



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:48 PM GMT

PDB ID : 4AKX
Title : Structure of the heterodimeric complex ExoU-SpcU from the type III secretion system (T3SS) of *Pseudomonas aeruginosa*
Authors : Gendrin, C.; Contreras-Martel, C.; Bouillot, S.; Elsen, S.; Lemaire, D.; Skoufias, D.A.; Huber, P.; Attree, I.; Dessen, A.
Deposited on : 2012-02-29
Resolution : 2.94 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

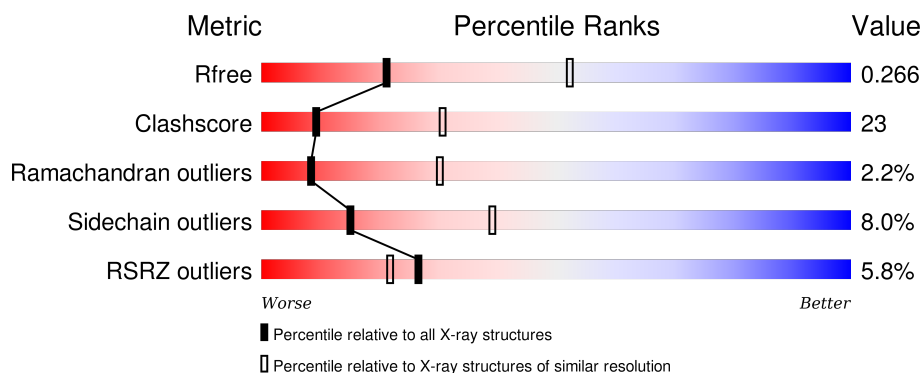
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

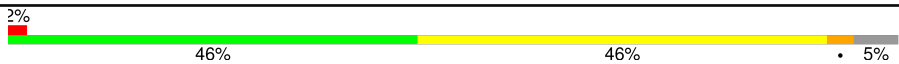

The reported resolution of this entry is 2.94 Å.

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	2057 (2.98-2.90)
Clashscore	102246	2308 (2.98-2.90)
Ramachandran outliers	100387	2245 (2.98-2.90)
Sidechain outliers	100360	2247 (2.98-2.90)
RSRZ outliers	91569	2065 (2.98-2.90)

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	127	
2	B	660	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5059 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 SPCU.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	121	Total	C	N	O	Se	0	1	0
			943	602	165	173	3			

- Molecule 2 is a protein called EXOU.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	534	Total	C	N	O	Se	0	1	0
			4078	2531	738	796	13			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	28	GLY	-	EXPRESSION TAG	UNP Q02IF1
B	29	SER	-	EXPRESSION TAG	UNP Q02IF1

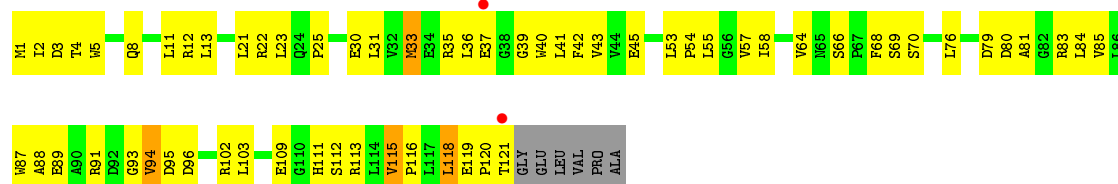
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	9	Total	O	0	0
			9	9		
3	B	29	Total	O	0	0
			29	29		

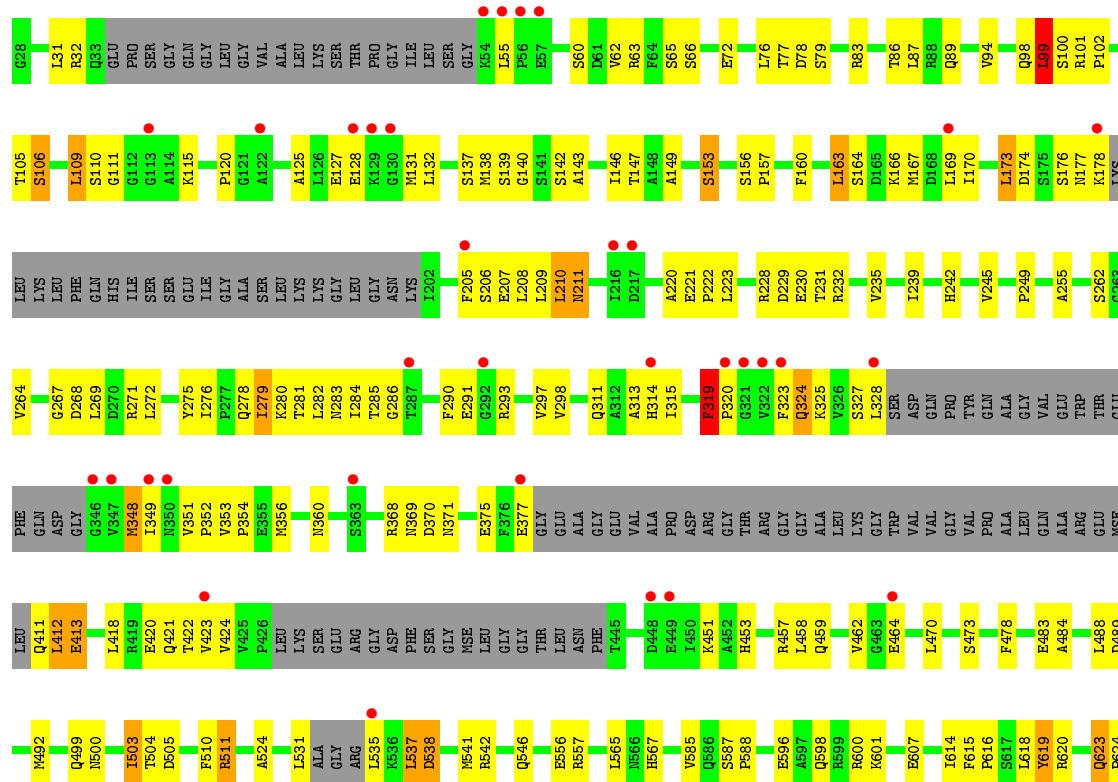
3 Residue-property plots

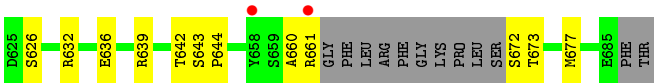
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according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



• Molecule 2: EXOU





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	154.89 Å 52.80 Å 121.81 Å 90.00° 127.18° 90.00°	Depositor
Resolution (Å)	97.05 – 2.94 48.54 – 2.94	Depositor EDS
% Data completeness (in resolution range)	98.7 (97.05-2.94) 98.8 (48.54-2.94)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.00 (at 2.96 Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.208 , 0.264 0.207 , 0.266	Depositor DCC
R_{free} test set	840 reflections (5.27%)	DCC
Wilson B-factor (Å ²)	81.0	Xtriage
Anisotropy	0.256	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 53.6	EDS
Estimated twinning fraction	0.018 for h,-k,-h-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 16783 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5059	wwPDB-VP
Average B, all atoms (Å ²)	98.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.39% 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

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	0.74	1/962 (0.1%)	0.71	0/1312
2	B	0.66	2/4119 (0.0%)	0.64	2/5545 (0.0%)
All	All	0.68	3/5081 (0.1%)	0.66	2/6857 (0.0%)

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
2	B	0	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	327	SER	CB-OG	14.86	1.61	1.42
2	B	377	GLU	C-O	7.65	1.37	1.23
1	A	87	TRP	CD2-CE2	5.94	1.48	1.41

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	565	LEU	CB-CG-CD2	-5.88	101.00	111.00
2	B	99	LEU	CA-CB-CG	5.09	127.01	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	319	PHE	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	943	0	961	58	0
2	B	4078	0	4115	178	0
3	A	9	0	0	2	0
3	B	29	0	0	3	0
All	All	5059	0	5076	230	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 23.

All (230) 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:1[A]:MSE:SE	1:A:1[A]:MSE:CE	2.14	1.45
1:A:21:LEU:HB2	1:A:33:MSE:HE3	1.21	1.20
1:A:119:GLU:HB3	1:A:120:PRO:HD3	1.35	1.08
2:B:221:GLU:HB3	2:B:222:PRO:HD2	1.35	1.06
2:B:453:HIS:CE1	2:B:457:ARG:HH11	1.73	1.05
2:B:489:ASP:H	2:B:492:MSE:HE3	1.18	1.04
2:B:489:ASP:H	2:B:492:MSE:CE	1.70	1.04
1:A:76:LEU:HD21	1:A:84:LEU:HD21	1.41	1.00
2:B:276:ILE:HB	2:B:279:ILE:HD11	1.46	0.97
2:B:206:SER:O	2:B:210:LEU:HD13	1.66	0.94
2:B:453:HIS:HE1	2:B:457:ARG:HH11	1.08	0.93
2:B:319:PHE:CE1	2:B:349:ILE:HD12	2.03	0.93
2:B:489:ASP:N	2:B:492:MSE:HE3	1.86	0.90
2:B:319:PHE:HE1	2:B:349:ILE:HD12	1.36	0.90
2:B:207:GLU:HA	2:B:210:LEU:HB2	1.54	0.88
2:B:164:SER:HA	2:B:167:MSE:HB3	1.55	0.88
2:B:488:LEU:O	2:B:511:ARG:NH2	2.07	0.87
2:B:276:ILE:HB	2:B:279:ILE:CD1	2.05	0.87
2:B:511:ARG:CG	2:B:511:ARG:HH11	1.86	0.86
2:B:76:LEU:HD11	2:B:89:GLN:HE21	1.40	0.86
2:B:499:GLN:HG2	2:B:500:ASN:HD22	1.39	0.86
2:B:76:LEU:HD11	2:B:89:GLN:NE2	1.93	0.84
2:B:146:ILE:HG13	2:B:147:THR:N	1.92	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:483:GLU:CG	2:B:556:GLU:HG2	2.10	0.82
2:B:453:HIS:HE1	2:B:457:ARG:NH1	1.79	0.81
1:A:119:GLU:HB3	1:A:120:PRO:CD	2.10	0.80
1:A:116:PRO:O	1:A:120:PRO:HD2	1.82	0.80
2:B:511:ARG:HH11	2:B:511:ARG:HG2	1.44	0.79
2:B:538:ASP:OD1	2:B:541:MSE:HG3	1.81	0.79
1:A:21:LEU:CB	1:A:33:MSE:HE3	2.10	0.79
2:B:351:VAL:HB	2:B:356:MSE:HE2	1.65	0.78
2:B:636:GLU:HA	2:B:636:GLU:OE1	1.85	0.77
1:A:80:ASP:OD2	2:B:32:ARG:HA	1.84	0.76
2:B:623:GLN:HA	2:B:623:GLN:HE21	1.49	0.76
2:B:360:ASN:HB2	2:B:368:ARG:HH21	1.51	0.76
2:B:285:THR:HG22	2:B:298:VAL:HG22	1.69	0.74
2:B:483:GLU:HG2	2:B:556:GLU:HG2	1.70	0.73
2:B:537:LEU:HA	2:B:541:MSE:HE3	1.71	0.73
2:B:607:GLU:OE1	2:B:607:GLU:HA	1.87	0.73
1:A:31:LEU:HD11	1:A:42:PHE:HB3	1.70	0.73
1:A:21:LEU:HB2	1:A:33:MSE:CE	2.12	0.72
2:B:293:ARG:HH11	2:B:673:THR:HG23	1.52	0.72
2:B:413:GLU:O	2:B:413:GLU:HG2	1.91	0.71
2:B:220:ALA:HB1	2:B:228:ARG:HH12	1.56	0.71
2:B:221:GLU:HB3	2:B:222:PRO:CD	2.19	0.69
1:A:76:LEU:CD2	1:A:84:LEU:HD21	2.19	0.69
1:A:115:VAL:HG12	1:A:116:PRO:HD3	1.73	0.69
2:B:453:HIS:CE1	2:B:457:ARG:HD2	2.29	0.68
2:B:319:PHE:HB3	2:B:320:PRO:HD2	1.74	0.68
2:B:323:PHE:C	2:B:325:LYS:H	1.98	0.67
1:A:79:ASP:HB2	2:B:31:LEU:O	1.94	0.67
2:B:99:LEU:HD23	2:B:100:SER:O	1.94	0.66
1:A:25:PRO:HA	2:B:62:VAL:HG12	1.76	0.66
2:B:499:GLN:HG2	2:B:500:ASN:ND2	2.09	0.66
2:B:83:ARG:HH21	2:B:99:LEU:HD21	1.61	0.66
2:B:77:THR:HG22	2:B:87:LEU:HB2	1.77	0.66
1:A:43:VAL:HG13	1:A:85:VAL:HG13	1.78	0.65
2:B:323:PHE:O	2:B:325:LYS:N	2.30	0.65
2:B:453:HIS:CE1	2:B:457:ARG:NH1	2.56	0.65
2:B:511:ARG:CG	2:B:511:ARG:NH1	2.51	0.65
2:B:101:ARG:HH12	2:B:420:GLU:HG2	1.63	0.64
2:B:128:GLU:OE1	2:B:459:GLN:NE2	2.32	0.63
1:A:23:LEU:HD13	1:A:33:MSE:HE2	1.81	0.63
2:B:94:VAL:HG13	2:B:478:PHE:O	1.99	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:83:ARG:NH2	2:B:99:LEU:HD21	2.14	0.62
2:B:411:GLN:O	2:B:411:GLN:HG2	1.99	0.62
2:B:221:GLU:CB	2:B:222:PRO:HD2	2.21	0.62
2:B:153:SER:HB3	2:B:231:THR:HG22	1.82	0.62
2:B:488:LEU:HA	2:B:492:MSE:HE3	1.80	0.62
1:A:45:GLU:HG3	1:A:83:ARG:HD3	1.81	0.61
2:B:102:PRO:HD2	2:B:370:ASP:OD1	2.01	0.61
2:B:110:SER:OG	2:B:375:GLU:HA	2.01	0.60
1:A:54:PRO:HG2	1:A:57:VAL:HG23	1.84	0.60
1:A:22:ARG:HD3	1:A:30:GLU:OE2	2.01	0.60
2:B:163:LEU:HG	2:B:167:MSE:HE3	1.83	0.60
2:B:231:THR:O	2:B:235:VAL:HG22	2.01	0.60
2:B:120:PRO:HG3	2:B:160:PHE:CE2	2.36	0.60
2:B:146:ILE:HG13	2:B:147:THR:H	1.67	0.60
2:B:286:GLY:O	2:B:297:VAL:HG22	2.02	0.59
2:B:83:ARG:NH2	3:B:2002:HOH:O	2.32	0.58
2:B:293:ARG:NH1	2:B:673:THR:HG23	2.19	0.58
2:B:146:ILE:O	2:B:149:ALA:HB3	2.04	0.58
2:B:511:ARG:NH1	2:B:511:ARG:HG3	2.19	0.57
2:B:208:LEU:HA	2:B:211:ASN:ND2	2.20	0.57
2:B:311:GLN:O	2:B:315:ILE:HG12	2.05	0.57
2:B:293:ARG:HH21	2:B:616:PRO:HB2	1.69	0.57
2:B:223:LEU:HD23	2:B:314:HIS:HD2	1.69	0.57
2:B:489:ASP:CG	2:B:492:MSE:HE2	2.26	0.56
2:B:242:HIS:O	2:B:245:VAL:HG22	2.06	0.56
1:A:12:ARG:HG2	1:A:12:ARG:HH11	1.70	0.55
2:B:166:LYS:O	2:B:170:ILE:HD12	2.07	0.55
1:A:70:SER:HB3	3:A:2007:HOH:O	2.07	0.55
2:B:348:MSE:HG3	2:B:349:ILE:N	2.22	0.55
2:B:77:THR:HG23	2:B:77:THR:O	2.06	0.55
2:B:78:ASP:O	2:B:86:THR:OG1	2.18	0.54
2:B:623:GLN:NE2	2:B:623:GLN:HA	2.22	0.54
2:B:106:SER:OG	2:B:137:SER:HB3	2.08	0.54
2:B:276:ILE:HG22	2:B:278:GLN:H	1.72	0.54
2:B:173:LEU:HA	2:B:176:SER:HB3	1.88	0.54
2:B:78:ASP:OD1	2:B:78:ASP:C	2.46	0.54
2:B:424:VAL:HG13	2:B:424:VAL:O	2.08	0.53
2:B:462:VAL:O	2:B:462:VAL:HG12	2.09	0.53
2:B:272:LEU:HB3	2:B:279:ILE:HD12	1.91	0.53
2:B:109:LEU:O	2:B:140:GLY:HA2	2.08	0.53
2:B:323:PHE:C	2:B:325:LYS:N	2.62	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:119:GLU:CB	1:A:120:PRO:HD3	2.22	0.53
2:B:489:ASP:H	2:B:492:MSE:HE2	1.66	0.53
2:B:72:GLU:HA	3:B:2001:HOH:O	2.07	0.53
2:B:619:TYR:N	2:B:619:TYR:CD1	2.76	0.52
2:B:488:LEU:CA	2:B:492:MSE:HE3	2.39	0.52
2:B:163:LEU:CD1	2:B:230:GLU:HG3	2.39	0.52
2:B:623:GLN:HE21	2:B:624:PRO:HD2	1.76	0.51
2:B:111:GLY:HA3	2:B:142:SER:HB3	1.92	0.51
1:A:13:LEU:HD11	1:A:40:TRP:CZ2	2.46	0.51
1:A:54:PRO:HG2	1:A:57:VAL:CG2	2.40	0.51
2:B:208:LEU:C	2:B:208:LEU:HD12	2.31	0.51
2:B:105:THR:OG1	2:B:369:ASN:HB2	2.10	0.51
2:B:146:ILE:CG1	2:B:147:THR:N	2.69	0.51
1:A:64:VAL:HG12	1:A:64:VAL:O	2.11	0.51
2:B:413:GLU:O	2:B:413:GLU:CG	2.60	0.50
1:A:111:HIS:O	1:A:115:VAL:HB	2.11	0.50
2:B:371:ASN:O	2:B:421:GLN:HB3	2.12	0.50
2:B:464:GLU:HB3	3:B:2015:HOH:O	2.11	0.50
1:A:1[B]:MSE:SE	1:A:4:THR:HG22	2.61	0.50
2:B:319:PHE:HB3	2:B:320:PRO:CD	2.38	0.50
2:B:169:LEU:HD23	2:B:173:LEU:HD23	1.93	0.50
1:A:1[B]:MSE:SE	1:A:3:ASP:HB3	2.62	0.50
2:B:143:ALA:O	2:B:146:ILE:HG12	2.12	0.49
2:B:228:ARG:O	2:B:232:ARG:HB2	2.11	0.49
2:B:360:ASN:ND2	2:B:368:ARG:HE	2.11	0.49
2:B:615:PHE:C	2:B:615:PHE:CD1	2.85	0.49
2:B:600:ARG:O	2:B:601:LYS:C	2.51	0.49
2:B:156:SER:HB2	2:B:157:PRO:HD2	1.93	0.49
1:A:5:TRP:O	1:A:8:GLN:HG2	2.12	0.49
2:B:488:LEU:HB3	2:B:492:MSE:HB2	1.94	0.49
2:B:615:PHE:N	2:B:616:PRO:HD2	2.28	0.49
1:A:84:LEU:HD23	1:A:85:VAL:N	2.28	0.48
2:B:77:THR:HG21	2:B:87:LEU:HD12	1.95	0.48
2:B:139:SER:HA	2:B:283:ASN:O	2.14	0.48
2:B:489:ASP:OD2	2:B:492:MSE:HE2	2.14	0.48
2:B:276:ILE:O	2:B:279:ILE:HG12	2.13	0.48
2:B:524:ALA:HB1	2:B:541:MSE:HG2	1.95	0.48
2:B:245:VAL:HG11	2:B:275:TYR:HB3	1.96	0.47
2:B:239:ILE:HD13	2:B:255:ALA:HB2	1.96	0.47
1:A:43:VAL:HG13	1:A:85:VAL:CG1	2.42	0.47
2:B:348:MSE:O	2:B:348:MSE:HE2	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:503:ILE:HG22	2:B:504:THR:HG23	1.96	0.47
1:A:81:ALA:HB1	2:B:636:GLU:HG2	1.95	0.47
1:A:93:GLY:O	1:A:94:VAL:C	2.54	0.47
2:B:324:GLN:HG3	2:B:324:GLN:O	2.14	0.47
2:B:127:GLU:HG2	2:B:132:LEU:HD22	1.97	0.46
1:A:35:ARG:HD2	1:A:36:LEU:H	1.80	0.46
2:B:285:THR:OG1	2:B:356:MSE:HG2	2.16	0.46
2:B:458:LEU:O	2:B:459:GLN:C	2.54	0.46
2:B:353:VAL:N	2:B:354:PRO:HD2	2.30	0.46
2:B:205:PHE:CZ	2:B:290:PHE:CD1	3.03	0.46
2:B:585:VAL:HG23	2:B:585:VAL:O	2.14	0.46
2:B:220:ALA:HB1	2:B:228:ARG:NH1	2.27	0.46
2:B:510:PHE:CD1	2:B:510:PHE:C	2.89	0.46
2:B:173:LEU:C	2:B:173:LEU:HD12	2.37	0.46
2:B:138:MSE:CE	2:B:280:LYS:HG3	2.46	0.46
1:A:109:GLU:OE2	1:A:113:ARG:NH2	2.49	0.46
2:B:268:ASP:HA	2:B:271:ARG:HB2	1.98	0.45
2:B:178:LYS:HZ3	2:B:319:PHE:HD2	1.64	0.45
1:A:1[B]:MSE:SE	1:A:4:THR:H	2.49	0.45
1:A:91:ARG:N	1:A:95:ASP:OD2	2.49	0.45
2:B:267:GLY:O	2:B:271:ARG:HG3	2.16	0.45
2:B:167:MSE:HE3	2:B:167:MSE:HB2	1.86	0.45
2:B:98:GLN:HG2	2:B:473:SER:OG	2.16	0.45
2:B:614:ILE:HD13	2:B:614:ILE:N	2.30	0.45
2:B:153:SER:HB2	2:B:231:THR:HA	1.99	0.45
2:B:174:ASP:HA	2:B:177:ASN:HB2	1.99	0.45
2:B:125:ALA:HA	2:B:459:GLN:HE22	1.81	0.44
1:A:53:LEU:HD23	1:A:121:THR:HG21	2.00	0.44
2:B:542:ARG:O	2:B:546:GLN:HG3	2.18	0.44
2:B:205:PHE:HA	2:B:208:LEU:HG	1.99	0.44
1:A:88:ALA:HB1	1:A:103:LEU:HD22	2.00	0.44
2:B:293:ARG:HE	2:B:616:PRO:HB2	1.82	0.44
2:B:478:PHE:CD1	2:B:484:ALA:HB2	2.52	0.44
2:B:127:GLU:HG2	2:B:132:LEU:CD2	2.47	0.44
2:B:412:LEU:HD12	2:B:412:LEU:O	2.18	0.43
2:B:615:PHE:HB3	2:B:616:PRO:CD	2.48	0.43
2:B:125:ALA:HA	2:B:459:GLN:NE2	2.34	0.43
1:A:116:PRO:O	1:A:120:PRO:CD	2.61	0.43
2:B:596:GLU:OE2	2:B:600:ARG:NE	2.49	0.43
1:A:36:LEU:HB2	1:A:39:GLY:O	2.19	0.43
2:B:282:LEU:HD21	2:B:284:ILE:HD11	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:418:LEU:HA	2:B:418:LEU:HD23	1.81	0.43
1:A:119:GLU:CB	1:A:120:PRO:CD	2.84	0.43
2:B:620:ARG:O	2:B:623:GLN:HB2	2.19	0.43
2:B:505:ASP:N	2:B:505:ASP:OD1	2.51	0.43
1:A:118:LEU:N	1:A:118:LEU:HD12	2.35	0.42
1:A:66:SER:C	1:A:68:PHE:H	2.23	0.42
2:B:229:ASP:OD1	2:B:232:ARG:NH2	2.53	0.42
1:A:94:VAL:HG12	1:A:95:ASP:OD1	2.20	0.42
2:B:163:LEU:O	2:B:167:MSE:HB2	2.20	0.42
2:B:503:ILE:HA	2:B:503:ILE:HD13	1.77	0.42
2:B:504:THR:O	2:B:505:ASP:C	2.57	0.42
1:A:88:ALA:CB	1:A:103:LEU:HD22	2.50	0.42
2:B:272:LEU:C	2:B:279:ILE:HG13	2.40	0.42
1:A:102:ARG:O	1:A:103:LEU:C	2.57	0.42
2:B:510:PHE:CE2	2:B:557:ARG:HD3	2.55	0.42
2:B:587:SER:HA	2:B:588:PRO:HD2	1.83	0.42
1:A:89:GLU:OE1	1:A:89:GLU:HA	2.20	0.42
2:B:101:ARG:HA	2:B:102:PRO:HD3	1.92	0.42
2:B:291:GLU:HG2	2:B:291:GLU:O	2.17	0.42
1:A:66:SER:C	1:A:68:PHE:N	2.74	0.41
1:A:45:GLU:CG	1:A:83:ARG:HD3	2.46	0.41
1:A:55:LEU:O	1:A:58:ILE:HB	2.20	0.41
3:A:2002:HOH:O	2:B:66:SER:HB2	2.19	0.41
1:A:1[B]:MSE:HE2	1:A:1[B]:MSE:HB2	1.61	0.41
1:A:25:PRO:CA	2:B:62:VAL:HG12	2.46	0.41
2:B:499:GLN:CG	2:B:500:ASN:HD22	2.22	0.41
1:A:11:LEU:HD11	2:B:62:VAL:O	2.21	0.41
2:B:618:LEU:HB3	2:B:619:TYR:CD1	2.56	0.41
2:B:352:PRO:C	2:B:354:PRO:HD2	2.41	0.41
1:A:84:LEU:HD23	1:A:85:VAL:H	1.85	0.41
2:B:412:LEU:HD22	2:B:639:ARG:HH22	1.85	0.41
2:B:643:SER:HB2	2:B:644:PRO:HD2	2.02	0.41
2:B:319:PHE:CB	2:B:320:PRO:CD	2.99	0.41
2:B:137:SER:HA	2:B:281:THR:OG1	2.21	0.41
2:B:483:GLU:HG3	2:B:556:GLU:HG2	1.96	0.40
2:B:221:GLU:CB	2:B:222:PRO:CD	2.91	0.40
2:B:78:ASP:H	2:B:499:GLN:NE2	2.19	0.40
2:B:146:ILE:HA	2:B:313:ALA:HB1	2.03	0.40
1:A:36:LEU:HB3	1:A:37:GLU:H	1.68	0.40
2:B:269:LEU:HD23	2:B:269:LEU:HA	1.64	0.40
1:A:76:LEU:HD23	1:A:76:LEU:C	2.41	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:101:ARG:NH1	2:B:420:GLU:HG2	2.34	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	119/127 (94%)	108 (91%)	9 (8%)	2 (2%)	11	37
2	B	519/660 (79%)	468 (90%)	39 (8%)	12 (2%)	8	29
All	All	638/787 (81%)	576 (90%)	48 (8%)	14 (2%)	8	30

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	94	VAL
2	B	319	PHE
2	B	153	SER
2	B	324	GLN
2	B	412	LEU
2	B	660	ALA
2	B	210	LEU
2	B	60	SER
2	B	115	LYS
2	B	537	LEU
1	A	2	ILE
2	B	249	PRO
2	B	538	ASP
2	B	567	HIS

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	102/103 (99%)	95 (93%)	7 (7%)	19	48
2	B	440/517 (85%)	404 (92%)	36 (8%)	14	38
All	All	542/620 (87%)	499 (92%)	43 (8%)	15	40

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	33	MSE
1	A	41	LEU
1	A	69	SER
1	A	96	ASP
1	A	112	SER
1	A	115	VAL
1	A	118	LEU
2	B	55	LEU
2	B	63	ARG
2	B	65	SER
2	B	79	SER
2	B	99	LEU
2	B	106	SER
2	B	109	LEU
2	B	131	MSE
2	B	163	LEU
2	B	173	LEU
2	B	209	LEU
2	B	211	ASN
2	B	262	SER
2	B	264	VAL
2	B	279	ILE
2	B	319	PHE
2	B	328	LEU
2	B	348	MSE
2	B	413	GLU
2	B	422	THR

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Mol	Chain	Res	Type
2	B	423	VAL
2	B	451	LYS
2	B	470	LEU
2	B	503	ILE
2	B	511	ARG
2	B	531	LEU
2	B	535	LEU
2	B	598	GLN
2	B	619	TYR
2	B	623	GLN
2	B	626	SER
2	B	632	ARG
2	B	642	THR
2	B	661	ARG
2	B	672	SER
2	B	677	MSE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	24	GLN
1	A	111	HIS
2	B	89	GLN
2	B	314	HIS
2	B	360	ASN
2	B	411	GLN
2	B	453	HIS
2	B	471	GLN
2	B	499	GLN
2	B	500	ASN
2	B	567	HIS
2	B	575	GLN
2	B	623	GLN
2	B	648	ASN

5.3.3 RNA ⓘ

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	119/127 (93%)	-0.25	2 (1%) 73 72	50, 76, 130, 190	0
2	B	521/660 (78%)	0.08	35 (6%) 21 16	41, 98, 175, 206	0
All	All	640/787 (81%)	0.02	37 (5%) 26 22	41, 94, 174, 206	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	347	VAL	15.8
2	B	346	GLY	8.0
1	A	121	THR	7.1
2	B	54	LYS	7.0
2	B	217	ASP	6.2
2	B	349	ILE	5.9
2	B	321	GLY	5.8
2	B	322	VAL	5.4
2	B	320	PRO	4.3
2	B	328	LEU	3.9
2	B	292	GLY	3.8
2	B	216	ILE	3.5
2	B	323	PHE	3.4
2	B	287	THR	3.3
2	B	658	TYR	3.2
2	B	55	LEU	3.1
2	B	377	GLU	2.9
2	B	122	ALA	2.8
2	B	178	LYS	2.8
2	B	205	PHE	2.8
2	B	314	HIS	2.7
2	B	350	ASN	2.7
2	B	423	VAL	2.6
2	B	56	PRO	2.5

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Mol	Chain	Res	Type	RSRZ
2	B	169	LEU	2.5
2	B	130	GLY	2.5
2	B	129	LYS	2.4
2	B	449	GLU	2.3
2	B	464	GLU	2.3
2	B	535	LEU	2.3
1	A	37	GLU	2.2
2	B	661	ARG	2.2
2	B	113	GLY	2.1
2	B	128	GLU	2.1
2	B	363	SER	2.1
2	B	57	GLU	2.0
2	B	448	ASP	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.