



# wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 1, 2016 – 07:28 PM GMT

PDB ID : 4OVN  
Title : Voltage-gated Sodium Channel 1.5 (Nav1.5) C-terminal domain in complex with Calmodulin poised for activation  
Authors : Gabelli, S.B.; Bianchet, M.A.; Boto, A.; Jakoncic, J.; Tomaselli, G.F.; Amzel, L.M.  
Deposited on : 2013-12-10  
Resolution : 2.80 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](mailto: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.

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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 : 1.7 (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



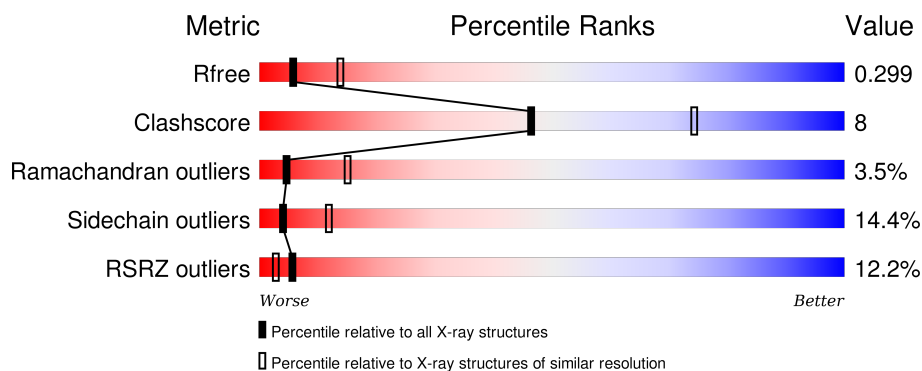
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.80 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	2393 (2.80-2.80)
Clashscore	102246	2827 (2.80-2.80)
Ramachandran outliers	100387	2782 (2.80-2.80)
Sidechain outliers	100360	2784 (2.80-2.80)
RSRZ outliers	91569	2404 (2.80-2.80)

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  $\geq 3$ , 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	149	<div> <div>12%</div> <div>73%</div> <div>23%</div> <div>..</div> </div>
1	B	149	<div> <div>19%</div> <div>73%</div> <div>21%</div> <div>..</div> </div>
1	C	149	<div> <div>7%</div> <div>77%</div> <div>19%</div> <div>..</div> </div>
1	D	149	<div> <div>5%</div> <div>72%</div> <div>21%</div> <div>...</div> </div>
1	E	149	<div> <div>21%</div> <div>69%</div> <div>25%</div> <div>..</div> </div>

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Mol	Chain	Length	Quality of chain
2	F	157	
2	G	157	
2	H	157	
2	I	157	
2	J	157	

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
3	MG	C	206	-	-	-	X



## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 23011 atoms, of which 11227 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 Calmodulin.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	145	Total	C	H	N	O	Se	0	0	0
			2213	701	1070	184	249	9			
1	B	144	Total	C	H	N	O	Se	0	0	0
			2197	696	1063	182	247	9			
1	C	145	Total	C	H	N	O	Se	0	0	0
			2213	701	1070	184	249	9			
1	D	145	Total	C	H	N	O	Se	0	0	0
			2213	701	1070	184	249	9			
1	E	144	Total	C	H	N	O	Se	0	0	0
			2197	696	1063	182	247	9			

- Molecule 2 is a protein called Sodium channel protein type 5 subunit alpha.

Mol	Chain	Residues	Atoms							ZeroOcc	AltConf	Trace
2	F	146	Total	C	H	N	O	S	Se	0	0	0
			2361	756	1174	198	224	1	8			
2	G	144	Total	C	H	N	O	S	Se	0	0	0
			2331	747	1158	196	221	1	8			
2	H	151	Total	C	H	N	O	S	Se	0	0	0
			2431	778	1207	203	234	1	8			
2	I	143	Total	C	H	N	O	S	Se	0	0	0
			2317	743	1151	195	219	1	8			
2	J	151	Total	C	H	N	O	S	Se	0	0	0
			2423	775	1201	203	235	1	8			

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	3	Total	Mg	0	0
			3	3		
3	A	4	Total	Mg	0	0
			4	4		

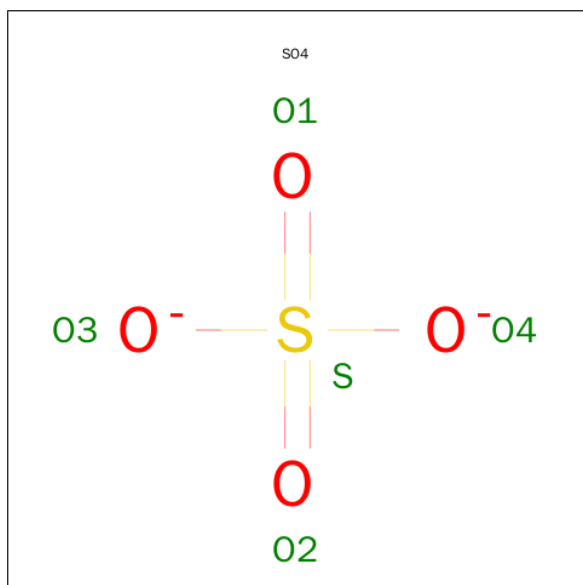
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	4	Total	Mg	0	0
			4	4		
3	C	5	Total	Mg	0	0
			5	5		
3	E	2	Total	Mg	0	0
			2	2		

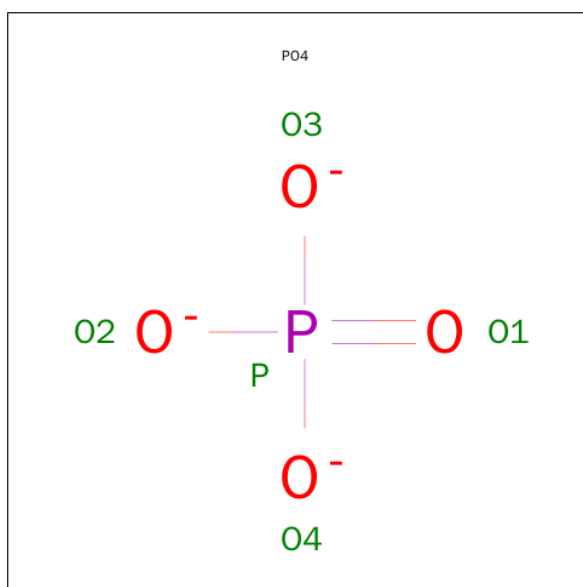
- Molecule 4 is SULFATE ION (three-letter code: SO<sub>4</sub>) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		
4	F	1	Total	O	S	0	0
			5	4	1		
4	G	1	Total	O	S	0	0
			5	4	1		
4	H	1	Total	O	S	0	0
			5	4	1		
4	J	1	Total	O	S	0	0
			5	4	1		

- Molecule 5 is PHOSPHATE ION (three-letter code: PO<sub>4</sub>) (formula: O<sub>4</sub>P).





Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	O	P	0	0
			5	4	1		
5	C	1	Total	O	P	0	0
			5	4	1		

- Molecule 6 is water.

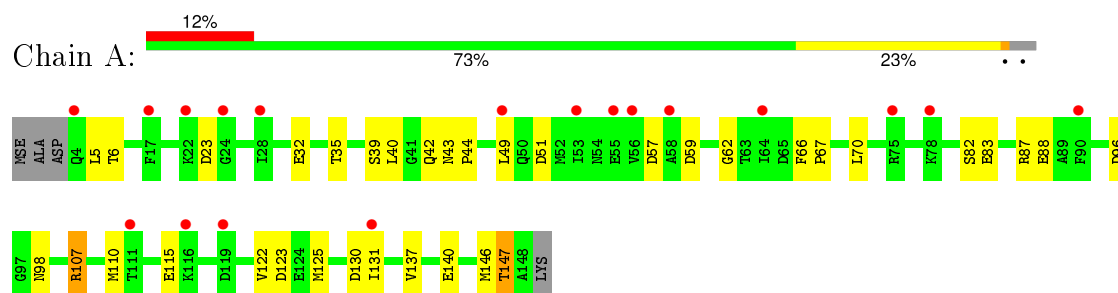
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	10	Total	O	0	0
			10	10		
6	F	3	Total	O	0	0
			3	3		
6	B	4	Total	O	0	0
			4	4		
6	G	3	Total	O	0	0
			3	3		
6	C	10	Total	O	0	0
			10	10		
6	H	4	Total	O	0	0
			4	4		
6	D	13	Total	O	0	0
			13	13		
6	I	3	Total	O	0	0
			3	3		
6	E	4	Total	O	0	0
			4	4		
6	J	8	Total	O	0	0
			8	8		



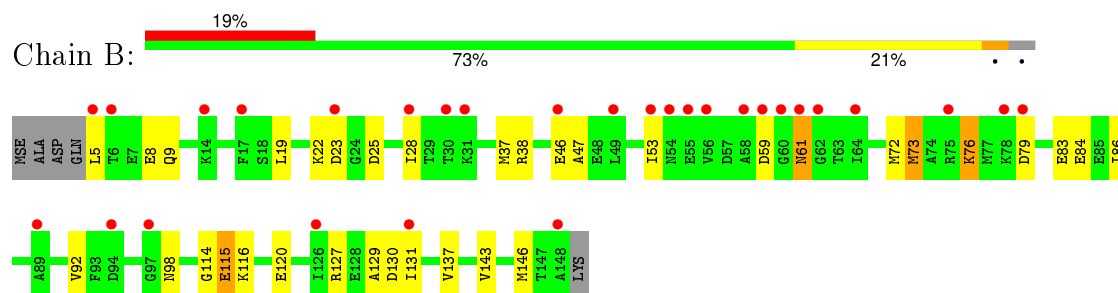
### 3 Residue-property plots [i](#)

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 ( $\text{RSRZ} > 2$ ). 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.

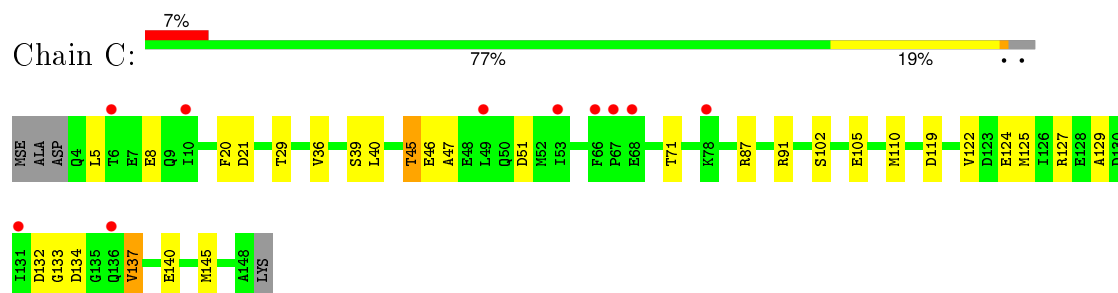
#### • Molecule 1: Calmodulin



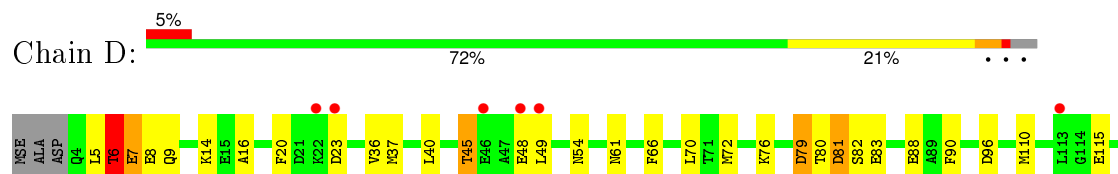
#### • Molecule 1: Calmodulin



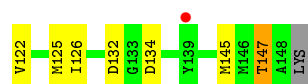
#### • Molecule 1: Calmodulin



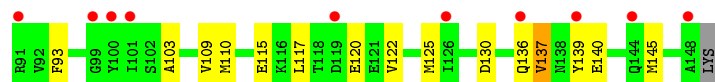
#### • Molecule 1: Calmodulin



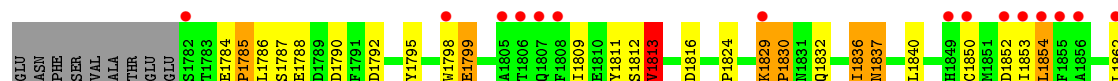




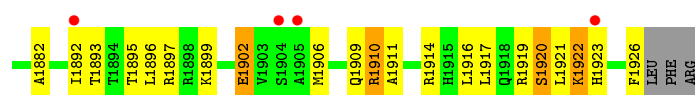
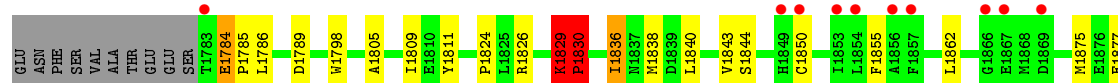
- Molecule 1: Calmodulin



- Molecule 2: Sodium channel protein type 5 subunit alpha



- Molecule 2: Sodium channel protein type 5 subunit alpha



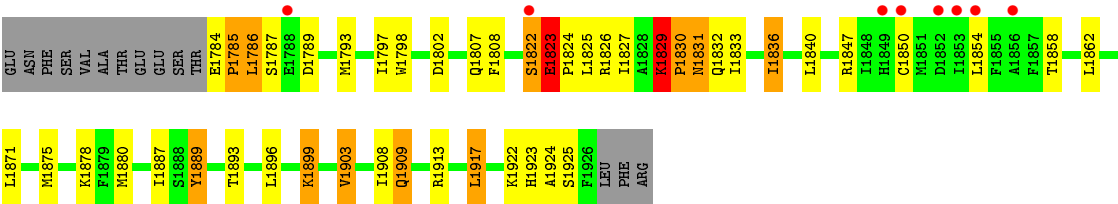
- Molecule 2: Sodium channel protein type 5 subunit alpha



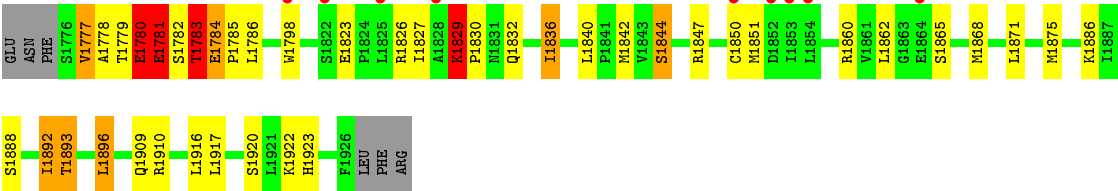
- Molecule 2: Sodium channel protein type 5 subunit alpha







● Molecule 2: Sodium channel protein type 5 subunit alpha





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	106.13 Å   99.01 Å   109.28 Å 90.00°   106.11°   90.00°	Depositor
Resolution (Å)	28.70 – 2.80 28.70 – 2.80	Depositor EDS
% Data completeness (in resolution range)	88.2 (28.70-2.80) 90.2 (28.70-2.80)	Depositor EDS
$R_{merge}$	0.12	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.48 (at 2.80 Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, $R_{free}$	0.211 , 0.282 0.245 , 0.299	Depositor DCC
$R_{free}$ test set	2390 reflections (5.19%)	DCC
Wilson B-factor (Å <sup>2</sup> )	78.3	Xtriage
Anisotropy	0.356	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 54.2	EDS
Estimated twinning fraction	0.021 for l,-k,h	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 48495 reflections	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	23011	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	93.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.71% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.



## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MG, PO4, 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	
		RMSZ	# Z  >5	RMSZ	# Z  >5
1	A	0.50	0/1146	0.66	0/1524
1	B	0.44	0/1137	0.58	0/1512
1	C	0.48	0/1146	0.82	3/1524 (0.2%)
1	D	0.49	0/1146	0.68	0/1524
1	E	0.42	0/1137	0.58	0/1512
2	F	0.44	0/1204	0.60	0/1608
2	G	0.45	0/1190	0.61	0/1589
2	H	0.49	0/1241	0.65	2/1659 (0.1%)
2	I	0.50	0/1183	0.67	1/1579 (0.1%)
2	J	0.51	0/1239	0.68	0/1656
All	All	0.47	0/11769	0.66	6/15687 (0.0%)

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
2	F	0	1
2	G	0	1
2	H	0	2
2	I	0	2
2	J	0	1
All	All	0	7

There are no bond length outliers.

The worst 5 of 6 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	21	ASP	CB-CG-OD2	-12.80	106.78	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	21	ASP	OD1-CG-OD2	10.61	143.45	123.30
1	C	21	ASP	CB-CG-OD1	-9.63	109.64	118.30
2	H	1829	LYS	C-N-CD	-6.77	105.70	120.60
2	I	1829	LYS	C-N-CD	-6.19	106.97	120.60

There are no chirality outliers.

5 of 7 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	F	1829	LYS	Peptide
2	G	1829	LYS	Peptide
2	H	1823	GLU	Peptide
2	H	1829	LYS	Peptide
2	I	1823	GLU	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	1143	1070	1070	17	0
1	B	1134	1063	1063	14	0
1	C	1143	1070	1070	8	0
1	D	1143	1070	1070	12	0
1	E	1134	1063	1063	18	0
2	F	1187	1174	1174	20	0
2	G	1173	1158	1158	21	0
2	H	1224	1207	1207	31	0
2	I	1166	1151	1151	28	0
2	J	1222	1201	1201	24	0
3	A	4	0	0	0	0
3	B	3	0	0	0	0
3	C	5	0	0	0	0
3	D	4	0	0	0	0
3	E	2	0	0	1	0
4	A	5	0	0	0	0
4	F	5	0	0	0	0
4	G	5	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	H	5	0	0	1	0
4	J	5	0	0	0	0
5	B	5	0	0	0	0
5	C	5	0	0	0	0
6	A	10	0	0	1	0
6	B	4	0	0	1	0
6	C	10	0	0	0	0
6	D	13	0	0	0	0
6	E	4	0	0	3	0
6	F	3	0	0	0	0
6	G	3	0	0	0	0
6	H	4	0	0	0	0
6	I	3	0	0	0	0
6	J	8	0	0	0	0
All	All	11784	11227	11227	175	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 8.

The worst 5 of 175 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:107:ARG:NH2	1:A:123:ASP:OD1	2.10	0.84
2:F:1927:LEU:HD23	2:F:1927:LEU:O	1.78	0.83
2:H:1925:SER:OG	2:H:1926:PHE:N	2.17	0.78
2:H:1925:SER:O	2:H:1926:PHE:HB3	1.84	0.76
2:I:1854:LEU:O	2:I:1858:THR:OG1	2.03	0.75

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	143/149 (96%)	127 (89%)	14 (10%)	2 (1%)	14	42
1	B	142/149 (95%)	122 (86%)	18 (13%)	2 (1%)	14	42
1	C	143/149 (96%)	126 (88%)	15 (10%)	2 (1%)	14	42
1	D	143/149 (96%)	124 (87%)	13 (9%)	6 (4%)	3	11
1	E	142/149 (95%)	120 (84%)	16 (11%)	6 (4%)	3	11
2	F	144/157 (92%)	118 (82%)	19 (13%)	7 (5%)	3	8
2	G	142/157 (90%)	118 (83%)	19 (13%)	5 (4%)	4	15
2	H	149/157 (95%)	125 (84%)	16 (11%)	8 (5%)	2	7
2	I	141/157 (90%)	119 (84%)	17 (12%)	5 (4%)	4	15
2	J	149/157 (95%)	126 (85%)	16 (11%)	7 (5%)	3	9
All	All	1438/1530 (94%)	1225 (85%)	163 (11%)	50 (4%)	4	15

5 of 50 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	F	1813	VAL
2	F	1830	PRO
1	B	115	GLU
2	G	1784	GLU
2	G	1829	LYS

### 5.3.2 Protein sidechains ⓘ

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	124/117 (106%)	111 (90%)	13 (10%)	8	24
1	B	123/117 (105%)	104 (85%)	19 (15%)	3	10
1	C	124/117 (106%)	110 (89%)	14 (11%)	7	22
1	D	124/117 (106%)	99 (80%)	25 (20%)	1	4
1	E	123/117 (105%)	111 (90%)	12 (10%)	10	28
2	F	132/134 (98%)	117 (89%)	15 (11%)	7	21

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	G	130/134 (97%)	108 (83%)	22 (17%)	2	7
2	H	136/134 (102%)	114 (84%)	22 (16%)	3	8
2	I	129/134 (96%)	108 (84%)	21 (16%)	3	8
2	J	136/134 (102%)	114 (84%)	22 (16%)	3	8
All	All	1281/1255 (102%)	1096 (86%)	185 (14%)	4	12

5 of 185 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	H	1780	GLU
2	H	1922	LYS
2	J	1836	ILE
2	H	1804	GLU
2	H	1862	LEU

Some 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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 25 ligands modelled in this entry, 18 are monoatomic - leaving 7 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	$\# Z  > 2$
4	SO4	A	205	-	4,4,4	0.12	0	6,6,6	0.20	0
5	PO4	B	204	-	4,4,4	0.33	0	6,6,6	0.27	0
5	PO4	C	205	-	4,4,4	0.23	0	6,6,6	0.28	0
4	SO4	F	2101	-	4,4,4	0.12	0	6,6,6	0.10	0
4	SO4	G	2101	-	4,4,4	0.07	0	6,6,6	0.20	0
4	SO4	H	2101	-	4,4,4	0.07	0	6,6,6	0.12	0
4	SO4	J	2101	-	4,4,4	0.12	0	6,6,6	0.10	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	SO4	A	205	-	-	0/0/0/0	0/0/0/0
5	PO4	B	204	-	-	0/0/0/0	0/0/0/0
5	PO4	C	205	-	-	0/0/0/0	0/0/0/0
4	SO4	F	2101	-	-	0/0/0/0	0/0/0/0
4	SO4	G	2101	-	-	0/0/0/0	0/0/0/0
4	SO4	H	2101	-	-	0/0/0/0	0/0/0/0
4	SO4	J	2101	-	-	0/0/0/0	0/0/0/0

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.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	G	2101	SO4	1	0
4	H	2101	SO4	1	0

## 5.7 Other polymers

There are no such residues in this entry.



## 5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.



## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

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	136/149 (91%)	0.69	18 (13%)	4 2	55, 89, 132, 165	0
1	B	135/149 (90%)	1.18	29 (21%)	1 1	65, 109, 181, 219	0
1	C	136/149 (91%)	0.41	10 (7%)	17 9	59, 87, 118, 138	0
1	D	136/149 (91%)	0.30	7 (5%)	32 21	57, 84, 119, 148	0
1	E	135/149 (90%)	1.14	31 (22%)	1 1	67, 122, 165, 200	0
2	F	138/157 (87%)	0.98	29 (21%)	1 1	55, 84, 138, 185	0
2	G	136/157 (86%)	0.69	14 (10%)	9 4	56, 80, 119, 153	0
2	H	143/157 (91%)	0.62	13 (9%)	11 6	54, 77, 119, 149	0
2	I	135/157 (85%)	0.44	8 (5%)	26 16	52, 67, 102, 115	0
2	J	143/157 (91%)	0.56	9 (6%)	23 14	53, 74, 100, 122	0
All	All	1373/1530 (89%)	0.70	168 (12%)	5 3	52, 84, 142, 219	0

The worst 5 of 168 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	53	ILE	10.3
1	E	54	ASN	9.2
1	B	62	GLY	8.2
1	B	54	ASN	8.0
1	B	56	VAL	6.8

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

There are no non-standard protein/DNA/RNA residues in this entry.



### 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

### 6.4 Ligands ⓘ

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
3	MG	C	206	1/1	0.92	1.05	13.15	66,66,66,66	0
3	MG	A	203	1/1	0.85	0.33	1.66	66,66,66,66	0
3	MG	B	202	1/1	0.88	0.38	1.25	66,66,66,66	0
3	MG	A	204	1/1	0.98	0.15	-0.55	70,70,70,70	0
3	MG	D	202	1/1	0.97	0.16	-0.67	71,71,71,71	0
3	MG	B	201	1/1	0.90	0.19	-0.79	113,113,113,113	0
3	MG	D	204	1/1	0.82	0.11	-0.85	91,91,91,91	0
3	MG	A	202	1/1	0.62	0.14	-0.95	115,115,115,115	0
3	MG	C	204	1/1	0.94	0.10	-1.06	102,102,102,102	0
3	MG	E	202	1/1	0.67	0.11	-1.26	83,83,83,83	0
3	MG	C	202	1/1	0.95	0.10	-1.37	68,68,68,68	0
3	MG	E	201	1/1	0.94	0.11	-1.44	113,113,113,113	0
3	MG	C	203	1/1	0.95	0.10	-1.63	80,80,80,80	0
3	MG	C	201	1/1	0.73	0.07	-1.64	94,94,94,94	0
3	MG	D	203	1/1	0.80	0.07	-1.68	96,96,96,96	0
3	MG	B	203	1/1	0.72	0.08	-1.81	95,95,95,95	0
3	MG	D	201	1/1	0.88	0.08	-1.94	89,89,89,89	0
4	SO4	F	2101	5/5	0.90	0.12	-2.12	66,82,101,114	0
3	MG	A	201	1/1	0.92	0.06	-2.42	89,89,89,89	0
5	PO4	B	204	5/5	0.75	0.35	-	103,108,111,135	0
4	SO4	J	2101	5/5	0.86	0.21	-	97,103,108,125	0
4	SO4	A	205	5/5	0.81	0.27	-	79,103,129,142	0
4	SO4	G	2101	5/5	0.72	0.35	-	98,105,111,140	0
4	SO4	H	2101	5/5	0.92	0.15	-	83,91,96,133	0
5	PO4	C	205	5/5	0.73	0.21	-	67,92,112,124	0

### 6.5 Other polymers ⓘ

There are no such residues in this entry.