



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 26, 2026 – 04:30 AM EDT

PDB ID : 7LV3 / pdb_00007lv3
Title : Crystal structure of human protein kinase G (PKG) R-C complex in inhibited state
Authors : Sharma, R.; Lying, Q.; Casteel, D.; Kim, C.
Deposited on : 2021-02-23
Resolution : 2.41 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

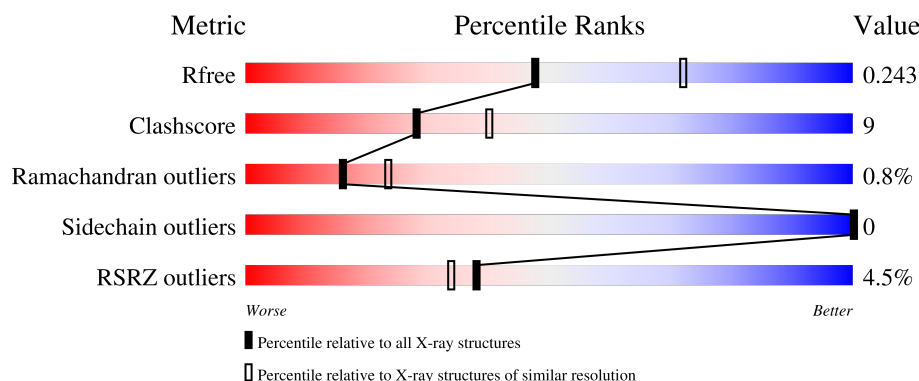
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.41 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	6062 (2.44-2.40)
Clashscore	190562	6562 (2.44-2.40)
Ramachandran outliers	187476	6481 (2.44-2.40)
Sidechain outliers	187428	6482 (2.44-2.40)
RSRZ outliers	180081	6066 (2.44-2.40)

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	617	
1	B	617	

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 9432 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.

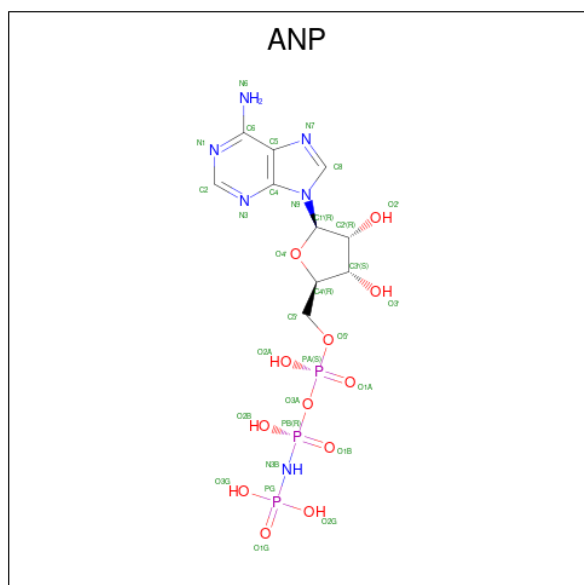
- Molecule 1 is a protein called Isoform Beta of cGMP-dependent protein kinase 1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	601	Total	C	N	O	P	S	0	0	0
			4653	2973	772	883	1	24			
1	B	594	Total	C	N	O	P	S	0	0	0
			4574	2926	752	871	1	24			

- Molecule 2 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Mn	0	0
			2	2		
2	B	2	Total	Mn	0	0
			2	2		

- Molecule 3 is PHOSPHOAMINOPHOSPHONIC ACID-ADENYLATE ESTER (CCD ID: ANP) (formula: C₁₀H₁₇N₆O₁₂P₃) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 4 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



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

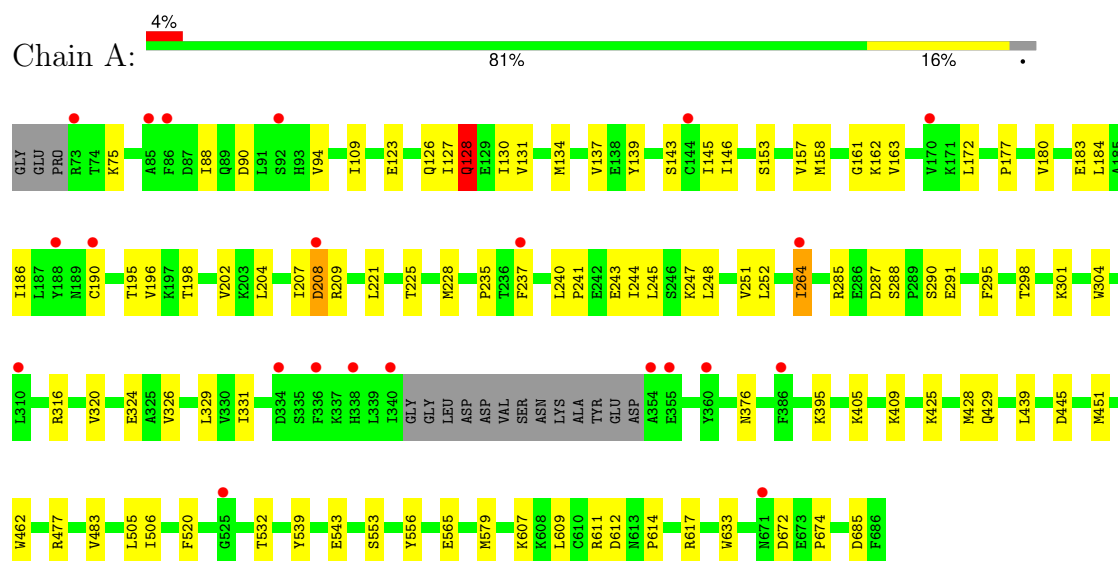
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	66	Total	O	0	0
			66	66		
5	B	61	Total	O	0	0
			61	61		

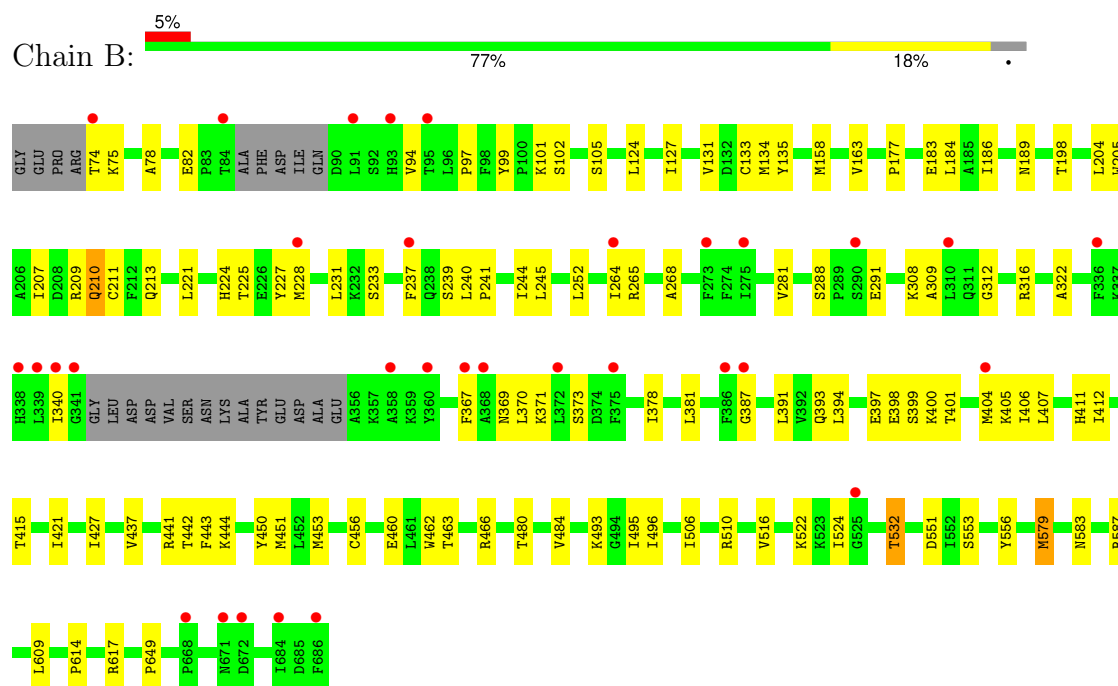
3 Residue-property plots

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: Isoform Beta of cGMP-dependent protein kinase 1



- Molecule 1: Isoform Beta of cGMP-dependent protein kinase 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	99.83Å 112.82Å 150.72Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.38 – 2.41 47.38 – 2.41	Depositor EDS
% Data completeness (in resolution range)	100.0 (47.38-2.41) 89.3 (47.38-2.41)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.53 (at 2.34Å)	Xtriage
Refinement program	PHENIX 1.17.1 _3660	Depositor
R, R_{free}	0.200 , 0.238 0.205 , 0.243	Depositor DCC
R_{free} test set	2000 reflections (2.73%)	wwPDB-VP
Wilson B-factor (Å ²)	57.8	Xtriage
Anisotropy	0.096	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 48.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	9432	wwPDB-VP
Average B, all atoms (Å ²)	87.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.72% 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: TPO, MN, EDO, ANP

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	0.27	3/4745 (0.1%)	0.49	6/6438 (0.1%)
1	B	0.22	0/4662	0.51	2/6324 (0.0%)
All	All	0.25	3/9407 (0.0%)	0.50	8/12762 (0.1%)

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	A	0	2
1	B	0	1
All	All	0	3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	674	PRO	CA-C	7.03	1.55	1.51
1	A	264	ILE	CB-CG2	5.38	1.70	1.52
1	A	208	ASP	CA-CB	5.27	1.62	1.53

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	207	ILE	CA-C-N	-9.32	106.55	122.67
1	A	207	ILE	C-N-CA	-9.32	106.55	122.67
1	B	369	ASN	CB-CA-C	-7.45	99.77	111.17
1	A	264	ILE	CG1-CB-CG2	6.22	129.37	110.70
1	B	493	LYS	CB-CA-C	-5.75	101.45	110.94
1	A	128	GLN	CA-CB-CG	5.56	125.22	114.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	674	PRO	O-C-N	5.48	123.83	121.31
1	A	208	ASP	CB-CA-C	5.14	119.34	110.45

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	128	GLN	Sidechain
1	A	290	SER	Peptide
1	B	444	LYS	Peptide

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	4653	0	4462	84	0
1	B	4574	0	4385	82	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	31	0	13	3	0
3	B	31	0	13	2	0
4	A	4	0	6	2	0
4	B	8	0	12	0	0
5	A	66	0	0	0	0
5	B	61	0	0	4	0
All	All	9432	0	8891	161	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 9.

All (161) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:264:ILE:HG22	1:A:320:VAL:HB	1.43	1.00
1:B:506:ILE:HD11	3:B:803:ANP:H2'	1.40	1.00

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:288:SER:OG	1:A:291:GLU:N	2.06	0.89
1:A:264:ILE:CG2	1:A:320:VAL:HB	2.03	0.86
1:A:146:ILE:HB	1:A:196:VAL:HG12	1.59	0.81
1:A:163:VAL:HG23	1:A:198:THR:HA	1.62	0.80
1:A:145:ILE:HG13	1:A:196:VAL:HG13	1.64	0.79
1:B:82:GLU:OE2	5:B:901:HOH:O	2.02	0.77
1:A:288:SER:HG	1:A:291:GLU:H	1.37	0.72
1:B:443:PHE:HB2	1:B:450:TYR:HB2	1.72	0.70
1:A:295:PHE:HZ	1:A:298:THR:HG23	1.57	0.69
1:A:184:LEU:HD23	1:B:579:MET:HE2	1.75	0.68
1:A:153:SER:HB2	1:A:208:ASP:OD2	1.95	0.66
1:B:124:LEU:HD12	1:B:124:LEU:H	1.61	0.66
1:A:429:GLN:HG3	1:A:439:LEU:HD22	1.78	0.66
1:B:412:ILE:HD13	1:B:421:ILE:HG13	1.77	0.65
1:B:213:GLN:OE1	5:B:901:HOH:O	2.15	0.65
1:B:407:LEU:HD11	1:B:451:MET:HE2	1.78	0.64
1:B:231:LEU:HD11	1:B:252:LEU:HD11	1.80	0.64
1:B:397:GLU:C	1:B:399:SER:H	2.08	0.60
1:B:397:GLU:O	1:B:399:SER:N	2.29	0.60
1:A:109:ILE:CG2	1:A:131:VAL:HG12	2.32	0.59
1:B:240:LEU:HG	1:B:245:LEU:HD11	1.84	0.59
1:A:208:ASP:OD2	1:A:209:ARG:N	2.34	0.59
1:A:264:ILE:CG2	1:A:320:VAL:CB	2.79	0.58
1:B:463:THR:HG23	1:B:466:ARG:HH21	1.68	0.58
1:A:75:LYS:HE3	1:B:462:TRP:CE3	2.39	0.58
1:A:506:ILE:HD11	3:A:803:ANP:H2'	1.86	0.58
1:A:153:SER:O	1:A:208:ASP:OD2	2.21	0.57
1:B:134:MET:HB3	1:B:207:ILE:HB	1.87	0.57
1:A:409:LYS:NZ	1:A:445:ASP:O	2.30	0.57
1:B:427:ILE:HG21	1:B:495:ILE:HG21	1.87	0.57
1:B:456:CYS:HB2	1:B:506:ILE:CG2	2.34	0.56
1:A:128:GLN:HA	1:A:131:VAL:CG2	2.35	0.56
1:B:237:PHE:O	1:B:239:SER:N	2.30	0.56
1:B:158:MET:HE1	1:B:163:VAL:HG13	1.88	0.56
1:A:161:GLY:HA3	1:A:202:VAL:HG12	1.88	0.56
1:B:189:ASN:OD1	1:B:209:ARG:NH2	2.37	0.55
1:B:227:TYR:O	1:B:231:LEU:HD12	2.05	0.55
1:B:158:MET:HG3	1:B:177:PRO:HA	1.88	0.55
1:A:88:ILE:HD12	1:A:88:ILE:H	1.72	0.55
1:A:543:GLU:OE2	1:A:617:ARG:NH2	2.40	0.54
1:A:251:VAL:HG23	1:A:331:ILE:HD13	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:123:GLU:HB2	1:A:126:GLN:HG3	1.91	0.52
1:B:506:ILE:HD11	3:B:803:ANP:C2'	2.27	0.52
1:B:288:SER:HB2	1:B:291:GLU:H	1.74	0.52
1:A:520:PHE:CE2	1:B:78:ALA:HB1	2.45	0.52
1:B:209:ARG:C	1:B:211:CYS:H	2.18	0.52
1:A:183:GLU:O	1:A:186:ILE:HG12	2.11	0.51
1:A:611:ARG:HE	4:A:804:EDO:H21	1.75	0.51
1:B:231:LEU:HD23	1:B:237:PHE:CE2	2.46	0.51
1:B:387:GLY:HA3	1:B:406:ILE:O	2.11	0.51
1:B:158:MET:HB3	1:B:204:LEU:HD23	1.93	0.51
1:B:240:LEU:HG	1:B:245:LEU:CD1	2.42	0.50
1:A:143:SER:N	1:A:198:THR:OG1	2.44	0.50
1:A:243:GLU:HG2	1:A:244:ILE:N	2.25	0.50
1:B:378:ILE:HD11	1:B:393:GLN:HB3	1.93	0.50
1:A:425:LYS:O	1:A:429:GLN:HB2	2.12	0.50
1:A:146:ILE:HB	1:A:196:VAL:CG1	2.38	0.50
1:A:184:LEU:CD1	1:A:190:CYS:HB3	2.42	0.49
1:A:607:LYS:O	1:A:611:ARG:NH1	2.46	0.49
1:B:441:ARG:HG2	1:B:442:THR:H	1.76	0.49
1:B:614:PRO:HA	1:B:617:ARG:HG3	1.94	0.49
1:B:97:PRO:HD2	1:B:133:CYS:SG	2.53	0.49
1:B:460:GLU:HA	1:B:506:ILE:HD13	1.95	0.49
1:B:264:ILE:HD12	1:B:265:ARG:H	1.78	0.48
1:A:243:GLU:O	1:A:244:ILE:HB	2.12	0.48
1:B:609:LEU:O	1:B:617:ARG:HD3	2.14	0.48
1:A:127:ILE:O	1:A:131:VAL:HG22	2.13	0.47
1:A:240:LEU:HD12	1:A:244:ILE:HG22	1.96	0.47
1:A:158:MET:HE1	1:A:162:LYS:HA	1.95	0.47
1:A:428:MET:HB3	1:A:439:LEU:HB2	1.96	0.47
1:A:128:GLN:HA	1:A:131:VAL:HG22	1.94	0.47
1:B:456:CYS:HB2	1:B:506:ILE:HG21	1.95	0.47
1:B:186:ILE:HA	1:B:209:ARG:HB2	1.96	0.47
1:B:397:GLU:C	1:B:399:SER:N	2.73	0.47
1:A:611:ARG:O	1:A:617:ARG:HD3	2.14	0.47
1:A:614:PRO:HA	1:A:617:ARG:HG3	1.95	0.47
1:B:309:ALA:HA	1:B:312:GLY:O	2.15	0.47
1:A:158:MET:HG2	1:A:204:LEU:HD23	1.96	0.47
1:B:231:LEU:C	1:B:233:SER:H	2.23	0.47
1:A:405:LYS:HB3	1:A:451:MET:HB2	1.96	0.47
1:B:456:CYS:HB2	1:B:506:ILE:HG22	1.96	0.47
1:A:109:ILE:HG22	1:A:131:VAL:HG12	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:295:PHE:CZ	1:A:298:THR:HG23	2.45	0.47
1:B:158:MET:HG2	5:B:906:HOH:O	2.14	0.47
1:A:506:ILE:CD1	3:A:803:ANP:H2'	2.45	0.46
1:A:543:GLU:CD	1:A:617:ARG:HH22	2.23	0.46
1:B:183:GLU:O	1:B:186:ILE:HG12	2.15	0.46
1:B:210:GLN:HA	1:B:213:GLN:HE21	1.80	0.46
1:B:466:ARG:NH1	5:B:911:HOH:O	2.42	0.46
1:A:90:ASP:O	1:A:94:VAL:HG13	2.16	0.46
1:A:128:GLN:OE1	1:A:131:VAL:CG2	2.64	0.46
1:A:237:PHE:HA	1:A:240:LEU:HD21	1.98	0.46
1:A:462:TRP:HZ2	1:B:74:THR:HB	1.81	0.46
1:A:539:TYR:OH	1:A:565:GLU:OE1	2.28	0.46
1:B:522:LYS:NZ	1:B:532:TPO:O1P	2.48	0.46
1:A:285:ARG:NH2	1:A:287:ASP:OD1	2.49	0.46
1:A:425:LYS:NZ	1:A:429:GLN:OE1	2.33	0.46
1:B:510:ARG:O	1:B:649:PRO:HG2	2.16	0.46
1:A:247:LYS:O	1:A:251:VAL:HG13	2.16	0.45
1:A:248:LEU:HA	1:A:251:VAL:HG22	1.99	0.45
1:A:252:LEU:HD13	1:A:329:LEU:HD13	1.97	0.45
1:A:264:ILE:HD12	1:A:264:ILE:HA	1.63	0.45
1:B:460:GLU:OE1	1:B:462:TRP:HB3	2.17	0.45
1:A:128:GLN:OE1	1:A:131:VAL:HG21	2.16	0.45
1:A:130:ILE:O	1:A:134:MET:HG3	2.16	0.45
1:B:224:HIS:CE1	1:B:228:MET:HE2	2.51	0.45
1:B:371:LYS:C	1:B:373:SER:H	2.25	0.45
1:B:496:ILE:HD12	1:B:551:ASP:O	2.17	0.44
1:A:579:MET:HE3	1:B:184:LEU:HA	2.00	0.44
1:B:381:LEU:HD21	1:B:391:LEU:HB2	1.98	0.44
1:B:394:LEU:HD23	1:B:394:LEU:HA	1.81	0.44
1:B:163:VAL:HG12	1:B:198:THR:HA	2.00	0.44
1:B:583:ASN:O	1:B:587:ARG:HG3	2.18	0.44
1:B:378:ILE:HD11	1:B:393:GLN:CB	2.48	0.44
1:A:161:GLY:O	1:A:177:PRO:HD3	2.18	0.43
1:B:127:ILE:O	1:B:131:VAL:HG23	2.17	0.43
1:A:221:LEU:O	1:A:225:THR:HG23	2.18	0.43
1:A:157:VAL:HG22	1:A:180:VAL:HG22	1.99	0.43
1:B:522:LYS:HG2	1:B:524:ILE:HG23	2.00	0.43
1:A:579:MET:HE2	1:B:184:LEU:HD13	1.99	0.43
1:B:240:LEU:HB2	1:B:241:PRO:HD2	2.01	0.43
1:B:99:TYR:CE2	1:B:135:TYR:HB3	2.53	0.43
1:B:102:SER:HG	1:B:105:SER:HG	1.67	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:268:ALA:O	1:B:316:ARG:HG2	2.19	0.43
1:B:367:PHE:HA	1:B:370:LEU:HD23	2.00	0.43
1:A:425:LYS:HE2	1:A:685:ASP:O	2.19	0.43
1:B:241:PRO:O	1:B:245:LEU:HD12	2.19	0.43
1:B:437:VAL:HG21	1:B:516:VAL:HG12	2.01	0.43
1:B:240:LEU:HD12	1:B:244:ILE:HG21	2.00	0.42
1:A:241:PRO:O	1:A:245:LEU:HB2	2.18	0.42
1:A:612:ASP:HB2	4:A:804:EDO:H12	2.00	0.42
1:A:131:VAL:HG22	1:A:131:VAL:H	1.54	0.42
1:A:376:ASN:ND2	1:A:395:LYS:HE2	2.34	0.42
1:A:137:VAL:HG13	1:A:204:LEU:HB2	2.02	0.42
1:B:411:HIS:O	1:B:415:THR:HG23	2.20	0.42
1:A:184:LEU:HD12	1:A:190:CYS:HB3	2.02	0.42
1:A:483:VAL:HG22	1:A:505:LEU:HD21	2.02	0.42
1:B:101:LYS:NZ	1:B:131:VAL:O	2.49	0.42
1:A:228:MET:HE2	1:A:228:MET:HB3	1.98	0.41
1:B:480:THR:O	1:B:484:VAL:HG23	2.20	0.41
1:B:391:LEU:HD11	1:B:401:THR:HB	2.02	0.41
1:A:609:LEU:O	1:A:617:ARG:HD2	2.19	0.41
1:B:101:LYS:HG2	1:B:205:TRP:CH2	2.56	0.41
1:A:672:ASP:N	1:A:672:ASP:OD1	2.53	0.41
1:B:281:VAL:HG12	1:B:322:ALA:HA	2.03	0.41
1:B:404:MET:HE3	1:B:450:TYR:HD2	1.86	0.41
1:A:146:ILE:O	1:A:195:THR:HG23	2.20	0.41
1:A:264:ILE:HD11	1:A:316:ARG:HG3	2.02	0.41
1:B:221:LEU:O	1:B:225:THR:HG23	2.21	0.41
1:B:553:SER:HA	1:B:556:TYR:CD2	2.56	0.41
1:A:139:TYR:CE2	1:A:145:ILE:HG22	2.56	0.41
1:A:324:GLU:O	1:A:326:VAL:HG23	2.21	0.41
1:A:553:SER:HA	1:A:556:TYR:CD2	2.56	0.41
1:B:400:LYS:HD3	1:B:400:LYS:HA	1.84	0.41
1:A:477:ARG:HA	1:A:633:TRP:CH2	2.56	0.40
1:A:243:GLU:HG2	1:A:244:ILE:H	1.87	0.40
3:A:803:ANP:H8	3:A:803:ANP:O5'	2.21	0.40
1:B:405:LYS:HB2	1:B:453:MET:HE2	2.04	0.40
1:A:235:PRO:HD3	1:A:304:TRP:CZ2	2.56	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	596/617 (97%)	566 (95%)	28 (5%)	2 (0%)	36	49
1	B	587/617 (95%)	551 (94%)	29 (5%)	7 (1%)	10	15
All	All	1183/1234 (96%)	1117 (94%)	57 (5%)	9 (1%)	16	23

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	94	VAL
1	B	398	GLU
1	B	308	LYS
1	A	172	LEU
1	A	301	LYS
1	B	579	MET
1	B	75	LYS
1	B	210	GLN
1	B	340	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	484/535 (90%)	484 (100%)	0	100	100
1	B	474/535 (89%)	474 (100%)	0	100	100
All	All	958/1070 (90%)	958 (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 (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	213	GLN
1	A	224	HIS
1	A	369	ASN
1	A	547	ASN
1	B	128	GLN
1	B	213	GLN
1	B	282	ASN
1	B	432	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
1	TPO	B	532	1	8,10,11	1.63	1 (12%)	10,14,16	2.25	2 (20%)
1	TPO	A	532	1	8,10,11	1.59	1 (12%)	10,14,16	2.12	2 (20%)

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
1	TPO	B	532	1	-	0/9/11/13	-
1	TPO	A	532	1	-	0/9/11/13	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	532	TPO	P-O1P	3.47	1.61	1.50
1	A	532	TPO	P-O1P	3.33	1.60	1.50

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	532	TPO	P-OG1-CB	-6.40	105.94	123.33
1	A	532	TPO	P-OG1-CB	-5.86	107.40	123.33
1	B	532	TPO	CG2-CB-CA	-2.34	108.71	113.26
1	A	532	TPO	CG2-CB-CA	-2.26	108.85	113.26

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	B	532	TPO	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 4 are monoatomic - leaving 5 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	$\# Z > 2$
4	EDO	B	804	-	3,3,3	0.44	0	2,2,2	0.39	0
4	EDO	B	805	-	3,3,3	0.43	0	2,2,2	0.38	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	ANP	B	803	2	33,33,33	3.24	11 (33%)	45,52,52	2.11	10 (22%)
3	ANP	A	803	2	33,33,33	3.11	11 (33%)	45,52,52	1.97	12 (26%)
4	EDO	A	804	-	3,3,3	0.44	0	2,2,2	0.32	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	EDO	B	804	-	-	1/1/1/1	-
4	EDO	B	805	-	-	0/1/1/1	-
3	ANP	B	803	2	-	2/18/38/38	0/3/3/3
3	ANP	A	803	2	-	1/18/38/38	0/3/3/3
4	EDO	A	804	-	-	1/1/1/1	-

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	803	ANP	PB-O3A	12.33	1.74	1.59
3	A	803	ANP	PB-O3A	11.17	1.73	1.59
3	A	803	ANP	PG-N3B	6.05	1.79	1.63
3	B	803	ANP	PG-O1G	5.68	1.54	1.46
3	B	803	ANP	PA-O5'	5.67	1.81	1.59
3	A	803	ANP	PA-O3A	-5.58	1.53	1.59
3	B	803	ANP	PG-N3B	5.58	1.77	1.63
3	A	803	ANP	PG-O1G	5.48	1.54	1.46
3	B	803	ANP	PA-O3A	-5.42	1.53	1.59
3	A	803	ANP	PA-O5'	5.30	1.80	1.59
3	B	803	ANP	C5-C4	3.37	1.45	1.39
3	A	803	ANP	C5-C4	3.19	1.44	1.39
3	A	803	ANP	PB-O2B	-2.98	1.48	1.56
3	B	803	ANP	O5'-C5'	-2.80	1.34	1.44
3	B	803	ANP	PB-O2B	-2.57	1.50	1.56
3	A	803	ANP	O5'-C5'	-2.46	1.35	1.44
3	B	803	ANP	PG-O2G	-2.41	1.50	1.56
3	A	803	ANP	PG-O2G	-2.38	1.50	1.56
3	A	803	ANP	PG-O3G	-2.32	1.50	1.56
3	A	803	ANP	C4-N3	2.13	1.38	1.34
3	B	803	ANP	C4-N3	2.12	1.38	1.34
3	B	803	ANP	PG-O3G	-2.07	1.51	1.56

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	803	ANP	O1G-PG-N3B	-8.68	98.99	111.77
3	A	803	ANP	O1G-PG-N3B	-6.45	102.28	111.77
3	B	803	ANP	O2B-PB-O1B	4.72	120.00	109.87
3	A	803	ANP	O2B-PB-O1B	4.68	119.90	109.87
3	A	803	ANP	O2A-PA-O3A	3.99	118.05	107.27
3	A	803	ANP	O2G-PG-O3G	3.74	117.63	107.59
3	B	803	ANP	O2A-PA-O3A	3.67	117.19	107.27
3	B	803	ANP	O2G-PG-O3G	3.65	117.41	107.59
3	B	803	ANP	N3-C2-N1	3.45	133.81	128.58
3	B	803	ANP	C2-N1-C6	-3.42	113.12	118.73
3	A	803	ANP	C2-N1-C6	-3.39	113.16	118.73
3	A	803	ANP	N3-C2-N1	3.36	133.67	128.58
3	A	803	ANP	O3A-PB-N3B	-2.65	99.25	106.59
3	B	803	ANP	O5'-PA-O1A	-2.43	99.31	108.94
3	A	803	ANP	O1B-PB-N3B	2.38	115.28	111.77
3	A	803	ANP	N3-C4-N9	2.35	131.16	127.17
3	B	803	ANP	N3-C4-N9	2.34	131.15	127.17
3	B	803	ANP	O3A-PA-O1A	2.34	117.74	110.70
3	A	803	ANP	O3A-PA-O1A	2.16	117.21	110.70
3	A	803	ANP	O2A-PA-O5'	-2.15	97.80	107.57
3	A	803	ANP	O5'-PA-O1A	-2.14	100.46	108.94
3	B	803	ANP	O2A-PA-O5'	-2.11	97.99	107.57

There are no chirality outliers.

All (5) torsion outliers are listed below:

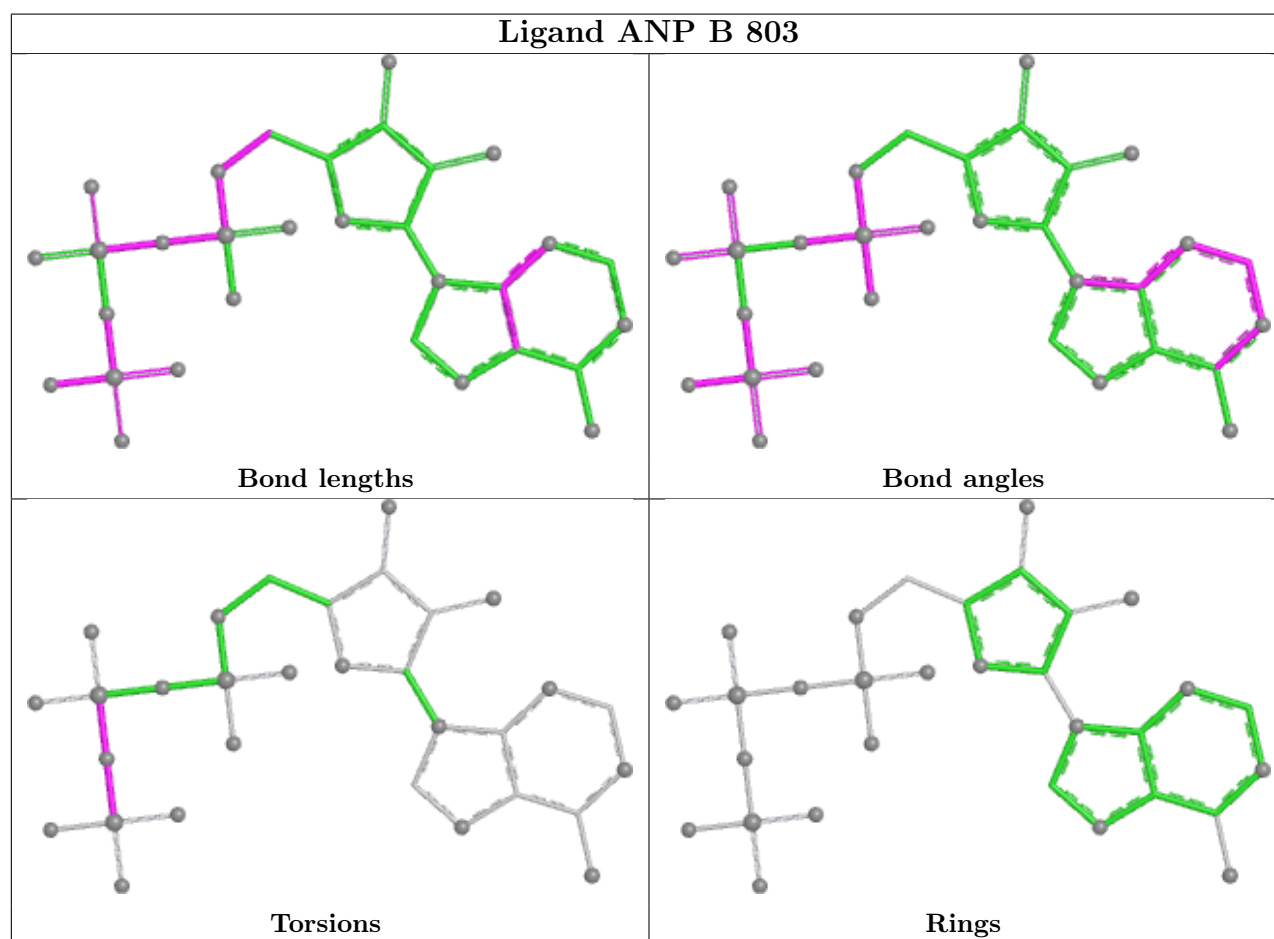
Mol	Chain	Res	Type	Atoms
3	A	803	ANP	PG-N3B-PB-O1B
3	B	803	ANP	PB-N3B-PG-O1G
3	B	803	ANP	PG-N3B-PB-O1B
4	A	804	EDO	O1-C1-C2-O2
4	B	804	EDO	O1-C1-C2-O2

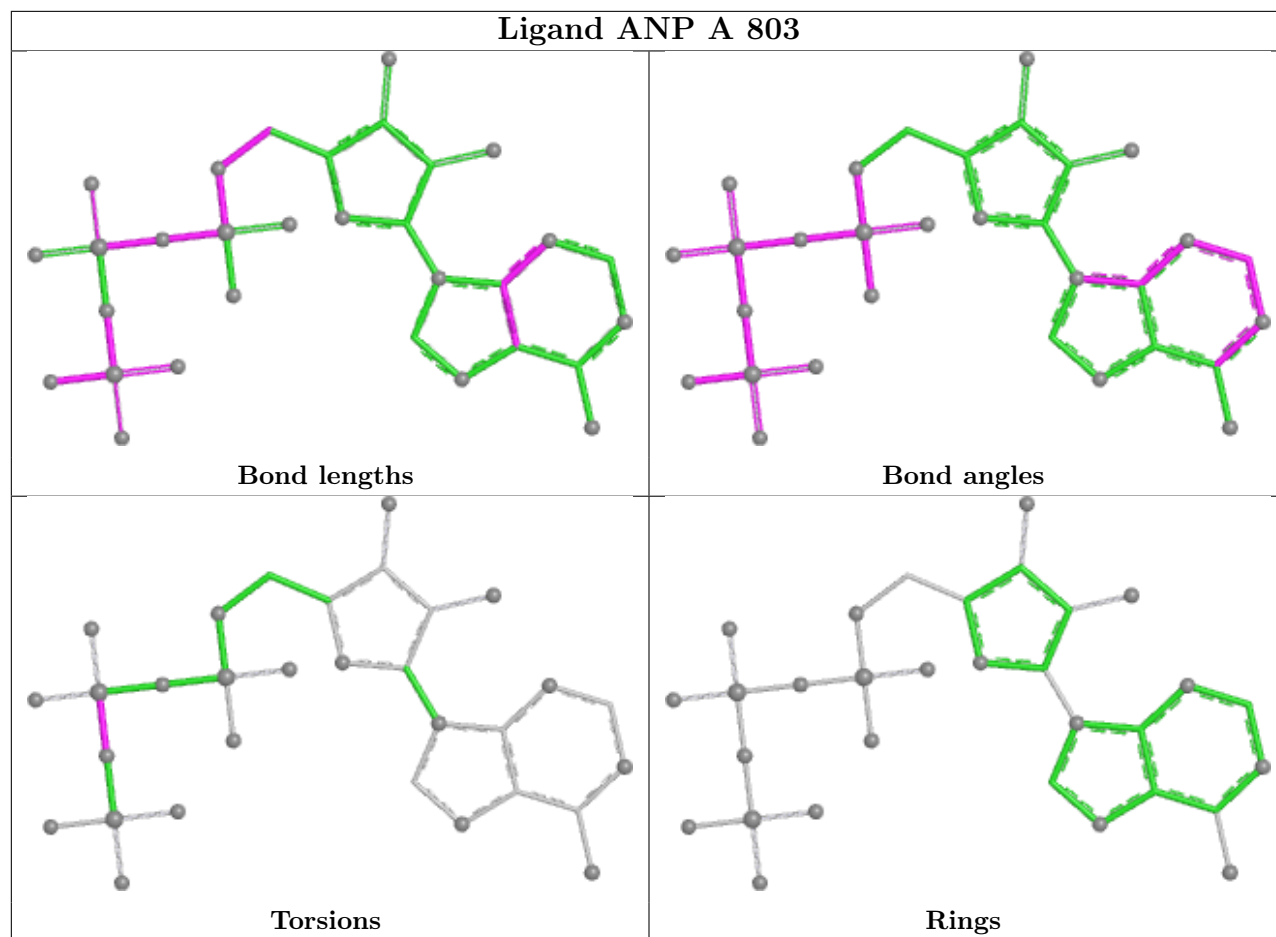
There are no ring outliers.

3 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	803	ANP	2	0
3	A	803	ANP	3	0
4	A	804	EDO	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	600/617 (97%)	0.22	22 (3%)	45 41	46, 82, 139, 173	0
1	B	593/617 (96%)	0.37	32 (5%)	31 28	48, 84, 144, 187	0
All	All	1193/1234 (96%)	0.30	54 (4%)	38 34	46, 83, 142, 187	0

All (54) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	386	PHE	4.7
1	B	84	THR	4.6
1	B	360	TYR	3.8
1	B	91	LEU	3.8
1	B	95	THR	3.7
1	A	86	PHE	3.7
1	B	375	PHE	3.6
1	B	684	ILE	3.6
1	A	340	ILE	3.6
1	B	338	HIS	3.4
1	B	339	LEU	3.3
1	A	525	GLY	3.3
1	A	188	TYR	3.2
1	B	273	PHE	3.2
1	A	354	ALA	3.1
1	A	237	PHE	3.1
1	B	372	LEU	3.1
1	B	686	PHE	3.0
1	B	310	LEU	2.9
1	B	672	ASP	2.9
1	A	338	HIS	2.9
1	A	190	CYS	2.9
1	B	275	ILE	2.8
1	A	671	ASN	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	336	PHE	2.8
1	A	85	ALA	2.8
1	B	341	GLY	2.7
1	A	360	TYR	2.7
1	B	387	GLY	2.7
1	B	93	HIS	2.6
1	B	368	ALA	2.6
1	B	237	PHE	2.6
1	B	671	ASN	2.5
1	B	358	ALA	2.5
1	A	144	CYS	2.4
1	A	355	GLU	2.4
1	B	264	ILE	2.4
1	B	340	ILE	2.4
1	A	92	SER	2.3
1	A	336	PHE	2.3
1	A	386	PHE	2.3
1	A	170	VAL	2.3
1	B	74	THR	2.3
1	B	290	SER	2.3
1	B	404	MET	2.2
1	A	73	ARG	2.2
1	A	208	ASP	2.1
1	B	367	PHE	2.1
1	B	668	PRO	2.1
1	A	310	LEU	2.0
1	B	525	GLY	2.0
1	A	264	ILE	2.0
1	B	228	MET	2.0
1	A	334	ASP	2.0

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

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
1	TPO	A	532	11/12	0.98	0.06	55,61,64,68	0
1	TPO	B	532	11/12	0.98	0.06	68,70,77,77	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

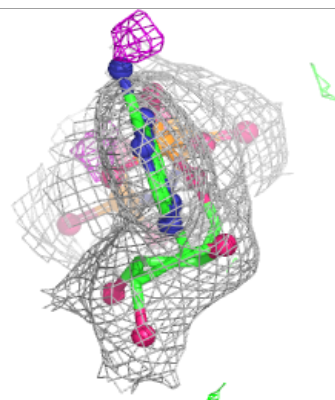
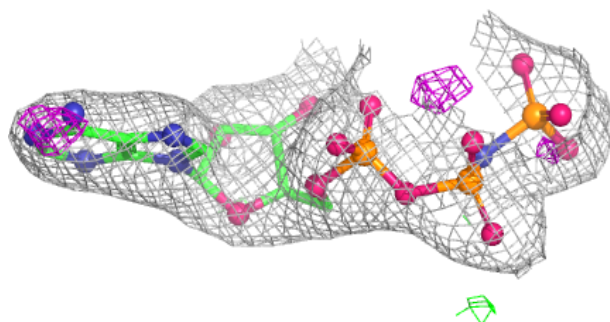
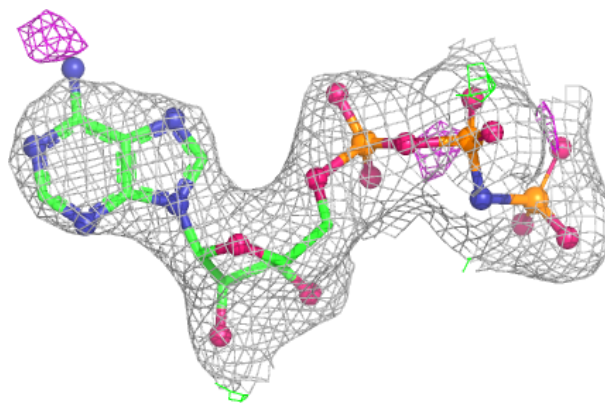
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	EDO	B	805	4/4	0.89	0.14	86,87,88,89	0
4	EDO	A	804	4/4	0.91	0.13	92,96,96,98	0
4	EDO	B	804	4/4	0.95	0.12	85,87,90,90	0
3	ANP	B	803	31/31	0.96	0.07	75,88,93,102	0
2	MN	B	802	1/1	0.98	0.06	74,74,74,74	0
3	ANP	A	803	31/31	0.98	0.06	56,61,67,70	0
2	MN	B	801	1/1	0.98	0.06	81,81,81,81	0
2	MN	A	802	1/1	1.00	0.03	64,64,64,64	0
2	MN	A	801	1/1	1.00	0.02	65,65,65,65	0

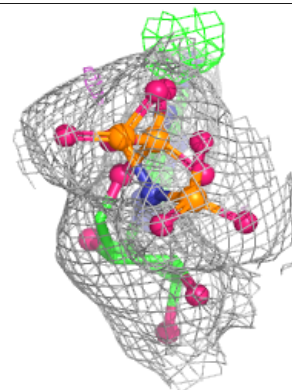
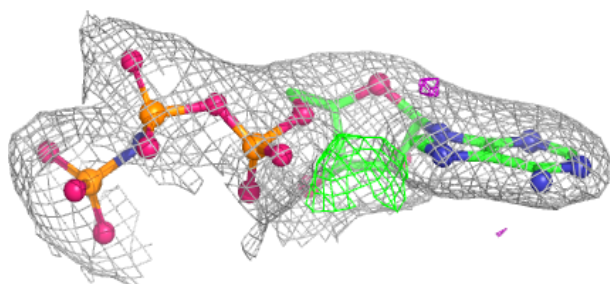
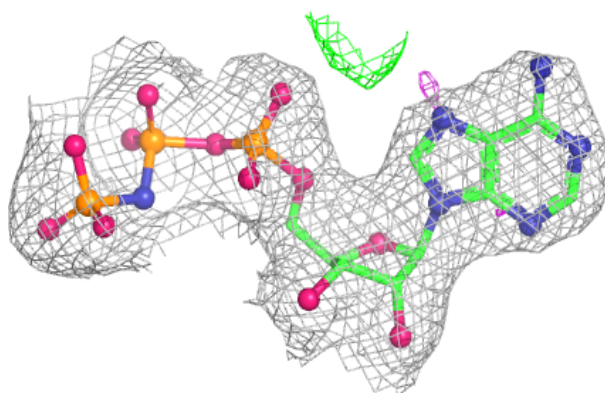
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around ANP B 803:

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

**Electron density around ANP A 803:**

$2mF_o-DF_c$ (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.