB2	2.21	0.41
1:B:255:ASN:ND2	1:B:258:LYS:CE	2.83	0.41
1:A:146:GLU:O	1:A:149:LYS:HG3	2.21	0.40
1:A:340:ILE:HD12	1:A:391:PHE:CZ	2.47	0.40
1:B:261:ASN:HB3	1:B:347:ILE:HG12	2.03	0.40
1:A:147:LEU:C	1:A:149:LYS:N	2.76	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	407/427 (95%)	395 (97%)	12 (3%)	0	100	100
1	B	398/427 (93%)	387 (97%)	11 (3%)	0	100	100
All	All	805/854 (94%)	782 (97%)	23 (3%)	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	376/392 (96%)	349 (93%)	27 (7%)	13	10

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	369/392 (94%)	342 (93%)	27 (7%)	13	9
All	All	745/784 (95%)	691 (93%)	54 (7%)	13	10

All (54) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	99	GLU
1	A	108	ASN
1	A	116	LYS
1	A	118	VAL
1	A	122	TYR
1	A	135	ILE
1	A	150	ASN
1	A	152	LYS
1	A	155	LYS
1	A	163	GLU
1	A	165	MET
1	A	170	SER
1	A	210	SER
1	A	213	PHE
1	A	248	LYS
1	A	252	LEU
1	A	253	GLN
1	A	287	SER
1	A	355	MET
1	A	385	GLN
1	A	429	LEU
1	A	455	LYS
1	A	458	GLU
1	A	459	ASN
1	A	464	MET
1	A	472	SER
1	A	477	LYS
1	B	99	GLU
1	B	104	GLU
1	B	108	ASN
1	B	116	LYS
1	B	118	VAL
1	B	122	TYR
1	B	135	ILE
1	B	150	ASN
1	B	151	ILE

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Mol	Chain	Res	Type
1	B	152	LYS
1	B	155	LYS
1	B	163	GLU
1	B	165	MET
1	B	170	SER
1	B	210	SER
1	B	213	PHE
1	B	220	ASN
1	B	252	LEU
1	B	283	LYS
1	B	287	SER
1	B	355	MET
1	B	385	GLN
1	B	429	LEU
1	B	437	TYR
1	B	458	GLU
1	B	459	ASN
1	B	477	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	150	ASN
1	A	185	GLN
1	A	220	ASN
1	A	253	GLN
1	A	255	ASN
1	A	260	ASN
1	A	328	ASN
1	A	335	HIS
1	A	341	HIS
1	A	392	HIS
1	A	420	ASN
1	A	433	ASN
1	B	108	ASN
1	B	150	ASN
1	B	197	ASN
1	B	253	GLN
1	B	255	ASN
1	B	260	ASN
1	B	284	ASN
1	B	328	ASN

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Mol	Chain	Res	Type
1	B	341	HIS
1	B	392	HIS
1	B	420	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](#)

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

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	NAP	A	502	-	50,52,52	2.73	21 (42%)	71,80,80	1.90	12 (16%)
4	1UQ	B	502	2	15,18,18	2.29	4 (26%)	16,25,25	1.05	0
4	1UQ	A	503	2	15,18,18	3.04	6 (40%)	16,25,25	1.13	1 (6%)

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	A	502	-	-	4/35/67/67	0/5/5/5
4	1UQ	B	502	2	-	3/15/19/19	0/1/1/1
4	1UQ	A	503	2	-	3/15/19/19	0/1/1/1

All (31) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	503	1UQ	C6-S7	8.74	1.91	1.81
3	A	502	NAP	C4N-C3N	8.40	1.52	1.39
3	A	502	NAP	C5N-C4N	6.73	1.50	1.38
3	A	502	NAP	O4D-C1D	6.42	1.49	1.40
3	A	502	NAP	C2N-C3N	-6.22	1.29	1.39
4	B	502	1UQ	C6-S7	5.80	1.87	1.81
3	A	502	NAP	PN-O3	5.43	1.65	1.59
4	A	503	1UQ	C6-C2	5.15	1.59	1.51
3	A	502	NAP	PA-O3	4.11	1.63	1.59
4	B	502	1UQ	P9-O12	3.52	1.60	1.54
3	A	502	NAP	C5B-C4B	3.32	1.61	1.51
3	A	502	NAP	C6N-C5N	-3.28	1.31	1.38
3	A	502	NAP	C1B-N9A	3.27	1.55	1.46
4	A	503	1UQ	C2-N3	3.17	1.39	1.35
4	A	503	1UQ	P9-O12	3.00	1.59	1.54
4	B	502	1UQ	P9-O10	-2.98	1.50	1.54
3	A	502	NAP	C3B-C2B	2.93	1.59	1.53
3	A	502	NAP	C4A-N9A	2.85	1.43	1.37
3	A	502	NAP	P2B-O1X	2.74	1.59	1.50
4	B	502	1UQ	C6-C2	2.64	1.55	1.51
3	A	502	NAP	PA-O1A	2.64	1.60	1.50
3	A	502	NAP	C2A-N1A	2.37	1.38	1.33
4	A	503	1UQ	C14-C13	2.33	1.42	1.39
3	A	502	NAP	C6N-N1N	-2.26	1.30	1.35
3	A	502	NAP	C6A-N6A	2.25	1.40	1.34
3	A	502	NAP	C3B-C4B	2.24	1.58	1.53
3	A	502	NAP	C2D-C3D	-2.15	1.47	1.53
3	A	502	NAP	PA-O5B	2.12	1.67	1.59
4	A	503	1UQ	P9-C8	-2.12	1.80	1.83
3	A	502	NAP	P2B-O2B	2.08	1.63	1.59
3	A	502	NAP	O4D-C4D	-2.02	1.40	1.45

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	502	NAP	C5N-C4N-C3N	-9.66	110.87	120.36
3	A	502	NAP	C5A-N7A-C8A	5.18	111.58	103.45
3	A	502	NAP	C4A-C5A-N7A	-4.51	105.42	110.58
3	A	502	NAP	N3A-C2A-N1A	-4.03	122.48	128.58
3	A	502	NAP	C3N-C7N-N7N	3.69	122.28	117.74
3	A	502	NAP	C5A-C4A-N3A	-3.29	122.19	126.72
3	A	502	NAP	C6A-C5A-C4A	2.85	121.07	117.18
3	A	502	NAP	N9A-C8A-N7A	-2.50	110.39	113.94
3	A	502	NAP	N6A-C6A-N1A	2.32	123.55	118.38
3	A	502	NAP	C2A-N1A-C6A	2.29	122.49	118.73
3	A	502	NAP	O2N-PN-O3	2.19	113.20	107.27
3	A	502	NAP	O7N-C7N-C3N	-2.12	117.00	119.60
4	A	503	1UQ	O1-C2-C6	-2.00	117.21	121.19

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	502	NAP	O4D-C1D-N1N-C2N
3	A	502	NAP	C3B-C2B-O2B-P2B
3	A	502	NAP	C1B-C2B-O2B-P2B
3	A	502	NAP	O4D-C1D-N1N-C6N
4	A	503	1UQ	O1-C2-N3-O4
4	B	502	1UQ	O1-C2-N3-O4
4	A	503	1UQ	N3-C2-C6-S7
4	B	502	1UQ	N3-C2-C6-S7
4	A	503	1UQ	O1-C2-C6-S7
4	B	502	1UQ	O1-C2-C6-S7

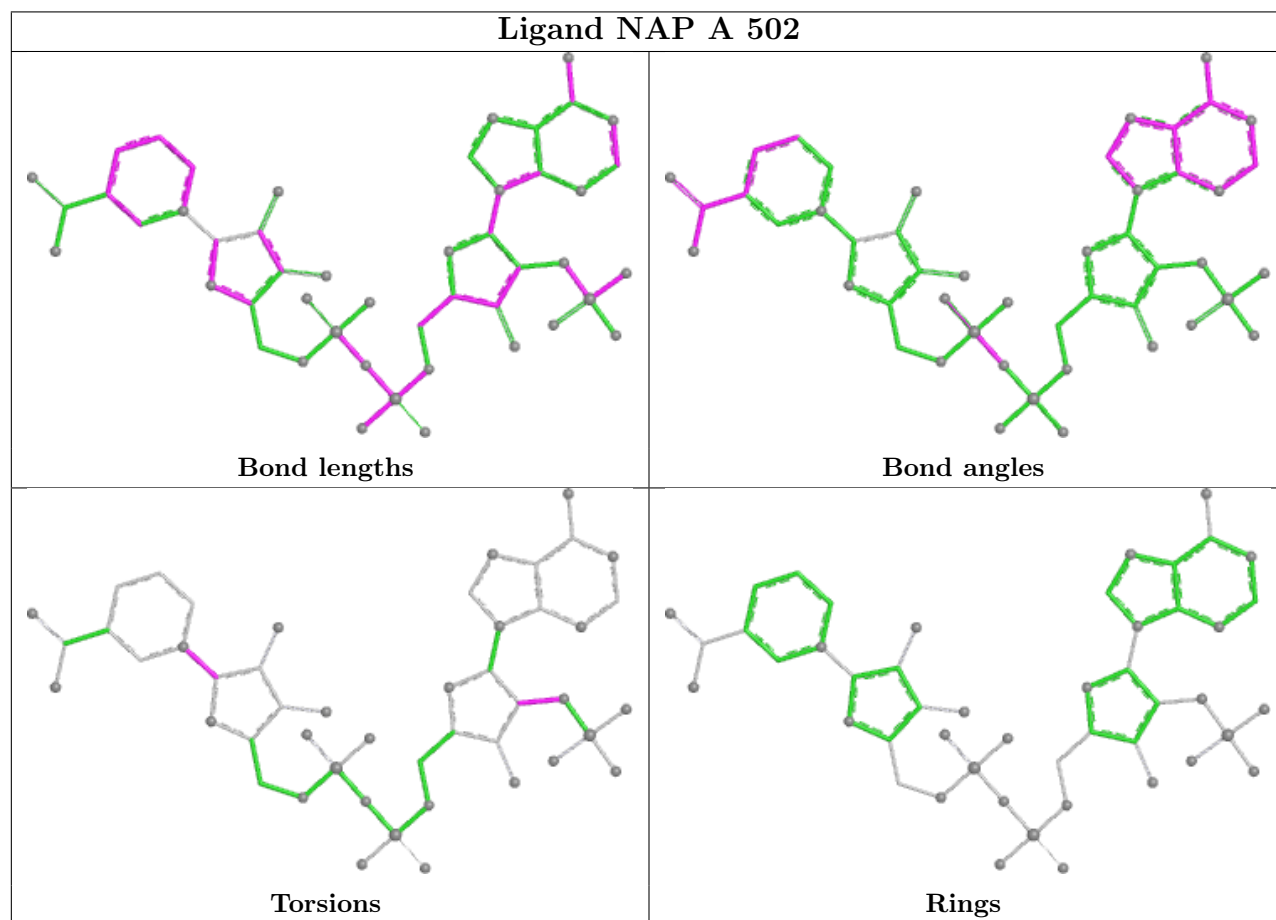
There are no ring outliers.

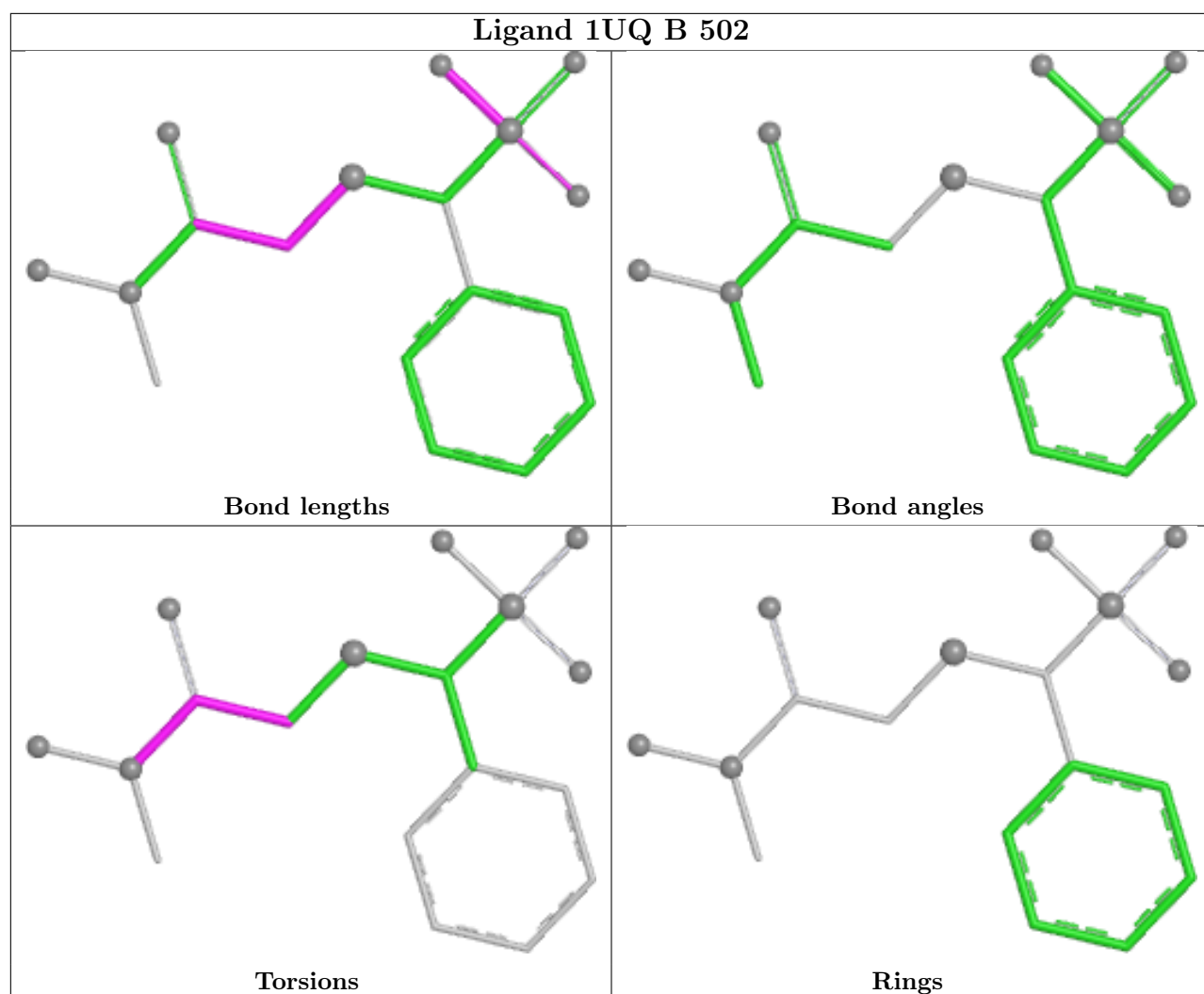
1 monomer is involved in 8 short contacts:

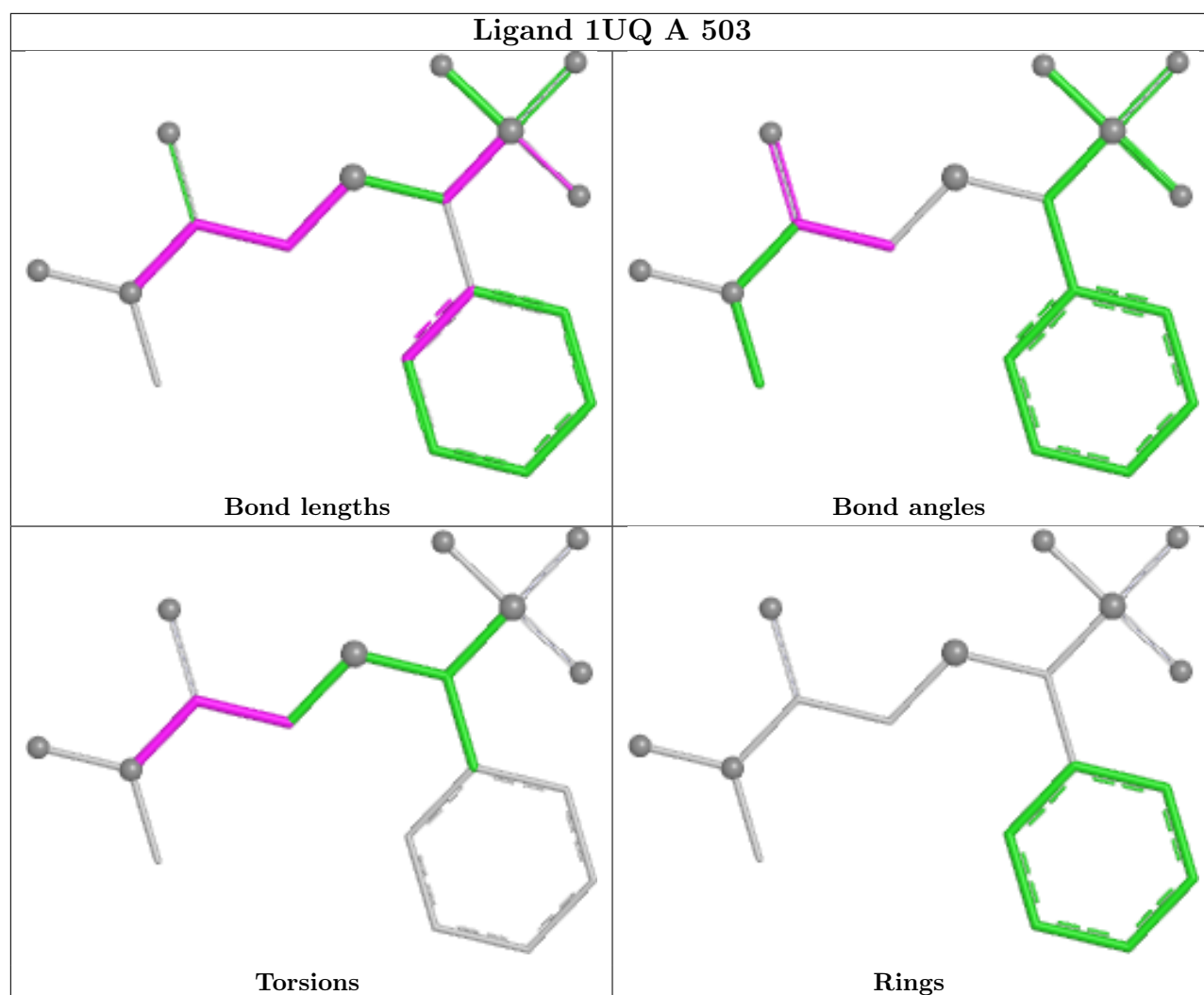
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	502	NAP	8	0

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	409/427 (95%)	-0.28	3 (0%) 84 84	16, 29, 62, 96	0
1	B	402/427 (94%)	-0.05	8 (1%) 65 64	15, 33, 79, 114	0
All	All	811/854 (94%)	-0.17	11 (1%) 73 73	15, 30, 75, 114	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	291	LEU	3.4
1	A	141	TYR	2.5
1	B	114	VAL	2.5
1	B	122	TYR	2.4
1	B	135	ILE	2.3
1	B	158	ILE	2.3
1	B	437	TYR	2.3
1	A	77	PRO	2.3
1	B	283	LYS	2.2
1	B	105	ASN	2.1
1	A	122	TYR	2.1

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](#)

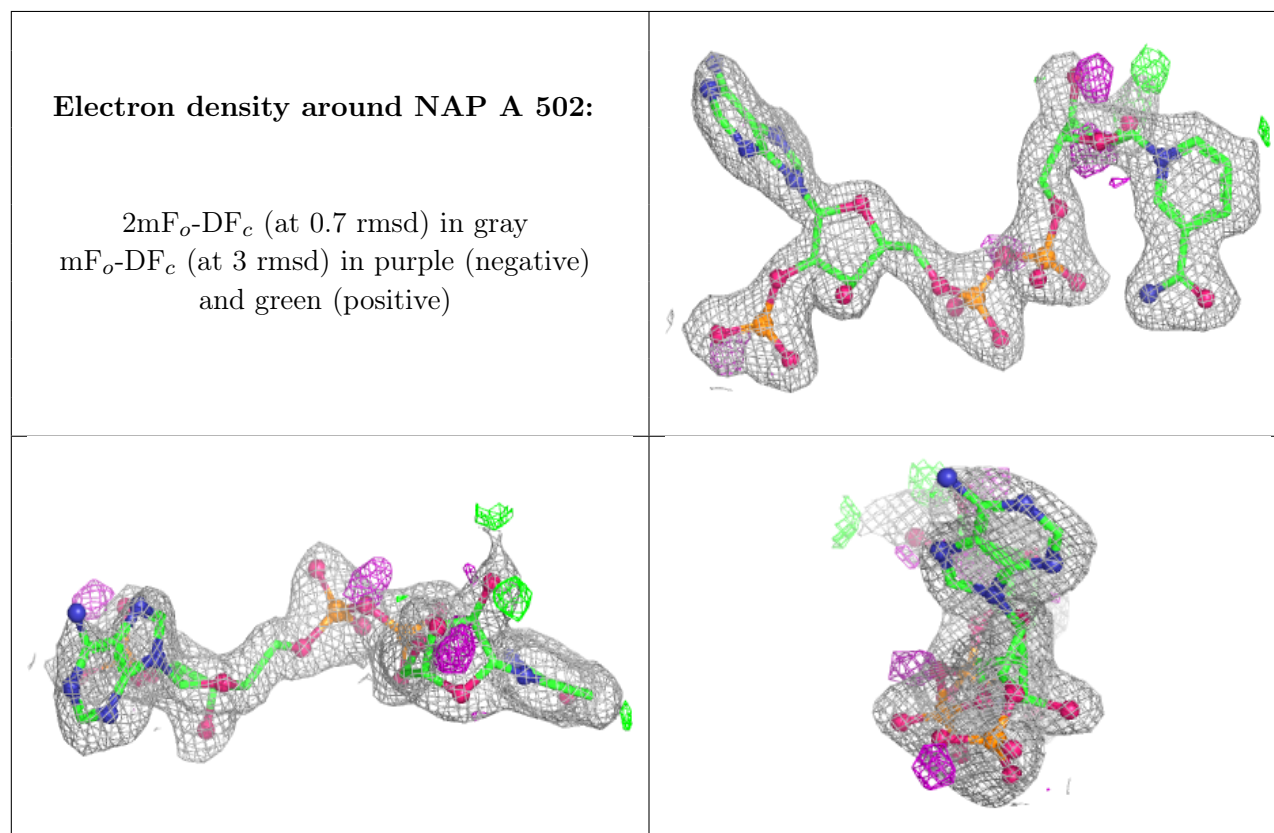
There are no oligosaccharides in this entry.

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.

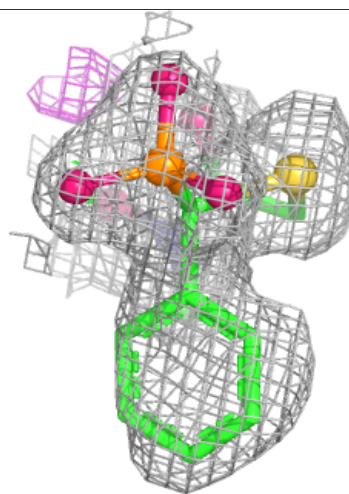
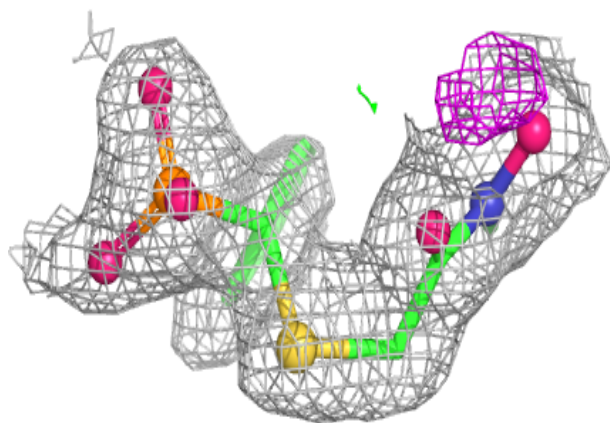
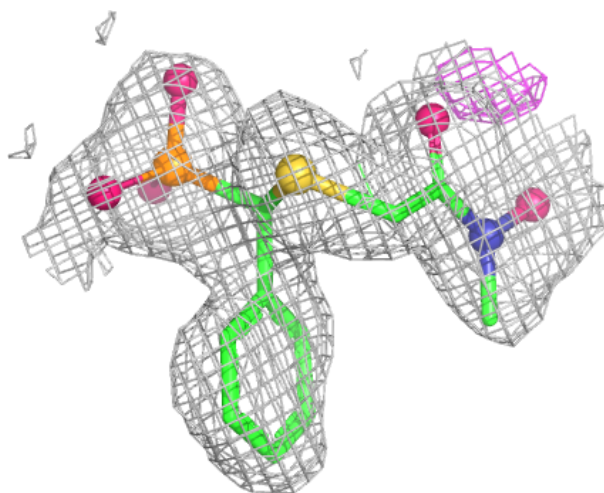
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	NAP	A	502	48/48	0.94	0.08	23,33,45,58	0
4	1UQ	B	502	18/18	0.97	0.06	30,37,44,45	0
4	1UQ	A	503	18/18	0.98	0.05	21,29,32,33	0
2	MN3	A	501	1/1	0.99	0.04	26,26,26,26	0
2	MN3	B	501	1/1	1.00	0.08	33,33,33,33	0

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.



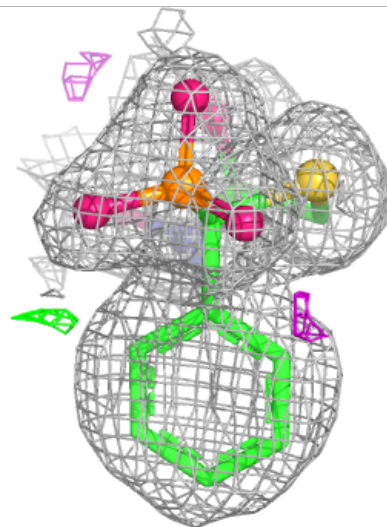
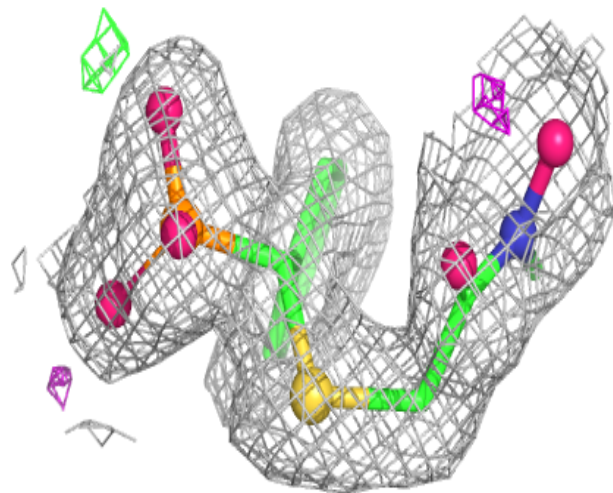
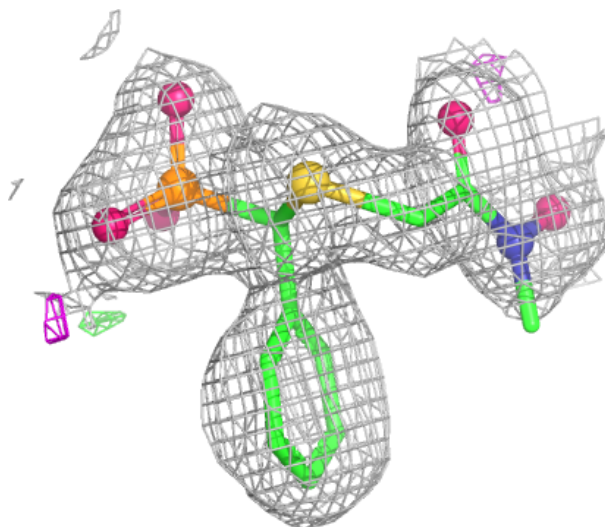
Electron density around 1UQ B 502:

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



Electron density around 1UQ A 503:

$2mF_o-DF_c$ (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.