



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:21 PM UTC

PDB ID : 6VT1 / pdb_00006vt1
Title : Naegleria gruberi RNA ligase D172A mutant apo
Authors : Unciuleac, M.C.; Goldgur, Y.; Shuman, S.
Deposited on : 2020-02-12
Resolution : 2.38 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

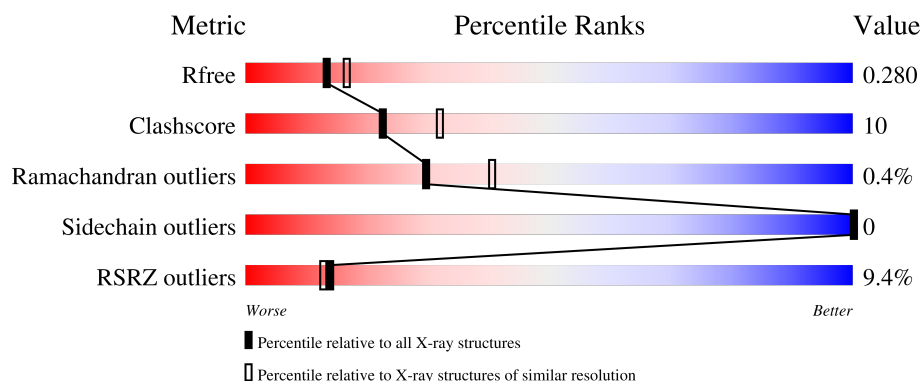
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.38 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	7164 (2.40-2.36)
Clashscore	190562	7722 (2.40-2.36)
Ramachandran outliers	187476	7626 (2.40-2.36)
Sidechain outliers	187428	7627 (2.40-2.36)
RSRZ outliers	180081	7170 (2.40-2.36)

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	340	4% 79% 20% .
1	B	340	8% 84% 15% .
1	C	340	16% 67% 31% .
1	D	340	8% 78% 20% ..

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 11143 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 RNA Ligase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	339	Total	C	N	O	S	0	0	0
			2708	1706	479	513	10			
1	B	339	Total	C	N	O	S	0	0	0
			2708	1706	479	513	10			
1	C	339	Total	C	N	O	S	0	0	0
			2708	1706	479	513	10			
1	D	339	Total	C	N	O	S	0	0	0
			2708	1706	479	513	10			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	SER	-	expression tag	UNP D2W2Z5
A	172	ALA	ASP	engineered mutation	UNP D2W2Z5
B	0	SER	-	expression tag	UNP D2W2Z5
B	172	ALA	ASP	engineered mutation	UNP D2W2Z5
C	0	SER	-	expression tag	UNP D2W2Z5
C	172	ALA	ASP	engineered mutation	UNP D2W2Z5
D	0	SER	-	expression tag	UNP D2W2Z5
D	172	ALA	ASP	engineered mutation	UNP D2W2Z5

- Molecule 2 is ADENOSINE MONOPHOSPHATE (CCD ID: AMP) (formula: C₁₀H₁₄N₅O₇P) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			22	10	5	6	1		
2	B	1	Total	C	N	O	P	0	0
			22	10	5	6	1		
2	C	1	Total	C	N	O	P	0	0
			22	10	5	6	1		
2	D	1	Total	C	N	O	P	0	0
			22	10	5	6	1		

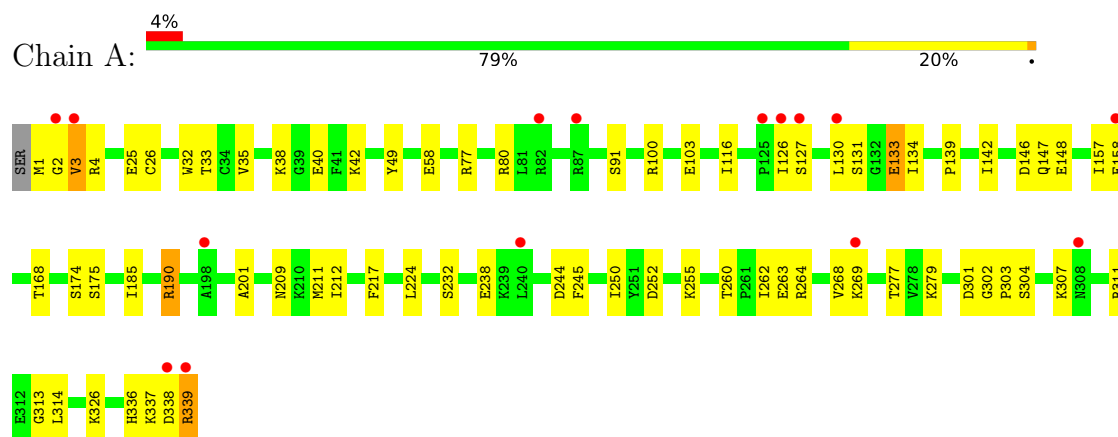
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	73	Total	O	0	0
			73	73		
3	B	48	Total	O	0	0
			48	48		
3	C	45	Total	O	0	0
			45	45		
3	D	57	Total	O	0	0
			57	57		

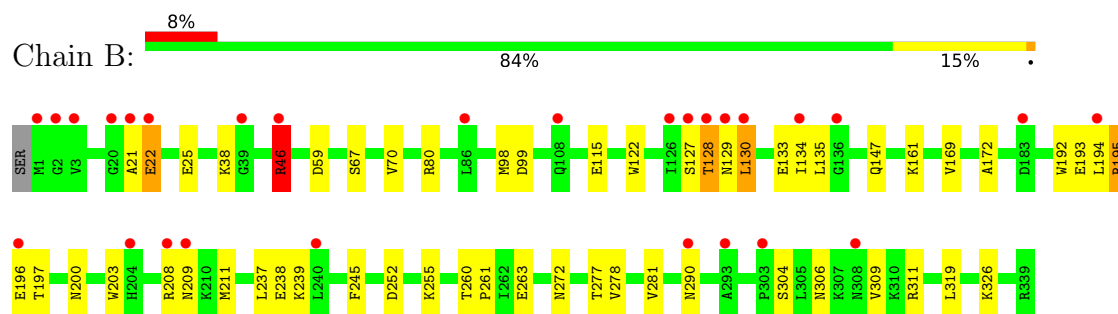
3 Residue-property plots [i](#)

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.

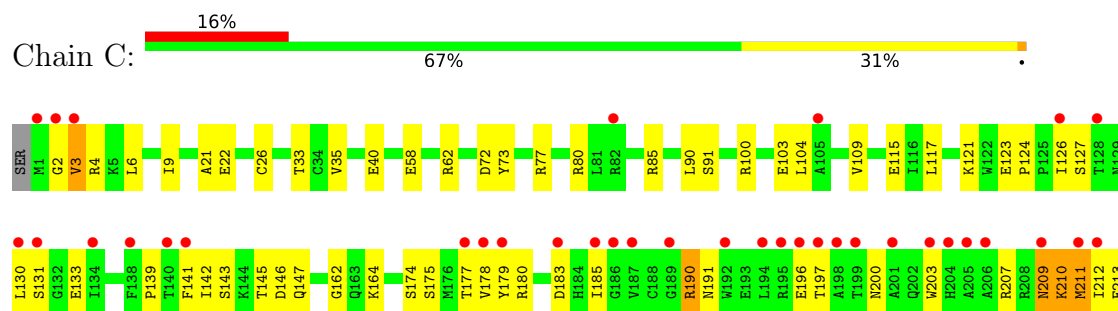
• Molecule 1: RNA Ligase

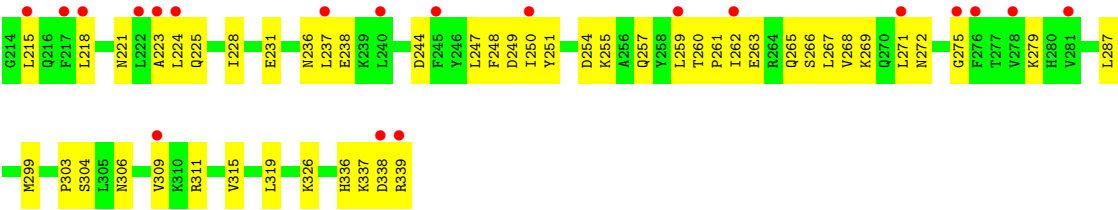


• Molecule 1: RNA Ligase

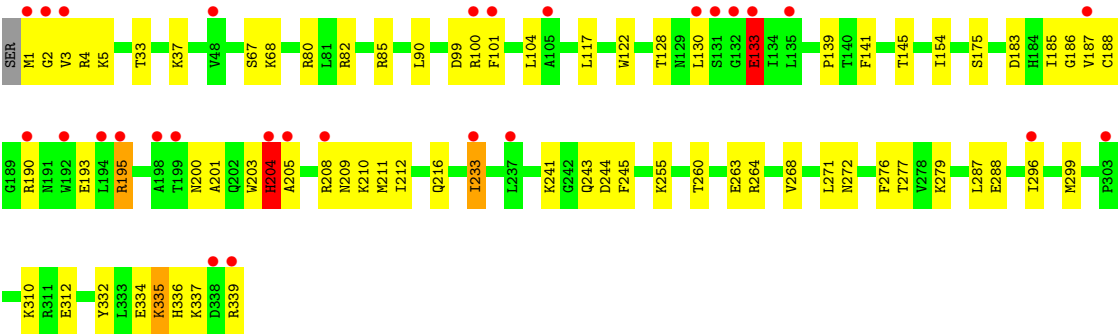
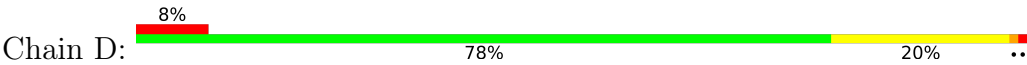


• Molecule 1: RNA Ligase





● Molecule 1: RNA Ligase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	125.27Å 104.18Å 120.56Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.04 – 2.38 49.04 – 2.38	Depositor EDS
% Data completeness (in resolution range)	94.2 (49.04-2.38) 91.9 (49.04-2.38)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.29 (at 2.37Å)	Xtriage
Refinement program	PHENIX 1.17.1-3660	Depositor
R, R_{free}	0.232 , 0.276 0.238 , 0.280	Depositor DCC
R_{free} test set	1974 reflections (3.16%)	wwPDB-VP
Wilson B-factor (Å ²)	47.2	Xtriage
Anisotropy	0.211	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 35.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.025 for l,-k,h	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	11143	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 29.94 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.4222e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² 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: 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	0.51	1/2755 (0.0%)	0.86	14/3710 (0.4%)
1	B	0.50	0/2755	0.84	8/3710 (0.2%)
1	C	0.53	0/2755	0.88	10/3710 (0.3%)
1	D	0.58	3/2755 (0.1%)	0.93	19/3710 (0.5%)
All	All	0.53	4/11020 (0.0%)	0.88	51/14840 (0.3%)

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	B	0	2
1	D	0	2
All	All	0	4

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	195	ARG	CZ-NH1	6.03	1.41	1.32
1	A	190	ARG	CD-NE	5.42	1.53	1.46
1	D	154	ILE	CA-CB	-5.21	1.47	1.54
1	D	133	GLU	CB-CG	5.12	1.67	1.52

All (51) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	115	GLU	CA-CB-CG	-11.79	90.51	114.10
1	C	22	GLU	CA-CB-CG	9.55	133.19	114.10
1	A	338	ASP	CB-CA-C	-8.99	99.81	111.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	46	ARG	CG-CD-NE	8.90	131.57	112.00
1	D	195	ARG	CB-CG-CD	-8.58	91.57	111.30
1	C	210	LYS	CA-CB-CG	-8.28	97.53	114.10
1	A	338	ASP	CA-C-N	-8.19	106.95	121.70
1	A	338	ASP	C-N-CA	-8.19	106.95	121.70
1	D	99	ASP	CA-C-N	-8.19	108.04	122.62
1	D	99	ASP	C-N-CA	-8.19	108.04	122.62
1	D	100	ARG	CG-CD-NE	8.03	129.65	112.00
1	D	204	HIS	N-CA-CB	-7.88	98.64	110.07
1	A	269	LYS	CB-CA-C	-7.86	98.86	110.96
1	A	339	ARG	CB-CG-CD	7.38	128.26	111.30
1	D	100	ARG	CB-CG-CD	-7.26	94.60	111.30
1	A	269	LYS	CB-CG-CD	-6.86	95.52	111.30
1	D	100	ARG	CA-CB-CG	6.80	127.70	114.10
1	B	128	THR	N-CA-C	6.72	120.05	109.96
1	D	195	ARG	CA-CB-CG	6.45	126.99	114.10
1	B	22	GLU	CB-CG-CD	-6.42	101.68	112.60
1	A	133	GLU	CA-CB-CG	-6.39	101.33	114.10
1	A	339	ARG	CA-CB-CG	-6.29	101.51	114.10
1	C	209	ASN	CA-C-N	6.14	131.12	122.46
1	C	209	ASN	C-N-CA	6.14	131.12	122.46
1	A	269	LYS	CA-CB-CG	5.96	126.02	114.10
1	D	190	ARG	CG-CD-NE	-5.93	98.95	112.00
1	C	211	MET	CB-CG-SD	-5.84	95.18	112.70
1	D	210	LYS	CA-CB-CG	5.81	125.72	114.10
1	D	128	THR	CA-C-N	-5.80	111.88	122.38
1	D	128	THR	C-N-CA	-5.80	111.88	122.38
1	C	133	GLU	CA-CB-CG	5.78	125.66	114.10
1	A	269	LYS	CG-CD-CE	5.77	124.57	111.30
1	D	100	ARG	CD-NE-CZ	5.74	132.44	124.40
1	C	21	ALA	CA-C-N	5.72	132.51	122.79
1	C	21	ALA	C-N-CA	5.72	132.51	122.79
1	B	208	ARG	CA-CB-CG	5.71	125.53	114.10
1	B	46	ARG	CB-CG-CD	5.57	124.10	111.30
1	B	46	ARG	CA-CB-CG	5.54	125.17	114.10
1	B	195	ARG	CG-CD-NE	5.49	124.07	112.00
1	C	183	ASP	CB-CA-C	-5.45	100.55	110.63
1	D	190	ARG	CB-CG-CD	-5.42	98.83	111.30
1	D	233	ILE	CG1-CB-CG2	-5.33	94.71	110.70
1	D	335	LYS	CG-CD-CE	5.29	123.48	111.30
1	A	307	LYS	CA-CB-CG	-5.28	103.54	114.10
1	A	307	LYS	CD-CE-NZ	5.27	128.76	111.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	133	GLU	CA-CB-CG	5.21	124.53	114.10
1	D	233	ILE	CA-CB-CG1	5.12	119.11	110.40
1	A	133	GLU	CA-C-N	5.12	128.97	122.37
1	A	133	GLU	C-N-CA	5.12	128.97	122.37
1	C	190	ARG	CB-CG-CD	5.09	123.00	111.30
1	D	204	HIS	CB-CG-ND1	-5.04	115.14	122.70

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	21	ALA	Peptide
1	B	46	ARG	Sidechain
1	D	133	GLU	Sidechain
1	D	204	HIS	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	2708	0	2709	52	0
1	B	2708	0	2709	39	0
1	C	2708	0	2708	84	0
1	D	2708	0	2709	60	0
2	A	22	0	12	0	0
2	B	22	0	12	1	0
2	C	22	0	12	2	0
2	D	22	0	12	0	0
3	A	73	0	0	10	0
3	B	48	0	0	6	0
3	C	45	0	0	5	0
3	D	57	0	0	6	0
All	All	11143	0	10883	229	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 10.

All (229) 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:D:85:ARG:NH1	1:D:334:GLU:CD	1.99	1.21
1:C:146:ASP:OD1	1:C:190:ARG:NH2	1.77	1.15
1:D:85:ARG:NH1	1:D:334:GLU:OE2	1.84	1.08
1:D:188:CYS:SG	3:D:554:HOH:O	2.13	1.04
1:D:85:ARG:NH1	1:D:334:GLU:OE1	1.85	1.02
1:B:25:GLU:OE1	3:B:501:HOH:O	1.84	0.94
1:B:195:ARG:NH1	1:B:197:THR:HG22	1.83	0.92
1:A:146:ASP:OD2	1:A:190:ARG:NH2	2.05	0.89
1:C:145:THR:O	2:C:401:AMP:O3'	1.94	0.85
1:D:85:ARG:HH12	1:D:334:GLU:CD	1.71	0.83
1:D:200:ASN:C	1:D:204:HIS:HD2	1.89	0.81
1:B:172:ALA:O	3:B:502:HOH:O	1.99	0.79
1:C:207:ARG:O	1:C:210:LYS:HG2	1.82	0.79
1:B:130:LEU:HB2	1:B:133:GLU:HG3	1.65	0.77
1:A:336:HIS:HB3	1:A:339:ARG:HD3	1.68	0.75
1:D:332:TYR:OH	1:D:339:ARG:NH1	2.20	0.75
1:D:85:ARG:HH11	1:D:334:GLU:CD	1.75	0.75
1:D:80:ARG:NH1	3:D:504:HOH:O	2.21	0.74
1:B:194:LEU:HD22	1:B:200:ASN:HD21	1.53	0.74
1:B:99:ASP:OD1	3:B:503:HOH:O	2.06	0.73
1:A:158:GLU:HG3	1:A:158:GLU:O	1.88	0.73
1:B:161:LYS:HA	3:B:506:HOH:O	1.88	0.72
1:C:275:GLY:O	3:C:501:HOH:O	2.07	0.72
1:C:126:ILE:HD13	1:C:131:SER:OG	1.90	0.72
1:B:195:ARG:HH11	1:B:197:THR:HG22	1.54	0.72
1:A:238:GLU:OE2	1:A:304:SER:OG	2.08	0.70
1:A:126:ILE:HD11	1:A:131:SER:OG	1.93	0.68
1:A:175:SER:O	3:A:502:HOH:O	2.09	0.68
1:D:145:THR:HG21	1:D:175:SER:OG	1.93	0.68
1:D:193:GLU:OE2	3:D:501:HOH:O	2.11	0.68
1:D:260:THR:OG1	1:D:263:GLU:HG3	1.94	0.67
1:D:5:LYS:O	3:D:502:HOH:O	2.12	0.67
1:C:124:PRO:HG2	1:C:339:ARG:NH1	2.09	0.67
1:A:133:GLU:CD	1:B:129:ASN:H	2.03	0.66
1:A:260:THR:OG1	3:A:501:HOH:O	1.92	0.66
1:A:262:ILE:N	3:A:501:HOH:O	2.27	0.66
1:C:250:ILE:HB	1:C:259:LEU:HD12	1.76	0.66
1:A:337:LYS:O	3:A:503:HOH:O	2.15	0.64
1:C:224:LEU:HD23	1:C:250:ILE:HG12	1.79	0.64
1:D:130:LEU:HD23	1:D:201:ALA:HB3	1.81	0.63
1:C:265:GLN:HA	1:C:268:VAL:HG12	1.78	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:335:LYS:HG2	1:D:336:HIS:CD2	2.33	0.63
1:D:264:ARG:O	1:D:268:VAL:HG23	1.98	0.63
1:C:259:LEU:HB3	1:C:263:GLU:HG3	1.79	0.63
1:A:133:GLU:HG3	1:A:134:ILE:N	2.14	0.62
1:A:42:LYS:NZ	3:A:509:HOH:O	2.31	0.62
1:D:200:ASN:C	1:D:204:HIS:CD2	2.76	0.62
1:A:133:GLU:OE1	1:B:130:LEU:N	2.33	0.61
1:A:133:GLU:HG3	1:A:134:ILE:H	1.65	0.61
1:B:239:LYS:HB2	1:B:306:ASN:HD22	1.66	0.61
1:A:232:SER:O	1:A:232:SER:OG	2.14	0.60
1:C:247:LEU:HG	1:C:248:PHE:N	2.15	0.60
1:A:263:GLU:N	3:A:501:HOH:O	1.96	0.60
1:D:82:ARG:NH2	3:D:504:HOH:O	2.35	0.60
1:B:22:GLU:O	1:B:38:LYS:HD3	2.02	0.59
1:C:80:ARG:HH22	1:C:338:ASP:C	2.11	0.59
1:C:115:GLU:CD	1:C:115:GLU:H	2.11	0.59
1:C:200:ASN:ND2	1:C:203:TRP:H	1.98	0.59
1:D:82:ARG:HH11	1:D:337:LYS:HB3	1.67	0.59
1:B:161:LYS:NZ	1:B:290:ASN:OD1	2.36	0.59
1:A:209:ASN:HB2	1:A:277:THR:O	2.04	0.58
1:C:209:ASN:ND2	1:C:211:MET:SD	2.76	0.58
1:C:250:ILE:HB	1:C:259:LEU:CD1	2.34	0.58
1:A:336:HIS:O	1:A:339:ARG:HD3	2.04	0.57
1:B:130:LEU:HB2	1:B:133:GLU:CG	2.35	0.57
1:B:260:THR:OG1	1:B:263:GLU:HG3	2.04	0.57
1:C:175:SER:HB3	1:C:190:ARG:HB2	1.87	0.57
1:A:130:LEU:HD23	1:A:201:ALA:HB3	1.87	0.57
1:A:224:LEU:HD23	1:A:250:ILE:HG12	1.86	0.57
1:A:252:ASP:CG	1:A:255:LYS:HG3	2.29	0.57
1:B:169:VAL:HG13	1:B:311:ARG:HD2	1.86	0.57
1:C:185:ILE:HG21	1:C:212:ILE:HG13	1.85	0.57
1:C:262:ILE:HD12	1:C:262:ILE:H	1.69	0.57
1:B:59:ASP:OD2	1:C:85:ARG:NH1	2.38	0.56
1:B:238:GLU:OE2	1:B:304:SER:OG	2.22	0.56
1:C:306:ASN:HB3	1:C:309:VAL:HG12	1.87	0.55
1:D:312:GLU:HG3	1:D:332:TYR:CE1	2.41	0.55
1:A:303:PRO:O	1:A:311:ARG:HD3	2.07	0.55
1:D:145:THR:HG23	1:D:145:THR:O	2.06	0.55
1:C:162:GLY:O	1:C:164:LYS:HE2	2.07	0.55
1:C:180:ARG:NH1	1:C:221:ASN:OD1	2.40	0.55
1:C:215:LEU:O	3:C:503:HOH:O	2.18	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:33:THR:HG23	1:D:90:LEU:O	2.07	0.55
1:D:244:ASP:OD2	1:D:279:LYS:HE3	2.06	0.54
1:D:209:ASN:HB2	1:D:277:THR:O	2.08	0.54
1:D:130:LEU:HD11	1:D:233:ILE:CD1	2.39	0.53
1:C:261:PRO:HG2	1:C:319:LEU:HD23	1.90	0.53
1:A:103:GLU:O	1:D:310:LYS:NZ	2.38	0.53
1:B:209:ASN:ND2	1:B:278:VAL:HG22	2.24	0.53
1:A:133:GLU:OE1	1:B:129:ASN:N	2.42	0.53
1:A:244:ASP:OD2	1:A:279:LYS:NZ	2.33	0.53
1:C:58:GLU:OE1	1:C:77:ARG:HD3	2.09	0.53
1:B:237:LEU:HG	1:B:309:VAL:HG11	1.91	0.53
1:C:231:GLU:HA	1:C:236:ASN:HB2	1.91	0.52
1:D:271:LEU:HG	1:D:276:PHE:HB2	1.91	0.52
1:A:1:MET:HG3	1:A:148:GLU:HB2	1.92	0.52
1:C:336:HIS:O	1:C:338:ASP:N	2.42	0.51
1:C:191:ASN:N	3:C:502:HOH:O	2.15	0.51
1:C:4:ARG:HB3	1:C:6:LEU:HD23	1.91	0.51
1:A:336:HIS:HB3	1:A:339:ARG:CD	2.39	0.51
1:C:147:GLN:O	1:C:326:LYS:HE2	2.11	0.51
1:C:223:ALA:HB3	1:C:251:TYR:HB3	1.91	0.51
1:B:272:ASN:OD1	1:B:277:THR:HA	2.11	0.50
1:C:237:LEU:HG	1:C:309:VAL:HG21	1.93	0.50
1:A:139:PRO:HG2	1:A:142:ILE:HD12	1.93	0.50
1:D:255:LYS:O	1:D:255:LYS:HE3	2.11	0.50
1:A:301:ASP:OD1	1:A:302:GLY:N	2.44	0.50
1:D:185:ILE:HG21	1:D:212:ILE:HG23	1.93	0.50
1:C:9:ILE:HG22	1:C:109:VAL:HA	1.94	0.50
1:D:2:GLY:C	1:D:4:ARG:H	2.19	0.50
1:C:145:THR:N	3:C:508:HOH:O	2.29	0.49
1:C:287:LEU:HD11	1:C:299:MET:HE1	1.94	0.49
1:A:168:THR:OG1	3:A:504:HOH:O	2.20	0.49
1:A:133:GLU:CG	1:A:134:ILE:H	2.26	0.49
1:A:264:ARG:O	1:A:268:VAL:HG23	2.12	0.49
1:D:201:ALA:C	1:D:204:HIS:HB2	2.37	0.49
1:D:183:ASP:OD1	1:D:183:ASP:N	2.40	0.49
1:D:139:PRO:HB3	1:D:141:PHE:CZ	2.48	0.49
1:C:197:THR:HB	1:C:200:ASN:HB2	1.95	0.48
1:C:244:ASP:OD2	1:C:279:LYS:HE3	2.13	0.48
1:D:130:LEU:HD11	1:D:233:ILE:HD12	1.95	0.48
1:D:200:ASN:O	1:D:204:HIS:CD2	2.66	0.48
1:D:272:ASN:OD1	1:D:277:THR:HA	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:335:LYS:HG2	1:D:336:HIS:CG	2.48	0.48
1:C:266:SER:O	1:C:269:LYS:HB3	2.13	0.48
1:D:201:ALA:HB1	1:D:243:GLN:OE1	2.12	0.48
1:B:161:LYS:HG2	3:B:506:HOH:O	2.14	0.48
1:A:116:ILE:HG13	3:A:527:HOH:O	2.13	0.48
1:C:178:VAL:HG21	1:C:211:MET:HE2	1.94	0.48
1:D:82:ARG:NH1	1:D:337:LYS:HB3	2.28	0.48
1:B:239:LYS:HB2	1:B:306:ASN:ND2	2.29	0.48
1:A:58:GLU:OE2	1:A:77:ARG:NH1	2.46	0.47
1:C:2:GLY:O	1:C:3:VAL:HB	2.13	0.47
1:C:85:ARG:HH21	1:C:90:LEU:HB2	1.78	0.47
1:C:174:SER:O	1:C:228:ILE:HG22	2.14	0.47
1:C:223:ALA:O	1:C:250:ILE:HA	2.14	0.47
1:D:67:SER:O	3:D:503:HOH:O	2.19	0.47
1:B:134:ILE:HG23	1:B:192:TRP:HB3	1.96	0.47
1:A:130:LEU:HD13	1:A:174:SER:OG	2.15	0.47
1:C:104:LEU:HD11	1:C:117:LEU:HD11	1.97	0.47
1:A:32:TRP:CG	1:A:49:TYR:HH	2.33	0.47
1:C:303:PRO:HD2	1:C:311:ARG:NH1	2.30	0.47
1:D:185:ILE:HD12	1:D:216:GLN:HG3	1.95	0.47
1:B:135:LEU:HB3	1:B:193:GLU:HG2	1.96	0.47
1:B:290:ASN:HA	3:B:506:HOH:O	2.14	0.47
1:A:35:VAL:HG23	1:A:91:SER:HB2	1.96	0.46
1:C:4:ARG:HH11	1:C:4:ARG:HG3	1.80	0.46
1:A:147:GLN:O	1:A:326:LYS:HE2	2.15	0.46
1:C:35:VAL:HG23	1:C:91:SER:HB2	1.97	0.46
1:D:201:ALA:HA	1:D:204:HIS:HB2	1.96	0.46
1:C:260:THR:OG1	1:C:263:GLU:HG2	2.14	0.46
1:C:268:VAL:O	1:C:272:ASN:ND2	2.27	0.46
1:C:221:ASN:HB2	1:C:254:ASP:OD2	2.16	0.46
1:C:178:VAL:CG2	1:C:211:MET:HE2	2.46	0.46
1:D:133:GLU:HB3	1:D:195:ARG:HH22	1.81	0.46
1:D:200:ASN:OD1	1:D:203:TRP:HD1	1.99	0.46
1:C:62:ARG:NH1	1:C:103:GLU:OE2	2.48	0.46
1:C:177:THR:OG1	1:C:225:GLN:HG2	2.16	0.45
1:D:205:ALA:HA	1:D:208:ARG:HG2	1.98	0.45
1:B:281:VAL:HG21	2:B:401:AMP:N6	2.31	0.45
1:C:40:GLU:OE2	1:C:100:ARG:NH2	2.49	0.45
1:C:218:LEU:HD12	1:C:218:LEU:HA	1.79	0.45
1:D:336:HIS:HB3	1:D:339:ARG:HG3	1.98	0.45
1:B:211:MET:HE1	1:B:245:PHE:CZ	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:262:ILE:HD11	1:C:319:LEU:HD22	1.99	0.45
1:A:133:GLU:CG	1:A:134:ILE:N	2.80	0.45
1:C:121:LYS:NZ	1:C:123:GLU:OE2	2.48	0.45
1:D:104:LEU:HD11	1:D:117:LEU:HD21	1.98	0.45
1:D:287:LEU:HD23	1:D:288:GLU:C	2.42	0.45
1:A:25:GLU:CD	1:A:38:LYS:HD2	2.41	0.45
1:A:127:SER:OG	1:A:130:LEU:HG	2.17	0.45
1:C:336:HIS:C	1:C:338:ASP:H	2.25	0.45
1:C:73:TYR:OH	3:C:504:HOH:O	2.21	0.45
1:A:260:THR:C	3:A:501:HOH:O	2.60	0.44
1:C:255:LYS:HG3	1:C:257:GLN:HG2	1.98	0.44
1:B:211:MET:HE1	1:B:245:PHE:CE1	2.51	0.44
1:C:139:PRO:HB3	1:C:141:PHE:CZ	2.52	0.44
1:D:80:ARG:HD2	1:D:122:TRP:CD2	2.53	0.44
1:C:267:LEU:C	1:C:271:LEU:HD12	2.42	0.44
1:D:1:MET:HG2	1:D:2:GLY:H	1.82	0.44
1:A:1:MET:HE1	1:A:4:ARG:HA	2.00	0.44
1:B:252:ASP:CG	1:B:255:LYS:HD2	2.43	0.44
1:C:260:THR:HB	1:C:319:LEU:O	2.18	0.44
1:D:101:PHE:HB3	1:D:104:LEU:HD12	2.00	0.43
1:C:200:ASN:HD22	1:C:203:TRP:HB2	1.82	0.43
1:D:187:VAL:HB	1:D:203:TRP:CE2	2.52	0.43
1:C:306:ASN:HB3	1:C:309:VAL:CG1	2.48	0.43
1:A:211:MET:HE1	1:A:245:PHE:CE1	2.53	0.43
1:C:267:LEU:O	1:C:271:LEU:HD12	2.19	0.43
1:A:185:ILE:HG21	1:A:212:ILE:HG23	2.01	0.43
1:C:263:GLU:HG2	1:C:263:GLU:H	1.71	0.43
1:C:162:GLY:O	1:C:164:LYS:CE	2.67	0.43
1:A:80:ARG:NH2	3:A:503:HOH:O	2.51	0.43
1:B:147:GLN:O	1:B:326:LYS:HE2	2.19	0.42
1:B:261:PRO:HG2	1:B:319:LEU:HD23	2.01	0.42
1:D:241:LYS:O	1:D:241:LYS:HG3	2.18	0.42
1:C:142:ILE:HG22	1:C:143:SER:O	2.19	0.42
1:C:265:GLN:HA	1:C:268:VAL:CG1	2.46	0.42
1:D:37:LYS:HE3	1:D:37:LYS:HB3	1.94	0.42
1:A:40:GLU:OE2	1:A:100:ARG:NH2	2.51	0.42
1:B:196:GLU:HG3	1:B:203:TRP:CZ2	2.55	0.42
1:C:315:VAL:HG21	2:C:401:AMP:C6	2.54	0.42
1:C:142:ILE:HD11	1:C:179:TYR:CD1	2.54	0.42
1:C:303:PRO:HD2	1:C:311:ARG:HH12	1.84	0.42
1:C:212:ILE:HG23	1:C:213:GLU:N	2.35	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:296:ILE:HD13	1:D:299:MET:HE3	2.02	0.42
1:A:313:GLY:O	1:A:314:LEU:HD23	2.19	0.41
1:B:127:SER:HB2	1:B:128:THR:H	1.53	0.41
1:C:200:ASN:HD22	1:C:203:TRP:H	1.66	0.41
1:C:4:ARG:HB3	1:C:6:LEU:CD2	2.50	0.41
1:C:247:LEU:HD11	1:C:249:ASP:O	2.20	0.41
1:C:304:SER:HB3	1:C:309:VAL:HG13	2.00	0.41
1:D:287:LEU:HD23	1:D:287:LEU:C	2.46	0.41
1:A:157:ILE:HD13	1:A:157:ILE:HA	1.93	0.41
1:B:260:THR:HB	1:B:319:LEU:O	2.21	0.41
1:B:46:ARG:HB3	1:B:98:MET:HG2	2.03	0.41
1:C:26:CYS:SG	1:C:33:THR:HG23	2.60	0.41
1:A:26:CYS:SG	1:A:33:THR:HG23	2.61	0.41
1:C:238:GLU:OE2	1:C:304:SER:OG	2.36	0.41
1:D:68:LYS:HD2	1:D:68:LYS:HA	1.87	0.41
1:D:260:THR:HG23	1:D:263:GLU:OE1	2.20	0.40
1:C:185:ILE:CG2	1:C:212:ILE:HG13	2.50	0.40
1:A:2:GLY:O	1:A:3:VAL:HB	2.22	0.40
1:A:217:PHE:CE2	1:C:72:ASP:HB2	2.57	0.40
1:B:67:SER:O	1:B:70:VAL:HG23	2.21	0.40
1:C:127:SER:HB2	1:C:130:LEU:HD13	2.04	0.40
1:D:186:GLY:HA3	1:D:193:GLU:OE1	2.22	0.40
1:D:211:MET:HE1	1:D:245:PHE:CZ	2.57	0.40
1:B:80:ARG:HD2	1:B:122:TRP:CE2	2.56	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	337/340 (99%)	325 (96%)	11 (3%)	1 (0%)	36 48

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	337/340 (99%)	318 (94%)	18 (5%)	1 (0%)	36	48
1	C	337/340 (99%)	321 (95%)	13 (4%)	3 (1%)	14	20
1	D	337/340 (99%)	319 (95%)	17 (5%)	1 (0%)	36	48
All	All	1348/1360 (99%)	1283 (95%)	59 (4%)	6 (0%)	30	40

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	130	LEU
1	C	196	GLU
1	C	337	LYS
1	D	3	VAL
1	A	3	VAL
1	C	3	VAL

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	292/293 (100%)	292 (100%)	0	100	100
1	B	292/293 (100%)	292 (100%)	0	100	100
1	C	292/293 (100%)	292 (100%)	0	100	100
1	D	292/293 (100%)	292 (100%)	0	100	100
All	All	1168/1172 (100%)	1168 (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 (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	92	GLN
1	A	107	ASN
1	A	129	ASN

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Mol	Chain	Res	Type
1	A	270	GLN
1	B	92	GLN
1	B	129	ASN
1	B	163	GLN
1	C	76	GLN
1	C	184	HIS
1	C	200	ASN
1	D	204	HIS
1	D	265	GLN
1	D	308	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

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	$\# Z > 2$
2	AMP	A	401	1	21,24,25	1.48	5 (23%)	30,35,38	2.18	8 (26%)
2	AMP	B	401	1	21,24,25	1.65	3 (14%)	30,35,38	2.63	8 (26%)
2	AMP	C	401	1	21,24,25	1.51	4 (19%)	30,35,38	2.20	7 (23%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	AMP	D	401	1	21,24,25	1.49	4 (19%)	30,35,38	2.37	9 (30%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	AMP	A	401	1	-	2/7/25/26	0/3/3/3
2	AMP	B	401	1	-	3/7/25/26	0/3/3/3
2	AMP	C	401	1	-	4/7/25/26	0/3/3/3
2	AMP	D	401	1	-	4/7/25/26	0/3/3/3

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	401	AMP	C5-C4	5.21	1.48	1.39
2	C	401	AMP	C5-C4	4.81	1.47	1.39
2	D	401	AMP	C5-C4	4.51	1.47	1.39
2	A	401	AMP	C5-C4	4.15	1.46	1.39
2	B	401	AMP	C5-C6	3.01	1.49	1.41
2	B	401	AMP	C5-N7	-2.86	1.33	1.39
2	D	401	AMP	C5-C6	2.86	1.48	1.41
2	C	401	AMP	C5-C6	2.74	1.48	1.41
2	A	401	AMP	C5-C6	2.62	1.48	1.41
2	A	401	AMP	C5-N7	-2.61	1.34	1.39
2	A	401	AMP	C8-N7	2.61	1.36	1.31
2	C	401	AMP	C8-N7	2.46	1.36	1.31
2	D	401	AMP	C8-N7	2.21	1.35	1.31
2	C	401	AMP	C5-N7	-2.09	1.35	1.39
2	D	401	AMP	C5-N7	-2.09	1.35	1.39
2	A	401	AMP	C4-N9	-2.01	1.33	1.37

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	AMP	C5-C4-N3	-8.25	115.36	126.72
2	D	401	AMP	C5-C4-N3	-7.22	116.78	126.72
2	C	401	AMP	C5-C4-N3	-6.82	117.33	126.72
2	B	401	AMP	N3-C4-N9	6.49	138.20	127.17
2	D	401	AMP	N3-C4-N9	5.59	136.68	127.17

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	AMP	C5-C4-N3	-5.54	119.09	126.72
2	C	401	AMP	N3-C4-N9	5.32	136.21	127.17
2	A	401	AMP	N3-C4-N9	4.98	135.63	127.17
2	C	401	AMP	C2-N3-C4	4.27	122.26	111.83
2	A	401	AMP	C4-N9-C8	4.19	110.14	105.74
2	D	401	AMP	C2-N3-C4	4.16	122.00	111.83
2	B	401	AMP	C4-C5-N7	-4.12	105.87	110.58
2	B	401	AMP	C2-N3-C4	4.03	121.68	111.83
2	B	401	AMP	C1'-N9-C8	-3.82	118.63	127.09
2	D	401	AMP	C4-C5-N7	-3.73	106.31	110.58
2	C	401	AMP	N3-C2-N1	-3.66	123.05	128.58
2	A	401	AMP	C2-N3-C4	3.57	120.55	111.83
2	B	401	AMP	C5-N7-C8	3.49	108.94	103.45
2	C	401	AMP	C4-C5-N7	-3.47	106.61	110.58
2	D	401	AMP	C1'-N9-C8	-3.41	119.52	127.09
2	A	401	AMP	N3-C2-N1	-3.27	123.63	128.58
2	A	401	AMP	N9-C8-N7	-3.25	109.33	113.94
2	B	401	AMP	C4-N9-C1'	3.10	133.88	126.63
2	D	401	AMP	N3-C2-N1	-2.89	124.21	128.58
2	A	401	AMP	C5-N7-C8	2.82	107.88	103.45
2	A	401	AMP	C4-C5-N7	-2.79	107.39	110.58
2	C	401	AMP	C5-N7-C8	2.70	107.69	103.45
2	D	401	AMP	C5-N7-C8	2.67	107.65	103.45
2	D	401	AMP	C4-N9-C1'	2.36	132.16	126.63
2	B	401	AMP	C6-C5-C4	2.24	120.24	117.18
2	D	401	AMP	C4-N9-C8	2.08	107.92	105.74
2	C	401	AMP	C6-C5-N7	2.01	135.96	132.09

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401	AMP	C4'-C5'-O5'-P
2	B	401	AMP	C4'-C5'-O5'-P
2	C	401	AMP	C4'-C5'-O5'-P
2	D	401	AMP	C4'-C5'-O5'-P
2	C	401	AMP	C2'-C1'-N9-C8
2	D	401	AMP	C2'-C1'-N9-C8
2	C	401	AMP	C3'-C4'-C5'-O5'
2	C	401	AMP	O4'-C1'-N9-C8
2	A	401	AMP	O4'-C1'-N9-C8
2	D	401	AMP	O4'-C1'-N9-C8

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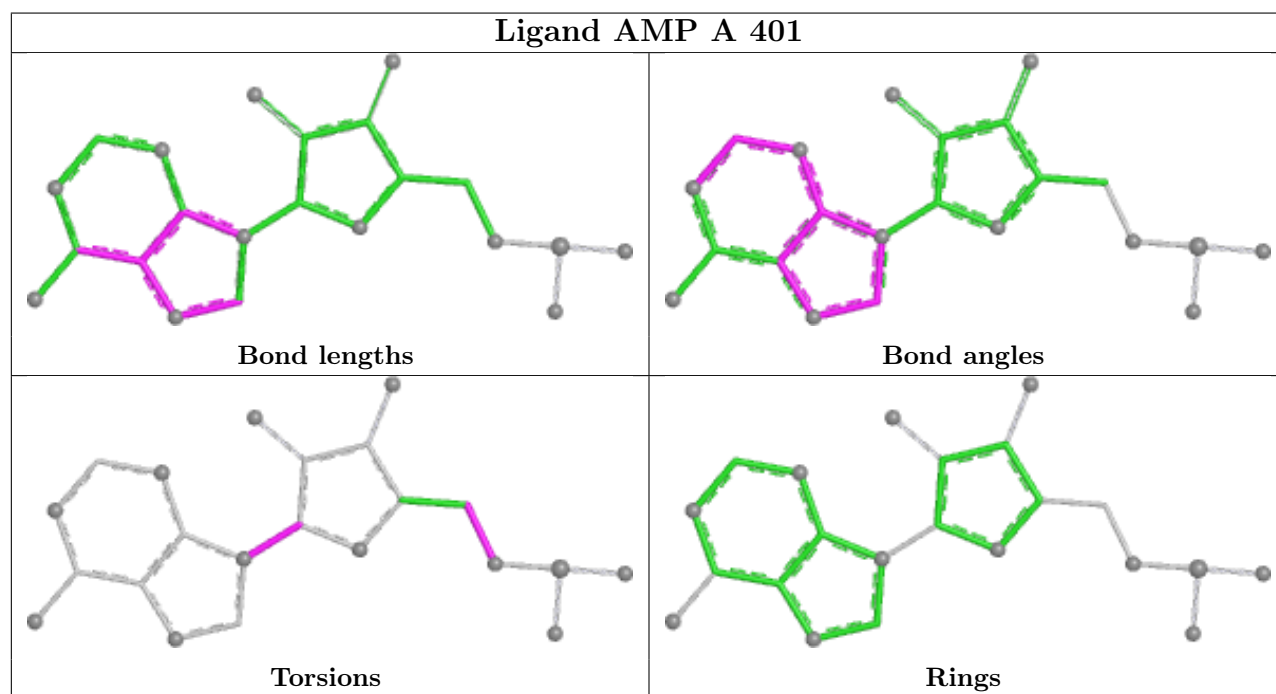
Mol	Chain	Res	Type	Atoms
2	B	401	AMP	C3'-C4'-C5'-O5'
2	B	401	AMP	O4'-C1'-N9-C8
2	D	401	AMP	C3'-C4'-C5'-O5'

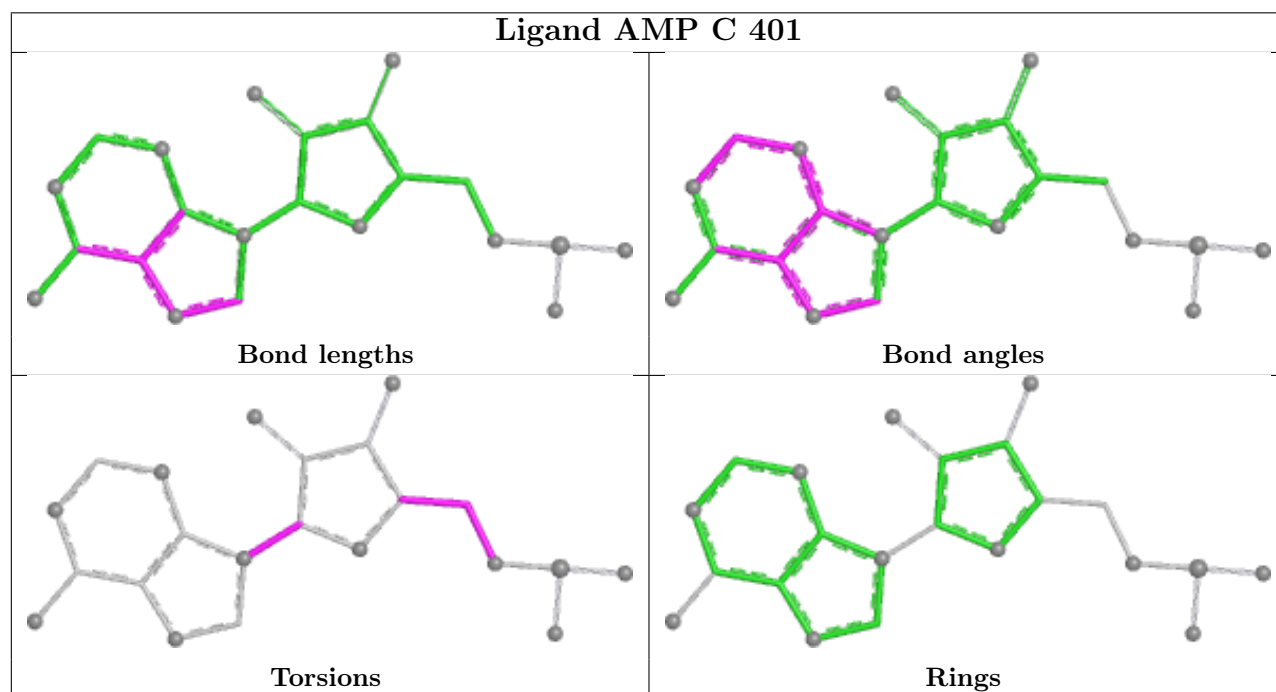
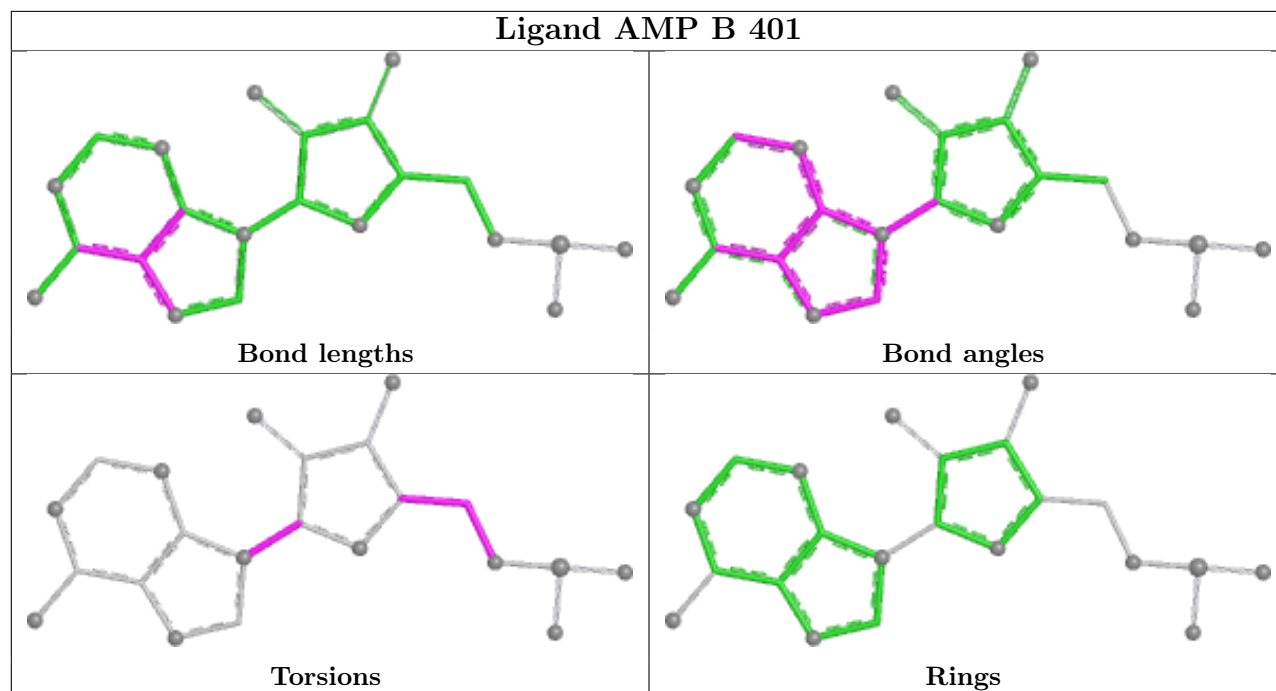
There are no ring outliers.

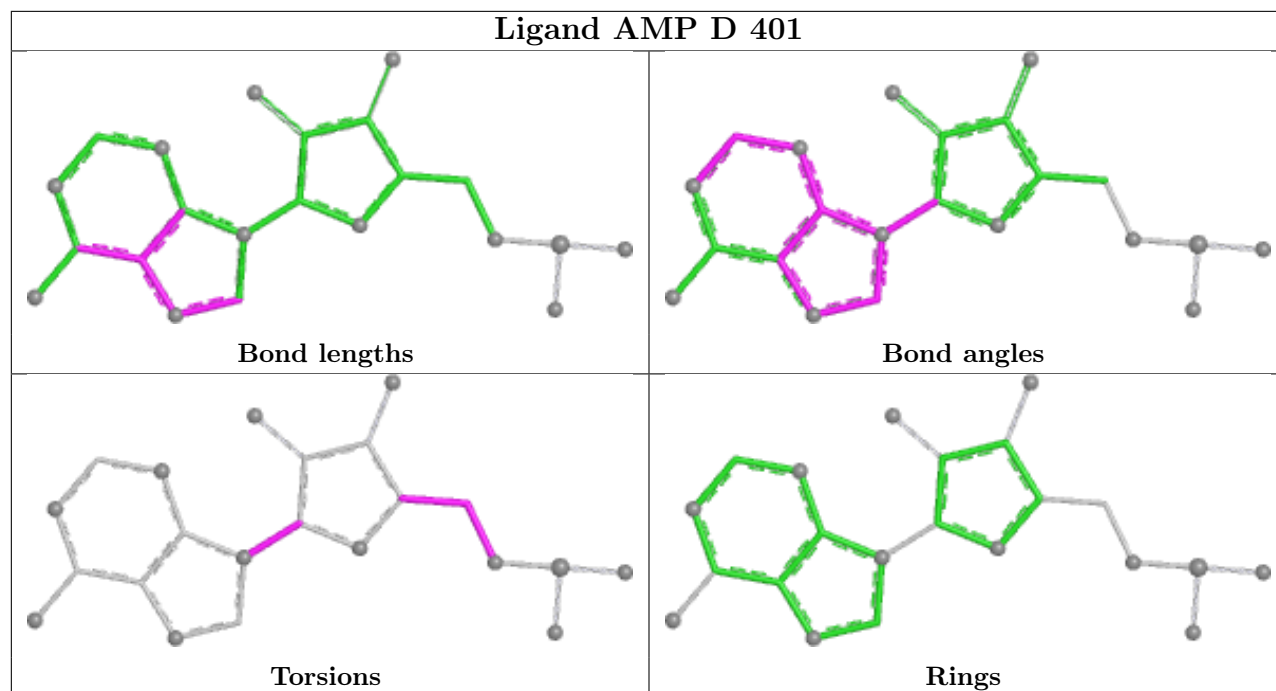
2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	401	AMP	1	0
2	C	401	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

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	339/340 (99%)	0.39	15 (4%) 39 38	29, 49, 71, 85	0
1	B	339/340 (99%)	0.71	28 (8%) 17 16	38, 59, 78, 100	0
1	C	339/340 (99%)	1.03	56 (16%) 4 4	33, 64, 87, 104	0
1	D	339/340 (99%)	0.65	28 (8%) 17 16	31, 53, 80, 90	0
All	All	1356/1360 (99%)	0.69	127 (9%) 14 13	29, 55, 81, 104	0

All (127) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	187	VAL	4.9
1	C	212	ILE	4.4
1	C	3	VAL	4.4
1	C	2	GLY	4.3
1	C	278	VAL	4.1
1	C	237	LEU	4.0
1	D	195	ARG	3.9
1	C	211	MET	3.8
1	C	217	PHE	3.7
1	C	185	ILE	3.7
1	B	128	THR	3.7
1	D	237	LEU	3.7
1	A	338	ASP	3.7
1	C	199	THR	3.7
1	D	2	GLY	3.7
1	C	198	ALA	3.7
1	B	1	MET	3.6
1	C	192	TRP	3.6
1	B	108	GLN	3.5
1	D	101	PHE	3.5
1	C	131	SER	3.5

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Mol	Chain	Res	Type	RSRZ
1	C	203	TRP	3.5
1	C	186	GLY	3.4
1	B	21	ALA	3.4
1	C	128	THR	3.4
1	C	224	LEU	3.4
1	D	339	ARG	3.2
1	A	339	ARG	3.2
1	C	215	LEU	3.1
1	D	130	LEU	3.1
1	B	126	ILE	3.1
1	C	178	VAL	3.1
1	C	276	PHE	3.0
1	D	198	ALA	3.0
1	A	127	SER	3.0
1	D	3	VAL	2.9
1	C	130	LEU	2.9
1	B	39	GLY	2.9
1	B	194	LEU	2.9
1	C	222	LEU	2.8
1	C	179	TYR	2.8
1	C	194	LEU	2.8
1	D	192	TRP	2.8
1	A	126	ILE	2.8
1	A	198	ALA	2.8
1	B	129	ASN	2.8
1	B	2	GLY	2.7
1	C	262	ILE	2.7
1	D	131	SER	2.7
1	C	138	PHE	2.7
1	C	338	ASP	2.7
1	D	194	LEU	2.7
1	C	196	GLU	2.6
1	B	183	ASP	2.6
1	D	303	PRO	2.6
1	C	126	ILE	2.6
1	D	204	HIS	2.6
1	C	218	LEU	2.6
1	C	339	ARG	2.6
1	C	141	PHE	2.6
1	D	199	THR	2.6
1	C	82	ARG	2.5
1	C	195	ARG	2.5

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Mol	Chain	Res	Type	RSRZ
1	C	183	ASP	2.5
1	B	303	PRO	2.5
1	C	209	ASN	2.5
1	A	3	VAL	2.5
1	D	1	MET	2.5
1	D	132	GLY	2.5
1	C	134	ILE	2.4
1	C	1	MET	2.4
1	D	100	ARG	2.4
1	C	105	ALA	2.4
1	C	250	ILE	2.4
1	B	86	LEU	2.4
1	D	135	LEU	2.4
1	B	196	GLU	2.4
1	B	3	VAL	2.4
1	C	240	LEU	2.4
1	D	205	ALA	2.3
1	D	233	ILE	2.3
1	A	2	GLY	2.3
1	C	189	GLY	2.3
1	C	140	THR	2.3
1	C	197	THR	2.3
1	B	46	ARG	2.3
1	D	296	ILE	2.3
1	B	136	GLY	2.3
1	C	275	GLY	2.3
1	D	133	GLU	2.3
1	B	240	LEU	2.2
1	C	259	LEU	2.2
1	D	187	VAL	2.3
1	B	22	GLU	2.2
1	A	130	LEU	2.2
1	D	48	VAL	2.2
1	C	223	ALA	2.2
1	C	281	VAL	2.2
1	A	87	ARG	2.2
1	B	308	ASN	2.2
1	D	208	ARG	2.2
1	A	125	PRO	2.2
1	B	127	SER	2.2
1	B	204	HIS	2.1
1	B	208	ARG	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	309	VAL	2.1
1	C	206	ALA	2.1
1	C	204	HIS	2.1
1	D	190	ARG	2.1
1	B	130	LEU	2.1
1	A	308	ASN	2.1
1	B	20	GLY	2.1
1	A	82	ARG	2.1
1	C	177	THR	2.0
1	A	158	GLU	2.0
1	C	201	ALA	2.0
1	C	205	ALA	2.0
1	A	240	LEU	2.0
1	B	209	ASN	2.0
1	C	271	LEU	2.0
1	D	338	ASP	2.0
1	C	245	PHE	2.0
1	A	269	LYS	2.0
1	B	293	ALA	2.0
1	D	105	ALA	2.0
1	B	290	ASN	2.0
1	B	134	ILE	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	AMP	D	401	22/23	0.90	0.10	34,47,63,67	0

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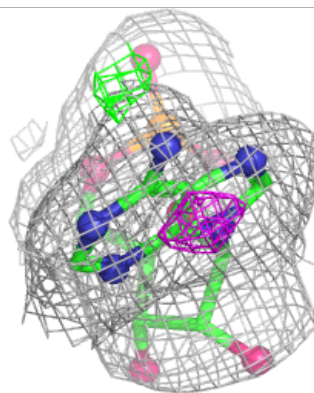
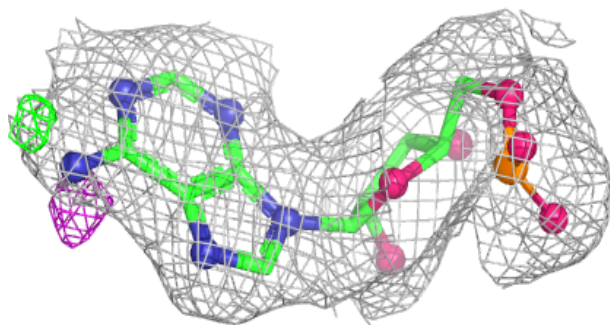
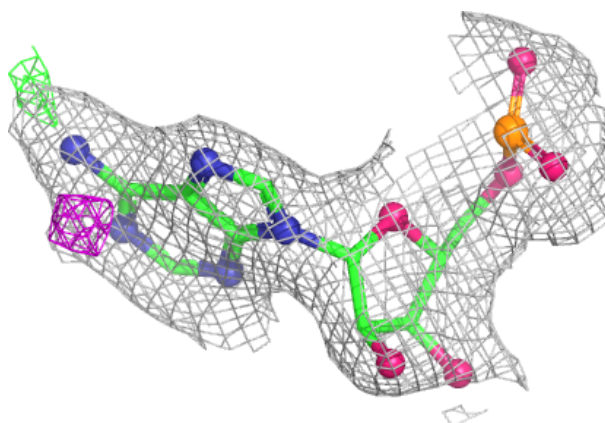
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	AMP	C	401	22/23	0.91	0.09	50,61,66,71	0
2	AMP	A	401	22/23	0.94	0.09	36,42,48,50	0
2	AMP	B	401	22/23	0.95	0.07	44,49,52,54	0

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.

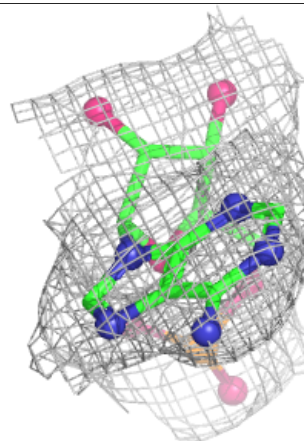
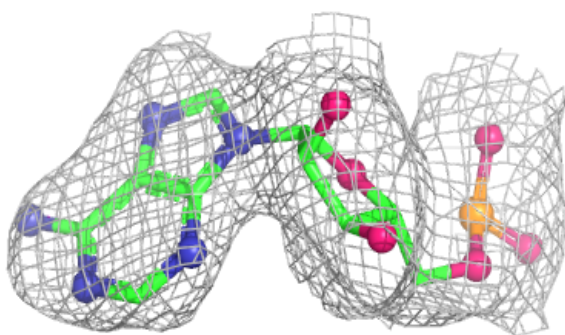
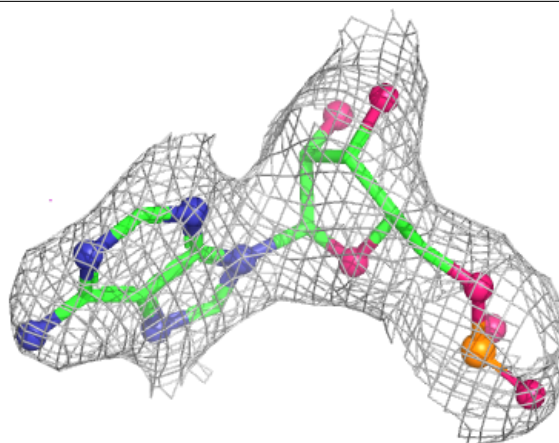
Electron density around AMP D 401:

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



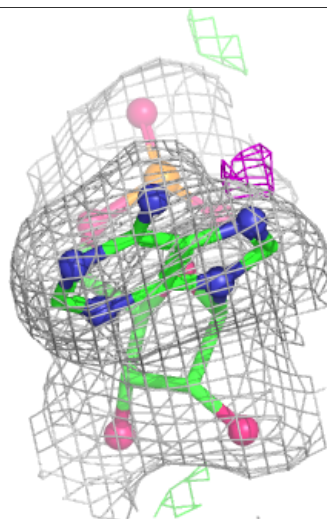
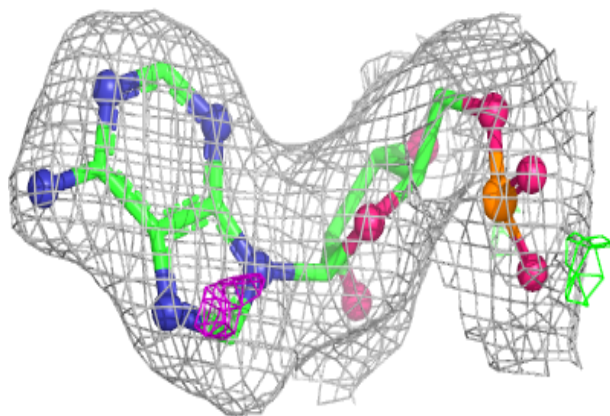
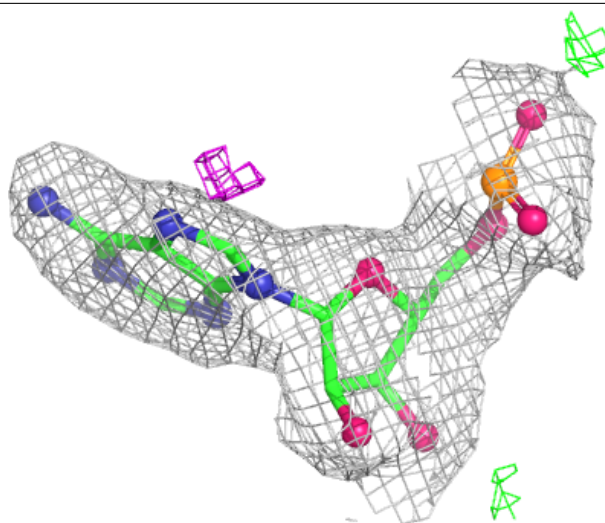
Electron density around AMP C 401:

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



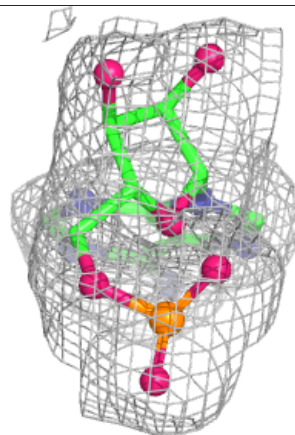
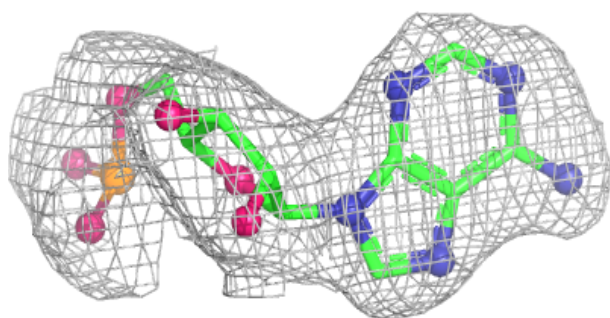
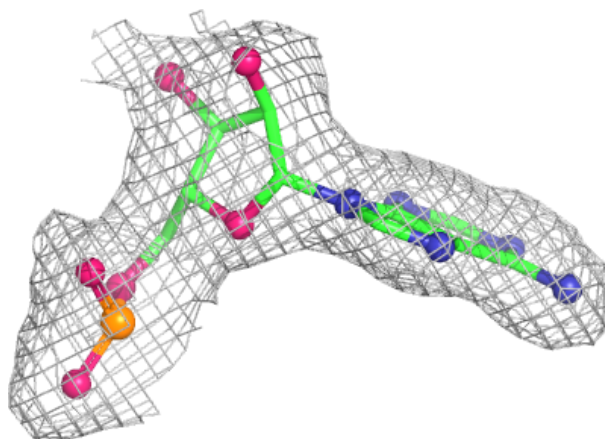
Electron density around AMP A 401:

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



Electron density around AMP B 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.