



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 9, 2023 – 11:02 PM EDT

PDB ID : 7RTG  
Title : Crystal Structure of the Human Adenosine Deaminase 1  
Authors : Ma, M.T.; Lieberman, R.L.; Blazeck, J.; Jennings, M.R.  
Deposited on : 2021-08-13  
Resolution : 2.59 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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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  
Xtriage (Phenix) : 1.13  
EDS : 2.35.1  
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

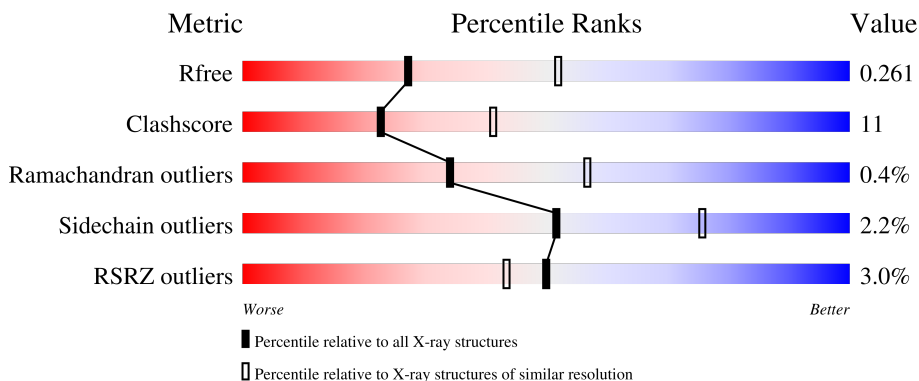
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.59 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	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

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  $\geq 3$ , 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	372	
1	B	372	

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5588 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 Adenosine deaminase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	349	2771	1764	475	519	13	0	1	0
1	B	352	2791	1776	478	524	13	0	1	0

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	364	GLU	-	expression tag	UNP P00813
A	365	ASN	-	expression tag	UNP P00813
A	366	LEU	-	expression tag	UNP P00813
A	367	TYR	-	expression tag	UNP P00813
A	368	PHE	-	expression tag	UNP P00813
A	369	GLN	-	expression tag	UNP P00813
A	370	SER	-	expression tag	UNP P00813
A	371	GLY	-	expression tag	UNP P00813
A	372	GLY	-	expression tag	UNP P00813
B	364	GLU	-	expression tag	UNP P00813
B	365	ASN	-	expression tag	UNP P00813
B	366	LEU	-	expression tag	UNP P00813
B	367	TYR	-	expression tag	UNP P00813
B	368	PHE	-	expression tag	UNP P00813
B	369	GLN	-	expression tag	UNP P00813
B	370	SER	-	expression tag	UNP P00813
B	371	GLY	-	expression tag	UNP P00813
B	372	GLY	-	expression tag	UNP P00813

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total 1	Zn 1	0	0
2	B	1	Total 1	Zn 1	0	0

- Molecule 3 is water.

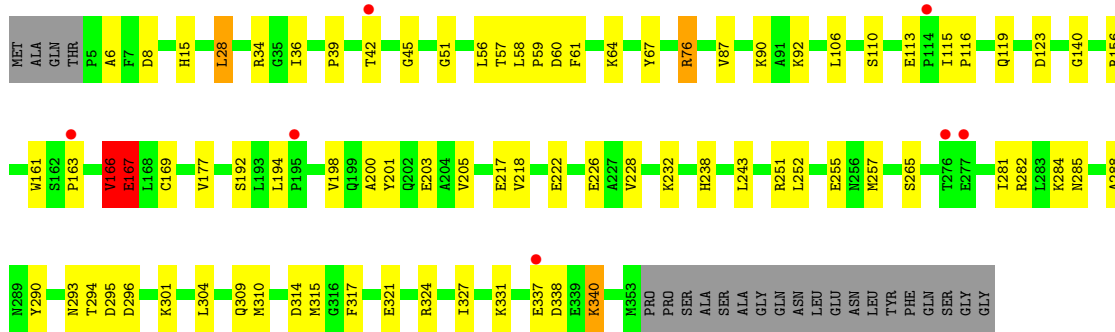
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	12	Total 12	O 12	0	0
3	B	12	Total 12	O 12	0	0

### 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.

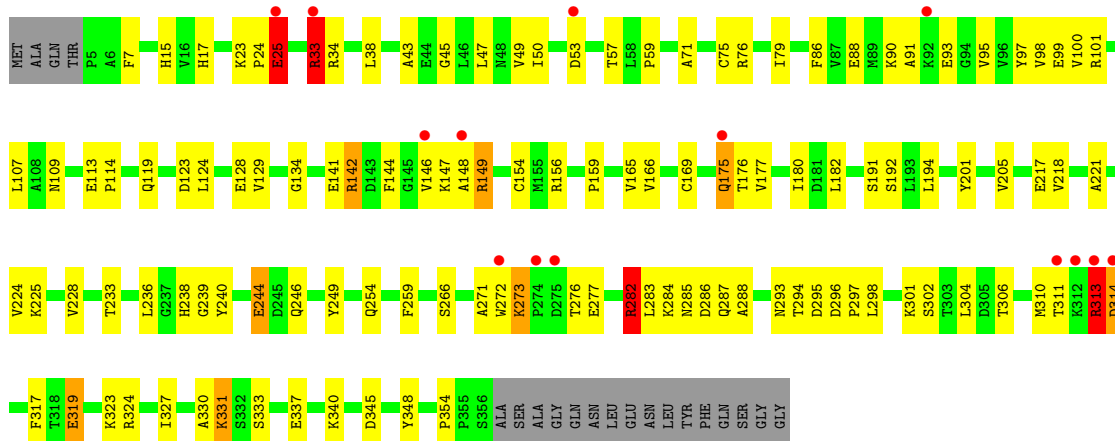
- Molecule 1: Adenosine deaminase

Chain A: 



- Molecule 1: Adenosine deaminase

Chain B: 



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	80.91Å 49.52Å 89.08Å 90.00° 96.18° 90.00°	Depositor
Resolution (Å)	42.17 – 2.59 43.22 – 2.59	Depositor EDS
% Data completeness (in resolution range)	56.6 (42.17-2.59) 70.3 (43.22-2.59)	Depositor EDS
$R_{merge}$	0.12	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.35 (at 2.58Å)	Xtrriage
Refinement program	PHENIX 1.14_3260	Depositor
R, $R_{free}$	0.201 , 0.261 0.202 , 0.261	Depositor DCC
$R_{free}$ test set	1682 reflections (10.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	42.0	Xtrriage
Anisotropy	0.036	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 43.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	5588	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	42.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.57% 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:  
ZN

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	0.48	2/2834 (0.1%)	0.70	6/3839 (0.2%)
1	B	0.51	4/2856 (0.1%)	1.14	30/3871 (0.8%)
All	All	0.49	6/5690 (0.1%)	0.95	36/7710 (0.5%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
1	B	1	3
All	All	1	5

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	76	ARG	CB-CG	-8.74	1.28	1.52
1	A	166	VAL	CB-CG1	-8.32	1.35	1.52
1	B	244	GLU	CB-CG	7.06	1.65	1.52
1	B	25	GLU	CG-CD	6.84	1.62	1.51
1	B	273	LYS	CE-NZ	6.36	1.65	1.49
1	B	25	GLU	CD-OE2	5.07	1.31	1.25

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	33	ARG	NE-CZ-NH2	-33.48	103.56	120.30
1	B	244	GLU	CA-CB-CG	15.18	146.80	113.40
1	B	282	ARG	CG-CD-NE	-14.01	82.38	111.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	33	ARG	NE-CZ-NH1	13.59	127.09	120.30
1	B	282	ARG	CA-CB-CG	13.30	142.66	113.40
1	B	313	ARG	NE-CZ-NH2	-13.11	113.75	120.30
1	A	76	ARG	NE-CZ-NH1	12.22	126.41	120.30
1	B	25	GLU	CG-CD-OE2	-12.22	93.86	118.30
1	A	76	ARG	NE-CZ-NH2	-11.41	114.59	120.30
1	B	282	ARG	CB-CA-C	-8.80	92.80	110.40
1	B	244	GLU	CB-CA-C	-8.61	93.18	110.40
1	B	244	GLU	N-CA-CB	8.38	125.68	110.60
1	A	76	ARG	CB-CG-CD	-8.32	89.97	111.60
1	B	25	GLU	CG-CD-OE1	8.15	134.61	118.30
1	B	175	GLN	N-CA-CB	8.09	125.16	110.60
1	B	282	ARG	N-CA-CB	7.98	124.96	110.60
1	B	313	ARG	NE-CZ-NH1	7.89	124.25	120.30
1	B	23	LYS	CG-CD-CE	7.74	135.11	111.90
1	B	33	ARG	NH1-CZ-NH2	7.36	127.50	119.40
1	B	273	LYS	CD-CE-NZ	-7.10	95.37	111.70
1	B	273	LYS	CG-CD-CE	7.02	132.97	111.90
1	A	28	LEU	CB-CG-CD2	6.75	122.48	111.00
1	B	306	THR	OG1-CB-CG2	6.74	125.49	110.00
1	B	273	LYS	CB-CG-CD	6.71	129.05	111.60
1	B	282	ARG	CB-CG-CD	6.61	128.77	111.60
1	B	314	ASP	CB-CG-OD1	6.46	124.12	118.30
1	B	175	GLN	CB-CA-C	-6.19	98.02	110.40
1	B	313	ARG	CD-NE-CZ	5.62	131.47	123.60
1	B	142	ARG	CB-CA-C	5.55	121.50	110.40
1	A	166	VAL	CG1-CB-CG2	5.50	119.70	110.90
1	B	254	GLN	CA-CB-CG	5.45	125.38	113.40
1	A	167	GLU	CG-CD-OE2	-5.29	107.73	118.30
1	B	314	ASP	CB-CG-OD2	-5.22	113.61	118.30
1	B	149	ARG	CB-CG-CD	5.15	124.98	111.60
1	B	23	LYS	CA-CB-CG	5.11	124.64	113.40
1	B	273	LYS	CA-CB-CG	5.07	124.56	113.40

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	B	306	THR	CB

All (5) planarity outliers are listed below:



Mol	Chain	Res	Type	Group
1	A	167	GLU	Sidechain
1	A	76	ARG	Sidechain
1	B	25	GLU	Sidechain
1	B	313	ARG	Peptide
1	B	33	ARG	Sidechain

## 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	2771	0	2742	46	2
1	B	2791	0	2761	81	2
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	12	0	0	0	0
3	B	12	0	0	0	0
All	All	5588	0	5503	127	2

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 11.

All (127) 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:310:MET:O	1:B:314:ASP:HB2	1.71	0.89
1:B:159:PRO:HG3	1:B:194:LEU:HD23	1.60	0.83
1:B:91:ALA:HA	1:B:146:VAL:HG21	1.62	0.80
1:B:97:TYR:CD1	1:B:147:LYS:HG2	2.20	0.77
1:B:282:ARG:O	1:B:286:ASP:N	2.20	0.75
1:B:97:TYR:HD1	1:B:147:LYS:HG2	1.53	0.73
1:A:6:ALA:HB2	1:A:309:GLN:HG2	1.70	0.72
1:B:57:THR:HG22	1:B:59:PRO:HD2	1.72	0.71
1:B:282:ARG:O	1:B:285:ASN:N	2.25	0.70
1:B:25:GLU:OE1	1:B:43:ALA:HB1	1.93	0.69
1:B:283:LEU:HD23	1:B:288:ALA:HB3	1.74	0.69
1:A:217:GLU:HB2	1:A:238:HIS:HB2	1.76	0.68
1:B:134:GLY:HA3	1:B:175:GLN:NE2	2.08	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:141:GLU:HG3	1:B:146:VAL:O	1.93	0.68
1:B:15:HIS:CG	1:B:101:ARG:HH12	2.12	0.67
1:A:42:THR:HG23	1:A:45:GLY:H	1.58	0.66
1:A:60:ASP:O	1:A:64:LYS:HE3	1.96	0.66
1:B:249:TYR:HE1	1:B:282:ARG:HH21	1.41	0.66
1:B:191:SER:HA	1:B:194:LEU:HD13	1.78	0.65
1:B:282:ARG:HB3	1:B:286:ASP:OD1	1.97	0.65
1:A:51:GLY:HA3	1:A:301:LYS:HE2	1.80	0.64
1:B:313:ARG:NH2	1:B:314:ASP:OD1	2.31	0.63
1:A:113:GLU:O	1:A:161:TRP:NE1	2.33	0.61
1:B:331:LYS:HA	1:B:340:LYS:HE2	1.82	0.61
1:A:87:VAL:HG11	1:A:140:GLY:HA3	1.82	0.61
1:B:93:GLU:OE2	1:B:304:LEU:HB3	2.00	0.60
1:B:24:PRO:HG3	1:B:298:LEU:HD13	1.81	0.60
1:B:201:TYR:O	1:B:205:VAL:HG23	2.02	0.59
1:B:154:CYS:HB2	1:B:182:LEU:HD12	1.85	0.58
1:A:282:ARG:HA	1:A:285:ASN:HB2	1.85	0.58
1:A:34:ARG:HB2	1:A:36:ILE:HD12	1.86	0.58
1:B:217:GLU:HB3	1:B:238:HIS:HB2	1.84	0.57
1:B:15:HIS:CG	1:B:101:ARG:NH1	2.72	0.57
1:A:58:LEU:HB3	1:A:59:PRO:HD3	1.87	0.56
1:B:345:ASP:HA	1:B:348:TYR:HB2	1.88	0.55
1:A:163:PRO:O	1:A:167:GLU:HG2	2.06	0.55
1:B:288:ALA:O	1:B:324:ARG:NH1	2.39	0.55
1:B:45:GLY:O	1:B:49:VAL:HG23	2.07	0.55
1:B:53:ASP:OD1	1:B:53:ASP:N	2.40	0.55
1:B:224:VAL:HG13	1:B:236:LEU:HD21	1.88	0.55
1:A:331:LYS:HA	1:A:340:LYS:HE2	1.87	0.55
1:B:24:PRO:HG3	1:B:298:LEU:CD1	2.37	0.55
1:B:76:ARG:NH1	1:B:128:GLU:OE2	2.40	0.55
1:A:166:VAL:HG11	1:A:200:ALA:HA	1.89	0.53
1:A:58:LEU:HG	1:A:218:VAL:HG21	1.91	0.53
1:B:24:PRO:HB3	1:B:47:LEU:HD11	1.90	0.53
1:B:272:TRP:NE1	1:B:276:THR:O	2.39	0.53
1:A:228:VAL:O	1:A:232:LYS:HA	2.09	0.53
1:B:134:GLY:HA2	1:B:176:THR:HG21	1.91	0.52
1:B:238:HIS:HB3	1:B:240:TYR:CE2	2.45	0.52
1:B:15:HIS:O	1:B:293:ASN:HB3	2.10	0.51
1:B:217:GLU:HB3	1:B:238:HIS:CB	2.41	0.51
1:A:290:TYR:OH	1:A:321:GLU:OE2	2.25	0.50
1:B:284:LYS:O	1:B:287:GLN:NE2	2.39	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:156:ARG:NH1	1:A:192:SER:OG	2.41	0.50
1:A:217:GLU:HB2	1:A:238:HIS:CB	2.40	0.49
1:A:327:ILE:O	1:A:331:LYS:HG2	2.11	0.49
1:B:17:HIS:HB2	1:B:296:ASP:OD1	2.11	0.49
1:A:166:VAL:O	1:A:169:CYS:HB2	2.12	0.49
1:B:277:GLU:H	1:B:277:GLU:CD	2.16	0.49
1:A:288:ALA:O	1:A:324:ARG:NH1	2.45	0.49
1:B:221:ALA:N	1:B:244:GLU:OE2	2.41	0.49
1:A:232:LYS:HB2	1:A:232:LYS:NZ	2.28	0.49
1:A:61:PHE:HZ	1:A:296:ASP:OD2	1.96	0.48
1:B:7:PHE:HE1	1:B:319:GLU:HG2	1.78	0.48
1:B:15:HIS:CD2	1:B:101:ARG:NH1	2.81	0.48
1:B:221:ALA:O	1:B:225:LYS:HG3	2.15	0.47
1:B:224:VAL:O	1:B:228:VAL:HG23	2.15	0.47
1:A:281:ILE:HD11	1:A:315:MET:HG2	1.96	0.47
1:B:98:VAL:O	1:B:148:ALA:HA	2.15	0.46
1:A:338:ASP:OD1	1:A:338:ASP:N	2.48	0.46
1:A:115:ILE:HG22	1:A:119:GLN:HB2	1.96	0.46
1:A:243:LEU:HD13	1:A:282:ARG:HD3	1.97	0.46
1:B:134:GLY:HA3	1:B:175:GLN:HE22	1.80	0.46
1:B:88:GLU:HB2	1:B:144:PHE:CE2	2.51	0.46
1:B:109:ASN:HB3	1:B:119:GLN:NE2	2.31	0.45
1:B:294:THR:HB	1:B:297:PRO:HD3	1.97	0.45
1:A:252:LEU:HB3	1:A:257:MET:HB3	1.98	0.45
1:B:91:ALA:CA	1:B:146:VAL:HG21	2.41	0.45
1:A:57:THR:OG1	1:A:59:PRO:HD2	2.17	0.44
1:B:228:VAL:HA	1:B:233:THR:HG22	1.98	0.44
1:B:148:ALA:O	1:B:149:ARG:HG3	2.16	0.44
1:B:240:TYR:CZ	1:B:266:SER:HB2	2.52	0.44
1:A:310:MET:O	1:A:314:ASP:HB2	2.17	0.44
1:B:313:ARG:HE	1:B:314:ASP:CG	2.21	0.43
1:B:283:LEU:HD23	1:B:288:ALA:CB	2.44	0.43
1:A:294:THR:HG23	1:A:304:LEU:HD13	2.00	0.43
1:A:106:LEU:HA	1:A:116:PRO:HG3	2.00	0.43
1:B:239:GLY:HA3	1:B:259:PHE:HD1	1.82	0.43
1:A:39:PRO:HG3	1:A:67:TYR:CD1	2.54	0.42
1:B:95:VAL:O	1:B:146:VAL:HG22	2.18	0.42
1:B:34:ARG:NH1	1:B:71:ALA:O	2.51	0.42
1:B:218:VAL:HG12	1:B:271:ALA:HA	2.02	0.42
1:A:15:HIS:O	1:A:293:ASN:HB3	2.18	0.42
1:A:201:TYR:O	1:A:205:VAL:HG23	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:50:ILE:HD11	1:B:298:LEU:HD11	2.01	0.42
1:B:169:CYS:HA	1:B:177:VAL:HG11	2.01	0.42
1:A:90:LYS:HD3	1:A:90:LYS:HA	1.57	0.42
1:B:297:PRO:HA	1:B:302:SER:HB3	2.01	0.42
1:A:92:LYS:HA	1:A:92:LYS:HD2	1.76	0.41
1:B:47:LEU:O	1:B:301:LYS:HE2	2.20	0.41
1:B:113:GLU:HA	1:B:114:PRO:HA	1.81	0.41
1:B:311:THR:HB	1:B:317:PHE:CG	2.55	0.41
1:A:251:ARG:O	1:A:255:GLU:HG3	2.20	0.41
1:B:323:LYS:O	1:B:327:ILE:HG13	2.20	0.41
1:B:310:MET:HA	1:B:314:ASP:OD2	2.21	0.41
1:B:123:ASP:N	1:B:123:ASP:OD1	2.53	0.41
1:B:156:ARG:NH1	1:B:192:SER:OG	2.53	0.41
1:B:166:VAL:HG22	1:B:180:ILE:HD12	2.01	0.41
1:A:169:CYS:HA	1:A:177:VAL:HG11	2.03	0.41
1:B:75:CYS:O	1:B:79:ILE:HG12	2.20	0.41
1:B:345:ASP:HA	1:B:348:TYR:CD2	2.56	0.41
1:A:56:LEU:HB3	1:A:57:THR:H	1.70	0.41
1:A:194:LEU:O	1:A:198:VAL:HG23	2.20	0.41
1:A:222:GLU:O	1:A:226:GLU:HG2	2.21	0.41
1:B:154:CYS:SG	1:B:165:VAL:HG21	2.61	0.41
1:B:330:ALA:O	1:B:333:SER:HB3	2.21	0.41
1:B:90:LYS:HA	1:B:90:LYS:HD3	1.67	0.40
1:B:99:GLU:HA	1:B:149:ARG:O	2.21	0.40
1:A:57:THR:HG23	1:A:60:ASP:H	1.86	0.40
1:A:282:ARG:HE	1:A:282:ARG:HB2	1.70	0.40
1:B:38:LEU:HD23	1:B:38:LEU:HA	1.86	0.40
1:A:123:ASP:N	1:A:123:ASP:OD1	2.52	0.40
1:B:124:LEU:HD12	1:B:124:LEU:HA	1.92	0.40
1:A:284:LYS:HD3	1:A:317:PHE:CE1	2.56	0.40
1:B:86:PHE:CE2	1:B:100:VAL:HB	2.56	0.40
1:B:107:LEU:HB2	1:B:129:VAL:HG11	2.03	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:8:ASP:OD1	1:B:337:GLU:OE2[2_655]	1.97	0.23
1:A:67:TYR:OH	1:B:354:PRO:O[1_455]	2.17	0.03

## 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	348/372 (94%)	334 (96%)	12 (3%)	2 (1%)	25	47
1	B	351/372 (94%)	337 (96%)	13 (4%)	1 (0%)	41	64
All	All	699/744 (94%)	671 (96%)	25 (4%)	3 (0%)	34	57

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	337	GLU
1	A	295	ASP
1	B	295	ASP

### 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	295/315 (94%)	289 (98%)	6 (2%)	55	78
1	B	299/315 (95%)	292 (98%)	7 (2%)	50	75
All	All	594/630 (94%)	581 (98%)	13 (2%)	52	76

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	LEU
1	A	110	SER
1	A	166	VAL

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	A	203	GLU
1	A	265	SER
1	A	340	LYS
1	B	33	ARG
1	B	142	ARG
1	B	246	GLN
1	B	273	LYS
1	B	282	ARG
1	B	319	GLU
1	B	331	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	135	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 [i](#)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 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, 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	349/372 (93%)	0.22	7 (2%) 65 60	27, 39, 52, 64	0
1	B	352/372 (94%)	0.45	14 (3%) 38 31	31, 43, 61, 76	0
All	All	701/744 (94%)	0.34	21 (2%) 50 43	27, 41, 57, 76	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	314	ASP	5.9
1	B	148	ALA	3.1
1	B	175	GLN	3.1
1	B	311	THR	3.0
1	B	313	ARG	3.0
1	A	114	PRO	2.9
1	A	276	THR	2.9
1	B	25	GLU	2.7
1	B	274	PRO	2.6
1	B	272	TRP	2.6
1	A	337	GLU	2.6
1	B	33	ARG	2.6
1	B	146	VAL	2.4
1	B	312	LYS	2.4
1	B	92	LYS	2.3
1	A	195	PRO	2.3
1	A	42	THR	2.3
1	B	275	ASP	2.3
1	B	53	ASP	2.2
1	A	163	PRO	2.1
1	A	277	GLU	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](#)

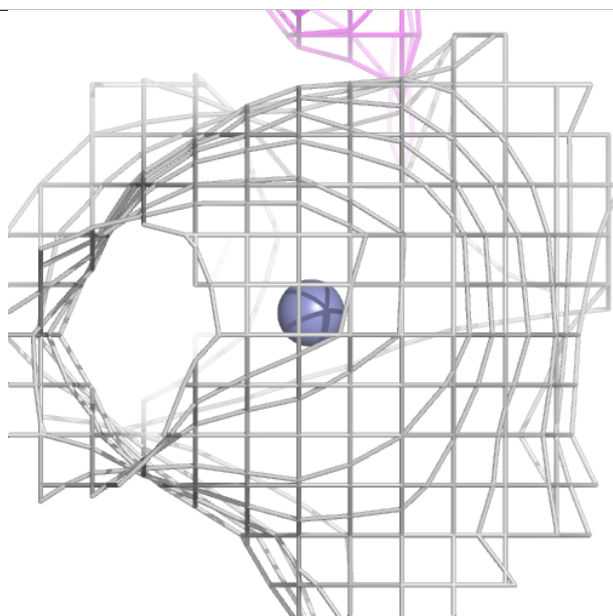
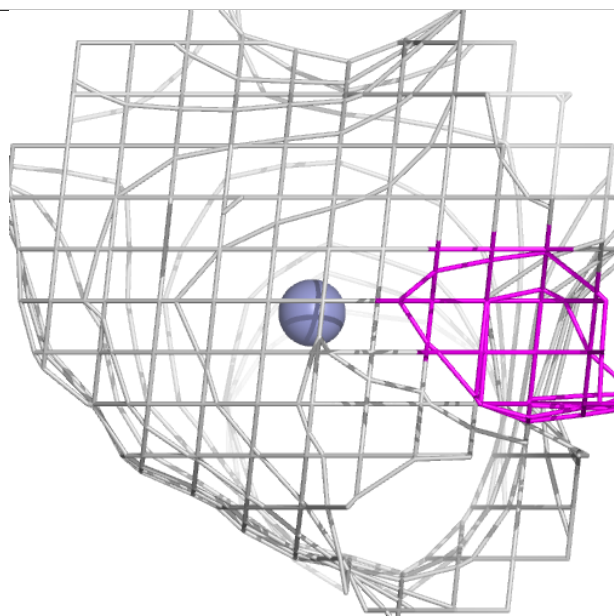
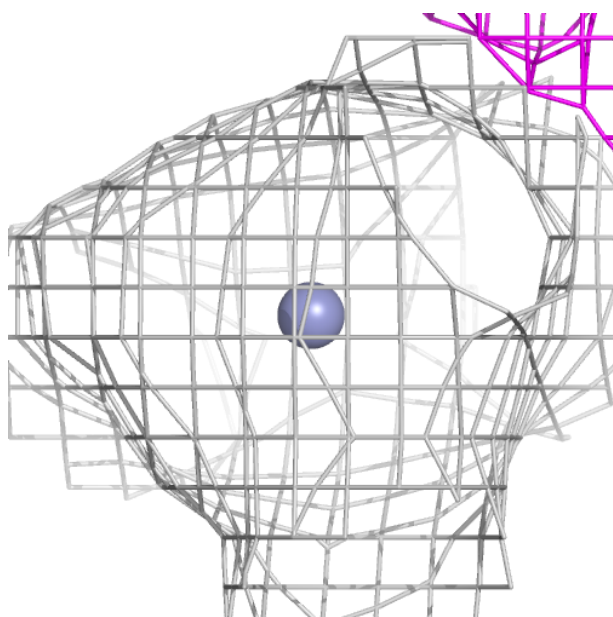
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

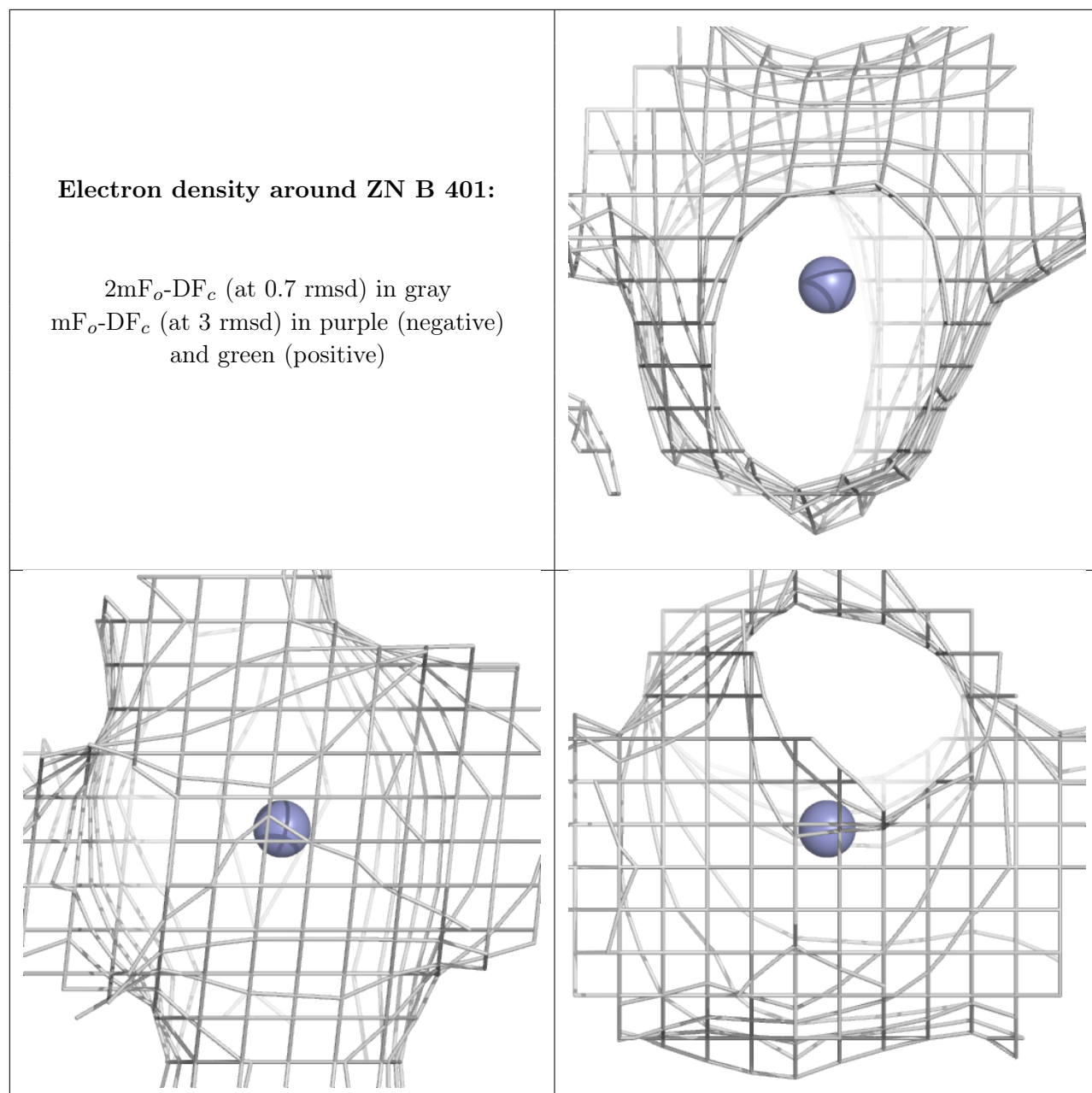
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	ZN	A	401	1/1	0.98	0.10	39,39,39,39	0
2	ZN	B	401	1/1	0.99	0.05	39,39,39,39	0

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 ZN A 401:**

$2mF_o-DF_c$  (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.