



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 11:57 PM GMT

PDB ID : 1Z6T  
Title : Structure of the apoptotic protease-activating factor 1 bound to ADP  
Authors : Riedl, S.J.; Li, W.; Chao, Y.; Schwarzenbacher, R.; Shi, Y.  
Deposited on : 2005-03-23  
Resolution : 2.21 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  
<http://wwpdb.org/validation/2016/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 : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

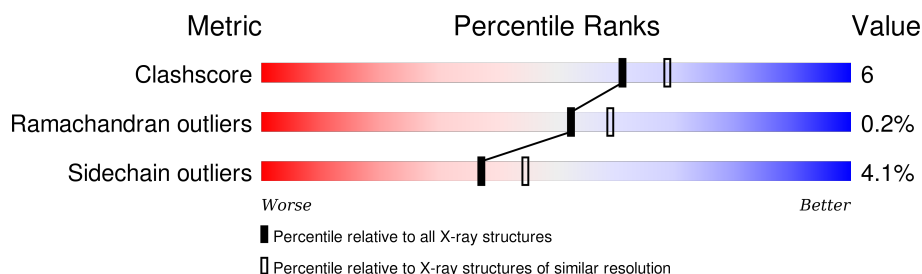
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.21 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))
Clashscore	102246	5146 (2.24-2.20)
Ramachandran outliers	100387	5065 (2.24-2.20)
Sidechain outliers	100360	5066 (2.24-2.20)

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. 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	591	 83% 13% ..
1	B	591	 85% 12% ..
1	C	591	 84% 14% ..
1	D	591	 84% 12% ..

## 2 Entry composition [i](#)

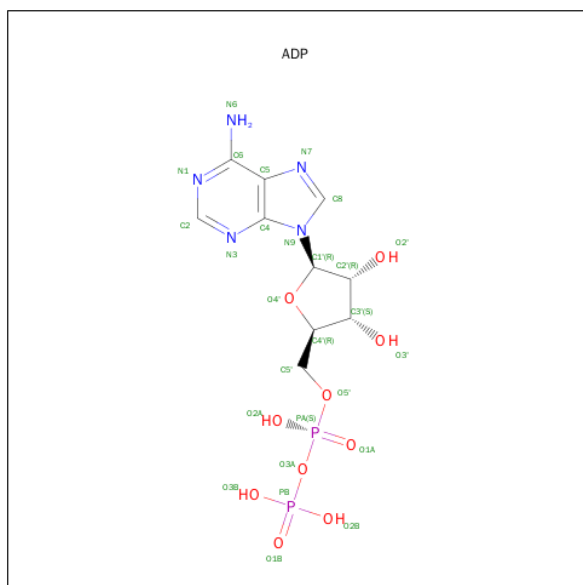
There are 3 unique types of molecules in this entry. The entry contains 19663 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Apoptotic protease activating factor 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	576	Total	C	N	O	S	0	0	0
			4656	2968	795	862	31			
1	B	586	Total	C	N	O	S	0	0	0
			4720	3003	806	880	31			
1	C	586	Total	C	N	O	S	0	0	0
			4720	3003	806	880	31			
1	D	576	Total	C	N	O	S	0	0	0
			4656	2968	795	862	31			

- Molecule 2 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula:  $C_{10}H_{15}N_5O_{10}P_2$ ).



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

- Molecule 3 is water.

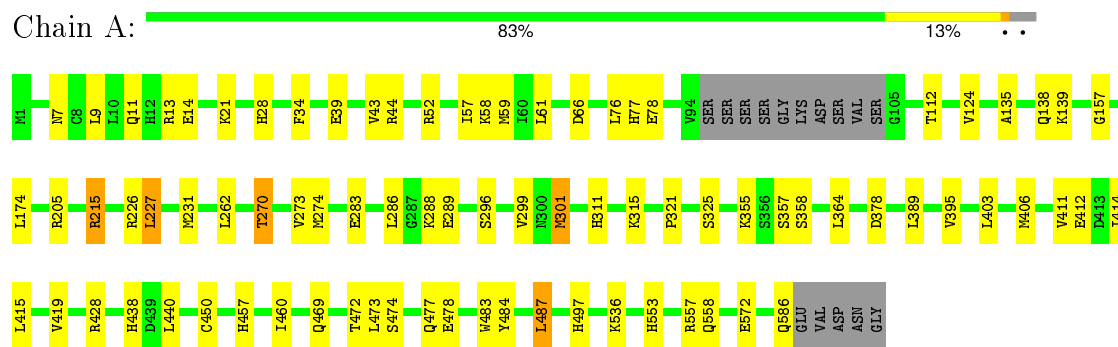
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	204	Total	O	0	0
			204	204		
3	B	220	Total	O	0	0
			220	220		
3	C	193	Total	O	0	0
			193	193		
3	D	186	Total	O	0	0
			186	186		

### 3 Residue-property plots

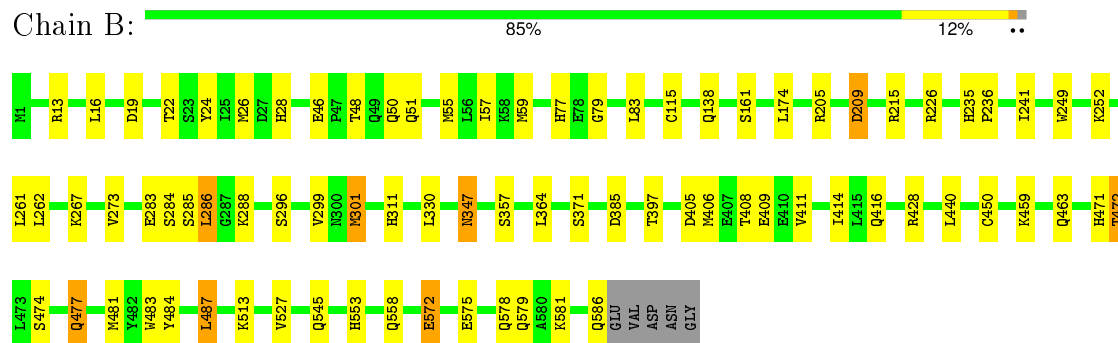
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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.

Note EDS was not executed.

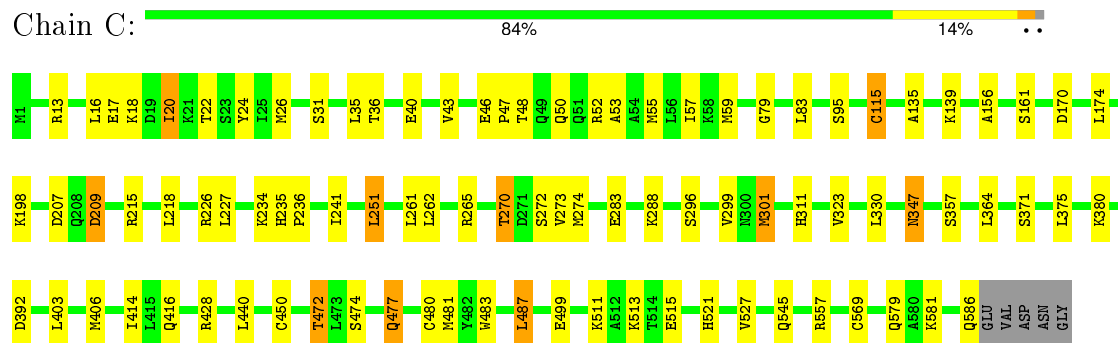
#### • Molecule 1: Apoptotic protease activating factor 1



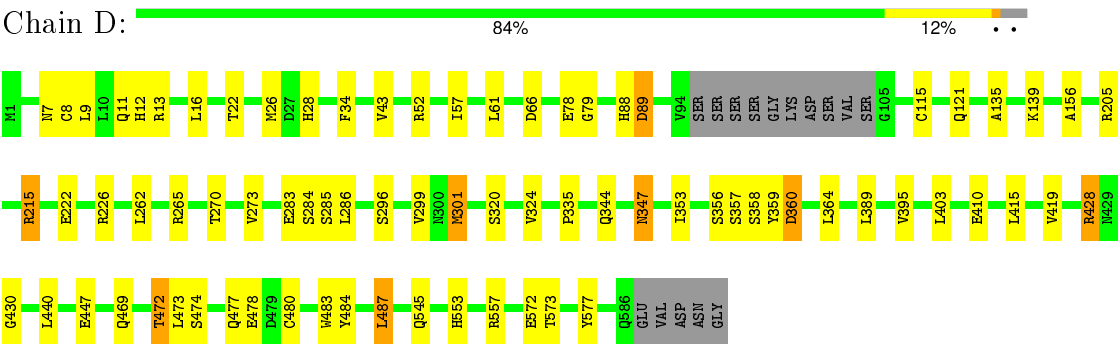
#### • Molecule 1: Apoptotic protease activating factor 1



#### • Molecule 1: Apoptotic protease activating factor 1



● Molecule 1: Apoptotic protease activating factor 1



## 4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	75.95Å 92.88Å 94.99Å 62.96° 89.99° 90.05°	Depositor
Resolution (Å)	15.00 – 2.21	Depositor
% Data completeness (in resolution range)	89.2 (15.00-2.21)	Depositor
$R_{merge}$	0.05	Depositor
$R_{sym}$	0.05	Depositor
Refinement program	REFMAC 5.2.0005	Depositor
R, $R_{free}$	0.189 , 0.244	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	19663	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	39.0	wwPDB-VP

## 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	0.63	0/4749	0.71	2/6405 (0.0%)
1	B	0.65	1/4814 (0.0%)	0.71	1/6493 (0.0%)
1	C	0.66	2/4814 (0.0%)	0.72	0/6493
1	D	0.65	0/4749	0.71	2/6405 (0.0%)
All	All	0.65	3/19126 (0.0%)	0.71	5/25796 (0.0%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	450	CYS	CB-SG	8.88	1.97	1.82
1	B	450	CYS	CB-SG	6.25	1.92	1.82
1	C	569	CYS	CB-SG	-5.42	1.73	1.81

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	205	ARG	NE-CZ-NH1	6.85	123.73	120.30
1	D	205	ARG	NE-CZ-NH2	-6.84	116.88	120.30
1	A	205	ARG	NE-CZ-NH2	-6.59	117.01	120.30
1	D	205	ARG	NE-CZ-NH1	5.61	123.10	120.30
1	B	205	ARG	NE-CZ-NH2	-5.50	117.55	120.30

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	4656	0	4696	54	1
1	B	4720	0	4756	51	1
1	C	4720	0	4756	59	0
1	D	4656	0	4696	51	2
2	A	27	0	12	1	0
2	B	27	0	12	0	0
2	C	27	0	12	0	0
2	D	27	0	12	0	0
3	A	204	0	0	13	0
3	B	220	0	0	11	0
3	C	193	0	0	8	2
3	D	186	0	0	10	0
All	All	19663	0	18952	208	3

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 6.

All (208) 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:26:MET:SD	1:C:35:LEU:HD21	1.72	1.29
1:A:157:GLY:HA2	3:A:1076:HOH:O	1.33	1.22
1:B:55:MET:HB2	3:B:1008:HOH:O	1.40	1.16
1:D:335:PRO:O	3:D:1050:HOH:O	1.82	0.96
1:D:121:GLN:OE1	3:D:912:HOH:O	1.84	0.94
1:C:26:MET:SD	1:C:35:LEU:CD2	2.55	0.93
1:A:450:CYS:SG	3:A:993:HOH:O	2.25	0.93
1:C:270:THR:HG21	3:C:1068:HOH:O	1.69	0.91
1:C:347:ASN:ND2	1:C:347:ASN:O	2.05	0.89
1:B:46:GLU:HG3	3:B:1008:HOH:O	1.73	0.86
1:C:472:THR:HG21	3:C:1047:HOH:O	1.78	0.84
1:D:270:THR:HG21	3:D:1000:HOH:O	1.81	0.80
1:B:55:MET:CB	3:B:1008:HOH:O	2.13	0.78
1:D:473:LEU:HD23	1:D:484:TYR:CE2	2.17	0.78
1:B:51:GLN:O	3:B:1008:HOH:O	1.99	0.78
1:C:579:GLN:OE1	3:C:945:HOH:O	2.03	0.76
1:B:579:GLN:OE1	3:B:930:HOH:O	2.03	0.76
1:B:13:ARG:NE	3:B:1009:HOH:O	2.06	0.75
1:A:44:ARG:NH2	3:A:1034:HOH:O	2.19	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:272:SER:OG	3:C:969:HOH:O	2.06	0.74
1:D:270:THR:HG23	3:D:1001:HOH:O	1.86	0.74
1:B:474:SER:O	1:B:477:GLN:HG3	1.88	0.74
1:A:473:LEU:HD23	1:A:484:TYR:CE2	2.22	0.74
1:B:347:ASN:O	1:B:347:ASN:ND2	2.22	0.73
1:D:474:SER:O	1:D:477:GLN:HG2	1.89	0.73
1:B:472:THR:CG2	3:B:947:HOH:O	2.37	0.72
1:A:270:THR:HG21	3:A:1071:HOH:O	1.89	0.71
1:A:270:THR:HG23	3:A:950:HOH:O	1.91	0.71
1:D:353:ILE:HB	1:D:447:GLU:OE2	1.91	0.71
1:A:438:HIS:NE2	3:A:1076:HOH:O	2.23	0.70
1:D:9:LEU:HD23	1:D:61:LEU:HD23	1.73	0.69
1:C:474:SER:O	1:C:477:GLN:HG2	1.94	0.68
1:A:536:LYS:H	1:A:536:LYS:HD2	1.57	0.68
1:A:321:PRO:O	3:A:970:HOH:O	2.11	0.67
1:C:357:SER:O	1:D:469:GLN:NE2	2.28	0.67
1:B:22:THR:HG23	3:B:1007:HOH:O	1.94	0.66
1:A:389:LEU:HD22	1:A:395:VAL:HG21	1.77	0.66
1:D:347:ASN:CG	1:D:347:ASN:O	2.34	0.66
1:A:9:LEU:HD23	1:A:61:LEU:HD23	1.77	0.66
1:D:222:GLU:OE2	1:D:557:ARG:NH2	2.29	0.65
1:A:536:LYS:H	1:A:536:LYS:CD	2.11	0.64
1:B:472:THR:HG22	3:B:947:HOH:O	2.00	0.62
1:B:22:THR:O	1:B:26:MET:HG3	2.00	0.62
1:C:79:GLY:HA2	1:C:115:CYS:SG	2.40	0.61
1:A:299:VAL:HG23	1:A:301:MET:HG3	1.81	0.61
1:A:215:ARG:HA	3:A:996:HOH:O	1.99	0.61
1:A:357:SER:OG	1:A:358:SER:N	2.34	0.61
1:A:76:LEU:HD22	1:A:112:THR:HG22	1.84	0.59
1:A:364:LEU:CD2	1:A:440:LEU:HB3	2.32	0.59
1:D:28:HIS:NE2	1:D:78:GLU:OE2	2.30	0.59
1:C:364:LEU:CD2	1:C:440:LEU:HB3	2.34	0.58
1:C:472:THR:CG2	3:C:931:HOH:O	2.51	0.58
1:D:357:SER:O	1:D:359:TYR:N	2.38	0.57
1:A:13:ARG:NH2	1:A:58:LYS:HG2	2.18	0.57
1:D:43:VAL:O	1:D:52:ARG:HG2	2.05	0.57
1:A:21:LYS:NZ	1:A:412:GLU:OE2	2.27	0.57
1:B:364:LEU:CD2	1:B:440:LEU:HB3	2.36	0.56
1:C:527:VAL:CG2	1:C:545:GLN:HE22	2.19	0.56
1:D:13:ARG:HG3	1:D:57:ILE:HG21	1.86	0.55
1:A:406:MET:HE2	1:A:411:VAL:HG22	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:573:THR:OG1	3:D:934:HOH:O	2.18	0.54
1:A:13:ARG:HH22	1:A:58:LYS:HG2	1.73	0.54
1:D:79:GLY:CA	1:D:115:CYS:SG	2.96	0.54
1:B:299:VAL:HG23	1:B:301:MET:HG2	1.90	0.53
1:B:24:TYR:CZ	1:B:416:GLN:OE1	2.62	0.53
1:C:209:ASP:C	1:C:209:ASP:OD1	2.46	0.53
1:C:20:ILE:HG22	1:C:83:LEU:CD1	2.39	0.53
1:D:215:ARG:HB3	3:D:1084:HOH:O	2.07	0.53
2:A:900:ADP:H5'1	3:A:1076:HOH:O	2.09	0.53
1:A:474:SER:N	1:A:477:GLN:OE1	2.36	0.53
1:A:76:LEU:HD22	1:A:112:THR:CG2	2.39	0.52
1:B:19:ASP:HB2	1:B:83:LEU:HD11	1.91	0.52
1:D:364:LEU:CD2	1:D:440:LEU:HB3	2.39	0.52
1:C:299:VAL:HG23	1:C:301:MET:CG	2.39	0.52
1:B:288:LYS:NZ	1:B:311:HIS:O	2.42	0.52
1:C:288:LYS:NZ	1:C:311:HIS:O	2.42	0.52
1:A:415:LEU:O	1:A:419:VAL:HG23	2.10	0.51
1:C:521:HIS:HB2	3:C:927:HOH:O	2.09	0.51
1:B:55:MET:O	1:B:59:MET:HG3	2.10	0.51
1:C:22:THR:HG21	1:C:43:VAL:HG11	1.93	0.51
1:D:156:ALA:HB2	1:D:265:ARG:CZ	2.41	0.51
1:C:207:ASP:HA	1:C:234:LYS:NZ	2.25	0.51
1:B:347:ASN:CG	1:B:347:ASN:O	2.49	0.51
1:C:17:GLU:O	1:C:50:GLN:HA	2.11	0.50
1:C:527:VAL:CG2	1:C:545:GLN:NE2	2.75	0.50
1:B:79:GLY:HA2	1:B:115:CYS:SG	2.51	0.50
1:D:415:LEU:O	1:D:419:VAL:HG23	2.12	0.50
1:B:463:GLN:NE2	3:B:948:HOH:O	2.44	0.50
1:D:473:LEU:HD23	1:D:484:TYR:CD2	2.47	0.49
1:B:299:VAL:HG23	1:B:301:MET:CG	2.42	0.49
1:B:385:ASP:OD2	1:B:459:LYS:NZ	2.46	0.49
1:D:572:GLU:HG3	1:D:577:TYR:CE1	2.47	0.49
1:C:24:TYR:CZ	1:C:416:GLN:OE1	2.65	0.49
1:D:34:PHE:CZ	1:D:66:ASP:HB3	2.48	0.49
1:D:553:HIS:O	1:D:557:ARG:HD2	2.13	0.49
1:C:26:MET:HG2	1:C:40:GLU:HG3	1.94	0.49
1:C:483:TRP:O	1:C:487:LEU:HB2	2.12	0.49
1:A:357:SER:O	1:B:471:HIS:NE2	2.45	0.49
1:C:48:THR:O	1:C:52:ARG:HG3	2.13	0.49
1:A:553:HIS:O	1:A:557:ARG:HD2	2.12	0.49
1:B:28:HIS:CE1	1:B:77:HIS:CE1	3.01	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:8:CYS:O	1:D:12:HIS:ND1	2.45	0.49
1:B:483:TRP:O	1:B:487:LEU:HB2	2.13	0.48
1:A:536:LYS:HD2	1:A:536:LYS:N	2.28	0.48
1:B:249:TRP:HA	1:B:252:LYS:HE2	1.94	0.48
1:B:406:MET:HE1	1:B:414:ILE:HD12	1.96	0.48
1:D:389:LEU:HD22	1:D:395:VAL:HG21	1.96	0.48
1:A:288:LYS:NZ	1:B:408:THR:HB	2.29	0.47
1:C:406:MET:HE1	1:C:414:ILE:HD12	1.96	0.47
1:D:359:TYR:CZ	1:D:440:LEU:HD22	2.49	0.47
1:C:13:ARG:HG3	1:C:57:ILE:HG21	1.97	0.47
1:B:553:HIS:H	1:B:553:HIS:CD2	2.32	0.47
1:C:55:MET:O	1:C:59:MET:HG3	2.15	0.47
1:A:406:MET:HE1	1:A:414:ILE:HD12	1.96	0.47
1:B:235:HIS:N	1:B:236:PRO:CD	2.77	0.47
1:B:527:VAL:CG2	1:B:545:GLN:HE22	2.28	0.47
1:D:299:VAL:HG23	1:D:301:MET:CG	2.45	0.47
1:D:428:ARG:HD3	1:D:428:ARG:N	2.30	0.47
1:A:364:LEU:HD21	1:A:440:LEU:HB3	1.96	0.47
1:C:207:ASP:HA	1:C:234:LYS:HZ2	1.80	0.47
1:A:13:ARG:HG3	1:A:57:ILE:HG21	1.97	0.47
1:A:43:VAL:O	1:A:52:ARG:HG2	2.14	0.47
1:D:477:GLN:HG3	3:D:909:HOH:O	2.15	0.46
1:D:88:HIS:O	1:D:89:ASP:C	2.53	0.46
1:C:299:VAL:HG23	1:C:301:MET:HG2	1.97	0.46
1:B:527:VAL:CG2	1:B:545:GLN:NE2	2.79	0.46
1:D:299:VAL:HG23	1:D:301:MET:HG3	1.97	0.46
1:C:22:THR:O	1:C:22:THR:HG22	2.14	0.46
1:A:124:VAL:CG2	3:A:1019:HOH:O	2.62	0.46
1:C:357:SER:C	1:D:469:GLN:HE22	2.19	0.46
1:C:170:ASP:O	1:C:174:LEU:HB2	2.16	0.46
1:D:283:GLU:HG3	1:D:285:SER:O	2.15	0.46
1:D:79:GLY:HA2	1:D:115:CYS:SG	2.56	0.46
1:C:364:LEU:HD22	1:C:440:LEU:HB3	1.99	0.45
1:C:20:ILE:CG2	1:C:83:LEU:HD13	2.46	0.45
1:A:288:LYS:NZ	1:B:409:GLU:OE1	2.50	0.45
1:A:473:LEU:HD23	1:A:484:TYR:CD2	2.51	0.45
1:A:315:LYS:NZ	1:B:405:ASP:OD1	2.49	0.45
1:A:13:ARG:NH2	1:A:58:LYS:CG	2.80	0.45
1:A:39:GLU:HG2	1:A:59:MET:HE2	1.99	0.45
1:C:241:ILE:HA	1:C:261:LEU:O	2.17	0.45
1:C:499:GLU:OE1	3:C:1004:HOH:O	2.21	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:480:CYS:HA	3:D:909:HOH:O	2.17	0.45
1:D:296:SER:HB2	1:D:301:MET:O	2.17	0.44
1:B:296:SER:HB2	1:B:301:MET:O	2.17	0.44
1:D:577:TYR:CD1	1:D:577:TYR:C	2.90	0.44
1:A:553:HIS:CD2	1:A:553:HIS:H	2.35	0.44
1:B:406:MET:HE2	1:B:411:VAL:HG22	2.00	0.44
1:D:320:SER:O	1:D:324:VAL:HG23	2.17	0.44
1:D:22:THR:O	1:D:26:MET:HG3	2.17	0.44
1:B:397:THR:HG21	1:B:408:THR:HG23	1.99	0.43
1:A:135:ALA:O	1:A:139:LYS:HG2	2.17	0.43
1:B:481:MET:HE2	1:B:481:MET:HA	2.00	0.43
1:B:209:ASP:OD1	1:B:209:ASP:C	2.56	0.43
1:D:359:TYR:CE2	1:D:440:LEU:HD22	2.53	0.43
1:A:227:LEU:O	1:A:231:MET:HG2	2.18	0.43
1:C:218:LEU:HD12	1:C:557:ARG:HE	1.83	0.43
1:A:457:HIS:O	1:A:460:ILE:HG22	2.18	0.43
1:C:296:SER:HB2	1:C:301:MET:O	2.19	0.43
1:D:364:LEU:HD22	1:D:440:LEU:HB3	2.00	0.43
1:B:138:GLN:HA	1:B:138:GLN:OE1	2.17	0.43
1:A:311:HIS:CD2	1:B:48:THR:CG2	3.02	0.43
1:B:330:LEU:HD21	1:B:371:SER:HA	2.00	0.43
1:A:406:MET:CE	1:A:411:VAL:HG22	2.48	0.43
1:C:20:ILE:CG2	1:C:83:LEU:CD1	2.96	0.43
1:C:375:LEU:O	1:C:380:LYS:HE3	2.18	0.43
1:C:330:LEU:HD21	1:C:371:SER:HA	2.00	0.43
1:A:325:SER:OG	3:A:970:HOH:O	2.17	0.42
1:A:483:TRP:O	1:A:487:LEU:HB2	2.19	0.42
1:A:34:PHE:CZ	1:A:66:ASP:HB3	2.54	0.42
1:B:572:GLU:O	1:B:578:GLN:NE2	2.42	0.42
1:A:296:SER:HB2	1:A:301:MET:O	2.18	0.42
1:A:289:GLU:HG3	1:B:50:GLN:HG3	2.01	0.42
1:A:28:HIS:NE2	1:A:78:GLU:OE2	2.35	0.42
1:C:156:ALA:HB2	1:C:265:ARG:CZ	2.49	0.42
1:D:78:GLU:O	1:D:430:GLY:N	2.42	0.42
1:C:17:GLU:HA	1:C:53:ALA:HB3	2.02	0.42
1:C:480:CYS:SG	1:C:481:MET:CE	3.08	0.42
1:C:513:LYS:NZ	1:C:521:HIS:HD2	2.18	0.42
1:A:124:VAL:HG22	3:A:1019:HOH:O	2.19	0.42
1:B:285:SER:OG	1:B:286:LEU:O	2.36	0.42
1:B:484:TYR:O	1:B:513:LYS:NZ	2.53	0.42
1:A:39:GLU:HA	1:A:59:MET:HE1	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:135:ALA:O	1:C:139:LYS:HG2	2.19	0.41
1:C:299:VAL:HG23	1:C:301:MET:HG3	2.01	0.41
1:D:428:ARG:N	3:D:912:HOH:O	2.50	0.41
1:D:473:LEU:CD1	3:D:909:HOH:O	2.69	0.41
1:B:16:LEU:HB3	1:B:57:ILE:HD11	2.01	0.41
1:C:198:LYS:NZ	1:C:392:ASP:O	2.47	0.41
1:C:511:LYS:O	1:C:515:GLU:HG3	2.21	0.41
1:C:235:HIS:N	1:C:236:PRO:CD	2.84	0.41
1:C:35:LEU:HG	1:C:36:THR:N	2.36	0.41
1:C:472:THR:HG22	3:C:931:HOH:O	2.20	0.41
1:B:79:GLY:CA	1:B:115:CYS:SG	3.09	0.41
1:C:480:CYS:SG	1:C:481:MET:HE2	2.61	0.41
1:D:472:THR:CG2	1:D:472:THR:O	2.69	0.41
1:D:7:ASN:O	1:D:11:GLN:HB2	2.21	0.41
1:C:16:LEU:HB3	1:C:57:ILE:HD11	2.03	0.41
1:D:135:ALA:O	1:D:139:LYS:HG2	2.21	0.41
1:B:241:ILE:HA	1:B:261:LEU:O	2.22	0.40
1:D:16:LEU:HB3	1:D:57:ILE:HD11	2.03	0.40
1:B:267:LYS:NZ	3:B:960:HOH:O	2.54	0.40
1:A:7:ASN:O	1:A:11:GLN:HB2	2.21	0.40
1:A:77:HIS:CE1	3:A:1020:HOH:O	2.73	0.40
1:C:251:LEU:HD12	1:C:251:LEU:HA	1.80	0.40
1:C:46:GLU:HA	1:C:47:PRO:HD3	1.95	0.40
1:D:483:TRP:O	1:D:487:LEU:HB2	2.22	0.40
1:C:323:VAL:HG22	1:C:364:LEU:HA	2.04	0.40

All (3) 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:D:360:ASP:OD1	3:C:1047:HOH:O[1_546]	1.97	0.23
1:D:360:ASP:OD2	3:C:1061:HOH:O[1_546]	2.10	0.10
1:A:469:GLN:NE2	1:B:357:SER:O[1_546]	2.17	0.03

## 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	572/591 (97%)	560 (98%)	11 (2%)	1 (0%)	52	59
1	B	584/591 (99%)	576 (99%)	8 (1%)	0	100	100
1	C	584/591 (99%)	571 (98%)	12 (2%)	1 (0%)	52	59
1	D	572/591 (97%)	558 (98%)	12 (2%)	2 (0%)	46	50
All	All	2312/2364 (98%)	2265 (98%)	43 (2%)	4 (0%)	52	59

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	355	LYS
1	D	358	SER
1	D	89	ASP
1	C	18	LYS

### 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	521/534 (98%)	498 (96%)	23 (4%)	35	41
1	B	530/534 (99%)	509 (96%)	21 (4%)	38	46
1	C	530/534 (99%)	506 (96%)	24 (4%)	34	40
1	D	521/534 (98%)	503 (96%)	18 (4%)	43	52
All	All	2102/2136 (98%)	2016 (96%)	86 (4%)	37	45

All (86) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	GLU
1	A	138	GLN

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Mol	Chain	Res	Type
1	A	174	LEU
1	A	215	ARG
1	A	226	ARG
1	A	227	LEU
1	A	262	LEU
1	A	270	THR
1	A	273	VAL
1	A	274	MET
1	A	283	GLU
1	A	286	LEU
1	A	301	MET
1	A	378	ASP
1	A	403	LEU
1	A	428	ARG
1	A	472	THR
1	A	478	GLU
1	A	487	LEU
1	A	497	HIS
1	A	558	GLN
1	A	572	GLU
1	A	586	GLN
1	B	161	SER
1	B	174	LEU
1	B	209	ASP
1	B	215	ARG
1	B	226	ARG
1	B	262	LEU
1	B	273	VAL
1	B	283	GLU
1	B	284	SER
1	B	286	LEU
1	B	301	MET
1	B	347	ASN
1	B	428	ARG
1	B	472	THR
1	B	477	GLN
1	B	487	LEU
1	B	558	GLN
1	B	572	GLU
1	B	575	GLU
1	B	581	LYS
1	B	586	GLN

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Mol	Chain	Res	Type
1	C	20	ILE
1	C	31	SER
1	C	95	SER
1	C	115	CYS
1	C	161	SER
1	C	209	ASP
1	C	215	ARG
1	C	226	ARG
1	C	227	LEU
1	C	251	LEU
1	C	262	LEU
1	C	270	THR
1	C	273	VAL
1	C	274	MET
1	C	283	GLU
1	C	301	MET
1	C	347	ASN
1	C	403	LEU
1	C	428	ARG
1	C	472	THR
1	C	477	GLN
1	C	487	LEU
1	C	581	LYS
1	C	586	GLN
1	D	215	ARG
1	D	226	ARG
1	D	262	LEU
1	D	273	VAL
1	D	284	SER
1	D	286	LEU
1	D	301	MET
1	D	344	GLN
1	D	347	ASN
1	D	356	SER
1	D	360	ASP
1	D	403	LEU
1	D	410	GLU
1	D	428	ARG
1	D	472	THR
1	D	478	GLU
1	D	487	LEU
1	D	545	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (32) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	65	ASN
1	A	77	HIS
1	A	134	ASN
1	A	153	HIS
1	A	344	GLN
1	A	420	ASN
1	A	429	ASN
1	A	463	GLN
1	A	553	HIS
1	A	586	GLN
1	B	11	GLN
1	B	134	ASN
1	B	153	HIS
1	B	347	ASN
1	B	521	HIS
1	B	524	HIS
1	B	543	ASN
1	B	553	HIS
1	B	586	GLN
1	C	11	GLN
1	C	134	ASN
1	C	153	HIS
1	C	477	GLN
1	C	521	HIS
1	C	543	ASN
1	C	586	GLN
1	D	7	ASN
1	D	73	ASN
1	D	134	ASN
1	D	153	HIS
1	D	347	ASN
1	D	521	HIS

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

4 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	$\# Z  > 2$
2	ADP	A	900	-	22,29,29	1.06	2 (9%)	27,45,45	2.03	4 (14%)
2	ADP	B	901	-	22,29,29	0.92	2 (9%)	27,45,45	2.04	6 (22%)
2	ADP	C	902	-	22,29,29	0.92	1 (4%)	27,45,45	1.95	3 (11%)
2	ADP	D	903	-	22,29,29	1.09	2 (9%)	27,45,45	1.85	5 (18%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	ADP	A	900	-	-	0/12/32/32	0/3/3/3
2	ADP	B	901	-	-	0/12/32/32	0/3/3/3
2	ADP	C	902	-	-	0/12/32/32	0/3/3/3
2	ADP	D	903	-	-	0/12/32/32	0/3/3/3

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	903	ADP	O4'-C1'	2.03	1.43	1.41
2	B	901	ADP	O4'-C1'	2.03	1.43	1.41
2	B	901	ADP	C5-C4	2.10	1.45	1.40
2	A	900	ADP	O4'-C1'	2.24	1.44	1.41
2	C	902	ADP	C5-C4	2.76	1.46	1.40
2	A	900	ADP	C5-C4	2.77	1.46	1.40
2	D	903	ADP	C5-C4	3.05	1.47	1.40

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	902	ADP	N3-C2-N1	-8.49	122.39	128.89
2	B	901	ADP	N3-C2-N1	-8.12	122.67	128.89
2	A	900	ADP	N3-C2-N1	-7.87	122.87	128.89
2	D	903	ADP	N3-C2-N1	-6.91	123.61	128.89
2	D	903	ADP	C4-C5-N7	-2.56	107.13	109.48
2	A	900	ADP	C4-C5-N7	-2.53	107.16	109.48
2	C	902	ADP	C1'-N9-C4	-2.35	123.40	126.94
2	B	901	ADP	C2'-C1'-N9	-2.22	110.90	114.29
2	B	901	ADP	PA-O3A-PB	-2.18	125.35	132.67
2	B	901	ADP	C4-C5-N7	-2.09	107.56	109.48
2	B	901	ADP	O2A-PA-O1A	2.06	123.67	112.53
2	D	903	ADP	O3B-PB-O1B	2.16	117.53	110.58
2	C	902	ADP	O3B-PB-O2B	2.18	115.68	107.38
2	A	900	ADP	C2-N1-C6	2.32	122.91	118.77
2	D	903	ADP	O3B-PB-O2B	2.33	116.27	107.38
2	B	901	ADP	O3B-PB-O2B	3.25	119.75	107.38
2	A	900	ADP	O3B-PB-O2B	3.30	119.93	107.38
2	D	903	ADP	C2'-C1'-N9	3.30	119.33	114.29

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	900	ADP	1	0

## 5.7 Other polymers ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

### 6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

### 6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

### 6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.