



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:45 PM UTC

PDB ID : 1RZU / pdb_00001rzu
Title : Crystal structure of the glycogen synthase from *A. tumefaciens* in complex with ADP
Authors : Buschiazzo, A.; Guerin, M.E.; Ugalde, J.E.; Ugalde, R.A.; Shepard, W.; Alzari, P.M.
Deposited on : 2003-12-29
Resolution : 2.30 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

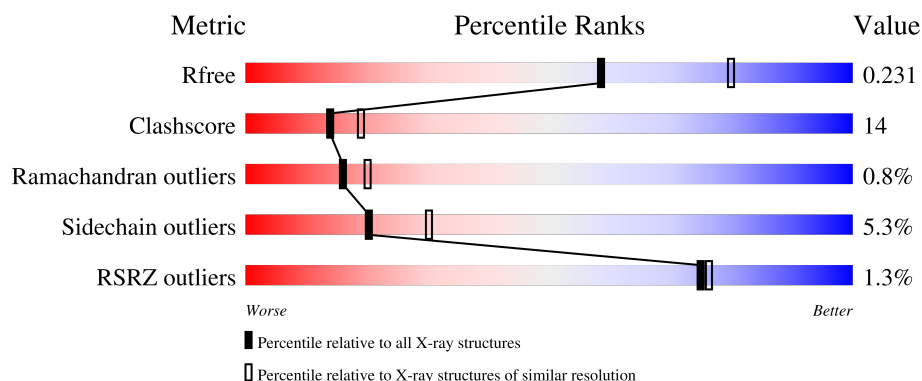
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	485	% 68% 27% ••
1	B	485	% 71% 22% 5% •

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7493 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 Glycogen synthase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	477	Total	C	N	O	S	0	0	0
			3618	2312	624	667	15			
1	B	478	Total	C	N	O	S	0	0	0
			3614	2310	622	668	14			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	481	HIS	-	expression tag	UNP P0A3F3
A	482	HIS	-	expression tag	UNP P0A3F3
A	483	HIS	-	expression tag	UNP P0A3F3
A	484	HIS	-	expression tag	UNP P0A3F3
A	485	HIS	-	expression tag	UNP P0A3F3
B	481	HIS	-	expression tag	UNP P0A3F3
B	482	HIS	-	expression tag	UNP P0A3F3
B	483	HIS	-	expression tag	UNP P0A3F3
B	484	HIS	-	expression tag	UNP P0A3F3
B	485	HIS	-	expression tag	UNP P0A3F3

- Molecule 2 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: C₁₀H₁₅N₅O₁₀P₂).



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

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	108	Total O 108 108	0	0
3	B	99	Total O 99 99	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	69.36Å 88.11Å 84.46Å 90.00° 98.19° 90.00°	Depositor
Resolution (Å)	50.00 – 2.30 50.00 – 2.30	Depositor EDS
% Data completeness (in resolution range)	90.7 (50.00-2.30) 90.0 (50.00-2.30)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.82 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.0	Depositor
R, R_{free}	0.180 , 0.223 0.185 , 0.231	Depositor DCC
R_{free} test set	4055 reflections (10.04%)	wwPDB-VP
Wilson B-factor (Å ²)	38.5	Xtriage
Anisotropy	0.015	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 50.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7493	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.70% 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

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	1.11	5/3702 (0.1%)	1.12	8/5040 (0.2%)
1	B	1.15	7/3698 (0.2%)	1.12	10/5037 (0.2%)
All	All	1.13	12/7400 (0.2%)	1.12	18/10077 (0.2%)

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	309	MET	SD-CE	-8.74	1.57	1.79
1	A	450	MET	SD-CE	-7.60	1.60	1.79
1	A	148	MET	SD-CE	-7.58	1.60	1.79
1	B	450	MET	SD-CE	-7.55	1.60	1.79
1	A	410	ALA	C-O	-7.02	1.15	1.24
1	A	188	MET	SD-CE	6.31	1.95	1.79
1	A	425	VAL	CA-C	-5.98	1.47	1.52
1	B	410	ALA	C-O	-5.92	1.17	1.24
1	B	212	THR	C-O	-5.66	1.16	1.23
1	B	436	ARG	C-O	-5.59	1.17	1.24
1	B	418	THR	C-O	-5.33	1.17	1.24
1	B	269	ASN	CB-CG	5.07	1.64	1.52

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	191	ILE	N-CA-C	7.35	119.20	112.29
1	B	99	GLN	N-CA-C	-7.25	104.96	114.31
1	B	100	THR	N-CA-C	-7.16	104.68	113.41
1	B	298	SER	N-CA-C	6.70	119.50	108.99
1	A	381	THR	N-CA-C	6.46	118.32	111.28
1	B	102	LYS	N-CA-C	-6.38	99.61	109.24
1	A	374	ARG	N-CA-C	-6.12	105.97	113.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	459	VAL	N-CA-C	-6.05	106.87	112.43
1	B	299	ARG	N-CA-C	-5.80	101.82	110.23
1	B	394	VAL	N-CA-C	5.72	118.18	108.86
1	A	373	SER	N-CA-C	5.58	117.94	110.35
1	B	291	SER	CA-C-N	5.35	125.61	119.83
1	B	291	SER	C-N-CA	5.35	125.61	119.83
1	A	52	THR	N-CA-C	5.34	117.61	108.90
1	B	316	ILE	CB-CA-C	-5.33	105.15	111.97
1	A	51	VAL	N-CA-C	5.30	116.85	109.80
1	A	213	VAL	N-CA-C	5.29	120.34	109.34
1	B	112	ALA	N-CA-C	-5.06	105.76	111.28

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	3618	0	3610	105	0
1	B	3614	0	3594	104	0
2	A	27	0	12	2	0
2	B	27	0	12	0	0
3	A	108	0	0	5	0
3	B	99	0	0	5	0
All	All	7493	0	7228	209	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 14.

All (209) 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:309:MET:HE1	1:B:370:ILE:HG21	1.34	1.09
1:A:382:GLN:HE21	1:A:382:GLN:H	1.05	1.00
1:B:382:GLN:HE21	1:B:382:GLN:H	1.09	0.97

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:102:LYS:HB2	1:A:102:LYS:NZ	1.89	0.88
1:B:14:ILE:HD11	1:B:95:PRO:HG2	1.59	0.83
1:B:309:MET:HE1	1:B:370:ILE:CG2	2.09	0.81
1:A:102:LYS:HB2	1:A:102:LYS:HZ2	1.49	0.77
1:B:144:THR:HB	1:B:145:PRO:HD3	1.67	0.76
1:B:86:PRO:HA	1:B:90:GLU:HG2	1.69	0.75
1:B:97:LEU:HB3	1:B:101:GLY:HA2	1.70	0.72
1:A:431:LYS:O	1:A:435:ARG:HG3	1.89	0.72
1:B:211:SER:HA	1:B:242:HIS:O	1.90	0.72
1:A:303:GLN:HG2	3:A:1591:HOH:O	1.91	0.70
1:B:402:ASP:OD1	3:B:2612:HOH:O	2.09	0.70
1:B:51:VAL:HG12	1:B:54:PRO:HG3	1.74	0.69
1:A:91:ARG:NH2	1:A:106:ASP:OD1	2.20	0.69
1:B:14:ILE:HD13	1:B:15:LYS:N	2.06	0.69
1:B:53:ASP:N	1:B:54:PRO:HD3	2.07	0.69
1:A:142:ALA:O	1:A:145:PRO:HD2	1.92	0.68
1:A:327:GLY:HA2	2:A:1531:ADP:HN61	1.57	0.68
1:A:44:TYR:HB2	1:A:47:VAL:HG23	1.75	0.68
1:B:382:GLN:HE21	1:B:382:GLN:N	1.88	0.67
1:A:314:ASP:OD2	1:A:343:ARG:NH1	2.28	0.67
1:B:123:ALA:HB1	1:B:153:THR:HG21	1.76	0.67
1:B:54:PRO:HA	1:B:73:VAL:HG22	1.77	0.66
1:A:102:LYS:HZ2	1:A:102:LYS:CB	2.07	0.66
1:B:51:VAL:HG13	1:B:73:VAL:HG11	1.78	0.66
1:A:13:LEU:HD21	1:A:75:HIS:CD2	2.31	0.65
1:B:419:GLY:N	1:B:450:MET:HE3	2.11	0.65
1:A:339:ALA:O	1:A:343:ARG:HG2	1.97	0.65
1:A:340:ALA:HA	1:A:343:ARG:HE	1.62	0.64
1:A:26:LEU:HB3	1:A:27:PRO:HD3	1.80	0.64
1:A:473:SER:HA	1:A:476:ILE:HG12	1.80	0.63
1:A:267:ALA:O	1:A:452:LYS:NZ	2.28	0.62
1:B:291:SER:HB2	1:B:292:PRO:HD2	1.81	0.62
1:A:144:THR:HB	1:A:145:PRO:HD3	1.81	0.62
1:A:475:LEU:C	1:A:477:SER:H	2.09	0.61
1:B:14:ILE:HD13	1:B:14:ILE:C	2.25	0.61
1:B:67:LYS:N	1:B:67:LYS:HD2	2.15	0.61
1:B:418:THR:C	1:B:450:MET:HE3	2.26	0.61
1:A:222:LEU:HD21	1:A:234:ILE:HG22	1.83	0.60
1:B:304:LYS:HE3	3:B:2577:HOH:O	2.00	0.60
1:A:45:PRO:HD3	1:A:89:TYR:O	2.01	0.60
1:A:382:GLN:HE21	1:A:382:GLN:N	1.87	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:382:GLN:H	1:A:382:GLN:NE2	1.88	0.60
1:A:148:MET:HE1	1:A:156:ILE:HG12	1.84	0.59
1:A:159:LEU:HD11	1:A:211:SER:HB3	1.84	0.59
1:A:53:ASP:N	1:A:54:PRO:HD3	2.18	0.59
1:A:97:LEU:N	1:A:97:LEU:HD22	2.16	0.59
1:A:54:PRO:HA	1:A:73:VAL:HG22	1.82	0.59
1:B:477:SER:O	1:B:478:LYS:HB2	2.02	0.59
1:A:327:GLY:HA2	2:A:1531:ADP:N6	2.18	0.58
1:A:340:ALA:HA	1:A:343:ARG:NE	2.18	0.58
1:A:200:LEU:O	1:A:204:LEU:HG	2.02	0.58
1:A:71:LEU:HD21	1:A:84:ASP:HB2	1.85	0.57
1:A:317:VAL:HG11	1:A:344:HIS:ND1	2.19	0.57
1:B:212:THR:HG21	1:B:218:ALA:N	2.20	0.57
1:B:41:ILE:H	1:B:41:ILE:HD13	1.69	0.57
1:B:200:LEU:O	1:B:204:LEU:HG	2.05	0.56
1:A:436:ARG:NH1	3:A:1577:HOH:O	2.25	0.56
1:B:143:MET:SD	1:B:146:VAL:HG21	2.45	0.56
1:A:307:ASP:CG	1:A:374:ARG:HH12	2.14	0.56
1:B:73:VAL:HG12	1:B:74:GLN:H	1.70	0.56
1:B:11:TYR:CZ	1:B:12:PRO:HB3	2.41	0.56
1:B:73:VAL:HG12	1:B:74:GLN:N	2.21	0.55
1:A:406:ASP:HB2	1:A:421:GLN:HE21	1.72	0.55
1:B:59:GLU:HA	1:B:68:ALA:O	2.06	0.55
1:B:382:GLN:H	1:B:382:GLN:NE2	1.92	0.54
1:B:19:LEU:C	1:B:19:LEU:HD23	2.32	0.54
1:B:129:TRP:CZ3	1:B:131:PRO:HD3	2.42	0.53
1:A:26:LEU:N	1:A:27:PRO:CD	2.71	0.53
1:A:262:HIS:ND1	1:A:263:ASP:OD1	2.41	0.53
1:B:262:HIS:ND1	1:B:263:ASP:OD1	2.38	0.53
1:A:47:VAL:O	1:A:51:VAL:HG23	2.09	0.53
1:A:289:ASP:OD2	1:A:291:SER:OG	2.26	0.53
1:A:102:LYS:HZ2	1:A:102:LYS:N	2.07	0.52
1:B:223:THR:OG1	1:B:225:GLU:HG2	2.09	0.52
1:A:120:ARG:HB3	1:A:125:VAL:HG21	1.91	0.52
1:B:40:LEU:HD21	1:B:83:LEU:HB2	1.91	0.52
1:A:102:LYS:HB2	1:A:102:LYS:HZ3	1.72	0.52
1:A:51:VAL:HG12	1:A:73:VAL:HG11	1.91	0.52
1:B:222:LEU:HD22	1:B:235:GLY:HA2	1.93	0.51
1:A:222:LEU:CD2	1:A:234:ILE:HG22	2.39	0.51
1:A:300:LEU:HA	1:A:306:ILE:HG13	1.93	0.50
1:B:89:TYR:O	1:B:91:ARG:N	2.44	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:70:LEU:HD12	1:A:82:ILE:O	2.12	0.50
1:B:269:ASN:HB2	3:B:2621:HOH:O	2.12	0.50
1:B:51:VAL:HG21	1:B:82:ILE:HD11	1.93	0.50
1:B:176:PHE:CE1	1:B:186:PHE:HB2	2.47	0.49
1:A:44:TYR:CZ	1:A:95:PRO:HB3	2.47	0.49
1:B:91:ARG:HH22	1:B:106:ASP:CG	2.20	0.49
1:B:99:GLN:HA	1:B:99:GLN:OE1	2.11	0.49
1:A:92:SER:O	1:A:99:GLN:NE2	2.45	0.49
1:A:76:GLU:O	1:A:77:ARG:HB2	2.12	0.49
1:B:148:MET:HA	3:B:2580:HOH:O	2.12	0.49
1:B:222:LEU:HD21	1:B:234:ILE:HG22	1.95	0.49
1:B:340:ALA:HA	1:B:343:ARG:CZ	2.42	0.49
1:A:267:ALA:HB1	1:A:456:LYS:HZ2	1.77	0.48
1:B:19:LEU:C	1:B:19:LEU:CD2	2.86	0.48
1:B:51:VAL:HG11	1:B:73:VAL:HG21	1.95	0.48
1:A:233:VAL:HG23	3:A:1592:HOH:O	2.12	0.48
1:A:277:LYS:HE2	1:A:322:ARG:HH22	1.79	0.48
1:B:51:VAL:CG1	1:B:54:PRO:HG3	2.40	0.48
1:A:33:HIS:NE2	1:A:462:GLU:O	2.39	0.48
1:A:340:ALA:HA	1:A:343:ARG:CG	2.43	0.48
1:B:61:THR:HG23	1:B:61:THR:O	2.14	0.48
1:B:11:TYR:CD1	1:B:12:PRO:HA	2.49	0.47
1:A:307:ASP:OD1	1:A:374:ARG:NH1	2.43	0.47
1:A:47:VAL:HG12	1:A:82:ILE:HD13	1.96	0.47
1:A:14:ILE:HG13	1:A:15:LYS:N	2.29	0.47
1:A:446:LEU:C	1:A:446:LEU:HD23	2.40	0.47
1:B:222:LEU:CD2	1:B:234:ILE:HG22	2.45	0.47
1:A:225:GLU:CD	1:A:225:GLU:H	2.22	0.47
1:A:324:VAL:HA	1:A:349:GLY:O	2.15	0.47
1:B:11:TYR:CE1	1:B:12:PRO:HB3	2.49	0.47
1:A:220:GLU:HB3	1:A:226:PHE:CD2	2.50	0.46
1:B:173:ALA:HB2	1:B:195:ASN:O	2.15	0.46
1:B:103:ASP:HB3	1:B:107:ASN:HD22	1.80	0.46
1:A:51:VAL:CG1	1:A:73:VAL:HG11	2.45	0.46
1:A:267:ALA:HB1	1:A:456:LYS:NZ	2.30	0.46
1:B:158:SER:OG	1:B:207:ALA:HA	2.15	0.46
1:A:254:ASN:OD1	1:A:255:PRO:HD2	2.15	0.46
1:A:298:SER:O	1:A:300:LEU:CD1	2.62	0.46
1:A:10:ILE:HG13	1:A:23:VAL:HG12	1.97	0.46
1:A:106:ASP:HB3	1:A:110:ARG:HD2	1.97	0.46
1:A:236:SER:C	1:A:237:ARG:HG2	2.41	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:74:GLN:OE1	1:A:78:LEU:C	2.59	0.46
1:A:298:SER:O	1:A:300:LEU:HD12	2.16	0.46
1:B:445:LYS:C	1:B:445:LYS:HD2	2.40	0.46
1:B:314:ASP:OD1	1:B:343:ARG:NH1	2.49	0.45
1:A:173:ALA:HA	1:A:197:VAL:HG23	1.98	0.45
1:B:123:ALA:HB1	1:B:153:THR:CG2	2.45	0.45
1:A:108:TRP:HB3	1:A:171:PHE:CE1	2.52	0.45
1:B:59:GLU:CD	1:B:59:GLU:C	2.83	0.45
1:A:334:GLU:HB3	1:A:352:ILE:CD1	2.46	0.45
1:A:380:LEU:HD21	3:A:1636:HOH:O	2.15	0.45
1:A:165:ILE:CD1	1:A:230:LEU:HD12	2.46	0.45
1:B:41:ILE:HD13	1:B:41:ILE:N	2.31	0.45
1:A:475:LEU:C	1:A:477:SER:N	2.73	0.45
1:A:423:SER:HA	1:A:424:PRO:C	2.41	0.45
1:A:104:TYR:C	1:A:106:ASP:H	2.25	0.45
1:B:106:ASP:HB3	1:B:110:ARG:HD2	1.98	0.45
1:B:26:LEU:N	1:B:27:PRO:CD	2.80	0.44
1:A:9:GLU:HB2	1:A:14:ILE:HG23	2.00	0.44
1:B:42:PRO:HB2	1:B:44:TYR:CE2	2.53	0.44
1:B:67:LYS:HD2	1:B:67:LYS:H	1.82	0.44
1:B:476:ILE:HG22	1:B:476:ILE:O	2.17	0.44
1:A:228:MET:HE2	3:A:1545:HOH:O	2.18	0.43
1:B:160:LEU:HD23	1:B:204:LEU:HD23	1.99	0.43
1:B:41:ILE:HD13	1:B:81:LEU:O	2.18	0.43
1:B:345:HIS:CD2	1:B:345:HIS:H	2.35	0.43
1:B:446:LEU:C	1:B:446:LEU:HD23	2.43	0.43
1:B:58:PHE:HB3	1:B:70:LEU:HB3	1.99	0.43
1:A:1:MET:HE1	1:A:473:SER:HB3	2.01	0.43
1:A:300:LEU:HG	1:A:306:ILE:HG21	2.01	0.43
1:B:176:PHE:CZ	1:B:186:PHE:HB2	2.54	0.43
1:B:267:ALA:O	1:B:452:LYS:NZ	2.52	0.43
1:B:58:PHE:CZ	1:B:125:VAL:CG1	3.01	0.43
1:B:294:PHE:CD1	1:B:370:ILE:HD11	2.53	0.43
1:B:309:MET:CE	1:B:370:ILE:CG2	2.88	0.43
1:A:93:GLY:C	1:A:94:GLY:O	2.61	0.43
1:B:53:ASP:N	1:B:54:PRO:CD	2.79	0.43
1:B:88:TYR:CE1	1:B:109:LYS:HG2	2.54	0.42
1:B:258:ASP:OD2	1:B:388:TYR:OH	2.35	0.42
1:A:98:GLY:C	1:A:100:THR:N	2.76	0.42
1:B:97:LEU:HD22	1:B:97:LEU:N	2.34	0.42
1:B:19:LEU:HD23	1:B:19:LEU:O	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:188:MET:C	1:A:190:GLY:H	2.27	0.42
1:A:295:CYS:HA	1:A:324:VAL:O	2.19	0.42
1:A:408:ASN:O	1:A:409:HIS:C	2.63	0.42
1:A:57:CYS:O	1:A:58:PHE:HB2	2.18	0.42
1:A:104:TYR:C	1:A:106:ASP:N	2.78	0.42
1:B:58:PHE:CZ	1:B:125:VAL:HG12	2.54	0.42
1:A:96:TYR:HB2	1:A:97:LEU:HD22	2.01	0.42
1:A:188:MET:C	1:A:190:GLY:N	2.75	0.42
1:B:44:TYR:CE1	1:B:95:PRO:HB3	2.55	0.41
1:A:23:VAL:HG22	1:A:137:HIS:CD2	2.55	0.41
1:B:23:VAL:HG22	1:B:137:HIS:CD2	2.56	0.41
1:B:71:LEU:O	1:B:81:LEU:HA	2.19	0.41
1:A:385:ALA:HB1	1:A:390:CYS:O	2.20	0.41
1:B:37:THR:O	1:B:79:ASP:HB3	2.20	0.41
1:B:220:GLU:HB3	1:B:226:PHE:CD2	2.55	0.41
1:A:269:ASN:C	1:A:269:ASN:ND2	2.75	0.41
1:B:103:ASP:HB3	1:B:107:ASN:ND2	2.36	0.41
1:B:355:ASN:OD1	1:B:357:PRO:HD2	2.20	0.41
1:B:44:TYR:CZ	1:B:95:PRO:HB3	2.56	0.41
1:A:165:ILE:HG23	1:A:228:MET:HG3	2.02	0.41
1:A:10:ILE:O	1:A:10:ILE:HG23	2.21	0.41
1:A:67:LYS:HD2	1:A:67:LYS:N	2.36	0.41
1:A:93:GLY:O	1:A:94:GLY:C	2.61	0.41
1:B:132:ASP:HB3	1:B:476:ILE:HD13	2.02	0.41
1:B:412:LEU:HD23	1:B:412:LEU:HA	1.90	0.41
1:A:270:LEU:O	1:A:271:LYS:C	2.63	0.41
1:B:45:PRO:O	1:B:49:ALA:N	2.50	0.41
1:B:175:ILE:HD13	1:B:176:PHE:N	2.35	0.41
1:A:81:LEU:HD23	1:A:81:LEU:HA	1.93	0.40
1:A:115:SER:HB3	1:A:144:THR:N	2.35	0.40
1:A:394:VAL:O	1:A:421:GLN:HA	2.22	0.40
1:B:11:TYR:HA	1:B:12:PRO:HA	1.83	0.40
1:B:40:LEU:HD22	1:B:41:ILE:N	2.35	0.40
1:B:60:PHE:O	1:B:62:ASP:CG	2.65	0.40
1:B:298:SER:HB2	3:B:2573:HOH:O	2.20	0.40
1:B:436:ARG:HD2	1:B:440:TYR:OH	2.21	0.40
1:A:334:GLU:HB3	1:A:352:ILE:HD13	2.03	0.40
1:B:52:THR:O	1:B:53:ASP:HB2	2.21	0.40
1:B:416:ALA:HA	1:B:453:LEU:HD21	2.03	0.40
1:A:51:VAL:HG12	1:A:52:THR:N	2.37	0.40
1:B:86:PRO:HB3	1:B:90:GLU:OE1	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:245:VAL:HG12	1:B:246:ASN:O	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	475/485 (98%)	454 (96%)	19 (4%)	2 (0%)	30	38
1	B	476/485 (98%)	444 (93%)	26 (6%)	6 (1%)	9	10
All	All	951/970 (98%)	898 (94%)	45 (5%)	8 (1%)	16	20

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	61	THR
1	B	62	ASP
1	B	98	GLY
1	A	330	ASP
1	B	90	GLU
1	B	477	SER
1	A	476	ILE
1	B	10	ILE

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	371/379 (98%)	351 (95%)	20 (5%)	20	29
1	B	369/379 (97%)	350 (95%)	19 (5%)	21	32
All	All	740/758 (98%)	701 (95%)	39 (5%)	20	30

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	LEU
1	A	40	LEU
1	A	52	THR
1	A	71	LEU
1	A	73	VAL
1	A	74	GLN
1	A	102	LYS
1	A	106	ASP
1	A	153	THR
1	A	165	ILE
1	A	189	GLU
1	A	192	GLU
1	A	211	SER
1	A	225	GLU
1	A	272	ASN
1	A	298	SER
1	A	309	MET
1	A	359	SER
1	A	382	GLN
1	A	432	GLN
1	B	12	PRO
1	B	14	ILE
1	B	19	LEU
1	B	40	LEU
1	B	41	ILE
1	B	60	PHE
1	B	153	THR
1	B	155	GLU
1	B	175	ILE
1	B	192	GLU
1	B	212	THR
1	B	225	GLU
1	B	285	ARG
1	B	298	SER
1	B	316	ILE

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Mol	Chain	Res	Type
1	B	356	GLU
1	B	382	GLN
1	B	424	PRO
1	B	445	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	75	HIS
1	A	195	ASN
1	A	269	ASN
1	A	272	ASN
1	A	382	GLN
1	A	421	GLN
1	B	75	HIS
1	B	382	GLN
1	B	421	GLN
1	B	442	HIS
1	B	449	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](#)

2 ligands 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	ADP	A	1531	-	28,29,29	1.46	5 (17%)	43,45,45	2.31	16 (37%)
2	ADP	B	2531	-	28,29,29	1.39	3 (10%)	43,45,45	2.39	17 (39%)

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	ADP	A	1531	-	-	2/16/32/32	0/3/3/3
2	ADP	B	2531	-	-	4/16/32/32	0/3/3/3

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1531	ADP	C5-C4	4.00	1.46	1.39
2	B	2531	ADP	C5-N7	-3.87	1.32	1.39
2	A	1531	ADP	C5-N7	-3.31	1.33	1.39
2	B	2531	ADP	C5-C4	3.30	1.45	1.39
2	B	2531	ADP	PA-O3A	-3.25	1.56	1.59
2	A	1531	ADP	C4-N9	-2.90	1.31	1.37
2	A	1531	ADP	C8-N7	2.34	1.36	1.31
2	A	1531	ADP	C5-C6	2.31	1.47	1.41

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1531	ADP	O5'-PA-O1A	-6.14	84.61	108.94
2	A	1531	ADP	C5-C4-N3	-5.37	119.32	126.72
2	B	2531	ADP	O2A-PA-O5'	-5.23	83.84	107.57
2	B	2531	ADP	C5-C4-N3	-5.23	119.52	126.72
2	B	2531	ADP	N3-C4-N9	5.19	135.99	127.17
2	B	2531	ADP	O4'-C1'-N9	5.00	117.70	108.09
2	B	2531	ADP	O2A-PA-O3A	4.98	120.72	107.27
2	A	1531	ADP	O2A-PA-O5'	-4.73	86.14	107.57
2	A	1531	ADP	N3-C4-N9	4.38	134.61	127.17

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1531	ADP	O2A-PA-O3A	4.30	118.90	107.27
2	A	1531	ADP	C2-N3-C4	3.62	120.66	111.83
2	B	2531	ADP	O5'-PA-O1A	-3.47	95.17	108.94
2	B	2531	ADP	N6-C6-N1	3.42	126.00	118.38
2	B	2531	ADP	O3B-PB-O2B	3.12	119.50	107.80
2	A	1531	ADP	N3-C2-N1	-3.10	123.89	128.58
2	B	2531	ADP	C4-N9-C8	2.88	108.76	105.74
2	A	1531	ADP	C4-N9-C8	2.84	108.72	105.74
2	A	1531	ADP	O3'-C3'-C4'	-2.65	103.48	111.08
2	B	2531	ADP	C2-N3-C4	2.55	118.06	111.83
2	A	1531	ADP	C4-C5-N7	-2.52	107.69	110.58
2	B	2531	ADP	N3-C2-N1	-2.49	124.81	128.58
2	B	2531	ADP	N9-C8-N7	-2.45	110.46	113.94
2	B	2531	ADP	C5-N7-C8	2.35	107.15	103.45
2	B	2531	ADP	C5-C6-N6	-2.30	117.58	123.29
2	A	1531	ADP	O3A-PA-O1A	2.30	117.63	110.70
2	A	1531	ADP	O3A-PB-O1B	-2.22	99.35	111.04
2	A	1531	ADP	N6-C6-N1	2.21	123.29	118.38
2	A	1531	ADP	O2B-PB-O1B	2.17	119.29	110.83
2	A	1531	ADP	O3B-PB-O3A	2.16	111.87	104.64
2	B	2531	ADP	O3A-PB-O1B	-2.09	100.05	111.04
2	A	1531	ADP	N9-C8-N7	-2.07	111.00	113.94
2	B	2531	ADP	C2-N1-C6	2.02	122.05	118.73
2	B	2531	ADP	O2B-PB-O3A	-2.01	97.89	104.64

There are no chirality outliers.

All (6) torsion outliers are listed below:

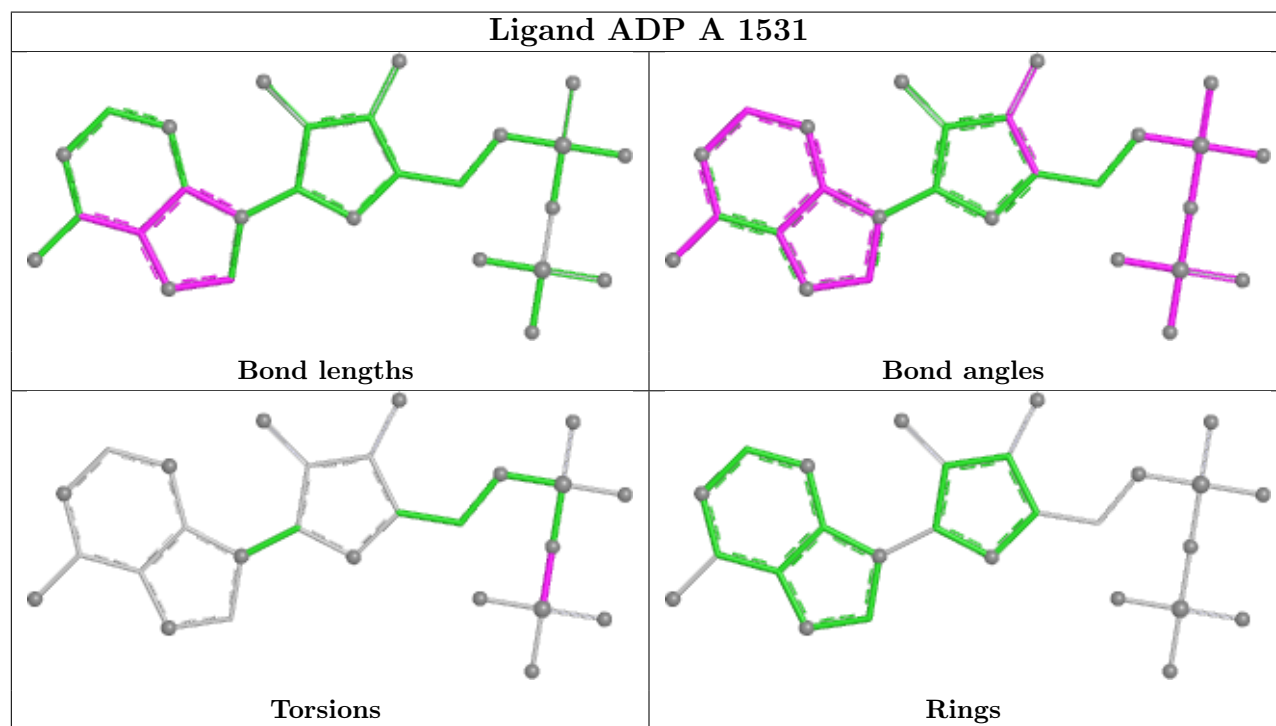
Mol	Chain	Res	Type	Atoms
2	A	1531	ADP	PA-O3A-PB-O3B
2	B	2531	ADP	C2'-C1'-N9-C8
2	B	2531	ADP	O4'-C4'-C5'-O5'
2	B	2531	ADP	O4'-C1'-N9-C8
2	B	2531	ADP	C3'-C4'-C5'-O5'
2	A	1531	ADP	PA-O3A-PB-O2B

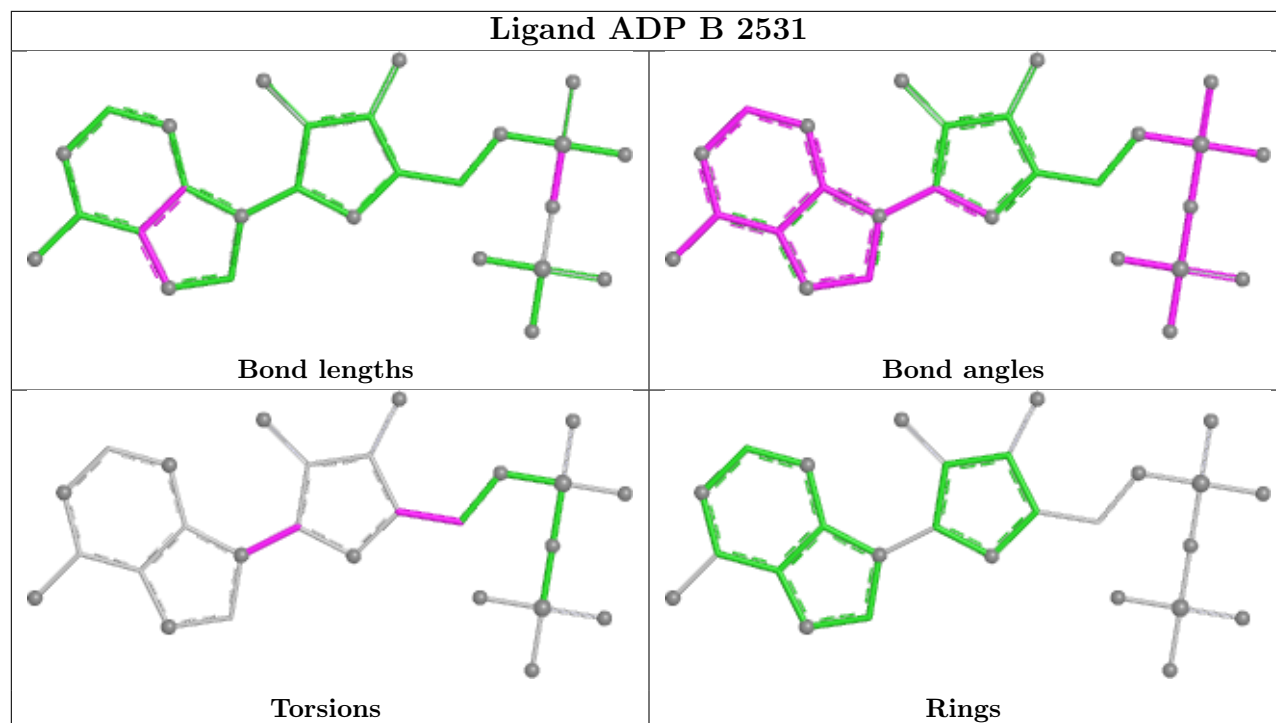
There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1531	ADP	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	477/485 (98%)	0.18	7 (1%) 72 73	12, 22, 31, 39	0
1	B	478/485 (98%)	0.20	5 (1%) 79 80	12, 22, 32, 66	0
All	All	955/970 (98%)	0.19	12 (1%) 75 76	12, 22, 32, 66	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	477	SER	3.7
1	B	476	ILE	3.5
1	A	413	ALA	3.3
1	A	416	ALA	3.1
1	B	197	VAL	2.6
1	A	409	HIS	2.5
1	A	417	ALA	2.3
1	A	100	THR	2.3
1	A	300	LEU	2.2
1	B	101	GLY	2.1
1	B	151	ALA	2.0
1	A	353	GLY	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 ⓘ

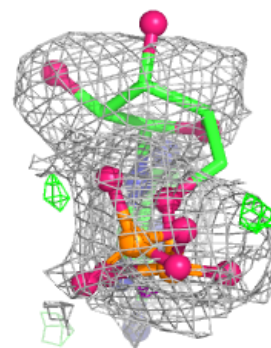
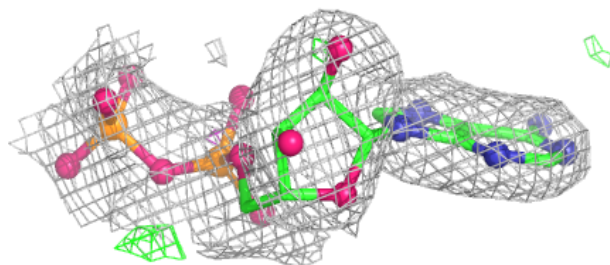
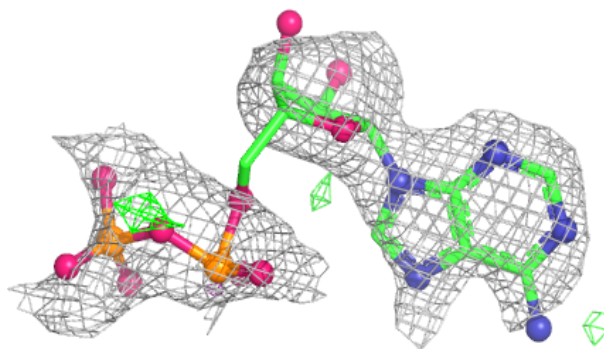
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists 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	ADP	A	1531	27/27	0.79	0.19	56,67,74,74	27
2	ADP	B	2531	27/27	0.94	0.08	32,36,60,64	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.

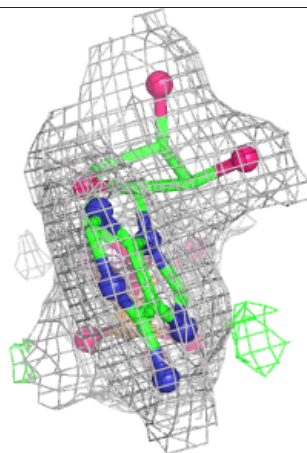
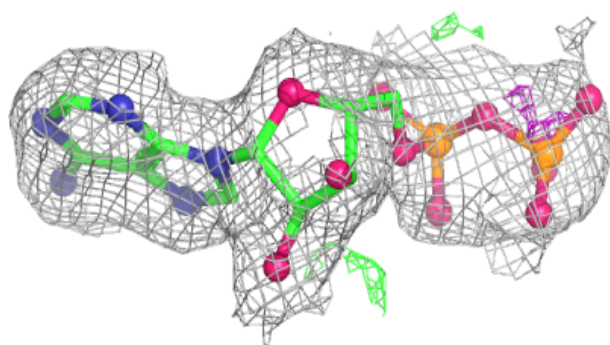
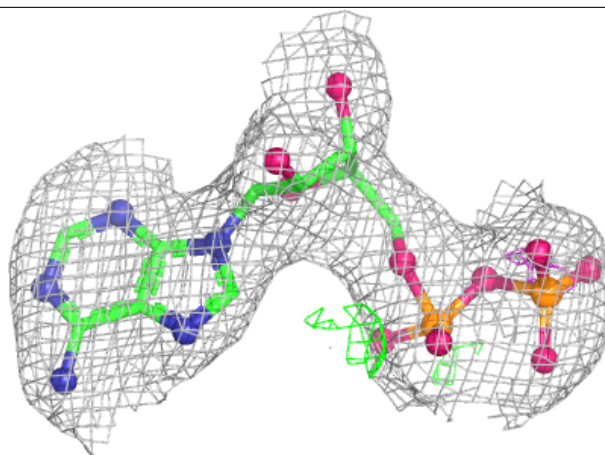
Electron density around ADP A 1531:

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 B 2531:

$2mF_o - DF_c$ (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.