



# Full wwPDB X-ray Structure Validation Report ⓘ

Sep 12, 2022 – 10:14 AM JST

PDB ID : 7X4N  
Title : Crystal Structure of C. elegans kinesin-4 KLP-12 complexed with tubulin and DARPin  
Authors : Taguchi, S.; Imasaki, T.; Saijo-Hamano, Y.; Sakai, N.; Nitta, R.  
Deposited on : 2022-03-03  
Resolution : 2.88 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](mailto: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>.

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

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.30
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.30

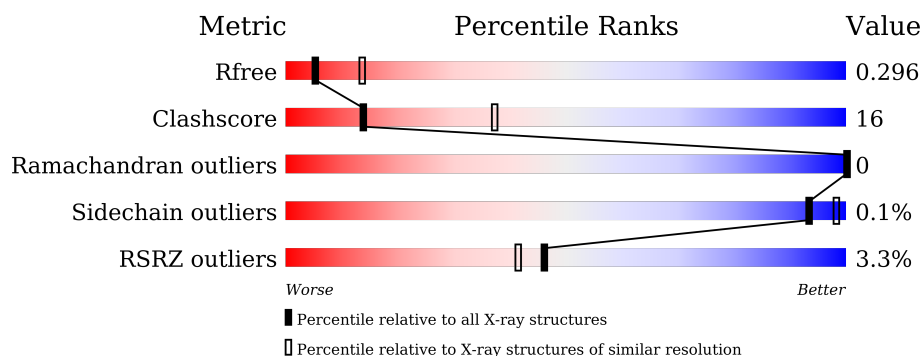
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.88 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))
$R_{free}$	130704	2691 (2.90-2.86)
Clashscore	141614	2947 (2.90-2.86)
Ramachandran outliers	138981	2868 (2.90-2.86)
Sidechain outliers	138945	2871 (2.90-2.86)
RSRZ outliers	127900	2629 (2.90-2.86)

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  $\geq 3$ , 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	451	<div> <div>3%</div> <div> <div></div> <div>65%</div> <div>31%</div> <div>..</div> </div> </div>
2	B	445	<div> <div>%</div> <div> <div></div> <div>61%</div> <div>35%</div> <div>..</div> </div> </div>
3	C	159	<div> <div>4%</div> <div> <div></div> <div>46%</div> <div>36%</div> <div>18%</div> </div> </div>
4	E	365	<div> <div>5%</div> <div> <div></div> <div>68%</div> <div>22%</div> <div>9%</div> </div> </div>

## 2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 10394 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 Tubulin alpha-1A chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	435	Total	C	N	O	S	0	0	0
			3394	2149	578	645	22			

- Molecule 2 is a protein called Tubulin beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	430	Total	C	N	O	S	6	1	0
			3382	2123	580	652	27			

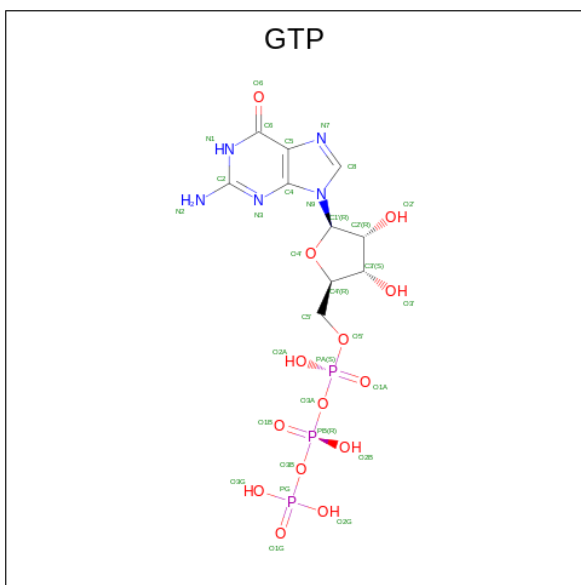
- Molecule 3 is a protein called DARPin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	131	Total	C	N	O	S	0	0	0
			965	606	170	186	3			

- Molecule 4 is a protein called Kinesin-like protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	E	332	Total	C	N	O	S	0	0	0
			2560	1587	460	497	16			

- Molecule 5 is GUANOSINE-5'-TRIPHOSPHATE (three-letter code: GTP) (formula:  $C_{10}H_{16}N_5O_{14}P_3$ ) (labeled as "Ligand of Interest" by depositor).

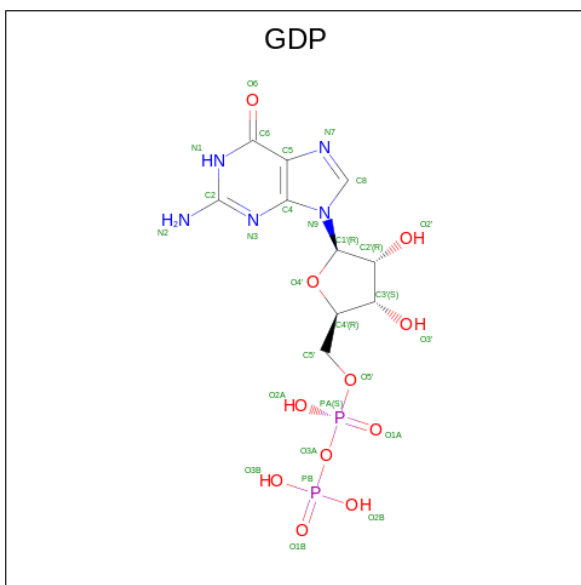


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	P	0	0
			32	10	5	14	3		

- Molecule 6 is MAGNESIUM ION (three-letter code: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

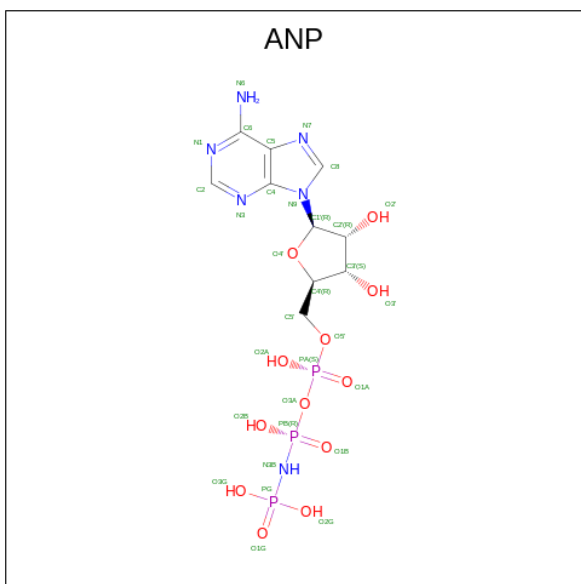
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	1	Total Mg 1 1	0	0
6	E	1	Total Mg 1 1	0	0

- Molecule 7 is GUANOSINE-5'-DIPHOSPHATE (three-letter code: GDP) (formula:  $\text{C}_{10}\text{H}_{15}\text{N}_5\text{O}_{11}\text{P}_2$ ) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	B	1	Total	C	N	O	P	0	0
			28	10	5	11	2		

- Molecule 8 is PHOSPHOAMINOPHOSPHONIC ACID-ADENYLATE ESTER (three-letter code: ANP) (formula:  $\text{C}_{10}\text{H}_{17}\text{N}_6\text{O}_{12}\text{P}_3$ ) (labeled as "Ligand of Interest" by depositor).

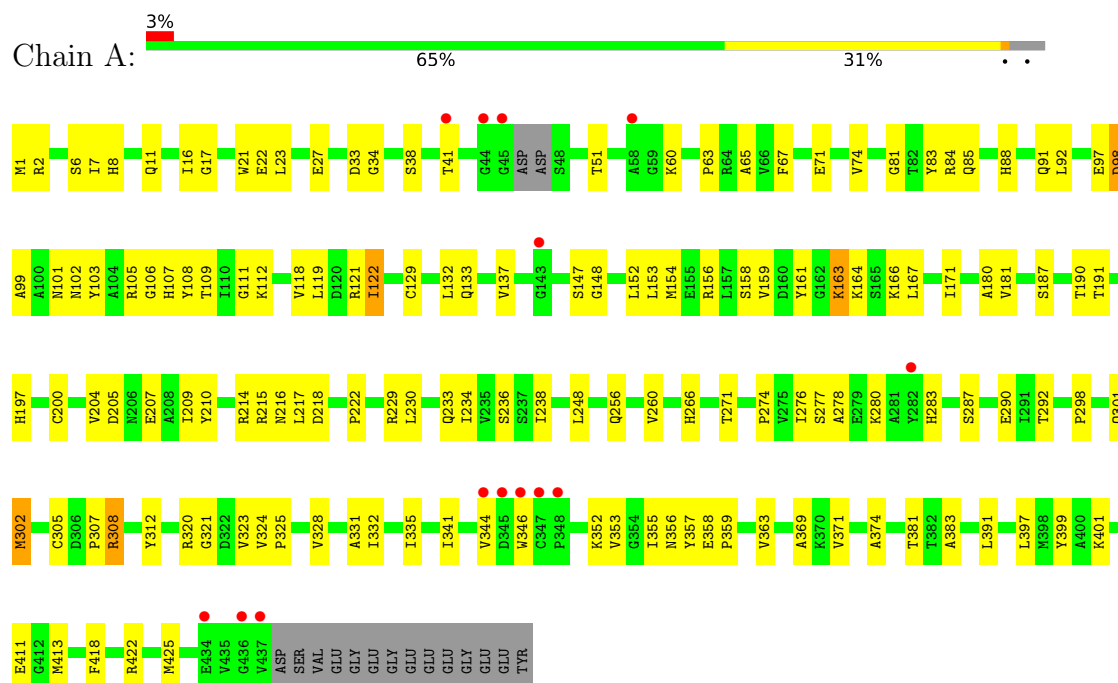


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
8	E	1	Total 31	C 10	N 6	O 12	P 3	0	0

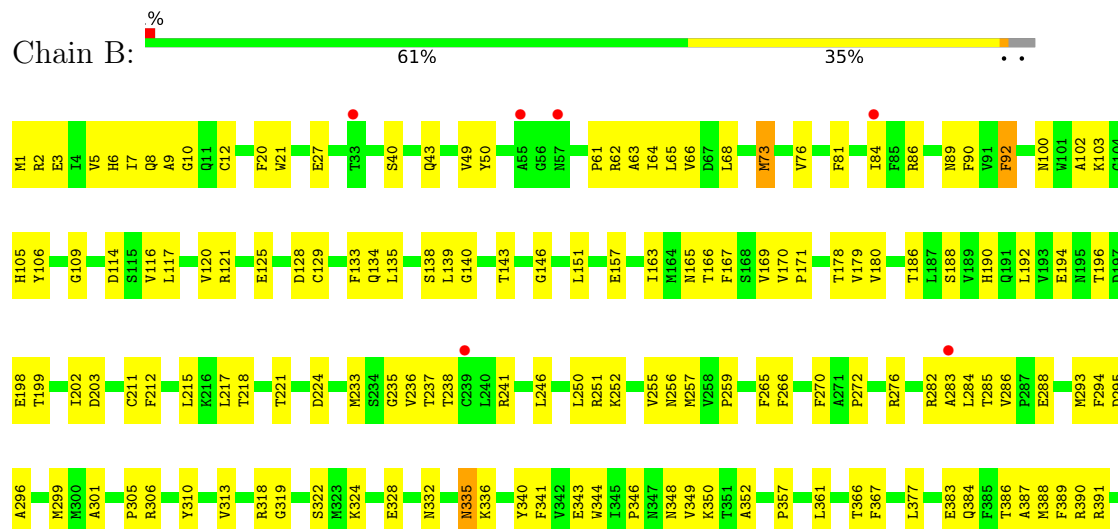
### 3 Residue-property plots

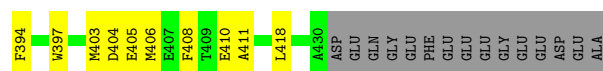
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density.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

#### • Molecule 1: Tubulin alpha-1A chain

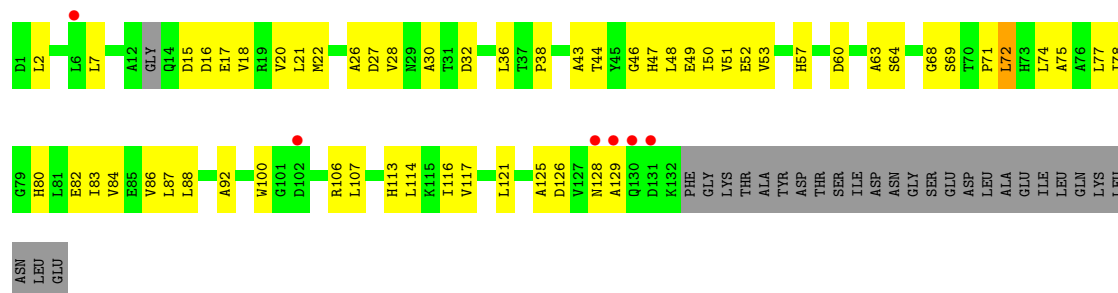


#### • Molecule 2: Tubulin beta chain

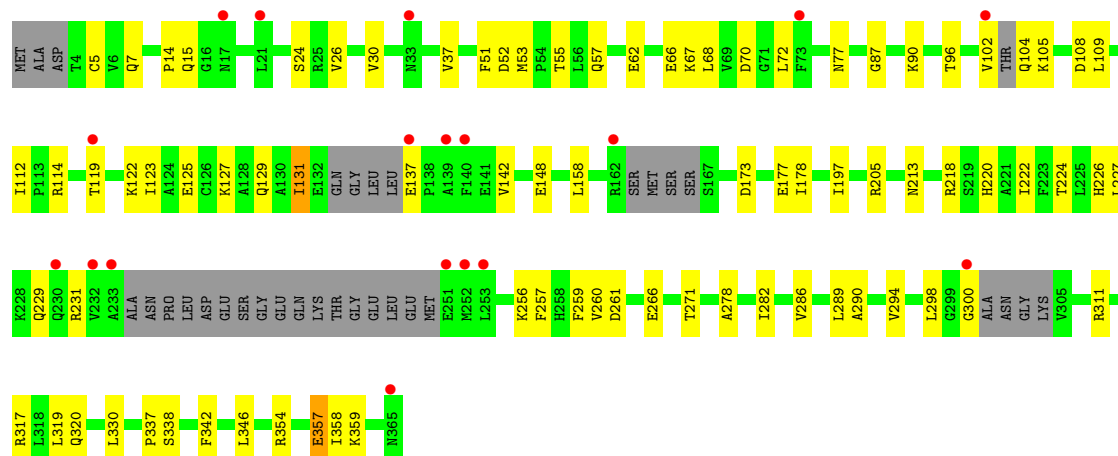




• Molecule 3: DARPin



• Molecule 4: Kinesin-like protein



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	82.33Å 80.98Å 117.75Å 90.00° 93.51° 90.00°	Depositor
Resolution (Å)	49.25 – 2.88 49.25 – 2.88	Depositor EDS
% Data completeness (in resolution range)	95.8 (49.25-2.88) 86.4 (49.25-2.88)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.27 (at 2.86Å)	Xtriage
Refinement program	PHENIX (1.18.2_3874)	Depositor
R, $R_{free}$	0.210 , 0.296 0.211 , 0.296	Depositor DCC
$R_{free}$ test set	1746 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	43.5	Xtriage
Anisotropy	0.740	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 46.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	10394	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.46% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: ANP, GTP, MG, GDP

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	0.49	0/3471	0.72	5/4710 (0.1%)
2	B	0.53	0/3457	0.80	5/4684 (0.1%)
3	C	0.45	0/977	0.74	2/1327 (0.2%)
4	E	0.47	0/2592	0.72	2/3490 (0.1%)
All	All	0.49	0/10497	0.75	14/14211 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
2	B	0	2
4	E	0	1
All	All	0	4

There are no bond length outliers.

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	293	MET	CB-CG-SD	-8.21	87.78	112.40
2	B	73	MET	CG-SD-CE	7.85	112.76	100.20
1	A	122	ILE	CG1-CB-CG2	-7.71	94.44	111.40
3	C	88	LEU	CB-CG-CD2	-6.67	99.67	111.00
4	E	289	LEU	CB-CG-CD2	6.66	122.33	111.00
1	A	121	ARG	CB-CG-CD	-6.23	95.41	111.60
2	B	217	LEU	CB-CG-CD2	5.99	121.17	111.00
4	E	131	ILE	CG1-CB-CG2	-5.82	98.60	111.40
1	A	98	ASP	CB-CG-OD1	5.81	123.53	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	72	LEU	CA-CB-CG	5.58	128.15	115.30
1	A	163	LYS	CD-CE-NZ	-5.35	99.39	111.70
2	B	92	PHE	CB-CG-CD2	5.32	124.53	120.80
2	B	157	GLU	N-CA-CB	5.29	120.12	110.60
1	A	302	MET	CB-CG-SD	-5.09	97.12	112.40

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	159	VAL	Peptide
2	B	284	LEU	Peptide
2	B	335	ASN	Peptide
4	E	357	GLU	Sidechain

## 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	3394	0	3308	113	0
2	B	3382	0	3262	126	1
3	C	965	0	982	40	0
4	E	2560	0	2538	57	1
5	A	32	0	12	0	0
6	A	1	0	0	0	0
6	E	1	0	0	0	0
7	B	28	0	12	2	0
8	E	31	0	13	4	0
All	All	10394	0	10127	324	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 16.

All (324) 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)
2:B:282:ARG:HE	2:B:283:ALA:H	1.09	0.98
1:A:180:ALA:HA	2:B:256:ASN:HD21	1.32	0.91
2:B:296:ALA:H	2:B:306:ARG:HH12	1.20	0.86
4:E:300:GLY:HA2	4:E:354:ARG:HH11	1.40	0.85
1:A:181:VAL:HG23	2:B:256:ASN:HD22	1.41	0.85
2:B:282:ARG:NE	2:B:283:ALA:H	1.76	0.83
1:A:344:VAL:HG12	1:A:346:TRP:H	1.48	0.77
3:C:69:SER:HB3	3:C:74:LEU:HD11	1.66	0.77
3:C:80:HIS:O	3:C:84:VAL:HG23	1.85	0.76
1:A:2:ARG:O	1:A:133:GLN:NE2	2.18	0.76
2:B:282:ARG:HE	2:B:283:ALA:N	1.85	0.75
2:B:324:LYS:O	2:B:328:GLU:HG3	1.85	0.75
4:E:72:LEU:HD21	4:E:257:PHE:HB2	1.68	0.75
1:A:11:GLN:HG3	1:A:74:VAL:HG21	1.68	0.75
2:B:335:ASN:OD1	2:B:335:ASN:O	2.05	0.74
2:B:170:VAL:HG11	2:B:377:LEU:HD21	1.69	0.74
3:C:71:PRO:HB2	3:C:87:LEU:HD21	1.68	0.74
2:B:116:VAL:HG11	2:B:151:LEU:HD11	1.69	0.73
3:C:121:LEU:HD23	3:C:125:ALA:HB2	1.70	0.72
2:B:306:ARG:HG3	2:B:340:TYR:CE2	2.25	0.72
1:A:88:HIS:H	1:A:91:GLN:NE2	1.87	0.72
3:C:7:LEU:HD12	3:C:38:PRO:HG3	1.71	0.72
2:B:140:GLY:HA2	2:B:171:PRO:HG3	1.72	0.71
4:E:173:ASP:N	4:E:177:GLU:O	2.23	0.69
1:A:88:HIS:H	1:A:91:GLN:HE21	1.41	0.69
1:A:102:ASN:HB3	1:A:105:ARG:HG2	1.75	0.68
1:A:397:LEU:HD12	2:B:346:PRO:HG3	1.74	0.68
2:B:169:VAL:HA	2:B:202:ILE:O	1.94	0.68
4:E:311:ARG:HG2	4:E:317:ARG:HD3	1.76	0.68
1:A:248:LEU:HD11	1:A:352:LYS:HD3	1.76	0.67
3:C:27:ASP:HB3	3:C:30:ALA:HB2	1.77	0.66
1:A:280:LYS:HD2	1:A:283:HIS:HB2	1.78	0.65
3:C:18:VAL:O	3:C:22:MET:HG3	1.97	0.65
4:E:282:ILE:O	4:E:286:VAL:HG23	1.97	0.65
2:B:6:HIS:O	2:B:63:ALA:HA	1.98	0.64
2:B:73:MET:HE2	2:B:92:PHE:HB3	1.78	0.64
2:B:114:ASP:H	2:B:117:LEU:HD12	1.62	0.64
1:A:21:TRP:CZ3	1:A:63:PRO:HB3	2.33	0.64
4:E:119:THR:O	4:E:123:ILE:HG13	1.97	0.64
4:E:226:HIS:HB2	4:E:256:LYS:NZ	2.14	0.63
4:E:5:CYS:SG	4:E:359:LYS:NZ	2.65	0.63
2:B:282:ARG:NH1	2:B:285:THR:O	2.31	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:386:THR:HG22	2:B:390:ARG:HD2	1.81	0.63
4:E:300:GLY:HA2	4:E:354:ARG:HD2	1.80	0.62
4:E:226:HIS:HB2	4:E:256:LYS:HZ3	1.64	0.62
1:A:167:LEU:CD1	1:A:200:CYS:HB3	2.30	0.62
2:B:21:TRP:CH2	2:B:61:PRO:HB3	2.35	0.62
2:B:319:GLY:HA2	2:B:357:PRO:HG3	1.81	0.62
1:A:88:HIS:HB3	1:A:91:GLN:HE21	1.65	0.61
3:C:52:GLU:HG2	3:C:86:VAL:HG11	1.81	0.61
2:B:246:LEU:HB3	2:B:352:ALA:HB2	1.83	0.61
3:C:17:GLU:O	3:C:21:LEU:HD23	2.00	0.61
3:C:72:LEU:HB3	3:C:87:LEU:HD23	1.80	0.61
2:B:296:ALA:H	2:B:306:ARG:NH1	1.95	0.61
1:A:328:VAL:HG11	1:A:353:VAL:HG11	1.83	0.61
2:B:12:CYS:SG	2:B:169:VAL:HG11	2.41	0.61
1:A:260:VAL:HG21	1:A:266:HIS:HB3	1.83	0.60
2:B:7:ILE:HG12	2:B:64:ILE:HB	1.82	0.60
2:B:102:ALA:HB2	2:B:403:MET:SD	2.40	0.60
1:A:132:LEU:HB3	1:A:164:LYS:HE3	1.83	0.60
1:A:229:ARG:HD2	1:A:363:VAL:HG21	1.84	0.60
1:A:274:PRO:HG2	1:A:371:VAL:HG11	1.84	0.60
1:A:97:GLU:HG2	2:B:1:MET:HG2	1.82	0.60
1:A:181:VAL:HG23	2:B:256:ASN:ND2	2.12	0.60
3:C:44:THR:HG23	3:C:74:LEU:HB3	1.83	0.60
1:A:99:ALA:O	1:A:101:ASN:N	2.35	0.59
2:B:65:LEU:HD22	2:B:90:PHE:CE2	2.36	0.59
2:B:143:THR:N	7:B:501:GDP:O2B	2.35	0.59
4:E:222:ILE:HG12	4:E:260:VAL:HG22	1.83	0.59
1:A:17:GLY:O	1:A:21:TRP:HD1	1.84	0.59
4:E:30:VAL:HG22	4:E:52:ASP:OD1	2.01	0.59
2:B:310:TYR:HE1	2:B:367:PHE:HZ	1.50	0.59
3:C:113:HIS:HB3	3:C:116:ILE:HD12	1.83	0.59
2:B:235:GLY:HA3	2:B:366:THR:HG21	1.84	0.58
1:A:312:TYR:CD1	1:A:341:ILE:HG23	2.38	0.58
4:E:77:ASN:OD1	4:E:256:LYS:N	2.36	0.58
1:A:106:GLY:O	1:A:111:GLY:HA3	2.03	0.58
1:A:209:ILE:HD11	1:A:302:MET:HG2	1.85	0.58
2:B:73:MET:HA	2:B:76:VAL:HG22	1.86	0.58
2:B:203:ASP:HB3	2:B:301:ALA:HA	1.84	0.58
2:B:165:ASN:HB3	2:B:198:GLU:HG3	1.86	0.58
3:C:69:SER:HB3	3:C:74:LEU:CD1	2.34	0.57
4:E:266:GLU:HB3	4:E:271:THR:HG21	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:381:THR:HG22	1:A:383:ALA:H	1.68	0.57
4:E:224:THR:HG22	4:E:256:LYS:HD2	1.85	0.57
3:C:48:LEU:HD22	3:C:82:GLU:HB3	1.87	0.57
1:A:21:TRP:CE3	1:A:63:PRO:HB3	2.40	0.57
1:A:217:LEU:HA	1:A:277:SER:HB3	1.86	0.57
2:B:332:ASN:OD1	2:B:336:LYS:HG3	2.05	0.57
1:A:312:TYR:CE1	1:A:341:ILE:HD12	2.39	0.56
1:A:401:LYS:HE2	2:B:344:TRP:CG	2.39	0.56
4:E:278:ALA:O	4:E:282:ILE:HG13	2.06	0.56
1:A:33:ASP:HA	1:A:85:GLN:CB	2.36	0.56
1:A:234:ILE:O	1:A:238:ILE:HG12	2.06	0.56
1:A:210:TYR:HE2	1:A:214:ARG:NH2	2.05	0.55
1:A:16:ILE:HD12	1:A:171:ILE:HD11	1.89	0.55
4:E:260:VAL:HG21	4:E:319:LEU:HD21	1.86	0.55
2:B:3:GLU:OE2	2:B:128:ASP:N	2.40	0.55
2:B:10:GLY:HA2	2:B:143:THR:HG23	1.88	0.55
3:C:60:ASP:HB3	3:C:63:ALA:HB2	1.89	0.55
1:A:163:LYS:HD2	1:A:164:LYS:HZ3	1.72	0.54
2:B:8:GLN:HE21	2:B:65:LEU:HG	1.72	0.54
1:A:22:GLU:HG3	1:A:83:TYR:OH	2.08	0.54
1:A:167:LEU:HD11	1:A:200:CYS:HB3	1.88	0.54
1:A:256:GLN:HA	1:A:260:VAL:HG12	1.88	0.54
1:A:328:VAL:HG21	1:A:355:ILE:HD11	1.88	0.54
4:E:357:GLU:O	4:E:359:LYS:HE3	2.08	0.54
2:B:335:ASN:OD1	2:B:335:ASN:C	2.46	0.54
3:C:49:GLU:O	3:C:53:VAL:HG23	2.07	0.54
1:A:204:VAL:HG13	1:A:302:MET:HE3	1.90	0.53
2:B:179:VAL:HG21	3:C:78:ILE:HG23	1.89	0.53
4:E:205:ARG:NH1	4:E:261:ASP:OD1	2.27	0.53
2:B:237:THR:O	2:B:241:ARG:HG3	2.08	0.53
1:A:356:ASN:O	1:A:358:GLU:N	2.42	0.53
1:A:166:LYS:HE2	1:A:197:HIS:O	2.08	0.53
1:A:33:ASP:HA	1:A:85:GLN:HB3	1.91	0.53
2:B:310:TYR:CE1	2:B:367:PHE:HZ	2.27	0.53
3:C:106:ARG:NH2	3:C:129:ALA:HA	2.24	0.53
1:A:274:PRO:HB2	1:A:276:ILE:HG12	1.90	0.53
4:E:67:LYS:HG3	4:E:68:LEU:HD22	1.91	0.53
1:A:88:HIS:CB	1:A:91:GLN:HE21	2.22	0.52
2:B:68:LEU:HD21	2:B:109:GLY:HA2	1.90	0.52
1:A:105:ARG:NH1	2:B:251:ARG:HD2	2.25	0.52
2:B:394:PHE:HD1	2:B:397:TRP:NE1	2.07	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:86:ARG:NH1	2:B:89:ASN:OD1	2.43	0.52
3:C:44:THR:HG23	3:C:74:LEU:HD23	1.91	0.52
1:A:312:TYR:HE1	1:A:341:ILE:HD12	1.75	0.52
2:B:7:ILE:O	2:B:135:LEU:HA	2.10	0.52
2:B:170:VAL:HB	2:B:203:ASP:HA	1.92	0.52
3:C:44:THR:CG2	3:C:74:LEU:HB3	2.38	0.52
2:B:12:CYS:HB2	7:B:501:GDP:C8	2.45	0.52
2:B:21:TRP:CZ2	2:B:61:PRO:HB3	2.45	0.52
3:C:44:THR:CG2	3:C:74:LEU:HD23	2.40	0.52
4:E:62:GLU:HG3	4:E:66:GLU:OE2	2.10	0.51
1:A:107:HIS:HD2	1:A:152:LEU:HB2	1.76	0.51
2:B:6:HIS:CD2	2:B:21:TRP:HE1	2.28	0.51
2:B:1:MET:HA	2:B:129:CYS:SG	2.51	0.51
4:E:87:GLY:H	8:E:401:ANP:HNB1	1.56	0.51
4:E:218:ARG:NH1	4:E:266:GLU:OE2	2.40	0.51
1:A:99:ALA:HB1	1:A:105:ARG:HB2	1.93	0.51
1:A:274:PRO:HD3	1:A:374:ALA:HA	1.92	0.51
2:B:64:ILE:HG22	2:B:66:VAL:HG23	1.91	0.51
4:E:30:VAL:HG12	4:E:37:VAL:HG23	1.92	0.51
2:B:190:HIS:CE1	2:B:410:GLU:HG3	2.45	0.50
4:E:129:GLN:HG3	4:E:131:ILE:H	1.76	0.50
2:B:313:VAL:HB	2:B:349:VAL:HG13	1.92	0.50
2:B:21:TRP:HE1	2:B:63:ALA:HB2	1.77	0.50
3:C:16:ASP:O	3:C:20:VAL:HG23	2.12	0.50
1:A:103:TYR:CE2	1:A:148:GLY:HA2	2.46	0.50
1:A:399:TYR:CE1	1:A:418:PHE:HB3	2.46	0.50
1:A:7:ILE:HB	1:A:137:VAL:HG22	1.94	0.50
1:A:298:PRO:O	1:A:301:GLN:HG2	2.12	0.50
1:A:88:HIS:N	1:A:91:GLN:HE21	2.10	0.50
3:C:43:ALA:HB1	3:C:75:ALA:HB2	1.93	0.50
4:E:24:SER:HB3	4:E:338:SER:HB3	1.92	0.50
4:E:102:VAL:HB	4:E:104:GLN:N	2.26	0.50
1:A:67:PHE:HB2	1:A:92:LEU:HD23	1.93	0.50
1:A:401:LYS:HE2	2:B:344:TRP:CD1	2.46	0.50
2:B:49:VAL:O	2:B:62:ARG:NH2	2.45	0.49
1:A:161:TYR:HA	1:A:163:LYS:HZ1	1.76	0.49
3:C:114:LEU:HA	3:C:117:VAL:HG22	1.94	0.49
1:A:8:HIS:CD2	1:A:67:PHE:CE1	3.00	0.49
1:A:321:GLY:O	1:A:357:TYR:HA	2.11	0.49
2:B:27:GLU:CD	2:B:318:ARG:HH22	2.16	0.49
2:B:138:SER:HA	2:B:169:VAL:HG12	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:230:LEU:O	1:A:234:ILE:HD12	2.13	0.49
1:A:1:MET:HG2	1:A:51:THR:HG22	1.94	0.49
4:E:109:LEU:O	4:E:114:ARG:NE	2.40	0.49
2:B:224:ASP:OD1	2:B:276:ARG:NH1	2.34	0.48
1:A:210:TYR:CE1	1:A:222:PRO:HD2	2.48	0.48
1:A:287:SER:OG	1:A:290:GLU:HG3	2.13	0.48
3:C:28:VAL:HG21	3:C:57:HIS:O	2.13	0.48
4:E:142:VAL:HG22	4:E:227:LEU:HD13	1.96	0.48
2:B:322:SER:C	2:B:324:LYS:H	2.16	0.48
4:E:213:ASN:HD22	8:E:401:ANP:HNB1	1.61	0.48
3:C:2:LEU:HB3	3:C:26:ALA:HB2	1.96	0.48
1:A:323:VAL:HG12	1:A:355:ILE:HD13	1.95	0.48
2:B:2:ARG:O	2:B:49:VAL:HG22	2.13	0.48
2:B:202:ILE:HD12	2:B:202:ILE:H	1.79	0.48
2:B:235:GLY:O	2:B:238:THR:HG22	2.13	0.48
2:B:343:GLU:OE2	2:B:343:GLU:N	2.47	0.47
4:E:70:ASP:OD1	4:E:122:LYS:HE3	2.14	0.47
4:E:122:LYS:HA	4:E:125:GLU:HB3	1.96	0.47
1:A:256:GLN:HA	1:A:260:VAL:CG1	2.44	0.47
2:B:199:THR:HB	2:B:265:PHE:CD2	2.49	0.47
4:E:15:GLN:NE2	4:E:26:VAL:HG23	2.28	0.47
2:B:192:LEU:HD22	2:B:196:THR:HG21	1.95	0.47
4:E:112:ILE:HG12	4:E:259:PHE:CG	2.50	0.47
1:A:129:CYS:SG	1:A:132:LEU:HD13	2.54	0.47
1:A:167:LEU:HD13	1:A:200:CYS:HB3	1.96	0.47
2:B:186:THR:O	2:B:411:ALA:HB1	2.15	0.47
2:B:341:PHE:HB3	2:B:348:ASN:HD21	1.78	0.46
4:E:52:ASP:OD2	4:E:53:MET:HG3	2.15	0.46
4:E:57:GLN:HB2	4:E:114:ARG:HH21	1.80	0.46
4:E:51:PHE:HB3	4:E:55:THR:HG21	1.95	0.46
2:B:20:PHE:CD1	2:B:233:MET:HE2	2.50	0.46
2:B:211:CYS:HA	2:B:215:LEU:HD12	1.98	0.46
2:B:294:PHE:O	2:B:306:ARG:NH2	2.48	0.46
3:C:72:LEU:HD22	3:C:92:ALA:HB1	1.97	0.46
1:A:214:ARG:O	1:A:218:ASP:HA	2.15	0.46
1:A:292:THR:HG22	1:A:335:ILE:HD11	1.97	0.46
4:E:158:LEU:HD21	4:E:197:ILE:HG23	1.98	0.46
2:B:198:GLU:HB2	2:B:266:PHE:CE2	2.51	0.46
4:E:72:LEU:CD2	4:E:257:PHE:HB2	2.42	0.46
2:B:212:PHE:CE2	2:B:218:THR:HA	2.51	0.46
4:E:122:LYS:HG3	4:E:125:GLU:OE1	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:134:GLN:HA	2:B:165:ASN:O	2.16	0.46
2:B:163:ILE:HG21	2:B:250:LEU:HB3	1.98	0.45
2:B:283:ALA:HA	2:B:286:VAL:HG22	1.97	0.45
1:A:119:LEU:HD12	1:A:156:ARG:NH2	2.32	0.45
2:B:387:ALA:HA	2:B:391:ARG:NH1	2.31	0.45
1:A:38:SER:HA	1:A:41:THR:HG23	1.98	0.45
2:B:221:THR:HG23	2:B:224:ASP:H	1.82	0.45
2:B:404:ASP:OD2	2:B:405:GLU:N	2.50	0.45
2:B:389:PHE:CE1	2:B:408:PHE:HB3	2.52	0.45
1:A:308:ARG:H	1:A:308:ARG:HD3	1.82	0.45
2:B:257:MET:O	2:B:259:PRO:HD3	2.17	0.45
2:B:265:PHE:CE1	2:B:418:LEU:HD21	2.51	0.45
1:A:23:LEU:O	1:A:27:GLU:HG3	2.17	0.45
1:A:422:ARG:HD2	1:A:422:ARG:O	2.17	0.45
3:C:28:VAL:O	3:C:38:PRO:HD2	2.17	0.45
3:C:64:SER:HB2	3:C:68:GLY:HA2	1.98	0.45
1:A:103:TYR:CD2	1:A:148:GLY:HA2	2.52	0.45
3:C:77:LEU:HD21	3:C:107:LEU:HD12	1.97	0.45
2:B:81:PHE:O	2:B:84:ILE:HG22	2.16	0.44
1:A:71:GLU:HB3	1:A:98:ASP:HB3	1.99	0.44
1:A:98:ASP:OD2	2:B:252:LYS:HE3	2.17	0.44
1:A:147:SER:HB2	1:A:190:THR:HB	2.00	0.44
2:B:288:GLU:O	2:B:288:GLU:HG3	2.18	0.44
1:A:229:ARG:CD	1:A:363:VAL:HG21	2.46	0.44
2:B:100:ASN:HB3	2:B:103:LYS:HB2	1.99	0.44
4:E:68:LEU:HD13	4:E:68:LEU:HA	1.83	0.44
1:A:215:ARG:NH1	1:A:216:ASN:OD1	2.50	0.44
1:A:271:THR:HA	1:A:301:GLN:HA	1.99	0.44
2:B:3:GLU:HG2	2:B:129:CYS:O	2.18	0.44
2:B:121:ARG:O	2:B:125:GLU:HG2	2.17	0.44
1:A:33:ASP:HA	1:A:85:GLN:HB2	1.99	0.44
4:E:15:GLN:HG3	4:E:337:PRO:HG2	1.98	0.44
1:A:305:CYS:O	1:A:307:PRO:HD3	2.18	0.44
2:B:86:ARG:HH11	2:B:89:ASN:HD21	1.65	0.44
2:B:114:ASP:N	2:B:117:LEU:HD12	2.31	0.44
3:C:32:ASP:OD2	3:C:36:LEU:N	2.50	0.44
2:B:9:ALA:HA	2:B:66:VAL:O	2.18	0.43
2:B:40:SER:HB3	2:B:43:GLN:HG3	1.99	0.43
2:B:73:MET:HE2	2:B:92:PHE:CB	2.44	0.43
3:C:15:ASP:OD1	3:C:15:ASP:N	2.50	0.43
4:E:96:THR:HG23	4:E:112:ILE:HD12	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:167:PHE:HE2	2:B:236:VAL:HG21	1.83	0.43
2:B:178:THR:HG22	2:B:180:VAL:HG12	1.99	0.43
2:B:165:ASN:CB	2:B:198:GLU:HG3	2.48	0.43
3:C:83:ILE:O	3:C:87:LEU:HB2	2.19	0.43
4:E:298:LEU:O	4:E:358:ILE:HG12	2.18	0.43
2:B:270:PHE:HE2	2:B:272:PRO:O	2.02	0.43
3:C:114:LEU:O	3:C:117:VAL:HG22	2.19	0.43
2:B:21:TRP:CH2	2:B:50:TYR:HB3	2.53	0.43
3:C:18:VAL:HG21	3:C:50:ILE:HD13	2.00	0.43
3:C:46:GLY:O	3:C:83:ILE:HD11	2.19	0.43
1:A:34:GLY:HA3	1:A:60:LYS:HG3	2.01	0.43
1:A:229:ARG:CZ	1:A:363:VAL:HG11	2.48	0.43
2:B:391:ARG:HD3	3:C:100:TRP:CZ3	2.53	0.43
1:A:118:VAL:O	1:A:122:ILE:HG13	2.19	0.42
2:B:190:HIS:ND1	2:B:411:ALA:HA	2.34	0.42
3:C:126:ASP:OD2	3:C:128:ASN:ND2	2.52	0.42
4:E:14:PRO:HD2	8:E:401:ANP:C6	2.49	0.42
1:A:328:VAL:HG12	1:A:332:ILE:HD12	2.01	0.42
1:A:187:SER:CB	1:A:391:LEU:HD21	2.49	0.42
2:B:250:LEU:HD23	2:B:250:LEU:HA	1.72	0.42
1:A:187:SER:HB3	1:A:391:LEU:HD21	2.01	0.42
2:B:135:LEU:HB3	2:B:166:THR:HG22	2.01	0.42
4:E:148:GLU:HA	4:E:220:HIS:O	2.20	0.42
1:A:2:ARG:C	1:A:133:GLN:HE21	2.23	0.42
1:A:229:ARG:HA	1:A:229:ARG:HD3	1.90	0.42
2:B:7:ILE:HD11	2:B:120:VAL:HG22	2.02	0.42
4:E:57:GLN:HG3	4:E:108:ASP:HA	2.02	0.42
4:E:105:LYS:HB2	4:E:105:LYS:HE3	1.87	0.42
4:E:290:ALA:O	4:E:294:VAL:HG23	2.20	0.42
1:A:331:ALA:O	1:A:335:ILE:HD12	2.19	0.42
2:B:5:VAL:O	2:B:133:PHE:HA	2.20	0.42
2:B:251:ARG:O	2:B:255:VAL:HG23	2.20	0.42
2:B:361:LEU:HD23	2:B:361:LEU:HA	1.74	0.42
4:E:178:ILE:HG13	4:E:320:GLN:HB3	2.02	0.42
3:C:47:HIS:O	3:C:51:VAL:HG23	2.20	0.41
4:E:224:THR:CG2	4:E:256:LYS:HD2	2.48	0.41
2:B:68:LEU:HD21	2:B:109:GLY:CA	2.50	0.41
2:B:105:HIS:HD2	2:B:146:GLY:O	2.02	0.41
2:B:383:GLU:HG3	2:B:384:GLN:N	2.35	0.41
4:E:90:LYS:NZ	8:E:401:ANP:O2B	2.38	0.41
4:E:7:GLN:HB2	4:E:330:LEU:HD12	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:E:257:PHE:CE2	4:E:259:PHE:CZ	3.07	0.41
1:A:341:ILE:HD13	1:A:341:ILE:HA	1.77	0.41
2:B:179:VAL:HG22	2:B:388:MET:CE	2.50	0.41
1:A:356:ASN:C	1:A:358:GLU:H	2.23	0.41
2:B:212:PHE:CD2	2:B:218:THR:HA	2.55	0.41
2:B:246:LEU:HD21	2:B:350:LYS:HB3	2.02	0.41
1:A:6:SER:O	1:A:65:ALA:HA	2.21	0.41
1:A:205:ASP:OD2	1:A:207:GLU:N	2.48	0.41
1:A:278:ALA:HA	1:A:369:ALA:HB2	2.02	0.41
2:B:299:MET:HE3	2:B:305:PRO:HG2	2.01	0.41
2:B:406:MET:HA	2:B:406:MET:HE2	2.02	0.41
1:A:154:MET:O	1:A:158:SER:HB2	2.21	0.41
1:A:233:GLN:O	1:A:236:SER:N	2.53	0.41
1:A:105:ARG:HD2	1:A:411:GLU:OE1	2.21	0.41
1:A:108:TYR:CE2	1:A:413:MET:HG2	2.56	0.41
1:A:209:ILE:CD1	1:A:302:MET:HG2	2.51	0.41
1:A:312:TYR:CE1	1:A:341:ILE:HG23	2.55	0.41
2:B:288:GLU:O	2:B:288:GLU:CG	2.69	0.41
1:A:17:GLY:O	1:A:21:TRP:CD1	2.70	0.41
2:B:394:PHE:CE2	3:C:78:ILE:HD12	2.56	0.41
4:E:229:GLN:NE2	4:E:231:ARG:HD3	2.36	0.41
1:A:7:ILE:HG21	1:A:153:LEU:HD21	2.02	0.40
1:A:81:GLY:O	1:A:84:ARG:HG2	2.21	0.40
1:A:191:THR:HB	1:A:425:MET:HE2	2.04	0.40
4:E:342:PHE:O	4:E:346:LEU:HD23	2.20	0.40
2:B:386:THR:HG22	2:B:390:ARG:CD	2.50	0.40
1:A:320:ARG:HE	1:A:320:ARG:HB3	1.56	0.40
1:A:358:GLU:OE1	1:A:359:PRO:HD2	2.21	0.40
4:E:127:LYS:HD2	4:E:137:GLU:OE1	2.21	0.40
2:B:12:CYS:HB3	2:B:138:SER:HB3	2.04	0.40
2:B:190:HIS:O	2:B:194:GLU:HG2	2.21	0.40
2:B:406:MET:HA	2:B:406:MET:CE	2.52	0.40
1:A:109:THR:O	1:A:112:LYS:HG2	2.22	0.40
1:A:324:VAL:HA	1:A:325:PRO:HD3	1.91	0.40
2:B:106:TYR:CE2	2:B:403:MET:HG2	2.57	0.40
2:B:139:LEU:HD22	2:B:188:SER:HB3	2.03	0.40

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

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:295:ASP:OD2	4:E:104:GLN:NE2[1_565]	2.15	0.05

## 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	431/451 (96%)	403 (94%)	28 (6%)	0	100	100
2	B	429/445 (96%)	396 (92%)	33 (8%)	0	100	100
3	C	127/159 (80%)	114 (90%)	13 (10%)	0	100	100
4	E	320/365 (88%)	297 (93%)	23 (7%)	0	100	100
All	All	1307/1420 (92%)	1210 (93%)	97 (7%)	0	100	100

There are no Ramachandran outliers to report.

### 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	364/377 (97%)	363 (100%)	1 (0%)	92	97
2	B	370/381 (97%)	370 (100%)	0	100	100
3	C	100/123 (81%)	100 (100%)	0	100	100
4	E	280/306 (92%)	280 (100%)	0	100	100
All	All	1114/1187 (94%)	1113 (100%)	1 (0%)	93	98

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	308	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	91	GLN
1	A	197	HIS
2	B	256	ASN
4	E	171	HIS

### 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 monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 5 ligands modelled in this entry, 2 are monoatomic - leaving 3 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	$\# Z  > 2$
5	GTP	A	501	6	26,34,34	1.01	3 (11%)	32,54,54	1.60	7 (21%)
7	GDP	B	501	-	24,30,30	0.95	1 (4%)	30,47,47	1.37	4 (13%)
8	ANP	E	401	6	29,33,33	1.25	4 (13%)	31,52,52	1.32	3 (9%)

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
5	GTP	A	501	6	-	3/18/38/38	0/3/3/3
7	GDP	B	501	-	-	3/12/32/32	0/3/3/3
8	ANP	E	401	6	-	4/14/38/38	0/3/3/3

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	E	401	ANP	PG-O1G	3.50	1.51	1.46
8	E	401	ANP	PB-O1B	3.15	1.51	1.46
5	A	501	GTP	C5-C6	-3.11	1.41	1.47
8	E	401	ANP	C8-N7	-2.41	1.30	1.34
7	B	501	GDP	O4'-C1'	2.27	1.44	1.41
8	E	401	ANP	PG-N3B	2.15	1.69	1.63
5	A	501	GTP	O4'-C4'	-2.10	1.40	1.45
5	A	501	GTP	C2-N3	2.05	1.38	1.33

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	E	401	ANP	O1G-PG-N3B	-4.49	105.15	111.77
5	A	501	GTP	PA-O3A-PB	-3.46	120.94	132.83
5	A	501	GTP	C8-N7-C5	3.25	109.17	102.99
8	E	401	ANP	O2G-PG-O1G	-3.18	105.45	113.45
5	A	501	GTP	PB-O3B-PG	-3.13	122.08	132.83
7	B	501	GDP	O3B-PB-O3A	3.11	115.06	104.64
5	A	501	GTP	O2G-PG-O3B	2.90	114.35	104.64
7	B	501	GDP	PA-O3A-PB	-2.89	122.92	132.83
5	A	501	GTP	C5-C6-N1	2.86	119.01	113.95
8	E	401	ANP	PB-O3A-PA	-2.62	123.39	132.62
5	A	501	GTP	C2-N1-C6	-2.33	120.81	125.10
7	B	501	GDP	C5-C6-N1	2.09	117.65	113.95
7	B	501	GDP	C8-N7-C5	2.08	106.95	102.99
5	A	501	GTP	O6-C6-N1	-2.05	118.23	120.65

There are no chirality outliers.

All (10) torsion outliers are listed below:

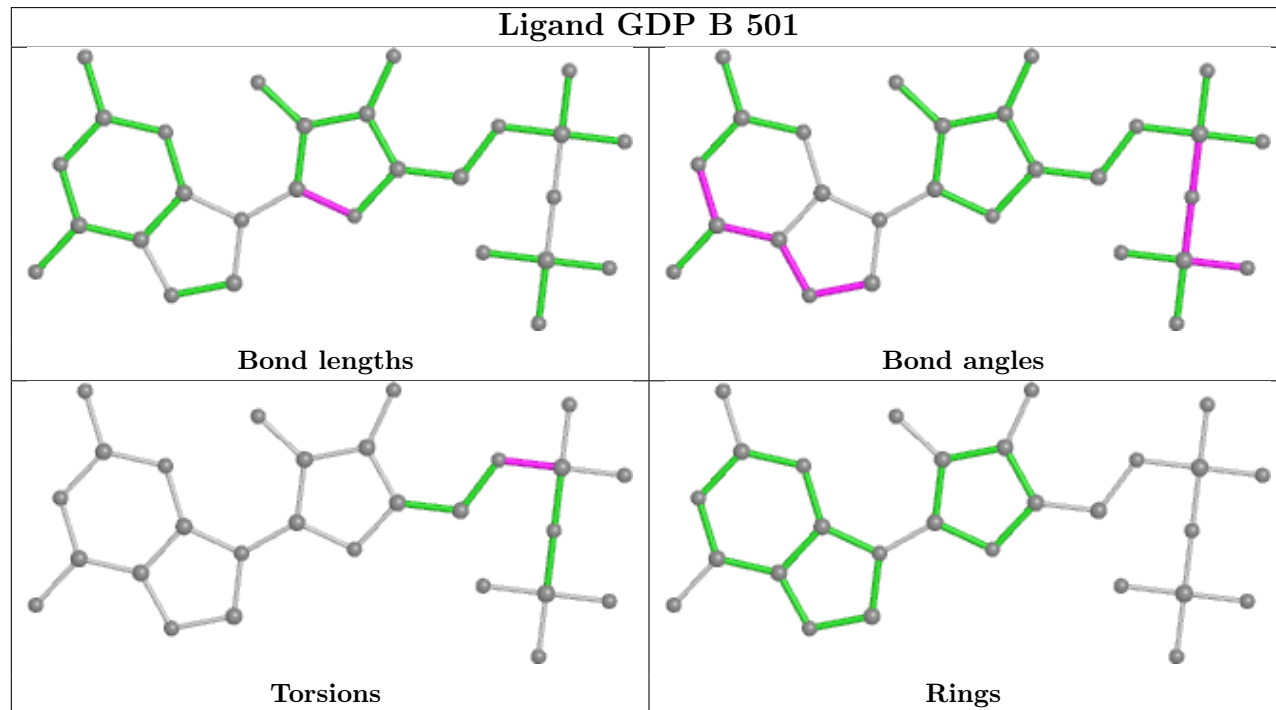
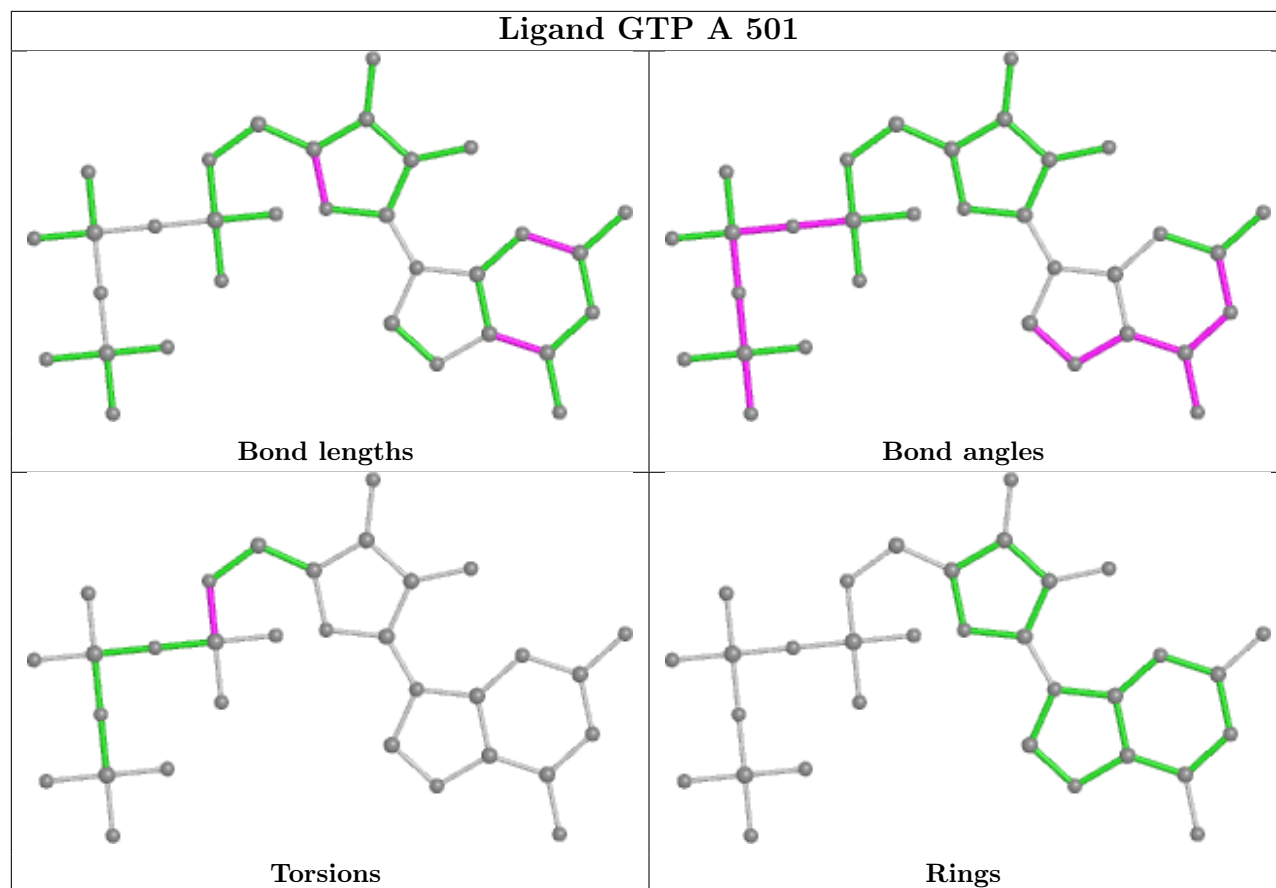
Mol	Chain	Res	Type	Atoms
5	A	501	GTP	C5'-O5'-PA-O1A
5	A	501	GTP	C5'-O5'-PA-O2A
7	B	501	GDP	C5'-O5'-PA-O1A
7	B	501	GDP	C5'-O5'-PA-O2A
8	E	401	ANP	PB-N3B-PG-O1G
8	E	401	ANP	PG-N3B-PB-O1B
8	E	401	ANP	PG-N3B-PB-O3A
8	E	401	ANP	C5'-O5'-PA-O3A
5	A	501	GTP	C5'-O5'-PA-O3A
7	B	501	GDP	C5'-O5'-PA-O3A

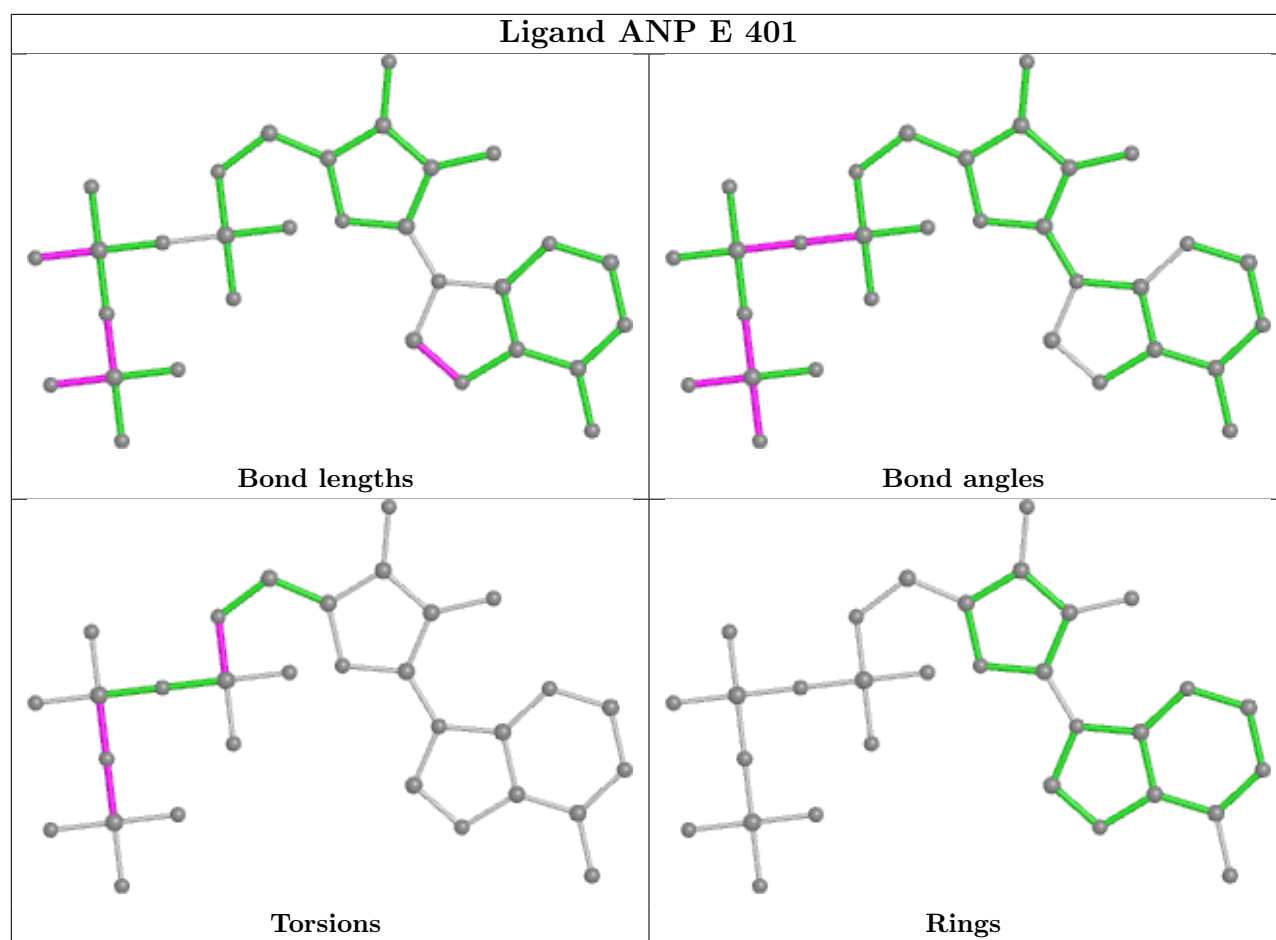
There are no ring outliers.

2 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	B	501	GDP	2	0
8	E	401	ANP	4	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 ⓘ

### 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	435/451 (96%)	0.20	14 (3%) 47 43	27, 48, 79, 112	0
2	B	430/445 (96%)	0.20	6 (1%) 75 75	28, 50, 80, 105	0
3	C	131/159 (82%)	0.53	6 (4%) 32 28	34, 54, 79, 113	0
4	E	332/365 (90%)	0.26	18 (5%) 25 22	31, 57, 89, 103	0
All	All	1328/1420 (93%)	0.25	44 (3%) 46 41	27, 51, 83, 113	0

All (44) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	C	129	ALA	5.2
2	B	283	ALA	4.7
3	C	128	ASN	4.5
3	C	130	GLN	4.3
2	B	57	ASN	3.8
1	A	346	TRP	3.6
4	E	253	LEU	3.6
1	A	58	ALA	3.5
4	E	21	LEU	3.2
4	E	252	MET	3.1
1	A	44	GLY	3.0
4	E	139	ALA	2.9
4	E	251	GLU	2.8
4	E	233	ALA	2.7
4	E	137	GLU	2.7
2	B	33	THR	2.7
1	A	437	VAL	2.7
4	E	300	GLY	2.7
4	E	232	VAL	2.7
1	A	344	VAL	2.6
3	C	102	ASP	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	41	THR	2.6
1	A	45	GLY	2.6
2	B	84	ILE	2.6
1	A	347	CYS	2.5
4	E	230	GLN	2.5
1	A	434	GLU	2.5
1	A	282	TYR	2.5
4	E	33	ASN	2.4
4	E	162	ARG	2.4
4	E	140	PHE	2.3
2	B	55	ALA	2.3
2	B	239	CYS	2.3
4	E	17	ASN	2.3
1	A	436	GLY	2.3
4	E	73	PHE	2.3
1	A	348	PRO	2.2
4	E	365	ASN	2.2
1	A	143	GLY	2.2
3	C	131	ASP	2.1
4	E	102	VAL	2.1
4	E	119	THR	2.1
1	A	345	ASP	2.1
3	C	6	LEU	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 monosaccharides 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, 95<sup>th</sup> 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.

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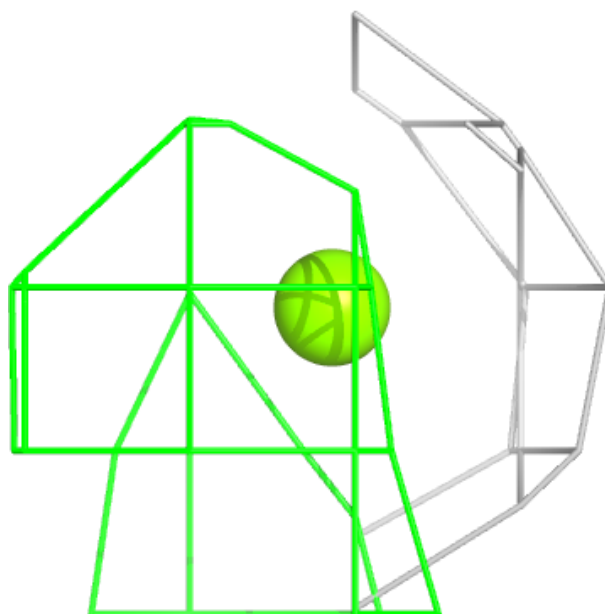
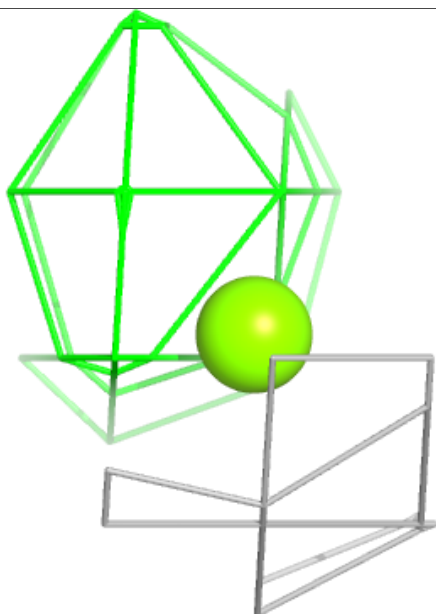
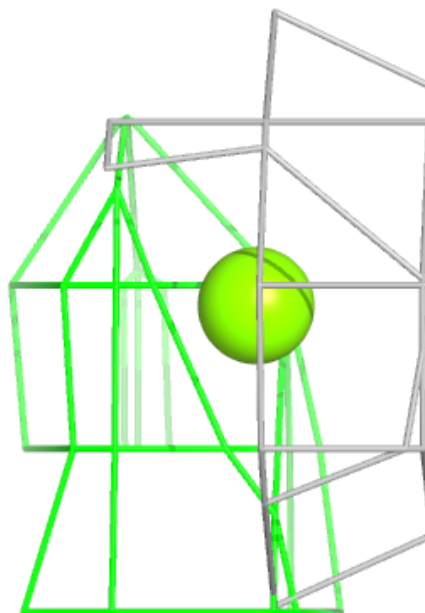
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
6	MG	A	502	1/1	0.93	0.39	46,46,46,46	0
8	ANP	E	401	31/31	0.94	0.16	30,44,52,60	0
6	MG	E	402	1/1	0.95	0.10	40,40,40,40	0
7	GDP	B	501	28/28	0.95	0.16	35,45,54,58	0
5	GTP	A	501	32/32	0.95	0.25	29,37,46,48	0

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

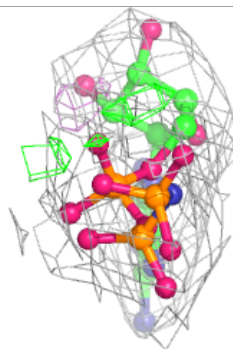
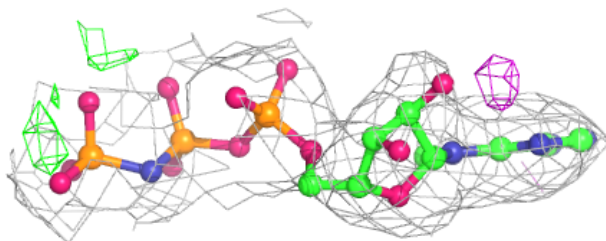
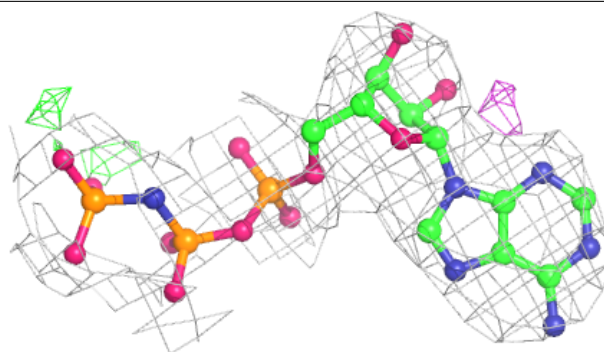
**Electron density around MG A 502:**

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



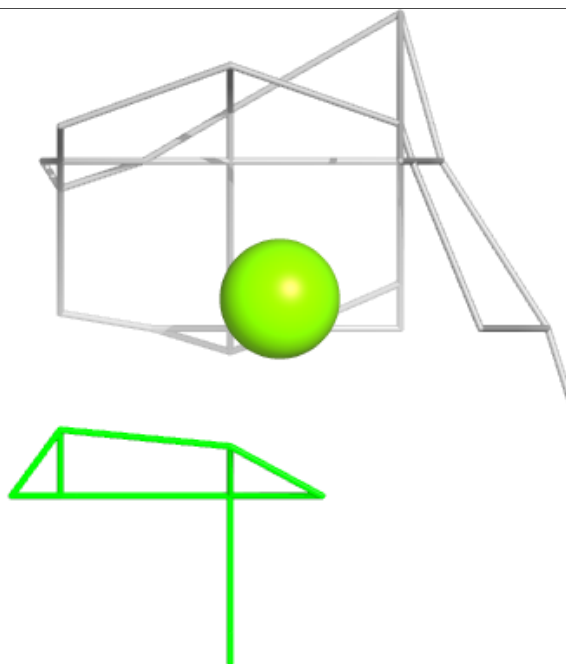
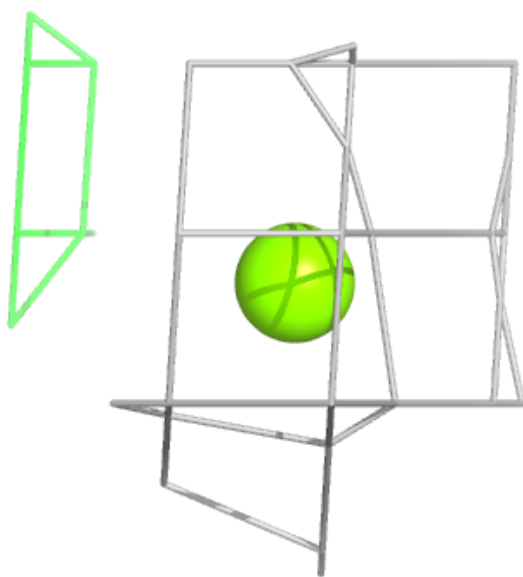
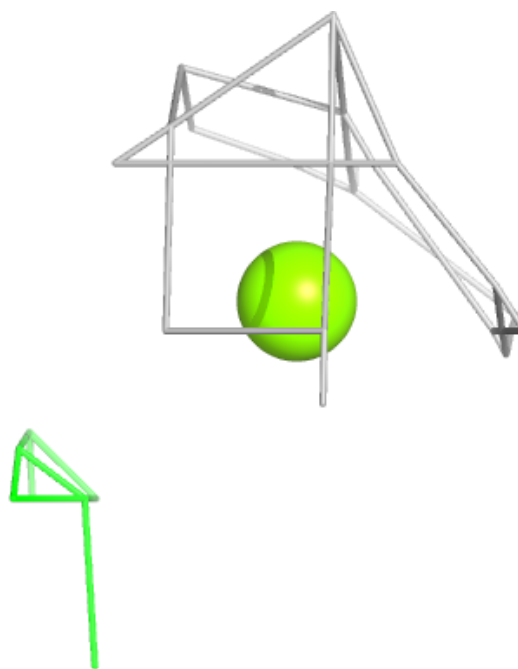
**Electron density around ANP E 401:**

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



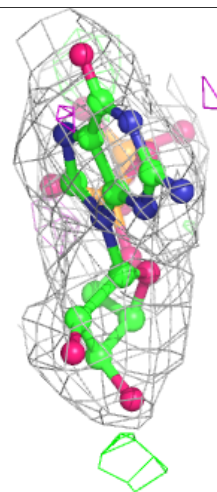
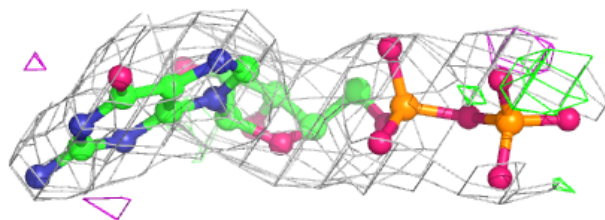
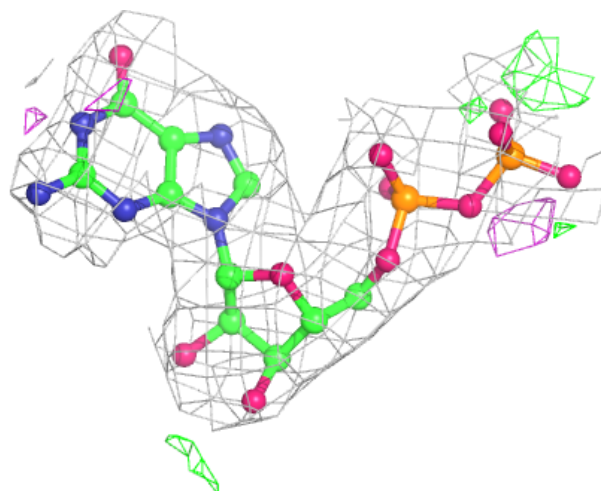
**Electron density around MG E 402:**

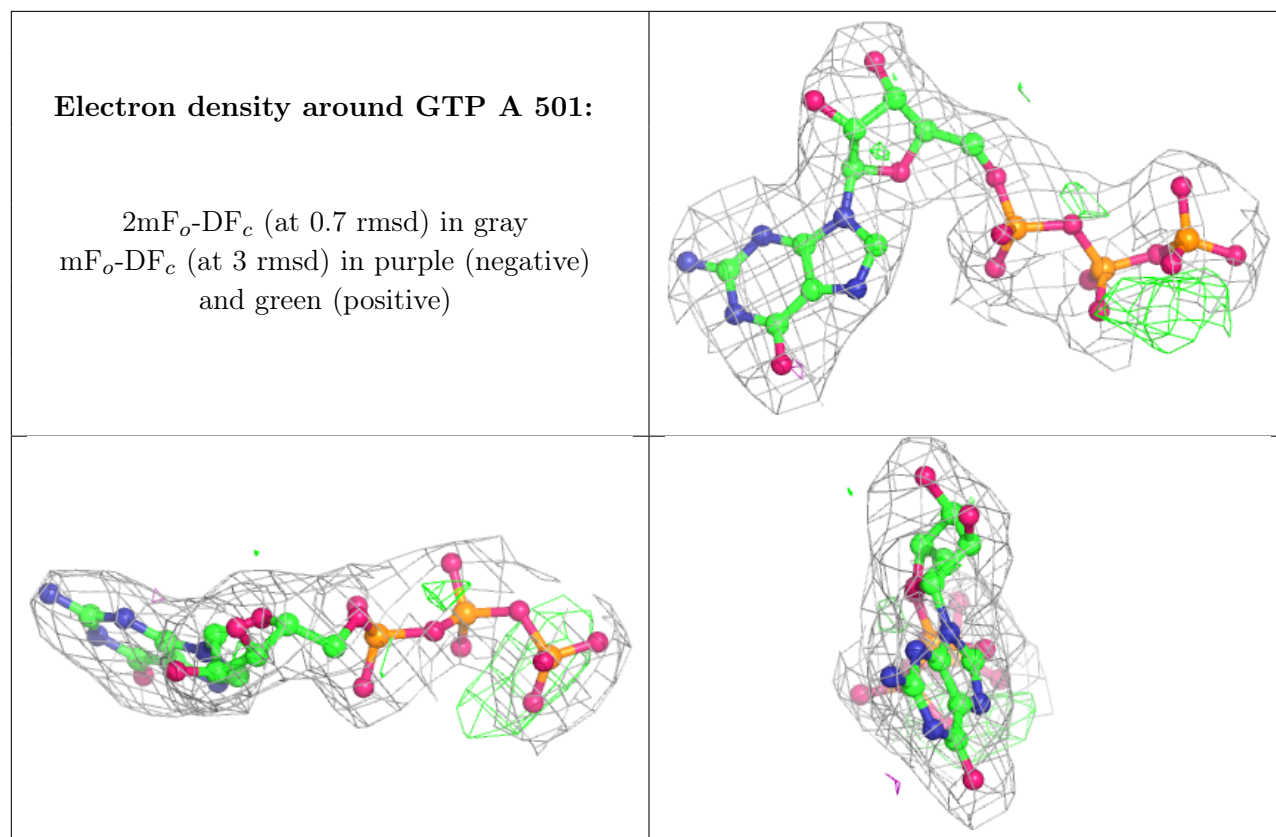
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



**Electron density around GDP B 501:**

$2mF_o-DF_c$  (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.