wwPDB X-ray Structure Validation Summary Report (i)

Aug 9, 2016 – 01:03 AM EDT

PDB ID	:	5EYA
Title	:	TRIM25 RING domain in complex with Ubc13-Ub conjugate
Authors	:	Pornillos, O.; Sanchez, J.G.
Deposited on	:	2015-11-24
Resolution	:	2.40 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

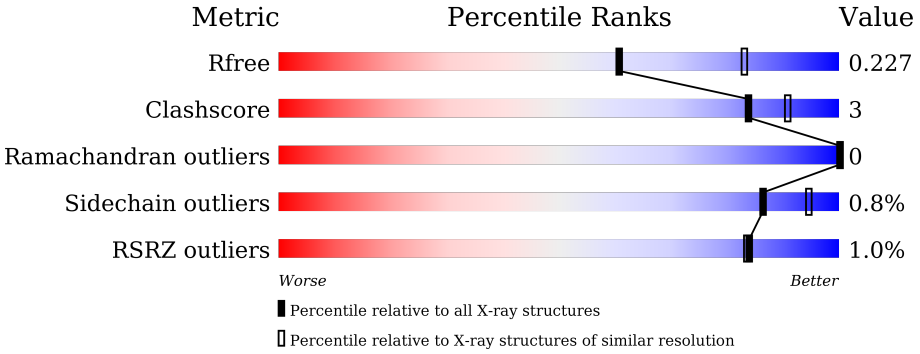
MolProbity	:	4.02b-467
Mogul	:	unknown
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20027939
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20027939

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	2919 (2.40-2.40)
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (2.40-2.40)

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 <=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	152	.% 93% 5% •
1	B	152	93% 5% •
2	F	86	2% 78% 10% 12%
2	G	86	.% 88% 12%
3	C	76	.% 91% 9%
3	D	76	97% •

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 9882 atoms, of which 4864 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 Ubiquitin-conjugating enzyme E2 N.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	149	2391	764	1202	206	216	3	0	0	0
1	B	149	2391	764	1202	206	216	3	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	87	LYS	CYS	engineered mutation	UNP P61088
B	87	LYS	CYS	engineered mutation	UNP P61088

- Molecule 2 is a protein called Tripartite motif-containing 25 variant.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
2	F	76	1165	378	570	99	110	8	0	1	0
2	G	86	1291	412	632	114	123	10	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
F	0	HIS	ALA	conflict	UNP Q59GW5
G	0	HIS	ALA	conflict	UNP Q59GW5

- Molecule 3 is a protein called Polyubiquitin-B.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
3	C	76	1230	378	629	105	117	1	0	0	0
3	D	76	1230	378	629	105	117	1	0	0	0

- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Zn		
4	G	2	2	2	0	0
4	F	2	2	2	0	0

- Molecule 5 is water.

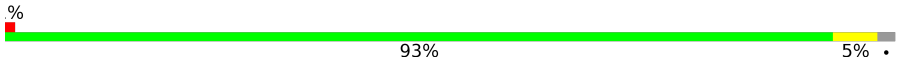
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	A	41	41	41	0	0
5	F	27	27	27	0	0
5	G	32	32	32	0	0
5	B	36	36	36	0	0
5	C	23	23	23	0	0
5	D	21	21	21	0	0

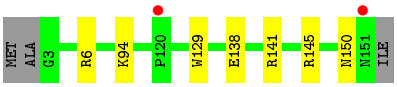
3 Residue-property plots (i)

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: Ubiquitin-conjugating enzyme E2 N

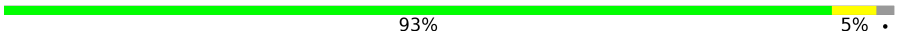
Chain A:

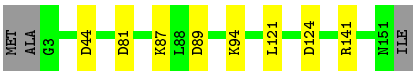




- Molecule 1: Ubiquitin-conjugating enzyme E2 N

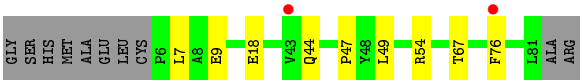
Chain B:





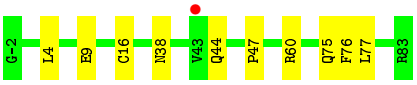
- Molecule 2: Tripartite motif-containing 25 variant

Chain F:



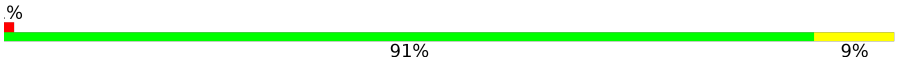
- Molecule 2: Tripartite motif-containing 25 variant

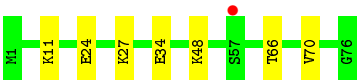
Chain G:



- Molecule 3: Polyubiquitin-B

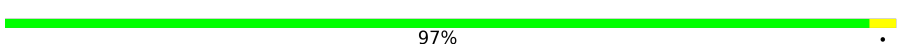
Chain C:

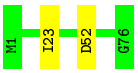




- Molecule 3: Polyubiquitin-B

Chain D:




M1
I23
D52
G76

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	53.05Å 75.78Å 169.10Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	28.97 – 2.40 45.23 – 2.21	Depositor EDS
% Data completeness (in resolution range)	89.7 (28.97-2.40) 79.0 (45.23-2.21)	Depositor EDS
R_{merge}	0.18	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.49 (at 2.22Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.189 , 0.228 0.192 , 0.227	Depositor DCC
R_{free} test set	1786 reflections (7.25%)	DCC
Wilson B-factor (Å^2)	26.8	Xtriage
Anisotropy	0.530	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.39 , 38.9	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	9882	wwPDB-VP
Average B, all atoms (Å^2)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 30.78 % of the origin peak, indicating pseudo translational symmetry.The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 1.2341e-03.The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: ZN

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	0.33	0/1219	0.49	0/1659
1	B	0.32	0/1219	0.50	0/1659
2	F	0.38	0/614	0.56	0/838
2	G	0.34	0/675	0.57	0/920
3	C	0.31	0/607	0.51	0/816
3	D	0.30	0/607	0.53	0/816
All	All	0.33	0/4941	0.52	0/6708

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	1189	1202	1202	6	0
1	B	1189	1202	1202	7	0
2	F	595	570	570	6	0
2	G	659	632	632	5	0
3	C	601	629	629	6	0
3	D	601	629	629	1	0
4	F	2	0	0	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	G	2	0	0	0	0
5	A	41	0	0	3	0
5	B	36	0	0	3	0
5	C	23	0	0	4	1
5	D	21	0	0	0	1
5	F	27	0	0	2	0
5	G	32	0	0	2	0
All	All	5018	4864	4864	29	1

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 3.

The worst 5 of 29 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:F:9:GLU:OE2	5:F:201:HOH:O	1.85	0.92
1:A:6:ARG:NE	2:F:18:GLU:OE2	2.21	0.72
1:B:89:ASP:OD2	5:B:201:HOH:O	2.08	0.71
2:G:16:CYS:O	5:G:201:HOH:O	2.07	0.70
2:G:38:ASN:OD1	2:G:60:ARG:NH2	2.30	0.63

All (1) 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)
5:C:119:HOH:O	5:D:101:HOH:O[1_455]	1.92	0.28

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	147/152 (97%)	144 (98%)	3 (2%)	0	100	100
1	B	147/152 (97%)	144 (98%)	3 (2%)	0	100	100
2	F	75/86 (87%)	75 (100%)	0	0	100	100
2	G	84/86 (98%)	84 (100%)	0	0	100	100
3	C	74/76 (97%)	74 (100%)	0	0	100	100
3	D	74/76 (97%)	74 (100%)	0	0	100	100
All	All	601/628 (96%)	595 (99%)	6 (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	127/129 (98%)	126 (99%)	1 (1%)	86	94
1	B	127/129 (98%)	127 (100%)	0	100	100
2	F	69/75 (92%)	69 (100%)	0	100	100
2	G	75/75 (100%)	73 (97%)	2 (3%)	52	73
3	C	68/68 (100%)	67 (98%)	1 (2%)	72	87
3	D	68/68 (100%)	68 (100%)	0	100	100
All	All	534/544 (98%)	530 (99%)	4 (1%)	86	95

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

Mol	Chain	Res	Type
1	A	129	TRP
2	G	4	LEU
2	G	75	GLN
3	C	70	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 4 ligands modelled in this entry, 4 are 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	149/152 (98%)	-0.27	2 (1%) 79 79	22, 41, 75, 87	0
1	B	149/152 (98%)	-0.32	0 100 100	20, 39, 72, 80	0
2	F	76/86 (88%)	-0.14	2 (2%) 59 58	19, 29, 71, 91	0
2	G	86/86 (100%)	-0.33	1 (1%) 81 81	16, 29, 58, 86	0
3	C	76/76 (100%)	-0.42	1 (1%) 79 79	20, 43, 63, 71	0
3	D	76/76 (100%)	-0.36	0 100 100	22, 44, 67, 77	0
All	All	612/628 (97%)	-0.30	6 (0%) 84 83	16, 38, 71, 91	0

The worst 5 of 6 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	F	43	VAL	3.4
1	A	120	PRO	3.1
1	A	151	ASN	3.0
3	C	57	SER	2.3
2	G	43	VAL	2.2

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
4	ZN	G	101	1/1	0.98	0.14	1.61	26,26,26,26	0
4	ZN	F	101	1/1	0.98	0.14	1.09	25,25,25,25	0
4	ZN	F	102	1/1	0.99	0.14	-0.12	27,27,27,27	0
4	ZN	G	102	1/1	0.99	0.14	-0.15	27,27,27,27	0

6.5 Other polymers (i)

There are no such residues in this entry.