Full wwPDB X-ray Structure Validation Report (i)

Aug 29, 2023 – 04:44 AM EDT

PDB ID	:	3OB6
Title	:	Structure of AdiC(N101A) in the open-to-out Arg+ bound conformation
Authors	:	Carpena, X.; Kowalczyk, L.; Ratera, M.; Valencia, E.; Vazquez-lbar, J.L.; Fita, I.; Palacin, M.
Deposited on	:	2010-08-06
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 (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:

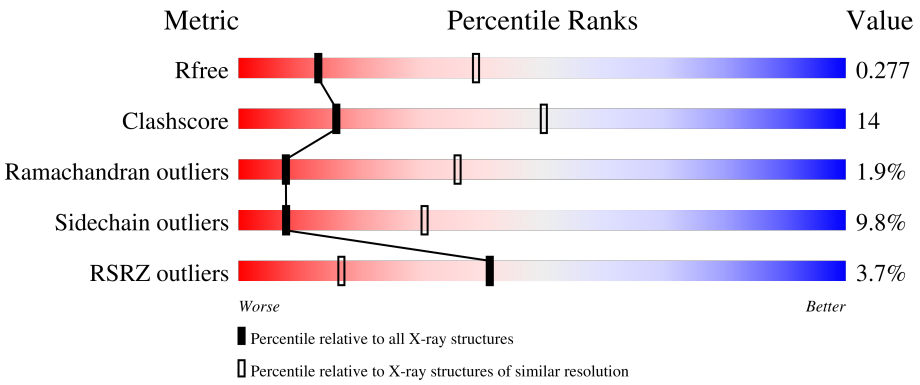
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, 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 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 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	A	445	3% 66% 26% 5% •
1	B	445	4% 66% 27% • •

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	ARG	B	450	-	-	-	X

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 6435 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.

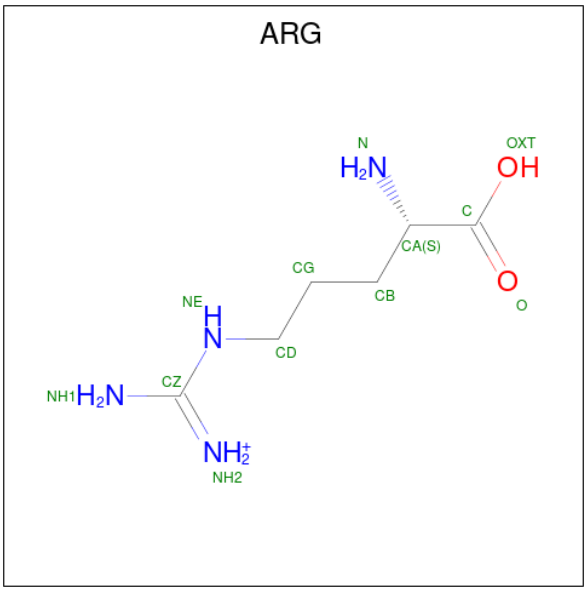
- Molecule 1 is a protein called AdiC arginine:agmatine antiporter.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	430	3181	2112	508	539	22	0	0	0
1	B	436	3230	2143	516	549	22	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	101	ALA	ASN	engineered mutation	UNP C5WBZ6
B	101	ALA	ASN	engineered mutation	UNP C5WBZ6

- Molecule 2 is ARGININE (three-letter code: ARG) (formula: $C_6H_{15}N_4O_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
2	A	1	12	6	4	2	0	0

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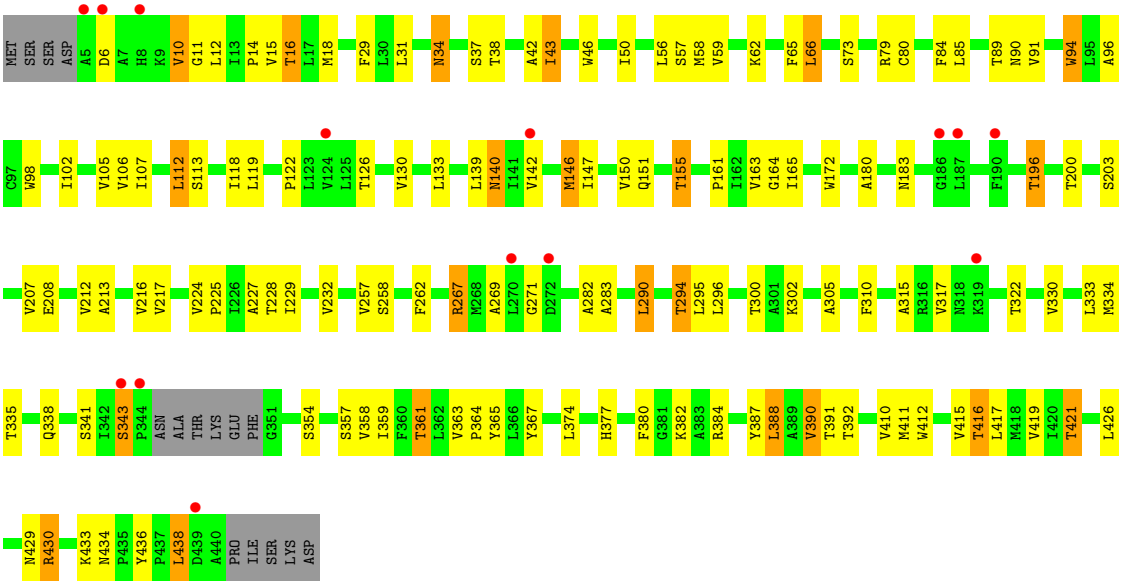
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
2	B	1	12	6	4	2	0	0

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.

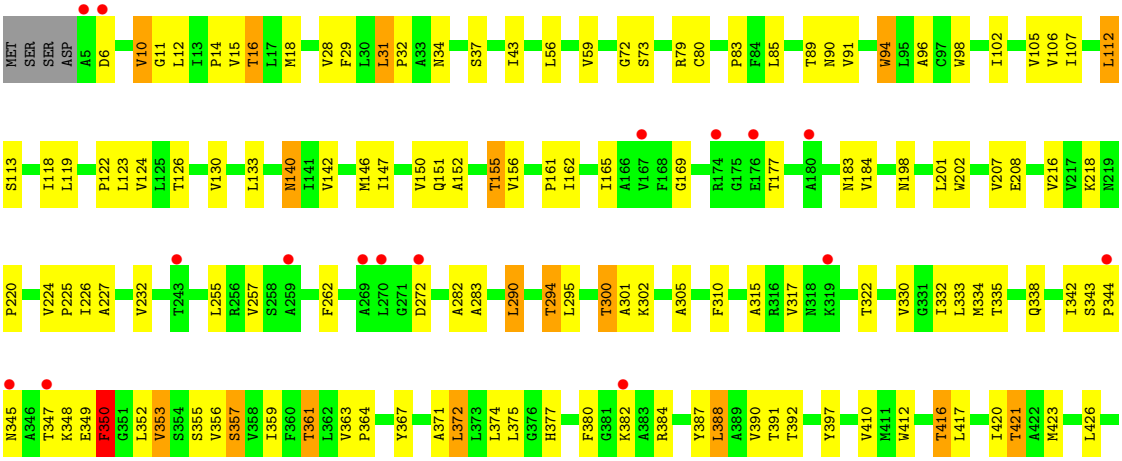
• Molecule 1: AdiC arginine:agmatine antiporter

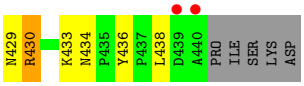
Chain A:



• Molecule 1: AdiC arginine:agmatine antiporter

Chain B:




N429
R430
K433
N434
P435
Y436
P437
L438
D439
A440
PRO
ILE
SER
LYS
ASP

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	86.07Å 77.20Å 104.45Å 90.00° 106.79° 90.00°	Depositor
Resolution (Å)	25.00 – 3.00 24.92 – 3.00	Depositor EDS
% Data completeness (in resolution range)	93.0 (25.00-3.00) 93.0 (24.92-3.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.57 (at 2.99Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.243 , 0.272 0.246 , 0.277	Depositor DCC
R_{free} test set	1255 reflections (5.10%)	wwPDB-VP
Wilson B-factor (Å^2)	99.9	Xtriage
Anisotropy	0.069	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.29 , 64.4	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	6435	wwPDB-VP
Average B, all atoms (Å^2)	93.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.94% 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)

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	0.65	0/3264	0.69	0/4467
1	B	0.64	0/3315	0.70	0/4537
All	All	0.64	0/6579	0.69	0/9004

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	3181	0	3300	95	0
1	B	3230	0	3347	107	0
2	A	12	0	12	0	0
2	B	12	0	12	0	0
All	All	6435	0	6671	190	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 14.

All (190) 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:85:LEU:HD21	1:B:85:LEU:HD21	1.21	1.11

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:11:GLY:O	1:B:15:VAL:HG23	1.66	0.95
1:B:349:GLU:HA	1:B:350:PHE:HB2	1.46	0.94
1:A:140:ASN:ND2	1:A:147:ILE:HD13	1.92	0.85
1:A:12:LEU:O	1:A:16:THR:HG22	1.77	0.84
1:B:140:ASN:ND2	1:B:147:ILE:HD13	1.93	0.84
1:B:151:GLN:O	1:B:155:THR:HG23	1.77	0.82
1:B:12:LEU:O	1:B:16:THR:HG22	1.80	0.82
1:A:85:LEU:CD2	1:B:85:LEU:HD21	2.10	0.80
1:A:11:GLY:O	1:A:15:VAL:HG23	1.82	0.80
1:A:438:LEU:HD12	1:B:79:ARG:NH1	1.97	0.79
1:A:85:LEU:HD21	1:B:85:LEU:CD2	2.08	0.79
1:A:151:GLN:O	1:A:155:THR:HG23	1.83	0.79
1:A:412:TRP:O	1:A:416:THR:HG23	1.82	0.78
1:A:436:TYR:OH	1:B:433:LYS:NZ	2.17	0.76
1:B:140:ASN:HD21	1:B:147:ILE:HD13	1.50	0.76
1:A:96:ALA:HB1	1:A:361:THR:HG23	1.69	0.74
1:B:349:GLU:HA	1:B:350:PHE:CB	2.20	0.72
1:B:412:TRP:O	1:B:416:THR:HG23	1.90	0.72
1:A:417:LEU:O	1:A:421:THR:HG23	1.91	0.69
1:B:133:LEU:HD21	1:B:334:MET:HE3	1.74	0.68
1:A:140:ASN:HD21	1:A:147:ILE:HD13	1.57	0.68
1:B:98:TRP:CZ2	1:B:333:LEU:HD13	2.29	0.68
1:A:34:ASN:HD22	1:A:34:ASN:N	1.92	0.67
1:A:433:LYS:NZ	1:B:436:TYR:OH	2.28	0.67
1:B:207:VAL:HG12	1:B:232:VAL:CG2	2.26	0.65
1:B:349:GLU:CA	1:B:350:PHE:HB2	2.25	0.64
1:B:130:VAL:HG22	1:B:335:THR:HG23	1.81	0.62
1:A:213:ALA:O	1:A:217:VAL:HG23	2.00	0.61
1:B:388:LEU:O	1:B:392:THR:HG23	2.00	0.61
1:A:426:LEU:HD21	1:B:392:THR:HG22	1.83	0.60
1:B:102:ILE:HD11	1:B:334:MET:HA	1.83	0.60
1:B:140:ASN:ND2	1:B:147:ILE:HG21	2.17	0.60
1:A:387:TYR:O	1:A:391:THR:HG23	2.01	0.60
1:B:96:ALA:HB1	1:B:361:THR:HG23	1.83	0.60
1:B:417:LEU:O	1:B:421:THR:HG23	2.01	0.59
1:B:122:PRO:O	1:B:126:THR:HG23	2.02	0.59
1:B:161:PRO:O	1:B:165:ILE:HD12	2.03	0.58
1:A:267:ARG:HA	1:A:271:GLY:HA2	1.86	0.58
1:A:65:PHE:HD1	1:A:66:LEU:HD13	1.68	0.58
1:A:388:LEU:O	1:A:392:THR:HG23	2.04	0.57
1:A:112:LEU:HD13	1:A:283:ALA:CB	2.34	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:12:LEU:HD21	1:B:226:ILE:CG2	2.35	0.57
1:A:133:LEU:HD21	1:A:334:MET:HE3	1.85	0.56
1:B:16:THR:HB	1:B:227:ALA:HA	1.86	0.56
1:A:377:HIS:HA	1:A:380:PHE:CD2	2.40	0.56
1:A:330:VAL:HG12	1:A:334:MET:CE	2.36	0.56
1:B:420:ILE:HA	1:B:423:MET:HE3	1.88	0.56
1:B:89:THR:HG23	1:B:364:PRO:HG3	1.87	0.56
1:A:12:LEU:C	1:A:12:LEU:HD13	2.26	0.55
1:B:147:ILE:CD1	1:B:295:LEU:HB2	2.36	0.55
1:A:290:LEU:O	1:A:290:LEU:HD23	2.06	0.55
1:A:65:PHE:CD1	1:A:66:LEU:HD13	2.41	0.55
1:B:12:LEU:HD21	1:B:226:ILE:HG21	1.89	0.55
1:A:94:TRP:CH2	1:A:330:VAL:HG22	2.42	0.54
1:B:208:GLU:OE1	1:B:208:GLU:N	2.29	0.54
1:A:392:THR:HG22	1:B:426:LEU:HD21	1.88	0.54
1:B:85:LEU:HD12	1:B:367:TYR:CE1	2.43	0.54
1:B:377:HIS:HA	1:B:380:PHE:CD2	2.42	0.54
1:B:387:TYR:O	1:B:391:THR:HG23	2.08	0.54
1:B:56:LEU:HD23	1:B:207:VAL:HG21	1.90	0.54
1:B:59:VAL:HG22	1:B:391:THR:HG22	1.89	0.53
1:B:112:LEU:HD13	1:B:283:ALA:CB	2.39	0.53
1:B:106:VAL:HG12	1:B:107:ILE:N	2.22	0.53
1:B:165:ILE:O	1:B:169:GLY:N	2.38	0.53
1:A:96:ALA:CB	1:A:361:THR:HG23	2.37	0.53
1:B:429:ASN:O	1:B:430:ARG:HB2	2.10	0.52
1:B:224:VAL:HB	1:B:225:PRO:HD3	1.91	0.52
1:B:29:PHE:HA	1:B:262:PHE:CE2	2.45	0.52
1:B:359:ILE:HD13	1:B:410:VAL:HG22	1.92	0.52
1:A:290:LEU:O	1:A:294:THR:CG2	2.58	0.52
1:B:34:ASN:HD22	1:B:34:ASN:N	2.08	0.52
1:A:34:ASN:N	1:A:34:ASN:ND2	2.57	0.51
1:A:290:LEU:CD2	1:A:290:LEU:C	2.79	0.51
1:B:352:LEU:O	1:B:355:SER:N	2.43	0.51
1:B:207:VAL:HG12	1:B:232:VAL:HG21	1.91	0.51
1:A:106:VAL:HG12	1:A:107:ILE:N	2.26	0.51
1:A:140:ASN:CG	1:A:147:ILE:HD13	2.30	0.51
1:A:207:VAL:HG12	1:A:232:VAL:CG2	2.41	0.50
1:B:330:VAL:HG12	1:B:334:MET:CE	2.41	0.50
1:A:113:SER:HB3	1:A:119:LEU:HD12	1.93	0.50
1:A:165:ILE:HD13	1:A:262:PHE:HD1	1.76	0.50
1:B:80:CYS:SG	1:B:371:ALA:HA	2.51	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:80:CYS:SG	1:B:374:LEU:HB2	2.52	0.50
1:A:359:ILE:HD13	1:A:410:VAL:HG22	1.93	0.50
1:B:363:VAL:HG12	1:B:364:PRO:HD3	1.92	0.49
1:A:122:PRO:O	1:A:126:THR:HG23	2.13	0.49
1:A:290:LEU:O	1:A:294:THR:HG23	2.12	0.49
1:B:146:MET:O	1:B:150:VAL:HG23	2.13	0.49
1:A:305:ALA:CB	1:A:315:ALA:HB2	2.42	0.49
1:A:43:ILE:O	1:A:43:ILE:HD13	2.13	0.49
1:A:147:ILE:CD1	1:A:295:LEU:HB2	2.43	0.49
1:A:85:LEU:HD12	1:A:367:TYR:CE1	2.48	0.48
1:B:80:CYS:SG	1:B:374:LEU:HD12	2.53	0.48
1:B:372:LEU:HD22	1:B:372:LEU:O	2.13	0.48
1:B:353:VAL:CG1	1:B:353:VAL:O	2.61	0.48
1:A:46:TRP:O	1:A:50:ILE:HD13	2.13	0.48
1:A:429:ASN:O	1:A:430:ARG:HB2	2.12	0.48
1:B:142:VAL:CG2	1:B:146:MET:HG2	2.43	0.48
1:B:152:ALA:O	1:B:156:VAL:HG12	2.14	0.48
1:B:12:LEU:C	1:B:12:LEU:HD13	2.34	0.48
1:A:56:LEU:HD23	1:A:207:VAL:HG21	1.96	0.47
1:B:353:VAL:O	1:B:353:VAL:HG13	2.14	0.47
1:B:147:ILE:CD1	1:B:295:LEU:HD22	2.44	0.47
1:A:172:TRP:CZ2	1:A:269:ALA:HA	2.50	0.47
1:A:80:CYS:SG	1:A:374:LEU:HD12	2.54	0.47
1:B:130:VAL:CG2	1:B:335:THR:HG23	2.45	0.47
1:A:12:LEU:O	1:A:16:THR:CG2	2.58	0.47
1:B:96:ALA:CB	1:B:361:THR:HG23	2.45	0.47
1:B:290:LEU:O	1:B:294:THR:CG2	2.63	0.46
1:A:62:LYS:HG3	1:A:66:LEU:HD22	1.96	0.46
1:B:133:LEU:HD21	1:B:334:MET:CE	2.43	0.46
1:B:113:SER:HB3	1:B:119:LEU:HD12	1.97	0.46
1:A:90:ASN:O	1:A:91:VAL:C	2.54	0.46
1:A:354:SER:O	1:A:358:VAL:HG23	2.15	0.46
1:B:34:ASN:N	1:B:34:ASN:ND2	2.63	0.46
1:A:257:VAL:HG22	1:A:257:VAL:O	2.16	0.46
1:A:290:LEU:HD23	1:A:290:LEU:C	2.36	0.46
1:A:434:ASN:OD1	1:B:80:CYS:HA	2.15	0.46
1:A:16:THR:HB	1:A:227:ALA:HA	1.97	0.46
1:A:56:LEU:HD21	1:A:365:TYR:CD1	2.51	0.46
1:A:161:PRO:O	1:A:165:ILE:HD12	2.16	0.46
1:B:147:ILE:HD11	1:B:295:LEU:HD22	1.98	0.46
1:A:89:THR:HG23	1:A:364:PRO:HG3	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:201:LEU:HD23	1:B:397:TYR:CE1	2.51	0.45
1:A:410:VAL:O	1:A:411:MET:C	2.54	0.45
1:B:290:LEU:O	1:B:294:THR:HG23	2.16	0.45
1:B:83:PRO:HG3	1:B:433:LYS:HE2	1.98	0.45
1:A:142:VAL:CG2	1:A:146:MET:HG2	2.46	0.45
1:A:415:VAL:O	1:A:419:VAL:HG23	2.17	0.45
1:B:184:VAL:HG12	1:B:184:VAL:O	2.15	0.45
1:B:257:VAL:HG22	1:B:257:VAL:O	2.17	0.45
1:A:224:VAL:HB	1:A:225:PRO:HD3	1.99	0.45
1:A:146:MET:O	1:A:150:VAL:HG23	2.17	0.45
1:A:29:PHE:CE2	1:A:282:ALA:HA	2.52	0.44
1:B:10:VAL:HG12	1:B:14:PRO:HB2	1.99	0.44
1:A:140:ASN:ND2	1:A:147:ILE:HG21	2.32	0.44
1:B:94:TRP:CH2	1:B:330:VAL:HG22	2.51	0.44
1:A:330:VAL:HG12	1:A:334:MET:HE1	2.00	0.44
1:B:133:LEU:HD21	1:B:334:MET:SD	2.58	0.44
1:B:140:ASN:CG	1:B:147:ILE:HD13	2.38	0.44
1:A:363:VAL:HG12	1:A:364:PRO:HD3	2.00	0.44
1:B:102:ILE:O	1:B:105:VAL:HG12	2.18	0.44
1:A:10:VAL:HG12	1:A:14:PRO:HB2	1.99	0.43
1:B:28:VAL:HG21	1:B:162:ILE:HD13	2.01	0.43
1:B:347:THR:HG22	1:B:348:LYS:HG3	1.99	0.43
1:A:79:ARG:NH1	1:B:438:LEU:HD23	2.33	0.43
1:A:212:VAL:CG1	1:A:296:LEU:HD22	2.49	0.43
1:B:90:ASN:O	1:B:91:VAL:C	2.56	0.43
1:A:228:THR:HG22	1:A:229:ILE:N	2.34	0.43
1:B:29:PHE:CE2	1:B:282:ALA:HA	2.54	0.43
1:B:356:VAL:HG23	1:B:357:SER:N	2.33	0.43
1:A:80:CYS:HA	1:B:434:ASN:OD1	2.18	0.43
1:A:302:LYS:NZ	1:A:317:VAL:HG11	2.34	0.42
1:B:198:ASN:O	1:B:202:TRP:CD1	2.72	0.42
1:B:342:ILE:O	1:B:343:SER:C	2.56	0.42
1:A:42:ALA:HA	1:A:196:THR:HG21	2.01	0.42
1:A:98:TRP:CZ2	1:A:333:LEU:HD13	2.53	0.42
1:B:207:VAL:HG12	1:B:232:VAL:HG22	2.00	0.42
1:A:85:LEU:CD2	1:B:85:LEU:CD2	2.86	0.42
1:B:302:LYS:NZ	1:B:317:VAL:HG11	2.34	0.42
1:A:57:SER:OG	1:A:58:MET:N	2.52	0.42
1:B:429:ASN:O	1:B:430:ARG:CB	2.67	0.42
1:A:102:ILE:O	1:A:105:VAL:HG12	2.20	0.42
1:A:207:VAL:HG12	1:A:232:VAL:HG22	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:10:VAL:HG23	1:A:216:VAL:HG13	2.02	0.42
1:B:85:LEU:HD12	1:B:367:TYR:CZ	2.55	0.42
1:A:390:VAL:HG12	1:A:391:THR:N	2.35	0.42
1:B:79:ARG:HB2	1:B:375:LEU:HD21	2.00	0.42
1:B:305:ALA:CB	1:B:315:ALA:HB2	2.50	0.42
1:A:163:VAL:O	1:A:164:GLY:C	2.58	0.42
1:B:123:LEU:C	1:B:123:LEU:HD13	2.40	0.42
1:B:372:LEU:HD22	1:B:380:PHE:HZ	1.85	0.42
1:B:420:ILE:HA	1:B:423:MET:CE	2.49	0.42
1:A:112:LEU:HD13	1:A:283:ALA:HB2	2.02	0.41
1:A:208:GLU:OE1	1:A:365:TYR:OH	2.24	0.41
1:A:15:VAL:HG21	1:A:217:VAL:HG13	2.02	0.41
1:B:31:LEU:N	1:B:32:PRO:HD2	2.35	0.41
1:A:139:LEU:O	1:A:142:VAL:HG22	2.20	0.41
1:A:200:THR:O	1:A:203:SER:HB2	2.20	0.41
1:B:377:HIS:HA	1:B:380:PHE:CE2	2.55	0.41
1:B:218:LYS:O	1:B:220:PRO:HD3	2.21	0.41
1:A:94:TRP:CH2	1:A:330:VAL:CG2	3.04	0.41
1:A:130:VAL:HG22	1:A:335:THR:HG23	2.03	0.41
1:B:300:THR:O	1:B:301:ALA:C	2.58	0.40
1:B:123:LEU:HD13	1:B:124:VAL:N	2.37	0.40
1:A:59:VAL:HG22	1:A:391:THR:HG22	2.04	0.40
1:B:216:VAL:O	1:B:216:VAL:HG13	2.20	0.40
1:A:84:PHE:CD2	1:A:85:LEU:HD22	2.57	0.40
1:B:330:VAL:HG12	1:B:334:MET:HE1	2.04	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	426/445 (96%)	392 (92%)	27 (6%)	7 (2%)	9 40

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	434/445 (98%)	388 (89%)	37 (8%)	9 (2%)	7	33
All	All	860/890 (97%)	780 (91%)	64 (7%)	16 (2%)	8	36

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	118	ILE
1	A	341	SER
1	A	343	SER
1	A	430	ARG
1	B	118	ILE
1	B	344	PRO
1	B	430	ARG
1	B	350	PHE
1	A	180	ALA
1	B	272	ASP
1	A	6	ASP
1	B	6	ASP
1	B	345	ASN
1	A	338	GLN
1	B	338	GLN
1	B	72	GLY

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	328/342 (96%)	294 (90%)	34 (10%)	7	27
1	B	333/342 (97%)	302 (91%)	31 (9%)	9	33
All	All	661/684 (97%)	596 (90%)	65 (10%)	8	30

All (65) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	VAL
1	A	16	THR
1	A	18	MET
1	A	31	LEU
1	A	34	ASN
1	A	37	SER
1	A	38	THR
1	A	43	ILE
1	A	66	LEU
1	A	73	SER
1	A	94	TRP
1	A	112	LEU
1	A	140	ASN
1	A	146	MET
1	A	155	THR
1	A	183	ASN
1	A	196	THR
1	A	258	SER
1	A	267	ARG
1	A	290	LEU
1	A	294	THR
1	A	300	THR
1	A	310	PHE
1	A	322	THR
1	A	343	SER
1	A	357	SER
1	A	361	THR
1	A	382	LYS
1	A	384	ARG
1	A	388	LEU
1	A	390	VAL
1	A	416	THR
1	A	421	THR
1	A	438	LEU
1	B	10	VAL
1	B	16	THR
1	B	18	MET
1	B	31	LEU
1	B	37	SER
1	B	43	ILE
1	B	73	SER
1	B	94	TRP
1	B	112	LEU

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Mol	Chain	Res	Type
1	B	140	ASN
1	B	155	THR
1	B	177	THR
1	B	183	ASN
1	B	255	LEU
1	B	290	LEU
1	B	294	THR
1	B	300	THR
1	B	310	PHE
1	B	322	THR
1	B	332	ILE
1	B	350	PHE
1	B	353	VAL
1	B	357	SER
1	B	361	THR
1	B	372	LEU
1	B	382	LYS
1	B	384	ARG
1	B	388	LEU
1	B	390	VAL
1	B	416	THR
1	B	421	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	22	ASN
1	A	34	ASN
1	A	140	ASN
1	A	151	GLN
1	A	183	ASN
1	A	338	GLN
1	B	22	ASN
1	B	34	ASN
1	B	140	ASN
1	B	151	GLN
1	B	183	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)

2 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	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	ARG	A	450	-	10,11,11	0.88	0	11,13,13	1.20	2 (18%)
2	ARG	B	450	-	10,11,11	0.82	1 (10%)	11,13,13	0.93	1 (9%)

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
2	ARG	A	450	-	-	3/11/11/11	-
2	ARG	B	450	-	-	4/11/11/11	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	450	ARG	OXT-C	-2.19	1.23	1.30

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	450	ARG	OXT-C-O	-2.69	117.97	124.09
2	A	450	ARG	OXT-C-CA	2.20	120.88	113.38
2	B	450	ARG	OXT-C-O	-2.20	119.09	124.09

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	450	ARG	NE-CD-CG-CB
2	A	450	ARG	OXT-C-CA-N
2	A	450	ARG	NE-CD-CG-CB
2	A	450	ARG	O-C-CA-N
2	B	450	ARG	O-C-CA-N
2	B	450	ARG	CG-CD-NE-CZ
2	B	450	ARG	OXT-C-CA-N

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, 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	430/445 (96%)	-0.29	14 (3%)	46	20	69, 88, 111, 133	0
1	B	436/445 (97%)	-0.23	18 (4%)	37	14	68, 93, 131, 154	0
All	All	866/890 (97%)	-0.26	32 (3%)	41	17	68, 90, 122, 154	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	440	ALA	5.5
1	A	272	ASP	5.2
1	B	176	GLU	4.7
1	B	270	LEU	4.6
1	B	439	ASP	4.3
1	A	5	ALA	4.1
1	B	167	VAL	4.0
1	B	269	ALA	3.9
1	A	8	HIS	3.5
1	B	5	ALA	3.4
1	B	259	ALA	3.4
1	A	319	LYS	3.4
1	B	319	LYS	3.3
1	A	343	SER	3.1
1	A	439	ASP	2.9
1	A	190	PHE	2.8
1	A	344	PRO	2.8
1	B	345	ASN	2.6
1	A	187	LEU	2.5
1	B	344	PRO	2.5
1	B	347	THR	2.5
1	A	142	VAL	2.5
1	B	174	ARG	2.5
1	B	382	LYS	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	6	ASP	2.4
1	A	186	GLY	2.3
1	A	6	ASP	2.2
1	B	272	ASP	2.2
1	B	243	THR	2.2
1	A	270	LEU	2.1
1	B	180	ALA	2.0
1	A	124	VAL	2.0

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, 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(Å^2)	Q<0.9
2	ARG	B	450	12/12	0.70	0.42	103,105,105,106	12
2	ARG	A	450	12/12	0.86	0.32	99,100,101,101	12

6.5 Other polymers (i)

There are no such residues in this entry.