



# wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 1, 2016 – 05:04 AM GMT

PDB ID : 2P9S  
Title : Structure of bovine Arp2/3 complex co-crystallized with ATP/Mg2+  
Authors : Nolen, B.J.; Pollard, T.D.  
Deposited on : 2007-03-26  
Resolution : 2.68 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

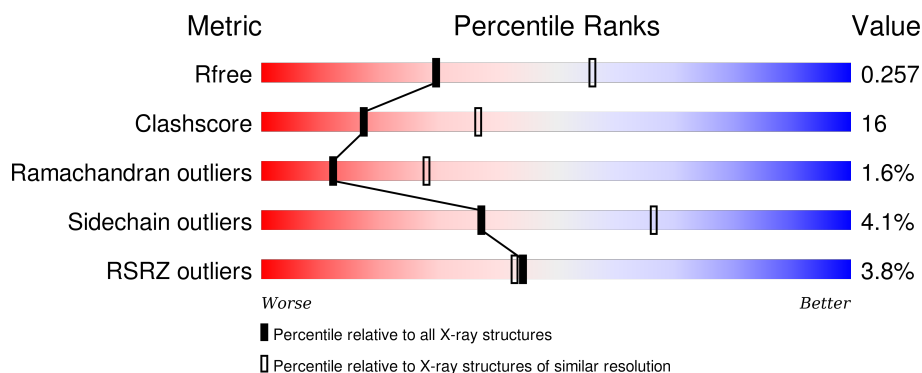
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.68 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))
$R_{free}$	91344	2780 (2.70-2.66)
Clashscore	102246	3138 (2.70-2.66)
Ramachandran outliers	100387	3089 (2.70-2.66)
Sidechain outliers	100360	3089 (2.70-2.66)
RSRZ outliers	91569	2789 (2.70-2.66)

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.

Mol	Chain	Length	Quality of chain
1	A	418	<div> <div>2%</div> <div>71% 21% 5%</div> </div>
2	B	394	<div> <div>7%</div> <div>30% 18% 50%</div> </div>
3	C	372	<div> <div>%</div> <div>61% 28% 8%</div> </div>
4	D	300	<div> <div>%</div> <div>70% 22% 7%</div> </div>
5	E	178	<div> <div>4%</div> <div>58% 34% . .</div> </div>

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Mol	Chain	Length	Quality of chain
6	F	168	<div><div>%</div><div><div></div><div>76%</div><div>21%</div><div>• •</div></div></div>
7	G	151	<div><div>10%</div><div><div></div><div>59%</div><div>27%</div><div>•</div><div>11%</div></div></div>

## 2 Entry composition [i](#)

There are 10 unique types of molecules in this entry. The entry contains 13608 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 Actin-like protein 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	397	Total	C	N	O	S	0	0	0
			3125	2006	517	587	15			

- Molecule 2 is a protein called Actin-like protein 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	197	Total	C	N	O	S	0	0	0
			1487	951	251	281	4			

- Molecule 3 is a protein called Actin-related protein 2/3 complex subunit 1B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	341	Total	C	N	O	S	0	0	0
			2641	1676	463	483	19			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	58	VAL	ILE	CONFLICT	UNP Q58CQ2

- Molecule 4 is a protein called Actin-related protein 2/3 complex subunit 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	280	Total	C	N	O	S	0	0	0
			2249	1430	392	419	8			

- Molecule 5 is a protein called Actin-related protein 2/3 complex subunit 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	E	173	Total	C	N	O	S	0	0	0
			1411	906	235	261	9			

- Molecule 6 is a protein called Actin-related protein 2/3 complex subunit 4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	F	167	Total	C	N	O	S	0	0	0
			1371	875	239	248	9			

- Molecule 7 is a protein called Actin-related protein 2/3 complex subunit 5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	G	135	Total	C	N	O	S	0	0	0
			1028	643	177	205	3			

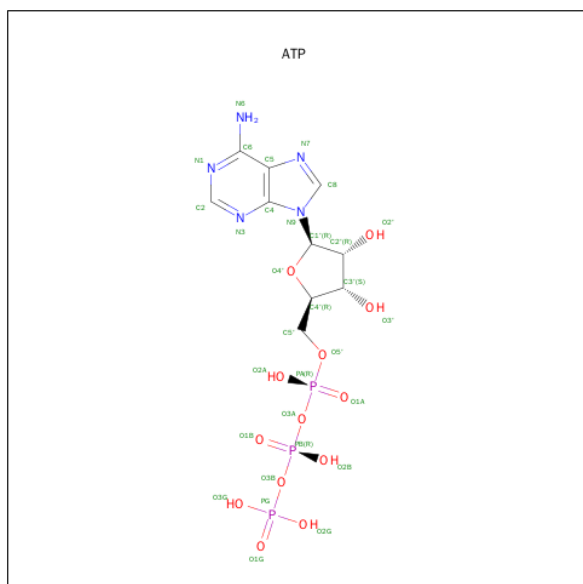
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
G	17	ASP	GLY	CONFLICT	UNP Q3SYX9
G	28	ASP	GLU	CONFLICT	UNP Q3SYX9

- Molecule 8 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	1	Total	Mg	0	0
			1	1		

- Molecule 9 is ADENOSINE-5'-TRIPHOSPHATE (three-letter code: ATP) (formula: C<sub>10</sub>H<sub>16</sub>N<sub>5</sub>O<sub>13</sub>P<sub>3</sub>).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
9	A	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
9	B	1	Total	C	N	O	P	0	0
			31	10	5	13	3		

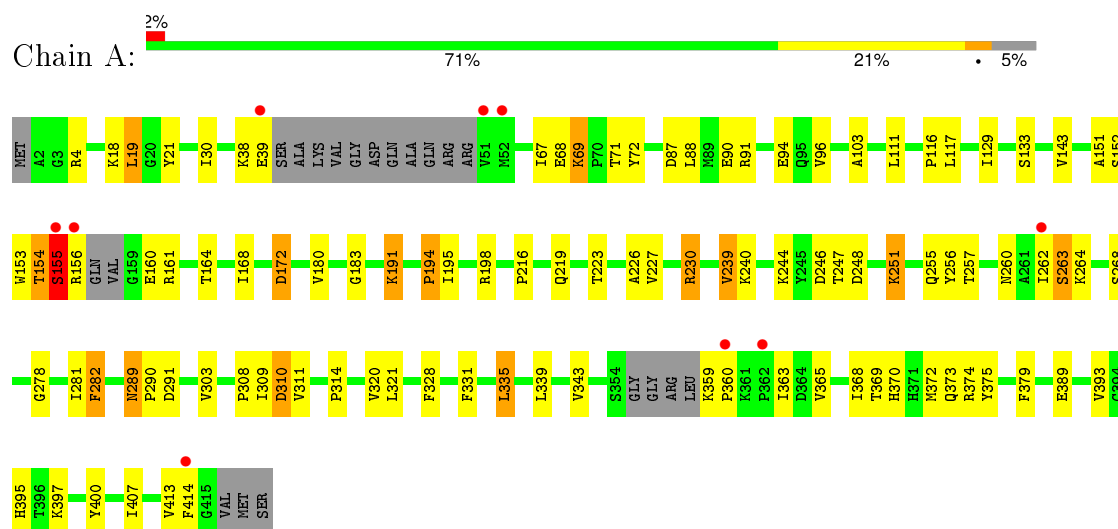
- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	50	Total	O	0	0
			50	50		
10	B	15	Total	O	0	0
			15	15		
10	C	61	Total	O	0	0
			61	61		
10	D	47	Total	O	0	0
			47	47		
10	E	14	Total	O	0	0
			14	14		
10	F	39	Total	O	0	0
			39	39		
10	G	7	Total	O	0	0
			7	7		

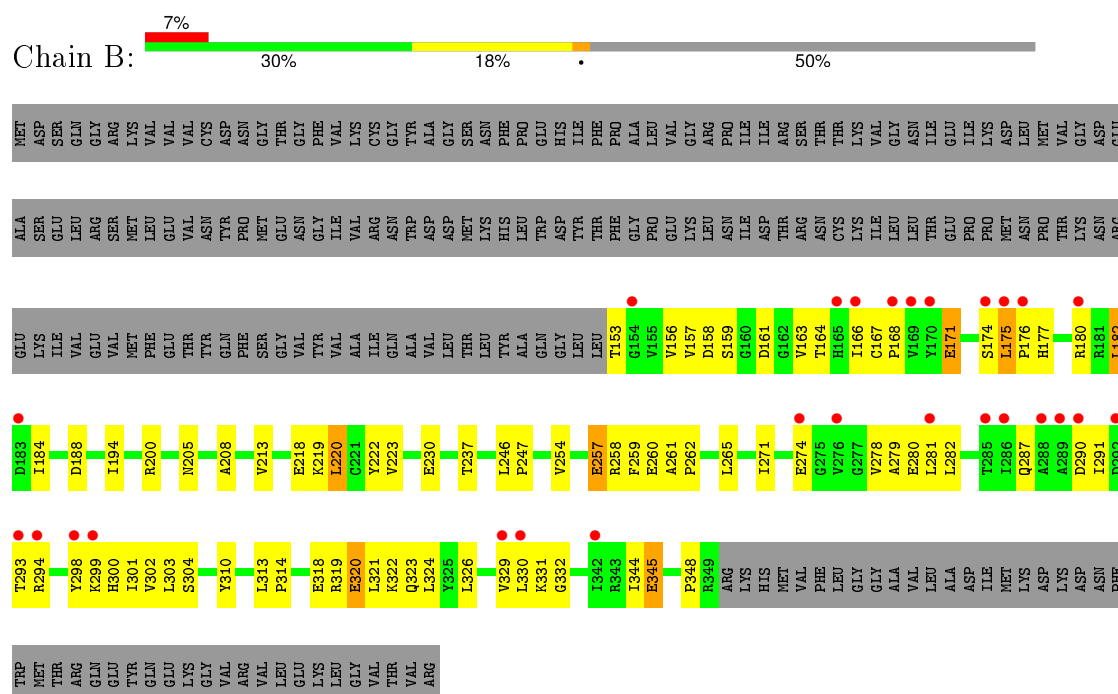
### 3 Residue-property plots

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.

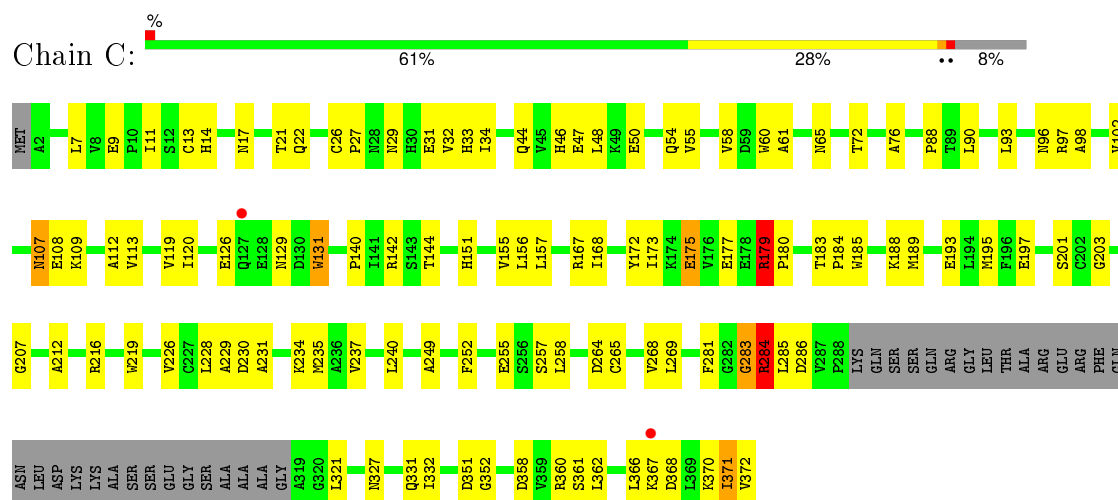
#### • Molecule 1: Actin-like protein 3



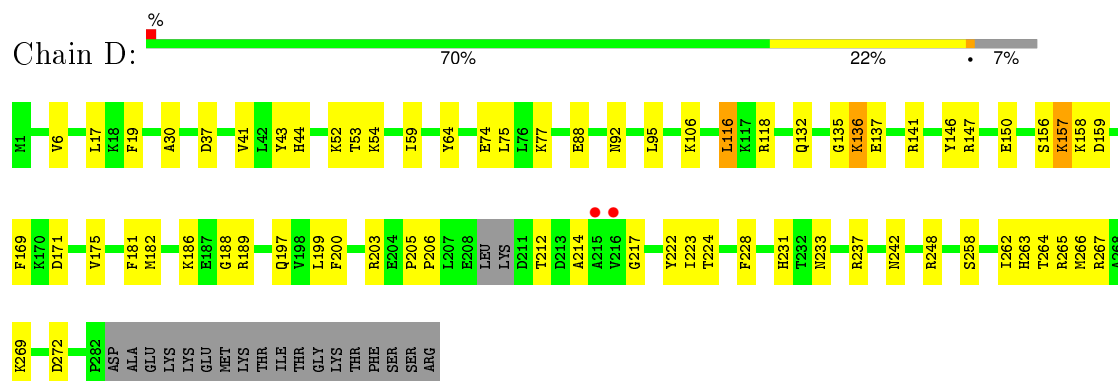
#### • Molecule 2: Actin-like protein 2



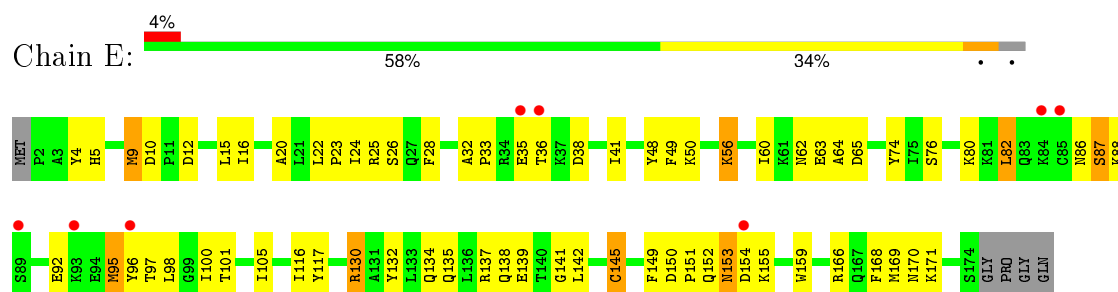
- Molecule 3: Actin-related protein 2/3 complex subunit 1B



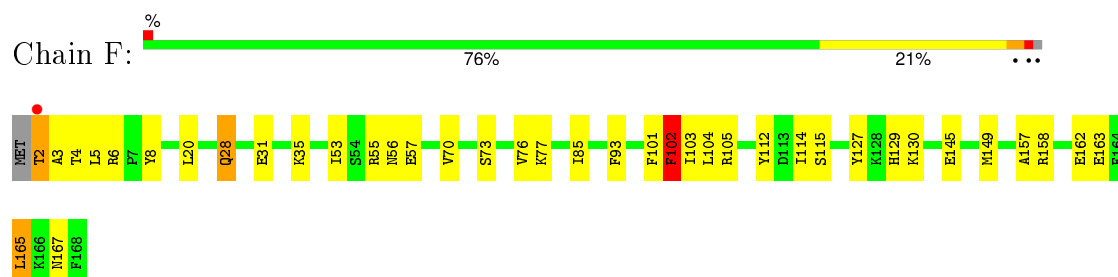
- Molecule 4: Actin-related protein 2/3 complex subunit 2



- Molecule 5: Actin-related protein 2/3 complex subunit 3

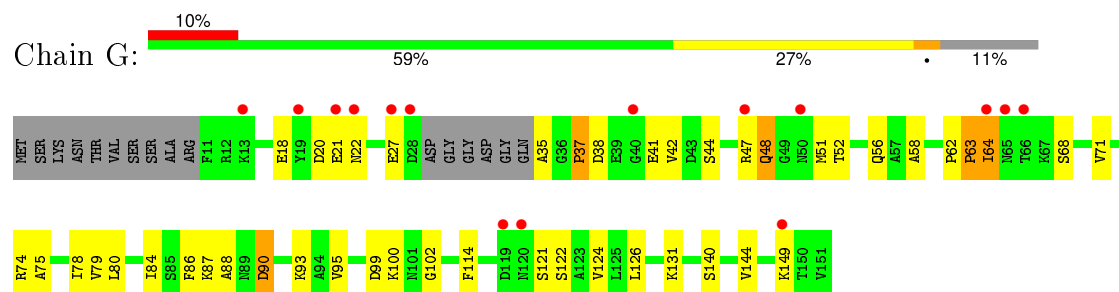


- Molecule 6: Actin-related protein 2/3 complex subunit 4





Chain G:



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	111.02Å 128.69Å 201.40Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.68 46.45 – 2.68	Depositor EDS
% Data completeness (in resolution range)	95.0 (30.00-2.68) 95.3 (46.45-2.68)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	10.40	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.70 (at 2.69Å)	Xtriage
Refinement program	CNS	Depositor
R, $R_{free}$	0.222 , 0.261 0.220 , 0.257	Depositor DCC
$R_{free}$ test set	3906 reflections (5.04%)	DCC
Wilson B-factor (Å <sup>2</sup> )	50.6	Xtriage
Anisotropy	0.431	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 44.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 80818 reflections	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	13608	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.78% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: MG, ATP

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	0.42	0/3205	0.63	1/4358 (0.0%)
2	B	0.34	0/1514	0.59	0/2059
3	C	0.40	0/2710	0.70	1/3680 (0.0%)
4	D	0.39	0/2297	0.59	0/3101
5	E	0.36	0/1445	0.62	0/1949
6	F	0.40	0/1393	0.61	0/1868
7	G	0.32	0/1040	0.55	0/1401
All	All	0.39	0/13604	0.62	2/18416 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	155	SER	N-CA-C	7.53	131.34	111.00
3	C	283	GLY	N-CA-C	6.30	128.86	113.10

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	3125	0	3001	87	0
2	B	1487	0	1440	56	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	2641	0	2589	100	0
4	D	2249	0	2206	62	0
5	E	1411	0	1413	71	0
6	F	1371	0	1410	34	0
7	G	1028	0	1029	31	0
8	A	1	0	0	0	0
9	A	31	0	12	2	0
9	B	31	0	12	3	0
10	A	50	0	0	2	0
10	B	15	0	0	0	0
10	C	61	0	0	1	0
10	D	47	0	0	1	0
10	E	14	0	0	0	0
10	F	39	0	0	0	0
10	G	7	0	0	0	0
All	All	13608	0	13112	425	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 16.

The worst 5 of 425 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)
3:C:216:ARG:HG2	3:C:230:ASP:HB3	1.32	1.10
3:C:183:THR:HG22	3:C:185:TRP:H	1.24	1.01
3:C:155:VAL:HG21	3:C:180:PRO:HG3	1.45	0.99
3:C:14:HIS:H	3:C:331:GLN:HE22	1.07	0.91
3:C:201:SER:HB3	7:G:149:LYS:NZ	1.90	0.87

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	389/418 (93%)	356 (92%)	29 (8%)	4 (1%)	19	42
2	B	195/394 (50%)	175 (90%)	14 (7%)	6 (3%)	5	11
3	C	337/372 (91%)	315 (94%)	18 (5%)	4 (1%)	16	37
4	D	276/300 (92%)	260 (94%)	15 (5%)	1 (0%)	39	67
5	E	171/178 (96%)	158 (92%)	9 (5%)	4 (2%)	8	19
6	F	165/168 (98%)	155 (94%)	8 (5%)	2 (1%)	16	37
7	G	131/151 (87%)	113 (86%)	13 (10%)	5 (4%)	4	8
All	All	1664/1981 (84%)	1532 (92%)	106 (6%)	26 (2%)	12	28

5 of 26 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	155	SER
2	B	171	GLU
2	B	345	GLU
3	C	371	ILE
5	E	87	SER

### 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	333/363 (92%)	317 (95%)	16 (5%)	31	59
2	B	151/345 (44%)	143 (95%)	8 (5%)	28	54
3	C	288/313 (92%)	277 (96%)	11 (4%)	40	69
4	D	242/264 (92%)	238 (98%)	4 (2%)	68	89
5	E	156/159 (98%)	149 (96%)	7 (4%)	34	62
6	F	154/155 (99%)	146 (95%)	8 (5%)	29	55
7	G	111/124 (90%)	106 (96%)	5 (4%)	34	62
All	All	1435/1723 (83%)	1376 (96%)	59 (4%)	37	66

5 of 59 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	C	131	TRP
3	C	367	LYS
7	G	18	GLU
3	C	140	PRO
3	C	179	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 32 such sidechains are listed below:

Mol	Chain	Res	Type
3	C	46	HIS
3	C	331	GLN
7	G	48	GLN
3	C	107	ASN
4	D	132	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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 1 is monoatomic - leaving 2 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
9	ATP	A	501	8	24,33,33	1.37	3 (12%)	31,52,52	2.47	5 (16%)
9	ATP	B	502	-	24,33,33	1.40	3 (12%)	31,52,52	2.46	4 (12%)

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
9	ATP	A	501	8	-	0/18/38/38	0/3/3/3
9	ATP	B	502	-	-	0/18/38/38	0/3/3/3

The worst 5 of 6 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	A	501	ATP	O4'-C1'	2.16	1.43	1.41
9	B	502	ATP	O4'-C1'	2.36	1.44	1.41
9	A	501	ATP	PG-O1G	3.14	1.61	1.51
9	B	502	ATP	PG-O1G	3.24	1.61	1.51
9	A	501	ATP	C2-N1	3.54	1.40	1.33

The worst 5 of 9 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	B	502	ATP	N3-C2-N1	-11.37	120.19	128.89
9	A	501	ATP	N3-C2-N1	-11.35	120.21	128.89
9	B	502	ATP	PA-O3A-PB	-4.12	121.16	132.73
9	A	501	ATP	PA-O3A-PB	-3.97	121.57	132.73
9	B	502	ATP	PB-O3B-PG	-3.88	119.67	132.67

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
9	A	501	ATP	2	0
9	B	502	ATP	3	0

## 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	397/418 (94%)	-0.06	9 (2%) 64 62	25, 48, 81, 96	0
2	B	197/394 (50%)	0.41	27 (13%) 4 3	34, 62, 96, 111	0
3	C	341/372 (91%)	-0.08	2 (0%) 90 91	27, 41, 69, 92	0
4	D	280/300 (93%)	-0.04	2 (0%) 89 89	27, 44, 78, 87	0
5	E	173/178 (97%)	0.12	8 (4%) 36 34	37, 58, 85, 100	0
6	F	167/168 (99%)	-0.17	1 (0%) 90 91	30, 43, 59, 89	0
7	G	135/151 (89%)	0.49	15 (11%) 7 5	35, 72, 96, 103	0
All	All	1690/1981 (85%)	0.05	64 (3%) 44 43	25, 49, 88, 111	0

The worst 5 of 64 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	51	VAL	8.5
5	E	154	ASP	4.7
2	B	290	ASP	4.3
1	A	52	MET	4.3
7	G	120	ASN	4.3

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

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

### 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
9	ATP	A	501	31/31	0.97	0.15	-0.12	41,45,48,51	0
9	ATP	B	502	31/31	0.94	0.12	-1.15	51,58,88,89	0
8	MG	A	500	1/1	0.93	0.43	-	43,43,43,43	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.