



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 11:10 PM UTC

PDB ID : 1BJ1 / pdb_00001bj1
Title : VASCULAR ENDOTHELIAL GROWTH FACTOR IN COMPLEX WITH A
NEUTRALIZING ANTIBODY
Authors : Muller, Y.A.; Christinger, H.W.; De Vos, A.M.
Deposited on : 1998-06-30
Resolution : 2.40 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)
Xtriage (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

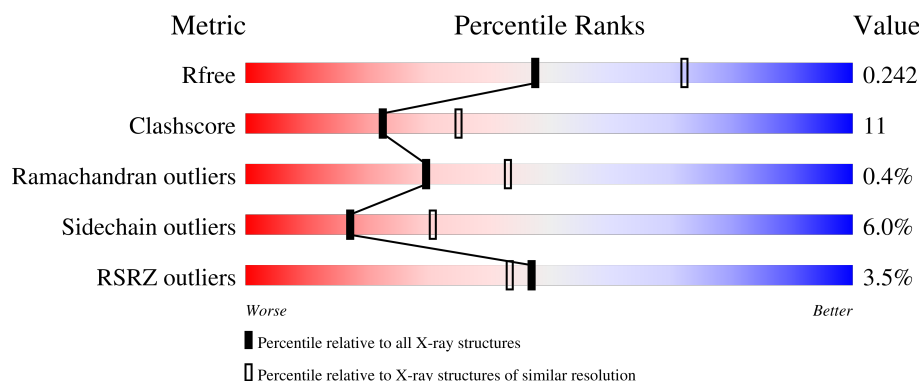
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	J	214	9% 64% 31% 5%
1	L	214	78% 18% .
2	H	231	71% 22% 6%
2	K	231	7% 66% 26% 6%
3	V	102	70% 20% 8%

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Mol	Chain	Length	Quality of chain
3	W	102	 A horizontal bar chart showing the quality of chain W. The bar is divided into three segments: green (62%), yellow (28%), and red (8%). The percentages are labeled below the bar.

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8694 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 Fab fragment, light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	213	Total	C	N	O	S	0	0	0
			1644	1031	272	336	5			
1	J	213	Total	C	N	O	S	0	0	0
			1644	1031	272	336	5			

- Molecule 2 is a protein called Fab fragment, heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	218	Total	C	N	O	S	0	0	0
			1663	1064	273	320	6			
2	K	218	Total	C	N	O	S	0	0	0
			1663	1064	273	320	6			

- Molecule 3 is a protein called Vascular endothelial growth factor A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	V	94	Total	C	N	O	S	0	0	0
			761	478	128	142	13			
3	W	94	Total	C	N	O	S	0	0	0
			761	478	128	142	13			

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	H	1	Total	O	S	0	0
			5	4	1		
4	K	1	Total	O	S	0	0
			5	4	1		

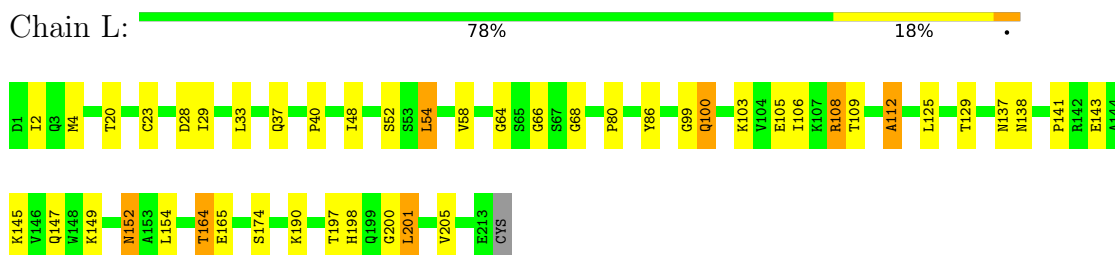
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	L	167	Total	O	0	0
			167	167		
5	H	190	Total	O	0	0
			190	190		
5	V	31	Total	O	0	0
			31	31		
5	W	34	Total	O	0	0
			34	34		
5	J	59	Total	O	0	0
			59	59		
5	K	67	Total	O	0	0
			67	67		

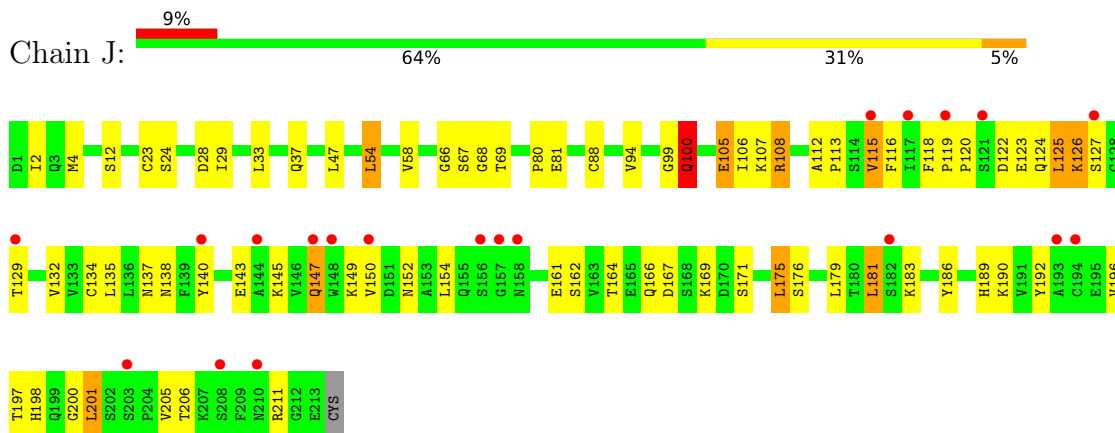
3 Residue-property plots

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.

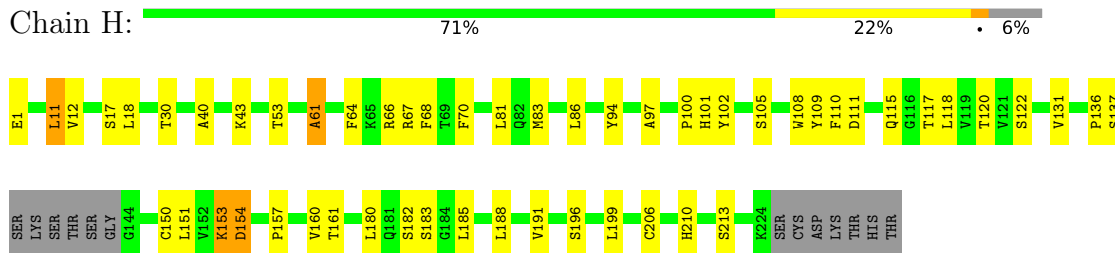
- Molecule 1: Fab fragment, light chain



- Molecule 1: Fab fragment, light chain

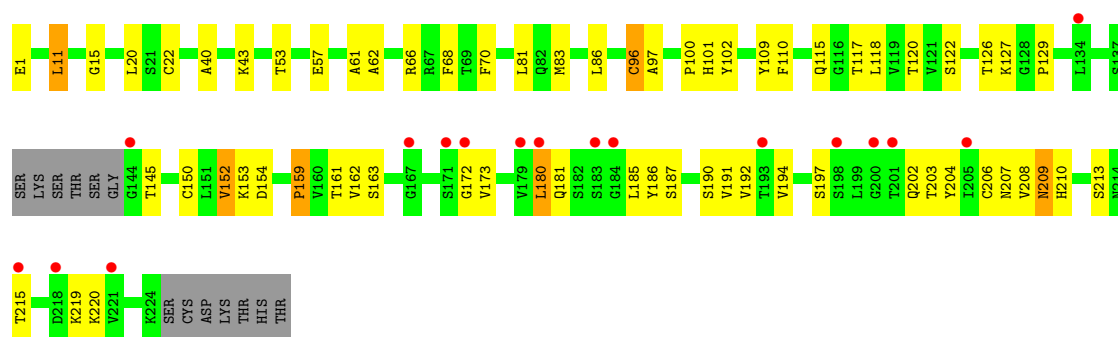


- Molecule 2: Fab fragment, heavy chain



- Molecule 2: Fab fragment, heavy chain





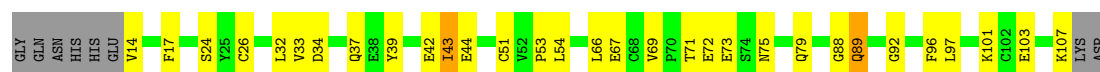
- Molecule 3: Vascular endothelial growth factor A

Chain V: 70% 20% 8%



- Molecule 3: Vascular endothelial growth factor A

Chain W: 62% 28% 8%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	89.86Å 66.98Å 140.51Å 90.00° 94.27° 90.00°	Depositor
Resolution (Å)	8.00 – 2.40 8.00 – 2.40	Depositor EDS
% Data completeness (in resolution range)	94.8 (8.00-2.40) 92.1 (8.00-2.40)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.13 (at 2.40Å)	Xtriage
Refinement program	X-PLOR 3.8	Depositor
R, R_{free}	0.196 , 0.266 0.185 , 0.242	Depositor DCC
R_{free} test set	5908 reflections (9.51%)	wwPDB-VP
Wilson B-factor (Å ²)	34.0	Xtriage
Anisotropy	0.639	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 87.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8694	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.68% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SO4

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	J	0.67	0/1682	0.98	6/2287 (0.3%)
1	L	0.81	0/1682	1.03	7/2287 (0.3%)
2	H	0.82	1/1712 (0.1%)	1.11	7/2338 (0.3%)
2	K	0.68	1/1712 (0.1%)	1.06	10/2338 (0.4%)
3	V	0.66	0/779	1.01	4/1050 (0.4%)
3	W	0.73	0/779	1.01	1/1050 (0.1%)
All	All	0.74	2/8346 (0.0%)	1.04	35/11350 (0.3%)

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
3	V	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	131	VAL	CA-CB	5.37	1.60	1.54
2	K	83	MET	SD-CE	-5.10	1.66	1.79

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	K	61	ALA	N-CA-C	-9.47	96.27	110.28
2	H	61	ALA	N-CA-C	-8.36	97.91	110.28
2	H	154	ASP	N-CA-C	8.08	121.12	111.02
2	H	111	ASP	N-CA-C	7.91	123.00	113.12
2	H	153	LYS	N-CA-C	7.58	121.99	109.95

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	K	66	ARG	N-CA-C	7.57	121.77	112.23
2	H	66	ARG	N-CA-C	7.04	119.03	111.36
2	K	62	ALA	N-CA-C	6.92	119.70	111.33
1	J	99	GLY	N-CA-C	-6.82	102.12	112.89
2	K	197	SER	N-CA-C	6.62	118.18	110.97
1	J	137	ASN	N-CA-C	6.39	119.92	109.76
3	V	93	GLU	N-CA-C	6.20	119.75	109.46
1	L	138	ASN	N-CA-C	6.19	118.76	111.02
2	H	199	LEU	N-CA-C	6.17	118.80	111.71
1	J	100	GLN	N-CA-C	6.05	120.52	113.20
1	J	143	GLU	N-CA-C	5.99	118.94	109.96
2	K	96	CYS	N-CA-C	-5.83	100.39	109.72
3	V	43	ILE	N-CA-C	5.75	118.23	111.05
3	W	43	ILE	N-CA-C	5.67	115.86	110.42
1	J	108	ARG	N-CA-C	5.57	115.76	108.07
2	K	209	ASN	N-CA-C	5.56	117.01	107.28
1	L	66	GLY	N-CA-C	5.53	118.48	111.85
2	K	57	GLU	CA-C-N	5.42	126.62	119.84
2	K	57	GLU	C-N-CA	5.42	126.62	119.84
1	L	112	ALA	CA-C-N	-5.34	115.08	120.31
1	L	112	ALA	C-N-CA	-5.34	115.08	120.31
3	V	27	HIS	N-CA-C	5.30	112.78	108.07
2	K	15	GLY	N-CA-C	-5.22	108.14	114.92
1	L	152	ASN	N-CA-C	5.21	120.81	113.56
3	V	37	GLN	N-CA-C	-5.13	105.76	112.23
2	H	101	HIS	N-CA-C	-5.07	101.70	109.76
1	L	137	ASN	N-CA-C	5.06	117.22	109.07
1	J	115	VAL	N-CA-C	5.05	116.52	109.55
1	L	99	GLY	N-CA-C	-5.03	104.75	112.85
2	K	86	LEU	N-CA-C	5.02	117.50	110.23

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	V	21	TYR	Sidechain

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	J	1644	0	1587	48	0
1	L	1644	0	1587	24	0
2	H	1663	0	1605	28	0
2	K	1663	0	1605	41	0
3	V	761	0	728	18	0
3	W	761	0	728	23	0
4	H	5	0	0	0	0
4	K	5	0	0	0	0
5	H	190	0	0	4	0
5	J	59	0	0	0	0
5	K	67	0	0	0	0
5	L	167	0	0	3	0
5	V	31	0	0	0	0
5	W	34	0	0	0	0
All	All	8694	0	7840	167	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 11.

All (167) 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:183:SER:HA	5:H:331:HOH:O	1.53	1.08
2:K:150:CYS:HG	2:K:206:CYS:HG	0.93	0.85
2:K:22:CYS:HG	2:K:96:CYS:HG	1.02	0.85
2:H:150:CYS:SG	2:H:206:CYS:SG	2.50	0.84
2:K:22:CYS:SG	2:K:96:CYS:SG	2.65	0.79
2:K:150:CYS:SG	2:K:206:CYS:SG	2.69	0.78
1:J:113:PRO:HD2	1:J:201:LEU:HD13	1.65	0.77
1:L:80:PRO:HA	1:L:106:ILE:HD13	1.64	0.77
2:K:206:CYS:SG	2:K:219:LYS:HB3	2.26	0.74
1:L:198:HIS:CD2	1:L:200:GLY:H	2.08	0.71
1:J:145:LYS:HB3	1:J:197:THR:HB	1.71	0.71
2:K:150:CYS:HG	2:K:206:CYS:CB	2.04	0.71
2:K:129:PRO:HB2	2:K:152:VAL:HG12	1.73	0.70
2:H:210:HIS:HD2	2:H:213:SER:OG	1.77	0.67
2:K:203:THR:HG23	2:K:220:LYS:HE3	1.77	0.66
3:V:89:GLN:HE22	2:K:53:THR:H	1.44	0.66
2:H:157:PRO:O	2:H:210:HIS:HE1	1.78	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:198:HIS:CD2	1:J:200:GLY:H	2.13	0.65
1:J:161:GLU:HB2	1:J:175:LEU:HD21	1.78	0.65
3:W:34:ASP:O	3:W:37:GLN:HG2	1.96	0.65
2:K:40:ALA:HB3	2:K:43:LYS:HE3	1.78	0.64
1:J:190:LYS:HA	1:J:211:ARG:HB3	1.78	0.63
2:K:20:LEU:HD12	2:K:81:LEU:HD23	1.80	0.63
2:K:127:LYS:NZ	2:K:154:ASP:HB3	2.14	0.63
1:J:162:SER:HB2	1:J:176:SER:OG	2.00	0.61
1:J:123:GLU:O	1:J:126:LYS:HG3	2.00	0.61
1:L:28:ASP:OD1	1:L:68:GLY:HA2	2.00	0.60
3:V:92:GLY:HA2	2:K:101:HIS:CE1	2.37	0.59
2:K:161:THR:HB	2:K:209:ASN:HB3	1.84	0.59
2:K:127:LYS:HZ2	2:K:154:ASP:HB3	1.66	0.59
3:V:48:LYS:O	3:W:17:PHE:HE1	1.86	0.58
1:J:107:LYS:HA	1:J:140:TYR:OH	2.03	0.58
1:J:189:HIS:O	1:J:211:ARG:HD3	2.04	0.58
3:W:44:GLU:H	3:W:44:GLU:CD	2.11	0.57
3:V:59:GLY:HA2	3:W:32:LEU:HD13	1.86	0.57
2:H:151:LEU:HG	2:H:153:LYS:HG3	1.85	0.57
3:V:60:CYS:HG	3:W:51:CYS:HG	1.53	0.57
1:J:113:PRO:HD3	1:J:198:HIS:CD2	2.38	0.57
1:J:115:VAL:CG2	1:J:196:VAL:HG21	2.35	0.57
1:L:54:LEU:HD22	1:L:58:VAL:HB	1.86	0.56
2:K:100:PRO:HD3	2:K:109:TYR:O	2.06	0.56
1:L:147:GLN:HG3	1:L:154:LEU:HD13	1.89	0.55
3:V:34:ASP:O	3:V:37:GLN:HG2	2.06	0.55
1:L:198:HIS:HD2	1:L:200:GLY:H	1.53	0.55
2:H:40:ALA:HB3	2:H:43:LYS:HE3	1.89	0.55
1:J:154:LEU:HD12	1:J:154:LEU:H	1.71	0.54
1:J:54:LEU:HD22	1:J:58:VAL:HB	1.90	0.54
1:J:205:VAL:HG12	1:J:206:THR:N	2.23	0.53
2:H:53:THR:H	3:W:89:GLN:HE22	1.56	0.53
2:K:129:PRO:HA	2:K:154:ASP:O	2.08	0.53
3:V:44:GLU:H	3:V:44:GLU:CD	2.17	0.53
2:K:153:LYS:HG2	2:K:187:SER:OG	2.09	0.53
2:K:173:VAL:HA	2:K:191:VAL:O	2.09	0.53
2:H:12:VAL:HG11	2:H:86:LEU:HD13	1.91	0.52
1:L:141:PRO:O	1:L:198:HIS:HE1	1.93	0.52
1:L:40:PRO:CG	1:L:165:GLU:HG2	2.40	0.52
1:J:28:ASP:OD1	1:J:68:GLY:HA2	2.09	0.52
2:H:161:THR:HG22	5:H:283:HOH:O	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:V:77:THR:CG2	3:W:14:VAL:HG22	2.40	0.51
1:L:103:LYS:HE2	5:L:379:HOH:O	2.11	0.51
1:L:112:ALA:HB1	1:L:201:LEU:HD13	1.92	0.51
3:W:73:GLU:HB3	3:W:97:LEU:HD11	1.92	0.51
1:J:2:ILE:HD13	1:J:29:ILE:HG22	1.92	0.50
1:J:80:PRO:HA	1:J:106:ILE:HD13	1.93	0.50
2:H:67:ARG:C	2:H:68:PHE:HD1	2.18	0.50
1:J:166:GLN:HE21	1:J:171:SER:HB3	1.75	0.50
2:K:145:THR:HG23	2:K:194:VAL:O	2.11	0.50
2:K:206:CYS:SG	2:K:206:CYS:O	2.71	0.49
2:K:11:LEU:HD21	2:K:122:SER:HB3	1.94	0.49
3:V:77:THR:HG22	3:W:14:VAL:HG22	1.95	0.49
1:J:4:MET:HE3	1:J:23:CYS:SG	2.52	0.49
3:V:71:THR:CG2	3:V:103:GLU:HB2	2.42	0.49
1:J:120:PRO:HD3	1:J:132:VAL:HG22	1.95	0.49
1:J:150:VAL:HG21	1:J:179:LEU:HD21	1.95	0.49
1:J:201:LEU:HG	1:J:205:VAL:HG23	1.95	0.49
2:K:180:LEU:HD23	2:K:186:TYR:CE1	2.48	0.49
1:J:24:SER:HA	1:J:69:THR:O	2.13	0.48
3:W:71:THR:CG2	3:W:103:GLU:HB2	2.42	0.48
3:W:69:VAL:HG22	3:W:103:GLU:O	2.13	0.48
2:K:202:GLN:HG2	2:K:204:TYR:CZ	2.48	0.48
2:H:100:PRO:HD3	2:H:109:TYR:O	2.13	0.48
3:W:79:GLN:HA	3:W:92:GLY:O	2.14	0.48
1:L:4:MET:HE3	1:L:23:CYS:SG	2.53	0.48
1:J:161:GLU:HA	1:J:176:SER:O	2.14	0.47
3:W:71:THR:HG21	3:W:103:GLU:HB2	1.95	0.47
1:J:167:ASP:OD1	1:J:169:LYS:HB2	2.15	0.47
1:J:190:LYS:HG2	1:J:211:ARG:N	2.29	0.47
1:L:20:THR:HG21	5:L:323:HOH:O	2.13	0.47
2:H:68:PHE:N	2:H:68:PHE:CD1	2.83	0.47
1:J:12:SER:HA	1:J:105:GLU:O	2.15	0.47
1:L:52:SER:HB3	1:L:64:GLY:O	2.15	0.47
2:K:210:HIS:HB3	2:K:215:THR:HB	1.96	0.46
2:K:163:SER:OG	2:K:207:ASN:HB2	2.15	0.46
1:L:147:GLN:HE21	1:L:149:LYS:HG3	1.81	0.46
2:H:136:PRO:O	2:H:137:SER:HB2	2.15	0.46
2:K:97:ALA:HB1	2:K:110:PHE:HB3	1.97	0.46
2:H:30:THR:HG22	2:H:53:THR:O	2.16	0.46
1:J:120:PRO:HG3	1:J:186:TYR:CE1	2.51	0.46
3:V:71:THR:HG21	3:V:103:GLU:HB2	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:68:PHE:HD1	2:H:68:PHE:N	2.14	0.46
1:J:150:VAL:HG22	1:J:192:TYR:CD1	2.51	0.46
1:J:66:GLY:O	1:J:67:SER:HB3	2.16	0.45
1:J:100:GLN:H	1:J:100:GLN:CD	2.25	0.45
1:J:181:LEU:HD13	1:J:186:TYR:HD1	1.82	0.45
2:K:210:HIS:HD2	2:K:213:SER:OG	1.98	0.45
2:H:11:LEU:HD21	2:H:122:SER:HB3	1.98	0.45
1:J:124:GLN:O	1:J:127:SER:HB2	2.17	0.45
1:J:23:CYS:SG	1:J:88:CYS:SG	3.10	0.45
2:H:83:MET:HE1	2:H:94:TYR:CZ	2.51	0.45
1:L:145:LYS:HB3	1:L:197:THR:HB	1.99	0.45
3:V:39:TYR:HB3	3:V:42:GLU:HG2	1.99	0.45
3:V:77:THR:O	3:W:14:VAL:HA	2.16	0.44
2:H:154:ASP:HB3	2:H:185:LEU:HD13	1.99	0.44
2:K:181:GLN:N	2:K:185:LEU:O	2.48	0.44
2:H:183:SER:CA	5:H:331:HOH:O	2.34	0.44
3:V:33:VAL:HG21	3:V:54:LEU:HD12	1.99	0.44
3:W:67:GLU:HG3	3:W:69:VAL:HG13	2.00	0.44
1:J:116:PHE:O	1:J:134:CYS:HA	2.18	0.44
1:J:149:LYS:HE2	1:J:154:LEU:HG	2.00	0.44
1:J:190:LYS:HG2	1:J:211:ARG:H	1.83	0.44
1:L:2:ILE:HD13	1:L:29:ILE:HG22	2.00	0.44
2:H:17:SER:HA	2:H:83:MET:O	2.18	0.44
2:K:129:PRO:HB2	2:K:152:VAL:CG1	2.44	0.44
3:W:33:VAL:HG21	3:W:54:LEU:HD12	2.00	0.44
1:L:147:GLN:NE2	1:L:149:LYS:HG3	2.33	0.43
2:H:102:TYR:HB2	2:H:105:SER:O	2.18	0.43
1:L:147:GLN:HG3	1:L:154:LEU:CD1	2.48	0.43
2:H:182:SER:O	5:H:331:HOH:O	2.20	0.43
3:W:39:TYR:HB3	3:W:42:GLU:HG2	2.00	0.43
1:J:37:GLN:HB2	1:J:47:LEU:HD11	2.00	0.43
1:L:108:ARG:HD3	1:L:109:THR:O	2.18	0.43
1:L:201:LEU:HG	1:L:205:VAL:HG23	2.01	0.43
3:W:75:ASN:HA	3:W:96:PHE:O	2.18	0.43
1:L:164:THR:CG2	5:L:244:HOH:O	2.67	0.43
2:H:97:ALA:HB1	2:H:110:PHE:HB3	2.01	0.43
1:L:37:GLN:HG3	1:L:86:TYR:CE2	2.54	0.43
1:J:135:LEU:HD22	2:K:191:VAL:HG11	2.01	0.43
1:J:147:GLN:NE2	1:J:149:LYS:HG3	2.34	0.43
2:K:191:VAL:CG1	2:K:192:VAL:N	2.82	0.42
1:J:115:VAL:HG21	1:J:196:VAL:HG21	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:V:82:ARG:HD2	2:K:102:TYR:CZ	2.55	0.42
2:K:70:PHE:CE1	2:K:81:LEU:HD13	2.55	0.42
1:J:183:LYS:O	1:J:186:TYR:HB3	2.21	0.41
3:V:101:LYS:HB2	3:V:101:LYS:HE2	1.83	0.41
1:J:115:VAL:HG22	1:J:196:VAL:HG21	2.03	0.41
2:K:172:GLY:O	2:K:192:VAL:HA	2.20	0.41
1:L:100:GLN:CD	1:L:100:GLN:H	2.29	0.41
1:J:125:LEU:HD12	1:J:129:THR:O	2.21	0.41
2:K:68:PHE:N	2:K:68:PHE:CD1	2.89	0.41
2:K:70:PHE:HE1	2:K:81:LEU:HD13	1.85	0.41
3:V:24:SER:HB3	3:W:53:PRO:HD3	2.03	0.41
2:K:153:LYS:HA	2:K:187:SER:OG	2.21	0.41
2:K:162:VAL:HG11	2:K:190:SER:CB	2.51	0.41
1:L:164:THR:HG22	1:L:174:SER:H	1.86	0.41
2:H:102:TYR:CD1	3:W:92:GLY:HA3	2.56	0.40
2:H:108:TRP:CZ3	3:W:88:GLY:HA3	2.55	0.40
3:W:72:GLU:HB3	3:W:101:LYS:HB2	2.02	0.40
2:H:61:ALA:HB3	2:H:64:PHE:HD2	1.86	0.40
2:H:210:HIS:CD2	2:H:213:SER:OG	2.66	0.40
3:W:66:LEU:O	3:W:107:LYS:HE3	2.21	0.40
1:J:112:ALA:HB1	1:J:201:LEU:CD1	2.50	0.40
2:K:208:VAL:HG12	2:K:209:ASN:N	2.36	0.40
2:H:70:PHE:CE1	2:H:81:LEU:HD13	2.57	0.40
3:V:89:GLN:HE21	3:V:89:GLN:HB3	1.74	0.40
1:J:118:PHE:HA	1:J:119:PRO:HD3	1.79	0.40
1:J:81:GLU:H	1:J:81:GLU:CD	2.28	0.40
1:J:135:LEU:CD2	2:K:191:VAL:HG11	2.51	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	J	211/214 (99%)	191 (90%)	19 (9%)	1 (0%)	24	37
1	L	211/214 (99%)	205 (97%)	6 (3%)	0	100	100
2	H	214/231 (93%)	209 (98%)	5 (2%)	0	100	100
2	K	214/231 (93%)	196 (92%)	17 (8%)	1 (0%)	24	37
3	V	92/102 (90%)	85 (92%)	6 (6%)	1 (1%)	11	18
3	W	92/102 (90%)	88 (96%)	3 (3%)	1 (1%)	11	18
All	All	1034/1094 (94%)	974 (94%)	56 (5%)	4 (0%)	30	43

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	J	138	ASN
3	V	26	CYS
3	W	26	CYS
2	K	159	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	J	190/191 (100%)	175 (92%)	15 (8%)	11	19
1	L	190/191 (100%)	177 (93%)	13 (7%)	14	25
2	H	182/194 (94%)	170 (93%)	12 (7%)	15	26
2	K	182/194 (94%)	172 (94%)	10 (6%)	19	34
3	V	89/96 (93%)	87 (98%)	2 (2%)	45	67
3	W	89/96 (93%)	86 (97%)	3 (3%)	32	54
All	All	922/962 (96%)	867 (94%)	55 (6%)	17	31

All (55) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	33	LEU

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Mol	Chain	Res	Type
1	L	48	ILE
1	L	54	LEU
1	L	100	GLN
1	L	105	GLU
1	L	108	ARG
1	L	125	LEU
1	L	129	THR
1	L	143	GLU
1	L	152	ASN
1	L	164	THR
1	L	190	LYS
1	L	201	LEU
2	H	1	GLU
2	H	11	LEU
2	H	18	LEU
2	H	115	GLN
2	H	117	THR
2	H	118	LEU
2	H	120	THR
2	H	160	VAL
2	H	180	LEU
2	H	188	LEU
2	H	191	VAL
2	H	196	SER
3	V	43	ILE
3	V	89	GLN
3	W	24	SER
3	W	43	ILE
3	W	89	GLN
1	J	33	LEU
1	J	54	LEU
1	J	94	VAL
1	J	100	GLN
1	J	105	GLU
1	J	108	ARG
1	J	122	ASP
1	J	125	LEU
1	J	126	LYS
1	J	147	GLN
1	J	152	ASN
1	J	164	THR
1	J	175	LEU

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Mol	Chain	Res	Type
1	J	181	LEU
1	J	201	LEU
2	K	1	GLU
2	K	11	LEU
2	K	115	GLN
2	K	117	THR
2	K	118	LEU
2	K	120	THR
2	K	126	THR
2	K	152	VAL
2	K	159	PRO
2	K	180	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	31	ASN
1	L	38	GLN
1	L	55	HIS
1	L	147	GLN
1	L	152	ASN
1	L	198	HIS
2	H	39	GLN
2	H	101	HIS
2	H	209	ASN
2	H	210	HIS
3	V	22	GLN
3	V	37	GLN
3	V	89	GLN
3	W	22	GLN
3	W	27	HIS
3	W	89	GLN
1	J	3	GLN
1	J	55	HIS
1	J	79	GLN
1	J	137	ASN
1	J	147	GLN
1	J	152	ASN
1	J	160	GLN
2	K	101	HIS
2	K	174	HIS
2	K	202	GLN

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Mol	Chain	Res	Type
2	K	210	HIS

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 ligands 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	$\# Z > 2$
4	SO4	K	232	-	4,4,4	0.76	0	6,6,6	0.55	0
4	SO4	H	232	-	4,4,4	0.59	0	6,6,6	0.60	0

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, 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	J	213/214 (99%)	0.29	20 (9%) 14 11	13, 52, 120, 132	0
1	L	213/214 (99%)	-0.92	0 100 100	7, 22, 43, 70	0
2	H	218/231 (94%)	-0.94	0 100 100	9, 21, 43, 60	0
2	K	218/231 (94%)	0.15	17 (7%) 19 15	13, 41, 119, 140	0
3	V	94/102 (92%)	-0.42	0 100 100	11, 40, 68, 76	0
3	W	94/102 (92%)	-0.39	0 100 100	12, 38, 63, 76	0
All	All	1050/1094 (95%)	-0.36	37 (3%) 47 43	7, 30, 114, 140	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	J	157	GLY	5.0
1	J	203	SER	4.1
1	J	144	ALA	3.7
2	K	184	GLY	3.6
1	J	182	SER	3.5
2	K	183	SER	3.5
2	K	201	THR	3.3
1	J	127	SER	3.1
1	J	148	TRP	3.1
2	K	221	VAL	3.0
1	J	194	CYS	3.0
1	J	121	SER	2.9
1	J	156	SER	2.9
2	K	215	THR	2.9
2	K	200	GLY	2.8
2	K	179	VAL	2.7
2	K	172	GLY	2.6
2	K	167	GLY	2.6
1	J	158	ASN	2.6

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Mol	Chain	Res	Type	RSRZ
1	J	147	GLN	2.6
1	J	119	PRO	2.5
2	K	171	SER	2.5
2	K	193	THR	2.5
1	J	117	ILE	2.5
1	J	150	VAL	2.4
2	K	144	GLY	2.4
2	K	218	ASP	2.4
1	J	210	ASN	2.4
2	K	205	ILE	2.4
1	J	115	VAL	2.3
2	K	180	LEU	2.3
1	J	208	SER	2.2
2	K	198	SER	2.1
1	J	193	ALA	2.1
1	J	129	THR	2.0
2	K	134	LEU	2.0
1	J	140	TYR	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 oligosaccharides 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, 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
4	SO4	K	232	5/5	0.96	0.09	58,61,64,66	0
4	SO4	H	232	5/5	0.99	0.05	37,42,48,50	0

6.5 Other polymers [i](#)

There are no such residues in this entry.