



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 01:18 PM UTC

PDB ID : 8XAK / pdb_00008xak
Title : Structure of Pif1-G4 complex
Authors : Hong, Z.; Song, H.
Deposited on : 2023-12-04
Resolution : 3.50 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

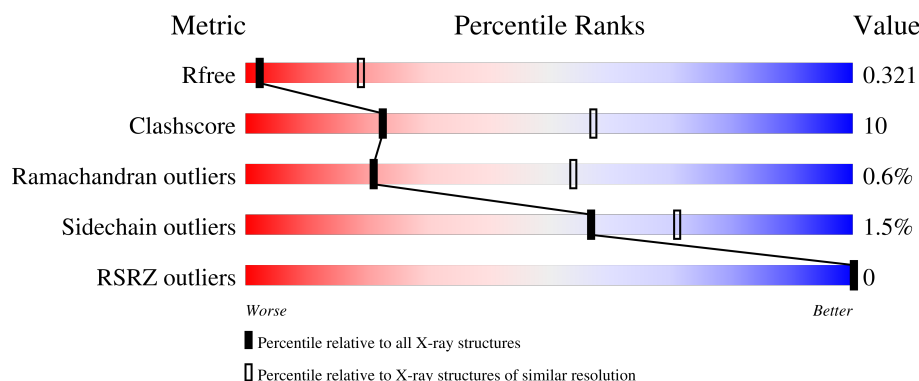
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.50 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	1085 (3.54-3.46)
Clashscore	190562	1140 (3.54-3.46)
Ramachandran outliers	187476	1113 (3.54-3.46)
Sidechain outliers	187428	1114 (3.54-3.46)
RSRZ outliers	180081	1084 (3.54-3.46)

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	545	
1	C	545	
2	B	34	
2	D	34	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	ALF	C	1002	-	-	X	-

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 9372 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.

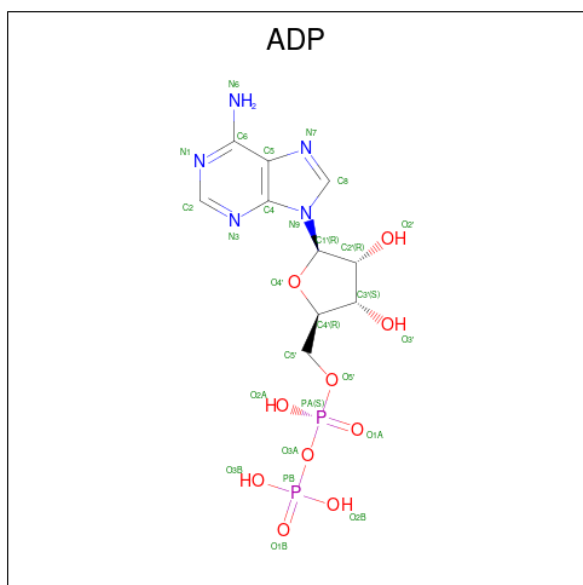
- Molecule 1 is a protein called ATP-dependent DNA helicase PIF1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	501	Total	C	N	O	S	0	0	0
			3997	2536	709	731	21			
1	C	490	Total	C	N	O	S	0	0	0
			3903	2478	689	715	21			

- Molecule 2 is a DNA chain called DNA (34-MER).

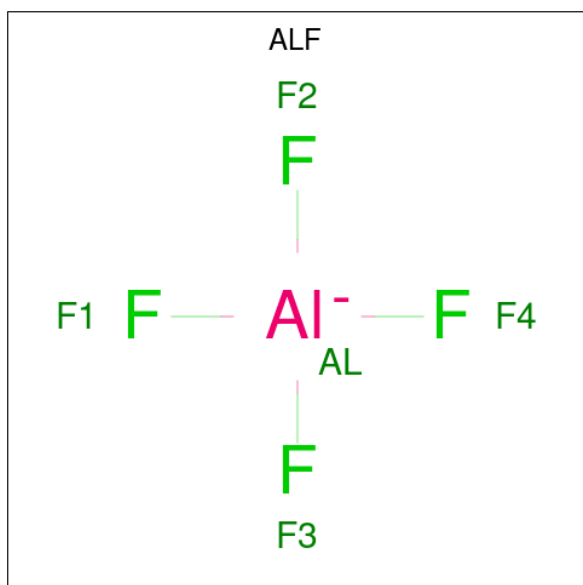
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	33	Total	C	N	O	P	0	0	0
			689	329	113	214	33			
2	B	34	Total	C	N	O	P	0	0	0
			711	340	116	221	34			

- Molecule 3 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	C	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

- Molecule 4 is TETRAFLUOROALUMINATE ION (CCD ID: ALF) (formula: AlF_4) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	Al	F	0	0
			5	1	4		
4	C	1	Total	Al	F	0	0
			5	1	4		

- Molecule 5 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Mg	0	0
			1	1		
5	C	1	Total	Mg	0	0
			1	1		

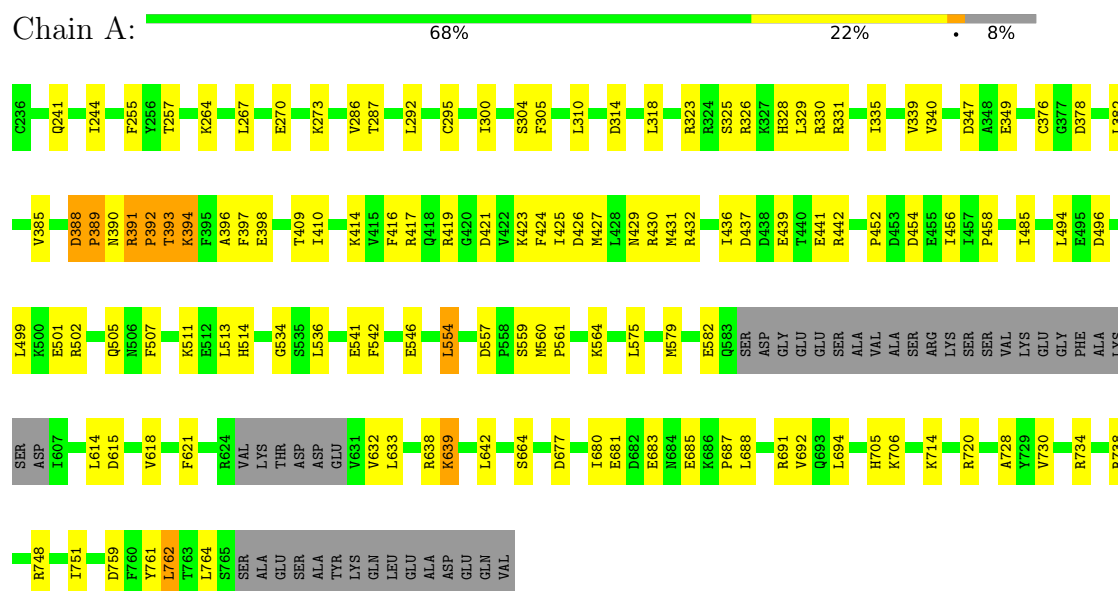
- Molecule 6 is POTASSIUM ION (CCD ID: K) (formula: K) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	D	3	Total 3	K 3	0	0
6	B	3	Total 3	K 3	0	0

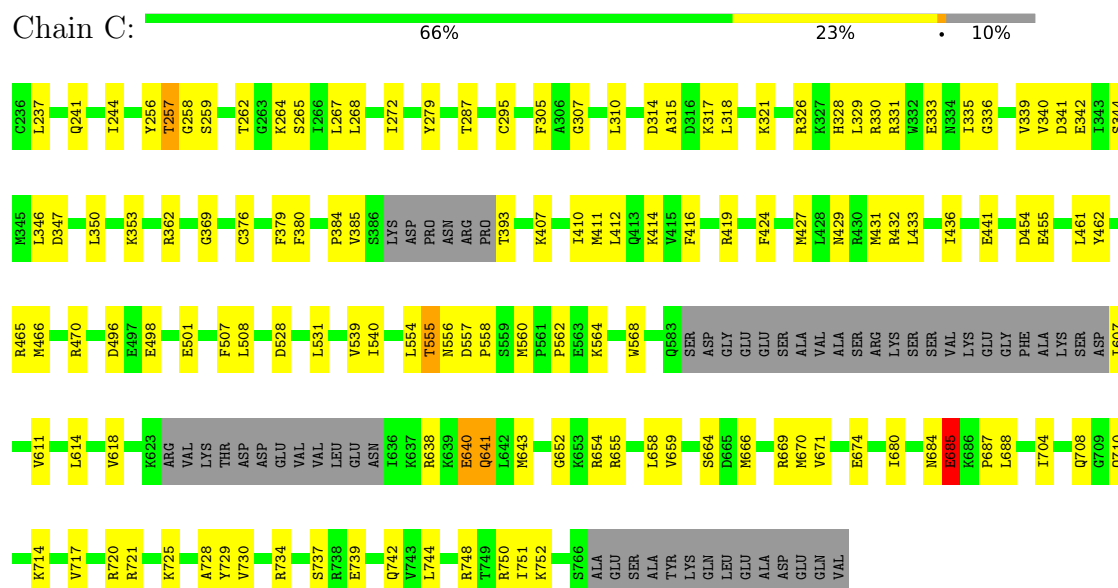
3 Residue-property plots

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: ATP-dependent DNA helicase PIF1



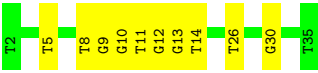
• Molecule 1: ATP-dependent DNA helicase PIF1



● Molecule 2: DNA (34-MER)



● Molecule 2: DNA (34-MER)



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	56.16Å 125.14Å 258.23Å 90.00° 91.47° 90.00°	Depositor
Resolution (Å)	48.00 – 3.50 48.00 – 3.50	Depositor EDS
% Data completeness (in resolution range)	99.4 (48.00-3.50) 92.3 (48.00-3.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.56 (at 3.48Å)	Xtriage
Refinement program	PHENIX (1.17.1_3660: ???)	Depositor
R, R_{free}	0.282 , 0.324 0.283 , 0.321	Depositor DCC
R_{free} test set	1999 reflections (8.85%)	wwPDB-VP
Wilson B-factor (Å ²)	49.8	Xtriage
Anisotropy	0.595	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 28.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.41$, $\langle L^2 \rangle = 0.23$	Xtriage
Estimated twinning fraction	0.149 for h,-k,-l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	9372	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 4.15% 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: ADP, MG, K, ALF

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	0.32	0/4065	0.79	12/5463 (0.2%)
1	C	0.36	0/3968	0.78	3/5329 (0.1%)
2	B	0.21	0/794	0.46	0/1230
2	D	0.21	0/769	0.48	0/1190
All	All	0.32	0/9596	0.74	15/13212 (0.1%)

There are no bond length outliers.

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	392	PRO	N-CA-C	-9.96	83.19	113.27
1	C	638	ARG	N-CA-C	-7.13	104.31	113.16
1	A	554	LEU	N-CA-C	-6.62	104.07	111.28
1	A	683	GLU	N-CA-C	-6.43	103.54	111.33
1	A	391	ARG	N-CA-C	6.26	119.95	108.94
1	C	562	PRO	CB-CA-C	-6.04	102.79	112.21
1	A	391	ARG	CB-CA-C	-5.91	101.92	111.14
1	C	555	THR	N-CA-C	-5.81	102.86	110.53
1	A	397	PHE	CA-CB-CG	5.59	119.39	113.80
1	A	762	LEU	CA-C-N	-5.45	113.55	122.54
1	A	762	LEU	C-N-CA	-5.45	113.55	122.54
1	A	685	GLU	CA-C-N	-5.27	113.60	120.67
1	A	685	GLU	C-N-CA	-5.27	113.60	120.67
1	A	394	LYS	N-CA-C	-5.08	102.78	110.24
1	A	638	ARG	N-CA-C	-5.01	106.43	112.54

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	3997	0	4105	85	1
1	C	3903	0	4005	91	0
2	B	711	0	393	8	1
2	D	689	0	377	10	0
3	A	27	0	12	1	0
3	C	27	0	12	0	0
4	A	5	0	0	1	0
4	C	5	0	0	2	0
5	A	1	0	0	0	0
5	C	1	0	0	0	0
6	B	3	0	0	0	0
6	D	3	0	0	0	0
All	All	9372	0	8904	185	1

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 10.

All (185) 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:466:MET:HE1	1:C:470:ARG:HD3	1.61	0.83
1:C:330:ARG:NH2	1:C:333:GLU:OE1	2.15	0.80
1:C:654:ARG:NH2	1:C:674:GLU:OE1	2.18	0.77
1:C:531:LEU:HD21	1:C:671:VAL:HG21	1.68	0.76
1:A:759:ASP:HA	1:A:762:LEU:HD12	1.66	0.75
1:C:244:ILE:HG12	1:C:410:ILE:HD12	1.67	0.74
1:C:721:ARG:HB3	2:D:2:DT:H3	1.50	0.74
1:C:314:ASP:HB2	1:C:317:LYS:HE2	1.70	0.71
1:A:423:LYS:HD2	1:A:738:ARG:HH22	1.55	0.71
1:C:658:LEU:HD11	1:C:670:MET:HG2	1.72	0.70
1:C:424:PHE:HA	1:C:427:MET:HE2	1.75	0.69
1:C:441:GLU:OE1	1:C:748:ARG:NH1	2.25	0.68
1:A:541:GLU:HG3	1:A:542:PHE:H	1.60	0.67
1:A:326:ARG:HD2	1:A:330:ARG:HE	1.60	0.66
2:B:8:DT:H2"	2:B:9:DG:H5"	1.78	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:557:ASP:HB2	1:A:560:MET:HB2	1.77	0.64
1:A:494:LEU:HD11	1:A:688:LEU:HG	1.80	0.63
1:C:431:MET:HE2	1:C:751:ILE:HD13	1.80	0.63
1:A:761:TYR:HA	1:A:764:LEU:HD13	1.81	0.62
1:A:323:ARG:HH11	1:A:323:ARG:HG2	1.64	0.62
1:A:680:ILE:HG22	1:A:681:GLU:H	1.65	0.62
1:C:684:ASN:O	1:C:685:GLU:HB2	2.01	0.61
1:A:728:ALA:HB3	1:A:751:ILE:HG22	1.83	0.60
1:C:540:ILE:HG22	1:C:618:VAL:HG12	1.83	0.60
1:A:287:THR:HA	1:A:300:ILE:O	2.01	0.60
1:A:244:ILE:HB	1:A:267:LEU:HD11	1.84	0.60
1:C:268:LEU:HD21	1:C:339:VAL:HG11	1.83	0.60
1:A:393:THR:O	1:A:394:LYS:C	2.45	0.59
1:C:461:LEU:HB2	1:C:717:VAL:HG12	1.84	0.59
1:A:431:MET:HE2	1:A:436:ILE:HD11	1.85	0.59
1:C:307:GLY:O	1:C:321:LYS:NZ	2.29	0.59
1:A:255:PHE:O	1:A:409:THR:HA	2.03	0.59
2:D:3:DT:H2''	2:D:4:DT:H5'	1.84	0.58
1:A:734:ARG:NH2	4:A:1002:ALF:F1	2.27	0.57
1:C:315:ALA:HB2	1:C:353:LYS:HG2	1.86	0.57
1:A:541:GLU:HB2	1:A:618:VAL:HG11	1.85	0.57
1:A:511:LYS:HA	1:A:691:ARG:HH12	1.70	0.57
1:C:687:PRO:HG2	2:D:18:DT:H5'	1.86	0.57
1:A:748:ARG:O	1:A:751:ILE:HG13	2.05	0.57
1:C:346:LEU:HG	1:C:350:LEU:HD23	1.87	0.56
1:C:607:ILE:HA	1:C:652:GLY:HA2	1.87	0.56
1:C:342:GLU:OE2	4:C:1002:ALF:F2	2.14	0.56
1:C:640:GLU:HG3	1:C:641:GLN:N	2.21	0.56
1:A:305:PHE:CE1	1:A:331:ARG:HG2	2.40	0.55
1:A:326:ARG:HD3	2:B:11:DT:H71	1.87	0.55
1:A:310:LEU:HD22	2:B:5:DT:C2	2.41	0.54
2:B:10:DG:H4'	2:B:11:DT:H72	1.90	0.54
1:C:385:VAL:HG13	2:D:3:DT:C2	2.43	0.54
1:A:295:CYS:O	1:A:664:SER:HB2	2.09	0.53
1:A:639:LYS:HA	1:A:642:LEU:HB2	1.90	0.53
1:A:417:ARG:NH1	3:A:1001:ADP:O2B	2.42	0.53
1:A:430:ARG:NH1	1:A:437:ASP:OD1	2.34	0.53
1:A:419:ARG:HA	1:A:425:ILE:HD11	1.92	0.52
1:C:555:THR:O	1:C:556:ASN:C	2.52	0.52
1:A:557:ASP:CB	1:A:560:MET:HB2	2.40	0.51
1:A:314:ASP:O	1:A:318:LEU:HD13	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:257:THR:HG23	1:C:411:MET:HG3	1.92	0.50
1:C:265:SER:OG	1:C:341:ASP:OD2	2.29	0.50
1:C:720:ARG:O	1:C:750:ARG:NH1	2.35	0.50
1:A:680:ILE:HG22	1:A:681:GLU:N	2.25	0.50
1:C:429:ASN:O	1:C:433:LEU:HD12	2.10	0.50
1:A:329:LEU:HD22	2:B:11:DT:C4	2.47	0.50
1:C:362:ARG:HD3	1:C:369:GLY:HA3	1.94	0.49
1:A:456:ILE:HB	1:A:714:LYS:HE3	1.95	0.49
1:A:264:LYS:HB3	1:A:376:CYS:SG	2.52	0.49
1:A:304:SER:O	1:A:328:HIS:HE1	1.96	0.49
1:A:382:LEU:HD13	1:A:705:HIS:ND1	2.28	0.49
2:B:13:DG:H2''	2:B:30:DG:H5'	1.94	0.49
1:A:730:VAL:O	1:A:734:ARG:NH1	2.45	0.48
1:A:680:ILE:O	1:A:687:PRO:HA	2.14	0.48
1:A:323:ARG:HA	1:A:329:LEU:HD12	1.95	0.48
1:C:350:LEU:HA	1:C:353:LYS:HE2	1.95	0.48
1:C:730:VAL:HG13	1:C:734:ARG:NH1	2.29	0.48
1:C:666:MET:O	1:C:666:MET:HG2	2.13	0.48
1:C:720:ARG:NH2	1:C:721:ARG:HD3	2.28	0.48
1:A:561:PRO:HG2	1:A:564:LYS:HD3	1.96	0.48
1:C:262:THR:HG22	1:C:414:LYS:O	2.13	0.48
1:C:326:ARG:NH1	1:C:330:ARG:HE	2.11	0.48
1:A:241:GLN:CA	1:A:267:LEU:HD12	2.44	0.47
1:A:513:LEU:HD23	1:A:694:LEU:HD12	1.95	0.47
2:D:16:DG:H4'	2:D:17:DT:H5''	1.96	0.47
1:A:286:VAL:HA	1:A:339:VAL:HB	1.95	0.47
1:A:507:PHE:CD2	1:A:691:ARG:HB3	2.50	0.47
1:A:546:GLU:HB3	1:A:575:LEU:HD21	1.96	0.47
1:C:305:PHE:O	1:C:328:HIS:HD2	1.98	0.47
1:A:507:PHE:CE2	1:A:691:ARG:HB3	2.50	0.47
1:C:507:PHE:CZ	1:C:680:ILE:HG12	2.49	0.47
1:C:557:ASP:HB2	1:C:560:MET:HB2	1.96	0.47
1:A:541:GLU:HG3	1:A:542:PHE:N	2.26	0.47
1:A:554:LEU:HG	1:A:621:PHE:HZ	1.80	0.47
1:C:432:ARG:HB2	1:C:729:TYR:CZ	2.50	0.47
1:A:615:ASP:O	1:A:618:VAL:HG22	2.15	0.47
1:A:677:ASP:HB3	1:A:692:VAL:HG22	1.97	0.47
1:C:441:GLU:CD	1:C:748:ARG:HH11	2.21	0.47
1:A:499:LEU:HD11	1:A:502:ARG:HH21	1.79	0.47
1:A:378:ASP:CG	1:A:432:ARG:HH21	2.23	0.46
1:C:607:ILE:HD12	1:C:652:GLY:HA3	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:429:ASN:O	1:A:432:ARG:HB3	2.15	0.46
1:C:264:LYS:HG2	1:C:412:LEU:HD12	1.97	0.46
1:C:432:ARG:HB2	1:C:729:TYR:CE2	2.51	0.46
1:A:325:SER:OG	1:A:328:HIS:HB2	2.16	0.46
1:A:496:ASP:HB3	1:A:499:LEU:HB2	1.98	0.46
1:C:244:ILE:HB	1:C:267:LEU:HD11	1.98	0.46
1:C:407:LYS:HD3	1:C:407:LYS:HA	1.45	0.46
1:A:409:THR:C	1:A:410:ILE:HG13	2.40	0.45
1:A:452:PRO:HB2	1:A:454:ASP:OD2	2.16	0.45
1:A:496:ASP:OD2	1:A:499:LEU:HD13	2.15	0.45
1:A:388:ASP:N	1:A:389:PRO:HD3	2.31	0.45
1:C:329:LEU:HD13	2:D:11:DT:C5	2.52	0.45
2:D:16:DG:H4'	2:D:17:DT:H3'	1.99	0.45
1:A:426:ASP:OD1	1:A:430:ARG:HD2	2.16	0.45
1:C:507:PHE:CE1	1:C:680:ILE:HG12	2.52	0.45
1:C:241:GLN:CB	1:C:267:LEU:HD12	2.47	0.45
1:C:734:ARG:NH2	4:C:1002:ALF:F1	2.40	0.45
1:C:643:MET:HB3	1:C:643:MET:HE3	1.78	0.45
1:C:380:PHE:HB3	1:C:432:ARG:HG3	1.99	0.44
1:C:728:ALA:HB3	1:C:751:ILE:HG12	1.99	0.44
1:C:328:HIS:CD2	1:C:331:ARG:HH11	2.35	0.44
1:C:611:VAL:HG22	1:C:655:ARG:HH12	1.81	0.44
1:A:391:ARG:C	1:A:393:THR:N	2.74	0.44
1:A:305:PHE:HZ	1:A:335:ILE:HD13	1.83	0.44
1:A:389:PRO:HB2	1:A:390:ASN:H	1.57	0.44
1:C:465:ARG:NH2	1:C:508:LEU:HD23	2.33	0.44
1:C:708:GLN:HA	1:C:734:ARG:HH11	1.81	0.44
1:A:421:ASP:HB3	1:A:424:PHE:HB3	1.99	0.44
1:A:541:GLU:HG2	1:A:614:LEU:HG	1.99	0.44
1:C:268:LEU:HD23	1:C:272:ILE:HD11	2.00	0.44
1:C:314:ASP:O	1:C:318:LEU:HD13	2.18	0.44
1:A:424:PHE:HA	1:A:427:MET:HE2	1.99	0.43
1:A:305:PHE:HE1	1:A:331:ARG:HG2	1.81	0.43
1:C:257:THR:OG1	1:C:258:GLY:N	2.50	0.43
1:C:539:VAL:HG22	1:C:659:VAL:HG22	2.01	0.43
1:C:462:TYR:O	1:C:704:ILE:HG13	2.18	0.43
1:C:310:LEU:HD23	2:D:5:DT:C4	2.54	0.43
1:C:564:LYS:HE3	1:C:568:TRP:NE1	2.34	0.43
1:C:455:GLU:H	1:C:455:GLU:CD	2.27	0.43
1:C:710:GLN:O	1:C:734:ARG:HD3	2.19	0.43
2:B:14:DT:H6	2:B:14:DT:H2'	1.67	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:305:PHE:HZ	1:C:335:ILE:HD13	1.83	0.43
1:C:344:SER:HA	1:C:379:PHE:CE1	2.53	0.43
1:A:392:PRO:O	1:A:393:THR:C	2.62	0.43
1:A:458:PRO:HB3	1:A:714:LYS:HG2	2.01	0.43
1:C:737:SER:HB2	1:C:739:GLU:OE1	2.18	0.43
1:C:241:GLN:HB3	1:C:267:LEU:HD12	2.00	0.42
1:A:255:PHE:CZ	1:A:257:THR:HB	2.55	0.42
1:C:614:LEU:HD23	1:C:618:VAL:HG21	2.01	0.42
1:A:579:MET:O	1:A:582:GLU:HG3	2.19	0.42
1:C:640:GLU:HG3	1:C:641:GLN:H	1.85	0.42
2:D:13:DG:H3'	2:D:29:DT:H2''	2.00	0.42
1:A:439:GLU:HG2	1:A:442:ARG:NH2	2.35	0.42
1:C:380:PHE:CB	1:C:432:ARG:HG3	2.50	0.42
1:A:347:ASP:OD1	1:A:349:GLU:HB2	2.19	0.42
1:A:270:GLU:HA	1:A:273:LYS:HE2	2.01	0.42
1:A:292:LEU:HD23	1:A:706:LYS:O	2.20	0.42
1:A:396:ALA:C	1:A:398:GLU:H	2.28	0.42
1:C:436:ILE:HD13	1:C:436:ILE:HA	1.91	0.42
1:C:742:GLN:HE21	1:C:744:LEU:HD21	1.84	0.42
1:C:295:CYS:O	1:C:664:SER:HB2	2.20	0.42
2:D:6:DT:H6	2:D:6:DT:H2'	1.65	0.42
1:C:730:VAL:O	1:C:734:ARG:HG2	2.18	0.42
1:A:485:ILE:HG12	1:A:514:HIS:CE1	2.55	0.41
1:C:416:PHE:O	1:C:419:ARG:NH1	2.52	0.41
1:A:499:LEU:CD1	1:A:502:ARG:HH21	2.33	0.41
1:A:534:GLY:O	1:A:536:LEU:HD22	2.20	0.41
1:C:725:LYS:HE3	1:C:752:LYS:HE3	2.02	0.41
1:A:323:ARG:NH1	2:B:12:DG:O5'	2.53	0.41
1:C:528:ASP:OD1	1:C:669:ARG:NH1	2.45	0.41
1:C:305:PHE:CZ	1:C:335:ILE:HD13	2.56	0.41
1:C:498:GLU:HG2	1:C:501:GLU:HB3	2.01	0.41
1:C:431:MET:HE3	1:C:431:MET:HB2	1.90	0.41
1:C:264:LYS:HB3	1:C:376:CYS:SG	2.60	0.41
1:A:436:ILE:HD12	1:A:751:ILE:HG12	2.02	0.41
1:C:326:ARG:HH12	1:C:330:ARG:HE	1.68	0.41
1:C:344:SER:O	1:C:384:PRO:HD3	2.21	0.41
1:C:454:ASP:OD1	1:C:714:LYS:NZ	2.36	0.41
1:A:241:GLN:HA	1:A:267:LEU:HD12	2.02	0.40
1:A:414:LYS:HD3	1:A:416:PHE:CZ	2.55	0.40
1:C:347:ASP:OD1	1:C:347:ASP:N	2.53	0.40
1:C:279:TYR:CD1	1:C:336:GLY:HA3	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:501:GLU:O	1:A:505:GLN:HG2	2.21	0.40
1:C:287:THR:O	1:C:340:VAL:HA	2.20	0.40
1:C:554:LEU:HD12	1:C:554:LEU:HA	1.84	0.40
1:A:287:THR:O	1:A:340:VAL:HA	2.21	0.40
1:A:441:GLU:OE2	1:A:748:ARG:HD2	2.20	0.40
1:C:256:TYR:HB3	1:C:410:ILE:HG13	2.03	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:720:ARG:NH2	2:B:26:DT:O2[1_455]	2.19	0.01

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	495/545 (91%)	472 (95%)	21 (4%)	2 (0%)	30	62
1	C	482/545 (88%)	466 (97%)	12 (2%)	4 (1%)	16	49
All	All	977/1090 (90%)	938 (96%)	33 (3%)	6 (1%)	21	54

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	259	SER
1	C	496	ASP
1	C	685	GLU
1	A	389	PRO
1	A	559	SER
1	C	558	PRO

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	438/474 (92%)	432 (99%)	6 (1%)	59	71
1	C	427/474 (90%)	420 (98%)	7 (2%)	55	70
All	All	865/948 (91%)	852 (98%)	13 (2%)	57	71

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	385	VAL
1	A	388	ASP
1	A	393	THR
1	A	632	VAL
1	A	633	LEU
1	A	639	LYS
1	C	237	LEU
1	C	257	THR
1	C	393	THR
1	C	640	GLU
1	C	641	GLN
1	C	685	GLU
1	C	688	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	334	ASN
1	A	514	HIS
1	A	647	HIS
1	A	755	GLN
1	C	252	HIS
1	C	303	HIS
1	C	328	HIS
1	C	487	ASN
1	C	514	HIS

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Mol	Chain	Res	Type
1	C	641	GLN
1	C	644	GLN
1	C	648	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 12 ligands modelled in this entry, 8 are monoatomic - leaving 4 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	ADP	A	1001	5	28,29,29	1.42	4 (14%)	43,45,45	1.82	8 (18%)
4	ALF	A	1002	-	4,4,4	1.35	0	-		
3	ADP	C	1001	5	28,29,29	1.42	5 (17%)	43,45,45	1.82	9 (20%)
4	ALF	C	1002	-	4,4,4	1.37	0	-		

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	ADP	A	1001	5	-	1/16/32/32	0/3/3/3
3	ADP	C	1001	5	-	2/16/32/32	0/3/3/3

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1001	ADP	C5-C4	4.78	1.47	1.39
3	C	1001	ADP	C5-C4	4.73	1.47	1.39
3	A	1001	ADP	C5-C6	2.77	1.48	1.41
3	C	1001	ADP	C5-C6	2.73	1.48	1.41
3	C	1001	ADP	C8-N7	2.36	1.36	1.31
3	A	1001	ADP	C8-N7	2.35	1.36	1.31
3	C	1001	ADP	C5-N7	-2.20	1.35	1.39
3	A	1001	ADP	C5-N7	-2.20	1.35	1.39
3	C	1001	ADP	PA-O3A	2.05	1.61	1.59

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	1001	ADP	C5-C4-N3	-5.82	118.70	126.72
3	A	1001	ADP	C5-C4-N3	-5.78	118.76	126.72
3	C	1001	ADP	N3-C4-N9	4.60	135.00	127.17
3	A	1001	ADP	N3-C4-N9	4.59	134.97	127.17
3	A	1001	ADP	C2-N3-C4	3.68	120.81	111.83
3	C	1001	ADP	C2-N3-C4	3.66	120.77	111.83
3	A	1001	ADP	C4-C5-N7	-3.53	106.55	110.58
3	C	1001	ADP	C4-C5-N7	-3.43	106.66	110.58
3	A	1001	ADP	N3-C2-N1	-3.23	123.70	128.58
3	C	1001	ADP	N3-C2-N1	-3.18	123.77	128.58
3	A	1001	ADP	C4-N9-C8	2.72	108.59	105.74
3	C	1001	ADP	C4-N9-C8	2.58	108.45	105.74
3	A	1001	ADP	C5-N7-C8	2.56	107.48	103.45
3	C	1001	ADP	C5-N7-C8	2.51	107.40	103.45
3	C	1001	ADP	C3'-C2'-C1'	2.18	105.58	101.46
3	A	1001	ADP	C6-C5-N7	2.13	136.19	132.09
3	C	1001	ADP	C6-C5-N7	2.03	136.00	132.09

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	1001	ADP	C5'-O5'-PA-O1A

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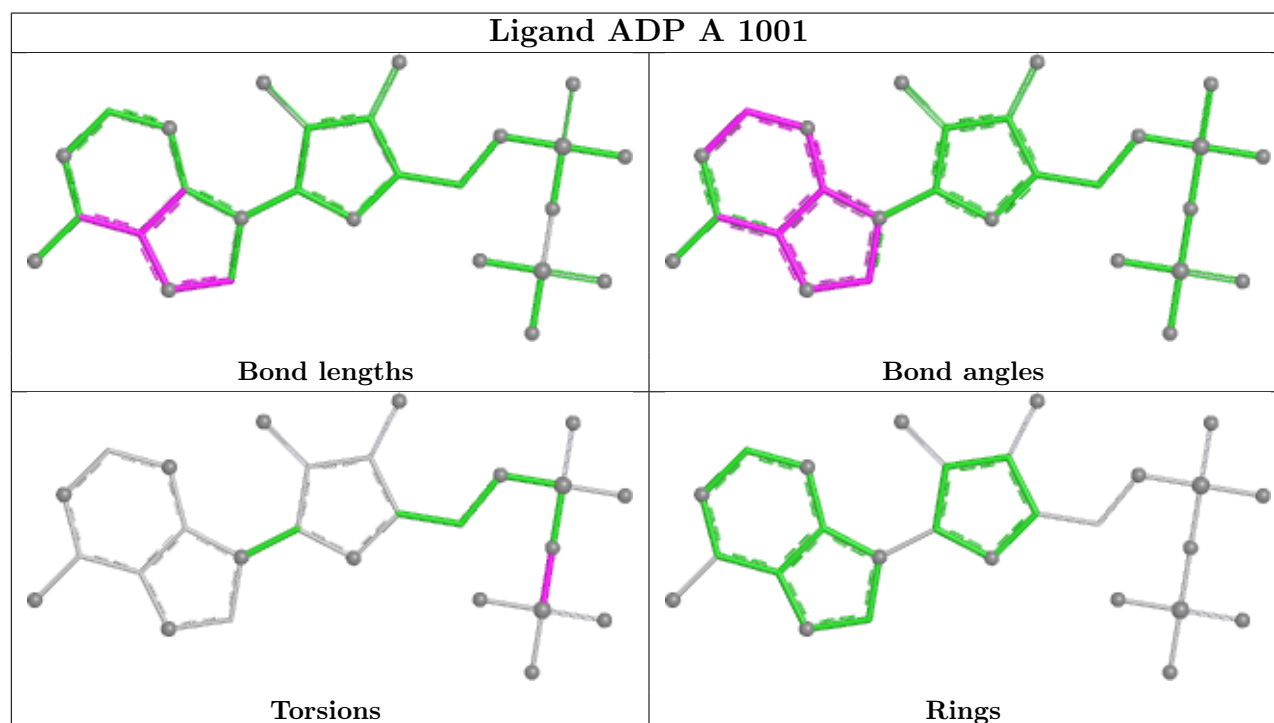
Mol	Chain	Res	Type	Atoms
3	C	1001	ADP	C5'-O5'-PA-O3A
3	A	1001	ADP	PA-O3A-PB-O1B

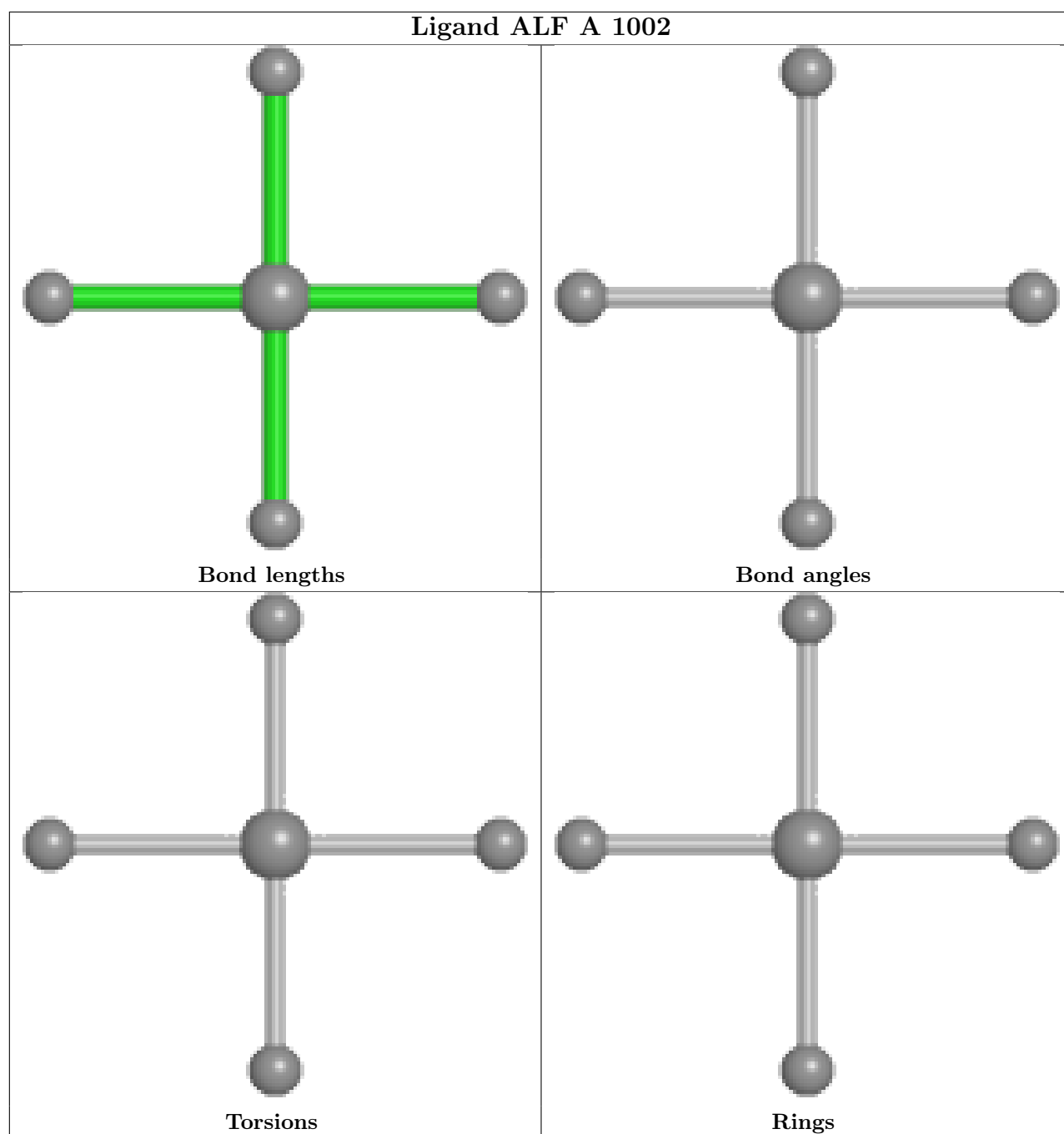
There are no ring outliers.

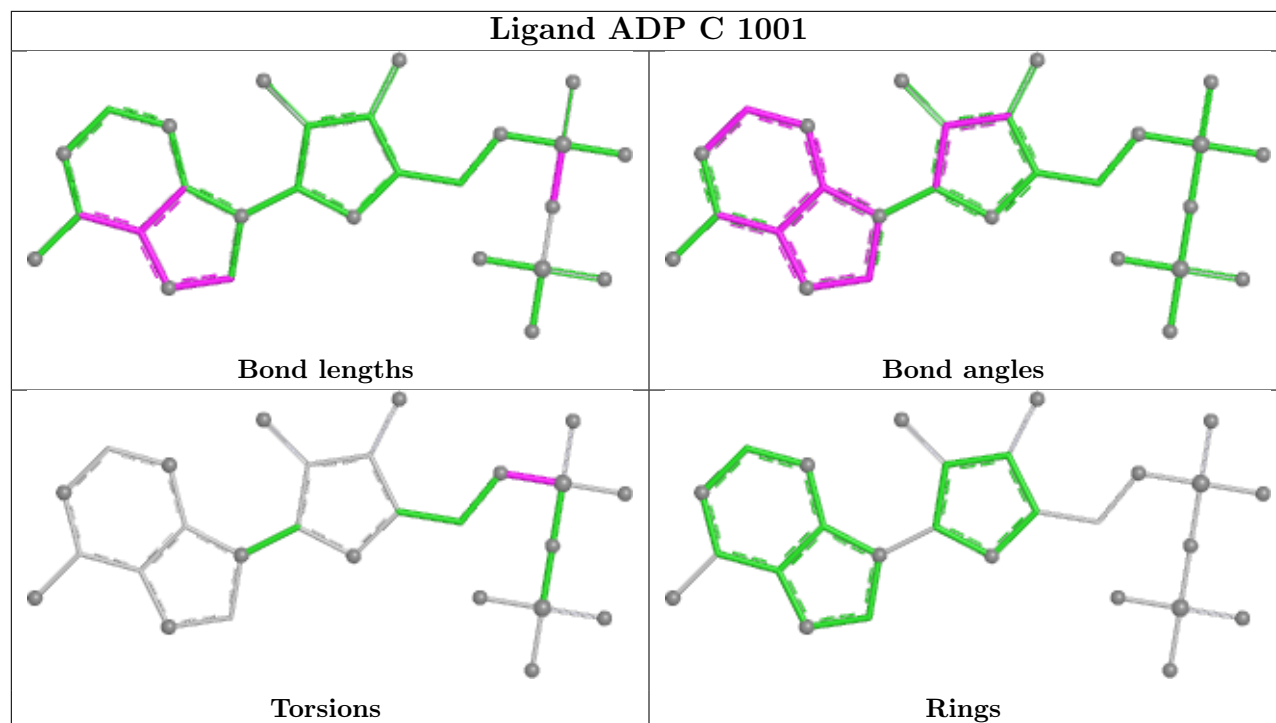
3 monomers are involved in 4 short contacts:

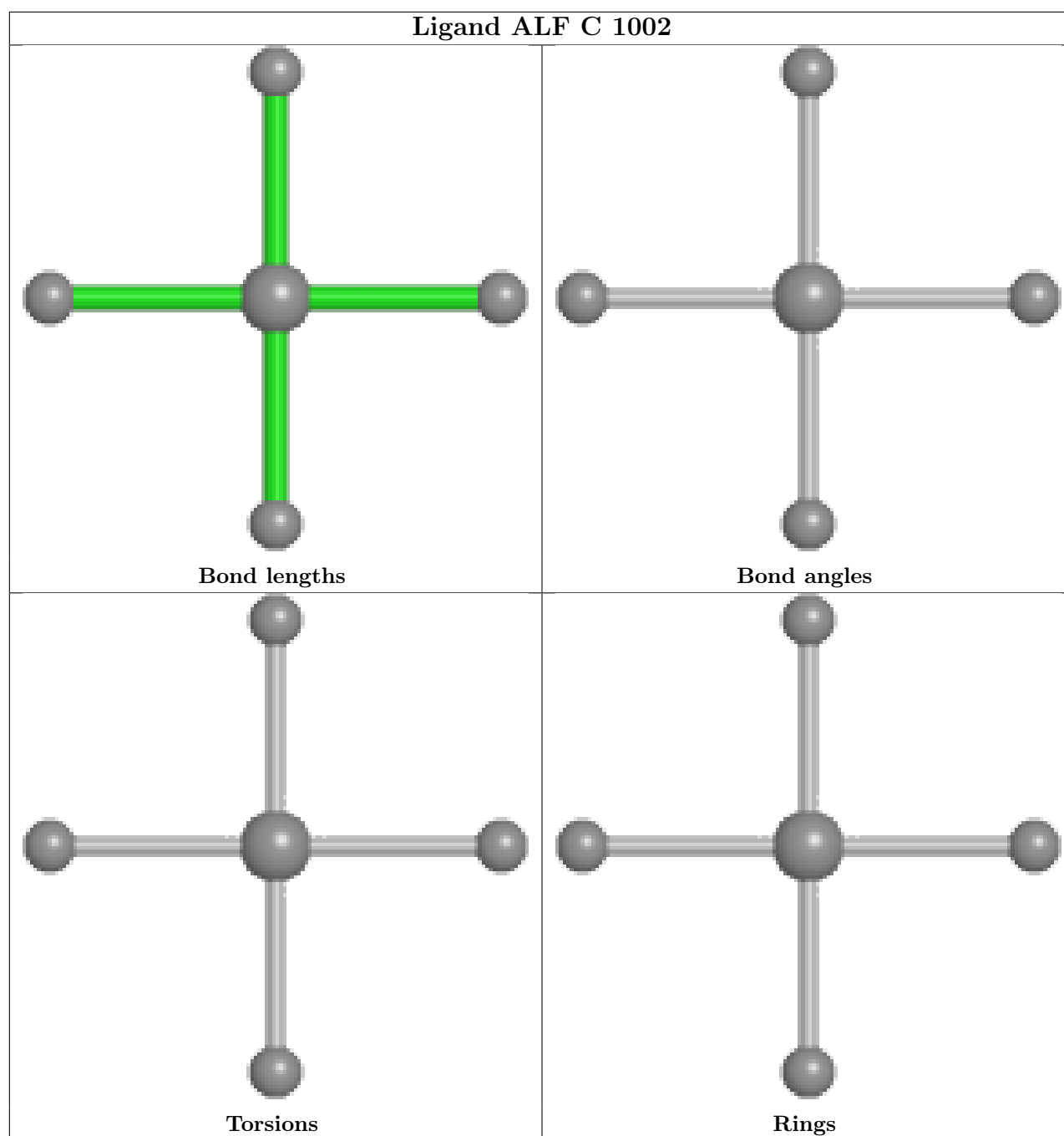
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1001	ADP	1	0
4	A	1002	ALF	1	0
4	C	1002	ALF	2	0

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	501/545 (91%)	-1.31	0 100 100	14, 35, 97, 122	0
1	C	490/545 (89%)	-1.30	0 100 100	18, 35, 98, 119	0
2	B	34/34 (100%)	-1.04	0 100 100	31, 52, 90, 108	0
2	D	33/34 (97%)	-1.07	0 100 100	23, 53, 99, 132	0
All	All	1058/1158 (91%)	-1.29	0 100 100	14, 36, 98, 132	0

There are no RSRZ outliers to report.

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(Å ²)	Q<0.9
5	MG	C	1003	1/1	0.99	0.10	30,30,30,30	0
3	ADP	C	1001	27/27	1.00	0.03	33,41,53,60	0
4	ALF	A	1002	5/5	1.00	0.03	20,21,23,23	0
4	ALF	C	1002	5/5	1.00	0.03	26,28,30,30	0

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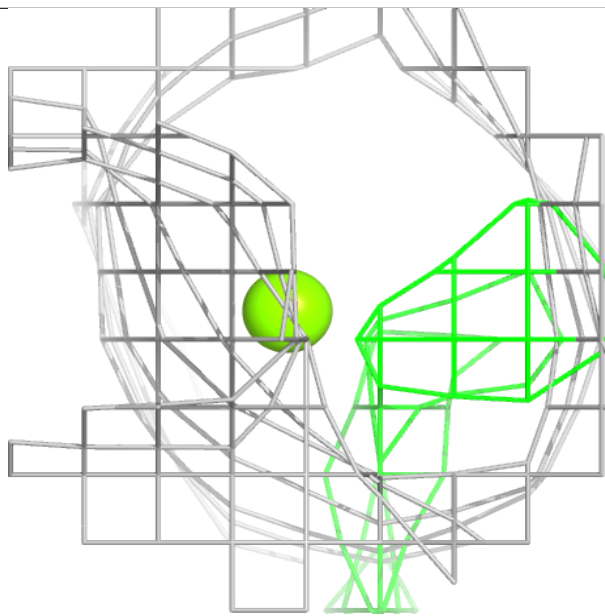
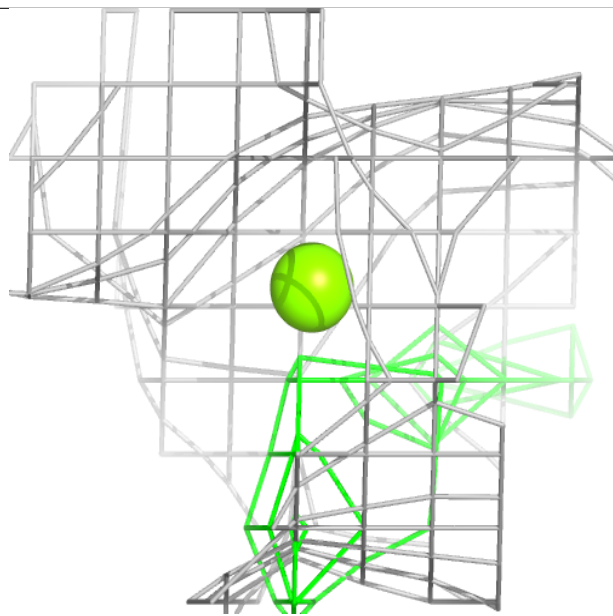
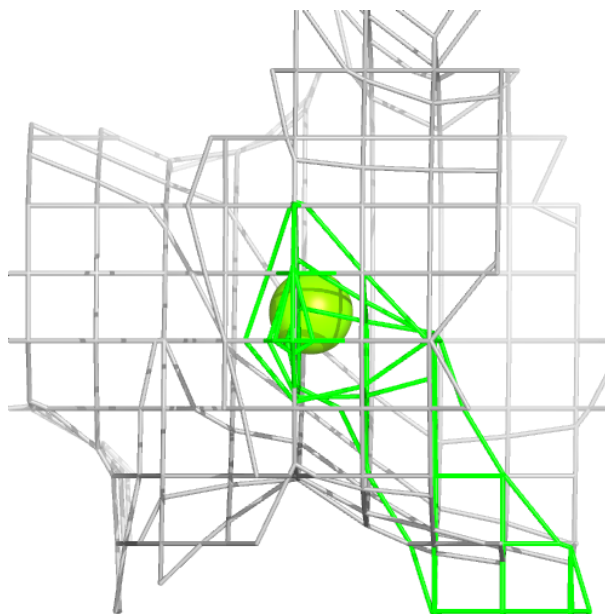
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	MG	A	1003	1/1	1.00	0.02	19,19,19,19	0
3	ADP	A	1001	27/27	1.00	0.03	29,37,50,56	0
6	K	D	2001	1/1	1.00	0.04	35,35,35,35	0
6	K	D	2002	1/1	1.00	0.02	50,50,50,50	0
6	K	D	2003	1/1	1.00	0.01	49,49,49,49	0
6	K	B	101	1/1	1.00	0.02	42,42,42,42	0
6	K	B	102	1/1	1.00	0.02	37,37,37,37	0
6	K	B	103	1/1	1.00	0.02	48,48,48,48	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.

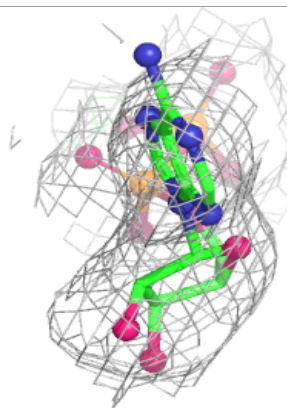
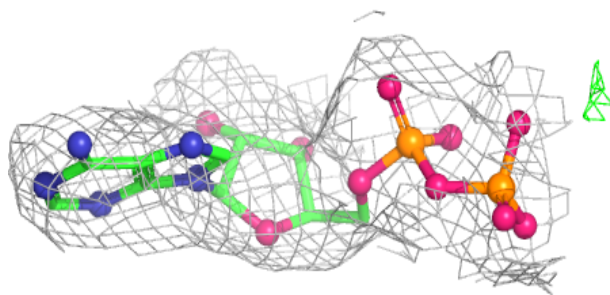
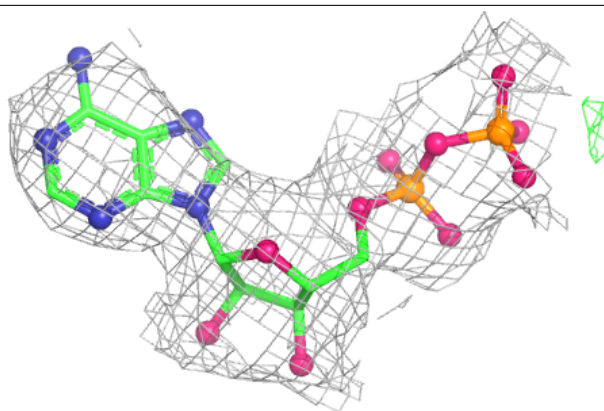
Electron density around MG C 1003:

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



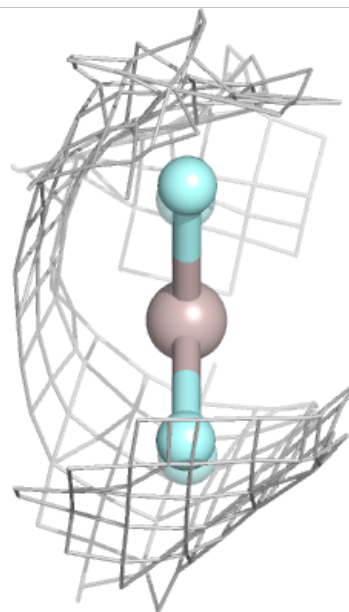
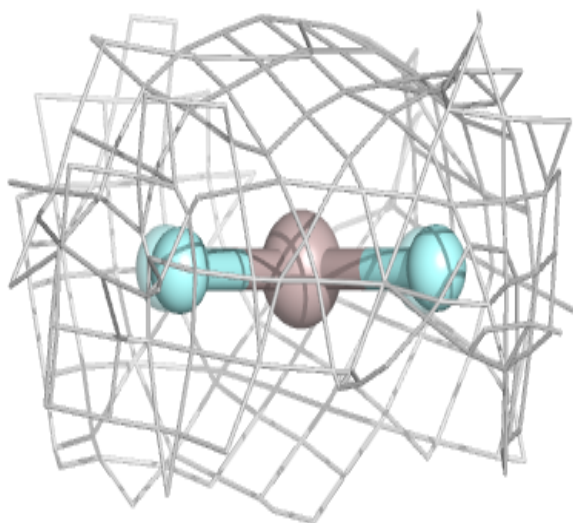
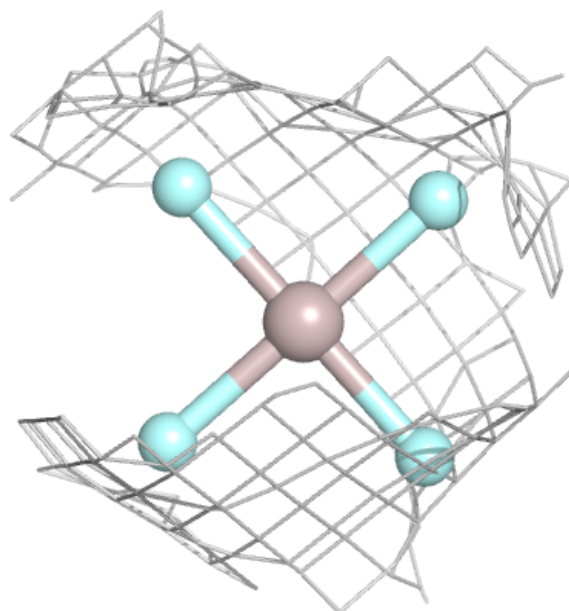
Electron density around ADP C 1001:

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



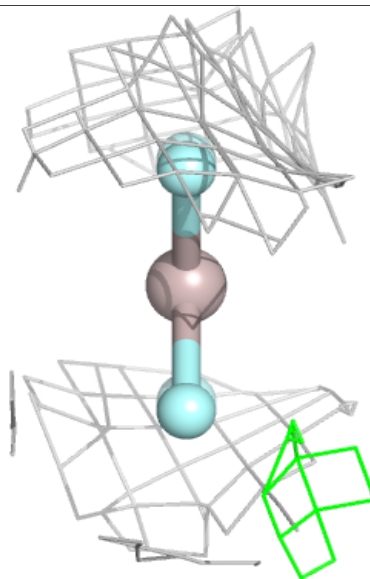
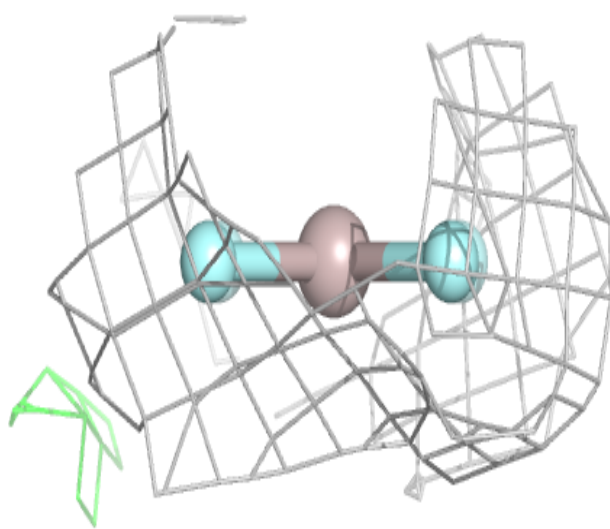
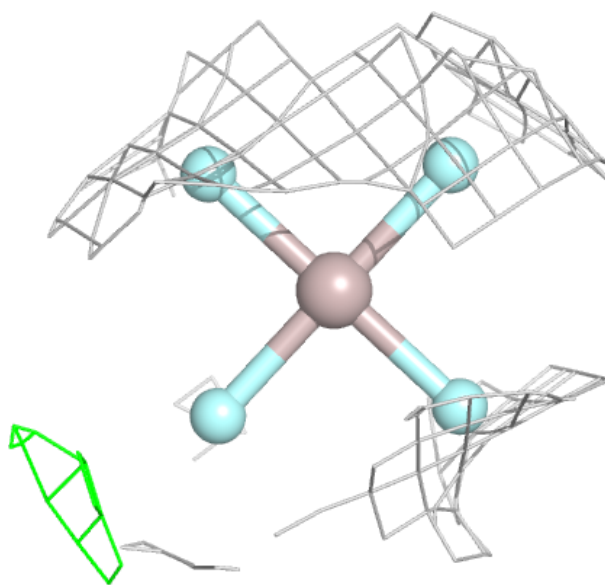
Electron density around ALF A 1002:

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



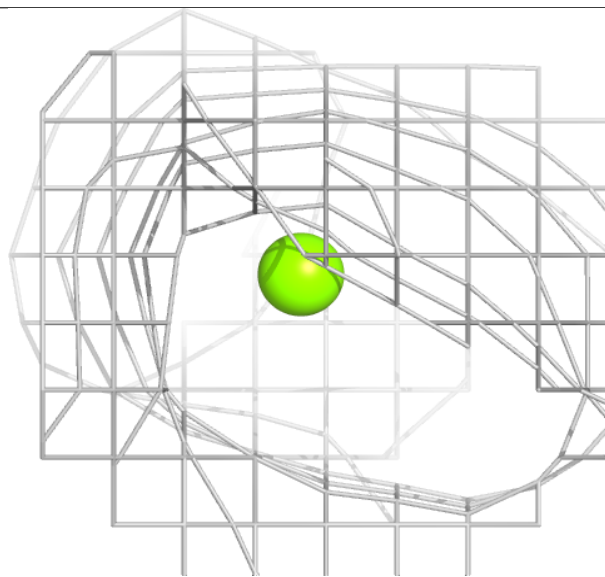
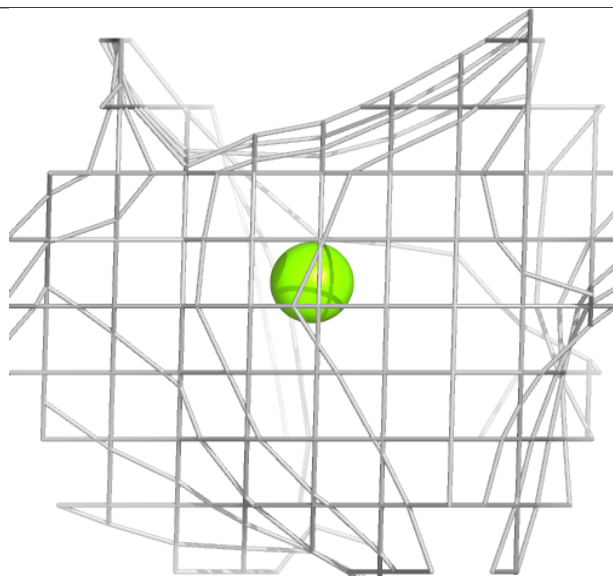
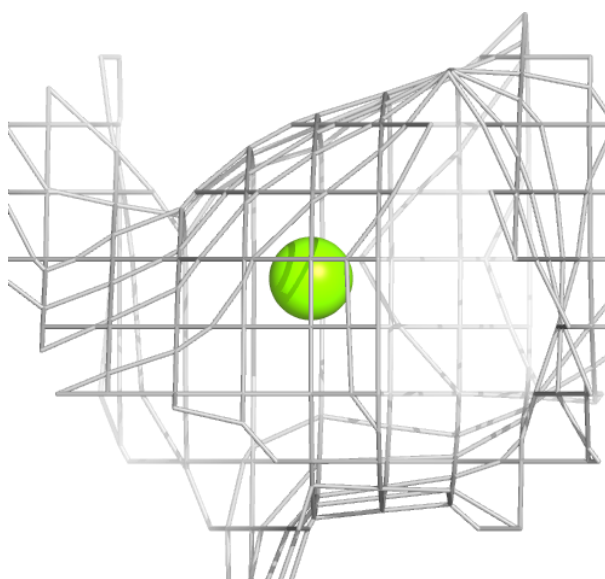
Electron density around ALF C 1002:

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



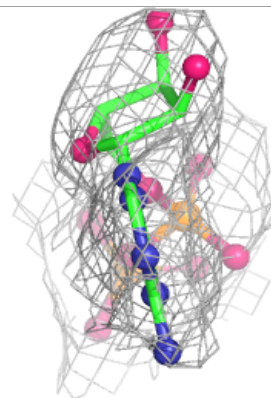
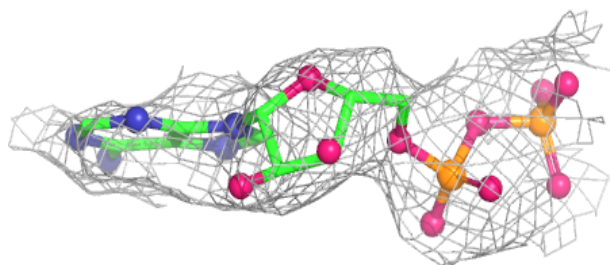
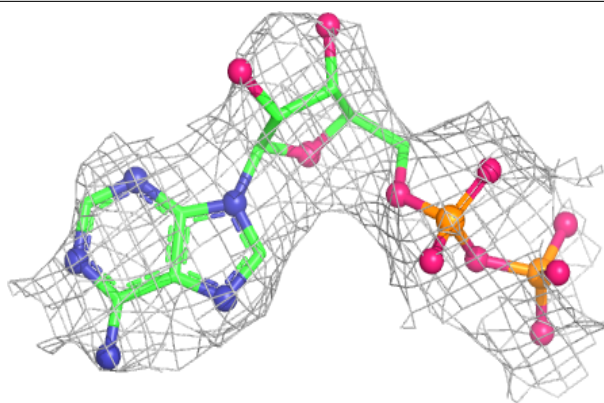
Electron density around MG A 1003:

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



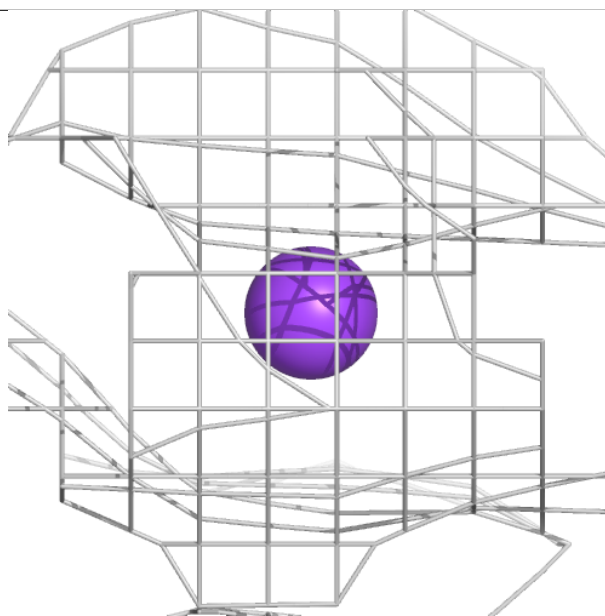
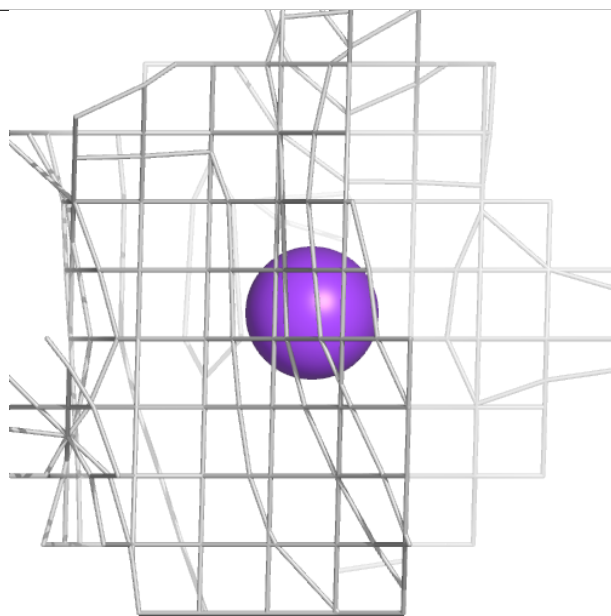
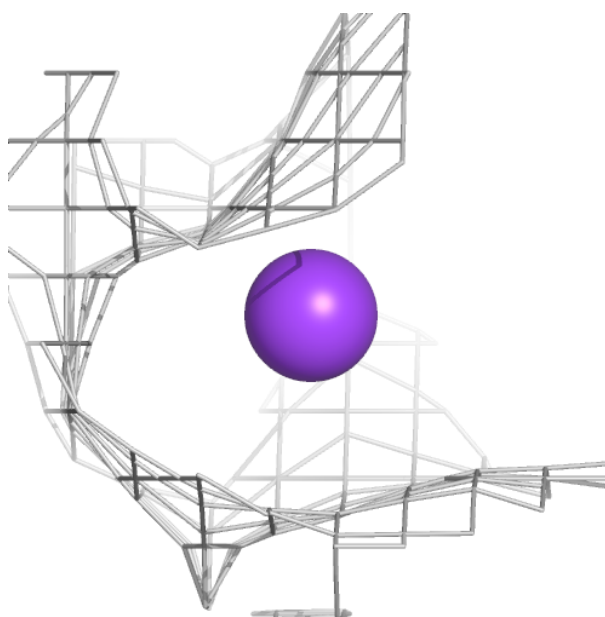
Electron density around ADP A 1001:

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



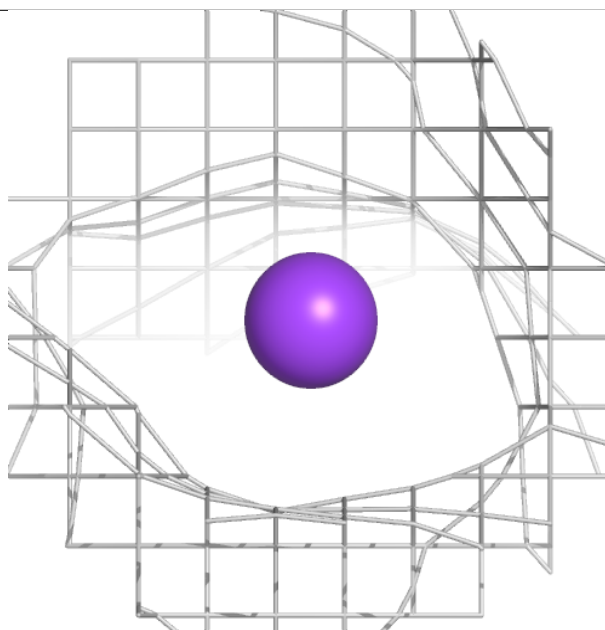
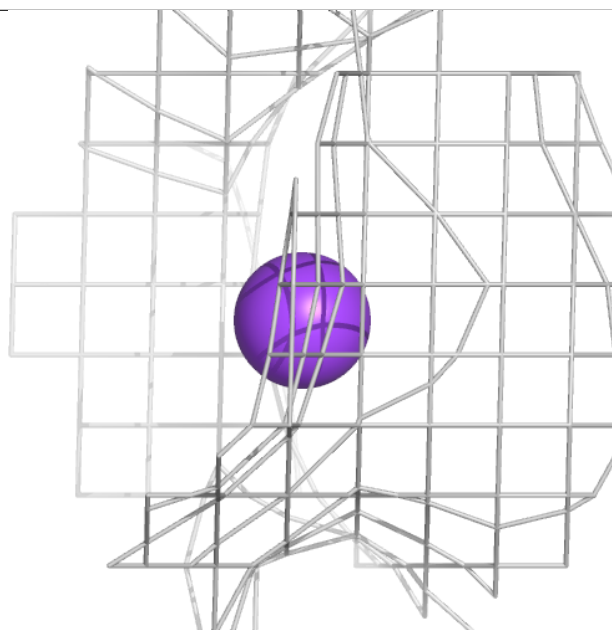
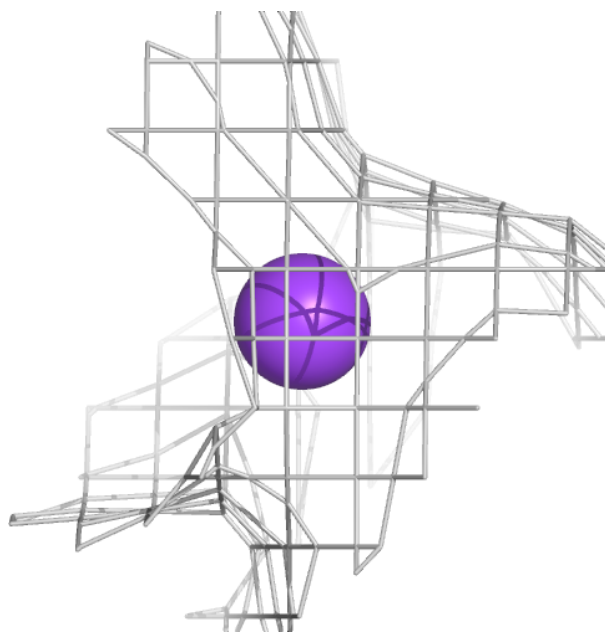
Electron density around K D 2001:

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



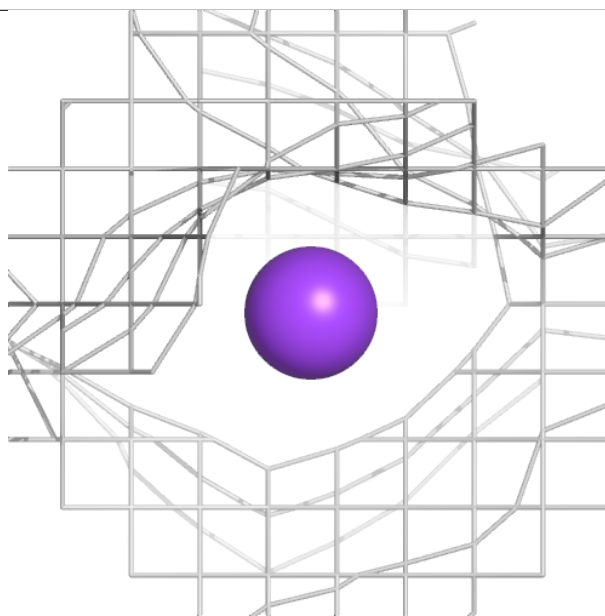
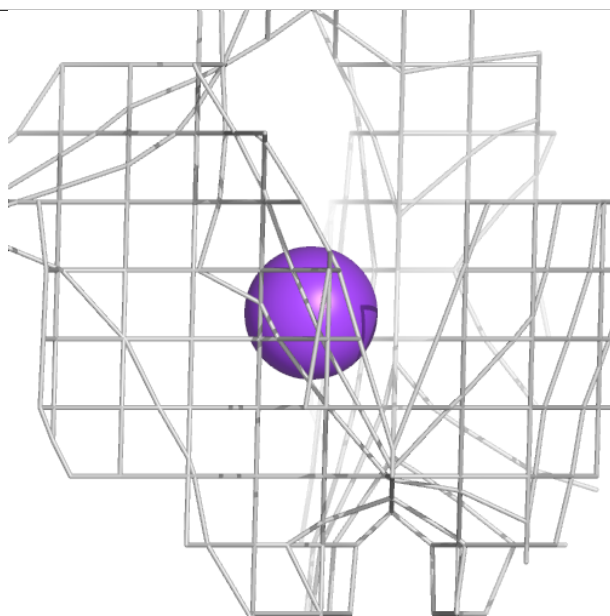
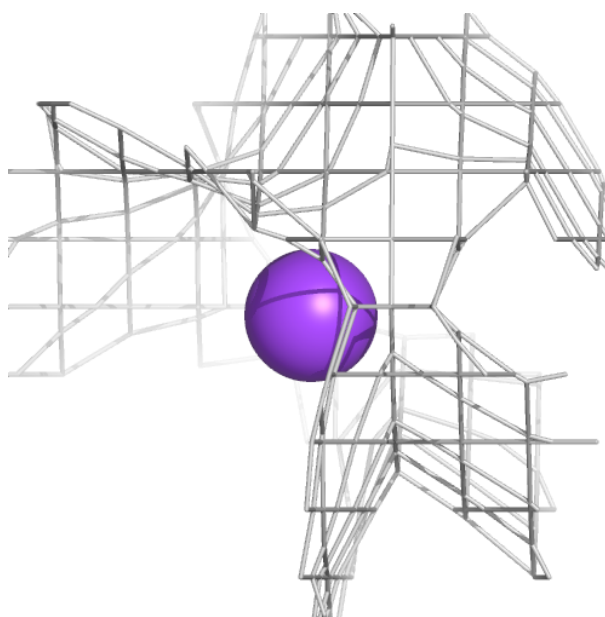
Electron density around K D 2002:

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



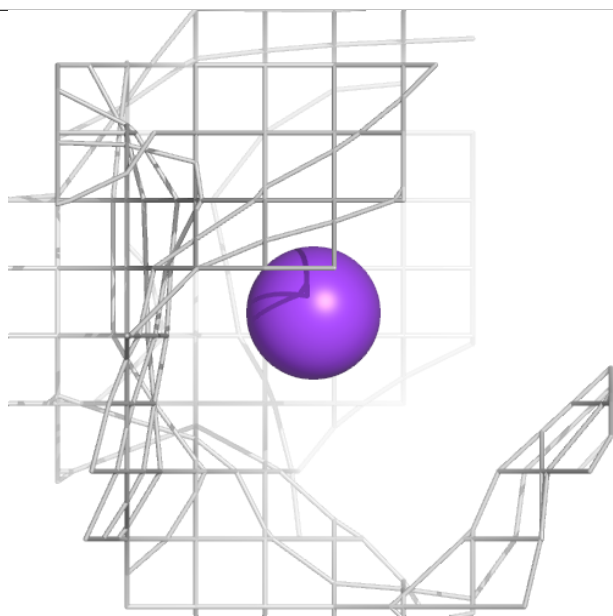
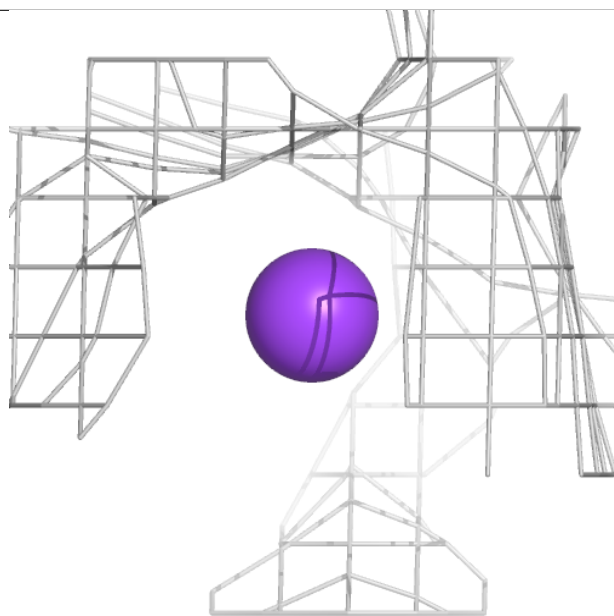
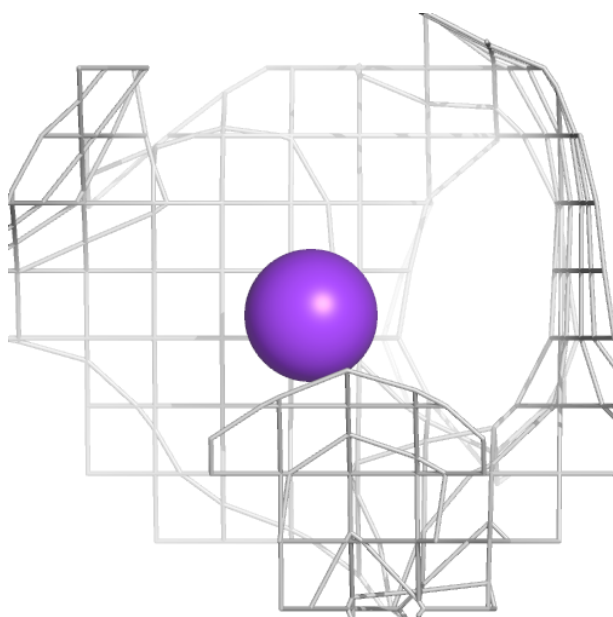
Electron density around K D 2003:

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



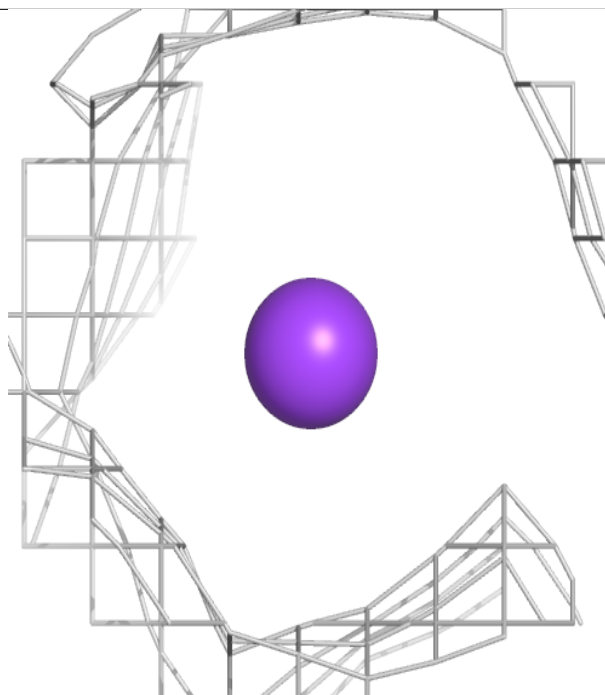
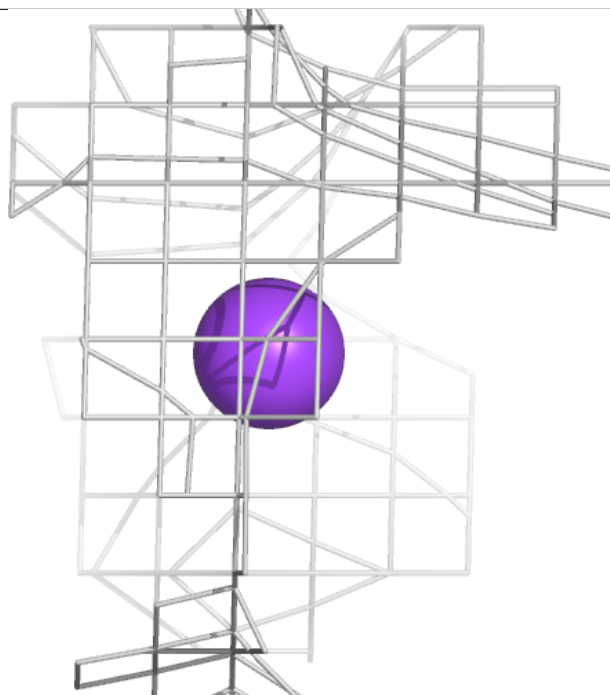
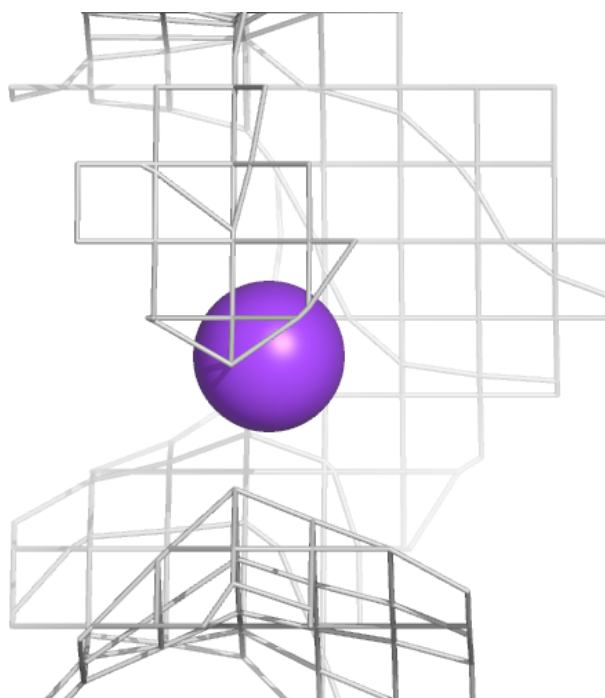
Electron density around K B 101:

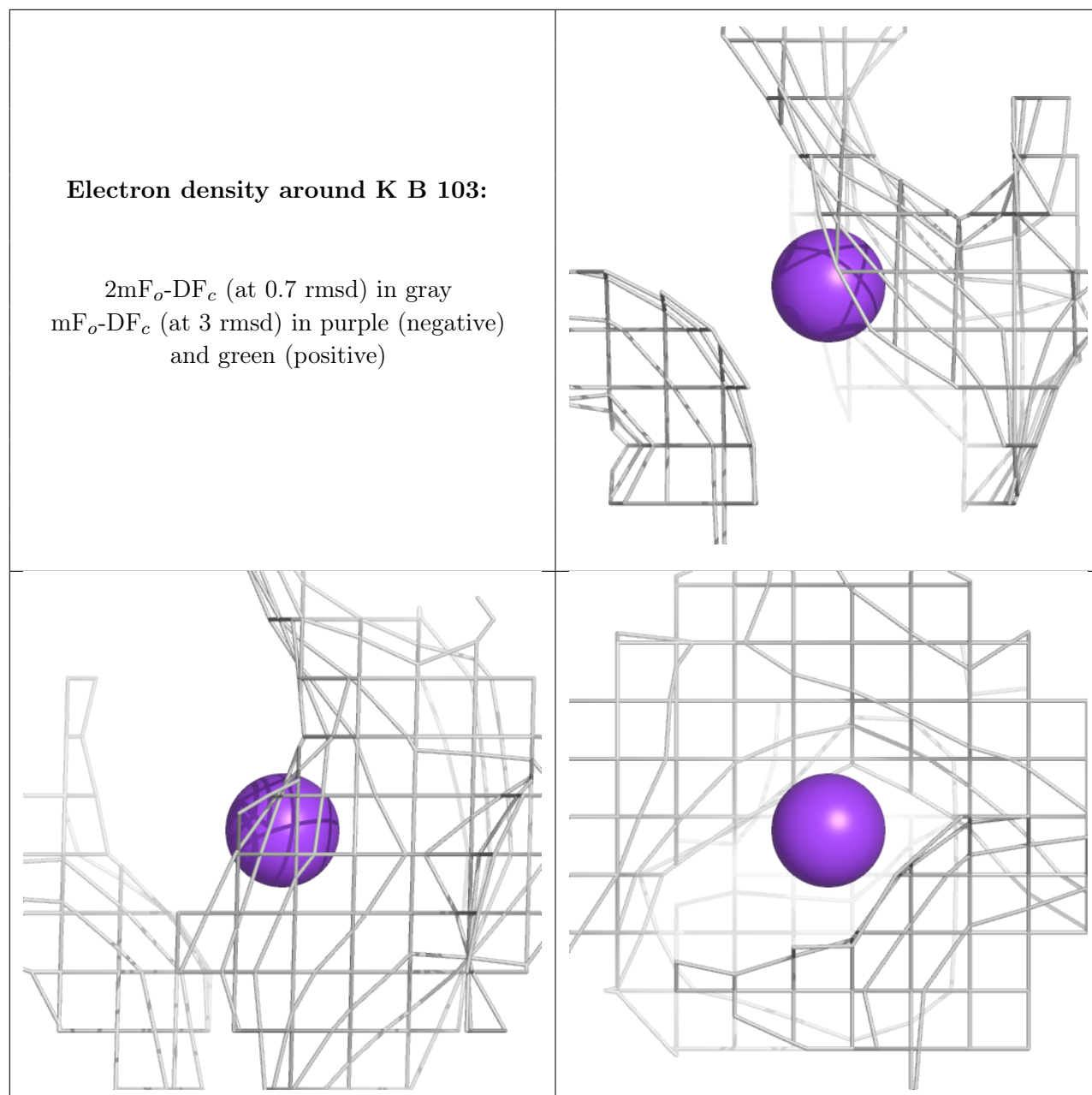
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around K B 102:

$2mF_o-DF_c$ (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.