



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 1, 2016 – 10:45 AM EDT

PDB ID : 5IO3
Title : Crystal structure of the legionella pneumophila effector protein RavZ - I422
Authors : Kwon, D.H.; Kim, L.; Kim, B.-W.; Hong, S.B.; Song, H.K.
Deposited on : 2016-03-08
Resolution : 2.74 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	:	unknown
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20028320
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)	:	rb-20028320

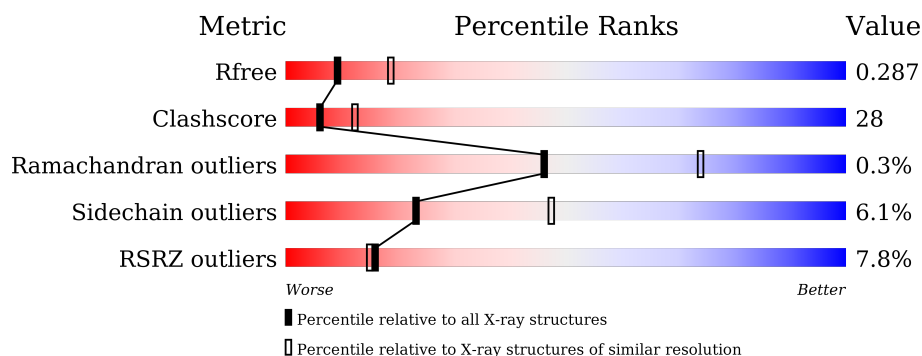
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.74 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	3050 (2.78-2.70)
Clashscore	102246	3424 (2.78-2.70)
Ramachandran outliers	100387	3367 (2.78-2.70)
Sidechain outliers	100360	3368 (2.78-2.70)
RSRZ outliers	91569	3055 (2.78-2.70)

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	502	<div> <div>6%</div> <div>37%</div> <div>37%</div> <div>•</div> <div>23%</div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3071 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 Uncharacterized protein RavZ.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	385	Total	C	N	O	S	0	0	0
			3038	1927	496	606	9			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	258	SER	CYS	engineered mutation	UNP Q5ZUV9

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	33	Total	O	0	0
			33	33		

4 Data and refinement statistics

Property	Value	Source
Space group	I 4 2 2	Depositor
Cell constants a, b, c, α , β , γ	222.04Å 222.04Å 71.30Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.25 – 2.74 42.97 – 2.74	Depositor EDS
% Data completeness (in resolution range)	97.9 (39.25-2.74) 97.9 (42.97-2.74)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.45 (at 2.73Å)	Xtriage
Refinement program	PHENIX (1.10.1 _2155: ???)	Depositor
R, R_{free}	0.224 , 0.287 0.236 , 0.287	Depositor DCC
R_{free} test set	1998 reflections (8.61%)	DCC
Wilson B-factor (Å ²)	51.2	Xtriage
Anisotropy	0.196	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.25 , 49.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.36$, $\langle L^2 \rangle = 0.19$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.83	EDS
Total number of atoms	3071	wwPDB-VP
Average B, all atoms (Å ²)	78.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.45% 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

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	0.55	1/3095 (0.0%)	0.74	3/4181 (0.1%)

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
1	A	0	2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	206	GLU	CB-CG	-5.06	1.42	1.52

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	217	ARG	NE-CZ-NH2	6.73	123.66	120.30
1	A	343	ARG	NE-CZ-NH1	5.96	123.28	120.30
1	A	351	MET	CB-CG-SD	5.02	127.45	112.40

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	118	SER	Peptide
1	A	253	LEU	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	3038	0	3017	171	1
2	A	33	0	0	9	0
All	All	3071	0	3017	171	1

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 28.

All (171) 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:376:ILE:HG12	1:A:382:ILE:HD11	1.54	0.85
1:A:88:VAL:HG11	1:A:111:ILE:HD11	1.62	0.81
1:A:107:ASN:O	2:A:601:HOH:O	1.99	0.79
1:A:273:SER:OG	2:A:602:HOH:O	2.01	0.77
1:A:308:ARG:NH2	1:A:331:GLU:OE1	2.18	0.77
1:A:95:PRO:HB2	1:A:149:PHE:HE1	1.50	0.76
1:A:387:TYR:CE1	1:A:423:LEU:HD13	2.19	0.76
1:A:118:SER:HA	1:A:120:TYR:H	1.50	0.76
1:A:329:LEU:HD11	1:A:373:LEU:HD23	1.69	0.75
1:A:53:PRO:HD3	1:A:309:ASN:HB2	1.68	0.75
1:A:221:ASP:O	1:A:225:THR:HG23	1.88	0.73
1:A:404:LYS:HA	1:A:407:GLN:HB2	1.70	0.73
1:A:410:LEU:HD21	1:A:421:LEU:HD22	1.71	0.72
1:A:287:ILE:HD12	1:A:288:THR:H	1.54	0.72
1:A:176:HIS:CE1	1:A:199:ILE:HD13	2.25	0.71
1:A:199:ILE:HA	1:A:251:ASN:HB2	1.72	0.71
1:A:176:HIS:HE1	1:A:199:ILE:HD13	1.57	0.69
1:A:54:GLU:OE2	1:A:67:ARG:NH1	2.26	0.69
1:A:176:HIS:HB2	1:A:254:THR:HG21	1.76	0.67
1:A:328:GLN:OE1	1:A:330:SER:N	2.21	0.67
1:A:248:ASN:ND2	1:A:251:ASN:OD1	2.28	0.67
1:A:429:ASP:OD2	2:A:604:HOH:O	2.13	0.65
1:A:397:LEU:HA	1:A:402:LEU:HD13	1.79	0.65
1:A:179:LEU:HB2	1:A:262:THR:HG23	1.78	0.65
1:A:204:TYR:CD1	1:A:207:SER:HB2	2.31	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:125:HIS:NE2	2:A:608:HOH:O	2.30	0.64
1:A:85:GLY:O	1:A:105:LYS:NZ	2.30	0.64
1:A:195:VAL:HG13	1:A:244:ARG:HD2	1.79	0.64
1:A:394:VAL:HG13	1:A:418:LEU:HD13	1.80	0.64
1:A:131:ASP:O	1:A:217:ARG:NH1	2.32	0.63
1:A:49:SER:O	1:A:312:VAL:HG23	2.01	0.60
1:A:142:ARG:HA	1:A:145:GLU:HG3	1.83	0.60
1:A:288:THR:HG23	1:A:290:ASN:H	1.66	0.60
1:A:376:ILE:CG1	1:A:382:ILE:HD11	2.29	0.60
1:A:369:LEU:O	1:A:373:LEU:HD12	2.01	0.60
1:A:142:ARG:NH2	1:A:145:GLU:OE1	2.36	0.59
1:A:94:LYS:HE3	1:A:99:ALA:O	2.03	0.59
1:A:124:TYR:HD2	1:A:220:CYS:HB3	1.67	0.59
1:A:93:ASP:HB2	1:A:95:PRO:HG3	1.85	0.58
1:A:111:ILE:HG22	1:A:158:THR:HB	1.85	0.58
1:A:203:THR:O	1:A:207:SER:OG	2.18	0.57
1:A:54:GLU:OE1	1:A:67:ARG:HD3	2.04	0.57
1:A:330:SER:O	1:A:334:VAL:HG13	2.05	0.57
1:A:419:GLN:O	1:A:423:LEU:HB2	2.05	0.56
1:A:247:TYR:OH	1:A:261:TYR:HB3	2.06	0.56
1:A:103:GLN:OE1	2:A:605:HOH:O	2.18	0.56
1:A:409:HIS:O	1:A:413:LEU:HG	2.05	0.56
1:A:136:ASN:HD22	1:A:138:LYS:H	1.53	0.56
1:A:287:ILE:CD1	1:A:288:THR:H	2.18	0.55
1:A:110:GLN:O	1:A:157:ALA:HA	2.07	0.55
1:A:394:VAL:O	1:A:398:MET:HG2	2.06	0.55
1:A:229:GLU:HG3	1:A:239:LEU:HD22	1.89	0.55
1:A:280:LYS:NZ	2:A:609:HOH:O	2.37	0.55
1:A:287:ILE:HD11	1:A:291:SER:HB3	1.90	0.54
1:A:179:LEU:HD23	1:A:196:TYR:HD1	1.71	0.54
1:A:56:SER:OG	1:A:61:LYS:NZ	2.40	0.54
1:A:126:VAL:HG12	1:A:169:ASP:HB2	1.90	0.54
1:A:265:THR:O	1:A:269:VAL:HG23	2.08	0.54
1:A:269:VAL:HG22	1:A:279:VAL:HG11	1.89	0.53
1:A:404:LYS:HD2	1:A:404:LYS:H	1.73	0.53
1:A:172:LYS:HB3	1:A:174:VAL:HG23	1.91	0.52
1:A:125:HIS:CE1	1:A:131:ASP:HB2	2.44	0.52
1:A:382:ILE:HD13	1:A:425:CYS:HB3	1.91	0.52
1:A:179:LEU:HD13	1:A:266:ILE:HG13	1.91	0.52
1:A:156:PRO:HA	1:A:186:PRO:HD3	1.92	0.52
1:A:184:TYR:HB2	1:A:191:LEU:CD2	2.41	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:309:ASN:ND2	1:A:334:VAL:HG21	2.26	0.51
1:A:72:TYR:CE1	1:A:113:LEU:HD21	2.45	0.51
1:A:400:SER:OG	1:A:401:SER:N	2.44	0.51
1:A:104:SER:OG	1:A:274:ALA:O	2.29	0.50
1:A:376:ILE:HD11	1:A:422:VAL:HG23	1.92	0.50
1:A:51:TYR:OH	1:A:69:HIS:ND1	2.39	0.50
1:A:198:SER:O	1:A:248:ASN:HB3	2.11	0.50
1:A:302:GLU:O	1:A:306:LYS:HB2	2.11	0.50
1:A:81:ALA:HB2	1:A:296:GLN:HG2	1.93	0.50
1:A:120:TYR:CD1	1:A:161:LEU:HD21	2.46	0.50
1:A:69:HIS:NE2	1:A:115:ARG:HD2	2.26	0.50
1:A:224:LEU:O	1:A:228:ILE:HG13	2.11	0.50
1:A:368:LEU:HD23	1:A:393:GLU:HG3	1.94	0.50
1:A:150:CYS:HA	1:A:153:ALA:HB2	1.93	0.50
1:A:369:LEU:CD1	1:A:373:LEU:HD11	2.41	0.49
1:A:225:THR:HG22	1:A:242:PHE:CD1	2.47	0.49
1:A:358:SER:OG	1:A:359:LYS:N	2.46	0.49
1:A:330:SER:HB3	1:A:377:LYS:HE3	1.94	0.49
1:A:431:THR:HA	1:A:433:ASP:HB3	1.95	0.49
1:A:54:GLU:O	1:A:306:LYS:HA	2.12	0.49
1:A:368:LEU:HD21	1:A:390:LEU:HA	1.93	0.49
1:A:111:ILE:HD12	1:A:270:ILE:HG23	1.95	0.49
1:A:194:THR:HG23	1:A:243:THR:HG23	1.94	0.49
1:A:410:LEU:HD23	1:A:410:LEU:HA	1.75	0.48
1:A:65:SER:HB3	1:A:118:SER:HB2	1.95	0.48
1:A:127:ASP:OD2	1:A:129:LYS:HG3	2.14	0.48
1:A:179:LEU:HD23	1:A:196:TYR:CD1	2.49	0.48
1:A:176:HIS:HB2	1:A:254:THR:CG2	2.42	0.48
1:A:170:ILE:HG22	1:A:172:LYS:H	1.77	0.48
1:A:328:GLN:OE1	1:A:329:LEU:N	2.46	0.48
1:A:407:GLN:NE2	2:A:607:HOH:O	2.25	0.47
1:A:369:LEU:HD22	1:A:397:LEU:HD13	1.95	0.47
1:A:311:GLY:O	1:A:315:ILE:HD12	2.15	0.47
1:A:142:ARG:O	1:A:146:ILE:HG13	2.15	0.47
1:A:287:ILE:HD12	1:A:288:THR:N	2.26	0.47
1:A:57:TRP:CZ3	1:A:305:ILE:HD12	2.49	0.47
1:A:398:MET:SD	1:A:418:LEU:HD11	2.55	0.47
1:A:125:HIS:HB2	1:A:131:ASP:O	2.15	0.47
1:A:164:VAL:CG1	1:A:175:GLN:HG3	2.45	0.47
1:A:381:GLU:C	1:A:382:ILE:HD12	2.36	0.46
1:A:203:THR:HG23	1:A:250:GLN:HG3	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:418:LEU:HD12	1:A:418:LEU:H	1.79	0.46
1:A:420:LYS:N	2:A:603:HOH:O	2.11	0.46
1:A:72:TYR:O	1:A:319:LEU:HD21	2.15	0.46
1:A:86:ASP:OD1	1:A:86:ASP:N	2.39	0.46
1:A:265:THR:HA	1:A:292:TYR:OH	2.15	0.46
1:A:308:ARG:HD3	1:A:331:GLU:OE2	2.16	0.46
1:A:255:GLU:OE2	1:A:257:ASN:CG	2.54	0.45
1:A:266:ILE:O	1:A:270:ILE:HG13	2.16	0.45
1:A:90:VAL:HG21	1:A:320:THR:O	2.16	0.45
1:A:153:ALA:HB3	1:A:184:TYR:CE2	2.52	0.45
1:A:252:ARG:HG3	1:A:255:GLU:HB2	1.98	0.45
1:A:69:HIS:O	1:A:72:TYR:HB2	2.16	0.45
1:A:68:LEU:HA	1:A:68:LEU:HD12	1.66	0.45
1:A:75:PHE:O	1:A:83:TYR:HA	2.17	0.45
1:A:75:PHE:HD1	1:A:84:LEU:HG	1.80	0.45
1:A:429:ASP:HA	1:A:430:ASP:HA	1.55	0.45
1:A:308:ARG:HH21	1:A:331:GLU:CD	2.19	0.45
1:A:76:PHE:CE2	1:A:322:GLY:HA2	2.52	0.45
1:A:208:LEU:HB3	1:A:213:LYS:HG2	1.98	0.44
1:A:76:PHE:CD2	1:A:322:GLY:HA2	2.53	0.44
1:A:390:LEU:O	1:A:394:VAL:HG23	2.18	0.44
1:A:106:LYS:C	2:A:612:HOH:O	2.55	0.44
1:A:96:LEU:HG	1:A:149:PHE:CD1	2.53	0.44
1:A:248:ASN:H	1:A:282:PRO:HB2	1.82	0.44
1:A:294:THR:H	1:A:297:HIS:CD2	2.36	0.44
1:A:397:LEU:O	1:A:402:LEU:HD22	2.19	0.43
1:A:164:VAL:HG11	1:A:175:GLN:HG3	1.99	0.43
1:A:184:TYR:CZ	1:A:186:PRO:HA	2.53	0.43
1:A:71:LEU:HB3	1:A:263:PHE:CE2	2.54	0.43
1:A:72:TYR:HA	1:A:75:PHE:CE2	2.54	0.43
1:A:172:LYS:HB3	1:A:174:VAL:CG2	2.49	0.43
1:A:222:GLU:O	1:A:226:GLN:HG3	2.18	0.42
1:A:107:ASN:HB2	1:A:274:ALA:O	2.19	0.42
1:A:208:LEU:HD21	1:A:216:TYR:HB2	2.01	0.42
1:A:194:THR:HA	1:A:243:THR:O	2.20	0.42
1:A:299:GLN:O	1:A:303:SER:OG	2.23	0.42
1:A:89:SER:HB2	1:A:105:LYS:HE3	2.01	0.42
1:A:146:ILE:HG21	1:A:146:ILE:HD13	1.83	0.42
1:A:373:LEU:HD12	1:A:373:LEU:H	1.84	0.42
1:A:423:LEU:HB3	1:A:424:PRO:HD3	2.00	0.42
1:A:208:LEU:HB3	1:A:213:LYS:CG	2.50	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:284:SER:O	1:A:284:SER:OG	2.32	0.42
1:A:362:LYS:O	1:A:366:VAL:HG23	2.19	0.42
1:A:170:ILE:HG22	1:A:172:LYS:N	2.35	0.41
1:A:296:GLN:CG	1:A:299:GLN:HE21	2.33	0.41
1:A:169:ASP:N	1:A:169:ASP:OD1	2.53	0.41
1:A:369:LEU:HD11	1:A:373:LEU:HD11	2.02	0.41
1:A:387:TYR:CZ	1:A:423:LEU:HD13	2.52	0.41
1:A:96:LEU:HA	1:A:96:LEU:HD23	1.81	0.41
1:A:257:ASN:OD1	1:A:258:SER:N	2.53	0.41
1:A:59:VAL:HG21	1:A:298:VAL:HG13	2.02	0.41
1:A:78:LYS:HG3	1:A:83:TYR:CE1	2.55	0.41
1:A:225:THR:O	1:A:229:GLU:HB2	2.20	0.41
1:A:196:TYR:CD2	1:A:245:ALA:HB3	2.56	0.41
1:A:409:HIS:O	1:A:413:LEU:N	2.51	0.41
1:A:184:TYR:OH	1:A:186:PRO:HA	2.21	0.41
1:A:196:TYR:HD2	1:A:245:ALA:HB3	1.87	0.40
1:A:432:ILE:HA	1:A:433:ASP:C	2.42	0.40
1:A:304:CYS:O	1:A:308:ARG:HB2	2.22	0.40
1:A:303:SER:HB3	1:A:408:PHE:CE2	2.56	0.40
1:A:418:LEU:O	1:A:422:VAL:HG12	2.21	0.40
1:A:71:LEU:HD21	1:A:305:ILE:HG13	2.03	0.40
1:A:209:SER:HB3	1:A:212:PHE:HD2	1.86	0.40
1:A:382:ILE:HD12	1:A:382:ILE:N	2.36	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:206:GLU:OE1	1:A:343:ARG:NH2[15_545]	2.15	0.05

5.3 Torsion angles

5.3.1 Protein backbone

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	383/502 (76%)	368 (96%)	14 (4%)	1 (0%)	46 74

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	119	ASP

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	346/449 (77%)	325 (94%)	21 (6%)	23 47

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	55	THR
1	A	78	LYS
1	A	89	SER
1	A	90	VAL
1	A	119	ASP
1	A	136	ASN
1	A	137	ASP
1	A	145	GLU
1	A	169	ASP
1	A	181	THR
1	A	183	SER
1	A	204	TYR
1	A	217	ARG
1	A	220	CYS
1	A	308	ARG
1	A	332	PHE
1	A	354	PHE
1	A	364	THR
1	A	375	ASP
1	A	385	SER

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Mol	Chain	Res	Type
1	A	425	CYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	136	ASN
1	A	248	ASN
1	A	275	GLN
1	A	299	GLN

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](#)

There are no ligands in this entry.

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 ⓘ

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	385/502 (76%)	0.47	30 (7%) 16 15	29, 73, 144, 204	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	353	ASN	7.6
1	A	354	PHE	7.2
1	A	348	GLU	6.8
1	A	352	LEU	4.7
1	A	350	SER	4.4
1	A	349	LYS	3.9
1	A	410	LEU	3.8
1	A	286	TYR	3.6
1	A	432	ILE	3.3
1	A	205	SER	3.2
1	A	377	LYS	3.2
1	A	201	ARG	3.2
1	A	289	SER	3.1
1	A	351	MET	3.0
1	A	325	LEU	3.0
1	A	249	HIS	2.8
1	A	82	PHE	2.8
1	A	228	ILE	2.7
1	A	261	TYR	2.6
1	A	287	ILE	2.5
1	A	378	GLY	2.5
1	A	356	GLY	2.5
1	A	161	LEU	2.5
1	A	372	ILE	2.4
1	A	433	ASP	2.4
1	A	84	LEU	2.4
1	A	202	ASP	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	270	ILE	2.1
1	A	301	ILE	2.0
1	A	206	GLU	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 carbohydrates in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.