



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:37 PM GMT

PDB ID : 4M63
Title : Crystal Structure of a Filament-Like Actin Trimer Bound to the Bacterial Effector VopL
Authors : Tomchick, D.R.; Zahm, J.A.; Rosen, M.K.
Deposited on : 2013-08-08
Resolution : 2.75 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
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

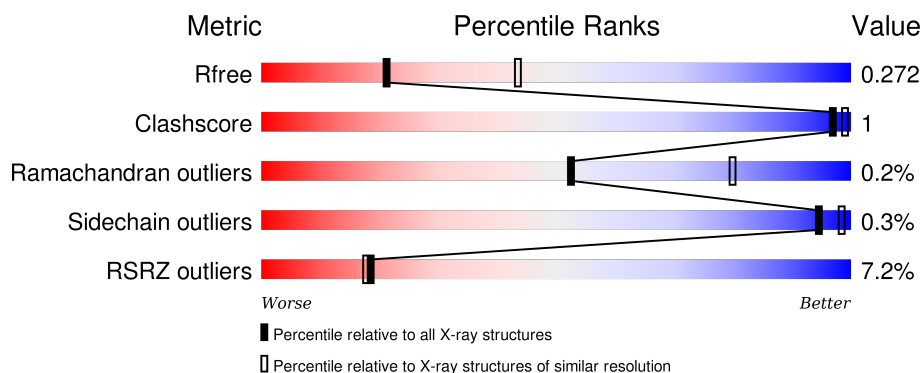
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.75 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	3050 (2.78-2.70)
Clashscore	102246	3424 (2.78-2.70)
Ramachandran outliers	100387	3367 (2.78-2.70)
Sidechain outliers	100360	3368 (2.78-2.70)
RSRZ outliers	91569	3055 (2.78-2.70)

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	241	<div> <div>12%</div> <div>89%</div> <div>7%</div> </div>
1	B	241	<div> <div>5%</div> <div>87%</div> <div>5%</div> <div>8%</div> </div>
2	C	377	<div> <div>7%</div> <div>92%</div> <div>•</div> <div>•</div> </div>
2	D	377	<div> <div>5%</div> <div>92%</div> <div>•</div> <div>•</div> </div>
2	E	377	<div> <div>6%</div> <div>90%</div> <div>•</div> <div>7%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	CA	C	402	-	-	-	X
4	CA	D	402	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 23742 atoms, of which 11782 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 T3SS2 effector VopL nucleation of actin polymerization.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	224	Total	C	H	N	O	S	0	0	0
			3508	1108	1737	302	353	8			
1	B	222	Total	C	H	N	O	S	0	0	0
			3478	1099	1720	300	351	8			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	244	GLY	-	EXPRESSION TAG	UNP L0I5A1
A	245	HIS	-	EXPRESSION TAG	UNP L0I5A1
A	246	MET	-	EXPRESSION TAG	UNP L0I5A1
B	244	GLY	-	EXPRESSION TAG	UNP L0I5A1
B	245	HIS	-	EXPRESSION TAG	UNP L0I5A1
B	246	MET	-	EXPRESSION TAG	UNP L0I5A1

- Molecule 2 is a protein called Actin-5C.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	C	361	Total	C	H	N	O	S	0	0	0
			5596	1779	2789	475	535	18			
2	D	361	Total	C	H	N	O	S	0	0	0
			5612	1783	2801	475	534	19			
2	E	350	Total	C	H	N	O	S	0	0	0
			5416	1723	2699	457	518	19			

There are 12 discrepancies between the modelled and reference sequences:

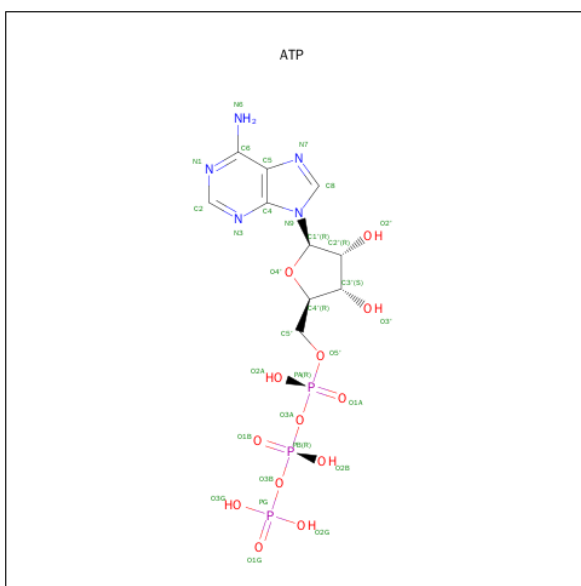
Chain	Residue	Modelled	Actual	Comment	Reference
C	-1	GLY	-	EXPRESSION TAG	UNP P10987
C	286	ALA	ASP	ENGINEERED MUTATION	UNP P10987
C	287	ALA	VAL	ENGINEERED MUTATION	UNP P10987
C	288	ALA	ASP	ENGINEERED MUTATION	UNP P10987

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-1	GLY	-	EXPRESSION TAG	UNP P10987
D	286	ALA	ASP	ENGINEERED MUTATION	UNP P10987
D	287	ALA	VAL	ENGINEERED MUTATION	UNP P10987
D	288	ALA	ASP	ENGINEERED MUTATION	UNP P10987
E	-1	GLY	-	EXPRESSION TAG	UNP P10987
E	286	ALA	ASP	EXPRESSION TAG	UNP P10987
E	287	ALA	VAL	ENGINEERED MUTATION	UNP P10987
E	288	ALA	ASP	ENGINEERED MUTATION	UNP P10987

- Molecule 3 is ADENOSINE-5'-TRIPHOSPHATE (three-letter code: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
3	C	1	Total 43	C 10	H 12	N 5	O 13	P 3	0	0
3	D	1	Total 43	C 10	H 12	N 5	O 13	P 3	0	0
3	E	1	Total 43	C 10	H 12	N 5	O 13	P 3	0	0

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	1	Total	Ca	0	0
			1	1		

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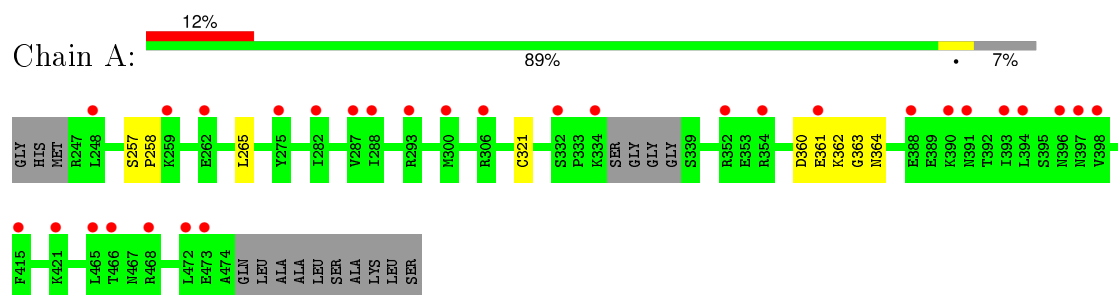
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	1	Total 1	Ca 1	0	0
4	E	1	Total 1	Ca 1	0	0

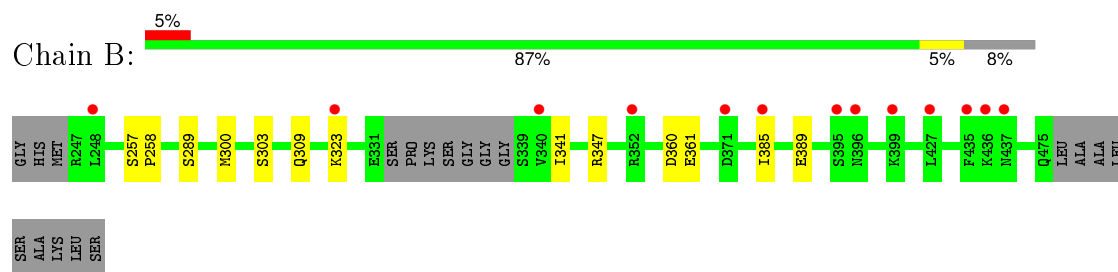
3 Residue-property plots [i](#)

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.

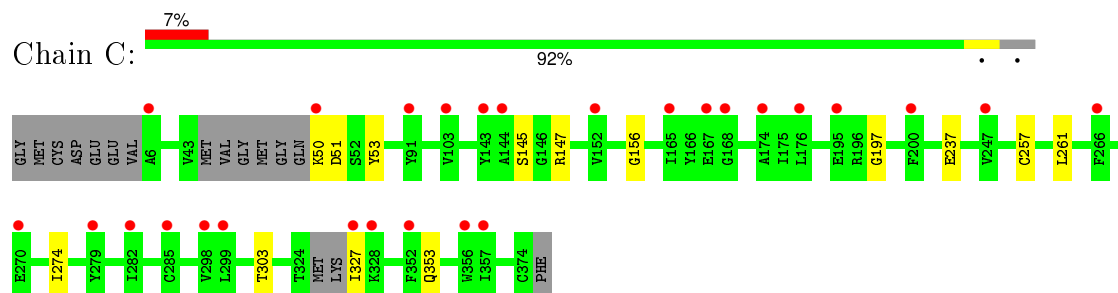
- Molecule 1: T3SS2 effector VopL nucleation of actin polymerization



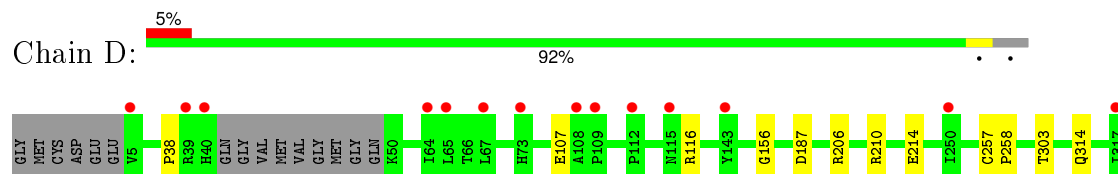
- Molecule 1: T3SS2 effector VopL nucleation of actin polymerization

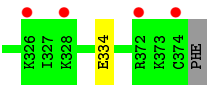


- Molecule 2: Actin-5C

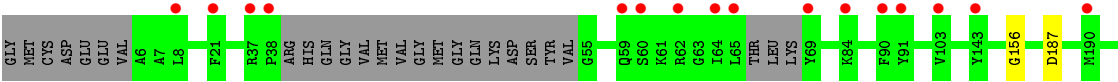
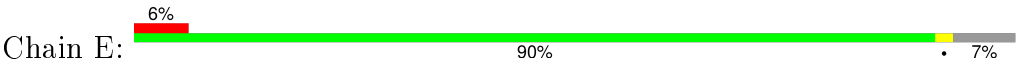


- Molecule 2: Actin-5C





● Molecule 2: Actin-5C



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	287.76Å 66.08Å 103.16Å 90.00° 94.01° 90.00°	Depositor
Resolution (Å)	44.59 – 2.75 44.59 – 2.75	Depositor EDS
% Data completeness (in resolution range)	98.1 (44.59-2.75) 98.2 (44.59-2.75)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.95 (at 2.77Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, R_{free}	0.219 , 0.272 0.219 , 0.272	Depositor DCC
R_{free} test set	2481 reflections (4.97%)	DCC
Wilson B-factor (Å ²)	75.3	Xtriage
Anisotropy	0.543	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 44.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	1 of 49988 reflections (0.002%)	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	23742	wwPDB-VP
Average B, all atoms (Å ²)	101.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.62% 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: CA, 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	0.22	0/1798	0.36	0/2419
1	B	0.22	0/1784	0.37	0/2400
2	C	0.21	0/2867	0.38	0/3885
2	D	0.21	0/2872	0.37	0/3892
2	E	0.21	0/2775	0.38	0/3760
All	All	0.21	0/12096	0.37	0/16356

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	1771	1737	1735	3	0
1	B	1758	1720	1718	8	0
2	C	2807	2789	2780	7	0
2	D	2811	2801	2792	7	0
2	E	2717	2699	2691	5	0
3	C	31	12	12	0	0
3	D	31	12	12	0	0
3	E	31	12	12	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	C	1	0	0	0	0
4	D	1	0	0	0	0
4	E	1	0	0	0	0
All	All	11960	11782	11752	30	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 1.

All (30) 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)
2:C:353:GLN:N	2:C:353:GLN:OE1	2.33	0.62
2:C:50:LYS:N	2:C:53:TYR:HH	1.99	0.59
2:C:145:SER:O	2:C:147:ARG:NH1	2.36	0.58
2:E:210:ARG:NH1	2:E:214:GLU:OE2	2.39	0.56
1:B:360:ASP:OD1	1:B:361:GLU:N	2.41	0.53
2:D:210:ARG:NH1	2:D:214:GLU:OE2	2.41	0.53
2:C:156:GLY:O	2:C:303:THR:OG1	2.25	0.52
2:C:237:GLU:N	2:C:237:GLU:OE1	2.43	0.51
2:D:314:GLN:OE1	2:D:314:GLN:N	2.43	0.51
2:E:156:GLY:O	2:E:303:THR:OG1	2.25	0.50
2:E:211:ASP:OD2	2:E:240:TYR:OH	2.30	0.49
2:D:156:GLY:O	2:D:303:THR:OG1	2.28	0.49
2:D:334:GLU:N	2:D:334:GLU:OE1	2.46	0.48
1:A:360:ASP:OD1	1:A:361:GLU:N	2.42	0.47
2:E:187:ASP:OD1	2:E:206:ARG:NH1	2.45	0.47
2:D:187:ASP:OD1	2:D:206:ARG:NH2	2.48	0.47
1:B:323:LYS:O	1:B:347:ARG:NH2	2.47	0.46
1:A:362:LYS:O	1:A:364:ASN:N	2.48	0.46
2:E:334:GLU:OE1	2:E:334:GLU:N	2.47	0.46
1:B:385:ILE:HG23	1:B:389:GLU:HB2	1.99	0.45
2:D:107:GLU:OE2	2:D:116:ARG:NH1	2.49	0.45
2:C:50:LYS:NZ	2:C:51:ASP:OD1	2.46	0.44
1:B:289:SER:HA	1:B:341:ILE:CD1	2.47	0.44
1:B:303:SER:O	1:B:309:GLN:NE2	2.47	0.43
1:B:289:SER:HA	1:B:341:ILE:HD11	2.01	0.43
2:D:257:CYS:HB3	2:D:258:PRO:HD3	2.00	0.43
1:B:257:SER:N	1:B:258:PRO:CD	2.83	0.42
1:A:257:SER:N	1:A:258:PRO:CD	2.82	0.42
1:B:300:MET:HE1	1:B:341:ILE:HG23	2.03	0.41
2:C:261:LEU:HG	2:C:274:ILE:HD11	2.03	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	220/241 (91%)	210 (96%)	9 (4%)	1 (0%)	34	62
1	B	218/241 (90%)	209 (96%)	9 (4%)	0	100	100
2	C	355/377 (94%)	339 (96%)	15 (4%)	1 (0%)	46	74
2	D	357/377 (95%)	346 (97%)	10 (3%)	1 (0%)	46	74
2	E	344/377 (91%)	336 (98%)	8 (2%)	0	100	100
All	All	1494/1613 (93%)	1440 (96%)	51 (3%)	3 (0%)	52	80

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	363	GLY
2	D	38	PRO
2	C	197	GLY

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	190/200 (95%)	188 (99%)	2 (1%)	80	93
1	B	188/200 (94%)	188 (100%)	0	100	100
2	C	302/315 (96%)	300 (99%)	2 (1%)	88	96

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	D	303/315 (96%)	303 (100%)	0	100	100
2	E	292/315 (93%)	292 (100%)	0	100	100
All	All	1275/1345 (95%)	1271 (100%)	4 (0%)	94	98

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	265	LEU
1	A	321	CYS
2	C	257	CYS
2	C	327	ILE

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

Mol	Chain	Res	Type
2	E	87	HIS

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 6 ligands modelled in this entry, 3 are monoatomic - leaving 3 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	$\# Z > 2$
3	ATP	C	401	4	24,33,33	0.96	1 (4%)	31,52,52	1.91	4 (12%)
3	ATP	D	401	4	24,33,33	0.98	1 (4%)	31,52,52	1.97	4 (12%)
3	ATP	E	401	4	24,33,33	0.95	1 (4%)	31,52,52	1.88	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
3	ATP	C	401	4	-	0/18/38/38	0/3/3/3
3	ATP	D	401	4	-	0/18/38/38	0/3/3/3
3	ATP	E	401	4	-	0/18/38/38	0/3/3/3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	401	ATP	C5-C4	2.96	1.47	1.40
3	C	401	ATP	C5-C4	3.05	1.47	1.40
3	D	401	ATP	C5-C4	3.08	1.47	1.40

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	401	ATP	N3-C2-N1	-7.50	123.15	128.89
3	E	401	ATP	N3-C2-N1	-7.47	123.18	128.89
3	C	401	ATP	N3-C2-N1	-7.39	123.24	128.89
3	C	401	ATP	PA-O3A-PB	-3.99	121.53	132.73
3	D	401	ATP	PA-O3A-PB	-3.74	122.23	132.73
3	E	401	ATP	PA-O3A-PB	-3.58	122.68	132.73
3	D	401	ATP	PB-O3B-PG	-3.50	120.92	132.67
3	E	401	ATP	C4-C5-N7	-3.40	106.35	109.48
3	C	401	ATP	PB-O3B-PG	-3.39	121.31	132.67
3	D	401	ATP	C4-C5-N7	-3.24	106.50	109.48
3	C	401	ATP	C4-C5-N7	-3.18	106.55	109.48
3	E	401	ATP	PB-O3B-PG	-2.91	122.92	132.67

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

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, 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	224/241 (92%)	0.75	30 (13%) 4 4	53, 103, 178, 231	0
1	B	222/241 (92%)	0.48	13 (5%) 26 26	48, 92, 149, 185	0
2	C	361/377 (95%)	0.59	27 (7%) 17 16	46, 89, 145, 188	0
2	D	361/377 (95%)	0.52	18 (4%) 32 32	47, 81, 125, 186	0
2	E	350/377 (92%)	0.46	22 (6%) 23 23	45, 75, 140, 183	0
All	All	1518/1613 (94%)	0.55	110 (7%) 18 18	45, 87, 148, 231	0

All (110) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	5	VAL	6.3
2	E	38	PRO	5.5
2	C	282	ILE	5.3
2	E	65	LEU	4.9
1	A	275	TYR	4.7
2	D	374	CYS	4.5
1	A	468	ARG	4.5
2	E	37	ARG	4.3
2	E	201	THR	4.2
1	A	334	LYS	4.2
1	A	472	LEU	4.2
1	B	352	ARG	3.8
2	E	64	ILE	3.8
2	C	328	LYS	3.8
2	E	60	SER	3.7
2	E	62	ARG	3.6
2	C	144	ALA	3.5
1	A	332	SER	3.5
2	E	59	GLN	3.5
1	A	397	ASN	3.4

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Mol	Chain	Res	Type	RSRZ
2	C	327	ILE	3.4
1	A	354	ARG	3.3
1	A	465	LEU	3.3
1	A	394	LEU	3.3
2	C	103	VAL	3.3
1	B	436	LYS	3.2
1	A	391	ASN	3.2
2	D	67	LEU	3.2
2	C	200	PHE	3.2
2	D	328	LYS	3.1
1	A	396	ASN	3.1
2	C	143	TYR	3.1
2	E	202	THR	3.0
2	E	8	LEU	3.0
1	A	288	ILE	3.0
1	A	282	ILE	2.9
1	B	427	LEU	2.9
1	A	361	GLU	2.9
1	A	473	GLU	2.9
1	A	287	VAL	2.9
2	C	279	TYR	2.9
2	E	197	GLY	2.9
2	C	298	VAL	2.8
1	A	421	LYS	2.8
2	C	168	GLY	2.8
2	C	285	CYS	2.7
2	C	352	PHE	2.7
2	D	73	HIS	2.7
2	D	250	ILE	2.7
1	B	435	PHE	2.7
2	D	40	HIS	2.6
2	E	199	SER	2.6
2	E	69	TYR	2.6
1	A	393	ILE	2.6
2	E	370	VAL	2.6
1	B	323	LYS	2.6
2	C	6	ALA	2.5
2	D	39	ARG	2.5
2	C	167	GLU	2.5
2	D	64	ILE	2.5
1	A	262	GLU	2.5
2	C	174	ALA	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	300	MET	2.5
2	D	317	ILE	2.4
2	E	190	MET	2.4
2	C	299	LEU	2.4
2	C	357	ILE	2.4
2	D	372	ARG	2.4
1	A	293	ARG	2.4
1	A	306	ARG	2.4
2	C	91	TYR	2.4
1	A	388	GLU	2.4
2	C	50	LYS	2.3
1	B	385	ILE	2.3
1	A	398	VAL	2.3
1	B	340	VAL	2.3
2	E	84	LYS	2.3
2	C	247	VAL	2.3
2	D	108	ALA	2.3
2	E	103	VAL	2.3
1	A	248	LEU	2.3
1	A	415	PHE	2.3
1	A	259	LYS	2.3
1	B	395	SER	2.3
2	C	165	ILE	2.3
2	E	143	TYR	2.3
2	E	21	PHE	2.2
2	D	326	LYS	2.2
1	B	399	LYS	2.2
1	B	371	ASP	2.2
2	C	270	GLU	2.2
2	C	356	TRP	2.1
1	B	248	LEU	2.1
2	C	152	VAL	2.1
2	C	176	LEU	2.1
2	E	91	TYR	2.1
1	B	437	ASN	2.1
2	D	115	ASN	2.1
2	E	90	PHE	2.1
2	C	195	GLU	2.1
1	B	396	ASN	2.1
1	A	390	LYS	2.1
2	D	112	PRO	2.0
2	D	143	TYR	2.0

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Mol	Chain	Res	Type	RSRZ
2	D	65	LEU	2.0
1	A	352	ARG	2.0
1	A	466	THR	2.0
2	E	194	THR	2.0
2	C	266	PHE	2.0
2	D	109	PRO	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 ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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
4	CA	D	402	1/1	0.96	0.37	8.52	74,74,74,74	0
4	CA	C	402	1/1	0.97	0.42	7.65	89,89,89,89	0
4	CA	E	402	1/1	0.94	0.22	1.26	69,69,69,69	0
3	ATP	E	401	31/31	0.95	0.16	-0.64	38,54,78,98	0
3	ATP	C	401	31/31	0.94	0.15	-0.81	54,69,82,89	0
3	ATP	D	401	31/31	0.95	0.15	-1.45	40,61,71,79	0

6.5 Other polymers ⓘ

There are no such residues in this entry.