



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 08:37 PM GMT

PDB ID : 1L4A
Title : X-RAY STRUCTURE OF THE NEURONAL COMPLEXIN/SNARE COMPLEX FROM THE SQUID LOLIGO PEALEI
Authors : Bracher, A.; Kadlec, J.; Betz, H.; Weissenhorn, W.
Deposited on : 2002-03-04
Resolution : 2.95 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

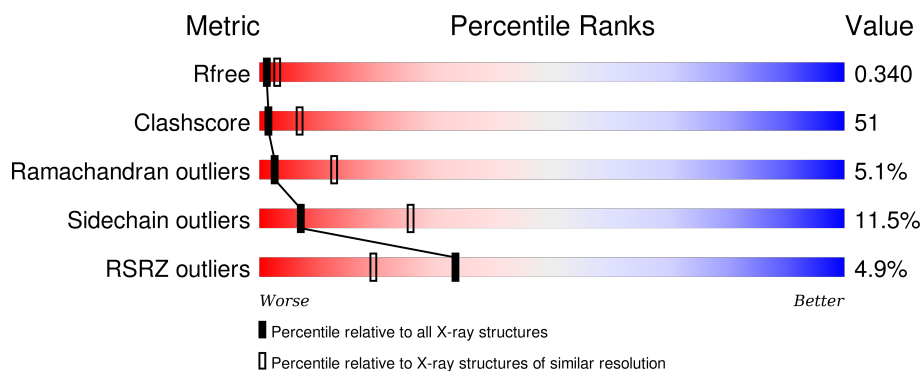
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.95 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	2184 (3.00-2.92)
Clashscore	102246	2552 (3.00-2.92)
Ramachandran outliers	100387	2468 (3.00-2.92)
Sidechain outliers	100360	2471 (3.00-2.92)
RSRZ outliers	91569	2201 (3.00-2.92)

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	80	<div> <div>10%</div> <div>25%</div> <div>43%</div> <div>15%</div> <div>18%</div> </div>
2	B	88	<div> <div>3%</div> <div>36%</div> <div>50%</div> <div>7%</div> <div>7%</div> </div>
3	C	83	<div> <div>%</div> <div>31%</div> <div>47%</div> <div>10%</div> <div>12%</div> </div>
4	D	87	<div> <div>2%</div> <div>40%</div> <div>39%</div> <div>7%</div> <div>14%</div> </div>
5	E	79	<div> <div>3%</div> <div>23%</div> <div>13%</div> <div>65%</div> </div>

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 2381 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 SYNAPTOBREVIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	66	Total	C	N	O	S	0	0	0
			485	294	89	101	1			

- Molecule 2 is a protein called S-SYNTAXIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	82	Total	C	N	O	S	0	0	0
			596	364	103	122	7			

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	178	GLY	-	CLONING ARTIFACT	UNP O46345
B	179	LYS	-	CLONING ARTIFACT	UNP O46345
B	180	SER	-	CLONING ARTIFACT	UNP O46345
B	181	ALA	-	CLONING ARTIFACT	UNP O46345
B	182	SER	-	CLONING ARTIFACT	UNP O46345

- Molecule 3 is a protein called S-SNAP25 fusion protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	73	Total	C	N	O	S	0	0	0
			535	314	97	118	6			

- Molecule 4 is a protein called S-SNAP25 fusion protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	75	Total	C	N	O	S	0	0	0
			547	321	102	119	5			

- Molecule 5 is a protein called SYNAPHIN A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	E	28	Total	C	N	O	S	0	0	0
			195	118	38	38	1			

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	20	GLY	-	CLONING ARTIFACT	UNP Q95PA1
E	21	LYS	-	CLONING ARTIFACT	UNP Q95PA1
E	22	SER	-	CLONING ARTIFACT	UNP Q95PA1
E	23	ALA	-	CLONING ARTIFACT	UNP Q95PA1
E	24	SER	-	CLONING ARTIFACT	UNP Q95PA1

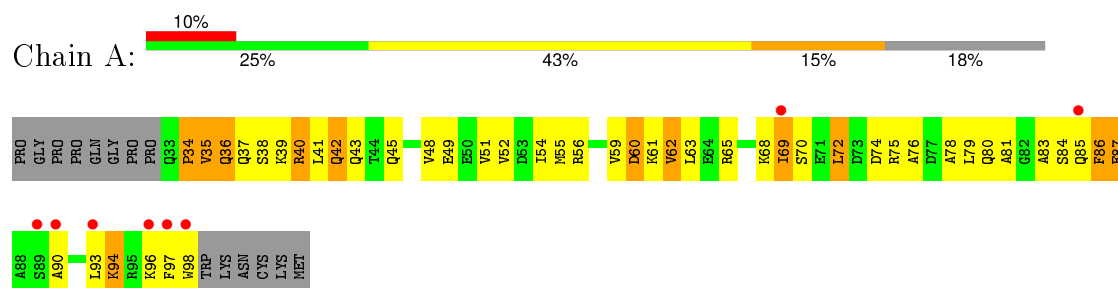
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	6	Total	O	0	0
			6	6		
6	B	3	Total	O	0	0
			3	3		
6	C	4	Total	O	0	0
			4	4		
6	D	10	Total	O	0	0
			10	10		

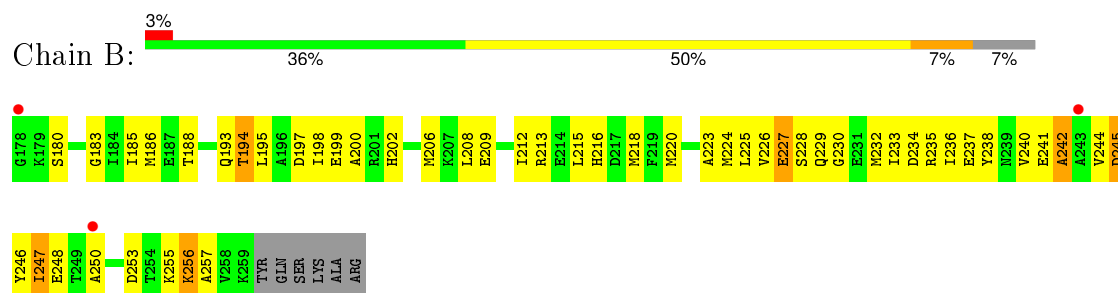
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

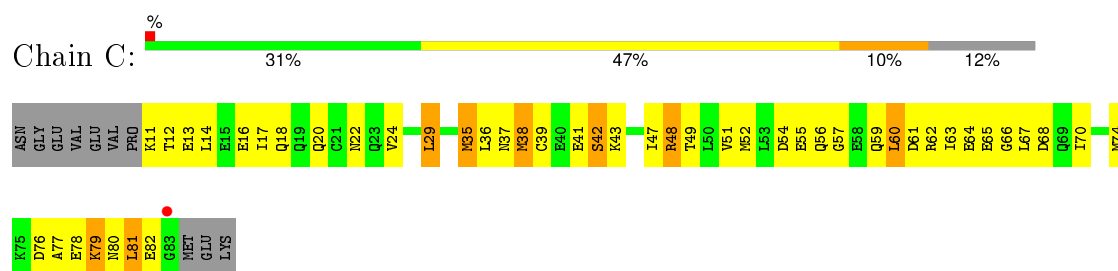
• Molecule 1: SYNAPTOBREVIN



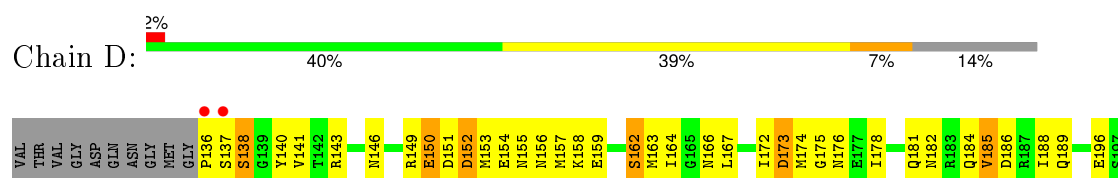
• Molecule 2: S-SYNTAXIN

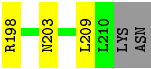


• Molecule 3: S-SNAP25 fusion protein

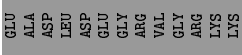
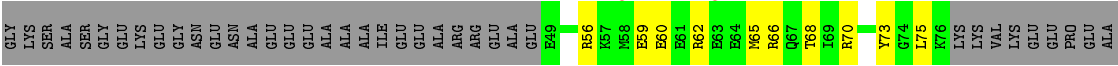


• Molecule 4: S-SNAP25 fusion protein





● Molecule 5: SYNAPHIN A



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	26.07Å 75.43Å 243.11Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 2.95 47.33 – 2.80	Depositor EDS
% Data completeness (in resolution range)	92.6 (15.00-2.95) 91.3 (47.33-2.80)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.23 (at 2.81Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.297 , 0.344 0.298 , 0.340	Depositor DCC
R_{free} test set	700 reflections (7.00%)	DCC
Wilson B-factor (Å ²)	67.3	Xtriage
Anisotropy	0.902	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 93.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	2 of 11610 reflections (0.017%)	Xtriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	2381	wwPDB-VP
Average B, all atoms (Å ²)	80.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.08% 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 [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	0/486	0.67	0/656
2	B	0.37	0/599	0.61	0/804
3	C	0.38	0/534	0.67	0/717
4	D	0.42	0/548	0.75	1/734 (0.1%)
5	E	0.34	0/195	0.51	0/261
All	All	0.38	0/2362	0.66	1/3172 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
4	D	138	SER	N-CA-C	-5.19	96.98	111.00

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	485	0	457	79	0
2	B	596	0	556	74	0
3	C	535	0	475	75	0
4	D	547	0	493	70	0
5	E	195	0	153	15	0
6	A	6	0	0	2	0
6	B	3	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	C	4	0	0	0	0
6	D	10	0	0	2	0
All	All	2381	0	2134	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 51.

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:94:LYS:HA	2:B:256:LYS:HE3	1.20	1.11
4:D:157:MET:HE2	4:D:157:MET:HA	1.35	1.07
5:E:66:ARG:HB3	5:E:70:ARG:HH12	1.35	0.91
1:A:65:ARG:HB2	4:D:182:ASN:OD1	1.71	0.90
1:A:75:ARG:HA	1:A:78:ALA:HB3	1.52	0.89
1:A:35:VAL:HG23	1:A:36:GLN:H	1.37	0.89
1:A:55:MET:O	1:A:59:VAL:HG23	1.73	0.89
2:B:233:ILE:HG22	3:C:59:GLN:HE21	1.36	0.88
3:C:22:ASN:ND2	4:D:149:ARG:HH22	1.72	0.86
5:E:62:ARG:O	5:E:66:ARG:HG3	1.76	0.86
3:C:22:ASN:HD21	4:D:149:ARG:HH22	1.21	0.85
4:D:157:MET:HE2	4:D:157:MET:CA	2.06	0.84
1:A:74:ASP:HA	6:A:19:HOH:O	1.78	0.84
2:B:233:ILE:HG22	3:C:59:GLN:NE2	1.93	0.83
3:C:36:LEU:HD23	4:D:156:ASN:HD22	1.44	0.83
1:A:48:VAL:CG2	4:D:164:ILE:HD13	2.08	0.83
4:D:157:MET:CE	4:D:157:MET:HA	2.10	0.82
1:A:38:SER:O	1:A:39:LYS:HG2	1.79	0.82
1:A:48:VAL:HG22	4:D:164:ILE:HD13	1.61	0.82
1:A:39:LYS:HB3	1:A:42:GLN:HG2	1.60	0.81
5:E:56:ARG:O	5:E:60:GLU:HG2	1.83	0.79
1:A:39:LYS:HB3	1:A:42:GLN:CG	2.13	0.78
3:C:36:LEU:CD2	4:D:156:ASN:HD22	1.97	0.77
1:A:69:ILE:HD11	2:B:233:ILE:HG13	1.68	0.76
1:A:69:ILE:HD11	2:B:233:ILE:CG1	2.14	0.76
3:C:36:LEU:HD11	4:D:159:GLU:OE2	1.87	0.75
3:C:80:ASN:C	3:C:82:GLU:H	1.92	0.72
2:B:212:ILE:HD12	3:C:35:MET:HG2	1.72	0.71
5:E:70:ARG:HG2	5:E:75:LEU:HB2	1.72	0.71
3:C:65:GLU:O	3:C:68:ASP:HB2	1.90	0.71
1:A:72:LEU:HD11	3:C:63:ILE:HD11	1.72	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:195:LEU:O	2:B:199:GLU:HG3	1.92	0.68
1:A:83:ALA:C	1:A:85:GLN:H	1.95	0.68
1:A:35:VAL:HG23	1:A:36:GLN:N	2.10	0.67
2:B:246:TYR:C	2:B:248:GLU:H	1.96	0.66
3:C:76:ASP:O	3:C:79:LYS:HG3	1.95	0.66
1:A:48:VAL:O	1:A:52:VAL:HG23	1.94	0.66
1:A:61:LYS:HB3	4:D:178:ILE:HG21	1.78	0.66
1:A:86:PHE:CE1	4:D:203:ASN:HA	2.31	0.66
2:B:197:ASP:O	2:B:200:ALA:HB3	1.97	0.65
2:B:229:GLN:O	2:B:233:ILE:HG13	1.96	0.65
1:A:61:LYS:HB3	4:D:178:ILE:CG2	2.27	0.65
3:C:63:ILE:O	3:C:67:LEU:HG	1.97	0.64
1:A:42:GLN:OE1	4:D:136:PRO:HB2	1.97	0.64
3:C:63:ILE:HG23	3:C:64:GLU:N	2.11	0.64
2:B:255:LYS:C	2:B:257:ALA:H	2.01	0.64
2:B:242:ALA:HA	2:B:245:ASP:HB3	1.79	0.64
3:C:43:LYS:O	3:C:47:ILE:HG12	1.98	0.63
2:B:228:SER:C	2:B:230:GLY:H	2.02	0.62
3:C:57:GLY:HA3	4:D:181:GLN:HE22	1.64	0.61
4:D:158:LYS:O	4:D:162:SER:HB2	2.00	0.61
2:B:202:HIS:CE1	2:B:206:MET:HE2	2.36	0.61
3:C:36:LEU:HD23	4:D:156:ASN:ND2	2.15	0.61
2:B:234:ASP:O	2:B:237:GLU:HB3	2.01	0.61
1:A:87:GLU:HA	2:B:250:ALA:CB	2.31	0.60
1:A:62:VAL:CG1	2:B:225:LEU:HD13	2.32	0.60
3:C:14:LEU:O	3:C:18:GLN:HG3	2.02	0.60
1:A:42:GLN:O	1:A:45:GLN:HB3	2.01	0.60
1:A:59:VAL:O	1:A:62:VAL:HG12	2.02	0.60
1:A:39:LYS:CE	4:D:136:PRO:HD2	2.31	0.60
1:A:94:LYS:HA	2:B:256:LYS:CE	2.14	0.60
1:A:35:VAL:CG2	1:A:36:GLN:H	2.08	0.60
4:D:137:SER:O	4:D:138:SER:HB3	2.03	0.59
2:B:244:VAL:HG13	3:C:70:ILE:HA	1.84	0.59
2:B:244:VAL:HG21	3:C:66:GLY:O	2.03	0.58
5:E:66:ARG:HG2	5:E:66:ARG:HH11	1.69	0.58
1:A:39:LYS:HE2	4:D:136:PRO:HD2	1.85	0.58
1:A:87:GLU:HA	2:B:250:ALA:HB2	1.84	0.58
2:B:247:ILE:HD11	3:C:74:MET:SD	2.44	0.57
2:B:202:HIS:CE1	2:B:206:MET:CE	2.87	0.57
2:B:228:SER:C	2:B:230:GLY:N	2.56	0.57
4:D:141:VAL:HA	6:D:12:HOH:O	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:22:ASN:ND2	4:D:149:ARG:NH2	2.50	0.57
2:B:244:VAL:HA	2:B:247:ILE:HG22	1.87	0.56
2:B:227:GLU:HG3	2:B:227:GLU:O	2.02	0.56
3:C:11:LYS:C	3:C:13:GLU:H	2.09	0.56
3:C:49:THR:HG21	4:D:174:MET:CE	2.35	0.56
1:A:54:ILE:HD13	4:D:172:ILE:HD11	1.88	0.56
3:C:52:MET:O	3:C:56:GLN:HG3	2.05	0.56
2:B:213:ARG:HG2	3:C:38:MET:SD	2.46	0.55
4:D:182:ASN:O	4:D:185:VAL:HG12	2.06	0.55
1:A:80:GLN:NE2	1:A:80:GLN:O	2.40	0.55
1:A:72:LEU:HD11	3:C:63:ILE:CD1	2.36	0.55
2:B:220:MET:HE1	3:C:41:GLU:O	2.07	0.55
1:A:83:ALA:C	1:A:85:GLN:N	2.60	0.54
3:C:12:THR:O	3:C:12:THR:HG22	2.07	0.54
2:B:230:GLY:HA2	3:C:56:GLN:NE2	2.22	0.54
1:A:60:ASP:O	1:A:63:LEU:HB2	2.08	0.54
1:A:86:PHE:HA	4:D:203:ASN:HD22	1.71	0.54
4:D:196:GLU:C	4:D:198:ARG:H	2.11	0.54
3:C:16:GLU:HG2	3:C:20:GLN:HE21	1.73	0.54
3:C:29:LEU:HD21	4:D:152:ASP:HB3	1.90	0.54
4:D:186:ASP:O	4:D:189:GLN:N	2.41	0.54
1:A:98:TRP:CB	2:B:256:LYS:HZ2	2.21	0.54
3:C:35:MET:CE	4:D:157:MET:CE	2.86	0.53
1:A:39:LYS:O	1:A:42:GLN:HG2	2.08	0.53
3:C:79:LYS:C	3:C:81:LEU:H	2.10	0.53
1:A:72:LEU:O	1:A:76:ALA:HB2	2.08	0.53
3:C:80:ASN:C	3:C:82:GLU:N	2.62	0.53
1:A:56:ARG:HG2	2:B:218:MET:HE1	1.91	0.53
1:A:38:SER:C	4:D:137:SER:CB	2.77	0.53
2:B:236:ILE:O	2:B:240:VAL:HG23	2.08	0.53
1:A:69:ILE:HD13	2:B:229:GLN:HB3	1.91	0.53
1:A:86:PHE:HE1	4:D:203:ASN:HA	1.72	0.53
5:E:65:MET:HA	5:E:68:THR:CG2	2.39	0.52
3:C:35:MET:CE	4:D:157:MET:HE1	2.40	0.52
2:B:228:SER:HB3	5:E:62:ARG:HH22	1.75	0.52
4:D:182:ASN:HB3	6:D:15:HOH:O	2.10	0.52
1:A:42:GLN:CD	1:A:42:GLN:H	2.10	0.52
3:C:39:CYS:HB3	4:D:163:MET:HE3	1.91	0.52
1:A:83:ALA:O	1:A:85:GLN:N	2.43	0.51
4:D:163:MET:O	4:D:167:LEU:HD13	2.10	0.51
4:D:184:GLN:O	4:D:188:ILE:HG12	2.09	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:74:MET:C	3:C:76:ASP:H	2.13	0.51
1:A:72:LEU:HD11	4:D:188:ILE:HG21	1.92	0.51
4:D:151:ASP:O	4:D:155:ASN:ND2	2.44	0.50
3:C:60:LEU:C	3:C:63:ILE:HG22	2.32	0.50
1:A:45:GLN:O	1:A:49:GLU:HG3	2.10	0.50
3:C:63:ILE:CG2	3:C:64:GLU:N	2.74	0.50
2:B:186:MET:C	2:B:188:THR:H	2.13	0.50
1:A:61:LYS:CB	4:D:178:ILE:HG21	2.39	0.50
1:A:45:GLN:HA	2:B:208:LEU:HD13	1.92	0.50
2:B:242:ALA:CA	2:B:245:ASP:HB3	2.41	0.49
1:A:90:ALA:HA	1:A:93:LEU:HB3	1.92	0.49
1:A:40:ARG:H	4:D:137:SER:CB	2.25	0.49
1:A:68:LYS:C	1:A:70:SER:H	2.15	0.49
1:A:42:GLN:OE1	4:D:136:PRO:CB	2.61	0.49
1:A:90:ALA:HB1	2:B:253:ASP:HB3	1.95	0.49
1:A:35:VAL:O	1:A:36:GLN:HB2	2.13	0.49
2:B:186:MET:C	2:B:188:THR:N	2.64	0.49
2:B:194:THR:O	2:B:198:ILE:HG13	2.13	0.49
1:A:49:GLU:HA	1:A:52:VAL:HG23	1.94	0.48
2:B:223:ALA:O	2:B:227:GLU:HB3	2.13	0.48
1:A:93:LEU:HG	2:B:256:LYS:HD2	1.96	0.48
3:C:22:ASN:ND2	4:D:149:ARG:HH12	2.12	0.47
1:A:79:LEU:HD11	4:D:196:GLU:HA	1.95	0.47
3:C:63:ILE:CD1	4:D:188:ILE:HG21	2.43	0.47
3:C:42:SER:HB3	4:D:167:LEU:HD21	1.97	0.47
3:C:66:GLY:C	3:C:68:ASP:H	2.18	0.47
1:A:62:VAL:HG11	2:B:225:LEU:HD13	1.96	0.47
2:B:244:VAL:HG22	3:C:70:ILE:CB	2.45	0.47
3:C:59:GLN:C	3:C:61:ASP:H	2.18	0.47
1:A:68:LYS:HD3	4:D:186:ASP:OD1	2.14	0.47
2:B:246:TYR:C	2:B:248:GLU:N	2.63	0.46
3:C:14:LEU:C	3:C:18:GLN:HE21	2.19	0.46
1:A:74:ASP:O	1:A:78:ALA:N	2.44	0.46
2:B:216:HIS:HA	3:C:42:SER:OG	2.15	0.46
3:C:76:ASP:C	3:C:78:GLU:H	2.18	0.46
3:C:64:GLU:O	3:C:68:ASP:CG	2.54	0.46
1:A:54:ILE:HD13	4:D:172:ILE:CD1	2.45	0.46
2:B:220:MET:O	2:B:224:MET:HG3	2.15	0.46
1:A:93:LEU:O	2:B:256:LYS:NZ	2.33	0.46
2:B:209:GLU:O	2:B:209:GLU:HG2	2.15	0.46
1:A:39:LYS:HB3	1:A:42:GLN:HG3	1.95	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:51:VAL:O	3:C:54:ASP:HB3	2.15	0.46
1:A:69:ILE:HG23	2:B:232:MET:HE3	1.97	0.46
1:A:39:LYS:N	4:D:137:SER:CB	2.79	0.46
1:A:40:ARG:HG2	4:D:137:SER:CB	2.45	0.46
3:C:63:ILE:HD13	4:D:188:ILE:HG21	1.97	0.46
1:A:69:ILE:HD11	2:B:233:ILE:HG12	1.95	0.45
3:C:77:ALA:O	3:C:80:ASN:ND2	2.49	0.45
2:B:255:LYS:C	2:B:257:ALA:N	2.68	0.45
4:D:137:SER:O	4:D:138:SER:CB	2.63	0.45
3:C:74:MET:C	3:C:76:ASP:N	2.70	0.45
4:D:153:MET:O	4:D:157:MET:HG2	2.16	0.45
4:D:159:GLU:O	4:D:163:MET:HG3	2.16	0.45
2:B:255:LYS:O	2:B:257:ALA:N	2.46	0.45
5:E:65:MET:HA	5:E:68:THR:HG22	1.99	0.45
2:B:183:GLY:C	2:B:185:ILE:H	2.20	0.45
3:C:57:GLY:CA	4:D:181:GLN:HE22	2.28	0.45
2:B:235:ARG:C	2:B:237:GLU:H	2.20	0.45
1:A:97:PHE:C	2:B:256:LYS:HZ3	2.20	0.45
2:B:233:ILE:O	2:B:233:ILE:HG22	2.17	0.44
3:C:48:ARG:O	3:C:52:MET:HG3	2.17	0.44
4:D:157:MET:HE2	4:D:157:MET:N	2.31	0.44
3:C:62:ARG:O	3:C:63:ILE:C	2.56	0.44
2:B:235:ARG:C	2:B:237:GLU:N	2.71	0.44
4:D:157:MET:C	4:D:159:GLU:N	2.70	0.44
4:D:150:GLU:O	4:D:151:ASP:C	2.56	0.44
1:A:81:ALA:HA	6:A:9:HOH:O	2.16	0.44
4:D:143:ARG:HD3	4:D:143:ARG:N	2.33	0.43
1:A:65:ARG:HG2	2:B:229:GLN:OE1	2.18	0.43
4:D:174:MET:O	4:D:178:ILE:HG12	2.18	0.43
1:A:69:ILE:O	1:A:69:ILE:HG12	2.19	0.43
1:A:75:ARG:O	1:A:79:LEU:HB2	2.17	0.43
5:E:60:GLU:HA	5:E:60:GLU:OE1	2.18	0.43
1:A:61:LYS:HB3	4:D:178:ILE:HG22	1.99	0.43
4:D:173:ASP:O	4:D:176:ASN:N	2.51	0.43
2:B:226:VAL:O	2:B:226:VAL:HG12	2.19	0.43
2:B:198:ILE:HG22	3:C:24:VAL:HG11	2.00	0.43
2:B:238:TYR:O	2:B:241:GLU:HB2	2.18	0.43
4:D:150:GLU:O	4:D:153:MET:N	2.51	0.43
4:D:164:ILE:C	4:D:166:ASN:N	2.71	0.43
3:C:76:ASP:C	3:C:78:GLU:N	2.72	0.43
2:B:232:MET:HE2	5:E:62:ARG:CZ	2.48	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:242:ALA:O	2:B:246:TYR:N	2.52	0.43
2:B:246:TYR:O	2:B:248:GLU:N	2.52	0.43
2:B:240:VAL:C	2:B:242:ALA:H	2.22	0.42
3:C:63:ILE:HG23	3:C:64:GLU:H	1.80	0.42
3:C:13:GLU:O	3:C:17:ILE:HG13	2.20	0.42
3:C:63:ILE:CG2	3:C:64:GLU:H	2.33	0.42
3:C:79:LYS:C	3:C:79:LYS:HD2	2.39	0.42
3:C:77:ALA:C	3:C:79:LYS:H	2.22	0.42
4:D:140:TYR:HD1	4:D:154:GLU:OE2	2.02	0.42
3:C:80:ASN:O	3:C:82:GLU:N	2.53	0.42
5:E:65:MET:C	5:E:68:THR:HG22	2.40	0.42
1:A:79:LEU:HA	1:A:79:LEU:HD12	1.76	0.41
3:C:77:ALA:HA	3:C:79:LYS:HE2	2.02	0.41
4:D:174:MET:HG2	4:D:175:GLY:N	2.35	0.41
4:D:149:ARG:O	4:D:153:MET:HG3	2.20	0.41
5:E:70:ARG:HA	5:E:75:LEU:HD12	2.02	0.41
1:A:90:ALA:HB1	2:B:253:ASP:CB	2.50	0.41
3:C:59:GLN:O	3:C:61:ASP:N	2.54	0.41
2:B:244:VAL:HG13	3:C:70:ILE:CA	2.50	0.41
2:B:218:MET:HG2	5:E:73:TYR:CE2	2.55	0.41
5:E:65:MET:CA	5:E:68:THR:HG22	2.51	0.41
1:A:69:ILE:O	1:A:69:ILE:HG23	2.20	0.41
1:A:51:VAL:HG12	2:B:215:LEU:HD11	2.01	0.41
2:B:212:ILE:HG22	3:C:38:MET:HE3	2.02	0.41
3:C:38:MET:O	3:C:42:SER:HB2	2.21	0.41
2:B:228:SER:O	2:B:230:GLY:N	2.54	0.41
3:C:78:GLU:C	3:C:80:ASN:H	2.23	0.41
3:C:79:LYS:C	3:C:81:LEU:N	2.72	0.41
5:E:65:MET:O	5:E:68:THR:HG22	2.21	0.41
3:C:59:GLN:C	3:C:61:ASP:N	2.74	0.40
3:C:64:GLU:HB2	4:D:188:ILE:CD1	2.51	0.40
2:B:193:GLN:NE2	6:B:10:HOH:O	2.53	0.40
1:A:41:LEU:HD12	1:A:41:LEU:HA	1.82	0.40
2:B:220:MET:CE	3:C:41:GLU:O	2.68	0.40
4:D:157:MET:O	4:D:159:GLU:N	2.55	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	64/80 (80%)	46 (72%)	11 (17%)	7 (11%)	0	2
2	B	80/88 (91%)	62 (78%)	14 (18%)	4 (5%)	3	13
3	C	71/83 (86%)	58 (82%)	11 (16%)	2 (3%)	6	28
4	D	73/87 (84%)	52 (71%)	18 (25%)	3 (4%)	3	17
5	E	26/79 (33%)	18 (69%)	8 (31%)	0	100	100
All	All	314/417 (75%)	236 (75%)	62 (20%)	16 (5%)	2	13

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	34	PRO
1	A	35	VAL
1	A	36	GLN
1	A	84	SER
1	A	96	LYS
2	B	180	SER
2	B	242	ALA
4	D	173	ASP
2	B	256	LYS
3	C	60	LEU
3	C	81	LEU
4	D	146	ASN
1	A	69	ILE
2	B	247	ILE
4	D	209	LEU
1	A	94	LYS

5.3.2 Protein sidechains [i](#)

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	48/69 (70%)	38 (79%)	10 (21%)	1	6
2	B	57/74 (77%)	54 (95%)	3 (5%)	28	65
3	C	54/76 (71%)	46 (85%)	8 (15%)	4	15
4	D	55/75 (73%)	51 (93%)	4 (7%)	17	50
5	E	13/64 (20%)	12 (92%)	1 (8%)	16	47
All	All	227/358 (63%)	201 (88%)	26 (12%)	7	26

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	34	PRO
1	A	37	GLN
1	A	40	ARG
1	A	42	GLN
1	A	43	GLN
1	A	60	ASP
1	A	62	VAL
1	A	72	LEU
1	A	86	PHE
1	A	87	GLU
2	B	194	THR
2	B	227	GLU
2	B	245	ASP
3	C	29	LEU
3	C	35	MET
3	C	37	ASN
3	C	38	MET
3	C	42	SER
3	C	48	ARG
3	C	55	GLU
3	C	79	LYS
4	D	150	GLU
4	D	152	ASP
4	D	162	SER
4	D	185	VAL
5	E	59	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	80	GLN
2	B	239	ASN
3	C	18	GLN
3	C	20	GLN
3	C	22	ASN
3	C	37	ASN
3	C	56	GLN
3	C	59	GLN
3	C	80	ASN
4	D	155	ASN
4	D	156	ASN
4	D	176	ASN
4	D	203	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](#)

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	66/80 (82%)	0.47	8 (12%) 6 3	42, 82, 116, 119	0
2	B	82/88 (93%)	0.34	3 (3%) 45 27	44, 78, 121, 125	0
3	C	73/83 (87%)	0.21	1 (1%) 78 59	35, 74, 112, 116	0
4	D	75/87 (86%)	0.30	2 (2%) 58 37	41, 72, 117, 119	0
5	E	28/79 (35%)	0.34	2 (7%) 19 10	80, 102, 132, 132	0
All	All	324/417 (77%)	0.33	16 (4%) 33 19	35, 81, 118, 132	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	250	ALA	3.5
1	A	89	SER	3.1
1	A	98	TRP	3.1
5	E	63	GLU	2.9
2	B	243	ALA	2.9
1	A	96	LYS	2.8
1	A	93	LEU	2.7
2	B	178	GLY	2.7
4	D	137	SER	2.6
5	E	58	MET	2.5
4	D	136	PRO	2.5
1	A	97	PHE	2.5
3	C	83	GLY	2.5
1	A	69	ILE	2.4
1	A	85	GLN	2.2
1	A	90	ALA	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.