



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 04:13 PM UTC

PDB ID : 4KH6 / pdb_00004kh6
Title : Toxoplasma gondii NTPDase1 C258S/C268S E493G crystallized with Mg and AMPNP
Authors : Krug, U.; Totzauer, R.; Strater, N.
Deposited on : 2013-04-30
Resolution : 2.40 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

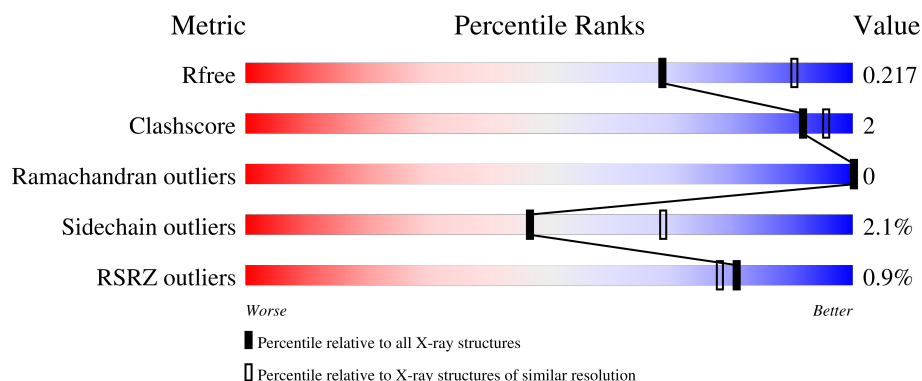
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	611	% 89% 8% .
1	B	611	% 88% 9% .

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9820 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.

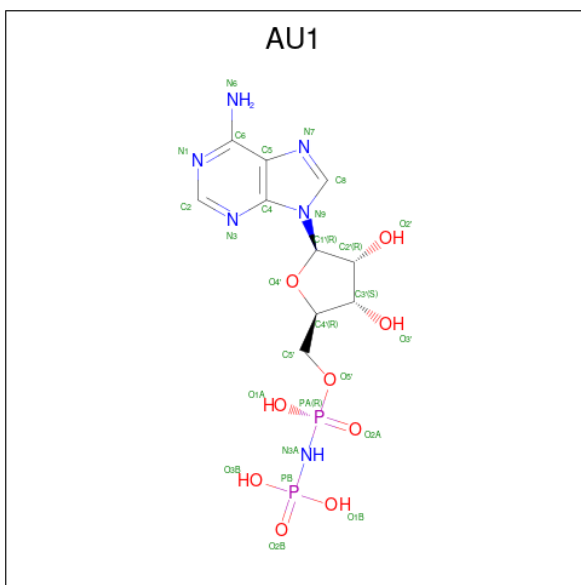
- Molecule 1 is a protein called Nucleoside-triphosphatase 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	594	Total	C	N	O	S	0	0	0
			4632	2910	816	881	25			
1	B	596	Total	C	N	O	S	0	0	0
			4645	2917	818	885	25			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	25	MET	-	initiating methionine	UNP Q27895
A	258	SER	CYS	engineered mutation	UNP Q27895
A	268	SER	CYS	engineered mutation	UNP Q27895
A	493	GLY	GLU	engineered mutation	UNP Q27895
A	629	GLU	-	expression tag	UNP Q27895
A	630	HIS	-	expression tag	UNP Q27895
A	631	HIS	-	expression tag	UNP Q27895
A	632	HIS	-	expression tag	UNP Q27895
A	633	HIS	-	expression tag	UNP Q27895
A	634	HIS	-	expression tag	UNP Q27895
A	635	HIS	-	expression tag	UNP Q27895
B	25	MET	-	initiating methionine	UNP Q27895
B	258	SER	CYS	engineered mutation	UNP Q27895
B	268	SER	CYS	engineered mutation	UNP Q27895
B	493	GLY	GLU	engineered mutation	UNP Q27895
B	629	GLU	-	expression tag	UNP Q27895
B	630	HIS	-	expression tag	UNP Q27895
B	631	HIS	-	expression tag	UNP Q27895
B	632	HIS	-	expression tag	UNP Q27895
B	633	HIS	-	expression tag	UNP Q27895
B	634	HIS	-	expression tag	UNP Q27895
B	635	HIS	-	expression tag	UNP Q27895

- Molecule 2 is 5'-O-[(R)-hydroxy(phosphonoamino)phosphoryl]adenosine (CCD ID: AU1) (formula: C₁₀H₁₆N₆O₉P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 27	C 10	N 6	O 9	P 2	27	0
2	B	1	Total 9	N 1	O 6	P 2		0	0

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

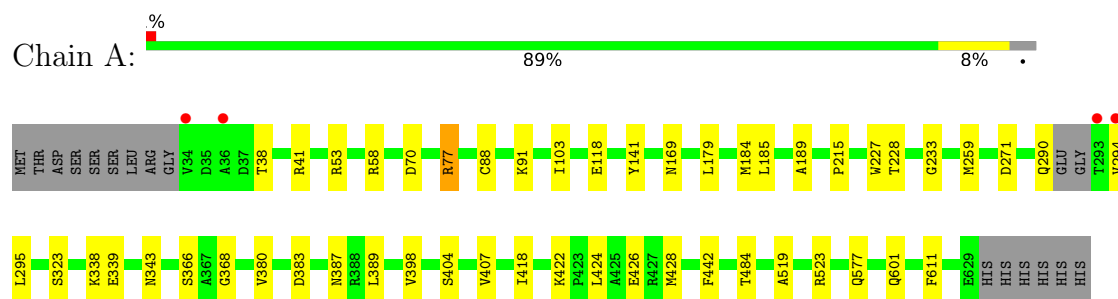
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	297	Total O 298 298	0	1
4	B	207	Total O 207 207	0	0

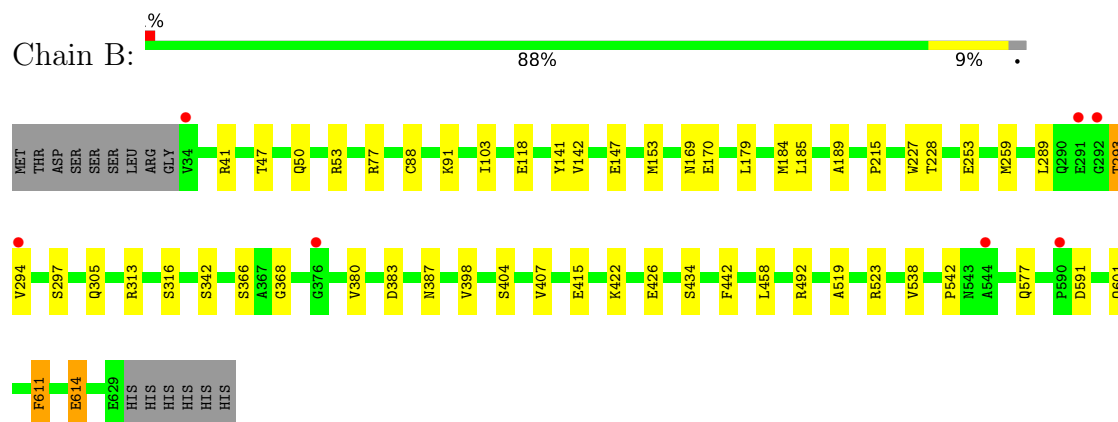
3 Residue-property plots [i](#)

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

• Molecule 1: Nucleoside-triphosphatase 2



• Molecule 1: Nucleoside-triphosphatase 2



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	73.71Å 150.24Å 241.56Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	24.19 – 2.40 24.19 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.7 (24.19-2.40) 99.4 (24.19-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.17 (at 2.39Å)	Xtriage
Refinement program	BUSTER 2.10.0	Depositor
R, R_{free}	0.170 , 0.213 0.175 , 0.217	Depositor DCC
R_{free} test set	1085 reflections (2.05%)	wwPDB-VP
Wilson B-factor (Å ²)	32.9	Xtriage
Anisotropy	0.157	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 56.7	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	0.95	EDS
Total number of atoms	9820	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.98% 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: MG, AU1

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	0.85	0/4723	1.36	17/6394 (0.3%)
1	B	0.84	1/4737 (0.0%)	1.37	27/6414 (0.4%)
All	All	0.84	1/9460 (0.0%)	1.37	44/12808 (0.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	142	VAL	CA-C	5.24	1.57	1.52

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	542	PRO	N-CA-C	7.07	124.07	113.81
1	A	189	ALA	N-CA-C	6.65	120.27	111.75
1	B	189	ALA	N-CA-C	6.58	120.17	111.75
1	A	383	ASP	CA-CB-CG	6.39	119.00	112.60
1	B	611	PHE	CA-C-N	6.29	130.26	120.82
1	B	611	PHE	C-N-CA	6.29	130.26	120.82
1	A	611	PHE	CA-C-N	6.26	130.21	120.82
1	A	611	PHE	C-N-CA	6.26	130.21	120.82
1	B	591	ASP	CA-CB-CG	6.03	118.63	112.60
1	B	179	LEU	N-CA-C	5.84	118.60	111.82
1	B	103	ILE	CB-CA-C	5.79	119.24	111.94
1	A	407	VAL	CA-C-N	5.70	128.23	120.54
1	A	407	VAL	C-N-CA	5.70	128.23	120.54
1	A	141	TYR	CA-C-N	5.58	124.79	120.33
1	A	141	TYR	C-N-CA	5.58	124.79	120.33
1	A	179	LEU	N-CA-C	5.55	118.25	111.82
1	B	41	ARG	CA-C-N	5.52	128.26	120.42
1	B	41	ARG	C-N-CA	5.52	128.26	120.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	407	VAL	CA-C-N	5.43	127.87	120.54
1	B	407	VAL	C-N-CA	5.43	127.87	120.54
1	B	601	GLN	N-CA-C	5.41	116.86	111.07
1	A	601	GLN	N-CA-C	5.40	116.85	111.07
1	A	215	PRO	CA-C-N	5.39	127.81	120.54
1	A	215	PRO	C-N-CA	5.39	127.81	120.54
1	B	141	TYR	CA-C-N	5.33	124.59	120.33
1	B	141	TYR	C-N-CA	5.33	124.59	120.33
1	B	383	ASP	CA-CB-CG	5.29	117.89	112.60
1	A	343	ASN	CA-CB-CG	5.21	117.81	112.60
1	B	170	GLU	N-CA-C	5.18	117.67	111.71
1	B	169	ASN	CA-CB-CG	5.17	117.77	112.60
1	B	53	ARG	CA-C-N	5.15	127.50	120.54
1	B	53	ARG	C-N-CA	5.15	127.50	120.54
1	A	271	ASP	CA-CB-CG	5.14	117.75	112.60
1	A	442	PHE	N-CA-C	5.14	116.88	111.28
1	B	442	PHE	N-CA-C	5.13	116.87	111.28
1	A	169	ASN	CA-CB-CG	5.12	117.72	112.60
1	A	58	ARG	N-CA-C	5.12	117.41	110.35
1	B	297	SER	CA-C-N	5.11	128.07	120.31
1	B	297	SER	C-N-CA	5.11	128.07	120.31
1	B	47	THR	CA-C-N	5.09	127.10	120.28
1	B	47	THR	C-N-CA	5.09	127.10	120.28
1	B	215	PRO	CA-C-N	5.05	127.36	120.54
1	B	215	PRO	C-N-CA	5.05	127.36	120.54
1	B	342	SER	N-CA-C	-5.03	106.57	113.30

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	4632	0	4584	15	0
1	B	4645	0	4594	15	0
2	A	27	0	16	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	9	0	3	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	298	0	0	0	0
4	B	207	0	0	1	0
All	All	9820	0	9197	30	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 2.

All (30) 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:415:GLU:HG3	1:B:458:LEU:HD13	1.52	0.91
1:B:289:LEU:HB3	1:B:293:THR:HG21	1.69	0.74
1:B:253:GLU:HG3	1:B:313:ARG:HG3	1.81	0.61
1:B:611:PHE:HB3	1:B:614:GLU:HG2	1.83	0.61
1:B:611:PHE:HB3	1:B:614:GLU:CG	2.33	0.58
1:B:184:MET:HG2	1:B:227:TRP:HB3	1.88	0.56
1:B:398:VAL:O	1:B:404:SER:HB3	2.05	0.55
1:A:290:GLN:HG3	1:A:484:THR:OG1	2.09	0.53
1:A:398:VAL:O	1:A:404:SER:HB3	2.09	0.53
1:A:366:SER:HB3	1:A:380:VAL:HG11	1.90	0.53
1:A:184:MET:HG2	1:A:227:TRP:HB3	1.90	0.52
1:A:38:THR:HG22	1:A:41:ARG:HH21	1.75	0.52
1:B:366:SER:HB3	1:B:380:VAL:HG11	1.93	0.50
1:A:422:LYS:HB3	1:A:426:GLU:HG3	1.95	0.49
1:A:338:LYS:HB2	1:A:389:LEU:HD11	1.95	0.48
1:A:185:LEU:HB3	1:A:228:THR:HG23	1.95	0.48
1:A:368:GLY:HA2	1:A:387:ASN:HA	1.97	0.47
1:B:88:CYS:HB2	1:B:91:LYS:HB2	1.97	0.47
1:B:368:GLY:HA2	1:B:387:ASN:HA	1.97	0.46
1:A:88:CYS:HB2	1:A:91:LYS:HB2	1.97	0.46
1:A:233:GLY:HA3	1:A:323:SER:HB3	1.97	0.46
1:B:422:LYS:HB3	1:B:426:GLU:HG3	1.97	0.45
1:B:153:MET:HE3	4:B:927:HOH:O	2.16	0.45
1:B:185:LEU:HB3	1:B:228:THR:HG23	1.98	0.43
1:A:424:LEU:HG	1:A:428:MET:HE2	2.01	0.43
1:B:519:ALA:O	1:B:523:ARG:HB2	2.19	0.43
1:A:70:ASP:HB3	1:A:77:ARG:HD2	2.01	0.42
1:A:519:ALA:O	1:A:523:ARG:HB2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:305:GLN:OE1	1:B:316:SER:HA	2.21	0.41
1:A:338:LYS:HE2	1:A:418:ILE:HG13	2.03	0.41

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	590/611 (97%)	580 (98%)	10 (2%)	0	100	100
1	B	594/611 (97%)	586 (99%)	8 (1%)	0	100	100
All	All	1184/1222 (97%)	1166 (98%)	18 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	511/526 (97%)	502 (98%)	9 (2%)	51	73
1	B	512/526 (97%)	500 (98%)	12 (2%)	44	66
All	All	1023/1052 (97%)	1002 (98%)	21 (2%)	47	69

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	53	ARG
1	A	77	ARG
1	A	103	ILE
1	A	118	GLU
1	A	259	MET
1	A	294	VAL
1	A	295	LEU
1	A	339	GLU
1	A	577	GLN
1	B	50	GLN
1	B	77	ARG
1	B	118	GLU
1	B	147	GLU
1	B	259	MET
1	B	293	THR
1	B	294	VAL
1	B	434	SER
1	B	492	ARG
1	B	538	VAL
1	B	577	GLN
1	B	614	GLU

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	270	ASN
1	A	284	GLN
1	A	290	GLN
1	A	363	GLN
1	A	412	GLN
1	A	448	GLN
1	B	448	GLN
1	B	577	GLN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 2 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
2	AU1	B	700	3	8,8,29	3.36	5 (62%)	8,13,45	3.35	2 (25%)
2	AU1	A	700	3	29,29,29	2.19	10 (34%)	40,45,45	2.18	14 (35%)

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	AU1	B	700	3	-	2/2/6/32	-
2	AU1	A	700	3	-	5/13/32/32	0/3/3/3

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	700	AU1	C5-C4	5.11	1.48	1.39
2	B	700	AU1	PB-N3A	4.82	1.76	1.63
2	A	700	AU1	PA-N3A	4.72	1.75	1.63
2	A	700	AU1	PB-N3A	4.68	1.75	1.63
2	B	700	AU1	PA-N3A	4.24	1.74	1.63
2	B	700	AU1	PB-O2B	4.20	1.52	1.46
2	B	700	AU1	PA-O2A	4.10	1.52	1.46
2	A	700	AU1	PA-O2A	3.59	1.51	1.46
2	A	700	AU1	PB-O2B	3.42	1.51	1.46
2	A	700	AU1	C5-C6	2.95	1.49	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	700	AU1	C8-N7	2.33	1.36	1.31
2	A	700	AU1	C5-N7	-2.29	1.34	1.39
2	B	700	AU1	PA-O5'	-2.10	1.51	1.56
2	A	700	AU1	PB-O1B	-2.08	1.51	1.56
2	A	700	AU1	PB-O3B	-2.05	1.51	1.56

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	700	AU1	O2B-PB-N3A	-7.12	101.28	111.77
2	A	700	AU1	C5-C4-N3	-5.92	118.56	126.72
2	B	700	AU1	O2A-PA-N3A	-5.69	103.39	111.77
2	A	700	AU1	N3-C4-N9	4.79	135.32	127.17
2	A	700	AU1	O2B-PB-N3A	-4.05	105.81	111.77
2	A	700	AU1	C2-N3-C4	3.87	121.28	111.83
2	A	700	AU1	O1A-PA-O2A	3.68	117.77	109.87
2	A	700	AU1	N1-C2-N3	-3.58	123.17	128.58
2	A	700	AU1	C4-C5-N7	-3.36	106.75	110.58
2	A	700	AU1	O2A-PA-N3A	-3.16	107.11	111.77
2	A	700	AU1	C5-N7-C8	2.47	107.34	103.45
2	A	700	AU1	C4-N9-C8	2.39	108.25	105.74
2	A	700	AU1	C3'-C2'-C1'	2.29	105.79	101.46
2	A	700	AU1	O1B-PB-O3B	2.20	113.51	107.59
2	A	700	AU1	C6-C5-N7	2.11	136.15	132.09
2	A	700	AU1	O5'-PA-O2A	-2.05	108.18	114.27

There are no chirality outliers.

All (7) torsion outliers are listed below:

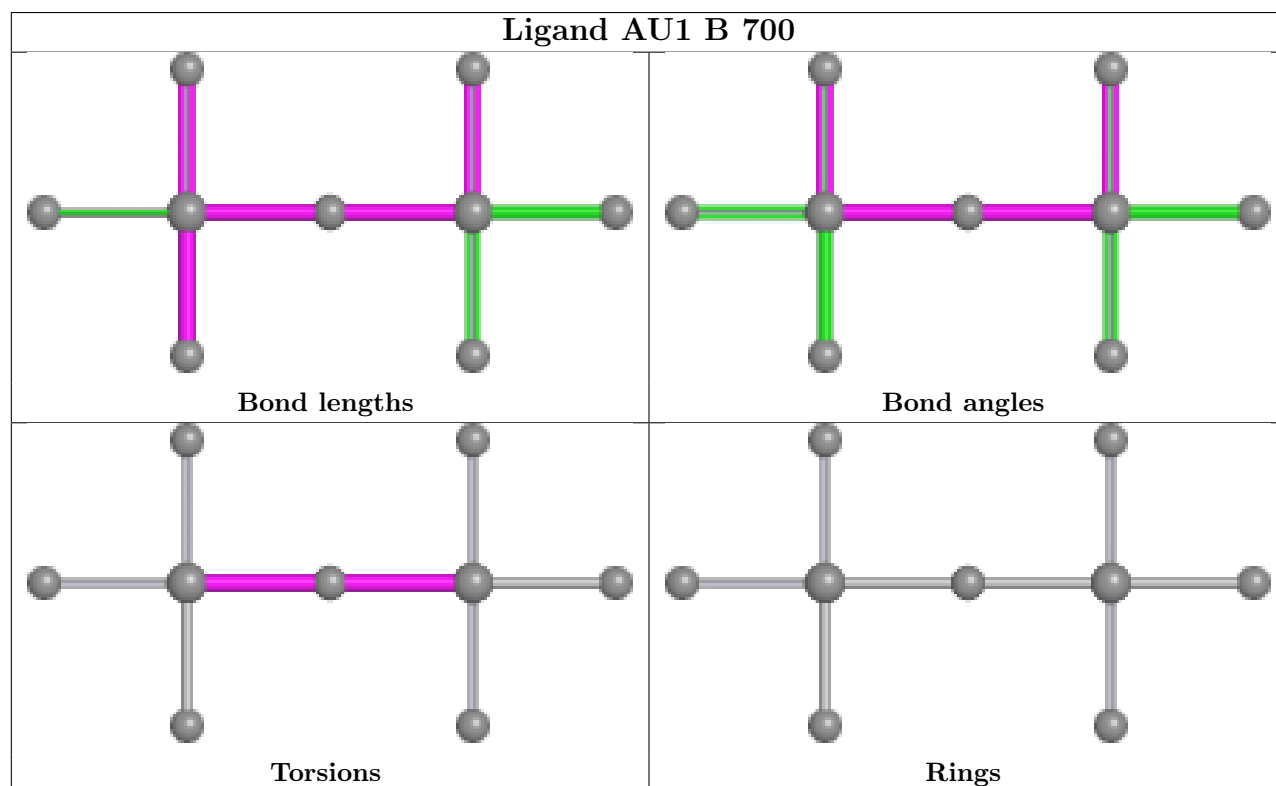
Mol	Chain	Res	Type	Atoms
2	A	700	AU1	PB-N3A-PA-O2A
2	A	700	AU1	C5'-O5'-PA-O1A
2	B	700	AU1	PB-N3A-PA-O2A
2	B	700	AU1	PA-N3A-PB-O2B
2	A	700	AU1	O4'-C4'-C5'-O5'
2	A	700	AU1	C3'-C4'-C5'-O5'
2	A	700	AU1	C5'-O5'-PA-N3A

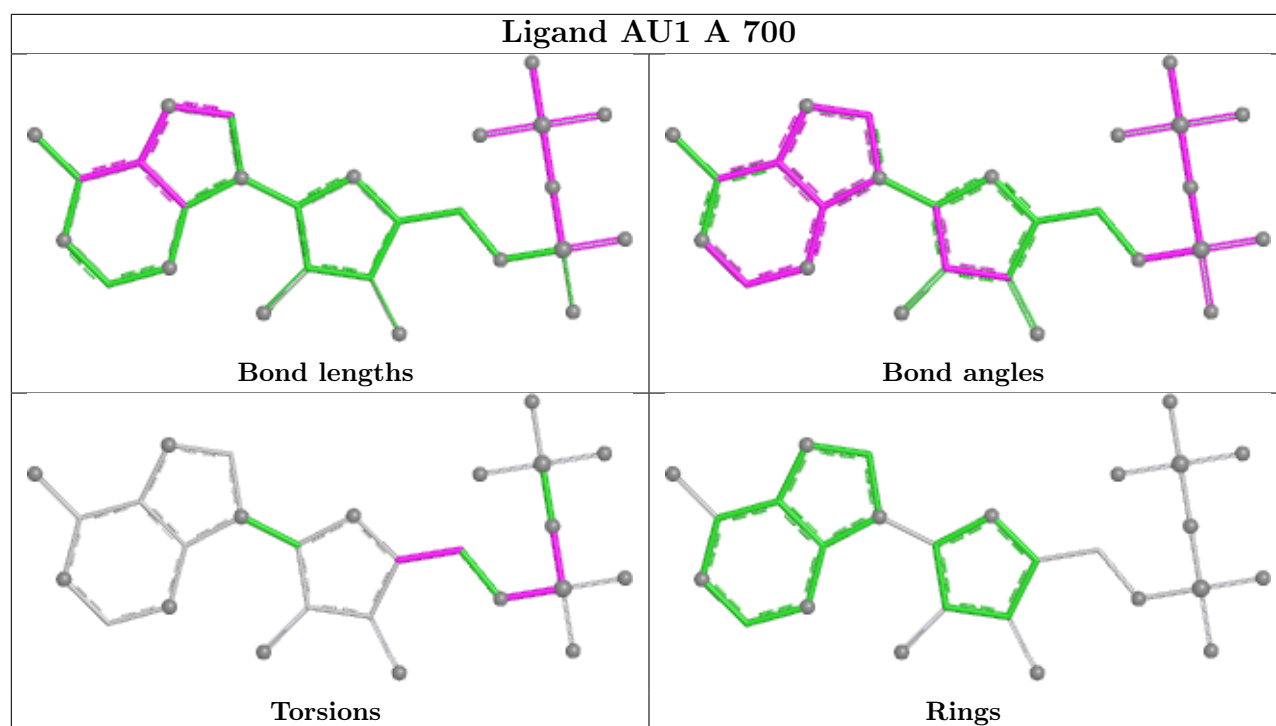
There are no ring outliers.

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 [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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	594/611 (97%)	-0.59	4 (0%) 84 81	20, 33, 60, 102	0
1	B	596/611 (97%)	-0.43	7 (1%) 76 73	21, 39, 80, 103	0
All	All	1190/1222 (97%)	-0.51	11 (0%) 81 78	20, 35, 71, 103	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	34	VAL	3.3
1	A	294	VAL	3.2
1	A	34	VAL	3.0
1	B	544	ALA	2.6
1	B	292	GLY	2.5
1	B	291	GLU	2.4
1	A	36	ALA	2.2
1	B	376	GLY	2.2
1	B	294	VAL	2.1
1	B	590	PRO	2.1
1	A	293	THR	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 ⓘ

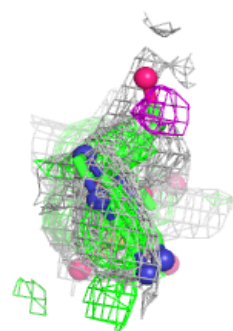
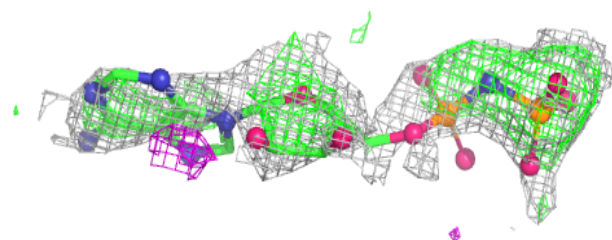
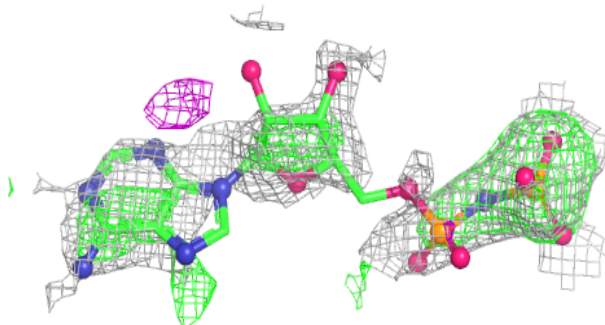
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	AU1	A	700	27/27	-	-	132,132,132,132	27
2	AU1	B	700	9/27	0.89	0.14	42,44,51,54	9
3	MG	A	701	1/1	0.92	0.05	42,42,42,42	0
3	MG	B	701	1/1	0.99	0.03	37,37,37,37	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.

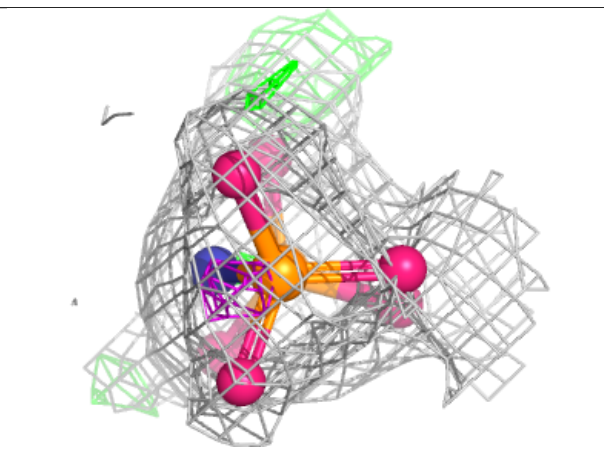
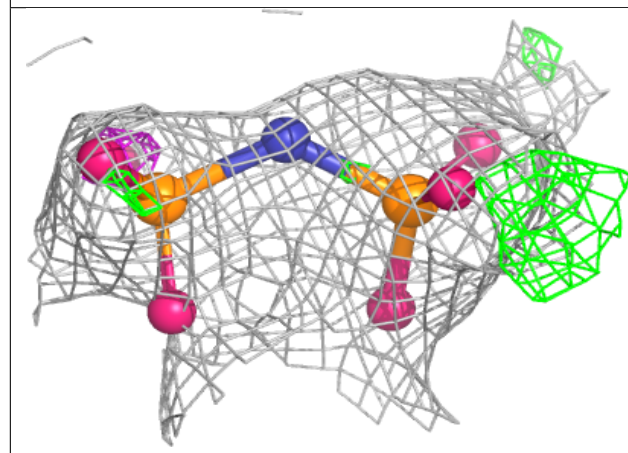
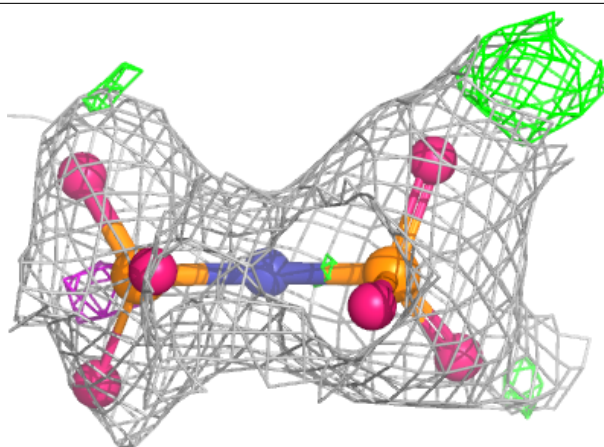
Electron density around AU1 A 700:

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 AU1 B 700:

$2mF_o - DF_c$ (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.