



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 07:39 AM UTC

PDB ID : 1EYZ / pdb_00001eyz
Title : STRUCTURE OF ESCHERICHIA COLI PURT-ENCODED GLYCINAMIDE RIBONUCLEOTIDE TRANSFORMYLASE COMPLEXED WITH MG AND AMPPNP
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Deposited on : 2000-05-09
Resolution : 1.75 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

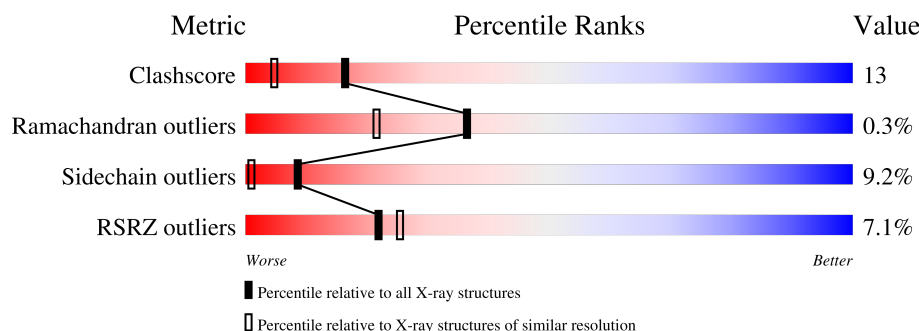
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.75 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))
Clashscore	190562	3299 (1.76-1.76)
Ramachandran outliers	187476	3274 (1.76-1.76)
Sidechain outliers	187428	3274 (1.76-1.76)
RSRZ outliers	180081	3183 (1.76-1.76)

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	392	
1	B	392	

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 6878 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 PHOSPHORIBOSYLGLYCINAMIDE FORMYLTRANSFERASE 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	389	Total	C	N	O	S	0	1	0
			2970	1870	529	558	13			
1	B	389	Total	C	N	O	S	0	2	0
			2979	1875	534	557	13			

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	3	Total	Mg	0	0
			3	3		
2	B	2	Total	Mg	0	0
			2	2		

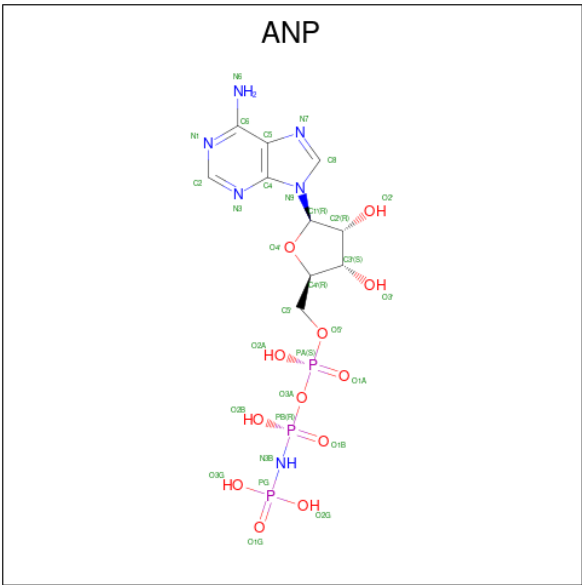
- Molecule 3 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Na	0	0
			1	1		
3	B	1	Total	Na	0	0
			1	1		

- Molecule 4 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

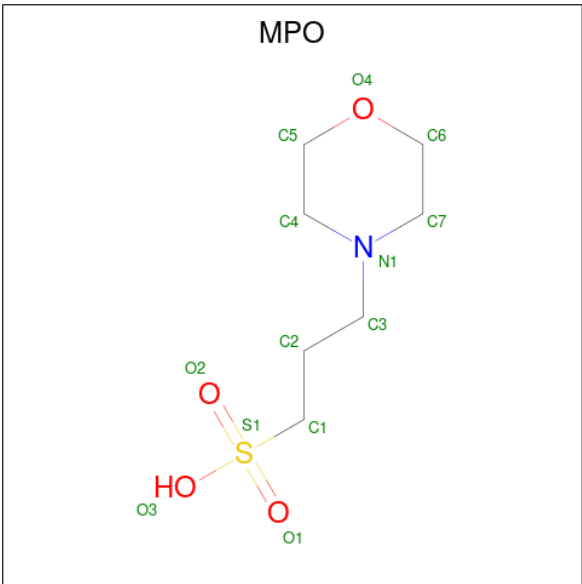
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Cl	0	0
			1	1		

- Molecule 5 is PHOSPHOAMINOPHOSPHONIC ACID-ADENYLATE ESTER (CCD ID: ANP) (formula: C₁₀H₁₇N₆O₁₂P₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	P	0	0
			31	10	6	12	3		
5	B	1	Total	C	N	O	P	0	0
			31	10	6	12	3		

- Molecule 6 is 3[N-MORPHOLINO]PROPANE SULFONIC ACID (CCD ID: MPO) (formula: $C_7H_{15}NO_4S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	A	1	Total	C	N	O	S	0	0
			13	7	1	4	1		

- Molecule 7 is water.

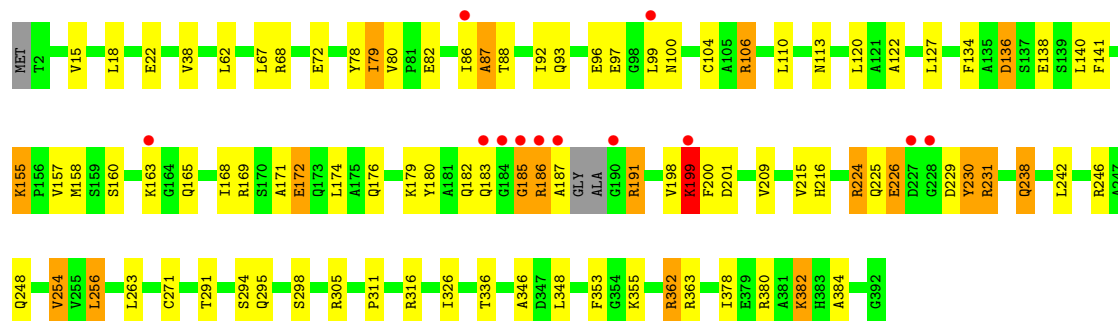
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	488	Total 488	O 488	0	0
7	B	358	Total 358	O 358	0	0

3 Residue-property plots

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

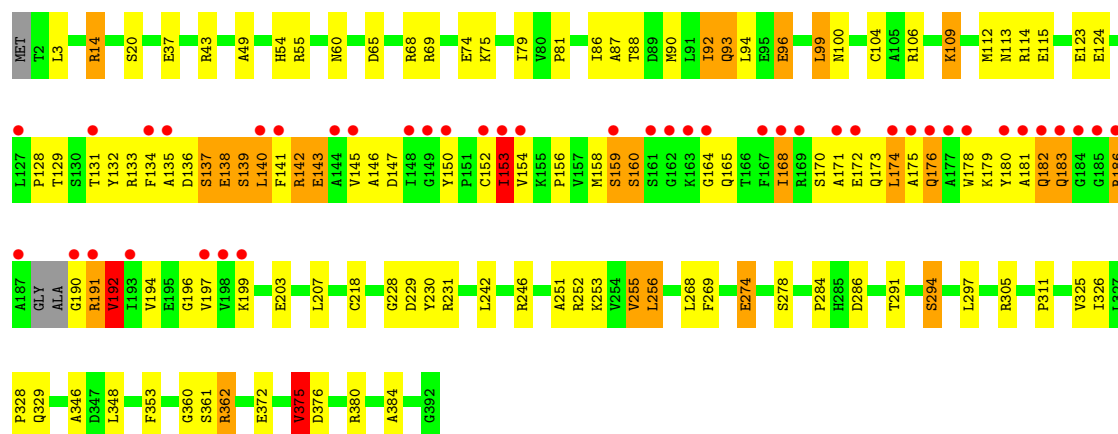
• Molecule 1: PHOSPHORIBOSYLGLYCINAMIDE FORMYLTRANSFERASE 2

Chain A: 



• Molecule 1: PHOSPHORIBOSYLGLYCINAMIDE FORMYLTRANSFERASE 2

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	62.30Å 179.50Å 75.70Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 1.75 30.00 – 1.75	Depositor EDS
% Data completeness (in resolution range)	96.1 (30.00-1.75) 96.4 (30.00-1.75)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.89 (at 1.70Å)	Xtriage
Refinement program	TNT	Depositor
R, R_{free}	0.185 , 0.228 0.177 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	16.4	Xtriage
Anisotropy	0.321	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 112.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6878	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.12% 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: NA, CL, MPO, ANP, MG

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	1.04	0/3023	1.47	23/4095 (0.6%)
1	B	1.04	5/3036 (0.2%)	1.49	25/4111 (0.6%)
All	All	1.04	5/6059 (0.1%)	1.48	48/8206 (0.6%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	284	PRO	CA-C	7.27	1.61	1.52
1	B	137	SER	N-CA	6.09	1.53	1.46
1	B	123	GLU	CD-OE2	5.58	1.35	1.25
1	B	284	PRO	N-CD	5.54	1.55	1.47
1	B	138	GLU	CD-OE2	5.18	1.35	1.25

All (48) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	60	ASN	CA-CB-CG	-9.48	103.12	112.60
1	A	353	PHE	N-CA-C	8.86	123.35	112.54
1	B	353	PHE	N-CA-C	8.57	123.00	112.54
1	A	238	GLN	N-CA-CB	-7.53	98.46	109.83
1	B	375	VAL	N-CA-CB	-7.38	101.92	110.55
1	B	168	ILE	N-CA-C	7.26	118.59	107.99
1	A	172	GLU	CB-CG-CD	-7.03	100.64	112.60
1	B	153	ILE	CA-C-N	-6.84	114.71	123.19
1	B	153	ILE	C-N-CA	-6.84	114.71	123.19
1	A	198	VAL	CA-C-N	-6.36	114.01	122.85
1	A	198	VAL	C-N-CA	-6.36	114.01	122.85
1	A	80	VAL	N-CA-C	6.36	114.44	108.15
1	A	294	SER	N-CA-C	6.26	120.91	113.28
1	B	192	VAL	N-CA-C	6.00	117.36	108.65

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	54	HIS	CA-CB-CG	-5.84	107.96	113.80
1	B	128	PRO	N-CA-CB	5.80	107.12	103.23
1	B	286	ASP	N-CA-C	5.79	118.34	111.33
1	B	284	PRO	N-CA-CB	5.74	108.14	103.32
1	B	294	SER	N-CA-C	5.73	119.99	112.89
1	A	186	ARG	CA-C-N	5.72	132.00	121.70
1	A	186	ARG	C-N-CA	5.72	132.00	121.70
1	A	168	ILE	N-CA-C	5.69	116.93	108.23
1	B	170	SER	N-CA-C	5.60	117.91	109.23
1	A	104	CYS	N-CA-CB	5.53	120.03	111.24
1	A	199	LYS	N-CA-CB	5.52	119.49	110.39
1	A	201	ASP	N-CA-C	-5.45	105.25	111.14
1	B	14	ARG	N-CA-C	5.45	117.84	109.07
1	A	326	ILE	N-CA-C	-5.43	99.13	106.85
1	B	199	LYS	N-CA-CB	5.43	118.48	110.49
1	A	79	ILE	N-CA-CB	5.42	118.69	111.64
1	A	185	GLY	CA-C-N	5.41	131.88	121.54
1	A	185	GLY	C-N-CA	5.41	131.88	121.54
1	B	136	ASP	CA-C-N	5.41	131.46	121.94
1	B	136	ASP	C-N-CA	5.41	131.46	121.94
1	B	326	ILE	N-CA-C	-5.41	99.17	106.85
1	A	38	VAL	N-CA-C	5.35	115.82	108.12
1	A	134	PHE	CA-CB-CG	-5.33	108.47	113.80
1	A	172	GLU	CB-CA-C	-5.27	102.03	110.79
1	A	136	ASP	N-CA-CB	5.25	119.64	110.98
1	A	230	TYR	N-CA-C	-5.23	102.54	110.28
1	B	360	GLY	N-CA-C	-5.15	106.36	111.56
1	B	176	GLN	CB-CA-C	-5.14	102.80	110.88
1	A	122	ALA	N-CA-C	5.13	116.68	111.14
1	B	175	ALA	CA-C-N	5.12	127.10	120.44
1	B	175	ALA	C-N-CA	5.12	127.10	120.44
1	B	60	ASN	N-CA-CB	5.07	118.00	110.49
1	B	325	VAL	N-CA-C	5.07	116.55	109.80
1	B	104	CYS	N-CA-CB	5.01	119.21	111.24

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	2970	0	2992	59	0
1	B	2979	0	3006	96	0
2	A	3	0	0	0	0
2	B	2	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	1	0	0	0	0
5	A	31	0	13	1	0
5	B	31	0	13	3	0
6	A	13	0	15	1	0
7	A	488	0	0	18	0
7	B	358	0	0	9	1
All	All	6878	0	6039	157	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 13.

All (157) 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:B:79:ILE:HD11	1:B:99:LEU:HD22	1.31	1.08
1:B:68:ARG:HB2	1:B:90:MET:HE3	1.30	1.08
1:B:115:GLU:HG3	1:B:134:PHE:CE1	2.01	0.95
1:B:191:ARG:HG3	1:B:192:VAL:N	1.82	0.94
1:B:173:GLN:HA	1:B:176:GLN:NE2	1.85	0.91
1:B:380[B]:ARG:HG2	1:B:380[B]:ARG:HH11	1.36	0.90
1:B:153:ILE:HD12	1:B:197:VAL:HG22	1.51	0.89
1:B:68:ARG:CB	1:B:90:MET:HE3	2.04	0.87
1:A:291:THR:HB	7:A:2002:HOH:O	1.80	0.81
1:B:156:PRO:HG2	1:B:159:SER:HB2	1.61	0.81
1:B:145:VAL:HG13	1:B:152:CYS:SG	2.22	0.79
1:B:141:PHE:O	1:B:145:VAL:HG23	1.83	0.78
1:B:79:ILE:HD11	1:B:99:LEU:CD2	2.12	0.76
1:A:295:GLN:HB2	7:A:2002:HOH:O	1.86	0.76
1:B:173:GLN:HA	1:B:176:GLN:HE21	1.49	0.76
1:B:153:ILE:CD1	1:B:197:VAL:HG22	2.15	0.76
1:B:132:TYR:O	1:B:133:ARG:HD2	1.89	0.73
1:B:143:GLU:O	1:B:146:ALA:HB3	1.89	0.72
1:B:65:ASP:O	1:B:69:ARG:HG3	1.91	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:298:SER:HA	7:A:2002:HOH:O	1.91	0.71
1:A:79:ILE:HD11	1:A:99:LEU:HD23	1.73	0.69
1:A:362:ARG:NH1	7:A:1375:HOH:O	2.26	0.69
1:B:178:TRP:O	1:B:182:GLN:NE2	2.26	0.68
1:A:113:ASN:HD22	1:A:158:MET:HE2	1.59	0.67
1:B:156:PRO:HG2	1:B:159:SER:CB	2.24	0.67
1:A:68:ARG:O	1:A:72:GLU:HG3	1.95	0.67
1:A:106:ARG:NH2	7:A:1297:HOH:O	2.28	0.66
1:B:153:ILE:HD12	1:B:197:VAL:CG2	2.24	0.65
1:A:180:TYR:O	1:A:183:GLN:HB2	1.96	0.65
1:A:199:LYS:NZ	7:A:1804:HOH:O	2.30	0.65
1:A:92:ILE:O	1:A:96:GLU:HG3	1.98	0.64
1:B:348:LEU:HB2	1:B:380[A]:ARG:NH2	2.13	0.63
1:B:228:GLY:HA2	7:B:1596:HOH:O	1.98	0.63
1:A:18:LEU:HD11	1:A:67:LEU:HD11	1.82	0.62
1:A:378:ILE:O	1:A:382:LYS:HG3	2.00	0.62
1:A:68:ARG:NH2	1:A:97:GLU:OE1	2.34	0.61
1:A:295:GLN:NE2	7:A:2002:HOH:O	2.28	0.61
1:B:93:GLN:O	1:B:96:GLU:HB3	2.01	0.60
1:A:169:ARG:NH1	7:A:1984:HOH:O	2.31	0.60
1:B:346:ALA:O	1:B:380[A]:ARG:NH2	2.34	0.60
1:B:362:ARG:NH2	7:B:1642:HOH:O	2.35	0.59
1:B:173:GLN:O	1:B:176:GLN:HB2	2.01	0.59
1:B:174:LEU:HD23	1:B:174:LEU:N	2.16	0.59
1:B:274:GLU:HB3	7:B:2046:HOH:O	2.01	0.59
1:A:79:ILE:HD11	1:A:99:LEU:CD2	2.32	0.59
1:A:113:ASN:ND2	1:A:158:MET:HE2	2.16	0.59
1:B:68:ARG:HE	1:B:90:MET:HE1	1.67	0.59
1:B:138:GLU:OE1	1:B:142:ARG:NH1	2.36	0.59
1:B:380[B]:ARG:HG2	1:B:380[B]:ARG:NH1	2.11	0.58
1:B:180:TYR:C	1:B:182:GLN:H	2.10	0.57
1:B:94:LEU:HB3	1:B:99:LEU:HD13	1.87	0.56
1:A:136:ASP:HB3	1:A:191:ARG:HD3	1.88	0.56
1:A:216:HIS:HD2	7:A:1811:HOH:O	1.89	0.56
1:A:183:GLN:HB3	7:A:1803:HOH:O	2.06	0.55
1:B:81:PRO:HB3	1:B:86:ILE:CD1	2.36	0.55
1:A:86:ILE:O	1:A:88:THR:HG23	2.07	0.55
1:A:185:GLY:C	1:A:187:ALA:H	2.15	0.54
1:B:88:THR:O	1:B:92:ILE:HD12	2.06	0.54
1:A:68:ARG:HD2	7:A:1316:HOH:O	2.07	0.54
1:A:171:ALA:HA	1:A:174:LEU:HG	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:115:GLU:HG3	1:B:134:PHE:CZ	2.43	0.53
1:A:229:ASP:HA	7:A:1394:HOH:O	2.09	0.53
1:B:380[A]:ARG:NH1	7:B:1479:HOH:O	2.21	0.53
1:B:79:ILE:CD1	1:B:99:LEU:HD22	2.23	0.52
1:B:113:ASN:ND2	1:B:158:MET:HE2	2.24	0.52
5:B:961:ANP:H8	5:B:961:ANP:O5'	2.10	0.51
1:B:291:THR:HA	1:B:294:SER:OG	2.10	0.51
1:B:135:ALA:CB	1:B:141:PHE:HA	2.41	0.51
1:A:209:VAL:HG22	1:A:263:LEU:HD12	1.92	0.51
1:A:348:LEU:HD11	1:A:384:ALA:CB	2.41	0.51
1:B:20:SER:HB3	1:B:49:ALA:HB3	1.92	0.50
1:B:156:PRO:HD3	1:B:164:GLY:O	2.11	0.50
1:B:179:LYS:O	1:B:182:GLN:HB2	2.10	0.50
1:A:86:ILE:HG22	1:A:87:ALA:N	2.26	0.50
1:B:231:ARG:NH2	7:B:1674:HOH:O	2.44	0.50
1:B:14:ARG:HG2	1:B:37:GLU:HB3	1.94	0.49
1:A:336:THR:HB	1:B:3:LEU:HD11	1.93	0.49
1:B:253:LYS:NZ	7:B:2103:HOH:O	2.46	0.49
1:B:171:ALA:O	1:B:174:LEU:HG	2.11	0.49
1:A:224:ARG:HG3	7:A:1717:HOH:O	2.12	0.49
1:A:15:VAL:HG22	1:A:78:TYR:HB2	1.95	0.48
1:B:138:GLU:O	1:B:142:ARG:HB2	2.13	0.48
1:A:127:LEU:CD1	1:A:254:VAL:HG23	2.43	0.48
5:A:400:ANP:O5'	5:A:400:ANP:H8	2.14	0.48
1:B:132:TYR:HA	1:B:194:VAL:O	2.14	0.48
1:B:141:PHE:CE1	1:B:145:VAL:CG2	2.96	0.48
1:B:131:THR:OG1	1:B:196:GLY:HA3	2.13	0.48
1:B:160:SER:HA	5:B:961:ANP:O1B	2.14	0.47
1:A:172:GLU:HG3	7:A:2092:HOH:O	2.13	0.47
1:B:93:GLN:OE1	1:B:93:GLN:HA	2.11	0.47
1:B:55:ARG:HE	1:B:74:GLU:CD	2.21	0.47
1:A:127:LEU:HD11	1:A:254:VAL:HG23	1.95	0.47
1:B:109:LYS:O	1:B:112:MET:HG3	2.15	0.47
1:B:203:GLU:HG3	1:B:269:PHE:CD1	2.49	0.47
1:B:180:TYR:C	1:B:182:GLN:N	2.73	0.47
1:A:110:LEU:HD13	1:A:120:LEU:HD22	1.96	0.47
1:B:114:ARG:HH22	5:B:961:ANP:PB	2.37	0.47
1:A:348:LEU:HD11	1:A:384:ALA:HB2	1.96	0.47
1:B:252:ARG:O	1:B:256:LEU:HB2	2.15	0.46
1:A:231:ARG:HA	1:A:231:ARG:HD3	1.70	0.46
1:B:154:VAL:O	1:B:165:GLN:HA	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:328:PRO:HD2	1:B:362:ARG:O	2.16	0.46
1:B:251:ALA:O	1:B:255:VAL:HG13	2.15	0.46
1:B:269:PHE:CE2	1:B:278:SER:HB2	2.51	0.46
1:B:88:THR:HG21	1:B:112:MET:HG2	1.98	0.45
1:A:62:LEU:HD23	1:A:62:LEU:HA	1.82	0.45
1:B:139:SER:HB3	7:B:2060:HOH:O	2.16	0.45
1:B:173:GLN:HA	1:B:176:GLN:HE22	1.72	0.45
1:B:180:TYR:CD2	1:B:183:GLN:HB2	2.51	0.45
1:B:156:PRO:CG	1:B:159:SER:CB	2.95	0.45
1:B:380[B]:ARG:NH1	1:B:380[B]:ARG:CG	2.79	0.45
1:A:18:LEU:HD23	1:A:18:LEU:HA	1.83	0.45
1:A:155:LYS:HG2	1:A:165:GLN:HG3	1.98	0.45
1:A:82:GLU:HB2	7:A:1103:HOH:O	2.17	0.45
1:B:181:ALA:C	1:B:182:GLN:NE2	2.75	0.45
1:A:138:GLU:O	1:A:141:PHE:HB3	2.16	0.44
1:A:180:TYR:CD2	1:A:183:GLN:NE2	2.86	0.44
1:A:200:PHE:HA	1:A:271:CYS:SG	2.57	0.44
1:B:203:GLU:HG3	1:B:269:PHE:HD1	1.82	0.44
1:A:185:GLY:O	1:A:187:ALA:N	2.50	0.44
1:B:135:ALA:HB1	1:B:141:PHE:N	2.33	0.44
1:A:256:LEU:HD12	1:A:256:LEU:HA	1.87	0.44
1:B:252:ARG:HG3	1:B:256:LEU:HD22	2.00	0.44
1:B:348:LEU:HD11	1:B:384:ALA:CB	2.48	0.44
1:A:225:GLN:NE2	7:A:1815:HOH:O	2.51	0.44
1:B:297:LEU:HA	7:B:1554:HOH:O	2.18	0.44
1:A:22:GLU:OE2	1:A:363:ARG:NH1	2.48	0.43
1:B:329:GLN:HG2	7:B:1565:HOH:O	2.19	0.43
1:A:100:ASN:C	1:A:100:ASN:HD22	2.26	0.43
1:B:115:GLU:HG3	1:B:134:PHE:CD1	2.52	0.43
1:B:141:PHE:CE1	1:B:145:VAL:HG21	2.53	0.43
1:B:305:ARG:HD2	1:B:311:PRO:O	2.19	0.43
1:A:157:VAL:HG22	1:A:191:ARG:O	2.19	0.43
1:A:209:VAL:O	1:A:215:VAL:HA	2.19	0.43
1:A:242:LEU:O	1:A:246:ARG:HG3	2.19	0.43
1:A:22:GLU:OE1	1:A:355:LYS:NZ	2.44	0.42
1:B:150:TYR:HE2	1:B:168:ILE:HG22	1.84	0.42
1:B:348:LEU:HD11	1:B:384:ALA:HB2	2.01	0.42
1:B:137:SER:O	1:B:140:LEU:N	2.51	0.42
1:B:242:LEU:O	1:B:246:ARG:HG3	2.19	0.42
1:B:141:PHE:CD2	1:B:178:TRP:HB2	2.54	0.42
1:A:305:ARG:HD3	1:A:311:PRO:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:207:LEU:HB2	1:B:218:CYS:HB2	2.02	0.42
1:A:68:ARG:NH2	1:A:97:GLU:CD	2.78	0.41
1:A:226:GLU:HB3	1:A:231:ARG:HG2	2.02	0.41
1:B:113:ASN:OD1	1:B:115:GLU:HB2	2.20	0.41
1:B:20:SER:CB	1:B:49:ALA:HB3	2.50	0.41
1:B:186:ARG:NH1	1:B:190:GLY:N	2.69	0.41
1:B:86:ILE:HG22	1:B:87:ALA:N	2.35	0.41
1:A:346:ALA:O	1:A:380:ARG:NH1	2.54	0.41
1:B:182:GLN:NE2	1:B:182:GLN:N	2.69	0.41
6:A:950:MPO:H72	7:A:1204:HOH:O	2.21	0.41
1:B:180:TYR:HD2	1:B:183:GLN:HB2	1.86	0.40
1:B:154:VAL:O	1:B:165:GLN:HG3	2.22	0.40
1:B:156:PRO:HB2	1:B:159:SER:HB3	2.03	0.40
1:B:375:VAL:HG12	1:B:376:ASP:N	2.35	0.40
1:A:248:GLN:NE2	7:A:2091:HOH:O	2.48	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)
7:B:1663:HOH:O	7:B:1663:HOH:O[2_655]	2.13	0.07

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	386/392 (98%)	371 (96%)	13 (3%)	2 (0%)	24	11
1	B	387/392 (99%)	371 (96%)	16 (4%)	0	100	100
All	All	773/784 (99%)	742 (96%)	29 (4%)	2 (0%)	36	21

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	186	ARG
1	A	87	ALA

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	312/312 (100%)	290 (93%)	22 (7%)	13	2
1	B	313/312 (100%)	276 (88%)	37 (12%)	5	0
All	All	625/624 (100%)	566 (91%)	59 (9%)	8	1

All (59) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	93[A]	GLN
1	A	93[B]	GLN
1	A	106	ARG
1	A	140	LEU
1	A	155	LYS
1	A	160	SER
1	A	163	LYS
1	A	176	GLN
1	A	179	LYS
1	A	182	GLN
1	A	191	ARG
1	A	199	LYS
1	A	224	ARG
1	A	226	GLU
1	A	230	TYR
1	A	231	ARG
1	A	238	GLN
1	A	254	VAL
1	A	256	LEU
1	A	316	ARG
1	A	362	ARG
1	A	382	LYS

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Mol	Chain	Res	Type
1	B	43[A]	ARG
1	B	43[B]	ARG
1	B	75	LYS
1	B	92	ILE
1	B	93	GLN
1	B	96	GLU
1	B	99	LEU
1	B	100	ASN
1	B	106	ARG
1	B	109	LYS
1	B	124	GLU
1	B	129	THR
1	B	139	SER
1	B	140	LEU
1	B	142	ARG
1	B	143	GLU
1	B	147	ASP
1	B	153	ILE
1	B	159	SER
1	B	160	SER
1	B	172	GLU
1	B	174	LEU
1	B	182	GLN
1	B	183	GLN
1	B	186	ARG
1	B	191	ARG
1	B	192	VAL
1	B	229	ASP
1	B	230	TYR
1	B	255	VAL
1	B	256	LEU
1	B	268	LEU
1	B	274	GLU
1	B	361	SER
1	B	362	ARG
1	B	372	GLU
1	B	375	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	100	ASN

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Mol	Chain	Res	Type
1	A	183	GLN
1	A	216	HIS
1	A	225	GLN
1	A	238	GLN
1	A	248	GLN
1	A	329	GLN
1	A	339	ASN
1	A	342	ASN
1	B	100	ASN
1	B	126	GLN
1	B	176	GLN
1	B	182	GLN
1	B	225	GLN
1	B	248	GLN
1	B	329	GLN
1	B	387	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 11 ligands modelled in this entry, 8 are monoatomic - leaving 3 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	ANP	B	961	2	33,33,33	1.36	5 (15%)	45,52,52	1.15	4 (8%)
5	ANP	A	400	2	33,33,33	1.39	8 (24%)	45,52,52	1.44	6 (13%)
6	MPO	A	950	-	13,13,13	1.95	2 (15%)	17,17,17	1.82	3 (17%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	ANP	B	961	2	-	2/18/38/38	0/3/3/3
5	ANP	A	400	2	-	2/18/38/38	0/3/3/3
6	MPO	A	950	-	-	0/7/15/15	0/1/1/1

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	950	MPO	C3-N1	4.44	1.57	1.47
6	A	950	MPO	C1-S1	4.35	1.83	1.77
5	B	961	ANP	C6-N6	-3.23	1.26	1.34
5	A	400	ANP	PG-O1G	3.05	1.50	1.46
5	A	400	ANP	C6-N6	-2.96	1.26	1.34
5	B	961	ANP	C2-N1	2.77	1.38	1.33
5	B	961	ANP	PB-O1B	2.69	1.50	1.46
5	B	961	ANP	PG-O2G	-2.66	1.49	1.56
5	A	400	ANP	PB-O1B	2.51	1.50	1.46
5	A	400	ANP	PG-O2G	-2.34	1.50	1.56
5	B	961	ANP	PG-O1G	2.33	1.49	1.46
5	A	400	ANP	PB-O2B	-2.31	1.50	1.56
5	A	400	ANP	C2-N1	2.14	1.37	1.33
5	A	400	ANP	PB-O3A	2.08	1.61	1.59
5	A	400	ANP	PA-O3A	-2.02	1.57	1.59

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	400	ANP	O1G-PG-N3B	-4.74	104.79	111.77
5	A	400	ANP	C2-N1-C6	4.48	126.09	118.73
6	A	950	MPO	O1-S1-C1	4.27	113.18	106.73
6	A	950	MPO	O3-S1-O1	-4.23	100.81	111.40
5	B	961	ANP	C2-N1-C6	3.84	125.04	118.73

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	400	ANP	N3-C2-N1	-3.20	123.73	128.58
5	B	961	ANP	N3-C2-N1	-2.56	124.71	128.58
5	A	400	ANP	C5-C6-N1	-2.49	111.19	117.51
6	A	950	MPO	O3-S1-C1	2.43	110.76	106.00
5	B	961	ANP	C5-C6-N1	-2.37	111.49	117.51
5	A	400	ANP	O3'-C3'-C2'	2.33	119.29	111.82
5	A	400	ANP	C5-C6-N6	2.33	129.06	123.29
5	B	961	ANP	C5-C6-N6	2.29	128.97	123.29

There are no chirality outliers.

All (4) torsion outliers are listed below:

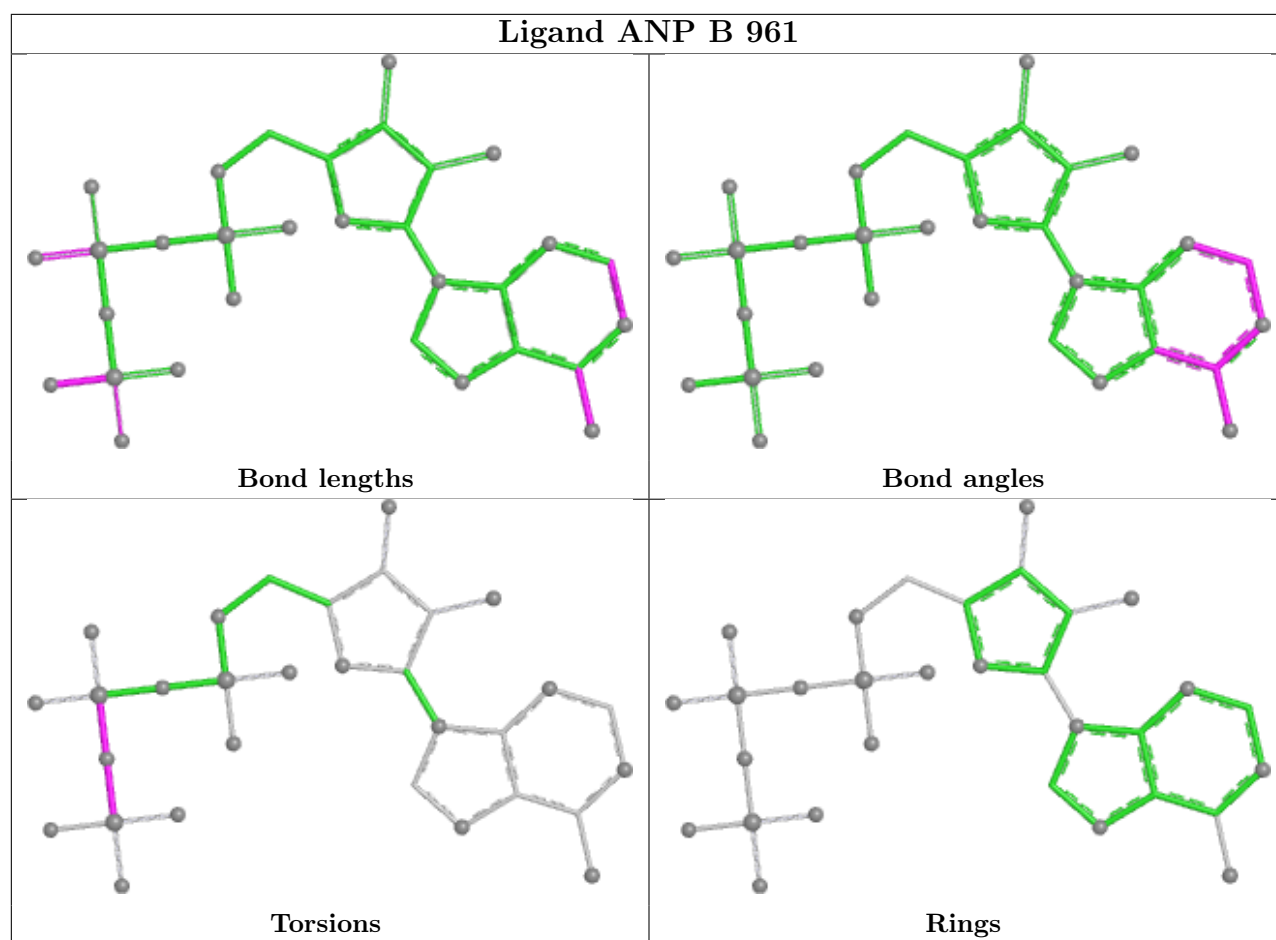
Mol	Chain	Res	Type	Atoms
5	A	400	ANP	PB-N3B-PG-O1G
5	A	400	ANP	PG-N3B-PB-O1B
5	B	961	ANP	PB-N3B-PG-O1G
5	B	961	ANP	PG-N3B-PB-O1B

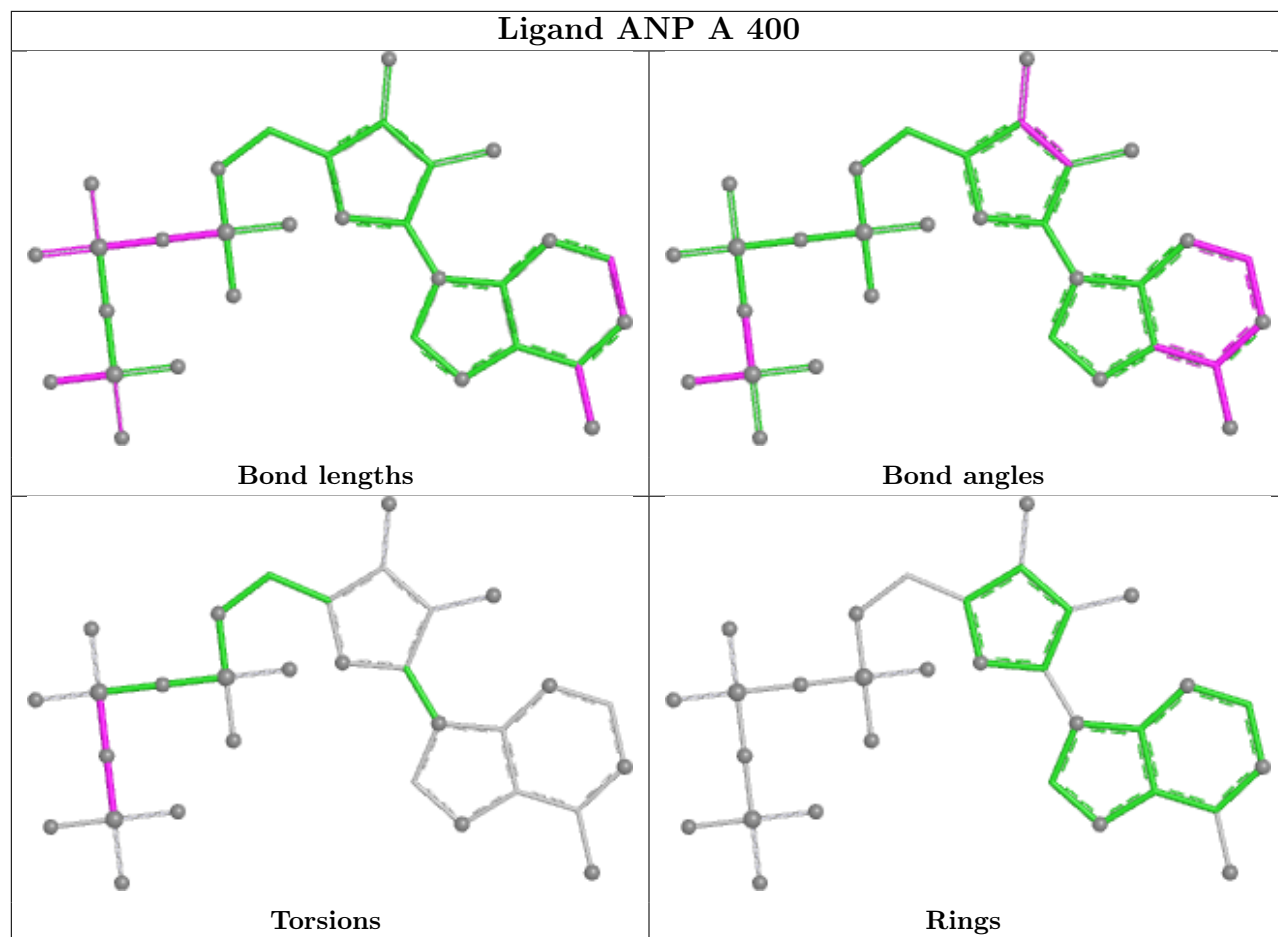
There are no ring outliers.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	961	ANP	3	0
5	A	400	ANP	1	0
6	A	950	MPO	1	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	389/392 (99%)	-0.24	12 (3%)	51	58	10, 17, 49, 97	1 (0%)
1	B	389/392 (99%)	0.23	43 (11%)	10	11	11, 22, 81, 100	2 (0%)
All	All	778/784 (99%)	-0.01	55 (7%)	22	25	10, 19, 70, 100	3 (0%)

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	187	ALA	5.3
1	B	135	ALA	5.2
1	B	190	GLY	5.0
1	B	187	ALA	5.0
1	B	175	ALA	5.0
1	A	190	GLY	4.7
1	B	149	GLY	4.5
1	B	180	TYR	4.1
1	A	185	GLY	4.0
1	B	141	PHE	3.9
1	B	181	ALA	3.9
1	B	176	GLN	3.9
1	B	131	THR	3.5
1	B	134	PHE	3.4
1	B	171	ALA	3.3
1	B	186	ARG	3.3
1	B	168	ILE	3.3
1	B	185	GLY	3.2
1	B	169	ARG	3.2
1	B	162	GLY	3.2
1	A	86	ILE	3.1
1	B	150	TYR	3.1
1	A	184	GLY	3.0
1	A	186	ARG	3.0

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Mol	Chain	Res	Type	RSRZ
1	B	191	ARG	3.0
1	B	167	PHE	3.0
1	B	199	LYS	2.9
1	B	164	GLY	2.9
1	B	198	VAL	2.8
1	B	140	LEU	2.8
1	B	178	TRP	2.8
1	B	197	VAL	2.8
1	B	177	ALA	2.7
1	B	163	LYS	2.7
1	B	145	VAL	2.7
1	B	154	VAL	2.7
1	B	127	LEU	2.6
1	B	159	SER	2.6
1	B	193	ILE	2.6
1	B	144	ALA	2.6
1	A	163	LYS	2.5
1	B	152	CYS	2.5
1	B	174	LEU	2.5
1	B	148	ILE	2.5
1	A	228	GLY	2.4
1	A	199	LYS	2.4
1	B	183	GLN	2.3
1	A	99	LEU	2.3
1	B	153	ILE	2.3
1	B	182	GLN	2.2
1	A	227	ASP	2.2
1	B	184	GLY	2.2
1	B	161	SER	2.2
1	A	183	GLN	2.1
1	B	172	GLU	2.1

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

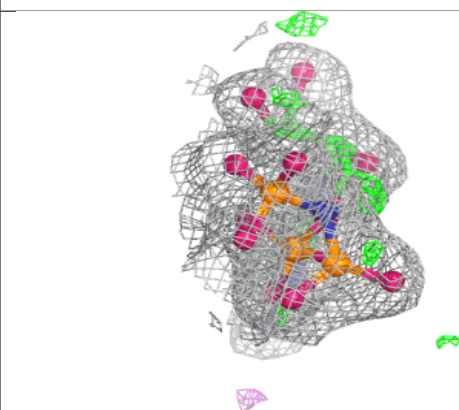
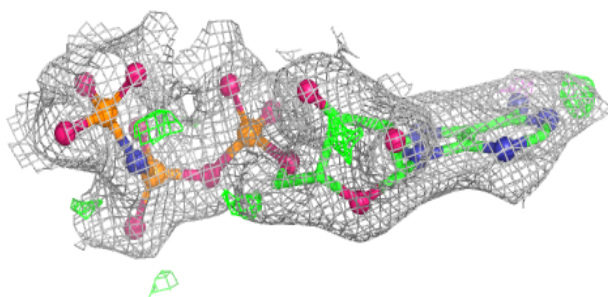
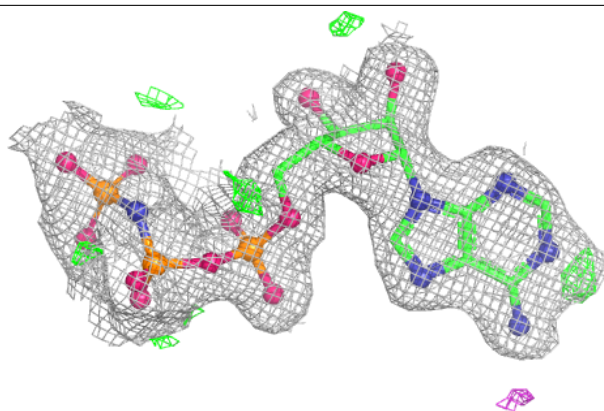
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	NA	B	960	1/1	0.62	0.30	37,37,37,37	0
3	NA	A	961	1/1	0.90	0.24	21,21,21,21	0
5	ANP	B	961	31/31	0.96	0.07	18,28,37,62	0
2	MG	A	403	1/1	0.97	0.06	29,29,29,29	0
2	MG	B	402	1/1	0.98	0.04	23,23,23,23	0
4	CL	A	970	1/1	0.98	0.11	26,26,26,26	0
2	MG	A	401	1/1	0.98	0.04	18,18,18,18	0
6	MPO	A	950	13/13	0.98	0.06	15,23,31,40	0
5	ANP	A	400	31/31	0.99	0.04	12,15,22,25	0
2	MG	A	402	1/1	0.99	0.04	15,15,15,15	0
2	MG	B	401	1/1	0.99	0.04	25,25,25,25	0

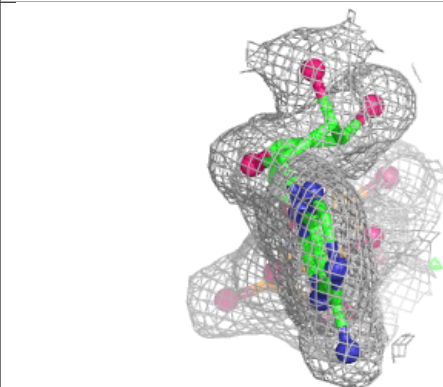
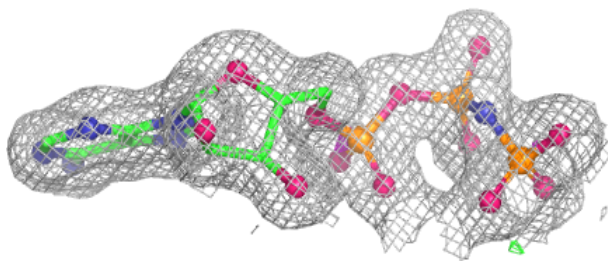
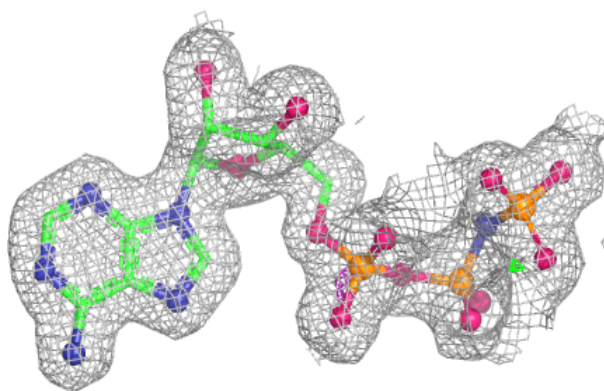
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around ANP B 961:

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

**Electron density around ANP A 400:**

$2mF_o-DF_c$ (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.