Full wwPDB X-ray Structure Validation Report ⓘ

Sep 20, 2023 – 02:29 PM EDT

PDB ID	:	5KJ8
Title	:	Structure of the Ca2+-bound synaptotagmin-1 SNARE complex (long unit cell form) - from synchrotron diffraction
Authors	:	Lyubimov, A.Y.; Uervirojnangkoorn, M.; Zhou, Q.; Zhao, M.; Sauter, N.K.; Brewster, A.S.; Weis, W.I.; Brunger, A.T.
Deposited on	:	2016-06-17
Resolution	:	4.10 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 types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references ⓘ) were used in the production of this report:

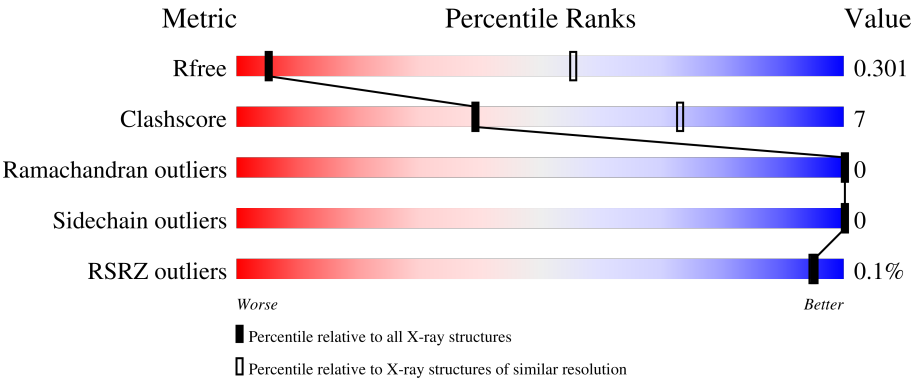
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

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 4.10 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	1193 (4.50-3.70)
Clashscore	141614	1003 (4.44-3.76)
Ramachandran outliers	138981	1005 (4.48-3.72)
Sidechain outliers	138945	1199 (4.50-3.70)
RSRZ outliers	127900	1034 (4.50-3.70)

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 of 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 <=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	63	86% 14%
1	G	63	79% 21%
2	B	66	79% 21%
2	H	66	73% 26% .
3	C	75	80% 19% .

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Mol	Chain	Length	Quality of chain
3	I	75	72% 27%
4	D	64	88% 11%
4	J	64	81% 17%
5	E	279	% 85% 15%
5	F	279	85% 13%
5	K	279	80% 20%

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 10515 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 Vesicle-associated membrane protein 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	63	Total	C	N	O	S	0	0	0
			502	305	95	101	1			
1	G	63	Total	C	N	O	S	0	0	0
			492	299	91	101	1			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	37	ALA	ASN	conflict	UNP P63025
G	37	ALA	ASN	conflict	UNP P63025

- Molecule 2 is a protein called Syntaxin-1A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	66	Total	C	N	O	S	1	0	0
			512	314	87	106	5			
2	H	65	Total	C	N	O	S	0	0	0
			510	314	84	107	5			

- Molecule 3 is a protein called Synaptosomal-associated protein 25.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	74	Total	C	N	O	S	0	0	0
			575	341	101	128	5			
3	I	74	Total	C	N	O	S	0	0	0
			549	325	98	121	5			

- Molecule 4 is a protein called Synaptosomal-associated protein 25.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	63	Total	C	N	O	S	0	0	0
			488	286	92	105	5			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	J	63	Total	C	N	O	S	0	0	0
			496	292	94	105	5			

- Molecule 5 is a protein called Synaptotagmin-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	E	279	Total	C	N	O	S	0	0	1
			2161	1389	363	402	7			
5	F	275	Total	C	N	O	S	2	0	1
			2064	1314	347	396	7			
5	K	279	Total	C	N	O	S	1	0	1
			2152	1384	359	402	7			

- Molecule 6 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	C	2	Total	Ca	0	0
			2	2		
6	E	4	Total	Ca	0	0
			4	4		
6	F	4	Total	Ca	0	0
			4	4		
6	K	4	Total	Ca	0	0
			4	4		

3 Residue-property plots (i)

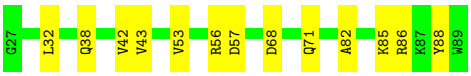
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Vesicle-associated membrane protein 3

Chain A:

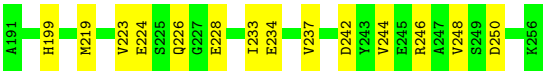
• Molecule 1: Vesicle-associated membrane protein 3

Chain G:



• Molecule 2: Syntaxin-1A

Chain B:



• Molecule 2: Syntaxin-1A

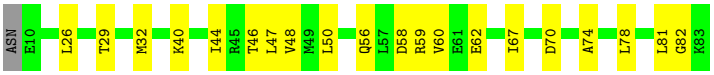
Chain H:

• Molecule 3: Synaptosomal-associated protein 25

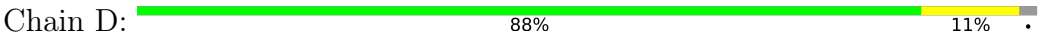
Chain C:

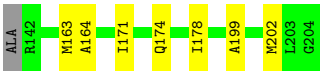
• Molecule 3: Synaptosomal-associated protein 25

Chain I:

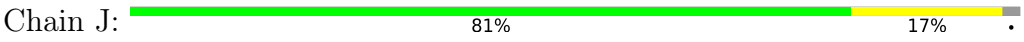


• Molecule 4: Synaptosomal-associated protein 25

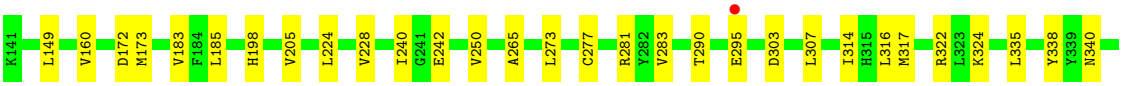


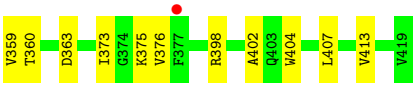


• Molecule 4: Synaptosomal-associated protein 25

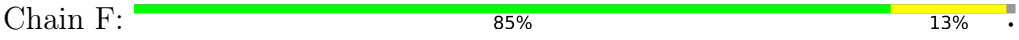


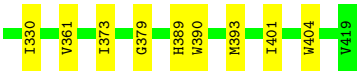
• Molecule 5: Synaptotagmin-1



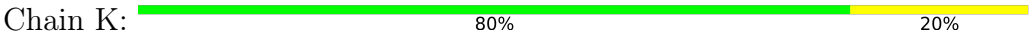


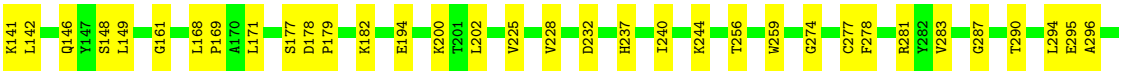
• Molecule 5: Synaptotagmin-1

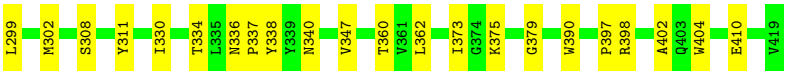




• Molecule 5: Synaptotagmin-1







4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	68.79Å 169.71Å 286.79Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.64 – 4.10 83.29 – 4.10	Depositor EDS
% Data completeness (in resolution range)	98.0 (49.64-4.10) 89.9 (83.29-4.10)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.84 (at 4.15Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.283 , 0.297 0.286 , 0.301	Depositor DCC
R_{free} test set	1338 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å^2)	134.5	Xtriage
Anisotropy	0.642	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.30 , 82.3	EDS
L-test for twinning[2]	$< \|L\| > = 0.45$, $< L^2 > = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.90	EDS
Total number of atoms	10515	wwPDB-VP
Average B, all atoms (Å^2)	155.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 17.57% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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

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	0.22	0/505	0.34	0/675
1	G	0.22	0/495	0.33	0/667
2	B	0.22	0/517	0.37	0/695
2	H	0.23	0/515	0.36	0/693
3	C	0.21	0/575	0.35	0/769
3	I	0.23	0/549	0.36	0/730
4	D	0.22	0/488	0.34	0/649
4	J	0.22	0/496	0.35	0/661
5	E	0.24	0/2209	0.41	1/2994 (0.0%)
5	F	0.22	0/2109	0.39	0/2860
5	K	0.22	0/2201	0.39	0/2983
All	All	0.23	0/10659	0.38	1/14376 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
5	E	303	ASP	CB-CA-C	-5.45	99.51	110.40

There are no chirality outliers.

There are no planarity outliers.

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	502	0	498	10	0
1	G	492	0	472	12	0
2	B	512	0	477	21	0
2	H	510	0	479	20	0
3	C	575	0	532	16	0
3	I	549	0	506	22	0
4	D	488	0	462	9	0
4	J	496	0	480	13	0
5	E	2161	0	2107	25	0
5	F	2064	0	1924	26	0
5	K	2152	0	2083	39	0
6	C	2	0	0	0	0
6	E	4	0	0	0	0
6	F	4	0	0	0	0
6	K	4	0	0	0	0
All	All	10515	0	10020	152	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 7.

All (152) 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)
2:H:248:VAL:HG21	3:I:70:ASP:HB3	1.63	0.78
3:C:40:LYS:NZ	5:E:295:GLU:OE2	2.16	0.77
5:F:281:ARG:HB3	5:F:290:THR:HB	1.74	0.70
2:B:234:GLU:OE2	3:C:59:ARG:NH1	2.22	0.69
5:K:281:ARG:HB3	5:K:290:THR:HB	1.75	0.69
2:B:248:VAL:HG21	3:C:70:ASP:HB3	1.74	0.69
2:B:224:GLU:HG2	5:F:286:ALA:HB2	1.77	0.67
2:H:244:VAL:HG21	3:I:67:ILE:HG23	1.78	0.66
5:F:312:VAL:HG22	5:F:361:VAL:HG22	1.78	0.65
5:K:148:SER:HB3	5:K:161:GLY:HA3	1.78	0.65
2:H:223:VAL:HG21	3:I:46:THR:HG23	1.78	0.65
5:E:242:GLU:HG2	5:E:265:ALA:HB2	1.80	0.64
1:A:57:ASP:HB2	2:B:226:GLN:HE21	1.63	0.62
2:B:223:VAL:O	3:C:53:GLN:NE2	2.29	0.62
1:A:64:ASP:HA	2:B:233:ILE:HG12	1.81	0.62
4:J:163:MET:HG2	5:K:338:TYR:CZ	2.35	0.62
2:H:237:VAL:HG21	3:I:60:VAL:HG13	1.82	0.61
3:C:48:VAL:HG21	5:E:402:ALA:HB2	1.84	0.60
5:E:172:ASP:OD1	5:E:173:MET:N	2.33	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:E:198:HIS:NE2	5:E:205:VAL:O	2.34	0.59
5:E:360:THR:HG23	5:E:375:LYS:HB3	1.84	0.59
4:D:163:MET:HG2	5:E:338:TYR:CZ	2.37	0.59
5:E:277:CYS:HB2	5:E:404:TRP:CD2	2.38	0.58
5:E:281:ARG:HB3	5:E:290:THR:HB	1.84	0.58
5:K:277:CYS:HB2	5:K:404:TRP:CD2	2.38	0.58
5:E:316:LEU:HB3	5:E:324:LYS:HB2	1.84	0.58
5:F:277:CYS:HB2	5:F:404:TRP:CD2	2.39	0.57
5:F:228:VAL:HG12	5:F:240:ILE:HD12	1.86	0.57
4:J:163:MET:HG2	5:K:338:TYR:CE1	2.39	0.57
2:B:237:VAL:HG21	3:C:60:VAL:HG13	1.87	0.57
4:D:163:MET:HG2	5:E:338:TYR:CE1	2.39	0.57
1:A:49:ASN:ND2	4:D:164:ALA:O	2.38	0.56
5:E:149:LEU:HD12	5:E:250:VAL:HG11	1.86	0.56
5:F:151:TYR:OH	5:F:156:ASN:OD1	2.23	0.56
1:A:53:VAL:HG23	4:D:171:ILE:HD13	1.87	0.56
2:B:242:ASP:OD2	1:G:85:LYS:NZ	2.23	0.56
1:A:56:ARG:NH1	2:B:226:GLN:OE1	2.32	0.56
5:F:146:GLN:NE2	5:F:257:GLU:OE2	2.39	0.56
2:H:248:VAL:HG22	3:I:74:ALA:HB2	1.87	0.55
5:K:274:GLY:HA3	5:K:299:LEU:HD23	1.88	0.55
5:E:283:VAL:HG22	5:E:398:ARG:HA	1.89	0.55
2:H:234:GLU:OE1	3:I:56:GLN:NE2	2.36	0.55
3:I:26:LEU:HD11	4:J:143:GLU:HG3	1.90	0.54
4:D:199:ALA:HA	4:D:202:MET:HE2	1.89	0.53
5:F:389:HIS:NE2	5:F:401:ILE:O	2.34	0.53
1:G:56:ARG:NH1	2:H:226:GLN:OE1	2.35	0.53
2:B:224:GLU:CG	5:F:286:ALA:HB2	2.39	0.53
5:F:148:SER:HB3	5:F:161:GLY:HA3	1.91	0.53
2:B:223:VAL:HG21	3:C:46:THR:HG23	1.90	0.52
5:E:363:ASP:HB2	5:E:373:ILE:HD11	1.90	0.52
1:G:32:LEU:HD11	2:H:202:ILE:HG13	1.91	0.52
5:K:302:MET:HG3	5:K:373:ILE:HG12	1.92	0.52
2:B:234:GLU:CD	3:C:59:ARG:HH12	2.12	0.52
5:E:314:ILE:HD12	5:E:359:VAL:HG22	1.92	0.52
3:I:47:LEU:HD13	5:K:340:ASN:ND2	2.24	0.52
5:K:146:GLN:HB2	5:K:259:TRP:CE2	2.44	0.52
5:F:311:TYR:HB3	5:F:330:ILE:HG12	1.91	0.52
3:C:26:LEU:O	3:C:29:THR:OG1	2.21	0.51
3:C:78:LEU:HD22	4:D:202:MET:HE1	1.92	0.51
3:I:58:ASP:OD1	4:J:177:GLN:NE2	2.36	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:244:VAL:HG21	3:C:67:ILE:HG23	1.92	0.50
5:K:308:SER:OG	5:K:334:THR:O	2.29	0.50
2:B:234:GLU:OE2	3:C:59:ARG:NH2	2.45	0.50
5:K:397:PRO:O	5:K:398:ARG:HG2	2.11	0.50
1:G:32:LEU:HD11	2:H:202:ILE:CG1	2.42	0.50
2:H:206:GLU:HA	2:H:209:ILE:HG22	1.94	0.49
4:J:160:LEU:HD23	4:J:163:MET:HE2	1.94	0.49
5:F:277:CYS:HB2	5:F:404:TRP:CE2	2.47	0.49
3:I:40:LYS:NZ	5:K:336:ASN:HD21	2.09	0.49
5:E:317:MET:HG2	5:E:322:ARG:HA	1.94	0.49
5:F:171:LEU:HD21	5:F:240:ILE:HG12	1.95	0.49
5:E:307:LEU:HD22	5:E:335:LEU:HD11	1.95	0.49
1:G:88:TYR:CD2	3:I:81:LEU:HD11	2.48	0.49
1:A:85:LYS:HD2	2:B:250:ASP:HA	1.96	0.48
3:I:26:LEU:HD23	4:J:146:MET:HB2	1.96	0.48
5:K:171:LEU:HD22	5:K:240:ILE:HG12	1.95	0.48
5:F:299:LEU:HD13	5:F:373:ILE:HD13	1.96	0.47
5:K:311:TYR:CE1	5:K:362:LEU:HB2	2.50	0.47
2:B:246:ARG:HE	2:H:246:ARG:HD3	1.79	0.47
5:K:283:VAL:HG22	5:K:398:ARG:HA	1.97	0.47
5:E:376:VAL:HG23	5:E:407:LEU:HD23	1.97	0.46
5:F:299:LEU:HB2	5:F:373:ILE:HG21	1.98	0.46
5:F:379:GLY:HA2	5:F:390:TRP:CE2	2.51	0.46
5:F:183:VAL:HG22	5:F:226:MET:HG2	1.98	0.46
1:G:43:VAL:HA	2:H:212:LEU:HD13	1.97	0.46
1:G:68:ASP:O	1:G:71:GLN:HG3	2.16	0.46
5:F:174:GLY:HA2	5:F:175:GLY:HA2	1.56	0.45
3:I:59:ARG:NH1	3:I:62:GLU:OE1	2.50	0.45
2:B:228:GLU:OE1	5:F:288:LYS:NZ	2.33	0.45
3:C:47:LEU:HD13	5:E:340:ASN:ND2	2.31	0.45
1:G:53:VAL:O	2:H:226:GLN:NE2	2.49	0.45
3:I:40:LYS:HE2	5:K:295:GLU:OE1	2.15	0.45
1:G:57:ASP:HB2	2:H:226:GLN:HE21	1.82	0.45
3:I:48:VAL:HG21	5:K:402:ALA:HB2	1.99	0.45
5:K:311:TYR:HB3	5:K:330:ILE:HG12	1.97	0.45
2:H:200:SER:O	2:H:203:ILE:HG13	2.17	0.45
5:F:228:VAL:O	5:F:240:ILE:N	2.48	0.44
5:K:141:LYS:HG3	5:K:142:LEU:HD12	1.99	0.44
4:J:159:ASN:ND2	5:K:336:ASN:O	2.36	0.44
2:B:199:HIS:CE1	3:C:24:GLU:HB3	2.52	0.44
5:K:375:LYS:HD3	5:K:410:GLU:HA	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:53:VAL:HG21	2:B:219:MET:SD	2.58	0.44
5:E:375:LYS:O	5:E:413:VAL:HG21	2.18	0.44
5:F:280:LEU:HB3	5:F:393:MET:SD	2.57	0.44
5:K:287:GLY:O	5:K:347:VAL:N	2.42	0.44
2:B:246:ARG:NE	2:H:246:ARG:HD3	2.33	0.43
5:E:228:VAL:HG12	5:E:240:ILE:HD12	2.00	0.43
1:G:82:ALA:HB1	1:G:86:ARG:NH1	2.33	0.43
5:K:379:GLY:HA2	5:K:390:TRP:CD2	2.53	0.43
5:E:277:CYS:HB2	5:E:404:TRP:CE2	2.53	0.43
3:I:26:LEU:CD2	4:J:146:MET:HB2	2.48	0.43
3:I:29:THR:HA	3:I:32:MET:HE2	2.00	0.43
5:K:177:SER:O	5:K:179:PRO:HD3	2.18	0.43
5:K:178:ASP:OD1	5:K:200:LYS:N	2.37	0.43
1:G:42:VAL:HG12	2:H:212:LEU:HD11	1.98	0.43
3:C:81:LEU:HD23	4:D:202:MET:SD	2.59	0.43
3:I:40:LYS:HZ1	5:K:336:ASN:CG	2.21	0.43
5:F:379:GLY:HA2	5:F:390:TRP:CD2	2.54	0.43
5:K:182:LYS:HG2	5:K:194:GLU:HB3	2.01	0.43
1:A:53:VAL:O	2:B:226:GLN:NE2	2.51	0.43
4:J:174:GLN:O	4:J:178:ILE:HG13	2.19	0.43
5:F:145:LEU:HD23	5:F:243:PHE:HE1	1.84	0.42
2:H:237:VAL:O	2:H:241:VAL:HG23	2.20	0.42
5:K:228:VAL:HG12	5:K:240:ILE:HD12	2.01	0.42
1:G:38:GLN:O	1:G:42:VAL:HG23	2.18	0.42
4:D:174:GLN:O	4:D:178:ILE:HG13	2.20	0.42
3:C:78:LEU:HD22	4:D:202:MET:CE	2.49	0.42
2:H:232:ARG:HD2	2:H:232:ARG:HA	1.77	0.42
5:K:149:LEU:HB3	5:K:256:THR:HB	2.01	0.42
5:K:232:ASP:H	5:K:237:HIS:CD2	2.38	0.42
5:F:229:TYR:HA	5:F:239:ILE:HA	2.02	0.42
3:I:82:GLY:HA2	4:J:202:MET:SD	2.59	0.42
1:A:32:LEU:HD23	1:A:32:LEU:HA	1.90	0.41
5:E:273:LEU:HD21	5:E:373:ILE:O	2.20	0.41
3:I:44:ILE:CG2	5:K:294:LEU:HD13	2.50	0.41
5:E:185:LEU:HD23	5:E:224:LEU:HA	2.02	0.41
1:A:38:GLN:O	1:A:42:VAL:HG23	2.19	0.41
5:K:277:CYS:HB2	5:K:404:TRP:CE2	2.54	0.41
5:K:225:VAL:HG13	5:K:244:LYS:HG2	2.01	0.41
5:K:177:SER:H	5:K:202:LEU:CD2	2.34	0.41
5:K:278:PHE:HA	5:K:294:LEU:HG	2.03	0.41
5:F:239:ILE:HD11	5:F:242:GLU:HB3	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:J:162:HIS:HB3	5:K:338:TYR:CD2	2.56	0.41
4:J:198:ARG:O	4:J:202:MET:HG2	2.21	0.41
5:K:168:LEU:HA	5:K:169:PRO:HD3	1.83	0.41
5:K:360:THR:HG23	5:K:375:LYS:HB3	2.03	0.41
3:I:50:LEU:HD23	3:I:50:LEU:HA	1.92	0.41
5:E:160:VAL:HG11	5:E:183:VAL:HG11	2.03	0.40
5:F:151:TYR:OH	5:F:214:VAL:O	2.39	0.40
2:H:229:MET:O	2:H:233:ILE:HG13	2.21	0.40
3:I:78:LEU:HD11	4:J:202:MET:HG3	2.04	0.40
5:K:296:ALA:HB3	5:K:337:PRO:HG2	2.04	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	61/63 (97%)	61 (100%)	0	0	100	100
1	G	61/63 (97%)	61 (100%)	0	0	100	100
2	B	64/66 (97%)	64 (100%)	0	0	100	100
2	H	63/66 (96%)	63 (100%)	0	0	100	100
3	C	72/75 (96%)	72 (100%)	0	0	100	100
3	I	72/75 (96%)	72 (100%)	0	0	100	100
4	D	61/64 (95%)	61 (100%)	0	0	100	100
4	J	61/64 (95%)	61 (100%)	0	0	100	100
5	E	277/279 (99%)	263 (95%)	14 (5%)	0	100	100
5	F	269/279 (96%)	257 (96%)	12 (4%)	0	100	100
5	K	277/279 (99%)	263 (95%)	14 (5%)	0	100	100
All	All	1338/1373 (98%)	1298 (97%)	40 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	53/54 (98%)	53 (100%)	0	100	100
1	G	51/54 (94%)	51 (100%)	0	100	100
2	B	55/60 (92%)	55 (100%)	0	100	100
2	H	56/60 (93%)	56 (100%)	0	100	100
3	C	61/67 (91%)	61 (100%)	0	100	100
3	I	57/67 (85%)	57 (100%)	0	100	100
4	D	52/55 (94%)	52 (100%)	0	100	100
4	J	54/55 (98%)	54 (100%)	0	100	100
5	E	229/249 (92%)	229 (100%)	0	100	100
5	F	212/249 (85%)	212 (100%)	0	100	100
5	K	227/249 (91%)	227 (100%)	0	100	100
All	All	1107/1219 (91%)	1107 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

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

Mol	Chain	Res	Type
5	E	356	GLN
2	H	213	HIS

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 monosaccharides in this entry.

5.6 Ligand geometry

Of 14 ligands modelled in this entry, 14 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^{th} 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	63/63 (100%)	0.08	0 100 100	118, 135, 162, 170	0
1	G	63/63 (100%)	-0.16	0 100 100	144, 168, 183, 193	0
2	B	66/66 (100%)	0.03	0 100 100	121, 136, 161, 171	1 (1%)
2	H	65/66 (98%)	-0.15	0 100 100	144, 167, 179, 181	0
3	C	74/75 (98%)	0.06	0 100 100	116, 140, 180, 191	0
3	I	74/75 (98%)	-0.15	0 100 100	150, 161, 182, 194	0
4	D	63/64 (98%)	-0.04	0 100 100	113, 138, 162, 167	0
4	J	63/64 (98%)	-0.27	0 100 100	154, 164, 184, 192	0
5	E	279/279 (100%)	-0.06	2 (0%) 87 82	119, 154, 177, 182	0
5	F	275/279 (98%)	-0.18	0 100 100	126, 156, 215, 228	1 (0%)
5	K	279/279 (100%)	-0.15	0 100 100	124, 156, 177, 193	1 (0%)
All	All	1364/1373 (99%)	-0.11	2 (0%) 95 95	113, 156, 185, 228	3 (0%)

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
5	E	377	PHE	2.6
5	E	295	GLU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains

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

6.3 Carbohydrates

There are no monosaccharides 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. The B-factors column lists the minimum, median, 95^{th} 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	B-factors(Å^2)	Q<0.9
6	CA	C	102	1/1	0.48	0.27	142,142,142,142	0
6	CA	K	501	1/1	0.56	0.10	192,192,192,192	0
6	CA	F	504	1/1	0.77	0.22	133,133,133,133	0
6	CA	E	501	1/1	0.83	0.18	157,157,157,157	0
6	CA	F	502	1/1	0.86	0.18	137,137,137,137	0
6	CA	C	101	1/1	0.92	0.34	110,110,110,110	0
6	CA	E	503	1/1	0.92	0.16	153,153,153,153	0
6	CA	F	501	1/1	0.92	0.33	81,81,81,81	0
6	CA	F	503	1/1	0.93	0.11	218,218,218,218	0
6	CA	E	504	1/1	0.94	0.20	146,146,146,146	0
6	CA	E	502	1/1	0.96	0.12	160,160,160,160	0
6	CA	K	504	1/1	0.96	0.20	136,136,136,136	0
6	CA	K	502	1/1	0.98	0.27	158,158,158,158	0
6	CA	K	503	1/1	0.99	0.18	134,134,134,134	0

6.5 Other polymers (i)

There are no such residues in this entry.