



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2016 – 11:21 PM GMT

PDB ID : 5EF5
Title : Crystal structure of Chaetomium thermophilum Raptor
Authors : Imseng, S.; Sauer, E.; Aylett, C.H.S.; Boehringer, D.; Hall, M.N.; Ban, N.; Maier, T.
Deposited on : 2015-10-23
Resolution : 4.30 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 : unknown
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026982
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-20026982

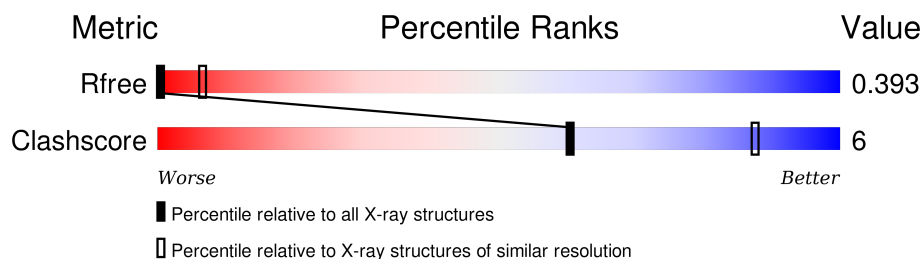
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 4.30 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	1059 (5.00-3.60)
Clashscore	102246	1166 (5.00-3.60)

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. 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	1029	 93% 7%
1	E	1029	 91% 6% •

2 Entry composition

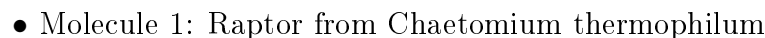
There is only 1 type of molecule in this entry. The entry contains 10164 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Raptor from Chaetomium thermophilum.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	E	1004	Total	C	N	O	0	0	0
			5019	3012	1004	1003			
1	A	1029	Total	C	N	O	0	0	0
			5145	3087	1029	1029			

- Molecule 1: Raptor from *Chaetomium thermophilum*



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	183.83Å 183.83Å 272.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.98 – 4.30 48.98 – 4.30	Depositor EDS
% Data completeness (in resolution range)	99.7 (48.98-4.30) 99.7 (48.98-4.30)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.75 (at 4.29Å)	Xtriage
Refinement program	PHENIX (phenix.refine: dev_1992)	Depositor
R, R_{free}	0.371 , 0.391 0.365 , 0.393	Depositor DCC
R_{free} test set	1593 reflections (4.92%)	DCC
Wilson B-factor (Å ²)	100.0	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 907.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 32407 reflections	Xtriage
F_o, F_c correlation	0.74	EDS
Total number of atoms	10164	wwPDB-VP
Average B, all atoms (Å ²)	77.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.91% 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

5.1 Standard geometry

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).

There are no protein, RNA or DNA chains available to summarize Z scores of covalent bonds and angles.

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	4730	UNK	Peptide
1	A	573	UNK	Peptide

5.2 Too-close contacts

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	5145	0	1096	35	0
1	E	5019	0	1054	34	0
All	All	10164	0	2150	69	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 6.

All (69) 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:48:UNK:HA	1:A:119:UNK:HA	1.64	0.79
1:E:107:UNK:N	1:E:108:UNK:HA	2.04	0.72
1:A:53:UNK:N	1:A:56:UNK:O	2.25	0.68
1:A:1551:UNK:O	1:A:1555:UNK:N	2.27	0.67
1:E:4548:UNK:N	1:E:4562:UNK:O	2.28	0.67
1:E:307:UNK:O	1:E:311:UNK:N	2.34	0.61
1:E:131:UNK:N	1:E:164:UNK:O	2.32	0.60
1:E:4685:UNK:O	1:E:4695:UNK:N	2.35	0.59
1:A:4654:UNK:HA	1:A:4661:UNK:HA	1.84	0.59
1:E:4525:UNK:O	1:E:4529:UNK:N	2.37	0.58
1:E:4692:UNK:O	1:E:4708:UNK:N	2.36	0.58
1:A:4672:UNK:O	1:A:4691:UNK:N	2.37	0.57
1:E:4593:UNK:N	1:E:4609:UNK:O	2.38	0.57
1:A:4762:UNK:O	1:A:4780:UNK:N	2.37	0.57
1:E:25:UNK:N	1:E:26:UNK:HA	2.19	0.56
1:E:40:UNK:O	1:E:74:UNK:N	2.39	0.56
1:A:200:UNK:O	1:A:204:UNK:N	2.39	0.56
1:E:82:UNK:O	1:E:86:UNK:N	2.40	0.55
1:E:1548:UNK:O	1:E:1552:UNK:N	2.40	0.54
1:A:4720:UNK:HA	1:A:4737:UNK:HA	1.88	0.54
1:A:4684:UNK:O	1:A:4700:UNK:N	2.40	0.54
1:A:23:UNK:N	1:A:24:UNK:HA	2.23	0.54
1:E:44:UNK:N	1:E:76:UNK:O	2.41	0.53
1:E:4686:UNK:HA	1:E:4694:UNK:HA	1.89	0.53
1:E:4558:UNK:HA	1:E:4572:UNK:HA	1.90	0.53
1:E:99:UNK:O	1:E:130:UNK:N	2.42	0.52
1:E:4681:UNK:HA	1:E:4698:UNK:HA	1.92	0.51
1:E:4799:UNK:HA	1:E:4800:UNK:C	2.41	0.51
1:A:199:UNK:O	1:A:203:UNK:N	2.43	0.51
1:A:40:UNK:O	1:A:86:UNK:N	2.44	0.50
1:A:1548:UNK:O	1:A:1552:UNK:N	2.44	0.50
1:E:212:UNK:O	1:E:216:UNK:N	2.45	0.49
1:E:261:UNK:O	1:E:265:UNK:N	2.46	0.49
1:E:25:UNK:H	1:E:26:UNK:HA	1.77	0.49
1:E:301:UNK:O	1:E:305:UNK:N	2.47	0.48
1:E:302:UNK:O	1:E:306:UNK:N	2.46	0.47
1:A:4539:UNK:N	1:A:4553:UNK:O	2.47	0.47
1:A:540:UNK:O	1:A:544:UNK:N	2.48	0.47
1:A:276:UNK:O	1:A:280:UNK:N	2.47	0.47
1:E:608:UNK:O	1:E:612:UNK:N	2.48	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:4831:UNK:HA	1:E:4845:UNK:O	2.15	0.47
1:A:46:UNK:O	1:A:92:UNK:N	2.47	0.47
1:A:4524:UNK:N	1:A:4830:UNK:O	2.48	0.46
1:A:513:UNK:HA	1:A:514:UNK:HA	1.83	0.46
1:E:4735:UNK:HA	1:E:4736:UNK:HA	1.63	0.46
1:E:42:UNK:N	1:E:74:UNK:O	2.49	0.46
1:A:4608:UNK:O	1:A:4621:UNK:N	2.49	0.45
1:E:4748:UNK:O	1:E:4765:UNK:HA	2.16	0.45
1:E:45:UNK:HA	1:E:78:UNK:O	2.16	0.45
1:A:110:UNK:N	1:A:145:UNK:O	2.50	0.44
1:A:4513:UNK:HA	1:A:4514:UNK:HA	1.79	0.44
1:E:4569:UNK:HA	1:E:4580:UNK:O	2.18	0.43
1:A:61:UNK:HA	1:A:62:UNK:C	2.48	0.43
1:E:137:UNK:HA	1:E:171:UNK:HA	2.00	0.43
1:A:243:UNK:O	1:A:247:UNK:N	2.52	0.43
1:A:455:UNK:O	1:A:459:UNK:N	2.52	0.43
1:A:128:UNK:CB	1:A:143:UNK:HA	2.48	0.42
1:E:493:UNK:HA	1:E:494:UNK:HA	1.84	0.42
1:A:489:UNK:O	1:A:490:UNK:C	2.67	0.42
1:A:406:UNK:N	1:A:407:UNK:HA	2.34	0.42
1:A:1552:UNK:O	1:A:1556:UNK:N	2.52	0.41
1:A:4763:UNK:HA	1:A:4779:UNK:HA	2.02	0.41
1:A:4524:UNK:O	1:A:4830:UNK:N	2.53	0.41
1:A:1505:UNK:O	1:A:1509:UNK:N	2.54	0.41
1:A:4815:UNK:HA	1:A:4825:UNK:O	2.20	0.41
1:E:4680:UNK:O	1:E:4699:UNK:N	2.53	0.41
1:E:4744:UNK:HA	1:E:4750:UNK:HA	2.03	0.40
1:A:4779:UNK:C	1:A:4781:UNK:N	2.81	0.40
1:A:303:UNK:HA	1:A:304:UNK:HA	1.32	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

There are no protein backbone outliers to report in this entry.

5.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

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](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	11
1	E	11

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	3030:UNK	C	4513:UNK	N	41.13
1	E	3040:UNK	C	4522:UNK	N	40.48
1	A	1557:UNK	C	3009:UNK	N	26.55
1	E	1565:UNK	C	3018:UNK	N	24.12
1	A	176:UNK	C	177:UNK	N	18.75
1	A	633:UNK	C	1505:UNK	N	17.54
1	E	49:UNK	C	50:UNK	N	16.92
1	E	612:UNK	C	1514:UNK	N	15.77
1	A	49:UNK	C	50:UNK	N	14.80
1	A	4630:UNK	C	4631:UNK	N	14.15

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Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	E	127:UNK	C	128:UNK	N	13.83
1	A	144:UNK	C	145:UNK	N	13.26
1	E	159:UNK	C	160:UNK	N	13.23
1	E	4638:UNK	C	4639:UNK	N	11.32
1	E	315:UNK	C	316:UNK	N	9.40
1	A	332:UNK	C	333:UNK	N	8.55
1	A	109:UNK	C	110:UNK	N	8.40
1	A	4799:UNK	C	4800:UNK	N	8.16
1	A	24:UNK	C	25:UNK	N	7.43
1	E	97:UNK	C	98:UNK	N	7.18
1	E	120:UNK	C	121:UNK	N	7.04
1	E	24:UNK	C	25:UNK	N	5.75

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	0/1029	-	-	-	-
1	E	0/1029	-	-	-	-
All	All	0/2058	-	-	-	-

There are no RSRZ outliers to report.

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 [i](#)

There are no such residues in this entry.