



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 29, 2020 – 09:54 PM BST

PDB ID : 6O73
Title : Crystal structure of apo Csm1-Csm4 cassette
Authors : Jia, N.; Patel, D.J.
Deposited on : 2019-03-07
Resolution : 3.00 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 ⓘ symbol.

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

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.13
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)	:	2.13

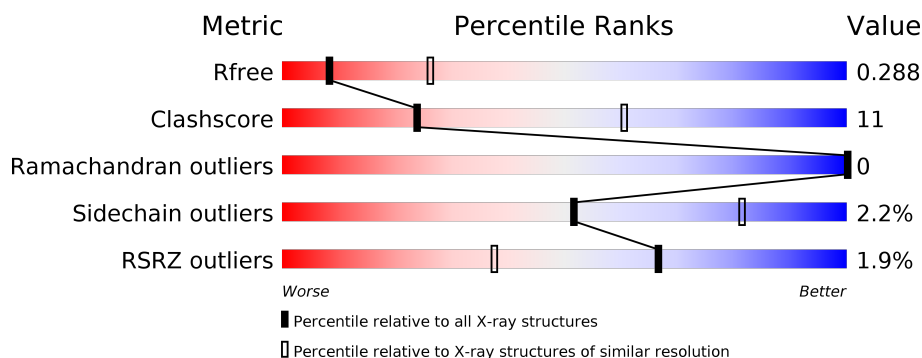
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2092 (3.00-3.00)
Clashscore	141614	2416 (3.00-3.00)
Ramachandran outliers	138981	2333 (3.00-3.00)
Sidechain outliers	138945	2336 (3.00-3.00)
RSRZ outliers	127900	1990 (3.00-3.00)

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 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	791	<div> <div>2%</div> <div> <div></div> <div>66%</div> <div>23%</div> <div>10%</div> </div> </div>
2	B	289	<div> <div>%</div> <div> <div></div> <div>63%</div> <div>20%</div> <div>16%</div> </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7663 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 Csm1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	713	5725	3685	983	1042	15	0	0	0

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-13	MET	-	initiating methionine	UNP B6YWB8
A	-12	GLY	-	expression tag	UNP B6YWB8
A	-11	SER	-	expression tag	UNP B6YWB8
A	-10	SER	-	expression tag	UNP B6YWB8
A	-9	HIS	-	expression tag	UNP B6YWB8
A	-8	HIS	-	expression tag	UNP B6YWB8
A	-7	HIS	-	expression tag	UNP B6YWB8
A	-6	HIS	-	expression tag	UNP B6YWB8
A	-5	HIS	-	expression tag	UNP B6YWB8
A	-4	HIS	-	expression tag	UNP B6YWB8
A	-3	SER	-	expression tag	UNP B6YWB8
A	-2	GLN	-	expression tag	UNP B6YWB8
A	-1	ASP	-	expression tag	UNP B6YWB8
A	0	PRO	-	expression tag	UNP B6YWB8

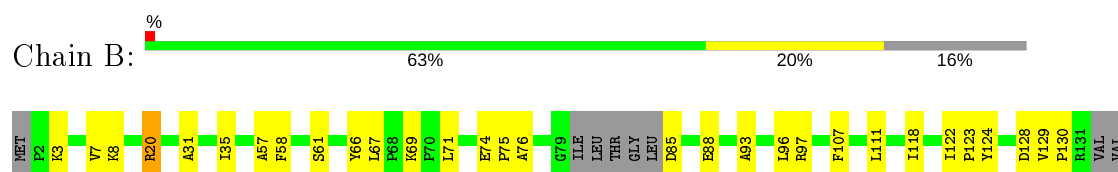
- Molecule 2 is a protein called Csm4.

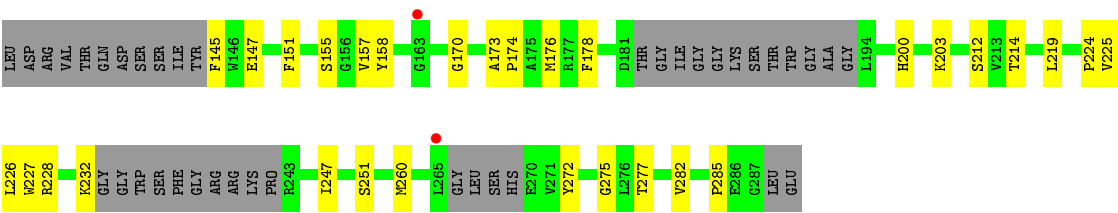
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	242	1937	1262	320	351	4	0	0	0

- Molecule 3 is NICKEL (II) ION (three-letter code: NI) (formula: Ni).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Ni	0	0
			1	1		

- Molecule 1: Csm1





4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	154.28 Å 154.28 Å 182.42 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.72 – 3.00 48.67 – 3.00	Depositor EDS
% Data completeness (in resolution range)	99.9 (48.72-3.00) 100.0 (48.67-3.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.14 (at 3.01 Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.245 , 0.293 0.244 , 0.288	Depositor DCC
R_{free} test set	2608 reflections (9.95%)	wwPDB-VP
Wilson B-factor (Å ²)	98.4	Xtriage
Anisotropy	0.098	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 56.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7663	wwPDB-VP
Average B, all atoms (Å ²)	108.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.44% 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, 0.333 respectively for untwinned datasets, and 0.375, 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: NI

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	0.63	0/5849	0.74	0/7888
2	B	0.60	0/1985	0.75	0/2682
All	All	0.62	0/7834	0.74	0/10570

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

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	5725	0	5730	128	0
2	B	1937	0	1935	45	0
3	A	1	0	0	0	0
All	All	7663	0	7665	170	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 11.

All (170) 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)
2:B:224:PRO:HG3	2:B:247:ILE:HD11	1.51	0.91
1:A:124:ASN:HB2	1:A:615:GLY:HA3	1.59	0.84
1:A:145:ASP:N	1:A:145:ASP:OD2	2.12	0.82
1:A:532:MET:HE3	1:A:537:LYS:HB3	1.69	0.73
1:A:712:ILE:HG12	1:A:728:LEU:CD2	2.18	0.72
1:A:26:TYR:HB2	1:A:33:GLN:HG3	1.72	0.72
1:A:36:ARG:O	1:A:40:ASP:HB2	1.89	0.72
2:B:214:THR:HG22	2:B:277:THR:HB	1.73	0.70
1:A:712:ILE:HD11	1:A:732:LEU:HG	1.74	0.68
2:B:57:ALA:HB1	2:B:157:VAL:HG11	1.76	0.68
2:B:224:PRO:HG3	2:B:247:ILE:CD1	2.24	0.67
1:A:183:LEU:HB3	1:A:198:LEU:HD23	1.75	0.67
2:B:224:PRO:HG2	2:B:227:TRP:CE3	2.32	0.65
1:A:26:TYR:H	1:A:33:GLN:NE2	1.94	0.64
2:B:96:LEU:HD23	2:B:118:ILE:HD11	1.80	0.63
1:A:138:LYS:O	1:A:139:GLU:OE2	2.15	0.63
2:B:85:ASP:HB3	2:B:88:GLU:HB2	1.82	0.61
1:A:179:LEU:HD22	1:A:202:MET:HE1	1.82	0.60
1:A:712:ILE:HG12	1:A:728:LEU:HD21	1.82	0.60
2:B:170:GLY:O	2:B:174:PRO:HG2	2.02	0.60
1:A:203:ARG:NH2	1:A:289:VAL:O	2.35	0.59
1:A:688:ILE:HG22	1:A:699:LEU:HD11	1.84	0.59
2:B:129:VAL:O	2:B:145:PHE:HB2	2.03	0.59
1:A:713:ARG:HG3	1:A:761:VAL:HG12	1.83	0.59
2:B:57:ALA:HB1	2:B:157:VAL:CG1	2.32	0.59
2:B:128:ASP:HB3	2:B:147:GLU:HB3	1.86	0.58
2:B:260:MET:HE1	2:B:272:TYR:HB3	1.85	0.57
1:A:460:PRO:HG2	1:A:477:PRO:HD3	1.85	0.57
1:A:775:VAL:HG12	1:A:775:VAL:O	2.05	0.57
1:A:134:LEU:HB2	1:A:184:THR:HG21	1.87	0.56
1:A:632:PRO:HB2	1:A:634:TYR:CE2	2.40	0.56
1:A:140:LEU:HB2	1:A:561:PHE:CE2	2.41	0.56
1:A:712:ILE:HG23	1:A:728:LEU:HD23	1.88	0.56
1:A:612:ALA:O	1:A:614:THR:O	2.23	0.55
1:A:580:ILE:HD11	1:A:598:GLN:HB3	1.87	0.55
2:B:130:PRO:HA	2:B:145:PHE:CB	2.36	0.55
1:A:345:ARG:NH1	1:A:533:ASP:OD1	2.39	0.55
2:B:76:ALA:HB2	2:B:285:PRO:HG3	1.89	0.55
1:A:575:PRO:HG2	1:A:578:PRO:HA	1.89	0.55
1:A:607:ARG:HG3	1:A:674:TRP:CE2	2.42	0.55
1:A:634:TYR:O	1:A:638:ASP:HB2	2.07	0.55
1:A:6:LEU:HD22	1:A:85:LEU:HD13	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:212:MET:HE2	1:A:227:CYS:HB2	1.89	0.54
1:A:264:ALA:O	1:A:267:GLU:HB2	2.08	0.54
1:A:627:PHE:HB2	1:A:636:MET:HE3	1.89	0.54
1:A:712:ILE:HG21	1:A:729:THR:HG22	1.90	0.53
1:A:78:LEU:O	1:A:80:PRO:HD3	2.08	0.53
1:A:180:GLU:OE2	1:A:547:TYR:OH	2.21	0.53
1:A:260:ARG:HH22	1:A:581:VAL:HG13	1.74	0.53
1:A:520:GLY:HA3	1:A:590:PHE:CZ	2.44	0.52
1:A:646:THR:O	1:A:650:GLU:HG3	2.09	0.52
1:A:522:VAL:HG11	1:A:525:LEU:HD13	1.91	0.52
1:A:668:HIS:HB3	1:A:762:TYR:CZ	2.44	0.52
1:A:660:ARG:HD2	1:A:765:ASP:OD2	2.10	0.51
1:A:712:ILE:HG12	1:A:728:LEU:HD23	1.91	0.51
2:B:225:VAL:HG12	2:B:226:LEU:HG	1.93	0.51
1:A:276:GLU:OE2	1:A:280:ARG:NE	2.43	0.51
2:B:151:PHE:HB3	2:B:155:SER:O	2.10	0.50
1:A:23:ALA:HB2	1:A:154:TYR:HB3	1.94	0.50
1:A:31:SER:O	1:A:57:SER:HB2	2.12	0.50
1:A:260:ARG:NH2	1:A:491:PHE:HB3	2.26	0.50
1:A:419:LEU:HA	1:A:422:LEU:HB3	1.92	0.49
1:A:635:ARG:HD2	2:B:145:PHE:HZ	1.77	0.49
1:A:204:MET:HE3	1:A:299:ILE:HD12	1.94	0.49
2:B:151:PHE:CE1	2:B:157:VAL:HG23	2.47	0.49
2:B:35:ILE:HD11	2:B:178:PHE:CG	2.46	0.49
1:A:212:MET:HE2	1:A:227:CYS:CB	2.42	0.49
1:A:570:SER:HB2	1:A:601:GLU:OE1	2.13	0.49
1:A:139:GLU:OE2	1:A:139:GLU:HA	2.12	0.49
2:B:130:PRO:HA	2:B:145:PHE:HB3	1.94	0.49
1:A:627:PHE:CD2	1:A:636:MET:HG2	2.48	0.49
1:A:687:ARG:HB3	1:A:741:LEU:HD11	1.95	0.48
1:A:557:ILE:O	1:A:578:PRO:HG3	2.13	0.48
2:B:66:TYR:O	2:B:158:TYR:OH	2.21	0.48
1:A:291:ASN:OD1	1:A:296:PHE:HB3	2.14	0.48
1:A:516:GLY:HA3	1:A:596:TRP:HA	1.95	0.48
1:A:607:ARG:HG3	1:A:674:TRP:CD2	2.49	0.47
1:A:204:MET:HE2	1:A:535:PRO:HB3	1.95	0.47
1:A:238:GLY:HA3	1:A:296:PHE:CZ	2.50	0.47
1:A:695:LEU:HD13	1:A:775:VAL:HG21	1.95	0.47
1:A:473:PRO:HA	1:A:482:PHE:CE1	2.49	0.47
1:A:725:TRP:CE3	1:A:761:VAL:HG21	2.49	0.47
1:A:140:LEU:HB2	1:A:561:PHE:CZ	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:477:PRO:HG2	1:A:480:ALA:HB2	1.97	0.47
1:A:302:ASN:O	1:A:302:ASN:ND2	2.38	0.46
1:A:469:ASN:HA	1:A:486:PHE:CD1	2.50	0.46
1:A:433:GLY:O	1:A:451:PHE:HA	2.16	0.46
1:A:41:LEU:HA	1:A:162:TRP:CH2	2.51	0.46
2:B:8:LYS:HD3	2:B:200:HIS:CE1	2.51	0.46
2:B:214:THR:HG21	2:B:277:THR:HG22	1.96	0.46
2:B:74:GLU:OE1	2:B:97:ARG:HG3	2.15	0.46
1:A:319:GLU:HA	1:A:374:ILE:HD11	1.97	0.46
1:A:503:GLU:O	1:A:507:MET:HG3	2.15	0.46
2:B:247:ILE:HG23	2:B:251:SER:CB	2.46	0.46
1:A:38:LEU:HD12	1:A:54:SER:HA	1.98	0.46
1:A:272:ASP:OD1	1:A:434:ARG:NH1	2.49	0.46
1:A:75:ILE:O	1:A:78:LEU:O	2.33	0.46
1:A:435:THR:HG21	1:A:461:VAL:HG12	1.99	0.45
1:A:724:ARG:O	1:A:727:TYR:HB3	2.16	0.45
1:A:272:ASP:OD1	1:A:449:SER:OG	2.33	0.45
1:A:708:LYS:O	1:A:712:ILE:HG13	2.16	0.45
1:A:491:PHE:CE2	1:A:493:LYS:HG2	2.52	0.45
1:A:545:MET:HE1	1:A:610:PHE:HE1	1.82	0.45
1:A:79:SER:OG	1:A:82:ARG:NE	2.48	0.45
2:B:74:GLU:HB2	2:B:75:PRO:HD3	1.98	0.45
1:A:179:LEU:HB3	1:A:202:MET:HE1	1.99	0.45
1:A:621:SER:HA	1:A:654:ARG:O	2.17	0.45
1:A:95:TRP:CE3	1:A:95:TRP:HA	2.51	0.44
2:B:35:ILE:HD11	2:B:178:PHE:CB	2.48	0.44
1:A:124:ASN:ND2	1:A:127:LYS:HG3	2.33	0.44
1:A:237:GLU:OE2	1:A:295:HIS:NE2	2.50	0.44
1:A:297:MET:HE1	1:A:356:ARG:HH11	1.83	0.44
1:A:709:LEU:HA	1:A:712:ILE:HD12	2.00	0.44
1:A:174:ARG:NH2	1:A:480:ALA:O	2.48	0.44
1:A:627:PHE:HB2	1:A:636:MET:CE	2.48	0.44
1:A:44:ASN:C	1:A:46:GLY:H	2.21	0.44
1:A:250:VAL:HB	1:A:633:ILE:HD11	1.99	0.44
2:B:7:VAL:HG11	2:B:176:MET:HG3	1.98	0.44
2:B:61:SER:HB2	2:B:123:PRO:CG	2.48	0.44
2:B:8:LYS:HD3	2:B:200:HIS:HE1	1.82	0.44
1:A:565:ILE:O	1:A:608:ARG:HD2	2.18	0.43
2:B:71:LEU:HD21	2:B:228:ARG:HH21	1.83	0.43
1:A:596:TRP:HB2	1:A:626:TYR:CD1	2.53	0.43
1:A:764:VAL:CG1	1:A:768:LEU:HG	2.47	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:11:GLY:HA2	1:A:97:VAL:HG13	1.99	0.43
1:A:174:ARG:C	1:A:177:PRO:HD2	2.39	0.43
1:A:176:LEU:HB2	1:A:177:PRO:HD3	1.99	0.43
1:A:503:GLU:HG3	1:A:629:GLU:OE1	2.18	0.43
1:A:712:ILE:HG21	1:A:729:THR:CG2	2.48	0.43
1:A:80:PRO:HB2	1:A:85:LEU:O	2.19	0.43
1:A:52:LEU:HD22	1:A:83:PHE:CD2	2.53	0.43
1:A:258:TYR:O	1:A:262:ARG:HG2	2.19	0.42
2:B:69:LYS:NZ	2:B:96:LEU:O	2.43	0.42
1:A:141:PRO:HD3	1:A:552:TYR:CZ	2.54	0.42
1:A:771:VAL:O	1:A:775:VAL:HG23	2.18	0.42
2:B:122:ILE:HA	2:B:123:PRO:HD3	1.90	0.42
2:B:219:LEU:HG	2:B:275:GLY:HA3	2.01	0.42
1:A:330:TYR:OH	1:A:366:ARG:HD2	2.19	0.42
2:B:93:ALA:HA	2:B:96:LEU:HD12	2.02	0.42
1:A:237:GLU:HB2	1:A:359:LEU:HD21	2.01	0.42
2:B:3:LYS:HB3	2:B:203:LYS:HG2	2.02	0.42
2:B:61:SER:HB2	2:B:123:PRO:HG3	2.02	0.42
1:A:5:GLU:O	1:A:9:LEU:HG	2.20	0.41
2:B:31:ALA:O	2:B:35:ILE:HG13	2.20	0.41
1:A:465:ILE:HG12	1:A:480:ALA:HB1	2.01	0.41
2:B:20:ARG:HD3	2:B:232:LYS:HE3	2.02	0.41
2:B:58:PHE:HB2	2:B:158:TYR:CE1	2.54	0.41
2:B:173:ALA:HB3	2:B:174:PRO:HD3	2.02	0.41
1:A:244:GLN:HG2	1:A:244:GLN:H	1.68	0.41
1:A:433:GLY:HA3	1:A:461:VAL:O	2.20	0.41
1:A:101:ASP:HA	1:A:205:THR:CG2	2.50	0.41
1:A:635:ARG:HD2	2:B:145:PHE:CZ	2.55	0.41
1:A:257:LYS:O	1:A:260:ARG:HG3	2.21	0.41
1:A:699:LEU:O	1:A:702:LYS:N	2.39	0.41
1:A:122:VAL:HA	1:A:195:ILE:HG23	2.03	0.41
1:A:588:ASP:OD2	1:A:589:ASP:HB2	2.21	0.41
2:B:67:LEU:HD13	2:B:107:PHE:CD2	2.56	0.41
1:A:631:THR:HG21	1:A:635:ARG:HH11	1.85	0.41
1:A:16:ILE:HG12	1:A:16:ILE:O	2.20	0.41
1:A:180:GLU:O	1:A:184:THR:HB	2.21	0.41
1:A:596:TRP:HB2	1:A:626:TYR:HD1	1.86	0.40
1:A:49:GLU:HG3	1:A:83:PHE:CE1	2.56	0.40
1:A:178:VAL:O	1:A:182:TYR:HD2	2.04	0.40
1:A:131:TRP:CZ2	1:A:157:LEU:HD13	2.56	0.40
1:A:327:SER:HA	1:A:371:PHE:CE2	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:310:LEU:HB3	1:A:335:TRP:CZ3	2.55	0.40
1:A:378:PHE:CE2	2:B:224:PRO:HD2	2.56	0.40
1:A:47:ARG:HD3	1:A:47:ARG:HA	1.91	0.40
1:A:525:LEU:CD2	1:A:588:ASP:HB3	2.52	0.40
2:B:111:LEU:HA	2:B:282:VAL:HG23	2.03	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	697/791 (88%)	639 (92%)	58 (8%)	0	100	100
2	B	230/289 (80%)	209 (91%)	21 (9%)	0	100	100
All	All	927/1080 (86%)	848 (92%)	79 (8%)	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	600/665 (90%)	585 (98%)	15 (2%)	47	79
2	B	204/240 (85%)	201 (98%)	3 (2%)	65	87
All	All	804/905 (89%)	786 (98%)	18 (2%)	52	81

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	ASP
1	A	95	TRP
1	A	102	ASN
1	A	105	SER
1	A	145	ASP
1	A	203	ARG
1	A	226	ARG
1	A	302	ASN
1	A	345	ARG
1	A	429	LEU
1	A	501	THR
1	A	588	ASP
1	A	638	ASP
1	A	641	SER
1	A	733	LEU
2	B	20	ARG
2	B	124	TYR
2	B	212	SER

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

Mol	Chain	Res	Type
2	B	200	HIS

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 monosaccharides in this entry.

5.6 Ligand geometry

Of 1 ligands modelled in this entry, 1 is 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 ⓘ

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, 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	713/791 (90%)	-0.11	16 (2%) 62 33	66, 104, 149, 204	0
2	B	242/289 (83%)	-0.02	2 (0%) 86 65	71, 102, 147, 172	0
All	All	955/1080 (88%)	-0.08	18 (1%) 66 37	66, 104, 149, 204	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	425	ASN	5.9
1	A	379	GLU	4.4
2	B	265	LEU	4.0
1	A	417	ASN	3.8
1	A	423	GLY	3.6
1	A	424	GLY	3.5
1	A	421	SER	3.1
1	A	216	GLY	2.6
1	A	228	ARG	2.6
1	A	341	ARG	2.4
1	A	56	PHE	2.4
2	B	163	GLY	2.3
1	A	258	TYR	2.3
1	A	54	SER	2.2
1	A	250	VAL	2.1
1	A	633	ILE	2.1
1	A	91	LEU	2.0
1	A	426	LEU	2.0

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 [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, 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	B-factors(\AA^2)	Q<0.9
3	NI	A	801	1/1	0.88	0.07	153,153,153,153	0

6.5 Other polymers [i](#)

There are no such residues in this entry.