



wwPDB X-ray Structure Validation Summary Report ⓘ

Dec 6, 2016 – 07:49 PM EST

PDB ID : 5I0J
Title : Crystal structure of DR2231_E47A mutant in complex with dUMP
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Deposited on : 2016-02-04
Resolution : 1.80 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 : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20028442
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-20028442

i

X-RAY DIFFRACTION

A.

 R_{free}

1	A	140
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2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4960 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.

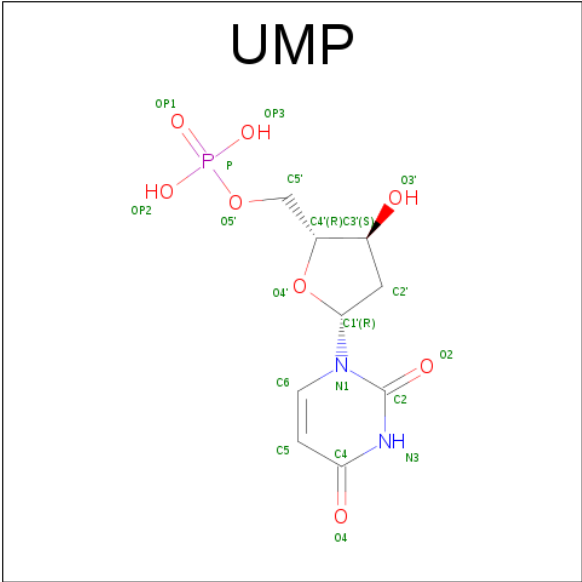
- Molecule 1 is a protein called DR2231.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	140	Total	C	N	O	S	0	0	0
			1069	666	202	200	1			
1	B	138	Total	C	N	O		0	0	0
			1056	658	200	198				
1	C	132	Total	C	N	O		0	0	0
			1010	632	188	190				
1	D	132	Total	C	N	O	S	0	0	0
			1005	629	186	189	1			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	47	ALA	GLU	engineered mutation	UNP Q9RS96
B	47	ALA	GLU	engineered mutation	UNP Q9RS96
C	47	ALA	GLU	engineered mutation	UNP Q9RS96
D	47	ALA	GLU	engineered mutation	UNP Q9RS96

- Molecule 2 is 2'-DEOXYURIDINE 5'-MONOPHOSPHATE (three-letter code: UMP) (formula: $C_9H_{13}N_2O_8P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			20	9	2	8	1		
2	B	1	Total	C	N	O	P	0	0
			20	9	2	8	1		

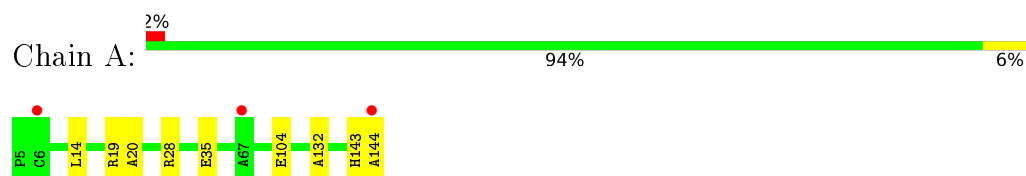
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	190	Total	O	0	0
			190	190		
3	B	185	Total	O	0	0
			185	185		
3	C	208	Total	O	0	0
			208	208		
3	D	197	Total	O	0	0
			197	197		

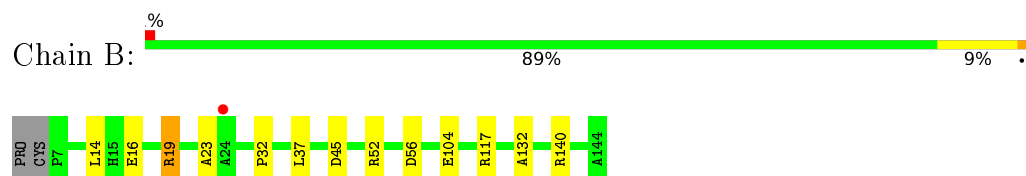
3 Residue-property plots [i](#)

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.

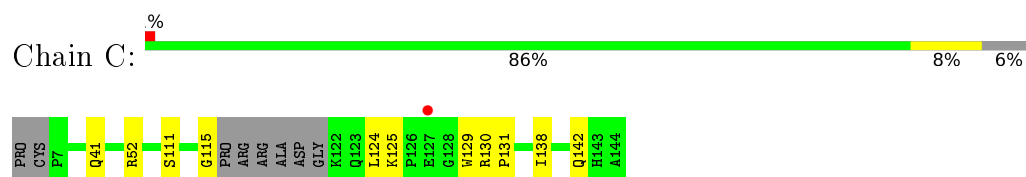
• Molecule 1: DR2231



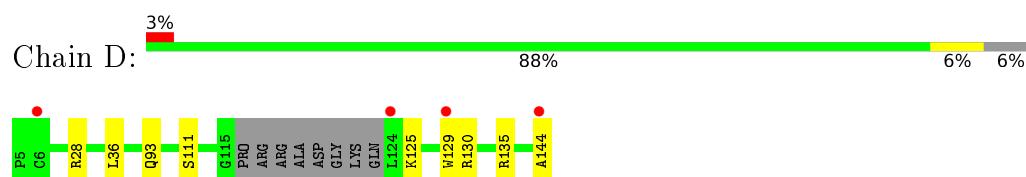
• Molecule 1: DR2231



• Molecule 1: DR2231



• Molecule 1: DR2231



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 ₁ 2 ₁ 2	Depositor
Cell constants a, b, c, α , β , γ	149.73 Å 78.07 Å 52.44 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.53 – 1.80 43.53 – 1.80	Depositor EDS
% Data completeness (in resolution range)	99.5 (43.53-1.80) 99.8 (43.53-1.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.83 (at 1.79 Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.173 , 0.211 0.169 , 0.207	Depositor DCC
R_{free} test set	2929 reflections (5.07%)	DCC
Wilson B-factor (Å ²)	19.1	Xtriage
Anisotropy	0.303	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 55.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4960	wwPDB-VP
Average B, all atoms (Å ²)	23.0	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 28.83 % 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 1.7209e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical 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: UMP

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	0.37	0/1093	0.50	0/1489
1	B	0.37	0/1079	0.53	1/1469 (0.1%)
1	C	0.37	0/1031	0.49	0/1403
1	D	0.37	0/1027	0.52	0/1400
All	All	0.37	0/4230	0.51	1/5761 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	19	ARG	NE-CZ-NH1	6.20	123.40	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	1069	0	1060	5	0
1	B	1056	0	1049	11	0
1	C	1010	0	1003	11	0
1	D	1005	0	993	11	0
2	A	20	0	11	0	0
2	B	20	0	11	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	190	0	0	0	0
3	B	185	0	0	6	0
3	C	208	0	0	3	0
3	D	197	0	0	2	0
All	All	4960	0	4127	36	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 4.

The worst 5 of 36 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)
1:D:135:ARG:HD2	1:D:135:ARG:H	1.36	0.89
1:D:28:ARG:HH12	1:D:144:ALA:HB2	1.51	0.76
1:D:130:ARG:NH2	3:D:201:HOH:O	2.20	0.75
1:D:135:ARG:N	1:D:135:ARG:HD2	2.03	0.74
1:B:16:GLU:HG3	1:B:19:ARG:HH22	1.53	0.73

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	138/140 (99%)	134 (97%)	4 (3%)	0	100	100
1	B	136/140 (97%)	133 (98%)	3 (2%)	0	100	100
1	C	128/140 (91%)	124 (97%)	4 (3%)	0	100	100
1	D	128/140 (91%)	125 (98%)	3 (2%)	0	100	100
All	All	530/560 (95%)	516 (97%)	14 (3%)	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	107/107 (100%)	106 (99%)	1 (1%)	84	80
1	B	105/107 (98%)	103 (98%)	2 (2%)	65	52
1	C	101/107 (94%)	101 (100%)	0	100	100
1	D	101/107 (94%)	101 (100%)	0	100	100
All	All	414/428 (97%)	411 (99%)	3 (1%)	88	86

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	LEU
1	B	14	LEU
1	B	140	ARG

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](#)

2 ligands are modelled in this entry.

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
2	UMP	A	201	-	17,21,21	4.05	10 (58%)	23,31,31	1.36	1 (4%)
2	UMP	B	201	-	17,21,21	3.85	9 (52%)	23,31,31	1.74	1 (4%)

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
2	UMP	A	201	-	-	0/6/22/22	0/2/2/2
2	UMP	B	201	-	-	0/6/22/22	0/2/2/2

The worst 5 of 19 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	201	UMP	C3'-C4'	-5.95	1.36	1.53
2	A	201	UMP	C3'-C4'	-5.90	1.36	1.53
2	A	201	UMP	O4'-C1'	-4.66	1.31	1.42
2	B	201	UMP	O4'-C1'	-4.48	1.32	1.42
2	B	201	UMP	O4-C4	-2.40	1.18	1.24

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	201	UMP	C4-N3-C2	5.73	120.24	114.21
2	B	201	UMP	C4-N3-C2	7.45	122.06	114.21

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	140/140 (100%)	-0.31	3 (2%) 67 62	9, 20, 44, 63	0
1	B	138/140 (98%)	-0.37	1 (0%) 89 87	10, 18, 38, 48	0
1	C	132/140 (94%)	-0.34	1 (0%) 87 85	9, 17, 40, 63	0
1	D	132/140 (94%)	-0.14	4 (3%) 54 48	9, 19, 48, 62	0
All	All	542/560 (96%)	-0.29	9 (1%) 73 69	9, 18, 43, 63	0

The worst 5 of 9 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	24	ALA	3.7
1	D	124	LEU	2.9
1	A	6	CYS	2.7
1	D	144	ALA	2.7
1	C	127	GLU	2.5

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(\AA^2)	Q<0.9
2	UMP	A	201	20/20	0.98	0.09	-0.09	8,11,13,16	0
2	UMP	B	201	20/20	0.98	0.08	-0.54	7,11,15,17	0

6.5 Other polymers [i](#)

There are no such residues in this entry.