



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 07:25 PM GMT

PDB ID : 1FKM
Title : CRYSTAL STRUCTURE OF THE YPT/RAB-GAP DOMAIN OF GYP1P
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Deposited on : 2000-08-09
Resolution : 1.90 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

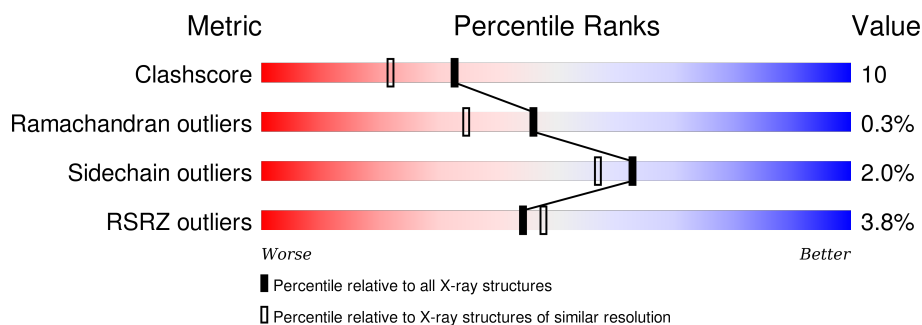
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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))
Clashscore	102246	5398 (1.90-1.90)
Ramachandran outliers	100387	5338 (1.90-1.90)
Sidechain outliers	100360	5339 (1.90-1.90)
RSRZ outliers	91569	4766 (1.90-1.90)

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	396	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2925 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 PROTEIN (GYP1P).

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	322	2759	1780	467	501	3	8	0	5	0

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	248	MSE	-	initiating met	UNP Q08484
A	413	MSE	MET	MODIFIED RESIDUE	UNP Q08484
A	484	MSE	MET	MODIFIED RESIDUE	UNP Q08484
A	489	MSE	MET	MODIFIED RESIDUE	UNP Q08484
A	494	MSE	MET	MODIFIED RESIDUE	UNP Q08484
A	500	MSE	MET	MODIFIED RESIDUE	UNP Q08484
A	518	MSE	MET	MODIFIED RESIDUE	UNP Q08484
A	554	MSE	MET	MODIFIED RESIDUE	UNP Q08484
A	566	MSE	MET	MODIFIED RESIDUE	UNP Q08484
A	591	MSE	MET	MODIFIED RESIDUE	UNP Q08484
A	593	MSE	MET	MODIFIED RESIDUE	UNP Q08484
A	617	MSE	MET	MODIFIED RESIDUE	UNP Q08484
A	638	HIS	-	EXPRESSION TAG	UNP Q08484
A	639	HIS	-	EXPRESSION TAG	UNP Q08484
A	640	HIS	-	EXPRESSION TAG	UNP Q08484
A	641	HIS	-	EXPRESSION TAG	UNP Q08484
A	642	HIS	-	EXPRESSION TAG	UNP Q08484
A	643	HIS	-	EXPRESSION TAG	UNP Q08484

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	166	Total	O	0	0
			166	166		

- Molecule 1: PROTEIN (GYP1P)



4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	74.06 Å 74.06 Å 277.75 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	6.00 – 1.90 6.00 – 1.90	Depositor EDS
% Data completeness (in resolution range)	94.7 (6.00-1.90) 94.6 (6.00-1.90)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.86 (at 1.90 Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.199 , 0.225 0.206 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	25.0	Xtriage
Anisotropy	0.316	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.58 , 71.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Outliers	0 of 33502 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	2925	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.95% 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](#)

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	0.32	0/2825	0.52	0/3821

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	2759	0	2707	57	0
2	A	166	0	0	7	0
All	All	2925	0	2707	57	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 10.

All (57) 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:A:341:ILE:HB	1:A:342:PRO:HD3	1.68	0.74
1:A:570:SER:HA	2:A:1186:HOH:O	1.88	0.72
1:A:608:LYS:HB2	1:A:608:LYS:NZ	2.06	0.70
1:A:430:THR:O	1:A:434:GLU:HG3	1.93	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:331:ILE:HB	1:A:332:PRO:HD3	1.77	0.67
1:A:265:LYS:HD3	1:A:265:LYS:O	1.96	0.66
1:A:458:VAL:HA	1:A:461:ILE:CD1	2.27	0.65
1:A:251:ILE:HD12	1:A:251:ILE:H	1.60	0.64
1:A:496:THR:HG23	1:A:615[A]:ILE:CD1	2.28	0.63
1:A:500:MSE:HG2	1:A:615[B]:ILE:HD11	1.80	0.63
1:A:319:LEU:HD21	1:A:362:ARG:HD2	1.80	0.61
1:A:615[A]:ILE:HD11	2:A:1130:HOH:O	2.02	0.60
1:A:328:SER:C	1:A:330:ASP:H	2.06	0.59
1:A:251:ILE:HD12	1:A:251:ILE:N	2.17	0.58
1:A:319:LEU:HA	1:A:322:THR:HG22	1.85	0.57
1:A:608:LYS:HB2	1:A:608:LYS:HZ3	1.71	0.56
1:A:580:ALA:O	1:A:584:ILE:HG12	2.06	0.55
1:A:604:ASN:N	1:A:605:PRO:HD3	2.22	0.55
1:A:341:ILE:HD11	1:A:364:LEU:HD12	1.88	0.55
1:A:330:ASP:HB3	1:A:365:TYR:OH	2.08	0.54
1:A:458:VAL:HA	1:A:461:ILE:HD13	1.89	0.53
1:A:266:THR:HG23	1:A:302:LYS:HD2	1.90	0.52
1:A:439:ASN:HD22	1:A:446:GLY:HA3	1.74	0.52
1:A:319:LEU:HG	1:A:323:PHE:CE2	2.45	0.51
1:A:251:ILE:CD1	1:A:251:ILE:H	2.24	0.50
1:A:452:LYS:HG3	2:A:1008:HOH:O	2.13	0.49
1:A:484:MSE:HG3	2:A:1172:HOH:O	2.12	0.49
1:A:265:LYS:HD2	1:A:267:ILE:O	2.14	0.48
1:A:400:GLN:OE1	1:A:407:LYS:NZ	2.47	0.48
1:A:500:MSE:HE2	2:A:1124:HOH:O	2.14	0.47
1:A:328:SER:HB2	1:A:330:ASP:OD1	2.14	0.47
1:A:494:MSE:O	1:A:498:ILE:HG12	2.14	0.47
1:A:447:ILE:HG21	1:A:481:PHE:CE1	2.49	0.47
1:A:448:LEU:O	1:A:452:LYS:HG3	2.14	0.47
1:A:447:ILE:HG21	1:A:481:PHE:CD1	2.50	0.46
1:A:389:GLU:OE1	1:A:494:MSE:HE3	2.16	0.46
1:A:450:GLN:HB3	1:A:484:MSE:HG2	1.98	0.45
1:A:408:ASP:HB3	1:A:411:THR:HG23	2.00	0.44
1:A:461:ILE:H	1:A:461:ILE:HD13	1.82	0.43
1:A:458:VAL:HA	1:A:461:ILE:HD11	2.01	0.43
1:A:285:HIS:HD2	2:A:1051:HOH:O	2.01	0.43
1:A:594:ASP:OD1	1:A:597:GLU:HG3	2.19	0.43
1:A:321:HIS:O	1:A:325:ASP:HB3	2.18	0.42
1:A:496:THR:O	1:A:500:MSE:HG3	2.19	0.42
1:A:496:THR:HG23	1:A:615[A]:ILE:HD13	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:307:PHE:O	1:A:311:LYS:HG2	2.19	0.42
1:A:311:LYS:HE3	1:A:427:TRP:CE2	2.55	0.42
1:A:472:GLU:O	1:A:473[A]:HIS:CG	2.73	0.42
1:A:505:LEU:O	1:A:509:SER:HB2	2.20	0.42
1:A:496:THR:HG23	1:A:615[A]:ILE:HG12	2.01	0.42
1:A:472:GLU:O	1:A:473[A]:HIS:ND1	2.52	0.42
1:A:482[B]:ARG:HD3	1:A:490:ARG:HB2	2.02	0.42
1:A:341:ILE:CB	1:A:342:PRO:HD3	2.45	0.41
1:A:450:GLN:CB	1:A:484:MSE:HG2	2.50	0.41
1:A:342:PRO:HA	2:A:1080:HOH:O	2.20	0.41
1:A:319:LEU:CA	1:A:322:THR:HG22	2.52	0.40
1:A:301:THR:HA	1:A:304:GLN:HG3	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	323/396 (82%)	309 (96%)	13 (4%)	1 (0%)	46	35

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	329	ARG

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	311/363 (86%)	305 (98%)	6 (2%)	65 59

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	438	ASP
1	A	457	LEU
1	A	461	ILE
1	A	484	MSE
1	A	494	MSE
1	A	572	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	249	ASN
1	A	253	GLN
1	A	285	HIS
1	A	435	GLN
1	A	439	ASN
1	A	453	ASN
1	A	467	ASN
1	A	572	ASN

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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	314/396 (79%)	-0.15	12 (3%)	44 48	16, 28, 52, 70	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	324	SER	8.7
1	A	325	ASP	5.3
1	A	327	HIS	5.2
1	A	328	SER	5.2
1	A	264	ASP	4.9
1	A	249	ASN	4.4
1	A	329	ARG	4.3
1	A	326	GLN	3.6
1	A	250	SER	2.8
1	A	265	LYS	2.6
1	A	322	THR	2.1
1	A	330	ASP	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 carbohydrates in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.