



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:06 AM GMT

PDB ID : 2ZJS  
Title : Crystal Structure of SecYE translocon from *Thermus thermophilus* with a Fab fragment  
Authors : Tsukazaki, T.; Mori, H.; Fukai, S.; Ishitani, R.; Perederina, A.; Vassylyev, D.G.; Ito, K.; Nureki, O.  
Deposited on : 2008-03-08  
Resolution : 3.20 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

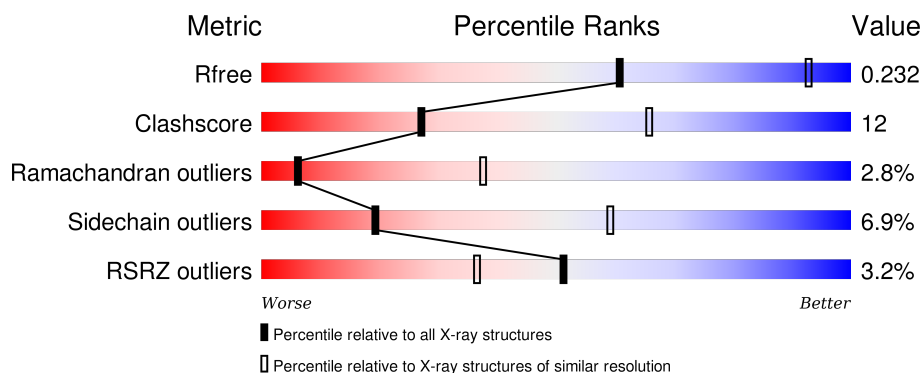
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.20 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	1124 (3.24-3.16)
Clashscore	102246	1024 (3.22-3.18)
Ramachandran outliers	100387	1004 (3.22-3.18)
Sidechain outliers	100360	1003 (3.22-3.18)
RSRZ outliers	91569	1129 (3.24-3.16)

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	Y	434	<div> <div>3%</div> <div>70% 22% . .</div> </div>
2	E	60	<div> <div>2%</div> <div>65% 12% 23%</div> </div>
3	H	221	<div> <div>4%</div> <div>69% 29% .</div> </div>
4	L	214	<div> <div>3%</div> <div>59% 34% 6% .</div> </div>

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6932 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 Preprotein translocase SecY subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	Y	415	Total	C	N	O	S	0	0	0
			3227	2155	529	537	6			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
Y	2	VAL	LEU	ENGINEERED	UNP Q8KZP3
Y	252	GLY	ARG	ENGINEERED	UNP Q8KZP3

- Molecule 2 is a protein called Preprotein translocase SecE subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	46	Total	C	N	O	S	0	0	0
			369	245	61	62	1			

- Molecule 3 is a protein called Fab56 (heavy chain).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	H	220	Total	C	N	O	S	0	0	0
			1667	1059	270	331	7			

- Molecule 4 is a protein called Fab56 (light chain).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	L	214	Total	C	N	O	S	0	0	0
			1667	1032	282	346	7			

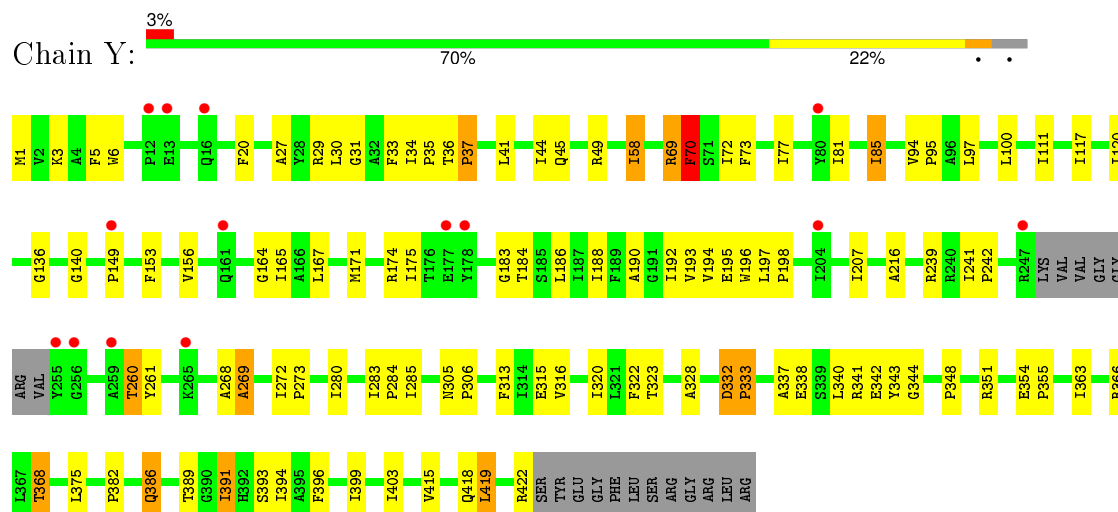
- Molecule 5 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	Y	1	Total 1	Zn 1	0	0
5	L	1	Total 1	Zn 1	0	0

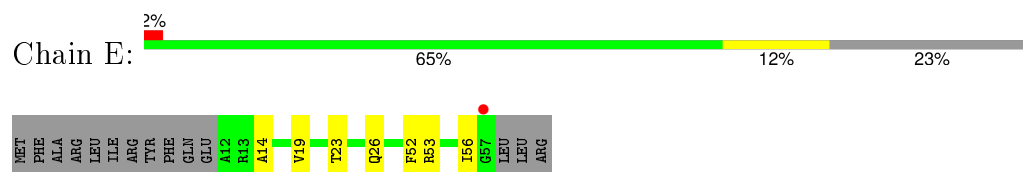
### 3 Residue-property plots

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.

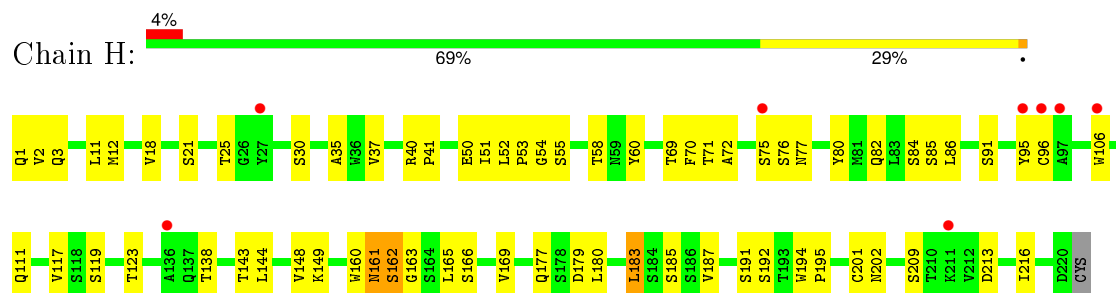
- Molecule 1: Preprotein translocase SecY subunit



- Molecule 2: Preprotein translocase SecE subunit

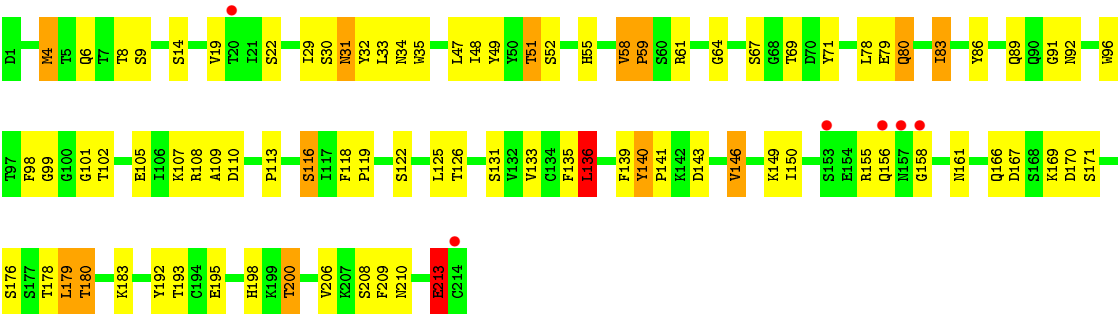


- Molecule 3: Fab56 (heavy chain)



- Molecule 4: Fab56 (light chain)





## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	189.10 Å   103.07 Å   78.00 Å 90.00°   105.18°   90.00°	Depositor
Resolution (Å)	48.43 – 3.20 48.43 – 3.19	Depositor EDS
% Data completeness (in resolution range)	98.6 (48.43-3.20) 93.6 (48.43-3.19)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.33 (at 3.19 Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.244   ,   0.280 0.244   ,   0.232	Depositor DCC
$R_{free}$ test set	1102 reflections (5.13%)	DCC
Wilson B-factor (Å <sup>2</sup> )	104.2	Xtriage
Anisotropy	0.374	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.26 , 91.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.28$	Xtriage
Outliers	0 of 22739 reflections	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	6932	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	127.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.03% 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ZN

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	Y	0.43	0/3306	0.54	0/4500
2	E	0.40	0/376	0.55	0/511
3	H	0.38	0/1716	0.56	0/2350
4	L	0.40	0/1702	0.62	0/2310
All	All	0.41	0/7100	0.56	0/9671

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
4	L	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
4	L	136	LEU	Peptide

### 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	Y	3227	0	3373	68	0
2	E	369	0	384	5	0
3	H	1667	0	1608	36	0
4	L	1667	0	1589	64	0
5	L	1	0	0	0	0
5	Y	1	0	0	0	0
All	All	6932	0	6954	166	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 12.

All (166) 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:Y:332:ASP:HB2	1:Y:333:PRO:HD3	1.24	1.15
1:Y:332:ASP:CB	1:Y:333:PRO:HD3	2.03	0.88
4:L:141:PRO:HD2	4:L:198:HIS:HE1	1.38	0.88
4:L:52:SER:HB2	4:L:64:GLY:HA3	1.58	0.84
1:Y:192:ILE:HD13	1:Y:403:ILE:HD11	1.59	0.84
1:Y:332:ASP:HB2	1:Y:333:PRO:CD	2.09	0.78
4:L:141:PRO:HD2	4:L:198:HIS:CE1	2.20	0.77
4:L:4:MET:HE1	4:L:29:ILE:HD11	1.65	0.77
3:H:106:TRP:HH2	4:L:96:TRP:CD1	2.02	0.76
4:L:136:LEU:HD11	4:L:146:VAL:HG13	1.65	0.76
1:Y:190:ALA:O	1:Y:194:VAL:HG23	1.86	0.75
1:Y:45:GLN:HB3	1:Y:49:ARG:HH21	1.51	0.75
1:Y:33:PHE:O	1:Y:35:PRO:HD3	1.89	0.72
1:Y:394:ILE:HD12	1:Y:394:ILE:H	1.54	0.72
4:L:167:ASP:HB3	4:L:170:ASP:O	1.91	0.70
3:H:71:THR:HB	3:H:80:TYR:HB2	1.72	0.70
4:L:108:ARG:HH12	4:L:170:ASP:HB2	1.59	0.67
3:H:37:VAL:HB	3:H:95:TYR:HB2	1.77	0.67
3:H:12:MET:CE	3:H:18:VAL:HG22	2.24	0.66
4:L:61:ARG:HH12	4:L:79:GLU:HB2	1.61	0.66
3:H:18:VAL:HG23	3:H:86:LEU:HD11	1.79	0.64
4:L:80:GLN:HE21	4:L:83:ILE:HG21	1.63	0.64
1:Y:175:ILE:HB	1:Y:183:GLY:HA2	1.80	0.63
4:L:195:GLU:HG2	4:L:206:VAL:HG22	1.79	0.63
4:L:55:HIS:HB3	4:L:58:VAL:HG12	1.81	0.62
4:L:198:HIS:ND1	4:L:200:THR:HB	2.14	0.62
1:Y:167:LEU:O	1:Y:171:MET:HG3	2.00	0.61
4:L:34:ASN:ND2	4:L:49:TYR:O	2.33	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:L:19:VAL:HG21	4:L:78:LEU:HD12	1.83	0.61
4:L:31:ASN:HB3	4:L:51:THR:OG1	2.00	0.60
4:L:6:GLN:HE21	4:L:99:GLY:HA3	1.66	0.60
4:L:198:HIS:HB3	4:L:200:THR:HG22	1.84	0.59
3:H:161:ASN:C	3:H:163:GLY:H	2.05	0.59
1:Y:272:ILE:N	1:Y:273:PRO:HD2	2.17	0.59
1:Y:29:ARG:CZ	1:Y:194:VAL:HG11	2.32	0.59
4:L:122:SER:HA	4:L:125:LEU:HD12	1.85	0.58
4:L:19:VAL:HG21	4:L:78:LEU:CD1	2.33	0.58
1:Y:338:GLU:O	1:Y:341:ARG:HB2	2.04	0.58
4:L:133:VAL:HG13	4:L:178:THR:HG22	1.86	0.57
4:L:80:GLN:O	4:L:83:ILE:HG22	2.04	0.57
4:L:140:TYR:O	4:L:141:PRO:C	2.42	0.57
3:H:12:MET:HE2	3:H:18:VAL:HG22	1.85	0.57
1:Y:366:ARG:HH11	2:E:14:ALA:HB3	1.70	0.57
2:E:52:PHE:O	2:E:56:ILE:HG13	2.05	0.56
1:Y:81:ILE:O	1:Y:85:ILE:HG22	2.06	0.56
1:Y:313:PHE:HA	1:Y:316:VAL:HG12	1.88	0.56
1:Y:332:ASP:CB	1:Y:333:PRO:CD	2.74	0.55
3:H:106:TRP:HH2	4:L:96:TRP:CG	2.25	0.55
4:L:105:GLU:HG3	4:L:166:GLN:NE2	2.22	0.54
1:Y:316:VAL:O	1:Y:320:ILE:HG12	2.07	0.54
1:Y:77:ILE:HG12	1:Y:184:THR:HG23	1.89	0.54
4:L:118:PHE:CE1	4:L:135:PHE:HD1	2.27	0.53
3:H:160:TRP:O	3:H:162:SER:N	2.42	0.53
3:H:30:SER:HA	3:H:53:PRO:HB2	1.90	0.53
3:H:106:TRP:CH2	4:L:96:TRP:CD1	2.92	0.53
3:H:91:SER:HB3	3:H:117:VAL:H	1.74	0.53
4:L:47:LEU:HD11	4:L:86:TYR:HE2	1.73	0.53
4:L:48:ILE:HG21	4:L:52:SER:C	2.29	0.53
1:Y:260:THR:HG22	1:Y:261:TYR:H	1.73	0.53
1:Y:136:GLY:HA2	1:Y:140:GLY:HA2	1.92	0.52
1:Y:58:ILE:HD11	1:Y:285:ILE:HD13	1.90	0.52
1:Y:241:ILE:HD11	1:Y:363:ILE:HG21	1.92	0.52
4:L:89:GLN:HB2	4:L:98:PHE:CD2	2.44	0.52
1:Y:216:ALA:HB1	1:Y:389:THR:HG23	1.93	0.51
4:L:193:THR:HA	4:L:208:SER:HB3	1.92	0.51
3:H:1:GLN:HG2	3:H:2:VAL:H	1.76	0.50
3:H:12:MET:HE3	3:H:18:VAL:HG22	1.91	0.50
1:Y:192:ILE:HD13	1:Y:403:ILE:CD1	2.36	0.50
4:L:136:LEU:HD11	4:L:146:VAL:CG1	2.38	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:L:58:VAL:HG22	4:L:59:PRO:HD2	1.92	0.50
4:L:150:ILE:CD1	4:L:155:ARG:HD3	2.42	0.50
1:Y:394:ILE:N	1:Y:394:ILE:HD12	2.25	0.50
4:L:118:PHE:HE1	4:L:135:PHE:HD1	1.59	0.49
1:Y:20:PHE:HZ	1:Y:174:ARG:HB3	1.78	0.49
3:H:149:LYS:HE2	3:H:177:GLN:OE1	2.12	0.49
1:Y:418:GLN:HE21	1:Y:422:ARG:HE	1.61	0.49
1:Y:354:GLU:HB3	1:Y:355:PRO:HD3	1.94	0.49
1:Y:386:GLN:HG2	1:Y:391:ILE:O	2.13	0.49
4:L:91:GLY:C	4:L:96:TRP:HZ3	2.16	0.49
3:H:53:PRO:HA	3:H:72:ALA:CB	2.43	0.49
1:Y:389:THR:HG22	1:Y:391:ILE:H	1.78	0.49
1:Y:415:VAL:O	1:Y:419:LEU:HB2	2.13	0.49
1:Y:72:ILE:O	1:Y:164:GLY:HA3	2.13	0.48
1:Y:328:ALA:HB2	1:Y:368:THR:HG21	1.95	0.48
4:L:149:LYS:HB2	4:L:193:THR:HB	1.96	0.48
3:H:165:LEU:HD13	3:H:187:VAL:HG21	1.95	0.48
1:Y:306:PRO:HB3	1:Y:315:GLU:OE1	2.14	0.48
3:H:53:PRO:HA	3:H:72:ALA:HB1	1.96	0.47
3:H:148:VAL:HB	3:H:183:LEU:HB3	1.96	0.47
4:L:110:ASP:HA	4:L:140:TYR:HB3	1.96	0.47
1:Y:193:VAL:O	1:Y:196:TRP:HB2	2.14	0.47
1:Y:382:PRO:O	1:Y:386:GLN:HB2	2.14	0.47
1:Y:337:ALA:HA	1:Y:340:LEU:HD12	1.97	0.47
3:H:70:PHE:HA	3:H:80:TYR:O	2.15	0.47
1:Y:100:LEU:HB2	1:Y:111:ILE:HD11	1.97	0.47
4:L:108:ARG:HG2	4:L:109:ALA:N	2.30	0.47
3:H:18:VAL:O	3:H:82:GLN:HA	2.15	0.47
4:L:67:SER:HA	4:L:71:TYR:CE2	2.50	0.47
1:Y:3:LYS:HA	1:Y:6:TRP:HB2	1.97	0.47
4:L:213:GLU:HG2	4:L:213:GLU:O	2.15	0.46
4:L:192:TYR:O	4:L:208:SER:HB2	2.15	0.46
1:Y:69:ARG:O	1:Y:70:PHE:C	2.53	0.46
1:Y:94:VAL:HA	1:Y:95:PRO:HD3	1.82	0.46
1:Y:41:LEU:HD22	1:Y:69:ARG:HA	1.98	0.46
3:H:194:TRP:O	3:H:195:PRO:C	2.53	0.46
3:H:160:TRP:HA	3:H:201:CYS:HA	1.97	0.45
1:Y:41:LEU:HD13	1:Y:69:ARG:HB2	1.99	0.45
1:Y:323:THR:HG21	1:Y:375:LEU:HD12	1.99	0.45
1:Y:20:PHE:CZ	1:Y:174:ARG:HB3	2.51	0.45
1:Y:394:ILE:CD1	1:Y:394:ILE:H	2.27	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:23:THR:HB	2:E:26:GLN:HB3	1.98	0.44
1:Y:27:ALA:HA	1:Y:30:LEU:HB3	1.99	0.44
4:L:150:ILE:HD12	4:L:155:ARG:HD3	1.99	0.44
3:H:161:ASN:C	3:H:163:GLY:N	2.70	0.44
4:L:29:ILE:HG23	4:L:92:ASN:HB2	1.98	0.44
1:Y:97:LEU:HA	1:Y:100:LEU:HD12	1.98	0.44
1:Y:94:VAL:HB	1:Y:97:LEU:HD12	1.99	0.44
4:L:33:LEU:HD13	4:L:71:TYR:CD1	2.53	0.44
1:Y:351:ARG:NH1	3:H:50:GLU:OE2	2.51	0.43
4:L:118:PHE:CE1	4:L:135:PHE:CD1	3.06	0.43
4:L:116:SER:HB2	4:L:118:PHE:CE1	2.54	0.43
4:L:119:PRO:HB3	4:L:209:PHE:CE1	2.54	0.43
3:H:202:ASN:OD1	3:H:213:ASP:HB3	2.17	0.43
1:Y:36:THR:HA	1:Y:37:PRO:HD2	1.80	0.43
3:H:58:THR:HG1	3:H:60:TYR:HE1	1.64	0.43
3:H:160:TRP:HZ3	3:H:216:ILE:HD11	1.83	0.43
4:L:158:GLY:O	4:L:179:LEU:HA	2.19	0.43
3:H:51:ILE:HG13	3:H:58:THR:HG22	2.00	0.43
1:Y:117:ILE:HA	1:Y:120:ILE:HG22	2.01	0.42
4:L:107:LYS:HD2	4:L:107:LYS:HA	1.92	0.42
4:L:35:TRP:CD1	4:L:52:SER:HB3	2.53	0.42
2:E:53:ARG:HH11	2:E:56:ILE:HD11	1.83	0.42
4:L:166:GLN:NE2	4:L:171:SER:HB2	2.35	0.42
4:L:86:TYR:O	4:L:101:GLY:HA2	2.20	0.42
1:Y:305:ASN:HA	1:Y:306:PRO:HD3	1.83	0.42
1:Y:393:SER:HG	1:Y:396:PHE:HD1	1.68	0.42
1:Y:280:ILE:HG13	1:Y:322:PHE:CZ	2.55	0.41
4:L:210:ASN:HB2	4:L:213:GLU:HB3	2.02	0.41
4:L:30:SER:O	4:L:32:TYR:CD1	2.73	0.41
1:Y:268:ALA:O	1:Y:269:ALA:C	2.59	0.41
1:Y:283:ILE:HB	1:Y:284:PRO:HD3	2.01	0.41
1:Y:167:LEU:HD22	1:Y:171:MET:SD	2.60	0.41
1:Y:272:ILE:N	1:Y:273:PRO:CD	2.82	0.41
1:Y:174:ARG:HA	1:Y:174:ARG:HD3	1.81	0.41
4:L:113:PRO:HG3	4:L:139:PHE:HB3	2.02	0.41
1:Y:188:ILE:HG22	1:Y:403:ILE:HD13	2.02	0.41
4:L:9:SER:O	4:L:102:THR:HA	2.20	0.41
4:L:155:ARG:HG3	4:L:156:GLN:N	2.35	0.41
4:L:140:TYR:HB3	4:L:141:PRO:HD3	2.02	0.41
3:H:166:SER:O	3:H:169:VAL:HG22	2.21	0.41
1:Y:239:ARG:HB2	2:E:19:VAL:HG22	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Y:153:PHE:HA	1:Y:156:VAL:HG12	2.01	0.41
3:H:179:ASP:OD1	3:H:180:LEU:HG	2.21	0.41
4:L:135:PHE:C	4:L:136:LEU:HD23	2.42	0.41
4:L:83:ILE:HG23	4:L:83:ILE:O	2.21	0.41
3:H:149:LYS:HE3	4:L:180:THR:HG21	2.02	0.41
4:L:118:PHE:HA	4:L:119:PRO:HD3	1.85	0.40
3:H:144:LEU:HB3	3:H:216:ILE:HD13	2.03	0.40
1:Y:31:GLY:HA2	1:Y:34:ILE:HG12	2.03	0.40
3:H:35:ALA:O	3:H:96:CYS:HA	2.21	0.40
4:L:125:LEU:HB3	4:L:183:LYS:HE3	2.04	0.40
3:H:191:SER:O	3:H:195:PRO:HD2	2.21	0.40
1:Y:393:SER:OG	1:Y:396:PHE:HD1	2.04	0.40
3:H:52:LEU:O	3:H:55:SER:O	2.38	0.40
1:Y:242:PRO:O	1:Y:348:PRO:HD2	2.21	0.40
1:Y:197:LEU:HB3	1:Y:198:PRO:HD3	2.03	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	Y	411/434 (95%)	373 (91%)	29 (7%)	9 (2%)	8	45
2	E	44/60 (73%)	39 (89%)	5 (11%)	0	100	100
3	H	218/221 (99%)	190 (87%)	19 (9%)	9 (4%)	3	27
4	L	212/214 (99%)	195 (92%)	10 (5%)	7 (3%)	5	32
All	All	885/929 (95%)	797 (90%)	63 (7%)	25 (3%)	6	37

All (25) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	Y	70	PHE
1	Y	269	ALA
1	Y	332	ASP
3	H	162	SER
4	L	213	GLU
1	Y	344	GLY
1	Y	368	THR
3	H	54	GLY
3	H	85	SER
3	H	161	ASN
3	H	209	SER
4	L	31	ASN
4	L	83	ILE
4	L	169	LYS
1	Y	149	PRO
3	H	41	PRO
3	H	75	SER
4	L	140	TYR
1	Y	37	PRO
1	Y	342	GLU
3	H	84	SER
4	L	51	THR
3	H	77	ASN
1	Y	333	PRO
4	L	59	PRO

### 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	Y	332/347 (96%)	314 (95%)	18 (5%)	27	68
2	E	38/51 (74%)	38 (100%)	0	100	100
3	H	187/188 (100%)	172 (92%)	15 (8%)	15	52
4	L	192/192 (100%)	173 (90%)	19 (10%)	10	38
All	All	749/778 (96%)	697 (93%)	52 (7%)	19	59

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

Mol	Chain	Res	Type
1	Y	1	MET
1	Y	5	PHE
1	Y	44	ILE
1	Y	58	ILE
1	Y	69	ARG
1	Y	70	PHE
1	Y	73	PHE
1	Y	85	ILE
1	Y	165	ILE
1	Y	186	LEU
1	Y	195	GLU
1	Y	207	ILE
1	Y	260	THR
1	Y	343	TYR
1	Y	386	GLN
1	Y	391	ILE
1	Y	399	ILE
1	Y	419	LEU
3	H	3	GLN
3	H	11	LEU
3	H	21	SER
3	H	25	THR
3	H	40	ARG
3	H	69	THR
3	H	76	SER
3	H	111	GLN
3	H	119	SER
3	H	123	THR
3	H	138	THR
3	H	143	THR
3	H	183	LEU
3	H	185	SER
3	H	192	SER
4	L	4	MET
4	L	8	THR
4	L	14	SER
4	L	22	SER
4	L	58	VAL
4	L	69	THR
4	L	80	GLN
4	L	116	SER
4	L	126	THR

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Mol	Chain	Res	Type
4	L	131	SER
4	L	136	LEU
4	L	143	ASP
4	L	146	VAL
4	L	161	ASN
4	L	176	SER
4	L	179	LEU
4	L	180	THR
4	L	200	THR
4	L	213	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	Y	302	ASN
1	Y	392	HIS
1	Y	418	GLN
3	H	161	ASN
4	L	6	GLN
4	L	77	ASN
4	L	80	GLN
4	L	190	ASN
4	L	210	ASN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

### 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

### 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

### 5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.



There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 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, 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	Y	415/434 (95%)	0.12	14 (3%) 49 34	105, 127, 156, 163	0
2	E	46/60 (76%)	-0.03	1 (2%) 65 50	133, 149, 163, 164	0
3	H	220/221 (99%)	0.03	8 (3%) 46 31	112, 125, 136, 145	0
4	L	214/214 (100%)	0.06	6 (2%) 56 42	104, 120, 128, 139	0
All	All	895/929 (96%)	0.08	29 (3%) 51 36	104, 124, 155, 164	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	E	57	GLY	4.6
4	L	157	ASN	4.6
1	Y	12	PRO	3.3
4	L	158	GLY	3.2
1	Y	178	TYR	2.9
4	L	214	CYS	2.8
1	Y	177	GLU	2.8
3	H	97	ALA	2.8
1	Y	16	GLN	2.8
1	Y	255	TYR	2.7
1	Y	13	GLU	2.6
3	H	136	ALA	2.6
3	H	75	SER	2.4
3	H	96	CYS	2.4
3	H	95	TYR	2.4
1	Y	259	ALA	2.3
3	H	106	TRP	2.3
1	Y	149	PRO	2.3
1	Y	247	ARG	2.3
1	Y	265	LYS	2.2
3	H	211	LYS	2.2

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Mol	Chain	Res	Type	RSRZ
1	Y	161	GLN	2.2
1	Y	80	TYR	2.1
1	Y	256	GLY	2.1
4	L	20	THR	2.1
4	L	156	GLN	2.1
3	H	27	TYR	2.1
1	Y	204	ILE	2.0
4	L	153	SER	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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
5	ZN	Y	435	1/1	0.99	0.10	-	89,89,89,89	1
5	ZN	L	223	1/1	0.98	0.20	-	97,97,97,97	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.