



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 5, 2023 – 03:17 AM EDT

PDB ID : 3V4R
Title : Crystal structure of a UvrB dimer-DNA complex
Authors : Webster, M.P.J.; Jukes, R.; Barrett, T.
Deposited on : 2011-12-15
Resolution : 3.25 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.35
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) : 2.35

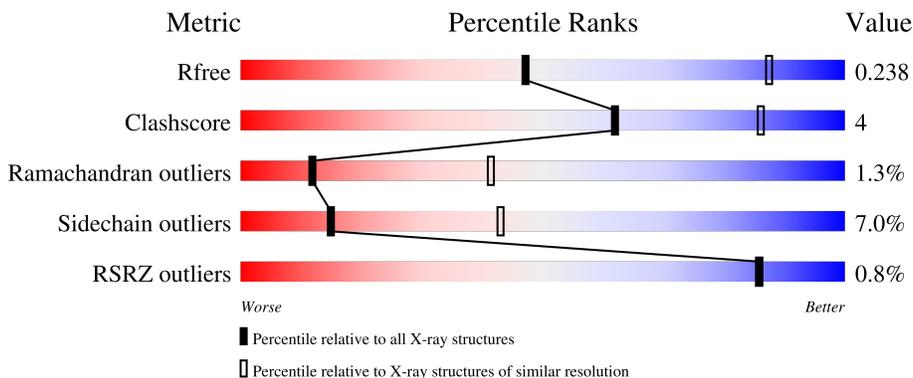
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.25 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	1191 (3.30-3.22)
Clashscore	141614	1251 (3.30-3.22)
Ramachandran outliers	138981	1229 (3.30-3.22)
Sidechain outliers	138945	1228 (3.30-3.22)
RSRZ outliers	127900	1154 (3.30-3.22)

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	667	
1	B	667	
2	C	8	
2	D	8	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 8815 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 UvrABC system protein B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	586	Total	C	N	O	S	0	0	0
			4321	2738	754	815	14			
1	B	579	Total	C	N	O	S	0	0	0
			4205	2663	739	790	13			

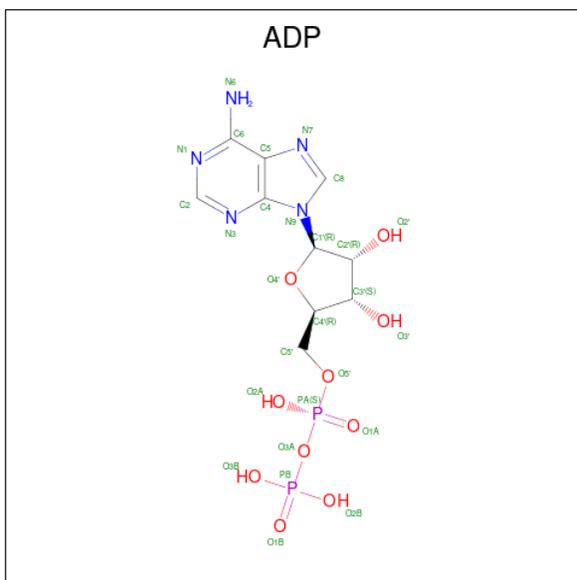
There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	HIS	-	expression tag	UNP P37954
A	-4	HIS	-	expression tag	UNP P37954
A	-3	HIS	-	expression tag	UNP P37954
A	-2	HIS	-	expression tag	UNP P37954
A	-1	HIS	-	expression tag	UNP P37954
A	0	HIS	-	expression tag	UNP P37954
B	-5	HIS	-	expression tag	UNP P37954
B	-4	HIS	-	expression tag	UNP P37954
B	-3	HIS	-	expression tag	UNP P37954
B	-2	HIS	-	expression tag	UNP P37954
B	-1	HIS	-	expression tag	UNP P37954
B	0	HIS	-	expression tag	UNP P37954

- Molecule 2 is a DNA chain called DNA: 5'-TACTGTTT-3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	C	7	Total	C	N	O	P	0	0	0
			131	64	20	41	6			
2	D	7	Total	C	N	O	P	0	0	0
			123	59	19	39	6			

- Molecule 3 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: C₁₀H₁₅N₅O₁₀P₂).

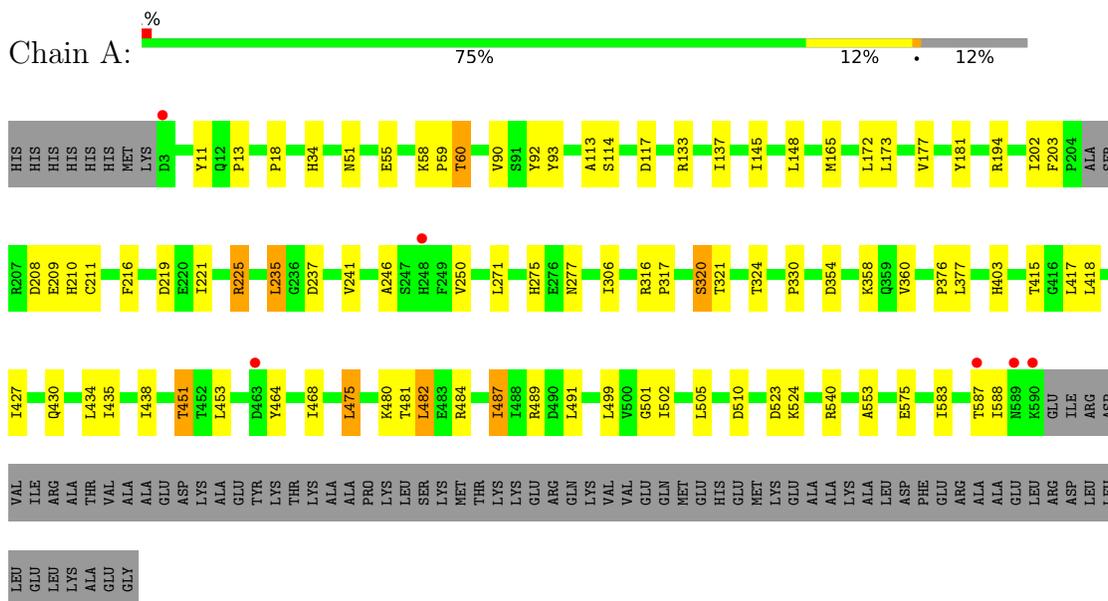


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
			Total	C	N	O			P
3	A	1	27	10	5	10	2	0	0
3	B	1	8	0	6	2	0	0	0

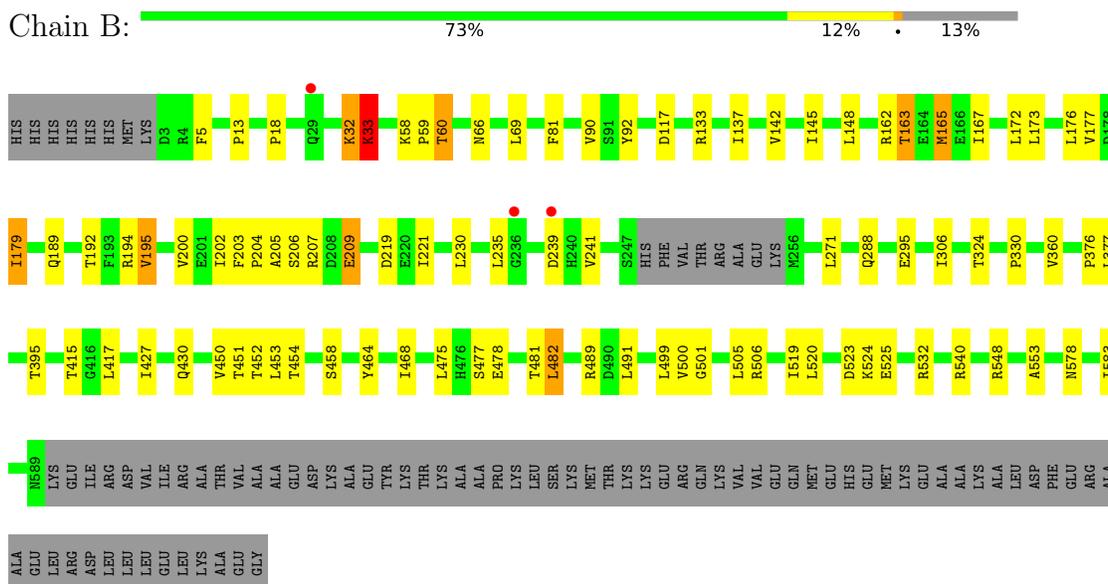
3 Residue-property plots

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

- Molecule 1: UvrABC system protein B

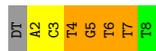


- Molecule 1: UvrABC system protein B

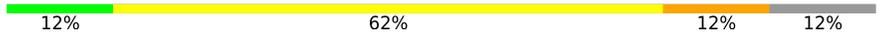


- Molecule 2: DNA: 5 -TACTGTTT-3

Chain C:  12% 25% 50% 12%



- Molecule 2: DNA: 5 -TACTGTTT-3

Chain D:  12% 62% 12% 12%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	95.94Å 100.31Å 163.63Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.91 – 3.25 19.91 – 3.25	Depositor EDS
% Data completeness (in resolution range)	100.0 (19.91-3.25) 100.0 (19.91-3.25)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.93 (at 3.22Å)	Xtrriage
Refinement program	BUSTER 2.10.0	Depositor
R, R_{free}	0.180 , 0.219 0.194 , 0.238	Depositor DCC
R_{free} test set	1297 reflections (5.10%)	wwPDB-VP
Wilson B-factor (Å ²)	75.2	Xtrriage
Anisotropy	0.173	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 65.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtrriage
Estimated twinning fraction	0.027 for k,h,-l	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	8815	wwPDB-VP
Average B, all atoms (Å ²)	68.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.70% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.52	0/4397	0.71	0/5996
1	B	0.50	0/4278	0.74	3/5843 (0.1%)
2	C	1.19	0/145	2.40	14/222 (6.3%)
2	D	1.22	0/136	2.50	12/209 (5.7%)
All	All	0.54	0/8956	0.85	29/12270 (0.2%)

There are no bond length outliers.

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	6	DT	O4'-C1'-N1	9.27	114.49	108.00
2	C	2	DA	O4'-C1'-N9	8.97	114.28	108.00
2	D	6	DT	O4'-C1'-N1	8.93	114.25	108.00
2	D	7	DT	C4-C5-C7	8.81	124.29	119.00
2	D	6	DT	P-O3'-C3'	8.37	129.75	119.70
2	D	3	DC	O4'-C1'-N1	7.70	113.39	108.00
2	D	2	DA	O4'-C1'-N9	7.60	113.32	108.00
2	C	3	DC	O4'-C1'-N1	6.95	112.86	108.00
2	C	6	DT	N3-C2-O2	-6.76	118.24	122.30
1	B	32	LYS	C-N-CA	6.65	138.32	121.70
2	C	6	DT	C6-C5-C7	-6.46	119.03	122.90
2	D	3	DC	N1-C2-O2	6.31	122.68	118.90
2	D	7	DT	C6-C5-C7	-6.31	119.12	122.90
2	C	6	DT	C4-C5-C7	5.97	122.58	119.00
2	D	4	DT	P-O3'-C3'	5.85	126.72	119.70
2	C	3	DC	N1-C2-O2	5.79	122.38	118.90
1	B	206	SER	C-N-CA	5.78	136.15	121.70
2	C	7	DT	O4'-C1'-N1	5.63	111.94	108.00
2	C	3	DC	C2-N1-C1'	5.61	124.97	118.80
2	C	4	DT	O4'-C1'-N1	5.58	111.90	108.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	4	DT	P-O3'-C3'	5.53	126.34	119.70
2	D	3	DC	P-O3'-C3'	5.49	126.29	119.70
2	D	2	DA	C3'-C2'-C1'	-5.30	96.14	102.50
1	B	33	LYS	N-CA-C	-5.29	96.72	111.00
2	C	5	DG	P-O3'-C3'	5.27	126.03	119.70
2	D	6	DT	C6-C5-C7	-5.11	119.83	122.90
2	C	4	DT	C6-C5-C7	-5.08	119.86	122.90
2	C	7	DT	P-O3'-C3'	5.07	125.78	119.70
2	D	3	DC	C1'-O4'-C4'	-5.04	105.06	110.10

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	4321	0	3986	38	0
1	B	4205	0	3844	35	0
2	C	131	0	78	3	0
2	D	123	0	70	1	0
3	A	27	0	12	0	0
3	B	8	0	0	0	0
All	All	8815	0	7990	74	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 4.

All (74) 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)
2:C:4:DT:H2''	2:C:5:DG:H5'	1.51	0.90
2:D:4:DT:H2''	2:D:5:DG:H5''	1.63	0.79
1:A:181:TYR:CE2	1:A:246:ALA:HB2	2.17	0.78
1:A:51:ASN:O	1:A:55:GLU:HG2	1.94	0.67
1:B:32:LYS:HA	1:B:33:LYS:HB2	1.79	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:450:VAL:HG22	1:B:500:VAL:HG13	1.84	0.60
1:A:435:ILE:HA	1:A:438:ILE:HD12	1.86	0.57
1:B:162:ARG:O	1:B:165:MET:HB2	2.04	0.57
1:A:481:THR:HG23	1:A:484:ARG:HH11	1.69	0.57
1:A:181:TYR:CD2	1:A:246:ALA:HB2	2.40	0.57
1:A:453:LEU:HD11	1:A:524:LYS:HB2	1.86	0.57
1:B:202:ILE:O	1:B:204:PRO:HD3	2.04	0.56
1:B:32:LYS:HA	1:B:33:LYS:CB	2.35	0.56
1:A:181:TYR:HE2	1:A:246:ALA:N	2.03	0.56
1:A:434:LEU:O	1:A:438:ILE:HG13	2.08	0.54
1:B:427:ILE:HA	1:B:430:GLN:HG3	1.89	0.53
2:C:6:DT:H2'	2:C:7:DT:C5	2.42	0.53
1:A:427:ILE:HA	1:A:430:GLN:HG3	1.89	0.53
1:A:475:LEU:HD11	1:A:487:ILE:HG21	1.89	0.53
1:B:5:PHE:HB3	1:B:81:PHE:HE2	1.74	0.53
1:A:145:ILE:HD12	1:A:376:PRO:HB2	1.91	0.52
1:B:173:LEU:O	1:B:177:VAL:HG23	2.10	0.51
1:A:173:LEU:O	1:A:177:VAL:HG23	2.11	0.49
1:A:208:ASP:O	1:A:210:HIS:N	2.45	0.49
1:A:354:ASP:OD1	1:A:358:LYS:NZ	2.44	0.49
1:B:60:THR:HG22	1:B:137:ILE:HG12	1.95	0.49
1:A:491:LEU:HD13	1:A:499:LEU:HG	1.94	0.49
1:A:202:ILE:O	1:A:211:CYS:HB2	2.13	0.48
1:B:477:SER:HA	1:B:506:ARG:HH22	1.78	0.48
1:B:505:LEU:HA	1:B:540:ARG:HH21	1.79	0.48
1:B:145:ILE:HD12	1:B:376:PRO:HB2	1.95	0.48
1:A:216:PHE:CE1	1:A:221:ILE:HG12	2.49	0.47
1:A:60:THR:HG22	1:A:137:ILE:HG12	1.94	0.47
1:A:464:TYR:CE1	1:A:468:ILE:HD11	2.50	0.47
1:B:5:PHE:HB3	1:B:81:PHE:CE2	2.50	0.47
1:B:451:THR:HG23	1:B:519:ILE:HA	1.97	0.47
1:A:482:LEU:HD12	1:B:482:LEU:HD23	1.97	0.46
1:B:464:TYR:CE1	1:B:468:ILE:HD11	2.49	0.46
1:A:316:ARG:NH2	1:A:321:THR:O	2.49	0.46
1:B:59:PRO:HG2	1:B:330:PRO:HG2	1.98	0.46
1:A:59:PRO:HG2	1:A:330:PRO:HG2	1.97	0.46
1:A:324:THR:HG21	1:A:377:LEU:HG	1.98	0.45
1:A:523:ASP:HB3	1:A:553:ALA:HB1	1.98	0.45
1:B:209:GLU:HG2	1:B:230:LEU:HD21	1.97	0.45
1:B:578:ASN:HA	1:B:583:ILE:HD12	1.99	0.45
1:B:195:VAL:HG22	1:B:200:VAL:HG22	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:451:THR:HA	1:B:501:GLY:O	2.16	0.45
1:B:395:THR:HB	1:B:532:ARG:HG2	1.97	0.45
1:A:317:PRO:O	1:A:320:SER:HB2	2.18	0.45
1:A:451:THR:HA	1:A:501:GLY:O	2.16	0.44
1:B:66:ASN:HA	2:C:5:DG:H5''	2.00	0.44
1:A:415:THR:HB	1:A:417:LEU:HD23	1.99	0.43
1:B:415:THR:HB	1:B:417:LEU:HD23	1.99	0.43
1:B:163:THR:O	1:B:221:ILE:O	2.37	0.43
1:A:13:PRO:HB2	1:A:18:PRO:HG3	2.01	0.43
1:A:275:HIS:C	1:A:277:ASN:H	2.21	0.43
1:B:452:THR:HG22	1:B:520:LEU:HD12	2.01	0.43
1:B:523:ASP:HB3	1:B:553:ALA:HB1	2.00	0.42
1:B:176:LEU:O	1:B:179:ILE:HG13	2.19	0.42
1:A:453:LEU:HD11	1:A:524:LYS:CB	2.48	0.42
1:B:453:LEU:HD11	1:B:524:LYS:HB2	2.01	0.42
1:A:418:LEU:HD22	1:A:575:GLU:HG3	2.01	0.42
1:B:163:THR:HG23	1:B:239:ASP:HA	2.02	0.42
1:B:165:MET:HG2	1:B:167:ILE:HG22	2.02	0.42
1:A:505:LEU:HA	1:A:540:ARG:HH21	1.85	0.41
1:B:13:PRO:HB2	1:B:18:PRO:HG3	2.02	0.41
1:A:34:HIS:CG	1:A:403:HIS:HB3	2.55	0.41
1:A:93:TYR:CD2	1:A:113:ALA:HB1	2.55	0.41
1:A:434:LEU:HG	1:A:438:ILE:HD11	2.03	0.41
1:B:324:THR:HG21	1:B:377:LEU:HG	2.01	0.41
1:A:489:ARG:HG3	1:B:489:ARG:HG3	2.03	0.41
1:B:491:LEU:HD13	1:B:499:LEU:HG	2.02	0.41
1:A:165:MET:HG2	1:A:221:ILE:HD12	2.02	0.40
1:A:225:ARG:HB3	1:A:237:ASP:HA	2.04	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	582/667 (87%)	546 (94%)	30 (5%)	6 (1%)	15	47
1	B	575/667 (86%)	543 (94%)	23 (4%)	9 (2%)	9	37
All	All	1157/1334 (87%)	1089 (94%)	53 (5%)	15 (1%)	12	41

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	209	GLU
1	A	587	THR
1	B	33	LYS
1	B	163	THR
1	B	165	MET
1	B	179	ILE
1	B	205	ALA
1	B	207	ARG
1	B	235	LEU
1	A	235	LEU
1	A	250	VAL
1	B	525	GLU
1	A	588	ILE
1	A	482	LEU
1	B	195	VAL

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	413/592 (70%)	386 (94%)	27 (6%)	17	46
1	B	392/592 (66%)	363 (93%)	29 (7%)	13	40
All	All	805/1184 (68%)	749 (93%)	56 (7%)	15	43

All (56) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	TYR

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Mol	Chain	Res	Type
1	A	58	LYS
1	A	60	THR
1	A	90	VAL
1	A	92	TYR
1	A	114	SER
1	A	117	ASP
1	A	133	ARG
1	A	148	LEU
1	A	172	LEU
1	A	194	ARG
1	A	203	PHE
1	A	219	ASP
1	A	225	ARG
1	A	235	LEU
1	A	241	VAL
1	A	271	LEU
1	A	306	ILE
1	A	320	SER
1	A	360	VAL
1	A	451	THR
1	A	475	LEU
1	A	480	LYS
1	A	487	ILE
1	A	502	ILE
1	A	510	ASP
1	A	583	ILE
1	B	58	LYS
1	B	60	THR
1	B	69	LEU
1	B	90	VAL
1	B	92	TYR
1	B	117	ASP
1	B	133	ARG
1	B	142	VAL
1	B	148	LEU
1	B	172	LEU
1	B	189	GLN
1	B	192	THR
1	B	194	ARG
1	B	203	PHE
1	B	209	GLU
1	B	219	ASP

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Mol	Chain	Res	Type
1	B	241	VAL
1	B	271	LEU
1	B	288	GLN
1	B	295	GLU
1	B	306	ILE
1	B	360	VAL
1	B	454	THR
1	B	458	SER
1	B	475	LEU
1	B	478	GLU
1	B	481	THR
1	B	482	LEU
1	B	548	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	355	GLN
1	A	364	HIS
1	B	277	ASN
1	B	355	GLN
1	B	359	GLN
1	B	364	HIS
1	B	439	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry

2 ligands 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
3	ADP	A	701	-	24,29,29	0.66	0	29,45,45	1.01	2 (6%)
3	ADP	B	701	-	3,7,29	0.36	0	6,10,45	0.94	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
3	ADP	A	701	-	-	0/12/32/32	0/3/3/3
3	ADP	B	701	-	-	1/2/5/32	-

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	701	ADP	O5'-C5'-C4'	2.62	118.02	108.99
3	A	701	ADP	C5-C6-N6	2.32	123.88	120.35

There are no chirality outliers.

All (1) torsion outliers are listed below:

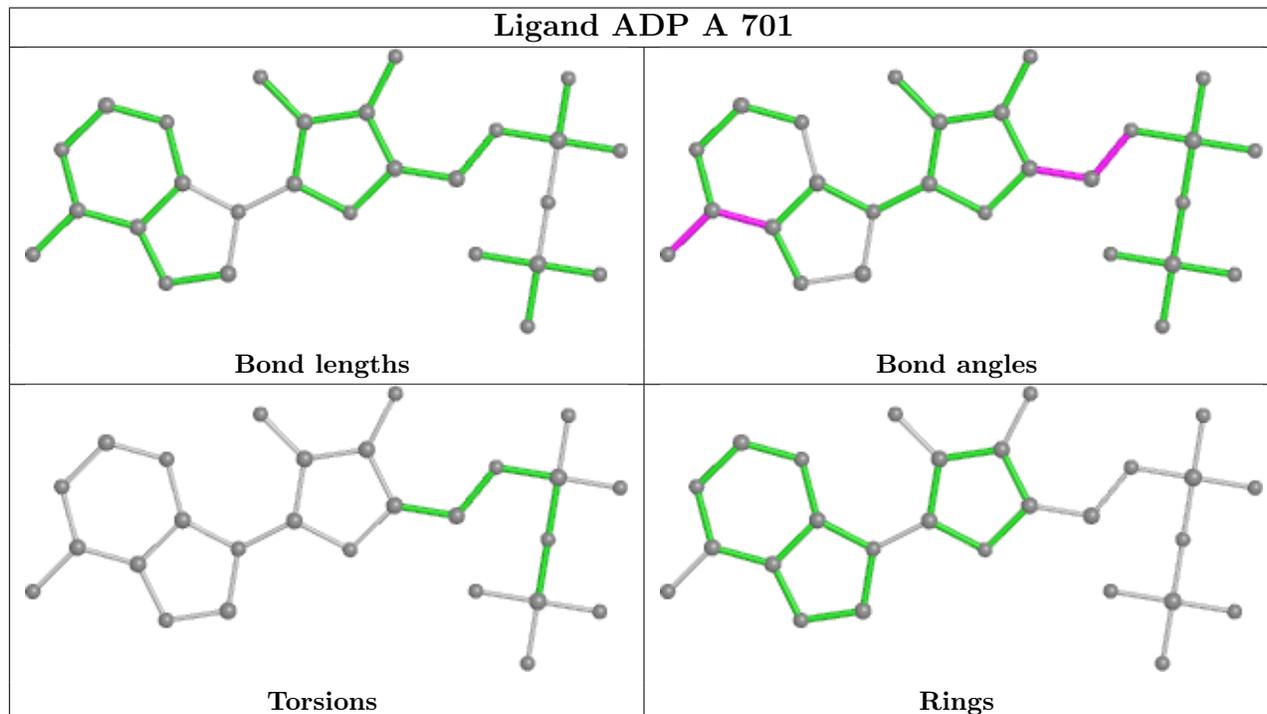
Mol	Chain	Res	Type	Atoms
3	B	701	ADP	PA-O3A-PB-O3B

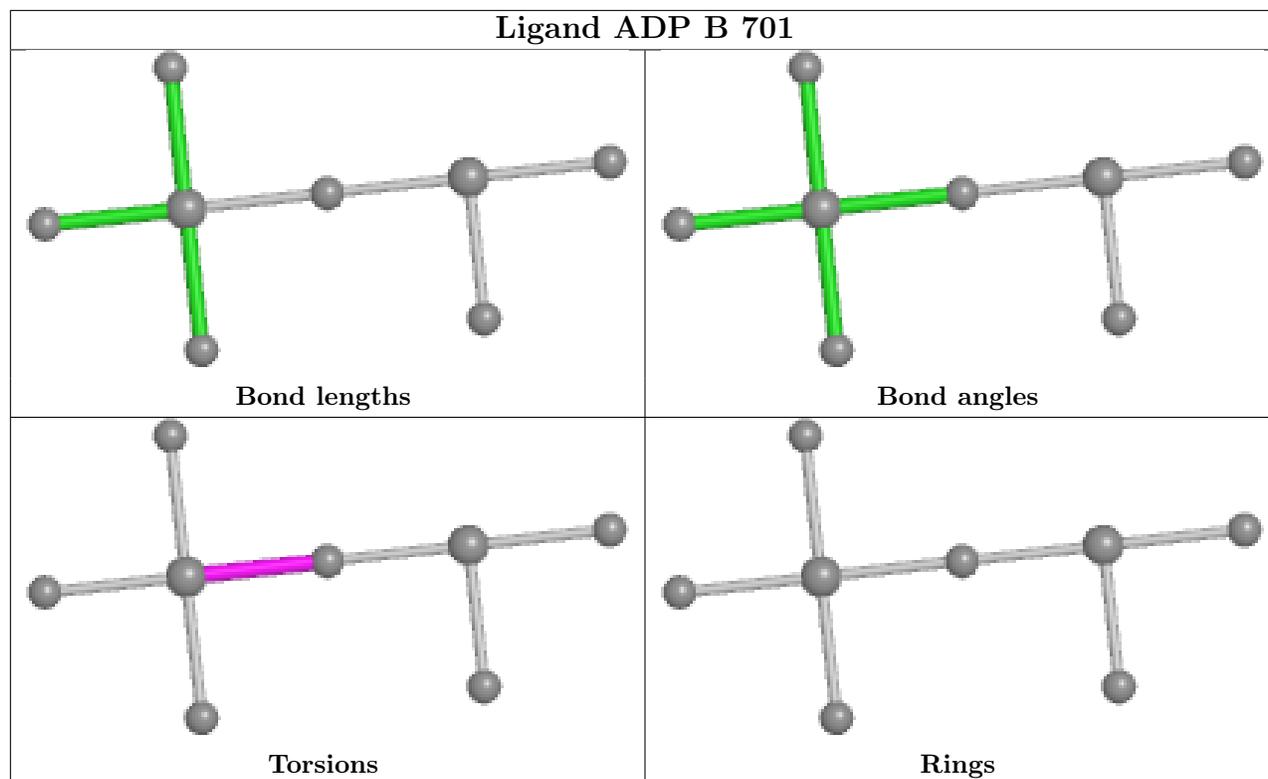
There are no ring outliers.

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	586/667 (87%)	-0.51	6 (1%) 82 82	32, 65, 94, 128	0
1	B	579/667 (86%)	-0.47	3 (0%) 91 90	36, 69, 98, 113	0
2	C	7/8 (87%)	0.17	0 100 100	79, 89, 95, 109	0
2	D	7/8 (87%)	-0.02	0 100 100	72, 75, 85, 88	0
All	All	1179/1350 (87%)	-0.48	9 (0%) 86 86	32, 68, 97, 128	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	589	ASN	3.3
1	A	587	THR	3.0
1	B	29	GLN	2.9
1	B	239	ASP	2.7
1	A	248	HIS	2.7
1	A	590	LYS	2.6
1	A	3	ASP	2.4
1	A	463	ASP	2.1
1	B	236	GLY	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](#)

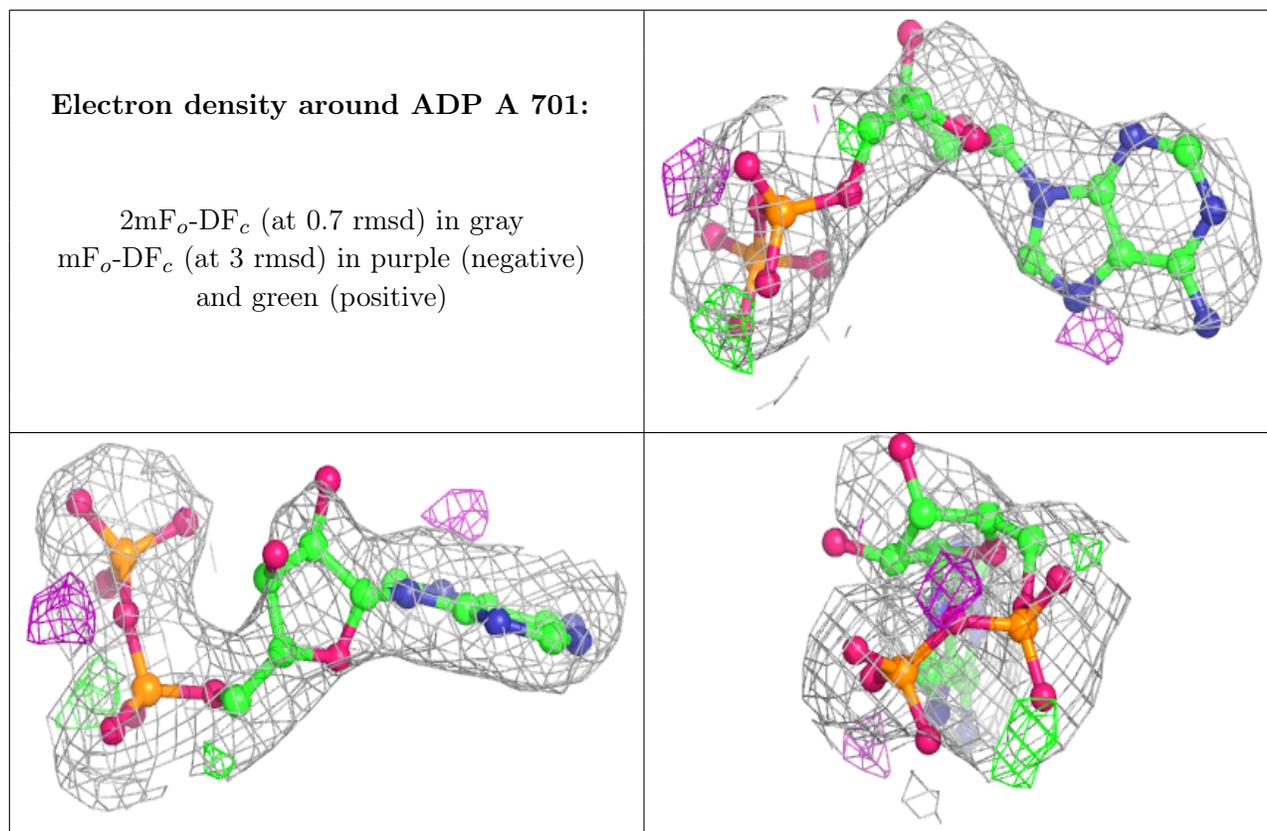
There are no monosaccharides in this entry.

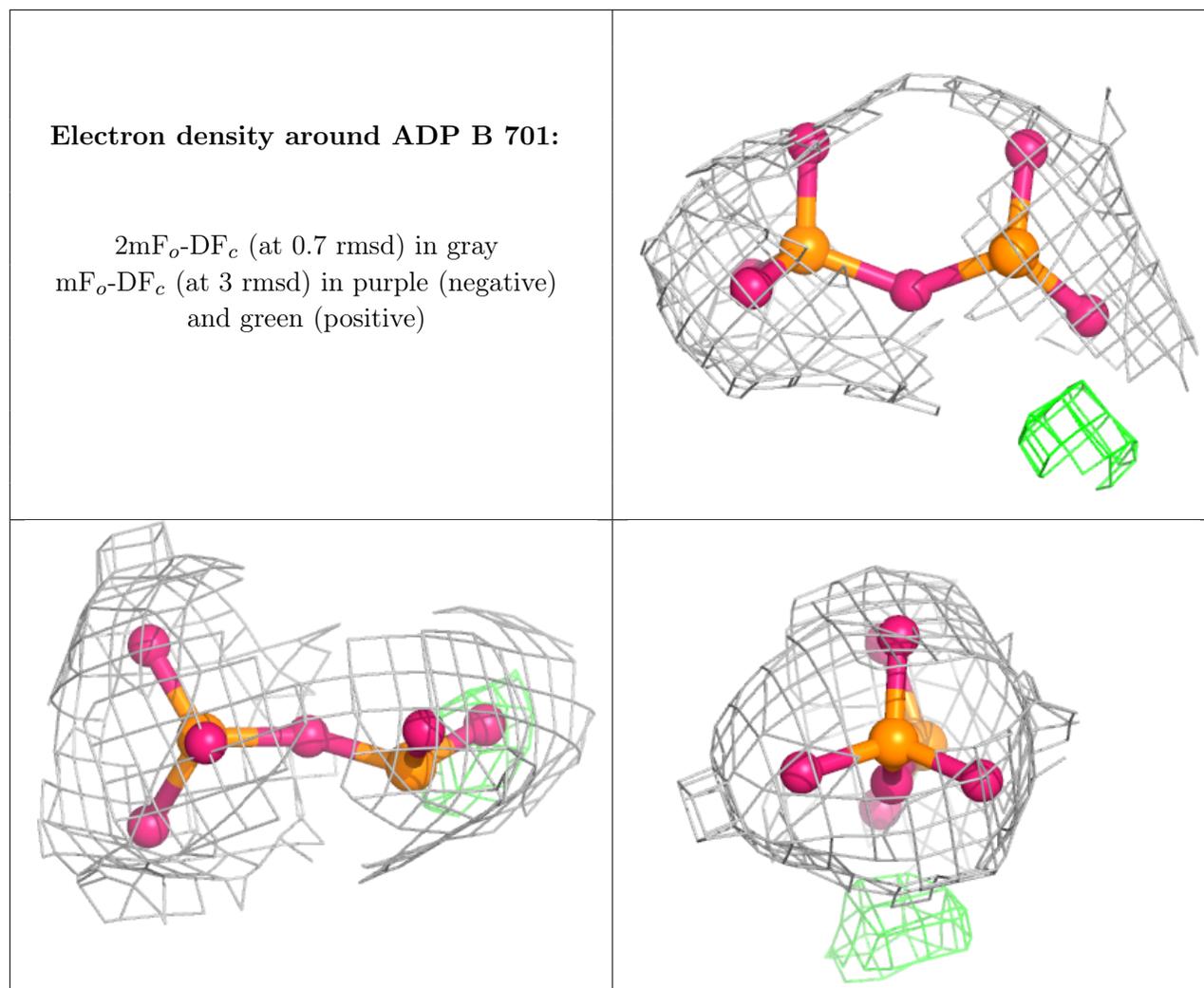
6.4 Ligands [i](#)

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(Å ²)	Q<0.9
3	ADP	A	701	27/27	0.87	0.26	100,105,109,110	0
3	ADP	B	701	8/27	0.94	0.13	117,119,121,122	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.





6.5 Other polymers [i](#)

There are no such residues in this entry.