



# Full wwPDB X-ray Structure Validation Report ⓘ

Sep 19, 2020 – 09:22 AM BST

PDB ID : 6VFN  
Title : Crystal structure of SpeG allosteric polyamine acetyltransferase from *Bacillus thuringiensis* in complex with spermine  
Authors : Tsimbalyuk, S.; Shornikov, A.; Le, V.T.B.; Kuhn, M.L.; Forwood, J.K.  
Deposited on : 2020-01-05  
Resolution : 2.50 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](mailto: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.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.14.6  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.14.6

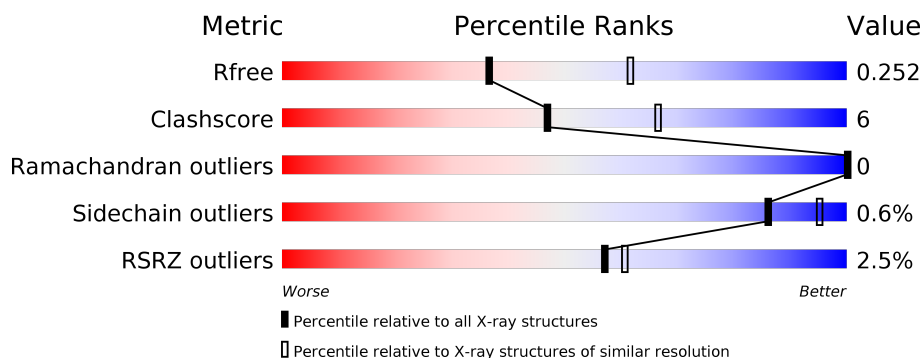
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.50 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))
$R_{free}$	130704	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	171	<div> <div></div> <div>77% 20% .</div> </div>
1	B	171	<div> <div>4%</div> <div>80% 18% .</div> </div>
1	C	171	<div> <div>%</div> <div>77% 20% .</div> </div>
1	D	171	<div> <div>4%</div> <div>84% 13% .</div> </div>
1	E	171	<div> <div>2%</div> <div>78% 19% .</div> </div>
1	F	171	<div> <div>%</div> <div>82% 16% .</div> </div>

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Mol	Chain	Length	Quality of chain
1	G	171	<div><div><div></div><div></div><div></div></div><div><div>2%</div><div>77%</div><div>20%</div><div>..</div></div></div>
1	H	171	<div><div><div></div><div></div><div></div></div><div><div>2%</div><div>85%</div><div>13%</div><div>.</div></div></div>
1	I	171	<div><div><div></div><div></div><div></div></div><div><div>4%</div><div>74%</div><div>24%</div><div>.</div></div></div>
1	J	171	<div><div><div></div><div></div><div></div></div><div><div>5%</div><div>82%</div><div>15%</div><div>..</div></div></div>
1	K	171	<div><div><div></div><div></div><div></div></div><div><div>3%</div><div>81%</div><div>16%</div><div>.</div></div></div>
1	L	171	<div><div><div></div><div></div><div></div></div><div><div>2%</div><div>85%</div><div>12%</div><div>.</div></div></div>

## 2 Entry composition

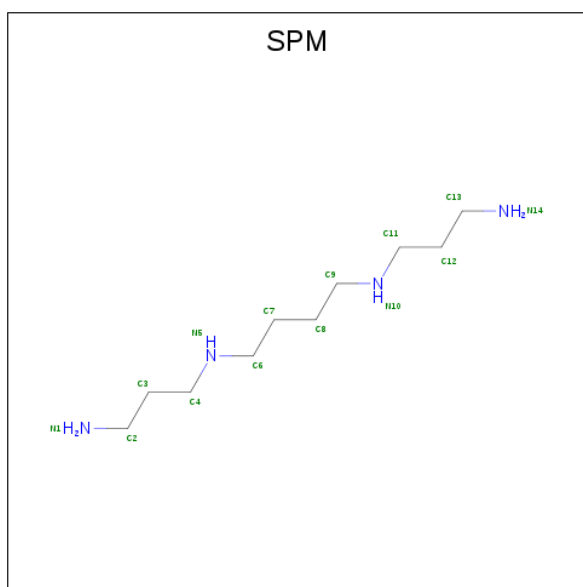
There are 2 unique types of molecules in this entry. The entry contains 16935 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Spermidine N1-acetyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	167	Total	C	N	O	S	0	0	0
			1403	903	229	263	8			
1	B	167	Total	C	N	O	S	0	0	0
			1403	904	230	261	8			
1	C	167	Total	C	N	O	S	0	0	0
			1411	908	231	264	8			
1	D	167	Total	C	N	O	S	0	0	0
			1391	896	228	259	8			
1	E	167	Total	C	N	O	S	0	0	0
			1411	909	231	263	8			
1	F	167	Total	C	N	O	S	0	0	0
			1392	898	230	256	8			
1	G	167	Total	C	N	O	S	0	0	0
			1381	888	228	257	8			
1	H	167	Total	C	N	O	S	0	0	0
			1403	904	229	262	8			
1	I	167	Total	C	N	O	S	0	0	0
			1399	901	229	261	8			
1	J	167	Total	C	N	O	S	0	0	0
			1399	899	228	264	8			
1	K	167	Total	C	N	O	S	0	0	0
			1379	889	226	256	8			
1	L	167	Total	C	N	O	S	0	0	0
			1395	898	228	261	8			

- Molecule 2 is SPERMINE (three-letter code: SPM) (formula: C<sub>10</sub>H<sub>26</sub>N<sub>4</sub>) (labeled as "Ligand of Interest" by author).



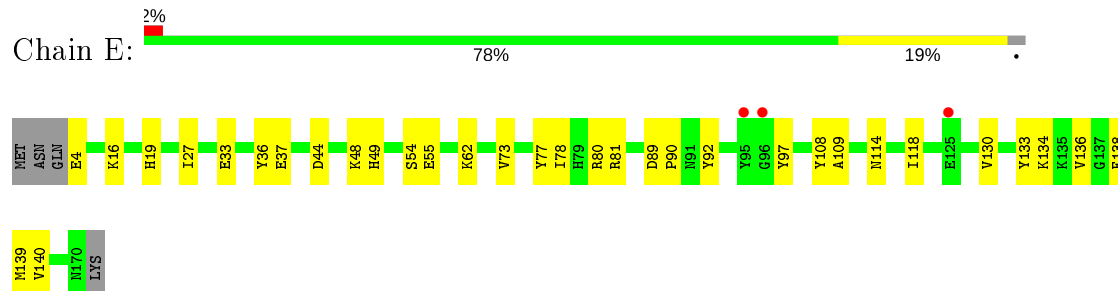
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total 14	C 10	N 4	0	0
2	B	1	Total 14	C 10	N 4	0	0
2	C	1	Total 14	C 10	N 4	0	0
2	D	1	Total 14	C 10	N 4	0	0
2	E	1	Total 14	C 10	N 4	0	0
2	E	1	Total 14	C 10	N 4	0	0
2	F	1	Total 14	C 10	N 4	0	0
2	G	1	Total 14	C 10	N 4	0	0
2	H	1	Total 14	C 10	N 4	0	0
2	J	1	Total 14	C 10	N 4	0	0
2	K	1	Total 14	C 10	N 4	0	0
2	L	1	Total 14	C 10	N 4	0	0



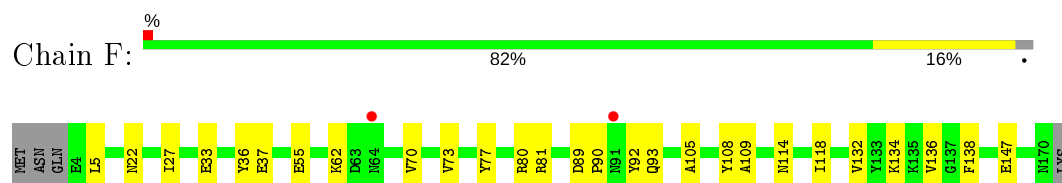
- Molecule 1: Spermidine N1-acetyltransferase



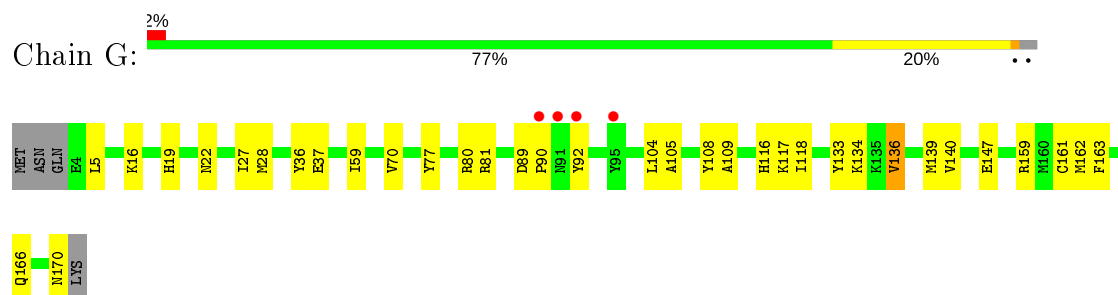
- Molecule 1: Spermidine N1-acetyltransferase



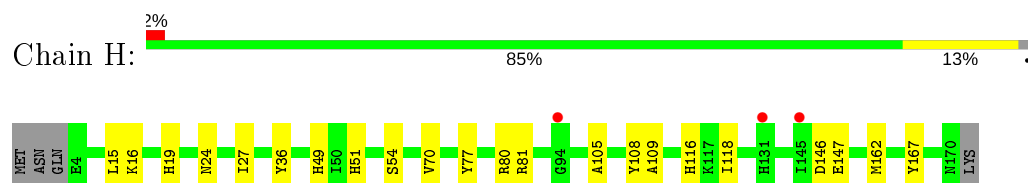
- Molecule 1: Spermidine N1-acetyltransferase



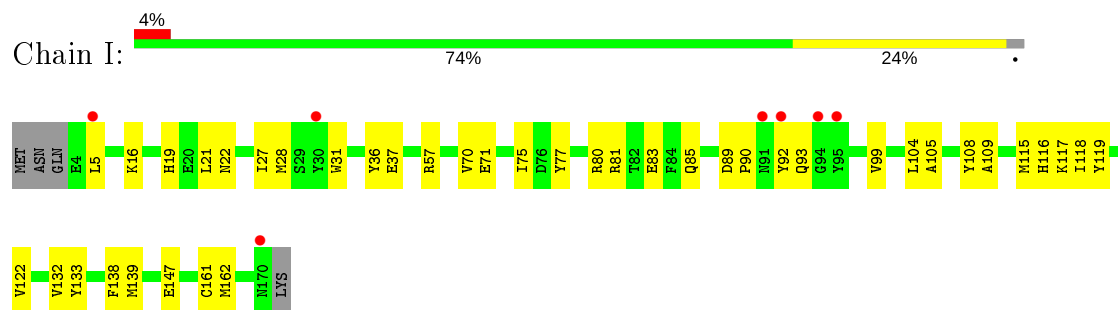
- Molecule 1: Spermidine N1-acetyltransferase



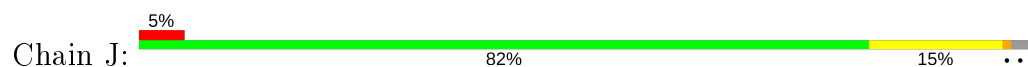
- Molecule 1: Spermidine N1-acetyltransferase

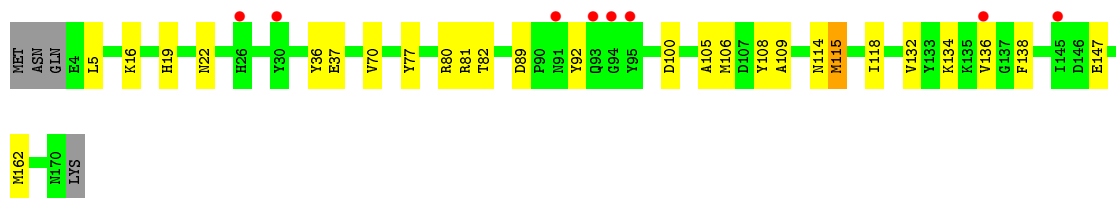


- Molecule 1: Spermidine N1-acetyltransferase

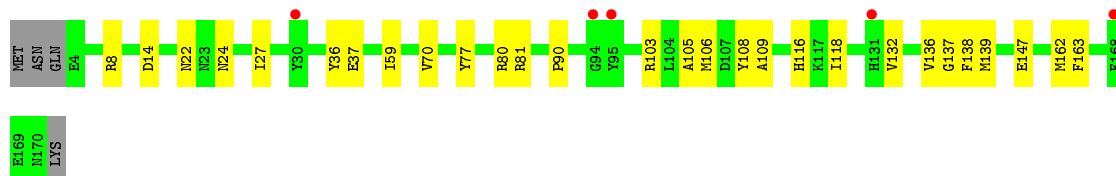
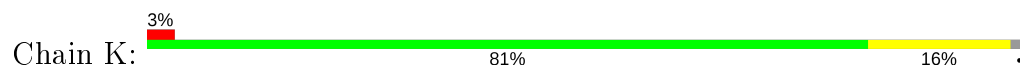


- Molecule 1: Spermidine N1-acetyltransferase

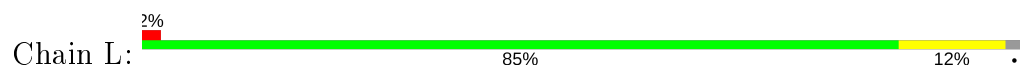




- Molecule 1: Spermidine N1-acetyltransferase



- Molecule 1: Spermidine N1-acetyltransferase





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	100.65Å 125.59Å 101.02Å 90.00° 106.50° 90.00°	Depositor
Resolution (Å)	29.87 – 2.50 29.87 – 2.50	Depositor EDS
% Data completeness (in resolution range)	100.0 (29.87-2.50) 100.0 (29.87-2.50)	Depositor EDS
$R_{merge}$	0.12	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.09 (at 2.51Å)	Xtriage
Refinement program	PHENIX refinement version: 1.13_2998	Depositor
R, $R_{free}$	0.226 , 0.252 0.226 , 0.252	Depositor DCC
$R_{free}$ test set	4214 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	39.5	Xtriage
Anisotropy	0.025	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 23.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.014 for l,k,h	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	16935	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	42.0	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.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

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

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	0.26	0/1438	0.44	0/1945
1	B	0.26	0/1438	0.43	0/1944
1	C	0.27	0/1446	0.43	0/1954
1	D	0.26	0/1426	0.44	0/1931
1	E	0.26	0/1446	0.43	0/1953
1	F	0.26	0/1427	0.44	0/1931
1	G	0.26	0/1415	0.44	0/1917
1	H	0.26	0/1438	0.43	0/1944
1	I	0.27	0/1434	0.44	0/1940
1	J	0.29	0/1434	0.46	0/1942
1	K	0.26	0/1414	0.44	0/1917
1	L	0.28	0/1430	0.45	0/1936
All	All	0.27	0/17186	0.44	0/23254

There are no bond length outliers.

There are no bond angle outliers.

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	1403	0	1307	23	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1403	0	1314	19	0
1	C	1411	0	1324	23	0
1	D	1391	0	1288	16	0
1	E	1411	0	1329	21	0
1	F	1392	0	1299	17	0
1	G	1381	0	1277	20	0
1	H	1403	0	1312	15	0
1	I	1399	0	1303	24	0
1	J	1399	0	1291	20	0
1	K	1379	0	1267	22	0
1	L	1395	0	1292	16	0
2	A	14	0	26	0	0
2	B	14	0	26	1	0
2	C	14	0	26	0	0
2	D	14	0	26	1	0
2	E	28	0	52	4	0
2	F	14	0	26	1	0
2	G	14	0	26	0	0
2	H	14	0	26	0	0
2	J	14	0	26	0	0
2	K	14	0	26	0	0
2	L	14	0	26	0	0
All	All	16935	0	15915	202	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 6.

All (202) 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:147:GLU:O	1:I:81:ARG:NH1	2.22	0.73
1:J:80:ARG:CZ	1:K:80:ARG:HD3	2.21	0.71
1:J:81:ARG:NH1	1:L:147:GLU:O	2.26	0.69
1:C:22:ASN:ND2	1:C:37:GLU:OE1	2.29	0.65
1:I:122:VAL:HG11	1:I:133:TYR:HE2	1.62	0.65
1:L:80:ARG:NH2	1:L:164:GLN:OE1	2.30	0.64
1:I:90:PRO:HA	1:I:93:GLN:HG3	1.82	0.62
1:E:27:ILE:HD11	1:E:90:PRO:HD3	1.80	0.62
1:J:89:ASP:HB3	1:J:92:TYR:HD2	1.65	0.62
1:J:77:TYR:HB3	1:K:77:TYR:HB3	1.83	0.61
1:C:109:ALA:HB3	1:C:118:ILE:HD11	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:166:GLN:O	1:C:170:ASN:ND2	2.32	0.60
1:F:109:ALA:HB3	1:F:118:ILE:HD11	1.84	0.60
1:A:109:ALA:HB3	1:A:118:ILE:HD11	1.84	0.59
1:F:147:GLU:O	1:K:81:ARG:NH1	2.36	0.59
1:A:89:ASP:HB3	1:A:92:TYR:HD1	1.68	0.59
1:D:109:ALA:HB3	1:D:118:ILE:HD11	1.83	0.59
1:K:109:ALA:HB3	1:K:118:ILE:HD11	1.85	0.58
1:G:109:ALA:HB3	1:G:118:ILE:HD11	1.86	0.58
1:J:22:ASN:ND2	1:J:37:GLU:OE1	2.27	0.58
1:F:89:ASP:HB3	1:F:92:TYR:HD2	1.68	0.58
1:A:22:ASN:ND2	1:A:37:GLU:OE1	2.34	0.57
1:A:103:ARG:HE	1:A:136:VAL:HG13	1.70	0.57
1:J:80:ARG:NH2	1:J:114:ASN:O	2.35	0.57
1:C:163:PHE:HB2	1:C:166:GLN:OE1	2.03	0.57
1:J:109:ALA:HB3	1:J:118:ILE:HD11	1.87	0.57
1:D:81:ARG:NH1	1:H:147:GLU:O	2.38	0.56
1:E:55:GLU:HG2	1:E:73:VAL:HG12	1.88	0.56
1:E:109:ALA:HB3	1:E:118:ILE:HD11	1.88	0.56
1:H:109:ALA:HB3	1:H:118:ILE:HD11	1.88	0.56
1:I:109:ALA:HB3	1:I:118:ILE:HD11	1.87	0.56
1:G:22:ASN:ND2	1:G:37:GLU:OE1	2.36	0.55
1:L:109:ALA:HB3	1:L:118:ILE:HD11	1.88	0.55
1:A:80:ARG:HB3	1:A:116:HIS:H	1.71	0.55
1:A:77:TYR:HB3	1:I:77:TYR:HB3	1.88	0.55
1:K:103:ARG:HA	1:K:106:MET:HE2	1.89	0.55
1:B:109:ALA:HB3	1:B:118:ILE:HD11	1.88	0.54
1:C:90:PRO:HA	1:C:93:GLN:HG3	1.90	0.54
1:D:77:TYR:HB3	1:L:77:TYR:HB3	1.88	0.54
1:D:28:MET:HE1	2:D:201:SPM:HN12	1.72	0.54
1:E:49:HIS:O	2:E:202:SPM:N14	2.41	0.53
1:F:90:PRO:HA	1:F:93:GLN:HG2	1.89	0.53
1:C:77:TYR:HB3	1:H:77:TYR:HB3	1.91	0.53
1:F:77:TYR:HB3	1:G:77:TYR:HB3	1.91	0.53
1:B:81:ARG:NH1	1:I:147:GLU:O	2.40	0.53
1:K:22:ASN:ND2	1:K:37:GLU:OE1	2.35	0.52
1:I:80:ARG:HB3	1:I:116:HIS:H	1.75	0.52
1:L:133:TYR:HA	1:L:136:VAL:HG22	1.91	0.52
1:K:70:VAL:HG11	1:K:105:ALA:HB2	1.91	0.51
1:J:147:GLU:O	1:L:81:ARG:NH1	2.42	0.51
1:F:70:VAL:HG11	1:F:105:ALA:HB2	1.92	0.51
1:B:55:GLU:HG2	1:B:73:VAL:HG12	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:80:ARG:HB3	1:G:116:HIS:H	1.76	0.51
1:F:132:VAL:O	1:F:136:VAL:HG23	2.11	0.50
1:F:5:LEU:HD12	1:F:62:LYS:HB2	1.92	0.50
1:F:27:ILE:HD11	1:F:90:PRO:HD3	1.92	0.50
1:D:147:GLU:O	1:H:81:ARG:NH1	2.45	0.50
1:A:5:LEU:HD12	1:A:62:LYS:HB2	1.92	0.50
1:B:80:ARG:HH22	1:B:164:GLN:HE22	1.59	0.50
1:C:89:ASP:HB3	1:C:92:TYR:HD1	1.75	0.50
1:E:133:TYR:HA	1:E:136:VAL:HG22	1.94	0.50
1:K:138:PHE:CD1	1:K:162:MET:HB2	2.47	0.50
1:B:113:LEU:HA	1:E:78:ILE:HD12	1.94	0.49
1:C:138:PHE:CD1	1:C:162:MET:HB2	2.47	0.49
1:D:108:TYR:OH	1:J:36:TYR:HB2	2.13	0.48
1:B:22:ASN:ND2	1:B:37:GLU:OE1	2.42	0.48
1:C:16:LYS:O	1:C:19:HIS:HB3	2.13	0.48
1:C:22:ASN:HA	1:C:28:MET:HE3	1.95	0.48
1:G:5:LEU:HD23	1:G:104:LEU:HD12	1.95	0.48
1:B:108:TYR:OH	1:C:36:TYR:HB2	2.13	0.48
1:C:134:LYS:HE3	1:C:140:VAL:HG23	1.95	0.48
1:E:130:VAL:HG12	1:E:134:LYS:HE3	1.95	0.48
1:E:89:ASP:HB3	1:E:92:TYR:HD1	1.77	0.48
1:F:55:GLU:HG2	1:F:73:VAL:HG12	1.96	0.48
1:I:70:VAL:HG11	1:I:105:ALA:HB2	1.95	0.48
1:F:81:ARG:NH1	1:K:147:GLU:O	2.44	0.48
1:B:165:LYS:NZ	1:B:169:GLU:OE2	2.38	0.48
1:C:27:ILE:HD11	1:C:90:PRO:HD3	1.95	0.48
1:B:77:TYR:HB3	1:E:77:TYR:HB3	1.96	0.48
1:H:80:ARG:HB3	1:H:116:HIS:H	1.78	0.48
1:B:89:ASP:HB3	1:B:92:TYR:HD1	1.79	0.47
1:G:108:TYR:OH	1:K:36:TYR:HB2	2.13	0.47
1:D:59:ILE:HG21	1:D:66:MET:HG2	1.96	0.47
1:E:33:GLU:OE2	2:E:201:SPM:H42	2.14	0.47
1:K:132:VAL:O	1:K:136:VAL:HG23	2.14	0.47
1:K:80:ARG:HB3	1:K:116:HIS:H	1.78	0.47
1:J:138:PHE:CD1	1:J:162:MET:HB2	2.49	0.47
1:L:80:ARG:HB3	1:L:116:HIS:H	1.79	0.47
1:C:70:VAL:HG11	1:C:105:ALA:HB2	1.96	0.47
1:K:27:ILE:HD11	1:K:90:PRO:HG3	1.97	0.47
1:B:134:LYS:HG3	1:B:160:MET:HE1	1.96	0.47
1:C:5:LEU:HD12	1:C:62:LYS:HB2	1.96	0.47
1:G:22:ASN:HA	1:G:28:MET:HE3	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:139:MET:HG2	1:I:161:CYS:O	2.14	0.47
1:J:109:ALA:O	1:J:115:MET:HB2	2.14	0.47
1:A:80:ARG:NE	1:A:114:ASN:O	2.44	0.47
1:D:70:VAL:HG11	1:D:105:ALA:HB2	1.96	0.47
1:L:138:PHE:CD1	1:L:162:MET:HB2	2.50	0.47
1:A:70:VAL:HG11	1:A:105:ALA:HB2	1.96	0.47
1:D:132:VAL:O	1:D:136:VAL:HG23	2.14	0.47
1:L:124:LYS:HZ2	1:L:156:ASN:HB2	1.80	0.47
1:A:5:LEU:HD11	1:A:60:VAL:HG12	1.96	0.47
1:A:22:ASN:HA	1:A:28:MET:HE3	1.97	0.46
1:A:108:TYR:OH	1:B:36:TYR:HB2	2.14	0.46
1:H:162:MET:HE1	1:H:167:TYR:HD1	1.79	0.46
1:I:5:LEU:HD23	1:I:104:LEU:HD12	1.97	0.46
1:K:103:ARG:HA	1:K:106:MET:CE	2.46	0.46
1:L:124:LYS:HB2	1:L:124:LYS:HE3	1.66	0.46
1:A:81:ARG:NH1	1:G:147:GLU:O	2.47	0.46
1:G:36:TYR:HB2	1:I:108:TYR:OH	2.15	0.46
1:B:139:MET:HG2	1:B:161:CYS:O	2.15	0.46
1:J:132:VAL:O	1:J:136:VAL:HG23	2.16	0.46
1:D:28:MET:HE3	1:D:33:GLU:HB2	1.96	0.46
1:D:61:GLU:HA	1:D:65:GLU:O	2.16	0.46
1:E:16:LYS:O	1:E:19:HIS:HB3	2.16	0.46
2:E:201:SPM:N14	1:H:49:HIS:O	2.49	0.46
1:G:70:VAL:HG11	1:G:105:ALA:HB2	1.98	0.46
1:E:36:TYR:HB2	1:H:108:TYR:OH	2.15	0.46
1:H:24:ASN:HB3	1:H:27:ILE:HB	1.98	0.46
1:I:22:ASN:ND2	1:I:37:GLU:OE1	2.37	0.46
1:F:36:TYR:HB2	1:J:108:TYR:OH	2.16	0.46
1:B:27:ILE:HD11	1:B:90:PRO:HG3	1.99	0.45
1:K:24:ASN:HB3	1:K:27:ILE:HB	1.97	0.45
1:F:80:ARG:HH21	1:F:114:ASN:HB3	1.81	0.45
1:J:16:LYS:O	1:J:19:HIS:HB3	2.17	0.45
1:A:55:GLU:HG2	1:A:73:VAL:HG12	1.98	0.45
1:E:44:ASP:OD1	1:H:51:HIS:NE2	2.47	0.45
1:H:36:TYR:HB2	1:L:108:TYR:OH	2.17	0.45
1:J:70:VAL:HG11	1:J:105:ALA:HB2	1.99	0.45
1:J:134:LYS:HA	1:J:138:PHE:O	2.16	0.44
1:K:139:MET:HE2	1:K:163:PHE:HE1	1.83	0.44
1:L:124:LYS:HZ1	1:L:156:ASN:HD22	1.64	0.44
1:A:124:LYS:HD2	1:A:158:ILE:HD11	2.00	0.44
1:E:48:LYS:HA	1:E:48:LYS:HD3	1.72	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:33:GLU:OE2	2:F:201:SPM:N1	2.50	0.44
1:A:24:ASN:HB3	1:A:27:ILE:HB	2.00	0.44
1:I:21:LEU:O	1:I:28:MET:HE3	2.18	0.44
1:G:16:LYS:O	1:G:19:HIS:HB3	2.18	0.44
1:C:5:LEU:HD11	1:C:60:VAL:HG12	2.00	0.44
1:I:27:ILE:HD11	1:I:90:PRO:HD3	1.99	0.44
1:A:103:ARG:HA	1:A:106:MET:HE2	2.00	0.43
1:I:16:LYS:O	1:I:19:HIS:HB3	2.18	0.43
1:F:134:LYS:HA	1:F:138:PHE:O	2.19	0.43
1:I:83:GLU:HA	1:I:119:TYR:O	2.18	0.43
1:I:81:ARG:HA	1:I:117:LYS:O	2.19	0.43
1:C:147:GLU:O	1:E:81:ARG:NH1	2.52	0.43
1:G:163:PHE:H	1:G:166:GLN:NE2	2.16	0.43
1:C:134:LYS:HA	1:C:138:PHE:O	2.19	0.43
1:E:62:LYS:HD3	1:E:97:TYR:CE1	2.54	0.43
1:K:108:TYR:OH	1:L:36:TYR:HB2	2.18	0.43
1:B:70:VAL:HG11	1:B:105:ALA:HB2	2.00	0.43
1:A:134:LYS:HA	1:A:138:PHE:O	2.19	0.43
1:C:123:ASP:OD1	1:C:125:GLU:HG2	2.19	0.43
1:G:134:LYS:HE2	1:G:140:VAL:HG23	2.00	0.43
1:C:75:ILE:HG23	1:C:115:MET:HE2	2.01	0.42
1:A:71:GLU:HB3	1:A:73:VAL:HG13	2.00	0.42
1:I:89:ASP:HB3	1:I:92:TYR:HD1	1.83	0.42
1:J:5:LEU:HD22	1:J:100:ASP:HB2	2.01	0.42
1:A:16:LYS:O	1:A:19:HIS:HB3	2.19	0.42
1:L:71:GLU:HB3	1:L:73:VAL:HG13	2.00	0.42
1:D:138:PHE:CD1	1:D:162:MET:HB2	2.55	0.42
1:L:55:GLU:HG2	1:L:73:VAL:HG12	2.01	0.42
1:B:33:GLU:OE1	2:B:201:SPM:H42	2.20	0.42
1:C:108:TYR:OH	1:D:36:TYR:HB2	2.19	0.42
1:G:133:TYR:O	1:G:136:VAL:HG12	2.19	0.42
1:E:37:GLU:OE2	2:E:201:SPM:H31	2.20	0.42
1:A:36:TYR:HB2	1:F:108:TYR:OH	2.19	0.42
1:A:138:PHE:CD2	1:A:162:MET:HB2	2.54	0.42
1:B:132:VAL:O	1:B:136:VAL:HG23	2.20	0.42
1:C:132:VAL:O	1:C:136:VAL:HG13	2.20	0.42
1:E:133:TYR:O	1:E:138:PHE:HB2	2.20	0.42
1:G:27:ILE:HD11	1:G:90:PRO:HG3	2.02	0.42
1:I:75:ILE:HG23	1:I:115:MET:HE2	2.01	0.42
1:A:147:GLU:OE2	1:G:159:ARG:NH2	2.53	0.42
1:J:82:THR:HG22	1:J:115:MET:HG2	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:59:ILE:HD13	1:K:59:ILE:HA	1.96	0.42
1:E:108:TYR:OH	1:I:36:TYR:HB2	2.20	0.41
1:H:15:LEU:HD23	1:H:15:LEU:HA	1.94	0.41
1:I:57:ARG:NE	1:I:71:GLU:OE1	2.51	0.41
1:C:24:ASN:HB3	1:C:27:ILE:HB	2.02	0.41
1:K:8:ARG:NH1	1:K:14:ASP:OD2	2.42	0.41
1:G:139:MET:HG2	1:G:161:CYS:O	2.21	0.41
1:H:16:LYS:O	1:H:19:HIS:HB3	2.20	0.41
1:B:138:PHE:CD1	1:B:162:MET:HB2	2.56	0.41
1:D:163:PHE:CE1	1:H:146:ASP:HB2	2.56	0.41
1:I:99:VAL:HG22	1:I:132:VAL:HG13	2.03	0.41
1:E:80:ARG:HH21	1:E:114:ASN:HB3	1.86	0.41
1:K:80:ARG:CZ	1:K:116:HIS:HB2	2.51	0.41
1:L:133:TYR:HB3	1:L:138:PHE:HB2	2.03	0.41
1:G:59:ILE:HD13	1:G:59:ILE:HA	1.96	0.41
1:I:138:PHE:CD1	1:I:162:MET:HB2	2.56	0.41
1:E:134:LYS:HE2	1:E:140:VAL:HG23	2.03	0.41
1:F:22:ASN:ND2	1:F:37:GLU:OE1	2.47	0.40
1:D:16:LYS:O	1:D:19:HIS:HB3	2.22	0.40
1:K:137:GLY:HA3	1:K:162:MET:HE2	2.03	0.40
1:J:80:ARG:NE	1:K:80:ARG:HD3	2.36	0.40
1:D:59:ILE:HD13	1:D:59:ILE:HA	1.96	0.40
1:G:81:ARG:HA	1:G:117:LYS:O	2.21	0.40
1:H:70:VAL:HG11	1:H:105:ALA:HB2	2.03	0.40
1:G:89:ASP:HB3	1:G:92:TYR:HD1	1.86	0.40
1:J:106:MET:HB2	1:J:106:MET:HE2	1.95	0.40
1:I:31:TRP:CE2	1:I:85:GLN:HB3	2.56	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	A	165/171 (96%)	160 (97%)	5 (3%)	0	100	100
1	B	165/171 (96%)	160 (97%)	5 (3%)	0	100	100
1	C	165/171 (96%)	160 (97%)	5 (3%)	0	100	100
1	D	165/171 (96%)	161 (98%)	4 (2%)	0	100	100
1	E	165/171 (96%)	161 (98%)	4 (2%)	0	100	100
1	F	165/171 (96%)	160 (97%)	5 (3%)	0	100	100
1	G	165/171 (96%)	160 (97%)	5 (3%)	0	100	100
1	H	165/171 (96%)	160 (97%)	5 (3%)	0	100	100
1	I	165/171 (96%)	160 (97%)	5 (3%)	0	100	100
1	J	165/171 (96%)	159 (96%)	6 (4%)	0	100	100
1	K	165/171 (96%)	161 (98%)	4 (2%)	0	100	100
1	L	165/171 (96%)	160 (97%)	5 (3%)	0	100	100
All	All	1980/2052 (96%)	1922 (97%)	58 (3%)	0	100	100

There are no Ramachandran outliers to report.

### 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	148/157 (94%)	148 (100%)	0	100	100
1	B	148/157 (94%)	148 (100%)	0	100	100
1	C	150/157 (96%)	150 (100%)	0	100	100
1	D	145/157 (92%)	144 (99%)	1 (1%)	84	94
1	E	150/157 (96%)	147 (98%)	3 (2%)	55	79
1	F	145/157 (92%)	145 (100%)	0	100	100
1	G	143/157 (91%)	140 (98%)	3 (2%)	53	78
1	H	148/157 (94%)	147 (99%)	1 (1%)	84	94
1	I	147/157 (94%)	147 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	J	147/157 (94%)	146 (99%)	1 (1%)	84	94
1	K	142/157 (90%)	142 (100%)	0	100	100
1	L	146/157 (93%)	144 (99%)	2 (1%)	67	86
All	All	1759/1884 (93%)	1748 (99%)	11 (1%)	86	95

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	54	SER
1	E	4	GLU
1	E	54	SER
1	E	139	MET
1	G	136	VAL
1	G	162	MET
1	G	170	ASN
1	H	54	SER
1	J	115	MET
1	L	139	MET
1	L	170	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	91	ASN
1	B	164	GLN
1	C	91	ASN
1	F	64	ASN
1	F	91	ASN
1	G	164	GLN
1	G	166	GLN
1	J	91	ASN
1	J	116	HIS
1	K	91	ASN
1	L	91	ASN
1	L	170	ASN

### 5.3.3 RNA ⓘ

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 monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

12 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	$\# Z  > 2$
2	SPM	B	201	-	13,13,13	0.33	0	12,12,12	0.70	0
2	SPM	A	201	-	13,13,13	0.34	0	12,12,12	0.71	0
2	SPM	D	201	-	13,13,13	0.33	0	12,12,12	0.71	0
2	SPM	K	201	-	13,13,13	0.33	0	12,12,12	0.71	0
2	SPM	F	201	-	13,13,13	0.34	0	12,12,12	0.72	0
2	SPM	E	201	-	13,13,13	0.33	0	12,12,12	0.73	0
2	SPM	C	201	-	13,13,13	0.33	0	12,12,12	0.71	0
2	SPM	E	202	-	13,13,13	0.34	0	12,12,12	0.71	0
2	SPM	H	201	-	13,13,13	0.33	0	12,12,12	0.71	0
2	SPM	J	201	-	13,13,13	0.34	0	12,12,12	0.81	0
2	SPM	L	201	-	13,13,13	0.34	0	12,12,12	0.72	0
2	SPM	G	201	-	13,13,13	0.34	0	12,12,12	0.73	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	SPM	B	201	-	-	2/11/11/11	-
2	SPM	A	201	-	-	2/11/11/11	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	SPM	D	201	-	-	2/11/11/11	-
2	SPM	K	201	-	-	1/11/11/11	-
2	SPM	F	201	-	-	1/11/11/11	-
2	SPM	E	201	-	-	1/11/11/11	-
2	SPM	C	201	-	-	2/11/11/11	-
2	SPM	E	202	-	-	2/11/11/11	-
2	SPM	H	201	-	-	2/11/11/11	-
2	SPM	J	201	-	-	2/11/11/11	-
2	SPM	L	201	-	-	1/11/11/11	-
2	SPM	G	201	-	-	2/11/11/11	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (20) torsion outliers are listed below:

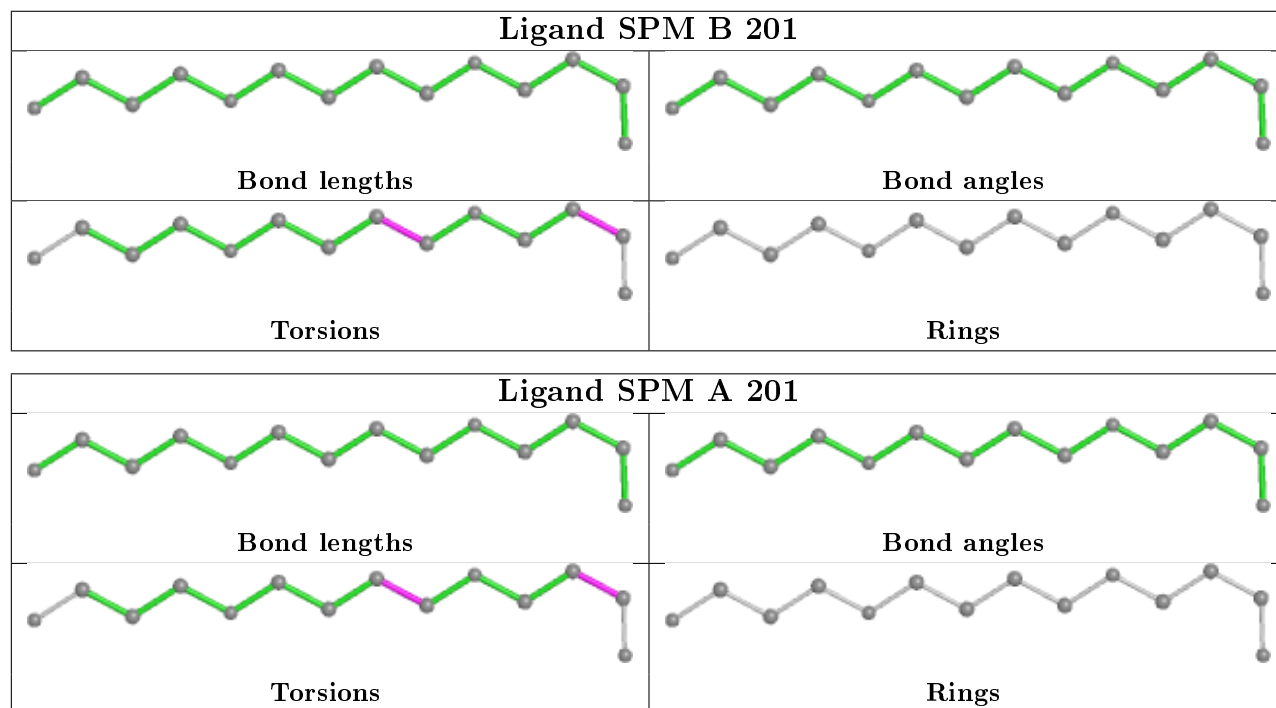
Mol	Chain	Res	Type	Atoms
2	B	201	SPM	N1-C2-C3-C4
2	A	201	SPM	N1-C2-C3-C4
2	D	201	SPM	N1-C2-C3-C4
2	K	201	SPM	N1-C2-C3-C4
2	F	201	SPM	N1-C2-C3-C4
2	E	201	SPM	N1-C2-C3-C4
2	C	201	SPM	N1-C2-C3-C4
2	E	202	SPM	N1-C2-C3-C4
2	H	201	SPM	N1-C2-C3-C4
2	J	201	SPM	N1-C2-C3-C4
2	L	201	SPM	N1-C2-C3-C4
2	G	201	SPM	N1-C2-C3-C4
2	B	201	SPM	N5-C6-C7-C8
2	H	201	SPM	N5-C6-C7-C8
2	G	201	SPM	N5-C6-C7-C8
2	A	201	SPM	N5-C6-C7-C8
2	D	201	SPM	N5-C6-C7-C8
2	J	201	SPM	N5-C6-C7-C8
2	E	202	SPM	N5-C6-C7-C8
2	C	201	SPM	N5-C6-C7-C8

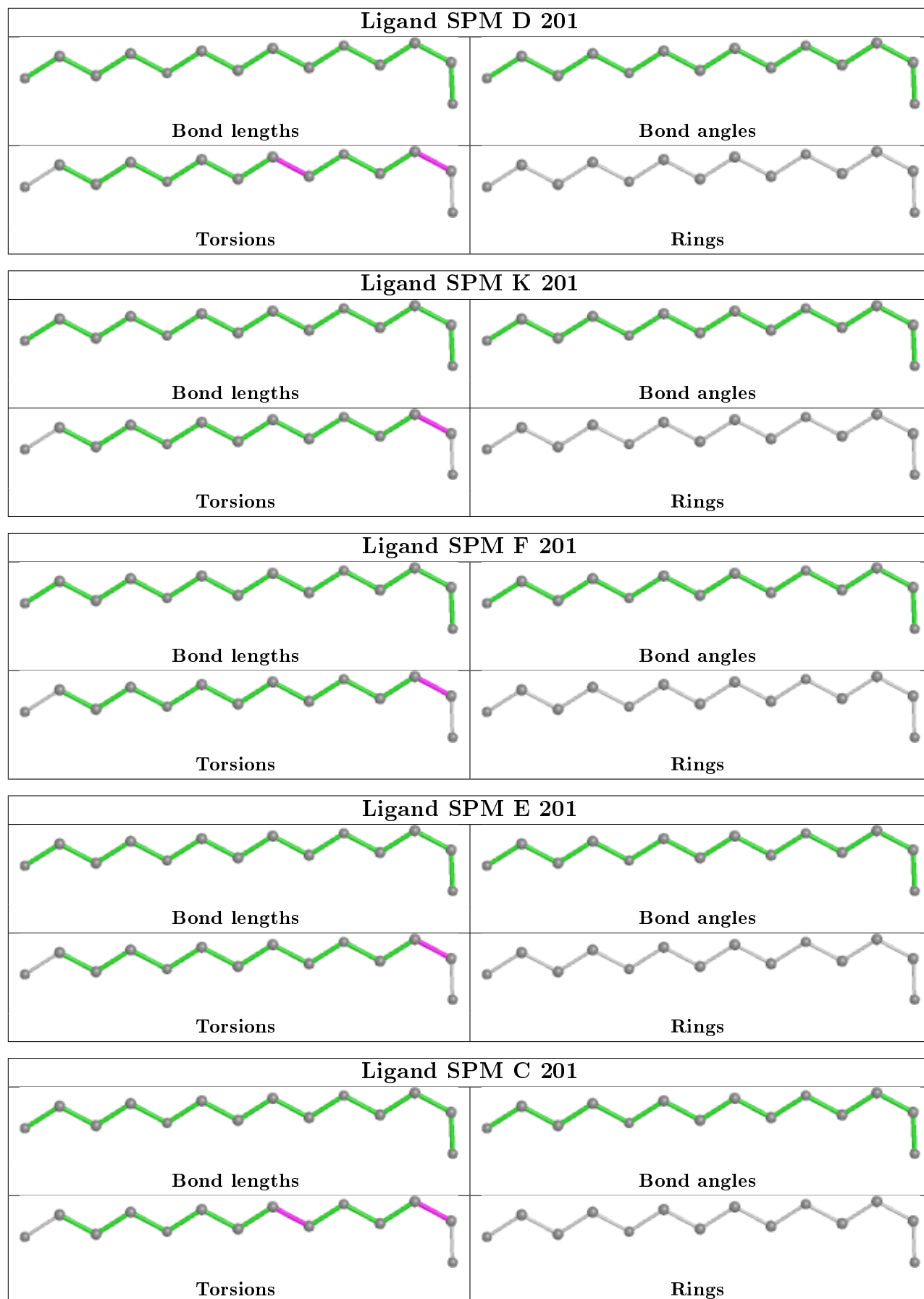
There are no ring outliers.

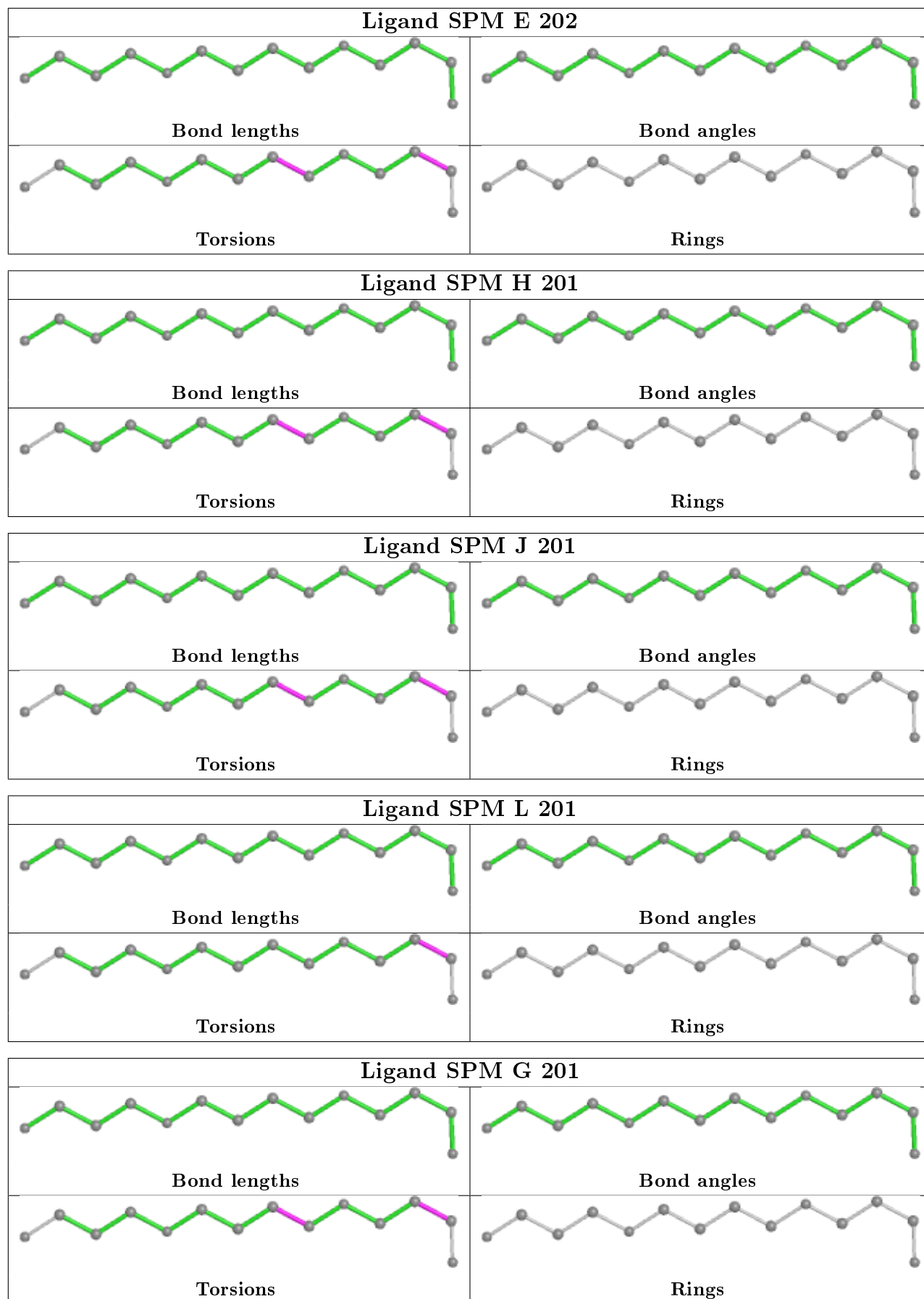
5 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	201	SPM	1	0
2	D	201	SPM	1	0
2	F	201	SPM	1	0
2	E	201	SPM	3	0
2	E	202	SPM	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.







## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	167/171 (97%)	-0.03	0 100 100	24, 35, 51, 57	0
1	B	167/171 (97%)	0.22	7 (4%) 36 39	25, 39, 56, 68	0
1	C	167/171 (97%)	0.04	1 (0%) 89 90	26, 38, 54, 65	0
1	D	167/171 (97%)	0.18	6 (3%) 42 46	29, 42, 63, 70	0
1	E	167/171 (97%)	0.20	3 (1%) 68 71	28, 40, 59, 67	0
1	F	167/171 (97%)	0.15	2 (1%) 79 80	25, 42, 57, 69	0
1	G	167/171 (97%)	0.17	4 (2%) 59 62	27, 39, 56, 68	0
1	H	167/171 (97%)	0.21	3 (1%) 68 71	30, 43, 58, 69	0
1	I	167/171 (97%)	0.26	7 (4%) 36 39	28, 40, 62, 75	0
1	J	167/171 (97%)	0.35	8 (4%) 30 32	30, 47, 64, 78	0
1	K	167/171 (97%)	0.31	5 (2%) 50 53	32, 45, 64, 71	0
1	L	167/171 (97%)	0.25	4 (2%) 59 62	32, 50, 66, 77	0
All	All	2004/2052 (97%)	0.19	50 (2%) 57 61	24, 41, 61, 78	0

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	J	94	GLY	4.9
1	E	96	GLY	4.4
1	L	131	HIS	4.1
1	D	30	TYR	3.9
1	B	125	GLU	3.8
1	G	95	TYR	3.8
1	K	131	HIS	3.7
1	D	131	HIS	3.6
1	K	30	TYR	3.4
1	B	30	TYR	3.3
1	E	95	TYR	3.3

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Mol	Chain	Res	Type	RSRZ
1	I	95	TYR	3.3
1	L	168	PHE	3.1
1	I	30	TYR	3.1
1	I	92	TYR	3.1
1	J	93	GLN	3.1
1	J	95	TYR	3.0
1	F	91	ASN	3.0
1	I	91	ASN	2.9
1	B	91	ASN	2.9
1	I	5	LEU	2.9
1	B	131	HIS	2.7
1	H	94	GLY	2.7
1	H	131	HIS	2.6
1	J	136	VAL	2.6
1	K	168	PHE	2.6
1	L	63	ASP	2.6
1	E	125	GLU	2.4
1	G	91	ASN	2.4
1	G	90	PRO	2.4
1	F	64	ASN	2.3
1	D	91	ASN	2.3
1	B	92	TYR	2.3
1	C	30	TYR	2.3
1	J	30	TYR	2.3
1	B	95	TYR	2.3
1	G	92	TYR	2.2
1	K	95	TYR	2.2
1	J	145	ILE	2.2
1	L	64	ASN	2.2
1	J	26	HIS	2.2
1	B	63	ASP	2.2
1	D	26	HIS	2.1
1	D	125	GLU	2.1
1	K	94	GLY	2.1
1	J	91	ASN	2.1
1	I	94	GLY	2.1
1	H	145	ILE	2.1
1	I	170	ASN	2.0
1	D	130	VAL	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

## 6.3 Carbohydrates ⓘ

There are no monosaccharides in this entry.

## 6.4 Ligands ⓘ

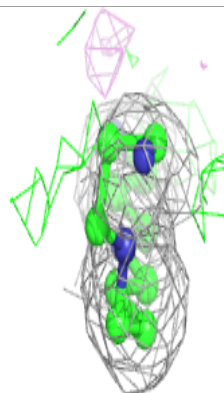
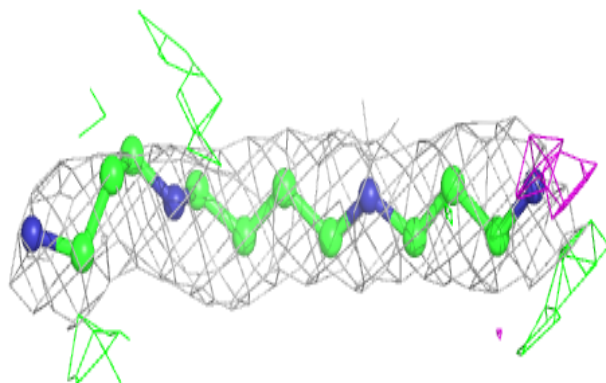
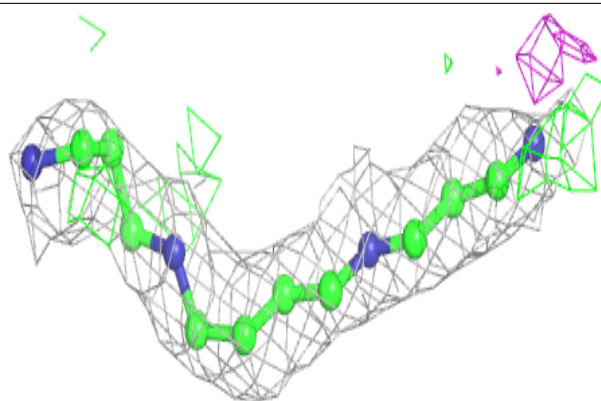
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 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	SPM	E	201	14/14	0.87	0.23	28,35,44,49	0
2	SPM	H	201	14/14	0.90	0.18	33,39,49,50	0
2	SPM	J	201	14/14	0.90	0.22	28,39,47,54	0
2	SPM	E	202	14/14	0.91	0.22	27,34,50,52	0
2	SPM	F	201	14/14	0.91	0.25	28,38,54,54	0
2	SPM	D	201	14/14	0.91	0.19	26,34,51,52	0
2	SPM	K	201	14/14	0.92	0.18	27,34,46,46	0
2	SPM	L	201	14/14	0.92	0.18	34,39,52,58	0
2	SPM	G	201	14/14	0.92	0.19	23,34,45,46	0
2	SPM	C	201	14/14	0.94	0.19	20,29,42,45	0
2	SPM	B	201	14/14	0.94	0.16	21,29,40,40	0
2	SPM	A	201	14/14	0.94	0.16	23,30,45,46	0

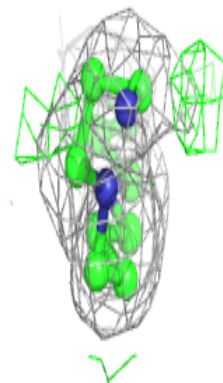
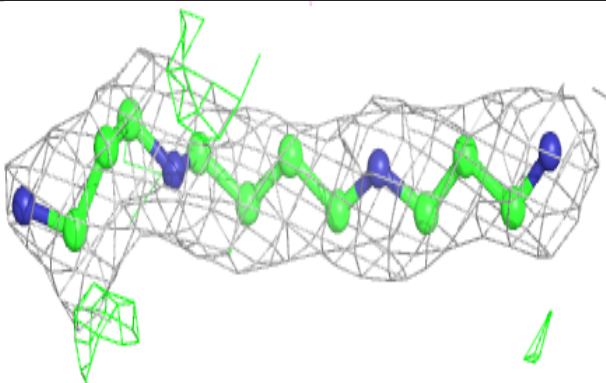
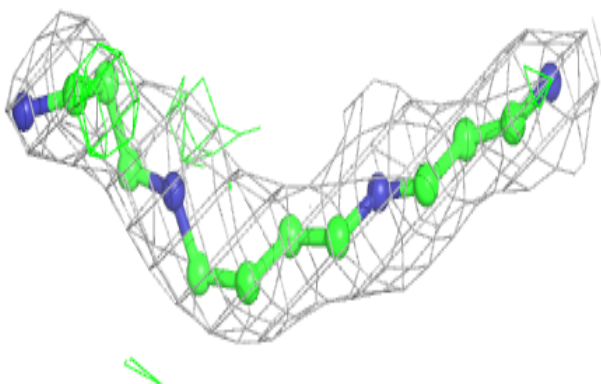
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

**Electron density around SPM E 201:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

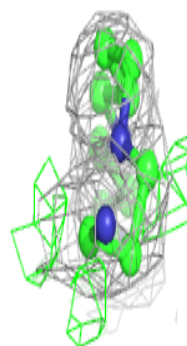
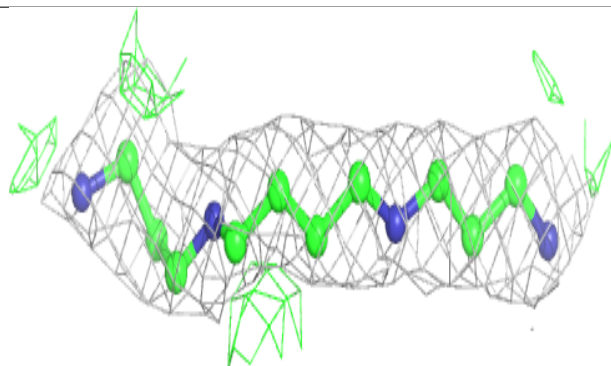
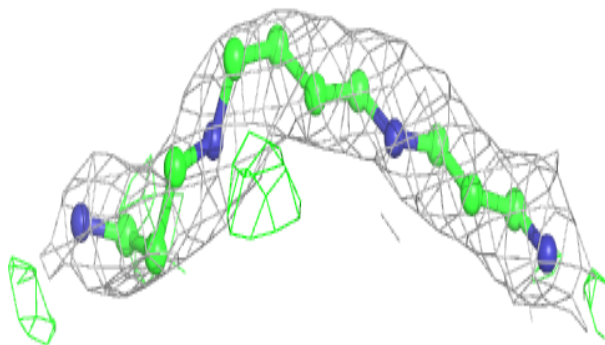
**Electron density around SPM H 201:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

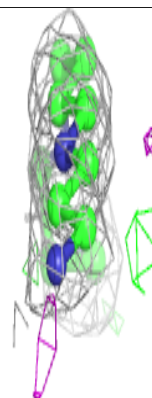
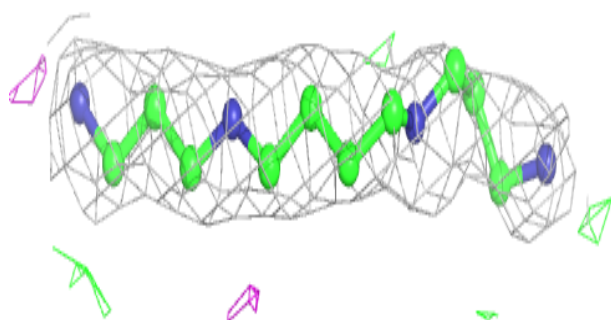
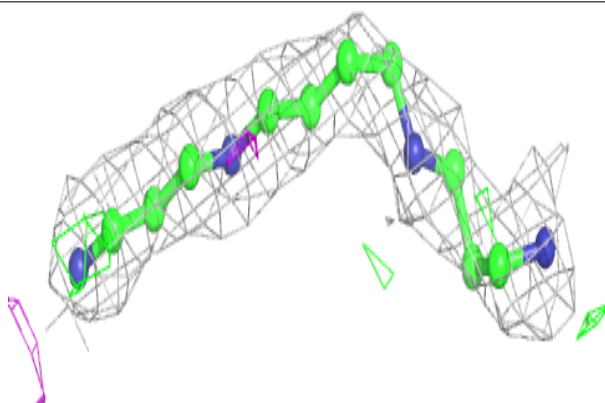


**Electron density around SPM J 201:**

$2mF_o - DF_c$  (at 0.7 rmsd) in gray  
 $mF_o - DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

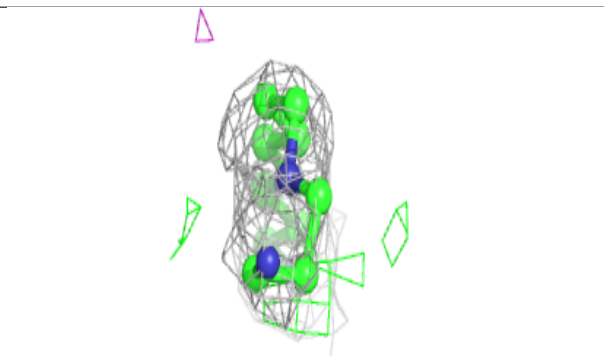
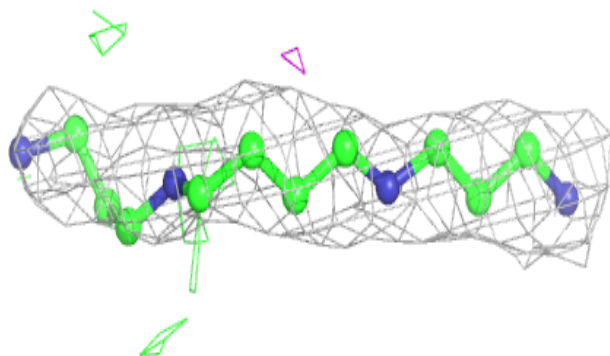
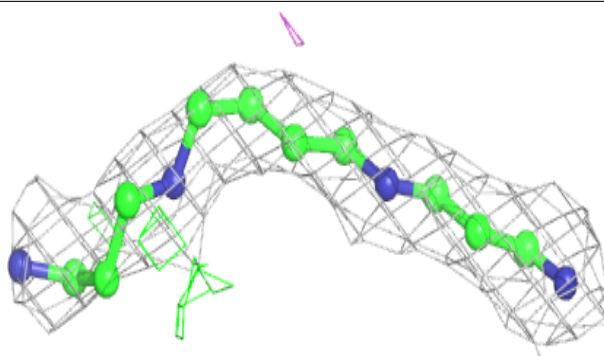
**Electron density around SPM E 202:**

$2mF_o - DF_c$  (at 0.7 rmsd) in gray  
 $mF_o - DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

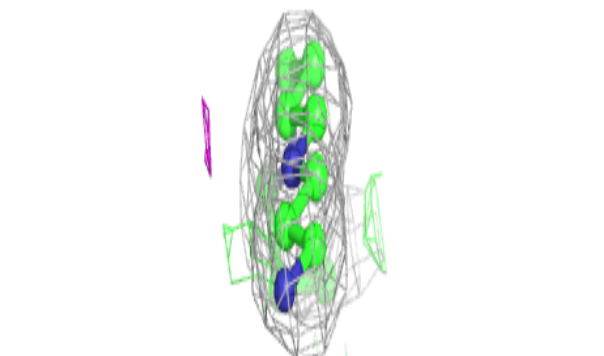
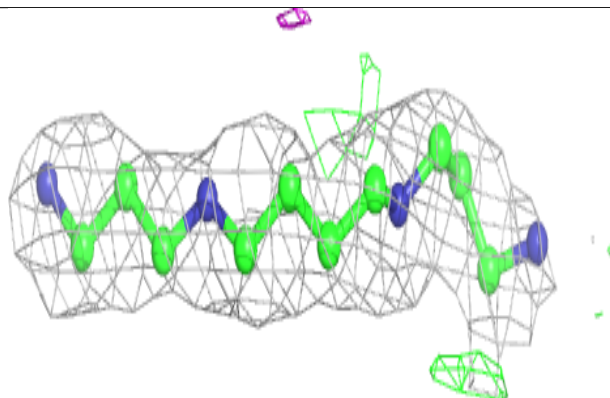
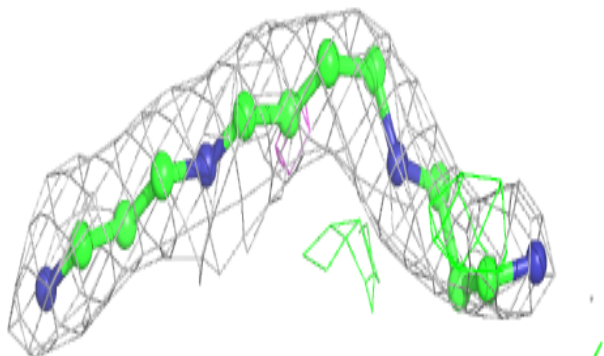


**Electron density around SPM F 201:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

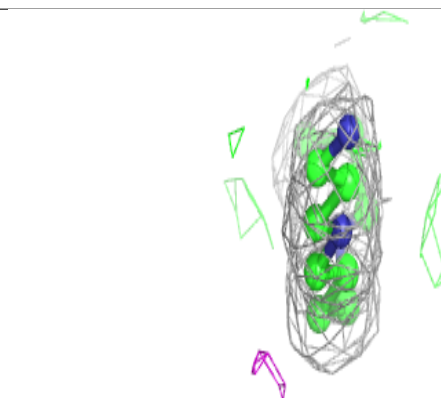
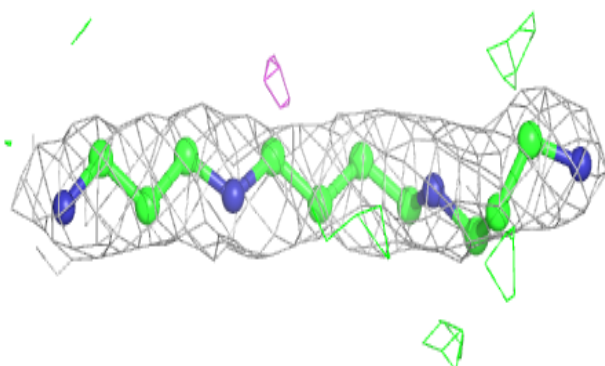
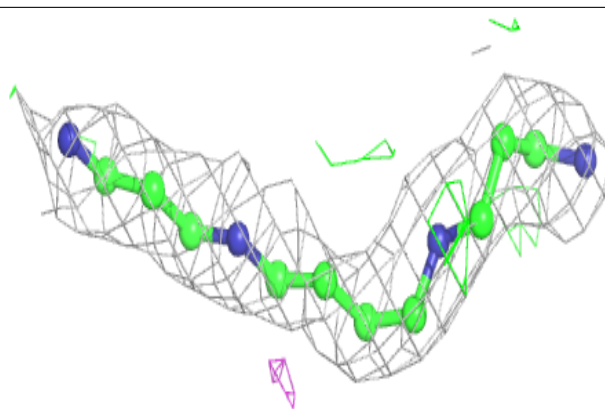
**Electron density around SPM D 201:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

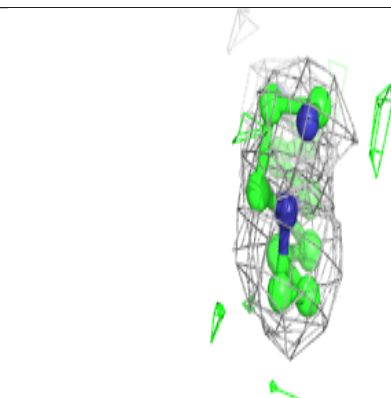
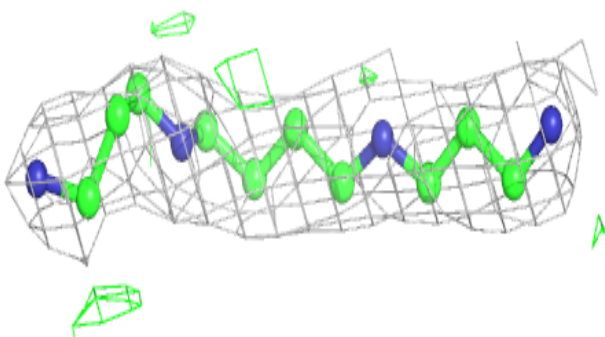
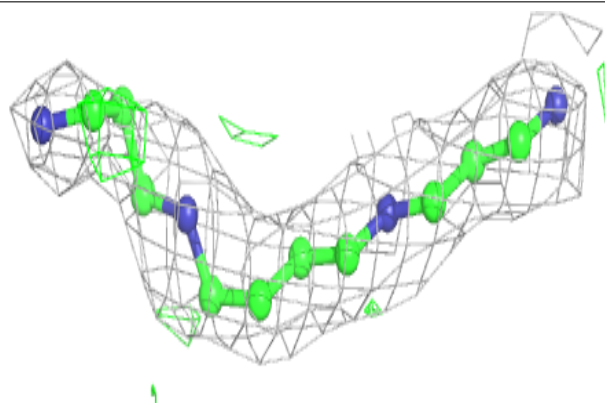


**Electron density around SPM K 201:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around SPM L 201:**

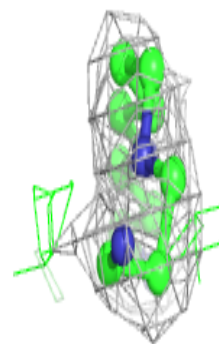
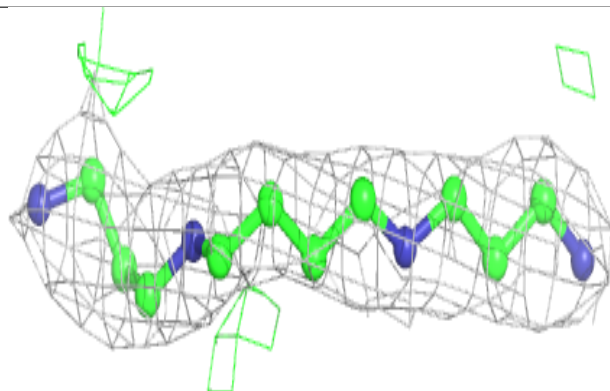
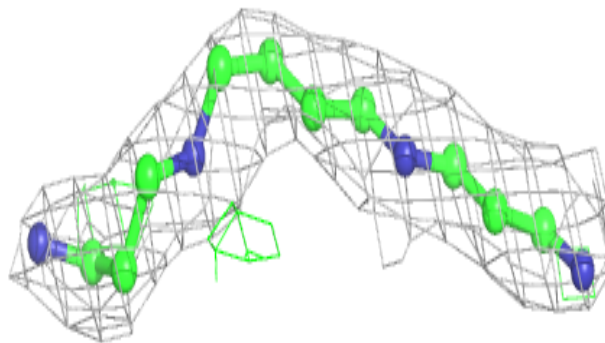
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



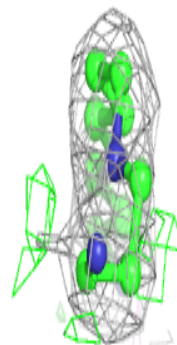
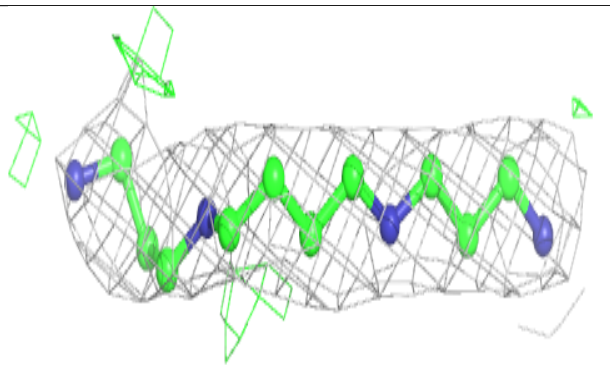
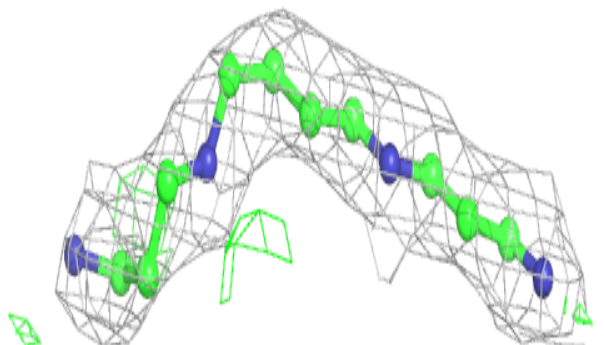


**Electron density around SPM G 201:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around SPM C 201:**

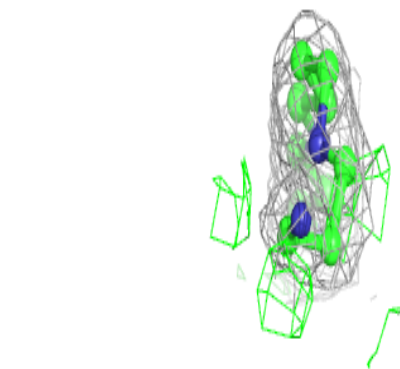
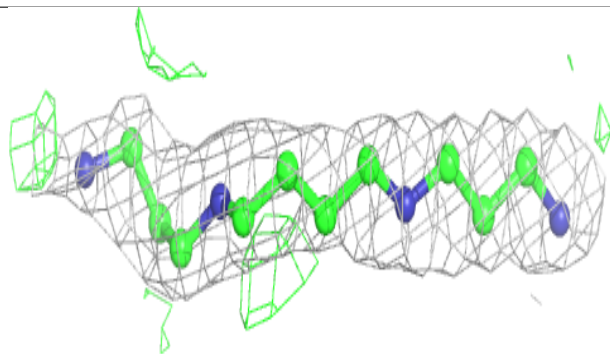
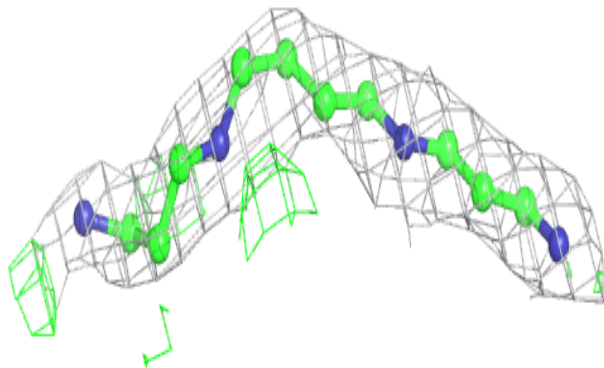
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



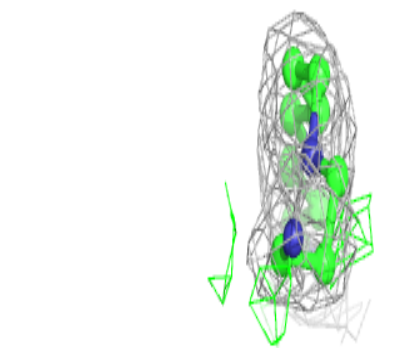
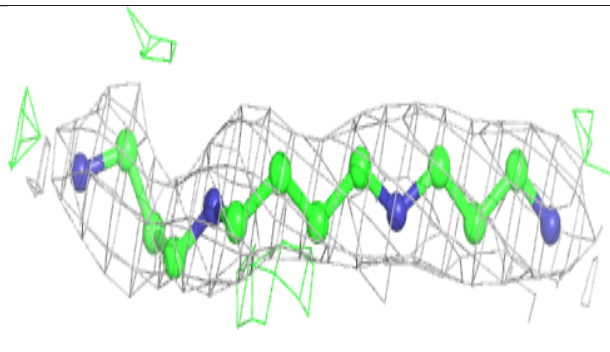
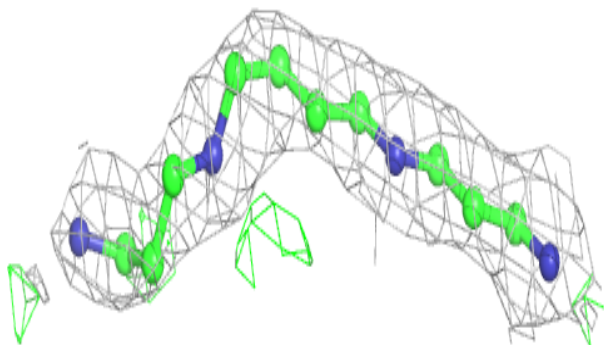


**Electron density around SPM B 201:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around SPM A 201:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



## 6.5 Other polymers

There are no such residues in this entry.