



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 19, 2020 – 12:43 AM EST

PDB ID : 6X1G
Title : Crystal structure of a GEF domain from the Orientia tsutsugamushi protein
OtDUB in complex with Rac1
Authors : Lim, C.S.; Xiong, Y.
Deposited on : 2020-05-18
Resolution : 1.60 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

<https://www.wwpdb.org/validation/2017/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
Xtriage (Phenix)	:	1.13
EDS	:	2.14.6
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)	:	2.14.6

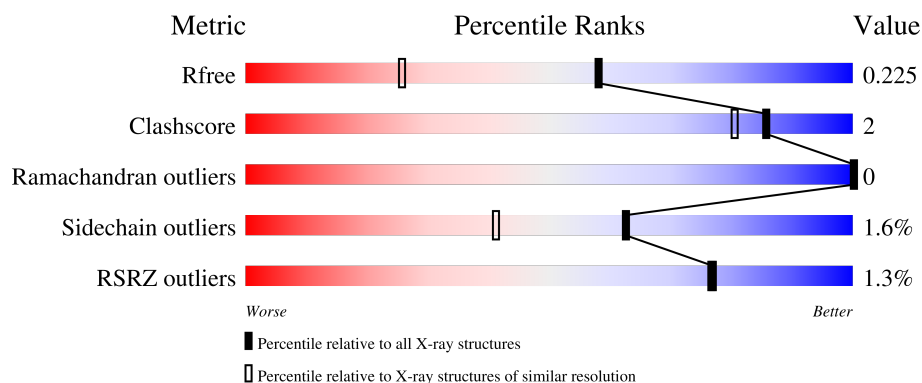
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.60 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	3398 (1.60-1.60)
Clashscore	141614	3665 (1.60-1.60)
Ramachandran outliers	138981	3564 (1.60-1.60)
Sidechain outliers	138945	3563 (1.60-1.60)
RSRZ outliers	127900	3321 (1.60-1.60)

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 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	219	<div> <div>2%</div> <div>90% 9%</div> </div>
1	C	219	<div> <div>94% 5%</div> </div>
2	B	183	<div> <div>2%</div> <div>86% 10%</div> </div>
2	D	183	<div> <div>2%</div> <div>90% 7%</div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6578 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 ULP_PROTEASE domain-containing protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	216	Total	C	N	O	S	0	0	0
			1740	1095	299	341	5			
1	C	216	Total	C	N	O	S	0	0	0
			1740	1095	299	341	5			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	547	MET	-	initiating methionine	UNP B3CVM3
A	760	HIS	-	expression tag	UNP B3CVM3
A	761	HIS	-	expression tag	UNP B3CVM3
A	762	HIS	-	expression tag	UNP B3CVM3
A	763	HIS	-	expression tag	UNP B3CVM3
A	764	HIS	-	expression tag	UNP B3CVM3
A	765	HIS	-	expression tag	UNP B3CVM3
C	547	MET	-	initiating methionine	UNP B3CVM3
C	760	HIS	-	expression tag	UNP B3CVM3
C	761	HIS	-	expression tag	UNP B3CVM3
C	762	HIS	-	expression tag	UNP B3CVM3
C	763	HIS	-	expression tag	UNP B3CVM3
C	764	HIS	-	expression tag	UNP B3CVM3
C	765	HIS	-	expression tag	UNP B3CVM3

- Molecule 2 is a protein called Ras-related C3 botulinum toxin substrate 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	178	Total	C	N	O	S	0	0	0
			1393	895	231	259	8			
2	D	178	Total	C	N	O	S	0	0	0
			1393	895	231	259	8			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	178	HIS	-	expression tag	UNP P63000
B	179	HIS	-	expression tag	UNP P63000
B	180	HIS	-	expression tag	UNP P63000
B	181	HIS	-	expression tag	UNP P63000
B	182	HIS	-	expression tag	UNP P63000
B	183	HIS	-	expression tag	UNP P63000
D	178	HIS	-	expression tag	UNP P63000
D	179	HIS	-	expression tag	UNP P63000
D	180	HIS	-	expression tag	UNP P63000
D	181	HIS	-	expression tag	UNP P63000
D	182	HIS	-	expression tag	UNP P63000
D	183	HIS	-	expression tag	UNP P63000

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	C	1	Total Mg 1 1	0	0

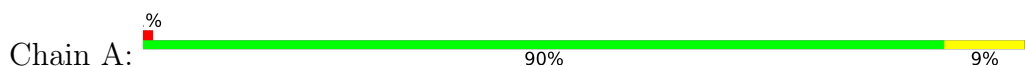
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	71	Total O 71 71	0	0
4	C	88	Total O 88 88	0	0
4	B	74	Total O 74 74	0	0
4	D	78	Total O 78 78	0	0

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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.

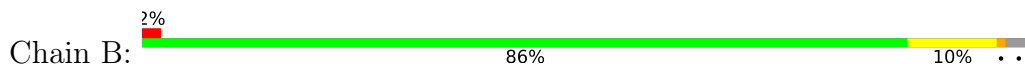
- Molecule 1: ULP_PROTEASE domain-containing protein



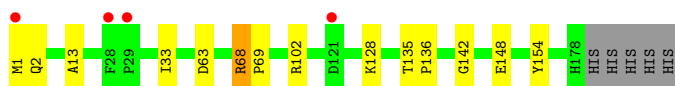
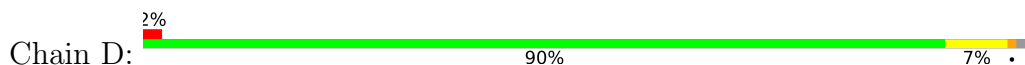
- Molecule 1: ULP_PROTEASE domain-containing protein



- Molecule 2: Ras-related C3 botulinum toxin substrate 1



- Molecule 2: Ras-related C3 botulinum toxin substrate 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	50.06Å 54.16Å 94.00Å 83.37° 76.34° 62.52°	Depositor
Resolution (Å)	48.05 – 1.60 48.05 – 1.60	Depositor EDS
% Data completeness (in resolution range)	81.5 (48.05-1.60) 81.5 (48.05-1.60)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.42 (at 1.60Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.181 , 0.218 0.190 , 0.225	Depositor DCC
R_{free} test set	4663 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	26.5	Xtriage
Anisotropy	0.085	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 45.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.146 for h,h-k,h-l 0.010 for -h,-h+k,-l 0.015 for -h,-k,-h+l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	6578	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.41% 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, 0.333 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: 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	0.85	1/1766 (0.1%)	0.87	0/2371
1	C	0.87	0/1766	0.87	0/2371
2	B	0.85	1/1424 (0.1%)	1.01	2/1937 (0.1%)
2	D	0.86	2/1424 (0.1%)	1.00	2/1937 (0.1%)
All	All	0.86	4/6380 (0.1%)	0.93	4/8616 (0.0%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	148	GLU	CD-OE1	-5.90	1.19	1.25
2	B	148	GLU	CD-OE1	-5.71	1.19	1.25
1	A	611	GLU	CD-OE2	-5.18	1.20	1.25
2	D	102	ARG	C-O	5.14	1.33	1.23

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	68	ARG	NE-CZ-NH1	9.95	125.27	120.30
2	B	68	ARG	NE-CZ-NH2	-9.28	115.66	120.30
2	D	68	ARG	NE-CZ-NH1	7.81	124.20	120.30
2	D	68	ARG	NE-CZ-NH2	-6.99	116.80	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	1740	0	1744	10	0
1	C	1740	0	1744	5	0
2	B	1393	0	1412	11	0
2	D	1393	0	1412	4	0
3	C	1	0	0	0	0
4	A	71	0	0	1	0
4	B	74	0	0	0	0
4	C	88	0	0	2	0
4	D	78	0	0	0	0
All	All	6578	0	6312	27	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 2.

All (27) 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:B:133:LYS:O	2:B:133:LYS:HG2	1.92	0.70
2:B:121:ASP:HA	2:B:126:ILE:HD11	1.81	0.62
1:A:598:LEU:O	1:A:608:SER:HA	2.09	0.52
1:C:681:GLU:O	1:C:685:GLU:HG2	2.09	0.52
2:B:132:LYS:O	2:B:133:LYS:HB3	2.10	0.52
1:A:676:LEU:HD13	2:B:45:MET:HG2	1.91	0.51
1:A:755:ILE:O	1:A:759:MET:HG2	2.11	0.51
1:A:680:SER:HB2	4:A:870:HOH:O	2.10	0.50
2:B:4:ILE:HG12	2:B:176:VAL:HG21	1.95	0.48
1:C:598:LEU:O	1:C:608:SER:HA	2.14	0.48
1:C:680:SER:HB2	4:C:983:HOH:O	2.12	0.48
1:C:755:ILE:O	1:C:759:MET:HG2	2.14	0.48
2:B:142:GLY:HA3	2:B:154:TYR:CZ	2.51	0.46
2:B:135:THR:HA	2:B:136:PRO:HD3	1.89	0.45
1:A:681:GLU:O	1:A:685:GLU:HG3	2.17	0.45
2:D:142:GLY:HA3	2:D:154:TYR:CZ	2.52	0.45
2:D:68:ARG:N	2:D:69:PRO:CD	2.79	0.45
1:A:666:LYS:HE3	1:A:691:VAL:HB	1.99	0.45
1:A:721:PHE:CE1	1:A:746:ALA:HB1	2.53	0.43
1:A:674:PRO:HD3	1:A:715:TYR:CD2	2.55	0.42
2:B:72:TYR:N	2:B:73:PRO:CD	2.83	0.42
1:A:707:HIS:NE2	2:B:3:ALA:HB1	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:671:ARG:HD3	2:B:40:TYR:HB3	2.01	0.41
2:B:68:ARG:N	2:B:69:PRO:CD	2.83	0.41
4:C:961:HOH:O	2:D:13:ALA:HB2	2.20	0.40
1:C:666:LYS:HE3	1:C:691:VAL:HB	2.03	0.40
2:D:135:THR:HA	2:D:136:PRO:HD3	1.96	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	214/219 (98%)	212 (99%)	2 (1%)	0	100	100
1	C	214/219 (98%)	212 (99%)	2 (1%)	0	100	100
2	B	176/183 (96%)	174 (99%)	2 (1%)	0	100	100
2	D	176/183 (96%)	173 (98%)	3 (2%)	0	100	100
All	All	780/804 (97%)	771 (99%)	9 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	196/199 (98%)	194 (99%)	2 (1%)	76	61

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	196/199 (98%)	195 (100%)	1 (0%)	88	80
2	B	154/159 (97%)	151 (98%)	3 (2%)	57	34
2	D	154/159 (97%)	149 (97%)	5 (3%)	39	15
All	All	700/716 (98%)	689 (98%)	11 (2%)	62	41

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	548	GLU
1	A	629	SER
1	C	583	ARG
2	B	45	MET
2	B	63	ASP
2	B	140	PRO
2	D	1	MET
2	D	2	GLN
2	D	33	ILE
2	D	63	ASP
2	D	128	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	592	GLN
1	C	592	GLN
1	C	749	GLN
2	B	2	GLN
2	B	103	HIS
2	D	103	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

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 [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	216/219 (98%)	-0.22	2 (0%) 84 84	24, 47, 82, 115	0
1	C	216/219 (98%)	-0.25	1 (0%) 91 90	22, 45, 74, 109	0
2	B	178/183 (97%)	-0.30	3 (1%) 70 69	25, 35, 81, 94	0
2	D	178/183 (97%)	-0.28	4 (2%) 62 60	24, 35, 82, 95	0
All	All	788/804 (98%)	-0.26	10 (1%) 77 77	22, 40, 80, 115	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	547	MET	13.1
2	D	28	PHE	5.1
1	A	547	MET	4.6
2	B	28	PHE	4.1
2	D	121	ASP	2.6
2	D	1	MET	2.5
2	D	29	PRO	2.3
2	B	13	ALA	2.2
2	B	46	VAL	2.1
1	A	548	GLU	2.1

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 monosaccharides 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, 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	B-factors(Å ²)	Q<0.9
3	MG	C	801	1/1	0.75	0.09	66,66,66,66	0

6.5 Other polymers [i](#)

There are no such residues in this entry.