



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 08:02 AM UTC

PDB ID : 4REW / pdb_00004rew
Title : Crystal structure of the non-phosphorylated human alpha1 beta2 gamma1
 holo-AMPK complex
Authors : Zhou, X.E.; Ke, J.; Li, X.; Wang, L.; Gu, X.; de Waal, P.W.; Tan, M.H.E.;
 Wang, D.; Wu, D.; Xu, H.E.; Melcher, K.
Deposited on : 2014-09-24
Resolution : 4.58 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

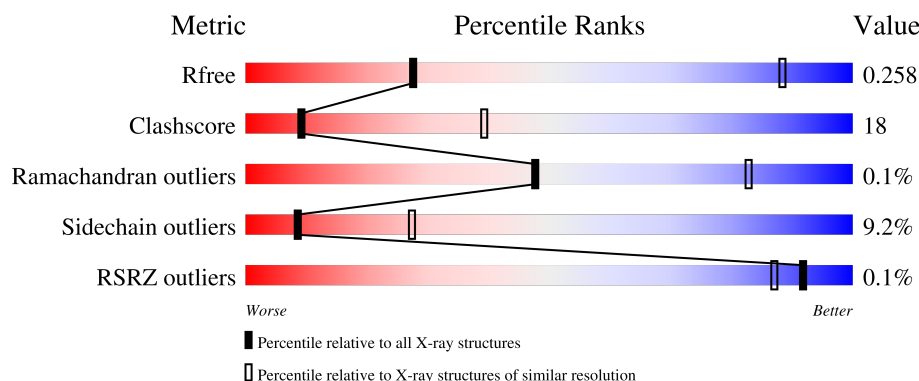
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 4.58 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	1003 (5.22-3.92)
Clashscore	190562	1019 (5.22-3.94)
Ramachandran outliers	187476	1068 (5.26-3.90)
Sidechain outliers	187428	1050 (5.26-3.90)
RSRZ outliers	180081	1139 (5.26-3.90)

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	540	
2	B	197	
3	G	304	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6499 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 5'-AMP-activated protein kinase catalytic subunit alpha-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	411	Total	C	N	O	S	0	0	0
			3326	2128	588	591	19			

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	12	SER	ARG	conflict	UNP Q13131
A	260	SER	THR	conflict	UNP Q13131
A	471	GLY	GLU	engineered mutation	UNP Q13131
A	474	ALA	GLU	engineered mutation	UNP Q13131
A	476	ALA	LYS	engineered mutation	UNP Q13131

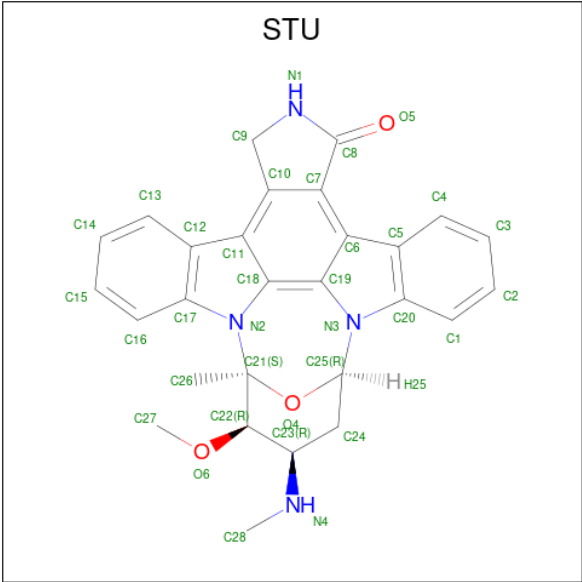
- Molecule 2 is a protein called 5'-AMP-activated protein kinase subunit beta-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	79	Total	C	N	O	S	0	0	0
			643	421	108	111	3			

- Molecule 3 is a protein called 5'-AMP-activated protein kinase subunit gamma-1.

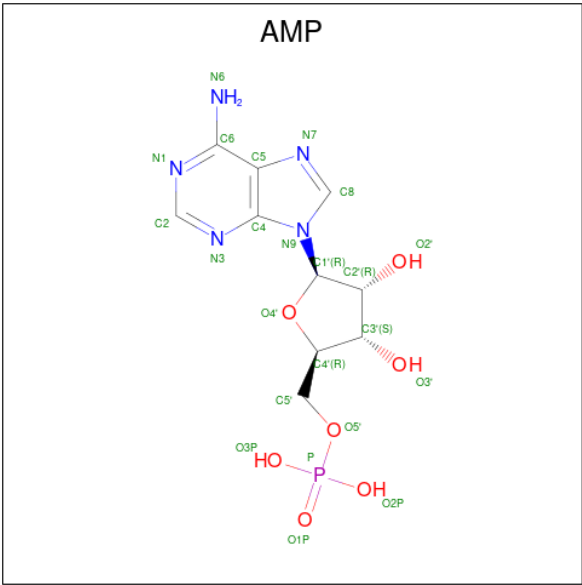
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	G	301	Total	C	N	O	S	0	0	0
			2426	1575	404	440	7			

- Molecule 4 is STAUROSPORINE (CCD ID: STU) (formula: C₂₈H₂₆N₄O₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			35	28	4	3		

- Molecule 5 is ADENOSINE MONOPHOSPHATE (CCD ID: AMP) (formula: C₁₀H₁₄N₅O₇P).

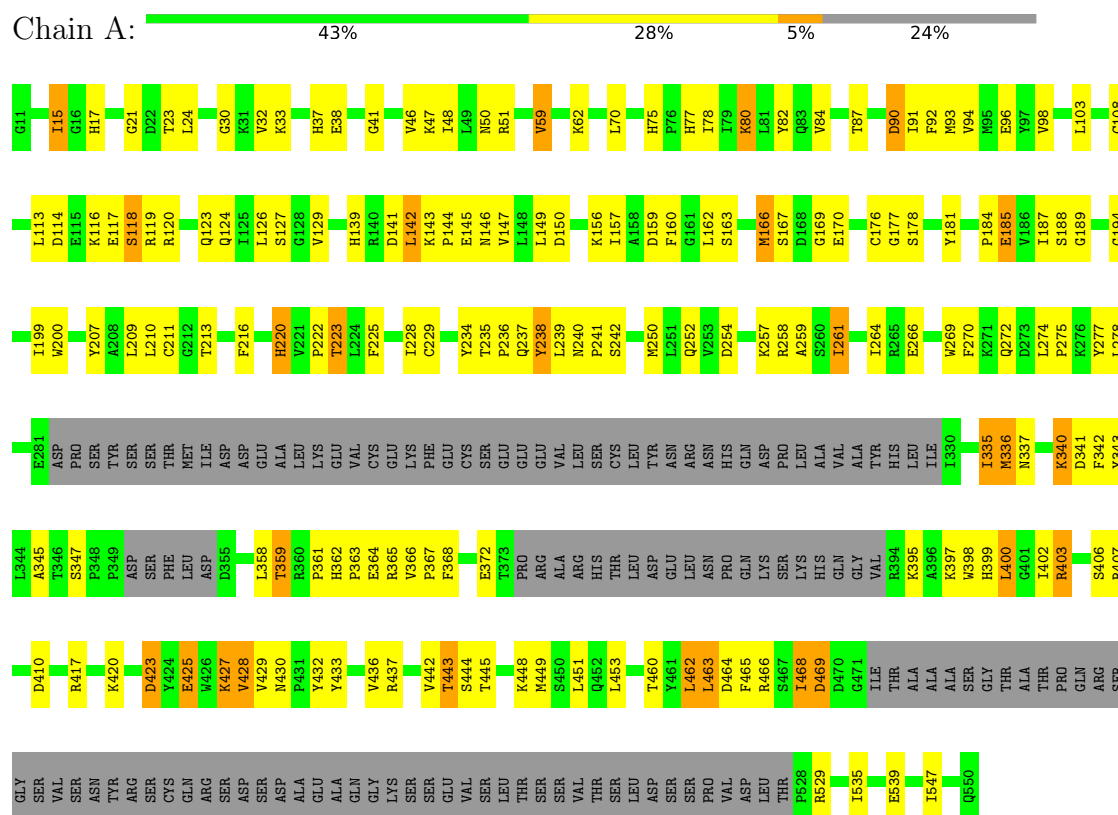


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	G	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
5	G	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
5	G	1	Total	C	N	O	P	0	0
			23	10	5	7	1		

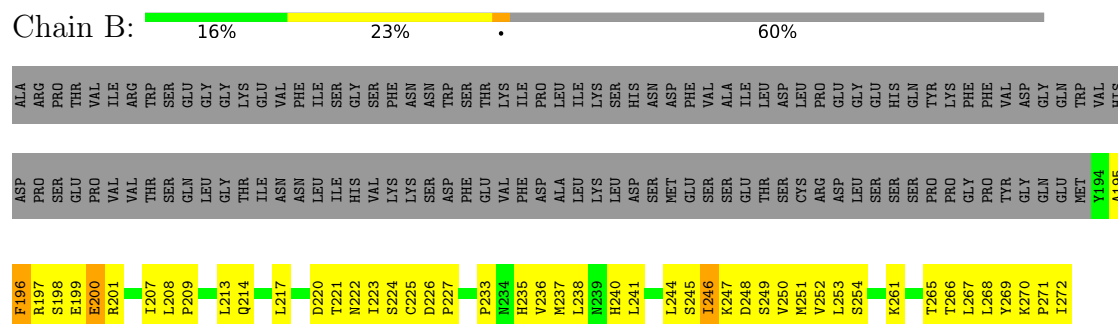
3 Residue-property plots

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: 5'-AMP-activated protein kinase catalytic subunit alpha-1



- Molecule 2: 5'-AMP-activated protein kinase subunit beta-2



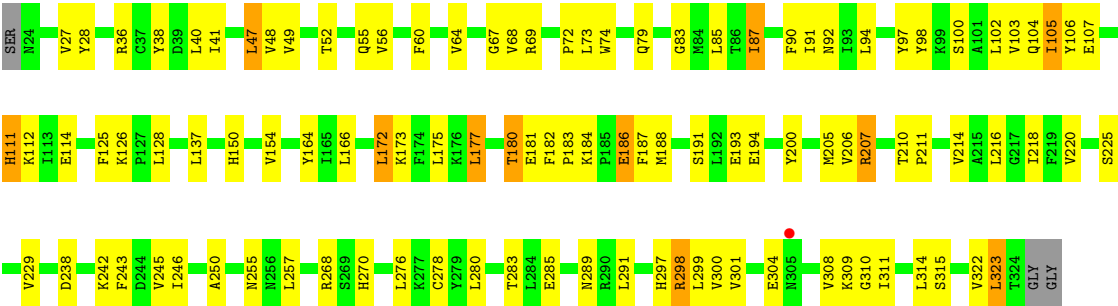
● Molecule 3: 5'-AMP-activated protein kinase subunit gamma-1

Chain G:

64%

31%

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4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	126.90Å 126.90Å 188.79Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.01 – 4.58 29.01 – 4.58	Depositor EDS
% Data completeness (in resolution range)	99.7 (29.01-4.58) 93.2 (29.01-4.58)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.98 (at 4.60Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, R_{free}	0.232 , 0.259 0.246 , 0.258	Depositor DCC
R_{free} test set	730 reflections (7.11%)	wwPDB-VP
Wilson B-factor (Å ²)	179.8	Xtriage
Anisotropy	0.104	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 167.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.044 for -h,-k,l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6499	wwPDB-VP
Average B, all atoms (Å ²)	217.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.70% 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: STU, AMP

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	0.35	0/3402	0.76	2/4589 (0.0%)
2	B	0.42	0/661	0.95	5/900 (0.6%)
3	G	0.34	0/2477	0.77	2/3363 (0.1%)
All	All	0.35	0/6540	0.78	9/8852 (0.1%)

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	G	184	LYS	N-CA-C	7.14	123.38	113.45
2	B	199	GLU	N-CA-C	7.07	120.11	110.35
3	G	182	PHE	N-CA-C	6.27	121.28	112.75
2	B	226	ASP	CA-C-N	5.24	125.54	119.47
2	B	226	ASP	C-N-CA	5.24	125.54	119.47
2	B	247	LYS	N-CA-C	-5.07	106.78	113.17
2	B	198	SER	N-CA-C	5.06	117.88	109.94
1	A	194	GLY	CA-C-N	-5.00	113.89	119.19
1	A	194	GLY	C-N-CA	-5.00	113.89	119.19

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	3326	0	3357	151	0
2	B	643	0	671	45	0
3	G	2426	0	2490	74	0
4	A	35	0	26	2	0
5	G	69	0	36	5	0
All	All	6499	0	6580	238	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 18.

All (238) 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:177:GLY:H	2:B:197:ARG:HG3	1.22	1.05
1:A:361:PRO:HB3	2:B:223:ILE:HG22	1.52	0.91
1:A:402:ILE:HB	1:A:462:LEU:HD23	1.63	0.80
1:A:177:GLY:N	2:B:197:ARG:HG3	1.97	0.80
3:G:27:VAL:HG11	3:G:183:PRO:HB2	1.62	0.79
3:G:69:ARG:HE	3:G:87:ILE:HD11	1.49	0.78
1:A:82:TYR:HB2	1:A:94:VAL:HG23	1.67	0.76
1:A:210:LEU:HD22	1:A:239:LEU:HD11	1.68	0.75
3:G:322:VAL:HG23	3:G:323:LEU:HG	1.67	0.75
1:A:50:ASN:HA	1:A:90:ASP:HB3	1.68	0.75
1:A:448:LYS:HB2	1:A:468:ILE:HG21	1.68	0.73
3:G:301:VAL:HG23	3:G:309:LYS:HB2	1.72	0.71
1:A:336:MET:O	1:A:340:LYS:N	2.24	0.70
1:A:400:LEU:HD13	1:A:464:ASP:HA	1.74	0.70
1:A:98:VAL:HB	1:A:150:ASP:HA	1.73	0.70
2:B:241:LEU:HD11	2:B:253:LEU:HD12	1.72	0.69
1:A:143:LYS:HG3	1:A:146:ASN:HB2	1.73	0.69
1:A:24:LEU:HB2	1:A:32:VAL:HG23	1.73	0.69
1:A:425:GLU:HB3	1:A:436:VAL:HG12	1.75	0.68
3:G:172:LEU:HD12	3:G:314:LEU:HD22	1.75	0.68
3:G:69:ARG:HB2	3:G:87:ILE:HD11	1.76	0.67
2:B:208:LEU:HD12	2:B:209:PRO:HD2	1.76	0.67
2:B:267:LEU:HB2	3:G:48:VAL:HG12	1.75	0.67
2:B:250:VAL:HA	2:B:271:PRO:HA	1.75	0.67
3:G:100:SER:HB3	3:G:103:VAL:HG22	1.76	0.67
1:A:46:VAL:HG12	1:A:94:VAL:HG12	1.76	0.66
3:G:83:GLY:HA2	3:G:128:LEU:HB2	1.78	0.66
3:G:105:ILE:HD13	3:G:105:ILE:H	1.60	0.65
1:A:37:HIS:HB3	1:A:41:GLY:H	1.62	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:188:SER:O	2:B:201:ARG:HD3	1.98	0.63
3:G:94:LEU:HG	3:G:257:LEU:HD11	1.82	0.62
1:A:77:HIS:CD2	1:A:124:GLN:HB3	2.35	0.61
1:A:254:ASP:HB3	1:A:257:LYS:HB2	1.81	0.61
3:G:92:ASN:HD22	3:G:220:VAL:HG12	1.65	0.61
3:G:97:TYR:HB3	3:G:107:GLU:HG3	1.82	0.61
3:G:91:ILE:HG23	3:G:216:LEU:HD22	1.83	0.61
1:A:368:PHE:CE2	3:G:67:GLY:HA2	2.36	0.60
1:A:397:LYS:HG2	2:B:244:LEU:HB2	1.84	0.59
1:A:432:TYR:HB3	1:A:453:LEU:HB3	1.84	0.59
1:A:59:VAL:HB	1:A:62:LYS:HB2	1.84	0.59
1:A:127:SER:HB3	1:A:278:LEU:HG	1.85	0.59
1:A:252:GLN:HG2	1:A:257:LYS:HB3	1.85	0.59
1:A:143:LYS:HB3	1:A:181:TYR:CE2	2.38	0.59
1:A:187:ILE:HD13	1:A:225:PHE:HB3	1.83	0.59
1:A:75:HIS:HB3	1:A:78:ILE:HB	1.83	0.58
3:G:105:ILE:HG12	3:G:105:ILE:O	2.03	0.58
2:B:217:LEU:HD21	2:B:265:THR:HG21	1.84	0.58
3:G:87:ILE:HD12	3:G:87:ILE:H	1.68	0.58
3:G:69:ARG:NH1	5:G:402:AMP:O1P	2.37	0.58
1:A:365:ARG:HB2	2:B:227:PRO:HG3	1.86	0.57
1:A:398:TRP:CD1	2:B:245:SER:HB2	2.40	0.57
2:B:195:ALA:HB1	2:B:196:PHE:HA	1.87	0.57
3:G:91:ILE:HG12	3:G:245:VAL:HG21	1.86	0.57
1:A:337:ASN:HA	1:A:340:LYS:HB2	1.87	0.56
1:A:160:PHE:O	1:A:163:SER:HB3	2.05	0.56
1:A:189:GLY:HA3	2:B:201:ARG:HB2	1.87	0.56
3:G:103:VAL:HG23	3:G:104:GLN:HG3	1.87	0.56
1:A:368:PHE:O	1:A:372:GLU:HG2	2.05	0.56
3:G:48:VAL:HG22	3:G:68:VAL:HG21	1.86	0.56
1:A:167:SER:HB2	1:A:170:GLU:HB2	1.87	0.56
3:G:188:MET:HE2	3:G:285:GLU:HG3	1.88	0.56
4:A:601:STU:H16	4:A:601:STU:H261	1.89	0.55
1:A:340:LYS:HD2	1:A:341:ASP:H	1.71	0.55
2:B:207:ILE:HD12	2:B:207:ILE:H	1.72	0.55
3:G:87:ILE:HG23	3:G:246:ILE:HD13	1.88	0.55
1:A:400:LEU:HD22	1:A:465:PHE:H	1.71	0.55
1:A:119:ARG:HH21	1:A:269:TRP:CD1	2.25	0.55
1:A:220:HIS:O	1:A:223:THR:OG1	2.26	0.54
3:G:193:GLU:HB2	3:G:280:LEU:HB3	1.89	0.54
3:G:98:TYR:CZ	3:G:257:LEU:HB2	2.43	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:185:GLU:OE1	1:A:258:ARG:NH2	2.40	0.54
1:A:361:PRO:HA	2:B:224:SER:HA	1.89	0.54
1:A:429:VAL:HG23	1:A:430:ASN:H	1.72	0.54
1:A:367:PRO:HG3	3:G:243:PHE:CE2	2.42	0.54
1:A:144:PRO:HA	1:A:147:VAL:HG22	1.90	0.53
1:A:119:ARG:HD3	1:A:272:GLN:O	2.09	0.53
1:A:451:LEU:HA	1:A:464:ASP:O	2.08	0.53
3:G:180:THR:OG1	3:G:181:GLU:N	2.40	0.53
1:A:365:ARG:HH11	2:B:223:ILE:C	2.17	0.53
3:G:52:THR:HB	3:G:114:GLU:HB2	1.90	0.53
1:A:365:ARG:NH1	2:B:225:CYS:H	2.07	0.53
3:G:297:HIS:HB3	3:G:298:ARG:HH21	1.74	0.53
1:A:184:PRO:HD3	1:A:200:TRP:CE2	2.45	0.52
1:A:118:SER:OG	1:A:209:LEU:HB3	2.10	0.52
1:A:236:PRO:HB2	1:A:238:TYR:HB2	1.90	0.52
3:G:150:HIS:HE2	5:G:403:AMP:P	2.33	0.52
3:G:300:VAL:HG12	3:G:311:ILE:HD12	1.92	0.51
1:A:406:SER:O	1:A:410:ASP:HB2	2.09	0.51
1:A:465:PHE:HB2	2:B:241:LEU:HB3	1.92	0.51
1:A:166:MET:HG2	2:B:235:HIS:CE1	2.46	0.51
1:A:345:ALA:HB1	2:B:221:THR:HB	1.91	0.51
1:A:442:VAL:HG21	3:G:79:GLN:HB3	1.91	0.51
1:A:445:THR:HG21	1:A:529:ARG:HD2	1.92	0.51
1:A:177:GLY:O	2:B:197:ARG:HA	2.11	0.51
1:A:342:PHE:HD2	1:A:343:TYR:HD1	1.59	0.51
1:A:539:GLU:HG2	2:B:270:LYS:HE2	1.93	0.51
1:A:187:ILE:HD11	1:A:229:CYS:SG	2.52	0.50
1:A:51:ARG:HE	1:A:91:ILE:HD11	1.76	0.50
1:A:361:PRO:HG2	1:A:366:VAL:HG12	1.92	0.50
1:A:222:PRO:O	1:A:225:PHE:N	2.45	0.50
1:A:162:LEU:HD23	1:A:176:CYS:SG	2.52	0.50
1:A:365:ARG:NH2	2:B:221:THR:OG1	2.45	0.50
3:G:278:CYS:SG	3:G:301:VAL:HG12	2.52	0.50
1:A:336:MET:HG3	1:A:340:LYS:HG2	1.92	0.49
1:A:123:GLN:HG2	1:A:274:LEU:HD12	1.94	0.49
4:A:601:STU:C18	4:A:601:STU:HN4	2.24	0.49
1:A:365:ARG:NH1	2:B:222:ASN:O	2.45	0.49
1:A:400:LEU:HD22	1:A:465:PHE:N	2.27	0.49
1:A:113:LEU:HB2	1:A:209:LEU:O	2.12	0.49
3:G:225:SER:OG	5:G:403:AMP:O2P	2.24	0.49
3:G:64:VAL:HG21	3:G:250:ALA:HA	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:154:VAL:HG22	3:G:164:TYR:O	2.12	0.48
1:A:442:VAL:HG21	3:G:79:GLN:CB	2.44	0.48
1:A:141:ASP:HB2	1:A:162:LEU:HD22	1.95	0.48
1:A:84:VAL:HG12	1:A:93:MET:HG2	1.96	0.48
1:A:535:ILE:HG21	3:G:74:TRP:CE2	2.48	0.48
2:B:235:HIS:HA	2:B:238:LEU:CD1	2.44	0.48
3:G:27:VAL:HG23	3:G:28:TYR:H	1.79	0.48
1:A:417:ARG:HD2	1:A:417:ARG:N	2.28	0.48
2:B:233:PRO:HG2	2:B:238:LEU:HD21	1.96	0.48
1:A:362:HIS:CE1	1:A:364:GLU:HB2	2.49	0.47
1:A:448:LYS:H	1:A:468:ILE:CD1	2.27	0.47
1:A:275:PRO:HB2	1:A:277:TYR:CD1	2.50	0.47
1:A:139:HIS:CE1	1:A:141:ASP:O	2.68	0.47
3:G:102:LEU:HD23	3:G:102:LEU:H	1.79	0.47
1:A:342:PHE:CD2	1:A:343:TYR:HD1	2.33	0.47
1:A:80:LYS:H	1:A:96:GLU:HG2	1.81	0.46
1:A:32:VAL:HG12	1:A:47:LYS:HA	1.97	0.46
1:A:359:THR:O	1:A:361:PRO:HD3	2.15	0.46
1:A:399:HIS:C	1:A:400:LEU:HG	2.40	0.46
3:G:285:GLU:O	3:G:289:ASN:ND2	2.41	0.46
1:A:143:LYS:HB3	1:A:181:TYR:HE2	1.81	0.46
1:A:400:LEU:CD2	1:A:465:PHE:H	2.29	0.46
1:A:437:ARG:HB2	1:A:448:LYS:HG2	1.98	0.46
3:G:297:HIS:O	3:G:314:LEU:HG	2.16	0.46
1:A:216:PHE:CD2	1:A:228:ILE:HA	2.51	0.46
3:G:206:VAL:O	3:G:229:VAL:HA	2.16	0.46
1:A:336:MET:HG3	1:A:340:LYS:CG	2.46	0.46
1:A:462:LEU:HD22	1:A:463:LEU:N	2.31	0.46
1:A:98:VAL:HG11	1:A:156:LYS:HE3	1.98	0.46
1:A:166:MET:HG2	2:B:235:HIS:HE1	1.79	0.46
2:B:252:VAL:HG12	2:B:269:TYR:HA	1.98	0.46
2:B:268:LEU:HD13	3:G:49:VAL:HG23	1.97	0.46
3:G:91:ILE:CG1	3:G:245:VAL:HG21	2.45	0.46
3:G:186:GLU:HG2	3:G:187:PHE:H	1.81	0.45
1:A:443:THR:HG22	1:A:444:SER:H	1.81	0.45
1:A:123:GLN:C	1:A:278:LEU:HD21	2.41	0.45
1:A:207:TYR:OH	1:A:234:TYR:O	2.28	0.45
1:A:449:MET:HE1	1:A:465:PHE:CD2	2.52	0.45
1:A:118:SER:OG	1:A:209:LEU:O	2.33	0.45
2:B:244:LEU:HD12	2:B:245:SER:N	2.31	0.45
3:G:173:LYS:O	3:G:177:LEU:HB2	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:139:HIS:NE2	1:A:159:ASP:O	2.48	0.45
1:A:340:LYS:CD	1:A:341:ASP:H	2.28	0.45
1:A:113:LEU:HD12	1:A:117:GLU:HG2	1.99	0.45
1:A:335:ILE:HD13	1:A:335:ILE:HA	1.86	0.45
1:A:126:LEU:HD23	1:A:270:PHE:CZ	2.52	0.45
3:G:200:TYR:HE2	3:G:301:VAL:HG21	1.82	0.45
1:A:141:ASP:OD2	1:A:143:LYS:HE2	2.17	0.44
1:A:30:GLY:HA3	1:A:48:ILE:O	2.17	0.44
1:A:169:GLY:HA2	2:B:236:VAL:O	2.18	0.44
1:A:261:ILE:H	1:A:261:ILE:HG13	1.62	0.44
1:A:466:ARG:HH11	2:B:240:HIS:CE1	2.35	0.44
2:B:261:LYS:HB2	3:G:38:TYR:OH	2.16	0.44
3:G:55:GLN:HA	3:G:112:LYS:HA	1.99	0.44
1:A:23:THR:HA	1:A:33:LYS:HG2	1.99	0.44
3:G:60:PHE:CE1	3:G:85:LEU:HD21	2.51	0.44
1:A:78:ILE:HD13	1:A:157:ILE:HB	2.00	0.44
1:A:448:LYS:O	1:A:468:ILE:HG12	2.17	0.44
2:B:200:GLU:OE2	2:B:200:GLU:N	2.46	0.44
3:G:47:LEU:HD11	3:G:72:PRO:HG2	2.00	0.44
3:G:90:PHE:O	3:G:94:LEU:HB2	2.18	0.44
1:A:275:PRO:HB2	1:A:277:TYR:HD1	1.82	0.44
1:A:427:LYS:H	1:A:427:LYS:HG3	1.56	0.44
1:A:178:SER:HA	2:B:196:PHE:O	2.18	0.44
3:G:191:SER:HA	3:G:283:THR:HA	2.00	0.44
1:A:210:LEU:HD13	1:A:239:LEU:HD12	2.00	0.43
2:B:254:SER:HB2	2:B:267:LEU:HA	2.00	0.43
1:A:363:PRO:O	1:A:366:VAL:HG22	2.19	0.43
3:G:315:SER:HB2	5:G:403:AMP:H5'1	1.99	0.43
2:B:222:ASN:OD1	2:B:223:ILE:N	2.47	0.43
1:A:139:HIS:CE1	1:A:159:ASP:O	2.72	0.43
1:A:462:LEU:HD22	1:A:463:LEU:H	1.84	0.43
1:A:21:GLY:O	1:A:33:LYS:HD2	2.19	0.43
1:A:147:VAL:HG12	1:A:157:ILE:HA	2.00	0.43
2:B:222:ASN:HB3	2:B:225:CYS:SG	2.59	0.43
3:G:242:LYS:O	3:G:245:VAL:HG22	2.19	0.43
1:A:103:LEU:HA	1:A:149:LEU:HD11	1.99	0.43
1:A:129:VAL:HG11	1:A:142:LEU:HD21	2.00	0.43
3:G:41:ILE:HD11	3:G:154:VAL:HG21	1.99	0.43
1:A:365:ARG:NH1	2:B:225:CYS:N	2.67	0.43
1:A:468:ILE:HB	1:A:469:ASP:H	1.54	0.43
3:G:200:TYR:CD2	3:G:310:GLY:HA3	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:40:LEU:HD21	3:G:137:LEU:HD13	2.00	0.43
1:A:437:ARG:HD2	1:A:448:LYS:HG2	2.01	0.43
3:G:27:VAL:HG23	3:G:28:TYR:N	2.33	0.42
3:G:56:VAL:HG22	3:G:111:HIS:O	2.19	0.42
3:G:205:MET:SD	3:G:308:VAL:HG21	2.58	0.42
1:A:114:ASP:O	1:A:118:SER:HB2	2.19	0.42
1:A:336:MET:C	1:A:340:LYS:HB2	2.45	0.42
1:A:364:GLU:HA	3:G:243:PHE:CZ	2.54	0.42
1:A:442:VAL:HG23	1:A:443:THR:OG1	2.18	0.42
3:G:243:PHE:HB3	5:G:402:AMP:H5'1	2.01	0.42
1:A:177:GLY:C	2:B:197:ARG:HA	2.45	0.42
1:A:340:LYS:O	1:A:342:PHE:N	2.52	0.42
2:B:248:ASP:OD1	2:B:249:SER:N	2.53	0.42
1:A:240:ASN:ND2	1:A:242:SER:HB3	2.35	0.42
1:A:428:VAL:HG22	1:A:433:TYR:CD1	2.55	0.42
1:A:17:HIS:O	1:A:38:GLU:HG3	2.20	0.41
1:A:399:HIS:O	1:A:400:LEU:HG	2.19	0.41
1:A:420:LYS:HE3	1:A:420:LYS:HB2	1.88	0.41
2:B:246:ILE:CG2	2:B:251:MET:HA	2.50	0.41
1:A:46:VAL:HA	1:A:93:MET:O	2.19	0.41
1:A:15:ILE:HD13	1:A:92:PHE:CG	2.55	0.41
3:G:207:ARG:HB2	3:G:210:THR:HG23	2.03	0.41
3:G:105:ILE:HG22	3:G:255:ASN:OD1	2.20	0.41
1:A:143:LYS:HD2	1:A:145:GLU:HB2	2.03	0.41
1:A:429:VAL:HG23	1:A:430:ASN:N	2.36	0.41
3:G:211:PRO:O	3:G:214:VAL:HG22	2.20	0.41
1:A:250:MET:HE2	1:A:264:ILE:HG12	2.02	0.41
1:A:120:ARG:O	1:A:124:GLN:HG3	2.20	0.41
1:A:240:ASN:HB2	1:A:241:PRO:HD2	2.01	0.41
1:A:362:HIS:CG	1:A:363:PRO:HD2	2.56	0.41
1:A:423:ASP:OD1	1:A:423:ASP:N	2.48	0.41
2:B:213:LEU:HD23	2:B:213:LEU:HA	1.83	0.41
3:G:73:LEU:HD21	3:G:85:LEU:HD12	2.03	0.41
3:G:291:LEU:HD13	3:G:299:LEU:HG	2.02	0.41
1:A:199:ILE:HG21	1:A:259:ALA:O	2.21	0.41
3:G:238:ASP:OD1	3:G:268:ARG:NH2	2.50	0.41
1:A:443:THR:HB	1:A:445:THR:HG23	2.03	0.40
3:G:214:VAL:O	3:G:218:ILE:HG13	2.22	0.40
1:A:211:CYS:C	1:A:213:THR:H	2.29	0.40
1:A:364:GLU:HA	3:G:243:PHE:CE2	2.56	0.40
1:A:403:ARG:O	1:A:460:THR:HG23	2.20	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:237:MET:HA	2:B:240:HIS:HD2	1.87	0.40
3:G:297:HIS:HB3	3:G:298:ARG:NH2	2.37	0.40
3:G:177:LEU:HD12	3:G:177:LEU:HA	1.89	0.40
1:A:463:LEU:HD23	1:A:463:LEU:HA	1.92	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	401/540 (74%)	373 (93%)	27 (7%)	1 (0%)	43	77
2	B	77/197 (39%)	71 (92%)	6 (8%)	0	100	100
3	G	299/304 (98%)	284 (95%)	15 (5%)	0	100	100
All	All	777/1041 (75%)	728 (94%)	48 (6%)	1 (0%)	48	83

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	428	VAL

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	369/485 (76%)	331 (90%)	38 (10%)	7	23
2	B	75/182 (41%)	68 (91%)	7 (9%)	8	26
3	G	276/278 (99%)	255 (92%)	21 (8%)	12	33
All	All	720/945 (76%)	654 (91%)	66 (9%)	8	27

All (66) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	ILE
1	A	59	VAL
1	A	70	LEU
1	A	80	LYS
1	A	87	THR
1	A	90	ASP
1	A	108	CYS
1	A	116	LYS
1	A	118	SER
1	A	142	LEU
1	A	166	MET
1	A	185	GLU
1	A	220	HIS
1	A	223	THR
1	A	235	THR
1	A	237	GLN
1	A	238	TYR
1	A	261	ILE
1	A	266	GLU
1	A	335	ILE
1	A	336	MET
1	A	340	LYS
1	A	347	SER
1	A	358	LEU
1	A	359	THR
1	A	395	LYS
1	A	400	LEU
1	A	403	ARG
1	A	407	ARG
1	A	423	ASP
1	A	425	GLU
1	A	427	LYS
1	A	443	THR
1	A	462	LEU

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Mol	Chain	Res	Type
1	A	463	LEU
1	A	468	ILE
1	A	469	ASP
1	A	547	ILE
2	B	196	PHE
2	B	200	GLU
2	B	214	GLN
2	B	220	ASP
2	B	246	ILE
2	B	266	THR
2	B	272	ILE
3	G	36	ARG
3	G	47	LEU
3	G	87	ILE
3	G	105	ILE
3	G	106	TYR
3	G	111	HIS
3	G	125	PHE
3	G	126	LYS
3	G	166	LEU
3	G	172	LEU
3	G	175	LEU
3	G	177	LEU
3	G	180	THR
3	G	186	GLU
3	G	194	GLU
3	G	207	ARG
3	G	270	HIS
3	G	276	LEU
3	G	298	ARG
3	G	304	GLU
3	G	323	LEU

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	83	GLN
3	G	79	GLN
3	G	196	GLN
3	G	319	GLN

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 ⓘ

4 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
4	STU	A	601	-	39,42,42	0.80	2 (5%)	50,68,68	2.49	12 (24%)
5	AMP	G	403	-	25,25,25	1.64	3 (12%)	37,38,38	1.80	10 (27%)
5	AMP	G	401	-	25,25,25	1.62	3 (12%)	37,38,38	1.84	10 (27%)
5	AMP	G	402	-	25,25,25	1.63	3 (12%)	37,38,38	1.82	10 (27%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	STU	A	601	-	-	2/4/42/42	-
5	AMP	G	403	-	-	0/10/26/26	0/3/3/3
5	AMP	G	401	-	-	0/10/26/26	0/3/3/3
5	AMP	G	402	-	-	0/10/26/26	0/3/3/3

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	G	402	AMP	C6-N6	5.94	1.49	1.34
5	G	403	AMP	C6-N6	5.91	1.49	1.34
5	G	401	AMP	C6-N6	5.82	1.49	1.34
5	G	401	AMP	P-O1P	3.15	1.60	1.50
5	G	403	AMP	P-O1P	3.13	1.60	1.50
5	G	402	AMP	P-O1P	3.09	1.60	1.50
5	G	402	AMP	P-O3P	-3.08	1.43	1.54
5	G	401	AMP	P-O3P	-3.06	1.43	1.54
5	G	403	AMP	P-O3P	-3.06	1.43	1.54
4	A	601	STU	O4-C25	-2.73	1.39	1.43
4	A	601	STU	C19-N3	-2.26	1.35	1.39

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	601	STU	C9-C10-C7	-9.21	106.50	109.88
4	A	601	STU	C28-N4-C23	-6.08	107.12	114.39
4	A	601	STU	C9-N1-C8	-5.86	108.52	113.86
5	G	402	AMP	C5-C4-N3	-5.72	118.85	126.72
5	G	403	AMP	C5-C4-N3	-5.68	118.90	126.72
5	G	401	AMP	C5-C4-N3	-5.62	118.98	126.72
4	A	601	STU	C9-C10-C11	4.89	132.44	128.73
4	A	601	STU	C27-O6-C22	-4.73	106.46	114.46
4	A	601	STU	C10-C9-N1	4.67	106.29	101.75
4	A	601	STU	C7-C8-N1	4.25	110.43	106.38
5	G	401	AMP	N3-C4-N9	4.14	134.21	127.17
5	G	403	AMP	N3-C4-N9	4.12	134.17	127.17
5	G	402	AMP	N3-C4-N9	4.09	134.13	127.17
5	G	401	AMP	N3-C2-N1	-3.53	123.25	128.58
5	G	403	AMP	N3-C2-N1	-3.51	123.27	128.58
5	G	402	AMP	N3-C2-N1	-3.50	123.29	128.58
5	G	401	AMP	N9-C8-N7	-3.09	109.55	113.94
5	G	402	AMP	C2-N3-C4	3.06	119.31	111.83
5	G	403	AMP	C2-N3-C4	3.03	119.24	111.83
4	A	601	STU	C16-C17-C12	-3.02	118.06	121.59
5	G	402	AMP	N9-C8-N7	-2.99	109.70	113.94
5	G	401	AMP	C2-N3-C4	2.98	119.10	111.83
4	A	601	STU	C13-C12-C17	2.88	122.57	119.24
5	G	401	AMP	C5-N7-C8	2.87	107.97	103.45
5	G	403	AMP	N9-C8-N7	-2.85	109.89	113.94
5	G	402	AMP	C5-N7-C8	2.81	107.87	103.45
5	G	402	AMP	C4-C5-N7	-2.77	107.42	110.58

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	G	403	AMP	C5-N7-C8	2.74	107.75	103.45
5	G	403	AMP	C4-C5-N7	-2.64	107.57	110.58
5	G	401	AMP	C4-C5-N7	-2.62	107.58	110.58
5	G	401	AMP	O2P-P-O1P	2.45	120.37	110.83
5	G	403	AMP	O2P-P-O1P	2.41	120.21	110.83
5	G	401	AMP	C6-C5-C4	2.34	120.37	117.18
5	G	402	AMP	O2P-P-O1P	2.32	119.89	110.83
5	G	401	AMP	C4-N9-C8	2.24	108.09	105.74
4	A	601	STU	O5-C8-C7	-2.24	125.22	128.47
5	G	402	AMP	C6-C5-C4	2.22	120.21	117.18
4	A	601	STU	C6-C7-C10	-2.18	119.24	120.76
5	G	403	AMP	C6-C5-C4	2.14	120.10	117.18
5	G	402	AMP	C4-N9-C8	2.14	107.98	105.74
5	G	403	AMP	C4-N9-C8	2.03	107.87	105.74
4	A	601	STU	C19-N3-C20	2.00	110.55	107.90

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	601	STU	C24-C23-N4-C28
4	A	601	STU	C22-C23-N4-C28

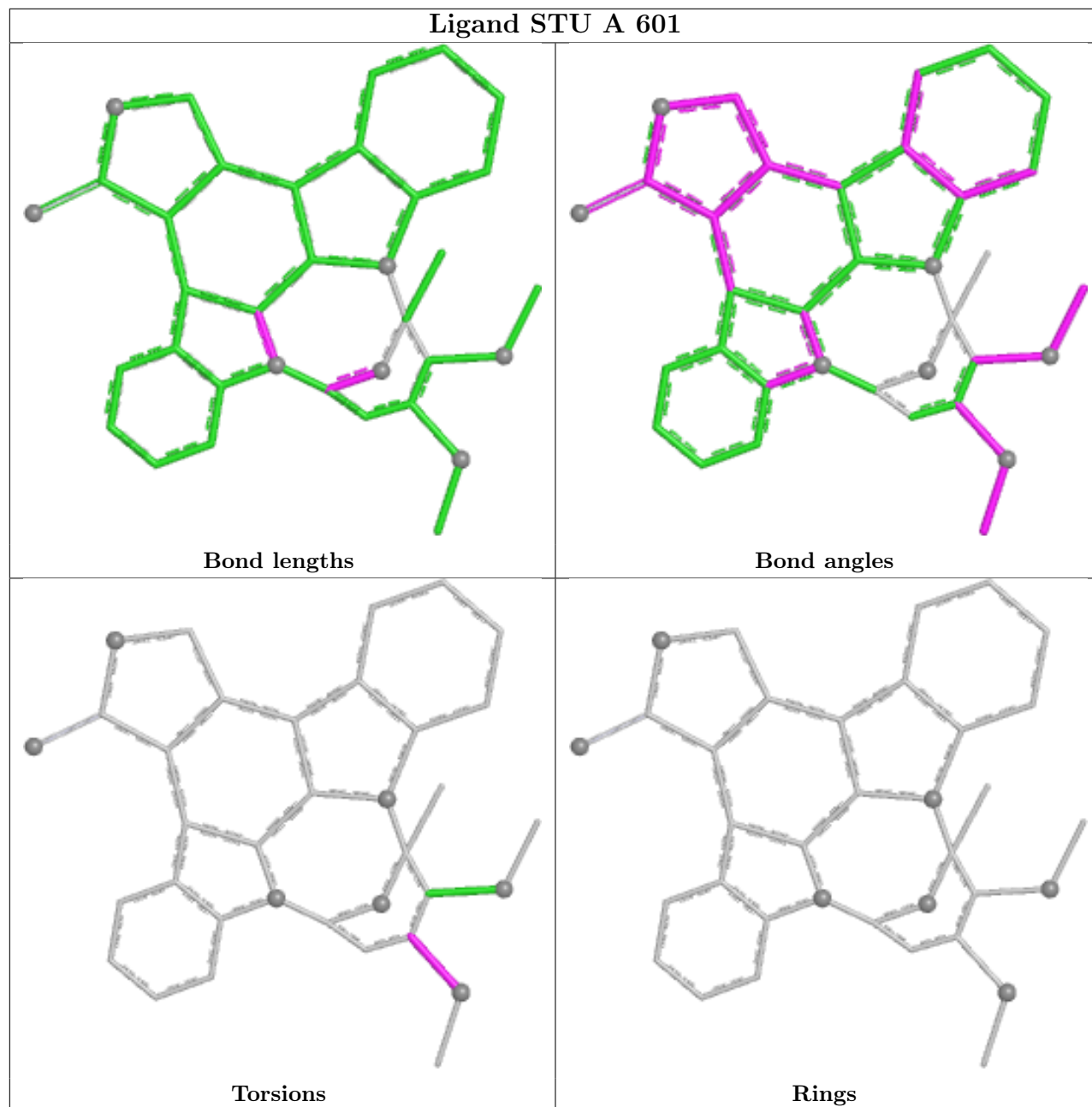
There are no ring outliers.

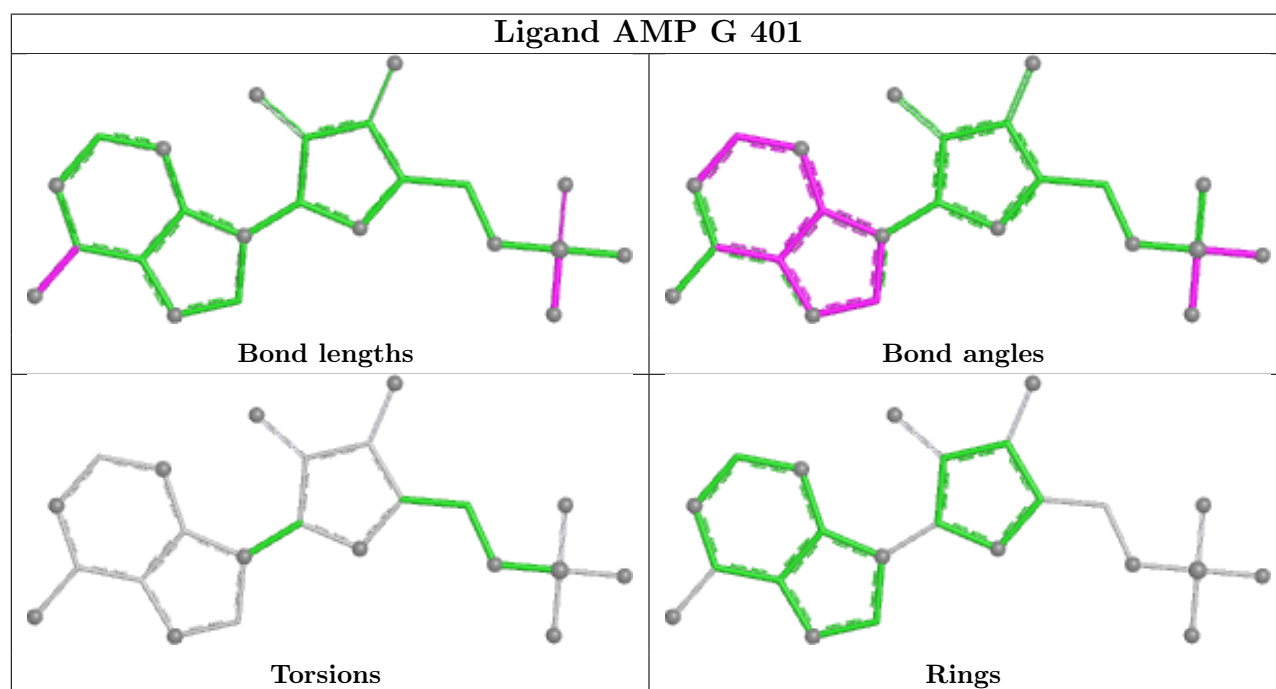
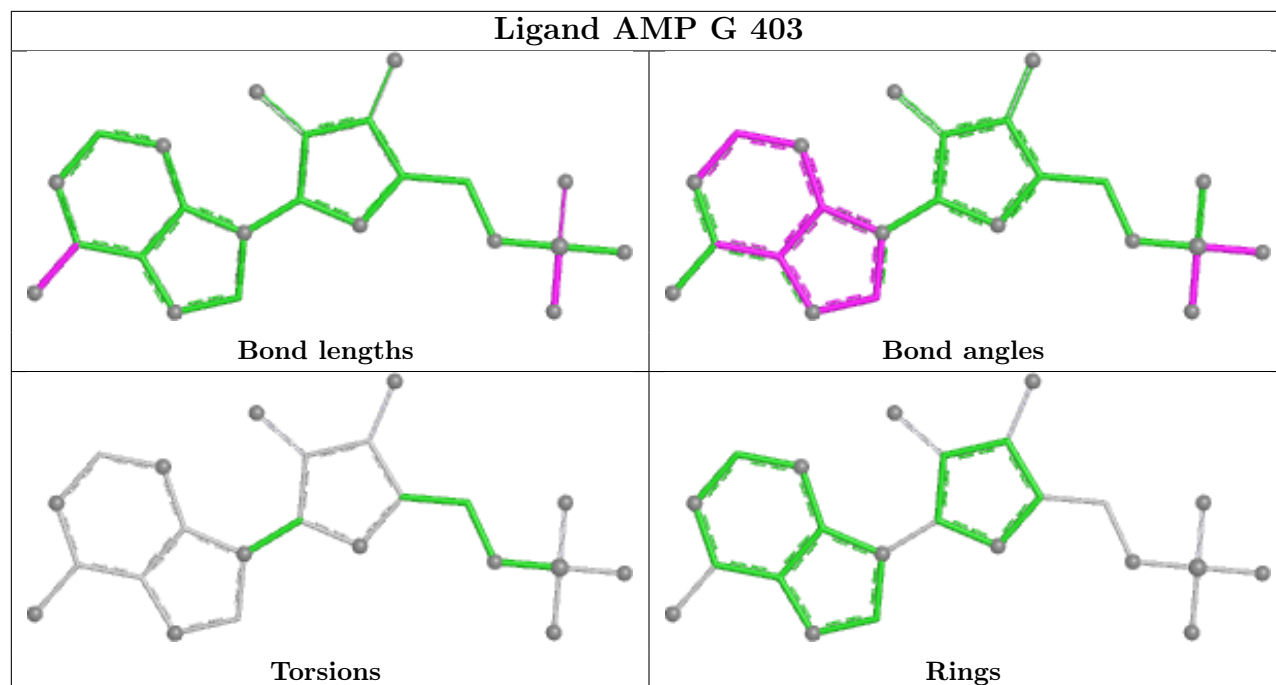
3 monomers are involved in 7 short contacts:

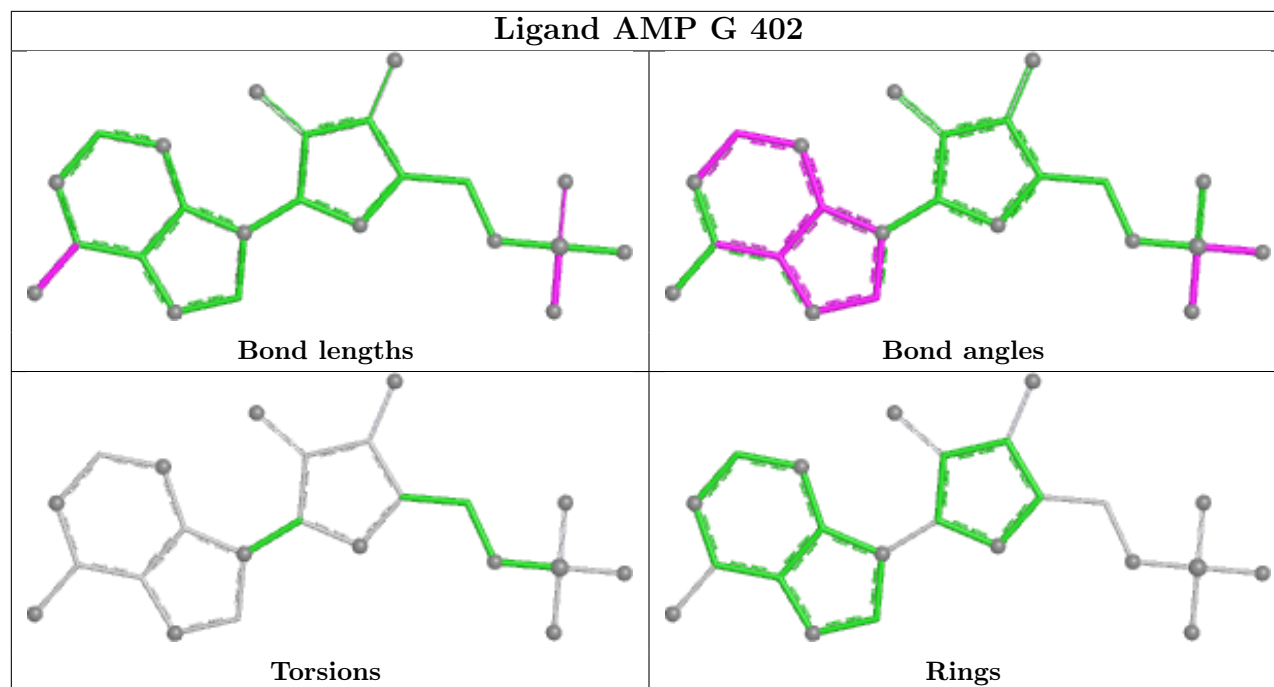
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	601	STU	2	0
5	G	403	AMP	3	0
5	G	402	AMP	2	0

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	411/540 (76%)	-0.29	0 100 100	187, 216, 263, 289	0
2	B	79/197 (40%)	-0.25	0 100 100	193, 207, 251, 259	0
3	G	301/304 (99%)	-0.32	1 (0%) 90 80	122, 211, 247, 293	0
All	All	791/1041 (75%)	-0.30	1 (0%) 92 87	122, 213, 256, 293	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	G	305	ASN	2.8

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(Å ²)	Q<0.9
5	AMP	G	403	23/23	0.90	0.09	196,202,208,212	0
4	STU	A	601	35/35	0.91	0.10	193,198,202,202	0

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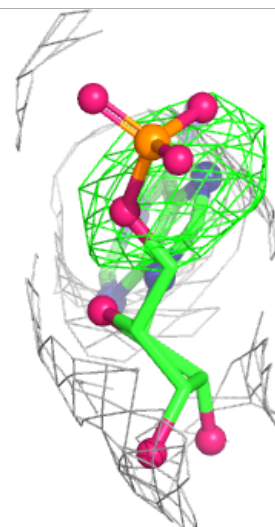
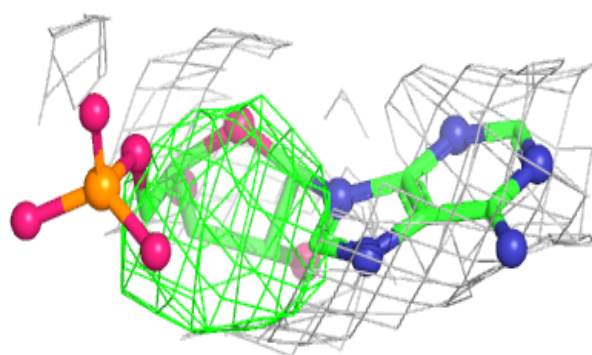
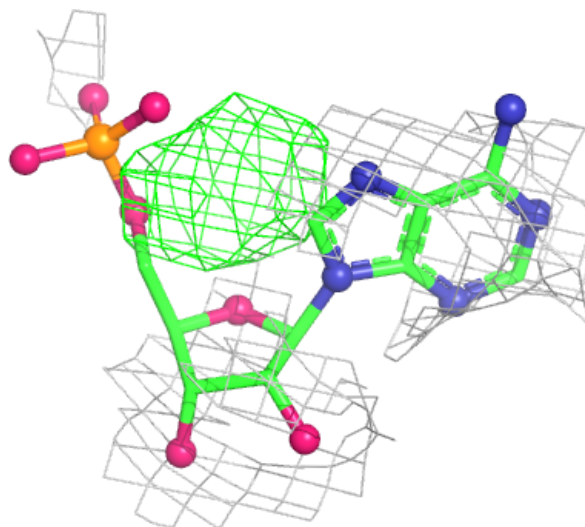
Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	AMP	G	401	23/23	0.93	0.07	201,211,216,218	0
5	AMP	G	402	23/23	0.95	0.11	186,188,194,195	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.

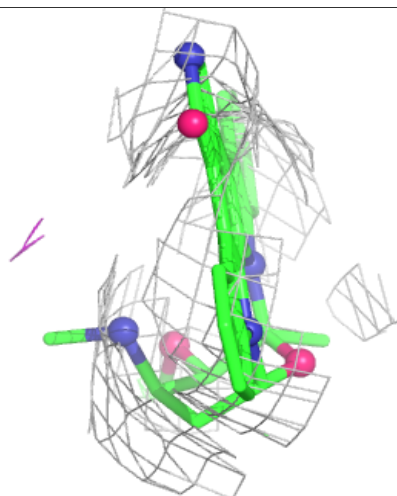
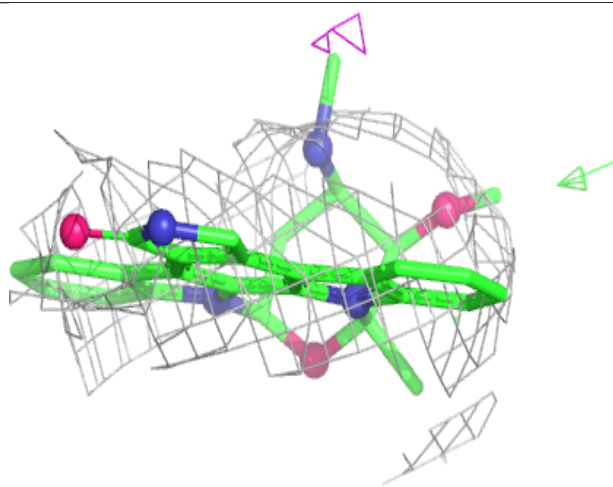
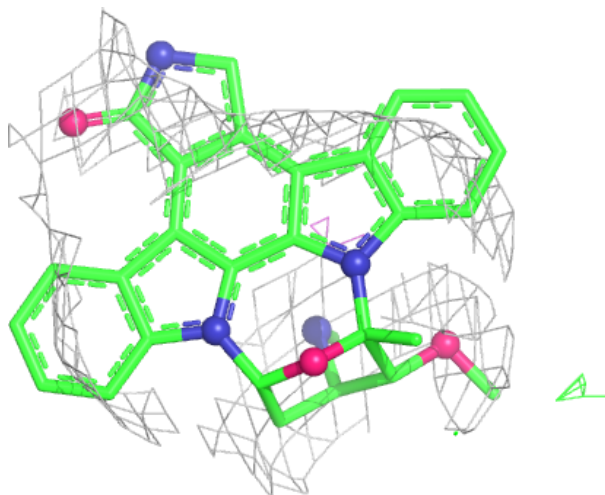
Electron density around AMP G 403:

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



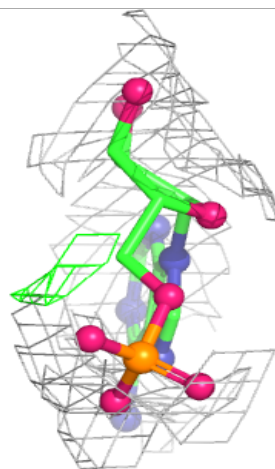
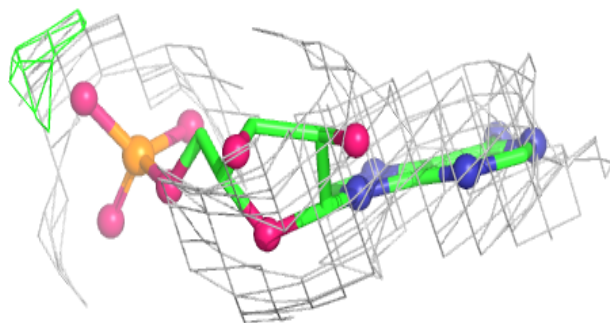
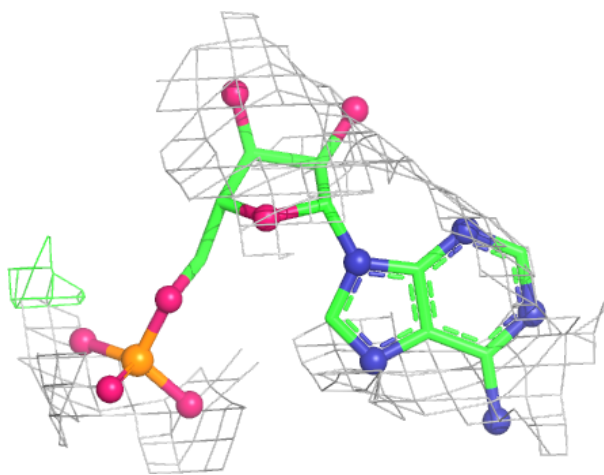
Electron density around STU A 601:

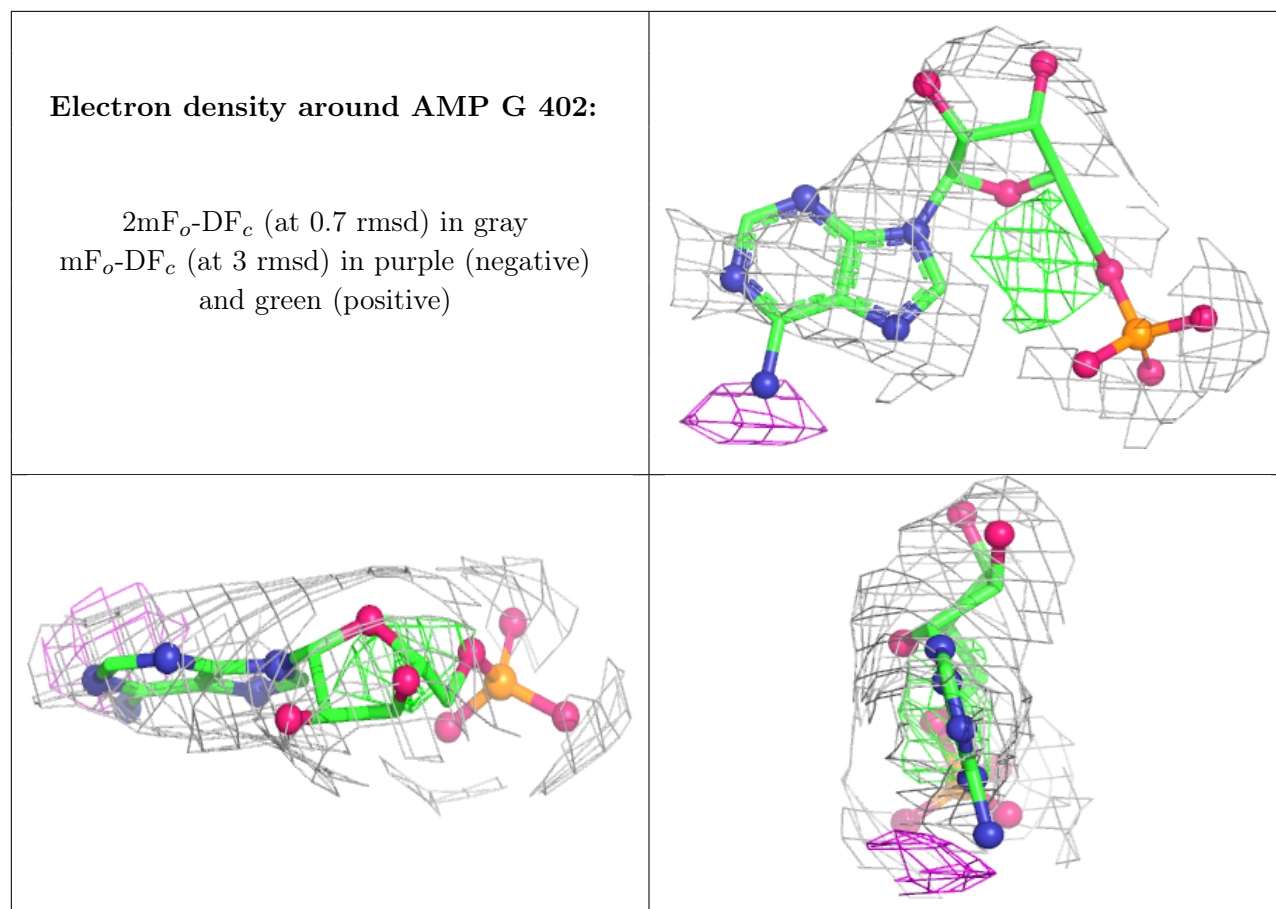
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around AMP G 401:

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





6.5 Other polymers [i](#)

There are no such residues in this entry.