



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:02 AM UTC

PDB ID : 2O1V / pdb_00002o1v
Title : Structure of full length GRP94 with ADP bound
Authors : Dollins, D.E.; Warren, J.J.; Immormino, R.M.; Gewirth, D.T.
Deposited on : 2006-11-29
Resolution : 2.45 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

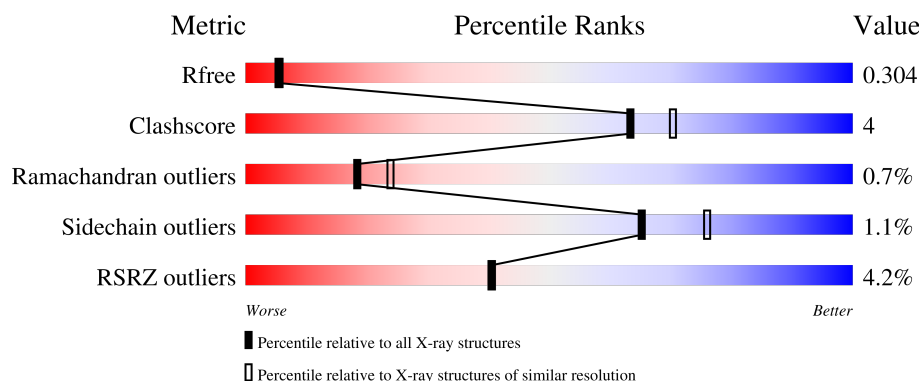
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

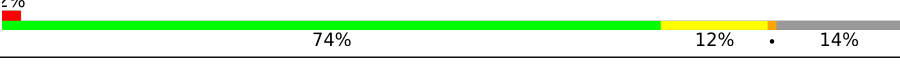
The reported resolution of this entry is 2.45 Å.

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	1190 (2.46-2.46)
Clashscore	190562	1229 (2.46-2.46)
Ramachandran outliers	187476	1218 (2.46-2.46)
Sidechain outliers	187428	1218 (2.46-2.46)
RSRZ outliers	180081	1190 (2.46-2.46)

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	666	 5% 77% 9% • 13%
1	B	666	 2% 74% 12% • 14%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9607 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 Endoplasmin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	578	Total	C	N	O	S	0	13	0
			4666	2978	785	886	17			
1	B	575	Total	C	N	O	S	0	0	0
			4623	2946	776	885	16			

There are 50 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	52	MET	-	expression tag	UNP P41148
A	53	GLY	-	expression tag	UNP P41148
A	54	SER	-	expression tag	UNP P41148
A	55	SER	-	expression tag	UNP P41148
A	56	HIS	-	expression tag	UNP P41148
A	57	HIS	-	expression tag	UNP P41148
A	58	HIS	-	expression tag	UNP P41148
A	59	HIS	-	expression tag	UNP P41148
A	60	HIS	-	expression tag	UNP P41148
A	61	HIS	-	expression tag	UNP P41148
A	62	SER	-	expression tag	UNP P41148
A	63	SER	-	expression tag	UNP P41148
A	64	GLY	-	expression tag	UNP P41148
A	65	LEU	-	expression tag	UNP P41148
A	66	VAL	-	expression tag	UNP P41148
A	67	PRO	-	expression tag	UNP P41148
A	68	ARG	-	expression tag	UNP P41148
A	69	GLY	-	expression tag	UNP P41148
A	70	SER	-	expression tag	UNP P41148
A	71	HIS	-	expression tag	UNP P41148
A	72	MET	-	expression tag	UNP P41148
A	287	GLY	-	see remark 999	UNP P41148
A	288	GLY	-	see remark 999	UNP P41148
A	289	GLY	-	see remark 999	UNP P41148
A	290	GLY	-	see remark 999	UNP P41148

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Chain	Residue	Modelled	Actual	Comment	Reference
B	52	MET	-	expression tag	UNP P41148
B	53	GLY	-	expression tag	UNP P41148
B	54	SER	-	expression tag	UNP P41148
B	55	SER	-	expression tag	UNP P41148
B	56	HIS	-	expression tag	UNP P41148
B	57	HIS	-	expression tag	UNP P41148
B	58	HIS	-	expression tag	UNP P41148
B	59	HIS	-	expression tag	UNP P41148
B	60	HIS	-	expression tag	UNP P41148
B	61	HIS	-	expression tag	UNP P41148
B	62	SER	-	expression tag	UNP P41148
B	63	SER	-	expression tag	UNP P41148
B	64	GLY	-	expression tag	UNP P41148
B	65	LEU	-	expression tag	UNP P41148
B	66	VAL	-	expression tag	UNP P41148
B	67	PRO	-	expression tag	UNP P41148
B	68	ARG	-	expression tag	UNP P41148
B	69	GLY	-	expression tag	UNP P41148
B	70	SER	-	expression tag	UNP P41148
B	71	HIS	-	expression tag	UNP P41148
B	72	MET	-	expression tag	UNP P41148
B	287	GLY	-	see remark 999	UNP P41148
B	288	GLY	-	see remark 999	UNP P41148
B	289	GLY	-	see remark 999	UNP P41148
B	290	GLY	-	see remark 999	UNP P41148

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	2	Total Mg 2 2	0	0
2	B	2	Total Mg 2 2	0	0

- Molecule 3 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: C₁₀H₁₅N₅O₁₀P₂).



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

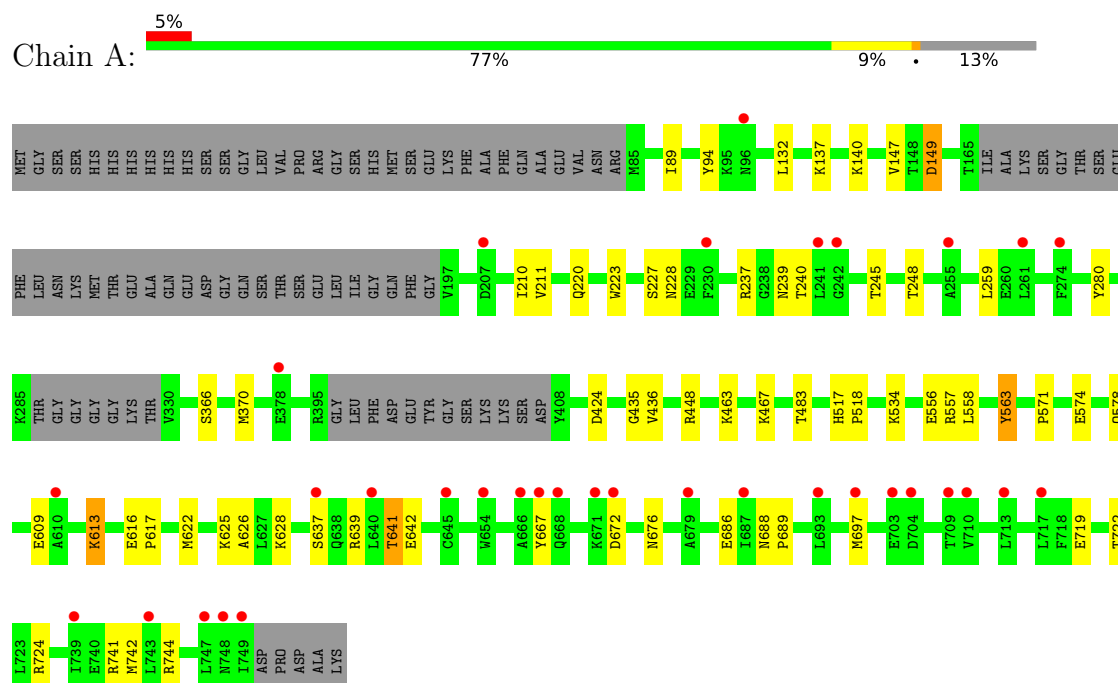
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	135	Total O 135 135	0	0
4	B	125	Total O 125 125	0	0

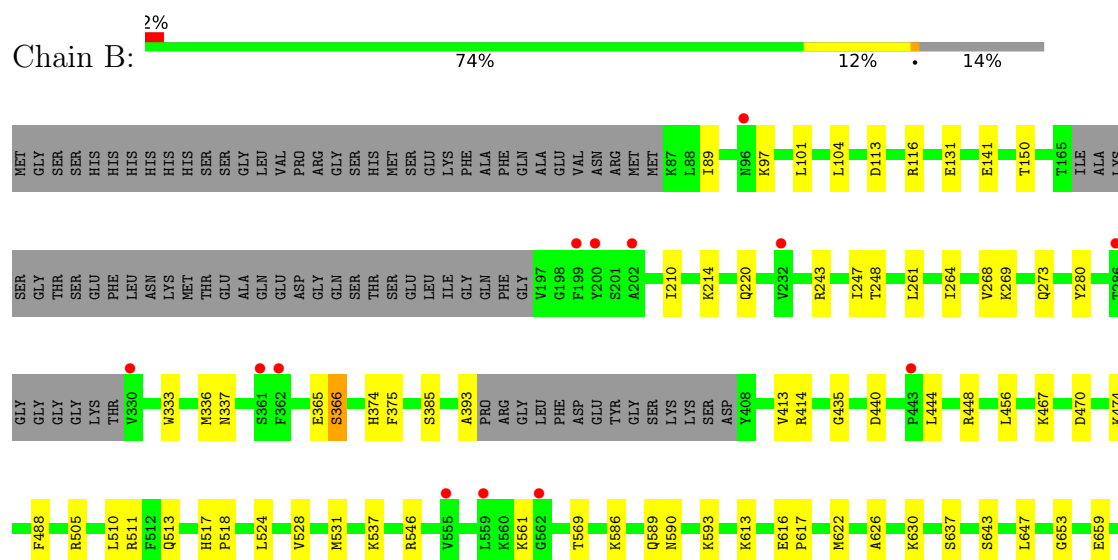
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Endoplasmic



• Molecule 1: Endoplasmic





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	99.08Å 108.85Å 148.98Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.45 20.00 – 2.45	Depositor EDS
% Data completeness (in resolution range)	99.6 (20.00-2.45) 99.4 (20.00-2.45)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.15 (at 2.45Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.248 , 0.299 0.264 , 0.304	Depositor DCC
R_{free} test set	2972 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	57.8	Xtriage
Anisotropy	0.357	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 61.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	9607	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.28% 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

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	0.66	9/4792 (0.2%)	0.99	6/6465 (0.1%)
1	B	0.52	0/4708	0.95	11/6351 (0.2%)
All	All	0.60	9/9500 (0.1%)	0.97	17/12816 (0.1%)

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	237	ARG	C-O	10.70	1.36	1.24
1	A	147	VAL	C-O	9.92	1.35	1.24
1	A	280	TYR	C-O	8.36	1.34	1.23
1	A	237	ARG	C-N	7.85	1.44	1.32
1	A	245	THR	C-O	6.07	1.31	1.23
1	A	239[A]	ASN	C-O	6.00	1.31	1.23
1	A	239[B]	ASN	C-O	6.00	1.31	1.23
1	A	132	LEU	C-N	5.87	1.41	1.33
1	A	132	LEU	C-O	5.67	1.30	1.23

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	97	LYS	N-CA-C	8.99	122.05	111.71
1	A	89	ILE	N-CA-C	-7.74	105.94	113.53
1	A	147	VAL	O-C-N	7.19	129.69	122.71
1	B	659	GLU	N-CA-C	-7.00	103.25	111.03
1	B	665	GLN	N-CA-C	-6.38	106.07	114.31
1	A	435	GLY	N-CA-C	6.25	119.00	110.69
1	B	435	GLY	N-CA-C	5.80	118.41	110.69
1	B	131	GLU	N-CA-C	5.75	118.12	110.53
1	B	691	HIS	CA-C-N	5.55	126.78	119.84
1	B	691	HIS	C-N-CA	5.55	126.78	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	643	SER	CA-C-N	5.39	124.91	119.19
1	B	643	SER	C-N-CA	5.39	124.91	119.19
1	B	666	ALA	N-CA-C	-5.29	106.67	113.23
1	A	563	TYR	N-CA-C	-5.28	103.42	110.55
1	B	670	GLY	N-CA-C	5.27	118.87	112.49
1	A	94	TYR	N-CA-C	-5.10	106.57	112.89
1	A	237	ARG	O-C-N	5.09	127.96	122.15

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	4666	0	4617	33	0
1	B	4623	0	4571	46	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	27	0	12	0	0
3	B	27	0	12	0	0
4	A	135	0	0	2	0
4	B	125	0	0	4	0
All	All	9607	0	9212	75	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 (75) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:637:SER:HB2	1:B:686:GLU:HB3	1.62	0.82
1:A:724:ARG:HD2	1:B:732:THR:HG23	1.67	0.76
1:A:741[B]:ARG:HH11	1:A:741[B]:ARG:CG	2.03	0.71
1:A:210:ILE:HG13	1:A:248:THR:HB	1.76	0.67
1:A:741[B]:ARG:HH11	1:A:741[B]:ARG:HG2	1.59	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:556:GLU:HB3	1:A:641:THR:HB	1.81	0.63
1:A:625:LYS:HA	1:A:628:LYS:HD2	1.82	0.62
1:B:104:LEU:HD22	1:B:247:ILE:HD13	1.82	0.61
1:A:637:SER:HB2	1:A:686:GLU:HB3	1.82	0.61
1:B:470:ASP:OD2	1:B:505:ARG:NH2	2.35	0.58
1:B:531:MET:HE1	1:B:586:LYS:HG2	1.86	0.58
1:B:590:ASN:HB3	1:B:593:LYS:HD2	1.87	0.57
1:B:653:GLY:HA2	1:B:678:TYR:HB2	1.88	0.56
1:A:370:MET:HE2	1:A:483:THR:HG22	1.87	0.56
1:A:140:LYS:HA	1:A:259:LEU:HD13	1.88	0.56
1:A:741[A]:ARG:HA	1:A:744[A]:ARG:HE	1.72	0.54
1:B:210:ILE:HB	1:B:248:THR:HB	1.90	0.54
1:B:414:ARG:HD2	1:B:444:LEU:HD22	1.89	0.53
1:B:413:VAL:HG23	1:B:444:LEU:HD13	1.91	0.52
1:B:666:ALA:O	1:B:671:LYS:N	2.39	0.52
1:B:474:LYS:HB2	4:B:872:HOH:O	2.10	0.52
1:B:517:HIS:CG	1:B:518:PRO:HD2	2.45	0.51
1:A:571:PRO:HG3	1:B:667:TYR:CE2	2.46	0.50
1:A:563:TYR:O	4:A:775:HOH:O	2.20	0.49
1:B:113:ASP:OD1	1:B:116:ARG:NH2	2.46	0.49
1:B:537:LYS:HD3	1:B:589:GLN:HB2	1.95	0.49
1:A:463:LYS:O	1:A:467:LYS:HG2	2.13	0.48
1:B:337:ASN:ND2	4:B:841:HOH:O	2.45	0.48
1:B:513:GLN:OE1	1:B:546:ARG:NH1	2.45	0.48
1:B:261:LEU:HA	1:B:264:ILE:HG22	1.96	0.47
1:B:719:GLU:OE2	1:B:735:TYR:HD1	1.97	0.47
1:A:667:TYR:HE1	1:B:569:THR:O	1.99	0.46
1:A:622:MET:HA	1:A:626:ALA:HB3	1.98	0.46
1:B:697:MET:HE2	1:B:713:LEU:HG	1.98	0.46
1:B:393:ALA:HA	1:B:488:PHE:HE2	1.80	0.46
1:B:622:MET:HA	1:B:626:ALA:HB3	1.98	0.46
1:A:220:GLN:HE22	1:A:240:THR:HB	1.81	0.45
1:B:647:LEU:HD23	1:B:647:LEU:HA	1.72	0.45
1:B:719:GLU:OE1	1:B:738:ARG:NH2	2.46	0.45
1:A:534:LYS:HA	1:A:534:LYS:HD2	1.84	0.45
1:A:517:HIS:CG	1:A:518:PRO:HD2	2.51	0.45
1:B:616:GLU:N	1:B:617:PRO:HD2	2.31	0.45
1:B:393:ALA:HA	1:B:488:PHE:CE2	2.51	0.45
1:A:149:ASP:OD2	1:A:149:ASP:C	2.60	0.45
1:B:375:PHE:CE1	1:B:467:LYS:HG3	2.51	0.44
1:A:719:GLU:O	1:A:722:THR:HB	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:336:MET:HA	1:B:336:MET:HE2	1.98	0.44
1:B:524:LEU:O	1:B:528:VAL:HG23	2.16	0.44
1:B:269:LYS:O	1:B:273:GLN:NE2	2.51	0.44
1:A:616:GLU:N	1:A:617:PRO:HD2	2.33	0.44
1:A:574:GLU:HB2	4:A:845:HOH:O	2.18	0.43
1:B:150:THR:O	1:B:243:ARG:NH2	2.48	0.43
1:B:630:LYS:NZ	4:B:828:HOH:O	2.48	0.43
1:B:561:LYS:HD3	1:B:561:LYS:HA	1.82	0.43
1:B:101:LEU:HD21	1:B:268:VAL:HG12	2.00	0.43
1:A:688:ASN:HA	1:A:689:PRO:HD2	1.93	0.42
1:A:211:VAL:HG22	1:A:223:TRP:HB3	2.00	0.42
1:B:613:LYS:HD2	1:B:613:LYS:HA	1.92	0.42
1:A:557:ARG:HB2	1:A:639:ARG:HA	2.00	0.42
1:B:141:GLU:H	1:B:141:GLU:CD	2.26	0.42
1:B:365:GLU:O	1:B:366:SER:HB3	2.19	0.42
1:A:424:ASP:OD1	1:A:424:ASP:N	2.53	0.42
1:A:574:GLU:O	1:A:578:GLN:HB2	2.20	0.42
1:B:505:ARG:HG3	4:B:779:HOH:O	2.19	0.41
1:B:666:ALA:HA	1:B:670:GLY:HA3	2.02	0.41
1:A:742:MET:SD	1:B:742:MET:HB3	2.61	0.41
1:A:697:MET:HE3	1:A:697:MET:HB2	1.70	0.41
1:B:214:LYS:HD3	1:B:220:GLN:HB2	2.02	0.41
1:A:641:THR:HG22	1:A:642:GLU:N	2.36	0.41
1:B:510:LEU:O	1:B:511:ARG:HD3	2.20	0.41
1:A:639:ARG:HD2	1:A:686:GLU:OE1	2.21	0.41
1:A:609:GLU:O	1:A:613:LYS:HD2	2.22	0.40
1:B:374:HIS:HA	1:B:385:SER:O	2.22	0.40
1:A:641:THR:HG22	1:A:642:GLU:HG3	2.03	0.40
1:B:280:TYR:HB3	1:B:333:TRP:HB3	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	583/666 (88%)	549 (94%)	29 (5%)	5 (1%)	14	17
1	B	565/666 (85%)	523 (93%)	39 (7%)	3 (0%)	24	32
All	All	1148/1332 (86%)	1072 (93%)	68 (6%)	8 (1%)	18	24

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	227	SER
1	A	366	SER
1	B	366	SER
1	A	228	ASN
1	B	89	ILE
1	A	676	ASN
1	B	677	TYR
1	A	672	ASP

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	508/598 (85%)	501 (99%)	7 (1%)	59	72
1	B	504/598 (84%)	500 (99%)	4 (1%)	73	83
All	All	1012/1196 (85%)	1001 (99%)	11 (1%)	65	76

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	137	LYS
1	A	149	ASP
1	A	436	VAL
1	A	448	ARG
1	A	558	LEU
1	A	613	LYS
1	A	641	THR
1	B	440	ASP

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Mol	Chain	Res	Type
1	B	448	ARG
1	B	456	LEU
1	B	732	THR

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	445	ASN
1	A	500	HIS
1	B	273	GLN
1	B	337	ASN
1	B	445	ASN
1	B	502	ASN
1	B	516	HIS
1	B	681	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 6 ligands modelled in this entry, 4 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
3	ADP	B	302	2	28,29,29	1.25	4 (14%)	43,45,45	1.68	8 (18%)
3	ADP	A	301	2	28,29,29	1.27	4 (14%)	43,45,45	1.76	8 (18%)

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	B	302	2	-	2/16/32/32	0/3/3/3
3	ADP	A	301	2	-	1/16/32/32	0/3/3/3

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	302	ADP	C5-N7	-3.17	1.33	1.39
3	A	301	ADP	PA-O3A	3.10	1.62	1.59
3	A	301	ADP	C5-N7	-2.86	1.33	1.39
3	B	302	ADP	PB-O3B	2.28	1.63	1.54
3	B	302	ADP	C1'-N9	2.23	1.52	1.46
3	A	301	ADP	C8-N9	-2.12	1.33	1.37
3	B	302	ADP	C4-N9	-2.11	1.33	1.37
3	A	301	ADP	C4-N9	-2.10	1.33	1.37

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	302	ADP	C5-C4-N3	-5.18	119.58	126.72
3	A	301	ADP	C5-C4-N3	-4.93	119.93	126.72
3	A	301	ADP	N3-C2-N1	-4.74	121.41	128.58
3	B	302	ADP	N3-C2-N1	-4.71	121.46	128.58
3	B	302	ADP	C2-N3-C4	3.54	120.49	111.83
3	A	301	ADP	C2-N3-C4	3.43	120.22	111.83
3	A	301	ADP	N3-C4-N9	3.33	132.83	127.17
3	B	302	ADP	N3-C4-N9	3.28	132.75	127.17
3	A	301	ADP	N9-C8-N7	-3.25	109.33	113.94
3	A	301	ADP	C4-C5-N7	-2.85	107.33	110.58
3	A	301	ADP	C5-N7-C8	2.84	107.91	103.45
3	B	302	ADP	C4-C5-N7	-2.72	107.47	110.58
3	B	302	ADP	N9-C8-N7	-2.61	110.23	113.94
3	B	302	ADP	C5-N7-C8	2.55	107.46	103.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	301	ADP	C4-N9-C8	2.34	108.19	105.74
3	B	302	ADP	O4'-C1'-N9	2.10	112.12	108.09

There are no chirality outliers.

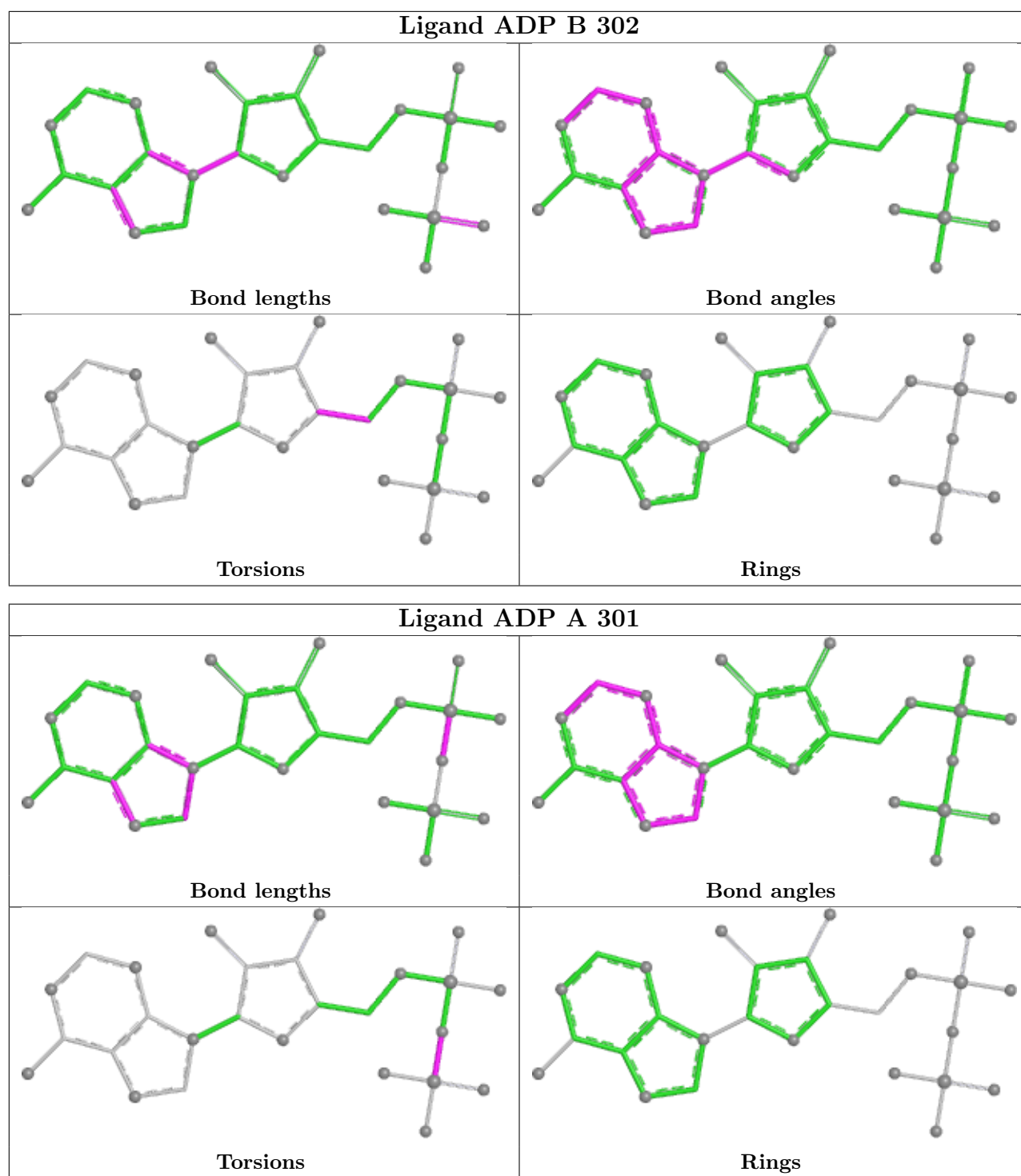
All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	302	ADP	O4'-C4'-C5'-O5'
3	B	302	ADP	C3'-C4'-C5'-O5'
3	A	301	ADP	PA-O3A-PB-O2B

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

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	578/666 (86%)	0.38	34 (5%)	28 25	13, 30, 43, 54	13 (2%)
1	B	575/666 (86%)	0.21	14 (2%)	59 62	10, 30, 40, 51	0
All	All	1153/1332 (86%)	0.29	48 (4%)	40 40	10, 30, 42, 54	13 (1%)

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	671	LYS	4.7
1	A	679	ALA	4.4
1	A	610	ALA	4.0
1	B	562	GLY	3.6
1	A	749	ILE	3.5
1	A	747	LEU	3.5
1	B	202	ALA	3.4
1	A	654	TRP	3.2
1	A	748	ASN	3.1
1	B	96	ASN	3.1
1	A	261	LEU	3.0
1	A	242	GLY	2.9
1	A	717	LEU	2.9
1	A	697	MET	2.9
1	A	709	THR	2.8
1	B	199	PHE	2.8
1	A	672	ASP	2.8
1	A	378	GLU	2.7
1	A	704	ASP	2.7
1	A	274	PHE	2.7
1	A	743	LEU	2.6
1	A	207	ASP	2.6
1	A	710	VAL	2.6
1	A	693	LEU	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	713	LEU	2.5
1	B	362	PHE	2.5
1	B	330	VAL	2.5
1	A	255	ALA	2.5
1	B	555	VAL	2.4
1	B	443	PRO	2.4
1	B	678	TYR	2.3
1	B	559	LEU	2.3
1	A	667	TYR	2.2
1	A	666	ALA	2.2
1	A	703	GLU	2.2
1	A	230	PHE	2.2
1	A	640	LEU	2.1
1	B	286	THR	2.1
1	A	739	ILE	2.1
1	A	668	GLN	2.1
1	A	241	LEU	2.1
1	A	637	SER	2.1
1	A	96	ASN	2.1
1	B	200	TYR	2.1
1	A	645	CYS	2.0
1	B	232	VAL	2.0
1	A	687	ILE	2.0
1	B	361	SER	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](#)

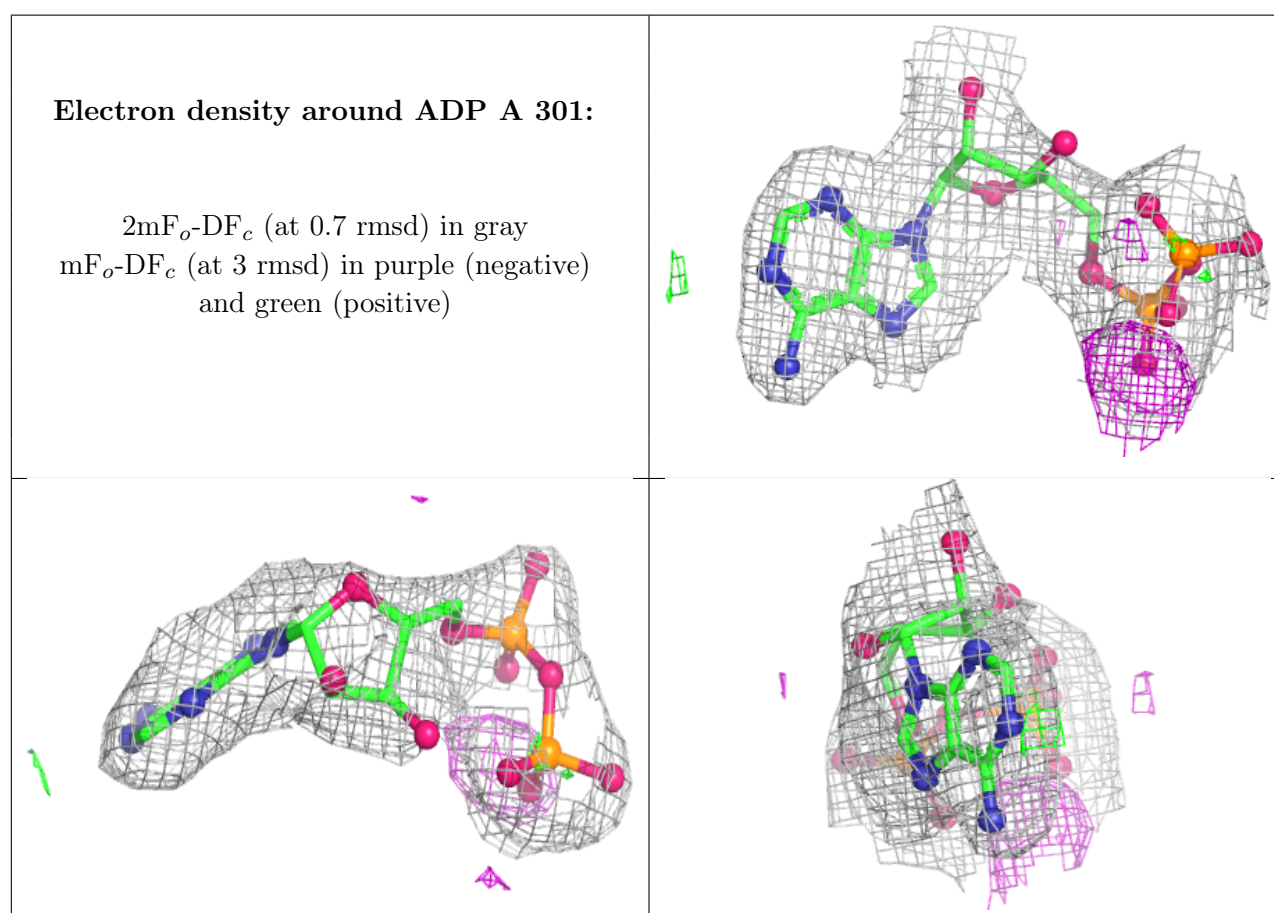
There are no oligosaccharides in this entry.

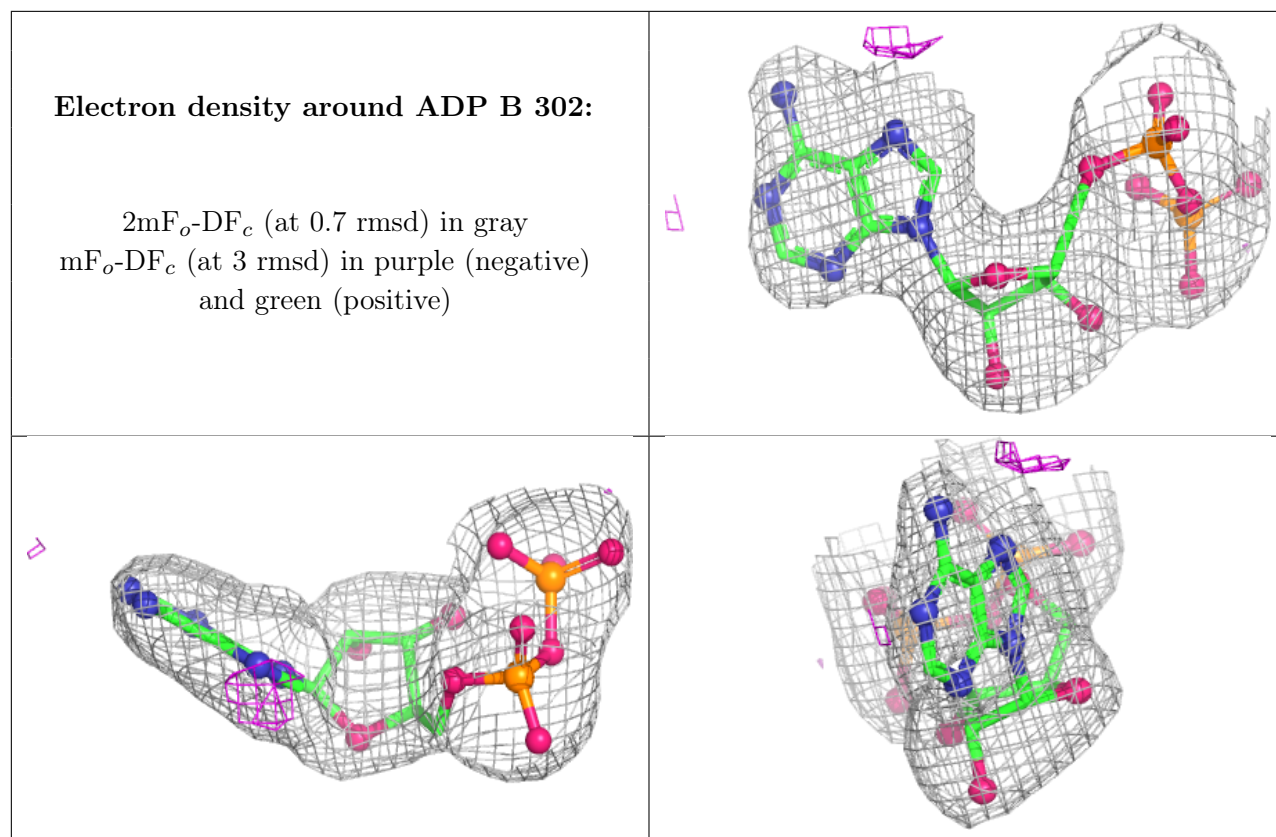
6.4 Ligands [i](#)

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
2	MG	B	756	1/1	0.85	0.08	34,34,34,34	0
2	MG	A	756	1/1	0.86	0.06	42,42,42,42	0
3	ADP	A	301	27/27	0.88	0.09	29,32,34,34	0
3	ADP	B	302	27/27	0.92	0.07	35,37,39,39	0
2	MG	B	755	1/1	0.96	0.09	37,37,37,37	0
2	MG	A	755	1/1	0.98	0.20	12,12,12,12	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.





6.5 Other polymers [i](#)

There are no such residues in this entry.