



# Full wwPDB X-ray Structure Validation Report ⓘ

Jun 15, 2020 – 11:00 pm BST

PDB ID : 2ZDH  
Title : Crystal structure of D-Alanine:D-Alanine Ligase with ADP and D-Alanine from *Thermus thermophilus* HB8  
Authors : Kitamuta, Y.; Yokoyama, S.; Kuramitsu, S.; RIKEN Structural Genomics/Proteomics Initiative (RSGI)  
Deposited on : 2007-11-22  
Resolution : 1.90 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.

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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

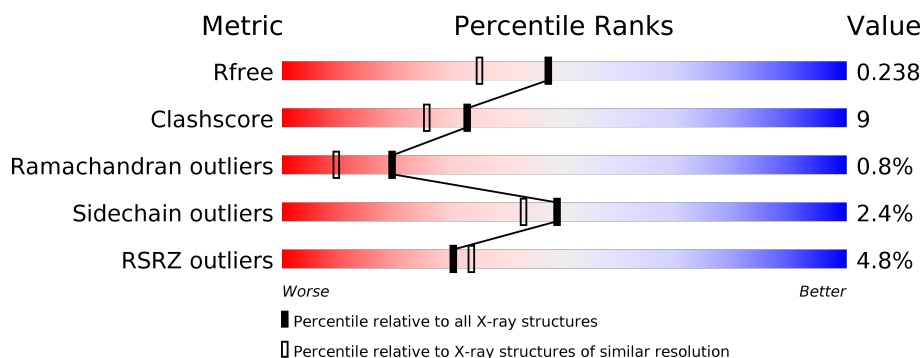
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.90 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	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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	319	<div> <div>83%</div> <div>16%</div> <div>•</div> </div>
1	B	319	<div> <div>4%</div> <div>78%</div> <div>18%</div> <div>••</div> </div>
1	C	319	<div> <div>6%</div> <div>78%</div> <div>19%</div> <div>••</div> </div>
1	D	319	<div> <div>8%</div> <div>80%</div> <div>19%</div> <div>•</div> </div>

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 10481 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 D-alanine–D-alanine ligase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	319	Total	C	N	O	S	0	0	0
			2453	1593	406	449	5			
1	B	312	Total	C	N	O	S	0	0	0
			2389	1552	398	434	5			
1	C	316	Total	C	N	O	S	0	0	0
			2416	1570	402	439	5			
1	D	319	Total	C	N	O	S	0	0	0
			2453	1593	406	449	5			

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

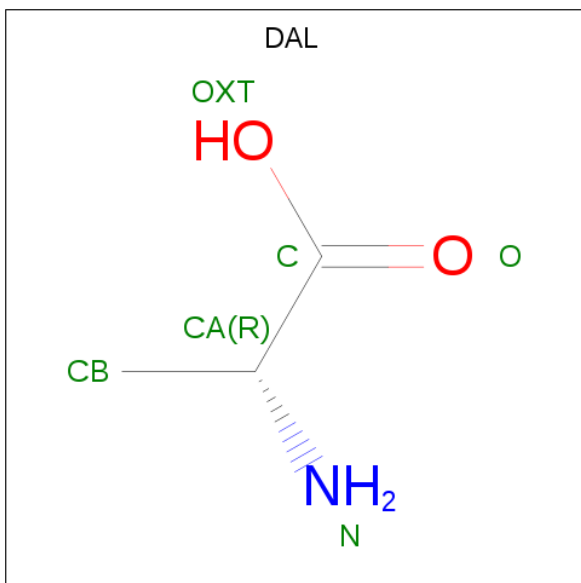
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	2	Total	Mg	0	0
			2	2		
2	A	2	Total	Mg	0	0
			2	2		
2	D	1	Total	Mg	0	0
			1	1		
2	C	2	Total	Mg	0	0
			2	2		

- Molecule 3 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: C<sub>10</sub>H<sub>15</sub>N<sub>5</sub>O<sub>10</sub>P<sub>2</sub>).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 27	C 10	N 5	O 10	P 2	0	0
3	B	1	Total 27	C 10	N 5	O 10	P 2	0	0
3	C	1	Total 27	C 10	N 5	O 10	P 2	0	0
3	D	1	Total 27	C 10	N 5	O 10	P 2	0	0

- Molecule 4 is D-ALANINE (three-letter code: DAL) (formula:  $\text{C}_3\text{H}_7\text{NO}_2$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			6	3	1	2		
4	D	1	Total	C	N	O	0	0
			6	3	1	2		

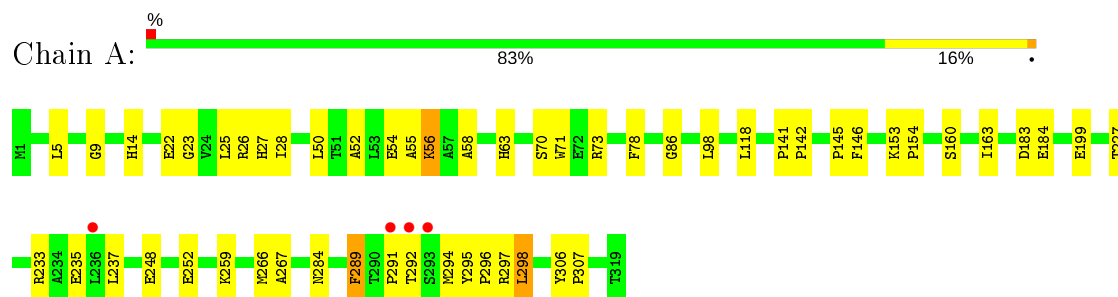
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	188	Total	O	0	0
			188	188		
5	B	174	Total	O	0	0
			174	174		
5	C	145	Total	O	0	0
			145	145		
5	D	136	Total	O	0	0
			136	136		

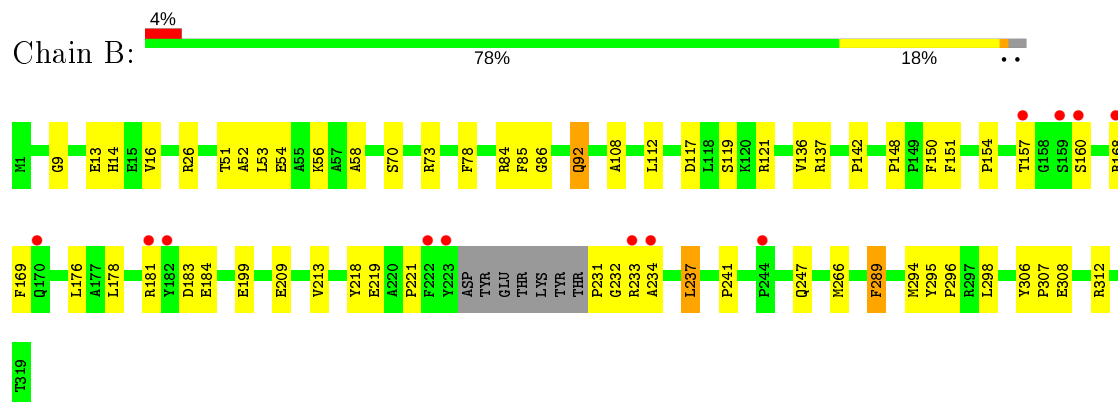
### 3 Residue-property plots [i](#)

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.

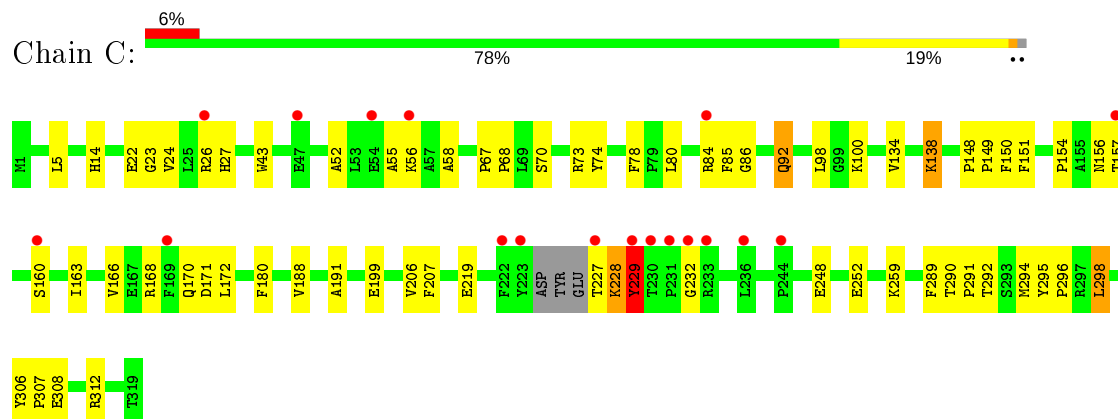
#### • Molecule 1: D-alanine–D-alanine ligase



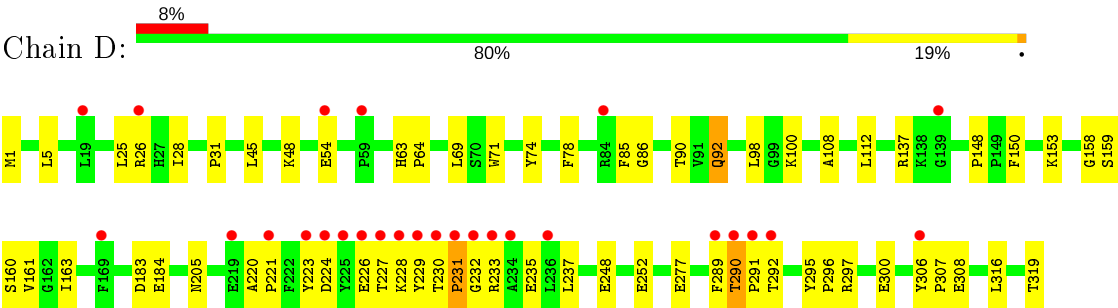
#### • Molecule 1: D-alanine–D-alanine ligase



#### • Molecule 1: D-alanine–D-alanine ligase



● Molecule 1: D-alanine–D-alanine ligase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	68.95Å 105.39Å 199.75Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.90 – 1.90 49.96 – 1.90	Depositor EDS
% Data completeness (in resolution range)	96.9 (47.90-1.90) 97.0 (49.96-1.90)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	0.08	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.79 (at 1.90Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.207 , 0.239 0.207 , 0.238	Depositor DCC
$R_{free}$ test set	11505 reflections (10.01%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	20.2	Xtriage
Anisotropy	0.380	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 47.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	10481	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.33% of the height of the origin peak. No significant pseudotranslation is detected.*

<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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: DAL, MG, ADP

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	0.33	0/2518	0.61	1/3434 (0.0%)
1	B	0.32	0/2451	0.61	1/3340 (0.0%)
1	C	0.32	0/2479	0.59	1/3380 (0.0%)
1	D	0.31	0/2518	0.58	0/3434
All	All	0.32	0/9966	0.60	3/13588 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	154	PRO	N-CA-C	-5.20	98.57	112.10
1	C	154	PRO	N-CA-C	-5.04	99.00	112.10
1	A	154	PRO	N-CA-C	-5.02	99.05	112.10

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2453	0	2470	40	0
1	B	2389	0	2415	38	0
1	C	2416	0	2429	52	0
1	D	2453	0	2470	47	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	1	0	0	0	0
3	A	27	0	12	0	0
3	B	27	0	12	0	0
3	C	27	0	12	0	0
3	D	27	0	12	0	0
4	A	6	0	6	0	0
4	D	6	0	6	0	0
5	A	188	0	0	1	0
5	B	174	0	0	2	0
5	C	145	0	0	1	0
5	D	136	0	0	1	0
All	All	10481	0	9844	172	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 (172) 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:292:THR:HB	1:A:297:ARG:HH11	1.37	0.88
1:C:5:LEU:HD11	1:C:80:LEU:HD12	1.62	0.81
1:D:108:ALA:O	1:D:112:LEU:HD23	1.83	0.79
1:C:14:HIS:HE1	1:C:58:ALA:H	1.34	0.76
1:D:290:THR:HG23	1:D:292:THR:H	1.51	0.74
1:D:295:TYR:HB3	1:D:296:PRO:HD3	1.69	0.74
1:C:229:TYR:O	1:C:229:TYR:HD2	1.70	0.73
1:A:14:HIS:HE1	1:A:58:ALA:H	1.33	0.73
1:D:230:THR:OG1	1:D:231:PRO:HD2	1.90	0.71
1:C:168:ARG:HG3	1:C:170:GLN:HG2	1.74	0.70
1:A:22:GLU:HG2	1:A:26:ARG:HH22	1.56	0.70
1:B:295:TYR:HB3	1:B:296:PRO:HD3	1.73	0.70
1:A:22:GLU:HG2	1:A:26:ARG:NH2	2.07	0.69
1:C:138:LYS:NZ	1:C:180:PHE:HB3	2.08	0.69
1:C:206:VAL:HG11	1:D:137:ARG:CD	2.22	0.69
1:D:220:ALA:HB1	1:D:221:PRO:HD2	1.75	0.68
1:D:289:PHE:O	1:D:290:THR:HB	1.93	0.68
1:D:290:THR:OG1	1:D:291:PRO:HD2	1.95	0.67
1:A:14:HIS:CE1	1:A:58:ALA:H	2.12	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:294:MET:HG3	1:A:298:LEU:HD22	1.75	0.66
1:C:248:GLU:O	1:C:252:GLU:HG3	1.95	0.66
1:D:160:SER:HA	1:D:163:ILE:HD12	1.78	0.65
1:B:14:HIS:HE1	1:B:58:ALA:H	1.45	0.64
1:C:295:TYR:HB3	1:C:296:PRO:HD3	1.77	0.64
1:D:316:LEU:HA	1:D:319:THR:HG22	1.78	0.64
1:D:45:LEU:HD22	1:D:69:LEU:HD13	1.78	0.63
1:B:14:HIS:CE1	1:B:58:ALA:H	2.16	0.63
1:C:14:HIS:CE1	1:C:58:ALA:H	2.14	0.63
1:D:227:THR:HA	1:D:233:ARG:HD3	1.79	0.63
1:A:295:TYR:HB3	1:A:296:PRO:HD3	1.81	0.62
1:C:229:TYR:O	1:C:229:TYR:CD2	2.51	0.62
1:C:160:SER:HA	1:C:163:ILE:HD12	1.79	0.62
1:A:248:GLU:O	1:A:252:GLU:HG3	2.00	0.62
1:B:108:ALA:O	1:B:112:LEU:HD13	2.01	0.61
1:C:290:THR:HG22	1:C:292:THR:H	1.65	0.61
1:D:48:LYS:HD3	1:D:48:LYS:H	1.65	0.61
1:D:45:LEU:O	1:D:48:LYS:HE2	2.01	0.60
1:C:199:GLU:OE1	1:C:294:MET:HG2	2.02	0.60
1:C:24:VAL:HG21	1:C:80:LEU:HD11	1.82	0.60
1:C:92:GLN:H	1:C:92:GLN:NE2	1.97	0.60
1:B:178:LEU:O	1:B:181:ARG:HG2	2.01	0.60
1:B:306:TYR:HB3	1:B:307:PRO:HD3	1.84	0.60
1:D:92:GLN:NE2	1:D:92:GLN:H	1.99	0.60
1:B:142:PRO:HG3	1:B:176:LEU:HD11	1.85	0.59
1:D:153:LYS:HB3	1:D:163:ILE:HG12	1.85	0.59
1:C:306:TYR:HB3	1:C:307:PRO:HD3	1.85	0.59
1:B:51:THR:O	1:B:54:GLU:HG2	2.04	0.58
1:C:148:PRO:HG3	1:C:150:PHE:CE2	2.37	0.58
1:A:23:GLY:HA3	1:A:291:PRO:HD3	1.85	0.58
1:B:26:ARG:NE	1:B:26:ARG:HA	2.18	0.58
1:B:148:PRO:HG3	1:B:150:PHE:CE2	2.38	0.57
1:A:306:TYR:HB3	1:A:307:PRO:HD3	1.86	0.57
1:D:248:GLU:O	1:D:252:GLU:HG3	2.05	0.57
1:D:289:PHE:O	1:D:290:THR:CB	2.52	0.57
1:A:153:LYS:HB3	1:A:163:ILE:HG12	1.87	0.57
1:D:48:LYS:N	1:D:48:LYS:HD3	2.20	0.57
1:C:156:ASN:O	1:C:157:THR:HG23	2.06	0.56
1:C:207:PHE:CD1	1:C:259:LYS:HE2	2.41	0.56
1:D:292:THR:HA	1:D:297:ARG:HD2	1.88	0.56
1:B:209:GLU:OE2	1:B:312:ARG:NH2	2.37	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:53:LEU:O	1:B:56:LYS:HD2	2.05	0.56
1:C:74:TYR:O	1:C:100:LYS:HE3	2.06	0.55
1:B:92:GLN:NE2	1:B:92:GLN:H	2.05	0.55
1:A:56:LYS:HE3	1:A:56:LYS:HA	1.88	0.55
1:B:52:ALA:HB1	1:B:58:ALA:HB2	1.88	0.55
1:B:237:LEU:HD13	1:B:241:PRO:HD3	1.88	0.54
1:C:168:ARG:HG3	1:C:170:GLN:CG	2.38	0.54
1:A:27:HIS:ND1	1:A:306:TYR:HD2	2.05	0.54
1:A:71:TRP:CD2	1:A:98:LEU:HD21	2.43	0.54
1:D:316:LEU:HA	1:D:319:THR:CG2	2.38	0.54
1:D:26:ARG:HD3	1:D:26:ARG:O	2.08	0.54
1:C:22:GLU:O	1:C:26:ARG:HG3	2.07	0.53
1:D:224:ASP:O	1:D:228:LYS:HB2	2.09	0.53
1:C:43:TRP:CE3	1:C:67:PRO:HG3	2.44	0.53
1:C:70:SER:HB2	1:C:73:ARG:HD3	1.90	0.53
1:D:306:TYR:HB3	1:D:307:PRO:HD3	1.92	0.52
1:A:259:LYS:HG2	1:B:137:ARG:NE	2.25	0.52
1:C:219:GLU:O	1:C:232:GLY:HA3	2.10	0.52
1:B:233:ARG:HG3	1:B:233:ARG:HH11	1.74	0.52
1:C:138:LYS:HB2	1:C:138:LYS:NZ	2.25	0.52
1:D:74:TYR:O	1:D:100:LYS:HE3	2.10	0.52
1:C:294:MET:HG3	1:C:298:LEU:HD22	1.93	0.51
1:A:27:HIS:HB2	1:A:306:TYR:CD2	2.45	0.51
1:C:149:PRO:HG2	1:C:191:ALA:HB3	1.93	0.51
1:A:50:LEU:O	1:A:54:GLU:HG2	2.10	0.51
1:B:168:ARG:HG2	1:B:168:ARG:HH11	1.76	0.50
1:D:159:SER:HB2	1:D:229:TYR:HE2	1.76	0.50
1:A:118:LEU:HD13	1:B:112:LEU:CD2	2.42	0.50
1:D:183:ASP:OD1	1:D:184:GLU:N	2.45	0.50
1:A:52:ALA:HB1	1:A:58:ALA:HB2	1.93	0.50
1:B:199:GLU:OE1	1:B:294:MET:HG2	2.11	0.49
1:D:71:TRP:CG	1:D:98:LEU:HD21	2.46	0.49
1:C:26:ARG:NH1	1:C:26:ARG:HB3	2.27	0.49
1:A:292:THR:O	1:A:297:ARG:NE	2.46	0.49
1:C:26:ARG:HH11	1:C:26:ARG:HB3	1.77	0.49
1:A:235:GLU:OE1	1:A:237:LEU:HD21	2.12	0.49
1:A:55:ALA:O	1:A:56:LYS:HB2	2.13	0.48
1:D:1:MET:O	1:D:31:PRO:HD2	2.13	0.48
1:D:48:LYS:CD	1:D:48:LYS:H	2.26	0.48
1:D:235:GLU:OE2	1:D:237:LEU:HD21	2.13	0.48
1:B:218:TYR:CE1	1:B:234:ALA:HB2	2.48	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:70:SER:HB2	1:A:73:ARG:HD3	1.94	0.48
1:B:266:MET:SD	1:B:289:PHE:HE1	2.37	0.48
1:C:134:VAL:HG22	1:C:188:VAL:HB	1.96	0.48
1:C:55:ALA:O	1:C:56:LYS:HG2	2.12	0.48
1:B:117:ASP:O	1:B:121:ARG:HG3	2.15	0.47
1:C:23:GLY:HA3	1:C:291:PRO:HD3	1.97	0.47
1:D:205:ASN:ND2	1:D:319:THR:HG23	2.30	0.47
1:C:26:ARG:HH11	1:C:26:ARG:CB	2.27	0.47
1:C:138:LYS:HB2	1:C:138:LYS:HZ2	1.80	0.46
1:A:227:THR:HA	1:A:233:ARG:HD3	1.97	0.46
1:C:100:LYS:HE2	5:C:950:HOH:O	2.15	0.46
1:C:55:ALA:C	1:C:56:LYS:HG2	2.36	0.46
1:C:67:PRO:HA	1:C:68:PRO:HD3	1.76	0.46
1:B:219:GLU:O	1:B:232:GLY:HA3	2.16	0.46
1:C:166:VAL:HG13	1:C:171:ASP:HB2	1.98	0.46
1:C:172:LEU:C	1:C:172:LEU:HD23	2.37	0.45
1:D:277:GLU:HG3	5:D:833:HOH:O	2.17	0.45
1:D:71:TRP:O	1:D:100:LYS:HE2	2.17	0.45
1:A:28:ILE:HG23	1:A:28:ILE:O	2.17	0.45
1:D:159:SER:HG	1:D:161:VAL:HG23	1.82	0.45
1:D:63:HIS:HA	1:D:64:PRO:HD3	1.89	0.45
1:D:5:LEU:HD23	1:D:25:LEU:HG	1.99	0.45
1:C:5:LEU:HD11	1:C:80:LEU:CD1	2.41	0.44
1:C:26:ARG:NH1	1:C:27:HIS:CE1	2.85	0.44
1:A:160:SER:HA	1:A:163:ILE:HD12	1.99	0.44
1:A:27:HIS:HD1	1:A:306:TYR:HD2	1.65	0.44
1:A:63:HIS:HD2	5:A:910:HOH:O	2.01	0.44
1:A:259:LYS:HG2	1:B:137:ARG:HE	1.82	0.44
1:A:27:HIS:ND1	1:A:306:TYR:CD2	2.85	0.44
1:D:148:PRO:HG3	1:D:150:PHE:CZ	2.52	0.44
1:C:22:GLU:HG2	1:C:26:ARG:HD2	2.00	0.44
1:A:183:ASP:OD2	1:A:184:GLU:N	2.50	0.43
1:A:26:ARG:C	1:A:27:HIS:HD2	2.22	0.43
1:B:183:ASP:OD2	1:B:184:GLU:N	2.51	0.43
1:C:52:ALA:HB1	1:C:58:ALA:HB2	2.01	0.43
1:D:226:GLU:O	1:D:230:THR:HB	2.18	0.43
1:B:136:VAL:HG22	1:B:142:PRO:HB3	2.00	0.43
1:A:145:PRO:HG2	1:A:146:PHE:CD1	2.54	0.43
1:D:296:PRO:O	1:D:300:GLU:HG3	2.18	0.43
1:C:206:VAL:HG11	1:D:137:ARG:HD2	1.97	0.43
1:D:85:PHE:HA	1:D:90:THR:OG1	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:213:VAL:H	1:B:247:GLN:HE21	1.65	0.43
1:B:308:GLU:O	1:B:312:ARG:HG3	2.18	0.42
1:A:267:ALA:HA	1:A:284:ASN:O	2.19	0.42
1:D:158:GLY:O	1:D:161:VAL:HG23	2.19	0.42
1:D:290:THR:HG23	1:D:292:THR:N	2.24	0.42
1:D:28:ILE:HG23	1:D:28:ILE:O	2.20	0.42
1:C:56:LYS:HE2	1:C:56:LYS:HB3	1.89	0.42
1:C:84:ARG:O	1:C:85:PHE:HB2	2.20	0.42
1:A:5:LEU:HD23	1:A:25:LEU:HG	2.02	0.42
1:A:9:GLY:O	1:A:14:HIS:HD2	2.03	0.42
1:B:112:LEU:O	1:B:119:SER:HB3	2.20	0.42
1:A:266:MET:SD	1:A:289:PHE:HE1	2.43	0.41
1:C:290:THR:HG23	1:C:291:PRO:HD2	2.03	0.41
1:B:84:ARG:NH2	1:B:85:PHE:CE2	2.88	0.41
1:B:157:THR:HB	5:B:896:HOH:O	2.20	0.41
1:C:27:HIS:HB2	1:C:306:TYR:CD2	2.55	0.41
1:B:221:PRO:HB3	1:B:231:PRO:HA	2.03	0.41
1:B:56:LYS:HG3	5:B:858:HOH:O	2.21	0.41
1:C:227:THR:O	1:C:228:LYS:CB	2.69	0.41
1:A:27:HIS:CG	1:A:306:TYR:HD2	2.39	0.41
1:C:308:GLU:O	1:C:312:ARG:HG3	2.21	0.41
1:A:199:GLU:OE1	1:A:294:MET:HG2	2.21	0.40
1:A:141:PRO:HA	1:A:142:PRO:HD3	1.84	0.40
1:B:181:ARG:HG3	1:B:181:ARG:HH21	1.86	0.40
1:D:71:TRP:CD2	1:D:98:LEU:HD21	2.55	0.40
1:B:13:GLU:O	1:B:16:VAL:HG22	2.22	0.40
1:D:296:PRO:HB3	1:D:306:TYR:CE1	2.56	0.40
1:B:9:GLY:O	1:B:14:HIS:HD2	2.03	0.40
1:B:70:SER:HB2	1:B:73:ARG:HD3	2.04	0.40
1:C:298:LEU:HD12	1:C:298:LEU:HA	1.83	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	317/319 (99%)	308 (97%)	8 (2%)	1 (0%)	41	31
1	B	308/319 (97%)	297 (96%)	9 (3%)	2 (1%)	25	15
1	C	312/319 (98%)	298 (96%)	11 (4%)	3 (1%)	15	6
1	D	317/319 (99%)	301 (95%)	12 (4%)	4 (1%)	12	4
All	All	1254/1276 (98%)	1204 (96%)	40 (3%)	10 (1%)	19	9

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	290	THR
1	B	86	GLY
1	C	229	TYR
1	D	86	GLY
1	D	232	GLY
1	B	160	SER
1	C	86	GLY
1	C	228	LYS
1	D	231	PRO
1	A	86	GLY

### 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	256/256 (100%)	252 (98%)	4 (2%)	62	60
1	B	249/256 (97%)	242 (97%)	7 (3%)	43	36
1	C	250/256 (98%)	242 (97%)	8 (3%)	39	30
1	D	256/256 (100%)	251 (98%)	5 (2%)	55	51
All	All	1011/1024 (99%)	987 (98%)	24 (2%)	49	43

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	56	LYS
1	A	78	PHE
1	A	289	PHE
1	A	298	LEU
1	B	78	PHE
1	B	92	GLN
1	B	151	PHE
1	B	169	PHE
1	B	237	LEU
1	B	289	PHE
1	B	298	LEU
1	C	78	PHE
1	C	92	GLN
1	C	98	LEU
1	C	138	LYS
1	C	151	PHE
1	C	229	TYR
1	C	289	PHE
1	C	298	LEU
1	D	54	GLU
1	D	78	PHE
1	D	92	GLN
1	D	223	TYR
1	D	308	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	14	HIS
1	A	27	HIS
1	A	63	HIS
1	A	156	ASN
1	B	14	HIS
1	B	92	GLN
1	B	247	GLN
1	C	14	HIS
1	C	27	HIS
1	C	63	HIS
1	C	92	GLN
1	C	170	GLN
1	D	63	HIS
1	D	92	GLN
1	D	156	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 ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 13 ligands modelled in this entry, 7 are monoatomic - leaving 6 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	ADP	B	702	2	24,29,29	1.16	2 (8%)	29,45,45	1.54	4 (13%)
3	ADP	C	703	2	24,29,29	1.24	2 (8%)	29,45,45	1.58	5 (17%)
3	ADP	D	704	2	24,29,29	1.29	2 (8%)	29,45,45	1.62	4 (13%)
3	ADP	A	701	2	24,29,29	1.14	2 (8%)	29,45,45	1.52	4 (13%)

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	ADP	B	702	2	-	2/12/32/32	0/3/3/3
3	ADP	C	703	2	-	0/12/32/32	0/3/3/3
3	ADP	D	704	2	-	2/12/32/32	0/3/3/3
3	ADP	A	701	2	-	1/12/32/32	0/3/3/3

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	703	ADP	O4'-C1'	4.40	1.47	1.41
3	D	704	ADP	O4'-C1'	4.25	1.47	1.41
3	B	702	ADP	O4'-C1'	4.01	1.46	1.41
3	A	701	ADP	O4'-C1'	3.92	1.46	1.41
3	D	704	ADP	C8-N7	-2.41	1.30	1.34
3	A	701	ADP	C8-N7	-2.13	1.30	1.34
3	B	702	ADP	C8-N7	-2.13	1.30	1.34
3	C	703	ADP	C8-N7	-2.09	1.31	1.34

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	704	ADP	PA-O3A-PB	-5.03	115.55	132.83
3	D	704	ADP	N3-C2-N1	-4.71	121.31	128.68
3	A	701	ADP	N3-C2-N1	-4.69	121.35	128.68
3	B	702	ADP	N3-C2-N1	-4.68	121.37	128.68
3	C	703	ADP	N3-C2-N1	-4.60	121.49	128.68
3	C	703	ADP	PA-O3A-PB	-4.46	117.51	132.83
3	A	701	ADP	PA-O3A-PB	-4.21	118.38	132.83
3	B	702	ADP	PA-O3A-PB	-3.96	119.24	132.83
3	B	702	ADP	O2B-PB-O3A	2.68	113.63	104.64
3	C	703	ADP	O2B-PB-O3A	2.67	113.58	104.64
3	C	703	ADP	PA-O5'-C5'	-2.42	107.47	121.68
3	A	701	ADP	O2B-PB-O3A	2.41	112.72	104.64
3	A	701	ADP	PA-O5'-C5'	-2.31	108.16	121.68
3	D	704	ADP	PA-O5'-C5'	-2.27	108.35	121.68
3	B	702	ADP	PA-O5'-C5'	-2.26	108.45	121.68
3	D	704	ADP	O2B-PB-O3A	2.13	111.78	104.64
3	C	703	ADP	C4-C5-N7	-2.01	107.31	109.40

There are no chirality outliers.

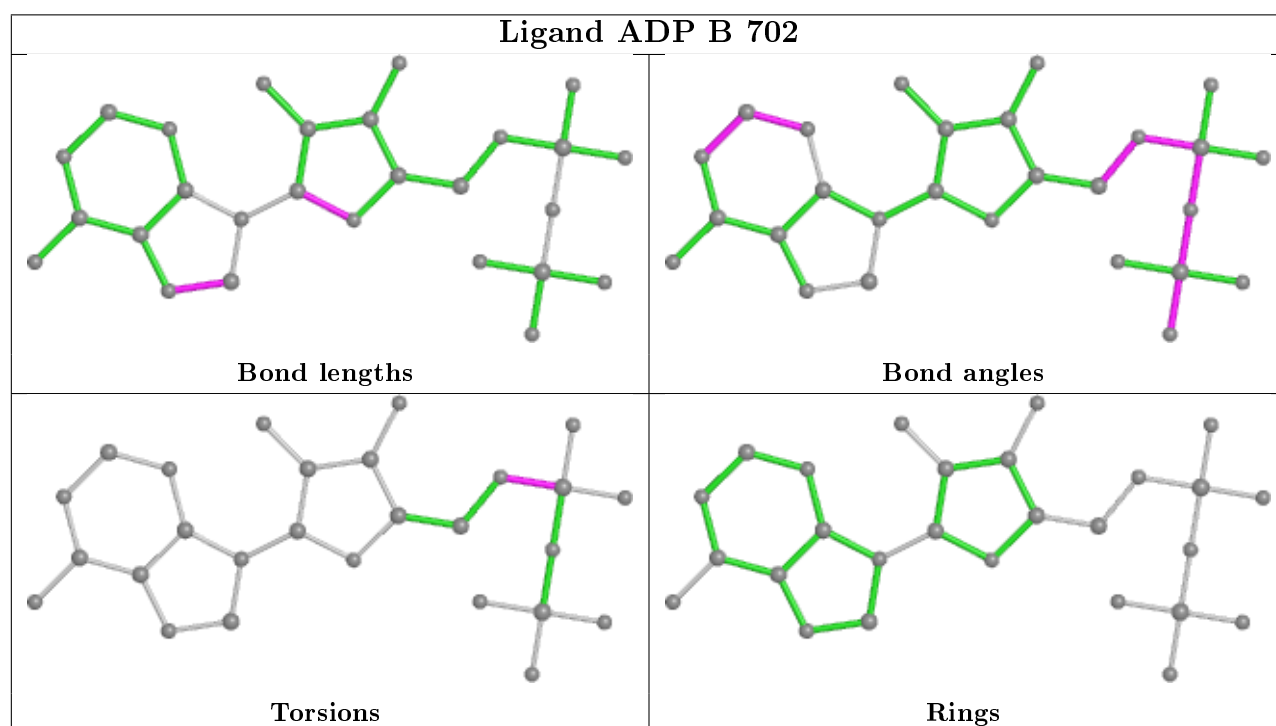
All (5) torsion outliers are listed below:

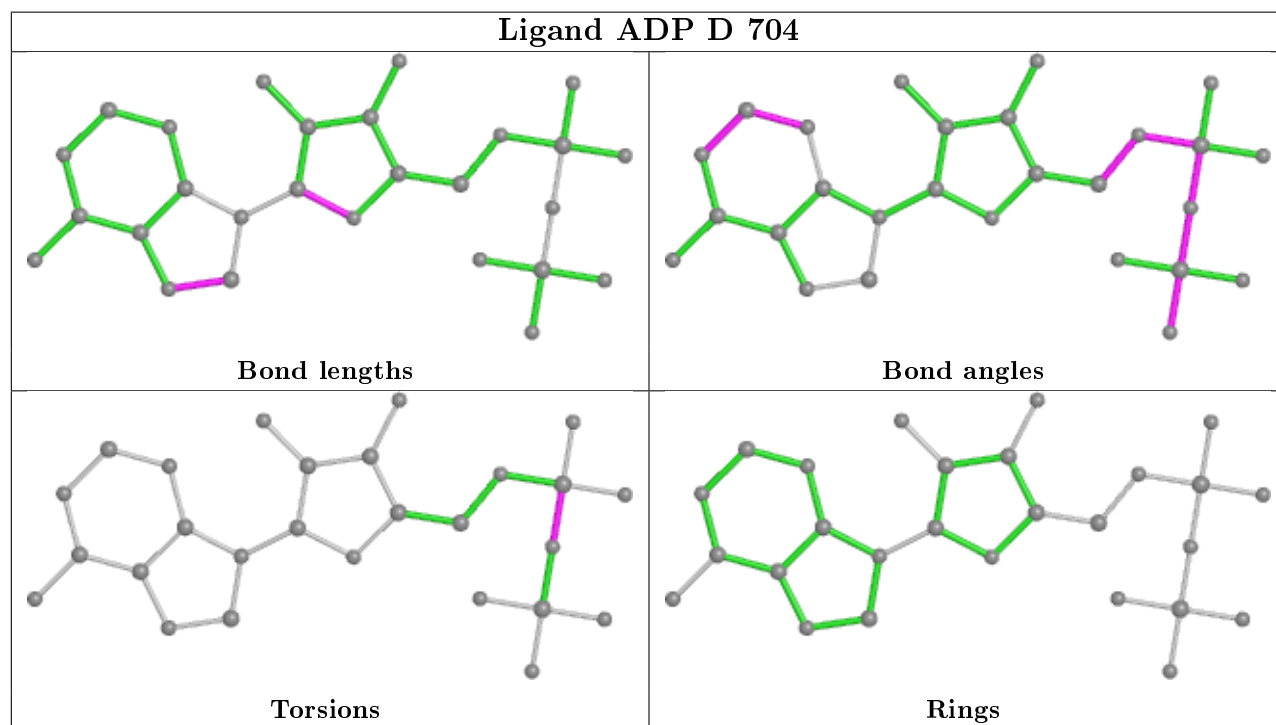
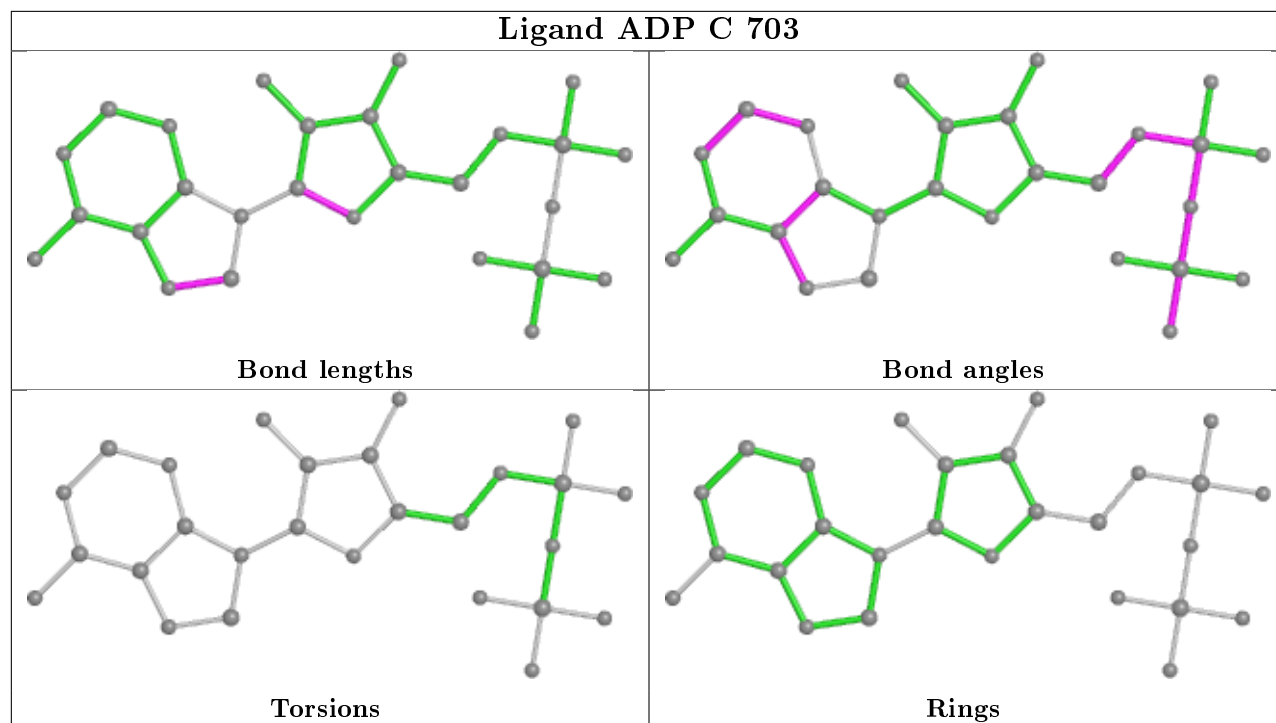
Mol	Chain	Res	Type	Atoms
3	B	702	ADP	C5'-O5'-PA-O3A
3	A	701	ADP	PB-O3A-PA-O2A
3	D	704	ADP	PB-O3A-PA-O2A
3	D	704	ADP	PB-O3A-PA-O1A
3	B	702	ADP	C5'-O5'-PA-O2A

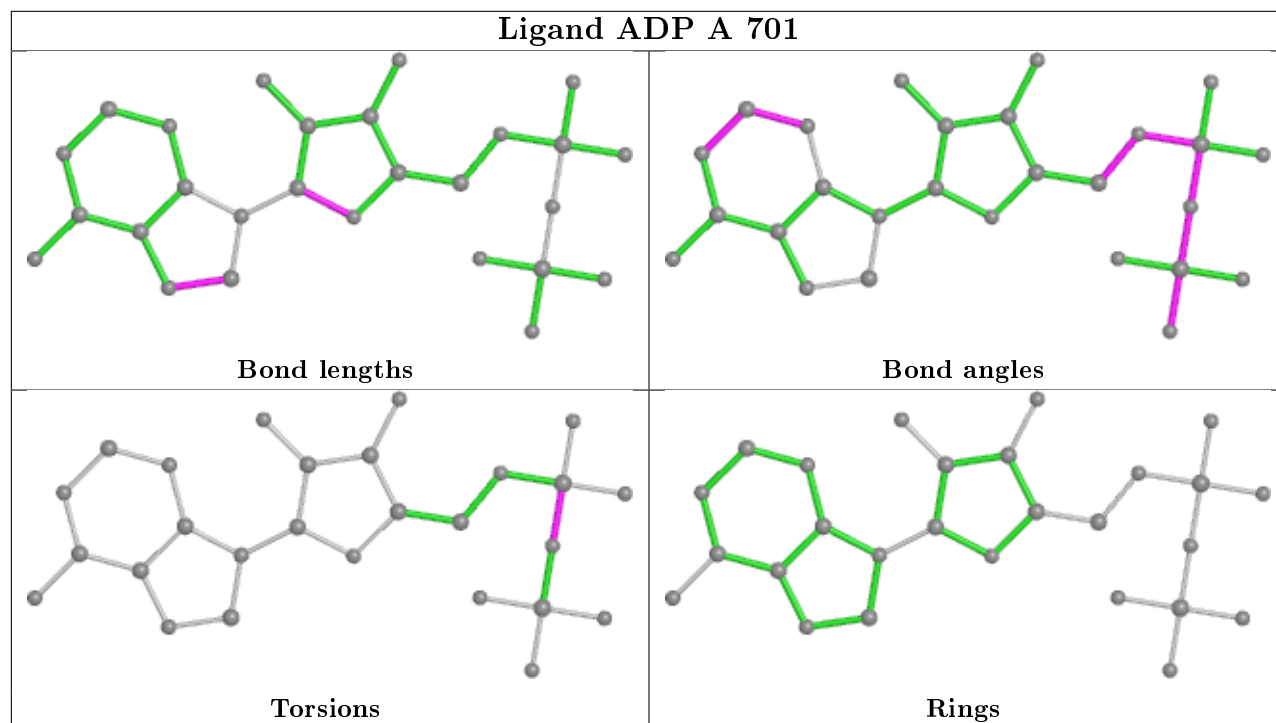
There are no ring outliers.

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 ⓘ

### 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	319/319 (100%)	0.13	4 (1%) 77 79	11, 18, 31, 44	0
1	B	312/319 (97%)	0.25	12 (3%) 40 43	10, 20, 39, 49	0
1	C	316/319 (99%)	0.36	18 (5%) 23 26	13, 24, 43, 54	0
1	D	319/319 (100%)	0.56	27 (8%) 10 12	12, 23, 48, 69	0
All	All	1266/1276 (99%)	0.33	61 (4%) 30 33	10, 21, 41, 69	0

All (61) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	231	PRO	13.0
1	C	229	TYR	8.2
1	D	291	PRO	7.4
1	D	229	TYR	7.3
1	C	223	TYR	7.2
1	D	225	TYR	6.5
1	A	292	THR	6.3
1	D	232	GLY	5.9
1	D	230	THR	5.6
1	C	222	PHE	5.0
1	D	292	THR	4.6
1	C	157	THR	4.3
1	D	228	LYS	4.3
1	D	290	THR	4.2
1	A	291	PRO	4.0
1	C	233	ARG	3.8
1	B	233	ARG	3.7
1	B	223	TYR	3.6
1	C	227	THR	3.5
1	C	230	THR	3.5
1	C	231	PRO	3.5

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Mol	Chain	Res	Type	RSRZ
1	D	227	THR	3.4
1	D	233	ARG	3.3
1	B	157	THR	3.3
1	B	222	PHE	3.2
1	B	181	ARG	3.2
1	D	226	GLU	3.2
1	C	160	SER	3.2
1	D	54	GLU	3.0
1	D	26	ARG	3.0
1	B	244	PRO	2.9
1	B	160	SER	2.8
1	C	26	ARG	2.7
1	D	221	PRO	2.7
1	C	169	PHE	2.7
1	D	19	LEU	2.7
1	C	236	LEU	2.6
1	A	293	SER	2.6
1	B	168	ARG	2.6
1	D	139	GLY	2.6
1	D	59	PRO	2.5
1	C	232	GLY	2.5
1	D	234	ALA	2.4
1	B	170	GLN	2.4
1	D	224	ASP	2.4
1	D	223	TYR	2.4
1	B	234	ALA	2.3
1	D	84	ARG	2.3
1	C	56	LYS	2.3
1	B	182	TYR	2.2
1	D	289	PHE	2.2
1	D	306	TYR	2.2
1	D	219	GLU	2.2
1	B	159	SER	2.2
1	C	244	PRO	2.2
1	D	236	LEU	2.1
1	C	54	GLU	2.1
1	C	84	ARG	2.1
1	D	169	PHE	2.1
1	C	47	GLU	2.1
1	A	236	LEU	2.0

## 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 carbohydrates 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, 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.

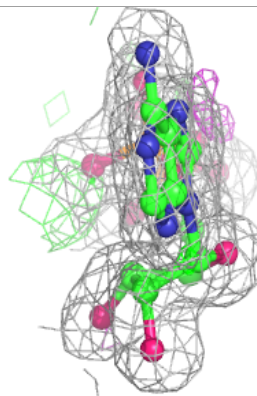
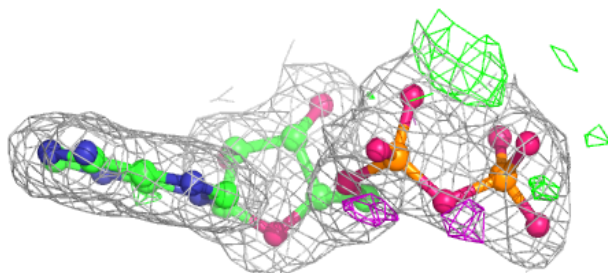
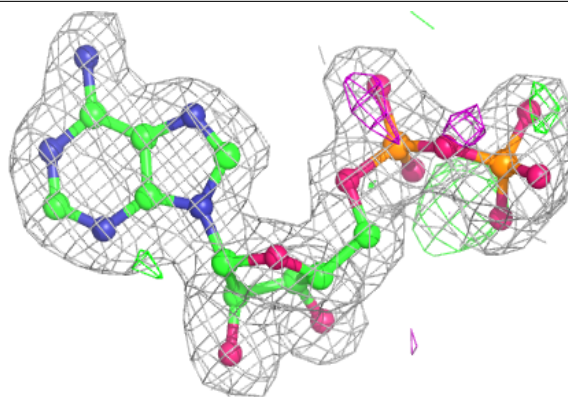
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	MG	D	817	1/1	0.53	0.30	55,55,55,55	0
4	DAL	D	802	6/6	0.69	0.21	30,33,35,37	0
4	DAL	A	801	6/6	0.79	0.18	14,21,26,28	0
2	MG	C	815	1/1	0.88	0.11	24,24,24,24	0
2	MG	B	813	1/1	0.93	0.12	18,18,18,18	0
2	MG	A	812	1/1	0.94	0.14	21,21,21,21	0
2	MG	A	811	1/1	0.96	0.18	26,26,26,26	0
3	ADP	D	704	27/27	0.96	0.12	14,17,23,24	0
3	ADP	C	703	27/27	0.96	0.10	18,20,22,25	0
2	MG	C	816	1/1	0.97	0.14	25,25,25,25	0
3	ADP	B	702	27/27	0.97	0.11	15,17,19,20	0
3	ADP	A	701	27/27	0.97	0.10	11,13,15,16	0
2	MG	B	814	1/1	0.97	0.14	21,21,21,21	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.

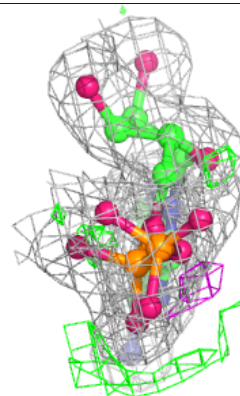
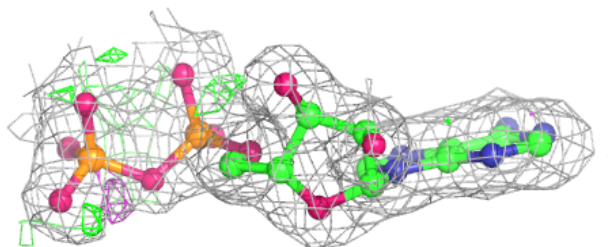
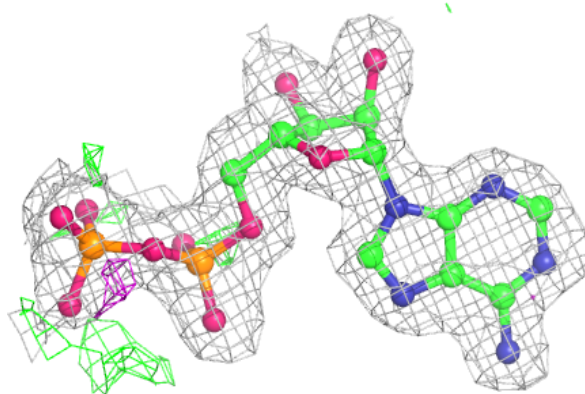


**Electron density around ADP D 704:**

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

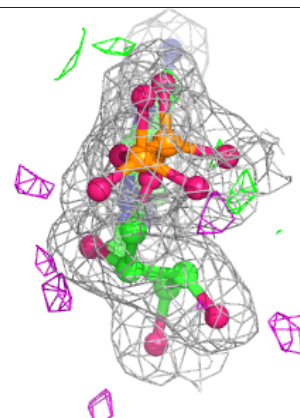
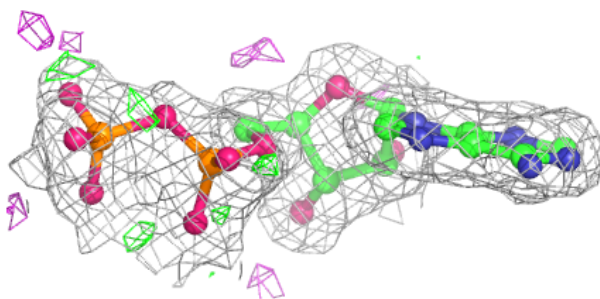
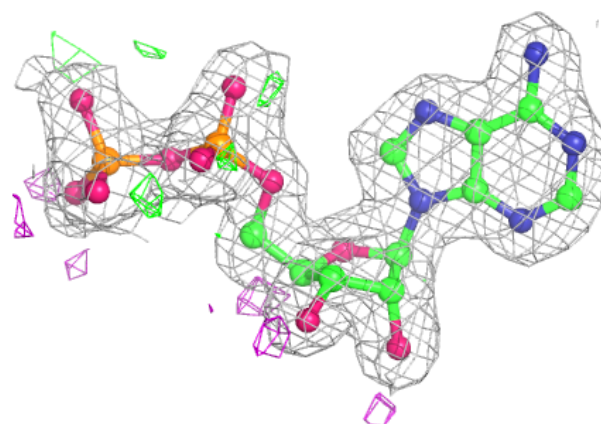
**Electron density around ADP C 703:**

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

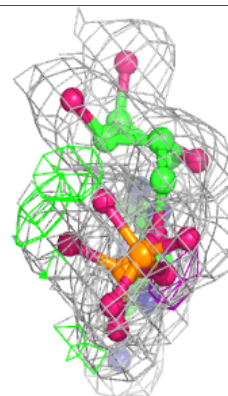
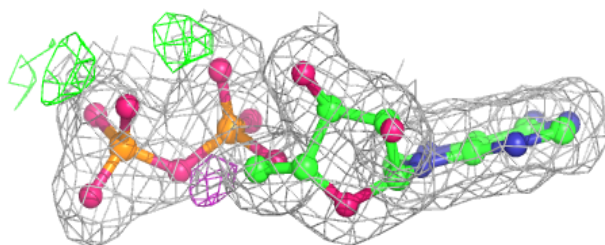
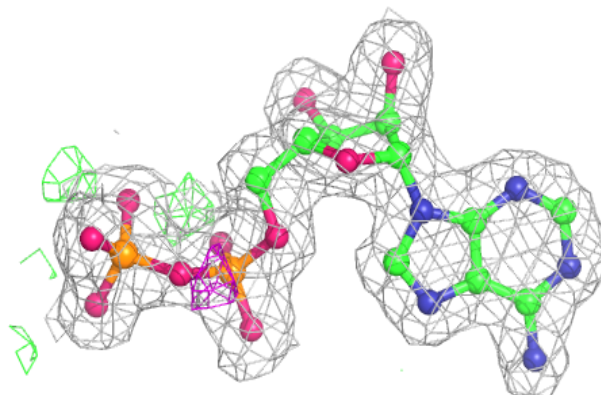


**Electron density around ADP B 702:**

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

**Electron density around ADP A 701:**

$2mF_o-DF_c$  (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.