



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 1, 2016 – 05:14 PM GMT

PDB ID : 4HMY
Title : Structural basis for recruitment and activation of the AP-1 clathrin adaptor complex by Arf1
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Deposited on : 2012-10-18
Resolution : 7.00 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
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

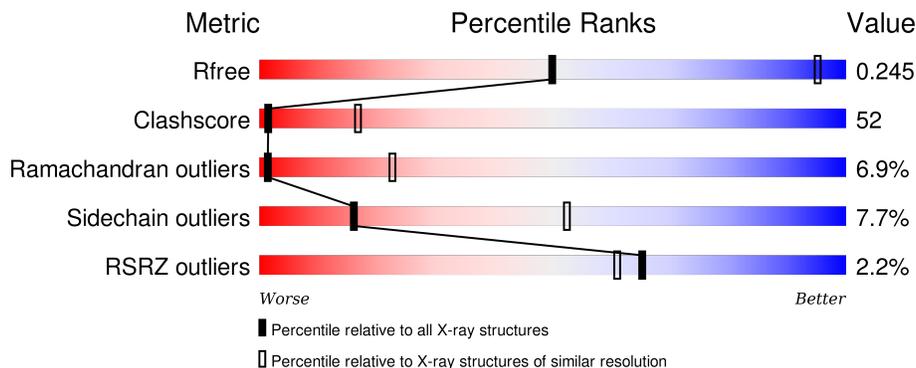
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 7.00 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	1014 (9.50-3.66)
Clashscore	102246	1063 (10.00-3.70)
Ramachandran outliers	100387	1035 (9.50-3.66)
Sidechain outliers	100360	1005 (9.50-3.66)
RSRZ outliers	91569	1013 (9.50-3.66)

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 ≥ 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	601	 4% 34% 56% 6%
2	B	586	 25% 58% 11%
3	M	423	 3% 25% 52% 12% 9%
4	S	154	 30% 60% 5% 6%
5	C	172	 2% 38% 52% 6%

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 14743 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 AP-1 complex subunit gamma-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	579	4583	2882	806	857	38	0	0	0

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	596	GLU	-	EXPRESSION TAG	UNP P22892
A	597	ASN	-	EXPRESSION TAG	UNP P22892
A	598	LEU	-	EXPRESSION TAG	UNP P22892
A	599	TYR	-	EXPRESSION TAG	UNP P22892
A	600	PHE	-	EXPRESSION TAG	UNP P22892
A	601	GLN	-	EXPRESSION TAG	UNP P22892

- Molecule 2 is a protein called AP-1 complex subunit beta-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	563	4458	2847	732	852	27	0	0	0

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-1	GLY	-	EXPRESSION TAG	UNP Q10567
B	0	SER	-	EXPRESSION TAG	UNP Q10567
B	359	ARG	LYS	VARIANT	UNP Q10567
B	476	LYS	GLU	VARIANT	UNP Q10567
B	488	PHE	ILE	ENGINEERED MUTATION	UNP Q10567

- Molecule 3 is a protein called AP-1 complex subunit mu-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	M	383	3119	2013	523	569	14	0	0	0

- Molecule 4 is a protein called AP-1 complex subunit sigma-3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
4	S	145	1219	795	200	219	5	0	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
S	148	CYS	SER	VARIANT	UNP Q96PC3

- Molecule 5 is a protein called ADP-ribosylation factor 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
5	C	165	1331	842	233	250	6	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

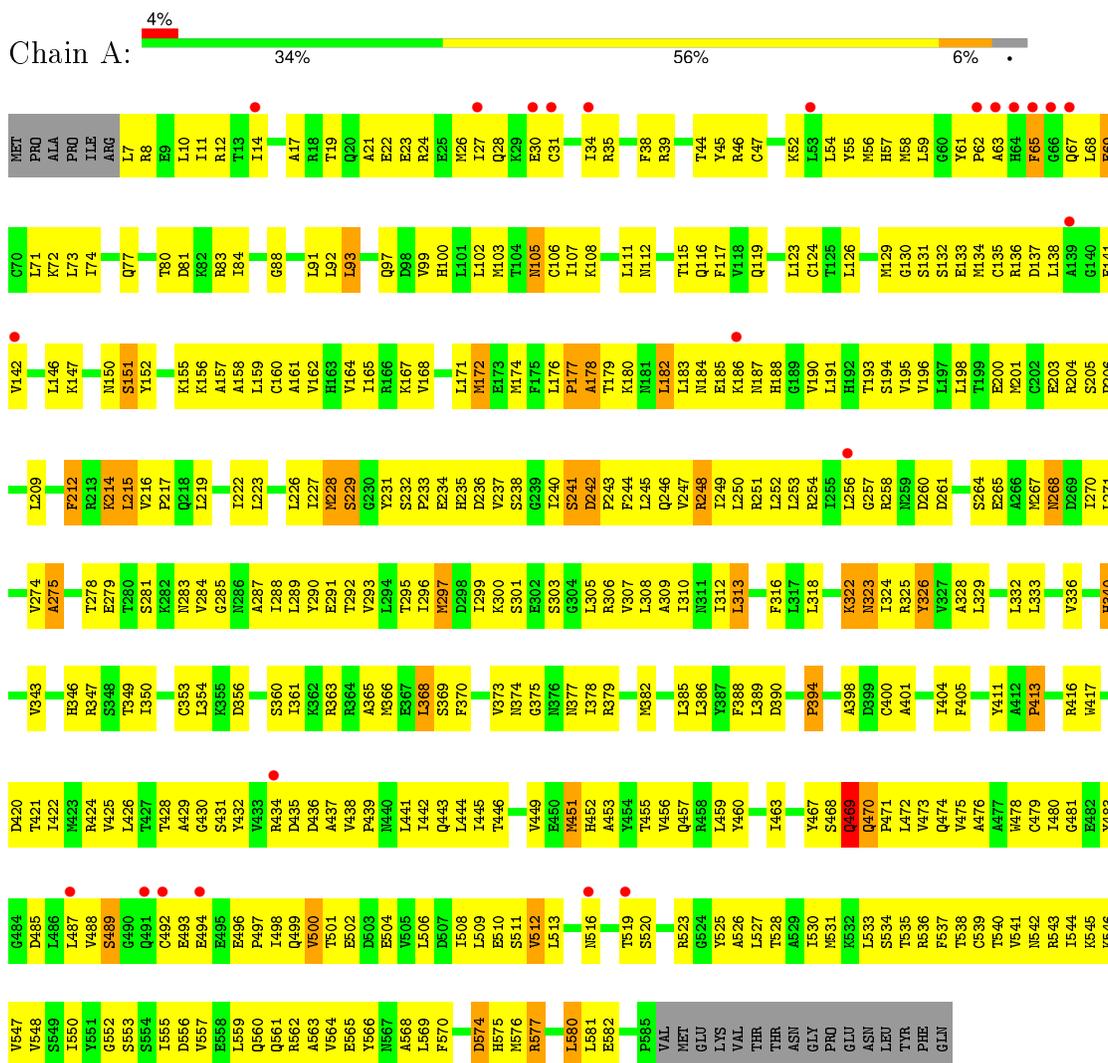
Chain	Residue	Modelled	Actual	Comment	Reference
C	10	MET	-	EXPRESSION TAG	UNP P84077
C	11	HIS	-	EXPRESSION TAG	UNP P84077
C	12	HIS	-	EXPRESSION TAG	UNP P84077
C	13	HIS	-	EXPRESSION TAG	UNP P84077
C	14	HIS	-	EXPRESSION TAG	UNP P84077
C	15	HIS	-	EXPRESSION TAG	UNP P84077
C	16	HIS	-	EXPRESSION TAG	UNP P84077
C	71	LEU	GLN	ENGINEERED MUTATION	UNP P84077

- Molecule 6 is GUANOSINE-5'-TRIPHOSPHATE (three-letter code: GTP) (formula: C₁₀H₁₆N₅O₁₄P₃).

3 Residue-property plots

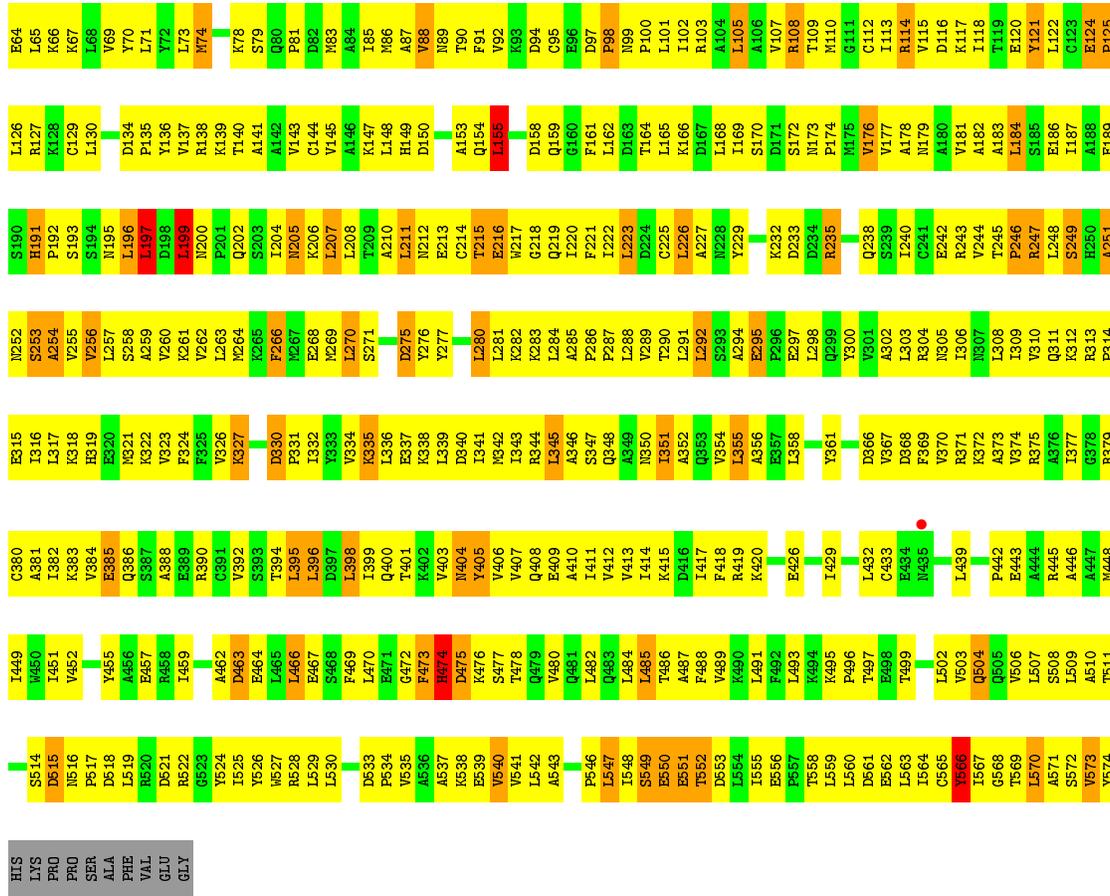
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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: AP-1 complex subunit gamma-1

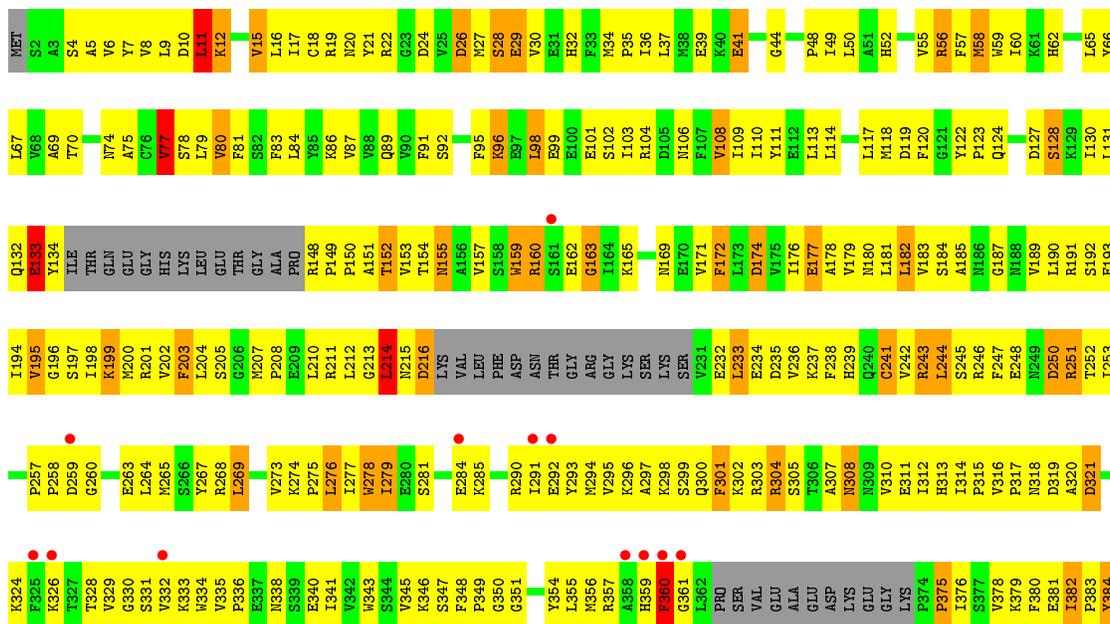


- Molecule 2: AP-1 complex subunit beta-1





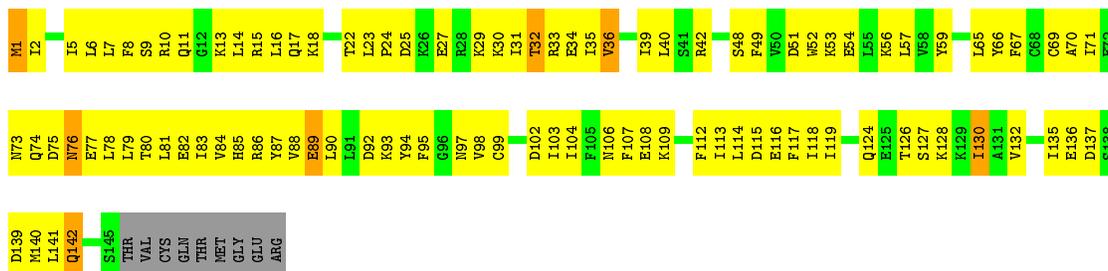
● Molecule 3: AP-1 complex subunit mu-1





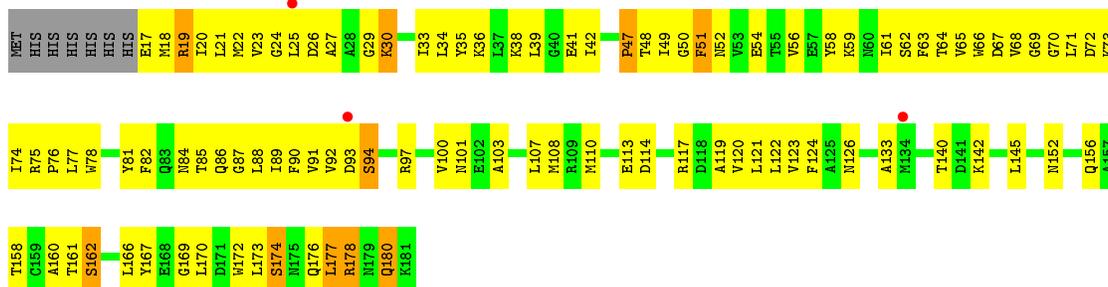
- Molecule 4: AP-1 complex subunit sigma-3

Chain S: 30% 60% 5% 6%



- Molecule 5: ADP-ribosylation factor 1

Chain C: 38% 52% 6% 2%



4 Data and refinement statistics

Property	Value	Source
Space group	P 64	Depositor
Cell constants a, b, c, α , β , γ	267.50Å 267.50Å 191.40Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 7.00 49.55 – 6.98	Depositor EDS
% Data completeness (in resolution range)	(Not available) (20.00-7.00) 93.6 (49.55-6.98)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.16	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.67 (at 6.68Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.200 , 0.250 0.203 , 0.245	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	357.1	Xtriage
Anisotropy	0.369	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 421.1	EDS
Estimated twinning fraction	0.046 for h,-h-k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 12305 reflections	Xtriage
F_o, F_c correlation	0.83	EDS
Total number of atoms	14743	wwPDB-VP
Average B, all atoms (Å ²)	321.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.11% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 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: GTP, MG

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	0.23	0/4646	0.50	0/6270
2	B	0.24	0/4525	0.54	0/6134
3	M	0.26	0/3191	0.53	0/4312
4	S	0.26	0/1242	0.47	0/1669
5	C	0.24	0/1354	0.47	0/1831
All	All	0.24	0/14958	0.51	0/20216

There are no bond length outliers.

There are no bond angle outliers.

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	4583	0	4700	439	0
2	B	4458	0	4593	589	0
3	M	3119	0	3129	382	0
4	S	1219	0	1249	133	0
5	C	1331	0	1328	109	0
6	C	32	0	12	3	0
7	C	1	0	0	0	0
All	All	14743	0	15011	1548	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 52.

The worst 5 of 1548 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:M:233:LEU:HD21	3:M:268:ARG:H	1.16	1.10
3:M:315:PRO:HA	3:M:340:GLU:HA	1.35	1.06
3:M:60:ILE:HD11	3:M:67:LEU:HD23	1.41	1.03
1:A:439:PRO:HG3	2:B:517:PRO:HD3	1.41	1.02
1:A:553:SER:HB3	2:B:546:PRO:HB2	1.48	0.95

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	577/601 (96%)	417 (72%)	122 (21%)	38 (7%)	1	24
2	B	561/586 (96%)	392 (70%)	125 (22%)	44 (8%)	1	20
3	M	375/423 (89%)	250 (67%)	94 (25%)	31 (8%)	1	18
4	S	143/154 (93%)	115 (80%)	23 (16%)	5 (4%)	4	39
5	C	163/172 (95%)	131 (80%)	24 (15%)	8 (5%)	3	31
All	All	1819/1936 (94%)	1305 (72%)	388 (21%)	126 (7%)	1	23

5 of 126 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	151	SER
1	A	184	ASN
1	A	214	LYS
1	A	215	LEU
1	A	229	SER

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	514/534 (96%)	496 (96%)	18 (4%)	43 74
2	B	503/523 (96%)	446 (89%)	57 (11%)	7 33
3	M	350/383 (91%)	312 (89%)	38 (11%)	8 35
4	S	135/143 (94%)	127 (94%)	8 (6%)	24 61
5	C	143/150 (95%)	138 (96%)	5 (4%)	43 74
All	All	1645/1733 (95%)	1519 (92%)	126 (8%)	16 52

5 of 126 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	355	LEU
2	B	551	GLU
4	S	74	GLN
2	B	385	GLU
2	B	466	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 48 such sidechains are listed below:

Mol	Chain	Res	Type
2	B	228	ASN
2	B	348	GLN
5	C	95	ASN
2	B	252	ASN
2	B	311	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 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 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
6	GTP	C	1001	7	25,34,34	1.13	2 (8%)	34,54,54	1.78	6 (17%)

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
6	GTP	C	1001	7	-	0/18/38/38	0/3/3/3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	C	1001	GTP	C2-N1	2.10	1.39	1.35
6	C	1001	GTP	C6-N1	3.94	1.40	1.33

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

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	C	1001	GTP	N3-C2-N1	-5.17	119.57	127.44
6	C	1001	GTP	C5-C6-N1	-3.79	118.41	123.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	C	1001	GTP	C1'-N9-C4	-3.46	121.73	126.94
6	C	1001	GTP	C6-C5-C4	-2.57	117.82	120.90
6	C	1001	GTP	C4-C5-N7	-2.27	107.39	109.48

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	C	1001	GTP	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, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ > 2	OWAB(Å ²)	Q < 0.9
1	A	579/601 (96%)	0.47	23 (3%) 42 40	356, 356, 356, 356	0
2	B	563/586 (96%)	-0.03	2 (0%) 93 90	293, 293, 293, 293	0
3	M	383/423 (90%)	0.21	12 (3%) 52 48	285, 320, 320, 320	0
4	S	145/154 (94%)	0.12	0 100 100	352, 352, 352, 352	0
5	C	165/172 (95%)	0.41	3 (1%) 71 66	299, 299, 299, 299	0
All	All	1835/1936 (94%)	0.23	40 (2%) 65 61	285, 320, 356, 356	0

The worst 5 of 40 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	62	PRO	4.5
1	A	63	ALA	3.9
1	A	494	GLU	3.4
1	A	66	GLY	3.4
1	A	491	GLN	3.3

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, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
6	GTP	C	1001	32/32	0.89	0.25	-0.87	299,299,299,299	0
7	MG	C	1002	1/1	0.99	0.18	-1.15	299,299,299,299	0

6.5 Other polymers [i](#)

There are no such residues in this entry.