



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 1, 2016 – 05:50 AM GMT

PDB ID : 2UZK
Title : Crystal structure of the human FOXO3a-DBD bound to DNA
Authors : Tsai, K.-L.; Sun, Y.-J.; Huang, C.-Y.; Yang, J.-Y.; Hung, M.-C.; Hsiao, C.-D.
Deposited on : 2007-04-30
Resolution : 2.70 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 (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

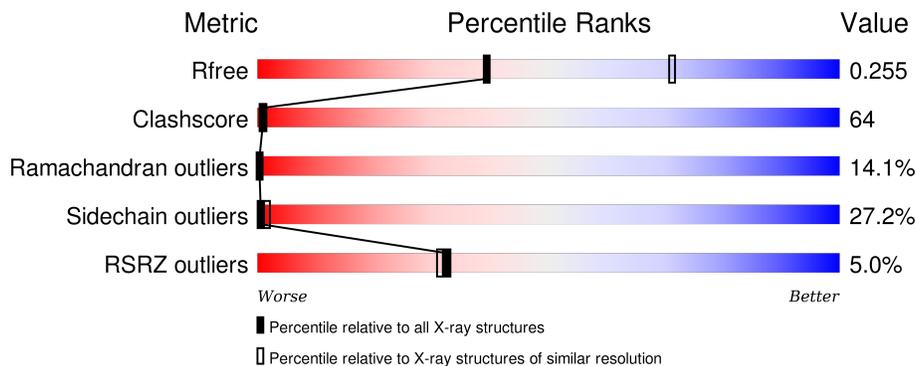
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	2103 (2.70-2.70)
Clashscore	102246	2422 (2.70-2.70)
Ramachandran outliers	100387	2382 (2.70-2.70)
Sidechain outliers	100360	2382 (2.70-2.70)
RSRZ outliers	91569	2107 (2.70-2.70)

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	97	
1	C	97	
2	B	13	
2	D	13	
3	E	13	

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Mol	Chain	Length	Quality of chain
3	F	13	 8% 92%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 2741 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 FORKHEAD BOX PROTEIN O3A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	97	Total 774	C 481	N 147	O 142	S 4	0	0	0
1	C	92	Total 704	C 437	N 131	O 133	S 3	0	0	1

- Molecule 2 is a DNA chain called 5'-D(*CP*TP*AP*TP*GP*TP*AP*AP*AP*CP*AP*AP*C)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	B	13	Total 262	C 127	N 50	O 73	P 12	0	0	0
2	D	13	Total 262	C 127	N 50	O 73	P 12	0	0	0

- Molecule 3 is a DNA chain called 5'-D(*GP*TP*TP*GP*TP*TP*TP*AP*CP*AP*TP*AP*G)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
3	E	13	Total 265	C 129	N 45	O 79	P 12	0	0	0
3	F	13	Total 265	C 129	N 45	O 79	P 12	0	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	46	Total 46 O 46	0	0
4	B	15	Total 15 O 15	0	0
4	C	72	Total 72 O 72	0	0

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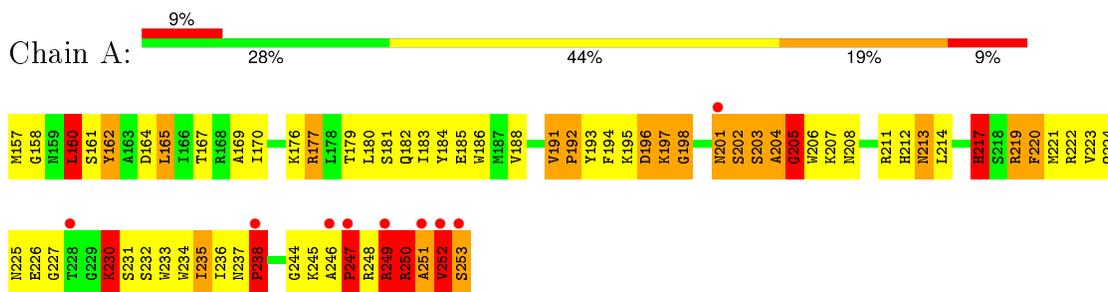
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	33	Total 33	O 33	0	0
4	E	22	Total 22	O 22	0	0
4	F	21	Total 21	O 21	0	0

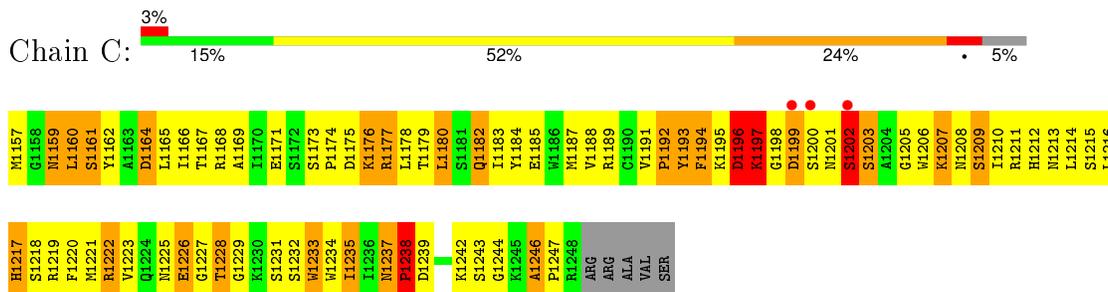
3 Residue-property plots [i](#)

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: FORKHEAD BOX PROTEIN O3A



- Molecule 1: FORKHEAD BOX PROTEIN O3A



- Molecule 2: 5'-D(*CP*TP*AP*TP*GP*TP*AP*AP*AP*CP*AP*AP*C)-3'



- Molecule 2: 5'-D(*CP*TP*AP*TP*GP*TP*AP*AP*AP*CP*AP*AP*C)-3'



- Molecule 3: 5'-D(*GP*TP*TP*GP*TP*TP*TP*AP*CP*AP*TP*AP*G)-3'



G25
T26
T27
G28
T29
T30
T31
A32
C33
A34
T35
A36
G37

- Molecule 3: 5'-D(*GP*TP*TP*GP*TP*TP*TP*AP*CP*AP*TP*AP*G)-3'

Chain F:  8% 92%

G1025
T1026
T1027
G1028
T1029
T1030
T1031
A1032
C1033
A1034
T1035
A1036
G1037

4 Data and refinement statistics

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α , β , γ	41.96Å 41.96Å 354.82Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	24.13 – 2.70 24.12 – 2.61	Depositor EDS
% Data completeness (in resolution range)	94.6 (24.13-2.70) 92.2 (24.12-2.61)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.84 (at 2.60Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.242 , 0.266 0.236 , 0.255	Depositor DCC
R_{free} test set	495 reflections (5.61%)	DCC
Wilson B-factor (Å ²)	24.2	Xtriage
Anisotropy	0.193	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.13 , 106.8	EDS
Estimated twinning fraction	0.399 for h,-h-k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.26$, $\langle L^2 \rangle = 0.11$	Xtriage
Outliers	0 of 9875 reflections	Xtriage
F_o, F_c correlation	0.86	EDS
Total number of atoms	2741	wwPDB-VP
Average B, all atoms (Å ²)	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 18.04% 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	0.75	2/792 (0.3%)	1.22	15/1064 (1.4%)
1	C	0.63	1/721 (0.1%)	1.50	10/972 (1.0%)
2	B	0.98	3/294 (1.0%)	1.20	6/451 (1.3%)
2	D	0.59	0/294	0.79	0/451
3	E	0.78	1/296 (0.3%)	0.87	0/456
3	F	0.49	0/296	0.83	0/456
All	All	0.71	7/2693 (0.3%)	1.18	31/3850 (0.8%)

The worst 5 of 7 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1	DC	O3'-P	10.65	1.74	1.61
3	E	37	DG	P-O5'	-8.27	1.51	1.59
1	C	1203	SER	CA-CB	7.60	1.64	1.52
1	A	247	PRO	C-N	-7.59	1.16	1.34
1	A	245	LYS	CB-CG	6.80	1.71	1.52

The worst 5 of 31 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	1237	ASN	C-N-CD	-21.96	72.29	120.60
1	C	1237	ASN	C-N-CA	12.43	174.22	122.00
1	C	1202	SER	C-N-CA	11.24	149.79	121.70
2	B	2	DT	O5'-P-OP1	-10.15	96.57	105.70
1	C	1196	ASP	N-CA-C	9.49	136.63	111.00

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	774	0	764	101	0
1	C	704	0	659	100	0
2	B	262	0	148	23	0
2	D	262	0	148	29	0
3	E	265	0	151	25	0
3	F	265	0	151	37	0
4	A	46	0	0	3	0
4	B	15	0	0	0	0
4	C	72	0	0	1	0
4	D	33	0	0	0	0
4	E	22	0	0	0	0
4	F	21	0	0	0	0
All	All	2741	0	2021	284	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 64.

The worst 5 of 284 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:D:1007:DA:H2''	2:D:1008:DA:H5'	1.26	1.14
1:C:1237:ASN:ND2	1:C:1238:PRO:HD3	1.68	1.07
1:A:237:ASN:HB2	1:A:238:PRO:HD2	1.11	1.05
1:A:160:LEU:HD12	1:A:165:LEU:HD22	1.39	1.05
2:B:5:DG:H2''	2:B:6:DT:H5''	1.37	1.03

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	95/97 (98%)	64 (67%)	17 (18%)	14 (15%)	0	0
1	C	90/97 (93%)	65 (72%)	13 (14%)	12 (13%)	0	0
All	All	185/194 (95%)	129 (70%)	30 (16%)	26 (14%)	0	0

5 of 26 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	198	GLY
1	A	204	ALA
1	A	217	HIS
1	A	238	PRO
1	A	247	PRO

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	84/84 (100%)	64 (76%)	20 (24%)	1	2
1	C	74/84 (88%)	51 (69%)	23 (31%)	0	0
All	All	158/168 (94%)	115 (73%)	43 (27%)	0	1

5 of 43 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	252	VAL
1	C	1177	ARG
1	C	1226	GLU
1	C	1159	ASN
1	C	1160	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 11 such sidechains are listed below:

Mol	Chain	Res	Type
1	A	217	HIS
1	C	1159	ASN
1	C	1208	ASN
1	A	213	ASN
1	C	1201	ASN

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](#)

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	97/97 (100%)	0.44	9 (9%) 11 8	0, 19, 61, 70	0
1	C	92/97 (94%)	0.34	3 (3%) 50 50	1, 18, 48, 58	0
2	B	13/13 (100%)	-0.58	0 100 100	3, 8, 18, 20	0
2	D	13/13 (100%)	-0.38	0 100 100	9, 14, 24, 27	0
3	E	13/13 (100%)	-0.56	0 100 100	3, 8, 23, 23	0
3	F	13/13 (100%)	-0.29	0 100 100	4, 17, 26, 26	0
All	All	241/246 (97%)	0.21	12 (4%) 32 31	0, 17, 56, 70	0

The worst 5 of 12 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	1200	SER	4.9
1	C	1199	ASP	4.8
1	A	228	THR	4.2
1	A	247	PRO	3.7
1	A	246	ALA	3.1

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.