Full wwPDB X-ray Structure Validation Report (i)

Aug 23, 2023 – 10:49 AM EDT

PDB ID	:	3D9J
Title	:	Snapshots of the RNA processing factor SCAF8 bound to different phosphorylated forms of the Carboxy-Terminal Domain of RNA-Polymerase II
Authors	:	Becker, R.; Loll, B.; Meinhart, A.
Deposited on	:	2008-05-27
Resolution	:	1.60 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
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

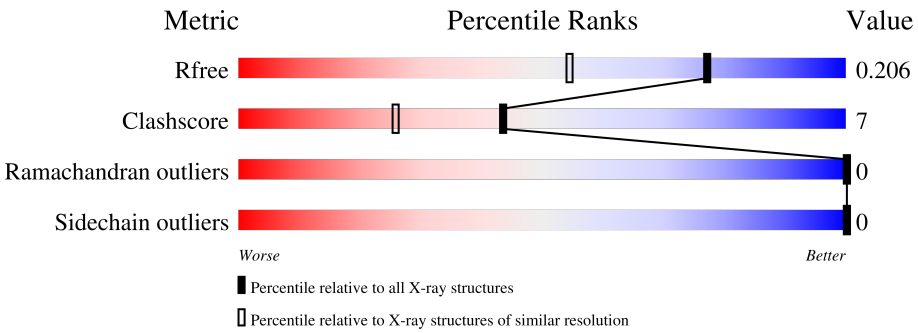
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.60 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	3398 (1.60-1.60)
Clashscore	141614	3665 (1.60-1.60)
Ramachandran outliers	138981	3564 (1.60-1.60)
Sidechain outliers	138945	3563 (1.60-1.60)

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 of the lower bar indicate the fraction of residues that contain outliers for >=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 <=5%

Mol	Chain	Length	Quality of chain
1	A	145	88% 7% 6%
1	B	145	83% 13% •

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	NH4	B	307	-	-	X	-

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 2669 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.

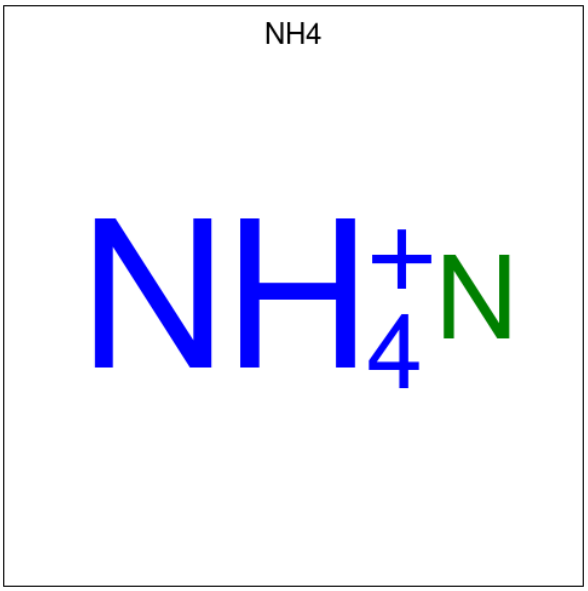
- Molecule 1 is a protein called RNA-binding protein 16.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	137	1156	753	197	202	4	0	9	0
1	B	140	1197	775	204	212	6	0	13	0

There are 18 discrepancies between the modelled and reference sequences:

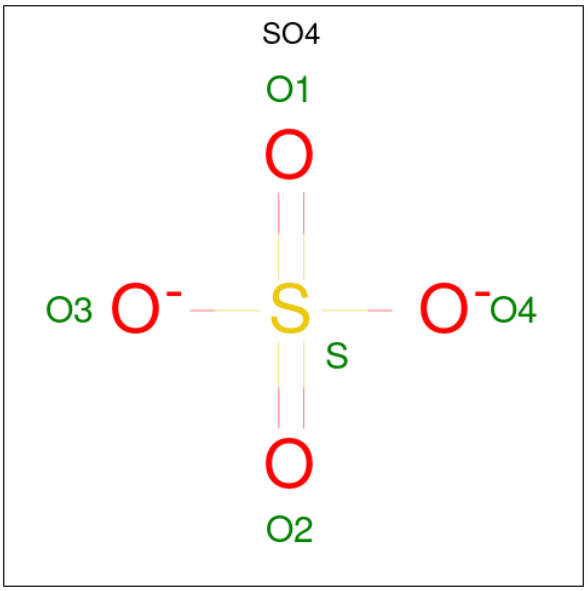
Chain	Residue	Modelled	Actual	Comment	Reference
A	137	ALA	-	expression tag	UNP Q9UPN6
A	138	LEU	-	expression tag	UNP Q9UPN6
A	139	GLU	-	expression tag	UNP Q9UPN6
A	140	HIS	-	expression tag	UNP Q9UPN6
A	141	HIS	-	expression tag	UNP Q9UPN6
A	142	HIS	-	expression tag	UNP Q9UPN6
A	143	HIS	-	expression tag	UNP Q9UPN6
A	144	HIS	-	expression tag	UNP Q9UPN6
A	145	HIS	-	expression tag	UNP Q9UPN6
B	137	ALA	-	expression tag	UNP Q9UPN6
B	138	LEU	-	expression tag	UNP Q9UPN6
B	139	GLU	-	expression tag	UNP Q9UPN6
B	140	HIS	-	expression tag	UNP Q9UPN6
B	141	HIS	-	expression tag	UNP Q9UPN6
B	142	HIS	-	expression tag	UNP Q9UPN6
B	143	HIS	-	expression tag	UNP Q9UPN6
B	144	HIS	-	expression tag	UNP Q9UPN6
B	145	HIS	-	expression tag	UNP Q9UPN6

- Molecule 2 is AMMONIUM ION (three-letter code: NH4) (formula: H_4N).



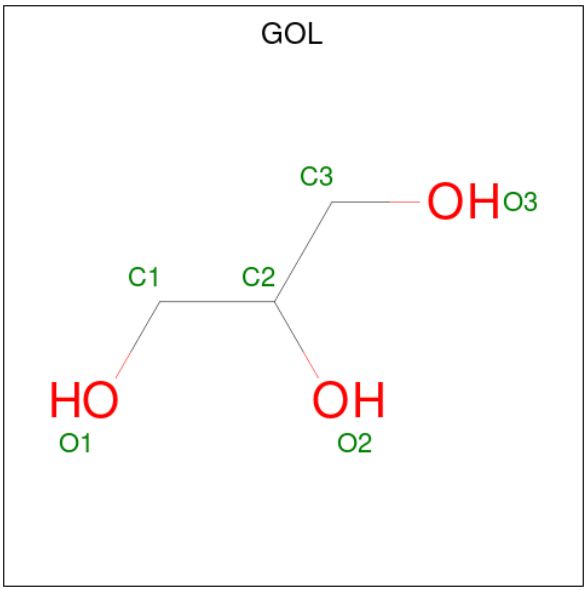
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total N 1 1	0	0
2	A	1	Total N 1 1	0	0
2	A	1	Total N 1 1	0	0
2	A	1	Total N 1 1	0	0
2	A	1	Total N 1 1	0	0
2	A	1	Total N 1 1	0	0
2	B	1	Total N 1 1	0	0
2	B	1	Total N 1 1	0	0
2	B	1	Total N 1 1	0	0
2	B	1	Total N 1 1	0	0
2	B	1	Total N 1 1	0	0
2	B	1	Total N 1 1	0	0

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
4	A	1	6	3	3	0	0
			Total	C	O		
4	A	1	6	3	3	0	0
			Total	C	O		
4	A	1	6	3	3	0	0
			Total	C	O		
4	B	1	6	3	3	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	A	115	115	115	0	0
			Total	O		
5	B	115	115	115	0	3

3 Residue-property plots (i)

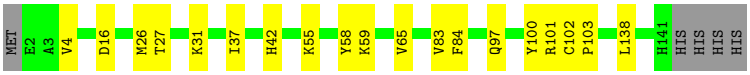
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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. 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: RNA-binding protein 16

Chain A:

- Molecule 1: RNA-binding protein 16

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 43	Depositor
Cell constants a, b, c, α, β, γ	57.62Å 57.62Å 107.15Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.76 – 1.60 40.74 – 1.50	Depositor EDS
% Data completeness (in resolution range)	98.7 (40.76-1.60) 98.5 (40.74-1.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.04	Depositor
$< I/\sigma(I) >$ [1]	2.63 (at 1.50Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.179 , 0.206 0.178 , 0.206	Depositor DCC
R_{free} test set	2625 reflections (4.76%)	wwPDB-VP
Wilson B-factor (Å^2)	20.0	Xtriage
Anisotropy	0.309	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 46.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	0.047 for h,-k,-l	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	2669	wwPDB-VP
Average B, all atoms (Å^2)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.89% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: GOL, SO4, NH4

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	0.64	0/1216	0.62	0/1636
1	B	0.60	0/1280	0.62	0/1724
All	All	0.62	0/2496	0.62	0/3360

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	1156	0	1207	12	0
1	B	1197	0	1224	21	0
2	A	6	0	0	0	0
2	B	6	0	0	2	0
3	A	30	0	0	0	0
3	B	20	0	0	0	0
4	A	18	0	24	4	0
4	B	6	0	8	0	0
5	A	115	0	0	5	0
5	B	115	0	0	2	0
All	All	2669	0	2463	33	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 7.

All (33) 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:112[A]:ARG:NH1	5:A:2120:HOH:O	2.04	0.89
1:B:4:VAL:HG13	1:B:42[B]:HIS:CD2	2.13	0.84
1:A:89:SER:O	4:A:1003:GOL:H12	1.84	0.77
1:B:102[B]:CYS:SG	1:B:103:PRO:HD2	2.30	0.72
1:B:55:LYS:HE2	5:B:2042:HOH:O	1.90	0.70
1:A:107:LYS:HE3	1:A:138:LEU:HD12	1.73	0.70
1:A:112[B]:ARG:NH1	5:A:2116:HOH:O	2.32	0.62
1:B:4:VAL:HG13	1:B:42[B]:HIS:NE2	2.14	0.62
1:A:127[A]:ILE:HD12	4:A:1003:GOL:H11	1.85	0.57
1:B:37[A]:ILE:HD12	1:B:84:PHE:CE2	2.39	0.56
1:A:121:ASN:OD1	5:A:2044:HOH:O	2.17	0.56
1:B:59:LYS:HB2	1:B:102[B]:CYS:SG	2.46	0.55
1:B:55:LYS:HE3	1:B:58:TYR:HE2	1.71	0.55
1:B:55:LYS:HE3	1:B:58:TYR:CE2	2.42	0.54
1:B:4:VAL:CG1	1:B:42[B]:HIS:CD2	2.91	0.52
1:B:102[B]:CYS:SG	1:B:103:PRO:CD	2.97	0.52
1:B:97[B]:GLN:HG2	1:B:101:ARG:HH21	1.74	0.52
1:A:34[B]:ILE:HD13	1:A:72:GLN:HG3	1.96	0.48
1:B:26[B]:MET:HE1	1:B:65:VAL:HA	1.97	0.47
1:A:124:LYS:NZ	5:A:2073:HOH:O	2.42	0.47
1:A:34[B]:ILE:CD1	1:A:72:GLN:HG3	2.46	0.45
1:B:100:TYR:HB3	1:B:138:LEU:HD11	1.98	0.45
1:B:37[A]:ILE:HD11	2:B:307:NH4:N	2.32	0.44
4:A:1000:GOL:H12	5:A:2100:HOH:O	2.17	0.44
1:A:127[A]:ILE:CD1	4:A:1003:GOL:H11	2.48	0.43
1:B:27:THR:CG2	1:B:31:LYS:HE2	2.49	0.42
1:A:34[B]:ILE:CD1	1:A:72:GLN:CG	2.98	0.41
1:A:22[B]:SER:HB2	1:B:16:ASP:OD2	2.21	0.41
1:B:42[A]:HIS:CE1	5:B:2068:HOH:O	2.72	0.41
1:B:37[A]:ILE:HD12	1:B:84:PHE:HE2	1.83	0.41
1:B:83:VAL:HG22	2:B:307:NH4:N	2.37	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	144/145 (99%)	143 (99%)	1 (1%)	0	100	100
1	B	152/145 (105%)	150 (99%)	2 (1%)	0	100	100
All	All	296/290 (102%)	293 (99%)	3 (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	134/133 (101%)	134 (100%)	0	100	100
1	B	142/133 (107%)	142 (100%)	0	100	100
All	All	276/266 (104%)	276 (100%)	0	100	100

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

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	121	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 26 ligands modelled in this entry, 12 are modelled with single atom - leaving 14 for Mogul analysis.

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	$\#\|Z\| > 2$
3	SO4	A	2003	-	4,4,4	0.28	0	6,6,6	0.44	0
3	SO4	B	2008	-	4,4,4	0.11	0	6,6,6	0.30	0
3	SO4	A	2000	-	4,4,4	0.26	0	6,6,6	0.49	0
4	GOL	A	1000	-	5,5,5	0.43	0	5,5,5	0.43	0
3	SO4	B	2006	-	4,4,4	0.17	0	6,6,6	0.19	0
3	SO4	A	2001	-	4,4,4	0.09	0	6,6,6	0.36	0
3	SO4	A	2002	-	4,4,4	0.15	0	6,6,6	0.39	0
3	SO4	B	2009	-	4,4,4	0.15	0	6,6,6	0.13	0
4	GOL	A	1003	-	5,5,5	0.34	0	5,5,5	0.47	0
4	GOL	B	1002	-	5,5,5	0.44	0	5,5,5	0.54	0
4	GOL	A	1001	-	5,5,5	0.34	0	5,5,5	0.27	0
3	SO4	A	2004	-	4,4,4	0.16	0	6,6,6	0.16	0
3	SO4	A	2005	-	4,4,4	0.20	0	6,6,6	0.11	0
3	SO4	B	2007	-	4,4,4	0.16	0	6,6,6	0.42	0

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
4	GOL	A	1000	-	-	3/4/4/4	-
4	GOL	A	1001	-	-	3/4/4/4	-
4	GOL	A	1003	-	-	1/4/4/4	-
4	GOL	B	1002	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	1001	GOL	C1-C2-C3-O3
4	B	1002	GOL	O1-C1-C2-C3
4	A	1000	GOL	O2-C2-C3-O3
4	A	1001	GOL	O2-C2-C3-O3
4	B	1002	GOL	O1-C1-C2-O2
4	A	1001	GOL	O1-C1-C2-C3
4	A	1000	GOL	C1-C2-C3-O3
4	A	1003	GOL	O1-C1-C2-C3
4	A	1000	GOL	O1-C1-C2-O2

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1000	GOL	1	0
4	A	1003	GOL	3	0

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)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.4 Ligands (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.5 Other polymers (i)

Unable to reproduce the depositors R factor - this section is therefore empty.