



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 26, 2026 – 08:09 AM EDT

PDB ID : 7T82 / pdb_00007t82
Title : Crystal Structure of LEUKOCIDIN E/CENTYRIN S26/FAB B438
Authors : Luo, J.; Malia, T.J.; Buckley, P.T.
Deposited on : 2021-12-15
Resolution : 3.50 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

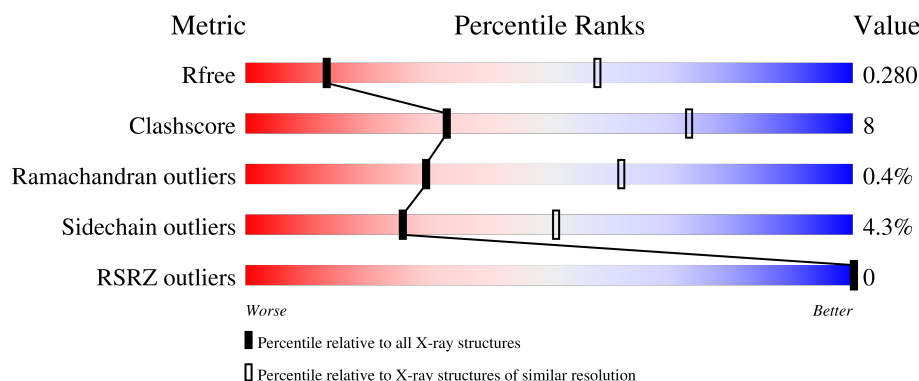
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.50 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	1085 (3.54-3.46)
Clashscore	190562	1140 (3.54-3.46)
Ramachandran outliers	187476	1113 (3.54-3.46)
Sidechain outliers	187428	1114 (3.54-3.46)
RSRZ outliers	180081	1084 (3.54-3.46)

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	A	294	 79% 19% •
1	B	294	 78% 21% ••
2	C	219	 81% 17% ••
2	L	219	 81% 16% ••
3	D	229	 70% 20% • 6%

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Mol	Chain	Length	Quality of chain
3	H	229	69%24%6%
4	E	98	85%12%••
4	F	98	81%16%••

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 12610 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 Leukocidin E.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	293	Total	C	N	O	S	0	0	0
			2331	1475	412	442	2			
1	B	292	Total	C	N	O	S	0	0	0
			2320	1467	407	444	2			

- Molecule 2 is a protein called Antibody Fab Light Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	217	Total	C	N	O	S	0	0	0
			1664	1048	276	335	5			
2	L	217	Total	C	N	O	S	0	0	0
			1664	1048	276	335	5			

- Molecule 3 is a protein called Antibody Fab Heavy Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	216	Total	C	N	O	S	0	0	0
			1615	1025	265	318	7			
3	H	216	Total	C	N	O	S	0	0	0
			1616	1025	265	319	7			

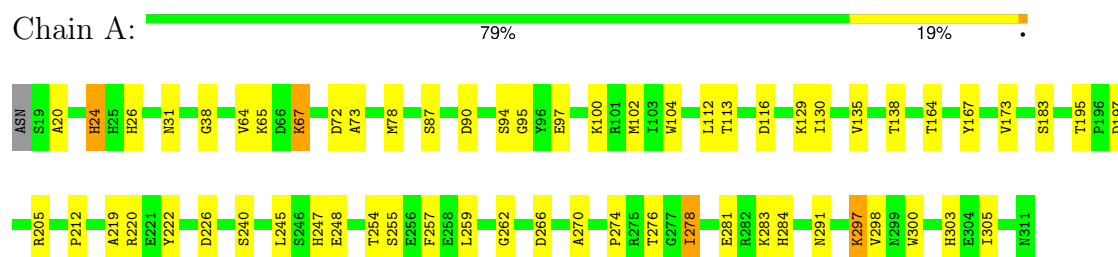
- Molecule 4 is a protein called Centyrin S26.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	E	96	Total	C	N	O	0	0	1
			700	450	116	134			
4	F	96	Total	C	N	O	0	0	1
			700	450	116	134			

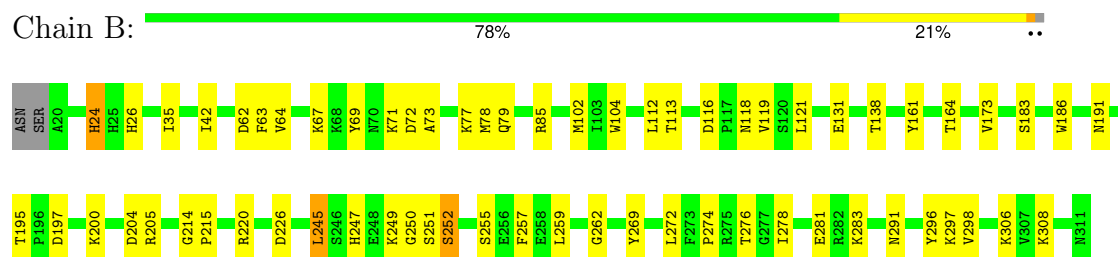
3 Residue-property plots

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.

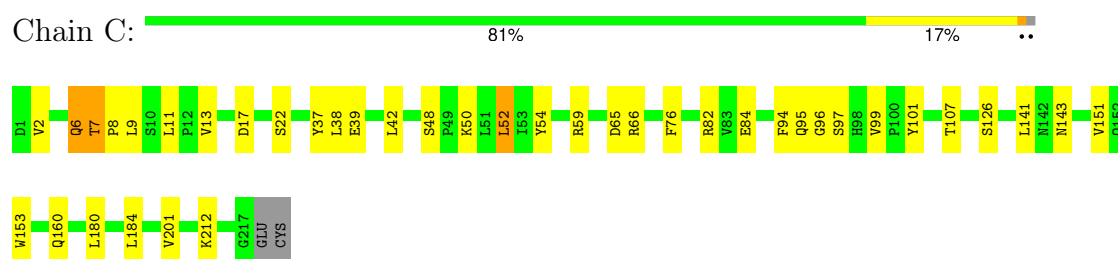
• Molecule 1: Leukocidin E



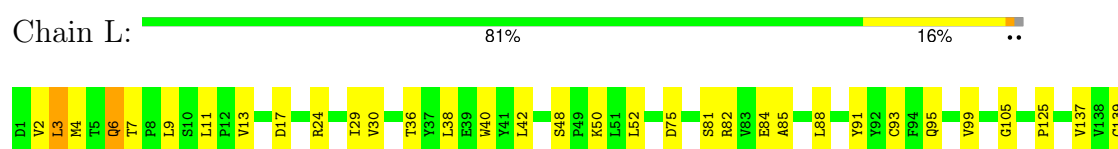
• Molecule 1: Leukocidin E



• Molecule 2: Antibody Fab Light Chain



• Molecule 2: Antibody Fab Light Chain

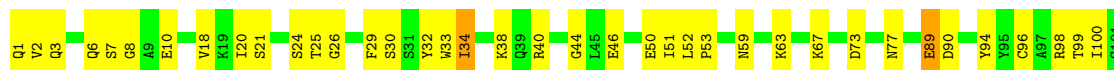




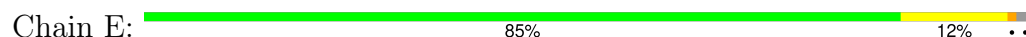
• Molecule 3: Antibody Fab Heavy Chain



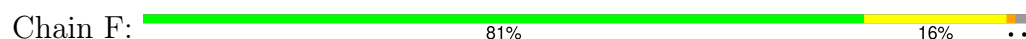
• Molecule 3: Antibody Fab Heavy Chain



• Molecule 4: Centyrin S26



• Molecule 4: Centyrin S26



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	344.26Å 54.77Å 115.73Å 90.00° 102.33° 90.00°	Depositor
Resolution (Å)	46.43 – 3.50 46.43 – 3.50	Depositor EDS
% Data completeness (in resolution range)	99.3 (46.43-3.50) 99.3 (46.43-3.50)	Depositor EDS
R_{merge}	0.24	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.62 (at 3.48Å)	Xtriage
Refinement program	PHENIX 1.8-3855	Depositor
R, R_{free}	0.229 , 0.279 0.232 , 0.280	Depositor DCC
R_{free} test set	1356 reflections (3.89%)	wwPDB-VP
Wilson B-factor (Å ²)	95.1	Xtriage
Anisotropy	0.365	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 59.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	12610	wwPDB-VP
Average B, all atoms (Å ²)	103.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.75% 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	A	0.19	0/2391	0.36	0/3233
1	B	0.17	0/2378	0.35	0/3215
2	C	0.13	0/1700	0.32	0/2308
2	L	0.13	0/1700	0.32	0/2308
3	D	0.16	0/1656	0.43	2/2256 (0.1%)
3	H	0.14	0/1657	0.35	0/2257
4	E	0.17	0/721	0.37	0/991
4	F	0.18	0/721	0.38	0/991
All	All	0.16	0/12924	0.36	2/17559 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	27	TYR	CA-CB-CG	-5.16	104.61	113.90
3	D	29	PHE	N-CA-C	-5.10	101.96	109.86

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2331	0	2252	43	0
1	B	2320	0	2245	43	1
2	C	1664	0	1621	27	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	L	1664	0	1621	17	0
3	D	1615	0	1562	35	1
3	H	1616	0	1565	31	0
4	E	700	0	656	8	0
4	F	700	0	656	13	0
All	All	12610	0	12178	194	1

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 8.

All (194) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:112:LEU:HD13	1:B:259:LEU:CD2	1.88	1.03
1:B:112:LEU:HD13	1:B:259:LEU:HD23	1.47	0.96
2:C:82:ARG:NH1	2:C:84:GLU:OE2	2.10	0.83
1:A:112:LEU:HD13	1:A:259:LEU:CD2	2.13	0.79
1:B:112:LEU:HD13	1:B:259:LEU:HD21	1.66	0.77
3:D:30:SER:HA	3:D:54:GLY:H	1.50	0.76
3:H:2:VAL:HG22	3:H:26:GLY:HA3	1.69	0.75
1:A:64:VAL:HB	1:A:73:ALA:HB3	1.71	0.71
2:L:82:ARG:NH1	2:L:84:GLU:OE2	2.24	0.70
1:B:72:ASP:HB2	1:B:247:HIS:HB3	1.73	0.69
3:H:121:ALA:O	3:H:124:SER:N	2.26	0.68
3:D:29:PHE:HD2	3:D:53:PRO:HG2	1.57	0.68
4:E:15:GLU:HG2	4:E:64:PRO:HB3	1.76	0.67
1:A:20:ALA:HB1	2:C:99:VAL:HG12	1.78	0.66
3:H:38:LYS:NZ	3:H:89:GLU:OE1	2.29	0.65
1:A:112:LEU:HD13	1:A:259:LEU:HD23	1.76	0.65
1:A:220:ARG:NH2	1:A:226:ASP:OD1	2.29	0.65
1:B:102:MET:HE2	1:B:104:TRP:HB3	1.77	0.65
3:H:34:ILE:HD11	3:H:96:CYS:HB2	1.79	0.64
3:D:177:ALA:HB2	3:D:187:LEU:HD23	1.80	0.63
1:A:38:GLY:HA3	1:A:67:LYS:HD3	1.81	0.63
2:C:48:SER:O	2:C:50:LYS:NZ	2.33	0.62
3:D:67:LYS:NZ	3:D:90:ASP:OD2	2.27	0.62
3:D:10:GLU:HG2	3:D:18:VAL:HG21	1.82	0.61
1:A:78:MET:SD	1:A:259:LEU:HD22	2.40	0.61
2:C:151:VAL:HG22	2:C:201:VAL:HG22	1.80	0.61
2:L:6:GLN:HB2	2:L:105:GLY:H	1.65	0.61
1:B:116:ASP:OD2	1:B:255:SER:OG	2.15	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:E:19:ARG:NE	4:E:58:ASP:OD1	2.34	0.61
4:F:37:ALA:HB2	4:F:45:ALA:HB2	1.83	0.60
1:B:220:ARG:NH2	1:B:226:ASP:OD1	2.29	0.60
1:A:112:LEU:HD13	1:A:259:LEU:HD21	1.84	0.60
2:L:48:SER:OG	2:L:50:LYS:NZ	2.32	0.59
3:H:6:GLN:NE2	3:H:94:TYR:O	2.25	0.59
3:H:202:THR:HG23	3:H:219:LYS:HE3	1.86	0.58
1:A:26:HIS:HD2	2:C:39:GLU:OE2	1.86	0.58
1:B:112:LEU:CD1	1:B:259:LEU:HD23	2.28	0.58
1:A:26:HIS:CE1	3:D:99:THR:HG21	2.38	0.58
3:H:136:SER:O	3:H:141:SER:N	2.37	0.57
1:A:72:ASP:HB2	1:A:247:HIS:HB3	1.86	0.57
1:A:247:HIS:HD1	1:A:300:TRP:NE1	2.02	0.57
2:C:7:THR:HG23	2:C:22:SER:HB2	1.86	0.57
1:A:24:HIS:HD2	3:D:33:TRP:NE1	2.02	0.57
1:B:121:LEU:HD23	1:B:245:LEU:HB3	1.87	0.57
1:A:129:LYS:HE3	1:A:173:VAL:HG11	1.87	0.56
3:D:3:GLN:HA	3:D:25:THR:HG23	1.87	0.56
3:H:132:PRO:HD3	3:H:218:LYS:HE2	1.87	0.55
3:H:29:PHE:HD2	3:H:53:PRO:HG2	1.71	0.55
1:A:257:PHE:HE2	1:A:298:VAL:HG13	1.72	0.55
4:F:80:ASN:N	4:F:80:ASN:OD1	2.40	0.54
3:H:10:GLU:HG2	3:H:18:VAL:HG21	1.88	0.54
3:D:132:PRO:HD3	3:D:218:LYS:HE2	1.90	0.54
1:A:95:GLY:C	4:F:80:ASN:HD21	2.16	0.54
1:B:64:VAL:HB	1:B:73:ALA:HB3	1.89	0.54
1:B:118:ASN:OD1	1:B:252:SER:OG	2.26	0.54
1:A:26:HIS:HB2	2:C:96:GLY:HA3	1.89	0.53
1:B:262:GLY:HA3	1:B:291:ASN:OD1	2.08	0.53
2:C:6:GLN:NE2	2:C:107:THR:OG1	2.41	0.53
4:F:32:PHE:HB2	4:F:50:VAL:HG22	1.91	0.53
3:H:7:SER:HB2	3:H:21:SER:H	1.74	0.53
3:D:32:TYR:CZ	3:D:100:ILE:HG23	2.44	0.53
1:B:24:HIS:HD2	3:H:33:TRP:CD1	2.28	0.52
3:D:202:THR:HG23	3:D:219:LYS:HE3	1.91	0.52
2:L:13:VAL:HG23	2:L:17:ASP:HB2	1.91	0.52
1:B:251:SER:HA	4:F:12:ARG:H	1.75	0.52
3:H:20:ILE:HG21	3:H:114:THR:HG21	1.91	0.52
1:B:42:ILE:HB	1:B:63:PHE:HB2	1.92	0.51
3:D:177:ALA:HA	3:D:187:LEU:HB3	1.92	0.51
1:B:131:GLU:HB2	1:B:173:VAL:HG22	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:87:SER:HB2	4:F:42:TRP:HE1	1.75	0.51
3:D:4:LEU:HD11	3:D:98:ARG:HB2	1.93	0.51
3:H:177:ALA:HB2	3:H:187:LEU:HD23	1.93	0.50
1:A:116:ASP:OD2	1:A:255:SER:OG	2.10	0.50
3:D:25:THR:OG1	3:D:26:GLY:N	2.43	0.49
3:H:67:LYS:NZ	3:H:90:ASP:OD2	2.45	0.49
2:L:30:VAL:HA	2:L:36:THR:HG22	1.93	0.49
3:H:147:LEU:HB2	3:H:220:VAL:HG11	1.93	0.49
2:C:99:VAL:HG11	3:D:59:ASN:ND2	2.28	0.49
3:D:28:LYS:HE2	3:D:28:LYS:HA	1.94	0.49
1:A:113:THR:HG22	1:A:183:SER:HB2	1.95	0.49
2:C:8:PRO:HG3	2:C:11:LEU:HD23	1.95	0.49
2:L:125:PRO:HD3	2:L:137:VAL:HG22	1.95	0.49
2:C:42:LEU:HB2	2:C:52:LEU:HD21	1.94	0.49
3:D:2:VAL:HG13	3:D:5:GLN:HB2	1.94	0.49
1:A:270:ALA:HB1	1:A:278:ILE:HD11	1.95	0.48
2:C:99:VAL:HG11	3:D:59:ASN:HD21	1.79	0.48
2:L:4:MET:HE3	2:L:4:MET:HB3	1.78	0.48
4:E:32:PHE:HB2	4:E:50:VAL:HG22	1.96	0.48
1:A:262:GLY:HA3	1:A:291:ASN:OD1	2.14	0.48
1:A:247:HIS:HD2	1:A:248:GLU:N	2.12	0.47
2:L:139:CYS:HB2	2:L:153:TRP:CH2	2.50	0.47
1:B:259:LEU:HD12	1:B:296:TYR:CE1	2.49	0.47
4:F:25:PRO:O	4:F:55:ARG:NH2	2.48	0.47
2:L:24:ARG:HD3	2:L:75:ASP:OD2	2.15	0.47
1:A:90:ASP:OD1	1:A:219:ALA:HB2	2.15	0.47
2:C:37:TYR:HB2	2:C:97:SER:HB3	1.97	0.47
1:B:24:HIS:CE1	3:H:50:GLU:OE2	2.68	0.46
1:B:138:THR:HG23	1:B:164:THR:HG22	1.96	0.46
1:A:197:ASP:OD1	1:A:197:ASP:N	2.48	0.46
2:L:29:ILE:HD13	2:L:38:LEU:HB2	1.96	0.46
3:D:29:PHE:HE1	3:D:77:ASN:CG	2.23	0.46
3:H:168:LEU:HD21	3:H:191:VAL:HG21	1.98	0.46
1:B:77:LYS:HE2	1:B:79:GLN:CD	2.41	0.46
2:C:126:SER:OG	3:D:131:PHE:HB3	2.16	0.46
4:E:24:ALA:HB3	4:E:55:ARG:CZ	2.46	0.46
1:A:94:SER:HA	1:A:97:GLU:OE2	2.16	0.46
1:B:205:ARG:HH21	1:B:281:GLU:CB	2.28	0.45
2:C:153:TRP:CE2	2:C:184:LEU:HB2	2.52	0.45
1:A:167:TYR:HD2	1:A:240:SER:HB3	1.80	0.45
1:B:78:MET:SD	1:B:259:LEU:HD22	2.56	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:13:VAL:HG23	2:C:17:ASP:HB2	1.98	0.45
3:D:61:ASN:OD1	3:D:62:GLU:N	2.50	0.45
1:A:205:ARG:HH21	1:A:281:GLU:CB	2.29	0.45
1:A:254:THR:HB	1:A:297:LYS:HE3	1.99	0.45
3:D:135:PRO:HD2	3:D:222:PRO:HA	1.99	0.45
1:A:205:ARG:HH21	1:A:281:GLU:HB3	1.82	0.45
1:B:173:VAL:HG23	1:B:191:ASN:HB2	1.99	0.45
1:A:283:LYS:NZ	4:F:85:SER:H	2.15	0.45
1:B:274:PRO:C	1:B:276:THR:H	2.25	0.45
3:D:55:SER:OG	3:D:57:SER:OG	2.32	0.44
3:H:24:SER:O	3:H:77:ASN:HB3	2.17	0.44
3:D:147:LEU:HB2	3:D:220:VAL:HG11	1.99	0.44
3:H:100:ILE:HB	3:H:107:PHE:HD1	1.82	0.44
3:H:102:THR:OG1	3:H:105:ASP:O	2.35	0.44
1:B:116:ASP:OD2	1:B:119:VAL:HG23	2.18	0.44
3:D:29:PHE:O	3:D:30:SER:OG	2.26	0.44
2:L:99:VAL:HG11	3:H:59:ASN:HD21	1.82	0.44
2:C:94:PHE:CE1	2:C:101:TYR:HB3	2.52	0.44
1:B:113:THR:HG22	1:B:183:SER:HB2	1.99	0.44
2:L:42:LEU:HD13	2:L:91:TYR:CZ	2.53	0.44
2:L:166:GLU:HB2	2:L:180:LEU:HD21	1.99	0.44
2:C:59:ARG:NH1	2:C:65:ASP:HA	2.33	0.44
2:C:38:LEU:HD22	2:C:76:PHE:CG	2.52	0.43
1:A:283:LYS:HZ3	4:F:85:SER:H	1.66	0.43
3:D:45:LEU:HD23	3:D:45:LEU:HA	1.88	0.43
1:B:283:LYS:NZ	4:E:85:SER:H	2.16	0.43
1:B:214:GLY:HA3	1:B:215:PRO:HA	1.77	0.43
1:B:257:PHE:HE2	1:B:298:VAL:HG13	1.84	0.43
3:H:46:GLU:OE1	3:H:63:LYS:NZ	2.45	0.43
1:A:100:LYS:HG3	1:A:222:TYR:CD1	2.54	0.43
4:F:25:PRO:HB2	4:F:28:ALA:HB2	2.00	0.43
1:A:130:ILE:HD13	1:A:135:VAL:HG22	2.01	0.43
1:A:65:LYS:HD3	1:A:303:HIS:NE2	2.34	0.43
1:B:306:LYS:HE3	1:B:308:LYS:HE3	2.00	0.43
1:B:250:GLY:C	1:B:251:SER:HG	2.26	0.42
2:C:66:ARG:HD2	2:C:82:ARG:O	2.19	0.42
1:A:281:GLU:OE1	4:F:82:THR:OG1	2.25	0.42
1:B:35:ILE:HG21	1:B:161:TYR:HB2	2.01	0.42
1:B:269:TYR:CZ	4:E:83:PRO:HG2	2.54	0.42
3:H:8:GLY:O	3:H:114:THR:HG23	2.19	0.42
1:B:200:LYS:HD3	1:B:204:ASP:OD2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:40:TRP:CZ3	2:L:93:CYS:HB3	2.54	0.42
3:H:106:TRP:CD1	3:H:106:TRP:C	2.97	0.42
1:B:26:HIS:CE1	3:H:99:THR:HG21	2.55	0.42
3:H:29:PHE:CD2	3:H:53:PRO:HG2	2.54	0.42
1:A:220:ARG:HD2	1:A:220:ARG:HA	1.80	0.42
1:B:69:TYR:CE1	1:B:71:LYS:HB2	2.55	0.42
1:A:138:THR:HG23	1:A:164:THR:HG22	2.01	0.42
1:B:62:ASP:OD2	1:B:161:TYR:OH	2.38	0.42
3:H:24:SER:OG	3:H:98:ARG:NH1	2.53	0.42
1:B:78:MET:HE1	1:B:186:TRP:CZ3	2.55	0.42
1:B:197:ASP:OD1	1:B:197:ASP:N	2.52	0.42
3:D:152:LYS:HG2	3:D:153:ASP:CG	2.45	0.42
1:B:205:ARG:HH21	1:B:281:GLU:HB3	1.85	0.41
1:B:283:LYS:HZ3	4:E:85:SER:H	1.67	0.41
4:F:32:PHE:O	4:F:49:THR:HA	2.19	0.41
1:A:102:MET:HE2	1:A:104:TRP:HB3	2.02	0.41
2:C:101:TYR:HB2	3:D:47:TRP:CE2	2.55	0.41
3:D:60:HIS:HB3	3:D:70:PHE:HE2	1.85	0.41
3:D:13:ASN:HA	3:D:119:SER:O	2.20	0.41
2:L:3:LEU:HD13	2:L:3:LEU:HA	1.92	0.41
1:A:274:PRO:C	1:A:276:THR:H	2.28	0.41
3:D:18:VAL:HG12	3:D:86:LEU:HD11	2.02	0.41
3:H:32:TYR:CZ	3:H:100:ILE:HG23	2.56	0.41
2:C:153:TRP:HB2	2:C:160:GLN:HB2	2.02	0.41
2:C:212:LYS:HE3	2:C:212:LYS:HB3	1.91	0.41
3:D:52:LEU:HD12	3:D:55:SER:HB3	2.02	0.41
3:H:40:ARG:NH1	3:H:44:GLY:HA3	2.36	0.41
3:H:157:GLU:OE2	3:H:177:ALA:HB3	2.20	0.41
1:A:266:ASP:OD2	1:A:284:HIS:CD2	2.74	0.41
2:C:94:PHE:CZ	2:C:101:TYR:HB3	2.55	0.41
2:C:141:LEU:HB2	2:C:180:LEU:HB3	2.03	0.41
3:D:125:THR:HG22	3:D:212:SER:HB3	2.03	0.41
2:L:144:PHE:N	2:L:177:THR:HB	2.36	0.41
2:C:39:GLU:HG2	2:C:54:TYR:HA	2.02	0.40
4:E:32:PHE:O	4:E:49:THR:HA	2.21	0.40
4:F:32:PHE:CE1	4:F:52:GLY:HA2	2.56	0.40
1:A:212:PRO:HG3	1:A:222:TYR:CZ	2.57	0.40
1:A:298:VAL:HG12	1:A:305:ILE:CD1	2.51	0.40
3:D:157:GLU:OE2	3:D:177:ALA:HB3	2.21	0.40
1:B:78:MET:HE3	1:B:78:MET:HB2	2.00	0.40
1:B:274:PRO:O	1:B:276:THR:N	2.54	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:101:TYR:HB2	3:D:47:TRP:CD2	2.55	0.40
2:L:85:ALA:O	2:L:88:LEU:HD22	2.21	0.40

All (1) 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:B:272:LEU:O	3:D:27:TYR:OH[1_554]	2.15	0.05

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	A	291/294 (99%)	266 (91%)	25 (9%)	0	100	100
1	B	290/294 (99%)	266 (92%)	24 (8%)	0	100	100
2	C	215/219 (98%)	205 (95%)	9 (4%)	1 (0%)	24	57
2	L	215/219 (98%)	205 (95%)	7 (3%)	3 (1%)	9	38
3	D	210/229 (92%)	190 (90%)	17 (8%)	3 (1%)	9	38
3	H	210/229 (92%)	191 (91%)	19 (9%)	0	100	100
4	E	94/98 (96%)	85 (90%)	9 (10%)	0	100	100
4	F	94/98 (96%)	85 (90%)	9 (10%)	0	100	100
All	All	1619/1680 (96%)	1493 (92%)	119 (7%)	7 (0%)	30	62

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	L	81	SER
2	L	143	ASN
2	C	143	ASN

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Mol	Chain	Res	Type
3	D	102	THR
2	L	2	VAL
3	D	3	GLN
3	D	27	TYR

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	254/259 (98%)	247 (97%)	7 (3%)	38	61
1	B	254/259 (98%)	245 (96%)	9 (4%)	32	57
2	C	190/194 (98%)	184 (97%)	6 (3%)	34	59
2	L	190/194 (98%)	182 (96%)	8 (4%)	26	52
3	D	178/193 (92%)	167 (94%)	11 (6%)	16	43
3	H	179/193 (93%)	168 (94%)	11 (6%)	17	43
4	E	70/80 (88%)	67 (96%)	3 (4%)	26	51
4	F	70/80 (88%)	66 (94%)	4 (6%)	18	45
All	All	1385/1452 (95%)	1326 (96%)	59 (4%)	26	51

All (59) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	HIS
1	A	31	ASN
1	A	67	LYS
1	A	195	THR
1	A	245	LEU
1	A	278	ILE
1	A	297	LYS
1	B	24	HIS
1	B	67	LYS
1	B	85	ARG
1	B	195	THR

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Mol	Chain	Res	Type
1	B	245	LEU
1	B	249	LYS
1	B	252	SER
1	B	278	ILE
1	B	297	LYS
2	C	2	VAL
2	C	6	GLN
2	C	7	THR
2	C	9	LEU
2	C	52	LEU
2	C	95	GLN
3	D	1	GLN
3	D	2	VAL
3	D	5	GLN
3	D	12	MET
3	D	25	THR
3	D	28	LYS
3	D	34	ILE
3	D	55	SER
3	D	77	ASN
3	D	89	GLU
3	D	98	ARG
4	E	15	GLU
4	E	72	ILE
4	E	87	ILE
4	F	46	ILE
4	F	72	ILE
4	F	80	ASN
4	F	87	ILE
2	L	3	LEU
2	L	6	GLN
2	L	7	THR
2	L	9	LEU
2	L	11	LEU
2	L	52	LEU
2	L	95	GLN
2	L	144	PHE
3	H	1	GLN
3	H	3	GLN
3	H	25	THR
3	H	30	SER
3	H	34	ILE

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Mol	Chain	Res	Type
3	H	51	ILE
3	H	52	LEU
3	H	73	ASP
3	H	89	GLU
3	H	106	TRP
3	H	107	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	24	HIS
1	A	26	HIS
1	A	118	ASN
1	B	24	HIS
1	B	58	ASN
2	C	6	GLN
2	C	27	GLN
2	C	58	ASN
2	C	95	GLN
2	C	143	ASN
3	D	3	GLN
3	D	180	GLN
2	L	129	GLN
2	L	143	ASN
2	L	204	GLN
3	H	3	GLN
3	H	164	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 [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
3	D	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	H	121:ALA	C	124:SER	N	3.27
1	D	121:ALA	C	124:SER	N	3.15

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	293/294 (99%)	-0.27	0 100 100	57, 76, 105, 138	0
1	B	292/294 (99%)	-0.24	0 100 100	60, 86, 123, 155	0
2	C	217/219 (99%)	-0.00	0 100 100	74, 122, 157, 170	0
2	L	217/219 (99%)	-0.14	0 100 100	99, 120, 140, 156	0
3	D	216/229 (94%)	-0.04	0 100 100	70, 108, 168, 180	0
3	H	216/229 (94%)	-0.02	0 100 100	96, 125, 152, 167	0
4	E	96/98 (97%)	-0.25	0 100 100	69, 86, 118, 150	0
4	F	96/98 (97%)	-0.26	0 100 100	64, 81, 101, 150	0
All	All	1643/1680 (97%)	-0.14	0 100 100	57, 104, 152, 180	0

There are no RSRZ outliers to report.

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.