



# Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2020 – 08:25 pm BST

PDB ID : 5ZQ6  
Title : SidE-Ubi-ADPr  
Authors : Wang, Y.; Gao, A.; Gao, P.  
Deposited on : 2018-04-17  
Resolution : 3.01 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

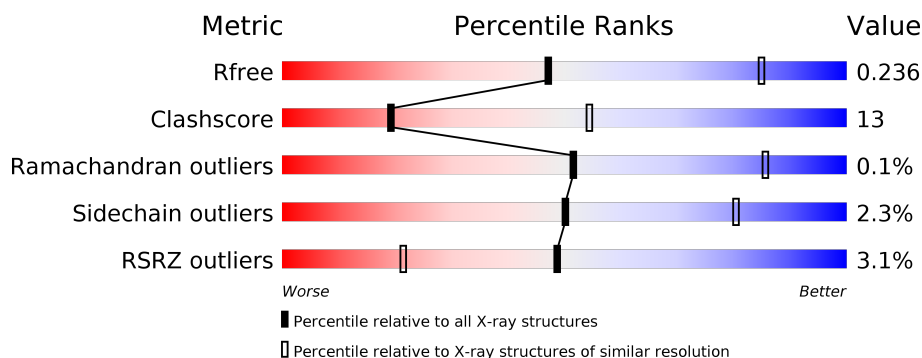
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.01 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	2092 (3.00-3.00)
Clashscore	141614	2416 (3.00-3.00)
Ramachandran outliers	138981	2333 (3.00-3.00)
Sidechain outliers	138945	2336 (3.00-3.00)
RSRZ outliers	127900	1990 (3.00-3.00)

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	845	<div> <div>0%</div> <div> <div></div> <div>72%</div> <div>25%</div> <div>••</div> </div> </div>
1	C	845	<div> <div>3%</div> <div> <div></div> <div>73%</div> <div>23%</div> <div>••</div> </div> </div>
2	B	79	<div> <div>6%</div> <div> <div></div> <div>68%</div> <div>28%</div> <div>•</div> </div> </div>
2	D	79	<div> <div>18%</div> <div> <div></div> <div>53%</div> <div>42%</div> <div>••</div> </div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 14395 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 Side.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	828	Total	C	N	O	S	Se	0	0	0
			6606	4164	1153	1267	6	16			
1	C	819	Total	C	N	O	S	Se	0	0	0
			6545	4132	1142	1249	6	16			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	221	MSE	-	initiating methionine	UNP Q6RCR1
A	1058	LEU	-	expression tag	UNP Q6RCR1
A	1059	GLU	-	expression tag	UNP Q6RCR1
A	1060	HIS	-	expression tag	UNP Q6RCR1
A	1061	HIS	-	expression tag	UNP Q6RCR1
A	1062	HIS	-	expression tag	UNP Q6RCR1
A	1063	HIS	-	expression tag	UNP Q6RCR1
A	1064	HIS	-	expression tag	UNP Q6RCR1
A	1065	HIS	-	expression tag	UNP Q6RCR1
C	221	MSE	-	initiating methionine	UNP Q6RCR1
C	1058	LEU	-	expression tag	UNP Q6RCR1
C	1059	GLU	-	expression tag	UNP Q6RCR1
C	1060	HIS	-	expression tag	UNP Q6RCR1
C	1061	HIS	-	expression tag	UNP Q6RCR1
C	1062	HIS	-	expression tag	UNP Q6RCR1
C	1063	HIS	-	expression tag	UNP Q6RCR1
C	1064	HIS	-	expression tag	UNP Q6RCR1
C	1065	HIS	-	expression tag	UNP Q6RCR1

- Molecule 2 is a protein called Ubiquitin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	76	Total	C	N	O	S	0	0	0
			595	375	102	117	1			

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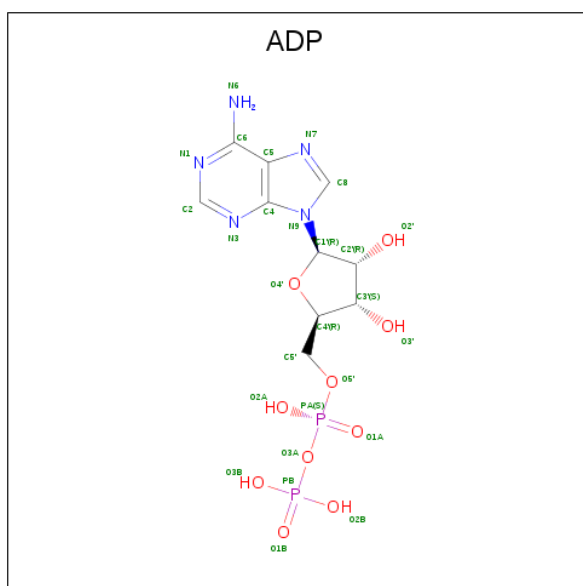
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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	76	Total	C	N	O	S	0	0	0
			595	375	102	117	1			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-2	GLY	-	expression tag	UNP P0CG48
B	-1	GLY	-	expression tag	UNP P0CG48
B	0	SER	-	expression tag	UNP P0CG48
B	42	ALA	ARG	engineered mutation	UNP P0CG48
D	-2	GLY	-	expression tag	UNP P0CG48
D	-1	GLY	-	expression tag	UNP P0CG48
D	0	SER	-	expression tag	UNP P0CG48
D	42	ALA	ARG	engineered mutation	UNP P0CG48

- Molecule 3 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula:  $C_{10}H_{15}N_5O_{10}P_2$ ).



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





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	92.32Å 128.00Å 191.06Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.94 – 3.01 47.77 – 3.01	Depositor EDS
% Data completeness (in resolution range)	99.0 (20.94-3.01) 99.1 (47.77-3.01)	Depositor EDS
$R_{merge}$	0.18	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.12 (at 3.01Å)	Xtriage
Refinement program	PHENIX (1.11.1 _2575: ???)	Depositor
R, $R_{free}$	0.180 , 0.232 0.184 , 0.236	Depositor DCC
$R_{free}$ test set	2286 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	68.4	Xtriage
Anisotropy	0.321	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 59.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	14395	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	83.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.00% 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: 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	0.54	0/6720	0.57	0/9053
1	C	0.46	0/6658	0.54	1/8968 (0.0%)
2	B	0.38	0/601	0.55	0/809
2	D	0.33	0/601	0.52	0/809
All	All	0.49	0/14580	0.56	1/19639 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	365	LEU	CB-CG-CD1	-6.63	99.73	111.00

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	6606	0	6542	167	0
1	C	6545	0	6498	170	0
2	B	595	0	621	16	0
2	D	595	0	621	33	0
3	A	27	0	12	3	0
3	C	27	0	12	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	14395	0	14306	371	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 13.

All (371) 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:C:366:ILE:CG2	1:C:367:PRO:HD3	1.22	1.55
1:C:366:ILE:HG23	1:C:367:PRO:CD	1.47	1.42
1:C:366:ILE:CG2	1:C:367:PRO:CD	2.07	1.24
1:A:401:GLN:O	1:A:405:VAL:HG23	1.52	1.07
1:C:258:GLU:HB2	1:C:337:GLU:HG3	1.11	1.07
1:C:366:ILE:HG22	1:C:367:PRO:HD3	1.39	1.05
1:A:245:TYR:HA	1:A:249:LEU:HB2	1.42	0.97
1:A:407:LEU:O	1:A:411:LYS:HG3	1.63	0.97
1:A:956:GLY:O	1:A:1000:PHE:HE1	1.52	0.93
1:C:365:LEU:HD12	1:C:366:ILE:N	1.83	0.93
1:C:366:ILE:HG22	1:C:367:PRO:CD	1.96	0.92
1:A:973:LEU:O	1:A:977:LEU:HB2	1.70	0.91
1:A:959:THR:HG23	1:A:960:PHE:N	1.90	0.86
1:C:685:THR:O	1:C:689:THR:HG22	1.77	0.84
1:A:956:GLY:O	1:A:1000:PHE:CE1	2.31	0.84
2:B:5:VAL:HG12	2:B:67:LEU:HB2	1.60	0.84
1:A:625:ASN:HB2	1:A:914:MSE:HG2	1.60	0.82
1:C:362:LYS:O	1:C:366:ILE:HG22	1.80	0.81
1:C:258:GLU:CB	1:C:337:GLU:HG3	2.05	0.80
1:C:257:VAL:HG12	1:C:337:GLU:HB2	1.63	0.79
1:A:249:LEU:HD13	1:A:276:LEU:HA	1.64	0.79
1:C:258:GLU:HB2	1:C:337:GLU:CG	2.06	0.79
1:A:410:VAL:HG13	1:A:411:LYS:HG2	1.63	0.78
2:D:5:VAL:HG12	2:D:67:LEU:HB2	1.65	0.77
1:A:404:MSE:SE	1:A:421:TYR:CD1	2.88	0.77
1:A:948:GLN:NE2	1:A:1008:MSE:SE	2.68	0.76
2:D:23:ILE:HD11	2:D:51:GLU:C	2.06	0.75
1:C:745:LYS:NZ	1:C:749:ASP:OD2	2.21	0.74
1:A:611:MSE:HE1	1:A:661:LEU:HD22	1.70	0.73
1:C:339:SER:OG	1:C:340:SER:N	2.21	0.73
1:C:412:GLN:O	1:C:464:LEU:HD23	1.88	0.72
1:A:926:VAL:HG13	1:C:919:LEU:HD12	1.72	0.72
1:A:411:LYS:O	1:A:417:TYR:CD2	2.44	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:594:ARG:NH2	1:C:600:GLU:OE2	2.21	0.71
1:A:959:THR:HG23	1:A:960:PHE:H	1.55	0.70
1:A:638:LEU:HD11	1:A:737:ALA:HA	1.73	0.70
1:C:411:LYS:HD2	1:C:417:TYR:CE1	2.27	0.70
1:C:412:GLN:O	1:C:464:LEU:CD2	2.39	0.69
2:D:45:PHE:HB2	2:D:67:LEU:HD22	1.73	0.68
1:A:271:ARG:NH2	1:A:334:ARG:O	2.26	0.68
1:A:482:ARG:NH2	1:A:522:GLU:OE2	2.26	0.68
1:A:745:LYS:NZ	1:A:749:ASP:OD2	2.25	0.68
1:C:976:LEU:HD11	1:C:1019:VAL:HG12	1.75	0.68
1:A:279:THR:HG22	1:A:283:MSE:HE2	1.77	0.67
1:C:641:CYS:O	1:C:811:ARG:NH1	2.28	0.67
1:A:953:THR:HG23	1:A:954:GLU:OE2	1.95	0.66
1:A:279:THR:O	1:A:283:MSE:HG3	1.96	0.66
1:C:513:ARG:HB2	1:C:516:GLU:HG3	1.78	0.66
1:A:278:HIS:ND1	1:A:406:ASP:OD2	2.28	0.65
1:A:411:LYS:NZ	3:A:1101:ADP:O3B	2.29	0.65
1:A:937:ARG:NH2	1:A:1011:GLN:OE1	2.29	0.65
1:C:365:LEU:C	1:C:365:LEU:HD12	2.13	0.65
1:A:404:MSE:HG3	3:A:1101:ADP:C8	2.32	0.65
1:C:366:ILE:HG23	1:C:367:PRO:HD3	0.65	0.65
1:A:358:VAL:HA	1:A:365:LEU:HD12	1.79	0.65
1:A:598:MSE:HE3	1:A:908:VAL:HG21	1.78	0.64
1:A:948:GLN:HE22	1:A:1008:MSE:SE	2.31	0.64
1:A:776:GLN:HB3	1:A:803:ASN:HB3	1.79	0.64
2:D:22:THR:HG22	2:D:55:THR:HG22	1.78	0.64
1:C:411:LYS:HD2	1:C:417:TYR:CD1	2.32	0.64
1:A:249:LEU:HD13	1:A:276:LEU:CA	2.28	0.63
1:C:941:LEU:HD23	1:C:969:LEU:HD21	1.80	0.63
1:C:321:ARG:HE	1:C:369:VAL:HG13	1.63	0.63
1:A:249:LEU:CD1	1:A:276:LEU:HA	2.28	0.62
1:C:319:GLU:O	1:C:323:ILE:HG12	1.99	0.62
1:A:253:TYR:CZ	1:A:271:ARG:HG2	2.35	0.62
1:A:959:THR:CG2	1:A:960:PHE:N	2.57	0.62
2:D:4:PHE:O	2:D:66:THR:HA	2.00	0.62
1:A:405:VAL:HG12	1:A:405:VAL:O	2.00	0.62
1:C:411:LYS:HD2	1:C:417:TYR:CZ	2.35	0.62
1:C:954:GLU:OE2	1:C:956:GLY:N	2.33	0.61
2:D:73:LEU:HD12	1:C:849:THR:HG21	1.83	0.61
1:A:614:PRO:HG2	1:A:660:THR:HG22	1.83	0.61
1:C:248:TYR:CZ	1:C:353:ALA:HB1	2.35	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:365:LEU:O	1:C:369:VAL:N	2.34	0.61
1:A:410:VAL:HG22	1:A:410:VAL:O	2.00	0.61
1:C:411:LYS:HG3	1:C:417:TYR:CG	2.35	0.61
1:A:959:THR:CG2	1:A:960:PHE:H	2.14	0.60
1:C:242:ARG:O	1:C:246:ASP:HB2	2.01	0.60
1:A:513:ARG:HB2	1:A:516:GLU:HG3	1.83	0.60
1:C:473:VAL:HG23	1:C:481:ILE:HD13	1.83	0.60
1:C:473:VAL:HG12	1:C:511:TYR:HD1	1.66	0.60
1:A:274:HIS:HB3	1:A:331:VAL:HG11	1.83	0.60
1:A:336:ASP:OD1	1:A:338:GLU:N	2.25	0.60
1:C:408:VAL:HG12	1:C:408:VAL:O	2.01	0.60
1:A:826:ILE:HD11	2:B:8:LEU:HG	1.84	0.60
1:A:980:ASN:OD1	1:A:983:ILE:N	2.29	0.60
1:C:543:GLU:O	1:C:547:LEU:HB2	2.01	0.60
1:C:672:SER:HB2	1:C:712:TRP:HA	1.84	0.60
1:A:287:GLU:HB2	1:A:324:MSE:HE1	1.85	0.59
1:C:280:LEU:HD23	1:C:283:MSE:HE3	1.84	0.59
1:C:463:HIS:HB2	1:C:514:VAL:HA	1.83	0.59
1:A:844:PRO:HB3	1:A:860:ILE:HG22	1.84	0.59
1:A:981:VAL:HG11	1:A:1026:GLU:HB2	1.85	0.59
1:A:482:ARG:HG2	1:A:496:LEU:HD23	1.84	0.58
1:A:404:MSE:SE	1:A:421:TYR:HD1	2.35	0.58
1:C:945:VAL:HG22	1:C:1008:MSE:HE3	1.86	0.58
1:C:323:ILE:HD12	1:C:395:ALA:HA	1.86	0.58
2:D:23:ILE:HD11	2:D:52:ASP:N	2.18	0.58
1:C:311:THR:HG23	1:C:313:ALA:H	1.68	0.57
1:A:1030:LEU:HD13	1:C:919:LEU:HD21	1.85	0.57
1:A:503:LYS:HE2	1:A:505:LEU:HD22	1.87	0.57
1:A:941:LEU:HD11	1:A:1008:MSE:HE1	1.87	0.57
2:D:5:VAL:HA	2:D:67:LEU:O	2.04	0.57
1:C:935:LEU:HD22	1:C:973:LEU:HD12	1.86	0.56
1:A:667:ASN:HA	1:A:681:LYS:HE2	1.86	0.56
1:A:287:GLU:HB2	1:A:324:MSE:CE	2.35	0.56
1:A:938:VAL:HG22	1:A:1015:ILE:HD12	1.87	0.56
1:C:257:VAL:CG1	1:C:337:GLU:HB2	2.35	0.56
1:A:480:PRO:HB3	1:A:512:MSE:HE3	1.86	0.56
1:A:672:SER:HB2	1:A:712:TRP:HA	1.87	0.56
1:A:619:ASP:HA	1:A:731:LYS:HE2	1.86	0.56
2:D:30:ILE:HG21	2:D:69:LEU:HD22	1.88	0.56
1:C:279:THR:HG22	1:C:283:MSE:HE2	1.87	0.56
1:A:1045:ILE:HG13	1:A:1046:ALA:N	2.21	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:366:ILE:HG22	1:C:367:PRO:HD2	1.84	0.55
1:A:400:ASN:OD1	3:A:1101:ADP:O2'	2.25	0.55
1:C:248:TYR:CE2	1:C:353:ALA:HB2	2.41	0.55
1:A:744:ASN:HB3	1:A:748:MSE:HE2	1.89	0.55
2:D:8:LEU:HD11	1:C:822:VAL:HG22	1.88	0.55
1:C:245:TYR:HA	1:C:249:LEU:HB2	1.89	0.55
1:C:424:GLN:OE1	3:C:1101:ADP:N6	2.40	0.54
1:C:878:LYS:NZ	1:C:884:ASP:OD1	2.33	0.54
1:A:543:GLU:O	1:A:547:LEU:HB2	2.06	0.54
1:C:1038:ASN:OD1	1:C:1042:LYS:NZ	2.35	0.54
1:A:283:MSE:HE2	1:A:328:ALA:HA	1.89	0.54
1:A:561:ASP:HB3	1:A:564:PHE:HB3	1.88	0.54
1:C:716:ARG:O	1:C:720:ILE:HG13	2.07	0.54
1:A:411:LYS:HB3	1:A:417:TYR:CE2	2.43	0.54
1:A:822:VAL:HB	1:A:823:PRO:HD3	1.89	0.54
1:C:627:GLN:HG2	1:C:913:LYS:HB2	1.89	0.54
1:A:336:ASP:C	1:A:336:ASP:OD1	2.47	0.53
1:A:356:LYS:HE2	1:A:360:GLU:OE1	2.09	0.53
1:C:411:LYS:HG3	1:C:417:TYR:CB	2.38	0.53
1:A:611:MSE:CE	1:A:738:LEU:HD13	2.39	0.53
1:C:248:TYR:CE2	1:C:353:ALA:CB	2.92	0.53
1:C:956:GLY:O	1:C:960:PHE:CB	2.57	0.53
1:A:611:MSE:HE2	1:A:661:LEU:HD13	1.90	0.53
1:C:365:LEU:HD12	1:C:366:ILE:H	1.71	0.53
1:C:406:ASP:O	1:C:409:ARG:NE	2.41	0.53
1:C:411:LYS:HD2	1:C:417:TYR:CG	2.43	0.53
2:B:39:ASP:OD1	2:B:39:ASP:N	2.39	0.52
1:C:544:TYR:CE1	1:C:545:GLN:HG3	2.45	0.52
1:C:351:ARG:NH2	1:C:352:ASP:OD1	2.42	0.52
1:C:245:TYR:OH	1:C:560:ASN:OD1	2.21	0.52
1:A:470:GLY:HA3	1:A:498:PHE:CE1	2.45	0.52
1:A:1053:ARG:NH1	1:C:1055:ILE:O	2.37	0.52
2:B:22:THR:HA	2:B:55:THR:HA	1.91	0.52
1:C:532:LYS:HG3	1:C:789:PHE:HB3	1.92	0.52
2:B:22:THR:HG22	2:B:25:ASN:CG	2.30	0.52
1:C:366:ILE:HD11	1:C:372:ASP:HA	1.91	0.52
1:A:641:CYS:CB	1:A:643:MSE:HE3	2.40	0.52
1:A:643:MSE:HE2	1:A:646:VAL:HG21	1.91	0.52
1:C:279:THR:O	1:C:283:MSE:HG3	2.10	0.52
2:D:14:THR:O	2:D:33:LYS:NZ	2.36	0.52
1:C:329:PHE:HA	1:C:332:THR:HG23	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:31:GLN:O	2:D:35:GLY:N	2.41	0.51
2:D:40:GLN:O	2:D:71:LEU:HA	2.10	0.51
1:C:400:ASN:ND2	3:C:1101:ADP:O2'	2.22	0.51
1:A:691:ARG:NH1	1:A:693:GLY:O	2.39	0.51
2:B:37:PRO:HB2	2:B:39:ASP:OD1	2.09	0.51
2:D:40:GLN:HG2	2:D:73:LEU:HD21	1.93	0.51
1:A:656:ARG:HG3	1:A:696:VAL:HG12	1.92	0.51
1:A:370:PHE:O	1:A:371:LYS:HB2	2.11	0.51
1:A:604:ALA:O	1:A:608:GLN:HG3	2.10	0.51
1:C:322:LYS:HD2	1:C:379:TYR:CE2	2.46	0.51
2:D:8:LEU:HG	1:C:826:ILE:HD11	1.93	0.51
1:C:644:ALA:O	1:C:748:MSE:HE1	2.11	0.51
2:B:5:VAL:HG23	2:B:13:ILE:HB	1.92	0.50
1:C:593:ARG:NH2	1:C:861:TYR:O	2.26	0.50
2:D:26:VAL:O	2:D:30:ILE:HG13	2.10	0.50
1:C:515:ASP:O	1:C:519:LYS:HG3	2.10	0.50
1:C:413:PRO:HB3	1:C:461:GLU:OE1	2.11	0.50
1:C:537:GLY:HA2	1:C:546:TYR:CZ	2.47	0.50
1:A:594:ARG:NH2	1:A:600:GLU:OE2	2.31	0.50
1:C:365:LEU:C	1:C:368:ASP:H	2.15	0.50
1:A:643:MSE:CE	1:A:646:VAL:HG21	2.41	0.50
1:A:841:LEU:HD13	1:A:894:LYS:HB2	1.94	0.50
1:C:954:GLU:OE2	1:C:957:TYR:N	2.31	0.50
2:D:2:GLN:HA	2:D:16:GLU:HA	1.93	0.50
1:A:405:VAL:O	1:A:441:GLN:OE1	2.30	0.49
1:A:716:ARG:O	1:A:720:ILE:HG13	2.12	0.49
1:C:411:LYS:HD2	1:C:417:TYR:CE2	2.47	0.49
2:D:74:ARG:NH2	1:C:690:ILE:O	2.42	0.49
1:A:611:MSE:HE3	1:A:738:LEU:HD13	1.94	0.49
1:A:780:ILE:HG23	1:A:800:ILE:HG12	1.94	0.49
2:B:4:PHE:O	2:B:66:THR:HA	2.12	0.49
1:C:348:GLU:HG2	1:C:351:ARG:HH12	1.77	0.49
1:A:655:ILE:HD13	1:A:742:LEU:HD22	1.93	0.49
1:A:948:GLN:HE21	1:A:1008:MSE:SE	2.46	0.49
1:C:278:HIS:O	1:C:282:THR:HG23	2.13	0.49
1:A:909:GLY:O	1:A:913:LYS:HG3	2.13	0.49
1:C:365:LEU:HA	1:C:368:ASP:HB2	1.94	0.49
1:A:941:LEU:HD23	1:A:969:LEU:HD21	1.94	0.49
1:C:323:ILE:HD13	1:C:379:TYR:HE1	1.78	0.49
1:C:638:LEU:HD21	1:C:736:THR:HG22	1.94	0.49
1:A:956:GLY:C	1:A:1000:PHE:HE1	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:411:LYS:HD3	1:A:417:TYR:CE1	2.49	0.48
1:C:293:ALA:HB2	1:C:429:ILE:HG22	1.95	0.48
2:D:73:LEU:HB2	1:C:849:THR:CG2	2.43	0.48
1:A:739:MSE:HE2	1:A:742:LEU:HD12	1.95	0.48
1:C:1006:ARG:HA	1:C:1009:LYS:HE2	1.95	0.48
1:A:341:LYS:O	1:A:342:ASN:HB2	2.13	0.48
1:C:572:ALA:O	1:C:576:THR:HG23	2.14	0.48
1:C:956:GLY:O	1:C:960:PHE:HB3	2.14	0.48
1:C:412:GLN:O	1:C:464:LEU:HD21	2.12	0.48
1:C:770:LYS:NZ	1:C:884:ASP:O	2.36	0.48
1:A:295:LYS:O	1:A:299:ARG:HG3	2.14	0.47
1:A:409:ARG:HG2	1:A:410:VAL:N	2.29	0.47
2:B:42:ALA:HB3	2:B:70:VAL:HG23	1.95	0.47
1:C:351:ARG:HG3	1:C:380:ALA:HB1	1.96	0.47
1:C:638:LEU:HD13	1:C:737:ALA:HA	1.96	0.47
1:A:329:PHE:O	1:A:350:SER:HB3	2.14	0.47
1:A:338:GLU:OE1	1:A:339:SER:N	2.45	0.47
1:A:911:LEU:HD12	1:A:914:MSE:HE2	1.95	0.47
1:C:411:LYS:HA	1:C:411:LYS:HD3	1.65	0.47
1:C:440:THR:HG23	1:C:576:THR:HG21	1.95	0.47
1:C:691:ARG:NH1	1:C:693:GLY:O	2.36	0.47
2:D:15:LEU:HD11	2:D:30:ILE:HG12	1.96	0.47
1:C:247:ASN:HB2	1:C:248:TYR:CD1	2.50	0.47
1:C:619:ASP:HA	1:C:731:LYS:HE3	1.97	0.47
1:C:505:LEU:HD12	1:C:505:LEU:HA	1.78	0.47
1:A:596:PRO:HG2	1:A:649:LEU:HD21	1.96	0.47
1:A:822:VAL:HA	1:A:826:ILE:HD13	1.97	0.47
1:A:1024:VAL:HG21	1:C:1045:ILE:HG21	1.97	0.47
1:C:248:TYR:CZ	1:C:353:ALA:CB	2.98	0.47
1:C:934:ASP:CG	1:C:937:ARG:HG3	2.35	0.47
1:A:455:PHE:HB3	1:A:534:LEU:HB2	1.96	0.46
1:A:652:ASP:HB3	1:A:696:VAL:HG11	1.96	0.46
1:C:662:MSE:HG3	1:C:686:TRP:CG	2.49	0.46
1:C:770:LYS:NZ	1:C:830:ASN:O	2.47	0.46
1:C:624:LEU:HD23	1:C:914:MSE:SE	2.65	0.46
2:B:45:PHE:HB3	2:B:50:LEU:HD21	1.97	0.46
1:C:1009:LYS:HB3	1:C:1009:LYS:HE2	1.67	0.46
1:C:954:GLU:HB3	1:C:957:TYR:HB3	1.97	0.46
1:C:970:GLU:O	1:C:974:ASN:ND2	2.49	0.46
1:A:397:VAL:O	1:A:401:GLN:HB2	2.15	0.46
2:D:27:LYS:HB3	2:D:38:PRO:HB3	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:22:THR:HG22	2:B:25:ASN:ND2	2.30	0.46
1:C:747:LEU:HB3	1:C:845:LYS:CB	2.46	0.46
1:A:970:GLU:O	1:A:974:ASN:ND2	2.48	0.46
1:C:243:TYR:CD2	1:C:357:TYR:HD1	2.33	0.46
1:C:473:VAL:HG12	1:C:511:TYR:CD1	2.48	0.46
1:C:780:ILE:HG23	1:C:800:ILE:HG12	1.96	0.46
1:A:935:LEU:HD12	1:C:912:LEU:HD21	1.97	0.45
1:A:308:ASP:OD2	1:A:310:ARG:HD3	2.16	0.45
1:A:597:ASN:O	1:A:601:ILE:HG13	2.17	0.45
2:D:51:GLU:HB3	2:D:54:ARG:HH21	1.81	0.45
1:A:253:TYR:CE1	1:A:271:ARG:HG2	2.51	0.45
1:A:278:HIS:O	1:A:282:THR:HG23	2.16	0.45
1:C:941:LEU:O	1:C:945:VAL:HG23	2.17	0.45
2:D:22:THR:O	2:D:26:VAL:HG23	2.17	0.45
1:A:941:LEU:O	1:A:945:VAL:HG23	2.16	0.45
1:C:411:LYS:HD2	1:C:417:TYR:CD2	2.52	0.45
1:C:411:LYS:CG	1:C:417:TYR:CG	3.00	0.45
1:C:841:LEU:HD13	1:C:894:LYS:HB2	1.99	0.45
1:A:252:PRO:HA	1:A:270:TYR:HB3	1.98	0.44
1:A:984:LEU:HA	1:A:987:VAL:HG12	1.99	0.44
2:B:15:LEU:HD11	2:B:30:ILE:HG12	1.98	0.44
1:C:243:TYR:CE1	1:C:247:ASN:OD1	2.70	0.44
1:C:247:ASN:C	1:C:248:TYR:CD1	2.90	0.44
2:B:21:ASP:HB2	2:B:56:LEU:HD12	1.99	0.44
1:A:1008:MSE:HE2	1:A:1012:MSE:CE	2.47	0.44
1:A:969:LEU:HD11	1:A:1012:MSE:HE1	1.99	0.44
1:C:711:ALA:HA	1:C:714:HIS:CD2	2.53	0.44
1:A:338:GLU:OE1	1:A:338:GLU:HA	2.17	0.44
1:A:772:LYS:HB2	1:A:772:LYS:HE3	1.77	0.44
1:A:986:LYS:HB2	1:A:986:LYS:HE3	1.85	0.44
2:D:42:ALA:HB3	2:D:70:VAL:HG23	2.00	0.44
1:A:846:GLN:HG3	1:A:850:HIS:NE2	2.32	0.44
1:A:980:ASN:OD1	1:A:982:THR:N	2.51	0.44
1:C:361:ASN:O	1:C:362:LYS:C	2.55	0.44
1:C:710:ASP:HB3	1:C:712:TRP:CE2	2.53	0.44
1:C:721:ALA:HB3	1:C:812:THR:HB	2.00	0.44
1:A:936:GLU:OE2	1:C:904:SER:HB2	2.18	0.44
1:A:1011:GLN:O	1:A:1015:ILE:HG12	2.18	0.44
1:C:397:VAL:O	1:C:401:GLN:HB2	2.18	0.44
1:A:366:ILE:HA	1:A:367:PRO:HA	1.70	0.43
1:A:650:LEU:O	1:A:745:LYS:HE2	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1025:LEU:HD23	1:C:1025:LEU:HA	1.89	0.43
1:C:604:ALA:O	1:C:608:GLN:HG3	2.18	0.43
1:A:336:ASP:OD1	1:A:338:GLU:CB	2.66	0.43
1:C:403:HIS:CD2	3:C:1101:ADP:H4'	2.53	0.43
1:C:710:ASP:HB3	1:C:712:TRP:NE1	2.33	0.43
1:C:846:GLN:HG3	1:C:850:HIS:NE2	2.33	0.43
1:A:582:LYS:HB3	1:A:582:LYS:HE2	1.79	0.43
1:C:972:SER:O	1:C:976:LEU:HD13	2.17	0.43
1:A:691:ARG:CZ	1:A:699:ILE:HG12	2.48	0.43
1:C:956:GLY:HA2	1:C:959:THR:OG1	2.19	0.43
1:C:961:LEU:HA	1:C:961:LEU:HD23	1.85	0.43
1:A:238:ARG:HD3	1:A:242:ARG:NH2	2.33	0.43
1:A:641:CYS:HB2	1:A:643:MSE:HE3	2.00	0.43
1:A:988:LEU:HD11	1:A:1020:GLU:HA	2.01	0.43
1:C:299:ARG:HD3	1:C:782:ALA:O	2.19	0.43
1:C:917:PRO:HB2	1:C:919:LEU:HG	2.00	0.43
1:C:955:SER:O	1:C:959:THR:OG1	2.30	0.43
1:A:997:TRP:CE3	1:A:1009:LYS:HB2	2.54	0.43
1:A:611:MSE:HE3	1:A:738:LEU:HB3	2.01	0.43
1:A:948:GLN:OE1	1:A:1004:GLU:HB2	2.19	0.43
1:A:658:ILE:HG12	1:A:739:MSE:HE1	1.99	0.43
1:A:739:MSE:O	1:A:743:THR:OG1	2.25	0.43
2:B:27:LYS:HB3	2:B:38:PRO:HB3	2.01	0.43
1:A:358:VAL:O	1:A:362:LYS:N	2.52	0.42
1:A:505:LEU:HD12	1:A:505:LEU:HA	1.72	0.42
1:C:846:GLN:HG3	1:C:850:HIS:CE1	2.54	0.42
1:C:824:ARG:NH1	1:C:876:ASP:OD2	2.46	0.42
1:A:343:TYR:C	1:A:343:TYR:CD1	2.93	0.42
1:A:820:ILE:O	1:A:823:PRO:HD2	2.19	0.42
1:C:415:GLU:HB2	1:C:514:VAL:HG21	2.02	0.42
1:A:596:PRO:HB3	1:A:649:LEU:HD11	2.01	0.42
1:A:770:LYS:NZ	1:A:830:ASN:O	2.48	0.42
1:C:396:HIS:O	1:C:400:ASN:HB2	2.20	0.42
1:A:532:LYS:HG3	1:A:789:PHE:HB3	2.00	0.42
1:A:924:ARG:O	1:A:927:GLU:HG2	2.20	0.42
1:A:404:MSE:SE	1:A:421:TYR:CE1	3.22	0.42
2:D:43:LEU:HA	2:D:43:LEU:HD23	1.90	0.42
1:C:411:LYS:HB3	1:C:417:TYR:CD2	2.55	0.42
1:C:412:GLN:HA	1:C:413:PRO:HA	1.74	0.42
1:C:333:GLY:HA3	1:C:350:SER:N	2.34	0.42
1:C:664:HIS:ND1	1:C:665:GLU:OE2	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:909:GLY:O	1:C:913:LYS:HG2	2.19	0.42
2:D:8:LEU:HD22	1:C:819:ASN:ND2	2.34	0.42
1:A:475:GLY:HA2	1:A:509:GLN:HG3	2.02	0.41
1:A:742:LEU:HD23	1:A:742:LEU:HA	1.95	0.41
1:C:584:ALA:HA	1:C:793:SER:HB3	2.02	0.41
1:A:483:GLU:OE2	1:A:497:LYS:HD2	2.20	0.41
1:C:830:ASN:O	1:C:885:ARG:HA	2.20	0.41
1:A:470:GLY:O	1:A:471:ARG:HD2	2.21	0.41
1:C:365:LEU:O	1:C:368:ASP:N	2.46	0.41
1:C:401:GLN:O	1:C:405:VAL:HG23	2.20	0.41
1:C:776:GLN:HB3	1:C:803:ASN:HB3	2.02	0.41
1:A:978:ASP:O	1:A:979:ASN:HB2	2.19	0.41
1:C:607:VAL:O	1:C:611:MSE:HG2	2.20	0.41
2:D:23:ILE:CD1	2:D:51:GLU:C	2.85	0.41
1:C:470:GLY:HA3	1:C:498:PHE:CE1	2.56	0.41
1:A:546:TYR:CE2	1:A:550:LEU:HD11	2.56	0.41
1:A:952:SER:HB3	1:A:958:LYS:HD2	2.03	0.41
1:A:978:ASP:O	1:A:979:ASN:CB	2.69	0.41
1:A:619:ASP:OD1	1:A:619:ASP:N	2.53	0.41
1:A:932:GLU:OE2	1:C:913:LYS:HD3	2.21	0.41
2:B:37:PRO:HA	2:B:38:PRO:HD3	1.92	0.41
1:C:385:ASP:OD1	1:C:388:HIS:HA	2.20	0.41
1:C:665:GLU:OE1	1:C:668:THR:OG1	2.27	0.41
1:C:689:THR:HG23	1:C:690:ILE:HG12	2.03	0.41
1:A:683:GLY:O	1:A:687:GLU:HG3	2.21	0.41
1:A:797:PHE:CZ	1:A:801:LYS:HD2	2.56	0.41
1:C:988:LEU:HD11	1:C:1020:GLU:HB2	2.01	0.41
2:D:23:ILE:CD1	2:D:51:GLU:O	2.69	0.41
2:D:71:LEU:O	1:C:854:SER:HA	2.21	0.41
1:A:293:ALA:HB2	1:A:429:ILE:HG22	2.03	0.41
1:A:404:MSE:HE2	1:A:404:MSE:HB3	1.75	0.41
1:A:405:VAL:CG1	1:A:405:VAL:O	2.68	0.41
1:A:583:ARG:NH2	1:A:786:GLU:OE1	2.47	0.41
1:A:370:PHE:O	1:A:372:ASP:N	2.49	0.40
2:B:33:LYS:HB3	2:B:33:LYS:HE3	1.88	0.40
1:C:956:GLY:O	1:C:960:PHE:HB2	2.21	0.40
1:C:986:LYS:HA	1:C:986:LYS:HD2	1.82	0.40
2:D:23:ILE:HD11	2:D:51:GLU:O	2.22	0.40
1:C:550:LEU:HD23	1:C:550:LEU:HA	1.86	0.40
1:C:624:LEU:HD12	1:C:624:LEU:HA	1.93	0.40
1:A:339:SER:HB2	1:A:340:SER:H	1.69	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:418:LEU:O	1:A:422:PHE:HB2	2.22	0.40
1:A:643:MSE:HE2	1:A:646:VAL:CG2	2.51	0.40
1:C:717:VAL:HG13	1:C:736:THR:HG23	2.03	0.40
1:C:748:MSE:HG2	1:C:845:LYS:HE3	2.02	0.40
2:D:17:VAL:HB	2:D:21:ASP:OD2	2.22	0.40
1:A:413:PRO:HA	1:A:414:PRO:HD3	1.90	0.40
1:A:961:LEU:HD23	1:A:961:LEU:HA	1.80	0.40
1:A:821:GLU:OE2	1:A:824:ARG:NH2	2.33	0.40
2:D:37:PRO:HA	2:D:38:PRO:HD3	1.94	0.40
2:D:23:ILE:CG1	2:D:52:ASP:HA	2.52	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	824/845 (98%)	798 (97%)	26 (3%)	0	100	100
1	C	813/845 (96%)	790 (97%)	22 (3%)	1 (0%)	51	85
2	B	74/79 (94%)	73 (99%)	1 (1%)	0	100	100
2	D	74/79 (94%)	72 (97%)	2 (3%)	0	100	100
All	All	1785/1848 (97%)	1733 (97%)	51 (3%)	1 (0%)	51	85

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	408	VAL

### 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	719/719 (100%)	698 (97%)	21 (3%)	42	76
1	C	713/719 (99%)	699 (98%)	14 (2%)	55	83
2	B	67/68 (98%)	67 (100%)	0	100	100
2	D	67/68 (98%)	66 (98%)	1 (2%)	65	87
All	All	1566/1574 (100%)	1530 (98%)	36 (2%)	50	80

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	339	SER
1	A	340	SER
1	A	363	SER
1	A	364	THR
1	A	372	ASP
1	A	386	LYS
1	A	392	ASP
1	A	404	MSE
1	A	409	ARG
1	A	411	LYS
1	A	471	ARG
1	A	544	TYR
1	A	615	ASP
1	A	619	ASP
1	A	670	PHE
1	A	950	LYS
1	A	954	GLU
1	A	978	ASP
1	A	986	LYS
1	A	1000	PHE
1	A	1002	SER
2	D	54	ARG
1	C	340	SER
1	C	364	THR

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Mol	Chain	Res	Type
1	C	392	ASP
1	C	412	GLN
1	C	480	PRO
1	C	544	TYR
1	C	615	ASP
1	C	619	ASP
1	C	627	GLN
1	C	643	MSE
1	C	670	PHE
1	C	707	MSE
1	C	752	ARG
1	C	928	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	347	HIS
1	A	441	GLN
1	A	446	HIS
1	A	974	ASN

### 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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

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	C	1101	-	24,29,29	1.00	1 (4%)	29,45,45	1.38	3 (10%)
3	ADP	A	1101	-	24,29,29	1.08	2 (8%)	29,45,45	1.28	3 (10%)

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	C	1101	-	-	2/12/32/32	0/3/3/3
3	ADP	A	1101	-	-	2/12/32/32	0/3/3/3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1101	ADP	C5-C4	2.87	1.48	1.40
3	C	1101	ADP	C5-C4	2.68	1.48	1.40
3	A	1101	ADP	C2-N3	2.24	1.35	1.32

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1101	ADP	C3'-C2'-C1'	3.70	106.55	100.98
3	C	1101	ADP	N3-C2-N1	-3.58	123.08	128.68
3	C	1101	ADP	C3'-C2'-C1'	3.49	106.24	100.98
3	A	1101	ADP	C4-C5-N7	-2.77	106.51	109.40
3	C	1101	ADP	PA-O3A-PB	-2.43	124.49	132.83
3	A	1101	ADP	N3-C2-N1	-2.40	124.93	128.68

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	1101	ADP	C3'-C4'-C5'-O5'
3	C	1101	ADP	O4'-C4'-C5'-O5'

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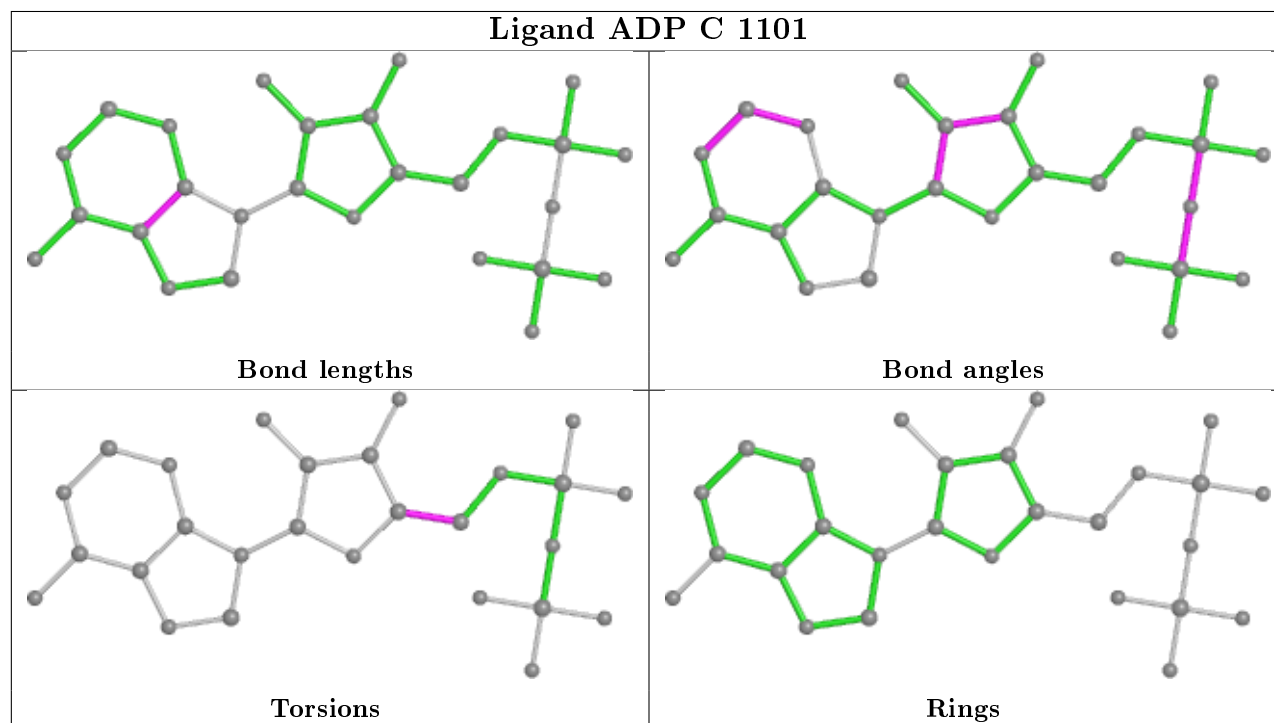
Mol	Chain	Res	Type	Atoms
3	A	1101	ADP	PB-O3A-PA-O1A
3	A	1101	ADP	PB-O3A-PA-O2A

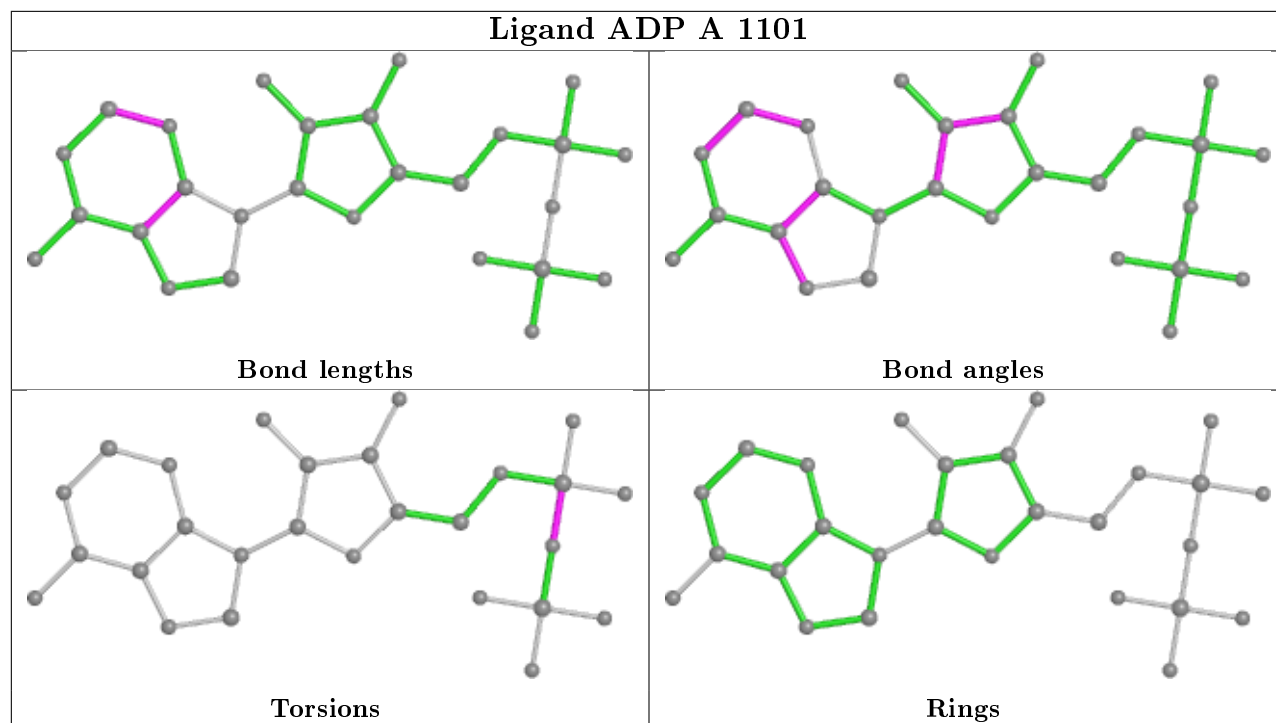
There are no ring outliers.

2 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	1101	ADP	3	0
3	A	1101	ADP	3	0

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	812/845 (96%)	-0.09	9 (1%) 80 56	33, 65, 122, 182	0
1	C	803/845 (95%)	0.11	27 (3%) 45 19	37, 85, 141, 225	0
2	B	76/79 (96%)	0.61	5 (6%) 18 5	59, 83, 114, 118	0
2	D	76/79 (96%)	1.09	14 (18%) 1 0	76, 131, 168, 183	0
All	All	1767/1848 (95%)	0.08	55 (3%) 49 21	33, 77, 139, 225	0

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	957	TYR	5.0
2	D	20	SER	4.9
2	D	32	ASP	4.8
2	B	20	SER	4.3
1	C	1001	ASN	3.9
1	A	223	VAL	3.7
1	C	1046	ALA	3.3
1	C	543	GLU	3.3
1	C	224	PRO	3.1
1	C	223	VAL	3.1
1	C	1058	LEU	3.1
2	B	2	GLN	3.1
1	A	1000	PHE	3.0
2	B	15	LEU	3.0
1	C	544	TYR	3.0
1	C	1045	ILE	2.9
1	A	222	SER	2.9
2	D	1	MET	2.9
1	C	266	GLY	2.9
2	D	15	LEU	2.8
1	C	988	LEU	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	953	THR	2.8
2	D	3	ILE	2.7
2	D	51	GLU	2.7
2	D	2	GLN	2.6
1	A	1056	PRO	2.5
1	C	983	ILE	2.5
2	B	17	VAL	2.5
1	C	344	GLU	2.5
1	A	1046	ALA	2.4
2	D	21	ASP	2.4
2	D	16	GLU	2.4
1	A	956	GLY	2.4
1	C	251	LYS	2.3
1	C	384	GLU	2.3
1	C	1021	LYS	2.3
1	C	343	TYR	2.3
1	C	979	ASN	2.3
2	B	32	ASP	2.3
1	C	1057	SER	2.3
1	C	985	GLY	2.2
1	C	984	LEU	2.2
2	D	53	GLY	2.2
1	C	1055	ILE	2.2
2	D	63	LYS	2.1
1	C	342	ASN	2.1
2	D	18	GLU	2.1
1	C	1028	GLN	2.1
1	C	1052	LEU	2.1
1	C	371	LYS	2.1
2	D	76	GLY	2.1
1	A	992	PRO	2.0
1	C	388	HIS	2.0
2	D	13	ILE	2.0
1	C	937	ARG	2.0

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

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

## 6.3 Carbohydrates [i](#)

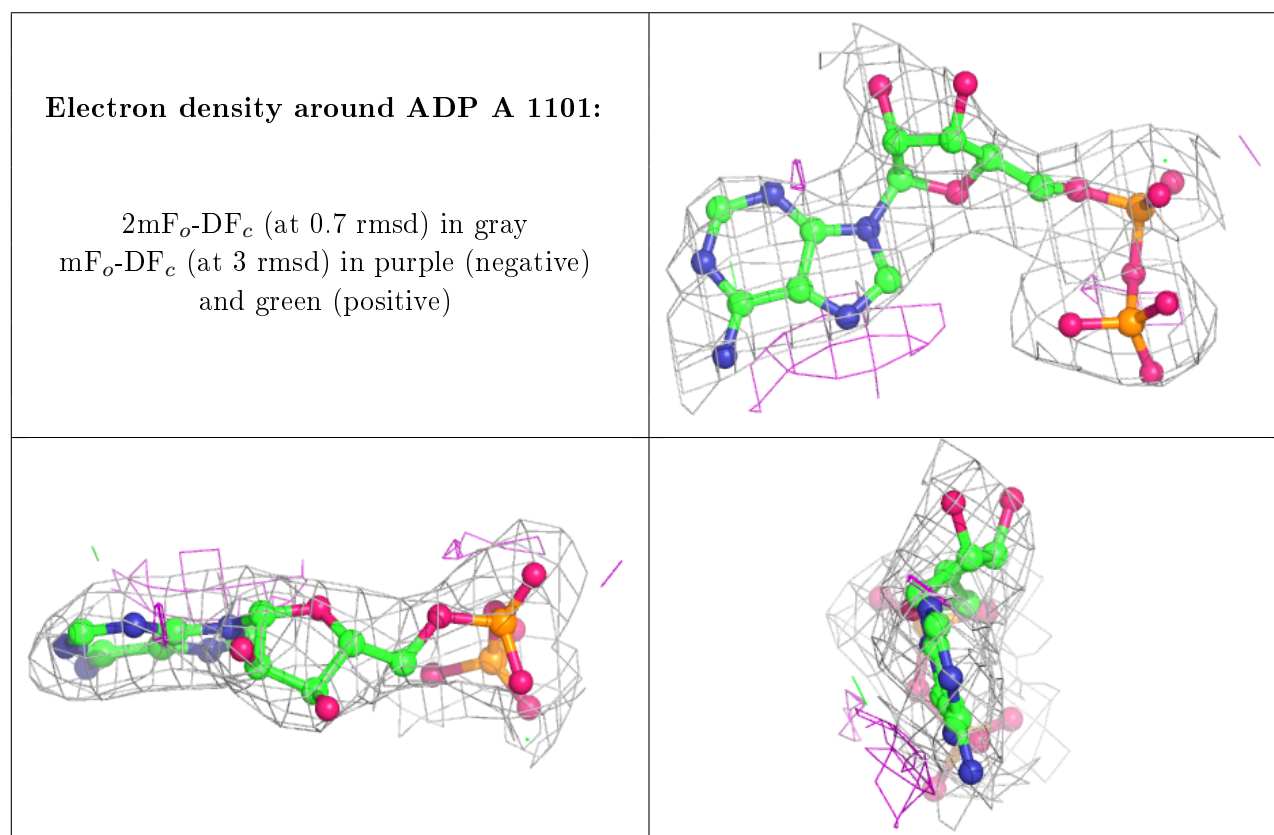
There are no carbohydrates in this entry.

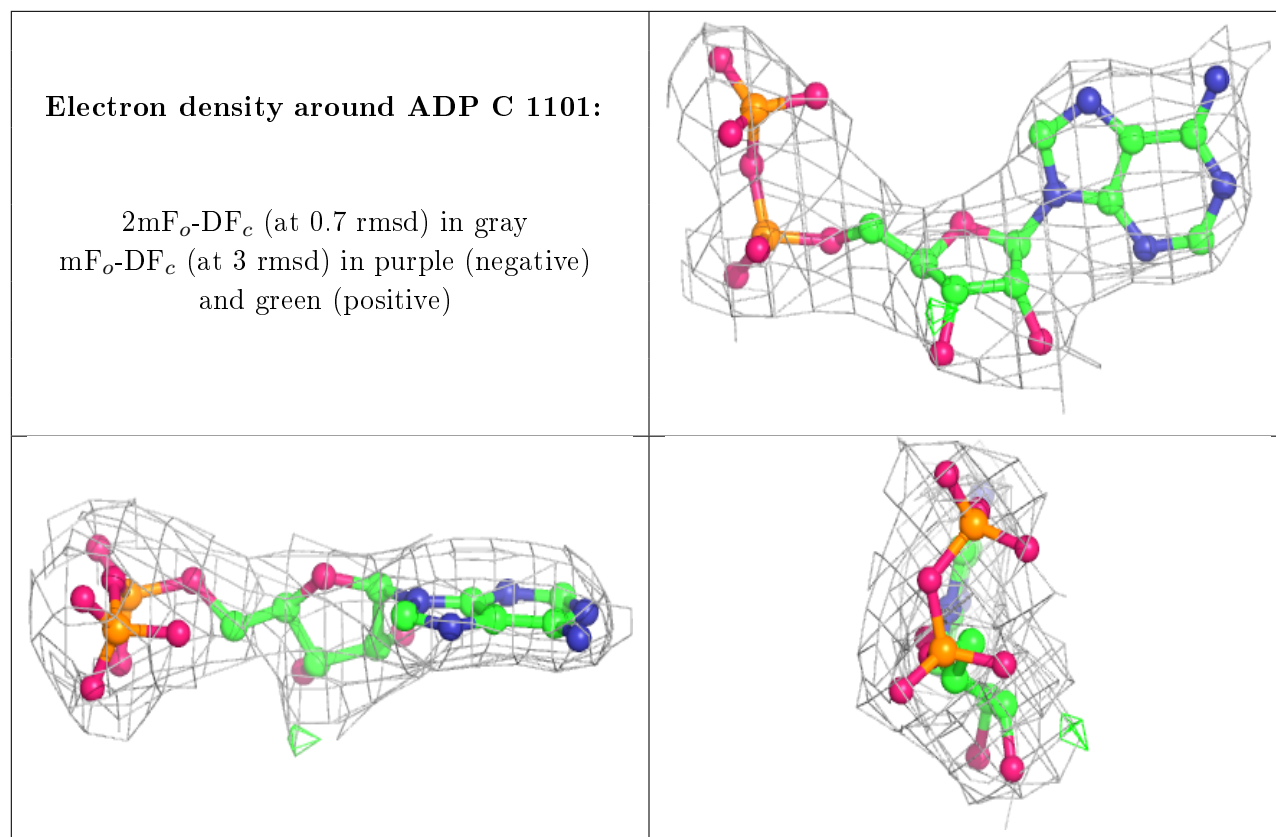
## 6.4 Ligands [i](#)

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
3	ADP	A	1101	27/27	0.84	0.28	52,82,168,171	0
3	ADP	C	1101	27/27	0.89	0.19	59,109,171,180	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.