



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 04:05 AM UTC

PDB ID : 2VVG / pdb_00002vvg
Title : Crystal Structure of the G.intestinalis Kinesin 2 GiKIN2a Motor Domain
Authors : Hoeng, J.C.; Loewe, J.; Dawson, S.C.; Cande, W.Z.; Sagolla, M.S.; Mancuso, J.J.
Deposited on : 2008-06-08
Resolution : 1.60 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

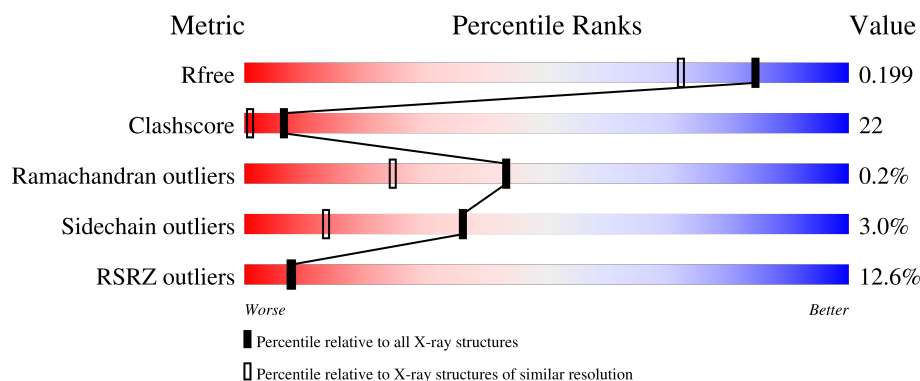
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.60 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	4673 (1.60-1.60)
Clashscore	190562	4931 (1.60-1.60)
Ramachandran outliers	187476	4831 (1.60-1.60)
Sidechain outliers	187428	4830 (1.60-1.60)
RSRZ outliers	180081	4672 (1.60-1.60)

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	350	9% 59% 27% 12%
1	B	350	13% 59% 27% 11%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 5420 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 KINESIN-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	308	Total	C	N	O	S	0	0	4
			2371	1477	422	461	11			
1	B	313	Total	C	N	O	S	0	0	4
			2411	1504	428	468	11			

- Molecule 2 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total 1	Mg 1	0	0
3	B	1	Total 1	Mg 1	0	0

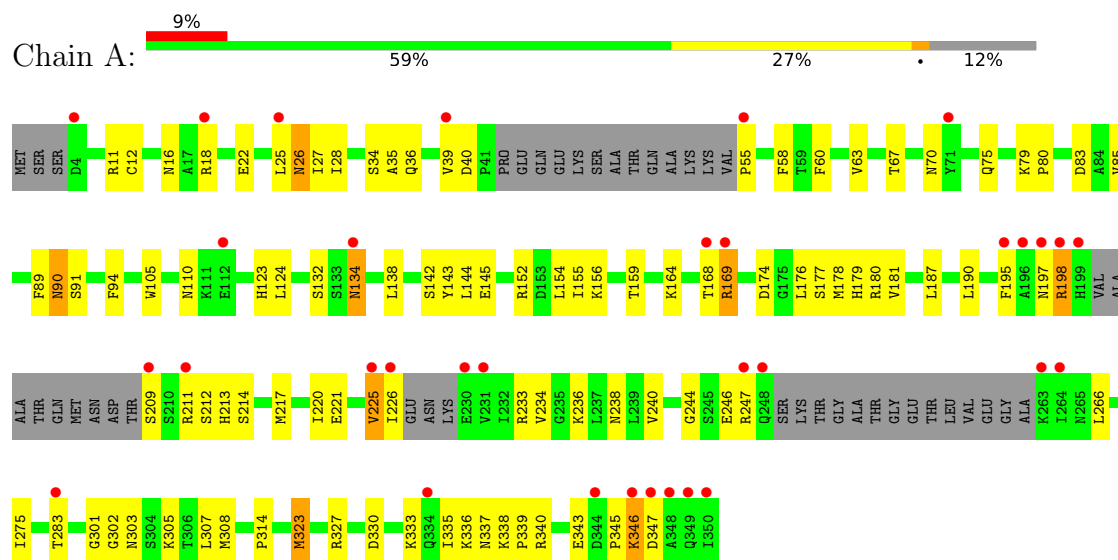
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	296	Total 296	O 296	0	0
4	B	286	Total 286	O 286	0	0

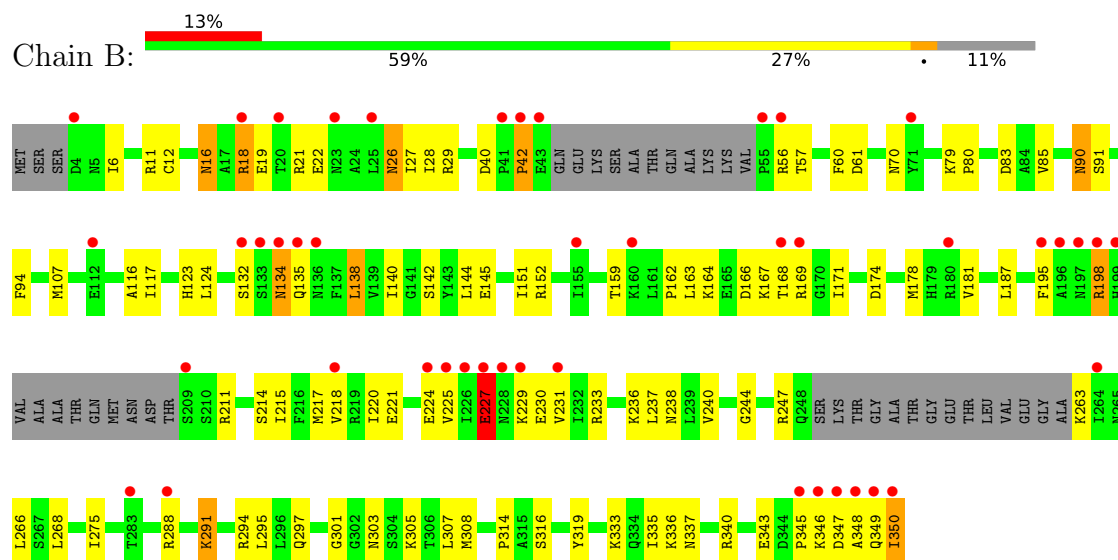
3 Residue-property plots

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: KINESIN-2



• Molecule 1: KINESIN-2



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	44.88Å 48.22Å 72.02Å 89.30° 87.52° 84.29°	Depositor
Resolution (Å)	50.00 – 1.60 50.00 – 1.60	Depositor EDS
% Data completeness (in resolution range)	95.9 (50.00-1.60) 95.9 (50.00-1.60)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.04 (at 1.60Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.203 , 0.224 0.198 , 0.199	Depositor DCC
R_{free} test set	1001 reflections (1.31%)	wwPDB-VP
Wilson B-factor (Å ²)	16.2	Xtriage
Anisotropy	0.138	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 51.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.018 for -h,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5420	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 70.44 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 3.1234e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.42	4/2402 (0.2%)	0.86	4/3237 (0.1%)
1	B	0.41	4/2444 (0.2%)	0.86	3/3295 (0.1%)
All	All	0.42	8/4846 (0.2%)	0.86	7/6532 (0.1%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	247	ARG	C-N	-5.88	1.25	1.33
1	B	198	ARG	C-N	-5.80	1.25	1.33
1	B	247	ARG	C-N	-5.79	1.25	1.33
1	A	40	ASP	C-N	-5.71	1.25	1.34
1	B	227	GLU	C-N	-5.67	1.25	1.33
1	B	42	PRO	C-N	-5.64	1.25	1.33
1	A	198	ARG	C-N	-5.35	1.25	1.33
1	A	225	VAL	C-N	-5.19	1.26	1.33

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	240	VAL	N-CA-C	6.86	117.77	108.17
1	A	246	GLU	N-CA-C	6.56	118.77	110.24
1	A	240	VAL	N-CA-C	6.23	117.27	108.36
1	B	90	ASN	N-CA-C	-6.04	100.72	110.32
1	B	305	LYS	N-CA-C	-5.38	100.93	109.76
1	A	305	LYS	N-CA-C	-5.06	100.99	109.24
1	A	90	ASN	N-CA-C	-5.01	101.55	109.96

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	2371	0	2371	99	0
1	B	2411	0	2415	112	0
2	A	27	0	12	0	0
2	B	27	0	12	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	296	0	0	23	0
4	B	286	0	0	17	0
All	All	5420	0	4810	210	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 22.

All (210) 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:A:323:MET:HA	1:A:323:MET:HE2	1.49	0.94
1:B:291:LYS:HA	4:B:2226:HOH:O	1.69	0.93
1:A:198:ARG:HH12	1:A:209:SER:HB3	1.34	0.92
1:A:58:PHE:CE1	1:A:323:MET:HE1	2.06	0.91
1:A:346:LYS:HD3	1:A:346:LYS:H	1.35	0.90
1:A:212:SER:HB2	4:A:2118:HOH:O	1.74	0.88
1:A:143:TYR:HB3	4:A:2115:HOH:O	1.75	0.85
1:A:58:PHE:HE1	1:A:323:MET:HE1	1.39	0.83
1:A:155:ILE:HG13	4:A:2115:HOH:O	1.77	0.83
1:B:117:ILE:HG13	4:B:2096:HOH:O	1.80	0.81
1:A:164:LYS:HG3	1:A:174:ASP:HB2	1.63	0.81
1:B:16:ASN:HD21	1:B:19:GLU:HG3	1.50	0.76
1:B:90:ASN:HD22	1:B:236:LYS:H	1.32	0.76
1:A:152:ARG:HG3	1:A:159:THR:HG22	1.69	0.75
1:A:26:ASN:ND2	1:A:28:ILE:H	1.84	0.75
1:A:346:LYS:HD3	1:A:346:LYS:N	2.01	0.75
1:B:163:LEU:HG	1:B:171:ILE:HD11	1.69	0.75
1:B:233:ARG:HD3	1:B:343:GLU:HG2	1.68	0.74
1:B:225:VAL:HG22	1:B:230:GLU:HG3	1.69	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:16:ASN:ND2	1:B:19:GLU:H	1.86	0.73
1:B:218:VAL:HG13	1:B:237:LEU:HB3	1.71	0.73
1:B:294:ARG:HD3	4:B:2226:HOH:O	1.87	0.72
1:B:116:ALA:HB3	4:B:2096:HOH:O	1.87	0.72
1:A:124:LEU:HD22	1:A:220:ILE:HD11	1.72	0.71
1:A:211:ARG:NE	1:A:211:ARG:HA	2.06	0.71
1:B:218:VAL:CG1	1:B:237:LEU:HB3	2.21	0.71
1:B:16:ASN:C	1:B:16:ASN:HD22	1.98	0.71
1:B:134:ASN:ND2	1:B:135:GLN:HG3	2.06	0.71
1:B:132:SER:HB2	1:B:134:ASN:ND2	2.06	0.70
1:B:198:ARG:HE	1:B:214:SER:HB3	1.57	0.70
1:A:90:ASN:HD22	1:A:236:LYS:H	1.39	0.70
1:B:152:ARG:HG3	1:B:159:THR:HG22	1.73	0.69
1:B:162:PRO:HG2	1:B:174:ASP:HB3	1.75	0.69
1:B:218:VAL:HG11	1:B:237:LEU:HD23	1.75	0.69
1:B:218:VAL:CG1	1:B:237:LEU:HD23	2.24	0.68
1:A:26:ASN:C	1:A:26:ASN:HD22	2.01	0.68
1:A:211:ARG:NH1	1:A:244:GLY:O	2.22	0.68
1:A:35:ALA:HB1	4:A:2034:HOH:O	1.93	0.67
1:A:145:GLU:HG3	4:A:2118:HOH:O	1.94	0.67
1:A:124:LEU:HD22	1:A:220:ILE:CD1	2.25	0.67
1:B:11:ARG:HH12	1:B:70:ASN:HD21	1.43	0.67
1:A:83:ASP:OD2	1:A:123:HIS:HE1	1.78	0.66
1:A:11:ARG:HH12	1:A:70:ASN:HD21	1.40	0.66
1:A:346:LYS:H	1:A:346:LYS:CD	1.99	0.66
1:B:124:LEU:HD22	1:B:220:ILE:HD11	1.78	0.66
1:B:263:LYS:HB3	1:B:266:LEU:HB2	1.78	0.66
1:A:134:ASN:HD22	1:A:134:ASN:N	1.94	0.65
1:B:83:ASP:OD2	1:B:123:HIS:HE1	1.79	0.65
1:A:323:MET:HA	1:A:323:MET:CE	2.26	0.65
1:A:11:ARG:HH12	1:A:70:ASN:ND2	1.95	0.64
1:A:26:ASN:HD22	1:A:28:ILE:H	1.44	0.63
1:B:227:GLU:HA	1:B:229:LYS:N	2.13	0.63
1:B:124:LEU:HD22	1:B:220:ILE:CD1	2.29	0.63
1:A:132:SER:HB3	1:A:134:ASN:ND2	2.14	0.63
1:B:168:THR:O	1:B:347:ASP:HB2	1.99	0.62
1:B:16:ASN:HD21	1:B:19:GLU:H	1.47	0.62
1:A:198:ARG:HE	1:A:214:SER:HB3	1.65	0.61
1:B:26:ASN:HD22	1:B:26:ASN:C	2.08	0.61
1:A:39:VAL:HG21	1:A:323:MET:HE3	1.83	0.61
1:A:134:ASN:HD22	1:A:134:ASN:H	1.49	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:163:LEU:HD22	1:B:291:LYS:HD3	1.83	0.61
1:B:18:ARG:HG3	1:B:18:ARG:HH11	1.65	0.60
1:A:303:ASN:ND2	1:A:340:ARG:H	1.99	0.60
1:A:198:ARG:HH12	1:A:209:SER:CB	2.11	0.60
1:B:27:ILE:HD13	1:B:316:SER:HA	1.84	0.60
1:A:169:ARG:HB2	1:A:169:ARG:NH1	2.16	0.60
1:A:181:VAL:HG21	1:A:187:LEU:HG	1.84	0.59
1:A:301:GLY:H	1:A:337:ASN:HD21	1.49	0.59
1:B:144:LEU:HD12	1:B:144:LEU:C	2.27	0.59
1:A:301:GLY:H	1:A:337:ASN:ND2	2.01	0.58
1:B:16:ASN:ND2	1:B:19:GLU:HG3	2.18	0.58
1:A:303:ASN:HD21	1:A:340:ARG:H	1.51	0.58
1:B:217:MET:HG2	1:B:238:ASN:HD22	1.69	0.57
1:A:63:VAL:HG13	4:A:2034:HOH:O	2.04	0.57
1:B:163:LEU:HD11	1:B:295:LEU:HD21	1.86	0.57
1:B:233:ARG:HH11	1:B:233:ARG:HG3	1.69	0.57
1:A:55:PRO:HG2	4:A:2042:HOH:O	2.04	0.57
1:A:145:GLU:OE1	1:A:198:ARG:HG3	2.05	0.57
1:B:217:MET:HG2	1:B:238:ASN:ND2	2.20	0.56
1:B:301:GLY:H	1:B:337:ASN:ND2	2.03	0.56
1:B:11:ARG:HH12	1:B:70:ASN:ND2	2.03	0.56
1:B:21:ARG:HD2	4:B:2016:HOH:O	2.04	0.56
1:A:142:SER:OG	1:A:178:MET:SD	2.64	0.56
1:A:221:GLU:HG2	1:A:234:VAL:HG22	1.88	0.56
1:A:233:ARG:HG2	1:A:343:GLU:HG2	1.87	0.56
1:B:60:PHE:HA	1:B:333:LYS:NZ	2.21	0.55
1:B:168:THR:HG23	1:B:346:LYS:HB3	1.87	0.55
1:B:301:GLY:H	1:B:337:ASN:HD21	1.53	0.55
1:A:198:ARG:NH2	4:A:2176:HOH:O	2.39	0.55
1:A:225:VAL:HG12	1:A:226:ILE:N	2.20	0.55
1:B:275:ILE:HG23	1:B:335:ILE:HD11	1.88	0.55
1:B:132:SER:HB2	1:B:134:ASN:HD21	1.71	0.55
1:B:29:ARG:HD3	1:B:40:ASP:OD2	2.07	0.54
1:B:233:ARG:CD	1:B:343:GLU:HG2	2.37	0.54
1:A:155:ILE:N	4:A:2115:HOH:O	2.41	0.54
1:B:61:ASP:OD1	1:B:333:LYS:HE2	2.08	0.54
1:B:233:ARG:HE	1:B:343:GLU:CD	2.16	0.53
1:B:85:VAL:CG2	1:B:91:SER:HB2	2.39	0.53
1:A:145:GLU:HG3	4:A:2175:HOH:O	2.08	0.53
1:A:327:ARG:NH2	4:A:2268:HOH:O	2.41	0.53
1:B:227:GLU:HG3	4:B:2189:HOH:O	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:67:THR:OG1	1:B:167:LYS:HE3	2.09	0.52
1:A:217:MET:HG2	1:A:238:ASN:ND2	2.24	0.52
1:B:303:ASN:ND2	1:B:340:ARG:H	2.07	0.52
1:A:134:ASN:ND2	1:A:134:ASN:H	2.08	0.52
1:B:291:LYS:HE3	4:B:2226:HOH:O	2.10	0.52
1:A:26:ASN:HD22	1:A:27:ILE:N	2.08	0.52
1:B:26:ASN:ND2	1:B:28:ILE:H	2.08	0.52
1:B:60:PHE:HA	1:B:333:LYS:HZ1	1.75	0.51
1:A:338:LYS:HB2	4:A:2242:HOH:O	2.11	0.51
1:B:138:LEU:HD22	1:B:140:ILE:HG13	1.91	0.51
1:A:275:ILE:HG23	1:A:335:ILE:HD11	1.93	0.51
1:A:180:ARG:HG3	1:A:180:ARG:HH11	1.75	0.51
1:B:168:THR:O	1:B:168:THR:HG22	2.11	0.50
1:A:60:PHE:HB2	4:A:2034:HOH:O	2.11	0.50
1:B:18:ARG:NE	1:B:22:GLU:OE1	2.44	0.50
1:A:330:ASP:O	1:A:333:LYS:HG2	2.12	0.50
1:B:164:LYS:HD2	4:B:2152:HOH:O	2.12	0.50
1:B:18:ARG:HG3	1:B:18:ARG:NH1	2.26	0.50
1:B:142:SER:OG	1:B:178:MET:SD	2.69	0.49
1:A:138:LEU:HD12	1:A:180:ARG:HH11	1.76	0.49
1:B:18:ARG:HE	1:B:22:GLU:CD	2.20	0.49
1:A:12:CYS:SG	1:A:314:PRO:HB3	2.53	0.49
1:A:156:LYS:HD3	4:A:2130:HOH:O	2.11	0.49
1:A:18:ARG:HB3	1:A:18:ARG:NH2	2.27	0.49
1:B:294:ARG:O	1:B:297:GLN:HG2	2.13	0.48
1:B:181:VAL:HG21	1:B:187:LEU:HG	1.95	0.48
1:B:221:GLU:CD	1:B:349:GLN:H	2.21	0.48
1:A:134:ASN:N	1:A:134:ASN:ND2	2.62	0.48
1:B:198:ARG:HG2	1:B:198:ARG:HH11	1.79	0.48
1:B:145:GLU:OE1	1:B:198:ARG:HG3	2.13	0.48
1:A:75:GLN:HE21	1:A:75:GLN:HA	1.79	0.48
1:B:94:PHE:CZ	1:B:308:MET:HB2	2.49	0.48
1:B:288:ARG:HG3	4:B:2223:HOH:O	2.13	0.48
4:A:2008:HOH:O	1:B:288:ARG:HD3	2.13	0.47
1:A:177:SER:OG	1:A:179:HIS:HE1	1.97	0.47
1:A:211:ARG:NE	1:A:211:ARG:CA	2.77	0.47
1:B:307:LEU:C	1:B:307:LEU:HD23	2.38	0.47
1:B:349:GLN:C	1:B:350:ILE:HG13	2.38	0.47
1:A:155:ILE:HD13	1:A:190:LEU:HD22	1.96	0.47
1:B:303:ASN:HD21	1:B:340:ARG:H	1.63	0.47
1:A:132:SER:HB3	1:A:134:ASN:HD21	1.80	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:211:ARG:NH1	1:B:244:GLY:O	2.48	0.47
1:B:227:GLU:HG3	1:B:229:LYS:N	2.30	0.47
1:A:75:GLN:HA	1:A:75:GLN:NE2	2.30	0.47
1:A:85:VAL:CG2	1:A:91:SER:HB2	2.44	0.47
1:B:42:PRO:HD3	1:B:319:TYR:CD2	2.50	0.47
1:B:12:CYS:SG	1:B:314:PRO:HB3	2.54	0.47
1:A:144:LEU:HD12	1:A:144:LEU:C	2.40	0.47
1:B:140:ILE:HG21	1:B:178:MET:HE3	1.96	0.46
1:B:349:GLN:O	1:B:350:ILE:OXT	2.33	0.46
1:A:176:LEU:HD21	1:A:217:MET:HE1	1.97	0.46
1:A:138:LEU:HD12	1:A:180:ARG:NH1	2.31	0.46
1:B:151:ILE:HG13	4:B:2124:HOH:O	2.15	0.46
1:B:134:ASN:HD22	1:B:134:ASN:C	2.24	0.46
1:B:198:ARG:NH2	4:B:2179:HOH:O	2.48	0.46
1:A:213:HIS:N	4:A:2118:HOH:O	2.49	0.46
1:B:56:ARG:HG2	1:B:57:THR:N	2.31	0.45
1:B:166:ASP:OD1	1:B:169:ARG:N	2.50	0.45
1:B:26:ASN:HD22	1:B:28:ILE:H	1.64	0.45
1:B:346:LYS:O	1:B:346:LYS:HG2	2.17	0.45
1:A:94:PHE:CZ	1:A:308:MET:HB2	2.52	0.45
1:B:233:ARG:HG3	4:B:2192:HOH:O	2.16	0.45
1:B:294:ARG:HB2	4:B:2226:HOH:O	2.15	0.45
1:A:197:ASN:ND2	4:A:2125:HOH:O	2.49	0.45
1:A:168:THR:O	1:A:169:ARG:HB2	2.17	0.44
1:A:283:THR:HA	4:A:2218:HOH:O	2.16	0.44
1:A:169:ARG:HB2	1:A:169:ARG:HH11	1.80	0.44
1:A:333:LYS:NZ	4:A:2272:HOH:O	2.48	0.44
1:A:25:LEU:HD12	1:A:27:ILE:HG23	1.99	0.44
1:B:229:LYS:N	4:B:2189:HOH:O	2.50	0.44
1:B:221:GLU:OE2	1:B:350:ILE:N	2.51	0.44
1:B:233:ARG:HG3	1:B:233:ARG:NH1	2.33	0.43
1:A:16:ASN:HD21	1:A:18:ARG:HH22	1.65	0.43
1:A:302:GLY:HA2	4:A:2242:HOH:O	2.18	0.43
1:B:85:VAL:HG21	1:B:237:LEU:HB2	2.00	0.43
1:A:154:LEU:HB2	4:A:2115:HOH:O	2.18	0.43
1:A:180:ARG:NH1	4:A:2159:HOH:O	2.52	0.43
1:A:26:ASN:ND2	1:A:26:ASN:C	2.72	0.43
1:B:124:LEU:CD2	1:B:220:ILE:HD11	2.47	0.43
1:A:79:LYS:HB3	1:A:80:PRO:HD3	2.00	0.43
1:A:345:PRO:C	1:A:347:ASP:H	2.27	0.43
1:B:134:ASN:ND2	1:B:134:ASN:C	2.76	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:215:ILE:HG22	1:B:217:MET:HG3	1.99	0.42
1:B:340:ARG:HD2	4:B:2115:HOH:O	2.19	0.42
1:A:58:PHE:CZ	1:A:323:MET:HE1	2.52	0.42
1:B:107:MET:HE3	4:B:2096:HOH:O	2.20	0.42
1:B:26:ASN:C	1:B:26:ASN:ND2	2.77	0.42
1:B:134:ASN:HD22	1:B:134:ASN:N	2.16	0.42
1:A:34:SER:HB2	1:A:36:GLN:HE21	1.85	0.42
1:A:307:LEU:C	1:A:307:LEU:HD23	2.45	0.41
1:A:336:LYS:HE3	1:A:336:LYS:HB2	1.89	0.41
1:A:105:TRP:CH2	1:A:110:ASN:ND2	2.89	0.41
1:B:268:LEU:HD12	1:B:268:LEU:HA	1.91	0.41
1:A:145:GLU:CG	4:A:2118:HOH:O	2.61	0.41
1:A:211:ARG:HA	1:A:211:ARG:HE	1.81	0.41
1:B:231:VAL:O	1:B:231:VAL:HG13	2.19	0.41
1:A:217:MET:HG2	1:A:238:ASN:HD22	1.85	0.41
1:B:26:ASN:HD22	1:B:27:ILE:N	2.18	0.41
1:B:79:LYS:HB3	1:B:80:PRO:HD3	2.02	0.41
1:B:233:ARG:NE	1:B:343:GLU:OE1	2.53	0.41
1:B:336:LYS:HE3	4:B:2265:HOH:O	2.21	0.40
1:B:345:PRO:HG2	1:B:348:ALA:HB2	2.02	0.40
1:B:6:ILE:HD12	1:B:335:ILE:HB	2.02	0.40
1:A:85:VAL:HA	1:A:89:PHE:O	2.21	0.40
1:B:90:ASN:ND2	1:B:236:LYS:H	2.09	0.40
1:A:18:ARG:O	1:A:22:GLU:HG3	2.21	0.40
1:A:303:ASN:HD22	1:A:339:PRO:HA	1.85	0.40
1:B:218:VAL:HG12	1:B:237:LEU:HB3	2.01	0.40
1:B:340:ARG:HG2	1:B:340:ARG:HH11	1.85	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	298/350 (85%)	287 (96%)	10 (3%)	1 (0%)	36	20
1	B	303/350 (87%)	293 (97%)	10 (3%)	0	100	100
All	All	601/700 (86%)	580 (96%)	20 (3%)	1 (0%)	43	24

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	169	ARG

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	262/300 (87%)	256 (98%)	6 (2%)	44	20
1	B	267/300 (89%)	257 (96%)	10 (4%)	30	10
All	All	529/600 (88%)	513 (97%)	16 (3%)	36	13

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	26	ASN
1	A	134	ASN
1	A	195	PHE
1	A	266	LEU
1	A	323	MET
1	A	346	LYS
1	B	16	ASN
1	B	18	ARG
1	B	26	ASN
1	B	134	ASN
1	B	138	LEU
1	B	195	PHE
1	B	224	GLU
1	B	227	GLU
1	B	291	LYS

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Mol	Chain	Res	Type
1	B	350	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (35) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	23	ASN
1	A	26	ASN
1	A	70	ASN
1	A	75	GLN
1	A	90	ASN
1	A	119	ASN
1	A	123	HIS
1	A	129	ASN
1	A	134	ASN
1	A	136	ASN
1	A	179	HIS
1	A	238	ASN
1	A	303	ASN
1	A	334	GLN
1	A	337	ASN
1	A	342	ASN
1	B	16	ASN
1	B	23	ASN
1	B	26	ASN
1	B	36	GLN
1	B	70	ASN
1	B	75	GLN
1	B	90	ASN
1	B	123	HIS
1	B	129	ASN
1	B	134	ASN
1	B	135	GLN
1	B	136	ASN
1	B	179	HIS
1	B	238	ASN
1	B	265	ASN
1	B	303	ASN
1	B	334	GLN
1	B	337	ASN
1	B	342	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 4 ligands modelled in this entry, 2 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$
2	ADP	A	1351	3	28,29,29	1.24	2 (7%)	43,45,45	1.14	3 (6%)
2	ADP	B	1351	3	28,29,29	1.29	3 (10%)	43,45,45	1.12	3 (6%)

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	1351	3	-	3/16/32/32	0/3/3/3
2	ADP	B	1351	3	-	3/16/32/32	0/3/3/3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1351	ADP	C2-N1	3.34	1.39	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1351	ADP	C2-N1	3.12	1.39	1.33
2	A	1351	ADP	PA-O3A	-2.09	1.57	1.59
2	B	1351	ADP	C1'-N9	2.04	1.52	1.46
2	B	1351	ADP	PA-O3A	-2.03	1.57	1.59

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1351	ADP	N3-C2-N1	-3.54	123.23	128.58
2	B	1351	ADP	N3-C2-N1	-3.49	123.30	128.58
2	B	1351	ADP	C2-N1-C6	2.84	123.40	118.73
2	A	1351	ADP	C2-N1-C6	2.81	123.34	118.73
2	B	1351	ADP	C3'-C2'-C1'	2.18	105.58	101.46
2	A	1351	ADP	C3'-C2'-C1'	2.15	105.54	101.46

There are no chirality outliers.

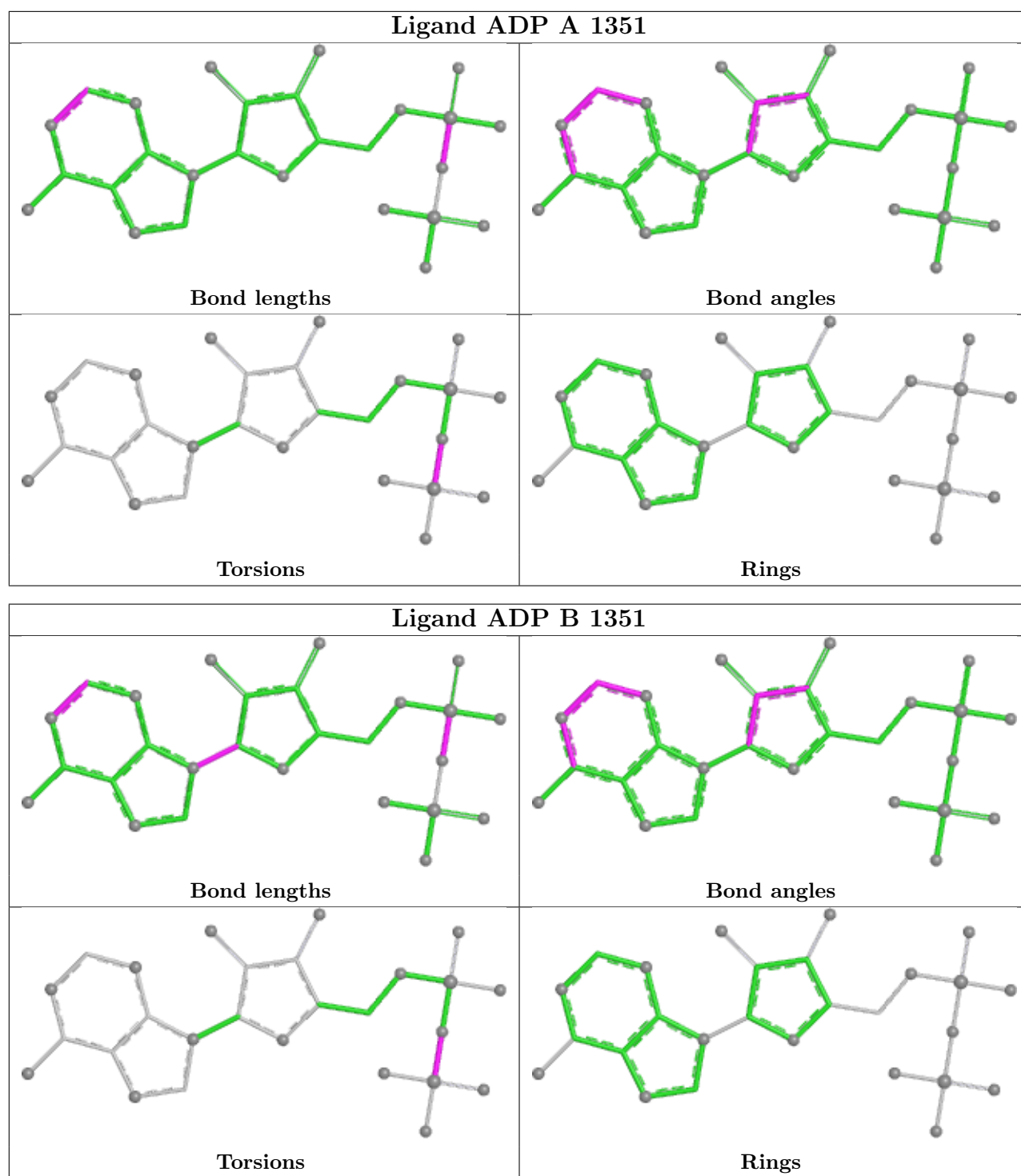
All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1351	ADP	PA-O3A-PB-O1B
2	B	1351	ADP	PA-O3A-PB-O1B
2	A	1351	ADP	PA-O3A-PB-O2B
2	A	1351	ADP	PA-O3A-PB-O3B
2	B	1351	ADP	PA-O3A-PB-O2B
2	B	1351	ADP	PA-O3A-PB-O3B

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	308/350 (88%)	1.04	33 (10%)	11 11	9, 17, 37, 51	1 (0%)
1	B	313/350 (89%)	1.10	45 (14%)	6 5	8, 18, 42, 55	0
All	All	621/700 (88%)	1.07	78 (12%)	8 7	8, 18, 40, 55	1 (0%)

All (78) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	248	GLN	12.7
1	A	199	HIS	9.3
1	B	350	ILE	8.7
1	B	199	HIS	8.2
1	A	226	ILE	7.8
1	B	228	ASN	6.2
1	A	198	ARG	5.9
1	A	195	PHE	5.6
1	B	195	PHE	5.3
1	B	168	THR	5.0
1	A	4	ASP	5.0
1	B	134	ASN	4.5
1	B	264	ILE	4.4
1	A	168	THR	4.4
1	B	226	ILE	4.3
1	B	133	SER	4.3
1	A	196	ALA	4.2
1	B	346	LYS	4.1
1	B	348	ALA	4.0
1	B	198	ARG	3.9
1	B	196	ALA	3.8
1	B	43	GLU	3.8
1	A	247	ARG	3.8
1	B	25	LEU	3.5

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Mol	Chain	Res	Type	RSRZ
1	B	227	GLU	3.5
1	B	56	ARG	3.5
1	B	347	ASP	3.5
1	A	197	ASN	3.3
1	B	349	GLN	3.1
1	B	71	TYR	3.1
1	A	25	LEU	3.1
1	B	197	ASN	3.1
1	A	112	GLU	3.1
1	B	209	SER	3.1
1	B	55	PRO	3.0
1	A	71	TYR	3.0
1	B	218	VAL	2.9
1	B	345	PRO	2.9
1	B	4	ASP	2.8
1	A	18	ARG	2.8
1	B	23	ASN	2.7
1	B	135	GLN	2.7
1	A	346	LYS	2.7
1	A	230	GLU	2.7
1	A	169	ARG	2.7
1	A	211	ARG	2.7
1	B	169	ARG	2.6
1	B	41	PRO	2.6
1	A	347	ASP	2.6
1	A	263	LYS	2.6
1	A	231	VAL	2.6
1	A	344	ASP	2.5
1	B	224	GLU	2.5
1	A	348	ALA	2.5
1	B	288	ARG	2.4
1	A	134	ASN	2.4
1	B	112	GLU	2.4
1	A	225	VAL	2.4
1	A	55	PRO	2.3
1	B	283	THR	2.3
1	B	132	SER	2.3
1	B	229	LYS	2.3
1	A	350	ILE	2.3
1	A	349	GLN	2.2
1	A	283	THR	2.2
1	A	264	ILE	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	42	PRO	2.2
1	A	209	SER	2.2
1	B	18	ARG	2.1
1	B	180	ARG	2.1
1	A	334	GLN	2.1
1	B	20	THR	2.1
1	A	39	VAL	2.1
1	B	231	VAL	2.1
1	B	160	LYS	2.1
1	B	136	ASN	2.0
1	B	155	ILE	2.0
1	B	225	VAL	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](#)

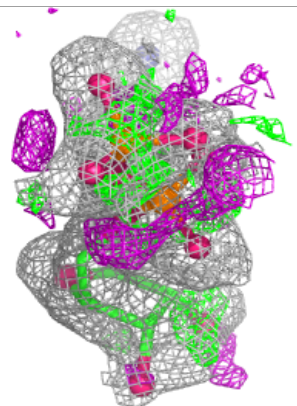
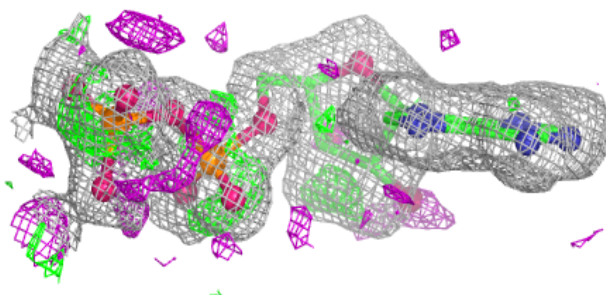
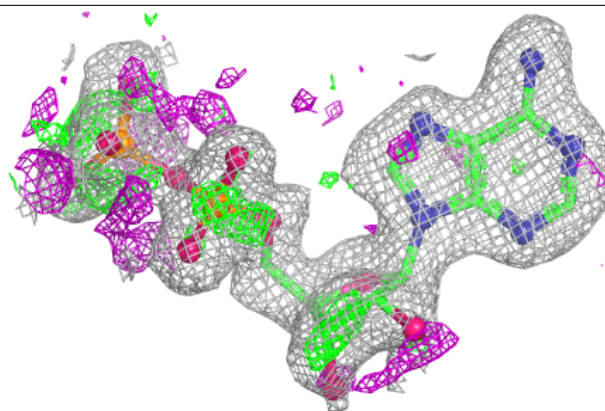
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
3	MG	A	1352	1/1	0.86	0.31	9,9,9,9	0
3	MG	B	1352	1/1	0.90	0.26	12,12,12,12	0
2	ADP	A	1351	27/27	0.93	0.10	9,14,26,36	0
2	ADP	B	1351	27/27	0.94	0.09	11,15,26,35	0

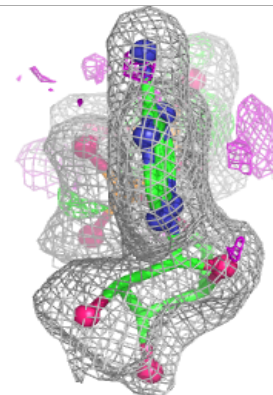
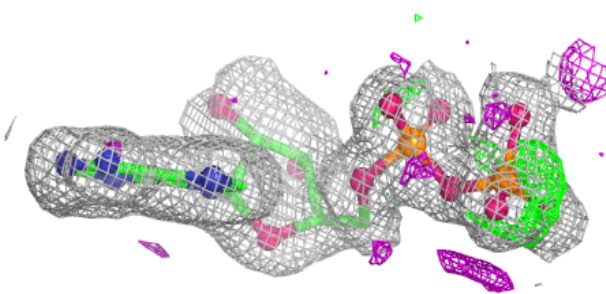
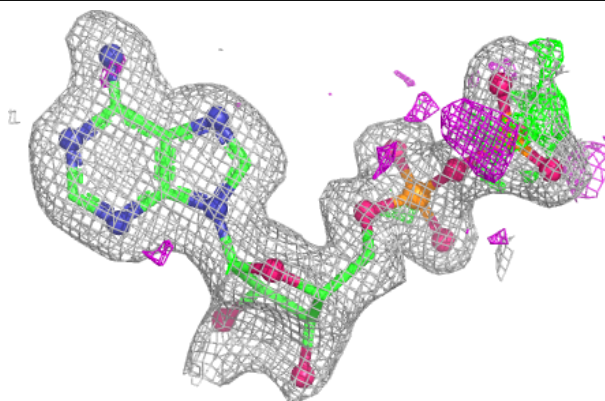
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around ADP A 1351:

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

**Electron density around ADP B 1351:**

$2mF_o-DF_c$ (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.