



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 2, 2026 – 01:15 AM UTC

PDB ID : 7MYJ / pdb_00007myj
Title : Structure of full length human AMPK (a2b1g1) in complex with a small molecule activator MSG011
Authors : Ovens, A.J.; Gee, Y.S.; Ling, N.X.Y.; Waters, N.J.; Yu, D.; Scott, J.W.; Parker, M.W.; Hoffman, N.J.; Kemp, B.E.; Baell, J.B.; Oakhill, J.S.; Langendorf, C.G.
Deposited on : 2021-05-21
Resolution : 2.95 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](#) ①) were used in the production of this report:

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

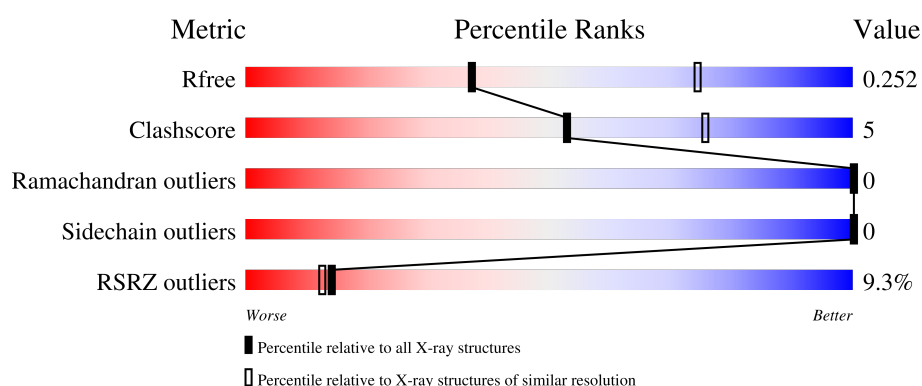
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.95 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	1130 (2.98-2.94)
Clashscore	190562	1157 (2.98-2.94)
Ramachandran outliers	187476	1101 (2.98-2.94)
Sidechain outliers	187428	1101 (2.98-2.94)
RSRZ outliers	180081	1130 (2.98-2.94)

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	565	
1	C	565	
2	B	270	
2	D	270	
3	E	336	

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Mol	Chain	Length	Quality of chain
3	F	336	8% 77% 13% 10%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 14740 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-2.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	435	Total	C	N	O	P	S	0	0	0
			3379	2172	571	612	1	23			
1	C	465	Total	C	N	O	P	S	1	0	0
			3601	2308	610	658	1	24			

There are 30 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-12	MET	-	initiating methionine	UNP P54646
A	-11	GLY	-	expression tag	UNP P54646
A	-10	SER	-	expression tag	UNP P54646
A	-9	SER	-	expression tag	UNP P54646
A	-8	HIS	-	expression tag	UNP P54646
A	-7	HIS	-	expression tag	UNP P54646
A	-6	HIS	-	expression tag	UNP P54646
A	-5	HIS	-	expression tag	UNP P54646
A	-4	HIS	-	expression tag	UNP P54646
A	-3	HIS	-	expression tag	UNP P54646
A	-2	SER	-	expression tag	UNP P54646
A	-1	GLN	-	expression tag	UNP P54646
A	0	ASP	-	expression tag	UNP P54646
A	1	PRO	-	expression tag	UNP P54646
A	271	GLY	ASP	conflict	UNP P54646
C	-12	MET	-	initiating methionine	UNP P54646
C	-11	GLY	-	expression tag	UNP P54646
C	-10	SER	-	expression tag	UNP P54646
C	-9	SER	-	expression tag	UNP P54646
C	-8	HIS	-	expression tag	UNP P54646
C	-7	HIS	-	expression tag	UNP P54646
C	-6	HIS	-	expression tag	UNP P54646
C	-5	HIS	-	expression tag	UNP P54646
C	-4	HIS	-	expression tag	UNP P54646
C	-3	HIS	-	expression tag	UNP P54646

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-2	SER	-	expression tag	UNP P54646
C	-1	GLN	-	expression tag	UNP P54646
C	0	ASP	-	expression tag	UNP P54646
C	1	PRO	-	expression tag	UNP P54646
C	271	GLY	ASP	conflict	UNP P54646

- Molecule 2 is a protein called 5'-AMP-activated protein kinase subunit beta-1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	184	Total	C	N	O	P	S	0	0	0
			1397	906	229	257	1	4			
2	D	187	Total	C	N	O	P	S	0	0	0
			1446	936	238	265	1	6			

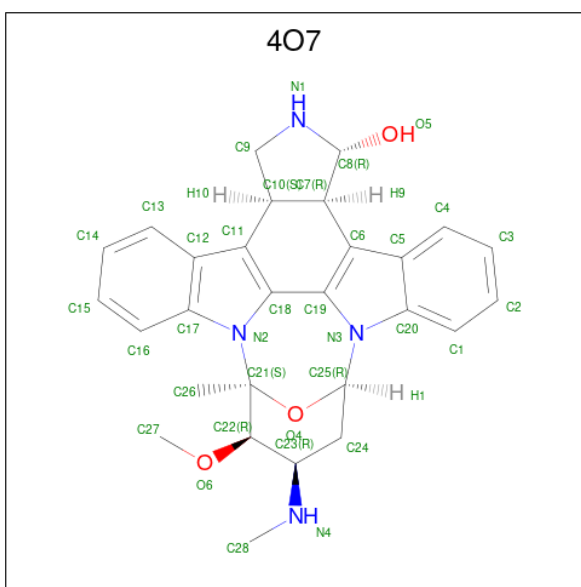
- Molecule 3 is a protein called 5'-AMP-activated protein kinase subunit gamma-1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	E	303	Total	C	N	O	S		0	0	0
			2342	1527	385	423	7				
3	F	303	Total	C	N	O	S		0	0	0
			2339	1524	386	422	7				

There are 12 discrepancies between the modelled and reference sequences:

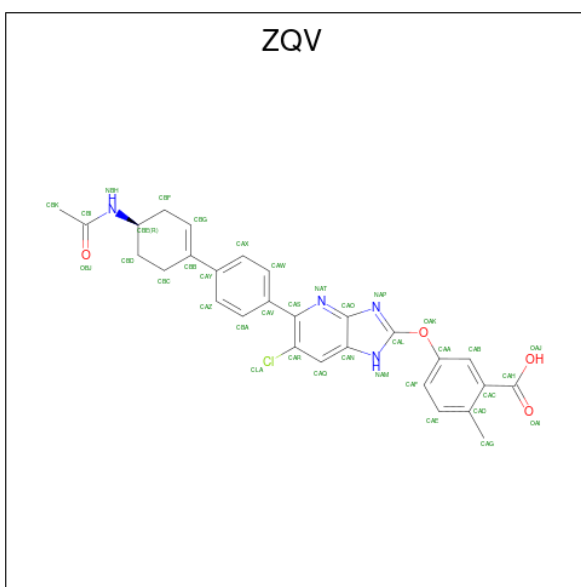
Chain	Residue	Modelled	Actual	Comment	Reference
E	-4	MET	-	initiating methionine	UNP P54619
E	-3	ALA	-	expression tag	UNP P54619
E	-2	ASP	-	expression tag	UNP P54619
E	-1	LEU	-	expression tag	UNP P54619
E	0	ASN	-	expression tag	UNP P54619
E	1	TRP	-	expression tag	UNP P54619
F	-4	MET	-	initiating methionine	UNP P54619
F	-3	ALA	-	expression tag	UNP P54619
F	-2	ASP	-	expression tag	UNP P54619
F	-1	LEU	-	expression tag	UNP P54619
F	0	ASN	-	expression tag	UNP P54619
F	1	TRP	-	expression tag	UNP P54619

- Molecule 4 is (5S,6R,7R,9R,13cR,14R,16aS)-6-methoxy-5-methyl-7-(methylamino)-6,7,8,9,14,15,16,16a-octahydro-5H,13cH-5,9-epoxy-4b,9a,1 5-triazadibenzo[b,h]cyclonona[1,2,3,4-jkl]cyclopenta[e]-as-indacen-14-ol (CCD ID: 4O7) (formula: C₂₈H₃₀N₄O₃).



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

- Molecule 5 is 5-(5-((4'R)-4'-acetamido-2',3',4',5'-tetrahydro[1,1'-biphenyl]-4-yl)-6-chloro-1H-imidazo[4,5-b]pyridin-2-yl)oxy)-2-methylbenzoic acid (CCD ID: ZQV) (formula: $C_{28}H_{25}ClN_4O_4$) (labeled as "Ligand of Interest" by depositor).



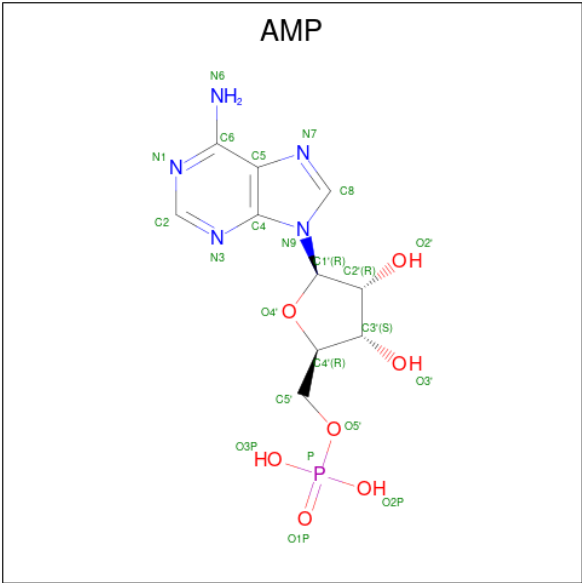
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	Cl	N	O	0
			37	28	1	4	4	

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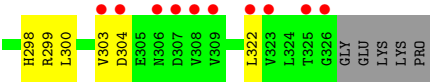
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	D	1	Total	C	Cl	N	O	0	0
			37	28	1	4	4		

- Molecule 6 is ADENOSINE MONOPHOSPHATE (CCD ID: AMP) (formula: C₁₀H₁₄N₅O₇P).



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

- Chain F:
-
- 8% 77% 13% 10%
- MET ALA ASP LEU ASN TRP GLU THR ILE VAL SER SER ASP SER SER PRO ALA VAL GLU ASN GLU HIS PRO GLN GLU THR PRO GLU S24 S25 Y29 M33 R37 C38 L41 V49 V50 F51 D52 T53 V57 F61 G68 P73 L74 F82 I94 T106



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	75.84Å 134.20Å 141.82Å 90.00° 92.55° 90.00°	Depositor
Resolution (Å)	48.72 – 2.95 48.72 – 2.95	Depositor EDS
% Data completeness (in resolution range)	99.0 (48.72-2.95) 98.9 (48.72-2.95)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.26 (at 2.96Å)	Xtriage
Refinement program	BUSTER 2.10.2	Depositor
R, R_{free}	0.213 , 0.243 (Not available) , 0.252	Depositor DCC
R_{free} test set	2992 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	70.8	Xtriage
Anisotropy	0.182	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 53.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.032 for h,-k,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	14740	wwPDB-VP
Average B, all atoms (Å ²)	75.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.41% 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: ZQV, TPO, SEP, 4O7, 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	0.78	0/3448	1.31	10/4678 (0.2%)
1	C	0.77	0/3674	1.28	4/4982 (0.1%)
2	B	0.79	0/1428	1.13	0/1959
2	D	0.79	0/1477	1.16	0/2018
3	E	0.79	0/2391	1.27	3/3260 (0.1%)
3	F	0.82	0/2388	1.26	3/3258 (0.1%)
All	All	0.79	0/14806	1.26	20/20155 (0.1%)

There are no bond length outliers.

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	221	VAL	N-CA-C	-6.86	104.50	110.74
1	A	159	GLY	N-CA-C	-6.64	102.81	111.37
1	A	218	HIS	CA-C-N	6.52	126.69	120.43
1	A	218	HIS	C-N-CA	6.52	126.69	120.43
1	C	275	TYR	N-CA-C	6.03	120.79	113.38
1	C	218	HIS	CA-C-N	5.98	126.17	120.43
1	C	218	HIS	C-N-CA	5.98	126.17	120.43
1	A	346	PRO	CB-CA-C	-5.82	103.82	110.92
1	A	224	LYS	N-CA-C	-5.62	105.24	111.36
1	A	160	LEU	N-CA-C	-5.44	104.72	111.90
3	E	106	ILE	N-CA-CB	5.39	117.80	111.66
1	C	224	LYS	N-CA-C	-5.35	105.53	111.36
1	A	175	GLY	CA-C-O	-5.33	118.48	122.37
3	E	139	PHE	N-CA-C	-5.33	105.47	111.28
3	F	106	ILE	CA-C-O	-5.24	118.07	122.63
1	A	346	PRO	N-CA-C	5.15	116.99	110.70
1	A	159	GLY	CA-C-O	-5.04	116.26	122.15
3	F	38	CYS	CA-C-N	5.03	127.27	120.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	38	CYS	C-N-CA	5.03	127.27	120.38
1	A	182	PRO	N-CA-C	5.02	121.22	113.75

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	3379	0	3216	42	0
1	C	3601	0	3435	23	0
2	B	1397	0	1305	23	0
2	D	1446	0	1360	17	0
3	E	2342	0	2341	25	0
3	F	2339	0	2316	27	0
4	A	35	0	0	0	0
4	C	35	0	0	0	0
5	A	37	0	0	0	0
5	D	37	0	0	0	0
6	E	46	0	24	0	0
6	F	46	0	24	0	0
All	All	14740	0	14021	143	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 5.

All (143) 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)
3:F:292:LEU:HD21	3:F:300:LEU:HG	1.18	1.16
3:E:41:LEU:HD13	3:E:138:LEU:HD21	1.26	1.09
2:B:125:TYR:HE1	2:B:153:ILE:HG12	1.15	1.06
3:E:41:LEU:CD1	3:E:138:LEU:HD21	1.86	1.06
3:F:292:LEU:CD2	3:F:300:LEU:HG	1.93	0.99
2:B:125:TYR:CE1	2:B:153:ILE:HG12	2.02	0.94

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:177:PRO:HB3	1:A:223:PHE:HE1	1.41	0.84
1:A:124:LEU:HD11	1:A:259:ILE:HG12	1.60	0.83
3:F:244:PHE:O	3:F:247:ILE:HG22	1.78	0.83
1:A:548:THR:HG21	2:B:270:ILE:HG23	1.62	0.81
3:E:41:LEU:HD13	3:E:138:LEU:CD2	2.08	0.80
3:E:244:PHE:O	3:E:247:ILE:HG22	1.82	0.80
3:E:216:ALA:O	3:E:220:PHE:CD2	2.40	0.75
3:F:292:LEU:HD21	3:F:300:LEU:CG	2.09	0.75
2:D:242:LEU:HD12	2:D:250:VAL:HG11	1.70	0.72
3:E:220:PHE:CZ	3:E:228:LEU:HD12	2.27	0.70
3:F:227:ALA:HB3	3:F:299:ARG:HD2	1.75	0.68
1:A:402:TRP:CH2	2:B:250:VAL:HG12	2.30	0.66
3:E:41:LEU:CD1	3:E:138:LEU:CD2	2.69	0.65
1:A:26:THR:HG21	1:A:160:LEU:HD13	1.79	0.65
1:A:124:LEU:HD23	1:A:268:PHE:CZ	2.31	0.64
1:A:137:HIS:ND1	1:A:140:LEU:HD23	2.12	0.64
1:A:548:THR:HG21	2:B:270:ILE:CG2	2.27	0.64
1:C:429:GLU:HB3	2:D:187:TYR:HB3	1.78	0.63
2:D:239:LEU:HD11	2:D:251:LEU:HD22	1.80	0.63
1:C:214:PHE:CG	1:C:226:ILE:HD13	2.36	0.60
1:A:185:ILE:HD12	1:A:223:PHE:CD1	2.35	0.60
1:C:72:ARG:NH2	1:C:290:ASP:OD1	2.34	0.60
1:C:543:CYS:SG	2:D:239:LEU:HD22	2.42	0.60
1:C:214:PHE:CB	1:C:226:ILE:HD13	2.33	0.59
1:A:190:TYR:CD1	1:A:195:VAL:HG21	2.37	0.59
1:A:407:ARG:NH2	1:A:460:VAL:HG11	2.17	0.59
1:C:222:LEU:O	1:C:226:ILE:HG12	2.03	0.58
1:A:440:VAL:HG12	1:A:542:MET:CE	2.33	0.58
1:A:293:VAL:HG23	1:A:293:VAL:O	2.03	0.57
2:D:97:PHE:CE1	2:D:98:ASN:ND2	2.73	0.57
3:E:220:PHE:CZ	3:E:228:LEU:CD1	2.87	0.57
2:B:125:TYR:CE1	2:B:153:ILE:CG1	2.83	0.57
2:B:206:LEU:HD21	2:B:211:LEU:HD11	1.87	0.57
1:C:428:PHE:HE1	1:C:442:ARG:HD3	1.69	0.57
1:A:26:THR:CG2	1:A:160:LEU:HD13	2.34	0.56
2:D:242:LEU:HB2	2:D:250:VAL:CG1	2.35	0.56
3:F:292:LEU:CD2	3:F:300:LEU:CG	2.78	0.55
1:A:456:GLN:HG2	1:A:458:TYR:CZ	2.42	0.55
3:F:82:PHE:O	3:F:129:LEU:HD22	2.07	0.55
2:B:206:LEU:CD2	2:B:211:LEU:HD11	2.36	0.55
3:F:57:VAL:HG22	3:F:61:PHE:CE2	2.42	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:171:ARG:HH21	1:C:189:LEU:HG	1.72	0.54
1:A:434:ASN:HB3	1:A:437:HIS:HB3	1.89	0.54
1:A:442:ARG:HH22	1:A:444:ASN:HD22	1.55	0.54
3:F:41:LEU:HD11	3:F:175:PHE:CD2	2.42	0.54
1:A:440:VAL:HG12	1:A:542:MET:HE1	1.90	0.53
1:A:26:THR:CB	1:A:160:LEU:HD13	2.38	0.53
3:F:53:THR:HB	3:F:115:GLU:HB3	1.90	0.53
3:E:285:LEU:HD21	3:E:322:LEU:HD11	1.91	0.53
1:C:171:ARG:HB2	1:C:189:LEU:HD23	1.90	0.52
3:F:37:ARG:HD2	3:F:135:ASN:HA	1.92	0.52
2:B:92:TYR:HB2	2:B:128:PHE:HB3	1.90	0.52
1:C:366:HIS:HB3	1:C:369:ARG:HG2	1.91	0.52
2:D:135:HIS:HB2	2:D:151:ASN:ND2	2.25	0.52
2:B:135:HIS:CD2	2:B:141:ILE:HD11	2.45	0.52
1:A:26:THR:HB	1:A:160:LEU:CD1	2.40	0.51
3:F:298:HIS:HB3	3:F:299:ARG:NH2	2.25	0.51
1:C:419:VAL:HG12	1:C:423:MET:HE2	1.92	0.51
3:E:220:PHE:CE2	3:E:228:LEU:HD12	2.45	0.51
1:A:140:LEU:HD13	1:A:200:CYS:SG	2.50	0.51
2:D:207:PRO:HD2	2:D:210:LEU:HD12	1.94	0.50
1:A:26:THR:HB	1:A:160:LEU:HD13	1.93	0.50
1:A:177:PRO:HB3	1:A:223:PHE:CE1	2.31	0.50
1:C:214:PHE:HB2	1:C:226:ILE:CD1	2.42	0.50
2:B:250:VAL:HG22	2:B:267:TYR:CD2	2.46	0.50
1:A:429:GLU:HB3	2:B:187:TYR:HB3	1.93	0.49
3:F:94:ILE:HG21	3:F:109:LEU:HD13	1.94	0.49
1:A:453:MET:HE1	1:A:469:PHE:HD2	1.78	0.48
1:A:419:VAL:HG12	1:A:423:MET:HE2	1.94	0.48
1:A:402:TRP:HB2	2:B:213:VAL:HG11	1.95	0.48
2:B:107:ARG:HE	2:B:110:ASN:H	1.61	0.48
3:F:280:TYR:CD1	3:F:303:VAL:HG23	2.49	0.48
1:C:214:PHE:CB	1:C:226:ILE:CD1	2.92	0.47
3:E:41:LEU:HD11	3:E:138:LEU:HD21	1.85	0.47
3:F:145:LEU:HD23	3:F:150:ILE:HG13	1.96	0.47
1:A:138:ARG:HH12	1:A:172:TPO:HB	1.79	0.47
1:A:138:ARG:HB2	1:A:160:LEU:O	2.14	0.47
2:B:81:VAL:HG22	2:B:115:ILE:HG12	1.95	0.47
1:C:136:VAL:HG11	1:C:192:GLY:HA3	1.97	0.47
3:E:206:MET:HG2	3:E:229:PRO:HD2	1.97	0.47
1:A:252:ASP:HB3	1:A:255:LYS:HB2	1.97	0.47
3:F:29:TYR:O	3:F:33:MET:HG2	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:265:LEU:HD22	3:F:49:VAL:HG22	1.97	0.46
3:E:194:GLU:HB2	3:E:281:LEU:HB3	1.98	0.46
1:C:370:MET:HG3	3:F:68:GLY:HA2	1.98	0.46
3:E:220:PHE:CE2	3:E:228:LEU:CD1	2.99	0.46
2:D:215:LEU:HD11	2:D:265:LEU:HD11	1.97	0.46
3:E:279:CYS:SG	3:E:302:VAL:HG22	2.56	0.45
1:A:96:VAL:HG11	1:A:154:LYS:HD2	1.99	0.45
3:F:50:VAL:HG22	3:F:73:PRO:HG2	1.99	0.45
2:D:200:PHE:O	2:D:200:PHE:CG	2.70	0.44
3:F:280:TYR:HD1	3:F:304:ASP:O	2.00	0.44
1:A:439:ARG:NH2	1:A:472:ILE:HD13	2.33	0.44
1:C:402:TRP:HB2	2:D:213:VAL:HG11	1.99	0.44
1:A:548:THR:CG2	2:B:270:ILE:CG2	2.94	0.44
2:D:267:TYR:HB2	3:F:51:PHE:HD1	1.83	0.44
1:C:173:SER:HB3	1:C:184:VAL:HG13	1.99	0.44
1:C:426:LEU:HD22	1:C:545:SER:HB3	2.00	0.44
2:D:200:PHE:O	2:D:200:PHE:CD1	2.70	0.43
3:E:106:ILE:O	3:E:110:GLU:HG2	2.18	0.43
3:E:201:TYR:HE1	3:E:302:VAL:HG11	1.82	0.43
3:E:303:VAL:HG12	3:E:309:VAL:HA	2.00	0.43
1:A:431:LYS:HB2	1:A:439:ARG:HB3	1.99	0.43
2:D:78:ARG:HE	2:D:119:PRO:HA	1.83	0.43
1:A:402:TRP:CH2	2:B:250:VAL:CG1	3.01	0.43
1:A:171:ARG:HB2	1:A:189:LEU:HD23	2.01	0.42
2:D:81:VAL:HG22	2:D:115:ILE:HG12	2.01	0.42
3:F:106:ILE:O	3:F:110:GLU:HG2	2.19	0.42
2:B:248:VAL:HA	2:B:269:PRO:HA	2.01	0.42
3:F:25:ASN:HB3	3:F:186:PRO:HA	2.00	0.42
1:A:202:VAL:HG13	1:A:213:PRO:HD2	2.00	0.42
1:A:338:SER:C	1:A:340:PHE:H	2.28	0.42
2:B:232:ASN:O	2:B:235:MET:HG3	2.19	0.42
2:B:207:PRO:HD2	2:B:210:LEU:HD12	2.01	0.42
1:C:214:PHE:HB2	1:C:226:ILE:HD11	2.01	0.42
1:C:374:ILE:HA	1:C:396:ALA:HB3	2.01	0.42
3:F:74:LEU:HD22	3:F:114:ILE:HG21	2.01	0.42
1:A:185:ILE:HD12	1:A:223:PHE:CE1	2.54	0.42
2:B:162:VAL:HG13	2:B:163:PHE:N	2.35	0.42
3:E:50:VAL:HG22	3:E:73:PRO:HG2	2.02	0.41
3:E:216:ALA:O	3:E:220:PHE:HD2	1.96	0.41
3:E:266:LEU:HD22	3:E:269:ARG:HE	1.85	0.41
3:F:277:LEU:CD1	3:F:295:ALA:HB1	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:447:THR:HG22	1:C:532:LEU:HB3	2.03	0.41
1:A:341:TYR:HD1	3:E:43:PRO:HA	1.84	0.41
1:A:47:LEU:HB3	1:A:52:ILE:HD11	2.03	0.41
3:E:41:LEU:HD11	3:E:138:LEU:CD2	2.47	0.41
3:E:280:TYR:HB2	3:E:283:GLU:HG3	2.02	0.41
1:C:252:ASP:HB3	1:C:255:LYS:HB2	2.03	0.41
2:D:76:GLN:OE1	2:D:76:GLN:N	2.54	0.41
2:B:142:VAL:HG23	2:B:152:ILE:HG21	2.03	0.41
2:B:250:VAL:HG21	2:B:267:TYR:CE2	2.56	0.40
3:F:138:LEU:HD11	3:F:172:ILE:HG12	2.04	0.40
3:E:74:LEU:HD22	3:E:114:ILE:HG21	2.02	0.40
1:A:216:ASP:OD2	1:A:221:THR:HB	2.21	0.40
1:C:397:VAL:HG22	1:C:398:LYS:N	2.36	0.40
3:F:285:LEU:HD21	3:F:322:LEU:HD21	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	422/565 (75%)	418 (99%)	4 (1%)	0	100	100
1	C	452/565 (80%)	447 (99%)	5 (1%)	0	100	100
2	B	177/270 (66%)	176 (99%)	1 (1%)	0	100	100
2	D	178/270 (66%)	175 (98%)	3 (2%)	0	100	100
3	E	301/336 (90%)	299 (99%)	2 (1%)	0	100	100
3	F	301/336 (90%)	295 (98%)	6 (2%)	0	100	100
All	All	1831/2342 (78%)	1810 (99%)	21 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	345/497 (69%)	345 (100%)	0	100	100
1	C	370/497 (74%)	370 (100%)	0	100	100
2	B	143/239 (60%)	143 (100%)	0	100	100
2	D	150/239 (63%)	150 (100%)	0	100	100
3	E	253/308 (82%)	253 (100%)	0	100	100
3	F	248/308 (80%)	248 (100%)	0	100	100
All	All	1509/2088 (72%)	1509 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	48	ASN
1	A	67	ASN
1	A	178	ASN
1	A	444	ASN
1	A	535	HIS
2	B	135	HIS
2	B	189	GLN
1	C	48	ASN
1	C	178	ASN
1	C	238	ASN
1	C	403	HIS
1	C	409	GLN
1	C	449	ASN
2	D	109	HIS
2	D	124	GLN
2	D	150	ASN
2	D	255	HIS
3	E	80	GLN
3	E	105	GLN

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Mol	Chain	Res	Type
3	E	135	ASN
3	E	248	ASN
3	E	282	HIS
3	E	290	ASN
3	F	80	GLN
3	F	112	HIS
3	F	257	ASN
3	F	320	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

4 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
2	SEP	D	108	2	8,9,10	0.66	0	7,12,14	0.66	0
1	TPO	A	172	1	8,10,11	0.75	0	10,14,16	1.00	1 (10%)
1	TPO	C	172	1	8,10,11	0.80	0	10,14,16	1.05	1 (10%)
2	SEP	B	108	2	8,9,10	0.89	0	7,12,14	2.07	2 (28%)

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 Torsions and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	SEP	D	108	2	-	3/6/8/10	-
1	TPO	A	172	1	-	0/9/11/13	-
1	TPO	C	172	1	-	4/9/11/13	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	SEP	B	108	2	-	4/6/8/10	-

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	108	SEP	OG-CB-CA	3.77	111.81	108.14
2	B	108	SEP	OG-P-O1P	3.60	116.17	106.44
1	C	172	TPO	O-C-CA	-2.77	117.64	124.77
1	A	172	TPO	O-C-CA	-2.47	118.43	124.77

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	108	SEP	CB-OG-P-O1P
2	B	108	SEP	CB-OG-P-O2P
2	B	108	SEP	CB-OG-P-O3P
1	C	172	TPO	N-CA-CB-CG2
1	C	172	TPO	N-CA-CB-OG1
1	C	172	TPO	C-CA-CB-CG2
2	D	108	SEP	CB-OG-P-O2P
2	D	108	SEP	CB-OG-P-O3P
2	D	108	SEP	CB-OG-P-O1P
2	B	108	SEP	CA-CB-OG-P
1	C	172	TPO	O-C-CA-CB

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	172	TPO	1	0

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

8 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
6	AMP	F	401	-	25,25,25	0.32	0	37,38,38	0.40	0
5	ZQV	D	701	-	39,41,41	2.49	10 (25%)	51,59,59	1.05	2 (3%)
6	AMP	E	401	-	25,25,25	0.34	0	37,38,38	0.41	0
4	4O7	A	601	-	38,42,42	3.26	10 (26%)	33,68,68	3.24	15 (45%)
6	AMP	E	402	-	25,25,25	0.35	0	37,38,38	0.43	0
6	AMP	F	402	-	25,25,25	0.34	0	37,38,38	0.40	0
5	ZQV	A	602	-	39,41,41	2.52	10 (25%)	51,59,59	1.26	7 (13%)
4	4O7	C	601	-	38,42,42	3.23	9 (23%)	33,68,68	3.24	15 (45%)

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
6	AMP	F	401	-	-	1/10/26/26	0/3/3/3
5	ZQV	D	701	-	-	8/18/30/30	0/5/5/5
6	AMP	E	401	-	-	4/10/26/26	0/3/3/3
4	4O7	A	601	-	-	2/4/58/58	-
6	AMP	E	402	-	-	1/10/26/26	0/3/3/3
6	AMP	F	402	-	-	4/10/26/26	0/3/3/3
5	ZQV	A	602	-	-	8/18/30/30	0/5/5/5
4	4O7	C	601	-	-	2/4/58/58	-

All (39) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	601	4O7	C18-C11	10.89	1.49	1.37
4	C	601	4O7	C18-C11	10.70	1.49	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	601	4O7	C19-C6	10.10	1.48	1.37
4	C	601	4O7	C19-C6	9.95	1.48	1.37
5	A	602	ZQV	CAV-CAS	-7.52	1.40	1.49
5	D	701	ZQV	CAV-CAS	-7.50	1.40	1.49
4	C	601	4O7	C8-N1	-6.84	1.35	1.45
4	A	601	4O7	C8-N1	-6.78	1.35	1.45
4	A	601	4O7	C12-C17	5.97	1.49	1.41
4	C	601	4O7	C12-C17	5.97	1.49	1.41
4	A	601	4O7	C5-C20	5.94	1.48	1.41
4	C	601	4O7	C5-C20	5.83	1.48	1.41
5	D	701	ZQV	CAG-CAD	-5.79	1.40	1.51
5	A	602	ZQV	CAG-CAD	-5.78	1.40	1.51
5	A	602	ZQV	CAY-CBB	-5.21	1.40	1.48
5	D	701	ZQV	CAN-CAO	-5.09	1.34	1.40
5	A	602	ZQV	CAN-CAO	-5.01	1.34	1.40
5	D	701	ZQV	CAY-CBB	-4.93	1.41	1.48
5	A	602	ZQV	CBF-CBG	-4.74	1.40	1.50
5	D	701	ZQV	CBF-CBG	-4.43	1.41	1.50
5	A	602	ZQV	CAC-CAH	-4.37	1.40	1.49
5	D	701	ZQV	CAC-CAH	-4.29	1.40	1.49
5	D	701	ZQV	CAQ-CAN	-3.41	1.34	1.39
5	A	602	ZQV	CAQ-CAN	-3.32	1.34	1.39
4	A	601	4O7	C10-C7	-3.24	1.48	1.54
4	C	601	4O7	C10-C7	-3.23	1.48	1.54
5	D	701	ZQV	CAO-NAP	-3.05	1.33	1.38
4	A	601	4O7	C25-N3	3.05	1.49	1.45
4	C	601	4O7	C25-N3	2.99	1.49	1.45
5	D	701	ZQV	CAN-NAM	-2.95	1.33	1.38
5	A	602	ZQV	CAN-NAM	-2.88	1.33	1.38
5	A	602	ZQV	CAO-NAP	-2.84	1.33	1.38
4	A	601	4O7	C21-N2	2.70	1.51	1.48
4	C	601	4O7	C21-N2	2.63	1.51	1.48
4	C	601	4O7	O5-C8	-2.62	1.28	1.40
5	D	701	ZQV	CBC-CBB	-2.58	1.42	1.50
4	A	601	4O7	O5-C8	-2.58	1.28	1.40
5	A	602	ZQV	CBC-CBB	-2.37	1.43	1.50
4	A	601	4O7	C17-N2	-2.12	1.38	1.41

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	601	4O7	C18-C19-C6	-10.07	114.74	122.87

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	601	4O7	C18-C19-C6	-9.62	115.10	122.87
4	A	601	4O7	C19-C18-C11	-6.85	117.34	122.87
4	A	601	4O7	C5-C6-C19	-6.58	102.26	107.63
4	C	601	4O7	C19-C18-C11	-6.27	117.81	122.87
4	C	601	4O7	C5-C6-C19	-6.10	102.66	107.63
4	C	601	4O7	C16-C17-C12	-4.81	116.59	121.95
4	A	601	4O7	C16-C17-C12	-4.47	116.96	121.95
4	C	601	4O7	C1-C20-C5	-4.43	117.00	121.95
4	C	601	4O7	C28-N4-C23	4.39	119.63	114.39
4	A	601	4O7	C28-N4-C23	4.35	119.58	114.39
4	A	601	4O7	C1-C20-C5	-4.27	117.19	121.95
4	A	601	4O7	C13-C12-C11	-4.09	127.55	134.06
4	C	601	4O7	C13-C12-C11	-4.01	127.68	134.06
4	A	601	4O7	C16-C17-N2	3.81	138.09	130.11
4	C	601	4O7	C16-C17-N2	3.75	137.96	130.11
4	A	601	4O7	C9-C10-C7	3.61	110.59	102.26
4	C	601	4O7	C9-C10-C7	3.46	110.25	102.26
5	A	602	ZQV	CBE-NBH-CBI	3.31	127.02	122.82
4	A	601	4O7	C4-C5-C6	-3.24	128.90	134.06
4	C	601	4O7	C12-C11-C18	-3.19	105.03	107.63
4	A	601	4O7	C12-C11-C18	-3.03	105.16	107.63
4	C	601	4O7	C4-C5-C6	-3.01	129.27	134.06
5	A	602	ZQV	CAQ-CAR-CAS	-3.00	118.84	121.88
5	D	701	ZQV	CAQ-CAR-CAS	-2.94	118.90	121.88
4	C	601	4O7	C12-C17-N2	-2.69	104.08	109.12
4	C	601	4O7	C4-C5-C20	2.63	122.28	119.24
5	A	602	ZQV	CAV-CAS-NAT	2.62	118.67	115.10
4	A	601	4O7	C12-C17-N2	-2.61	104.24	109.12
4	A	601	4O7	C4-C5-C20	2.57	122.20	119.24
5	D	701	ZQV	CAV-CAS-NAT	2.40	118.38	115.10
5	A	602	ZQV	CAY-CBB-CBG	-2.35	117.99	122.19
4	C	601	4O7	C1-C20-N3	2.33	135.33	130.55
4	C	601	4O7	C13-C12-C17	2.28	121.87	119.24
4	A	601	4O7	C1-C20-N3	2.21	135.07	130.55
5	A	602	ZQV	OBJ-CBI-NBH	2.20	125.87	121.98
5	A	602	ZQV	CBC-CBB-CBG	2.08	122.36	120.63
4	A	601	4O7	C13-C12-C17	2.03	121.58	119.24
5	A	602	ZQV	CBD-CBE-NBH	2.02	114.61	110.57

There are no chirality outliers.

All (30) torsion outliers are listed below:

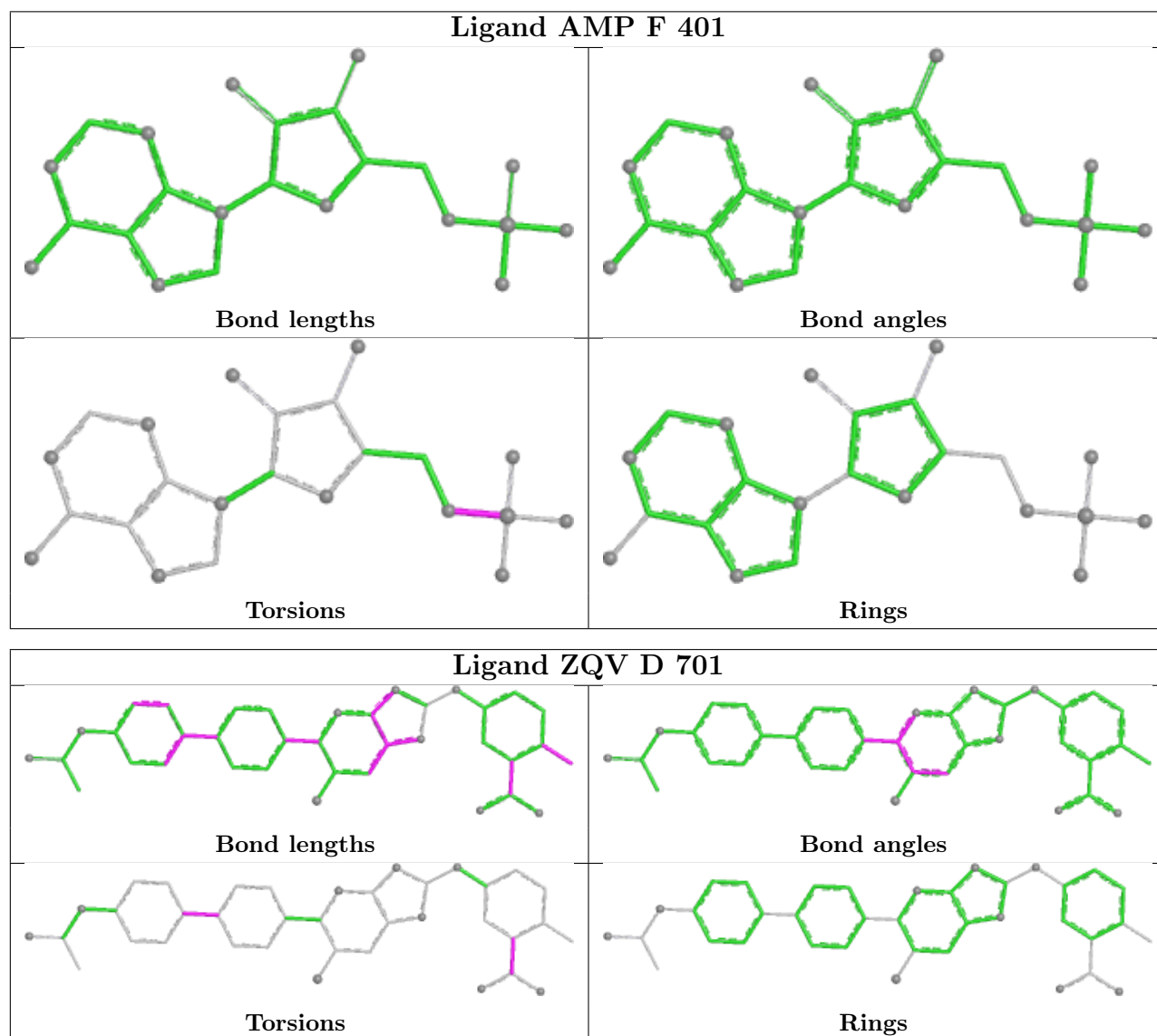
Mol	Chain	Res	Type	Atoms
6	E	401	AMP	C5'-O5'-P-O2P
6	E	401	AMP	C5'-O5'-P-O3P
6	F	402	AMP	C5'-O5'-P-O2P
6	F	402	AMP	C5'-O5'-P-O3P
6	F	402	AMP	O4'-C4'-C5'-O5'
6	F	402	AMP	C3'-C4'-C5'-O5'
5	A	602	ZQV	CAX-CAY-CBB-CBG
5	A	602	ZQV	CAZ-CAY-CBB-CBG
5	A	602	ZQV	CAX-CAY-CBB-CBC
5	A	602	ZQV	CAZ-CAY-CBB-CBC
4	A	601	4O7	C24-C23-N4-C28
4	C	601	4O7	C24-C23-N4-C28
5	D	701	ZQV	CAZ-CAY-CBB-CBG
6	E	401	AMP	O4'-C4'-C5'-O5'
5	D	701	ZQV	CAX-CAY-CBB-CBG
5	A	602	ZQV	CAB-CAC-CAH-OAI
5	D	701	ZQV	CAX-CAY-CBB-CBC
5	D	701	ZQV	CAB-CAC-CAH-OAJ
5	A	602	ZQV	CAB-CAC-CAH-OAJ
5	D	701	ZQV	CAB-CAC-CAH-OAI
4	A	601	4O7	C22-C23-N4-C28
4	C	601	4O7	C22-C23-N4-C28
5	D	701	ZQV	CAZ-CAY-CBB-CBC
5	D	701	ZQV	CAD-CAC-CAH-OAJ
6	E	402	AMP	C5'-O5'-P-O1P
6	F	401	AMP	C5'-O5'-P-O1P
5	A	602	ZQV	CAD-CAC-CAH-OAI
5	D	701	ZQV	CAD-CAC-CAH-OAI
6	E	401	AMP	C3'-C4'-C5'-O5'
5	A	602	ZQV	CAD-CAC-CAH-OAJ

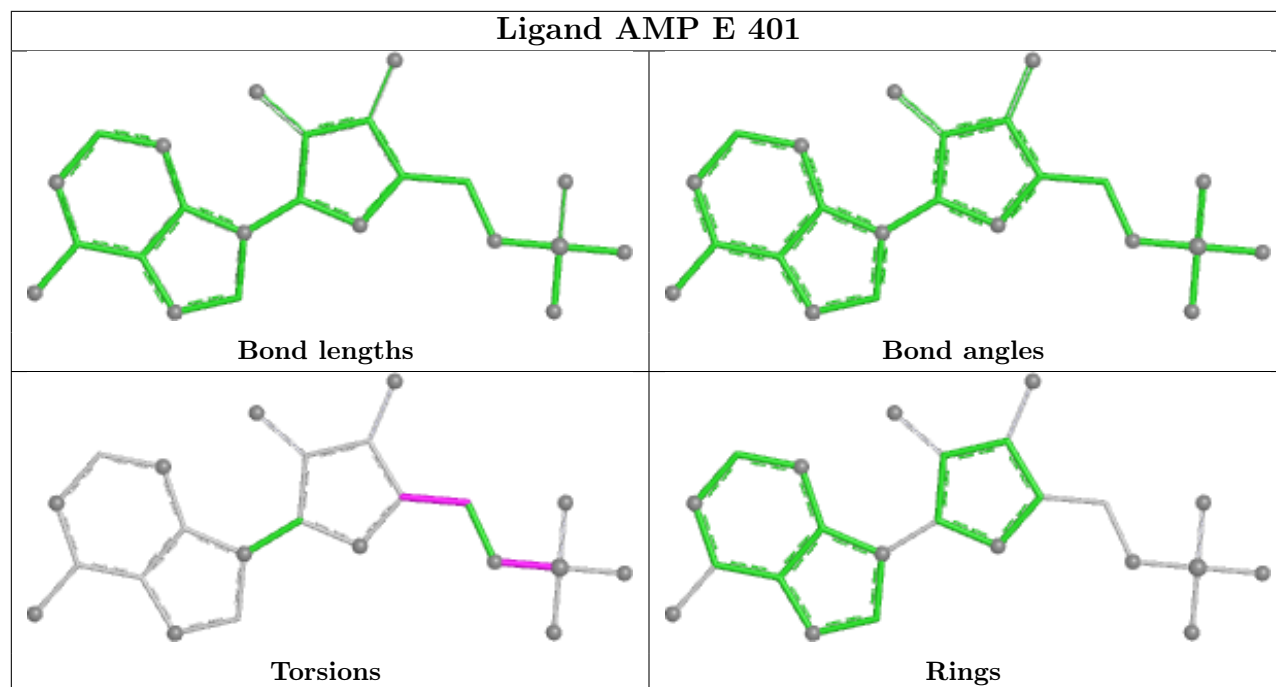
There are no ring outliers.

No monomer is involved in short contacts.

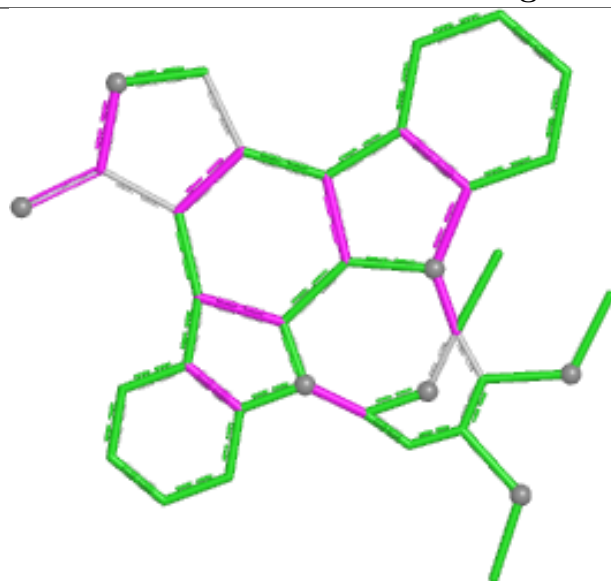
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

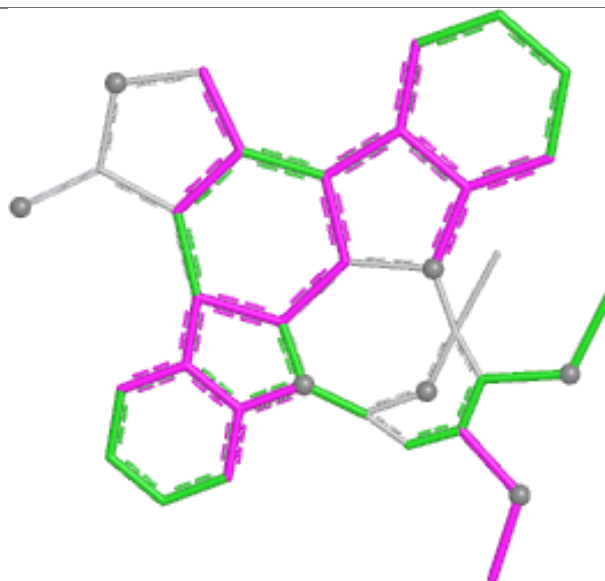




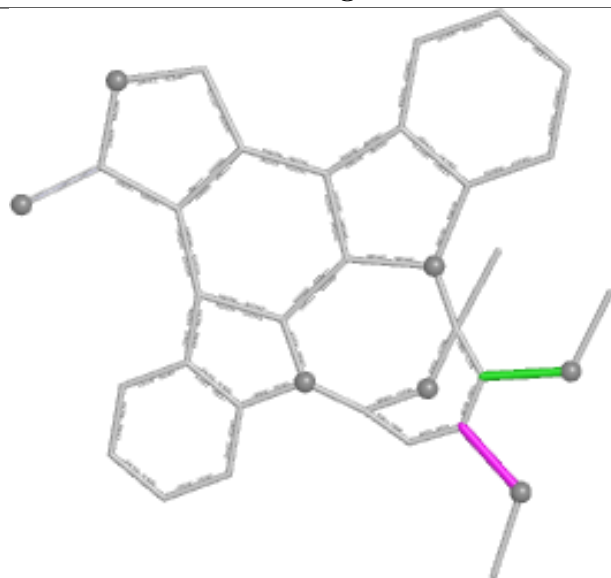
Ligand 4O7 A 601



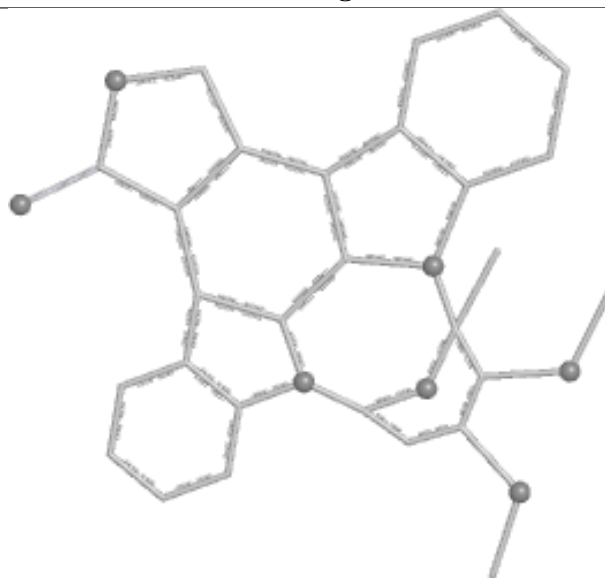
Bond lengths



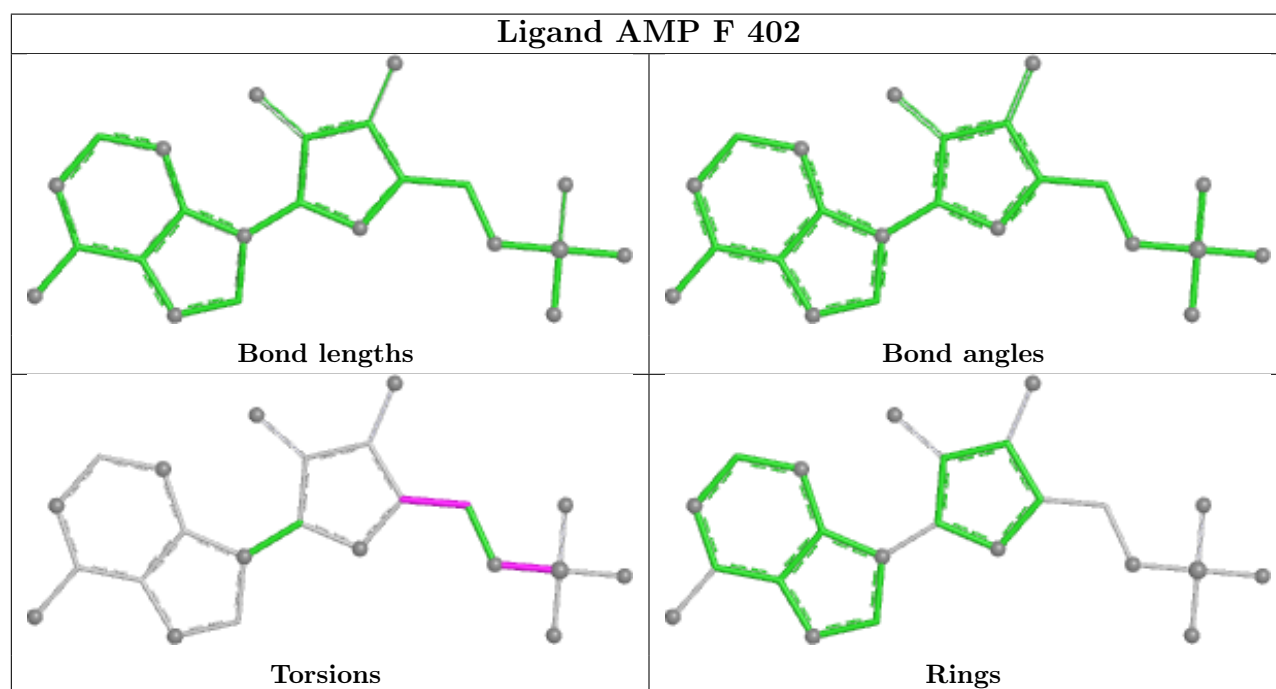
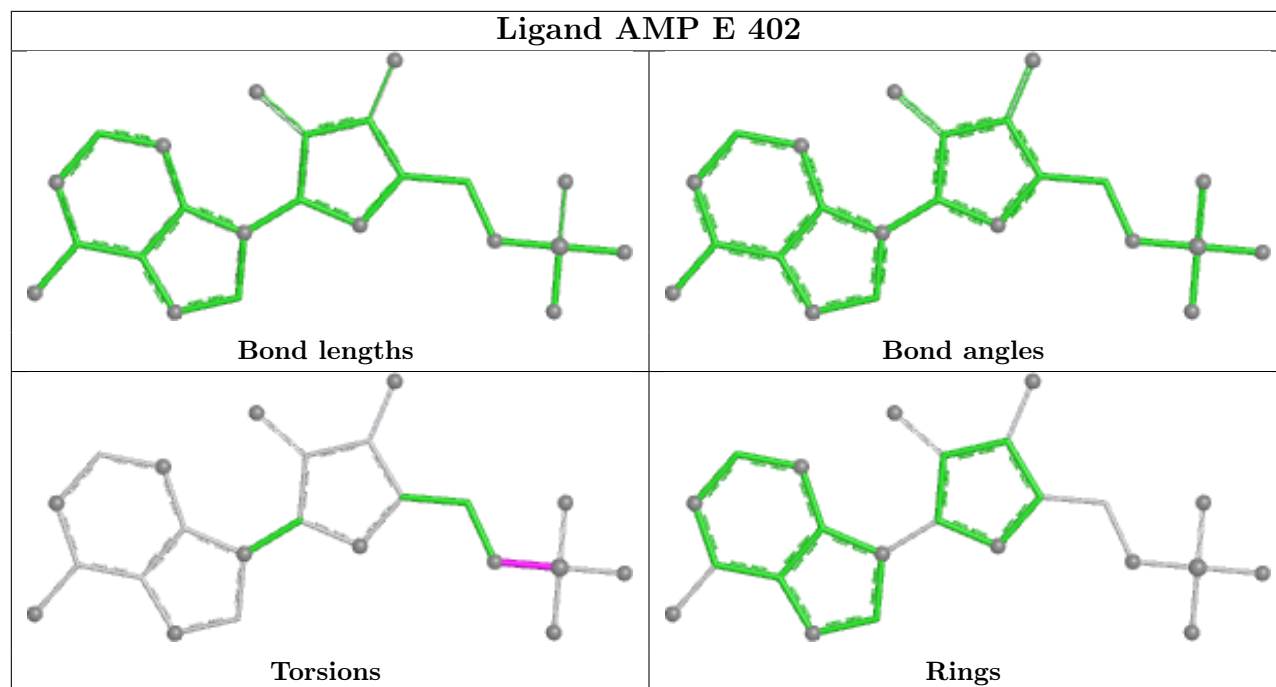
Bond angles

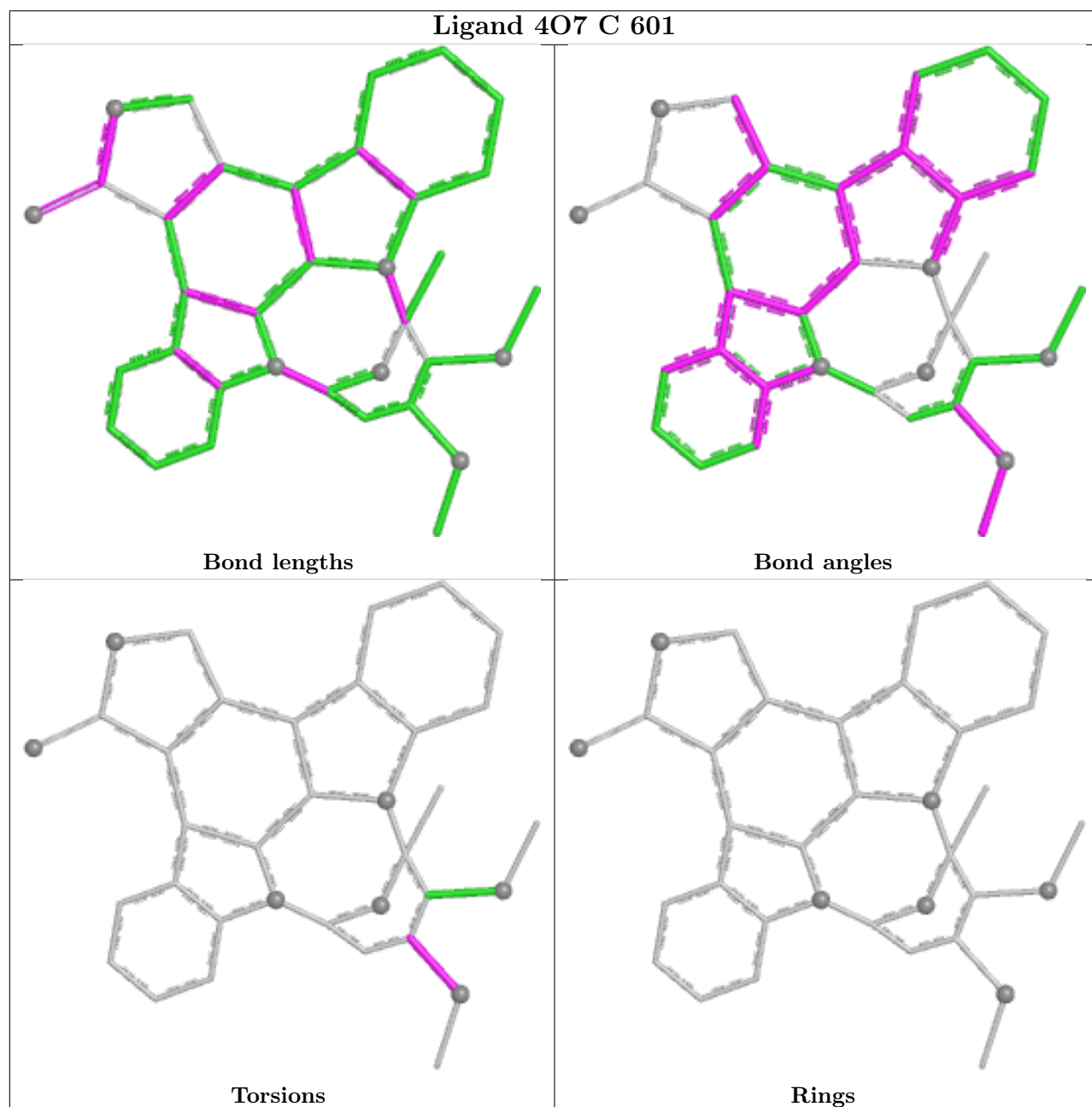
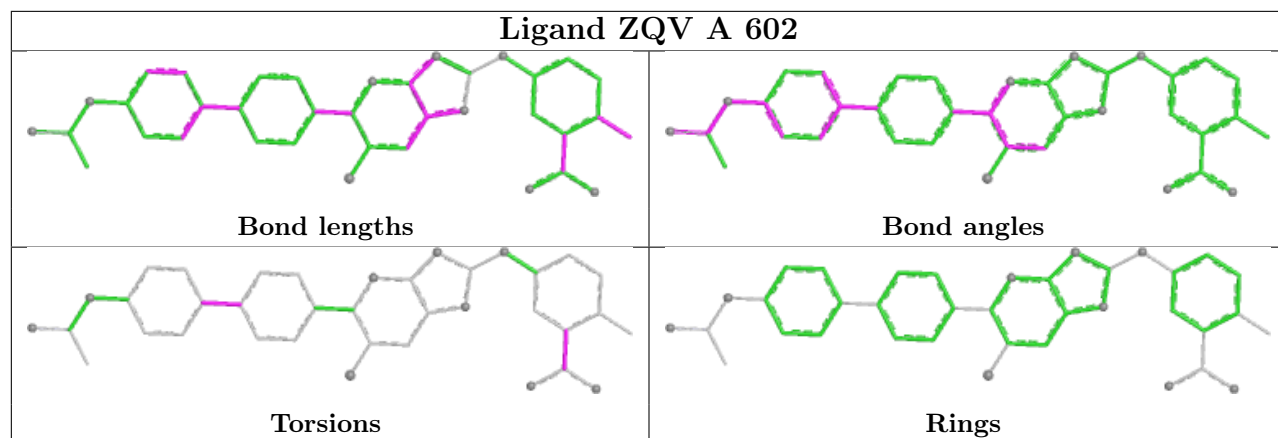


Torsions



Rings





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	434/565 (76%)	0.82	43 (9%) 13 11	52, 78, 115, 159	0
1	C	464/565 (82%)	0.59	37 (7%) 18 16	43, 69, 112, 139	0
2	B	183/270 (67%)	1.03	27 (14%) 5 5	52, 88, 162, 207	0
2	D	186/270 (68%)	0.79	22 (11%) 9 8	41, 66, 138, 173	0
3	E	303/336 (90%)	0.54	18 (5%) 28 23	44, 68, 103, 134	0
3	F	303/336 (90%)	0.64	27 (8%) 15 14	41, 66, 96, 116	0
All	All	1873/2342 (79%)	0.71	174 (9%) 14 12	41, 72, 116, 207	0

All (174) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	312	TYR	6.5
3	F	325	THR	5.3
1	A	293	VAL	5.0
3	F	107	TYR	4.9
3	E	26	ASN	4.7
3	F	307	ASP	4.4
2	D	226	ALA	4.4
1	A	551	ALA	4.3
1	C	303	THR	4.2
3	F	306	ASN	4.2
1	C	357	MET	4.1
2	D	224	ASP	4.1
3	F	326	GLY	4.0
3	F	303	VAL	4.0
2	D	225	PRO	3.9
3	F	308	VAL	3.9
1	A	311	LEU	3.8
3	F	304	ASP	3.8
2	B	184	PRO	3.8

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Mol	Chain	Res	Type	RSRZ
2	D	191	PRO	3.7
1	A	281	PRO	3.7
3	F	233	GLU	3.7
1	A	347	PRO	3.6
3	F	309	VAL	3.6
1	A	160	LEU	3.6
3	F	271	HIS	3.6
2	D	199	ARG	3.6
3	E	326	GLY	3.5
2	B	148	THR	3.5
1	C	173	SER	3.5
3	F	108	GLU	3.5
3	E	124	ASP	3.4
2	B	173	CYS	3.4
1	C	313	SER	3.4
2	D	270	ILE	3.4
1	A	323	ALA	3.4
1	C	171	ARG	3.4
2	B	175	ASP	3.4
2	B	209	HIS	3.3
2	B	270	ILE	3.3
1	A	171	ARG	3.3
1	A	398	LYS	3.3
1	C	377	SER	3.3
1	A	530	PRO	3.3
1	C	156	ALA	3.2
3	E	126	PHE	3.2
1	C	348	SER	3.1
1	A	27	PHE	3.1
1	A	174	CYS	3.1
1	A	176	SER	3.1
1	C	471	SER	3.1
1	C	336	GLN	3.1
1	A	532	LEU	3.0
3	E	323	VAL	3.0
1	A	157	ASP	3.0
2	B	193	VAL	3.0
1	C	362	GLY	2.9
3	E	304	ASP	2.9
1	C	474	ASP	2.9
1	A	307	VAL	2.8
3	E	220	PHE	2.8

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Mol	Chain	Res	Type	RSRZ
2	D	195	LYS	2.8
1	A	397	VAL	2.8
2	D	235	MET	2.8
3	F	236	ARG	2.7
3	F	126	PHE	2.7
2	B	146	LEU	2.7
3	F	112	HIS	2.7
3	F	223	HIS	2.7
2	D	146	LEU	2.7
2	B	133	TRP	2.7
2	B	172	LYS	2.7
3	E	306	ASN	2.7
1	A	405	GLY	2.7
1	C	304	GLU	2.7
2	D	194	CYS	2.6
1	C	109	GLY	2.6
1	C	530	PRO	2.6
2	D	111	ASN	2.6
2	D	200	PHE	2.6
1	A	191	ALA	2.6
2	D	77	ALA	2.6
1	A	310	SER	2.6
3	F	272	TYR	2.6
2	B	198	GLU	2.6
1	A	432	VAL	2.5
2	B	223	CYS	2.5
1	A	223	PHE	2.5
1	A	304	GLU	2.5
2	B	111	ASN	2.5
2	B	221	ILE	2.5
2	D	174	SER	2.5
3	E	233	GLU	2.5
1	A	473	ASP	2.4
1	C	305	SER	2.4
2	D	176	VAL	2.4
1	C	396	ALA	2.4
1	A	531	ARG	2.4
3	F	24	SER	2.4
1	C	306	GLU	2.4
1	C	309	ASN	2.4
3	F	232	ASP	2.4
2	B	187	TYR	2.4

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Mol	Chain	Res	Type	RSRZ
2	D	110	ASN	2.4
1	A	290	ASP	2.4
2	D	75	ALA	2.4
3	E	24	SER	2.4
3	F	125	SER	2.4
3	F	124	ASP	2.3
2	D	187	TYR	2.3
3	E	272	TYR	2.3
3	E	325	THR	2.3
1	A	399	LYS	2.3
1	A	362	GLY	2.3
1	A	308	MET	2.3
1	C	312	TYR	2.3
3	E	198	ILE	2.3
2	B	191	PRO	2.3
1	A	475	GLU	2.3
1	A	288	ILE	2.3
1	C	406	ILE	2.3
3	E	128	PRO	2.3
1	C	157	ASP	2.2
1	A	309	ASN	2.2
3	E	322	LEU	2.2
2	B	153	ILE	2.2
2	D	76	GLN	2.2
2	D	175	ASP	2.2
2	B	196	PRO	2.2
3	F	323	VAL	2.2
1	C	400	ALA	2.2
1	C	551	ALA	2.2
3	F	256	ASN	2.2
1	C	290	ASP	2.2
1	C	316	PRO	2.2
1	A	345	SER	2.2
2	B	192	TYR	2.2
3	E	121	TYR	2.2
1	A	400	ALA	2.2
2	D	237	ASN	2.1
1	C	347	PRO	2.1
1	C	308	MET	2.1
1	C	27	PHE	2.1
2	B	150	ASN	2.1
1	A	365	PRO	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	446	VAL	2.1
1	A	289	ASP	2.1
1	A	474	ASP	2.1
2	B	131	GLY	2.1
2	B	219	THR	2.1
3	E	183	PHE	2.1
3	F	322	LEU	2.1
2	B	190	GLU	2.1
1	C	177	PRO	2.1
1	C	361	PRO	2.1
3	E	234	LYS	2.1
1	A	291	GLU	2.1
2	D	192	TYR	2.1
3	F	106	ILE	2.1
1	A	374	ILE	2.1
1	C	108	HIS	2.0
2	B	145	GLN	2.0
1	C	132	ARG	2.0
3	F	147	ARG	2.0
2	B	197	GLU	2.0
1	C	450	TYR	2.0
1	A	305	SER	2.0
1	C	338	SER	2.0
1	C	358	HIS	2.0
3	F	237	VAL	2.0
2	B	110	ASN	2.0
1	A	8	ASP	2.0
1	C	339	GLU	2.0
2	B	113	VAL	2.0

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

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
1	TPO	C	172	11/12	0.54	0.30	92,95,96,97	1
1	TPO	A	172	11/12	0.70	0.19	90,92,94,94	0
2	SEP	D	108	10/11	0.83	0.14	69,73,77,78	0
2	SEP	B	108	10/11	0.88	0.11	80,81,86,86	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

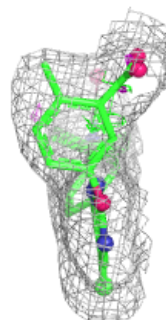
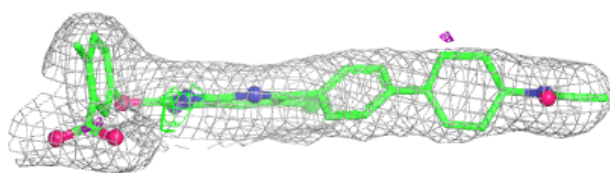
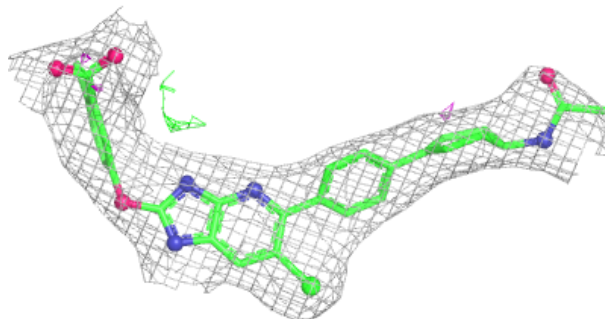
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
5	ZQV	A	602	37/37	0.91	0.12	70,72,78,78	0
5	ZQV	D	701	37/37	0.93	0.11	54,60,76,78	0
6	AMP	E	402	23/23	0.95	0.08	61,63,70,70	0
6	AMP	F	401	23/23	0.95	0.09	55,57,61,62	0
4	4O7	A	601	35/35	0.96	0.08	45,46,48,49	0
6	AMP	E	401	23/23	0.96	0.07	59,62,64,65	0
6	AMP	F	402	23/23	0.96	0.08	53,57,60,61	0
4	4O7	C	601	35/35	0.97	0.08	44,46,48,50	0

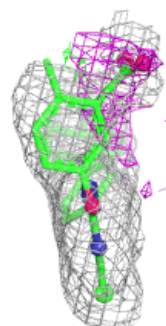
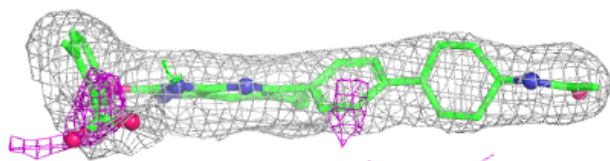
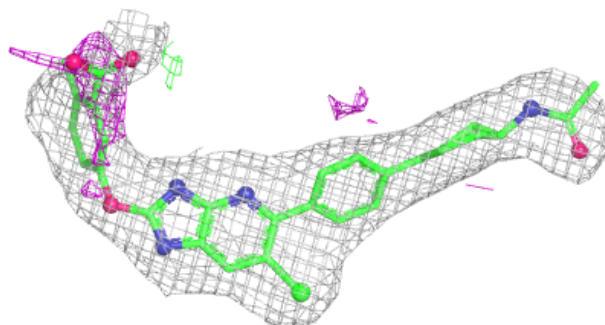
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around ZQV A 602:

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

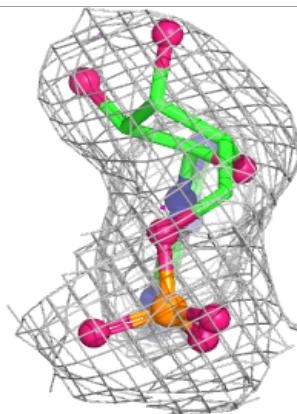
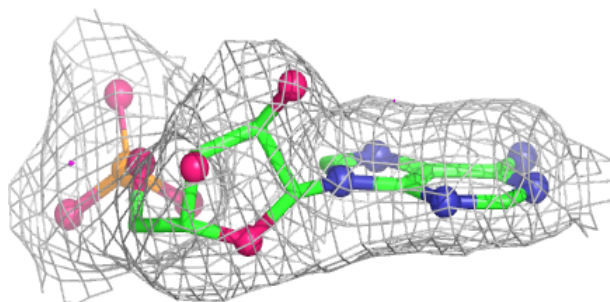
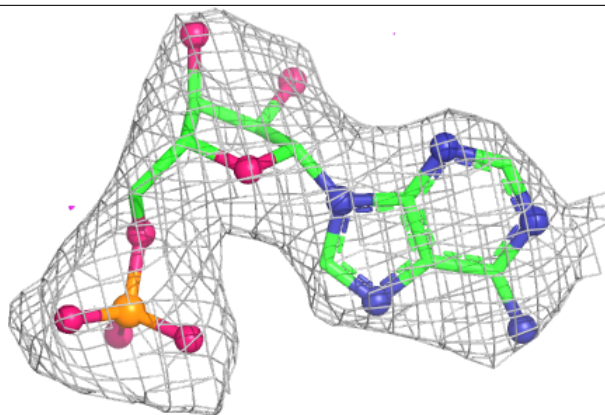
**Electron density around ZQV D 701:**

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



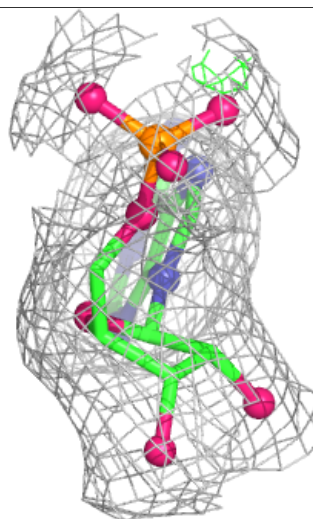
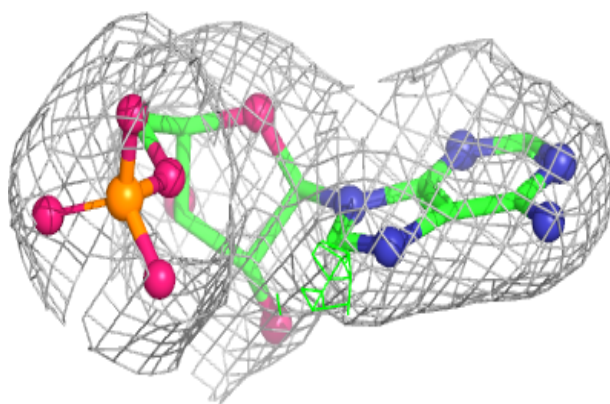
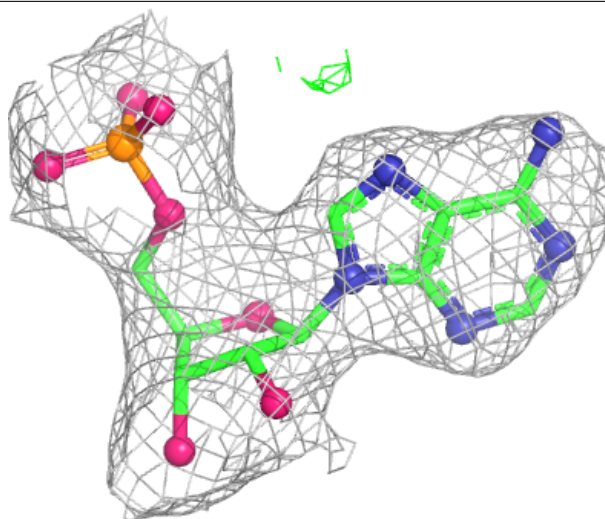
Electron density around AMP E 402:

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



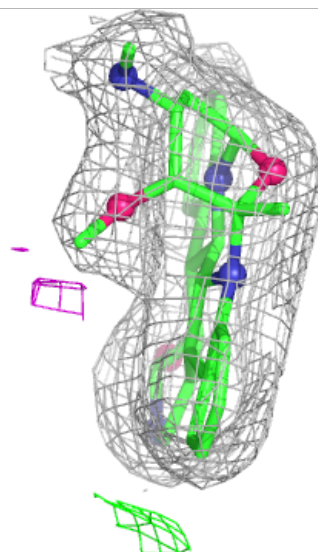
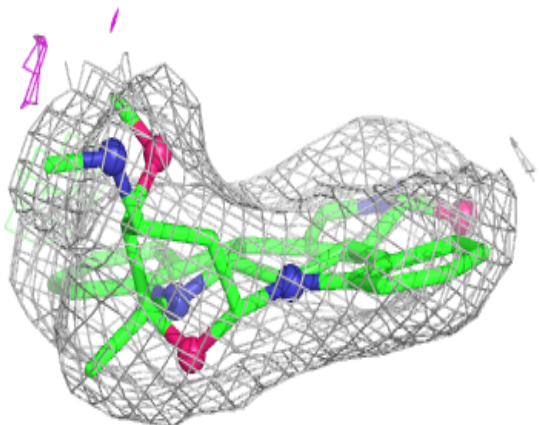
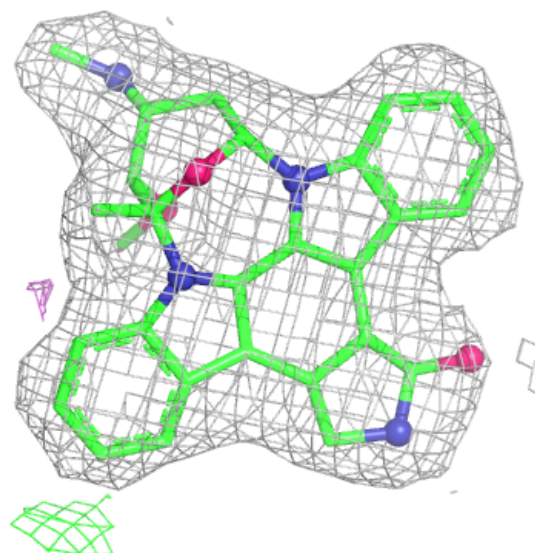
Electron density around AMP F 401:

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



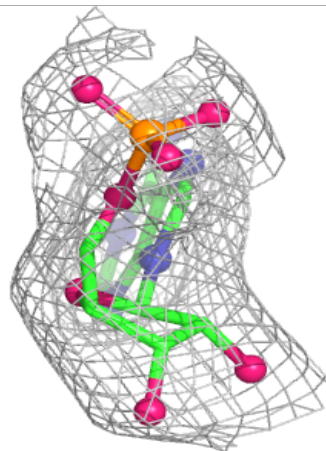
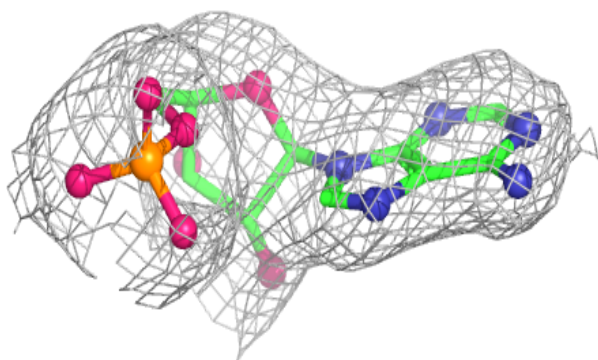
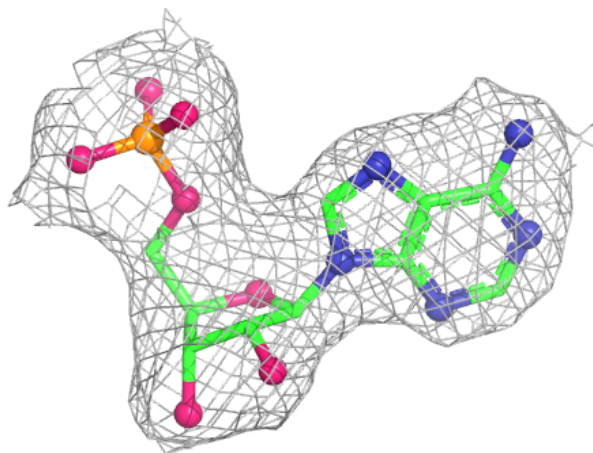
Electron density around 4O7 A 601:

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



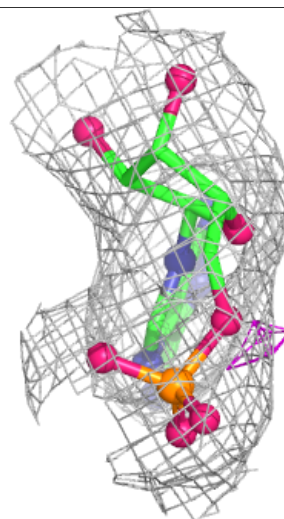
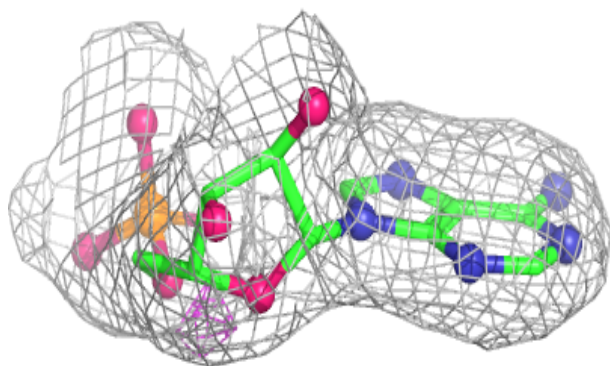
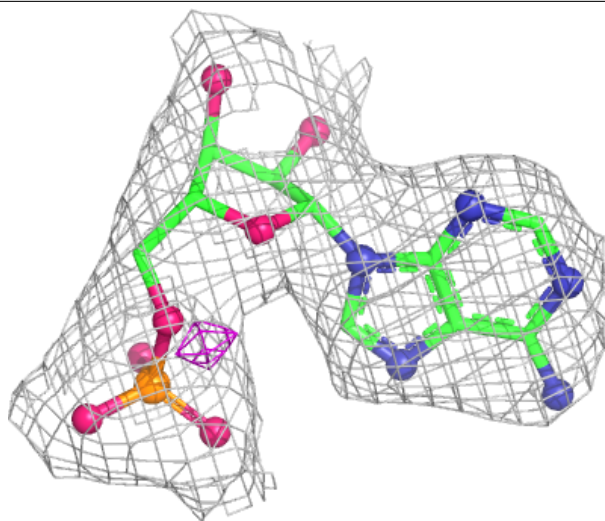
Electron density around AMP E 401:

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



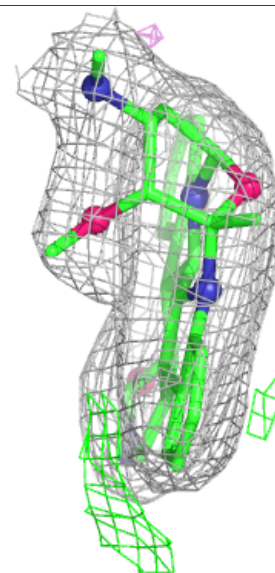
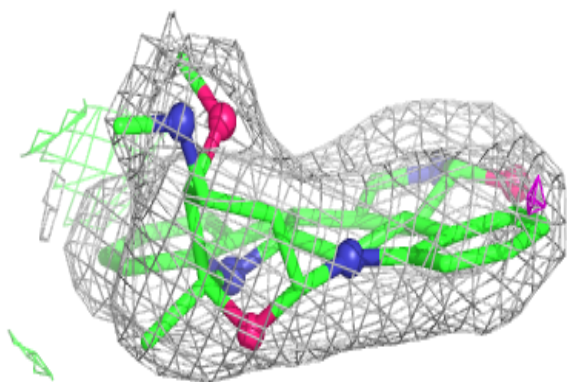
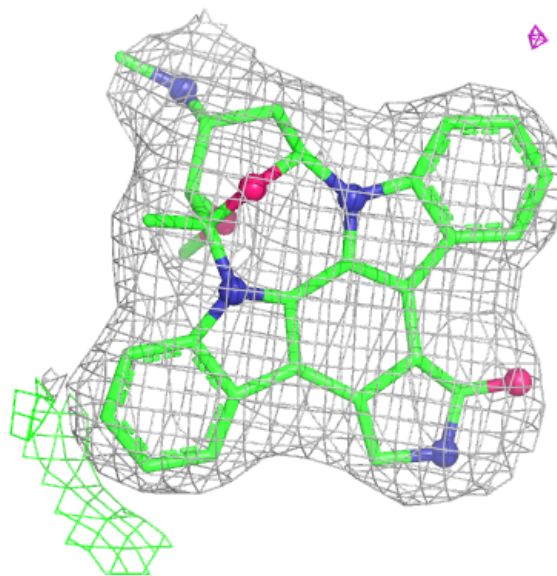
Electron density around AMP F 402:

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



Electron density around 4O7 C 601:

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.