



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 08:59 AM UTC

PDB ID : 3U9P / pdb_00003u9p
Title : Crystal Structure of Murine Siderocalin in Complex with an Fab Fragment
Authors : Correnti, C.; Strong, R.K.
Deposited on : 2011-10-19
Resolution : 2.80 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
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

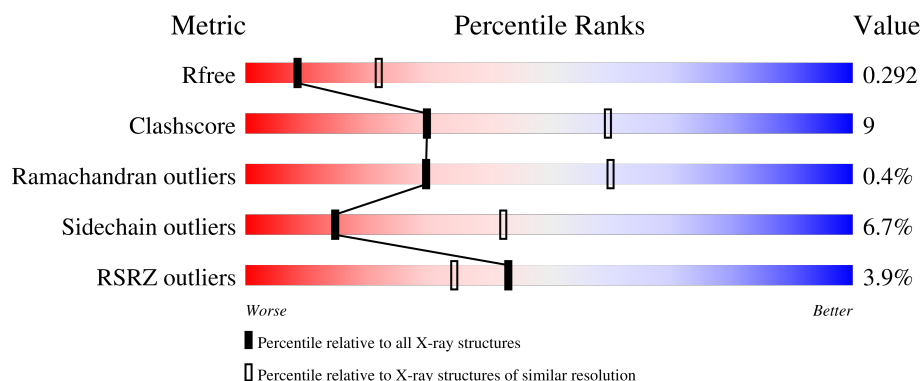
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	C	182	0% 77% 14% 7%
1	D	182	5% 73% 17% 9%
2	H	216	11% 67% 17% 15%
2	K	216	2% 69% 21% 7%
3	L	212	0% 82% 16% 2%

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
3	M	212	% 84% 15%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8339 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 Neutrophil gelatinase-associated lipocalin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	169	Total	C	N	O	S	0	0	0
			1266	826	209	226	5			
1	D	166	Total	C	N	O	S	0	0	0
			1229	801	206	217	5			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-7	GLY	-	expression tag	UNP P11672
C	-6	SER	-	expression tag	UNP P11672
D	-7	GLY	-	expression tag	UNP P11672
D	-6	SER	-	expression tag	UNP P11672

- Molecule 2 is a protein called Monoclonal Fab Fragment Heavy Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	184	Total	C	N	O	S	0	0	0
			1229	780	209	233	7			
2	K	200	Total	C	N	O	S	0	0	0
			1414	902	237	268	7			

- Molecule 3 is a protein called Monoclonal Fab Fragment Light Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	L	212	Total	C	N	O	S	0	0	0
			1561	986	254	314	7			
3	M	211	Total	C	N	O	S	0	0	0
			1579	997	257	318	7			

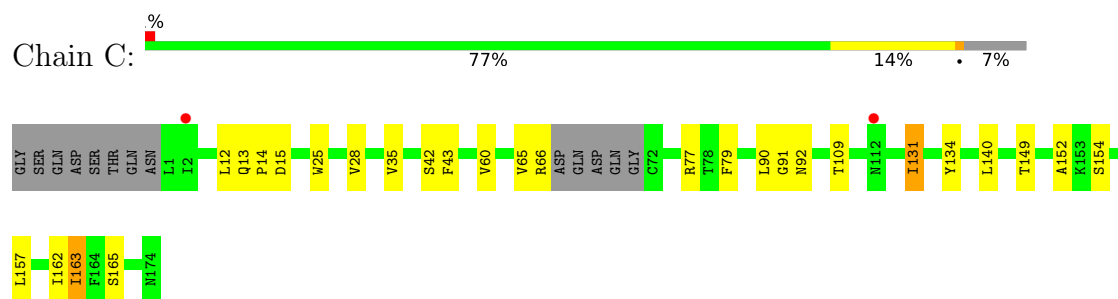
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	4	Total 4	O 4	0	0
4	D	5	Total 5	O 5	0	0
4	H	5	Total 5	O 5	0	0
4	K	15	Total 15	O 15	0	0
4	L	18	Total 18	O 18	0	0
4	M	14	Total 14	O 14	0	0

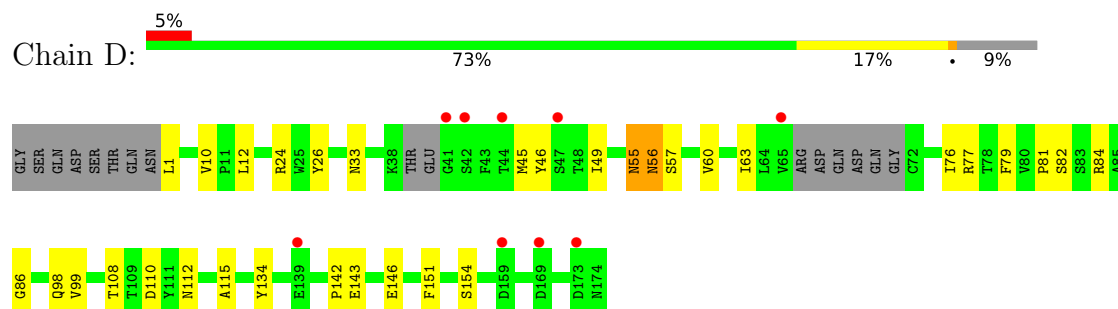
3 Residue-property plots [i](#)

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.

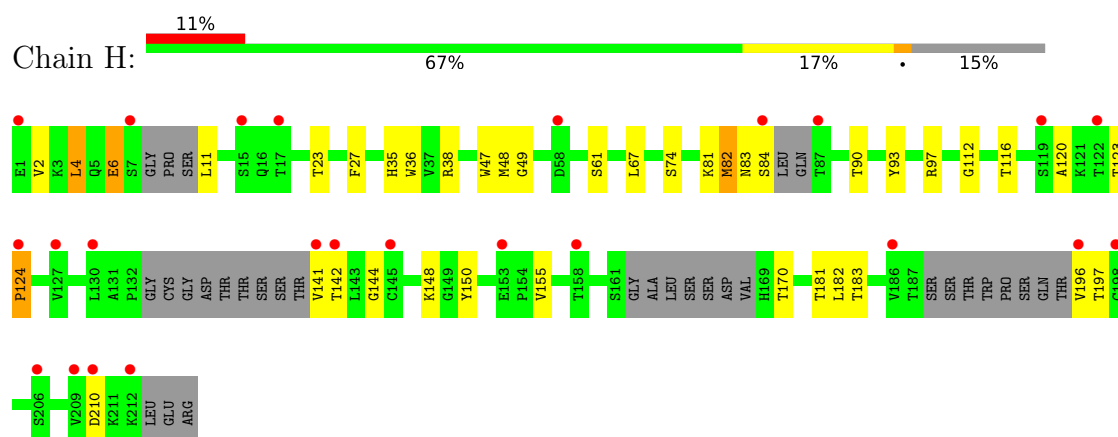
- Molecule 1: Neutrophil gelatinase-associated lipocalin



- Molecule 1: Neutrophil gelatinase-associated lipocalin

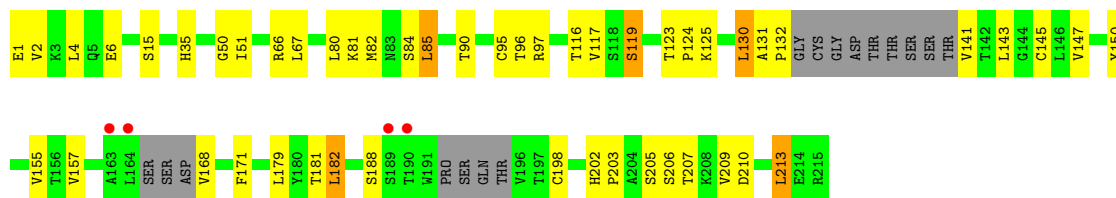


- Molecule 2: Monoclonal Fab Fragment Heavy Chain



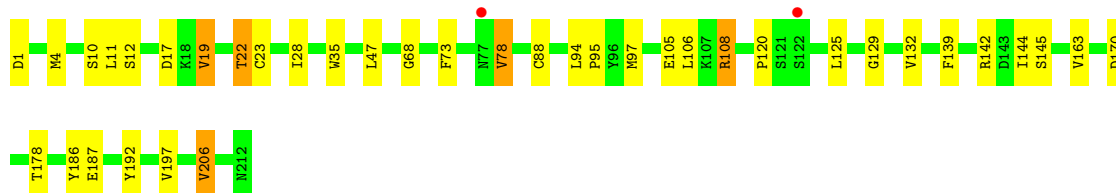
- Molecule 2: Monoclonal Fab Fragment Heavy Chain

Chain K: 



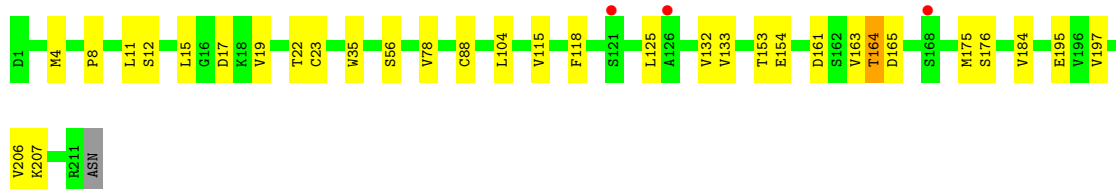
• Molecule 3: Monoclonal Fab Fragment Light Chain

Chain L: 



• Molecule 3: Monoclonal Fab Fragment Light Chain

Chain M: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	82.13Å 124.09Å 147.49Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.51 – 2.80 49.51 – 2.80	Depositor EDS
% Data completeness (in resolution range)	93.1 (49.51-2.80) 93.1 (49.51-2.80)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.12 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.5.0072	Depositor
R, R_{free}	0.219 , 0.292 0.222 , 0.292	Depositor DCC
R_{free} test set	1758 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	46.6	Xtriage
Anisotropy	0.198	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 62.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	8339	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.89% 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](#)

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	C	0.63	0/1299	0.85	0/1771
1	D	0.59	0/1261	0.86	2/1720 (0.1%)
2	H	0.66	0/1255	0.83	1/1726 (0.1%)
2	K	0.70	0/1447	0.89	0/1989
3	L	0.62	0/1596	0.83	0/2178
3	M	0.67	0/1614	0.84	0/2200
All	All	0.65	0/8472	0.85	3/11584 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	10	VAL	CA-C-N	-6.20	114.26	120.21
1	D	10	VAL	C-N-CA	-6.20	114.26	120.21
2	H	144	GLY	N-CA-C	5.74	119.67	111.12

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	C	1266	0	1135	19	0
1	D	1229	0	1078	17	1
2	H	1229	0	992	25	1
2	K	1414	0	1269	36	1

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	L	1561	0	1405	25	0
3	M	1579	0	1460	20	0
4	C	4	0	0	0	0
4	D	5	0	0	0	0
4	H	5	0	0	0	0
4	K	15	0	0	1	0
4	L	18	0	0	1	0
4	M	14	0	0	0	0
All	All	8339	0	7339	133	2

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 9.

All (133) 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:123:THR:HG22	2:H:124:PRO:CD	2.04	0.86
2:K:143:LEU:HB3	2:K:213:LEU:HD11	1.62	0.80
2:K:130:LEU:CD2	3:M:118:PHE:CG	2.68	0.77
2:H:90:THR:HG23	2:H:116:THR:HA	1.73	0.71
3:L:4:MET:HE3	3:L:23:CYS:SG	2.30	0.71
3:L:17:ASP:O	3:L:78:VAL:HG23	1.93	0.69
2:H:6:GLU:OE1	2:H:112:GLY:N	2.27	0.68
1:D:115:ALA:HB3	1:D:134:TYR:HB2	1.75	0.68
2:H:123:THR:HG22	2:H:124:PRO:HD2	1.76	0.66
3:M:161:ASP:HB3	3:M:175:MET:HE2	1.77	0.66
2:H:67:LEU:HD13	2:H:82:MET:HG3	1.76	0.65
1:C:77:ARG:HG3	1:C:90:LEU:HD11	1.78	0.65
2:K:130:LEU:HD23	3:M:118:PHE:CD2	2.31	0.65
2:K:130:LEU:HD23	3:M:118:PHE:CG	2.31	0.65
3:M:4:MET:HE3	3:M:23:CYS:SG	2.37	0.65
2:K:6:GLU:HG3	2:K:95:CYS:HB2	1.79	0.65
2:H:123:THR:HG22	2:H:124:PRO:HD3	1.79	0.64
1:C:60:VAL:HG21	1:C:79:PHE:CD1	2.31	0.64
3:M:164:THR:HG23	3:M:165:ASP:O	1.97	0.64
3:L:12:SER:HA	3:L:105:GLU:O	1.99	0.63
3:L:142:ARG:NH2	3:L:163:VAL:HG21	2.15	0.61
2:H:36:TRP:HB3	2:H:48:MET:HE3	1.83	0.60
2:K:66:ARG:NH2	2:K:82:MET:CE	2.64	0.60
2:H:182:LEU:O	2:H:183:THR:HG23	2.01	0.59
2:K:90:THR:HG23	2:K:116:THR:HA	1.83	0.59

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:125:LEU:HD23	3:L:129:GLY:O	2.03	0.59
1:C:149:THR:HG23	1:C:162:ILE:CD1	2.34	0.58
3:L:206:VAL:O	3:L:206:VAL:HG22	2.02	0.58
3:L:142:ARG:CZ	3:L:163:VAL:HG21	2.34	0.57
2:H:155:VAL:HG22	2:H:182:LEU:HD13	1.87	0.56
1:C:43:PHE:CE2	1:C:163:ILE:HG22	2.40	0.56
2:K:80:LEU:C	2:K:80:LEU:HD23	2.32	0.55
1:D:26:TYR:HB3	1:D:45:MET:HE2	1.89	0.55
3:M:8:PRO:HG3	3:M:11:LEU:HD13	1.87	0.55
3:L:1:ASP:HB3	3:L:97:MET:HE1	1.89	0.54
3:M:206:VAL:O	3:M:206:VAL:HG23	2.08	0.54
1:C:43:PHE:HE2	1:C:163:ILE:HG22	1.72	0.54
2:K:66:ARG:CZ	2:K:82:MET:HE3	2.39	0.53
2:H:47:TRP:CZ2	2:H:49:GLY:HA2	2.42	0.53
2:K:66:ARG:NH2	2:K:82:MET:HE2	2.24	0.53
3:L:142:ARG:NE	3:L:163:VAL:HG11	2.24	0.53
1:D:84:ARG:HD3	3:L:94:LEU:HD12	1.90	0.53
3:L:108:ARG:NH1	3:L:170:ASP:O	2.40	0.52
3:M:35:TRP:CZ3	3:M:88:CYS:HB3	2.44	0.52
1:D:55:ASN:O	1:D:56:ASN:CB	2.56	0.52
2:H:155:VAL:CG2	2:H:182:LEU:HD13	2.40	0.52
2:H:83:ASN:O	2:H:84:SER:CB	2.58	0.52
3:L:206:VAL:O	3:L:206:VAL:CG2	2.54	0.52
2:H:38:ARG:HB3	2:H:48:MET:HE2	1.93	0.51
1:C:25:TRP:CE3	1:C:134:TYR:HB3	2.45	0.51
1:D:76:ILE:O	1:D:77:ARG:HD2	2.10	0.51
2:K:168:VAL:N	4:K:218:HOH:O	2.44	0.51
3:L:23:CYS:SG	3:L:88:CYS:SG	3.01	0.51
2:K:130:LEU:CD2	3:M:118:PHE:CB	2.88	0.51
3:L:1:ASP:HB3	3:L:97:MET:CE	2.42	0.51
2:H:150:TYR:CZ	2:H:155:VAL:HG13	2.47	0.50
2:K:124:PRO:HG3	2:K:205:SER:HB2	1.94	0.49
2:K:130:LEU:CD2	3:M:118:PHE:CD2	2.93	0.49
1:C:28:VAL:HG21	1:C:140:LEU:HD12	1.94	0.49
1:C:77:ARG:CG	1:C:90:LEU:HD11	2.43	0.49
1:C:152:ALA:HB1	1:C:157:LEU:HD12	1.94	0.49
1:D:46:TYR:HB2	1:D:63:ILE:O	2.13	0.49
1:D:151:PHE:O	1:D:154:SER:CB	2.61	0.48
3:M:115:VAL:O	3:M:207:LYS:HE2	2.12	0.48
3:L:145:SER:HB3	3:L:197:VAL:HG23	1.95	0.48
1:D:60:VAL:HG21	1:D:79:PHE:CD1	2.48	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:120:PRO:HD3	3:L:132:VAL:HG22	1.95	0.48
2:K:130:LEU:HD23	3:M:118:PHE:CB	2.43	0.48
2:K:124:PRO:HG3	2:K:205:SER:CB	2.44	0.47
2:K:157:VAL:HG21	2:K:182:LEU:HD21	1.95	0.47
3:M:195:GLU:CG	3:M:206:VAL:HG12	2.45	0.47
2:H:4:LEU:HD12	2:H:4:LEU:N	2.29	0.47
2:H:123:THR:CG2	2:H:124:PRO:HD3	2.44	0.47
2:H:196:VAL:O	2:H:196:VAL:HG23	2.15	0.47
2:K:67:LEU:HD11	2:K:80:LEU:HG	1.97	0.47
2:K:130:LEU:HD11	3:M:133:VAL:HG21	1.96	0.47
2:K:150:TYR:CE2	2:K:155:VAL:HG13	2.50	0.47
3:L:35:TRP:CE2	3:L:73:PHE:HB2	2.49	0.47
1:C:43:PHE:CZ	1:C:165:SER:HB3	2.50	0.47
1:D:1:LEU:HD22	1:D:33:ASN:ND2	2.30	0.47
1:D:142:PRO:O	1:D:146:GLU:N	2.48	0.47
3:L:139:PHE:CD2	3:L:144:ILE:HD12	2.50	0.47
2:K:202:HIS:HB3	2:K:207:THR:HB	1.95	0.46
2:H:148:LYS:HA	2:H:181:THR:HG23	1.98	0.46
1:C:14:PRO:O	1:C:15:ASP:C	2.59	0.46
2:H:48:MET:HE1	2:H:93:TYR:CD2	2.50	0.46
2:K:141:VAL:N	2:K:188:SER:O	2.49	0.46
3:L:22:THR:HG22	4:L:228:HOH:O	2.16	0.46
3:L:4:MET:HE3	3:L:23:CYS:HG	1.81	0.46
1:D:110:ASP:C	1:D:112:ASN:H	2.24	0.45
2:H:123:THR:CG2	2:H:124:PRO:CD	2.88	0.45
2:K:209:VAL:CG1	2:K:210:ASP:N	2.79	0.45
2:K:80:LEU:HD23	2:K:81:LYS:N	2.32	0.45
2:H:150:TYR:CE2	2:H:155:VAL:HG13	2.51	0.45
3:L:19:VAL:HG13	3:L:78:VAL:HG22	1.98	0.44
1:C:149:THR:HG23	1:C:162:ILE:HD12	1.99	0.44
2:H:67:LEU:HD12	2:H:81:LYS:O	2.18	0.44
1:C:35:VAL:HG13	1:C:163:ILE:HD12	2.00	0.44
2:K:4:LEU:HD23	2:K:95:CYS:SG	2.57	0.44
2:K:66:ARG:NH2	2:K:82:MET:HE3	2.32	0.44
1:D:12:LEU:HD11	1:D:86:GLY:HA3	1.99	0.44
2:H:48:MET:HE1	2:H:93:TYR:HD2	1.83	0.43
1:C:28:VAL:HG21	1:C:140:LEU:CD1	2.48	0.43
1:D:60:VAL:HG21	1:D:79:PHE:CG	2.54	0.43
3:L:11:LEU:HD12	3:L:12:SER:H	1.83	0.43
1:C:13:GLN:HG2	1:C:109:THR:HG22	2.00	0.43
2:K:85:LEU:HD13	2:K:117:VAL:HG21	2.00	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:24:ARG:HA	1:D:49:ILE:HA	1.99	0.43
2:K:202:HIS:O	2:K:203:PRO:C	2.61	0.43
2:K:125:LYS:CB	2:K:147:VAL:HG13	2.49	0.42
3:L:28:ILE:HG23	3:L:68:GLY:HA2	2.01	0.42
3:M:163:VAL:HG22	3:M:175:MET:HG3	2.01	0.42
2:H:35:HIS:CD2	2:H:47:TRP:HE1	2.37	0.42
2:K:35:HIS:ND1	2:K:50:GLY:HA3	2.35	0.42
2:K:51:ILE:HG23	2:K:51:ILE:O	2.20	0.42
3:M:17:ASP:O	3:M:78:VAL:HG23	2.20	0.42
3:M:153:THR:O	3:M:154:GLU:C	2.63	0.42
1:C:91:GLY:O	1:C:92:ASN:C	2.63	0.42
2:K:198:CYS:O	2:K:210:ASP:HA	2.19	0.42
1:C:12:LEU:O	1:C:13:GLN:C	2.63	0.42
3:L:186:TYR:O	3:L:192:TYR:OH	2.37	0.42
1:C:65:VAL:O	1:C:66:ARG:CB	2.68	0.42
2:K:131:ALA:HB1	2:K:132:PRO:HD2	2.03	0.41
1:D:110:ASP:O	1:D:112:ASN:N	2.42	0.41
2:K:143:LEU:HD13	2:K:213:LEU:HD11	2.03	0.41
1:D:57:SER:HA	1:D:81:PRO:HD3	2.02	0.41
3:M:11:LEU:HB3	3:M:104:LEU:HD12	2.02	0.41
2:K:205:SER:O	2:K:206:SER:C	2.63	0.41
1:D:98:GLN:HG2	1:D:99:VAL:HG23	2.03	0.41
2:H:2:VAL:HG13	2:H:27:PHE:CD2	2.56	0.41
2:K:171:PHE:CD1	3:M:176:SER:HB2	2.56	0.41
1:C:131:ILE:HD11	1:C:152:ALA:HB2	2.03	0.40
3:L:94:LEU:HA	3:L:95:PRO:C	2.46	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:143:GLU:OE1	2:H:210:ASP:O[3_455]	2.02	0.18
2:K:1:GLU:N	2:K:119:SER:OG[4_446]	2.18	0.02

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries

of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	165/182 (91%)	155 (94%)	9 (6%)	1 (1%)	21	51
1	D	160/182 (88%)	149 (93%)	10 (6%)	1 (1%)	21	51
2	H	172/216 (80%)	162 (94%)	8 (5%)	2 (1%)	10	34
2	K	192/216 (89%)	183 (95%)	9 (5%)	0	100	100
3	L	210/212 (99%)	203 (97%)	7 (3%)	0	100	100
3	M	209/212 (99%)	203 (97%)	6 (3%)	0	100	100
All	All	1108/1220 (91%)	1055 (95%)	49 (4%)	4 (0%)	30	60

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	H	124	PRO
1	D	56	ASN
2	H	120	ALA
1	C	42	SER

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	C	116/165 (70%)	113 (97%)	3 (3%)	40	75
1	D	107/165 (65%)	104 (97%)	3 (3%)	38	73
2	H	98/190 (52%)	86 (88%)	12 (12%)	5	16
2	K	135/190 (71%)	121 (90%)	14 (10%)	7	22
3	L	157/189 (83%)	147 (94%)	10 (6%)	16	44
3	M	165/189 (87%)	155 (94%)	10 (6%)	17	46
All	All	778/1088 (72%)	726 (93%)	52 (7%)	15	42

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

Mol	Chain	Res	Type
1	C	131	ILE
1	C	154	SER
1	C	163	ILE
1	D	55	ASN
1	D	82	SER
1	D	108	THR
2	H	4	LEU
2	H	6	GLU
2	H	11	LEU
2	H	23	THR
2	H	61	SER
2	H	74	SER
2	H	82	MET
2	H	97	ARG
2	H	141	VAL
2	H	142	THR
2	H	170	THR
2	H	197	THR
2	K	2	VAL
2	K	15	SER
2	K	84	SER
2	K	85	LEU
2	K	96	THR
2	K	97	ARG
2	K	119	SER
2	K	123	THR
2	K	130	LEU
2	K	145	CYS
2	K	179	LEU
2	K	181	THR
2	K	182	LEU
2	K	213	LEU
3	L	10	SER
3	L	19	VAL
3	L	22	THR
3	L	47	LEU
3	L	78	VAL
3	L	106	LEU
3	L	108	ARG
3	L	178	THR
3	L	187	GLU
3	L	206	VAL

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
3	M	12	SER
3	M	15	LEU
3	M	19	VAL
3	M	22	THR
3	M	56	SER
3	M	125	LEU
3	M	132	VAL
3	M	164	THR
3	M	184	VAL
3	M	197	VAL

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

Mol	Chain	Res	Type
1	C	36	GLN
1	C	55	ASN
1	C	56	ASN
1	C	87	GLN
1	C	100	GLN
1	D	92	ASN
2	K	13	GLN
2	K	39	GLN
3	L	37	GLN
3	L	124	GLN
3	L	137	ASN
3	M	38	GLN
3	M	93	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 oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

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	C	169/182 (92%)	0.10	2 (1%) 76 68	24, 44, 65, 74	0
1	D	166/182 (91%)	0.48	9 (5%) 31 24	34, 57, 92, 99	0
2	H	184/216 (85%)	0.76	24 (13%) 7 6	35, 67, 100, 114	0
2	K	200/216 (92%)	0.06	4 (2%) 65 56	21, 35, 62, 76	0
3	L	212/212 (100%)	-0.03	2 (0%) 81 74	22, 38, 74, 101	0
3	M	211/212 (99%)	-0.07	3 (1%) 73 64	19, 37, 54, 64	0
All	All	1142/1220 (93%)	0.20	44 (3%) 43 34	19, 45, 86, 114	0

All (44) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	K	164	LEU	4.7
2	H	87	THR	4.6
2	H	141	VAL	4.4
2	H	212	LYS	3.8
2	H	186	VAL	3.5
2	H	153	GLU	3.5
2	H	210	ASP	3.4
2	H	145	CYS	3.2
2	K	189	SER	3.2
2	K	163	ALA	3.2
1	D	65	VAL	3.0
2	H	206	SER	3.0
3	L	77	ASN	2.8
2	K	190	THR	2.8
2	H	209	VAL	2.7
1	D	47	SER	2.7
2	H	84	SER	2.7
3	L	122	SER	2.7
1	D	169	ASP	2.6

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
2	H	142	THR	2.6
2	H	158	THR	2.5
2	H	15	SER	2.5
2	H	130	LEU	2.5
1	D	42	SER	2.5
1	D	44	THR	2.5
3	M	168	SER	2.4
2	H	17	THR	2.3
1	D	139	GLU	2.3
2	H	1	GLU	2.3
1	D	159	ASP	2.3
2	H	127	VAL	2.3
2	H	58	ASP	2.2
1	C	112	ASN	2.2
1	D	173	ASP	2.2
2	H	124	PRO	2.2
2	H	122	THR	2.2
2	H	119	SER	2.2
1	C	2	ILE	2.1
2	H	198	CYS	2.1
3	M	126	ALA	2.1
1	D	41	GLY	2.1
2	H	196	VAL	2.0
2	H	7	SER	2.0
3	M	121	SER	2.0

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

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

6.3 Carbohydrates [i](#)

There are no 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.