



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 01:22 AM UTC

PDB ID : 6BIT / pdb_00006bit
Title : SIRPalha antibody complex
Authors : Ring, N.G.; Herndler-Brandstetter, D.; Weiskopf, K.; Shan, L.; Volkmer, J.P.; George, B.M.; Lietzenmayer, M.; McKenna, K.M.; Naik, T.J.; McCarty, A.; Zheng, Y.; Ring, A.M.; Flavell, R.A.; Weissman, I.L.
Deposited on : 2017-11-03
Resolution : 2.19 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 : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtrriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

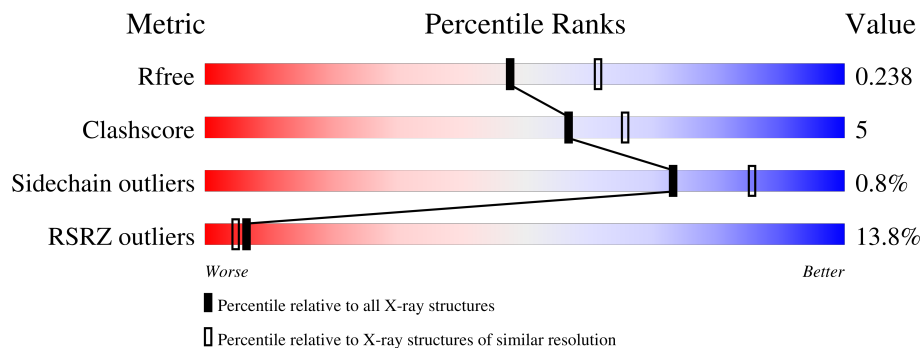
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.19 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}	180053	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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 $\leq 5\%$. 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	I	213	
1	J	213	
2	K	212	
2	L	212	
3	G	117	
3	H	117	

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 8398 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 KWAR23 Fab heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	I	213	Total 1600	C 1014	N 265	O 316	S 5	0	0	0
1	J	207	Total 1559	C 992	N 253	O 309	S 5	0	0	0

- Molecule 2 is a protein called KWAR23 Fab light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	K	212	Total 1618	C 1013	N 267	O 331	S 7	0	0	0
2	L	211	Total 1608	C 1008	N 263	O 330	S 7	0	0	0

- Molecule 3 is a protein called Tyrosine-protein phosphatase non-receptor type substrate 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	H	115	Total 884	C 556	N 157	O 168	S 3	0	1	0
3	G	115	Total 884	C 555	N 156	O 170	S 3	0	1	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	I	9	Total 9 9	0	0
4	K	17	Total 17 17	0	0
4	H	55	Total 55 55	0	0
4	J	60	Total 60 60	0	0

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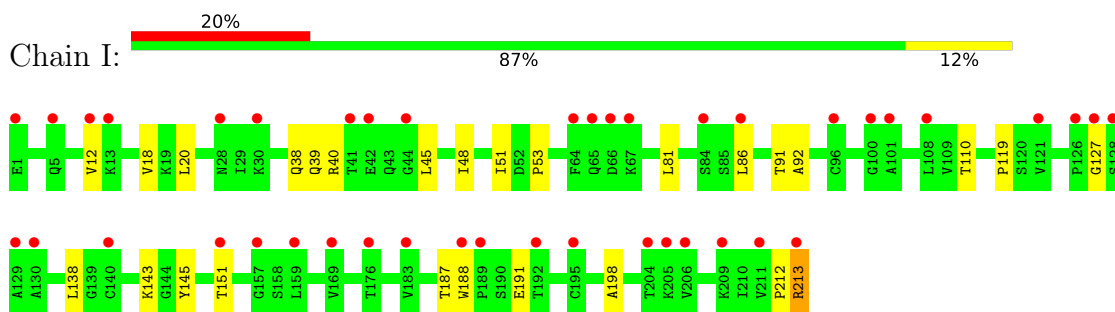
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	L	36	Total	O	0	0
			36	36		
4	G	68	Total	O	0	0
			68	68		

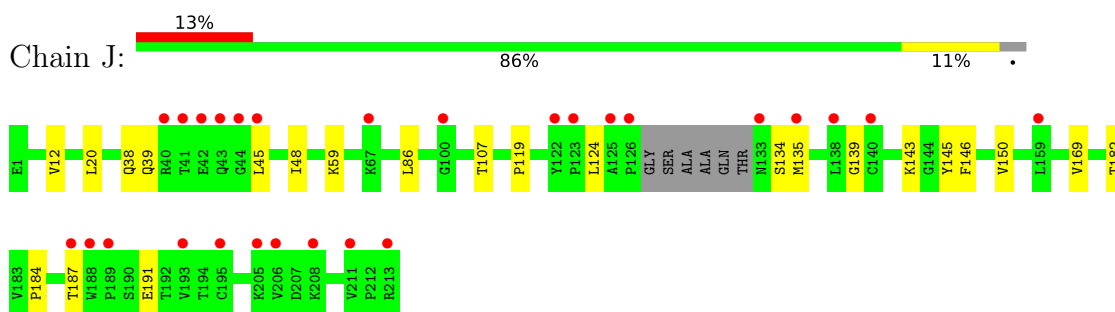
3 Residue-property plots [i](#)

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.

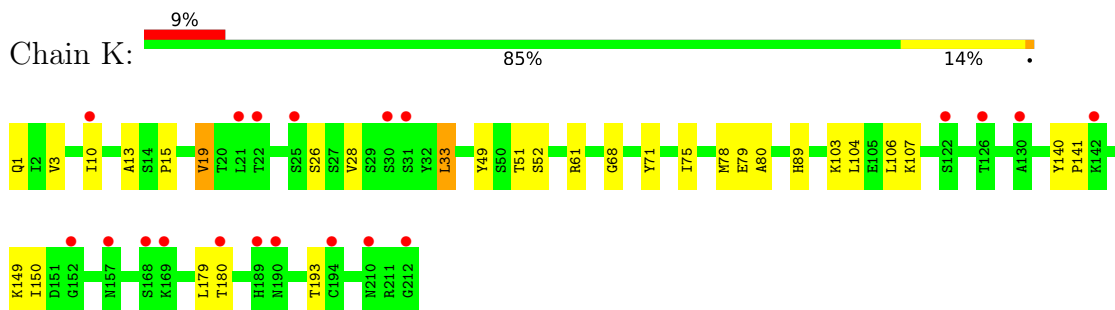
- Molecule 1: KWAR23 Fab heavy chain



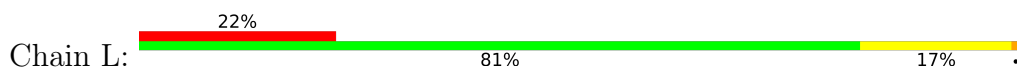
- Molecule 1: KWAR23 Fab heavy chain

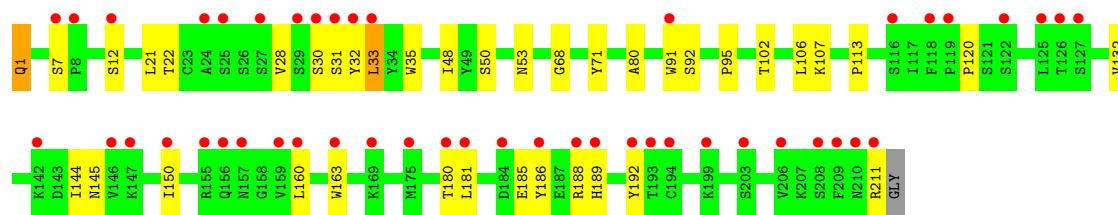


- Molecule 2: KWAR23 Fab light chain

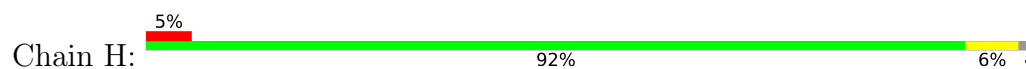


- Molecule 2: KWAR23 Fab light chain

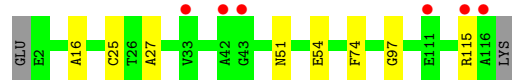
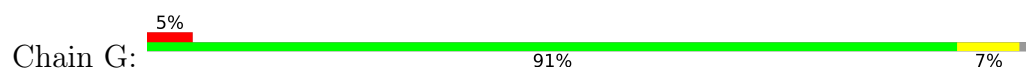




- Molecule 3: Tyrosine-protein phosphatase non-receptor type substrate 1



- Molecule 3: Tyrosine-protein phosphatase non-receptor type substrate 1



4 Data and refinement statistics i

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	164.97Å 164.97Å 96.72Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	47.62 – 2.19 47.62 – 2.19	Depositor EDS
% Data completeness (in resolution range)	98.8 (47.62-2.19) 95.8 (47.62-2.19)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.82 (at 2.20Å)	Xtrriage
Refinement program	PHENIX 1.11.1_2575	Depositor
R, R_{free}	0.210 , 0.240 (Not available) , 0.238	Depositor DCC
R_{free} test set	3856 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	41.8	Xtrriage
Anisotropy	0.179	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 43.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.021 for -h,-k,l	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8398	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.41% 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](#)

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

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	I	0.10	0/1643	0.31	0/2254
1	J	0.43	2/1601 (0.1%)	0.33	0/2196
2	K	0.44	4/1652 (0.2%)	0.35	0/2248
2	L	0.10	0/1642	0.35	0/2236
3	G	0.10	0/904	0.30	0/1225
3	H	0.40	1/904 (0.1%)	0.58	4/1224 (0.3%)
All	All	0.31	7/8346 (0.1%)	0.37	4/11383 (0.0%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	H	29	SER	C-N	-11.10	1.17	1.33
2	K	51	THR	C-O	-6.82	1.15	1.24
2	K	52	SER	CA-C	-6.72	1.44	1.52
1	J	146	PHE	C-O	-6.38	1.17	1.24
2	K	52	SER	C-O	-6.20	1.15	1.24
1	J	150	VAL	C-O	-5.90	1.17	1.24
2	K	49	TYR	C-O	-5.46	1.16	1.23

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	H	29	SER	O-C-N	-13.13	110.46	123.46
3	H	28	THR	O-C-N	-6.12	114.49	122.39
3	H	29	SER	CA-C-N	5.77	132.06	122.54
3	H	29	SER	C-N-CA	5.77	132.06	122.54

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	I	1600	0	1552	16	0
1	J	1559	0	1505	12	0
2	K	1618	0	1542	17	0
2	L	1608	0	1528	25	0
3	G	884	0	881	6	0
3	H	884	0	887	3	0
4	G	68	0	0	0	0
4	H	55	0	0	1	0
4	I	9	0	0	0	0
4	J	60	0	0	0	0
4	K	17	0	0	0	0
4	L	36	0	0	0	0
All	All	8398	0	7895	74	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 5.

All (74) 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)
3:H:111:GLU:OE1	4:H:201:HOH:O	2.15	0.65
2:K:149:LYS:HB2	2:K:193:THR:HB	1.79	0.63
2:L:28:VAL:HG13	2:L:68:GLY:HA2	1.81	0.62
1:J:119:PRO:HB3	1:J:145:TYR:HB3	1.83	0.61
1:J:143:LYS:HE3	2:L:180:THR:HG21	1.84	0.59
2:L:189:HIS:O	2:L:211:ARG:NH1	2.37	0.58
2:L:21:LEU:HD22	2:L:102:THR:HG21	1.86	0.57
2:L:113:PRO:HG3	2:L:144:ILE:HD11	1.86	0.56
1:J:135:MET:HE3	1:J:182:THR:HG22	1.86	0.56
3:H:32:PRO:O	3:H:93:LYS:NZ	2.38	0.56
2:L:32:TYR:HD2	2:L:91:TRP:CD2	2.23	0.56
2:L:7:SER:HB3	2:L:22:THR:HB	1.87	0.55
2:K:78:MET:HE1	2:K:104:LEU:HD21	1.87	0.55
3:G:16:ALA:HA	3:G:115:ARG:HG3	1.89	0.55
1:J:59:LYS:NZ	3:G:97:GLY:O	2.40	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:40:ARG:HG2	1:I:92:ALA:HB2	1.89	0.54
2:K:15:PRO:HG3	2:K:106:LEU:HD11	1.90	0.53
1:I:12:VAL:HG11	1:I:86:LEU:HD12	1.90	0.52
2:K:28:VAL:HG13	2:K:68:GLY:HA2	1.91	0.52
1:J:39:GLN:HB2	1:J:45:LEU:HD23	1.93	0.51
2:K:80:ALA:HA	2:K:106:LEU:HD13	1.91	0.51
1:J:187:THR:O	1:J:191:GLU:N	2.38	0.51
2:K:13:ALA:HB3	2:K:78:MET:HE3	1.92	0.51
1:I:38:GLN:HB3	1:I:48:ILE:HD11	1.92	0.51
1:J:38:GLN:HB2	1:J:48:ILE:HD11	1.92	0.51
1:I:39:GLN:HB2	1:I:45:LEU:HD23	1.94	0.50
2:L:188:ARG:HG3	2:L:189:HIS:CD2	2.47	0.49
2:L:80:ALA:HA	2:L:106:LEU:HG	1.94	0.49
1:I:119:PRO:HB3	1:I:145:TYR:HB3	1.95	0.49
2:L:50:SER:OG	2:L:53:ASN:ND2	2.43	0.48
2:L:1:PCA:HB3	2:L:95:PRO:HD2	1.94	0.48
1:I:91:THR:HG23	1:I:110:THR:HA	1.96	0.47
1:I:127:GLY:HA2	1:I:213:ARG:N	2.30	0.47
2:L:32:TYR:HB3	2:L:91:TRP:H	1.81	0.46
1:I:143:LYS:HZ1	2:K:180:THR:HG21	1.81	0.46
1:J:12:VAL:HG11	1:J:86:LEU:HD12	1.97	0.45
1:I:213:ARG:HH21	2:L:145:ASN:HB2	1.80	0.45
2:K:33:LEU:HA	2:K:89:HIS:O	2.16	0.45
2:L:33:LEU:HG	2:L:71:TYR:CG	2.50	0.45
2:L:186:TYR:O	2:L:192:TYR:OH	2.34	0.45
2:L:32:TYR:HD2	2:L:91:TRP:CE2	2.35	0.44
3:G:25:CYS:SG	3:G:74:PHE:HB2	2.57	0.44
2:K:150:ILE:HD11	2:K:179:LEU:HD21	2.00	0.44
2:L:144:ILE:HG22	2:L:163:TRP:CH2	2.53	0.44
1:I:212:PRO:O	1:I:213:ARG:HB2	2.17	0.43
3:G:16:ALA:HB2	3:G:115:ARG:HE	1.82	0.43
2:K:3:VAL:HB	2:K:26:SER:HB3	2.00	0.43
2:L:181:LEU:HB3	2:L:185:GLU:HG3	2.01	0.43
1:J:20:LEU:HD22	1:J:107:THR:HG21	2.00	0.43
1:J:184:PRO:O	1:J:187:THR:OG1	2.34	0.43
3:H:36:ILE:HD13	3:H:93:LYS:HA	1.99	0.43
1:I:18:VAL:HG12	1:I:86:LEU:HD11	2.00	0.43
2:K:61:ARG:HG3	2:K:75:ILE:HG23	2.01	0.43
3:G:27:ALA:HB2	3:G:74:PHE:CE2	2.54	0.43
2:L:30:SER:HB3	2:L:32:TYR:HD1	1.83	0.43
2:L:150:ILE:HD11	2:L:189:HIS:CD2	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:32:TYR:CD1	2:L:92:SER:HB3	2.53	0.42
2:K:19:VAL:HG11	2:K:78:MET:HE2	2.01	0.42
2:K:33:LEU:HD22	2:K:71:TYR:CD2	2.54	0.42
1:J:124:LEU:HB2	1:J:139:GLY:HA3	2.01	0.42
2:L:12:SER:OG	2:L:107:LYS:HB2	2.18	0.42
2:K:140:TYR:CG	2:K:141:PRO:HA	2.54	0.42
1:J:169:VAL:HG11	2:L:160:LEU:HD22	2.01	0.42
1:I:51:ILE:O	1:I:53:PRO:HD3	2.20	0.42
1:I:187:THR:O	1:I:191:GLU:N	2.51	0.41
2:L:35:TRP:HB2	2:L:48:ILE:HB	2.03	0.41
1:I:138:LEU:HD11	1:I:188:TRP:CD1	2.56	0.41
1:I:20:LEU:HD12	1:I:81:LEU:HD23	2.02	0.41
2:K:107:LYS:HA	2:K:140:TYR:OH	2.21	0.41
1:I:151:THR:HB	1:I:198:ALA:HB3	2.02	0.40
3:G:51:ASN:HB3	3:G:54:GLU:O	2.20	0.40
2:K:10:ILE:HD12	2:K:103:LYS:O	2.22	0.40
2:K:61:ARG:CZ	2:K:79:GLU:HG3	2.52	0.40
2:L:120:PRO:HD3	2:L:132:VAL:HG22	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

There are no protein backbone outliers to report in this entry.

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	I	177/181 (98%)	176 (99%)	1 (1%)	78 89
1	J	173/181 (96%)	172 (99%)	1 (1%)	78 89
2	K	183/184 (100%)	181 (99%)	2 (1%)	65 79

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	L	182/184 (99%)	180 (99%)	2 (1%)	65	79
3	G	96/97 (99%)	96 (100%)	0	100	100
3	H	96/97 (99%)	95 (99%)	1 (1%)	68	81
All	All	907/924 (98%)	900 (99%)	7 (1%)	73	85

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	I	213	ARG
2	K	19	VAL
2	K	33	LEU
3	H	13	VAL
1	J	134	SER
2	L	31	SER
2	L	33	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	I	3	GLN
1	I	39	GLN
1	I	43	GLN
2	K	37	GLN
2	K	38	GLN
2	K	53	ASN
2	K	137	ASN
2	L	157	ASN
2	L	161	ASN
2	L	190	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

2 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul

statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	PCA	K	1	2	7,8,9	1.92	1 (14%)	9,10,12	2.16	5 (55%)
2	PCA	L	1	2	7,8,9	1.93	1 (14%)	9,10,12	2.17	5 (55%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	PCA	K	1	2	-	0/0/11/13	0/1/1/1
2	PCA	L	1	2	-	0/0/11/13	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	L	1	PCA	CD-N	5.00	1.46	1.34
2	K	1	PCA	CD-N	4.96	1.46	1.34

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	K	1	PCA	OE-CD-CG	-3.06	121.25	126.72
2	L	1	PCA	OE-CD-CG	-3.02	121.33	126.72
2	L	1	PCA	CA-N-CD	-2.87	103.74	113.58
2	K	1	PCA	CA-N-CD	-2.78	104.05	113.58
2	L	1	PCA	CB-CA-N	2.76	110.82	103.24
2	K	1	PCA	CB-CA-N	2.71	110.70	103.24
2	K	1	PCA	CG-CD-N	2.64	114.85	108.39
2	L	1	PCA	CG-CD-N	2.59	114.72	108.39
2	K	1	PCA	CB-CA-C	-2.50	109.23	112.66
2	L	1	PCA	CB-CA-C	-2.45	109.30	112.66

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	L	1	PCA	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides 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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
3	H	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	H	29:SER	C	30:LEU	N	1.17

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	I	213/213 (100%)	1.36	42 (19%) 3 2	46, 70, 87, 98	0
1	J	207/213 (97%)	0.82	27 (13%) 7 5	28, 48, 93, 109	0
2	K	211/212 (99%)	0.91	20 (9%) 14 11	43, 60, 89, 95	0
2	L	210/212 (99%)	1.17	47 (22%) 2 1	32, 55, 97, 110	0
3	G	115/117 (98%)	0.19	6 (5%) 33 29	23, 38, 57, 76	1 (0%)
3	H	115/117 (98%)	0.36	6 (5%) 33 29	29, 42, 57, 81	1 (0%)
All	All	1071/1084 (98%)	0.90	148 (13%) 6 5	23, 56, 90, 110	2 (0%)

All (148) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	L	31	SER	9.6
2	L	32	TYR	8.3
2	L	30	SER	7.0
1	J	133	ASN	6.7
2	K	212	GLY	6.5
1	I	129	ALA	6.4
1	J	100	GLY	6.2
2	K	31	SER	5.9
1	I	189	PRO	5.8
2	L	91	TRP	5.7
1	J	43	GLN	5.5
1	J	42	GLU	5.4
2	L	7	SER	5.3
2	L	33	LEU	5.2
1	J	44	GLY	5.1
2	L	210	ASN	4.8
1	I	100	GLY	4.8
2	L	211	ARG	4.8
2	L	150	ILE	4.7

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Mol	Chain	Res	Type	RSRZ
2	L	29	SER	4.5
1	J	45	LEU	4.3
1	J	213	ARG	4.2
2	L	157	ASN	4.1
1	I	127	GLY	4.1
1	I	213	ARG	4.0
2	L	8	PRO	3.7
2	L	155	ARG	3.7
1	I	44	GLY	3.6
1	I	211	VAL	3.6
1	I	126	PRO	3.5
1	J	205	LYS	3.5
2	K	30	SER	3.4
2	L	25	SER	3.3
1	J	211	VAL	3.2
1	J	187	THR	3.2
2	K	168	SER	3.1
2	L	209	PHE	3.1
3	G	115	ARG	3.1
2	L	125	LEU	3.1
3	G	43	GLY	3.0
1	I	204	THR	3.0
3	G	116	ALA	3.0
1	J	40	ARG	3.0
1	J	126	PRO	3.0
1	I	188	TRP	3.0
1	I	96	CYS	2.8
2	L	194	CYS	2.8
2	L	193	THR	2.8
1	I	205	LYS	2.8
3	H	117	LYS	2.8
3	H	70	ASN	2.7
1	J	188	TRP	2.7
1	I	67	LYS	2.7
3	H	3	GLU	2.7
2	K	25	SER	2.7
1	I	151	THR	2.6
1	I	159	LEU	2.6
2	L	181	LEU	2.6
1	J	206	VAL	2.6
2	L	169	LYS	2.6
2	L	122	SER	2.6

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Mol	Chain	Res	Type	RSRZ
2	L	147	LYS	2.6
1	J	208	LYS	2.6
1	J	140	CYS	2.6
1	I	84	SER	2.6
1	I	30	LYS	2.5
2	L	184	ASP	2.5
1	J	123	PRO	2.5
3	G	42	ALA	2.5
2	K	122	SER	2.5
2	L	12	SER	2.5
2	L	119	PRO	2.5
2	L	203	SER	2.5
1	I	108	LEU	2.5
2	L	189	HIS	2.5
1	I	5	GLN	2.5
2	L	126	THR	2.5
1	I	121	VAL	2.5
1	I	192	THR	2.5
2	K	22	THR	2.4
2	L	175	MET	2.4
2	K	142	LYS	2.4
1	I	65	GLN	2.4
1	I	13	LYS	2.4
1	I	130	ALA	2.4
2	K	210	ASN	2.4
2	L	188	ARG	2.4
2	L	186	TYR	2.4
1	J	125	ALA	2.4
1	I	28	ASN	2.4
1	J	135	MET	2.4
1	J	189	PRO	2.4
2	K	169	LYS	2.3
1	I	12	VAL	2.3
1	I	183	VAL	2.3
3	H	113	SER	2.3
2	L	199	LYS	2.3
1	J	41	THR	2.3
2	K	157	ASN	2.3
2	K	130	ALA	2.3
1	I	86	LEU	2.3
1	J	138	LEU	2.3
1	J	67	LYS	2.3

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Mol	Chain	Res	Type	RSRZ
2	L	192	TYR	2.3
1	I	41	THR	2.2
2	K	180	THR	2.2
2	L	142	LYS	2.2
1	I	195	CYS	2.2
2	K	189	HIS	2.2
2	K	190	ASN	2.2
2	K	152	GLY	2.2
1	I	1	GLU	2.2
1	I	128	SER	2.2
1	I	176	THR	2.2
2	L	180	THR	2.2
1	J	193	VAL	2.2
2	L	146	VAL	2.2
1	J	159	LEU	2.2
1	I	206	VAL	2.1
1	I	101	ALA	2.1
2	L	24	ALA	2.1
2	L	116	SER	2.1
2	K	10	ILE	2.1
3	H	27	ALA	2.1
2	K	21	LEU	2.1
2	L	127	SER	2.1
1	J	122	TYR	2.1
1	I	157	GLY	2.1
1	I	169	VAL	2.1
3	H	13	VAL	2.1
3	G	33	VAL	2.1
1	I	209	LYS	2.1
2	L	118	PHE	2.1
2	L	206	VAL	2.1
2	L	163	TRP	2.1
1	I	42	GLU	2.0
2	L	27	SER	2.1
3	G	111	GLU	2.0
1	I	66	ASP	2.0
1	I	140	CYS	2.0
1	J	195	CYS	2.0
2	L	156	GLN	2.0
2	K	126	THR	2.0
1	I	64	PHE	2.0
2	L	159	VAL	2.0

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Mol	Chain	Res	Type	RSRZ
2	L	208	SER	2.0
2	L	160	LEU	2.0
2	K	194	CYS	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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, 95th 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
2	PCA	K	1	8/9	0.88	0.13	57,64,70,77	0
2	PCA	L	1	8/9	0.95	0.09	40,48,54,63	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides 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.