



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:13 PM GMT

PDB ID : 4O5P  
Title : Crystal structure of an uncharacterized protein from Pseudomonas aeruginosa  
Authors : Hu, H.D.; Gao, Z.Q.; Zhang, H.; Dong, Y.H.  
Deposited on : 2013-12-19  
Resolution : 2.00 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](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

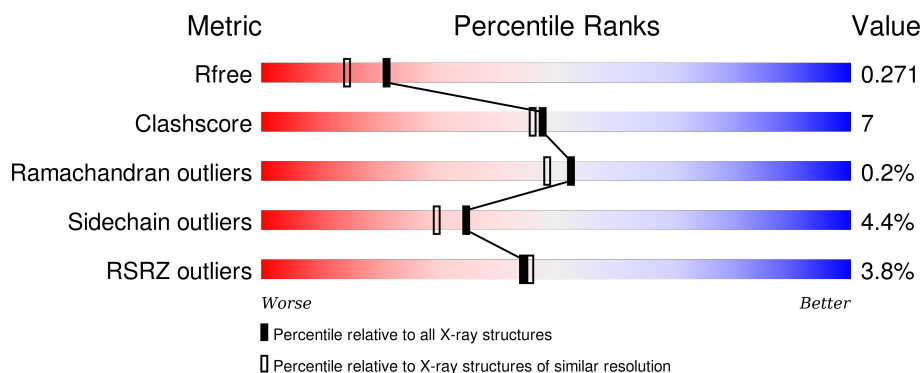
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.00 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))
$R_{free}$	91344	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (2.00-2.00)

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  $\geq 3$ , 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	884	<div> <div>3%</div> <div>70%</div> <div>13%</div> <div>•</div> <div>16%</div> </div>
1	B	884	<div> <div>3%</div> <div>70%</div> <div>13%</div> <div>•</div> <div>16%</div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 12543 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.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	742	Total	C	N	O	S	Se	0	0	0
			5753	3626	1025	1080	8	14			
1	B	741	Total	C	N	O	S	Se	0	1	0
			5765	3636	1025	1082	8	14			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	EXPRESSION TAG	UNP Q9HYV3
A	0	SER	-	EXPRESSION TAG	UNP Q9HYV3
B	-1	GLY	-	EXPRESSION TAG	UNP Q9HYV3
B	0	SER	-	EXPRESSION TAG	UNP Q9HYV3

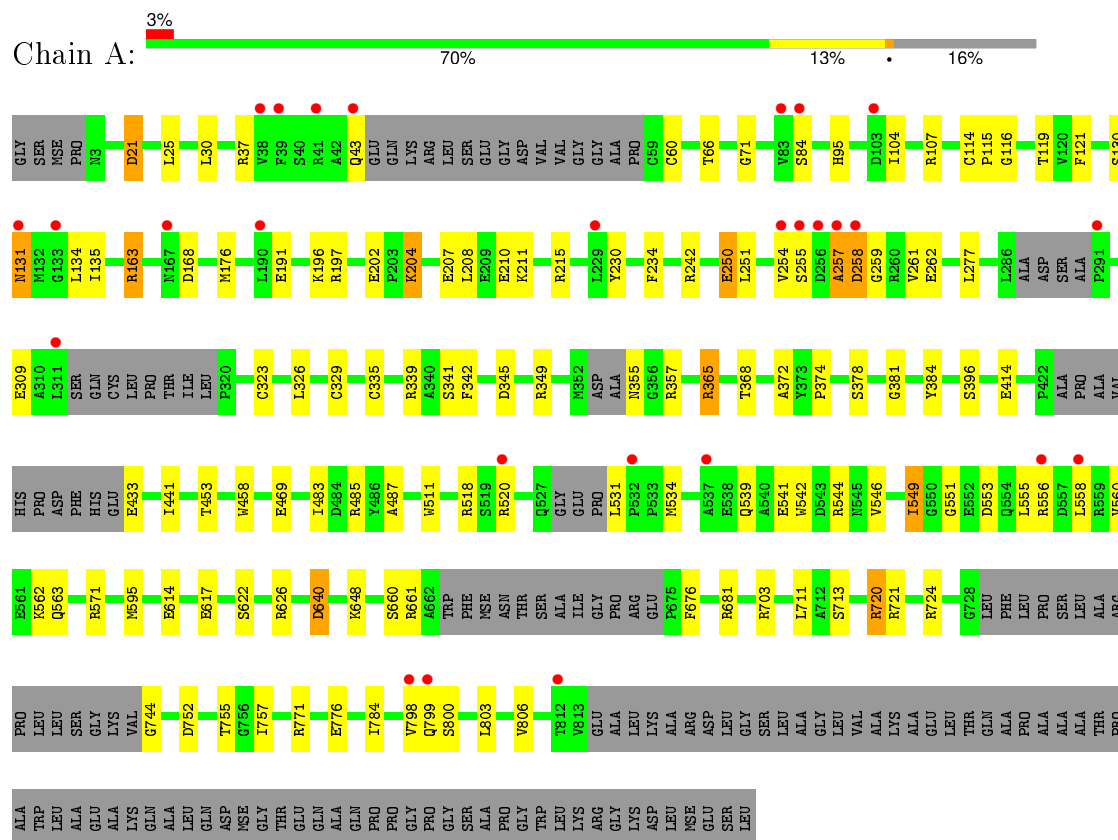
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	520	Total	O	0	0
			520	520		
2	B	505	Total	O	0	0
			505	505		

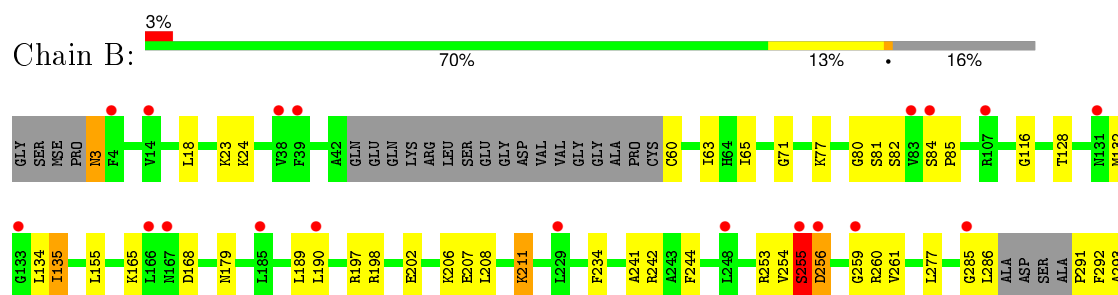
### 3 Residue-property plots

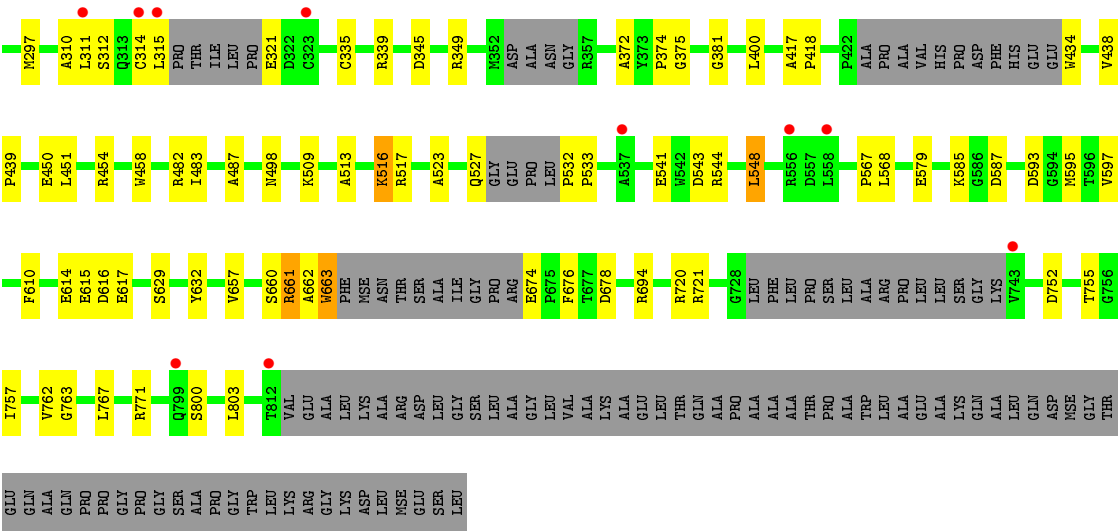
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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: Uncharacterized protein



#### • Molecule 1: Uncharacterized protein





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	61.60Å 81.08Å 92.14Å 97.83° 89.96° 90.01°	Depositor
Resolution (Å)	28.30 – 2.00 28.30 – 2.00	Depositor EDS
% Data completeness (in resolution range)	89.0 (28.30-2.00) 88.0 (28.30-2.00)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.57 (at 2.00Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, $R_{free}$	0.226 , 0.269 0.230 , 0.271	Depositor DCC
$R_{free}$ test set	5464 reflections (5.26%)	DCC
Wilson B-factor (Å <sup>2</sup> )	23.3	Xtriage
Anisotropy	0.300	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 45.8	EDS
Estimated twinning fraction	0.480 for h,-k,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.29$	Xtriage
Outliers	0 of 109303 reflections	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	12543	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.18% 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 [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	0.44	1/5859 (0.0%)	0.60	1/7902 (0.0%)
1	B	0.45	0/5872	0.62	0/7921
All	All	0.44	1/11731 (0.0%)	0.61	1/15823 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	258	ASP	CB-CG	-5.12	1.41	1.51

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	720	ARG	NE-CZ-NH1	5.15	122.88	120.30

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	5753	0	5631	77	0
1	B	5765	0	5639	92	0
2	A	520	0	0	13	0
2	B	505	0	0	16	1
All	All	12543	0	11270	169	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 7.

All (169) 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:B:255:SER:HB3	1:B:261:VAL:HA	1.18	1.17
1:B:255:SER:CB	1:B:261:VAL:HA	1.81	1.09
1:B:255:SER:HB3	1:B:260:ARG:O	1.71	0.90
1:B:255:SER:HA	1:B:256:ASP:O	1.74	0.88
1:A:755:THR:HG23	1:A:757:ILE:H	1.39	0.84
1:B:255:SER:HB3	1:B:261:VAL:CA	2.07	0.80
1:A:721:ARG:NH2	1:A:744:GLY:O	2.16	0.78
1:B:255:SER:HB2	1:B:261:VAL:HG22	1.66	0.78
1:B:255:SER:CB	1:B:261:VAL:HG22	2.19	0.72
1:B:255:SER:OG	1:B:261:VAL:HG13	1.89	0.71
1:B:771:ARG:NH2	2:B:1208:HOH:O	2.22	0.71
1:B:755:THR:HG23	1:B:757:ILE:H	1.56	0.71
1:B:721:ARG:NH1	2:B:1377:HOH:O	2.25	0.69
1:B:253:ARG:NH1	2:B:1172:HOH:O	2.22	0.69
1:A:255:SER:HB3	1:A:261:VAL:HG22	1.76	0.68
1:B:616:ASP:OD2	1:B:694:ARG:HD3	1.94	0.67
1:A:215:ARG:NH1	2:A:1071:HOH:O	2.27	0.67
1:B:197:ARG:NH2	2:B:1111:HOH:O	2.19	0.67
1:A:703:ARG:NH2	1:A:784:ILE:O	2.28	0.66
1:B:579:GLU:OE2	1:B:661:ARG:NH2	2.30	0.64
1:A:595:MSE:HE2	1:A:721:ARG:O	1.97	0.64
1:A:771:ARG:NH2	2:A:1390:HOH:O	2.30	0.63
1:B:450:GLU:OE2	1:B:454:ARG:NE	2.30	0.63
1:B:595:MSE:HE2	1:B:721:ARG:O	1.99	0.63
1:A:485:ARG:NH1	2:A:1138:HOH:O	2.30	0.63
1:A:355:ASN:HB3	1:A:357:ARG:H	1.64	0.63
1:A:131:ASN:ND2	1:A:131:ASN:O	2.33	0.62
1:B:314:CYS:SG	1:B:315:LEU:N	2.73	0.61
1:B:255:SER:OG	1:B:261:VAL:HA	2.01	0.60
1:B:615:GLU:OE2	1:B:721:ARG:NH1	2.35	0.59
1:B:254:VAL:O	1:B:255:SER:OG	2.16	0.59
1:B:80:GLY:HA2	1:B:85:PRO:HB3	1.85	0.58
1:A:309:GLU:OE2	1:A:365:ARG:NH2	2.35	0.58
1:A:617:GLU:OE1	1:A:720:ARG:NH2	2.25	0.58
1:B:135:ILE:H	1:B:285:GLY:HA3	1.70	0.57
1:B:375:GLY:HA2	1:B:657:VAL:HG13	1.86	0.57
1:A:752:ASP:OD2	1:A:755:THR:HG22	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:752:ASP:OD2	1:B:755:THR:HG22	2.05	0.56
1:B:516:LYS:HG2	1:B:517:ARG:N	2.19	0.56
1:A:254:VAL:HG13	1:A:255:SER:H	1.71	0.56
1:B:255:SER:CB	1:B:261:VAL:CA	2.70	0.56
1:B:617:GLU:OE1	1:B:720:ARG:NH2	2.30	0.55
1:B:291:PRO:HB3	1:B:597:VAL:HG11	1.86	0.55
1:B:71:GLY:O	1:B:116:GLY:HA3	2.07	0.55
1:B:242:ARG:HD2	1:B:277:LEU:HD13	1.89	0.55
1:A:511:TRP:CZ2	1:A:549:ILE:HG21	2.41	0.54
1:B:77:LYS:O	1:B:81:SER:HB3	2.07	0.54
1:A:130:SER:O	1:A:563:GLN:HG3	2.08	0.54
1:A:163:ARG:NH1	2:A:1395:HOH:O	2.39	0.54
1:A:250:GLU:HG2	1:A:251:LEU:N	2.24	0.53
1:B:255:SER:HA	1:B:256:ASP:C	2.30	0.53
1:B:482:ARG:NH2	2:B:940:HOH:O	2.27	0.53
1:B:614:GLU:HG3	1:B:720:ARG:NH1	2.24	0.52
1:A:595:MSE:HE3	1:A:720:ARG:NE	2.25	0.52
1:A:541:GLU:HG2	1:A:544:ARG:HH21	1.74	0.52
1:A:30:LEU:HA	1:A:37:ARG:HH22	1.73	0.52
1:A:798:VAL:HG23	1:A:799:GLN:HG3	1.91	0.51
1:B:255:SER:CB	1:B:260:ARG:O	2.52	0.51
1:B:168:ASP:HA	2:B:1392:HOH:O	2.11	0.50
1:A:242:ARG:HD2	1:A:277:LEU:HD13	1.93	0.50
1:B:80:GLY:HA2	1:B:85:PRO:CB	2.40	0.50
1:B:155:LEU:HB3	1:B:208:LEU:HD13	1.94	0.50
1:A:511:TRP:CE2	1:A:549:ILE:HD12	2.47	0.49
1:A:30:LEU:HD22	1:A:37:ARG:NH2	2.27	0.49
1:A:384:TYR:O	2:A:1020:HOH:O	2.20	0.49
1:B:198:ARG:NH2	2:B:1213:HOH:O	2.37	0.49
1:A:676:PHE:O	1:A:681:ARG:NH2	2.45	0.49
1:A:204:LYS:O	1:A:208:LEU:HG	2.13	0.49
1:B:63:ILE:HG22	1:B:65:ILE:HD11	1.95	0.48
1:A:134:LEU:HD12	1:A:134:LEU:HA	1.60	0.48
1:A:43:GLN:O	2:A:1172:HOH:O	2.20	0.48
1:B:614:GLU:HG2	1:B:721:ARG:O	2.13	0.48
1:B:23:LYS:HB3	1:B:24:LYS:HZ2	1.78	0.48
1:B:483:ILE:O	1:B:487:ALA:HB3	2.14	0.48
1:B:509:LYS:HG3	2:B:1217:HOH:O	2.14	0.48
1:B:513:ALA:HA	1:B:516:LYS:HD3	1.95	0.48
1:B:595:MSE:HE3	1:B:720:ARG:NE	2.29	0.48
1:A:378:SER:HB2	1:A:384:TYR:CE2	2.49	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:257:ALA:C	1:A:259:GLY:H	2.17	0.47
1:B:3:ASN:HD22	1:B:3:ASN:N	2.11	0.47
1:B:662:ALA:O	1:B:663:TRP:HE3	1.96	0.47
1:A:339:ARG:NH1	2:A:1128:HOH:O	2.46	0.47
1:B:253:ARG:HH22	1:B:310:ALA:HB2	1.80	0.47
1:B:134:LEU:HA	1:B:134:LEU:HD12	1.51	0.47
1:A:469:GLU:HG3	2:A:1008:HOH:O	2.14	0.47
1:A:341:SER:HG	1:A:342:PHE:HD1	1.59	0.47
1:B:629:SER:O	1:B:632:TYR:HB3	2.15	0.47
1:A:71:GLY:O	1:A:116:GLY:HA3	2.15	0.47
1:B:165:LYS:HE3	1:B:165:LYS:HB2	1.74	0.46
1:B:211:LYS:HD2	2:B:1102:HOH:O	2.14	0.46
1:B:567:PRO:O	1:B:568:LEU:HD23	2.15	0.46
1:B:543:ASP:O	2:B:1402:HOH:O	2.20	0.46
1:B:255:SER:OG	1:B:261:VAL:HG22	2.16	0.46
1:B:234:PHE:CE1	1:B:381:GLY:HA3	2.50	0.46
1:A:483:ILE:O	1:A:487:ALA:HB3	2.16	0.45
1:A:255:SER:OG	1:A:261:VAL:HG13	2.15	0.45
1:A:66:THR:HA	1:A:230:TYR:O	2.16	0.45
1:A:30:LEU:CA	1:A:37:ARG:HH22	2.29	0.45
1:B:763:GLY:H	1:B:767:LEU:HD22	1.82	0.45
1:A:614:GLU:HG2	1:A:721:ARG:O	2.17	0.45
1:A:595:MSE:HE3	1:A:720:ARG:HE	1.80	0.45
1:B:523:ALA:O	1:B:527:GLN:NE2	2.50	0.45
1:B:255:SER:HB3	1:B:260:ARG:C	2.35	0.45
1:B:374:PRO:HD3	1:B:458:TRP:CZ3	2.52	0.45
1:B:674:GLU:N	2:B:1371:HOH:O	2.50	0.44
1:A:323:CYS:SG	2:A:1400:HOH:O	2.61	0.44
1:B:259:GLY:O	1:B:312:SER:HB2	2.18	0.44
1:B:674:GLU:HG2	1:B:676:PHE:O	2.17	0.44
1:B:579:GLU:CD	1:B:661:ARG:HH22	2.20	0.44
1:B:293:ALA:HB3	1:B:610:PHE:HB2	1.99	0.44
1:A:534:MSE:HE3	1:A:539:GLN:OE1	2.17	0.44
1:A:640:ASP:OD2	2:A:1138:HOH:O	2.20	0.44
1:A:30:LEU:HD22	1:A:37:ARG:HH21	1.80	0.44
1:B:335:CYS:SG	1:B:372:ALA:HB1	2.58	0.44
1:A:433:GLU:N	2:A:1309:HOH:O	2.50	0.43
1:B:3:ASN:ND2	1:B:3:ASN:N	2.66	0.43
1:A:339:ARG:CZ	1:A:660:SER:HB2	2.48	0.43
1:A:622:SER:O	1:A:626:ARG:HG3	2.19	0.43
1:B:339:ARG:NE	1:B:660:SER:HB3	2.33	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:532:PRO:HA	1:B:533:PRO:HD2	1.78	0.43
1:A:104:ILE:HD13	1:A:107:ARG:HH12	1.84	0.43
1:A:549:ILE:O	1:A:551:GLY:N	2.44	0.42
1:A:191:GLU:OE1	1:A:196:LYS:NZ	2.50	0.42
1:A:234:PHE:CE1	1:A:381:GLY:HA3	2.53	0.42
1:A:374:PRO:HD3	1:A:458:TRP:CZ3	2.54	0.42
1:A:21:ASP:O	1:A:25:LEU:HG	2.18	0.42
1:B:179:ASN:O	1:B:189:LEU:HD23	2.19	0.42
1:B:291:PRO:HA	1:B:292:PHE:HA	1.84	0.42
1:B:207:GLU:O	1:B:211:LYS:HG2	2.19	0.42
1:A:84:SER:O	1:A:84:SER:OG	2.26	0.42
1:A:335:CYS:SG	1:A:372:ALA:HB1	2.59	0.42
1:A:541:GLU:HG2	1:A:544:ARG:NH2	2.33	0.42
1:A:207:GLU:O	1:A:211:LYS:HG2	2.20	0.42
1:A:711:LEU:HD12	2:A:1164:HOH:O	2.19	0.42
1:B:190:LEU:HD12	1:B:190:LEU:HA	1.83	0.42
1:B:241:ALA:O	1:B:244:PHE:HB3	2.20	0.42
1:A:595:MSE:HE3	1:A:720:ARG:HG3	2.02	0.42
1:A:204:LYS:HD3	1:A:204:LYS:HA	1.83	0.42
1:A:95:HIS:HB3	1:A:441:ILE:HD12	2.01	0.42
1:A:755:THR:HG21	1:A:806:VAL:HG22	2.02	0.41
1:B:755:THR:HG21	1:B:757:ILE:HD12	2.02	0.41
1:B:544:ARG:O	1:B:548:LEU:HD22	2.20	0.41
1:A:800:SER:HB2	1:A:803:LEU:HB2	2.01	0.41
1:A:453:THR:HG21	1:A:648:LYS:HE3	2.01	0.41
1:B:498:ASN:HA	2:B:1240:HOH:O	2.18	0.41
1:B:297:MSE:SE	2:B:1160:HOH:O	2.87	0.41
1:B:417:ALA:HA	1:B:418:PRO:HD3	1.90	0.41
1:A:176:MSE:HG2	1:A:197:ARG:HA	2.01	0.41
1:B:434:TRP:N	2:B:1254:HOH:O	2.52	0.41
1:B:438:VAL:HG13	1:B:439:PRO:HD2	2.03	0.41
1:B:206:LYS:HA	1:B:206:LYS:HD3	1.60	0.41
1:A:562:LYS:NZ	2:A:1271:HOH:O	2.53	0.41
1:B:800:SER:HB3	1:B:803:LEU:HB2	2.02	0.41
1:B:374:PRO:HB3	1:B:458:TRP:CE2	2.56	0.41
1:A:520:ARG:CB	1:A:531:LEU:HD11	2.51	0.41
1:B:254:VAL:O	1:B:255:SER:CB	2.68	0.40
1:B:661:ARG:HD3	1:B:678:ASP:HB2	2.03	0.40
1:A:614:GLU:HG3	1:A:720:ARG:NH1	2.37	0.40
1:A:326:LEU:HB3	1:A:365:ARG:NH2	2.36	0.40
1:A:542:TRP:O	1:A:546:VAL:HG23	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:555:LEU:O	1:A:558:LEU:HB2	2.20	0.40
1:A:114:CYS:HA	1:A:115:PRO:HD2	1.97	0.40
1:B:541:GLU:HG2	2:B:1123:HOH:O	2.20	0.40
1:A:119:THR:HG22	1:A:562:LYS:HG2	2.03	0.40
1:B:400:LEU:HA	1:B:451:LEU:HD11	2.04	0.40
1:B:60:CYS:N	2:B:990:HOH:O	2.54	0.40
1:A:614:GLU:OE2	1:A:720:ARG:HB2	2.22	0.40
1:B:595:MSE:HE3	1:B:720:ARG:HE	1.85	0.40
1:A:329:CYS:HB3	1:A:368:THR:HG22	2.04	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)
2:B:1264:HOH:O	2:B:1387:HOH:O[1_455]	2.17	0.03

## 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	724/884 (82%)	697 (96%)	26 (4%)	1 (0%)	56	53
1	B	724/884 (82%)	702 (97%)	20 (3%)	2 (0%)	46	41
All	All	1448/1768 (82%)	1399 (97%)	46 (3%)	3 (0%)	52	48

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	257	ALA
1	B	255	SER
1	B	256	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	596/687 (87%)	567 (95%)	29 (5%)	31	25
1	B	598/687 (87%)	575 (96%)	23 (4%)	40	36
All	All	1194/1374 (87%)	1142 (96%)	52 (4%)	35	30

All (52) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	21	ASP
1	A	60	CYS
1	A	121	PHE
1	A	131	ASN
1	A	135	ILE
1	A	163	ARG
1	A	168	ASP
1	A	202	GLU
1	A	204	LYS
1	A	210	GLU
1	A	250	GLU
1	A	258	ASP
1	A	262	GLU
1	A	345	ASP
1	A	349	ARG
1	A	365	ARG
1	A	396	SER
1	A	414	GLU
1	A	518	ARG
1	A	549	ILE
1	A	553	ASP
1	A	556	ARG
1	A	560	VAL
1	A	571	ARG
1	A	640	ASP
1	A	661	ARG
1	A	713	SER

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Mol	Chain	Res	Type
1	A	724	ARG
1	A	776	GLU
1	B	3	ASN
1	B	18	LEU
1	B	82	SER
1	B	84	SER
1	B	128	THR
1	B	132	MSE
1	B	135	ILE
1	B	202	GLU
1	B	211	LYS
1	B	255	SER
1	B	286	LEU
1	B	311	LEU
1	B	321	GLU
1	B	345	ASP
1	B	349	ARG
1	B	516	LYS
1	B	548	LEU
1	B	585	LYS
1	B	587	ASP
1	B	593	ASP
1	B	661	ARG
1	B	663	TRP
1	B	762	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	131	ASN
1	A	646	GLN

### 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 [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, 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	728/884 (82%)	0.14	27 (3%)	45 47	11, 24, 42, 65	2 (0%)
1	B	727/884 (82%)	0.15	29 (3%)	42 44	11, 24, 41, 56	0
All	All	1455/1768 (82%)	0.15	56 (3%)	44 45	11, 24, 42, 65	2 (0%)

All (56) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	83	VAL	6.9
1	B	83	VAL	6.2
1	A	311	LEU	5.0
1	B	131	ASN	5.0
1	A	812	THR	4.9
1	B	314	CYS	4.1
1	B	256	ASP	3.8
1	B	255	SER	3.8
1	A	556	ARG	3.7
1	A	84	SER	3.6
1	B	190	LEU	3.6
1	B	84	SER	3.6
1	A	256	ASP	3.6
1	B	311	LEU	3.5
1	A	131	ASN	3.5
1	B	556	ARG	3.4
1	B	743	VAL	3.3
1	A	257	ALA	3.2
1	A	39	PHE	3.2
1	B	812	THR	3.0
1	B	133	GLY	2.9
1	A	190	LEU	2.9
1	B	799	GLN	2.8
1	A	43	GLN	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	798	VAL	2.7
1	B	14	VAL	2.6
1	B	323	CYS	2.6
1	B	39	PHE	2.5
1	A	254	VAL	2.5
1	A	167	ASN	2.5
1	A	133	GLY	2.5
1	A	255	SER	2.5
1	A	537	ALA	2.5
1	A	229	LEU	2.5
1	B	315	LEU	2.5
1	A	558	LEU	2.4
1	B	185	LEU	2.4
1	B	259	GLY	2.4
1	A	291	PRO	2.4
1	B	285	GLY	2.4
1	A	799	GLN	2.4
1	A	41	ARG	2.3
1	A	532	PRO	2.3
1	A	258	ASP	2.3
1	B	558	LEU	2.3
1	B	166	LEU	2.2
1	B	537	ALA	2.2
1	B	107	ARG	2.2
1	B	248	LEU	2.2
1	B	38	VAL	2.2
1	B	229	LEU	2.1
1	A	520	ARG	2.1
1	A	103	ASP	2.1
1	B	4	PHE	2.0
1	A	38	VAL	2.0
1	B	167	ASN	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

There are no ligands in this entry.

## 6.5 Other polymers

There are no such residues in this entry.