



# Full wwPDB X-ray Structure Validation Report ⓘ

May 29, 2020 – 05:51 am BST

PDB ID : 1F0L  
Title : 1.55 ANGSTROM CRYSTAL STRUCTURE OF WILD TYPE DIPHTHERIA TOXIN  
Authors : Steere, B.; Eisenberg, D.  
Deposited on : 2000-05-16  
Resolution : 1.55 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](mailto: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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.11

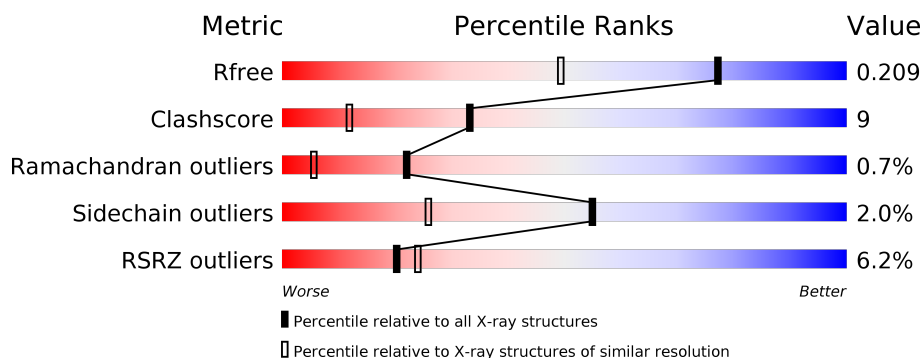
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.55 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))
$R_{free}$	130704	1483 (1.56-1.56)
Clashscore	141614	1529 (1.56-1.56)
Ramachandran outliers	138981	1498 (1.56-1.56)
Sidechain outliers	138945	1495 (1.56-1.56)
RSRZ outliers	127900	1465 (1.56-1.56)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	535	<div> <div>7%</div> <div> <div></div> <div>83%</div> <div>14%</div> <div>.</div> </div> </div>
1	B	535	<div> <div>5%</div> <div> <div></div> <div>85%</div> <div>11%</div> <div>..</div> </div> </div>

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 10204 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.

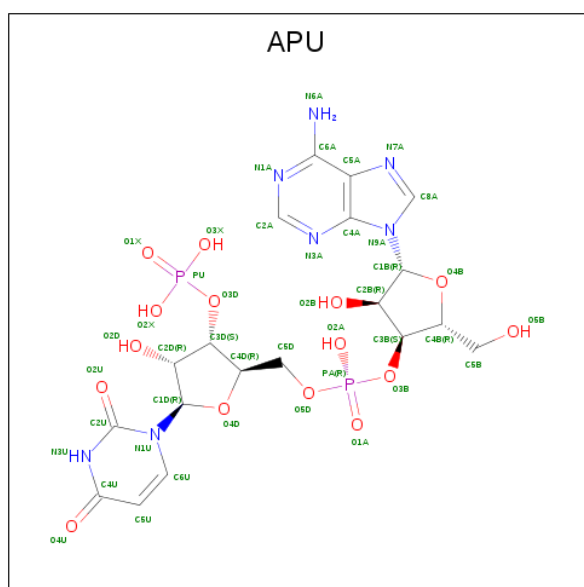
- Molecule 1 is a protein called DIPHTHERIA TOXIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	520	Total	C	N	O	S	0	35	0
			4112	2596	692	808	16			
1	B	522	Total	C	N	O	S	0	43	0
			4188	2636	707	830	15			

- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	5	Total	Cl	0	0
			5	5		
2	A	4	Total	Cl	0	0
			4	4		

- Molecule 3 is ADENYLYL-3'-5'-PHOSPHO-URIDINE-3'-MONOPHOSPHATE (three-letter code: APU) (formula: C<sub>19</sub>H<sub>25</sub>N<sub>7</sub>O<sub>15</sub>P<sub>2</sub>).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 43	C 19	N 7	O 15	P 2	0	0
3	B	1	Total 43	C 19	N 7	O 15	P 2	0	0

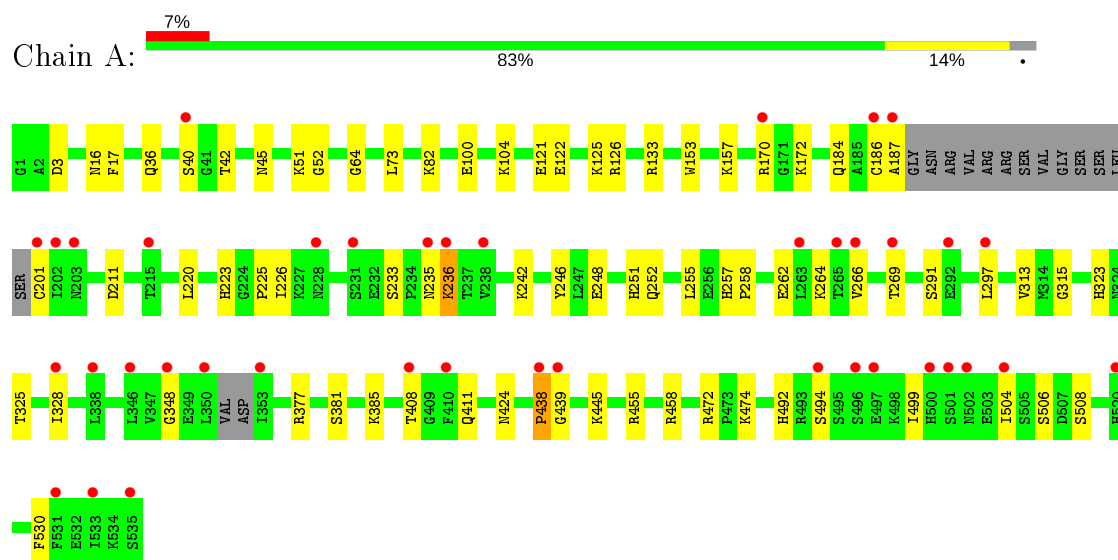
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	910	Total 910	O 910	0	0
4	B	899	Total 899	O 899	0	0

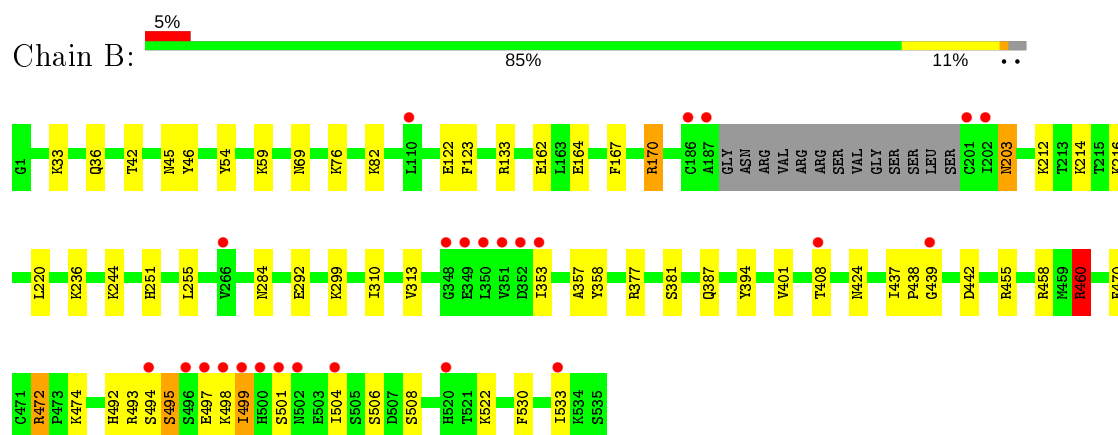
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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: DIPHTHERIA TOXIN



#### • Molecule 1: DIPHTHERIA TOXIN



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	167.77 Å   134.88 Å   46.87 Å 90.00°   90.00°   90.00°	Depositor
Resolution (Å)	40.00 – 1.55 37.51 – 1.55	Depositor EDS
% Data completeness (in resolution range)	97.0 (40.00-1.55) 94.6 (37.51-1.55)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.37 (at 1.55 Å)	Xtriage
Refinement program	SHELXL, SHELXL-97	Depositor
R, $R_{free}$	0.188   ,   0.238 0.190   ,   0.209	Depositor DCC
$R_{free}$ test set	7212 reflections (4.83%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	18.3	Xtriage
Anisotropy	0.263	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 80.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	10204	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 31.11 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.1698e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality ⓘ

### 5.1 Standard geometry ⓘ

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

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	0.43	0/4328	1.05	7/5853 (0.1%)
1	B	0.44	0/4425	1.10	15/5983 (0.3%)
All	All	0.44	0/8753	1.08	22/11836 (0.2%)

There are no bond length outliers.

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	460	ARG	NE-CZ-NH2	14.50	127.55	120.30
1	B	455	ARG	NE-CZ-NH1	-9.57	115.51	120.30
1	B	460	ARG	CD-NE-CZ	8.44	135.42	123.60
1	B	455	ARG	NE-CZ-NH2	-8.09	116.25	120.30
1	B	455	ARG	NH1-CZ-NH2	8.02	128.22	119.40
1	B	377	ARG	NE-CZ-NH1	7.96	124.28	120.30
1	A	458	ARG	NE-CZ-NH2	-7.54	116.53	120.30
1	B	54	TYR	CB-CG-CD2	-7.24	116.65	121.00
1	B	493	ARG	NE-CZ-NH2	-6.92	116.84	120.30
1	A	377	ARG	NE-CZ-NH1	6.82	123.71	120.30
1	B	54	TYR	CB-CG-CD1	6.33	124.80	121.00
1	B	458	ARG	NE-CZ-NH2	-6.24	117.18	120.30
1	A	126	ARG	CD-NE-CZ	6.22	132.31	123.60
1	A	17	PHE	CB-CG-CD2	-5.79	116.75	120.80
1	B	358	TYR	CG-CD1-CE1	5.76	125.91	121.30
1	A	126	ARG	NE-CZ-NH1	5.75	123.17	120.30
1	A	170	ARG	NE-CZ-NH1	5.72	123.16	120.30
1	A	377	ARG	NE-CZ-NH2	-5.55	117.52	120.30
1	B	394	TYR	CB-CG-CD2	5.40	124.24	121.00
1	B	472[A]	ARG	NE-CZ-NH1	5.11	122.85	120.30
1	B	472[B]	ARG	NE-CZ-NH1	5.11	122.85	120.30
1	B	123	PHE	CB-CG-CD1	5.08	124.36	120.80

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	4112	0	4023	80	0
1	B	4188	0	4086	64	0
2	A	4	0	0	0	0
2	B	5	0	0	0	0
3	A	43	0	22	0	0
3	B	43	0	22	0	0
4	A	910	0	0	56	0
4	B	899	0	0	40	0
All	All	10204	0	8153	144	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 (144) 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:313[B]:VAL:HG23	4:A:1521:HOH:O	1.27	1.28
1:A:122[A]:GLU:CD	4:A:850:HOH:O	1.66	1.27
1:B:284[B]:ASN:ND2	4:B:2253:HOH:O	1.68	1.23
1:A:122[B]:GLU:OE1	4:A:851:HOH:O	1.52	1.21
1:A:122[A]:GLU:CG	4:A:850:HOH:O	1.91	1.14
1:B:36[A]:GLN:NE2	4:B:1720:HOH:O	1.79	1.13
1:A:408[A]:THR:HG22	4:A:1595:HOH:O	1.49	1.10
1:A:313[B]:VAL:HG22	4:A:1691:HOH:O	1.52	1.08
1:B:162[A]:GLU:OE1	4:B:2566:HOH:O	1.70	1.08
1:A:36[B]:GLN:OE1	4:A:1094:HOH:O	1.71	1.07
1:B:353[B]:ILE:HB	4:B:1811:HOH:O	1.53	1.05
1:B:353[B]:ILE:CB	4:B:1811:HOH:O	2.03	1.03
1:A:122[A]:GLU:HG2	4:A:850:HOH:O	1.53	1.03
1:A:408[B]:THR:HG23	4:A:1594:HOH:O	1.61	1.01
1:B:442[B]:ASP:CG	4:B:2348:HOH:O	1.96	1.01

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:408[A]:THR:CG2	4:A:1595:HOH:O	2.02	1.00
1:A:125[A]:LYS:HD2	4:A:847:HOH:O	1.60	0.98
1:A:122[A]:GLU:OE2	4:A:849:HOH:O	1.79	0.97
1:A:313[B]:VAL:CG2	4:A:1691:HOH:O	2.08	0.96
1:B:313[A]:VAL:HG23	4:B:2408:HOH:O	1.67	0.93
1:A:474[A]:LYS:NZ	4:A:1146:HOH:O	2.00	0.93
1:B:162[A]:GLU:CD	4:B:2566:HOH:O	2.07	0.91
1:B:442[A]:ASP:OD2	4:B:1809:HOH:O	1.94	0.86
1:B:442[A]:ASP:OD1	4:B:1808:HOH:O	1.92	0.86
1:B:442[B]:ASP:OD1	4:B:2348:HOH:O	1.88	0.85
1:A:36[B]:GLN:NE2	4:A:1292:HOH:O	2.08	0.84
1:A:211[A]:ASP:OD1	4:A:1584:HOH:O	1.98	0.82
1:B:82[A]:LYS:CE	4:B:2566:HOH:O	2.29	0.80
1:B:36[B]:GLN:HG2	4:B:1719:HOH:O	1.81	0.79
1:A:211[A]:ASP:OD1	4:A:812:HOH:O	2.02	0.77
1:B:212:LYS:HE2	4:B:1756:HOH:O	1.84	0.76
1:B:82[A]:LYS:HE2	4:B:2566:HOH:O	1.84	0.76
1:A:122[A]:GLU:HG3	4:A:849:HOH:O	1.88	0.72
1:B:313[A]:VAL:HG22	4:B:2407:HOH:O	1.89	0.72
1:A:122[A]:GLU:CG	4:A:849:HOH:O	2.38	0.70
1:B:408[A]:THR:HG21	4:B:2531:HOH:O	1.91	0.70
1:B:313[A]:VAL:CG2	4:B:2407:HOH:O	2.40	0.70
1:A:104:LYS:NZ	4:A:1649:HOH:O	2.25	0.69
1:B:164[A]:GLU:OE2	4:B:2303:HOH:O	2.10	0.69
1:B:460:ARG:NH1	1:B:474[A]:LYS:HA	2.09	0.67
1:A:172[A]:LYS:NZ	4:A:908:HOH:O	1.66	0.66
1:B:357:ALA:HB3	4:B:2537:HOH:O	1.97	0.64
1:B:408[B]:THR:HG23	4:B:2227:HOH:O	1.96	0.63
1:A:266:VAL:O	1:A:269:THR:HB	1.99	0.63
1:A:100[A]:GLU:CG	4:A:1496:HOH:O	2.46	0.62
1:A:438[A]:PRO:HB2	4:A:1624:HOH:O	1.98	0.62
1:A:104:LYS:HE3	1:A:411:GLN:OE1	1.99	0.62
1:A:220:LEU:HD23	1:A:257:HIS:CD2	2.35	0.61
1:A:445[C]:LYS:NZ	4:A:1273:HOH:O	2.32	0.61
1:B:460:ARG:NH1	1:B:474[B]:LYS:HA	2.16	0.61
1:B:76:LYS:HE2	4:B:2518:HOH:O	2.02	0.59
1:A:211[A]:ASP:CG	4:A:1584:HOH:O	2.39	0.59
1:A:313[B]:VAL:HG23	4:A:1691:HOH:O	1.91	0.58
1:A:223:HIS:CD2	1:A:226:ILE:HD12	2.38	0.58
1:A:251:HIS:CE1	1:A:255:LEU:HD22	2.39	0.58
1:A:408[A]:THR:HG22	4:A:1594:HOH:O	2.03	0.58

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:498:LYS:HA	4:B:2162:HOH:O	2.04	0.57
1:A:121:GLU:HG3	4:A:1339:HOH:O	2.03	0.56
1:B:442[B]:ASP:OD1	1:B:492:HIS:HB3	2.05	0.56
1:B:522:LYS:HE3	4:B:2151:HOH:O	2.05	0.56
1:B:251:HIS:CE1	1:B:255:LEU:HD22	2.41	0.55
1:A:40:SER:HA	4:A:1442:HOH:O	2.06	0.55
1:A:385:LYS:NZ	4:A:1660:HOH:O	2.38	0.55
1:A:125[A]:LYS:CD	4:A:846:HOH:O	2.54	0.54
1:B:33[B]:LYS:NZ	4:B:2128:HOH:O	2.40	0.54
1:A:408[A]:THR:HG23	1:A:499:ILE:O	2.08	0.54
1:B:310:ILE:O	1:B:313[B]:VAL:HG12	2.08	0.53
1:B:408[B]:THR:HG21	4:B:2531:HOH:O	2.07	0.53
1:B:82[A]:LYS:NZ	4:B:2566:HOH:O	2.42	0.52
1:A:408[B]:THR:CG2	4:A:1594:HOH:O	2.39	0.51
1:B:442[B]:ASP:OD1	1:B:492:HIS:O	2.29	0.50
1:B:504:ILE:HG12	1:B:506:SER:O	2.12	0.50
1:A:439[B]:GLY:O	1:A:494:SER:OG	2.30	0.49
1:B:439[A]:GLY:O	1:B:495:SER:HB2	2.12	0.49
1:A:100[A]:GLU:CD	4:A:1496:HOH:O	2.49	0.49
1:B:216:LYS:O	1:B:220:LEU:HG	2.11	0.49
1:A:408[B]:THR:OG1	1:A:499:ILE:O	2.29	0.49
1:A:125[A]:LYS:HD2	4:A:846:HOH:O	2.12	0.48
1:A:125[A]:LYS:HD3	4:A:846:HOH:O	2.14	0.48
1:B:437[B]:ILE:HD12	1:B:499:ILE:HD13	1.95	0.48
1:A:233:SER:HB2	1:A:246:TYR:CZ	2.49	0.48
1:A:121:GLU:HG2	4:A:1240:HOH:O	2.13	0.47
1:A:172[A]:LYS:HE2	4:A:1426:HOH:O	2.15	0.47
1:A:186[A]:CYS:O	1:A:187[A]:ALA:HB2	2.15	0.47
1:A:297:LEU:HD21	4:A:1663:HOH:O	2.14	0.47
1:B:214:LYS:HE2	4:B:2319:HOH:O	2.14	0.47
1:B:522:LYS:HE2	4:B:2500:HOH:O	2.14	0.46
1:B:244[A]:LYS:NZ	4:B:2237:HOH:O	2.49	0.46
1:A:348:GLY:HA3	4:A:1406:HOH:O	2.15	0.46
1:B:437[A]:ILE:HD12	1:B:499:ILE:HD13	1.96	0.46
1:B:442[A]:ASP:HB2	4:B:2348:HOH:O	2.15	0.46
1:B:494:SER:HB3	4:B:2348:HOH:O	2.16	0.46
1:A:184:GLN:NE2	4:A:1316:HOH:O	2.49	0.46
1:B:470:PHE:CZ	1:B:472[B]:ARG:HG3	2.51	0.45
1:A:64:GLY:HA2	4:A:1465:HOH:O	2.15	0.45
1:A:187[A]:ALA:HB3	1:A:201[A]:CYS:SG	2.56	0.45
1:A:223:HIS:CD2	1:A:225:PRO:HD2	2.50	0.45

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:46:TYR:OH	1:B:442[B]:ASP:OD2	2.30	0.45
1:B:495:SER:OG	1:B:497:GLU:HG3	2.17	0.45
1:A:122[A]:GLU:OE2	4:A:850:HOH:O	2.08	0.45
1:A:411:GLN:HG2	4:A:1646:HOH:O	2.16	0.45
1:A:82[A]:LYS:NZ	4:A:1325:HOH:O	2.49	0.45
1:A:16:ASN:ND2	4:A:1255:HOH:O	2.49	0.45
1:A:236:LYS:HE3	4:A:1679:HOH:O	2.16	0.44
1:B:59:LYS:NZ	4:B:2303:HOH:O	2.49	0.44
1:A:157:LYS:NZ	4:A:830:HOH:O	2.50	0.44
1:A:262:GLU:OE1	1:A:262:GLU:N	2.50	0.44
1:B:203:ASN:ND2	4:B:1932:HOH:O	2.50	0.44
1:A:439[A]:GLY:O	1:A:494:SER:OG	2.30	0.44
1:B:236:LYS:NZ	4:B:2475:HOH:O	2.50	0.44
1:A:252:GLN:NE2	4:A:1713:HOH:O	2.50	0.44
1:A:411:GLN:NE2	4:A:1645:HOH:O	2.50	0.44
1:B:36[B]:GLN:NE2	4:B:1720:HOH:O	2.50	0.44
1:B:284[B]:ASN:ND2	4:B:1973:HOH:O	2.49	0.43
1:A:242:LYS:NZ	4:A:1254:HOH:O	2.50	0.43
1:B:508:SER:HB3	1:B:530:PHE:CD1	2.53	0.43
1:B:387:GLN:HB3	4:B:2535:HOH:O	2.18	0.43
1:A:51[B]:LYS:NZ	4:A:1613:HOH:O	2.51	0.43
1:A:264:LYS:HE2	4:A:910:HOH:O	2.17	0.43
1:B:167:PHE:O	1:B:170:ARG:HB2	2.19	0.42
1:A:257:HIS:ND1	1:A:258[B]:PRO:HD2	2.34	0.42
1:A:492:HIS:HE1	4:A:1120:HOH:O	2.02	0.42
1:B:408[B]:THR:OG1	1:B:499:ILE:O	2.29	0.42
1:B:353[B]:ILE:CG2	4:B:1811:HOH:O	2.55	0.41
1:A:315:GLY:HA3	1:A:323:HIS:CD2	2.56	0.41
1:B:292:GLU:OE2	1:B:299:LYS:HE2	2.20	0.41
1:A:508:SER:HB3	1:A:530:PHE:CD1	2.56	0.41
1:B:401:VAL:HB	4:B:2377:HOH:O	2.20	0.41
1:A:328:ILE:HG23	4:A:1682:HOH:O	2.21	0.41
1:A:172[A]:LYS:CE	4:A:908:HOH:O	2.47	0.41
1:A:504:ILE:HG12	1:A:506:SER:O	2.21	0.41
1:A:52:GLY:HA2	1:A:153:TRP:CD1	2.56	0.40
1:A:472:ARG:HD3	4:A:1261:HOH:O	2.22	0.40
1:B:504:ILE:O	1:B:504:ILE:HG23	2.21	0.40
1:A:325:THR:OG1	1:A:328:ILE:HD13	2.21	0.40
1:A:73:LEU:HD23	1:A:474[B]:LYS:O	2.21	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	550/535 (103%)	534 (97%)	12 (2%)	4 (1%)	22	5
1	B	559/535 (104%)	543 (97%)	11 (2%)	5 (1%)	17	3
All	All	1109/1070 (104%)	1077 (97%)	23 (2%)	9 (1%)	22	4

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	381	SER
1	B	381	SER
1	B	438[A]	PRO
1	B	438[B]	PRO
1	B	501	SER
1	A	424	ASN
1	A	438[A]	PRO
1	A	438[B]	PRO
1	B	424	ASN

### 5.3.2 Protein sidechains [i](#)

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	464/452 (103%)	457 (98%)	7 (2%)	65	37
1	B	480/452 (106%)	470 (98%)	10 (2%)	53	24
All	All	944/904 (104%)	927 (98%)	17 (2%)	55	31

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	ASP
1	A	42	THR
1	A	45	ASN
1	A	133	ARG
1	A	235	ASN
1	A	236	LYS
1	A	291	SER
1	B	42	THR
1	B	45	ASN
1	B	69	ASN
1	B	133	ARG
1	B	170	ARG
1	B	203	ASN
1	B	460	ARG
1	B	495	SER
1	B	499	ILE
1	B	533	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	184	GLN
1	A	203	ASN
1	A	223	HIS
1	A	492	HIS
1	B	16	ASN
1	B	184	GLN
1	B	203	ASN
1	B	223	HIS
1	B	257	HIS
1	B	373	ASN
1	B	502	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 11 ligands modelled in this entry, 9 are monoatomic - leaving 2 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	$\# Z  > 2$
3	APU	B	602	-	40,47,47	1.36	5 (12%)	45,72,72	1.34	6 (13%)
3	APU	A	601	-	40,47,47	1.46	4 (10%)	45,72,72	1.41	5 (11%)

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
3	APU	B	602	-	-	4/20/58/58	0/5/5/5
3	APU	A	601	-	-	5/20/58/58	0/5/5/5

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	601	APU	C4U-N3U	4.99	1.41	1.33
3	B	602	APU	C4U-N3U	4.36	1.40	1.33
3	A	601	APU	C2D-C1D	-3.46	1.48	1.53
3	A	601	APU	C8A-N7A	-2.70	1.29	1.34
3	B	602	APU	C2D-C1D	-2.27	1.50	1.53
3	B	602	APU	O4B-C1B	2.26	1.44	1.41
3	B	602	APU	C8A-N7A	-2.18	1.30	1.34
3	B	602	APU	C6U-C5U	-2.04	1.33	1.38
3	A	601	APU	PU-O3D	2.01	1.63	1.59

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	601	APU	C5U-C4U-N3U	-4.58	113.23	123.31
3	B	602	APU	C5U-C4U-N3U	-4.39	113.64	123.31
3	A	601	APU	O4B-C4B-C5B	3.36	116.47	109.21
3	B	602	APU	O4D-C1D-C2D	-3.29	102.12	106.93
3	A	601	APU	O2B-C2B-C3B	2.96	119.56	111.17
3	B	602	APU	O2B-C2B-C3B	2.91	119.42	111.17
3	A	601	APU	O4D-C1D-C2D	-2.91	102.68	106.93
3	B	602	APU	O4B-C4B-C5B	2.78	115.23	109.21
3	A	601	APU	O3B-C3B-C2B	-2.73	101.78	111.68
3	B	602	APU	O3B-C3B-C2B	-2.39	103.00	111.68
3	B	602	APU	O3X-PU-O1X	2.12	118.99	110.68

There are no chirality outliers.

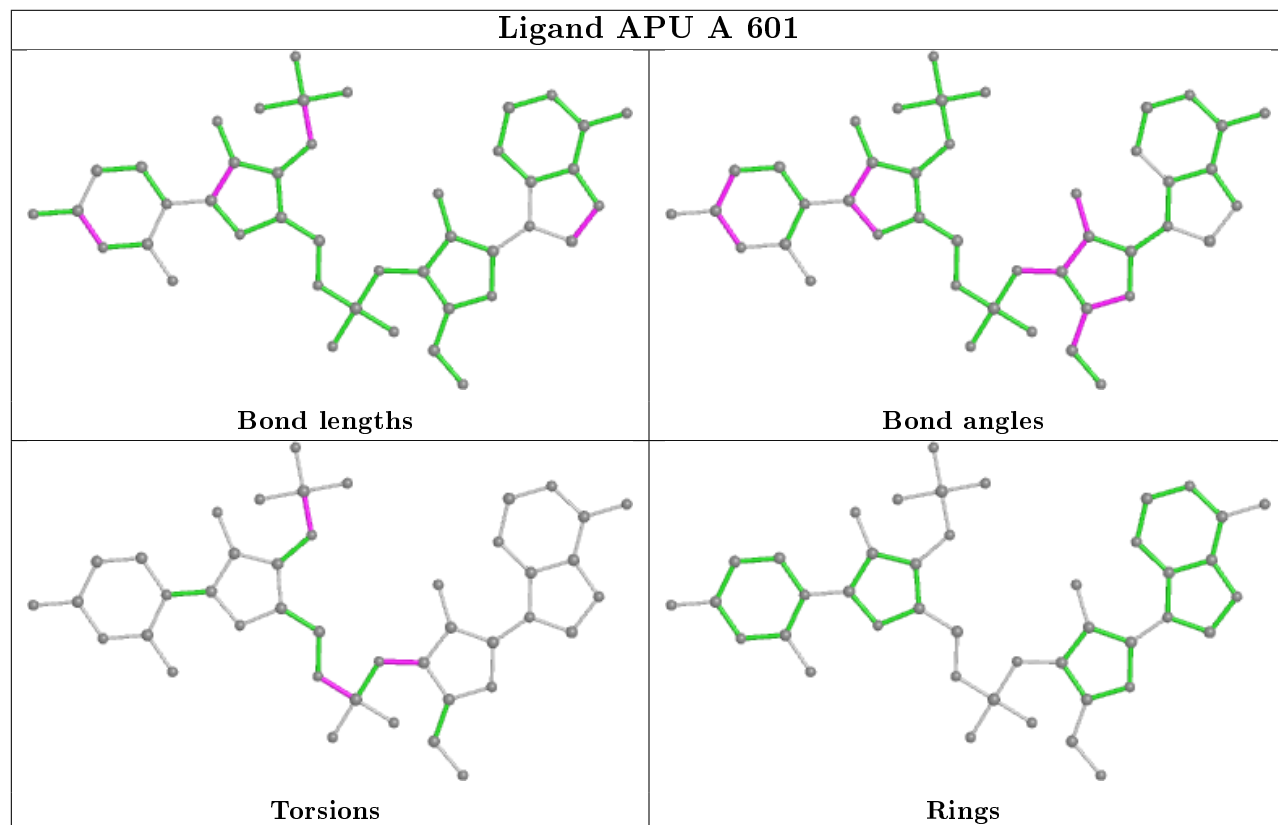
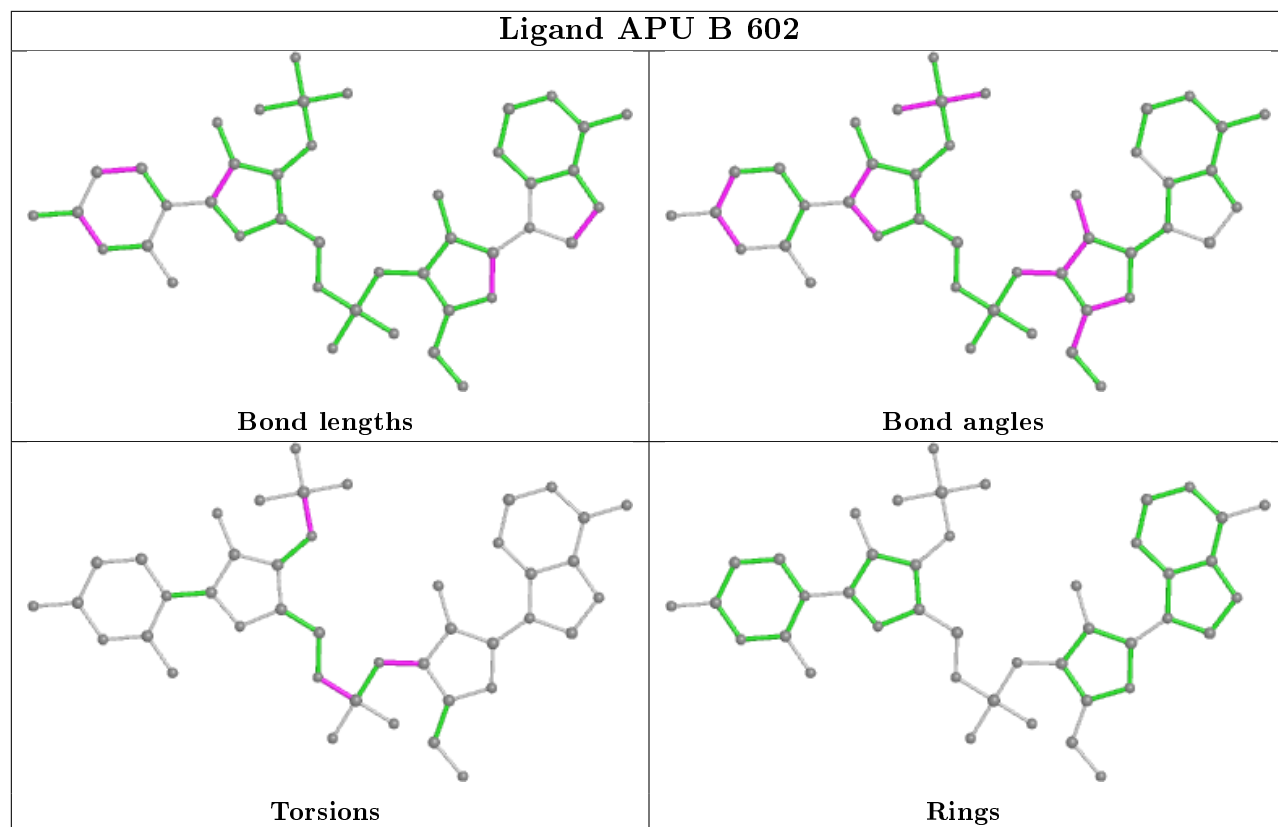
All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	602	APU	C3D-O3D-PU-O1X
3	B	602	APU	C4B-C3B-O3B-PA
3	A	601	APU	C4B-C3B-O3B-PA
3	B	602	APU	C2B-C3B-O3B-PA
3	A	601	APU	C2B-C3B-O3B-PA
3	A	601	APU	C3D-O3D-PU-O1X
3	B	602	APU	C5D-O5D-PA-O3B
3	A	601	APU	C5D-O5D-PA-O3B
3	A	601	APU	C5D-O5D-PA-O1A

There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



## 5.7 Other polymers

There are no such residues in this entry.

## 5.8 Polymer linkage issues

There are no chain breaks in this entry.

## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> 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 <sup>2</sup> )	Q < 0.9
1	A	520/535 (97%)	0.35	40 (7%)	13 16	15, 24, 45, 66	0
1	B	522/535 (97%)	0.18	25 (4%)	30 35	14, 22, 39, 60	1 (0%)
All	All	1042/1070 (97%)	0.27	65 (6%)	20 24	14, 23, 43, 66	1 (0%)

All (65) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	353	ILE	8.1
1	A	350	LEU	6.1
1	A	187[A]	ALA	5.8
1	B	501	SER	5.7
1	A	501	SER	5.2
1	B	353[A]	ILE	5.2
1	B	350	LEU	5.2
1	B	502	ASN	5.0
1	A	348	GLY	4.9
1	B	202[A]	ILE	4.7
1	B	504	ILE	4.7
1	A	201[A]	CYS	4.5
1	B	201[A]	CYS	4.4
1	B	439[A]	GLY	4.4
1	B	348	GLY	4.3
1	A	439[A]	GLY	4.2
1	A	228	ASN	4.0
1	B	499	ILE	3.8
1	A	40	SER	3.8
1	A	496	SER	3.8
1	A	266	VAL	3.7
1	A	186[A]	CYS	3.7
1	B	187[A]	ALA	3.6
1	A	269	THR	3.4

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	B	496	SER	3.4
1	B	497	GLU	3.3
1	A	502	ASN	3.3
1	B	266	VAL	3.2
1	A	235	ASN	3.2
1	A	497	GLU	3.2
1	A	231	SER	3.2
1	B	500	HIS	3.1
1	A	438[A]	PRO	3.0
1	A	170	ARG	3.0
1	B	186[A]	CYS	3.0
1	A	292	GLU	2.8
1	B	494	SER	2.8
1	A	504	ILE	2.8
1	A	531	PHE	2.8
1	A	328	ILE	2.8
1	A	346	LEU	2.6
1	A	533	ILE	2.6
1	B	349	GLU	2.6
1	B	351[A]	VAL	2.6
1	A	203	ASN	2.5
1	A	338	LEU	2.5
1	A	215	THR	2.5
1	B	408[A]	THR	2.5
1	A	410	PHE	2.4
1	A	202[A]	ILE	2.4
1	A	408[A]	THR	2.3
1	A	238	VAL	2.3
1	A	535[A]	SER	2.3
1	A	297	LEU	2.3
1	B	352[A]	ASP	2.3
1	B	533	ILE	2.2
1	A	494	SER	2.2
1	A	236	LYS	2.2
1	A	265	THR	2.2
1	B	520	HIS	2.1
1	A	500	HIS	2.1
1	A	263	LEU	2.1
1	B	110	LEU	2.1
1	B	498	LYS	2.0
1	A	520	HIS	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 ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands ⓘ

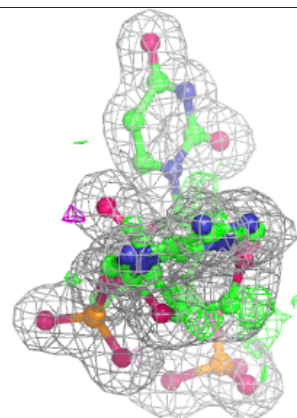
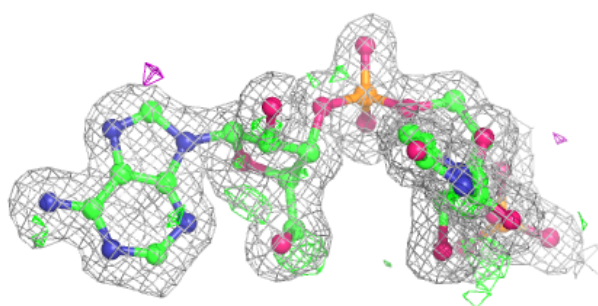
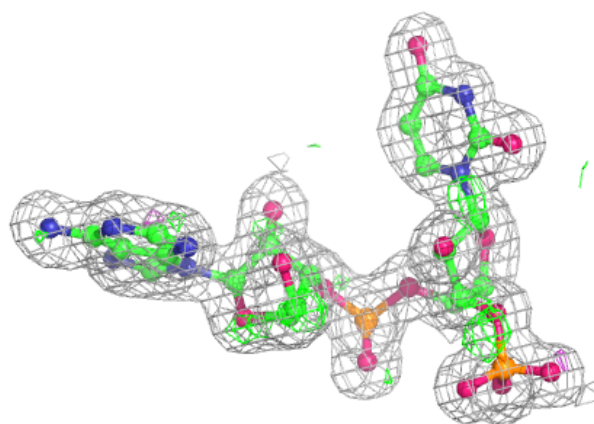
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

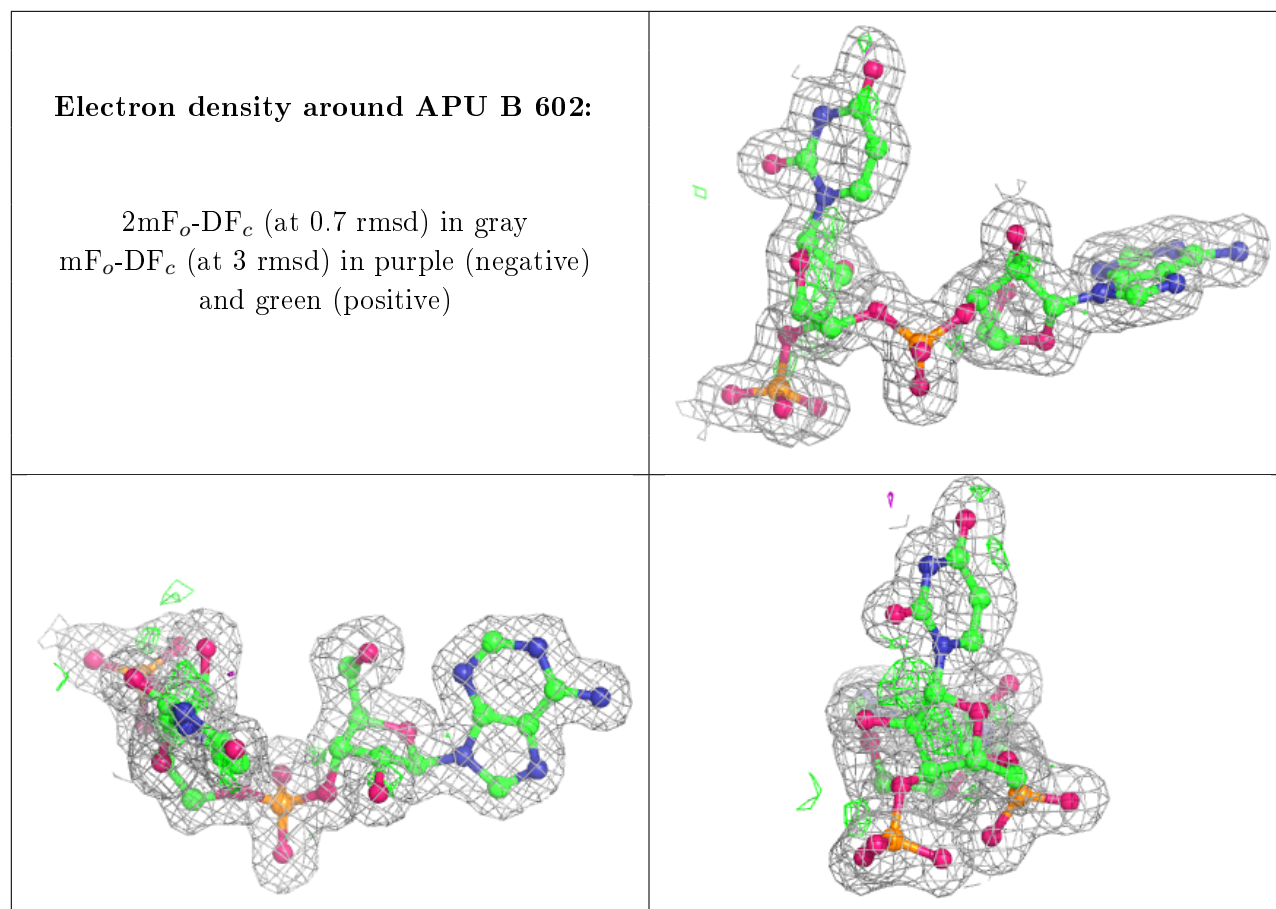
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	CL	A	703	1/1	0.94	0.05	38,38,38,38	0
3	APU	A	601	43/43	0.96	0.07	15,18,21,21	0
2	CL	B	708	1/1	0.96	0.09	34,34,34,34	0
2	CL	A	702	1/1	0.96	0.05	24,24,24,24	0
3	APU	B	602	43/43	0.97	0.07	14,17,21,22	0
2	CL	B	706	1/1	0.97	0.06	28,28,28,28	0
2	CL	A	704	1/1	0.98	0.07	31,31,31,31	0
2	CL	B	709	1/1	0.99	0.06	28,28,28,28	0
2	CL	A	701	1/1	0.99	0.03	20,20,20,20	0
2	CL	B	707	1/1	0.99	0.04	27,27,27,27	0
2	CL	B	705	1/1	1.00	0.04	19,19,19,19	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 APU 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)





## 6.5 Other polymers ⓘ

There are no such residues in this entry.