



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 23, 2026 – 07:25 PM JST

PDB ID : 9KPV / pdb_00009kpv
Title : Crystal structure of enzyme
Authors : Li, L.; Lin, L.Y.; Liu, C.; Hao, Z.Y.; Zhang, J.; Xiang, W.S.; Yuchi, Z.G.
Deposited on : 2024-11-24
Resolution : 2.51 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 : 1.8.5 (274361), CSD as541be (2020)
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

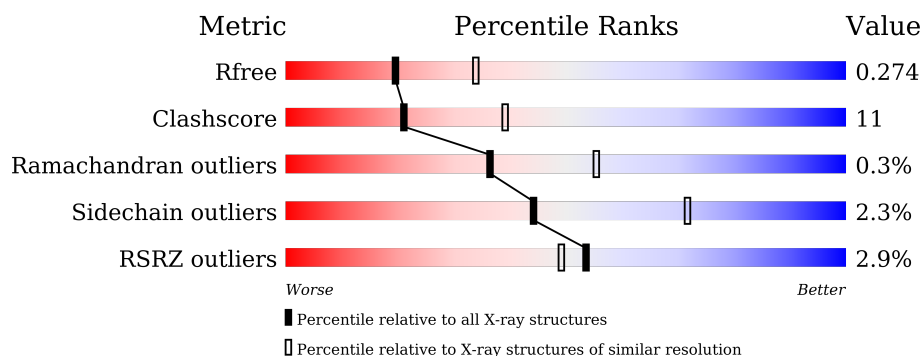
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.51 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	525	3% 71% 21% 7%
1	B	525	2% 67% 24% 8%
1	C	525	3% 67% 23% 8%
1	D	525	2% 71% 22% 6%

2 Entry composition [i](#)

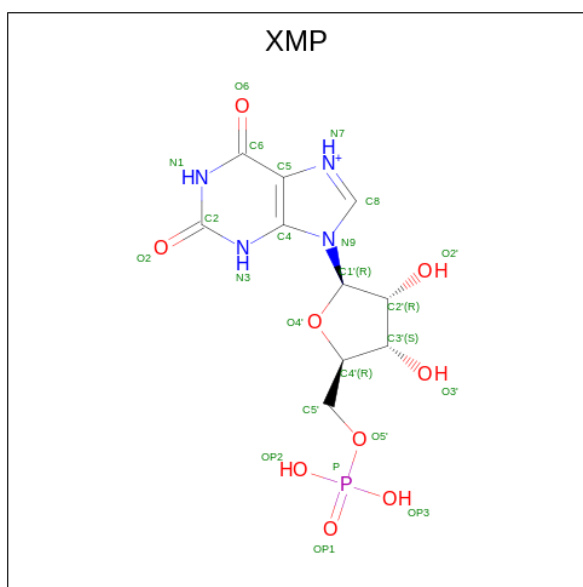
There are 4 unique types of molecules in this entry. The entry contains 15089 atoms, of which 48 are hydrogens and 0 are deuteriums.

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 GMP synthase [glutamine-hydrolyzing].

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	487	Total	C	N	O	S	0	0	0
			3702	2371	635	678	18			
1	B	481	Total	C	N	O	S	0	0	0
			3708	2379	636	673	20			
1	C	483	Total	C	N	O	S	0	0	0
			3684	2357	634	675	18			
1	D	494	Total	C	N	O	S	0	0	0
			3822	2445	659	698	20			

- Molecule 2 is XANTHOSINE-5'-MONOPHOSPHATE (CCD ID: XMP) (formula: $C_{10}H_{14}N_4O_9P$).



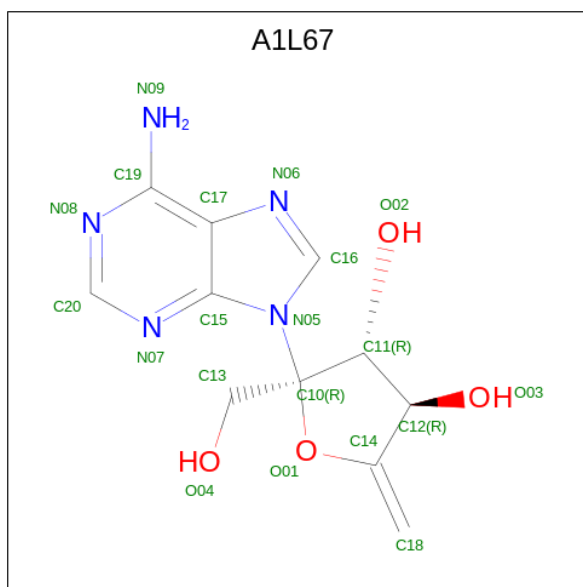
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	H	N	O	P	0
			36	10	12	4	9	1	
2	B	1	Total	C	H	N	O	P	0
			36	10	12	4	9	1	

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
2	C	1	Total	C	H	N	O	P	0	0
			36	10	12	4	9	1		
2	D	1	Total	C	H	N	O	P	0	0
			36	10	12	4	9	1		

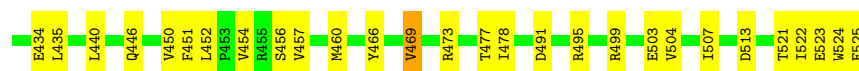
- Molecule 3 is (2 {R},3 {R},4 {R})-2-(6-aminopurin-9-yl)-2-(hydroxymethyl)-5-methylidene-oxolane-3,4-diol (CCD ID: A1L67) (formula: C₁₁H₁₃N₅O₄) (labeled as "Ligand of Interest" by depositor).



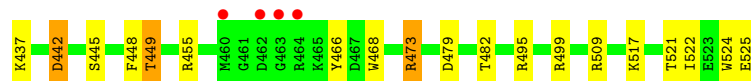
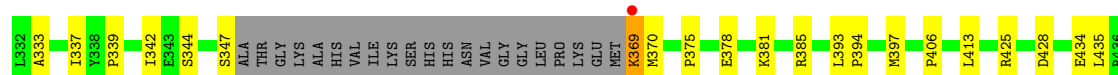
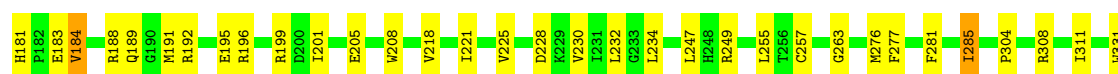
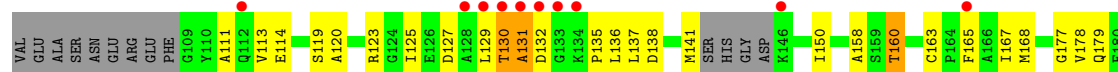
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	C	1	Total	C	N	O	0	0
			20	11	5	4		

- Molecule 4 is water.

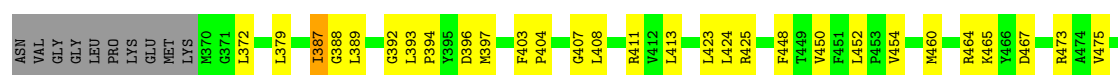
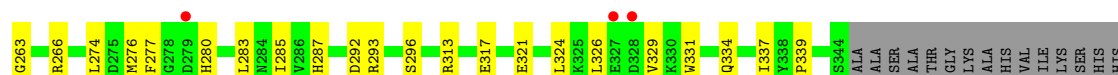
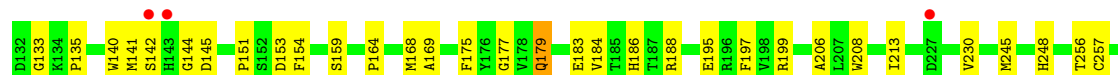
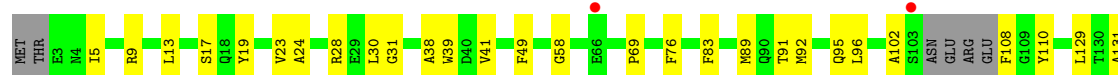
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	3	Total	O	0	0
			3	3		
4	B	3	Total	O	0	0
			3	3		
4	C	2	Total	O	0	0
			2	2		
4	D	1	Total	O	0	0
			1	1		



• Molecule 1: GMP synthase [glutamine-hydrolyzing]



• Molecule 1: GMP synthase [glutamine-hydrolyzing]



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	204.34Å 78.80Å 167.11Å 90.00° 123.68° 90.00°	Depositor
Resolution (Å)	48.94 – 2.51 48.94 – 2.51	Depositor EDS
% Data completeness (in resolution range)	92.2 (48.94-2.51) 92.6 (48.94-2.51)	Depositor EDS
R_{merge}	0.21	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.49 (at 2.51Å)	Xtriage
Refinement program	PHENIX (1.20.1_4487: ???)	Depositor
R, R_{free}	0.205 , 0.274 0.206 , 0.274	Depositor DCC
R_{free} test set	1986 reflections (2.60%)	wwPDB-VP
Wilson B-factor (Å ²)	37.8	Xtriage
Anisotropy	0.038	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 22.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	15089	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.19% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: XMP, A1L67

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	0.41	0/3779	0.57	0/5138
1	B	0.41	0/3786	0.55	0/5141
1	C	0.41	0/3762	0.58	0/5112
1	D	0.42	0/3905	0.57	0/5301
All	All	0.41	0/15232	0.57	0/20692

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	3702	0	3556	69	0
1	B	3708	0	3634	92	0
1	C	3684	0	3530	92	0
1	D	3822	0	3734	81	1
2	A	24	12	12	3	1
2	B	24	12	12	5	0
2	C	24	12	12	2	0
2	D	24	12	12	1	0
3	C	20	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	3	0	0	0	0
4	B	3	0	0	0	0
4	C	2	0	0	0	0
4	D	1	0	0	0	0
All	All	15041	48	14502	324	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 11.

All (324) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:263:GLY:O	1:D:425:ARG:NH2	2.06	0.89
1:C:92:MET:HE3	1:C:168:MET:HE3	1.64	0.79
1:D:89:MET:HG3	1:D:177:GLY:HA3	1.68	0.75
1:C:225:VAL:HG23	1:C:331:TRP:CD1	2.20	0.74
1:B:43:GLU:HG3	1:B:74:TYR:HB2	1.70	0.74
1:D:39:TRP:CE2	1:D:69:PRO:HB3	2.24	0.72
1:B:167:ILE:HG12	1:B:178:VAL:HG12	1.72	0.70
1:C:131:ALA:HB2	1:C:136:LEU:HD21	1.74	0.69
1:B:69:PRO:O	1:B:87:TYR:OH	2.07	0.69
1:C:167:ILE:HG12	1:C:178:VAL:HG12	1.75	0.69
1:C:369:LYS:HD2	1:C:370:MET:H	1.58	0.69
1:C:263:GLY:O	1:C:425:ARG:NH2	2.26	0.69
1:A:43:GLU:HG3	1:A:74:TYR:CG	2.29	0.68
1:C:129:LEU:O	1:C:136:LEU:HD23	1.94	0.67
1:C:232:LEU:HB2	1:C:333:ALA:HB3	1.76	0.67
1:D:230:VAL:HG22	1:D:331:TRP:HB2	1.75	0.66
1:C:69:PRO:O	1:C:87:TYR:OH	2.11	0.66
1:B:454:VAL:HG23	1:B:469:VAL:HG13	1.78	0.66
1:A:304:PRO:HB3	1:A:525:GLU:HB3	1.76	0.65
1:A:276:MET:HE2	1:A:394:PRO:HG2	1.79	0.65
1:C:218:VAL:HG21	1:C:249:ARG:HD3	1.80	0.64
1:C:199:ARG:HG3	1:C:199:ARG:HH11	1.63	0.63
1:C:413:LEU:HD11	1:C:524:TRP:HE3	1.63	0.63
1:D:248:HIS:HB2	1:D:283:LEU:HD21	1.80	0.63
1:A:230:VAL:HG22	1:A:331:TRP:HB2	1.79	0.63
1:C:428:ASP:OD1	1:C:449:THR:HG22	1.99	0.62
1:B:41:VAL:HG23	1:B:72:PRO:HG2	1.81	0.62
1:B:42:THR:H	1:B:45:GLN:HE21	1.47	0.62
1:B:118:ASP:OD2	1:B:123:ARG:NH1	2.31	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:89:MET:HG3	1:C:177:GLY:HA3	1.80	0.62
1:C:413:LEU:HD11	1:C:524:TRP:CE3	2.34	0.62
1:C:150:ILE:CB	1:C:165:PHE:HE2	2.12	0.62
1:B:450:VAL:HG21	1:B:524:TRP:CD1	2.34	0.62
1:C:381:LYS:HG3	1:C:385:ARG:HH11	1.65	0.62
1:C:230:VAL:HG22	1:C:331:TRP:HB2	1.80	0.61
1:C:276:MET:HE2	1:C:394:PRO:HG2	1.82	0.61
1:D:413:LEU:HD21	1:D:524:TRP:HE3	1.66	0.61
1:A:308:ARG:NH2	1:A:525:GLU:O	2.34	0.61
1:C:221:ILE:O	1:C:225:VAL:HG12	2.01	0.61
1:C:93:ALA:O	1:C:95:GLN:N	2.34	0.60
1:B:37:TRP:CH2	1:B:49:PHE:HB2	2.37	0.60
1:C:130:THR:C	1:C:132:ASP:H	2.09	0.59
1:B:413:LEU:HD11	1:B:524:TRP:CE3	2.38	0.59
1:D:9:ARG:HD3	1:D:49:PHE:CZ	2.37	0.59
1:A:467:ASP:HB2	1:A:508:SER:HB3	1.84	0.59
1:C:9:ARG:HD3	1:C:49:PHE:CE2	2.38	0.59
1:B:280:HIS:CE1	1:D:135:PRO:HG3	2.38	0.59
1:C:89:MET:HG3	1:C:177:GLY:CA	2.33	0.59
1:C:308:ARG:HH21	2:C:601:XMP:H7	1.51	0.58
1:C:522:ILE:HD12	2:C:601:XMP:H5'1	1.85	0.58
1:B:37:TRP:HH2	1:B:49:PHE:HB2	1.67	0.58
1:B:276:MET:O	1:B:280:HIS:HB3	2.04	0.57
1:B:504:VAL:HG11	1:B:507:ILE:HD12	1.86	0.57
1:C:455:ARG:NH1	1:C:466:TYR:O	2.37	0.57
1:D:102:ALA:HA	1:D:145:ASP:OD2	2.05	0.57
1:D:76:PHE:CE1	1:D:92:MET:HG3	2.39	0.57
1:A:413:LEU:HD11	1:A:524:TRP:HE3	1.69	0.56
1:B:9:ARG:HH21	1:B:50:ASN:HB3	1.71	0.56
1:D:393:LEU:HD13	1:D:397:MET:HE1	1.86	0.56
1:C:30:LEU:HD22	1:C:195:GLU:HG3	1.88	0.56
1:C:76:PHE:CE1	1:C:92:MET:HG3	2.41	0.56
1:C:111:ALA:HB3	1:C:141:MET:HG3	1.87	0.56
1:A:87:TYR:HA	1:A:90:GLN:HG3	1.87	0.56
1:A:413:LEU:HD11	1:A:524:TRP:CE3	2.41	0.56
1:C:9:ARG:HD3	1:C:49:PHE:CZ	2.41	0.56
1:D:83:PHE:HB2	1:D:197:PHE:CE1	2.41	0.55
1:B:277:PHE:HB2	1:B:285:ILE:HD11	1.87	0.55
1:B:393:LEU:HD13	1:B:397:MET:HE1	1.88	0.55
1:D:277:PHE:HB3	1:D:283:LEU:HD12	1.88	0.55
1:D:473:ARG:NH2	1:D:521:THR:O	2.39	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:408:LEU:HD22	1:D:424:LEU:HD23	1.88	0.55
1:D:337:ILE:HG13	1:D:339:PRO:HD2	1.89	0.54
1:B:395:TYR:CD1	1:D:188:ARG:HD3	2.43	0.54
1:B:460:MET:HA	1:B:460:MET:HE2	1.90	0.54
1:D:110:TYR:OH	1:D:186:HIS:ND1	2.32	0.54
1:B:522:ILE:HD12	2:B:601:XMP:H5'1	1.88	0.54
1:A:321:GLU:O	1:A:325:LYS:HG3	2.08	0.53
1:B:522:ILE:HD12	2:B:601:XMP:C5'	2.38	0.53
1:C:311:ILE:HG21	1:C:406:PRO:HA	1.90	0.53
1:C:434:GLU:OE1	1:C:495:ARG:HD2	2.08	0.53
1:C:448:PHE:CZ	1:C:473:ARG:HG3	2.43	0.53
1:C:455:ARG:HH21	1:C:468:TRP:NE1	2.07	0.53
1:C:113:VAL:HG22	1:C:137:LEU:HB2	1.91	0.53
1:D:17:SER:HB2	1:D:58:GLY:C	2.33	0.53
1:B:39:TRP:CE2	1:B:69:PRO:HB3	2.43	0.53
1:A:393:LEU:HB2	1:A:398:LEU:HD11	1.91	0.53
1:B:46:ILE:HD11	1:B:72:PRO:HG2	1.90	0.53
1:B:413:LEU:HD11	1:B:524:TRP:HE3	1.74	0.53
1:D:30:LEU:HD13	1:D:195:GLU:HA	1.91	0.53
1:C:225:VAL:HG21	1:C:230:VAL:CG2	2.39	0.52
1:A:22:LEU:O	1:A:26:ARG:HG2	2.09	0.52
1:B:8:HIS:CD2	1:B:205:GLU:HG3	2.43	0.52
1:A:265:LEU:O	1:A:425:ARG:HD3	2.10	0.52
1:B:184:VAL:HG12	1:B:186:HIS:H	1.74	0.52
1:C:381:LYS:HG3	1:C:385:ARG:NH1	2.25	0.52
1:C:234:LEU:HB2	1:C:257:CYS:HB3	1.91	0.52
1:A:110:TYR:CD2	1:A:140:TRP:HZ3	2.28	0.51
1:D:183:GLU:OE1	1:D:183:GLU:N	2.30	0.51
1:C:127:ASP:OD2	1:C:138:ASP:N	2.43	0.51
1:A:68:SER:OG	1:A:70:ARG:NH2	2.44	0.51
1:A:71:ALA:N	1:A:95:GLN:OE1	2.43	0.51
1:C:337:ILE:HG13	1:C:339:PRO:HD2	1.92	0.51
1:D:480:PHE:CE2	1:D:481:MET:HG3	2.46	0.51
1:A:308:ARG:HH21	1:A:525:GLU:C	2.18	0.51
1:C:304:PRO:HB3	1:C:525:GLU:HB3	1.92	0.51
1:D:140:TRP:CD1	1:D:184:VAL:HG21	2.46	0.51
1:A:115:VAL:HB	1:A:135:PRO:HB2	1.93	0.51
1:B:452:LEU:HB2	1:B:469:VAL:HG22	1.92	0.50
1:B:422:ASP:OD1	1:B:425:ARG:NH1	2.43	0.50
1:C:130:THR:C	1:C:132:ASP:N	2.69	0.50
1:D:89:MET:HG3	1:D:177:GLY:CA	2.40	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:276:MET:HE2	1:D:394:PRO:HG2	1.93	0.50
1:A:8:HIS:CE1	1:A:205:GLU:HG3	2.46	0.50
1:B:308:ARG:HH21	2:B:601:XMP:H7	1.59	0.50
1:A:308:ARG:NH2	1:A:410:VAL:HG21	2.27	0.50
1:B:317:GLU:O	1:B:321:GLU:HG3	2.12	0.50
1:C:442:ASP:OD1	1:C:442:ASP:N	2.44	0.50
1:A:154:PHE:CE2	1:A:168:MET:HE3	2.47	0.49
1:B:139:VAL:HA	1:B:187:THR:HG23	1.94	0.49
1:D:154:PHE:CD2	1:D:168:MET:HG3	2.47	0.49
1:B:499:ARG:O	1:B:503:GLU:HB2	2.12	0.49
1:C:129:LEU:HA	1:C:135:PRO:HA	1.93	0.49
1:C:160:THR:HG22	1:C:163:CYS:HB3	1.92	0.49
1:D:257:CYS:O	1:D:285:ILE:HA	2.12	0.49
1:A:458:GLY:O	1:A:464:ARG:HA	2.12	0.49
1:D:96:LEU:O	1:D:151:PRO:HB3	2.12	0.49
1:A:9:ARG:NH1	1:A:49:PHE:O	2.43	0.49
1:B:311:ILE:HG21	1:B:406:PRO:HA	1.95	0.49
1:D:292:ASP:O	1:D:296:SER:HB2	2.11	0.49
1:B:125:ILE:O	1:B:137:LEU:HD21	2.13	0.49
1:D:423:LEU:HD21	1:D:505:ASN:O	2.12	0.48
1:D:108:PHE:CE1	1:D:142:SER:HB3	2.48	0.48
1:A:184:VAL:HG12	1:A:186:HIS:H	1.78	0.48
1:A:247:LEU:HB2	1:A:255:LEU:HD22	1.95	0.48
1:B:42:THR:N	1:B:45:GLN:HE21	2.11	0.48
1:C:8:HIS:CE1	1:C:205:GLU:HG2	2.49	0.48
1:A:119:SER:O	1:A:123:ARG:HB2	2.13	0.48
1:B:37:TRP:CE3	1:B:41:VAL:HG11	2.48	0.48
1:B:434:GLU:OE1	1:B:495:ARG:NH1	2.46	0.48
1:D:387:ILE:HG13	1:D:388:GLY:N	2.28	0.48
1:B:274:LEU:HD11	1:B:287:HIS:CG	2.49	0.48
1:C:221:ILE:HG21	1:C:247:LEU:HD13	1.94	0.48
1:D:482:THR:HA	1:D:517:LYS:O	2.13	0.48
1:B:37:TRP:HE3	1:B:41:VAL:HG11	1.77	0.48
1:C:26:ARG:CD	1:C:191:MET:HE1	2.44	0.48
1:A:225:VAL:HG22	1:A:331:TRP:CD1	2.48	0.48
1:B:435:LEU:HD23	1:B:440:LEU:HB2	1.94	0.48
1:B:395:TYR:CG	1:D:188:ARG:HD3	2.49	0.48
1:C:120:ALA:O	1:C:196:ARG:NH1	2.44	0.48
1:B:15:PHE:HE2	1:B:56:LEU:HB3	1.79	0.47
1:A:5:ILE:HD12	1:A:208:TRP:CE3	2.48	0.47
1:A:158:ALA:HB3	1:A:167:ILE:HG13	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:234:LEU:HD12	1:A:236:GLY:H	1.78	0.47
1:D:266:ARG:HH21	1:D:403:PHE:CB	2.27	0.47
1:C:71:ALA:HB3	1:C:95:GLN:HE22	1.79	0.47
1:B:117:ASN:HB2	1:B:156:THR:O	2.13	0.47
1:B:466:TYR:CE2	1:C:304:PRO:HG2	2.50	0.47
1:C:26:ARG:NE	1:C:191:MET:HE1	2.30	0.47
1:B:153:ASP:OD1	1:B:153:ASP:N	2.44	0.47
1:D:293:ARG:NH2	1:D:321:GLU:OE2	2.43	0.47
1:A:9:ARG:HD3	1:A:49:PHE:CZ	2.49	0.47
1:B:208:TRP:CD1	1:B:213:ILE:HD11	2.50	0.47
1:D:396:ASP:OD1	1:D:397:MET:N	2.48	0.47
1:B:473:ARG:HD3	1:B:513:ASP:HB3	1.96	0.47
1:B:28:ARG:HB3	1:B:383:GLU:OE2	2.14	0.47
1:B:424:LEU:HD22	1:B:451:PHE:HB2	1.97	0.47
1:B:304:PRO:HB3	1:B:525:GLU:HB3	1.95	0.47
1:B:385:ARG:HD3	1:B:399:TYR:CE2	2.49	0.47
1:C:22:LEU:O	1:C:26:ARG:HG2	2.15	0.46
1:D:245:MET:HE3	1:D:245:MET:HA	1.97	0.46
1:A:30:LEU:HG	1:A:191:MET:HE1	1.97	0.46
1:B:135:PRO:HG3	1:D:280:HIS:NE2	2.31	0.46
1:B:331:TRP:CE3	1:B:371:GLY:HA3	2.51	0.46
1:D:169:ALA:HA	1:D:175:PHE:O	2.15	0.46
1:C:86:CYS:HG	1:C:181:HIS:CE1	2.33	0.46
1:A:39:TRP:CE2	1:A:69:PRO:HB3	2.50	0.46
1:B:24:ALA:O	1:B:28:ARG:HG2	2.16	0.46
1:B:39:TRP:CD2	1:B:69:PRO:HB3	2.51	0.46
1:B:58:GLY:HA2	1:B:86:CYS:HB3	1.97	0.46
1:D:413:LEU:HD21	1:D:524:TRP:CE3	2.48	0.46
1:A:53:GLY:HA3	1:A:202:CYS:SG	2.56	0.46
1:A:63:THR:HA	1:A:70:ARG:HH22	1.81	0.46
1:B:99:HIS:HB2	1:B:149:ALA:HB3	1.98	0.46
1:A:408:LEU:HD22	1:A:424:LEU:HD23	1.98	0.46
1:C:35:GLU:OE1	1:C:37:TRP:NE1	2.47	0.46
1:C:138:ASP:HB2	1:C:188:ARG:HB3	1.98	0.46
1:A:137:LEU:HD23	1:A:137:LEU:HA	1.74	0.46
1:A:151:PRO:HB2	1:A:154:PHE:CD1	2.51	0.45
1:A:125:ILE:O	1:A:137:LEU:HD21	2.17	0.45
1:D:38:ALA:O	1:D:41:VAL:HG13	2.15	0.45
1:D:39:TRP:CD2	1:D:69:PRO:HB3	2.51	0.45
1:D:452:LEU:O	1:D:454:VAL:N	2.43	0.45
1:A:395:TYR:O	1:A:398:LEU:N	2.49	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:405:GLY:H	2:A:601:XMP:C2	2.29	0.45
1:C:221:ILE:HD11	1:C:375:PRO:HB3	1.99	0.45
1:A:83:PHE:HB2	1:A:197:PHE:CE1	2.52	0.45
1:A:276:MET:O	1:A:280:HIS:HB3	2.17	0.45
1:D:452:LEU:C	1:D:454:VAL:H	2.23	0.45
1:B:9:ARG:HA	1:B:33:TYR:O	2.17	0.45
1:C:89:MET:SD	1:C:179:GLN:HG2	2.57	0.45
1:C:397:MET:HE3	1:C:397:MET:HB3	1.56	0.45
1:D:31:GLY:O	1:D:206:ALA:HA	2.17	0.45
1:B:15:PHE:CE2	1:B:56:LEU:HB3	2.52	0.44
1:B:295:LEU:HD13	1:B:418:LYS:HB2	1.99	0.44
1:C:247:LEU:HB3	1:C:255:LEU:HD22	1.99	0.44
1:D:188:ARG:HG3	1:D:188:ARG:HH11	1.81	0.44
1:A:238:VAL:O	1:A:242:VAL:HG23	2.17	0.44
1:A:418:LYS:HD3	1:A:422:ASP:OD2	2.17	0.44
1:A:391:LEU:HD23	1:A:391:LEU:HA	1.76	0.44
1:D:313:ARG:O	1:D:317:GLU:HG3	2.17	0.44
1:A:244:ALA:HB1	1:A:283:LEU:HD21	1.99	0.44
1:D:141:MET:HE2	1:D:179:GLN:HB3	1.99	0.44
1:A:337:ILE:HA	1:A:376:LEU:O	2.17	0.44
1:B:22:LEU:HD12	1:B:22:LEU:HA	1.76	0.44
1:B:153:ASP:OD2	1:B:170:ASN:ND2	2.51	0.44
1:B:119:SER:O	1:B:123:ARG:HB2	2.18	0.44
1:B:230:VAL:HG22	1:B:331:TRP:HB2	1.98	0.44
1:D:153:ASP:OD1	1:D:153:ASP:N	2.41	0.44
1:D:274:LEU:HD21	1:D:287:HIS:HB2	1.99	0.44
1:C:125:ILE:O	1:C:137:LEU:HD21	2.18	0.44
1:B:151:PRO:HB2	1:B:154:PHE:CD1	2.53	0.44
1:B:421:CYS:HB3	1:B:425:ARG:NH2	2.32	0.44
1:C:21:GLN:HG2	1:C:36:LEU:HD13	2.00	0.44
1:C:437:LYS:HB3	1:C:437:LYS:HE2	1.68	0.44
1:B:477:THR:O	1:B:478:ILE:HD13	2.18	0.43
1:A:522:ILE:HB	2:A:601:XMP:H5'1	1.99	0.43
1:B:220:ARG:HH12	1:B:378:GLU:CD	2.27	0.43
1:A:37:TRP:CH2	1:A:49:PHE:HB2	2.54	0.43
1:C:281:PHE:CE2	1:C:393:LEU:HD23	2.54	0.43
1:D:131:ALA:C	1:D:133:GLY:H	2.26	0.43
1:D:408:LEU:HD11	1:D:425:ARG:HG3	2.00	0.43
1:C:247:LEU:HD13	1:C:247:LEU:HA	1.86	0.43
1:A:333:ALA:HA	1:A:373:VAL:O	2.18	0.43
1:C:196:ARG:HG3	1:C:201:ILE:HD12	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:424:LEU:HD12	1:D:424:LEU:HA	1.72	0.43
1:C:183:GLU:HG2	1:C:184:VAL:N	2.34	0.43
1:D:450:VAL:HG21	1:D:524:TRP:CD1	2.53	0.43
1:A:54:ILE:HD13	1:A:75:VAL:HG22	1.99	0.43
1:B:310:ILE:O	1:B:314:VAL:HG23	2.19	0.43
1:D:256:THR:HG21	1:D:326:LEU:HD11	2.00	0.43
1:D:473:ARG:HD3	1:D:513:ASP:HB3	2.00	0.43
1:B:523:GLU:HG2	2:B:601:XMP:H2'	2.01	0.43
1:D:5:ILE:HD12	1:D:208:TRP:CE3	2.54	0.43
1:D:326:LEU:HB2	1:D:329:VAL:CG2	2.49	0.43
1:D:460:MET:HB2	1:D:465:LYS:HB3	2.01	0.43
1:C:13:LEU:HB2	1:C:56:LEU:HD23	2.01	0.43
1:D:324:LEU:HD23	1:D:324:LEU:HA	1.70	0.43
1:B:333:ALA:HA	1:B:373:VAL:O	2.19	0.42
1:B:521:THR:H	1:C:509:ARG:HH22	1.66	0.42
1:C:167:ILE:HA	1:C:177:GLY:O	2.19	0.42
1:A:424:LEU:HD11	1:A:449:THR:O	2.19	0.42
1:C:32:VAL:HG12	1:C:33:TYR:O	2.19	0.42
1:C:37:TRP:CH2	1:C:49:PHE:HB2	2.53	0.42
1:D:403:PHE:CE2	1:D:407:GLY:HA2	2.54	0.42
1:A:5:ILE:HD12	1:A:208:TRP:CZ3	2.54	0.42
1:D:334:GLN:OE1	1:D:372:LEU:HD11	2.18	0.42
1:A:111:ALA:O	1:A:139:VAL:HG22	2.20	0.42
1:A:225:VAL:HG13	1:A:228:ASP:HB2	2.02	0.42
1:C:76:PHE:HE2	1:C:95:GLN:NE2	2.17	0.42
1:C:199:ARG:HG3	1:C:199:ARG:NH1	2.33	0.42
1:A:86:CYS:O	1:A:89:MET:N	2.52	0.42
1:A:460:MET:HB2	1:A:460:MET:HE3	1.61	0.42
1:A:501:ILE:HG13	1:A:510:VAL:HG23	2.01	0.42
1:B:192:ARG:NH2	1:D:389:LEU:O	2.52	0.42
1:B:477:THR:C	1:B:478:ILE:HD13	2.45	0.42
1:B:313:ARG:HH11	1:B:317:GLU:CD	2.28	0.42
1:A:243:THR:OG1	1:A:335:GLY:HA2	2.19	0.42
1:D:372:LEU:HD12	1:D:372:LEU:HA	1.66	0.42
1:B:31:GLY:C	1:B:208:TRP:HB3	2.45	0.42
1:B:274:LEU:HD23	1:B:274:LEU:HA	1.82	0.42
1:B:276:MET:HE2	1:D:129:LEU:HD13	2.02	0.42
1:C:5:ILE:HD12	1:C:208:TRP:CE3	2.55	0.42
1:D:24:ALA:O	1:D:28:ARG:HG2	2.20	0.42
1:B:124:GLY:O	1:D:392:GLY:HA2	2.20	0.41
1:B:522:ILE:HB	2:B:601:XMP:H5'1	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:434:GLU:OE2	1:C:499:ARG:NH1	2.50	0.41
1:C:6:HIS:HE1	1:C:378:GLU:OE1	2.03	0.41
1:C:482:THR:HA	1:C:517:LYS:O	2.20	0.41
1:D:199:ARG:HD3	1:D:206:ALA:HB2	2.02	0.41
1:C:13:LEU:HD23	1:C:13:LEU:HA	1.87	0.41
1:A:517:LYS:HA	1:A:520:ALA:O	2.19	0.41
1:D:213:ILE:HG21	1:D:379:LEU:HD21	2.03	0.41
1:D:287:HIS:CD2	1:D:287:HIS:C	2.97	0.41
1:B:313:ARG:O	1:B:317:GLU:HG3	2.20	0.41
1:D:475:VAL:HB	1:D:483:ALA:HB1	2.02	0.41
1:B:172:GLU:HA	1:B:172:GLU:OE2	2.21	0.41
1:D:91:THR:OG1	1:D:95:GLN:OE1	2.30	0.41
1:B:194:LEU:HD23	1:B:194:LEU:HA	1.82	0.41
1:C:339:PRO:HA	1:C:342:ILE:HD12	2.02	0.41
1:D:266:ARG:HH21	1:D:403:PHE:HB2	1.85	0.41
1:B:402:PRO:HD2	1:B:446:GLN:HG3	2.01	0.41
1:C:141:MET:HE2	1:C:141:MET:HB3	1.89	0.41
1:D:404:PRO:HA	2:D:601:XMP:N7	2.35	0.41
1:A:308:ARG:HH12	2:A:601:XMP:H7	1.68	0.41
1:A:387:ILE:HG13	1:A:388:GLY:N	2.35	0.41
1:B:8:HIS:CD2	1:B:205:GLU:H	2.39	0.41
1:B:210:PRO:HA	1:B:387:ILE:HG22	2.03	0.41
1:C:119:SER:O	1:C:123:ARG:HB3	2.20	0.41
1:C:129:LEU:C	1:C:136:LEU:HD23	2.45	0.41
1:D:19:TYR:O	1:D:23:VAL:HG23	2.20	0.41
1:A:18:GLN:HE21	1:A:18:GLN:HB3	1.66	0.41
1:C:17:SER:HB2	1:C:58:GLY:C	2.46	0.41
1:C:435:LEU:HD12	1:C:435:LEU:HA	1.93	0.41
1:C:517:LYS:HG3	1:C:521:THR:HA	2.03	0.41
1:B:265:LEU:HD13	1:B:269:GLU:HG2	2.02	0.40
1:B:304:PRO:HG2	1:C:466:TYR:CE2	2.56	0.40
1:C:114:GLU:O	1:C:158:ALA:HA	2.21	0.40
1:A:31:GLY:C	1:A:208:TRP:HB3	2.45	0.40
1:D:423:LEU:HA	1:D:423:LEU:HD23	1.81	0.40
1:A:267:LEU:HA	1:A:425:ARG:HD2	2.03	0.40
1:A:499:ARG:O	1:A:503:GLU:HB2	2.21	0.40
1:B:19:TYR:O	1:B:23:VAL:HG23	2.21	0.40
1:B:53:GLY:HA3	1:B:202:CYS:SG	2.62	0.40
1:B:457:VAL:HB	1:B:466:TYR:CE2	2.57	0.40
1:C:277:PHE:HB2	1:C:285:ILE:HD11	2.02	0.40
1:A:423:LEU:HD23	1:A:426:ARG:NH1	2.36	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:231:ILE:HG12	1:B:232:LEU:N	2.37	0.40
1:C:189:GLN:OE1	1:C:192:ARG:HD3	2.20	0.40
1:C:479:ASP:OD1	1:C:482:THR:N	2.54	0.40
1:D:159:SER:HB3	1:D:164:PRO:O	2.20	0.40
1:D:411:ARG:HD3	1:D:448:PHE:CD2	2.56	0.40
1:A:63:THR:HG23	1:A:70:ARG:NH1	2.37	0.40
1:D:9:ARG:HD3	1:D:49:PHE:CE2	2.57	0.40
1:D:230:VAL:HG22	1:D:331:TRP:CB	2.46	0.40

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

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:464:ARG:NH2	2:A:601:XMP:OP1[2_556]	1.46	0.74

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	477/525 (91%)	453 (95%)	23 (5%)	1 (0%)	43	63
1	B	471/525 (90%)	455 (97%)	16 (3%)	0	100	100
1	C	473/525 (90%)	438 (93%)	32 (7%)	3 (1%)	21	38
1	D	488/525 (93%)	470 (96%)	17 (4%)	1 (0%)	43	63
All	All	1909/2100 (91%)	1816 (95%)	88 (5%)	5 (0%)	36	55

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	144	GLY
1	C	94	MET

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Mol	Chain	Res	Type
1	C	130	THR
1	C	131	ALA
1	A	478	ILE

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	372/442 (84%)	367 (99%)	5 (1%)	61	82
1	B	382/442 (86%)	370 (97%)	12 (3%)	35	62
1	C	369/442 (84%)	355 (96%)	14 (4%)	29	56
1	D	395/442 (89%)	391 (99%)	4 (1%)	68	86
All	All	1518/1768 (86%)	1483 (98%)	35 (2%)	44	72

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	75	VAL
1	A	159	SER
1	A	185	THR
1	A	326	LEU
1	A	415	GLU
1	B	41	VAL
1	B	68	SER
1	B	75	VAL
1	B	91	THR
1	B	99	HIS
1	B	205	GLU
1	B	247	LEU
1	B	316	VAL
1	B	373	VAL
1	B	456	SER
1	B	469	VAL
1	B	491	ASP
1	C	7	LYS

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Mol	Chain	Res	Type
1	C	22	LEU
1	C	65	GLU
1	C	160	THR
1	C	184	VAL
1	C	228	ASP
1	C	285	ILE
1	C	344	SER
1	C	347	SER
1	C	369	LYS
1	C	442	ASP
1	C	445	SER
1	C	449	THR
1	C	473	ARG
1	D	13	LEU
1	D	179	GLN
1	D	387	ILE
1	D	467	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	18	GLN
1	A	248	HIS
1	A	287	HIS
1	B	8	HIS
1	B	45	GLN
1	B	112	GLN
1	B	186	HIS
1	C	6	HIS
1	C	73	GLN
1	C	170	ASN
1	C	224	GLN
1	C	284	ASN
1	C	446	GLN
1	D	50	ASN
1	D	90	GLN
1	D	287	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

5 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	XMP	A	601	-	26,26,26	0.67	0	35,40,40	0.99	3 (8%)
2	XMP	C	601	-	26,26,26	0.84	1 (3%)	35,40,40	1.05	2 (5%)
3	A1L67	C	602	-	16,22,22	5.14	8 (50%)	14,34,34	6.85	11 (78%)
2	XMP	D	601	-	26,26,26	0.71	1 (3%)	35,40,40	1.16	3 (8%)
2	XMP	B	601	-	26,26,26	0.75	0	35,40,40	1.07	3 (8%)

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	XMP	A	601	-	-	5/10/26/26	0/3/3/3
2	XMP	C	601	-	-	0/10/26/26	0/3/3/3
3	A1L67	C	602	-	-	0/3/28/28	0/3/3/3
2	XMP	D	601	-	-	2/10/26/26	0/3/3/3
2	XMP	B	601	-	-	4/10/26/26	0/3/3/3

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	602	A1L67	C17-C15	13.46	1.64	1.39
3	C	602	A1L67	C15-N05	11.66	1.51	1.38
3	C	602	A1L67	C17-N06	6.68	1.51	1.39
3	C	602	A1L67	C19-N09	4.44	1.45	1.34
3	C	602	A1L67	C20-N08	-3.83	1.26	1.33
3	C	602	A1L67	C15-N07	3.05	1.40	1.34
3	C	602	A1L67	C19-N08	-2.85	1.27	1.35
2	D	601	XMP	C4-N3	-2.27	1.33	1.37
3	C	602	A1L67	C18-C14	2.21	1.34	1.31
2	C	601	XMP	C4-N3	-2.13	1.33	1.37

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	602	A1L67	N07-C15-N05	14.90	145.00	126.58
3	C	602	A1L67	C17-C15-N07	-11.26	112.06	126.75
3	C	602	A1L67	C19-C17-N06	7.93	146.81	132.02
3	C	602	A1L67	N08-C20-N07	-7.71	116.53	128.60
3	C	602	A1L67	C20-N07-C15	6.74	127.66	111.75
3	C	602	A1L67	C19-C17-C15	-6.11	108.97	117.18
3	C	602	A1L67	C17-C19-N08	5.43	129.04	117.39
3	C	602	A1L67	C15-C17-N06	-5.27	104.20	110.62
3	C	602	A1L67	C17-C19-N09	-4.39	113.88	123.43
3	C	602	A1L67	C20-N08-C19	4.06	125.73	118.77
2	C	601	XMP	C8-N7-C5	3.04	109.74	104.24
2	B	601	XMP	C8-N7-C5	2.88	109.46	104.24
3	C	602	A1L67	O01-C14-C12	-2.76	102.19	107.97
2	A	601	XMP	C8-N7-C5	2.73	109.18	104.24
2	C	601	XMP	N9-C8-N7	-2.63	108.43	113.39
2	D	601	XMP	O3'-C3'-C4'	-2.63	103.44	111.05
2	B	601	XMP	N9-C8-N7	-2.55	108.59	113.39
2	D	601	XMP	C8-N7-C5	2.53	108.81	104.24
2	A	601	XMP	N9-C8-N7	-2.47	108.74	113.39
2	D	601	XMP	N9-C8-N7	-2.40	108.87	113.39
2	A	601	XMP	C8-N9-C4	2.14	108.49	106.71
2	B	601	XMP	C8-N9-C4	2.07	108.43	106.71

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	601	XMP	C5'-O5'-P-OP3

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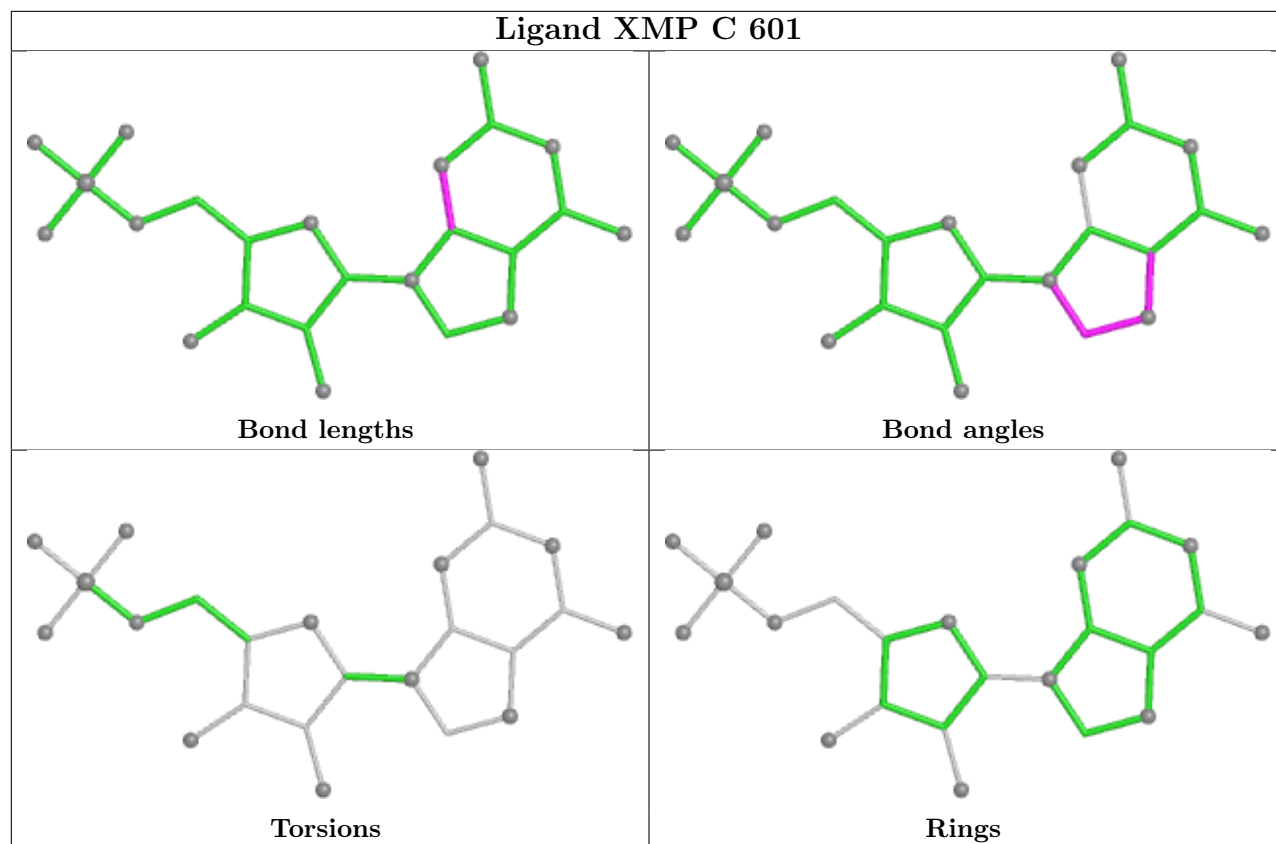
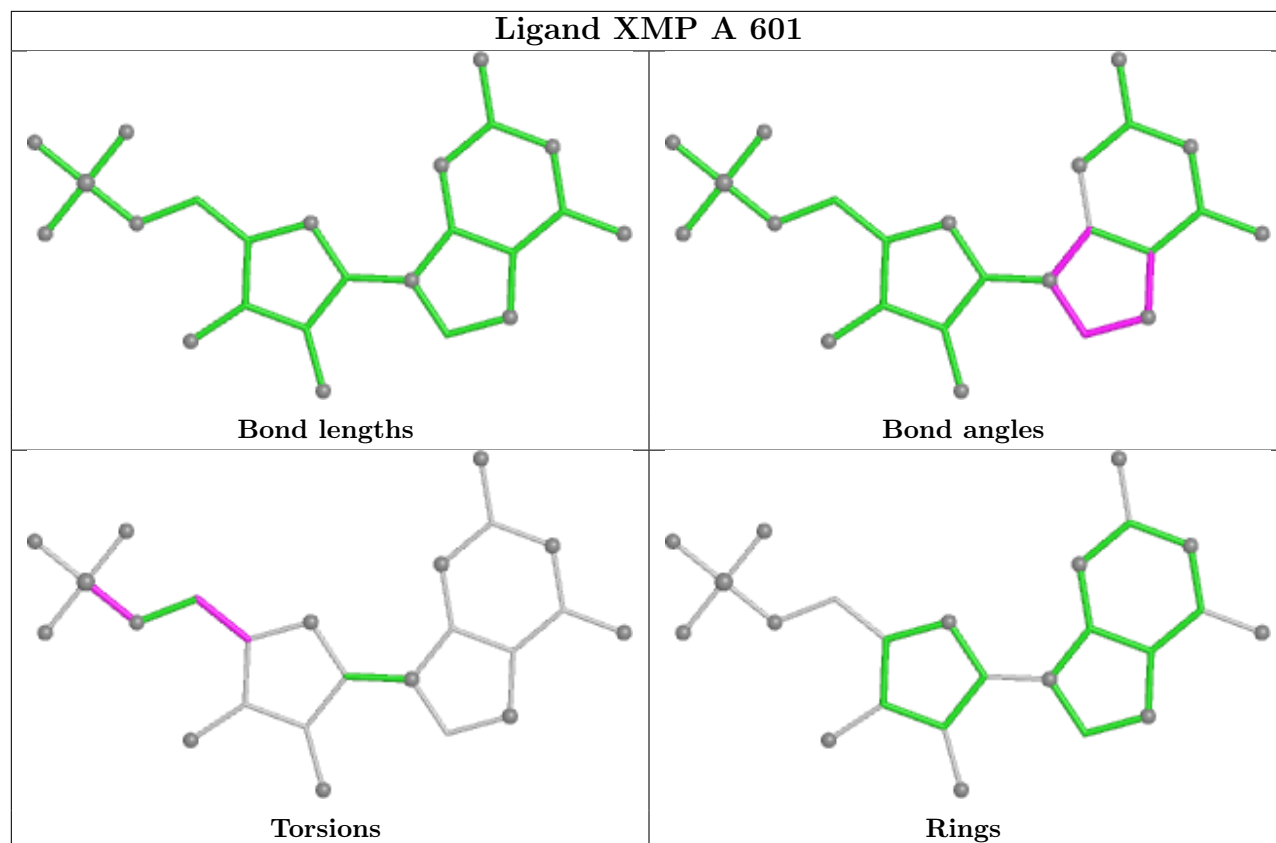
Mol	Chain	Res	Type	Atoms
2	A	601	XMP	C5'-O5'-P-OP2
2	B	601	XMP	C5'-O5'-P-OP1
2	B	601	XMP	O4'-C4'-C5'-O5'
2	B	601	XMP	C3'-C4'-C5'-O5'
2	A	601	XMP	O4'-C4'-C5'-O5'
2	D	601	XMP	O4'-C4'-C5'-O5'
2	D	601	XMP	C3'-C4'-C5'-O5'
2	A	601	XMP	C3'-C4'-C5'-O5'
2	A	601	XMP	C5'-O5'-P-OP1
2	B	601	XMP	C5'-O5'-P-OP2

There are no ring outliers.

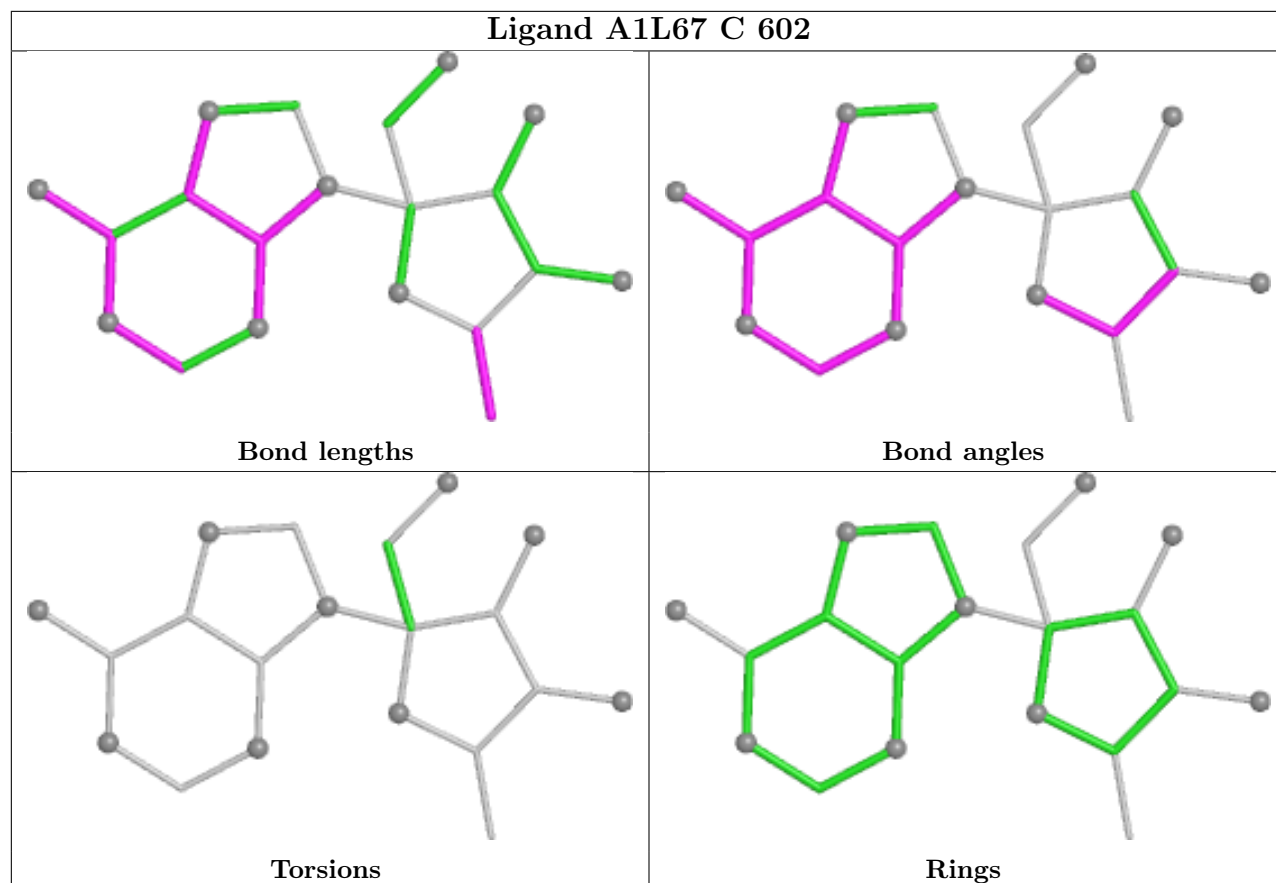
4 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601	XMP	3	1
2	C	601	XMP	2	0
2	D	601	XMP	1	0
2	B	601	XMP	5	0

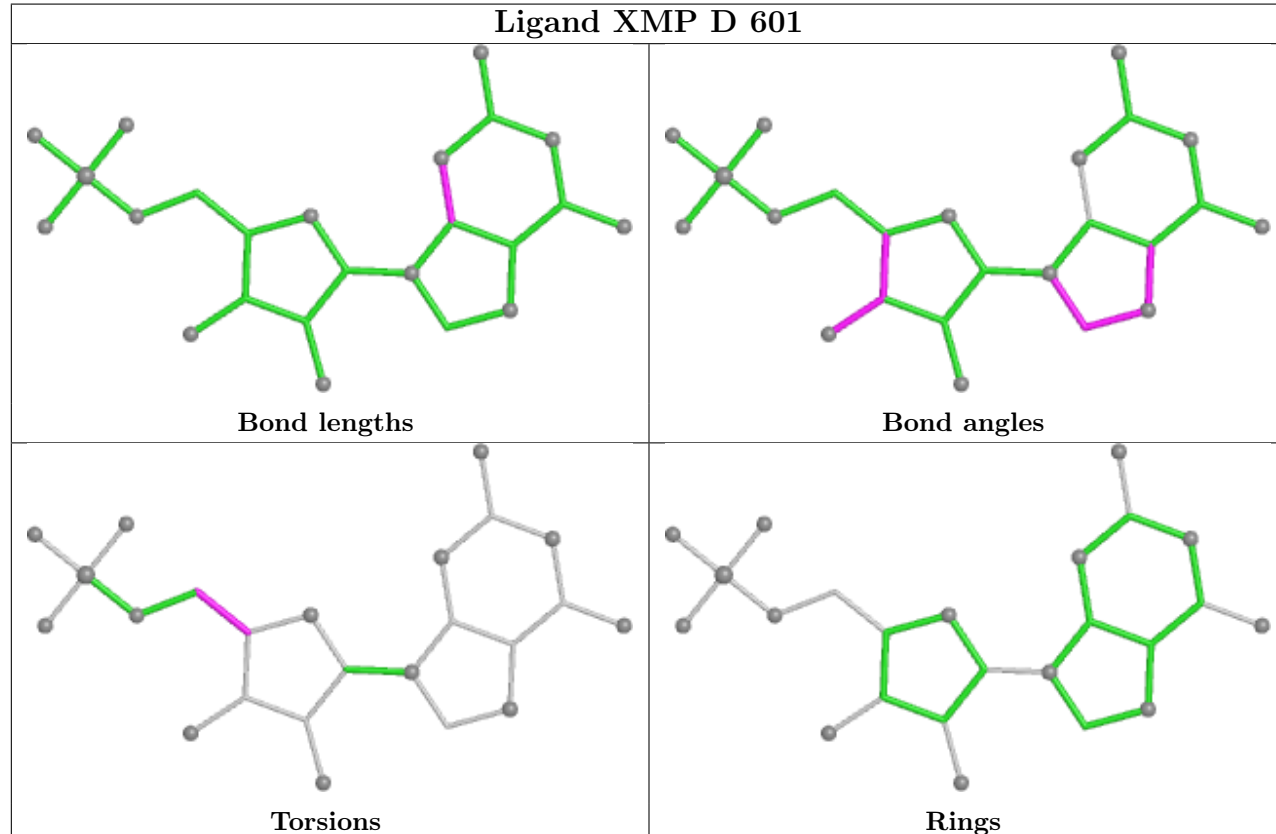
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

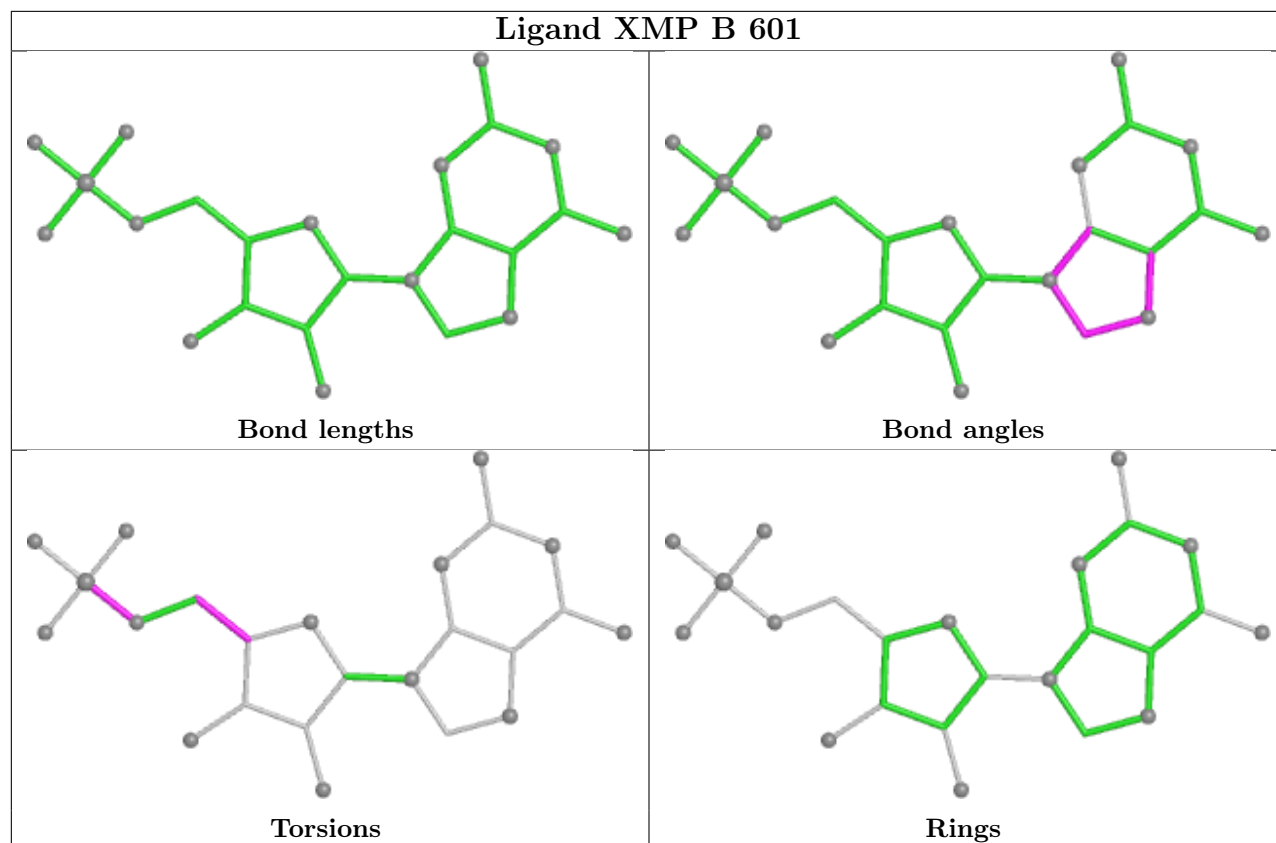


Ligand A1L67 C 602



Ligand XMP D 601





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	487/525 (92%)	0.12	18 (3%) 45 40	21, 36, 55, 75	0
1	B	481/525 (91%)	0.02	11 (2%) 61 57	18, 34, 53, 70	0
1	C	483/525 (92%)	0.17	18 (3%) 45 40	18, 37, 64, 80	0
1	D	494/525 (94%)	-0.07	9 (1%) 67 64	18, 33, 50, 63	0
All	All	1945/2100 (92%)	0.06	56 (2%) 53 49	18, 34, 56, 80	0

All (56) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	129	LEU	5.6
1	C	462	ASP	4.9
1	B	102	ALA	4.9
1	C	130	THR	4.8
1	C	464	ARG	4.5
1	C	63	THR	4.1
1	A	104	ASN	4.0
1	A	142	SER	3.6
1	A	108	PHE	3.5
1	A	327	GLU	3.5
1	D	328	ASP	3.3
1	B	346	ALA	3.3
1	C	133	GLY	3.1
1	C	131	ALA	3.1
1	A	103	SER	3.1
1	A	328	ASP	3.1
1	B	345	ALA	3.0
1	C	94	MET	2.8
1	C	460	MET	2.8
1	A	326	LEU	2.7
1	C	165	PHE	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	99	HIS	2.7
1	C	463	GLY	2.6
1	A	67	ASN	2.6
1	C	112	GLN	2.6
1	C	128	ALA	2.6
1	A	64	THR	2.5
1	A	370	MET	2.5
1	B	93	ALA	2.5
1	C	62	SER	2.5
1	B	109	GLY	2.4
1	A	57	SER	2.4
1	C	369	LYS	2.4
1	D	279	ASP	2.4
1	B	200	ASP	2.4
1	A	66	GLU	2.4
1	C	134	LYS	2.4
1	C	146	LYS	2.4
1	D	103	SER	2.4
1	C	132	ASP	2.4
1	D	142	SER	2.3
1	B	94	MET	2.3
1	A	330	LYS	2.3
1	A	459	VAL	2.3
1	A	160	THR	2.3
1	A	217	ALA	2.3
1	A	289	PRO	2.3
1	B	68	SER	2.3
1	D	143	HIS	2.2
1	A	77	GLU	2.2
1	D	327	GLU	2.2
1	D	66	GLU	2.2
1	D	227	ASP	2.1
1	B	19	TYR	2.1
1	B	161	GLU	2.1
1	D	480	PHE	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

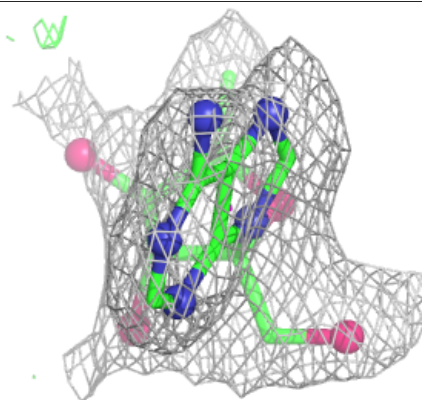
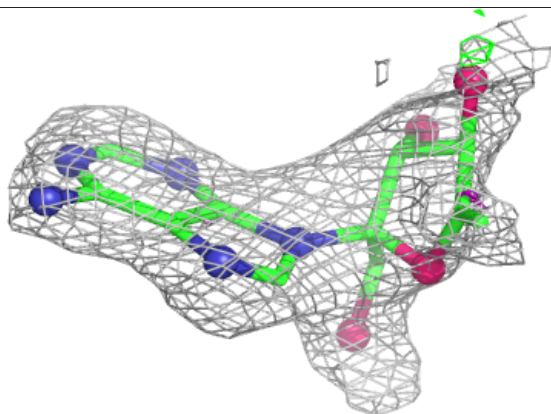
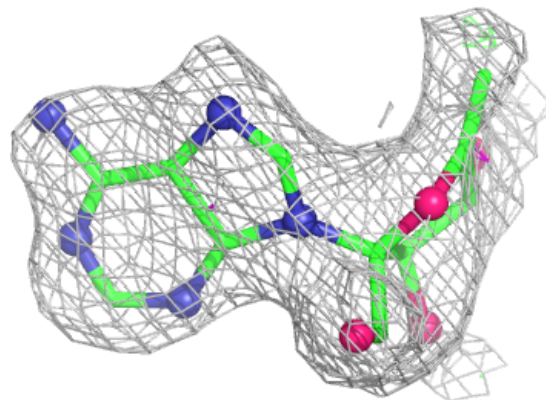
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	A1L67	C	602	20/20	0.78	0.14	33,47,61,62	0
2	XMP	D	601	24/24	0.81	0.13	20,20,20,20	0
2	XMP	A	601	24/24	0.91	0.09	20,20,20,20	0
2	XMP	B	601	24/24	0.93	0.08	20,20,20,20	0
2	XMP	C	601	24/24	0.94	0.07	20,20,20,20	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.

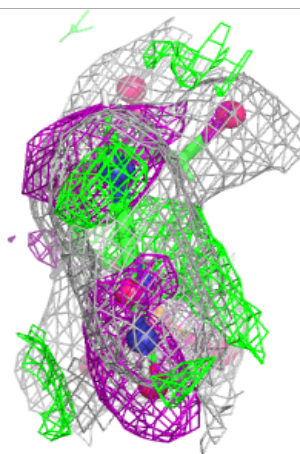
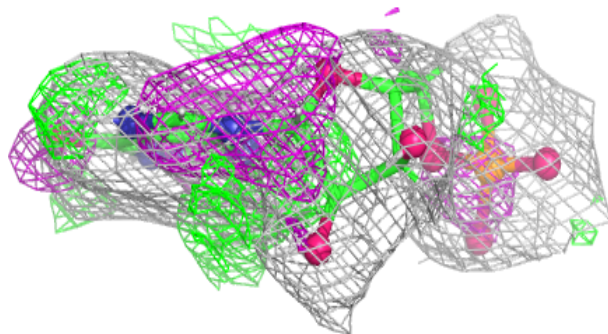
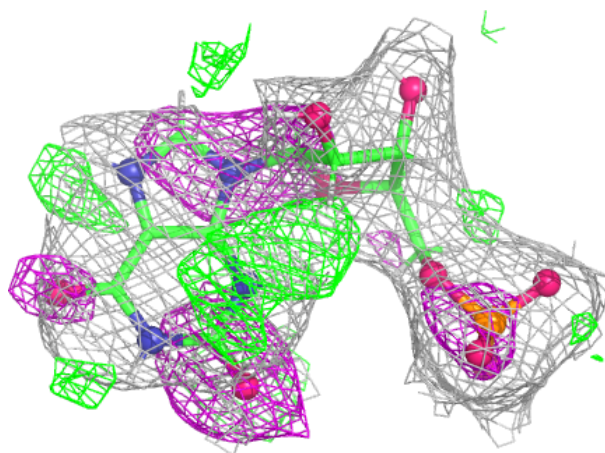
Electron density around A1L67 C 602:

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



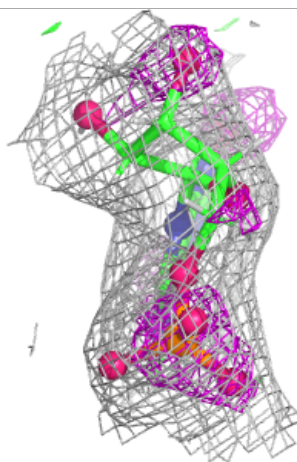
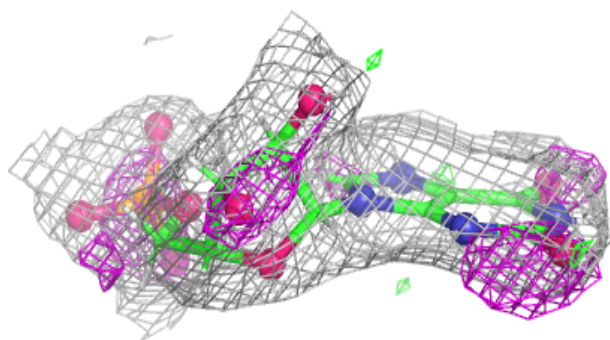
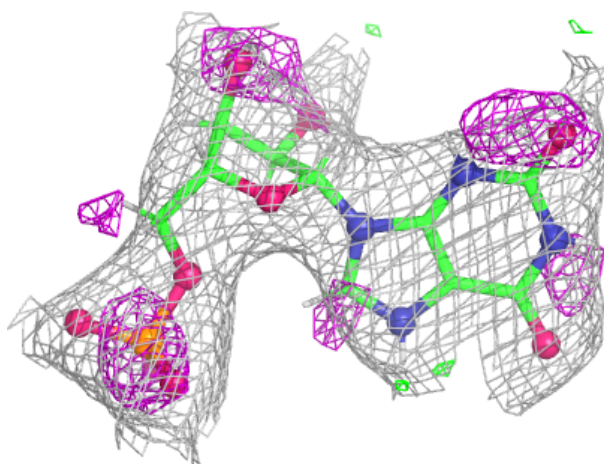
Electron density around XMP D 601:

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



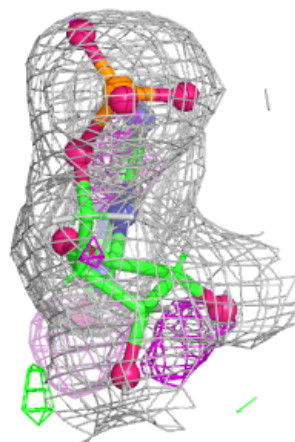
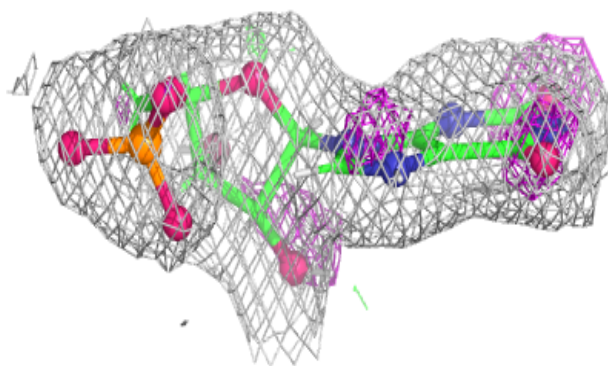
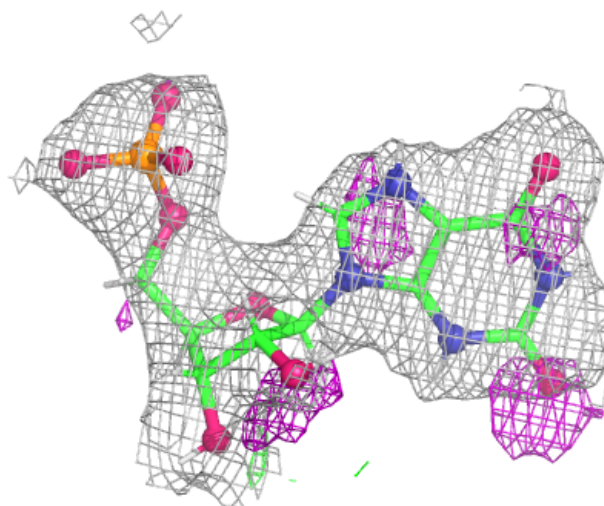
Electron density around XMP A 601:

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



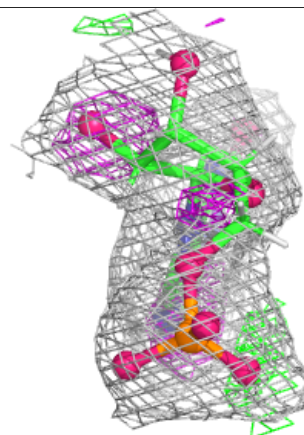
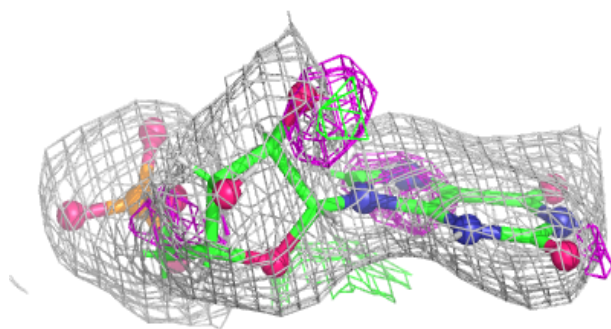
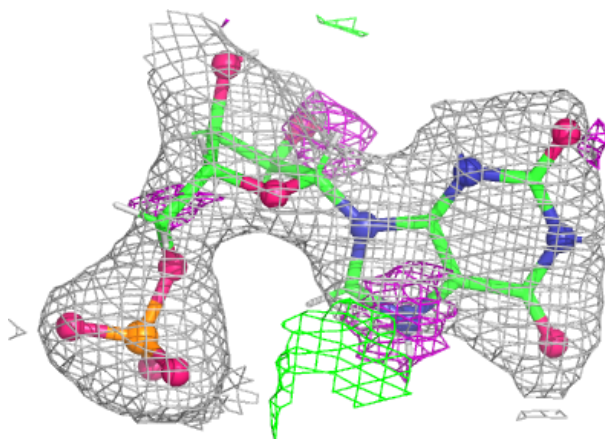
Electron density around XMP B 601:

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



Electron density around XMP 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.