



Full wwPDB X-ray Structure Validation Report ⓘ

May 26, 2020 – 05:50 pm BST

PDB ID : 1L2W
Title : Crystal Structure of the Yersinia Virulence Effector YopE Chaperone-binding Domain in Complex with its Secretion Chaperone, SycE
Authors : Birtalan, S.C.; Phillips, R.M.; Ghosh, P.
Deposited on : 2002-02-25
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

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:

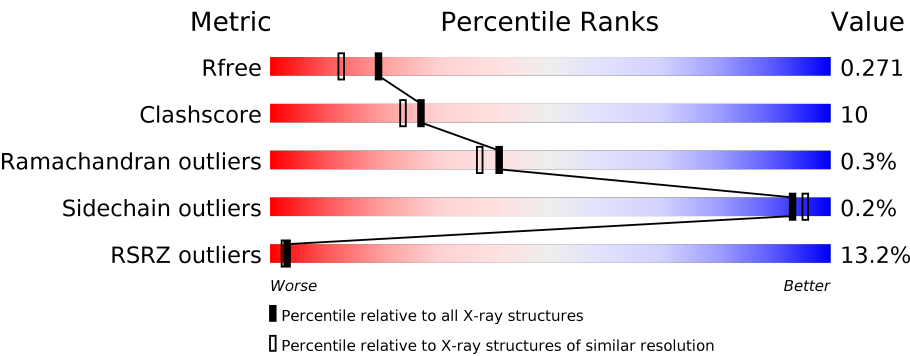
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

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 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}	130704	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (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 ≥ 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	123	<div><div>7%</div><div>80%19%.</div></div>
1	B	123	<div><div>5%</div><div>76%20%5%</div></div>
1	C	123	<div><div>8%</div><div>83%12%5%</div></div>
1	D	123	<div><div>11%</div><div>76%22%. .</div></div>
1	E	123	<div><div>11%</div><div>76%19%. 5%</div></div>
1	F	123	<div><div>15%</div><div>73%24%. .</div></div>

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Mol	Chain	Length	Quality of chain
1	G	123	
1	H	123	
2	I	69	
2	J	69	
2	K	69	
2	L	69	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9594 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 YopE regulator.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	122	Total	C	N	O	S	0	0	0
			965	614	158	189	4			
1	B	117	Total	C	N	O	S	0	0	0
			934	597	153	181	3			
1	C	117	Total	C	N	O	S	0	0	0
			934	597	153	181	3			
1	D	121	Total	C	N	O	S	0	0	0
			959	611	157	187	4			
1	E	117	Total	C	N	O	S	0	0	0
			934	597	153	181	3			
1	F	119	Total	C	N	O	S	0	0	0
			946	604	155	183	4			
1	G	119	Total	C	N	O	S	0	0	0
			946	604	155	183	4			
1	H	117	Total	C	N	O	S	0	0	0
			934	597	153	181	3			

- Molecule 2 is a protein called Outer membrane virulence protein yopE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	I	57	Total	C	N	O	S	0	0	0
			428	258	81	87	2			
2	J	56	Total	C	N	O	S	0	0	0
			422	255	80	85	2			
2	K	55	Total	C	N	O	S	0	0	0
			415	250	79	84	2			
2	L	55	Total	C	N	O	S	0	0	0
			415	250	79	84	2			

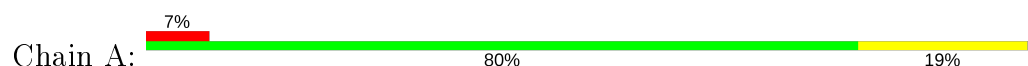
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	59	Total 59	O 59	0	0
3	B	42	Total 42	O 42	0	0
3	C	38	Total 38	O 38	0	0
3	D	48	Total 48	O 48	0	0
3	E	40	Total 40	O 40	0	0
3	F	36	Total 36	O 36	0	0
3	G	38	Total 38	O 38	0	0
3	H	25	Total 25	O 25	0	0
3	I	13	Total 13	O 13	0	0
3	J	6	Total 6	O 6	0	0
3	K	10	Total 10	O 10	0	0
3	L	7	Total 7	O 7	0	0

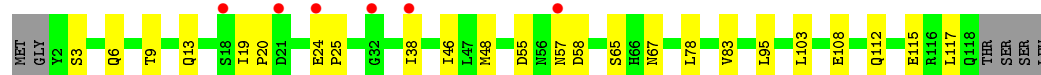
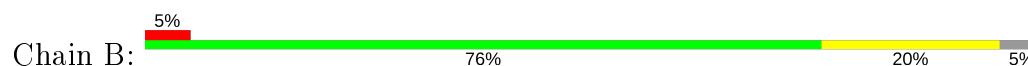
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

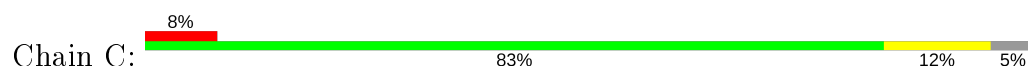
- Molecule 1: YopE regulator



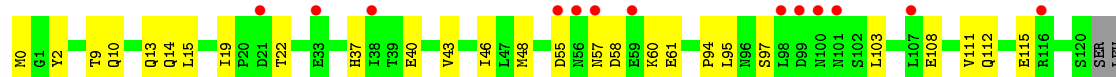
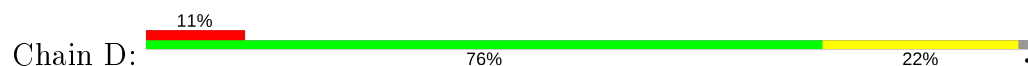
- Molecule 1: YopE regulator



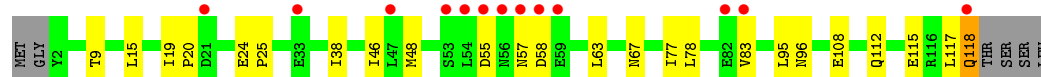
- Molecule 1: YopE regulator



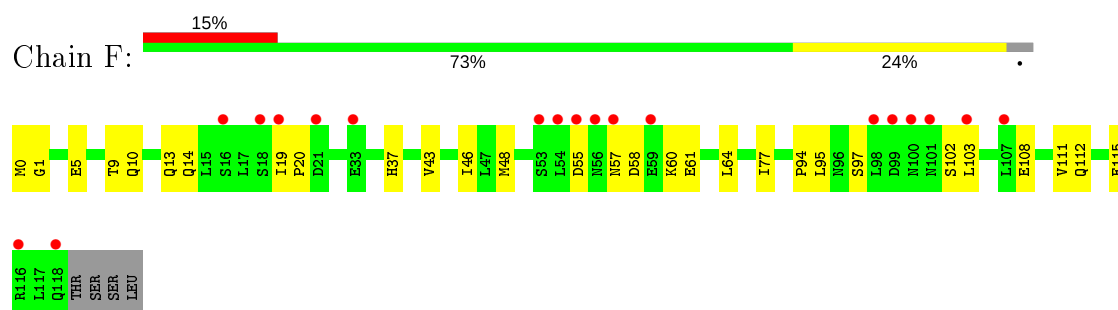
- Molecule 1: YopE regulator



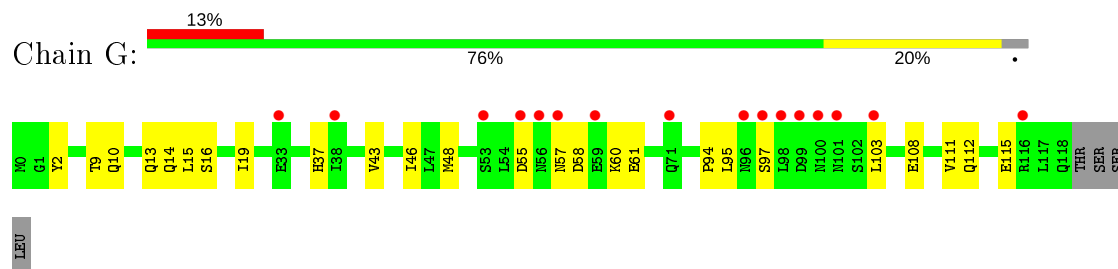
- Molecule 1: YopE regulator



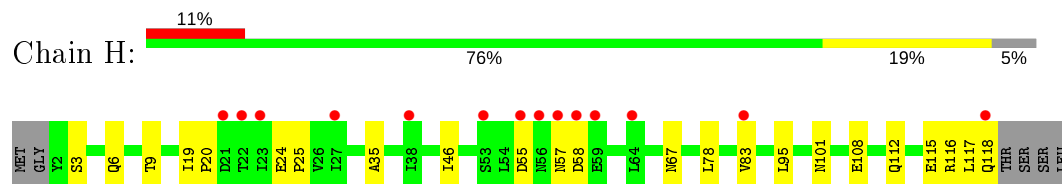
- Molecule 1: YopE regulator



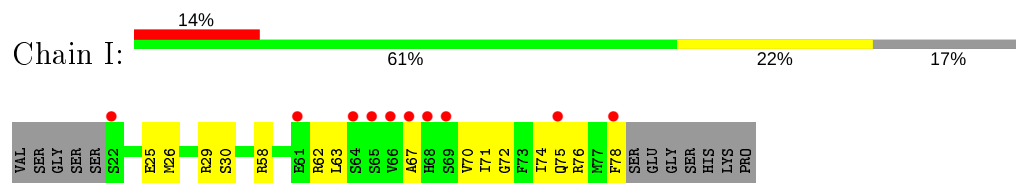
- Molecule 1: YopE regulator



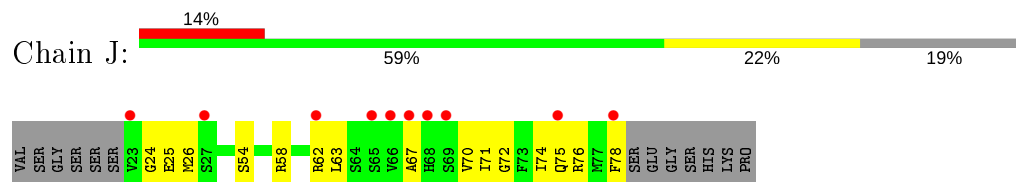
- Molecule 1: YopE regulator



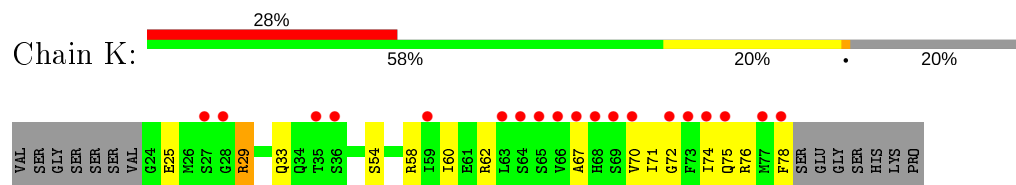
- Molecule 2: Outer membrane virulence protein yopE



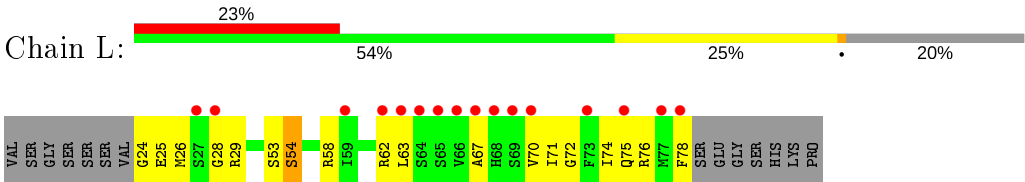
- Molecule 2: Outer membrane virulence protein yopE



- Molecule 2: Outer membrane virulence protein yopE



- Molecule 2: Outer membrane virulence protein yopE



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	72.84Å 73.35Å 74.26Å 103.37° 109.18° 107.36°	Depositor
Resolution (Å)	20.00 – 2.00 19.98 – 1.96	Depositor EDS
% Data completeness (in resolution range)	97.3 (20.00-2.00) 96.3 (19.98-1.96)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.59 (at 1.96Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.251 , 0.275 0.248 , 0.271	Depositor DCC
R_{free} test set	4252 reflections (4.79%)	wwPDB-VP
Wilson B-factor (Å ²)	36.3	Xtriage
Anisotropy	0.188	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 46.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.010 for -h,-l,-k	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9594	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.47% 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 [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	0.38	0/986	0.57	0/1343
1	B	0.37	0/955	0.59	0/1302
1	C	0.36	0/955	0.59	0/1302
1	D	0.35	0/980	0.57	0/1335
1	E	0.36	0/955	0.59	0/1302
1	F	0.33	0/967	0.55	0/1317
1	G	0.34	0/967	0.55	0/1317
1	H	0.34	0/955	0.59	0/1302
2	I	0.35	0/432	0.55	0/579
2	J	0.29	0/426	0.50	0/571
2	K	0.32	0/419	0.51	0/561
2	L	0.32	0/419	0.50	0/561
All	All	0.35	0/9416	0.57	0/12792

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	965	0	948	21	0
1	B	934	0	916	18	0
1	C	934	0	916	16	0
1	D	959	0	943	21	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	934	0	916	20	0
1	F	946	0	931	22	0
1	G	946	0	931	20	0
1	H	934	0	916	20	0
2	I	428	0	419	16	0
2	J	422	0	414	12	0
2	K	415	0	405	15	0
2	L	415	0	405	20	0
3	A	59	0	0	1	0
3	B	42	0	0	0	0
3	C	38	0	0	0	0
3	D	48	0	0	1	0
3	E	40	0	0	1	0
3	F	36	0	0	2	0
3	G	38	0	0	1	0
3	H	25	0	0	0	0
3	I	13	0	0	0	0
3	J	6	0	0	0	0
3	K	10	0	0	0	0
3	L	7	0	0	0	0
All	All	9594	0	9060	186	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 10.

All (186) 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:67:ASN:HD21	1:B:78:LEU:H	1.07	0.98
1:C:67:ASN:HD21	1:C:78:LEU:H	1.08	0.98
1:H:67:ASN:HD21	1:H:78:LEU:H	1.14	0.92
1:E:67:ASN:HD21	1:E:78:LEU:H	1.21	0.83
1:B:67:ASN:ND2	1:B:78:LEU:H	1.75	0.83
1:C:67:ASN:ND2	1:C:78:LEU:H	1.80	0.80
1:A:108:GLU:HA	2:I:26:MET:CE	2.13	0.79
2:L:25:GLU:HG2	2:L:26:MET:H	1.45	0.79
1:H:67:ASN:ND2	1:H:78:LEU:H	1.80	0.79
1:A:0:MET:HA	1:A:40:GLU:OE1	1.85	0.76
1:A:108:GLU:HA	2:I:26:MET:HE1	1.68	0.75
1:E:67:ASN:ND2	1:E:78:LEU:H	1.85	0.73
1:F:43:VAL:HG23	3:F:126:HOH:O	1.88	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:9:THR:O	1:A:13:GLN:HG3	1.92	0.69
2:L:25:GLU:HG2	2:L:26:MET:N	2.07	0.69
2:I:25:GLU:HG2	2:I:30:SER:HA	1.75	0.69
1:F:9:THR:O	1:F:13:GLN:HG3	1.93	0.68
1:C:101:ASN:HB3	1:G:2:TYR:HB3	1.75	0.68
1:G:9:THR:O	1:G:13:GLN:HG3	1.94	0.67
1:C:57:ASN:HD22	1:C:57:ASN:N	1.93	0.67
1:G:43:VAL:HG23	3:G:146:HOH:O	1.93	0.67
1:D:9:THR:O	1:D:13:GLN:HG3	1.95	0.66
1:C:20:PRO:HG3	2:J:62:ARG:CZ	2.26	0.65
1:D:22:THR:HG23	2:L:53:SER:O	1.97	0.65
1:B:20:PRO:HG3	2:I:62:ARG:CZ	2.27	0.65
1:D:0:MET:HA	1:D:40:GLU:OE1	1.96	0.65
1:B:67:ASN:HD21	1:B:78:LEU:N	1.89	0.64
2:L:25:GLU:CD	2:L:28:GLY:H	2.02	0.64
1:G:108:GLU:O	1:G:112:GLN:HG3	1.99	0.63
1:H:83:VAL:HB	2:L:76:ARG:NH1	2.13	0.63
1:E:20:PRO:HG3	2:K:62:ARG:CZ	2.29	0.62
1:D:2:TYR:HB3	1:H:101:ASN:HB3	1.82	0.61
1:C:83:VAL:HB	2:J:76:ARG:NH1	2.17	0.60
1:A:32:GLY:HA3	2:I:29:ARG:HD3	1.83	0.59
2:L:26:MET:O	2:L:29:ARG:HD2	2.01	0.59
1:F:108:GLU:O	1:F:112:GLN:HG3	2.02	0.59
1:D:108:GLU:O	1:D:112:GLN:HG3	2.03	0.58
1:E:83:VAL:HB	2:K:76:ARG:NH1	2.19	0.58
1:G:111:VAL:O	1:G:115:GLU:HG3	2.04	0.57
1:H:108:GLU:O	1:H:112:GLN:HG3	2.04	0.57
1:A:108:GLU:HA	2:I:26:MET:HE3	1.88	0.56
2:J:54:SER:O	2:J:58:ARG:HG3	2.05	0.56
1:H:57:ASN:HD22	1:H:57:ASN:N	2.03	0.56
1:B:57:ASN:HD22	1:B:57:ASN:N	2.03	0.56
1:F:0:MET:CE	1:F:1:GLY:H	2.18	0.56
1:A:43:VAL:HG23	3:A:134:HOH:O	2.06	0.56
1:D:111:VAL:O	1:D:115:GLU:HG3	2.07	0.55
1:C:108:GLU:O	1:C:112:GLN:HG3	2.08	0.54
1:B:108:GLU:O	1:B:112:GLN:HG3	2.07	0.54
1:D:46:ILE:HD11	1:D:95:LEU:HD13	1.89	0.54
1:E:57:ASN:N	1:E:57:ASN:HD22	2.04	0.54
1:D:55:ASP:HB3	1:D:58:ASP:OD2	2.08	0.54
2:I:67:ALA:O	2:I:71:ILE:HG13	2.08	0.54
1:A:33:GLU:CD	1:A:118:GLN:HE22	2.11	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:9:THR:HG23	1:F:19:ILE:HD13	1.89	0.54
1:H:67:ASN:HD21	1:H:78:LEU:N	1.96	0.53
1:A:108:GLU:O	1:A:112:GLN:HG3	2.09	0.53
2:I:74:ILE:O	2:I:78:PHE:HB2	2.08	0.53
1:F:55:ASP:HB3	1:F:58:ASP:OD2	2.09	0.53
1:B:46:ILE:HD11	1:B:95:LEU:HD13	1.91	0.52
2:J:67:ALA:HB3	2:J:70:VAL:HG23	1.92	0.52
1:H:46:ILE:HD11	1:H:95:LEU:HD13	1.92	0.52
2:J:67:ALA:O	2:J:71:ILE:HG13	2.10	0.52
2:K:67:ALA:HB3	2:K:70:VAL:HG23	1.91	0.52
1:E:46:ILE:HD11	1:E:95:LEU:HD13	1.92	0.52
2:I:67:ALA:HB3	2:I:70:VAL:HG23	1.92	0.52
1:C:57:ASN:N	1:C:57:ASN:ND2	2.56	0.52
2:L:67:ALA:O	2:L:71:ILE:HG13	2.10	0.52
1:A:94:PRO:HB2	1:A:97:SER:HB3	1.92	0.51
1:D:9:THR:HG23	1:D:19:ILE:HD13	1.91	0.51
2:I:72:GLY:O	2:I:76:ARG:HB2	2.11	0.51
2:K:67:ALA:O	2:K:71:ILE:HG13	2.11	0.51
2:L:72:GLY:O	2:L:76:ARG:HB2	2.10	0.51
1:G:55:ASP:HB3	1:G:58:ASP:OD2	2.10	0.51
1:H:35:ALA:HB1	2:L:63:LEU:HD11	1.92	0.51
1:C:67:ASN:HD21	1:C:78:LEU:N	1.92	0.51
1:D:43:VAL:HG23	3:D:164:HOH:O	2.10	0.51
1:C:101:ASN:ND2	1:G:2:TYR:HD2	2.09	0.51
1:G:15:LEU:O	2:L:24:GLY:HA2	2.11	0.50
1:F:111:VAL:O	1:F:115:GLU:HG3	2.11	0.50
1:G:9:THR:HG23	1:G:19:ILE:HD13	1.93	0.50
2:J:74:ILE:O	2:J:78:PHE:HB2	2.11	0.50
2:L:67:ALA:HB3	2:L:70:VAL:HG23	1.92	0.50
1:H:55:ASP:HB3	1:H:58:ASP:OD2	2.12	0.50
2:K:74:ILE:O	2:K:78:PHE:HB2	2.12	0.50
1:E:63:LEU:HD13	1:F:64:LEU:HD11	1.94	0.50
1:G:94:PRO:HB2	1:G:97:SER:HB3	1.94	0.50
1:E:108:GLU:O	1:E:112:GLN:HG3	2.11	0.50
2:K:72:GLY:O	2:K:76:ARG:HB2	2.12	0.50
2:K:25:GLU:HG3	2:K:29:ARG:O	2.12	0.49
1:B:9:THR:HG23	1:B:19:ILE:HD13	1.94	0.49
1:D:15:LEU:O	2:J:24:GLY:HA3	2.12	0.49
1:F:102:SER:HB3	3:F:156:HOH:O	2.11	0.49
1:G:46:ILE:HD11	1:G:95:LEU:HD13	1.95	0.49
2:J:72:GLY:O	2:J:76:ARG:HB2	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:24:GLU:HB3	1:E:25:PRO:HD2	1.94	0.49
2:L:74:ILE:O	2:L:78:PHE:HB2	2.13	0.48
1:E:55:ASP:HB3	1:E:58:ASP:OD2	2.13	0.48
1:E:77:ILE:HG21	1:F:77:ILE:HD12	1.94	0.48
1:H:24:GLU:HB3	1:H:25:PRO:HD2	1.94	0.48
1:E:117:LEU:O	1:E:118:GLN:CB	2.62	0.48
1:G:55:ASP:OD1	1:G:57:ASN:HB2	2.13	0.48
1:E:15:LEU:HD22	2:K:60:ILE:HD13	1.95	0.48
1:A:111:VAL:O	1:A:115:GLU:HG3	2.13	0.48
1:A:9:THR:HG23	1:A:19:ILE:HD13	1.94	0.48
1:B:55:ASP:HB3	1:B:58:ASP:OD2	2.13	0.48
1:C:24:GLU:HB3	1:C:25:PRO:HD2	1.95	0.48
1:C:55:ASP:HB3	1:C:58:ASP:OD2	2.13	0.48
1:H:20:PRO:HG3	2:L:62:ARG:CZ	2.44	0.48
1:F:94:PRO:HB2	1:F:97:SER:HB3	1.96	0.48
1:A:60:LYS:HG3	1:A:61:GLU:CD	2.35	0.47
1:F:46:ILE:HD11	1:F:95:LEU:HD13	1.95	0.47
1:D:55:ASP:OD1	1:D:57:ASN:HB2	2.14	0.47
1:D:94:PRO:HB2	1:D:97:SER:HB3	1.96	0.47
1:G:60:LYS:HG3	1:G:61:GLU:CD	2.35	0.47
1:A:46:ILE:HD11	1:A:95:LEU:HD13	1.96	0.47
1:B:24:GLU:HB3	1:B:25:PRO:HD2	1.97	0.47
1:G:46:ILE:HG21	1:G:103:LEU:HD22	1.97	0.47
1:F:55:ASP:OD1	1:F:57:ASN:HB2	2.15	0.46
1:C:57:ASN:H	1:C:57:ASN:ND2	2.13	0.46
1:E:57:ASN:N	1:E:57:ASN:ND2	2.64	0.46
1:A:60:LYS:HG3	1:A:61:GLU:H	1.81	0.46
1:F:60:LYS:HG3	1:F:61:GLU:CD	2.36	0.46
1:B:57:ASN:ND2	1:B:57:ASN:N	2.64	0.46
2:L:54:SER:HB2	2:L:58:ARG:CZ	2.47	0.45
1:D:46:ILE:HG21	1:D:103:LEU:HD22	1.99	0.45
1:D:60:LYS:HG3	1:D:61:GLU:CD	2.37	0.45
1:H:9:THR:HG23	1:H:19:ILE:HD13	1.98	0.45
1:H:57:ASN:N	1:H:57:ASN:ND2	2.64	0.45
1:A:60:LYS:HG3	1:A:61:GLU:N	2.32	0.45
1:A:32:GLY:CA	2:I:29:ARG:HB3	2.48	0.44
1:H:116:ARG:C	1:H:118:GLN:H	2.19	0.44
1:A:46:ILE:HG21	1:A:103:LEU:HD22	1.99	0.44
1:G:16:SER:HB2	2:L:24:GLY:N	2.33	0.44
1:C:101:ASN:ND2	1:G:2:TYR:CD2	2.85	0.44
1:F:46:ILE:HG21	1:F:103:LEU:HD22	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:111:VAL:HG21	2:I:26:MET:HE2	2.00	0.44
1:D:60:LYS:HG3	1:D:61:GLU:H	1.83	0.44
1:B:115:GLU:OE1	2:I:58:ARG:NH2	2.51	0.43
1:F:20:PRO:HG3	2:K:33:GLN:OE1	2.18	0.43
1:G:60:LYS:HG3	1:G:61:GLU:N	2.33	0.43
2:K:70:VAL:O	2:K:74:ILE:HG13	2.18	0.43
1:E:67:ASN:HD21	1:E:78:LEU:N	2.02	0.43
1:C:35:ALA:HB1	2:J:63:LEU:HD11	2.01	0.43
1:D:60:LYS:HG3	1:D:61:GLU:N	2.34	0.43
1:D:2:TYR:HD2	1:H:101:ASN:ND2	2.17	0.43
2:K:54:SER:HB2	2:K:58:ARG:CZ	2.48	0.43
1:E:115:GLU:OE1	2:K:58:ARG:NH2	2.52	0.42
2:I:71:ILE:CG2	2:I:75:GLN:HE21	2.32	0.42
2:J:71:ILE:CG2	2:J:75:GLN:HE21	2.32	0.42
1:B:38:ILE:HG12	1:B:48:MET:HG2	2.00	0.42
1:F:10:GLN:O	1:F:14:GLN:HG3	2.19	0.42
1:E:38:ILE:HG12	1:E:48:MET:HG2	2.00	0.42
2:L:71:ILE:CG2	2:L:75:GLN:HE21	2.32	0.42
1:A:55:ASP:OD2	1:A:120:SER:HA	2.19	0.42
2:I:63:LEU:HD23	2:I:63:LEU:HA	1.90	0.42
1:F:60:LYS:HG3	1:F:61:GLU:N	2.35	0.42
1:H:83:VAL:HB	2:L:76:ARG:HH11	1.84	0.42
1:D:10:GLN:O	1:D:14:GLN:HG3	2.20	0.42
1:H:19:ILE:HA	1:H:20:PRO:HD3	1.90	0.42
1:H:3:SER:OG	1:H:6:GLN:HG3	2.18	0.42
1:G:60:LYS:HG3	1:G:61:GLU:H	1.85	0.41
1:B:46:ILE:HG21	1:B:103:LEU:CD2	2.50	0.41
1:E:96:ASN:ND2	3:E:156:HOH:O	2.52	0.41
1:F:0:MET:HE3	1:F:1:GLY:H	1.85	0.41
1:F:37:HIS:O	1:F:48:MET:HA	2.20	0.41
2:K:29:ARG:HD2	2:K:29:ARG:H	1.84	0.41
2:J:70:VAL:O	2:J:74:ILE:HG13	2.20	0.41
1:D:46:ILE:CG1	1:D:95:LEU:HD13	2.50	0.41
1:G:37:HIS:O	1:G:48:MET:HA	2.21	0.41
2:J:25:GLU:HG2	2:J:26:MET:N	2.35	0.41
2:L:70:VAL:O	2:L:74:ILE:HG13	2.21	0.41
1:E:9:THR:HG23	1:E:19:ILE:HD13	2.03	0.41
1:B:13:GLN:CG	1:B:19:ILE:HD11	2.51	0.41
2:K:71:ILE:CG2	2:K:75:GLN:HE21	2.33	0.41
1:F:0:MET:HE1	1:F:5:GLU:HG3	2.03	0.41
1:A:82:GLU:HG2	1:B:65:SER:OG	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:3:SER:OG	1:B:6:GLN:HG3	2.21	0.40
1:C:9:THR:HG23	1:C:19:ILE:HD13	2.03	0.40
1:D:37:HIS:O	1:D:48:MET:HA	2.20	0.40
1:E:20:PRO:HG3	2:K:62:ARG:NE	2.36	0.40
1:G:10:GLN:O	1:G:14:GLN:HG3	2.22	0.40
2:L:63:LEU:HD23	2:L:63:LEU:HA	1.90	0.40
1:F:60:LYS:HG3	1:F:61:GLU:H	1.86	0.40
1:H:115:GLU:OE2	2:L:58:ARG:CZ	2.69	0.40
1:B:83:VAL:HB	2:I:76:ARG:NH1	2.37	0.40

There are no symmetry-related clashes.

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	120/123 (98%)	117 (98%)	3 (2%)	0	100	100
1	B	115/123 (94%)	112 (97%)	2 (2%)	1 (1%)	17	11
1	C	115/123 (94%)	113 (98%)	2 (2%)	0	100	100
1	D	119/123 (97%)	117 (98%)	2 (2%)	0	100	100
1	E	115/123 (94%)	114 (99%)	1 (1%)	0	100	100
1	F	117/123 (95%)	114 (97%)	3 (3%)	0	100	100
1	G	117/123 (95%)	115 (98%)	2 (2%)	0	100	100
1	H	115/123 (94%)	112 (97%)	2 (2%)	1 (1%)	17	11
2	I	55/69 (80%)	51 (93%)	4 (7%)	0	100	100
2	J	54/69 (78%)	50 (93%)	4 (7%)	0	100	100
2	K	53/69 (77%)	48 (91%)	5 (9%)	0	100	100
2	L	53/69 (77%)	47 (89%)	5 (9%)	1 (2%)	8	3
All	All	1148/1260 (91%)	1110 (97%)	35 (3%)	3 (0%)	41	37

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	H	117	LEU
1	B	117	LEU
2	L	54	SER

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	112/113 (99%)	112 (100%)	0	100	100
1	B	108/113 (96%)	108 (100%)	0	100	100
1	C	108/113 (96%)	108 (100%)	0	100	100
1	D	111/113 (98%)	111 (100%)	0	100	100
1	E	108/113 (96%)	107 (99%)	1 (1%)	78	83
1	F	109/113 (96%)	109 (100%)	0	100	100
1	G	109/113 (96%)	109 (100%)	0	100	100
1	H	108/113 (96%)	108 (100%)	0	100	100
2	I	48/58 (83%)	48 (100%)	0	100	100
2	J	47/58 (81%)	47 (100%)	0	100	100
2	K	46/58 (79%)	45 (98%)	1 (2%)	52	55
2	L	46/58 (79%)	46 (100%)	0	100	100
All	All	1060/1136 (93%)	1058 (100%)	2 (0%)	93	95

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	118	GLN
2	K	29	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	118	GLN
1	B	57	ASN
1	B	67	ASN
1	B	101	ASN
1	C	57	ASN
1	C	67	ASN
1	C	101	ASN
1	C	118	GLN
1	D	57	ASN
1	E	57	ASN
1	E	67	ASN
1	E	101	ASN
1	E	118	GLN
1	F	57	ASN
1	G	37	HIS
1	G	57	ASN
1	G	118	GLN
1	H	57	ASN
1	H	67	ASN
1	H	101	ASN
2	I	75	GLN
2	J	75	GLN
2	K	75	GLN
2	L	75	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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

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	122/123 (99%)	0.26	9 (7%) 14 13	21, 34, 59, 79	0
1	B	117/123 (95%)	0.35	6 (5%) 28 27	23, 40, 67, 76	0
1	C	117/123 (95%)	0.35	10 (8%) 10 10	27, 41, 70, 84	0
1	D	121/123 (98%)	0.41	13 (10%) 6 5	26, 38, 68, 83	0
1	E	117/123 (95%)	0.47	13 (11%) 5 4	25, 38, 75, 89	0
1	F	119/123 (96%)	0.77	19 (15%) 1 1	28, 48, 80, 87	0
1	G	119/123 (96%)	0.60	16 (13%) 3 2	26, 45, 79, 84	0
1	H	117/123 (95%)	0.85	14 (11%) 4 3	24, 48, 82, 97	0
2	I	57/69 (82%)	1.10	10 (17%) 1 1	26, 52, 90, 93	0
2	J	56/69 (81%)	1.05	10 (17%) 1 1	32, 51, 91, 97	0
2	K	55/69 (79%)	2.09	19 (34%) 0 0	30, 64, 91, 93	0
2	L	55/69 (79%)	1.63	16 (29%) 0 0	32, 68, 99, 101	0
All	All	1172/1260 (93%)	0.69	155 (13%) 3 2	21, 43, 84, 101	0

All (155) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	K	64	SER	10.5
2	K	78	PHE	9.7
2	L	64	SER	9.3
2	K	70	VAL	9.1
2	K	66	VAL	8.8
2	K	67	ALA	8.1
2	I	64	SER	8.1
2	K	65	SER	8.0
2	J	23	VAL	7.9
2	J	75	GLN	7.7
1	H	21	ASP	7.6

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Mol	Chain	Res	Type	RSRZ
2	L	66	VAL	7.5
2	I	65	SER	7.3
1	H	57	ASN	7.0
2	I	67	ALA	6.9
1	E	57	ASN	6.8
2	I	66	VAL	6.8
2	K	68	HIS	6.7
2	K	75	GLN	6.6
1	G	100	ASN	6.3
1	H	55	ASP	6.2
2	L	78	PHE	5.9
2	J	78	PHE	5.8
1	E	53	SER	5.8
2	L	68	HIS	5.8
2	K	72	GLY	5.7
2	L	75	GLN	5.4
1	F	56	ASN	5.4
1	F	33	GLU	5.3
2	K	69	SER	5.3
1	G	116	ARG	5.1
1	H	22	THR	5.1
1	H	56	ASN	5.1
1	F	101	ASN	4.8
1	D	98	LEU	4.8
2	L	28	GLY	4.7
1	A	121	SER	4.5
2	L	65	SER	4.5
2	J	68	HIS	4.4
1	D	100	ASN	4.4
1	A	120	SER	4.4
2	J	66	VAL	4.3
1	A	57	ASN	4.2
1	F	57	ASN	4.2
2	K	74	ILE	4.2
1	E	55	ASP	4.1
1	C	57	ASN	4.1
1	E	56	ASN	4.1
1	G	33	GLU	4.0
1	B	21	ASP	4.0
2	L	67	ALA	3.9
1	F	16	SER	3.9
1	B	57	ASN	3.9

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Mol	Chain	Res	Type	RSRZ
1	F	55	ASP	3.9
1	F	99	ASP	3.9
2	K	77	MET	3.8
1	C	53	SER	3.7
2	J	65	SER	3.7
1	F	19	ILE	3.7
1	H	53	SER	3.6
2	K	28	GLY	3.6
1	C	21	ASP	3.6
2	K	27	SER	3.5
1	D	33	GLU	3.5
1	D	57	ASN	3.5
2	I	75	GLN	3.5
2	I	69	SER	3.5
1	E	54	LEU	3.4
2	L	69	SER	3.4
1	C	56	ASN	3.3
2	J	69	SER	3.3
1	G	56	ASN	3.3
1	C	55	ASP	3.3
1	F	116	ARG	3.3
1	G	57	ASN	3.3
1	F	100	ASN	3.2
1	C	54	LEU	3.2
1	H	83	VAL	3.2
1	F	59	GLU	3.2
1	G	96	ASN	3.2
1	E	59	GLU	3.1
2	J	67	ALA	3.1
1	G	98	LEU	3.1
1	G	55	ASP	3.1
1	F	103	LEU	3.1
1	A	100	ASN	3.0
1	B	38	ILE	2.9
1	G	101	ASN	2.9
1	F	53	SER	2.8
1	D	99	ASP	2.8
1	G	97	SER	2.8
1	H	118	GLN	2.8
1	F	98	LEU	2.8
1	E	82	GLU	2.7
1	G	38	ILE	2.7

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Mol	Chain	Res	Type	RSRZ
1	E	33	GLU	2.7
2	L	27	SER	2.7
1	F	18	SER	2.7
1	H	27	ILE	2.7
2	L	59	ILE	2.7
2	L	63	LEU	2.7
1	E	58	ASP	2.6
1	G	59	GLU	2.6
2	L	77	MET	2.6
1	A	99	ASP	2.6
1	D	101	ASN	2.6
1	B	24	GLU	2.6
1	F	107	LEU	2.6
1	C	24	GLU	2.6
1	E	21	ASP	2.6
1	A	56	ASN	2.5
2	K	59	ILE	2.5
2	I	78	PHE	2.5
1	D	38	ILE	2.5
1	H	59	GLU	2.5
2	L	73	PHE	2.4
1	D	55	ASP	2.4
1	D	21	ASP	2.4
1	E	47	LEU	2.4
2	K	36	SER	2.4
2	L	70	VAL	2.3
2	I	22	SER	2.3
1	D	107	LEU	2.3
1	H	38	ILE	2.3
1	D	116	ARG	2.3
1	F	118	GLN	2.2
1	A	110	LEU	2.2
1	H	64	LEU	2.2
2	I	61	GLU	2.2
1	E	83	VAL	2.2
2	I	68	HIS	2.2
2	K	73	PHE	2.2
2	J	27	SER	2.2
2	L	62	ARG	2.2
2	K	35	THR	2.2
2	K	63	LEU	2.2
1	B	18	SER	2.2

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Mol	Chain	Res	Type	RSRZ
1	G	53	SER	2.2
1	C	116	ARG	2.2
1	C	115	GLU	2.2
2	J	62	ARG	2.2
1	A	98	LEU	2.1
1	H	58	ASP	2.1
1	C	22	THR	2.1
1	D	59	GLU	2.1
1	G	103	LEU	2.1
1	G	71	GLN	2.1
1	D	56	ASN	2.1
1	H	23	ILE	2.1
1	B	32	GLY	2.1
1	E	118	GLN	2.1
1	A	101	ASN	2.0
1	F	21	ASP	2.0
1	F	54	LEU	2.0
1	G	99	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.