

# Full wwPDB X-ray Structure Validation Report (i)

#### Aug 23, 2023 - 10:20 AM EDT

PDB ID	:	3DN9
Title	:	Carboxysome Subunit, CcmK1 C-terminal deletion mutant
Authors	:	Tanaka, S.; Sawaya, M.R.; Yeates, T.O.
Deposited on		
Resolution	:	2.28 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 (1)) were used in the production of this report:

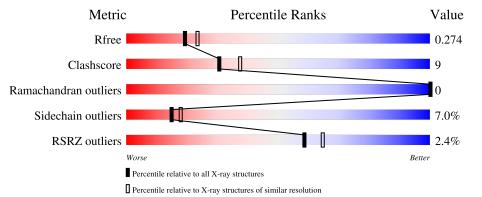
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
$\mathrm{EDS}$	:	2.35
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber $(2001)$
Ideal 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:  $X\text{-}RAY \, DIFFRACTION$ 

The reported resolution of this entry is 2.28 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	$\begin{array}{c} \textbf{Whole archive} \\ (\#\textbf{Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
$R_{free}$	130704	$6980 \ (2.30-2.26)$
Clashscore	141614	7711 (2.30-2.26)
Ramachandran outliers	138981	7597 (2.30-2.26)
Sidechain outliers	138945	7598 (2.30-2.26)
RSRZ outliers	127900	6849 (2.30-2.26)

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% 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	А	99	75%	12%	·	11%
1	В	99	74%	13%	•	9%
1	С	99	.% 75%	13%	•	11%
1	D	99	73%	15%	•	9%
1	Е	99	72%	16%	•	11%



Mol	Chain	Length	Quality of chain			
			4%			
1	F	99	71%	16%	•	9%



# 2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 4066 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.

Mol	Chain	Residues		At	oms			ZeroOcc	AltConf	Trace
1	А	88	Total	С	Ν	0	S	0	2	0
	A	00	661	417	117	125	2	0		0
1	В	90	Total	С	Ν	0	S	0	1	0
	D	90	675	426	120	127	2	0	1	U
1	С	88	Total	С	Ν	0	S	0	0	0
	U	88	653	410	117	124	2	0		0
1	D	90	Total	С	Ν	0	S	0	1	0
	D	90	677	427	122	126	2	0	L	0
1	Е	88	Total	С	Ν	Ο	S	0	0	0
	Ľ	88	653	410	117	124	2	0	0	0
1	F	90	Total	С	Ν	0	S	0	0	0
	Г	90	669	422	119	126	2		0	0

• Molecule 1 is a protein called CcmK1 C-terminal deletion mutant.

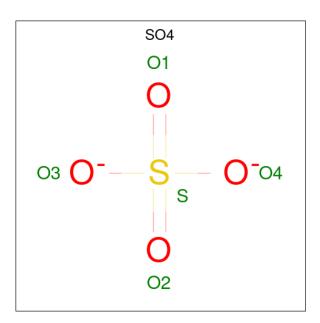
There are 48 discrepancies between the modelled and reference sequences:

	Modelled	Actual	Comment	Reference
92	LEU	-	expression tag	UNP P72760
93	GLU	-	expression tag	UNP P72760
94	HIS	-	expression tag	UNP P72760
95	HIS	-	expression tag	UNP P72760
96	HIS	-	expression tag	UNP P72760
97	HIS	-	expression tag	UNP P72760
98	HIS	-	expression tag	UNP P72760
99	HIS	-	expression tag	UNP P72760
92	LEU	-	expression tag	UNP P72760
93	GLU	-	expression tag	UNP P72760
94	HIS	-	expression tag	UNP P72760
95	HIS	-	expression tag	UNP P72760
96	HIS	-	expression tag	UNP P72760
97	HIS	-	expression tag	UNP P72760
98	HIS	-	expression tag	UNP P72760
99	HIS	-	expression tag	UNP P72760
92	LEU	-	expression tag	UNP P72760
	93         94         95         96         97         98         99         92         93         94         95         96         97         98         99         92         93         94         95         96         97         98         99	93         GLU           94         HIS           95         HIS           96         HIS           97         HIS           98         HIS           99         HIS           92         LEU           93         GLU           94         HIS           95         HIS           99         HIS           99         HIS           91         HIS           92         LEU           93         GLU           94         HIS           95         HIS           96         HIS           97         HIS           98         HIS           99         HIS	93       GLU       -         94       HIS       -         95       HIS       -         96       HIS       -         97       HIS       -         97       HIS       -         98       HIS       -         99       HIS       -         93       GLU       -         94       HIS       -         95       HIS       -         96       HIS       -         97       HIS       -         98       HIS       -         97       HIS       -         98       HIS       -         99       HIS       -         99       HIS       -	93GLU-expression tag94HIS-expression tag95HIS-expression tag96HIS-expression tag97HIS-expression tag98HIS-expression tag99HIS-expression tag93GLU-expression tag93GLU-expression tag93HIS-expression tag94HIS-expression tag95HIS-expression tag96HIS-expression tag97HIS-expression tag98HIS-expression tag99HIS-expression tag98HIS-expression tag99HIS-expression tag99HIS-expression tag



Chain	Residue	Modelled	Actual	Comment	Reference
С	93	GLU	-	expression tag	UNP P72760
С	94	HIS	-	expression tag	UNP P72760
С	95	HIS	-	expression tag	UNP P72760
С	96	HIS	-	expression tag	UNP P72760
С	97	HIS	-	expression tag	UNP P72760
С	98	HIS	-	expression tag	UNP P72760
С	99	HIS	-	expression tag	UNP P72760
D	92	LEU	-	expression tag	UNP P72760
D	93	GLU	-	expression tag	UNP P72760
D	94	HIS	-	expression tag	UNP P72760
D	95	HIS	-	expression tag	UNP P72760
D	96	HIS	-	expression tag	UNP P72760
D	97	HIS	-	expression tag	UNP P72760
D	98	HIS	-	expression tag	UNP P72760
D	99	HIS	-	expression tag	UNP P72760
Е	92	LEU	-	expression tag	UNP P72760
E	93	GLU	-	expression tag	UNP P72760
Е	94	HIS	-	expression tag	UNP P72760
E	95	HIS	-	expression tag	UNP P72760
Ε	96	HIS	-	expression tag	UNP P72760
Е	97	HIS	-	expression tag	UNP P72760
Е	98	HIS	-	expression tag	UNP P72760
Ε	99	HIS	-	expression tag	UNP P72760
F	92	LEU	-	expression tag	UNP P72760
F	93	GLU	-	expression tag	UNP P72760
F	94	HIS	-	expression tag	UNP P72760
F	95	HIS	-	expression tag	UNP P72760
F	96	HIS	-	expression tag	UNP P72760
F	97	HIS	-	expression tag	UNP P72760
F	98	HIS	-	expression tag	UNP P72760
F	99	HIS	-	expression tag	UNP P72760





Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	А	1	Total O S	0	0
		1	5 4 1		0
2	E	1	Total O S	0	0
2	Ľ	1	5 4 1	0	0
0	E	1	Total O S	0	0
	Ľ	1	$5 \ 4 \ 1$	0	0
0	Б	1	Total O S	0	0
	Г	1	5 4 1	0	0

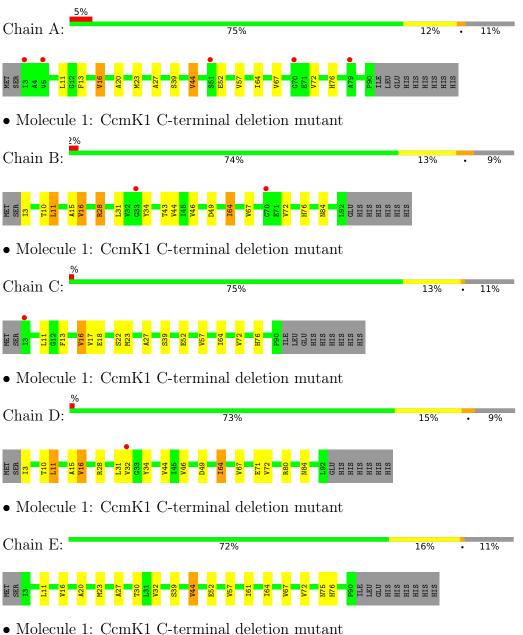
• Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	А	3	Total O 3 3	0	0
3	В	16	Total         O           16         16	0	0
3	С	11	Total O 11 11	0	0
3	D	7	Total O 7 7	0	0
3	Е	9	Total O 9 9	0	0
3	F	12	TotalO1212	0	0



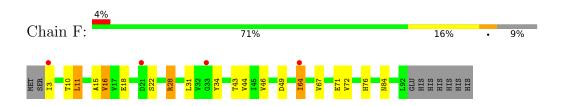
# 3 Residue-property plots (i)

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: CcmK1 C-terminal deletion mutant







#### Data and refinement statistics (i) 4

Property	Value	Source
Space group	P 1	Depositor
Cell constants	49.20Å 49.19Å 49.21Å	Denesiten
a, b, c, $\alpha$ , $\beta$ , $\gamma$	$90.65^{\circ}$ $91.41^{\circ}$ $91.43^{\circ}$	Depositor
Resolution (Å)	17.61 - 2.28	Depositor
Resolution (A)	17.61 - 2.28	EDS
% Data completeness	77.4 (17.61-2.28)	Depositor
(in resolution range)	$76.1 \ (17.61-2.28)$	EDS
R <sub>merge</sub>	0.11	Depositor
R <sub>sym</sub>	(Not available)	Depositor
$< I/\sigma(I) > 1$	1.14 (at 2.27Å)	Xtriage
Refinement program	REFMAC	Depositor
D D	0.222 , $0.274$	Depositor
$R, R_{free}$	0.224 , $0.274$	DCC
$R_{free}$ test set	816 reflections $(5.02\%)$	wwPDB-VP
Wilson B-factor $(Å^2)$	32.1	Xtriage
Anisotropy	0.074	Xtriage
Bulk solvent $k_{sol}(e/Å^3), B_{sol}(Å^2)$	0.37, $32.9$	EDS
L-test for twinning <sup>2</sup>	$<  L  > = 0.50, < L^2 > = 0.33$	Xtriage
	Continue	ed on next page

<sup>&</sup>lt;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.



Property	Value	Source	
	0.000 for l,k,-h		
	0.000 for -l,k,h		
	0.000 for h,-l,k		
	0.000 for h,l,-k		
	0.000 for -k,h,l		
	0.000 for k,-h,l		
	0.309 for l,h,k		
	0.309 for k,l,h		
	0.000 for -k,l,-h		
	0.000 for -l,-h,k		
	0.000 for -k,-l,h		
Estimated twinning fraction	0.000 for l,-h,-k	Xtriage	
-	0.000 for -l,h,-k		
	0.000 for k,-l,-h		
	0.016 for -h,k,-l		
	0.020 for h,-k,-l		
	0.022 for -h,-k,l		
	0.022 for k,h,-l		
	0.035 for -k,-h,-l		
	0.008 for -h,l,k		
	0.026 for -h,-l,-k		
	0.018 for l,-k,h		
	0.029 for -l,-k,-h		
$\mathbf{F}_o, \mathbf{F}_c$ correlation	0.93	EDS	
Total number of atoms	4066	wwPDB-V	
Average B, all atoms $(Å^2)$	34.0	wwPDB-V	

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 12.31% of the height of the origin peak. No significant pseudotranslation is detected.



# 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:  $\mathrm{SO4}$ 

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		
WIOI	Ullalli	RMSZ	# Z  > 5	RMSZ	# Z  > 5	
1	А	0.65	0/674	0.57	0/915	
1	В	0.65	0/685	0.59	1/930~(0.1%)	
1	С	0.64	0/660	0.58	0/896	
1	D	0.65	0/687	0.61	0/933	
1	Е	0.67	0/660	0.58	0/896	
1	F	0.64	0/676	0.60	1/918~(0.1%)	
All	All	0.65	0/4042	0.59	2/5488~(0.0%)	

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Ζ	$Observed(^{o})$	$Ideal(^{o})$
1	F	28	ARG	NE-CZ-NH1	5.13	122.86	120.30
1	В	28	ARG	NE-CZ-NH1	5.01	122.80	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	А	661	0	690	7	0
1	В	675	0	704	18	0
1	С	653	0	674	7	0



Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	677	0	709	20	0
1	Е	653	0	674	8	0
1	F	669	0	696	20	0
2	А	5	0	0	0	0
2	Е	10	0	0	0	0
2	F	5	0	0	0	0
3	А	3	0	0	0	0
3	В	16	0	0	0	0
3	С	11	0	0	0	0
3	D	7	0	0	1	0
3	Е	9	0	0	0	0
3	F	12	0	0	0	0
All	All	4066	0	4147	75	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 9.

All (75) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom 1	A + a	Interatomic	Clash
Atom-1	Atom-2	distance $(\text{\AA})$	overlap (Å)
1:D:10:THR:HG21	1:D:15:ALA:HB3	1.30	1.10
1:F:16:VAL:HG22	1:F:44:VAL:HG23	1.31	1.08
1:D:16:VAL:HG22	1:D:44:VAL:HG23	1.22	1.07
1:B:10:THR:HG21	1:B:15:ALA:HB3	1.37	1.05
1:B:16:VAL:HG22	1:B:44:VAL:HG23	1.31	1.04
1:F:10:THR:HG21	1:F:15:ALA:HB3	1.40	1.02
1:D:16:VAL:HG22	1:D:44:VAL:CG2	2.03	0.88
1:D:64:ILE:C	1:D:64:ILE:HD12	2.00	0.82
1:D:64:ILE:HG12	1:D:72:VAL:HG23	1.60	0.81
1:F:64:ILE:HG12	1:F:72:VAL:HG23	1.64	0.80
1:D:16:VAL:CG2	1:D:44:VAL:HG23	2.10	0.79
1:B:64:ILE:HG12	1:B:72:VAL:HG23	1.64	0.79
1:B:64:ILE:C	1:B:64:ILE:HD12	2.06	0.76
1:D:28:ARG:HG2	1:D:28:ARG:HH11	1.50	0.75
1:F:64:ILE:C	1:F:64:ILE:HD12	2.08	0.74
1:B:16:VAL:HG22	1:B:44:VAL:CG2	2.13	0.73
1:B:28:ARG:HH11	1:B:28:ARG:HG2	1.53	0.73
1:D:10:THR:HG21	1:D:15:ALA:CB	2.13	0.72
1:F:16:VAL:HG22	1:F:44:VAL:CG2	2.15	0.72
1:F:28:ARG:HH11	1:F:28:ARG:HG2	1.54	0.72
1:C:64:ILE:HD13	1:C:72:VAL:HG22	1.78	0.66



Continued from previous page Interatomic Clash								
Atom-1	Atom-2	distance (Å)	overlap (Å)					
1:B:10:THR:HG21	1:B:15:ALA:CB	2.21						
1:F:16:VAL:CG2	1:F:44:VAL:HG23	2.21 2.19	0.65					
1:D:64:ILE:HD12	1:D:64:ILE:O	1.99	0.63					
1:B:16:VAL:CG2	1:B:44:VAL:HG23	2.19	0.63					
1:D:28:ARG:HH11	1:D:28:ARG:CG	2.19	0.59					
1:D:10:THR:CG2	1:D:15:ALA:HB3	2.10	0.59					
1:D:34:TYR:HE2	3:D:105:HOH:O	1.86	0.58					
1:F:10:THR:HG21	1:F:15:ALA:CB	2.27	0.57					
1:B:67:VAL:HG13	1:F:15:ALA:CB 1:C:76:HIS:HB2	1.86	0.57					
1:B:28:ARG:HH11	1:B:28:ARG:CG	2.17	0.56					
1:B:10:THR:CG2	1:B:15:ALA:HB3	2.23	0.55					
1:D:64:ILE:C	1:D:64:ILE:CD1	2.72	0.55					
1:F:31:LEU:HA	1:F:46:VAL:HG12	1.89	0.54					
1:E:64:ILE:HD13	1:E:72:VAL:HG22	1.91	0.53					
1:F:28:ARG:HH11	1:F:28:ARG:CG	2.19	0.53					
1:A:23:MET:HE3	1:A:57:VAL:HG22	1.91	0.53					
1:A:27:ALA:HB1	1:A:52:GLU:OE2	2.09	0.52					
1:E:23:MET:HE3	1:E:57:VAL:HG22	1.91	0.52					
1:E:61:ILE:HD11	1:E:75:ASN:CG	2.30	0.52					
1:C:23:MET:HE3	1:C:57:VAL:HG22	1.91	0.52					
1:B:64:ILE:HD12	1:B:64:ILE:O	2.10	0.52					
1:C:27:ALA:HB1	1:C:52:GLU:OE2	2.10	0.52					
1:B:64:ILE:C	1:B:64:ILE:CD1	2.78	0.51					
1:F:10:THR:CG2	1:F:15:ALA:HB3	2.28	0.51					
1:B:31:LEU:HA	1:B:46:VAL:HG12	1.92	0.51					
1:A:67:VAL:HG13	1:B:76:HIS:HB2	1.91	0.50					
1:E:27:ALA:HB1	1:E:52:GLU:OE2	2.11	0.50					
1:D:31:LEU:HA	1:D:46:VAL:HG12	1.94	0.49					
1:F:64:ILE:C	1:F:64:ILE:CD1	2.79	0.48					
1:F:10:THR:HG23	1:F:71:GLU:O	2.13	0.48					
1:D:67:VAL:HG13	1:E:76:HIS:HB2	1.96	0.47					
1:C:13:PHE:O	1:C:16:VAL:HG12	2.14	0.47					
1:C:18:GLU:OE2	1:C:22:SER:OG	2.32	0.47					
1:F:31:LEU:HD11	1:F:44:VAL:HG13	1.98	0.46					
1:A:20:ALA:HB2	1:A:44:VAL:HG21	1.96	0.46					
1:A:64:ILE:HD13	1:A:72:VAL:HG22	1.97	0.46					
1:E:30:THR:O	1:E:32:VAL:HG13	2.16	0.46					
1:E:67:VAL:HG13	1:F:76:HIS:HB2	1.98	0.46					
1:F:34:TYR:HA	1:F:43:THR:O	2.16	0.46					
1:A:76:HIS:HB2	1:F:67:VAL:HG13	1.98	0.45					
1:A:13:PHE:O	1:A:16:VAL:HG12	2.16	0.45					



Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:13:PHE:O	1:C:17:VAL:HG23	2.17	0.45
1:F:64:ILE:HD12	1:F:64:ILE:O	2.17	0.44
1:F:18:GLU:OE2	1:F:22:SER:OG	2.37	0.43
1:B:34:TYR:HA	1:B:43:THR:O	2.18	0.43
1:D:32:VAL:O	1:D:32:VAL:HG23	2.18	0.43
1:D:10:THR:HG23	1:D:71:GLU:O	2.19	0.43
1:D:28:ARG:CG	1:D:28:ARG:NH1	2.80	0.42
1:D:64:ILE:O	1:D:64:ILE:CD1	2.67	0.42
1:E:20:ALA:HB2	1:E:44:VAL:HG21	2.02	0.41
1:B:10:THR:HG22	1:B:11:LEU:C	2.41	0.41
1:F:10:THR:HG22	1:F:11:LEU:C	2.41	0.41
1:B:28:ARG:CG	1:B:28:ARG:NH1	2.82	0.40
1:D:10:THR:HG22	1:D:11:LEU:C	2.42	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	А	88/99~(89%)	87~(99%)	1 (1%)	0	100 100
1	В	89/99~(90%)	87~(98%)	2(2%)	0	100 100
1	С	86/99~(87%)	85~(99%)	1 (1%)	0	100 100
1	D	89/99~(90%)	86~(97%)	3~(3%)	0	100 100
1	Е	86/99~(87%)	85~(99%)	1 (1%)	0	100 100
1	F	88/99~(89%)	86~(98%)	2(2%)	0	100 100
All	All	526/594~(89%)	516 (98%)	10 (2%)	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 side chain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric Outliers		Percentiles
1	А	72/81~(89%)	68~(94%)	4 (6%)	21 27
1	В	73/81~(90%)	67~(92%)	6 (8%)	11 13
1	$\mathbf{C}$	70/81~(86%)	67~(96%)	3~(4%)	29 38
1	D	73/81~(90%)	65~(89%)	8 (11%)	6 6
1	Ε	70/81~(86%)	66~(94%)	4 (6%)	20 26
1	F	72/81~(89%)	66~(92%)	6 (8%)	11 12
All	All	430/486~(88%)	399~(93%)	31~(7%)	15 17

All (31) residues with a non-rotameric side chain are listed below:

Mol	Chain	Res	Type
1	А	11	LEU
1	А	16	VAL
1	А	39	SER
1	А	44	VAL
1	В	3	ILE
1	В	11	LEU
1	В	16	VAL
1	В	49	ASP
1	В	64	ILE
1	В	84	ASN
1	C C C	11	LEU
1	С	16	VAL
1		39	SER
1	D	3	ILE
1	D	11	LEU
1	D	16	VAL
1	D	49	ASP
1	D	64	ILE
1	D	80[A]	ARG
1	D	80[B]	ARG
1	D	84	ASN
1	E	11	LEU



COULL	Commuted from previous page								
Mol	Chain	$\operatorname{Res}$	Type						
1	Е	16	VAL						
1	Е	39	SER						
1	Ε	44	VAL						
1	F	3	ILE						
1	F	11	LEU						
1	F	16	VAL						
1	F	49	ASP						
1	F	64	ILE						
1	F	84	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)

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)

4 ligands 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	B	ond leng	$\operatorname{gths}$	В	ond ang	gles
IVIOI	туре	Chain	nes	LIIIK	Counts	RMSZ	# Z >2	Counts	RMSZ	# Z >2
2	SO4	F	100	-	4,4,4	0.24	0	$6,\!6,\!6$	0.16	0
2	SO4	Е	101	-	4,4,4	0.23	0	$6,\!6,\!6$	0.38	0



Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
	туре			LIIIK	Counts	RMSZ	# Z >2	Counts	RMSZ	# Z >2
2	SO4	А	100	-	4,4,4	0.16	0	$6,\!6,\!6$	0.41	0
2	SO4	Е	100	-	4,4,4	0.15	0	$6,\!6,\!6$	0.17	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

### 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,  $95^{th}$  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	$\langle RSRZ \rangle$	#RSRZ>2	$OWAB(A^2)$	Q<0.9
1	А	88/99~(88%)	0.35	5 (5%) 23 28	18, 32, 48, 57	0
1	В	90/99~(90%)	0.22	2 (2%) 62 68	24, 33, 43, 48	0
1	С	88/99~(88%)	0.13	1 (1%) 80 84	18, 32, 48, 57	0
1	D	90/99~(90%)	0.18	1 (1%) 80 84	24, 33, 42, 48	0
1	Ε	88/99~(88%)	0.18	0 100 100	18, 32, 48, 57	0
1	F	90/99~(90%)	0.24	4 (4%) 34 40	24, 33, 42, 48	0
All	All	534/594~(89%)	0.22	13 (2%) 59 65	18, 33, 45, 57	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	В	70	GLY	4.8
1	А	70	GLY	3.3
1	В	33	GLY	3.1
1	А	51	SER	2.9
1	F	3	ILE	2.4
1	С	3	ILE	2.4
1	F	33	GLY	2.4
1	D	32	VAL	2.2
1	F	64	ILE	2.2
1	А	5	VAL	2.2
1	А	3	ILE	2.2
1	F	21	ASP	2.2
1	А	79	ALA	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 monosaccharides in this entry.

### 6.4 Ligands (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,  $95^{th}$  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	$\mathbf{B} ext{-factors}(\mathrm{\AA}^2)$	Q < 0.9
2	SO4	F	100	5/5	0.85	0.16	66,69,69,71	0
2	SO4	Е	100	5/5	0.93	0.11	71,71,72,73	0
2	SO4	А	100	5/5	0.94	0.11	$65,\!65,\!66,\!67$	0
2	SO4	Е	101	5/5	0.96	0.12	$50,\!51,\!55,\!55$	0

### 6.5 Other polymers (i)

There are no such residues in this entry.

