



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 21, 2021 – 02:39 am BST

PDB ID : 6IC9  
Title : Crystal structure of the SPOC domain of human PHF3 in complex with RNA polymerase II CTD diheptapeptide phosphorylated on Ser2Ser7  
Authors : Grishkovskaya, I.; Djinovic-Carugo, K.; Slade, D.  
Deposited on : 2018-12-03  
Resolution : 1.75 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](mailto:validation@mail.wwpdb.org)

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.23.2  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0267  
CCP4 : 7.1.010 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.23.2

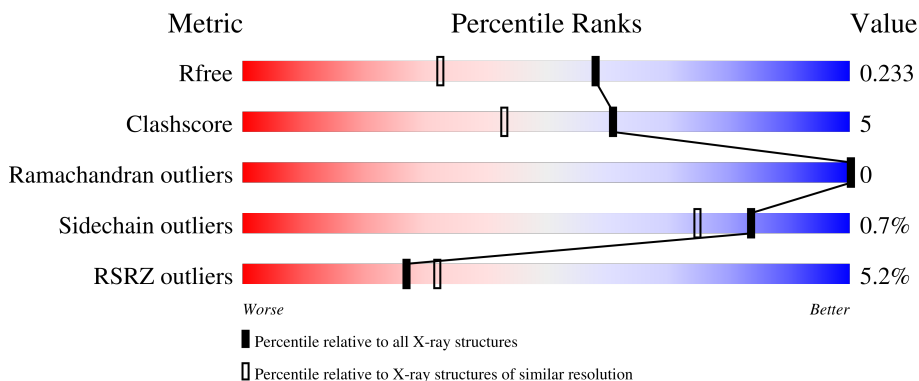
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.75 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))
$R_{free}$	130704	3764 (1.76-1.72)
Clashscore	141614	3923 (1.76-1.72)
Ramachandran outliers	138981	3878 (1.76-1.72)
Sidechain outliers	138945	3878 (1.76-1.72)
RSRZ outliers	127900	3705 (1.76-1.72)

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 of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. 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	162	 2% (Poor fit), 82% (0 outliers), 8% (1 outlier), 10% (2+ outliers)
1	B	162	 6% (Poor fit), 77% (0 outliers), 13% (1 outlier), 10% (2+ outliers)
2	C	14	 21% (Poor fit), 64% (0 outliers), 29% (1 outlier), 7% (2+ outliers)
2	D	14	 50% (0 outliers), 14% (1 outlier), 7% (2+ outliers), 29% (Not modelled)

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 2710 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 PHD finger protein 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	146	1177	770	193	210	4	0	2	0
1	B	146	1167	761	192	210	4	0	1	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1195	GLY	-	expression tag	UNP Q92576
A	1196	ALA	-	expression tag	UNP Q92576
A	1197	MET	-	expression tag	UNP Q92576
A	1198	GLY	-	expression tag	UNP Q92576
B	1195	GLY	-	expression tag	UNP Q92576
B	1196	ALA	-	expression tag	UNP Q92576
B	1197	MET	-	expression tag	UNP Q92576
B	1198	GLY	-	expression tag	UNP Q92576

- Molecule 2 is a protein called TYR-SEP-PRO-THR-SER-PRO-SEP-TYR-SEP-PRO-THR-SER-PRO.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	C	13	108	61	13	31	3	0	0	0
2	D	10	83	44	10	26	3	0	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	86	Total	O	0	0
			86	86		
3	B	60	Total	O	0	0
			60	60		

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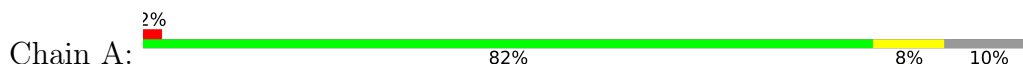
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<b>Mol</b>	<b>Chain</b>	<b>Residues</b>	<b>Atoms</b>	<b>ZeroOcc</b>	<b>AltConf</b>
3	C	21	Total O 21 21	0	0
3	D	8	Total O 8 8	0	0

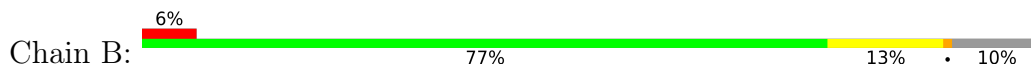
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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: PHD finger protein 3



- Molecule 1: PHD finger protein 3



- Molecule 2: TYR-SEP-PRO-THR-SER-PRO-SEP-TYR-SEP-PRO-THR-SER-PRO



- Molecule 2: TYR-SEP-PRO-THR-SER-PRO-SEP-TYR-SEP-PRO-THR-SER-PRO



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	56.22Å 68.11Å 100.87Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.10 – 1.75 49.10 – 1.75	Depositor EDS
% Data completeness (in resolution range)	98.7 (49.10-1.75) 99.0 (49.10-1.75)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.09 (at 1.75Å)	Xtrriage
Refinement program	PHENIX dev_3150	Depositor
R, $R_{free}$	0.195 , 0.232 0.195 , 0.233	Depositor DCC
$R_{free}$ test set	1953 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	32.7	Xtrriage
Anisotropy	0.191	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	(Not available) , (Not available)	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	2710	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	40.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.97% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>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: SEP

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	0.72	0/1208	0.84	2/1642 (0.1%)
1	B	0.67	1/1198 (0.1%)	0.74	1/1628 (0.1%)
2	C	0.69	0/80	0.66	0/106
2	D	0.53	0/54	0.54	0/71
All	All	0.69	1/2540 (0.0%)	0.79	3/3447 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	1270	CYS	CB-SG	-6.24	1.71	1.82

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1348	ARG	NE-CZ-NH2	-6.76	116.92	120.30
1	A	1338	ARG	NE-CZ-NH1	-6.13	117.24	120.30
1	B	1248	ARG	NE-CZ-NH2	-5.03	117.78	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	1177	0	1205	6	0
1	B	1167	0	1187	16	0
2	C	108	0	79	1	0
2	D	83	0	62	1	0
3	A	86	0	0	1	0
3	B	60	0	0	2	0
3	C	21	0	0	1	0
3	D	8	0	0	0	0
All	All	2710	0	2533	23	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 5.

All (23) 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:1206:ASN:N	3:A:1401:HOH:O	2.15	0.80
1:A:1268:GLU:HG2	1:A:1350:LYS:HG2	1.63	0.78
2:C:14:SER:OG	3:C:101:HOH:O	2.02	0.77
1:A:1265:GLY:O	1:A:1350:LYS:NZ	2.15	0.73
1:B:1248:ARG:HD2	2:D:11:SEP:O3P	2.02	0.59
1:B:1324:HIS:CG	1:B:1325:PRO:HD3	2.40	0.57
1:B:1337:HIS:NE2	3:B:1401:HOH:O	2.03	0.55
1:A:1304:ASN:HB3	1:A:1308:VAL:HG22	1.90	0.53
1:B:1209:TRP:CZ2	1:B:1211:GLY:HA3	2.44	0.53
1:B:1307:GLN:CD	1:B:1307:GLN:H	2.11	0.53
1:B:1296:LYS:HA	1:B:1296:LYS:HE2	1.92	0.52
1:B:1296:LYS:HD3	3:B:1439:HOH:O	2.10	0.50
1:B:1305:MET:HB3	1:B:1307:GLN:OE1	2.11	0.50
1:B:1234:TYR:HA	1:B:1237:GLU:OE2	2.12	0.49
1:A:1267:LYS:HE3	1:A:1349:GLN:OE1	2.13	0.48
1:B:1255:TRP:HA	1:B:1255:TRP:CE3	2.48	0.48
1:B:1306:LYS:HD2	1:B:1306:LYS:H	1.81	0.45
1:B:1258:VAL:O	1:B:1262:LYS:HG3	2.18	0.43
1:B:1268:GLU:OE1	1:B:1350:LYS:HE3	2.19	0.43
1:A:1209:TRP:CZ2	1:A:1211:GLY:HA3	2.53	0.43
1:B:1321:LYS:N	1:B:1321:LYS:HD2	2.33	0.43
1:B:1261:ILE:HG23	1:B:1267:LYS:HB2	2.01	0.42
1:B:1306:LYS:H	1:B:1306:LYS:CD	2.33	0.41

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	146/162 (90%)	144 (99%)	2 (1%)	0	100	100
1	B	145/162 (90%)	143 (99%)	2 (1%)	0	100	100
2	C	8/14 (57%)	8 (100%)	0	0	100	100
2	D	6/14 (43%)	6 (100%)	0	0	100	100
All	All	305/352 (87%)	301 (99%)	4 (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	131/141 (93%)	129 (98%)	2 (2%)	65	47
1	B	130/141 (92%)	129 (99%)	1 (1%)	81	72
2	C	10/10 (100%)	10 (100%)	0	100	100
2	D	7/10 (70%)	7 (100%)	0	100	100
All	All	278/302 (92%)	275 (99%)	3 (1%)	84	59

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

Mol	Chain	Res	Type
1	A	1313[A]	LEU
1	A	1313[B]	LEU
1	B	1206	ASN

Sometimes 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](#)

6 non-standard protein/DNA/RNA residues 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	SEP	C	4	2	8,9,10	1.02	0	8,12,14	1.29	1 (12%)
2	SEP	D	4	2,1	8,9,10	1.75	3 (37%)	8,12,14	1.16	1 (12%)
2	SEP	D	11	2	8,9,10	1.57	1 (12%)	8,12,14	1.78	3 (37%)
2	SEP	C	9	2	8,9,10	1.74	1 (12%)	8,12,14	1.33	2 (25%)
2	SEP	D	9	2	8,9,10	1.71	1 (12%)	8,12,14	1.63	2 (25%)
2	SEP	C	11	2	8,9,10	1.47	1 (12%)	8,12,14	2.35	5 (62%)

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	SEP	C	4	2	-	2/5/8/10	-
2	SEP	D	4	2,1	-	3/5/8/10	-
2	SEP	D	11	2	-	0/5/8/10	-
2	SEP	C	9	2	-	2/5/8/10	-
2	SEP	D	9	2	-	0/5/8/10	-
2	SEP	C	11	2	-	1/5/8/10	-

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	9	SEP	P-O1P	3.66	1.62	1.50
2	C	9	SEP	P-O1P	3.58	1.62	1.50
2	D	4	SEP	P-O1P	3.53	1.61	1.50
2	D	11	SEP	P-O1P	3.40	1.61	1.50
2	C	11	SEP	P-O1P	3.13	1.60	1.50
2	D	4	SEP	P-O2P	2.22	1.63	1.54
2	D	4	SEP	P-O3P	2.01	1.62	1.54

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	11	SEP	OG-CB-CA	4.03	112.06	108.14
2	D	9	SEP	P-OG-CB	-3.58	108.44	118.30
2	D	11	SEP	OG-CB-CA	3.31	111.37	108.14
2	C	11	SEP	P-OG-CB	-3.06	109.85	118.30
2	D	11	SEP	O3P-P-OG	2.61	113.67	106.73
2	C	4	SEP	P-OG-CB	-2.60	111.13	118.30
2	C	9	SEP	O3P-P-OG	2.44	113.21	106.73
2	D	4	SEP	P-OG-CB	-2.27	112.05	118.30
2	D	11	SEP	P-OG-CB	-2.24	112.14	118.30
2	D	9	SEP	OG-CB-CA	2.18	110.27	108.14
2	C	11	SEP	OG-P-O1P	2.14	112.49	106.47
2	C	11	SEP	O2P-P-OG	-2.13	101.06	106.73
2	C	11	SEP	O3P-P-OG	2.10	112.33	106.73
2	C	9	SEP	P-OG-CB	-2.04	112.67	118.30

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	4	SEP	CB-OG-P-O3P
2	C	9	SEP	CB-OG-P-O2P
2	C	9	SEP	CB-OG-P-O3P
2	C	11	SEP	N-CA-CB-OG
2	D	4	SEP	CB-OG-P-O2P
2	C	4	SEP	N-CA-CB-OG
2	D	4	SEP	N-CA-CB-OG
2	C	4	SEP	CB-OG-P-O3P

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	11	SEP	1	0

## 5.5 Carbohydrates [i](#)

There are no monosaccharides 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	146/162 (90%)	0.13	3 (2%) 63 70	25, 33, 53, 62	0
1	B	146/162 (90%)	0.25	10 (6%) 17 21	27, 40, 65, 84	0
2	C	10/14 (71%)	1.04	3 (30%) 0 0	29, 37, 66, 80	0
2	D	7/14 (50%)	0.11	0 100 100	46, 49, 53, 53	0
All	All	309/352 (87%)	0.22	16 (5%) 27 32	25, 37, 61, 84	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	1351	LEU	7.2
2	C	15	PRO	4.4
1	B	1306	LYS	4.2
1	B	1349	GLN	3.5
1	B	1233	GLU	3.5
1	A	1324	HIS	3.4
1	B	1234	TYR	2.9
1	B	1266	THR	2.9
1	A	1327	VAL	2.7
1	B	1255	TRP	2.7
1	B	1206	ASN	2.6
2	C	14	SER	2.6
1	B	1350	LYS	2.4
1	A	1321	LYS	2.3
1	B	1305	MET	2.3
2	C	3	TYR	2.0

### 6.2 Non-standard residues in protein, DNA, RNA chains [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. The B-factors column lists the minimum,

median, 95<sup>th</sup> 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	B-factors(Å <sup>2</sup> )	Q<0.9
2	SEP	D	4	10/11	0.60	0.23	59,70,84,87	0
2	SEP	D	9	10/11	0.80	0.18	50,66,89,96	0
2	SEP	C	9	10/11	0.90	0.10	34,39,73,74	0
2	SEP	D	11	10/11	0.96	0.07	43,50,64,70	0
2	SEP	C	11	10/11	0.98	0.12	34,38,46,47	0
2	SEP	C	4	10/11	0.98	0.12	35,40,46,48	0

### 6.3 Carbohydrates [i](#)

There are no monosaccharides 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.