



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 06:11 AM UTC

PDB ID : 8W49 / pdb_00008w49
Title : mouse isocitrate dehydrogenase with isocitrate and magnesium
Authors : Kang, B.S.; Cho, H.Y.
Deposited on : 2023-08-23
Resolution : 1.77 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

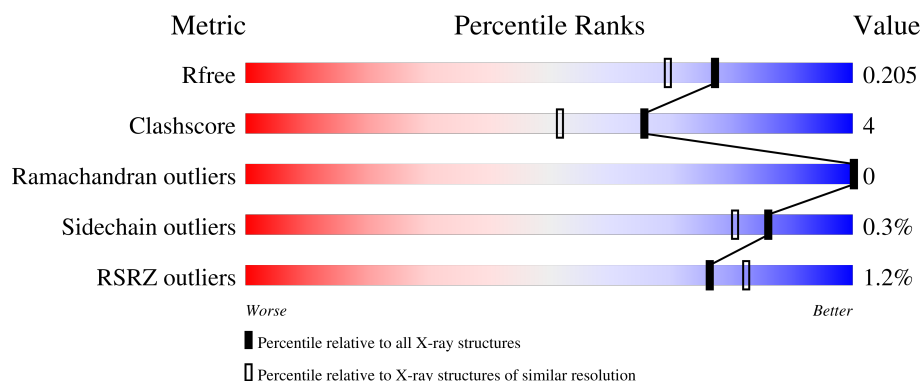
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.77 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	1365 (1.78-1.78)
Clashscore	190562	1395 (1.78-1.78)
Ramachandran outliers	187476	1382 (1.78-1.78)
Sidechain outliers	187428	1382 (1.78-1.78)
RSRZ outliers	180081	1365 (1.78-1.78)

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	1% 92% 7%
1	B	414	92% 7%
1	C	414	1% 92% 7%
1	D	414	2% 89% 11%

2 Entry composition ⓘ

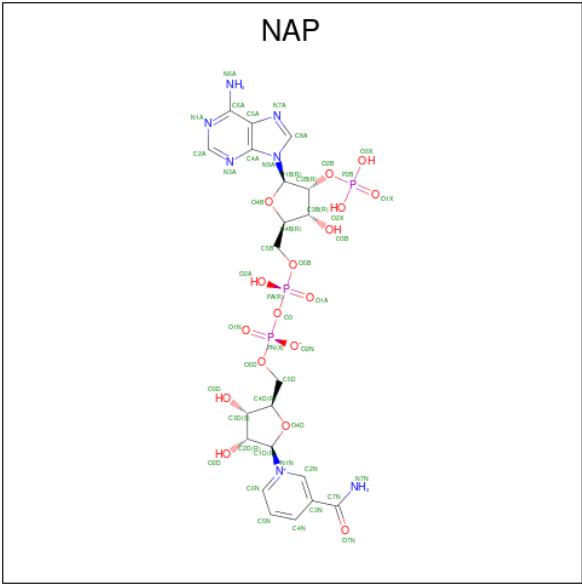
There are 6 unique types of molecules in this entry. The entry contains 14744 atoms, of which 0 are hydrogens and 0 are deuteriums.

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	6	0	0
			3248	2062	549	616	21			
1	B	412	Total	C	N	O	S	2	0	0
			3250	2062	549	618	21			
1	C	411	Total	C	N	O	S	5	1	0
			3248	2063	548	616	21			
1	D	411	Total	C	N	O	S	2	2	0
			3250	2063	548	618	21			

- Molecule 2 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (CCD ID: NAP) (formula: C₂₁H₂₈N₇O₁₇P₃) (labeled as "Ligand of Interest" by depositor).



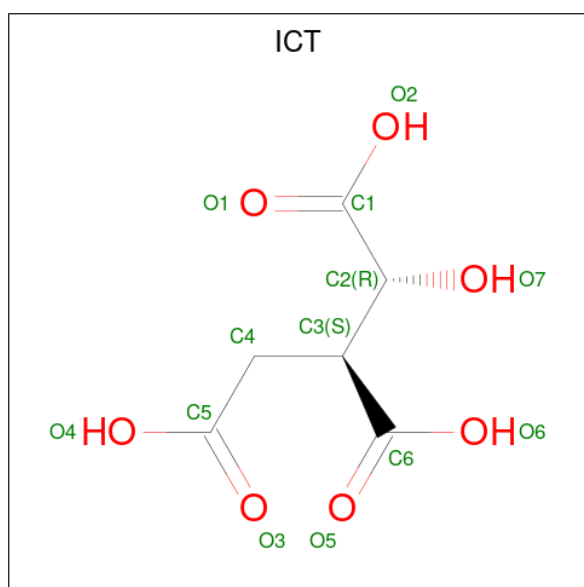
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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	C	1	Total	C	N	O	P	0	0
			48	21	7	17	3		

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

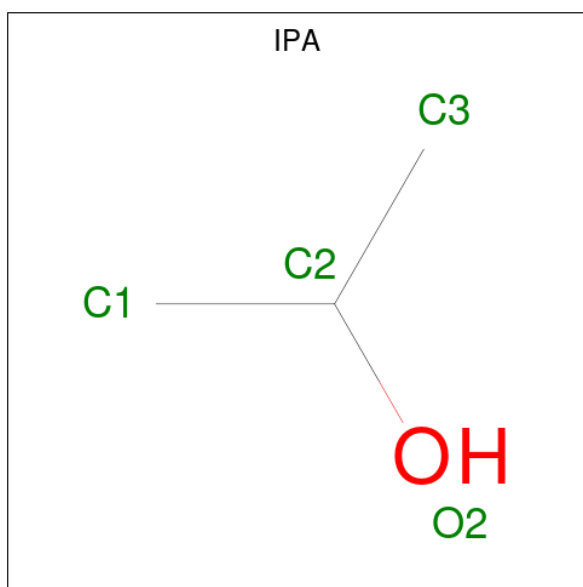
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Mg	0	0
			2	2		
3	C	2	Total	Mg	0	0
			2	2		

- Molecule 4 is ISOCITRIC ACID (CCD ID: ICT) (formula: C₆H₈O₇) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			13	6	7		
4	B	1	Total	C	O	0	0
			13	6	7		
4	C	1	Total	C	O	0	0
			13	6	7		
4	D	1	Total	C	O	0	0
			13	6	7		

- Molecule 5 is ISOPROPYL ALCOHOL (CCD ID: IPA) (formula: C₃H₈O).



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

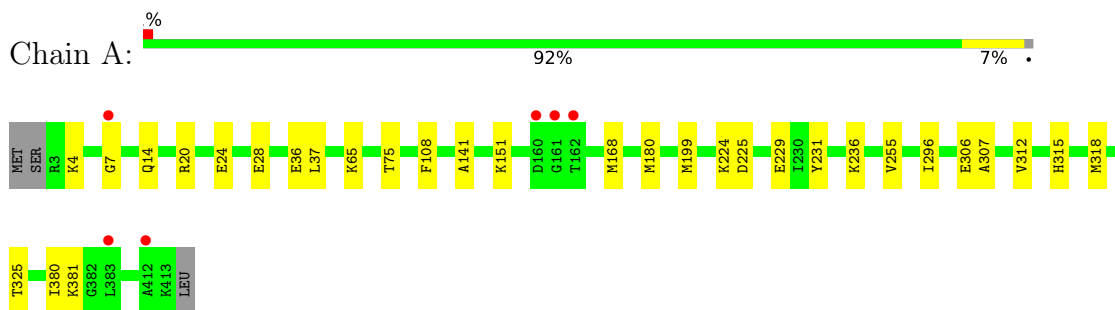
- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	372	Total O 372 372	0	0
6	B	460	Total O 460 460	0	0
6	C	350	Total O 350 350	0	0
6	D	338	Total O 338 338	0	0

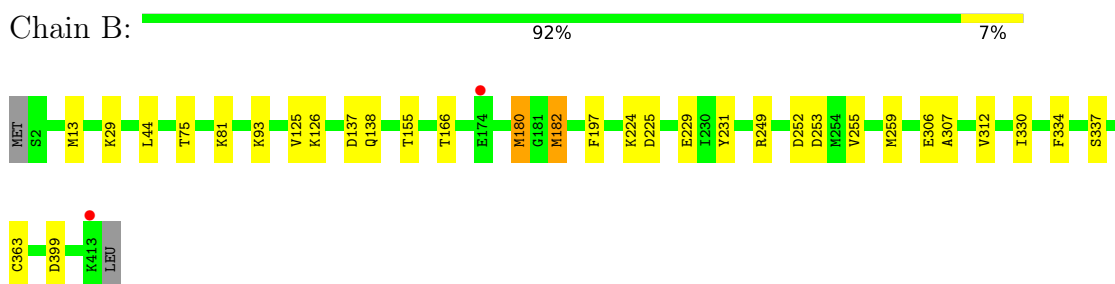
3 Residue-property plots

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.

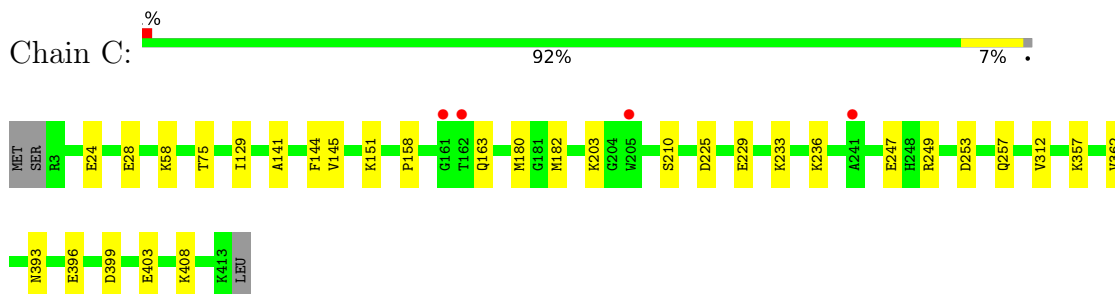
- Molecule 1: Isocitrate dehydrogenase [NADP] cytoplasmic



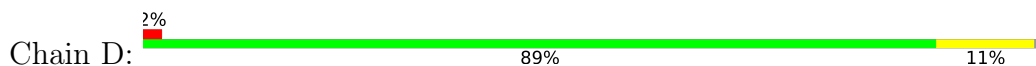
- Molecule 1: Isocitrate dehydrogenase [NADP] cytoplasmic

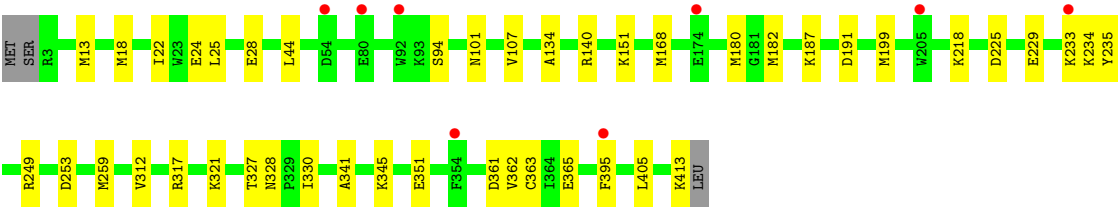


- Molecule 1: Isocitrate dehydrogenase [NADP] cytoplasmic



- Molecule 1: Isocitrate dehydrogenase [NADP] cytoplasmic





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	98.63Å 93.45Å 111.28Å 90.00° 113.90° 90.00°	Depositor
Resolution (Å)	33.90 – 1.77 33.90 – 1.77	Depositor EDS
% Data completeness (in resolution range)	96.0 (33.90-1.77) 96.2 (33.90-1.77)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.43 (at 1.77Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.170 , 0.205 0.170 , 0.205	Depositor DCC
R_{free} test set	1946 reflections (1.08%)	wwPDB-VP
Wilson B-factor (Å ²)	17.3	Xtriage
Anisotropy	0.038	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 41.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.017 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	14744	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 29.93 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.4249e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: NAP, MG, ICT, IPA

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	0.35	0/3313	0.55	0/4464
1	B	0.37	0/3315	0.60	3/4468 (0.1%)
1	C	0.32	0/3316	0.56	0/4470
1	D	0.32	0/3321	0.54	0/4476
All	All	0.34	0/13265	0.56	3/17878 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	180	MET	CG-SD-CE	-6.23	87.20	100.90
1	B	137	ASP	CB-CA-C	-6.21	109.40	116.54
1	B	182	MET	CG-SD-CE	5.76	113.58	100.90

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	3248	0	3223	18	0
1	B	3250	0	3217	23	0
1	C	3248	0	3221	25	0
1	D	3250	0	3222	32	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	48	0	25	3	0
2	B	48	0	25	2	0
2	C	48	0	25	3	0
3	A	2	0	0	0	0
3	C	2	0	0	0	0
4	A	13	0	4	2	0
4	B	13	0	4	1	0
4	C	13	0	4	2	0
4	D	13	0	4	1	0
5	A	8	0	16	1	0
5	B	8	0	16	3	0
5	C	4	0	8	0	0
5	D	8	0	16	1	0
6	A	372	0	0	0	0
6	B	460	0	0	2	0
6	C	350	0	0	2	0
6	D	338	0	0	2	0
All	All	14744	0	13030	99	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 4.

All (99) 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:C:144:PHE:HA	1:D:218:LYS:HE3	1.43	1.01
1:B:180:MET:HE1	1:B:182:MET:HE2	1.48	0.96
1:C:180:MET:HE2	1:C:182:MET:HE2	1.48	0.95
1:C:180:MET:CE	1:C:182:MET:HE2	2.09	0.82
1:D:180:MET:HE2	1:D:182:MET:HE2	1.63	0.81
1:D:187:LYS:NZ	1:D:191:ASP:OD1	2.16	0.78
2:C:501:NAP:C4N	4:C:504:ICT:H2	2.18	0.73
1:D:259:MET:HE2	5:D:503:IPA:H12	1.71	0.71
1:A:315:HIS:HA	1:A:318:MET:HE2	1.73	0.70
1:B:180:MET:HE1	1:B:182:MET:CE	2.19	0.70
2:A:501:NAP:C4N	4:A:504:ICT:H2	2.22	0.69
1:B:138:GLN:HG3	1:B:182:MET:HE1	1.75	0.68
2:B:501:NAP:C4N	4:B:502:ICT:H2	2.25	0.67
1:D:101:ASN:OD1	1:D:140:ARG:NH1	2.28	0.66
1:D:24:GLU:O	1:D:28:GLU:HG2	1.96	0.66
1:C:24:GLU:O	1:C:28:GLU:HG2	1.97	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:253:ASP:O	1:C:257:GLN:HG3	1.98	0.64
1:C:180:MET:HE2	1:C:182:MET:CE	2.26	0.63
1:B:225:ASP:O	1:B:229:GLU:HG2	1.99	0.62
1:B:259:MET:HE2	5:B:504:IPA:H11	1.80	0.61
1:B:29:LYS:NZ	1:B:399:ASP:OD1	2.32	0.60
1:A:7:GLY:HA3	1:A:37:LEU:HD23	1.86	0.58
1:A:306:GLU:HG3	1:A:307:ALA:O	2.04	0.58
1:D:341:ALA:O	1:D:345:LYS:HG2	2.04	0.57
1:B:306:GLU:HG3	1:B:307:ALA:O	2.04	0.57
1:C:182:MET:HE3	1:D:182:MET:HE3	1.85	0.56
1:C:225:ASP:O	1:C:229:GLU:HG3	2.07	0.55
1:D:229:GLU:HA	1:D:233:LYS:HZ3	1.72	0.55
1:C:58:LYS:HE3	6:C:610:HOH:O	2.08	0.54
1:D:187:LYS:HZ1	1:D:191:ASP:CG	2.11	0.53
1:C:362:VAL:HG23	1:C:408:LYS:HD2	1.90	0.53
1:C:145:VAL:H	1:D:218:LYS:NZ	2.07	0.52
1:C:393:ASN:OD1	1:C:396:GLU:HG3	2.09	0.52
1:B:255:VAL:O	5:B:504:IPA:H12	2.11	0.51
1:C:247:GLU:OE1	1:C:249:ARG:NH2	2.41	0.51
1:D:151:LYS:HD3	1:D:168:MET:HE1	1.93	0.51
1:A:141:ALA:HB1	1:A:180:MET:HE3	1.93	0.50
1:A:75:THR:O	2:A:501:NAP:H2N	2.12	0.50
1:C:210:SER:HA	1:C:249:ARG:O	2.12	0.50
1:B:81:LYS:N	1:B:81:LYS:HD3	2.27	0.49
1:B:75:THR:O	2:B:501:NAP:H2N	2.12	0.49
1:A:224:LYS:HZ2	1:B:81:LYS:HE3	1.77	0.49
1:A:380:ILE:HG22	1:A:381:LYS:HD3	1.95	0.49
1:A:225:ASP:O	1:A:229:GLU:HG3	2.12	0.49
1:B:180:MET:CE	1:B:182:MET:CE	2.88	0.49
2:C:501:NAP:H4N	4:C:504:ICT:H2	1.95	0.48
1:A:318:MET:HE3	1:A:325:THR:HG22	1.95	0.48
1:B:224:LYS:NZ	6:B:608:HOH:O	2.44	0.48
1:C:144:PHE:CA	1:D:218:LYS:HE3	2.29	0.47
1:C:129:ILE:CD1	1:C:203:LYS:HE2	2.45	0.47
1:A:65:LYS:O	1:A:65:LYS:HD2	2.13	0.47
2:A:501:NAP:H4N	4:A:504:ICT:H2	1.96	0.47
1:B:81:LYS:H	1:B:81:LYS:HE2	1.79	0.47
1:A:14:GLN:HG3	1:A:20:ARG:NH1	2.29	0.47
5:B:503:IPA:H12	6:B:1004:HOH:O	2.14	0.47
1:D:413:LYS:HB3	1:D:413:LYS:HE3	1.63	0.47
1:C:158:PRO:HD2	1:C:163:GLN:O	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:231:TYR:CE1	1:A:236:LYS:HG3	2.51	0.46
1:D:107:VAL:HG23	1:D:134:ALA:HB2	1.98	0.46
1:B:330:ILE:HD12	1:B:363:CYS:HB3	1.97	0.45
1:A:151:LYS:HG2	1:A:168:MET:HE1	1.99	0.45
1:D:317:ARG:O	1:D:321:LYS:HG3	2.17	0.45
1:D:361:ASP:O	1:D:365:GLU:HG3	2.17	0.45
1:A:4:LYS:HD2	1:A:36:GLU:HG3	1.99	0.44
1:C:236:LYS:HB3	1:C:236:LYS:HE3	1.66	0.44
1:B:197:PHE:CZ	1:B:231:TYR:HB2	2.53	0.44
1:C:233:LYS:HE3	1:C:233:LYS:HB3	1.55	0.43
1:D:18:MET:HE2	6:D:602:HOH:O	2.17	0.43
1:D:199:MET:HE2	1:D:199:MET:HB2	1.88	0.43
1:C:141:ALA:HB1	1:C:180:MET:HE3	2.01	0.43
1:D:22:ILE:HD11	1:D:327:THR:HB	2.01	0.43
1:B:13:MET:HB3	1:B:44:LEU:HG	2.01	0.43
1:B:93:LYS:HD2	1:B:93:LYS:HA	1.85	0.43
1:B:334:PHE:HA	1:B:337:SER:OG	2.19	0.42
1:C:75:THR:O	2:C:501:NAP:H2N	2.19	0.42
1:D:94:SER:OG	4:D:501:ICT:O3	2.32	0.42
1:D:225:ASP:O	1:D:229:GLU:HG3	2.18	0.42
1:B:249:ARG:HD3	1:B:253:ASP:OD2	2.19	0.42
1:D:330:ILE:HD12	1:D:363:CYS:HB3	2.01	0.42
1:A:108:PHE:CD2	1:A:296:ILE:HD12	2.55	0.42
1:D:229:GLU:HA	1:D:233:LYS:NZ	2.35	0.42
1:D:25:LEU:HD11	1:D:395:PHE:CD1	2.54	0.42
1:C:399:ASP:O	1:C:403:GLU:HG3	2.20	0.42
1:D:13:MET:HB3	1:D:44:LEU:HG	2.01	0.42
1:D:362:VAL:HG21	1:D:405:LEU:HA	2.01	0.42
1:A:255:VAL:HB	5:A:506:IPA:H31	2.02	0.42
1:C:151:LYS:HB2	1:C:151:LYS:HE2	1.87	0.42
1:A:24:GLU:O	1:A:28:GLU:HG2	2.20	0.41
1:D:187:LYS:NZ	6:D:606:HOH:O	2.52	0.41
1:C:357:LYS:NZ	6:C:624:HOH:O	2.52	0.41
1:D:328:ASN:OD1	1:D:330:ILE:HG12	2.21	0.41
1:D:351:GLU:OE2	1:D:351:GLU:N	2.42	0.41
1:B:155:THR:HG22	1:B:166:THR:HG23	2.03	0.41
1:D:234:LYS:HD3	1:D:235:TYR:CZ	2.55	0.41
1:D:249:ARG:HD3	1:D:253:ASP:OD2	2.21	0.40
1:B:252:ASP:OD1	1:B:252:ASP:N	2.54	0.40
1:C:393:ASN:CG	1:C:396:GLU:HG3	2.47	0.40
1:A:108:PHE:CE1	1:A:199:MET:HG2	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:125:VAL:HG23	1:B:126:LYS:HG3	2.03	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	409/414 (99%)	396 (97%)	13 (3%)	0	100	100
1	B	410/414 (99%)	397 (97%)	13 (3%)	0	100	100
1	C	410/414 (99%)	401 (98%)	9 (2%)	0	100	100
1	D	411/414 (99%)	401 (98%)	10 (2%)	0	100	100
All	All	1640/1656 (99%)	1595 (97%)	45 (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	346/350 (99%)	345 (100%)	1 (0%)	86	80
1	B	346/350 (99%)	345 (100%)	1 (0%)	86	80
1	C	346/350 (99%)	345 (100%)	1 (0%)	86	80
1	D	347/350 (99%)	346 (100%)	1 (0%)	86	80

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	1385/1400 (99%)	1381 (100%)	4 (0%)	86 80

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	312	VAL
1	B	312	VAL
1	C	312	VAL
1	D	312	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	68	ASN
1	A	228	GLN
1	A	242	GLN
1	A	283	GLN
1	B	53	ASN
1	B	68	ASN
1	C	228	GLN
1	C	238	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](#)

Of 18 ligands modelled in this entry, 4 are monoatomic - leaving 14 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	$\# Z > 2$
5	IPA	D	503	-	3,3,3	0.45	0	3,3,3	0.46	0
5	IPA	D	502	-	3,3,3	0.61	0	3,3,3	0.26	0
4	ICT	D	501	3	12,12,12	1.21	0	13,16,16	1.74	4 (30%)
5	IPA	B	504	-	3,3,3	0.46	0	3,3,3	0.52	0
5	IPA	A	506	-	3,3,3	0.51	0	3,3,3	0.63	0
2	NAP	A	501	-	50,52,52	0.58	0	71,80,80	0.74	2 (2%)
4	ICT	A	504	3	12,12,12	1.31	1 (8%)	13,16,16	1.49	2 (15%)
5	IPA	B	503	-	3,3,3	0.49	0	3,3,3	0.36	0
2	NAP	B	501	-	50,52,52	0.66	2 (4%)	71,80,80	0.68	2 (2%)
4	ICT	C	504	3	12,12,12	1.34	1 (8%)	13,16,16	1.77	4 (30%)
5	IPA	A	505	-	3,3,3	0.59	0	3,3,3	0.31	0
5	IPA	C	505	-	3,3,3	0.47	0	3,3,3	0.43	0
4	ICT	B	502	3	12,12,12	1.33	1 (8%)	13,16,16	1.88	6 (46%)
2	NAP	C	501	-	50,52,52	0.61	1 (2%)	71,80,80	0.76	1 (1%)

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	ICT	D	501	3	-	6/16/16/16	-
2	NAP	A	501	-	-	4/35/67/67	0/5/5/5
4	ICT	A	504	3	-	6/16/16/16	-
2	NAP	B	501	-	-	4/35/67/67	0/5/5/5
4	ICT	C	504	3	-	5/16/16/16	-
4	ICT	B	502	3	-	4/16/16/16	-
2	NAP	C	501	-	-	5/35/67/67	0/5/5/5

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	504	ICT	C3-C2	-2.65	1.51	1.54
4	B	502	ICT	C3-C2	-2.61	1.51	1.54
2	B	501	NAP	C2N-N1N	2.08	1.37	1.35
4	A	504	ICT	C3-C2	-2.05	1.51	1.54
2	C	501	NAP	P2B-O2B	2.04	1.63	1.59
2	B	501	NAP	P2B-O2B	2.03	1.63	1.59

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	501	ICT	O6-C6-C3	3.06	121.87	113.93
4	B	502	ICT	C3-C4-C5	-2.86	108.60	113.95
2	C	501	NAP	C6N-N1N-C2N	-2.85	119.45	121.88
4	C	504	ICT	O7-C2-C1	2.78	116.64	110.69
4	C	504	ICT	O2-C1-C2	2.60	120.53	113.31
4	B	502	ICT	O6-C6-C3	2.51	120.45	113.93
4	D	501	ICT	O6-C6-O5	-2.51	118.39	124.08
4	D	501	ICT	O2-C1-C2	2.45	120.12	113.31
4	C	504	ICT	C3-C4-C5	-2.41	109.45	113.95
2	A	501	NAP	C6N-N1N-C2N	-2.34	119.89	121.88
4	B	502	ICT	C4-C3-C6	-2.30	107.59	111.25
4	B	502	ICT	O4-C5-C4	2.26	121.03	114.00
4	B	502	ICT	O7-C2-C1	2.23	115.45	110.69
2	B	501	NAP	C6N-N1N-C2N	-2.21	120.00	121.88
4	D	501	ICT	C3-C4-C5	-2.15	109.92	113.95
2	B	501	NAP	P2B-O2B-C2B	-2.14	117.72	123.43
4	A	504	ICT	O2-C1-C2	2.08	119.09	113.31
4	C	504	ICT	O6-C6-C3	2.04	119.24	113.93
4	A	504	ICT	C4-C3-C6	-2.04	108.00	111.25
2	A	501	NAP	O3B-C3B-C2B	2.04	116.89	111.19
4	B	502	ICT	O5-C6-C3	-2.03	118.21	123.01

There are no chirality outliers.

All (34) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	NAP	O4D-C1D-N1N-C2N
2	A	501	NAP	O4D-C1D-N1N-C6N
2	A	501	NAP	C2D-C1D-N1N-C2N
2	B	501	NAP	O4D-C1D-N1N-C2N
2	B	501	NAP	O4D-C1D-N1N-C6N
2	B	501	NAP	C2D-C1D-N1N-C2N

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Mol	Chain	Res	Type	Atoms
2	C	501	NAP	O4D-C1D-N1N-C2N
2	C	501	NAP	O4D-C1D-N1N-C6N
2	C	501	NAP	C2D-C1D-N1N-C2N
4	A	504	ICT	O1-C1-C2-C3
4	A	504	ICT	O2-C1-C2-C3
4	B	502	ICT	O1-C1-C2-C3
4	B	502	ICT	O2-C1-C2-C3
4	B	502	ICT	O1-C1-C2-O7
4	B	502	ICT	O2-C1-C2-O7
4	A	504	ICT	O1-C1-C2-O7
4	A	504	ICT	O2-C1-C2-O7
4	A	504	ICT	C4-C3-C6-O5
4	A	504	ICT	C4-C3-C6-O6
4	D	501	ICT	C4-C3-C6-O5
4	D	501	ICT	C4-C3-C6-O6
4	D	501	ICT	O2-C1-C2-C3
4	C	504	ICT	C1-C2-C3-C6
2	C	501	NAP	C5B-O5B-PA-O1A
4	D	501	ICT	O1-C1-C2-C3
2	A	501	NAP	C2D-C1D-N1N-C6N
2	B	501	NAP	C2D-C1D-N1N-C6N
2	C	501	NAP	C2D-C1D-N1N-C6N
4	D	501	ICT	O2-C1-C2-O7
4	D	501	ICT	O1-C1-C2-O7
4	C	504	ICT	C3-C4-C5-O4
4	C	504	ICT	C3-C4-C5-O3
4	C	504	ICT	O2-C1-C2-C3
4	C	504	ICT	O1-C1-C2-C3

There are no ring outliers.

11 monomers are involved in 14 short contacts:

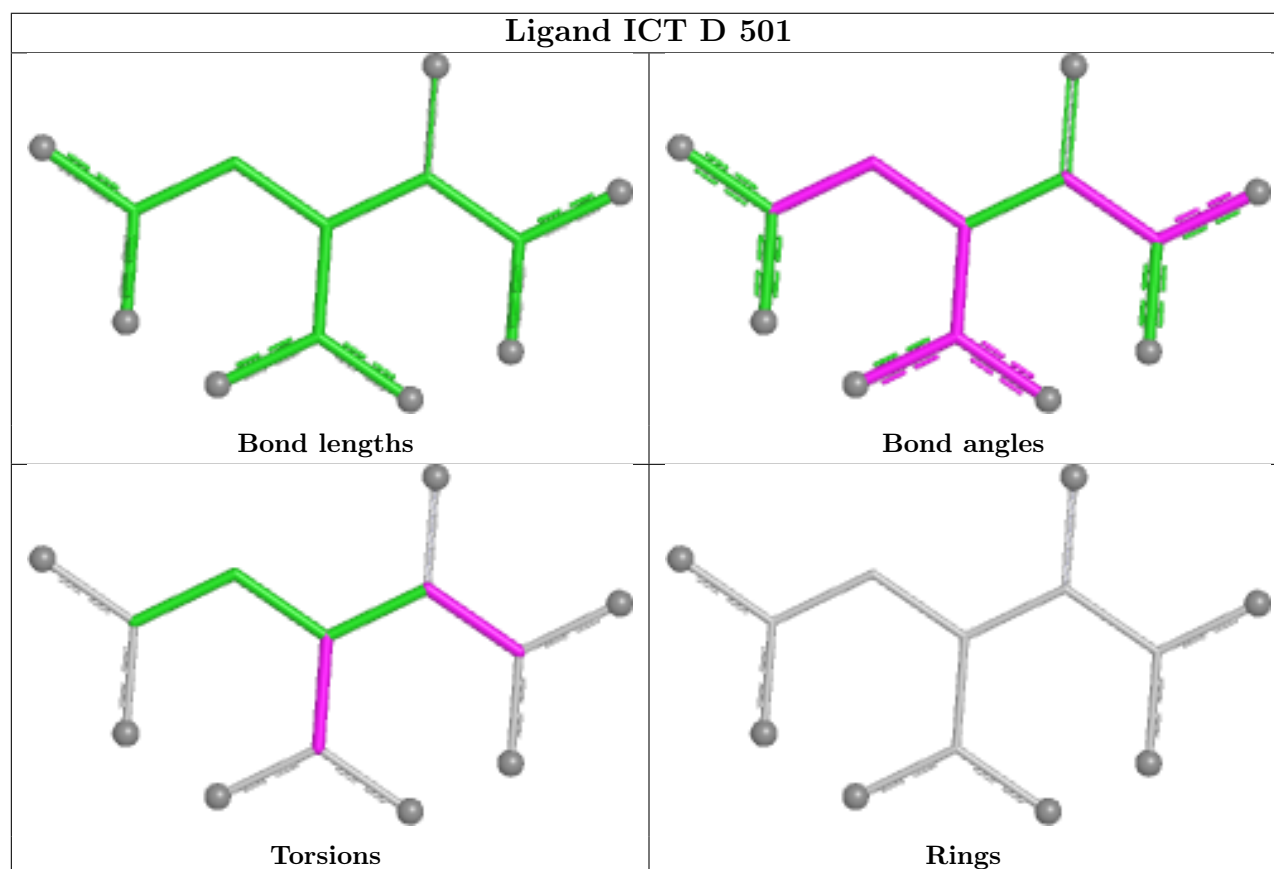
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	D	503	IPA	1	0
4	D	501	ICT	1	0
5	B	504	IPA	2	0
5	A	506	IPA	1	0
2	A	501	NAP	3	0
4	A	504	ICT	2	0
5	B	503	IPA	1	0
2	B	501	NAP	2	0
4	C	504	ICT	2	0

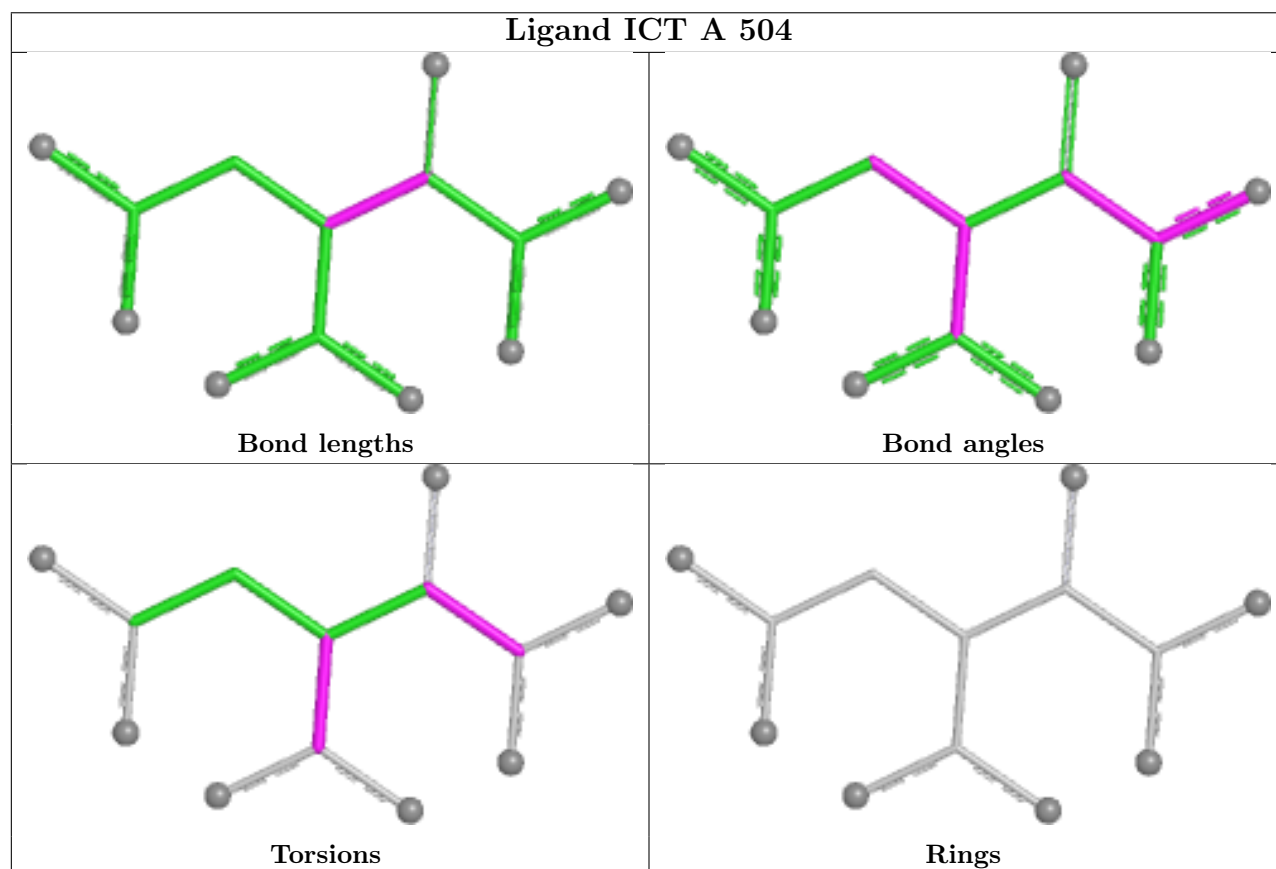
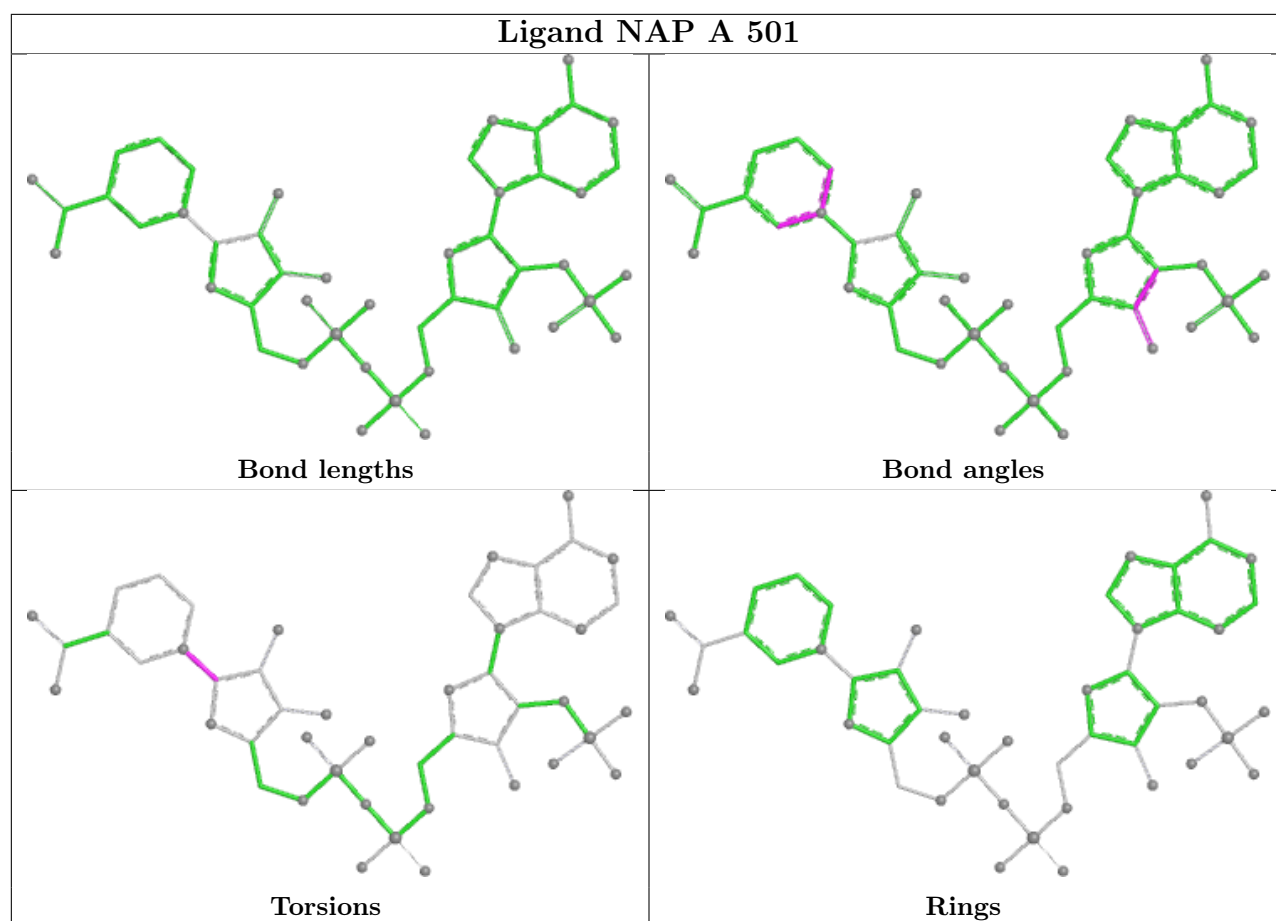
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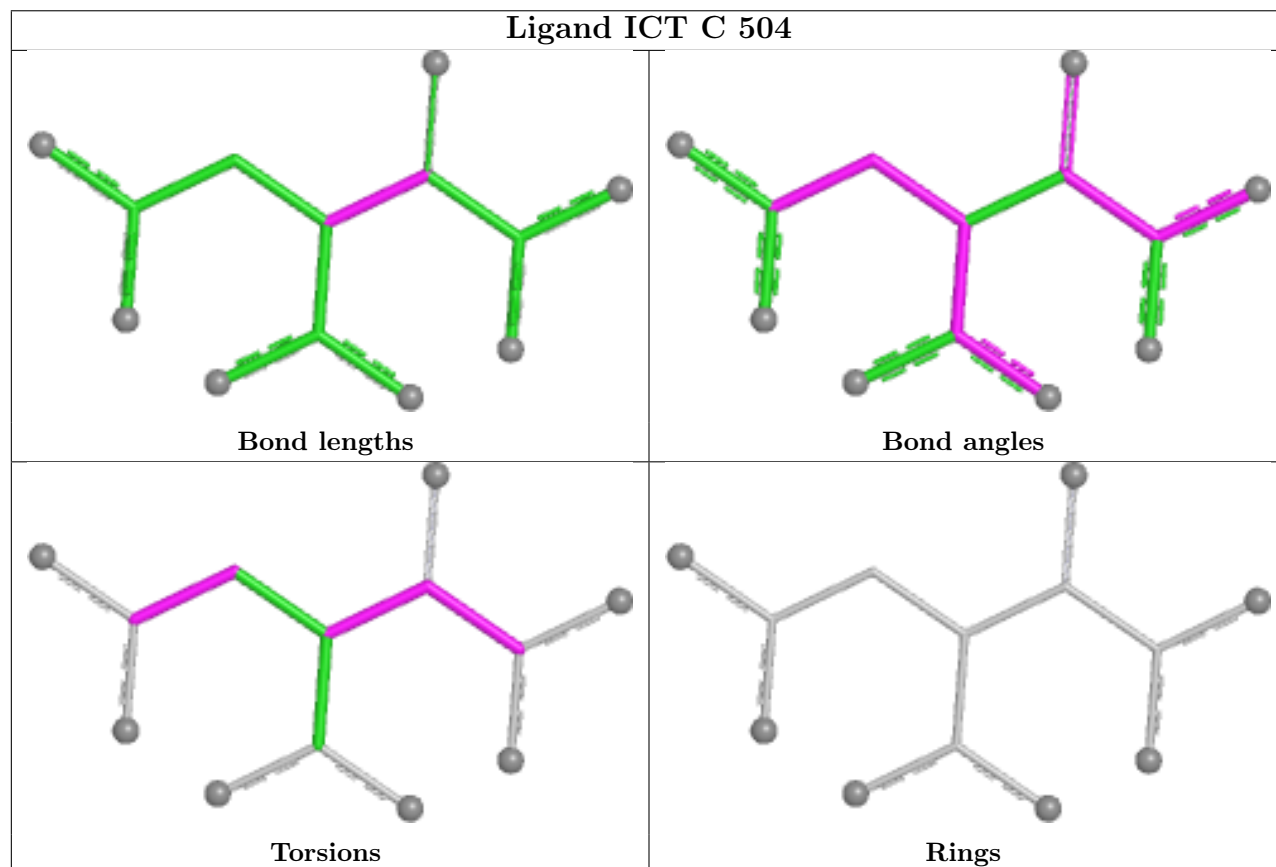
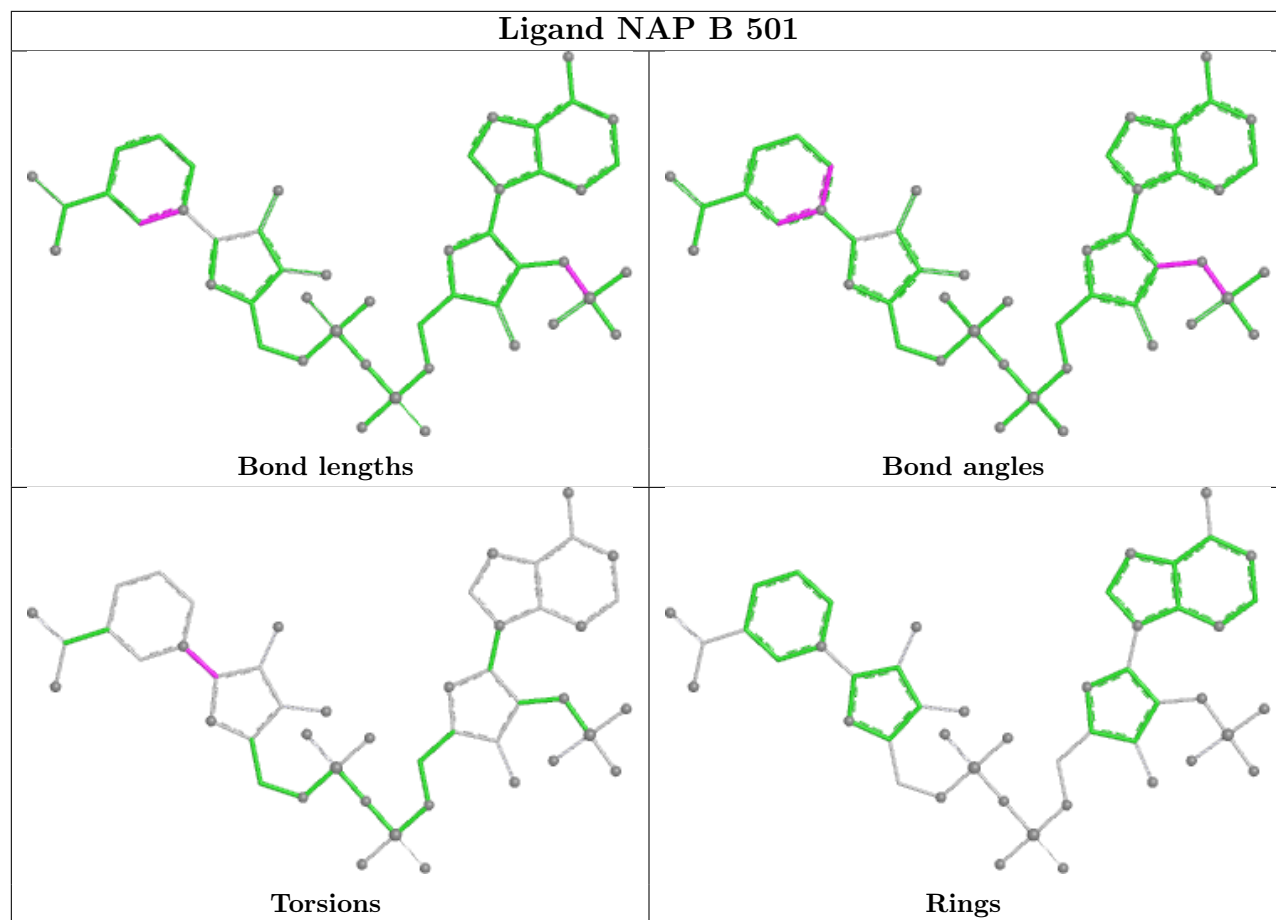
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	502	ICT	1	0
2	C	501	NAP	3	0

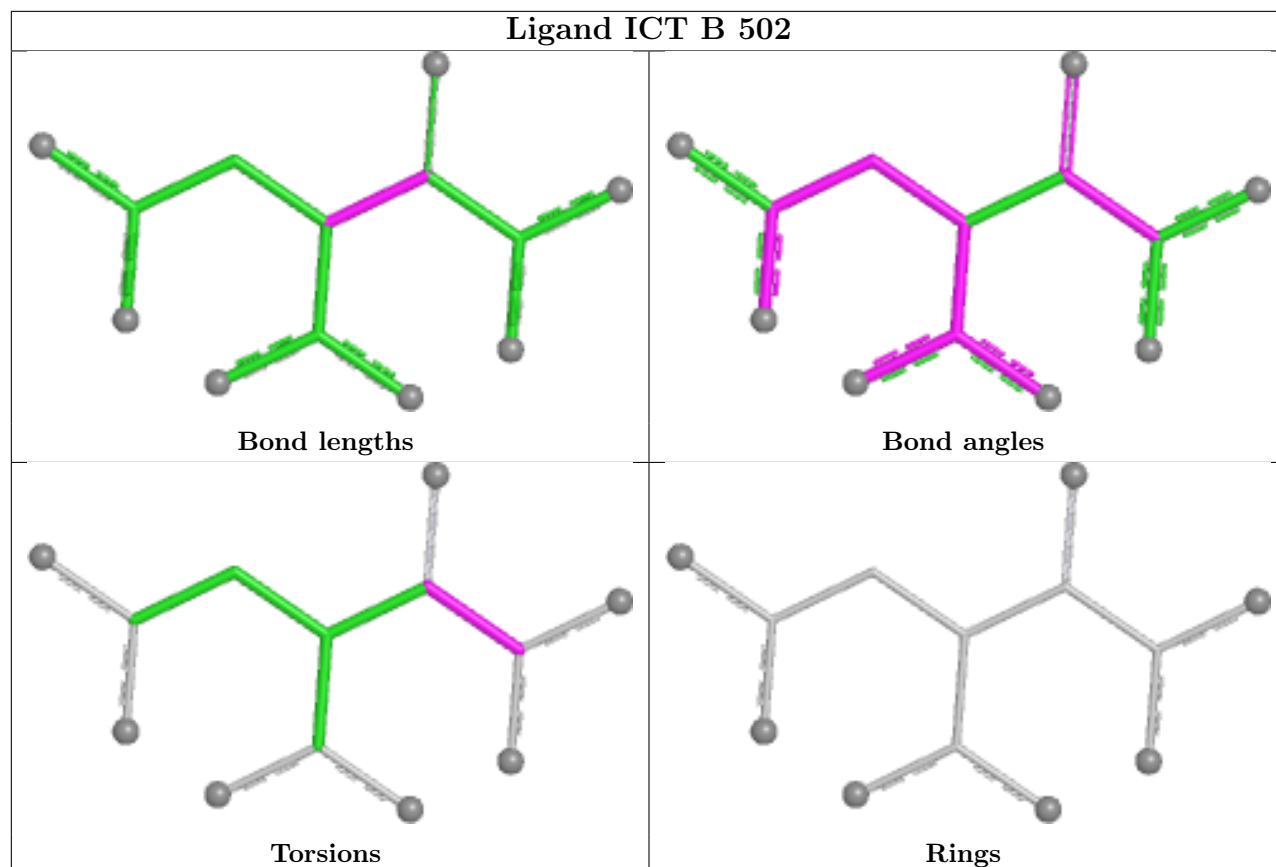
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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



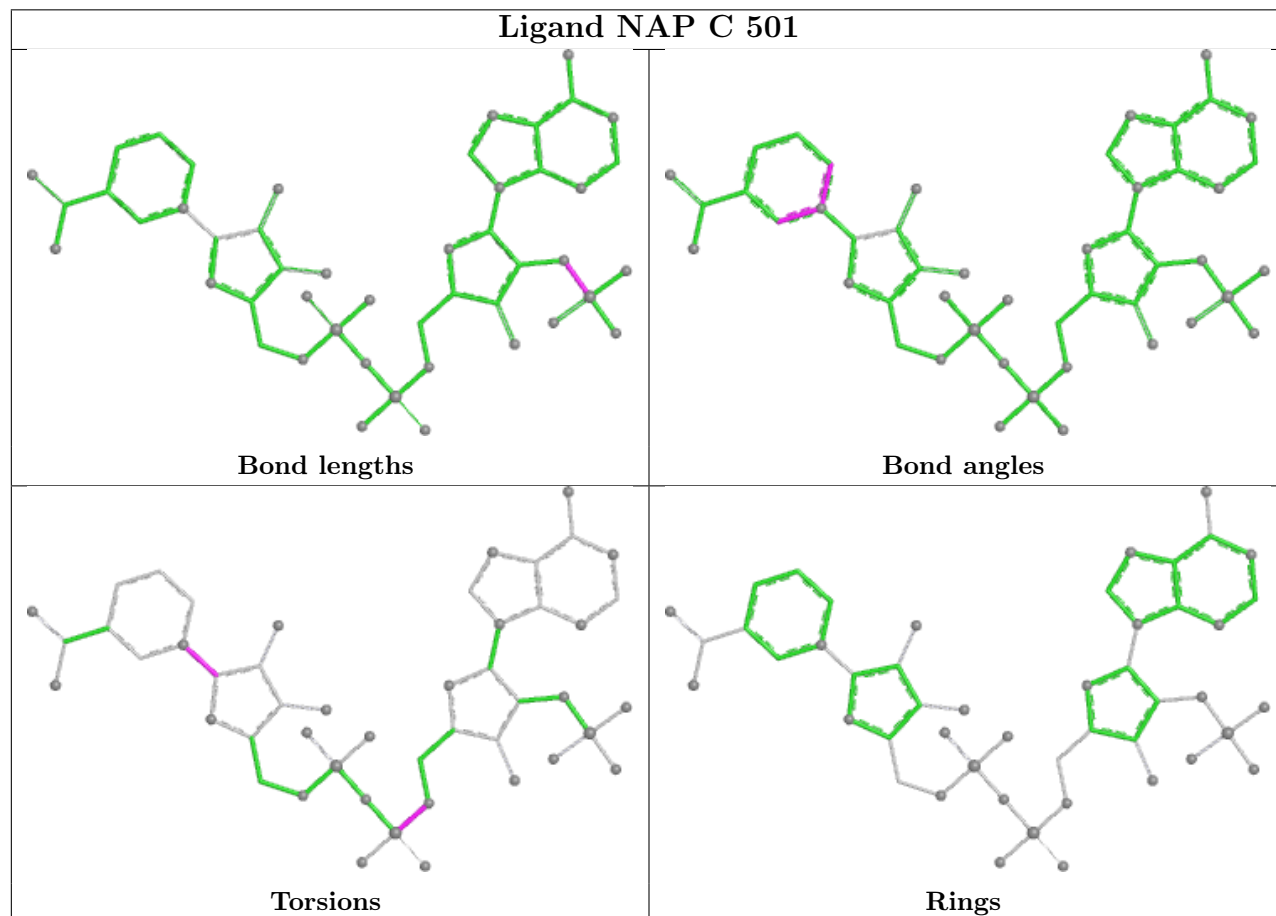




Ligand ICT B 502



Ligand NAP C 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

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	411/414 (99%)	-0.01	6 (1%) 72 78	9, 19, 39, 67	6 (1%)
1	B	412/414 (99%)	-0.21	2 (0%) 87 91	8, 16, 34, 59	2 (0%)
1	C	411/414 (99%)	0.04	4 (0%) 79 85	8, 20, 39, 68	9 (2%)
1	D	411/414 (99%)	0.22	8 (1%) 66 74	9, 23, 38, 62	4 (0%)
All	All	1645/1656 (99%)	0.01	20 (1%) 76 82	8, 20, 38, 68	21 (1%)

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	174	GLU	4.0
1	B	174	GLU	3.8
1	A	161	GLY	3.7
1	D	395	PHE	3.6
1	A	7	GLY	3.5
1	D	354	PHE	3.1
1	A	162	THR	3.0
1	C	205	TRP	2.7
1	C	161	GLY	2.7
1	D	92	TRP	2.7
1	D	205	TRP	2.6
1	C	241	ALA	2.6
1	A	412	ALA	2.4
1	D	54	ASP	2.3
1	A	383	LEU	2.2
1	B	413	LYS	2.2
1	D	80	GLU	2.2
1	C	162	THR	2.2
1	D	233	LYS	2.2
1	A	160	ASP	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 [i](#)

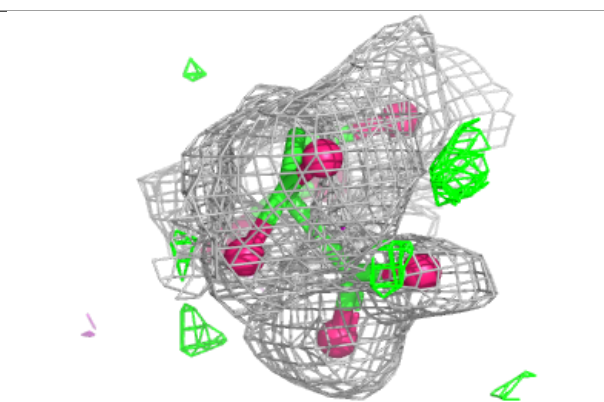
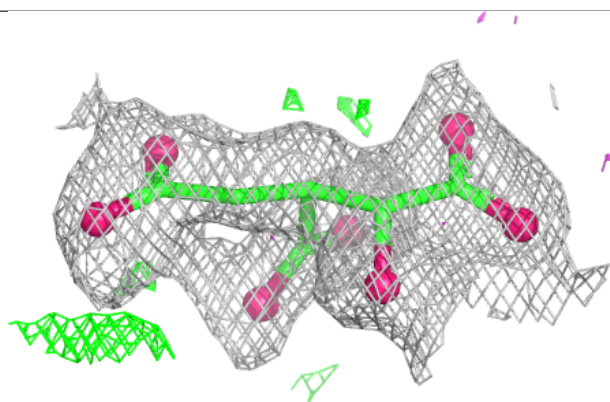
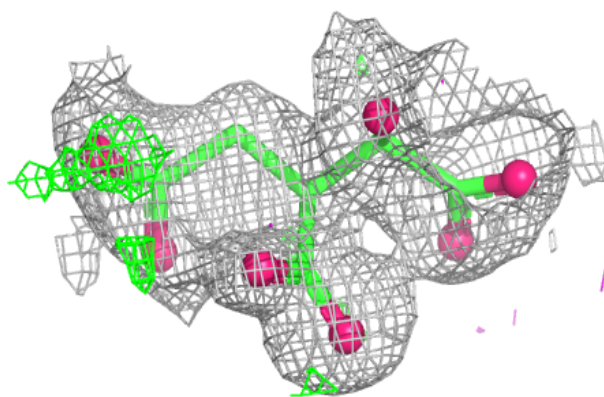
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
4	ICT	D	501	13/13	0.80	0.14	31,47,60,60	0
5	IPA	C	505	4/4	0.85	0.15	33,36,37,40	0
4	ICT	A	504	13/13	0.86	0.12	26,35,47,48	0
4	ICT	B	502	13/13	0.86	0.12	23,35,41,48	0
5	IPA	A	506	4/4	0.87	0.13	22,24,31,36	0
5	IPA	A	505	4/4	0.87	0.15	22,31,39,41	0
5	IPA	D	503	4/4	0.87	0.14	23,35,37,43	0
4	ICT	C	504	13/13	0.88	0.12	26,35,46,52	0
5	IPA	D	502	4/4	0.89	0.12	23,28,28,38	0
5	IPA	B	504	4/4	0.92	0.12	18,24,29,32	0
5	IPA	B	503	4/4	0.93	0.10	24,26,27,31	0
3	MG	C	503	1/1	0.98	0.04	21,21,21,21	0
2	NAP	A	501	48/48	0.98	0.05	7,11,21,27	0
2	NAP	B	501	48/48	0.98	0.04	8,14,22,26	0
2	NAP	C	501	48/48	0.99	0.04	7,11,18,28	0
3	MG	A	502	1/1	0.99	0.05	10,10,10,10	0
3	MG	C	502	1/1	0.99	0.04	8,8,8,8	0
3	MG	A	503	1/1	1.00	0.01	10,10,10,10	0

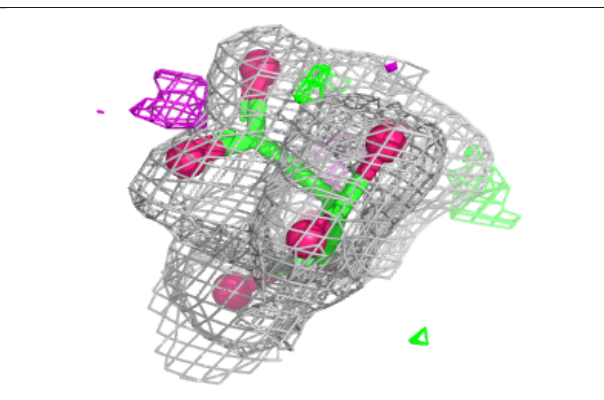
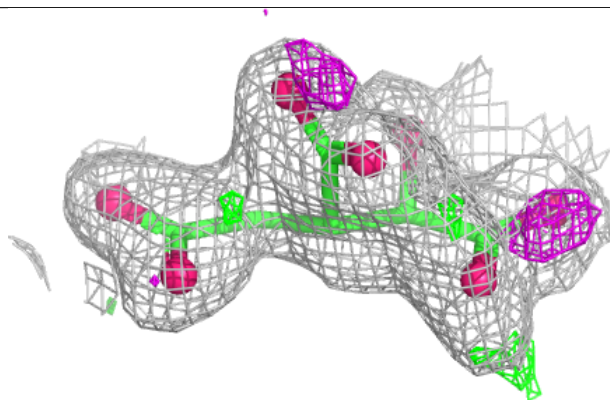
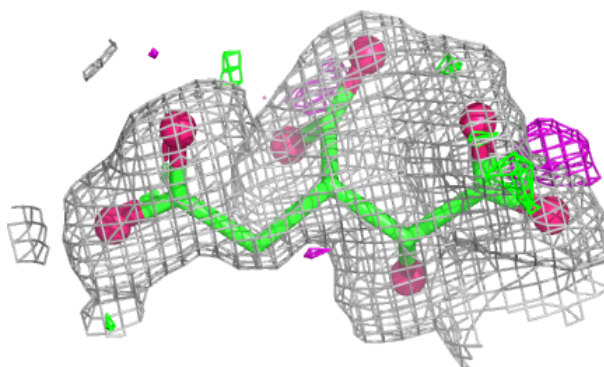
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around ICT D 501:

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

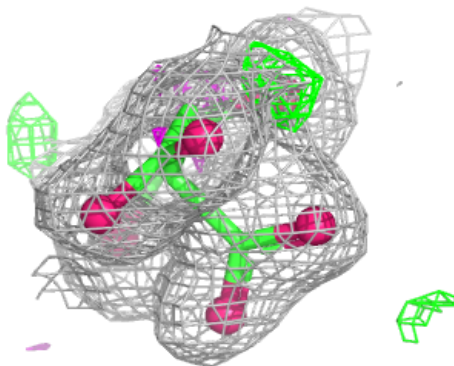
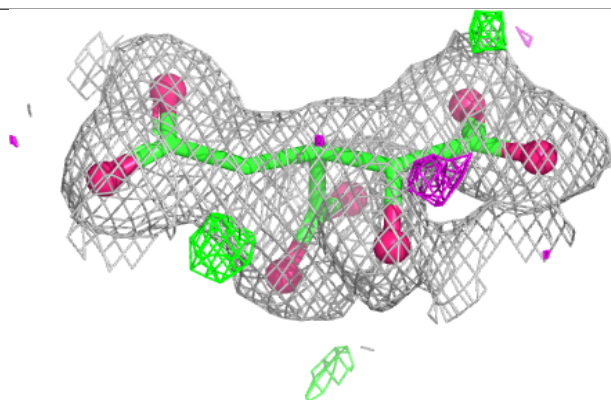
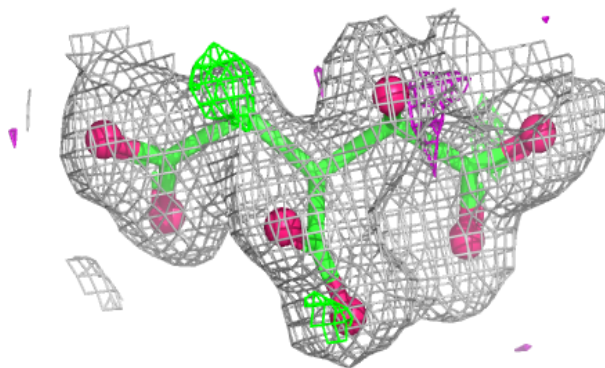
**Electron density around ICT A 504:**

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

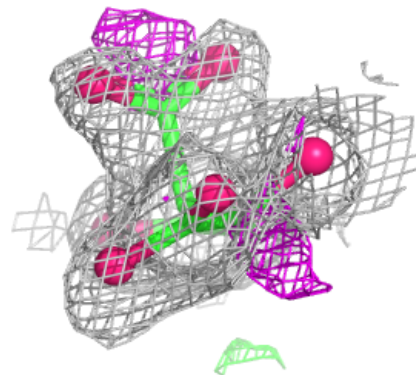
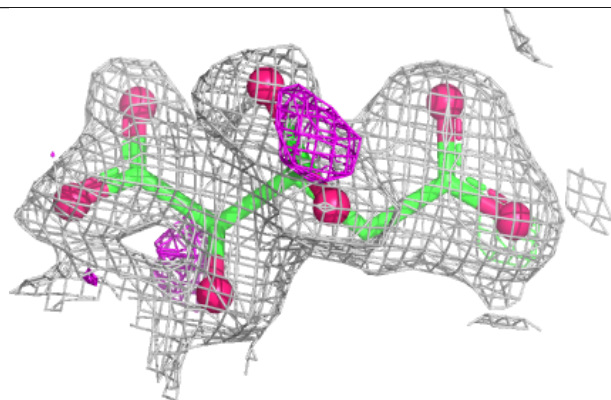
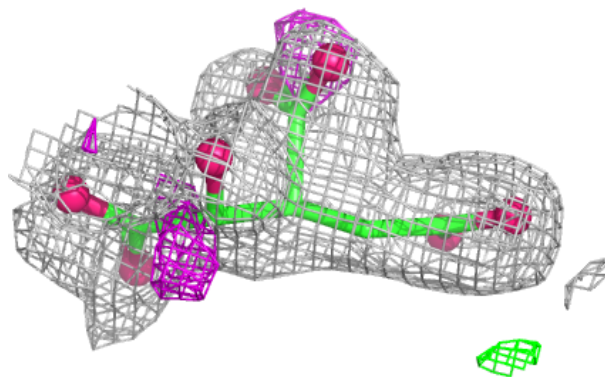


Electron density around ICT B 502:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

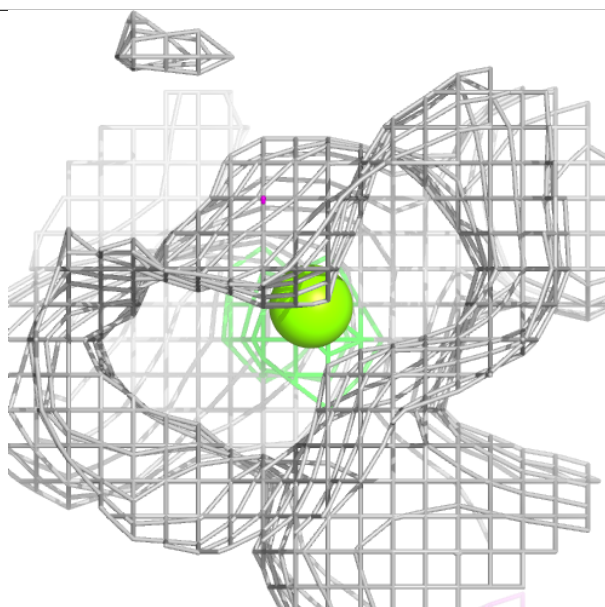
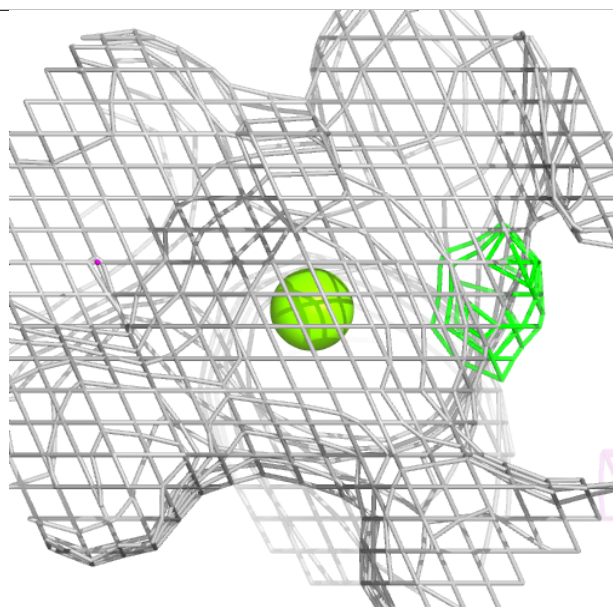
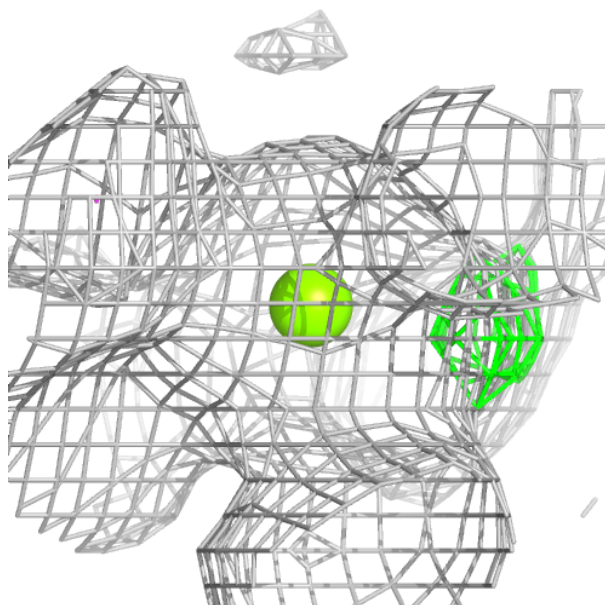
**Electron density around ICT C 504:**

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



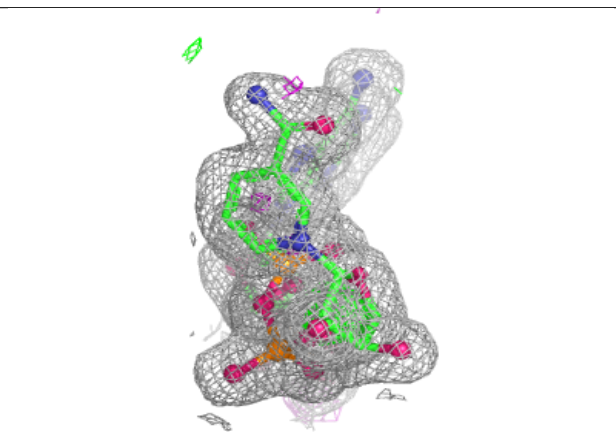
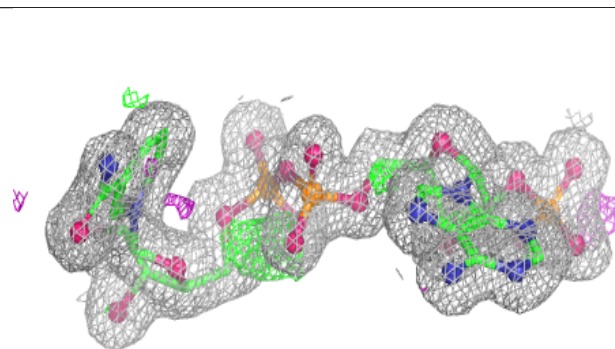
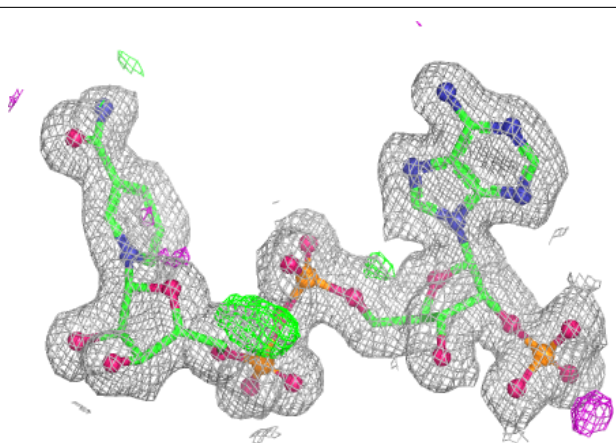
Electron density around MG C 503:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

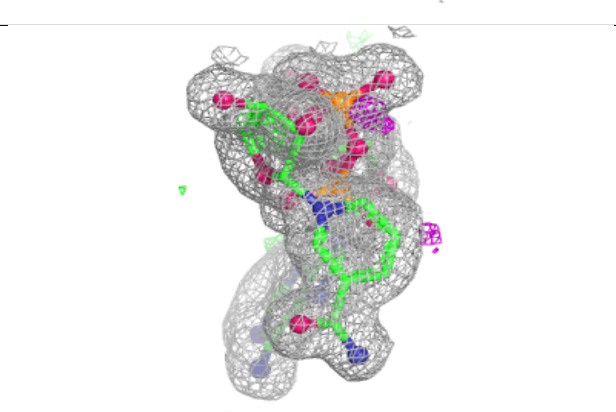
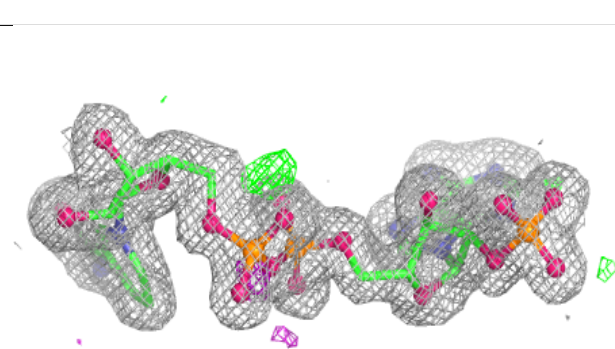
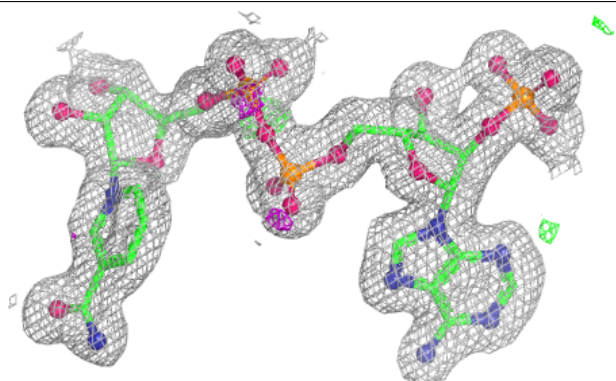


Electron density around NAP A 501:

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

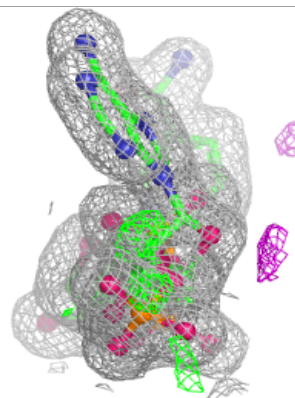
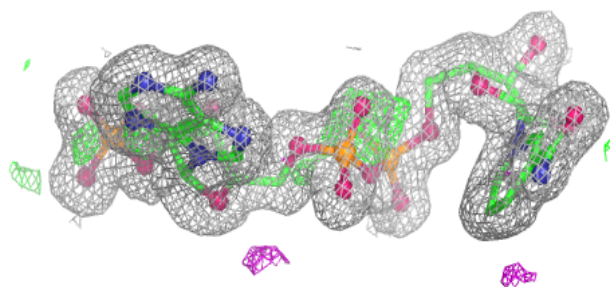
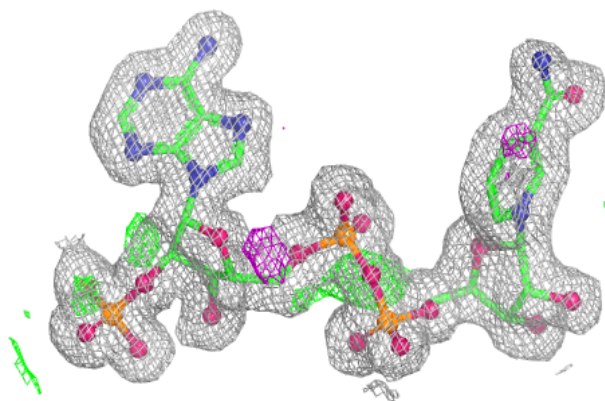
**Electron density around NAP B 501:**

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



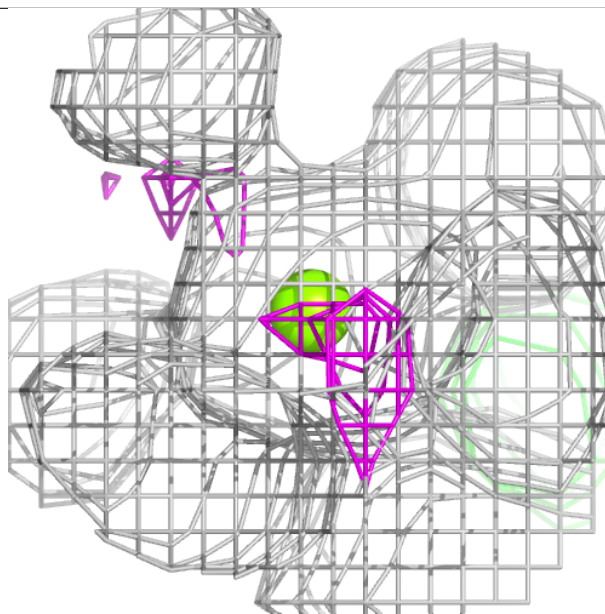
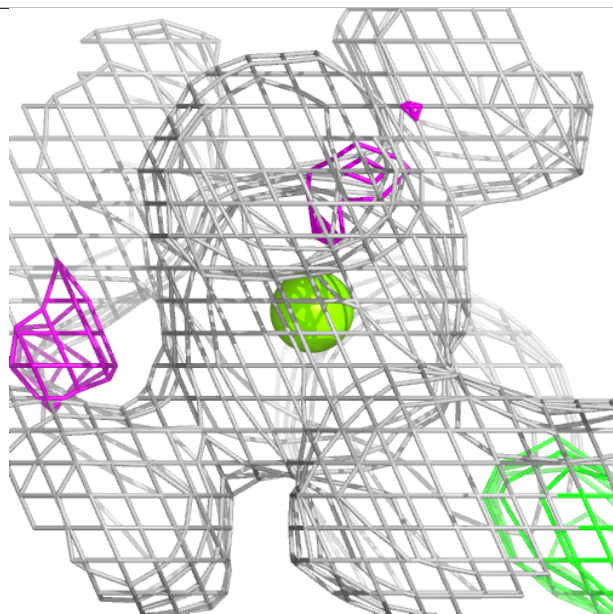
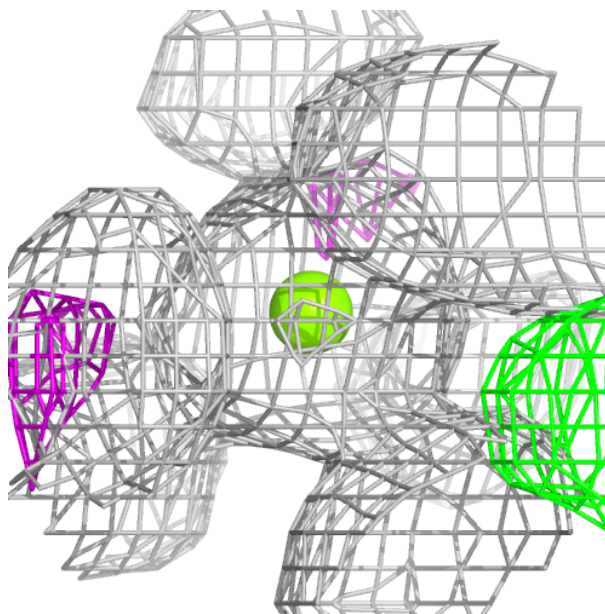
Electron density around NAP C 501:

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



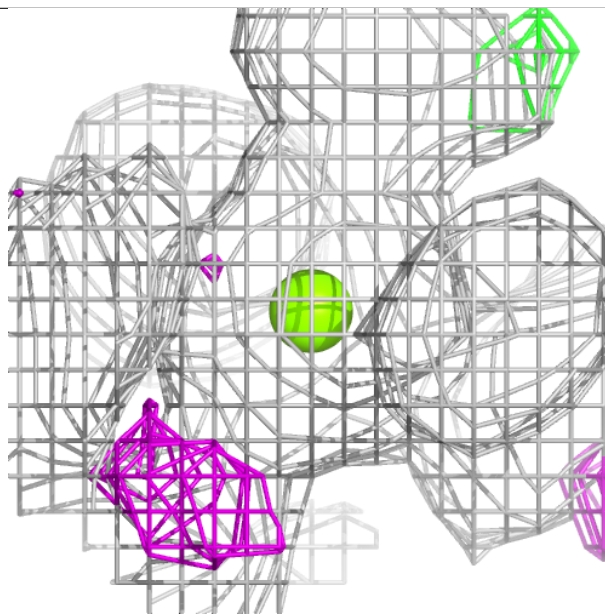
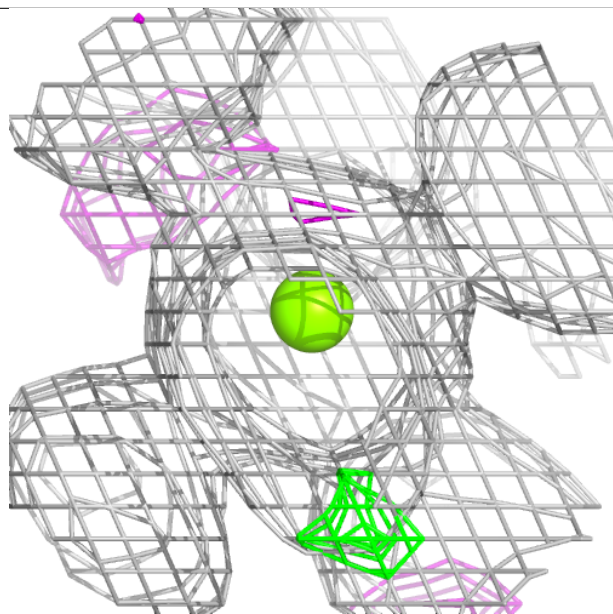
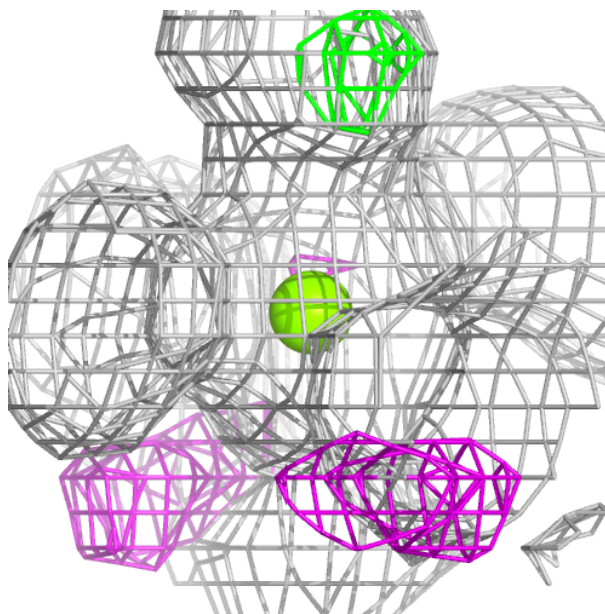
Electron density around MG A 502:

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



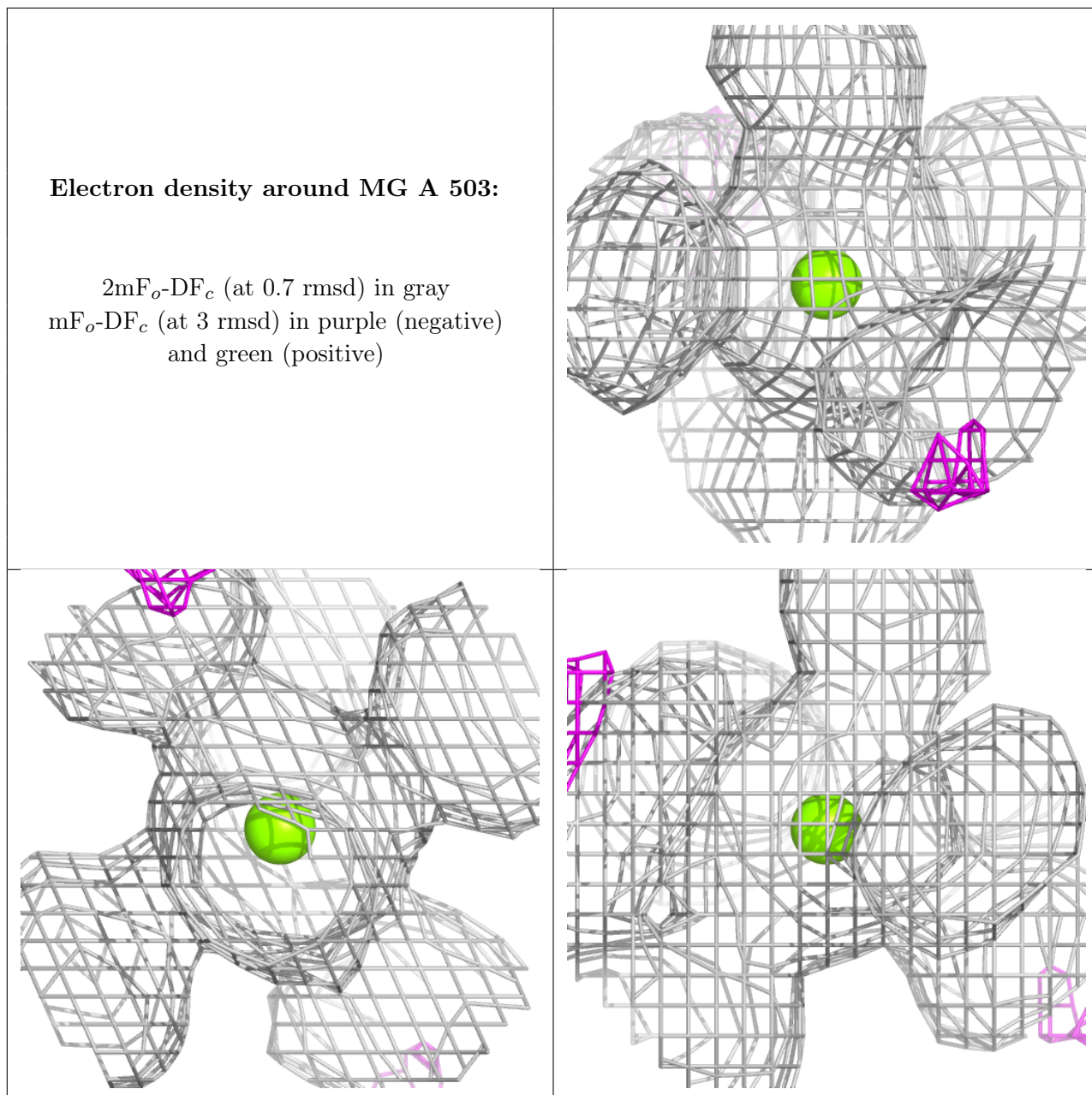
Electron density around MG C 502:

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



Electron density around MG 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 ⓘ

There are no such residues in this entry.