



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 04:37 AM UTC

PDB ID : 2GA0 / pdb_00002ga0
Title : Variable Small Protein 1 of Borrelia turicatae (VspA or Vsp1)
Authors : Lawson, C.L.; Yung, B.H.; Barbour, A.G.; Zuckert, W.R.
Deposited on : 2006-03-07
Resolution : 2.70 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

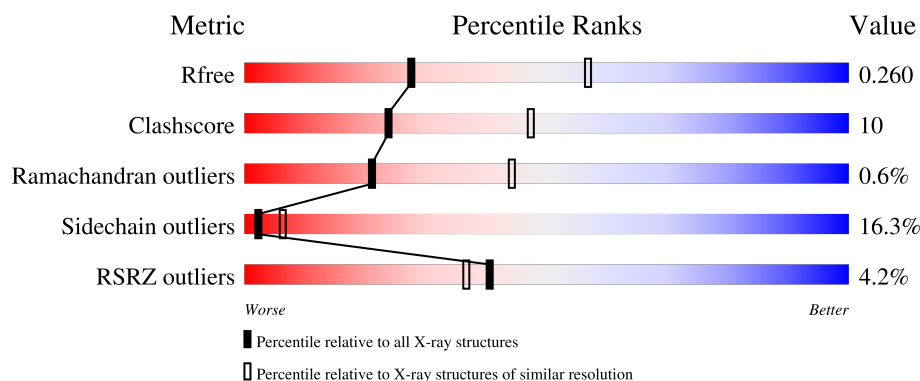
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. 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	181	
1	B	181	
1	C	181	
1	D	181	
1	E	181	

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Mol	Chain	Length	Quality of chain
1	F	181	
1	G	181	
1	H	181	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 9606 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 surface protein VspA.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	A	164	Total	C	N	O	0	0	0
			1182	740	196	246			
1	B	164	Total	C	N	O	0	0	0
			1182	740	196	246			
1	C	164	Total	C	N	O	0	0	0
			1182	740	196	246			
1	D	164	Total	C	N	O	0	0	0
			1182	740	196	246			
1	E	164	Total	C	N	O	0	0	0
			1182	740	196	246			
1	F	164	Total	C	N	O	0	0	0
			1182	740	196	246			
1	G	164	Total	C	N	O	0	0	0
			1182	740	196	246			
1	H	164	Total	C	N	O	0	0	0
			1182	740	196	246			

- Molecule 2 is NICKEL (II) ION (CCD ID: NI) (formula: Ni).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Ni	0	0
			2	2		
2	B	2	Total	Ni	0	0
			2	2		
2	C	1	Total	Ni	0	0
			1	1		
2	D	2	Total	Ni	0	0
			2	2		
2	E	2	Total	Ni	0	0
			2	2		
2	F	1	Total	Ni	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	G	2	Total 2	Ni 2	0	0
2	H	2	Total 2	Ni 2	0	0

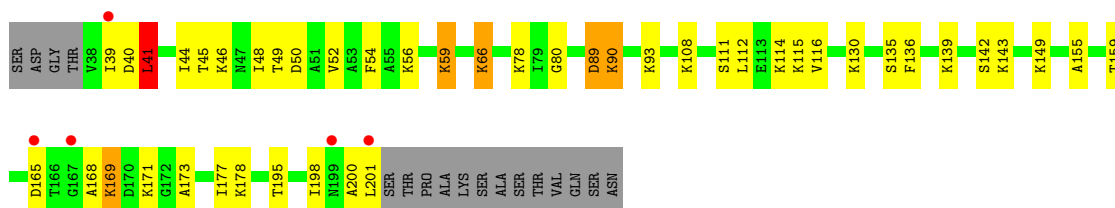
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	20	Total 20	O 20	0	0
3	B	16	Total 16	O 16	0	0
3	C	11	Total 11	O 11	0	0
3	D	13	Total 13	O 13	0	0
3	E	16	Total 16	O 16	0	0
3	F	24	Total 24	O 24	0	0
3	G	18	Total 18	O 18	0	0
3	H	18	Total 18	O 18	0	0

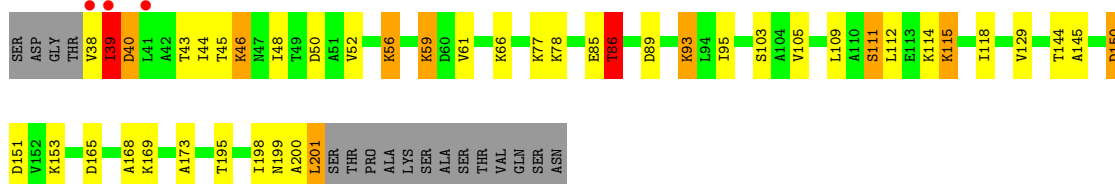
3 Residue-property plots [i](#)

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.

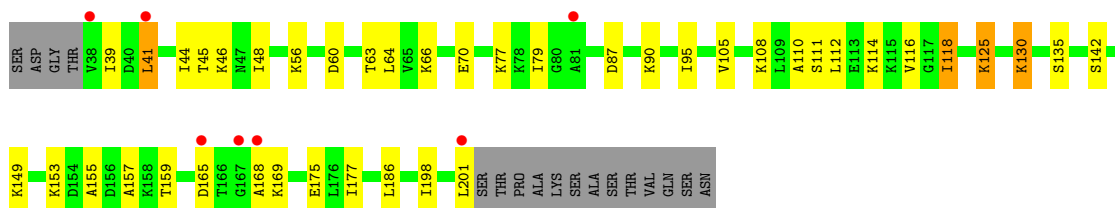
- Molecule 1: surface protein VspA



- Molecule 1: surface protein VspA

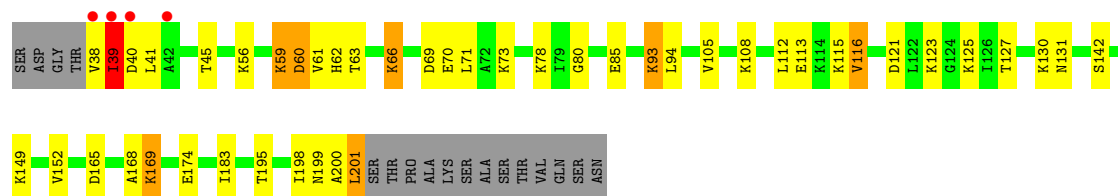


- Molecule 1: surface protein VspA

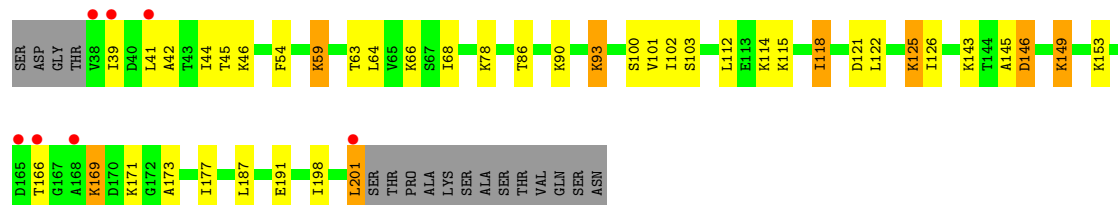


- Molecule 1: surface protein VspA

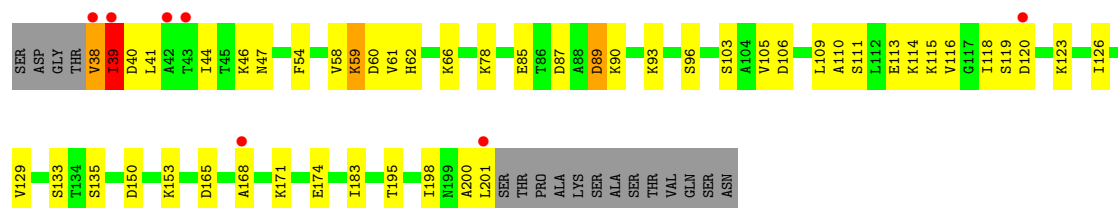




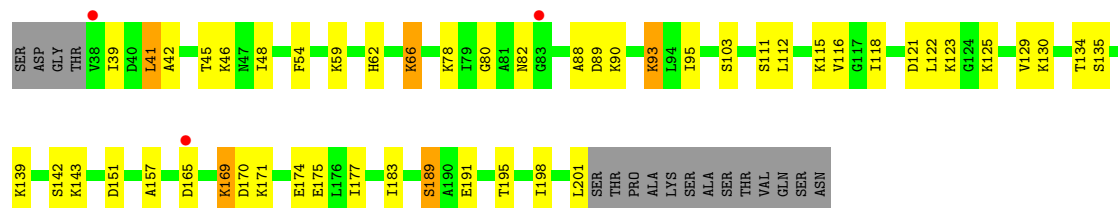
- Molecule 1: surface protein VspA



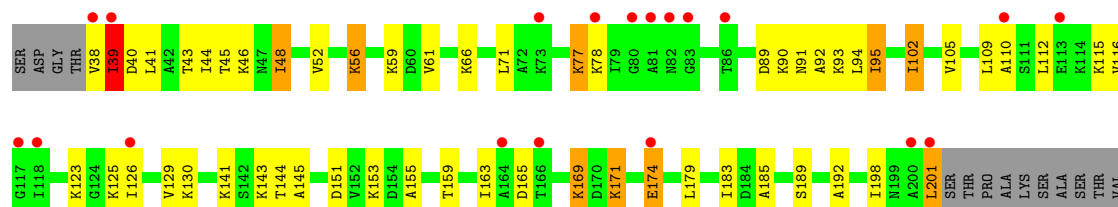
- Molecule 1: surface protein VspA



- Molecule 1: surface protein VspA



- Molecule 1: surface protein VspA



GLN
SER
ASN

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	241.84Å 69.11Å 87.58Å 90.00° 104.95° 90.00°	Depositor
Resolution (Å)	50.00 – 2.70 50.00 – 2.71	Depositor EDS
% Data completeness (in resolution range)	98.7 (50.00-2.70) 98.7 (50.00-2.71)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.08 (at 2.69Å)	Xtrriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.216 , 0.272 0.213 , 0.260	Depositor DCC
R_{free} test set	1902 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å ²)	39.9	Xtrriage
Anisotropy	0.428	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 36.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	9606	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 32.90 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 8.7245e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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:
NI

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	0.97	0/1185	1.15	0/1598
1	B	0.89	1/1185 (0.1%)	1.14	2/1598 (0.1%)
1	C	0.86	1/1185 (0.1%)	1.10	0/1598
1	D	0.85	1/1185 (0.1%)	1.12	0/1598
1	E	0.83	1/1185 (0.1%)	1.07	0/1598
1	F	0.89	0/1185	1.19	1/1598 (0.1%)
1	G	0.89	1/1185 (0.1%)	1.12	1/1598 (0.1%)
1	H	0.80	0/1185	1.13	0/1598
All	All	0.87	5/9480 (0.1%)	1.13	4/12784 (0.0%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	101	VAL	CA-CB	5.62	1.60	1.54
1	B	40	ASP	CA-C	5.46	1.60	1.52
1	D	116	VAL	CA-CB	5.40	1.61	1.54
1	G	129	VAL	CA-CB	5.32	1.60	1.54
1	C	105	VAL	CA-CB	5.12	1.60	1.54

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	88	ALA	N-CA-C	5.63	117.55	110.24
1	F	47	ASN	N-CA-C	5.29	117.46	111.11
1	B	86	THR	OG1-CB-CG2	-5.21	98.88	109.30
1	B	150	ASP	N-CA-C	5.17	118.36	111.75

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1182	0	1257	41	0
1	B	1182	0	1257	35	0
1	C	1182	0	1257	21	0
1	D	1182	0	1257	32	0
1	E	1182	0	1257	23	0
1	F	1182	0	1257	25	0
1	G	1182	0	1257	27	0
1	H	1182	0	1257	30	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	1	0	0	0	0
2	D	2	0	0	0	0
2	E	2	0	0	0	0
2	F	1	0	0	0	0
2	G	2	0	0	0	0
2	H	2	0	0	0	0
3	A	20	0	0	4	0
3	B	16	0	0	4	0
3	C	11	0	0	0	0
3	D	13	0	0	2	0
3	E	16	0	0	2	0
3	F	24	0	0	2	0
3	G	18	0	0	2	0
3	H	18	0	0	4	0
All	All	9606	0	10056	198	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 10.

All (198) 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:A:165:ASP:HB3	1:A:169:LYS:NZ	1.58	1.18
1:H:112:LEU:HA	3:H:224:HOH:O	1.67	0.94
1:C:56:LYS:HE3	1:D:59:LYS:HZ1	1.32	0.93

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:165:ASP:HB3	1:A:169:LYS:HZ1	1.29	0.93
1:E:59:LYS:HE3	3:F:238:HOH:O	1.67	0.92
3:E:227:HOH:O	1:F:59:LYS:HG3	1.68	0.91
1:A:143:LYS:HD2	1:A:171:LYS:HG2	1.52	0.89
1:G:42:ALA:O	1:G:45:THR:HG22	1.74	0.86
1:A:165:ASP:HB3	1:A:169:LYS:HZ3	1.39	0.86
1:C:56:LYS:HE3	1:D:59:LYS:NZ	1.90	0.85
1:H:198:ILE:O	1:H:201:LEU:HB2	1.76	0.85
1:G:165:ASP:O	1:G:169:LYS:HE3	1.76	0.85
1:H:112:LEU:HD23	3:H:224:HOH:O	1.74	0.85
1:A:59:LYS:HG3	3:B:228:HOH:O	1.77	0.83
1:B:198:ILE:O	1:B:201:LEU:HB2	1.78	0.83
1:G:198:ILE:O	1:G:201:LEU:HB2	1.80	0.81
1:E:198:ILE:O	1:E:201:LEU:HB2	1.79	0.80
1:A:46:LYS:HE2	1:A:50:ASP:OD1	1.82	0.78
1:F:38:VAL:O	1:F:39:ILE:HG13	1.84	0.78
1:F:165:ASP:OD1	1:F:168:ALA:O	2.02	0.77
1:E:42:ALA:HA	1:E:45:THR:HG22	1.67	0.76
1:A:165:ASP:CB	1:A:169:LYS:NZ	2.47	0.76
1:A:66:LYS:HD2	3:A:217:HOH:O	1.86	0.75
1:A:149:LYS:HD2	3:A:231:HOH:O	1.88	0.74
1:D:169:LYS:NZ	1:D:169:LYS:HB3	2.04	0.72
1:A:59:LYS:HE2	3:B:228:HOH:O	1.89	0.72
1:D:198:ILE:O	1:D:201:LEU:HB2	1.89	0.72
1:A:59:LYS:NZ	1:B:56:LYS:HZ2	1.89	0.71
1:A:165:ASP:O	1:A:169:LYS:HE2	1.91	0.70
1:H:185:ALA:HA	3:H:215:HOH:O	1.91	0.70
1:A:90:LYS:HE3	1:B:150:ASP:OD1	1.92	0.70
1:G:115:LYS:HD3	1:G:118:ILE:HD11	1.74	0.70
1:D:38:VAL:O	1:D:39:ILE:HG13	1.91	0.69
1:B:38:VAL:C	1:B:39:ILE:HG13	2.17	0.68
1:D:199:ASN:HA	3:D:224:HOH:O	1.91	0.68
1:A:165:ASP:CB	1:A:169:LYS:HZ3	2.05	0.68
1:E:54:PHE:CE1	1:E:118:ILE:HD11	2.29	0.68
1:H:38:VAL:C	1:H:39:ILE:HG13	2.18	0.68
1:B:86:THR:HG22	3:B:223:HOH:O	1.94	0.67
1:F:38:VAL:C	1:F:39:ILE:HG13	2.20	0.66
1:A:59:LYS:NZ	1:B:56:LYS:NZ	2.42	0.66
1:G:80:GLY:HA2	1:H:144:THR:OG1	1.95	0.66
1:F:198:ILE:O	1:F:201:LEU:HB2	1.96	0.65
1:H:192:ALA:HA	3:H:221:HOH:O	1.95	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:66:LYS:HD2	3:G:222:HOH:O	1.95	0.64
1:A:168:ALA:HA	1:C:111:SER:HB2	1.79	0.64
1:F:110:ALA:O	1:F:113:GLU:HG2	1.97	0.64
1:A:165:ASP:O	1:A:169:LYS:CE	2.45	0.64
1:A:198:ILE:O	1:A:201:LEU:HB2	1.98	0.63
1:D:169:LYS:HB3	1:D:169:LYS:HZ1	1.64	0.63
1:G:201:LEU:HD13	1:H:45:THR:OG1	1.98	0.62
1:H:143:LYS:HG3	1:H:171:LYS:HG2	1.80	0.62
1:D:165:ASP:OD1	1:D:168:ALA:O	2.18	0.61
1:D:70:GLU:OE2	1:D:73:LYS:HE3	2.01	0.60
1:A:59:LYS:HZ3	1:B:56:LYS:HZ2	1.47	0.60
1:D:59:LYS:HE3	3:D:215:HOH:O	2.02	0.59
1:E:93:LYS:NZ	1:F:87:ASP:O	2.33	0.59
1:A:59:LYS:HZ3	1:B:56:LYS:NZ	2.01	0.58
1:D:66:LYS:HE3	1:D:69:ASP:OD2	2.03	0.58
1:C:56:LYS:CE	1:D:59:LYS:HZ1	2.12	0.57
1:E:201:LEU:HD21	1:F:201:LEU:HD11	1.86	0.57
1:F:115:LYS:HD3	1:F:118:ILE:HD11	1.86	0.57
1:C:110:ALA:HB2	1:C:130:LYS:HE3	1.87	0.56
1:F:61:VAL:HG11	1:F:109:LEU:HD21	1.87	0.56
1:A:56:LYS:NZ	1:B:59:LYS:HZ3	2.04	0.56
1:G:93:LYS:HE3	1:H:89:ASP:O	2.05	0.56
1:E:42:ALA:HA	1:E:45:THR:CG2	2.35	0.56
1:D:38:VAL:C	1:D:39:ILE:HG13	2.31	0.56
1:H:89:ASP:HB3	1:H:151:ASP:OD2	2.06	0.56
1:H:71:LEU:HB3	1:H:94:LEU:HD11	1.89	0.55
1:H:109:LEU:HD11	1:H:129:VAL:HG12	1.89	0.55
1:A:56:LYS:NZ	1:B:59:LYS:NZ	2.55	0.55
1:E:145:ALA:O	1:E:149:LYS:HG2	2.06	0.55
1:B:89:ASP:HB3	1:B:151:ASP:OD2	2.07	0.54
1:E:121:ASP:OD2	1:E:125:LYS:NZ	2.41	0.54
1:G:165:ASP:HB3	1:G:169:LYS:NZ	2.22	0.54
1:B:169:LYS:HB3	1:B:173:ALA:HB3	1.90	0.54
1:F:66:LYS:HG3	3:F:237:HOH:O	2.08	0.54
1:A:198:ILE:HG22	1:B:45:THR:HG23	1.89	0.53
1:F:123:LYS:HA	1:F:126:ILE:HD12	1.91	0.53
1:E:173:ALA:O	1:E:177:ILE:HG13	2.08	0.53
1:G:95:ILE:HD11	1:G:157:ALA:HA	1.91	0.53
1:E:93:LYS:HD3	1:F:89:ASP:O	2.09	0.52
1:D:121:ASP:OD2	1:D:125:LYS:NZ	2.41	0.52
1:A:56:LYS:HZ1	1:B:59:LYS:NZ	2.08	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:165:ASP:OD1	1:B:168:ALA:O	2.29	0.51
1:D:62:HIS:HA	1:D:183:ILE:HG12	1.92	0.51
1:B:46:LYS:HD2	1:B:50:ASP:OD1	2.10	0.50
1:C:70:GLU:OE2	1:C:70:GLU:HA	2.11	0.50
1:E:187:LEU:O	1:E:191:GLU:HG3	2.12	0.50
1:A:80:GLY:HA2	1:B:144:THR:OG1	2.11	0.50
1:H:169:LYS:HB2	1:H:174:GLU:HG2	1.92	0.50
1:D:80:GLY:HA3	1:D:85:GLU:OE2	2.11	0.50
1:C:198:ILE:O	1:C:201:LEU:HB2	2.12	0.49
1:B:115:LYS:HB3	1:B:118:ILE:HD11	1.94	0.49
1:B:48:ILE:O	1:B:52:VAL:HG23	2.12	0.49
1:C:198:ILE:HG22	1:D:45:THR:HG23	1.93	0.49
1:H:110:ALA:HB2	1:H:130:LYS:HE2	1.94	0.49
1:B:61:VAL:HG13	1:B:105:VAL:HG13	1.95	0.49
1:C:56:LYS:CE	1:D:59:LYS:NZ	2.69	0.49
1:D:113:GLU:OE1	1:D:130:LYS:HG2	2.13	0.49
1:E:68:ILE:HG13	1:E:102:ILE:HD11	1.95	0.49
1:E:143:LYS:HG3	1:E:171:LYS:HG2	1.94	0.48
1:H:179:LEU:O	1:H:183:ILE:HG22	2.12	0.48
1:A:90:LYS:CE	1:B:150:ASP:OD1	2.59	0.48
1:A:59:LYS:HZ1	1:B:56:LYS:NZ	2.08	0.48
1:D:71:LEU:HB3	1:D:94:LEU:HD11	1.96	0.48
1:A:56:LYS:HZ1	1:B:59:LYS:HZ2	1.62	0.48
1:C:87:ASP:O	1:D:93:LYS:NZ	2.40	0.47
1:F:62:HIS:HD2	1:F:183:ILE:HG23	1.78	0.47
1:D:40:ASP:OD1	1:D:40:ASP:C	2.57	0.47
1:H:61:VAL:HG13	1:H:105:VAL:HG13	1.94	0.47
1:A:89:ASP:O	1:B:93:LYS:HD3	2.14	0.47
1:G:122:LEU:HD21	1:G:189:SER:HB3	1.97	0.47
1:A:56:LYS:HZ2	1:B:59:LYS:HZ3	1.61	0.47
1:A:78:LYS:NZ	3:A:221:HOH:O	2.48	0.47
1:D:169:LYS:NZ	1:D:169:LYS:CB	2.76	0.47
1:C:41:LEU:HD11	1:D:201:LEU:HD22	1.97	0.47
1:C:118:ILE:HD12	1:C:118:ILE:HA	1.54	0.46
1:E:63:THR:OG1	1:F:60:ASP:OD2	2.32	0.46
1:H:155:ALA:O	1:H:159:THR:HG23	2.15	0.46
3:E:230:HOH:O	1:F:66:LYS:HE3	2.15	0.46
1:F:93:LYS:O	1:F:96:SER:HB2	2.15	0.46
1:F:61:VAL:HG13	1:F:105:VAL:HG13	1.97	0.46
1:B:46:LYS:HB3	1:B:46:LYS:HE3	1.50	0.45
1:D:39:ILE:HG22	1:D:40:ASP:H	1.80	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:82:ASN:C	1:H:141:LYS:HD3	2.41	0.45
1:A:200:ALA:O	1:A:201:LEU:C	2.60	0.45
1:B:38:VAL:O	1:B:39:ILE:HG13	2.16	0.45
1:F:38:VAL:C	1:F:39:ILE:CG1	2.88	0.45
1:E:146:ASP:OD1	1:E:146:ASP:N	2.50	0.45
1:A:165:ASP:CB	1:A:169:LYS:HZ1	2.15	0.45
1:A:169:LYS:HB3	1:A:173:ALA:HB3	1.99	0.45
1:C:175:GLU:OE1	1:C:175:GLU:N	2.50	0.44
1:D:200:ALA:O	1:D:201:LEU:C	2.60	0.44
1:E:169:LYS:HB3	1:E:173:ALA:HB3	1.99	0.44
1:G:54:PHE:HE1	1:G:112:LEU:HD22	1.82	0.44
1:G:121:ASP:OD2	1:G:125:LYS:NZ	2.43	0.44
1:C:165:ASP:OD1	1:C:168:ALA:O	2.35	0.44
1:B:40:ASP:O	1:B:44:ILE:HD12	2.17	0.44
1:G:59:LYS:HG3	3:G:218:HOH:O	2.17	0.44
1:B:86:THR:CG2	3:B:223:HOH:O	2.62	0.44
1:C:64:LEU:HD23	1:C:64:LEU:HA	1.69	0.44
1:G:198:ILE:HG22	1:H:45:THR:HG23	1.99	0.44
1:C:125:LYS:HB3	1:C:186:LEU:HD13	2.00	0.44
1:G:143:LYS:HD2	1:G:171:LYS:HG2	1.99	0.44
1:B:56:LYS:HE3	1:B:56:LYS:HB3	1.40	0.43
1:C:112:LEU:HD23	1:C:112:LEU:HA	1.88	0.43
1:G:93:LYS:HD2	1:H:77:LYS:O	2.18	0.43
1:A:66:LYS:HE2	3:A:234:HOH:O	2.18	0.43
1:B:111:SER:O	1:B:114:LYS:HB2	2.19	0.43
1:C:155:ALA:O	1:C:159:THR:HG23	2.19	0.43
1:H:40:ASP:OD1	1:H:40:ASP:C	2.61	0.43
1:H:125:LYS:HE2	1:H:189:SER:OG	2.18	0.43
1:E:115:LYS:HB3	1:E:118:ILE:HD12	2.01	0.43
1:G:59:LYS:HE2	1:H:56:LYS:HE3	2.00	0.43
1:C:60:ASP:OD1	1:C:108:LYS:NZ	2.49	0.42
1:F:54:PHE:O	1:F:58:VAL:HG23	2.19	0.42
1:D:60:ASP:OD1	1:D:108:LYS:HD3	2.19	0.42
1:A:48:ILE:O	1:A:52:VAL:HG23	2.19	0.42
1:G:139:LYS:NZ	1:G:175:GLU:OE2	2.52	0.42
1:A:54:PHE:HE1	1:A:112:LEU:HD22	1.85	0.42
1:F:200:ALA:O	1:F:201:LEU:C	2.62	0.42
1:D:149:LYS:HG3	1:D:152:VAL:HG22	2.02	0.42
1:E:122:LEU:O	1:E:126:ILE:HG13	2.20	0.42
1:F:40:ASP:O	1:F:44:ILE:HD12	2.19	0.42
1:H:183:ILE:O	1:H:183:ILE:HG13	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:63:THR:HB	1:D:63:THR:HG21	2.01	0.42
1:E:64:LEU:HD23	1:E:64:LEU:HA	1.81	0.42
1:B:59:LYS:HA	1:B:59:LYS:HD2	1.77	0.42
1:F:111:SER:O	1:F:114:LYS:HB2	2.18	0.42
1:H:91:ASN:O	1:H:95:ILE:HG12	2.20	0.42
1:E:42:ALA:CA	1:E:45:THR:HG22	2.44	0.41
1:H:144:THR:HG23	1:H:145:ALA:N	2.35	0.41
1:B:200:ALA:O	1:B:201:LEU:C	2.63	0.41
1:B:109:LEU:CD1	1:B:129:VAL:HG12	2.51	0.41
1:E:54:PHE:CE2	1:E:122:LEU:HD13	2.54	0.41
1:E:100:SER:O	1:E:103:SER:HB2	2.20	0.41
1:A:173:ALA:O	1:A:177:ILE:HG13	2.20	0.41
1:G:165:ASP:HB3	1:G:169:LYS:HZ1	1.85	0.41
1:G:198:ILE:O	1:G:201:LEU:CB	2.61	0.41
1:H:91:ASN:O	1:H:92:ALA:C	2.63	0.41
1:H:102:ILE:HD12	1:H:102:ILE:HA	1.93	0.41
1:G:174:GLU:HA	1:G:177:ILE:HD12	2.03	0.41
1:D:61:VAL:HG13	1:D:105:VAL:HG13	2.03	0.41
1:D:127:THR:HG22	1:D:131:ASN:ND2	2.36	0.41
1:A:40:ASP:O	1:A:41:LEU:C	2.64	0.40
1:B:144:THR:HG23	1:B:145:ALA:N	2.36	0.40
1:A:136:PHE:CD2	1:A:136:PHE:C	3.00	0.40
1:D:38:VAL:O	1:D:38:VAL:HG23	2.22	0.40
1:F:103:SER:O	1:F:106:ASP:HB3	2.21	0.40
1:G:89:ASP:HB3	1:G:151:ASP:OD2	2.21	0.40
1:A:155:ALA:O	1:A:159:THR:HG23	2.21	0.40
1:F:119:SER:O	1:F:120:ASP:C	2.64	0.40
1:G:42:ALA:C	1:G:45:THR:HG22	2.41	0.40
1:G:62:HIS:HD2	1:G:183:ILE:HG23	1.87	0.40
1:C:95:ILE:HD11	1:C:157:ALA:HA	2.03	0.40
1:G:198:ILE:HG12	1:H:48:ILE:HD11	2.03	0.40

There are no symmetry-related clashes.

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	162/181 (90%)	154 (95%)	6 (4%)	2 (1%)	10	27
1	B	162/181 (90%)	152 (94%)	9 (6%)	1 (1%)	21	44
1	C	162/181 (90%)	155 (96%)	7 (4%)	0	100	100
1	D	162/181 (90%)	155 (96%)	6 (4%)	1 (1%)	21	44
1	E	162/181 (90%)	155 (96%)	6 (4%)	1 (1%)	21	44
1	F	162/181 (90%)	154 (95%)	7 (4%)	1 (1%)	21	44
1	G	162/181 (90%)	155 (96%)	6 (4%)	1 (1%)	21	44
1	H	162/181 (90%)	155 (96%)	6 (4%)	1 (1%)	21	44
All	All	1296/1448 (90%)	1235 (95%)	53 (4%)	8 (1%)	21	44

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	39	ILE
1	D	39	ILE
1	E	41	LEU
1	F	39	ILE
1	H	39	ILE
1	A	41	LEU
1	A	89	ASP
1	G	41	LEU

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	126/140 (90%)	105 (83%)	21 (17%)	2	6
1	B	126/140 (90%)	106 (84%)	20 (16%)	2	7
1	C	126/140 (90%)	105 (83%)	21 (17%)	2	6
1	D	126/140 (90%)	109 (86%)	17 (14%)	4	9

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	E	126/140 (90%)	107 (85%)	19 (15%)	3	7
1	F	126/140 (90%)	108 (86%)	18 (14%)	3	8
1	G	126/140 (90%)	105 (83%)	21 (17%)	2	6
1	H	126/140 (90%)	99 (79%)	27 (21%)	1	3
All	All	1008/1120 (90%)	844 (84%)	164 (16%)	2	6

All (164) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	39	ILE
1	A	41	LEU
1	A	44	ILE
1	A	45	THR
1	A	49	THR
1	A	59	LYS
1	A	66	LYS
1	A	90	LYS
1	A	93	LYS
1	A	108	LYS
1	A	111	SER
1	A	114	LYS
1	A	115	LYS
1	A	116	VAL
1	A	130	LYS
1	A	135	SER
1	A	139	LYS
1	A	142	SER
1	A	169	LYS
1	A	178	LYS
1	A	195	THR
1	B	39	ILE
1	B	43	THR
1	B	46	LYS
1	B	56	LYS
1	B	59	LYS
1	B	66	LYS
1	B	77	LYS
1	B	78	LYS
1	B	85	GLU
1	B	86	THR
1	B	93	LYS

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Mol	Chain	Res	Type
1	B	95	ILE
1	B	103	SER
1	B	111	SER
1	B	112	LEU
1	B	115	LYS
1	B	153	LYS
1	B	195	THR
1	B	199	ASN
1	B	201	LEU
1	C	39	ILE
1	C	41	LEU
1	C	44	ILE
1	C	45	THR
1	C	46	LYS
1	C	48	ILE
1	C	66	LYS
1	C	77	LYS
1	C	79	ILE
1	C	90	LYS
1	C	114	LYS
1	C	116	VAL
1	C	118	ILE
1	C	125	LYS
1	C	130	LYS
1	C	135	SER
1	C	142	SER
1	C	149	LYS
1	C	153	LYS
1	C	169	LYS
1	C	177	ILE
1	D	39	ILE
1	D	41	LEU
1	D	56	LYS
1	D	59	LYS
1	D	60	ASP
1	D	66	LYS
1	D	78	LYS
1	D	93	LYS
1	D	112	LEU
1	D	115	LYS
1	D	116	VAL
1	D	123	LYS

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Mol	Chain	Res	Type
1	D	142	SER
1	D	169	LYS
1	D	174	GLU
1	D	195	THR
1	D	201	LEU
1	E	39	ILE
1	E	44	ILE
1	E	46	LYS
1	E	59	LYS
1	E	66	LYS
1	E	78	LYS
1	E	86	THR
1	E	90	LYS
1	E	93	LYS
1	E	112	LEU
1	E	114	LYS
1	E	118	ILE
1	E	125	LYS
1	E	146	ASP
1	E	149	LYS
1	E	153	LYS
1	E	166	THR
1	E	169	LYS
1	E	201	LEU
1	F	38	VAL
1	F	39	ILE
1	F	41	LEU
1	F	46	LYS
1	F	59	LYS
1	F	78	LYS
1	F	85	GLU
1	F	89	ASP
1	F	90	LYS
1	F	116	VAL
1	F	129	VAL
1	F	133	SER
1	F	135	SER
1	F	150	ASP
1	F	153	LYS
1	F	171	LYS
1	F	174	GLU
1	F	195	THR

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Mol	Chain	Res	Type
1	G	39	ILE
1	G	41	LEU
1	G	46	LYS
1	G	48	ILE
1	G	66	LYS
1	G	78	LYS
1	G	90	LYS
1	G	93	LYS
1	G	103	SER
1	G	111	SER
1	G	116	VAL
1	G	123	LYS
1	G	130	LYS
1	G	134	THR
1	G	135	SER
1	G	142	SER
1	G	169	LYS
1	G	170	ASP
1	G	189	SER
1	G	191	GLU
1	G	195	THR
1	H	39	ILE
1	H	41	LEU
1	H	43	THR
1	H	44	ILE
1	H	46	LYS
1	H	48	ILE
1	H	52	VAL
1	H	56	LYS
1	H	59	LYS
1	H	66	LYS
1	H	77	LYS
1	H	78	LYS
1	H	90	LYS
1	H	93	LYS
1	H	95	ILE
1	H	102	ILE
1	H	115	LYS
1	H	116	VAL
1	H	123	LYS
1	H	126	ILE
1	H	153	LYS

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Mol	Chain	Res	Type
1	H	163	ILE
1	H	165	ASP
1	H	169	LYS
1	H	171	LYS
1	H	174	GLU
1	H	201	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	47	ASN
1	B	199	ASN
1	C	47	ASN
1	E	47	ASN
1	G	47	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](#)

Of 14 ligands modelled in this entry, 14 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

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	A	164/181 (90%)	-0.13	5 (3%)	52	49	15, 19, 21, 22	0
1	B	164/181 (90%)	0.02	3 (1%)	67	65	17, 19, 21, 22	0
1	C	164/181 (90%)	0.00	7 (4%)	40	36	17, 19, 21, 22	0
1	D	164/181 (90%)	0.04	4 (2%)	59	56	16, 19, 21, 23	0
1	E	164/181 (90%)	0.06	7 (4%)	40	36	17, 19, 21, 23	0
1	F	164/181 (90%)	-0.05	7 (4%)	40	36	16, 19, 21, 22	0
1	G	164/181 (90%)	-0.10	3 (1%)	67	65	16, 19, 21, 22	0
1	H	164/181 (90%)	0.77	19 (11%)	9	8	16, 19, 21, 22	0
All	All	1312/1448 (90%)	0.08	55 (4%)	40	37	15, 19, 21, 23	0

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	38	VAL	7.0
1	B	38	VAL	6.8
1	A	165	ASP	5.4
1	H	38	VAL	5.2
1	H	81	ALA	4.6
1	E	38	VAL	4.5
1	G	165	ASP	4.1
1	D	39	ILE	4.0
1	C	201	LEU	4.0
1	B	39	ILE	3.9
1	C	168	ALA	3.8
1	F	38	VAL	3.8
1	H	201	LEU	3.4
1	A	199	ASN	3.4
1	C	38	VAL	3.3
1	H	113	GLU	3.3

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Mol	Chain	Res	Type	RSRZ
1	D	40	ASP	3.2
1	E	168	ALA	3.2
1	F	43	THR	3.0
1	H	83	GLY	2.9
1	F	168	ALA	2.9
1	F	120	ASP	2.9
1	H	118	ILE	2.8
1	E	165	ASP	2.8
1	A	201	LEU	2.7
1	H	82	ASN	2.7
1	H	117	GLY	2.7
1	H	39	ILE	2.6
1	C	167	GLY	2.5
1	B	41	LEU	2.5
1	A	167	GLY	2.4
1	E	201	LEU	2.4
1	E	39	ILE	2.3
1	H	80	GLY	2.3
1	C	41	LEU	2.3
1	H	110	ALA	2.3
1	H	166	THR	2.3
1	D	42	ALA	2.3
1	E	166	THR	2.3
1	F	39	ILE	2.2
1	H	200	ALA	2.2
1	H	78	LYS	2.2
1	G	83	GLY	2.2
1	C	81	ALA	2.2
1	H	86	THR	2.2
1	F	201	LEU	2.2
1	H	174	GLU	2.2
1	E	41	LEU	2.2
1	C	165	ASP	2.2
1	H	73	LYS	2.1
1	F	42	ALA	2.1
1	A	39	ILE	2.1
1	H	126	ILE	2.1
1	H	164	ALA	2.1
1	G	38	VAL	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($\text{\AA}^2$)	Q<0.9
2	NI	H	13	1/1	0.85	0.09	80,80,80,80	0
2	NI	D	8	1/1	0.90	0.09	67,67,67,67	0
2	NI	E	5	1/1	0.90	0.07	48,48,48,48	0
2	NI	E	12	1/1	0.90	0.13	74,74,74,74	0
2	NI	G	11	1/1	0.90	0.09	67,67,67,67	0
2	NI	A	14	1/1	0.90	0.09	61,61,61,61	0
2	NI	H	10	1/1	0.91	0.07	49,49,49,49	0
2	NI	B	4	1/1	0.91	0.07	65,65,65,65	0
2	NI	C	3	1/1	0.94	0.06	43,43,43,43	0
2	NI	F	6	1/1	0.96	0.05	43,43,43,43	0
2	NI	A	7	1/1	0.96	0.06	41,41,41,41	0
2	NI	D	1	1/1	0.97	0.04	42,42,42,42	0
2	NI	B	2	1/1	0.97	0.04	42,42,42,42	0
2	NI	G	9	1/1	0.98	0.04	47,47,47,47	0

6.5 Other polymers [i](#)

There are no such residues in this entry.