



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 6, 2025 – 10:16 AM JST

PDB ID : 9JVT / pdb_00009jvt
Title : Structure of the C-terminal 4 domains (V4-V6-HP) villin bound to an actin
Authors : Robinson, R.C.
Deposited on : 2024-10-09
Resolution : 3.29 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 2.0
EDS : 3.0
buster-report : 1.1.7 (2018)
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
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) : 2.46

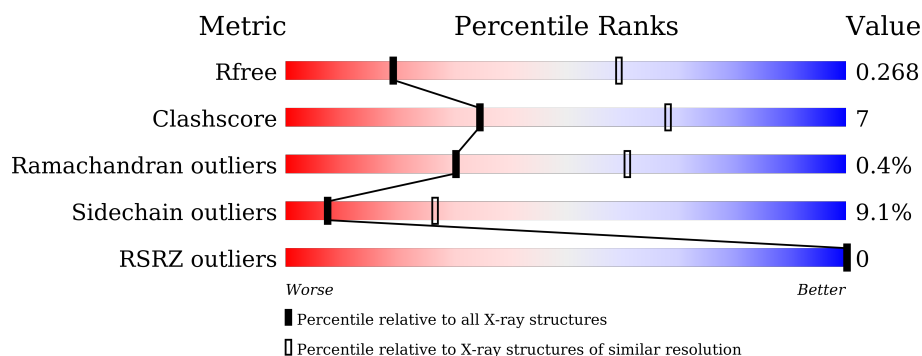
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.29 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	164625	1085 (3.32-3.28)
Clashscore	180529	1128 (3.32-3.28)
Ramachandran outliers	177936	1125 (3.32-3.28)
Sidechain outliers	177891	1124 (3.32-3.28)
RSRZ outliers	164620	1085 (3.32-3.28)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	V	485	<div> <div>58%</div> <div>19%</div> <div>•</div> <div>21%</div> </div>
1	v	485	<div> <div>68%</div> <div>20%</div> <div>•</div> <div>10%</div> </div>
2	G	377	<div> <div>72%</div> <div>21%</div> <div>•</div> <div>5%</div> </div>
2	g	377	<div> <div>71%</div> <div>19%</div> <div>•</div> <div>7%</div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 12102 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 Villin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	v	435	Total	C	N	O	S	0	0	0
			3470	2195	577	680	18			
1	V	381	Total	C	N	O	S	0	0	0
			3018	1897	510	595	16			

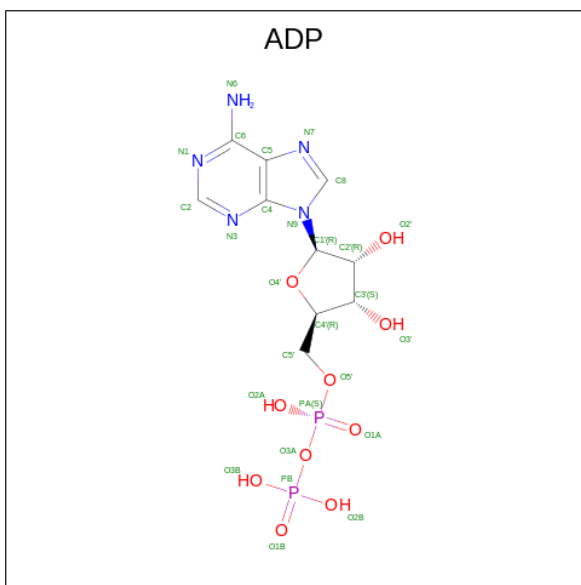
- Molecule 2 is a protein called Actin, alpha skeletal muscle.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	g	351	Total	C	N	O	S	0	0	0
			2743	1739	451	535	18			
2	G	359	Total	C	N	O	S	0	0	0
			2807	1779	469	540	19			

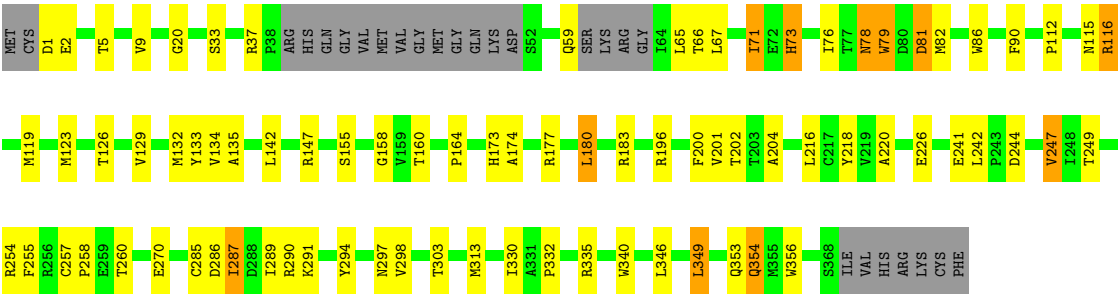
- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	v	4	Total	Ca	0	0
			4	4		
3	g	1	Total	Ca	0	0
			1	1		
3	V	4	Total	Ca	0	0
			4	4		
3	G	1	Total	Ca	0	0
			1	1		

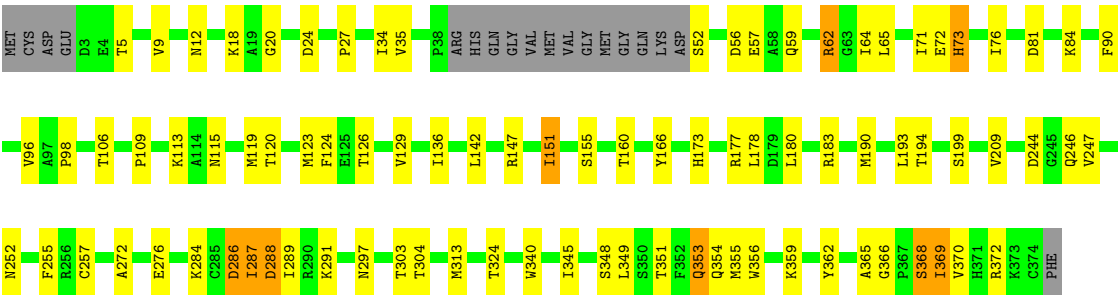
- Molecule 4 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: C₁₀H₁₅N₅O₁₀P₂) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	g	1	Total 27	C 10	N 5	O 10	P 2	0	0
4	G	1	Total 27	C 10	N 5	O 10	P 2	0	0



● Molecule 2: Actin, alpha skeletal muscle



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	131.12Å 97.69Å 134.31Å 90.00° 103.95° 90.00°	Depositor
Resolution (Å)	46.62 – 3.29 46.62 – 3.29	Depositor EDS
% Data completeness (in resolution range)	99.4 (46.62-3.29) 99.3 (46.62-3.29)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.35 (at 3.32Å)	Xtriage
Refinement program	PHENIX (1.17.1_3660: ???)	Depositor
R, R_{free}	0.241 , 0.268 0.242 , 0.268	Depositor DCC
R_{free} test set	2000 reflections (7.97%)	wwPDB-VP
Wilson B-factor (Å ²)	98.3	Xtriage
Anisotropy	0.349	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 80.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	12102	wwPDB-VP
Average B, all atoms (Å ²)	125.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.36% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: HIC, CA, ADP

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	V	0.13	0/3086	0.34	0/4163
1	v	0.11	0/3549	0.32	0/4778
2	G	0.13	0/2854	0.37	0/3868
2	g	0.14	0/2788	0.37	0/3781
All	All	0.13	0/12277	0.35	0/16590

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	V	3018	0	2879	45	0
1	v	3470	0	3326	52	0
2	G	2807	0	2774	44	0
2	g	2743	0	2694	42	0
3	G	1	0	0	0	0
3	V	4	0	0	0	0
3	g	1	0	0	0	0
3	v	4	0	0	0	0
4	G	27	0	12	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	g	27	0	12	0	0
All	All	12102	0	11697	176	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 7.

All (176) 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:V:413:LEU:HG	1:V:465:LYS:HD3	1.56	0.86
1:v:381:MET:HA	1:v:384:HIS:HB2	1.65	0.79
2:g:155:SER:HA	2:g:160:THR:HG22	1.65	0.77
2:g:285:CYS:HB3	2:g:289:ILE:HD11	1.68	0.76
1:v:630:CYS:HB2	1:v:655:VAL:HG23	1.71	0.73
2:G:155:SER:HA	2:G:160:THR:HG22	1.70	0.72
2:g:332:PRO:HG2	2:g:335:ARG:HB3	1.73	0.71
1:V:673:ARG:HB2	1:V:676:GLU:HB2	1.75	0.68
2:G:35:VAL:HG11	2:G:81:ASP:HB3	1.76	0.68
2:g:86:TRP:HH2	2:g:119:MET:HG2	1.59	0.67
2:g:353:GLN:HA	2:g:356:TRP:CD1	2.32	0.65
1:v:673:ARG:NH1	1:v:675:ASP:OD1	2.29	0.64
1:V:566:LYS:HB2	1:V:616:LYS:HA	1.79	0.64
2:g:76:ILE:HD13	2:g:82:MET:HG2	1.80	0.63
2:g:353:GLN:HA	2:g:356:TRP:HD1	1.64	0.63
1:V:450:LEU:HD22	1:V:459:ARG:HG2	1.81	0.63
1:v:551:VAL:HG12	1:v:564:ALA:HA	1.81	0.62
2:G:72:GLU:O	2:G:183:ARG:NH1	2.30	0.62
2:G:353:GLN:HA	2:G:356:TRP:HD1	1.65	0.62
1:v:635:GLY:HA3	1:V:373:THR:HG21	1.80	0.62
1:V:548:SER:HB2	1:V:615:LEU:HB2	1.83	0.61
2:g:142:LEU:HD22	2:g:147:ARG:HB2	1.83	0.60
1:V:623:ARG:HD3	1:V:624:PRO:HD2	1.84	0.59
1:V:733:GLU:HA	1:V:736:LYS:HE2	1.85	0.59
2:g:180:LEU:HD11	2:g:260:THR:HG22	1.83	0.58
1:v:526:VAL:HB	1:v:551:VAL:HG23	1.85	0.58
2:g:116:ARG:HG3	2:g:134:VAL:HG21	1.87	0.57
1:V:408:ILE:HG21	1:V:458:GLU:HA	1.87	0.56
1:v:547:ASN:HA	1:v:610:ALA:HB3	1.87	0.54
2:G:288:ASP:OD1	2:G:288:ASP:N	2.38	0.54
2:g:123:MET:HE3	2:g:129:VAL:HG11	1.88	0.54
2:G:190:MET:HG2	2:G:209:VAL:HG21	1.89	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:g:79:TRP:CH2	2:g:119:MET:HG3	2.44	0.53
1:V:470:ASP:OD2	1:V:479:GLN:NE2	2.41	0.53
1:V:549:ASN:H	1:V:615:LEU:HB3	1.72	0.53
2:G:123:MET:HE3	2:G:129:VAL:HG11	1.90	0.53
1:V:526:VAL:O	1:V:550:ASP:HB3	2.09	0.53
2:g:79:TRP:HH2	2:g:119:MET:HG3	1.73	0.52
2:g:242:LEU:HB2	2:g:244:ASP:OD1	2.10	0.52
1:v:758:MET:HB3	1:v:762:ASP:HB2	1.91	0.52
1:v:428:ASP:CG	1:v:506:LYS:H	2.17	0.52
1:v:406:TYR:HB3	1:v:413:LEU:HD22	1.92	0.52
1:v:739:GLY:O	1:v:741:LYS:N	2.36	0.52
1:V:407:ARG:HG2	1:V:409:GLU:HG3	1.92	0.51
1:V:745:MET:SD	1:V:745:MET:N	2.84	0.51
1:v:573:ARG:O	1:v:577:LYS:HG3	2.10	0.51
1:v:687:TYR:HD2	1:V:380:ALA:HB1	1.76	0.51
1:V:645:PHE:H	1:V:645:PHE:HD1	1.59	0.50
1:v:734:LEU:HB3	1:v:743:MET:SD	2.52	0.50
1:V:660:ALA:O	1:V:662:HIS:N	2.44	0.50
2:G:287:ILE:HD13	2:G:288:ASP:OD1	2.12	0.50
1:V:707:GLY:N	1:V:709:GLU:OE1	2.41	0.50
1:v:626:ARG:HH12	1:V:386:ASN:HB3	1.77	0.50
1:v:452:ARG:HD2	1:v:724:LEU:HD11	1.94	0.50
1:V:417:GLU:HB2	1:V:420:TYR:HD2	1.76	0.49
1:v:688:LEU:HD11	1:v:697:PRO:HA	1.92	0.49
2:g:133:TYR:HE2	2:g:135:ALA:HB2	1.78	0.49
1:V:503:GLN:H	1:V:538:GLN:HB3	1.78	0.49
2:G:52:SER:HB2	2:G:84:LYS:HD3	1.95	0.49
2:G:142:LEU:HD22	2:G:147:ARG:HB2	1.93	0.49
1:v:492:MET:HE2	1:v:533:ASN:HA	1.93	0.49
2:g:216:LEU:O	2:g:254:ARG:NE	2.33	0.49
1:V:610:ALA:HB1	1:V:615:LEU:HD13	1.95	0.49
2:G:120:THR:OG1	2:G:370:VAL:HG11	2.12	0.49
1:v:526:VAL:O	1:v:550:ASP:HB3	2.12	0.49
1:V:497:GLY:O	1:V:534:THR:OG1	2.31	0.49
1:v:402:LYS:HZ3	1:v:435:THR:HG22	1.78	0.48
2:G:178:LEU:HG	2:G:180:LEU:H	1.78	0.48
2:G:194:THR:HG22	2:G:199:SER:HA	1.95	0.48
2:G:365:ALA:O	2:G:372:ARG:NH2	2.46	0.48
2:G:353:GLN:HA	2:G:356:TRP:CD1	2.47	0.48
2:g:286:ASP:O	2:g:290:ARG:HG3	2.14	0.48
2:g:78:ASN:N	2:g:78:ASN:OD1	2.45	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:g:90:PHE:CE2	2:g:123:MET:HE1	2.49	0.47
1:V:543:ALA:HB3	1:V:607:THR:HG22	1.96	0.47
2:G:173:HIS:O	2:G:284:LYS:NZ	2.40	0.47
1:v:731:TYR:CE1	1:v:735:LYS:HD2	2.50	0.47
2:g:20:GLY:N	2:g:340:TRP:HE1	2.12	0.47
1:v:521:THR:HA	1:v:555:PHE:O	2.15	0.47
1:v:791:THR:O	1:v:795:PHE:HB2	2.14	0.47
2:G:24:ASP:HB2	2:G:340:TRP:HH2	1.78	0.47
2:G:252:ASN:HA	2:G:255:PHE:CZ	2.50	0.47
2:g:112:PRO:HB2	2:g:115:ASN:HB2	1.97	0.46
1:V:434:TYR:CD2	1:V:494:MET:HE3	2.50	0.46
1:v:759:SER:O	1:v:763:VAL:HG23	2.15	0.46
1:v:770:VAL:HG13	1:v:778:PRO:HD2	1.95	0.46
2:G:109:PRO:HA	2:G:136:ILE:HG23	1.97	0.46
1:v:407:ARG:NH2	1:v:428:ASP:OD2	2.49	0.46
1:v:543:ALA:HB3	1:v:607:THR:HG22	1.97	0.46
2:g:257:CYS:HB3	2:g:258:PRO:HD3	1.98	0.46
1:V:405:ILE:HG12	1:V:432:ILE:HG23	1.98	0.46
2:G:155:SER:HB3	2:G:304:THR:HG23	1.96	0.46
2:G:18:LYS:HE2	2:G:18:LYS:HB2	1.72	0.46
1:v:408:ILE:HD12	1:v:461:VAL:HG12	1.98	0.46
2:G:365:ALA:HB1	2:G:372:ARG:NH2	2.31	0.46
2:G:59:GLN:O	2:G:62:ARG:NE	2.49	0.45
1:v:627:LEU:HD11	1:v:656:PHE:HB3	1.98	0.45
2:G:20:GLY:N	2:G:340:TRP:HE1	2.15	0.45
1:v:701:ILE:HD13	1:v:748:VAL:HB	1.98	0.45
1:v:448:TYR:HB3	1:v:466:THR:HG21	1.98	0.45
2:G:12:ASN:ND2	2:G:106:THR:O	2.49	0.45
1:V:561:TYR:HB2	1:V:599:PHE:HE2	1.81	0.45
2:g:354:GLN:H	2:g:354:GLN:HG2	1.58	0.45
2:g:241:GLU:HG2	2:g:247:VAL:HG12	1.99	0.44
2:G:313:MET:HE3	2:G:313:MET:HB3	1.87	0.44
1:v:540:PRO:O	1:v:542:ARG:N	2.47	0.44
2:g:37:ARG:HB2	2:g:66:THR:OG1	2.18	0.44
1:V:537:GLU:HB2	1:V:719:VAL:HG11	1.99	0.44
1:V:623:ARG:HG3	1:V:695:ARG:HH22	1.82	0.44
2:G:27:PRO:HB3	2:G:340:TRP:CD1	2.53	0.44
2:g:202:THR:O	2:g:204:ALA:N	2.46	0.44
2:G:90:PHE:CZ	2:G:123:MET:HE1	2.53	0.44
1:V:629:GLN:O	1:V:639:ALA:HA	2.18	0.44
1:v:488:PRO:HG2	1:v:491:PHE:HB2	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:g:220:ALA:HB1	2:g:226:GLU:HG3	2.00	0.44
2:g:73:HIC:HB2	2:g:177:ARG:NH1	2.34	0.43
2:g:132:MET:HE2	2:g:132:MET:HB2	1.92	0.43
1:V:667:TRP:HA	1:V:704:VAL:HG13	1.99	0.43
2:G:124:PHE:CD1	2:G:359:LYS:HG2	2.53	0.43
2:g:287:ILE:H	2:g:287:ILE:HG12	1.62	0.43
1:v:521:THR:HG22	1:v:556:SER:HA	2.00	0.43
2:g:313:MET:HE3	2:g:313:MET:HB3	1.89	0.43
1:V:622:HIS:ND1	1:V:622:HIS:C	2.76	0.43
2:G:73:HIC:HB2	2:G:177:ARG:HH12	1.84	0.43
2:G:119:MET:O	2:G:123:MET:HG2	2.18	0.43
1:v:781:ILE:HA	1:v:789:HIS:CE1	2.54	0.43
1:v:763:VAL:HG21	1:v:788:LYS:HE2	2.01	0.43
1:v:432:ILE:HG21	1:v:494:MET:HE1	2.01	0.43
1:v:416:LEU:HD23	1:v:416:LEU:HA	1.83	0.43
2:G:286:ASP:OD1	2:G:286:ASP:N	2.50	0.43
1:v:512:HIS:HB3	1:v:513:ASN:H	1.67	0.42
2:g:71:ILE:HG23	2:g:76:ILE:HG12	2.01	0.42
1:V:395:MET:SD	1:V:395:MET:N	2.87	0.42
2:G:362:TYR:O	2:G:366:GLY:N	2.51	0.42
2:g:298:VAL:HA	2:g:330:ILE:HB	2.01	0.42
2:G:166:TYR:CD1	2:G:289:ILE:HG23	2.54	0.42
2:g:155:SER:OG	2:g:303:THR:HB	2.18	0.42
1:v:434:TYR:CE1	1:v:436:TYR:HB3	2.54	0.42
1:V:457:ASP:HA	2:G:345:ILE:HG12	2.02	0.42
1:V:542:ARG:NH1	1:V:730:SER:HB3	2.35	0.42
1:v:378:PHE:CE1	1:V:639:ALA:HB3	2.55	0.42
2:g:86:TRP:CH2	2:g:119:MET:HG2	2.47	0.42
2:g:287:ILE:HD13	2:g:290:ARG:CZ	2.50	0.42
1:v:405:ILE:HD11	1:v:495:PHE:HE1	1.84	0.42
2:G:354:GLN:HG3	2:G:355:MET:HG3	2.02	0.42
2:g:164:PRO:HG2	2:g:174:ALA:HB3	2.02	0.42
1:V:630:CYS:HB2	1:V:655:VAL:HG23	2.02	0.42
2:G:369:ILE:HA	2:G:372:ARG:HB2	2.01	0.42
2:g:346:LEU:O	2:g:349:LEU:HB2	2.20	0.42
1:V:407:ARG:HD3	1:V:430:TYR:CZ	2.55	0.42
2:g:78:ASN:HB3	2:g:81:ASP:HB2	2.02	0.41
1:V:411:MET:HE2	2:G:348:SER:HB2	2.02	0.41
1:V:619:GLU:HA	1:V:622:HIS:CD2	2.55	0.41
1:v:527:ARG:HG3	1:v:528:GLY:N	2.35	0.41
1:v:675:ASP:OD2	1:v:675:ASP:N	2.52	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:90:PHE:CG	2:G:98:PRO:HG3	2.55	0.41
1:v:444:HIS:CG	1:v:473:LEU:HD13	2.55	0.41
1:v:547:ASN:OD1	1:v:548:SER:N	2.52	0.41
2:G:151:ILE:HG23	2:G:297:ASN:HA	2.03	0.41
1:v:381:MET:HG3	1:V:690:SER:HB3	2.02	0.41
1:v:770:VAL:O	1:v:773:GLN:HG2	2.21	0.41
2:G:272:ALA:HB1	2:G:276:GLU:HB2	2.02	0.41
1:v:790:LEU:HD23	1:v:790:LEU:HA	1.86	0.41
1:v:819:LYS:HB2	1:v:819:LYS:HE3	1.73	0.41
1:V:516:ASP:OD1	1:V:516:ASP:N	2.54	0.41
1:V:548:SER:C	1:V:550:ASP:H	2.28	0.41
2:G:34:ILE:HD11	2:G:59:GLN:HE21	1.86	0.41
1:V:381:MET:O	1:V:385:ASN:N	2.54	0.41
2:G:52:SER:CB	2:G:84:LYS:HD3	2.51	0.41
2:g:218:TYR:O	2:g:255:PHE:HA	2.21	0.40
1:V:416:LEU:HD22	1:V:420:TYR:HB2	2.03	0.40
1:v:691:ASP:C	1:v:693:SER:H	2.29	0.40
2:G:193:LEU:HD23	2:G:193:LEU:HA	1.94	0.40
1:v:413:LEU:HA	1:v:413:LEU:HD23	1.84	0.40
2:g:158:GLY:HA3	2:g:183:ARG:NH1	2.36	0.40
2:g:290:ARG:O	2:g:294:TYR:N	2.52	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	V	379/485 (78%)	347 (92%)	30 (8%)	2 (0%)	25	56
1	v	429/485 (88%)	402 (94%)	25 (6%)	2 (0%)	25	56
2	G	354/377 (94%)	341 (96%)	12 (3%)	1 (0%)	37	66
2	g	344/377 (91%)	332 (96%)	11 (3%)	1 (0%)	37	66

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	1506/1724 (87%)	1422 (94%)	78 (5%)	6 (0%)	30 61

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	v	512	HIS
2	g	5	THR
2	G	368	SER
1	v	718	GLY
1	V	718	GLY
1	V	389	VAL

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	V	320/414 (77%)	285 (89%)	35 (11%)	5 20
1	v	370/414 (89%)	342 (92%)	28 (8%)	11 34
2	G	304/319 (95%)	275 (90%)	29 (10%)	7 25
2	g	297/319 (93%)	271 (91%)	26 (9%)	8 28
All	All	1291/1466 (88%)	1173 (91%)	118 (9%)	7 27

All (118) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	v	398	ASP
1	v	402	LYS
1	v	455	THR
1	v	516	ASP
1	v	531	GLN
1	v	539	VAL
1	v	551	VAL
1	v	578	LYS
1	v	620	ASN

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Mol	Chain	Res	Type
1	v	636	VAL
1	v	641	GLU
1	v	643	VAL
1	v	655	VAL
1	v	661	ASP
1	v	662	HIS
1	v	663	THR
1	v	668	LEU
1	v	675	ASP
1	v	688	LEU
1	v	699	THR
1	v	701	ILE
1	v	702	MET
1	v	743	MET
1	v	749	ARG
1	v	757	GLU
1	v	760	PHE
1	v	777	LEU
1	v	791	THR
2	g	1	ASP
2	g	2	GLU
2	g	9	VAL
2	g	33	SER
2	g	59	GLN
2	g	65	LEU
2	g	67	LEU
2	g	71	ILE
2	g	78	ASN
2	g	79	TRP
2	g	81	ASP
2	g	116	ARG
2	g	126	THR
2	g	173	HIS
2	g	180	LEU
2	g	196	ARG
2	g	200	PHE
2	g	201	VAL
2	g	247	VAL
2	g	249	THR
2	g	270	GLU
2	g	287	ILE
2	g	291	LYS

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Mol	Chain	Res	Type
2	g	297	ASN
2	g	349	LEU
2	g	354	GLN
1	V	374	VAL
1	V	391	LYS
1	V	392	GLU
1	V	393	THR
1	V	397	ASP
1	V	398	ASP
1	V	402	LYS
1	V	419	ARG
1	V	438	VAL
1	V	480	VAL
1	V	483	VAL
1	V	512	HIS
1	V	515	THR
1	V	539	VAL
1	V	556	SER
1	V	568	CYS
1	V	583	SER
1	V	609	TYR
1	V	612	ASP
1	V	615	LEU
1	V	622	HIS
1	V	636	VAL
1	V	645	PHE
1	V	655	VAL
1	V	663	THR
1	V	685	ILE
1	V	695	ARG
1	V	698	ASP
1	V	699	THR
1	V	702	MET
1	V	704	VAL
1	V	738	LEU
1	V	745	MET
1	V	747	GLN
1	V	749	ARG
2	G	5	THR
2	G	9	VAL
2	G	56	ASP
2	G	57	GLU

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Mol	Chain	Res	Type
2	G	62	ARG
2	G	64	ILE
2	G	65	LEU
2	G	71	ILE
2	G	76	ILE
2	G	96	VAL
2	G	113	LYS
2	G	115	ASN
2	G	126	THR
2	G	151	ILE
2	G	244	ASP
2	G	246	GLN
2	G	247	VAL
2	G	257	CYS
2	G	286	ASP
2	G	287	ILE
2	G	288	ASP
2	G	291	LYS
2	G	303	THR
2	G	324	THR
2	G	349	LEU
2	G	351	THR
2	G	353	GLN
2	G	368	SER
2	G	369	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
2	g	88	HIS
1	V	439	ASN
2	G	12	ASN
2	G	59	GLN
2	G	353	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

2 non-standard protein/DNA/RNA residues are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	HIC	G	73	2	8,11,12	1.68	2 (25%)	6,14,16	0.84	0
2	HIC	g	73	2	8,11,12	1.66	2 (25%)	6,14,16	0.98	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
2	HIC	G	73	2	-	1/5/6/8	0/1/1/1
2	HIC	g	73	2	-	0/5/6/8	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	G	73	HIC	CD2-CG	3.60	1.41	1.36
2	g	73	HIC	CD2-CG	3.60	1.41	1.36
2	G	73	HIC	CZ-NE2	-2.03	1.42	1.48
2	g	73	HIC	CZ-NE2	-2.02	1.42	1.48

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	G	73	HIC	O-C-CA-CB

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	G	73	HIC	1	0
2	g	73	HIC	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 12 ligands modelled in this entry, 10 are monoatomic - leaving 2 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$
4	ADP	G	401	3	24,29,29	0.93	1 (4%)	29,45,45	1.40	4 (13%)
4	ADP	g	401	3	24,29,29	0.94	1 (4%)	29,45,45	1.41	4 (13%)

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	ADP	G	401	3	-	0/12/32/32	0/3/3/3
4	ADP	g	401	3	-	0/12/32/32	0/3/3/3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	g	401	ADP	C5-C4	2.46	1.47	1.40
4	G	401	ADP	C5-C4	2.46	1.47	1.40

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	G	401	ADP	PA-O3A-PB	-3.29	121.52	132.83
4	g	401	ADP	N3-C2-N1	-3.11	123.81	128.68
4	g	401	ADP	C3'-C2'-C1'	2.98	105.46	100.98
4	G	401	ADP	N3-C2-N1	-2.97	124.03	128.68
4	g	401	ADP	PA-O3A-PB	-2.96	122.66	132.83
4	G	401	ADP	C3'-C2'-C1'	2.78	105.16	100.98
4	g	401	ADP	C4-C5-N7	-2.77	106.51	109.40
4	G	401	ADP	C4-C5-N7	-2.69	106.60	109.40

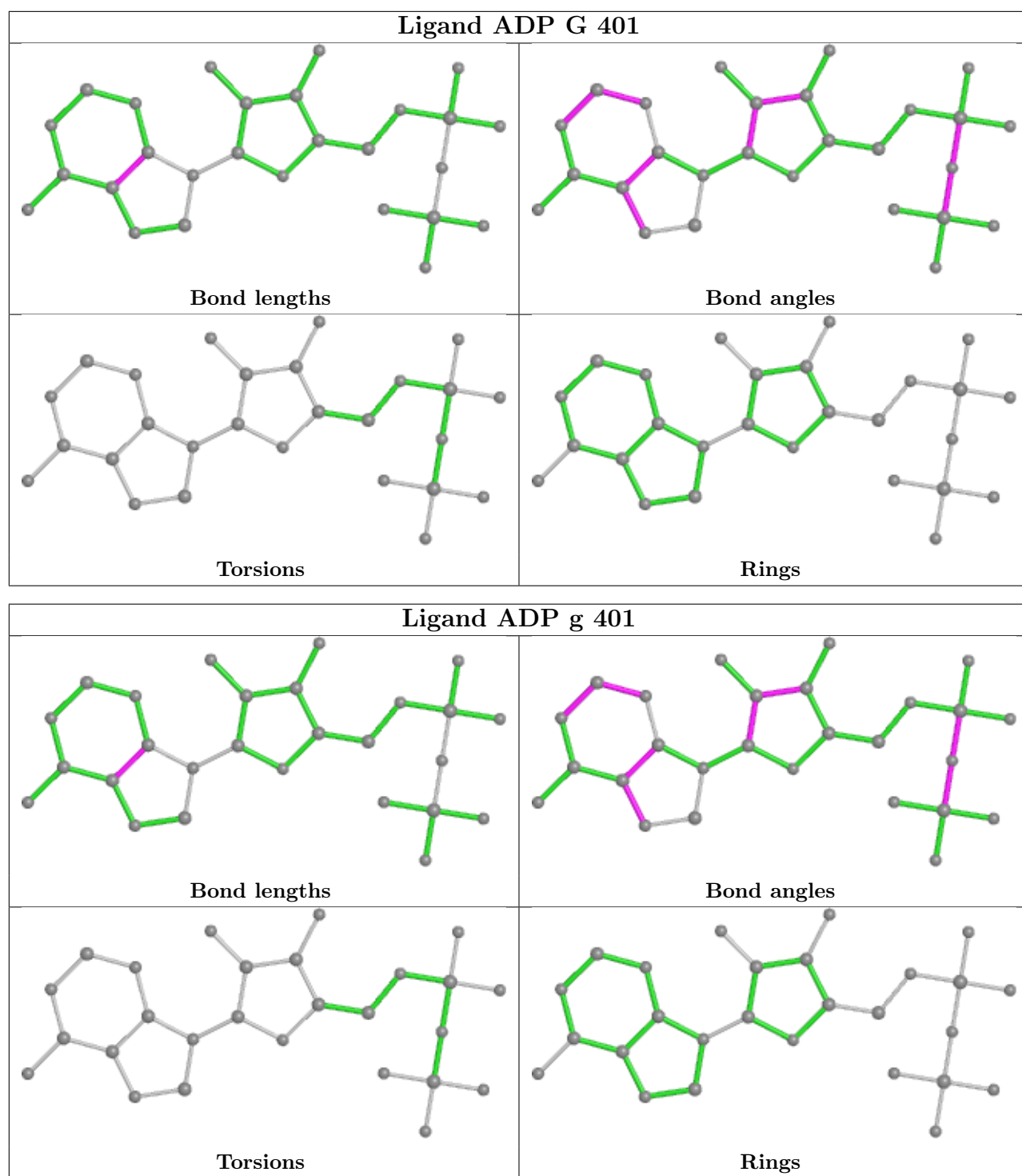
There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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 ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	V	381/485 (78%)	-0.67	0 100 100	69, 129, 196, 244	0
1	v	435/485 (89%)	-0.76	0 100 100	67, 115, 167, 196	0
2	G	358/377 (94%)	-0.62	0 100 100	82, 128, 178, 221	0
2	g	350/377 (92%)	-0.71	0 100 100	68, 113, 190, 255	0
All	All	1524/1724 (88%)	-0.69	0 100 100	67, 122, 186, 255	0

There are no RSRZ outliers to report.

6.2 Non-standard residues in protein, DNA, RNA chains [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
2	HIC	G	73	11/12	0.81	0.07	120,137,139,141	0
2	HIC	g	73	11/12	0.87	0.07	135,145,154,156	0

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

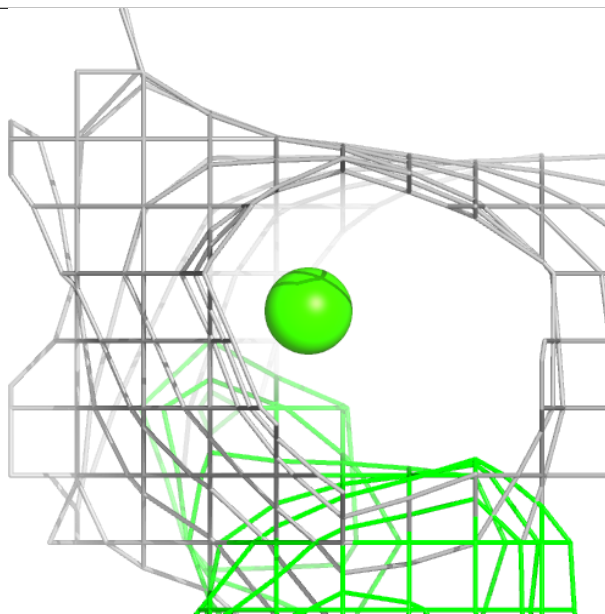
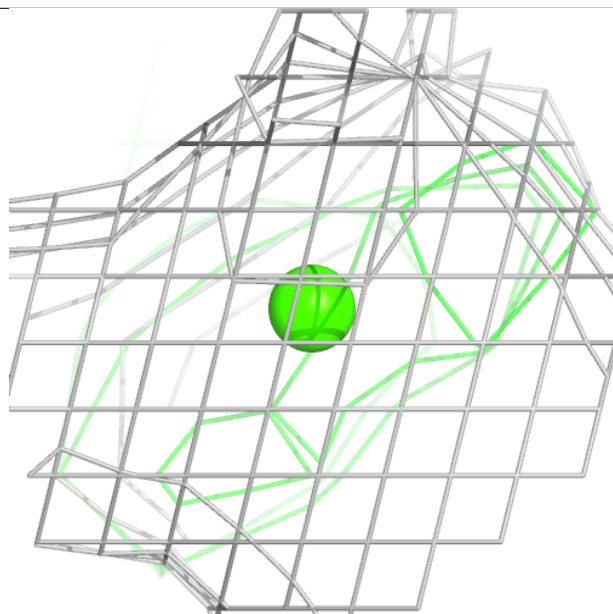
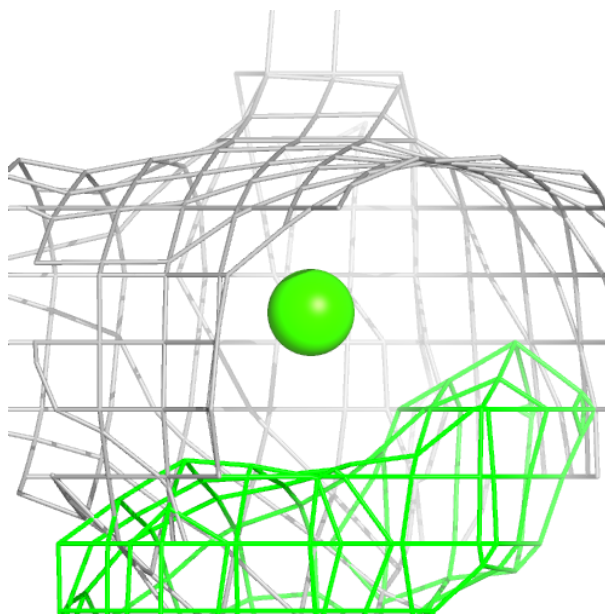
labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	CA	G	402	1/1	0.81	0.11	87,87,87,87	0
3	CA	V	901	1/1	0.84	0.09	203,203,203,203	0
4	ADP	G	401	27/27	0.85	0.09	82,97,106,106	0
3	CA	v	902	1/1	0.89	0.08	84,84,84,84	0
3	CA	v	901	1/1	0.91	0.08	100,100,100,100	0
3	CA	g	402	1/1	0.93	0.05	82,82,82,82	0
4	ADP	g	401	27/27	0.94	0.06	82,84,92,99	0
3	CA	v	903	1/1	0.94	0.05	105,105,105,105	0
3	CA	V	904	1/1	0.95	0.07	82,82,82,82	0
3	CA	V	902	1/1	0.96	0.05	102,102,102,102	0
3	CA	V	903	1/1	0.96	0.05	92,92,92,92	0
3	CA	v	904	1/1	0.98	0.06	82,82,82,82	0

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

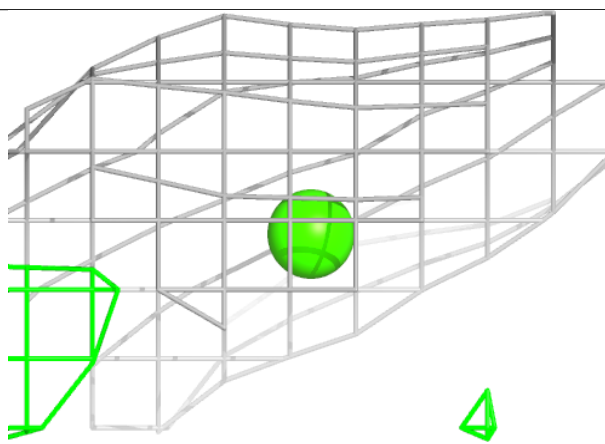
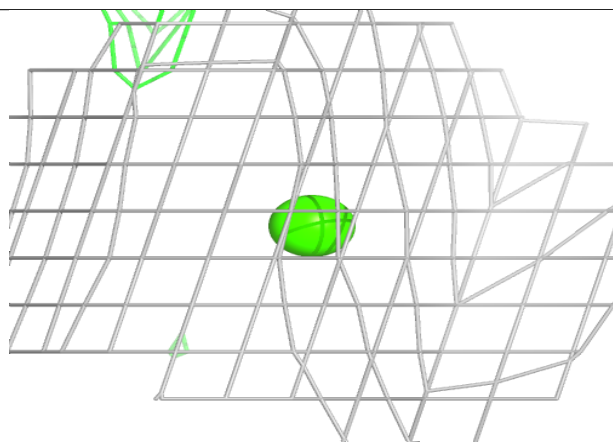
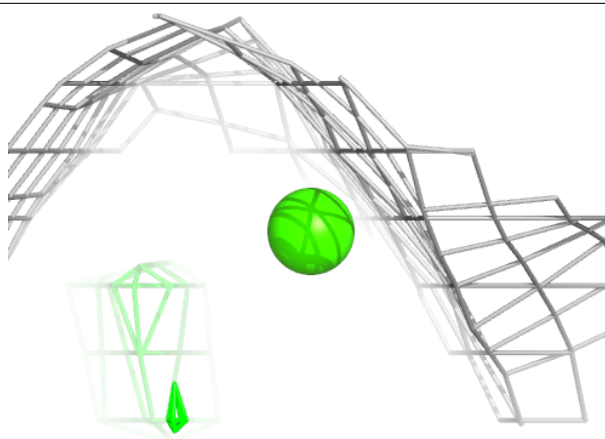
Electron density around CA G 402:

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

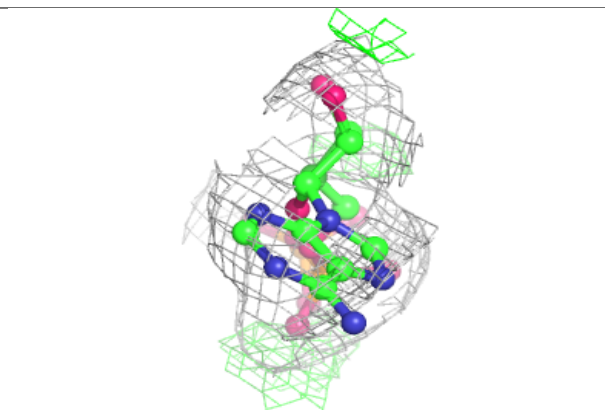
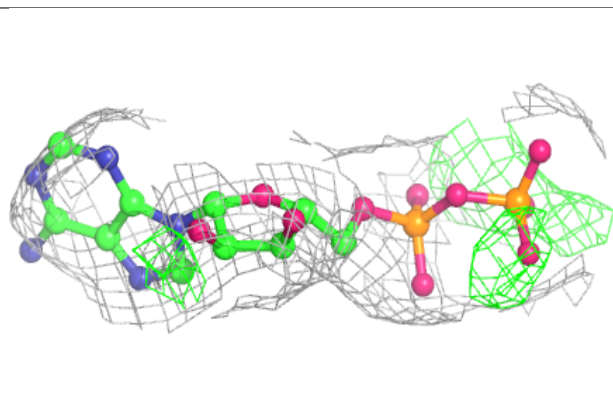
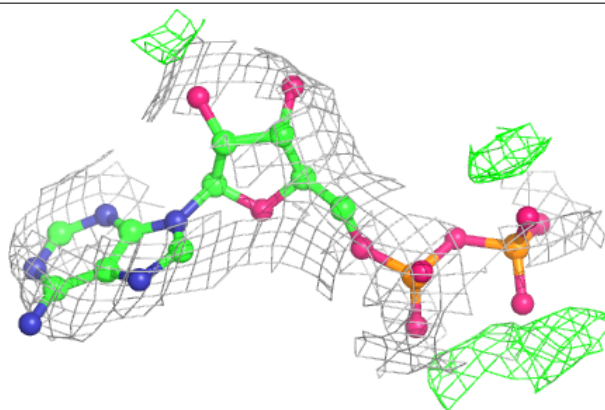


Electron density around CA V 901:

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

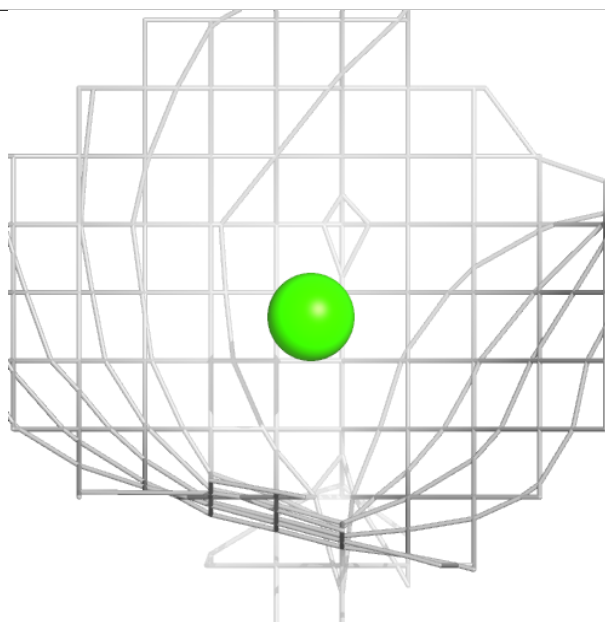
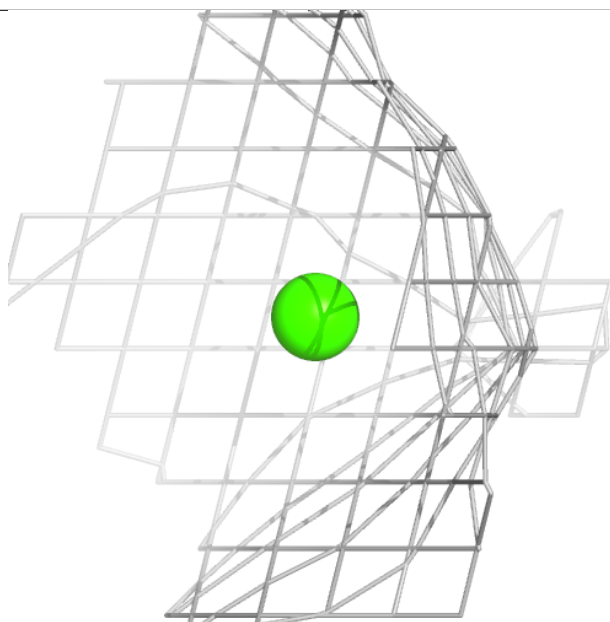
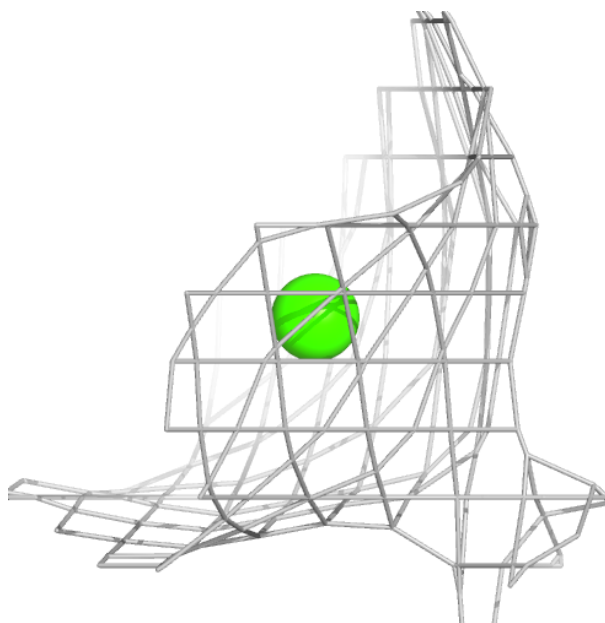
**Electron density around ADP G 401:**

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



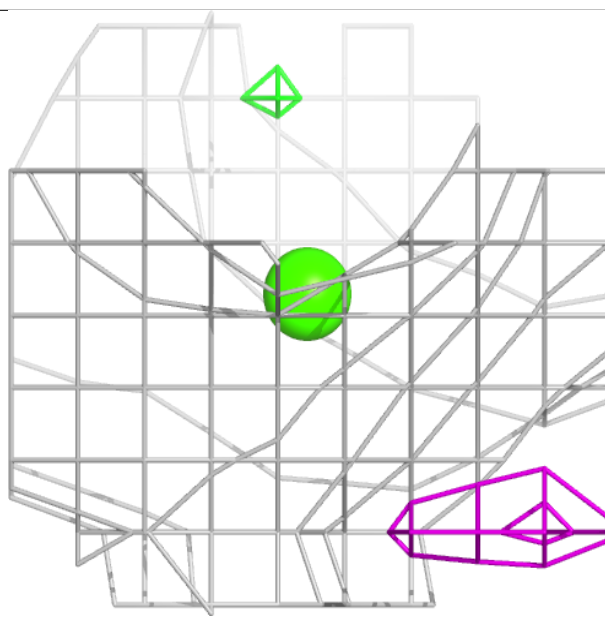
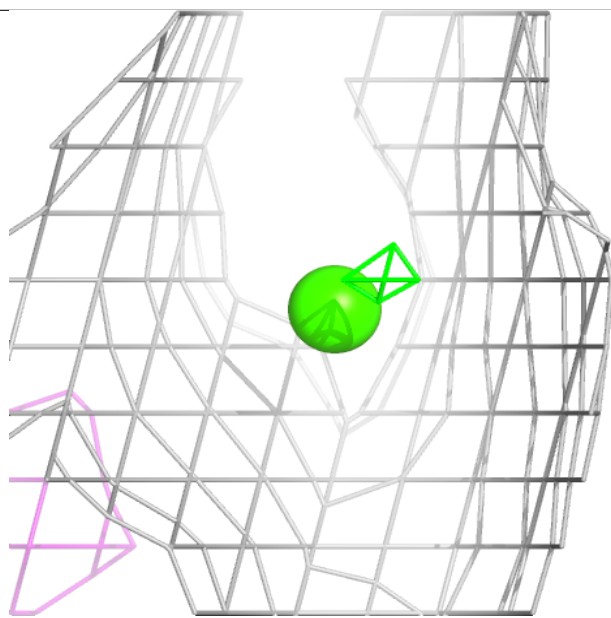
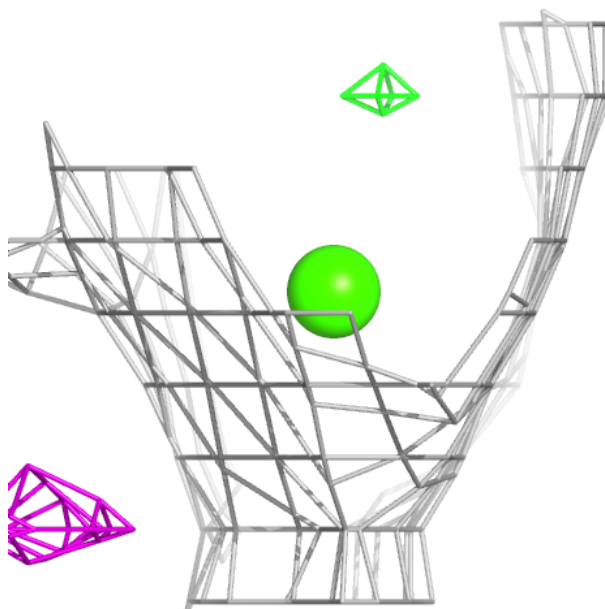
Electron density around CA v 902:

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



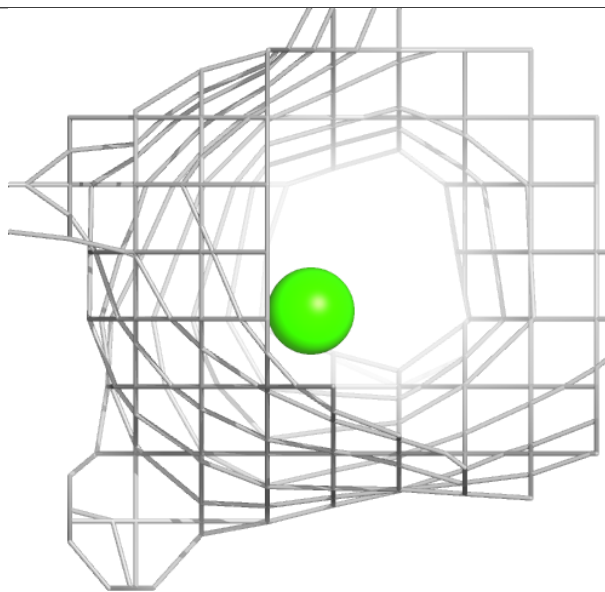
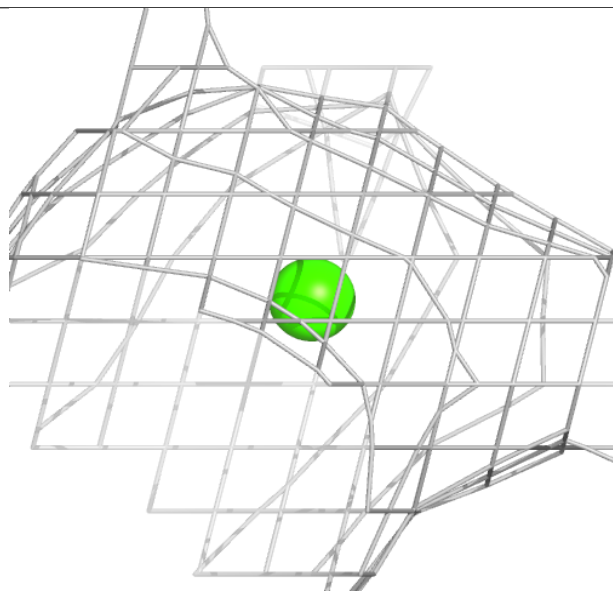
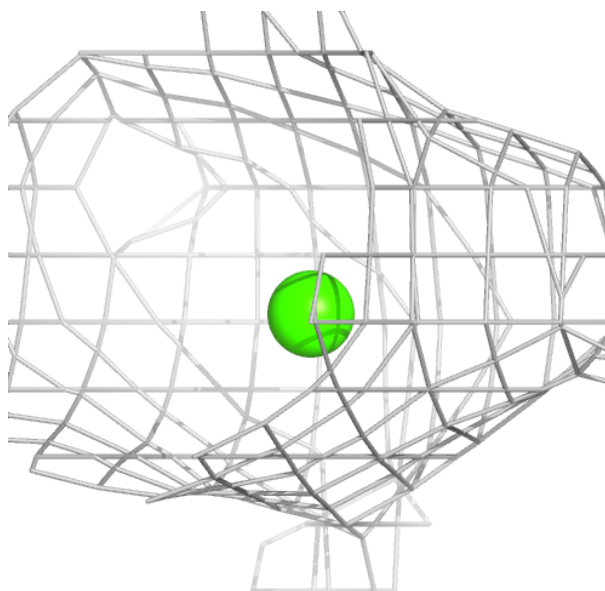
Electron density around CA v 901:

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



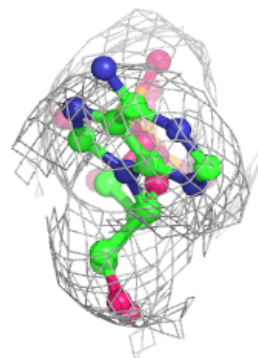
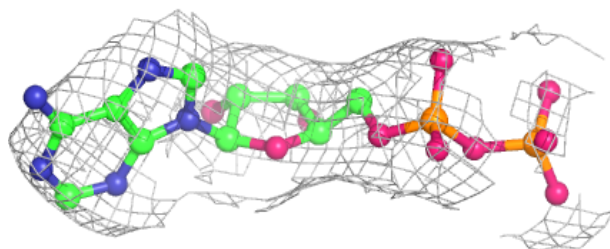
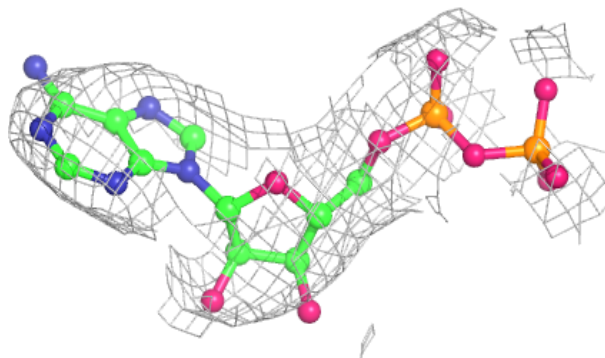
Electron density around CA g 402:

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



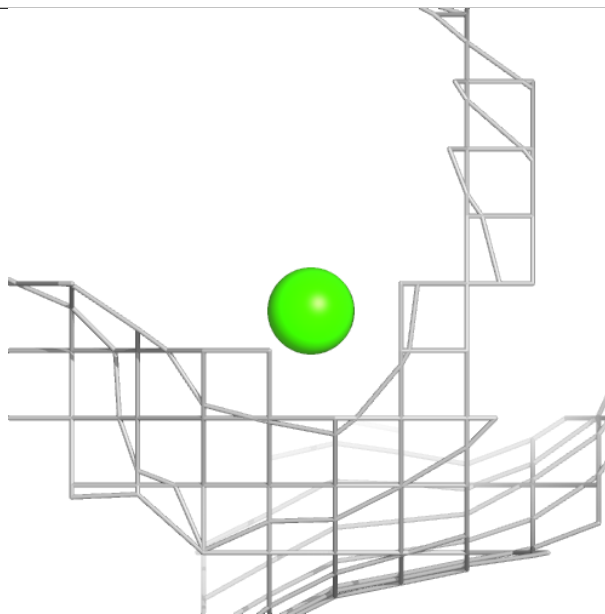
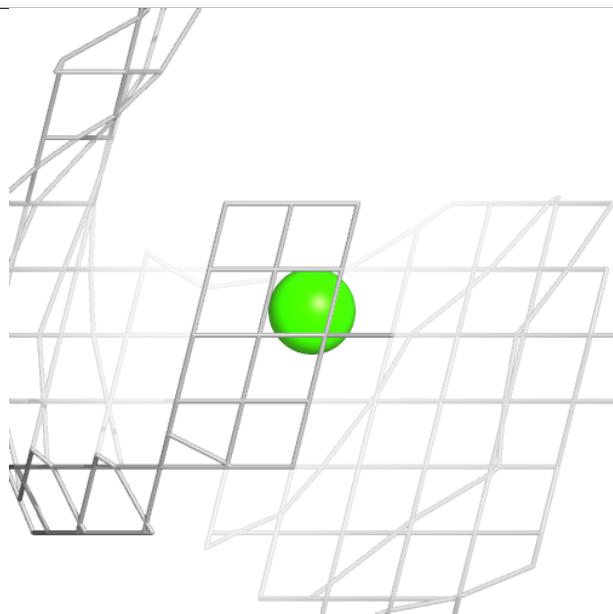
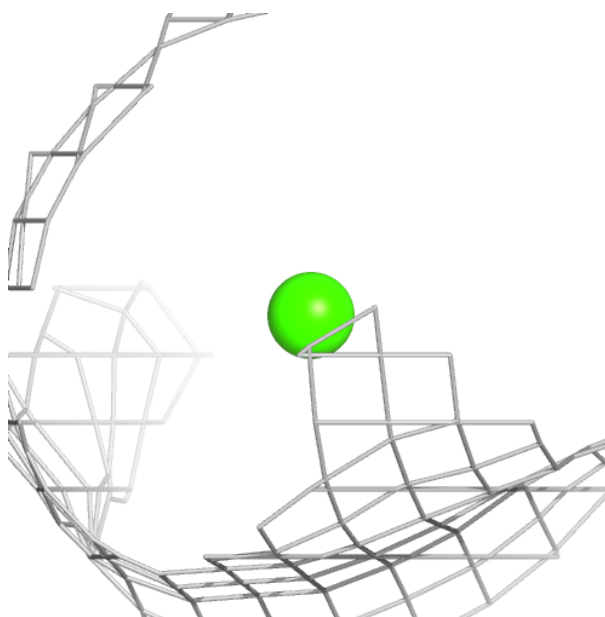
Electron density around ADP g 401:

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



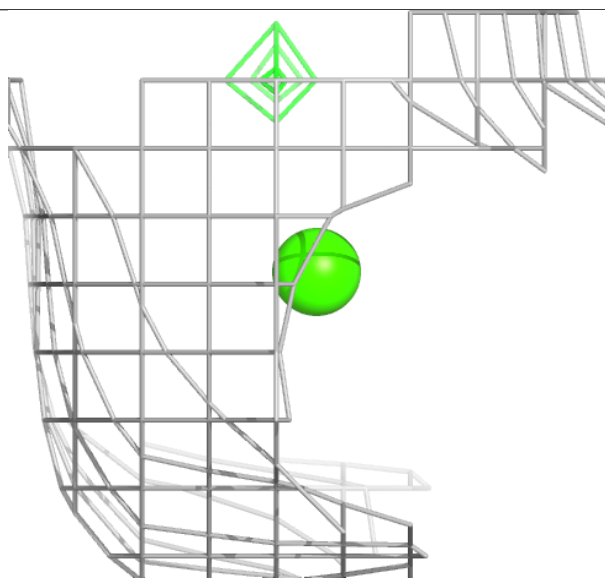
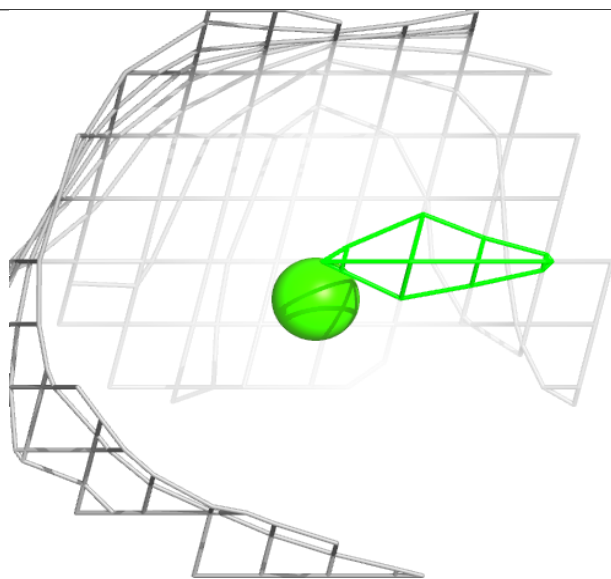
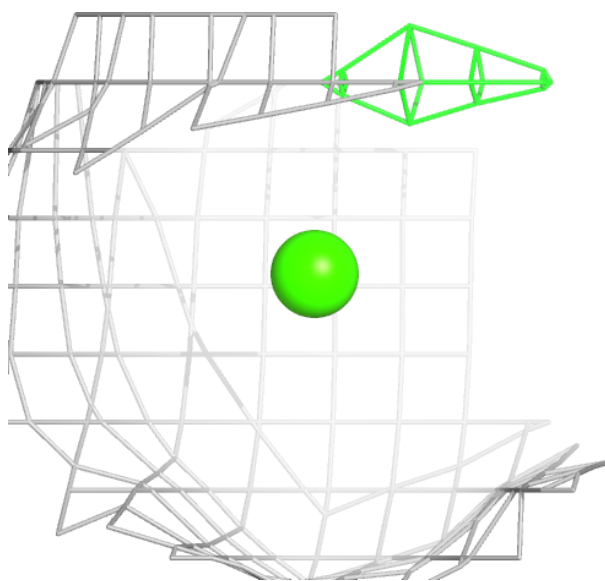
Electron density around CA v 903:

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



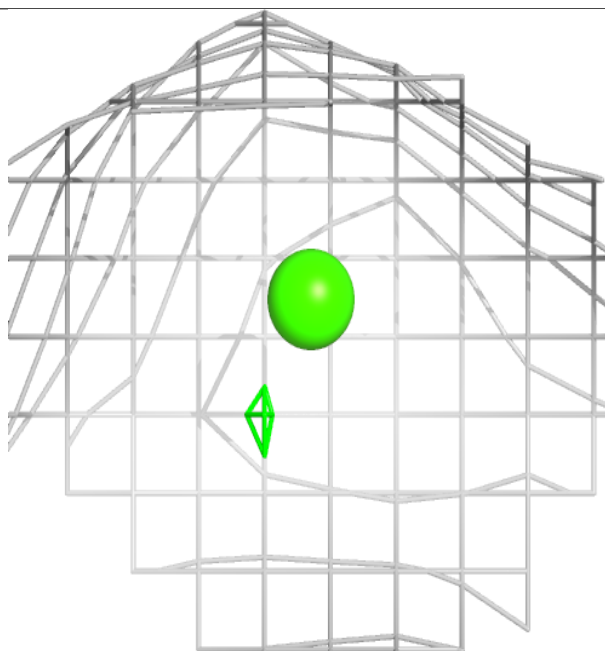
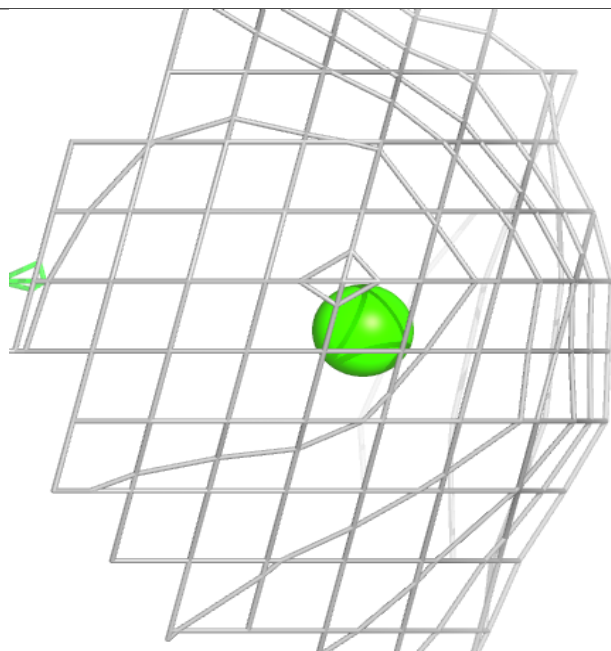
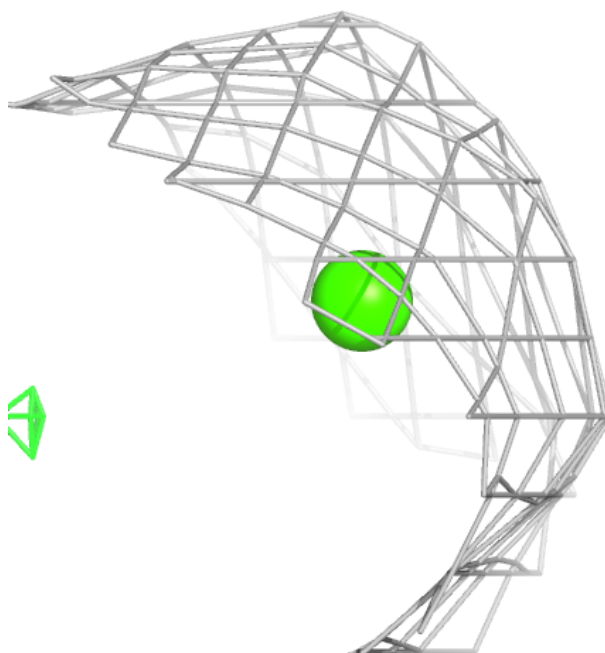
Electron density around CA V 904:

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



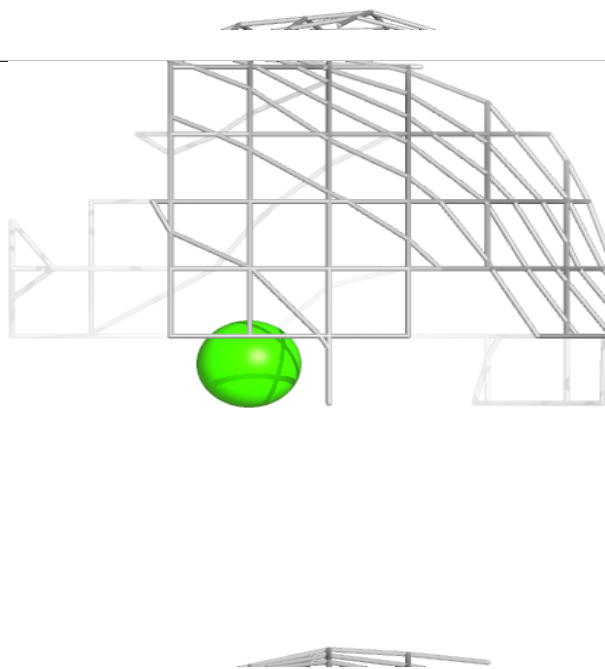
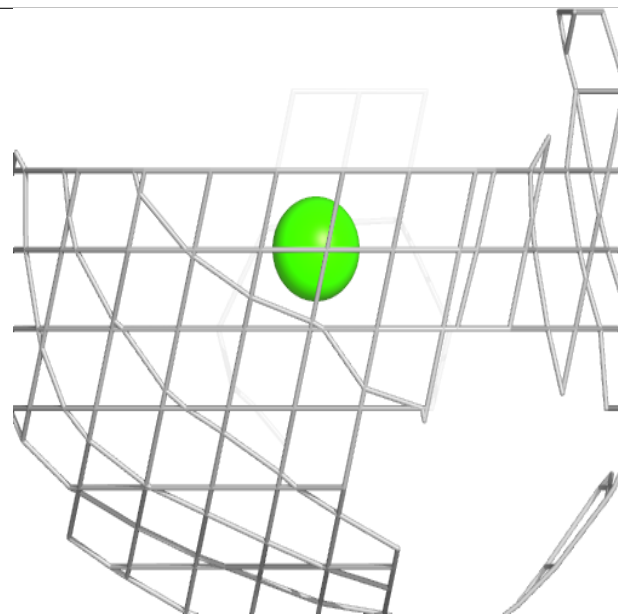
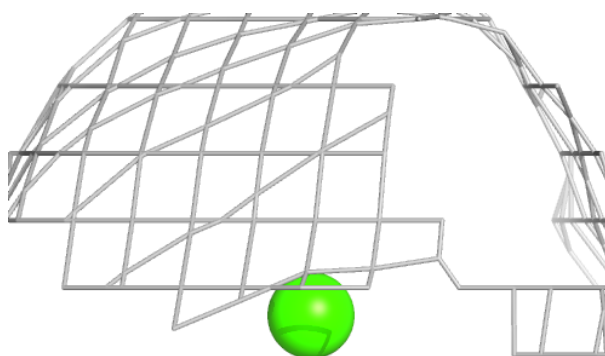
Electron density around CA V 902:

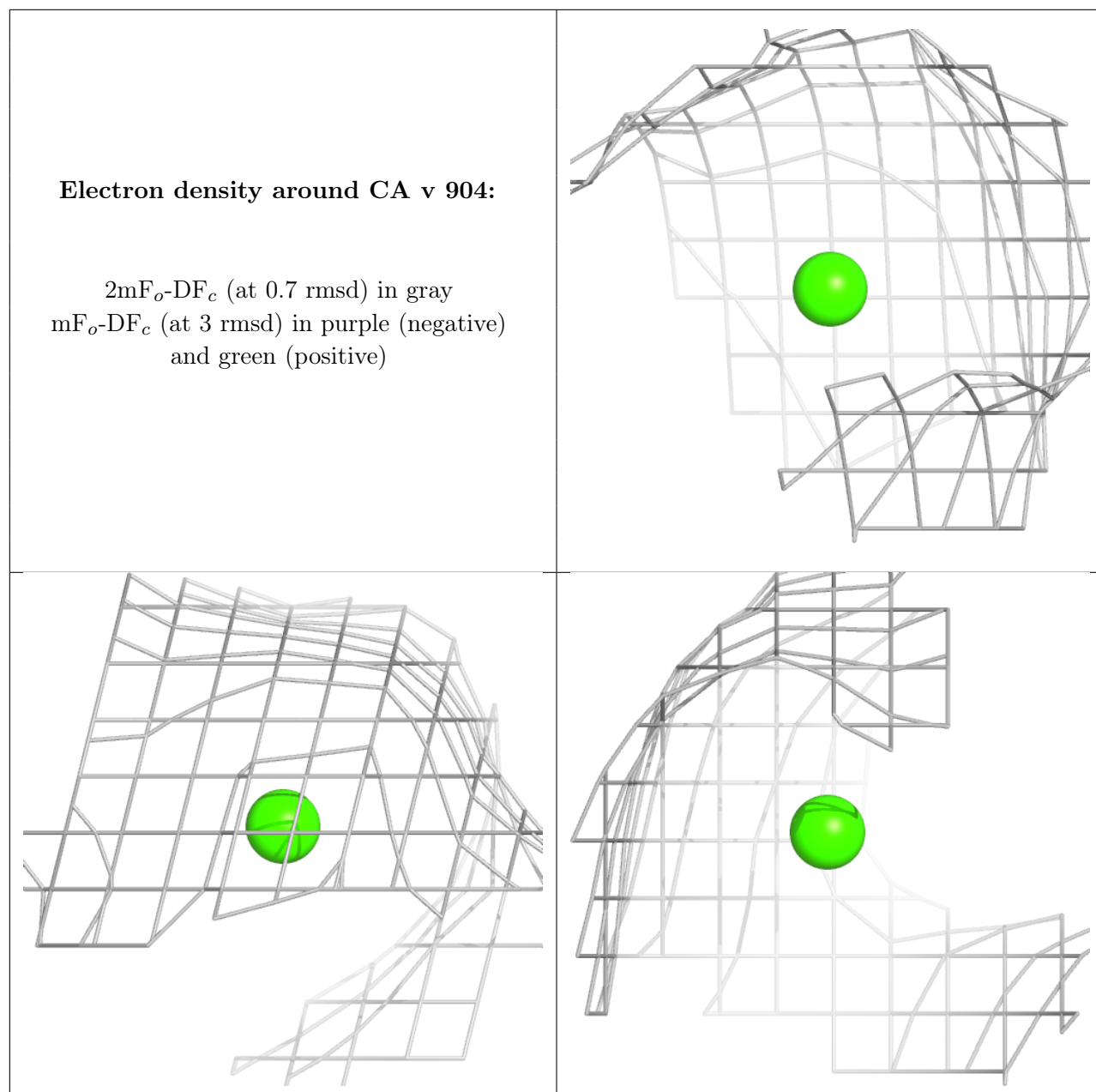
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around CA V 903:

$2mF_o-DF_c$ (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.