



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:49 PM GMT

PDB ID : 3V5S
Title : Structure of Sodium/Calcium Exchanger from *Methanococcus jannaschii*
Authors : Liao, J.; Li, H.; Zeng, W.; Sauer, D.B.; Belmares, R.; Jiang, Y.
Deposited on : 2011-12-16
Resolution : 3.50 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
<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

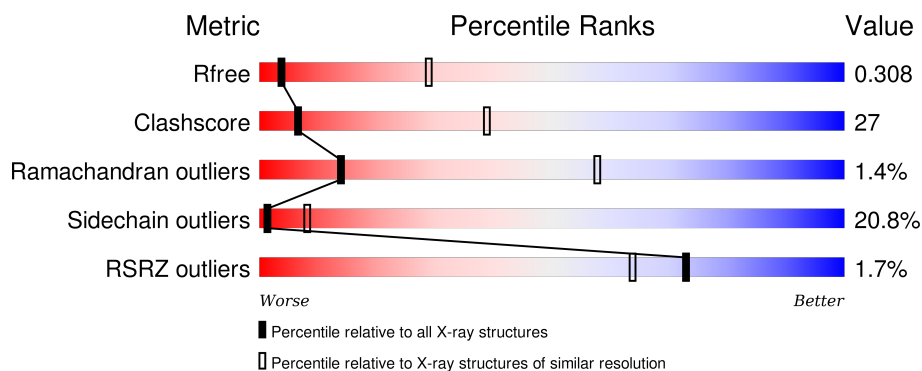
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.50 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	1051 (3.60-3.40)
Clashscore	102246	1157 (3.60-3.40)
Ramachandran outliers	100387	1120 (3.60-3.40)
Sidechain outliers	100360	1121 (3.60-3.40)
RSRZ outliers	91569	1058 (3.60-3.40)

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	320	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2172 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 Sodium/Calcium Exchanger.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	289	Total	C	N	O	S	0	0	0
			2170	1453	340	366	11			

There are 19 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	2	VAL	LEU	ENGINEERED MUTATION	UNP Q57556
A	303	GLY	-	EXPRESSION TAG	UNP Q57556
A	304	GLY	-	EXPRESSION TAG	UNP Q57556
A	305	SER	-	EXPRESSION TAG	UNP Q57556
A	306	LEU	-	EXPRESSION TAG	UNP Q57556
A	307	VAL	-	EXPRESSION TAG	UNP Q57556
A	308	PRO	-	EXPRESSION TAG	UNP Q57556
A	309	ARG	-	EXPRESSION TAG	UNP Q57556
A	310	GLY	-	EXPRESSION TAG	UNP Q57556
A	311	SER	-	EXPRESSION TAG	UNP Q57556
A	312	ARG	-	EXPRESSION TAG	UNP Q57556
A	313	SER	-	EXPRESSION TAG	UNP Q57556
A	314	HIS	-	EXPRESSION TAG	UNP Q57556
A	315	HIS	-	EXPRESSION TAG	UNP Q57556
A	316	HIS	-	EXPRESSION TAG	UNP Q57556
A	317	HIS	-	EXPRESSION TAG	UNP Q57556
A	318	HIS	-	EXPRESSION TAG	UNP Q57556
A	319	HIS	-	EXPRESSION TAG	UNP Q57556
A	320	HIS	-	EXPRESSION TAG	UNP Q57556

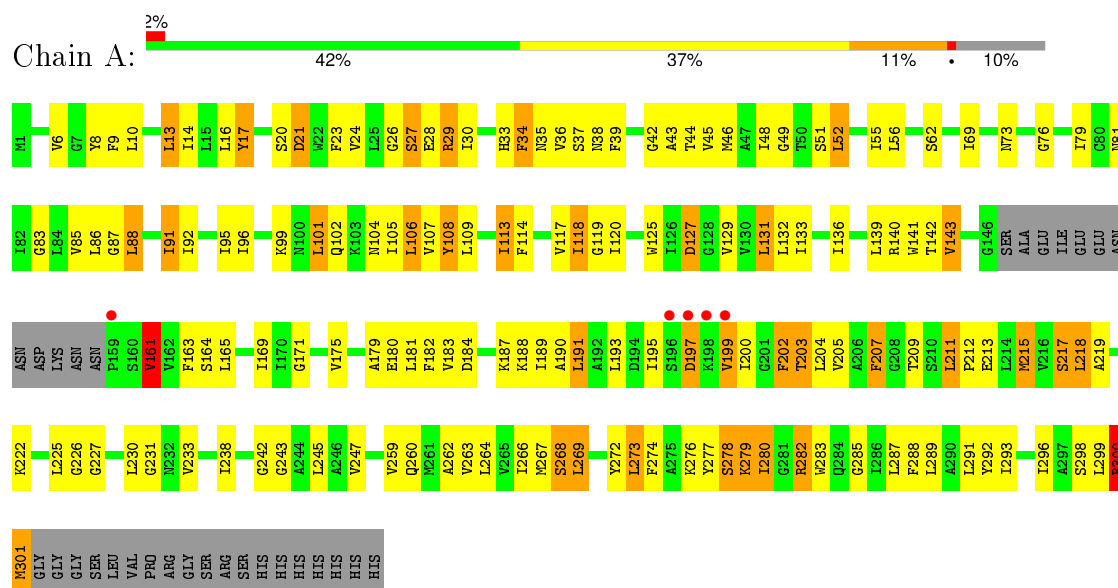
- Molecule 2 is CADMIUM ION (three-letter code: CD) (formula: Cd).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Cd	0	0
			2	2		

3 Residue-property plots

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: Sodium/Calcium Exchanger



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	84.99 Å 84.99 Å 181.63 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	37.20 – 3.50 42.65 – 3.50	Depositor EDS
% Data completeness (in resolution range)	98.3 (37.20-3.50) 98.3 (42.65-3.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.86 (at 3.48 Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.2_869)	Depositor
R, R_{free}	0.281 , 0.303 0.271 , 0.308	Depositor DCC
R_{free} test set	412 reflections (4.69%)	DCC
Wilson B-factor (Å ²)	141.5	Xtriage
Anisotropy	0.163	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 113.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	1 of 8815 reflections (0.011%)	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	2172	wwPDB-VP
Average B, all atoms (Å ²)	131.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.12% 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: CD

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	0.35	0/2210	0.69	3/3002 (0.1%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	279	LYS	N-CA-CB	-5.80	100.17	110.60
1	A	282	ARG	N-CA-CB	-5.56	100.60	110.60
1	A	278	SER	CB-CA-C	-5.52	99.61	110.10

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	2170	0	2327	123	0
2	A	2	0	0	0	0
All	All	2172	0	2327	123	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 27.

All (123) 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:260:GLN:CG	1:A:299:LEU:HD11	1.61	1.29
1:A:260:GLN:HG2	1:A:299:LEU:CD1	1.74	1.16
1:A:81:ASN:HD21	1:A:213:GLU:HG2	1.39	0.88
1:A:17:TYR:HE1	1:A:222:LYS:HZ1	1.22	0.86
1:A:35:ASN:O	1:A:35:ASN:OD1	1.98	0.82
1:A:260:GLN:HG2	1:A:299:LEU:HD11	0.83	0.79
1:A:299:LEU:HG	1:A:300:ARG:N	2.01	0.74
1:A:69:ILE:O	1:A:73:ASN:ND2	2.20	0.73
1:A:91:ILE:HA	1:A:282:ARG:HA	1.71	0.72
1:A:102:GLN:HG2	1:A:276:LYS:HA	1.72	0.71
1:A:215:MET:HA	1:A:218:LEU:HG	1.74	0.70
1:A:299:LEU:O	1:A:301:MET:N	2.24	0.70
1:A:96:ILE:HG13	1:A:279:LYS:HE2	1.73	0.70
1:A:299:LEU:HG	1:A:300:ARG:H	1.56	0.69
1:A:203:THR:O	1:A:207:PHE:HB3	1.93	0.69
1:A:197:ASP:HB2	1:A:200:ILE:HG13	1.79	0.62
1:A:132:LEU:O	1:A:136:ILE:HG12	1.99	0.62
1:A:202:PHE:HZ	1:A:300:ARG:CZ	2.14	0.60
1:A:288:PHE:O	1:A:292:TYR:N	2.28	0.60
1:A:299:LEU:O	1:A:301:MET:HG3	2.02	0.59
1:A:42:GLY:HA2	1:A:46:MET:HB2	1.85	0.59
1:A:99:LYS:O	1:A:99:LYS:HG2	2.01	0.59
1:A:108:TYR:OH	1:A:268:SER:OG	2.21	0.58
1:A:88:LEU:HD12	1:A:91:ILE:HG13	1.86	0.57
1:A:199:VAL:HA	1:A:202:PHE:HB3	1.85	0.57
1:A:79:ILE:HG23	1:A:292:TYR:HE2	1.69	0.57
1:A:23:PHE:O	1:A:27:SER:HB2	2.04	0.57
1:A:79:ILE:HD11	1:A:263:VAL:HG12	1.86	0.56
1:A:238:ILE:O	1:A:242:GLY:N	2.38	0.56
1:A:9:PHE:HE1	1:A:182:PHE:CE2	2.23	0.56
1:A:44:THR:OG1	1:A:45:VAL:N	2.37	0.56
1:A:260:GLN:CG	1:A:299:LEU:CD1	2.56	0.56
1:A:30:ILE:O	1:A:34:PHE:HB2	2.05	0.56
1:A:114:PHE:O	1:A:118:ILE:HG23	2.06	0.56
1:A:125:TRP:O	1:A:129:VAL:HG23	2.07	0.55
1:A:197:ASP:HB3	1:A:200:ILE:H	1.71	0.55
1:A:95:ILE:HG12	1:A:226:GLY:HA3	1.89	0.55
1:A:171:GLY:O	1:A:175:VAL:HG23	2.07	0.55
1:A:287:LEU:O	1:A:291:LEU:N	2.29	0.55
1:A:218:LEU:HD12	1:A:219:ALA:N	2.22	0.54
1:A:165:LEU:O	1:A:169:ILE:HG12	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:105:ILE:HG22	1:A:272:TYR:HD1	1.73	0.54
1:A:107:VAL:HG21	1:A:141:TRP:CH2	2.42	0.54
1:A:99:LYS:HA	1:A:102:GLN:OE1	2.08	0.53
1:A:274:PHE:HB2	1:A:280:ILE:HG22	1.91	0.53
1:A:102:GLN:HG2	1:A:276:LYS:CA	2.38	0.53
1:A:188:LYS:HA	1:A:191:LEU:HD23	1.90	0.52
1:A:88:LEU:HA	1:A:91:ILE:HG13	1.90	0.52
1:A:180:GLU:O	1:A:184:ASP:HB2	2.10	0.52
1:A:211:LEU:HB3	1:A:212:PRO:HD3	1.92	0.52
1:A:79:ILE:HD11	1:A:267:MET:HG3	1.90	0.52
1:A:34:PHE:O	1:A:35:ASN:HB3	2.10	0.51
1:A:29:ARG:HE	1:A:164:SER:HB3	1.75	0.51
1:A:218:LEU:HD13	1:A:222:LYS:HE3	1.93	0.51
1:A:23:PHE:HZ	1:A:46:MET:HA	1.75	0.50
1:A:262:ALA:O	1:A:266:ILE:HG12	2.12	0.50
1:A:33:HIS:HB3	1:A:161:VAL:HG13	1.94	0.50
1:A:6:VAL:HA	1:A:189:ILE:HD11	1.93	0.50
1:A:81:ASN:HA	1:A:85:VAL:HB	1.95	0.49
1:A:207:PHE:CD2	1:A:207:PHE:C	2.86	0.48
1:A:109:LEU:HG	1:A:113:ILE:HD11	1.95	0.48
1:A:16:LEU:HD21	1:A:179:ALA:N	2.28	0.48
1:A:199:VAL:O	1:A:203:THR:OG1	2.29	0.48
1:A:86:LEU:HB2	1:A:233:VAL:HG21	1.96	0.48
1:A:213:GLU:O	1:A:217:SER:HB2	2.14	0.48
1:A:200:ILE:O	1:A:204:LEU:HD13	2.15	0.47
1:A:91:ILE:HD13	1:A:282:ARG:O	2.14	0.47
1:A:106:LEU:HB2	1:A:272:TYR:CE1	2.50	0.47
1:A:83:GLY:O	1:A:289:LEU:HD22	2.15	0.47
1:A:101:LEU:HA	1:A:104:ASN:HD22	1.79	0.47
1:A:264:LEU:O	1:A:268:SER:HB2	2.15	0.47
1:A:10:LEU:O	1:A:14:ILE:HD12	2.15	0.47
1:A:26:GLY:O	1:A:30:ILE:HG12	2.15	0.46
1:A:8:TYR:HB3	1:A:181:LEU:O	2.16	0.46
1:A:129:VAL:O	1:A:133:ILE:HG12	2.15	0.46
1:A:29:ARG:HB3	1:A:164:SER:HB3	1.98	0.46
1:A:269:LEU:HD12	1:A:269:LEU:HA	1.66	0.46
1:A:183:VAL:O	1:A:187:LYS:HD3	2.16	0.46
1:A:21:ASP:O	1:A:24:VAL:N	2.49	0.46
1:A:292:TYR:CZ	1:A:296:ILE:HD11	2.52	0.45
1:A:43:ALA:HA	1:A:231:GLY:O	2.17	0.45
1:A:273:LEU:O	1:A:277:TYR:HB2	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:202:PHE:O	1:A:205:VAL:N	2.49	0.45
1:A:29:ARG:CB	1:A:164:SER:HB3	2.48	0.45
1:A:49:GLY:HA2	1:A:52:LEU:HD12	1.99	0.45
1:A:102:GLN:HB3	1:A:276:LYS:HB2	1.98	0.44
1:A:13:LEU:HA	1:A:16:LEU:HD12	1.98	0.44
1:A:48:ILE:O	1:A:52:LEU:N	2.49	0.44
1:A:118:ILE:HD11	1:A:131:LEU:HD21	2.00	0.44
1:A:299:LEU:CG	1:A:300:ARG:N	2.79	0.44
1:A:289:LEU:O	1:A:293:ILE:HG12	2.17	0.44
1:A:109:LEU:O	1:A:113:ILE:HG13	2.18	0.44
1:A:139:LEU:O	1:A:143:VAL:HB	2.18	0.44
1:A:79:ILE:HG23	1:A:292:TYR:CE2	2.52	0.43
1:A:87:GLY:O	1:A:285:GLY:HA3	2.17	0.43
1:A:289:LEU:HD13	1:A:289:LEU:HA	1.89	0.43
1:A:20:SER:OG	1:A:215:MET:HG2	2.18	0.43
1:A:42:GLY:HA2	1:A:46:MET:HG3	2.01	0.43
1:A:69:ILE:HD13	1:A:69:ILE:HA	1.92	0.43
1:A:243:GLY:O	1:A:247:VAL:HG23	2.19	0.43
1:A:190:ALA:HA	1:A:193:LEU:HD12	2.01	0.43
1:A:95:ILE:O	1:A:279:LYS:HD2	2.19	0.43
1:A:202:PHE:CD2	1:A:202:PHE:C	2.93	0.42
1:A:269:LEU:O	1:A:272:TYR:N	2.52	0.42
1:A:298:SER:O	1:A:298:SER:OG	2.30	0.42
1:A:88:LEU:HA	1:A:88:LEU:HD12	1.74	0.42
1:A:76:GLY:HA2	1:A:79:ILE:HG22	2.01	0.42
1:A:289:LEU:O	1:A:293:ILE:N	2.34	0.42
1:A:197:ASP:HB2	1:A:200:ILE:CG1	2.46	0.42
1:A:38:ASN:HD22	1:A:225:LEU:HD22	1.85	0.42
1:A:106:LEU:O	1:A:109:LEU:HB3	2.20	0.41
1:A:202:PHE:CZ	1:A:300:ARG:CZ	2.99	0.41
1:A:101:LEU:O	1:A:104:ASN:N	2.53	0.41
1:A:230:LEU:HA	1:A:230:LEU:HD12	1.91	0.41
1:A:29:ARG:NH2	1:A:163:PHE:HB3	2.36	0.41
1:A:88:LEU:HD13	1:A:289:LEU:CD2	2.50	0.41
1:A:205:VAL:O	1:A:209:THR:HG23	2.21	0.41
1:A:39:PHE:HA	1:A:227:GLY:O	2.21	0.41
1:A:117:VAL:C	1:A:119:GLY:H	2.25	0.40
1:A:299:LEU:CG	1:A:300:ARG:H	2.30	0.40
1:A:120:ILE:H	1:A:120:ILE:HG13	1.62	0.40
1:A:102:GLN:N	1:A:102:GLN:OE1	2.42	0.40
1:A:274:PHE:CB	1:A:280:ILE:HG22	2.51	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	285/320 (89%)	221 (78%)	60 (21%)	4 (1%)	14 58

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	127	ASP
1	A	161	VAL
1	A	300	ARG
1	A	195	ILE

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	231/257 (90%)	183 (79%)	48 (21%)	1 8

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	LEU
1	A	17	TYR
1	A	21	ASP
1	A	27	SER

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Mol	Chain	Res	Type
1	A	28	GLU
1	A	29	ARG
1	A	34	PHE
1	A	36	VAL
1	A	37	SER
1	A	51	SER
1	A	52	LEU
1	A	55	ILE
1	A	56	LEU
1	A	62	SER
1	A	88	LEU
1	A	91	ILE
1	A	92	ILE
1	A	101	LEU
1	A	106	LEU
1	A	108	TYR
1	A	113	ILE
1	A	118	ILE
1	A	127	ASP
1	A	131	LEU
1	A	140	ARG
1	A	142	THR
1	A	143	VAL
1	A	161	VAL
1	A	191	LEU
1	A	197	ASP
1	A	199	VAL
1	A	202	PHE
1	A	203	THR
1	A	207	PHE
1	A	211	LEU
1	A	215	MET
1	A	217	SER
1	A	218	LEU
1	A	245	LEU
1	A	259	VAL
1	A	268	SER
1	A	269	LEU
1	A	273	LEU
1	A	278	SER
1	A	280	ILE
1	A	283	TRP

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Mol	Chain	Res	Type
1	A	300	ARG
1	A	301	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	81	ASN
1	A	104	ASN
1	A	145	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

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 ⓘ

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 [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	289/320 (90%)	-0.35	5 (1%) 73 64	97, 125, 166, 191	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	196	SER	5.5
1	A	198	LYS	5.4
1	A	197	ASP	2.8
1	A	159	PRO	2.5
1	A	199	VAL	2.3

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](#)

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, 95th 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 ²)	Q<0.9
2	CD	A	322	1/1	0.97	0.21	0.47	118,118,118,118	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	CD	A	321	1/1	0.89	0.17	-	136,136,136,136	1

6.5 Other polymers [i](#)

There are no such residues in this entry.