



# Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 02:27 AM GMT

PDB ID : 4DMU  
Title : Crystal structure of the von Willebrand factor A3 domain in complex with a collagen III derived triple-helical peptide  
Authors : Brondijk, T.H.C.; Huizinga, E.G.  
Deposited on : 2012-02-08  
Resolution : 2.80 Å(reported)

This is a full wwPDB 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 <http://wwpdb.org/ValidationPDFNotes.html>

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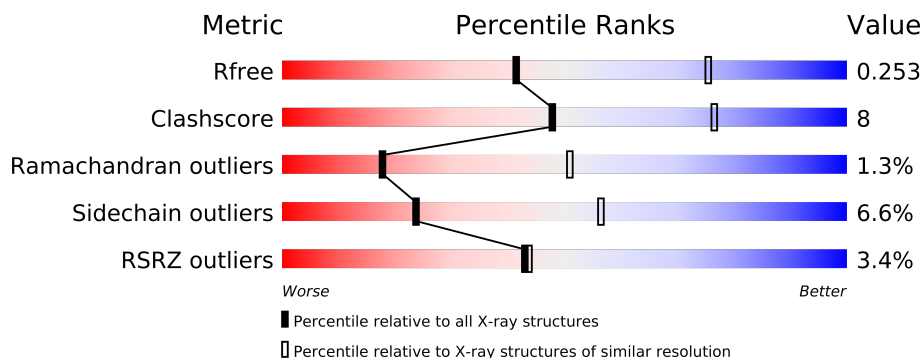
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 2.80 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	66092	1799 (2.80-2.80)
Clashscore	79885	2295 (2.80-2.80)
Ramachandran outliers	78287	2252 (2.80-2.80)
Sidechain outliers	78261	2254 (2.80-2.80)
RSRZ outliers	66119	1802 (2.80-2.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	87	
1	C	87	
1	E	87	
1	G	87	
1	I	87	
1	K	87	
2	B	192	
2	D	192	
2	F	192	
2	H	192	
2	J	192	
2	L	192	

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 11651 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Collagen III derived triple-helical peptide.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	79	Total	C	N	O	S	0	0	0
			529	332	91	103	3			
1	C	79	Total	C	N	O	S	0	0	0
			529	332	91	103	3			
1	E	79	Total	C	N	O	S	0	0	0
			529	332	91	103	3			
1	G	79	Total	C	N	O	S	0	0	0
			529	332	91	103	3			
1	I	79	Total	C	N	O	S	0	0	0
			529	332	91	103	3			
1	K	80	Total	C	N	O	S	0	0	0
			536	337	92	104	3			

- Molecule 2 is a protein called von Willebrand factor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	188	Total	C	N	O	S	0	0	0
			1406	890	241	269	6			
2	D	188	Total	C	N	O	S	0	0	0
			1406	890	241	269	6			
2	F	188	Total	C	N	O	S	0	0	0
			1406	890	241	269	6			
2	H	187	Total	C	N	O	S	0	0	0
			1400	887	240	268	5			
2	J	191	Total	C	N	O	S	0	0	0
			1426	902	244	274	6			
2	L	188	Total	C	N	O	S	0	0	0
			1406	890	241	269	6			

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	C	1	Total	O	S	0	0
			5	4	1		
3	F	1	Total	O	S	0	0
			5	4	1		
3	I	1	Total	O	S	0	0
			5	4	1		
3	K	1	Total	O	S	0	0
			5	4	1		

### 3 Residue-property plots i

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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 ( $\text{RSRZ} > 2$ ). 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: Collagen III derived triple-helical peptide

Chain A:



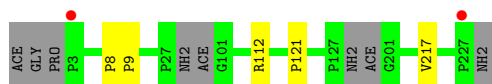
- Molecule 1: Collagen III derived triple-helical peptide

Chain C:



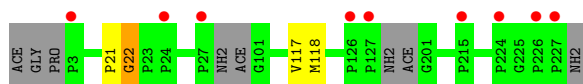
- Molecule 1: Collagen III derived triple-helical peptide

Chain E:



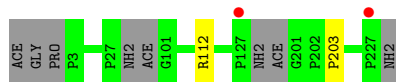
- Molecule 1: Collagen III derived triple-helical peptide

Chain G:



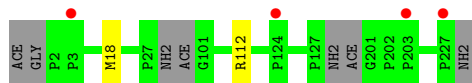
- Molecule 1: Collagen III derived triple-helical peptide

Chain I:



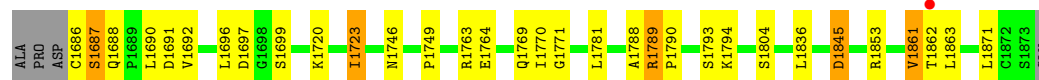
- Molecule 1: Collagen III derived triple-helical peptide

Chain K:



- Molecule 2: von Willebrand factor

Chain B:



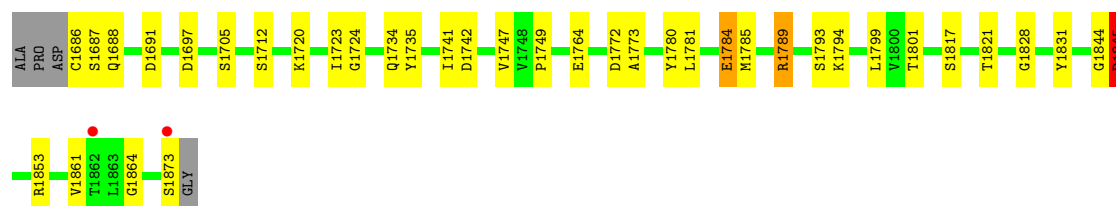
- Molecule 2: von Willebrand factor

Chain D:



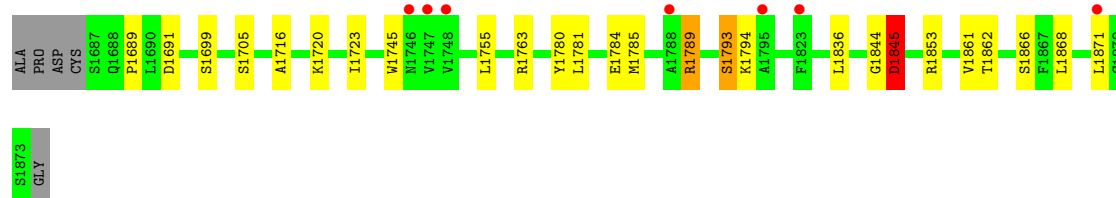
- Molecule 2: von Willebrand factor

Chain F:



- Molecule 2: von Willebrand factor

Chain H:



- Molecule 2: von Willebrand factor

Chain J:



- Molecule 2: von Willebrand factor

Chain L:



L1871
C1872
S1873
GLY

## 4 Data and refinement statistics

Property	Value	Source
Space group	P 6	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	187.61Å 187.61Å 89.37Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	46.32 – 2.80 46.32 – 2.80	Depositor EDS
% Data completeness (in resolution range)	92.2 (46.32-2.80) 92.4 (46.32-2.80)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.03 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.5.0110	Depositor
R, $R_{free}$	0.202 , 0.253 0.203 , 0.253	Depositor DCC
$R_{free}$ test set	2097 reflections (5.39%)	DCC
Wilson B-factor (Å <sup>2</sup> )	54.9	Xtriage
Anisotropy	0.112	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 8.5	EDS
Estimated twinning fraction	0.518 for H, K, L 0.482 for H+K, -K, -L 0.477 for h,-h-k,-l	Xtriage
Reported twinning fraction	0.518 for H, K, L 0.482 for H+K, -K, -L	Depositor
L-test for twinning	$\langle  L  \rangle = 0.43$ , $\langle L^2 \rangle = 0.26$	Xtriage
Outliers	2 of 41013 reflections (0.005%)	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	11651	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 33.33 % 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 8.1368e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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



## 5 Model quality

### 5.1 Standard geometry

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

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z  > 5$	RMSZ	# $ Z  > 5$
1	A	0.61	0/361	0.66	0/460
1	C	0.75	0/361	0.68	0/460
1	E	0.61	0/361	0.71	0/460
1	G	0.54	0/361	0.65	0/460
1	I	0.71	0/361	0.72	0/460
1	K	0.68	0/368	0.69	0/468
2	B	0.64	0/1431	0.73	0/1946
2	D	0.72	0/1431	0.79	1/1946 (0.1%)
2	F	0.64	0/1431	0.75	0/1946
2	H	0.57	0/1425	0.65	0/1938
2	J	0.88	0/1452	0.87	1/1976 (0.1%)
2	L	0.78	0/1431	0.80	0/1946
All	All	0.70	0/10774	0.75	2/14466 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	J	1789	ARG	NE-CZ-NH2	5.09	122.85	120.30
2	D	1692	VAL	CB-CA-C	-5.07	101.78	111.40

There are no chirality outliers.

There are no planarity outliers.

### 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit,

and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	529	0	0	1	0
1	C	529	0	0	4	0
1	E	529	0	0	3	0
1	G	529	0	0	3	0
1	I	529	0	0	2	0
1	K	536	0	0	2	0
2	B	1406	0	0	16	0
2	D	1406	0	0	17	0
2	F	1406	0	0	15	0
2	H	1400	0	0	10	0
2	J	1426	0	0	12	0
2	L	1406	0	0	16	0
3	C	5	0	0	0	0
3	F	5	0	0	0	0
3	I	5	0	0	0	0
3	K	5	0	0	0	0
All	All	11651	0	0	92	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 8.

All (92) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:D:1781:LEU:O	2:D:1794:LYS:NZ	2.18	0.75
2:H:1781:LEU:O	2:H:1794:LYS:NZ	2.22	0.72
2:J:1861:VAL:CG2	2:J:1862:THR:N	2.53	0.72
1:A:112:ARG:NH1	2:B:1764:GLU:OE2	2.25	0.69
2:D:1845:ASP:OD1	2:D:1845:ASP:N	2.28	0.66
2:D:1691:ASP:OD1	2:D:1789:ARG:NH2	2.29	0.65
2:F:1686:CYS:O	2:F:1688:GLN:N	2.30	0.65
1:C:112:ARG:NH1	2:D:1764:GLU:OE2	2.31	0.64
2:B:1691:ASP:OD1	2:B:1789:ARG:NH2	2.31	0.63
2:F:1781:LEU:O	2:F:1794:LYS:NZ	2.32	0.63
2:L:1861:VAL:CG2	2:L:1862:THR:N	2.64	0.61
2:B:1781:LEU:O	2:B:1794:LYS:NZ	2.34	0.61
2:F:1697:ASP:OD2	2:F:1801:THR:OG1	2.18	0.60
2:F:1845:ASP:N	2:F:1845:ASP:OD1	2.34	0.60
1:I:112:ARG:NH2	2:J:1764:GLU:OE2	2.35	0.59
2:B:1861:VAL:O	2:B:1863:LEU:N	2.35	0.59

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:J:1745:TRP:O	2:J:1747:VAL:N	2.35	0.59
2:J:1845:ASP:OD1	2:J:1845:ASP:N	2.36	0.58
2:L:1716:ALA:CB	2:L:1861:VAL:CG2	2.81	0.58
2:L:1691:ASP:OD1	2:L:1789:ARG:NH2	2.37	0.57
1:E:112:ARG:NH2	2:F:1764:GLU:OE2	2.37	0.57
2:L:1734:GLN:NE2	2:L:1742:ASP:OD2	2.38	0.57
2:B:1723:ILE:CD1	2:B:1723:ILE:N	2.68	0.56
2:B:1861:VAL:CG2	2:B:1862:THR:N	2.68	0.56
2:H:1745:TRP:CD1	2:H:1780:TYR:CZ	2.94	0.56
2:F:1691:ASP:OD1	2:F:1789:ARG:NH2	2.39	0.56
2:L:1871:LEU:N	2:L:1871:LEU:CD2	2.68	0.55
2:D:1735:TYR:OH	2:D:1799:LEU:O	2.25	0.55
2:H:1784:GLU:OE2	2:H:1784:GLU:N	2.40	0.55
2:H:1716:ALA:CB	2:H:1861:VAL:CG2	2.86	0.54
2:D:1804:SER:OG	2:D:1806:ASP:O	2.27	0.53
2:B:1686:CYS:O	2:B:1687:SER:CB	2.57	0.53
2:F:1741:ILE:CD1	2:F:1780:TYR:CE1	2.91	0.53
2:F:1828:GLY:O	2:F:1831:TYR:CE2	2.62	0.53
2:D:1716:ALA:CB	2:D:1861:VAL:CG2	2.88	0.52
2:F:1686:CYS:C	2:F:1688:GLN:N	2.64	0.51
1:K:112:ARG:NH1	2:L:1764:GLU:OE2	2.44	0.51
2:J:1776:PHE:CD1	2:J:1776:PHE:C	2.83	0.51
2:J:1861:VAL:O	2:J:1863:LEU:N	2.44	0.51
2:L:1686:CYS:O	2:L:1687:SER:CB	2.58	0.51
2:H:1691:ASP:OD1	2:H:1789:ARG:NH2	2.44	0.51
2:D:1861:VAL:CG2	2:D:1862:THR:N	2.75	0.49
2:H:1689:PRO:O	2:H:1793:SER:N	2.46	0.49
2:D:1853:ARG:NH1	2:D:1856:ASP:OD2	2.46	0.48
1:G:21:HYP:O	1:G:22:GLY:O	2.31	0.48
2:B:1690:LEU:CD2	2:B:1692:VAL:CG2	2.91	0.48
2:B:1845:ASP:N	2:B:1845:ASP:OD1	2.46	0.47
2:F:1844:GLY:O	2:F:1845:ASP:C	2.53	0.47
2:F:1735:TYR:OH	2:F:1799:LEU:O	2.33	0.47
2:D:1697:ASP:OD1	2:D:1699:SER:OG	2.33	0.47
2:F:1784:GLU:OE2	2:F:1784:GLU:N	2.48	0.47
2:L:1741:ILE:CD1	2:L:1780:TYR:CE1	2.98	0.46
2:J:1684:PRO:CB	2:J:1685:ASP:CB	2.93	0.46
2:J:1697:ASP:OD2	2:J:1801:THR:OG1	2.33	0.46
2:L:1784:GLU:N	2:L:1784:GLU:OE2	2.49	0.46
2:B:1696:LEU:CD2	2:B:1696:LEU:N	2.79	0.45
2:B:1769:GLN:OE1	2:D:1865:ASN:ND2	2.49	0.45
2:D:1697:ASP:OD2	2:D:1801:THR:OG1	2.35	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:F:1723:ILE:CG2	2:F:1724:GLY:N	2.81	0.45
2:J:1704:ALA:O	2:J:1705:SER:C	2.56	0.44
1:G:21:HYP:C	1:G:22:GLY:O	2.66	0.44
2:D:1808:VAL:O	2:D:1809:ASP:C	2.56	0.43
1:G:117:VAL:CG1	1:G:118:MET:N	2.80	0.43
2:H:1723:ILE:N	2:H:1723:ILE:CD1	2.82	0.43
1:C:117:VAL:CG1	1:C:118:MET:N	2.81	0.43
2:J:1784:GLU:OE2	2:J:1784:GLU:N	2.51	0.43
1:C:8:PRO:O	1:C:9:HYP:C	2.67	0.43
2:H:1836:LEU:N	2:H:1836:LEU:CD1	2.82	0.43
2:H:1844:GLY:O	2:H:1845:ASP:C	2.57	0.43
2:D:1776:PHE:C	2:D:1776:PHE:CD1	2.92	0.42
2:J:1744:PRO:O	2:J:1747:VAL:CG2	2.67	0.42
2:B:1770:ILE:O	2:B:1771:GLY:C	2.56	0.42
2:B:1781:LEU:O	2:B:1788:ALA:CB	2.67	0.42
2:B:1836:LEU:CD1	2:B:1836:LEU:N	2.82	0.42
2:D:1747:VAL:CG1	2:D:1753:HIS:CE1	3.02	0.42
1:C:18:MET:O	2:D:1786:HIS:NE2	2.53	0.42
1:E:8:PRO:O	1:E:9:HYP:C	2.68	0.42
2:F:1772:ASP:O	2:F:1773:ALA:C	2.57	0.42
2:L:1829:ASP:N	2:L:1829:ASP:OD2	2.53	0.42
2:L:1735:TYR:OH	2:L:1799:LEU:O	2.38	0.42
2:L:1861:VAL:O	2:L:1863:LEU:N	2.53	0.41
1:E:121:HYP:OD1	1:I:203:HYP:OD1	2.38	0.41
2:H:1862:THR:O	2:H:1862:THR:CG2	2.68	0.41
2:B:1861:VAL:C	2:B:1863:LEU:N	2.74	0.41
2:L:1866:SER:OG	2:L:1867:PHE:N	2.54	0.41
2:L:1774:LEU:O	2:L:1775:GLY:C	2.59	0.41
2:J:1691:ASP:OD1	2:J:1729:GLN:NE2	2.54	0.41
2:B:1697:ASP:OD1	2:B:1699:SER:N	2.54	0.41
2:D:1832:ASP:O	2:D:1835:GLN:N	2.54	0.41
1:K:18:MET:O	2:L:1786:HIS:NE2	2.54	0.41
2:L:1856:ASP:O	2:L:1857:LEU:C	2.60	0.41
2:F:1734:GLN:NE2	2:F:1742:ASP:OD2	2.55	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	56/87 (64%)	52 (93%)	4 (7%)	0	100	100
1	C	56/87 (64%)	51 (91%)	4 (7%)	1 (2%)	13	39
1	E	56/87 (64%)	53 (95%)	3 (5%)	0	100	100
1	G	56/87 (64%)	51 (91%)	4 (7%)	1 (2%)	13	39
1	I	56/87 (64%)	54 (96%)	2 (4%)	0	100	100
1	K	56/87 (64%)	53 (95%)	3 (5%)	0	100	100
2	B	186/192 (97%)	172 (92%)	9 (5%)	5 (3%)	8	25
2	D	186/192 (97%)	163 (88%)	21 (11%)	2 (1%)	21	57
2	F	186/192 (97%)	168 (90%)	14 (8%)	4 (2%)	10	32
2	H	185/192 (96%)	169 (91%)	15 (8%)	1 (0%)	38	76
2	J	189/192 (98%)	168 (89%)	19 (10%)	2 (1%)	21	57
2	L	186/192 (97%)	169 (91%)	14 (8%)	3 (2%)	14	44
All	All	1454/1674 (87%)	1323 (91%)	112 (8%)	19 (1%)	18	51

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	1687	SER
2	F	1687	SER
2	D	1687	SER
2	F	1864	GLY
1	G	22	GLY
2	J	1746	ASN
2	L	1687	SER
2	H	1845	ASP
2	B	1746	ASN
2	F	1845	ASP
2	J	1784	GLU
2	L	1746	ASN
2	B	1688	GLN

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Mol	Chain	Res	Type
2	B	1790	PRO
2	L	1862	THR
1	C	26	PRO
2	D	1725	PRO
2	B	1749	PRO
2	F	1749	PRO

### 5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	32/33 (97%)	32 (100%)	0	100	100
1	C	32/33 (97%)	32 (100%)	0	100	100
1	E	32/33 (97%)	31 (97%)	1 (3%)	52	86
1	G	32/33 (97%)	32 (100%)	0	100	100
1	I	32/33 (97%)	32 (100%)	0	100	100
1	K	33/33 (100%)	33 (100%)	0	100	100
2	B	155/157 (99%)	145 (94%)	10 (6%)	24	57
2	D	155/157 (99%)	144 (93%)	11 (7%)	21	51
2	F	155/157 (99%)	141 (91%)	14 (9%)	14	37
2	H	154/157 (98%)	141 (92%)	13 (8%)	16	41
2	J	157/157 (100%)	141 (90%)	16 (10%)	11	29
2	L	155/157 (99%)	146 (94%)	9 (6%)	28	63
All	All	1124/1140 (99%)	1050 (93%)	74 (7%)	24	56

All (74) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	1720	LYS
2	B	1723	ILE
2	B	1763	ARG
2	B	1789	ARG
2	B	1793	SER

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Mol	Chain	Res	Type
2	B	1804	SER
2	B	1845	ASP
2	B	1853	ARG
2	B	1861	VAL
2	B	1871	LEU
2	D	1705	SER
2	D	1720	LYS
2	D	1738	ILE
2	D	1747	VAL
2	D	1785	MET
2	D	1789	ARG
2	D	1793	SER
2	D	1845	ASP
2	D	1853	ARG
2	D	1866	SER
2	D	1871	LEU
1	E	217	VAL
2	F	1705	SER
2	F	1712	SER
2	F	1720	LYS
2	F	1747	VAL
2	F	1784	GLU
2	F	1785	MET
2	F	1789	ARG
2	F	1793	SER
2	F	1817	SER
2	F	1821	THR
2	F	1845	ASP
2	F	1853	ARG
2	F	1861	VAL
2	F	1873	SER
2	H	1699	SER
2	H	1705	SER
2	H	1720	LYS
2	H	1755	LEU
2	H	1763	ARG
2	H	1785	MET
2	H	1789	ARG
2	H	1793	SER
2	H	1845	ASP
2	H	1853	ARG
2	H	1866	SER

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Mol	Chain	Res	Type
2	H	1868	LEU
2	H	1871	LEU
2	J	1685	ASP
2	J	1686	CYS
2	J	1719	SER
2	J	1720	LYS
2	J	1738	ILE
2	J	1763	ARG
2	J	1776	PHE
2	J	1785	MET
2	J	1789	ARG
2	J	1793	SER
2	J	1845	ASP
2	J	1853	ARG
2	J	1861	VAL
2	J	1866	SER
2	J	1871	LEU
2	J	1873	SER
2	L	1705	SER
2	L	1720	LYS
2	L	1738	ILE
2	L	1763	ARG
2	L	1782	THR
2	L	1789	ARG
2	L	1793	SER
2	L	1845	ASP
2	L	1853	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

126 non-standard protein/DNA/RNA residues 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
1	HYP	A	103	1	8,8,9	5.97	2 (25%)	8,10,12	2.04	3 (37%)
1	HYP	A	106	1	8,8,9	6.42	2 (25%)	8,10,12	1.50	2 (25%)
1	HYP	A	109	1	8,8,9	5.91	2 (25%)	8,10,12	1.42	1 (12%)
1	HYP	A	115	1	8,8,9	6.87	2 (25%)	8,10,12	1.54	2 (25%)
1	HYP	A	121	1	8,8,9	6.61	2 (25%)	8,10,12	1.40	1 (12%)
1	HYP	A	124	1	8,8,9	6.53	2 (25%)	8,10,12	1.56	3 (37%)
1	HYP	A	127	1	8,8,9	6.63	2 (25%)	8,10,12	1.63	3 (37%)
1	HYP	A	15	1	8,8,9	7.59	2 (25%)	8,10,12	1.55	3 (37%)
1	HYP	A	203	1	8,8,9	6.25	2 (25%)	8,10,12	1.35	1 (12%)
1	HYP	A	206	1	8,8,9	7.23	2 (25%)	8,10,12	0.79	0
1	HYP	A	209	1	8,8,9	6.87	2 (25%)	8,10,12	1.33	1 (12%)
1	HYP	A	21	1	8,8,9	7.10	2 (25%)	8,10,12	1.51	1 (12%)
1	HYP	A	215	1	8,8,9	6.09	2 (25%)	8,10,12	1.42	2 (25%)
1	HYP	A	221	1	8,8,9	6.48	2 (25%)	8,10,12	1.27	2 (25%)
1	HYP	A	224	1	8,8,9	6.80	2 (25%)	8,10,12	1.23	1 (12%)
1	HYP	A	227	1	8,8,9	7.26	2 (25%)	8,10,12	1.66	3 (37%)
1	HYP	A	24	1	8,8,9	6.36	2 (25%)	8,10,12	1.37	3 (37%)
1	HYP	A	27	1	8,8,9	6.72	2 (25%)	8,10,12	1.20	1 (12%)
1	HYP	A	3	1	8,8,9	6.47	2 (25%)	8,10,12	1.85	3 (37%)
1	HYP	A	6	1	8,8,9	6.56	2 (25%)	8,10,12	1.36	1 (12%)
1	HYP	A	9	1	8,8,9	6.49	2 (25%)	8,10,12	1.14	1 (12%)
1	HYP	C	103	1	8,8,9	6.02	2 (25%)	8,10,12	1.43	1 (12%)
1	HYP	C	106	1	8,8,9	6.15	2 (25%)	8,10,12	1.29	0
1	HYP	C	109	1	8,8,9	6.57	2 (25%)	8,10,12	1.68	2 (25%)
1	HYP	C	115	1	8,8,9	6.36	2 (25%)	8,10,12	1.81	3 (37%)
1	HYP	C	121	1	8,8,9	6.07	2 (25%)	8,10,12	1.09	1 (12%)
1	HYP	C	124	1	8,8,9	6.62	2 (25%)	8,10,12	1.90	3 (37%)
1	HYP	C	127	1	8,8,9	7.33	2 (25%)	8,10,12	1.73	3 (37%)
1	HYP	C	15	1	8,8,9	6.56	3 (37%)	8,10,12	1.83	3 (37%)
1	HYP	C	203	1	8,8,9	6.09	1 (12%)	8,10,12	1.25	0
1	HYP	C	206	1	8,8,9	6.71	2 (25%)	8,10,12	2.39	2 (25%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
1	HYP	C	209	1	8,8,9	5.65	2 (25%)	8,10,12	2.06	5 (62%)
1	HYP	C	21	1	8,8,9	4.62	1 (12%)	8,10,12	1.63	2 (25%)
1	HYP	C	215	1	8,8,9	5.85	1 (12%)	8,10,12	0.97	0
1	HYP	C	221	1	8,8,9	6.85	2 (25%)	8,10,12	1.21	1 (12%)
1	HYP	C	224	1	8,8,9	6.29	2 (25%)	8,10,12	1.35	1 (12%)
1	HYP	C	227	1	8,8,9	6.55	2 (25%)	8,10,12	0.97	0
1	HYP	C	24	1	8,8,9	6.65	2 (25%)	8,10,12	1.60	3 (37%)
1	HYP	C	27	1	8,8,9	6.39	2 (25%)	8,10,12	1.56	1 (12%)
1	HYP	C	3	1	8,8,9	6.36	2 (25%)	8,10,12	2.44	5 (62%)
1	HYP	C	6	1	8,8,9	6.69	1 (12%)	8,10,12	1.31	1 (12%)
1	HYP	C	9	1	8,8,9	4.79	2 (25%)	8,10,12	1.35	2 (25%)
1	HYP	E	103	1	8,8,9	6.22	1 (12%)	8,10,12	1.10	0
1	HYP	E	106	1	8,8,9	6.39	2 (25%)	8,10,12	1.26	1 (12%)
1	HYP	E	109	1	8,8,9	6.42	1 (12%)	8,10,12	1.09	1 (12%)
1	HYP	E	115	1	8,8,9	5.86	1 (12%)	8,10,12	1.41	2 (25%)
1	HYP	E	121	1	8,8,9	7.05	2 (25%)	8,10,12	1.33	1 (12%)
1	HYP	E	124	1	8,8,9	6.62	2 (25%)	8,10,12	1.63	3 (37%)
1	HYP	E	127	1	8,8,9	7.38	2 (25%)	8,10,12	1.77	4 (50%)
1	HYP	E	15	1	8,8,9	6.04	2 (25%)	8,10,12	1.88	3 (37%)
1	HYP	E	203	1	8,8,9	5.97	1 (12%)	8,10,12	1.20	1 (12%)
1	HYP	E	206	1	8,8,9	5.41	2 (25%)	8,10,12	1.41	2 (25%)
1	HYP	E	209	1	8,8,9	4.92	1 (12%)	8,10,12	1.69	2 (25%)
1	HYP	E	21	1	8,8,9	6.30	1 (12%)	8,10,12	1.91	2 (25%)
1	HYP	E	215	1	8,8,9	7.06	2 (25%)	8,10,12	0.99	1 (12%)
1	HYP	E	221	1	8,8,9	5.93	1 (12%)	8,10,12	0.98	0
1	HYP	E	224	1	8,8,9	6.64	2 (25%)	8,10,12	0.90	0
1	HYP	E	227	1	8,8,9	6.29	1 (12%)	8,10,12	1.27	1 (12%)
1	HYP	E	24	1	8,8,9	6.63	2 (25%)	8,10,12	1.20	2 (25%)
1	HYP	E	27	1	8,8,9	6.17	2 (25%)	8,10,12	1.91	1 (12%)
1	HYP	E	3	1	8,8,9	6.04	2 (25%)	8,10,12	1.88	4 (50%)
1	HYP	E	6	1	8,8,9	6.83	2 (25%)	8,10,12	1.84	3 (37%)
1	HYP	E	9	1	8,8,9	5.72	2 (25%)	8,10,12	0.83	0
1	HYP	G	103	1	8,8,9	6.95	2 (25%)	8,10,12	1.53	2 (25%)
1	HYP	G	106	1	8,8,9	6.52	2 (25%)	8,10,12	1.20	1 (12%)
1	HYP	G	109	1	8,8,9	6.32	2 (25%)	8,10,12	0.93	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
1	HYP	G	115	1	8,8,9	6.55	2 (25%)	8,10,12	1.61	2 (25%)
1	HYP	G	121	1	8,8,9	6.67	2 (25%)	8,10,12	1.37	1 (12%)
1	HYP	G	124	1	8,8,9	6.45	2 (25%)	8,10,12	1.89	3 (37%)
1	HYP	G	127	1	8,8,9	8.16	2 (25%)	8,10,12	1.64	2 (25%)
1	HYP	G	15	1	8,8,9	6.22	1 (12%)	8,10,12	1.71	2 (25%)
1	HYP	G	203	1	8,8,9	6.20	1 (12%)	8,10,12	0.97	0
1	HYP	G	206	1	8,8,9	6.73	2 (25%)	8,10,12	1.40	1 (12%)
1	HYP	G	209	1	8,8,9	6.12	1 (12%)	8,10,12	1.47	2 (25%)
1	HYP	G	21	1	8,8,9	6.77	2 (25%)	8,10,12	1.54	2 (25%)
1	HYP	G	215	1	8,8,9	6.90	2 (25%)	8,10,12	1.10	1 (12%)
1	HYP	G	221	1	8,8,9	6.87	2 (25%)	8,10,12	1.61	2 (25%)
1	HYP	G	224	1	8,8,9	6.55	2 (25%)	8,10,12	1.29	1 (12%)
1	HYP	G	227	1	8,8,9	6.70	2 (25%)	8,10,12	1.64	3 (37%)
1	HYP	G	24	1	8,8,9	6.69	2 (25%)	8,10,12	1.44	1 (12%)
1	HYP	G	27	1	8,8,9	6.51	2 (25%)	8,10,12	1.31	2 (25%)
1	HYP	G	3	1	8,8,9	6.41	2 (25%)	8,10,12	1.96	3 (37%)
1	HYP	G	6	1	8,8,9	6.64	2 (25%)	8,10,12	1.64	2 (25%)
1	HYP	G	9	1	8,8,9	6.47	2 (25%)	8,10,12	1.57	2 (25%)
1	HYP	I	103	1	8,8,9	6.88	2 (25%)	8,10,12	1.63	3 (37%)
1	HYP	I	106	1	8,8,9	5.68	1 (12%)	8,10,12	1.13	1 (12%)
1	HYP	I	109	1	8,8,9	6.32	2 (25%)	8,10,12	1.46	2 (25%)
1	HYP	I	115	1	8,8,9	6.00	2 (25%)	8,10,12	1.65	2 (25%)
1	HYP	I	121	1	8,8,9	7.88	2 (25%)	8,10,12	1.08	0
1	HYP	I	124	1	8,8,9	6.05	2 (25%)	8,10,12	1.19	1 (12%)
1	HYP	I	127	1	8,8,9	7.54	2 (25%)	8,10,12	1.94	4 (50%)
1	HYP	I	15	1	8,8,9	5.76	1 (12%)	8,10,12	1.57	2 (25%)
1	HYP	I	203	1	8,8,9	6.60	2 (25%)	8,10,12	1.31	1 (12%)
1	HYP	I	206	1	8,8,9	6.84	2 (25%)	8,10,12	1.85	2 (25%)
1	HYP	I	209	1	8,8,9	6.48	2 (25%)	8,10,12	1.99	4 (50%)
1	HYP	I	21	1	8,8,9	5.00	1 (12%)	8,10,12	1.80	3 (37%)
1	HYP	I	215	1	8,8,9	7.17	2 (25%)	8,10,12	1.31	1 (12%)
1	HYP	I	221	1	8,8,9	6.56	2 (25%)	8,10,12	0.90	0
1	HYP	I	224	1	8,8,9	6.76	2 (25%)	8,10,12	1.53	2 (25%)
1	HYP	I	227	1	8,8,9	6.45	2 (25%)	8,10,12	1.37	1 (12%)
1	HYP	I	24	1	8,8,9	6.95	2 (25%)	8,10,12	1.91	3 (37%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
1	HYP	I	27	1	8,8,9	6.55	2 (25%)	8,10,12	1.84	1 (12%)
1	HYP	I	3	1	8,8,9	6.66	2 (25%)	8,10,12	1.96	2 (25%)
1	HYP	I	6	1	8,8,9	6.40	1 (12%)	8,10,12	0.99	0
1	HYP	I	9	1	8,8,9	6.70	2 (25%)	8,10,12	0.91	0
1	HYP	K	103	1	8,8,9	6.42	2 (25%)	8,10,12	1.48	1 (12%)
1	HYP	K	106	1	8,8,9	5.55	2 (25%)	8,10,12	1.64	2 (25%)
1	HYP	K	109	1	8,8,9	6.27	2 (25%)	8,10,12	1.16	0
1	HYP	K	115	1	8,8,9	5.99	2 (25%)	8,10,12	1.72	2 (25%)
1	HYP	K	121	1	8,8,9	6.72	1 (12%)	8,10,12	1.40	1 (12%)
1	HYP	K	124	1	8,8,9	6.59	2 (25%)	8,10,12	1.82	3 (37%)
1	HYP	K	127	1	8,8,9	7.22	2 (25%)	8,10,12	1.49	2 (25%)
1	HYP	K	15	1	8,8,9	6.15	2 (25%)	8,10,12	1.43	1 (12%)
1	HYP	K	203	1	8,8,9	6.13	1 (12%)	8,10,12	1.16	1 (12%)
1	HYP	K	206	1	8,8,9	6.34	2 (25%)	8,10,12	1.47	2 (25%)
1	HYP	K	209	1	8,8,9	5.65	3 (37%)	8,10,12	1.68	2 (25%)
1	HYP	K	21	1	8,8,9	6.77	2 (25%)	8,10,12	1.18	1 (12%)
1	HYP	K	215	1	8,8,9	6.88	2 (25%)	8,10,12	1.22	1 (12%)
1	HYP	K	221	1	8,8,9	6.44	2 (25%)	8,10,12	0.93	0
1	HYP	K	224	1	8,8,9	6.70	2 (25%)	8,10,12	1.06	0
1	HYP	K	227	1	8,8,9	7.10	2 (25%)	8,10,12	0.91	0
1	HYP	K	24	1	8,8,9	6.64	2 (25%)	8,10,12	1.67	2 (25%)
1	HYP	K	27	1	8,8,9	6.27	2 (25%)	8,10,12	1.63	3 (37%)
1	HYP	K	3	1	8,8,9	6.68	1 (12%)	8,10,12	1.14	1 (12%)
1	HYP	K	6	1	8,8,9	7.35	2 (25%)	8,10,12	1.45	1 (12%)
1	HYP	K	9	1	8,8,9	5.86	2 (25%)	8,10,12	1.88	3 (37%)

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
1	HYP	A	103	1	-	0/0/11/13	0/1/1/1
1	HYP	A	106	1	-	0/0/11/13	0/1/1/1
1	HYP	A	109	1	-	0/0/11/13	0/1/1/1
1	HYP	A	115	1	-	0/0/11/13	0/1/1/1
1	HYP	A	121	1	-	0/0/11/13	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	HYP	A	124	1	-	0/0/11/13	0/1/1/1
1	HYP	A	127	1	-	0/0/11/13	0/1/1/1
1	HYP	A	15	1	-	0/0/11/13	0/1/1/1
1	HYP	A	203	1	-	0/0/11/13	0/1/1/1
1	HYP	A	206	1	-	0/0/11/13	0/1/1/1
1	HYP	A	209	1	-	0/0/11/13	0/1/1/1
1	HYP	A	21	1	-	0/0/11/13	0/1/1/1
1	HYP	A	215	1	-	0/0/11/13	0/1/1/1
1	HYP	A	221	1	-	0/0/11/13	0/1/1/1
1	HYP	A	224	1	-	0/0/11/13	0/1/1/1
1	HYP	A	227	1	-	0/0/11/13	0/1/1/1
1	HYP	A	24	1	-	0/0/11/13	0/1/1/1
1	HYP	A	27	1	-	0/0/11/13	0/1/1/1
1	HYP	A	3	1	-	0/0/11/13	0/1/1/1
1	HYP	A	6	1	-	0/0/11/13	0/1/1/1
1	HYP	A	9	1	-	0/0/11/13	0/1/1/1
1	HYP	C	103	1	-	0/0/11/13	0/1/1/1
1	HYP	C	106	1	-	0/0/11/13	0/1/1/1
1	HYP	C	109	1	-	0/0/11/13	0/1/1/1
1	HYP	C	115	1	-	0/0/11/13	0/1/1/1
1	HYP	C	121	1	-	0/0/11/13	0/1/1/1
1	HYP	C	124	1	-	0/0/11/13	0/1/1/1
1	HYP	C	127	1	-	0/0/11/13	0/1/1/1
1	HYP	C	15	1	-	0/0/11/13	0/1/1/1
1	HYP	C	203	1	-	0/0/11/13	0/1/1/1
1	HYP	C	206	1	-	0/0/11/13	0/1/1/1
1	HYP	C	209	1	-	0/0/11/13	0/1/1/1
1	HYP	C	21	1	-	0/0/11/13	0/1/1/1
1	HYP	C	215	1	-	0/0/11/13	0/1/1/1
1	HYP	C	221	1	-	0/0/11/13	0/1/1/1
1	HYP	C	224	1	-	0/0/11/13	0/1/1/1
1	HYP	C	227	1	-	0/0/11/13	0/1/1/1
1	HYP	C	24	1	-	0/0/11/13	0/1/1/1
1	HYP	C	27	1	-	0/0/11/13	0/1/1/1
1	HYP	C	3	1	-	0/0/11/13	0/1/1/1
1	HYP	C	6	1	-	0/0/11/13	0/1/1/1
1	HYP	C	9	1	-	0/0/11/13	0/1/1/1
1	HYP	E	103	1	-	0/0/11/13	0/1/1/1
1	HYP	E	106	1	-	0/0/11/13	0/1/1/1
1	HYP	E	109	1	-	0/0/11/13	0/1/1/1
1	HYP	E	115	1	-	0/0/11/13	0/1/1/1
1	HYP	E	121	1	-	0/0/11/13	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	HYP	E	124	1	-	0/0/11/13	0/1/1/1
1	HYP	E	127	1	-	0/0/11/13	0/1/1/1
1	HYP	E	15	1	-	0/0/11/13	0/1/1/1
1	HYP	E	203	1	-	0/0/11/13	0/1/1/1
1	HYP	E	206	1	-	0/0/11/13	0/1/1/1
1	HYP	E	209	1	-	0/0/11/13	0/1/1/1
1	HYP	E	21	1	-	0/0/11/13	0/1/1/1
1	HYP	E	215	1	-	0/0/11/13	0/1/1/1
1	HYP	E	221	1	-	0/0/11/13	0/1/1/1
1	HYP	E	224	1	-	0/0/11/13	0/1/1/1
1	HYP	E	227	1	-	0/0/11/13	0/1/1/1
1	HYP	E	24	1	-	0/0/11/13	0/1/1/1
1	HYP	E	27	1	-	0/0/11/13	0/1/1/1
1	HYP	E	3	1	-	0/0/11/13	0/1/1/1
1	HYP	E	6	1	-	0/0/11/13	0/1/1/1
1	HYP	E	9	1	-	0/0/11/13	0/1/1/1
1	HYP	G	103	1	-	0/0/11/13	0/1/1/1
1	HYP	G	106	1	-	0/0/11/13	0/1/1/1
1	HYP	G	109	1	-	0/0/11/13	0/1/1/1
1	HYP	G	115	1	-	0/0/11/13	0/1/1/1
1	HYP	G	121	1	-	0/0/11/13	0/1/1/1
1	HYP	G	124	1	-	0/0/11/13	0/1/1/1
1	HYP	G	127	1	-	0/0/11/13	0/1/1/1
1	HYP	G	15	1	-	0/0/11/13	0/1/1/1
1	HYP	G	203	1	-	0/0/11/13	0/1/1/1
1	HYP	G	206	1	-	0/0/11/13	0/1/1/1
1	HYP	G	209	1	-	0/0/11/13	0/1/1/1
1	HYP	G	21	1	-	0/0/11/13	0/1/1/1
1	HYP	G	215	1	-	0/0/11/13	0/1/1/1
1	HYP	G	221	1	-	0/0/11/13	0/1/1/1
1	HYP	G	224	1	-	0/0/11/13	0/1/1/1
1	HYP	G	227	1	-	0/0/11/13	0/1/1/1
1	HYP	G	24	1	-	0/0/11/13	0/1/1/1
1	HYP	G	27	1	-	0/0/11/13	0/1/1/1
1	HYP	G	3	1	-	0/0/11/13	0/1/1/1
1	HYP	G	6	1	-	0/0/11/13	0/1/1/1
1	HYP	G	9	1	-	0/0/11/13	0/1/1/1
1	HYP	I	103	1	-	0/0/11/13	0/1/1/1
1	HYP	I	106	1	-	0/0/11/13	0/1/1/1
1	HYP	I	109	1	-	0/0/11/13	0/1/1/1
1	HYP	I	115	1	-	0/0/11/13	0/1/1/1
1	HYP	I	121	1	-	0/0/11/13	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	HYP	I	124	1	-	0/0/11/13	0/1/1/1
1	HYP	I	127	1	-	0/0/11/13	0/1/1/1
1	HYP	I	15	1	-	0/0/11/13	0/1/1/1
1	HYP	I	203	1	-	0/0/11/13	0/1/1/1
1	HYP	I	206	1	-	0/0/11/13	0/1/1/1
1	HYP	I	209	1	-	0/0/11/13	0/1/1/1
1	HYP	I	21	1	-	0/0/11/13	0/1/1/1
1	HYP	I	215	1	-	0/0/11/13	0/1/1/1
1	HYP	I	221	1	-	0/0/11/13	0/1/1/1
1	HYP	I	224	1	-	0/0/11/13	0/1/1/1
1	HYP	I	227	1	-	0/0/11/13	0/1/1/1
1	HYP	I	24	1	-	0/0/11/13	0/1/1/1
1	HYP	I	27	1	-	0/0/11/13	0/1/1/1
1	HYP	I	3	1	-	0/0/11/13	0/1/1/1
1	HYP	I	6	1	-	0/0/11/13	0/1/1/1
1	HYP	I	9	1	-	0/0/11/13	0/1/1/1
1	HYP	K	103	1	-	0/0/11/13	0/1/1/1
1	HYP	K	106	1	-	0/0/11/13	0/1/1/1
1	HYP	K	109	1	-	0/0/11/13	0/1/1/1
1	HYP	K	115	1	-	0/0/11/13	0/1/1/1
1	HYP	K	121	1	-	0/0/11/13	0/1/1/1
1	HYP	K	124	1	-	0/0/11/13	0/1/1/1
1	HYP	K	127	1	-	0/0/11/13	0/1/1/1
1	HYP	K	15	1	-	0/0/11/13	0/1/1/1
1	HYP	K	203	1	-	0/0/11/13	0/1/1/1
1	HYP	K	206	1	-	0/0/11/13	0/1/1/1
1	HYP	K	209	1	-	0/0/11/13	0/1/1/1
1	HYP	K	21	1	-	0/0/11/13	0/1/1/1
1	HYP	K	215	1	-	0/0/11/13	0/1/1/1
1	HYP	K	221	1	-	0/0/11/13	0/1/1/1
1	HYP	K	224	1	-	0/0/11/13	0/1/1/1
1	HYP	K	227	1	-	0/0/11/13	0/1/1/1
1	HYP	K	24	1	-	0/0/11/13	0/1/1/1
1	HYP	K	27	1	-	0/0/11/13	0/1/1/1
1	HYP	K	3	1	-	0/0/11/13	0/1/1/1
1	HYP	K	6	1	-	0/0/11/13	0/1/1/1
1	HYP	K	9	1	-	0/0/11/13	0/1/1/1

All (232) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	G	127	HYP	O-C	22.78	1.27	1.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	I	121	HYP	O-C	22.03	1.26	1.11
1	A	15	HYP	O-C	21.22	1.26	1.11
1	I	127	HYP	O-C	21.06	1.25	1.11
1	E	127	HYP	O-C	20.61	1.25	1.11
1	K	6	HYP	O-C	20.52	1.25	1.11
1	C	127	HYP	O-C	20.49	1.25	1.11
1	A	206	HYP	O-C	20.29	1.25	1.11
1	A	227	HYP	O-C	20.17	1.25	1.11
1	K	127	HYP	O-C	20.13	1.25	1.11
1	I	215	HYP	O-C	19.92	1.25	1.11
1	K	227	HYP	O-C	19.85	1.25	1.11
1	A	21	HYP	O-C	19.75	1.25	1.11
1	E	215	HYP	O-C	19.70	1.25	1.11
1	E	121	HYP	O-C	19.69	1.25	1.11
1	G	103	HYP	O-C	19.47	1.24	1.11
1	G	215	HYP	O-C	19.34	1.24	1.11
1	I	24	HYP	O-C	19.25	1.24	1.11
1	I	103	HYP	O-C	19.23	1.24	1.11
1	A	209	HYP	O-C	19.20	1.24	1.11
1	K	215	HYP	O-C	19.18	1.24	1.11
1	C	221	HYP	O-C	19.18	1.24	1.11
1	G	221	HYP	O-C	19.16	1.24	1.11
1	A	115	HYP	O-C	19.12	1.24	1.11
1	I	206	HYP	O-C	19.03	1.24	1.11
1	E	6	HYP	O-C	19.02	1.24	1.11
1	A	224	HYP	O-C	18.99	1.24	1.11
1	G	21	HYP	O-C	18.94	1.24	1.11
1	I	224	HYP	O-C	18.87	1.24	1.11
1	K	121	HYP	O-C	18.86	1.24	1.11
1	K	21	HYP	O-C	18.86	1.24	1.11
1	C	206	HYP	O-C	18.82	1.24	1.11
1	G	206	HYP	O-C	18.78	1.24	1.11
1	A	27	HYP	O-C	18.77	1.24	1.11
1	I	9	HYP	O-C	18.76	1.24	1.11
1	K	224	HYP	O-C	18.75	1.24	1.11
1	K	3	HYP	O-C	18.75	1.24	1.11
1	C	6	HYP	O-C	18.74	1.24	1.11
1	G	227	HYP	O-C	18.69	1.24	1.11
1	G	121	HYP	O-C	18.65	1.24	1.11
1	G	24	HYP	O-C	18.63	1.24	1.11
1	E	224	HYP	O-C	18.61	1.24	1.11
1	C	24	HYP	O-C	18.55	1.24	1.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	I	3	HYP	O-C	18.51	1.24	1.11
1	A	121	HYP	O-C	18.48	1.24	1.11
1	A	127	HYP	O-C	18.47	1.24	1.11
1	K	24	HYP	O-C	18.45	1.24	1.11
1	E	124	HYP	O-C	18.45	1.24	1.11
1	E	24	HYP	O-C	18.44	1.24	1.11
1	I	203	HYP	O-C	18.44	1.24	1.11
1	K	124	HYP	O-C	18.42	1.24	1.11
1	C	124	HYP	O-C	18.42	1.24	1.11
1	G	6	HYP	O-C	18.39	1.24	1.11
1	I	221	HYP	O-C	18.35	1.24	1.11
1	C	227	HYP	O-C	18.30	1.24	1.11
1	C	109	HYP	O-C	18.29	1.24	1.11
1	I	27	HYP	O-C	18.25	1.24	1.11
1	A	6	HYP	O-C	18.24	1.24	1.11
1	A	124	HYP	O-C	18.21	1.23	1.11
1	G	115	HYP	O-C	18.20	1.23	1.11
1	C	15	HYP	O-C	18.20	1.23	1.11
1	G	224	HYP	O-C	18.20	1.23	1.11
1	A	3	HYP	O-C	18.16	1.23	1.11
1	G	106	HYP	O-C	18.15	1.23	1.11
1	G	9	HYP	O-C	18.13	1.23	1.11
1	G	27	HYP	O-C	18.12	1.23	1.11
1	A	221	HYP	O-C	18.08	1.23	1.11
1	E	109	HYP	O-C	18.06	1.23	1.11
1	A	9	HYP	O-C	18.05	1.23	1.11
1	I	227	HYP	O-C	18.04	1.23	1.11
1	K	221	HYP	O-C	18.01	1.23	1.11
1	G	124	HYP	O-C	18.00	1.23	1.11
1	K	103	HYP	O-C	17.94	1.23	1.11
1	G	3	HYP	O-C	17.92	1.23	1.11
1	I	209	HYP	O-C	17.89	1.23	1.11
1	A	106	HYP	O-C	17.89	1.23	1.11
1	I	6	HYP	O-C	17.89	1.23	1.11
1	E	106	HYP	O-C	17.83	1.23	1.11
1	C	27	HYP	O-C	17.82	1.23	1.11
1	C	3	HYP	O-C	17.74	1.23	1.11
1	A	24	HYP	O-C	17.71	1.23	1.11
1	K	206	HYP	O-C	17.69	1.23	1.11
1	G	109	HYP	O-C	17.69	1.23	1.11
1	C	115	HYP	O-C	17.65	1.23	1.11
1	E	227	HYP	O-C	17.64	1.23	1.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	I	109	HYP	O-C	17.63	1.23	1.11
1	E	21	HYP	O-C	17.57	1.23	1.11
1	C	224	HYP	O-C	17.57	1.23	1.11
1	K	109	HYP	O-C	17.54	1.23	1.11
1	K	27	HYP	O-C	17.48	1.23	1.11
1	E	103	HYP	O-C	17.44	1.23	1.11
1	G	15	HYP	O-C	17.41	1.23	1.11
1	A	203	HYP	O-C	17.40	1.23	1.11
1	G	203	HYP	O-C	17.38	1.23	1.11
1	E	27	HYP	O-C	17.21	1.23	1.11
1	K	203	HYP	O-C	17.17	1.23	1.11
1	G	209	HYP	O-C	17.15	1.23	1.11
1	C	203	HYP	O-C	17.13	1.23	1.11
1	C	106	HYP	O-C	17.12	1.23	1.11
1	K	15	HYP	O-C	17.09	1.23	1.11
1	A	215	HYP	O-C	17.01	1.23	1.11
1	C	121	HYP	O-C	16.99	1.23	1.11
1	I	124	HYP	O-C	16.88	1.23	1.11
1	E	3	HYP	O-C	16.81	1.23	1.11
1	E	15	HYP	O-C	16.80	1.23	1.11
1	C	103	HYP	O-C	16.79	1.22	1.11
1	E	203	HYP	O-C	16.70	1.22	1.11
1	K	115	HYP	O-C	16.68	1.22	1.11
1	E	221	HYP	O-C	16.62	1.22	1.11
1	A	103	HYP	O-C	16.53	1.22	1.11
1	A	109	HYP	O-C	16.53	1.22	1.11
1	I	115	HYP	O-C	16.42	1.22	1.11
1	C	215	HYP	O-C	16.39	1.22	1.11
1	E	115	HYP	O-C	16.35	1.22	1.11
1	K	9	HYP	O-C	16.22	1.22	1.11
1	I	15	HYP	O-C	16.01	1.22	1.11
1	E	9	HYP	O-C	15.99	1.22	1.11
1	I	106	HYP	O-C	15.94	1.22	1.11
1	C	209	HYP	O-C	15.66	1.22	1.11
1	K	209	HYP	O-C	15.55	1.22	1.11
1	K	106	HYP	O-C	15.42	1.22	1.11
1	E	206	HYP	O-C	15.09	1.21	1.11
1	I	21	HYP	O-C	13.81	1.20	1.11
1	E	209	HYP	O-C	13.75	1.20	1.11
1	C	9	HYP	O-C	13.23	1.20	1.11
1	C	21	HYP	O-C	12.96	1.20	1.11
1	I	24	HYP	CA-C	3.86	1.55	1.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	G	6	HYP	CA-C	3.65	1.55	1.48
1	A	227	HYP	CA-C	3.60	1.55	1.48
1	G	127	HYP	CA-C	3.47	1.54	1.48
1	I	209	HYP	CA-C	3.45	1.54	1.48
1	A	21	HYP	CA-C	3.35	1.54	1.48
1	K	127	HYP	CA-C	3.33	1.54	1.48
1	I	3	HYP	CA-C	3.32	1.54	1.48
1	I	127	HYP	CA-C	3.31	1.54	1.48
1	A	9	HYP	CA-C	3.29	1.54	1.48
1	A	103	HYP	CA-C	3.27	1.54	1.48
1	I	215	HYP	CA-C	3.27	1.54	1.48
1	E	24	HYP	CA-C	3.26	1.54	1.48
1	G	224	HYP	CA-C	3.26	1.54	1.48
1	G	115	HYP	CA-C	3.25	1.54	1.48
1	K	24	HYP	CA-C	3.24	1.54	1.48
1	C	124	HYP	CA-C	3.23	1.54	1.48
1	I	115	HYP	CA-C	3.21	1.54	1.48
1	A	127	HYP	CA-C	3.13	1.54	1.48
1	G	27	HYP	CA-C	3.08	1.54	1.48
1	E	127	HYP	CA-C	3.07	1.54	1.48
1	I	206	HYP	CA-C	3.06	1.54	1.48
1	C	109	HYP	CA-C	3.05	1.54	1.48
1	G	24	HYP	CA-C	3.05	1.54	1.48
1	A	115	HYP	CA-C	3.03	1.54	1.48
1	G	221	HYP	CA-C	3.02	1.54	1.48
1	I	224	HYP	CA-C	3.01	1.54	1.48
1	C	127	HYP	CA-C	3.01	1.54	1.48
1	A	6	HYP	CA-C	3.00	1.53	1.48
1	G	227	HYP	CA-C	3.00	1.53	1.48
1	G	206	HYP	CA-C	2.98	1.53	1.48
1	K	6	HYP	CA-C	2.94	1.53	1.48
1	E	124	HYP	CA-C	2.92	1.53	1.48
1	A	24	HYP	CA-C	2.91	1.53	1.48
1	C	24	HYP	CA-C	2.91	1.53	1.48
1	G	106	HYP	CA-C	2.91	1.53	1.48
1	I	121	HYP	CA-C	2.90	1.53	1.48
1	I	27	HYP	CA-C	2.90	1.53	1.48
1	K	227	HYP	CA-C	2.87	1.53	1.48
1	A	124	HYP	CA-C	2.85	1.53	1.48
1	A	221	HYP	CA-C	2.84	1.53	1.48
1	K	9	HYP	CA-C	2.83	1.53	1.48
1	G	124	HYP	CA-C	2.81	1.53	1.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	6	HYP	CA-C	2.80	1.53	1.48
1	E	3	HYP	CA-C	2.80	1.53	1.48
1	A	27	HYP	CA-C	2.79	1.53	1.48
1	K	27	HYP	CA-C	2.78	1.53	1.48
1	A	209	HYP	CA-C	2.77	1.53	1.48
1	A	224	HYP	CA-C	2.76	1.53	1.48
1	C	115	HYP	CA-C	2.75	1.53	1.48
1	A	106	HYP	CA-C	2.73	1.53	1.48
1	K	15	HYP	CA-C	2.73	1.53	1.48
1	A	203	HYP	CA-C	2.70	1.53	1.48
1	C	227	HYP	CA-C	2.70	1.53	1.48
1	C	221	HYP	CA-C	2.66	1.53	1.48
1	A	15	HYP	CA-C	2.66	1.53	1.48
1	C	27	HYP	CA-C	2.65	1.53	1.48
1	I	103	HYP	CA-C	2.65	1.53	1.48
1	K	124	HYP	CA-C	2.65	1.53	1.48
1	G	21	HYP	CA-C	2.64	1.53	1.48
1	K	209	HYP	CA-C	2.63	1.53	1.48
1	G	121	HYP	CA-C	2.61	1.53	1.48
1	E	215	HYP	CA-C	2.59	1.53	1.48
1	C	9	HYP	CA-C	2.58	1.53	1.48
1	E	27	HYP	CA-C	2.56	1.53	1.48
1	I	221	HYP	CA-C	2.55	1.53	1.48
1	C	106	HYP	CA-C	2.53	1.53	1.48
1	K	21	HYP	CA-C	2.52	1.53	1.48
1	C	3	HYP	CA-C	2.45	1.52	1.48
1	I	124	HYP	CA-C	2.45	1.52	1.48
1	I	203	HYP	CA-C	2.45	1.52	1.48
1	G	3	HYP	CA-C	2.44	1.52	1.48
1	I	227	HYP	CA-C	2.44	1.52	1.48
1	I	9	HYP	CA-C	2.42	1.52	1.48
1	I	109	HYP	CA-C	2.41	1.52	1.48
1	K	106	HYP	CA-C	2.40	1.52	1.48
1	C	224	HYP	CA-C	2.40	1.52	1.48
1	K	221	HYP	CA-C	2.38	1.52	1.48
1	G	103	HYP	CA-C	2.38	1.52	1.48
1	E	15	HYP	CA-C	2.38	1.52	1.48
1	C	121	HYP	CA-C	2.37	1.52	1.48
1	A	121	HYP	CA-C	2.37	1.52	1.48
1	K	115	HYP	CA-C	2.36	1.52	1.48
1	K	206	HYP	CA-C	2.35	1.52	1.48
1	K	224	HYP	CA-C	2.34	1.52	1.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	206	HYP	CA-C	2.33	1.52	1.48
1	C	15	HYP	CA-C	2.32	1.52	1.48
1	K	103	HYP	CA-C	2.32	1.52	1.48
1	E	121	HYP	CA-C	2.32	1.52	1.48
1	A	109	HYP	CA-C	2.30	1.52	1.48
1	E	106	HYP	CA-C	2.29	1.52	1.48
1	K	109	HYP	CA-C	2.25	1.52	1.48
1	C	15	HYP	CB-CA	-2.23	1.48	1.55
1	G	9	HYP	CA-C	2.21	1.52	1.48
1	K	215	HYP	CA-C	2.20	1.52	1.48
1	G	109	HYP	CA-C	2.19	1.52	1.48
1	E	206	HYP	CA-C	2.18	1.52	1.48
1	C	209	HYP	CA-C	2.18	1.52	1.48
1	K	209	HYP	CB-CA	-2.15	1.49	1.55
1	A	3	HYP	CA-C	2.12	1.52	1.48
1	G	215	HYP	CA-C	2.10	1.52	1.48
1	E	224	HYP	CA-C	2.08	1.52	1.48
1	A	215	HYP	CA-C	2.08	1.52	1.48
1	C	103	HYP	CA-C	2.06	1.52	1.48
1	C	206	HYP	CA-C	2.05	1.52	1.48
1	E	9	HYP	CA-C	2.02	1.52	1.48

All (210) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	206	HYP	CG-CD-N	-5.36	96.84	105.47
1	E	27	HYP	CG-CD-N	-4.14	98.80	105.47
1	I	27	HYP	CG-CD-N	-3.81	99.33	105.47
1	I	206	HYP	CG-CD-N	-3.78	99.38	105.47
1	C	27	HYP	CG-CD-N	-3.72	99.49	105.47
1	C	15	HYP	CG-CD-N	-3.65	99.59	105.47
1	I	3	HYP	CG-CD-N	-3.65	99.60	105.47
1	A	103	HYP	CD-N-CA	3.61	114.97	108.15
1	K	124	HYP	CD-N-CA	3.47	114.69	108.15
1	C	124	HYP	CD-N-CA	3.44	114.64	108.15
1	C	3	HYP	C-CA-N	-3.43	104.52	110.86
1	C	206	HYP	CB-CG-CD	-3.43	98.59	103.46
1	G	124	HYP	CD-N-CA	3.39	114.55	108.15
1	E	21	HYP	CG-CD-N	-3.38	100.02	105.47
1	K	106	HYP	C-CA-N	3.34	117.04	110.86
1	C	115	HYP	CD-N-CA	3.28	114.35	108.15
1	A	121	HYP	CG-CD-N	-3.27	100.20	105.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	109	HYP	CG-CD-N	-3.21	100.31	105.47
1	I	209	HYP	CD-N-CA	3.20	114.19	108.15
1	E	6	HYP	C-CA-N	3.18	116.74	110.86
1	G	124	HYP	CG-CB-CA	3.14	108.81	104.22
1	K	115	HYP	CB-CG-CD	3.13	107.91	103.46
1	I	24	HYP	CD-N-CA	3.12	114.04	108.15
1	K	24	HYP	CB-CG-CD	3.11	107.89	103.46
1	G	3	HYP	C-CA-N	-3.11	105.11	110.86
1	G	24	HYP	CB-CG-CD	3.10	107.87	103.46
1	I	24	HYP	CB-CG-CD	3.10	107.87	103.46
1	C	124	HYP	CG-CB-CA	3.08	108.72	104.22
1	A	3	HYP	CB-CG-CD	3.08	107.85	103.46
1	C	209	HYP	CD-N-CA	3.05	113.91	108.15
1	C	3	HYP	CG-CB-CA	3.05	108.68	104.22
1	K	215	HYP	CG-CD-N	-3.04	100.57	105.47
1	G	127	HYP	CG-CB-CA	2.97	108.55	104.22
1	I	21	HYP	CD-N-CA	2.95	113.72	108.15
1	E	206	HYP	CG-CD-N	-2.95	100.72	105.47
1	G	221	HYP	CD-N-CA	2.93	113.69	108.15
1	G	9	HYP	CD-N-CA	2.93	113.68	108.15
1	A	103	HYP	CB-CG-CD	2.93	107.62	103.46
1	I	127	HYP	CB-CG-CD	2.93	107.62	103.46
1	C	21	HYP	C-CA-N	-2.92	105.46	110.86
1	K	24	HYP	CD-N-CA	2.92	113.66	108.15
1	G	227	HYP	CB-CG-CD	2.91	107.59	103.46
1	A	103	HYP	CG-CB-CA	2.91	108.46	104.22
1	I	209	HYP	CG-CB-CA	2.88	108.43	104.22
1	G	115	HYP	C-CA-N	2.88	116.19	110.86
1	I	224	HYP	CB-CG-CD	2.87	107.55	103.46
1	E	209	HYP	CG-CB-CA	2.87	108.41	104.22
1	I	21	HYP	CG-CD-N	-2.86	100.86	105.47
1	A	227	HYP	CB-CG-CD	2.86	107.53	103.46
1	E	127	HYP	CB-CG-CD	2.85	107.51	103.46
1	E	15	HYP	OD1-CG-CB	-2.84	102.47	110.16
1	K	209	HYP	CD-N-CA	2.84	113.50	108.15
1	E	21	HYP	C-CA-N	-2.81	105.66	110.86
1	K	127	HYP	CG-CB-CA	2.81	108.32	104.22
1	C	24	HYP	CD-N-CA	2.80	113.44	108.15
1	C	3	HYP	CB-CA-N	-2.80	100.30	104.59
1	C	209	HYP	CG-CB-CA	2.79	108.30	104.22
1	E	124	HYP	CD-N-CA	2.79	113.42	108.15
1	E	3	HYP	CB-CA-N	-2.78	100.32	104.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	K	9	HYP	CG-CB-CA	-2.77	100.18	104.22
1	G	15	HYP	CB-CG-CD	2.77	107.40	103.46
1	K	27	HYP	CG-CB-CA	2.77	108.26	104.22
1	A	106	HYP	C-CA-N	2.75	115.95	110.86
1	G	15	HYP	CG-CD-N	-2.75	101.05	105.47
1	C	127	HYP	CB-CG-CD	2.75	107.37	103.46
1	I	127	HYP	CD-N-CA	2.74	113.33	108.15
1	I	15	HYP	CG-CB-CA	2.73	108.21	104.22
1	K	9	HYP	CB-CG-CD	2.73	107.35	103.46
1	A	127	HYP	CG-CB-CA	2.73	108.21	104.22
1	G	221	HYP	CG-CB-CA	2.72	108.19	104.22
1	K	21	HYP	CG-CD-N	-2.70	101.12	105.47
1	K	27	HYP	CD-N-CA	2.69	113.23	108.15
1	I	127	HYP	C-CA-N	2.69	115.84	110.86
1	A	115	HYP	CD-N-CA	2.69	113.22	108.15
1	C	224	HYP	CD-N-CA	2.68	113.22	108.15
1	C	127	HYP	CG-CB-CA	2.68	108.13	104.22
1	E	115	HYP	CD-N-CA	2.68	113.20	108.15
1	C	3	HYP	CD-N-CA	2.65	113.15	108.15
1	C	9	HYP	CD-N-CA	2.64	113.14	108.15
1	K	124	HYP	CB-CG-CD	2.64	107.22	103.46
1	I	115	HYP	CB-CG-CD	2.64	107.22	103.46
1	G	124	HYP	CB-CG-CD	2.63	107.20	103.46
1	A	127	HYP	CB-CG-CD	2.62	107.19	103.46
1	E	15	HYP	CG-CB-CA	2.61	108.03	104.22
1	K	206	HYP	C-CA-N	2.61	115.69	110.86
1	K	103	HYP	CG-CB-CA	2.61	108.03	104.22
1	G	121	HYP	CB-CG-CD	2.58	107.13	103.46
1	G	115	HYP	CB-CG-CD	2.58	107.13	103.46
1	I	3	HYP	CB-CG-CD	2.57	107.12	103.46
1	G	21	HYP	CG-CB-CA	2.57	107.97	104.22
1	E	209	HYP	CD-N-CA	2.57	113.00	108.15
1	K	124	HYP	CG-CB-CA	2.56	107.96	104.22
1	A	209	HYP	CD-N-CA	2.56	112.99	108.15
1	A	124	HYP	CD-N-CA	2.55	112.96	108.15
1	A	15	HYP	CG-CB-CA	2.55	107.94	104.22
1	I	127	HYP	CG-CB-CA	2.54	107.93	104.22
1	A	115	HYP	CB-CG-CD	2.54	107.08	103.46
1	G	206	HYP	CD-N-CA	2.54	112.95	108.15
1	G	227	HYP	CG-CB-CA	2.54	107.92	104.22
1	I	106	HYP	CG-CD-N	-2.53	101.39	105.47
1	G	21	HYP	CD-N-CA	2.53	112.93	108.15

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	21	HYP	CD-N-CA	2.53	112.93	108.15
1	A	3	HYP	C-CA-N	-2.52	106.19	110.86
1	A	227	HYP	C-CA-N	2.52	115.52	110.86
1	I	103	HYP	CB-CG-CD	2.51	107.03	103.46
1	I	224	HYP	CD-N-CA	2.49	112.86	108.15
1	G	127	HYP	CB-CG-CD	2.48	107.00	103.46
1	C	115	HYP	CB-CG-CD	2.49	107.00	103.46
1	G	6	HYP	CD-N-CA	2.48	112.83	108.15
1	A	215	HYP	CB-CG-CD	2.46	106.97	103.46
1	C	124	HYP	CB-CG-CD	2.46	106.96	103.46
1	E	127	HYP	C-CA-N	2.44	115.38	110.86
1	K	121	HYP	OD1-CG-CB	-2.44	103.56	110.16
1	A	109	HYP	CB-CG-CD	2.42	106.91	103.46
1	A	127	HYP	CD-N-CA	2.41	112.70	108.15
1	K	6	HYP	C-CA-N	2.41	115.32	110.86
1	G	27	HYP	CG-CB-CA	2.41	107.73	104.22
1	I	206	HYP	OD1-CG-CB	-2.39	103.68	110.16
1	K	127	HYP	CD-N-CA	2.39	112.67	108.15
1	A	224	HYP	CD-N-CA	2.39	112.65	108.15
1	E	3	HYP	CG-CB-CA	2.39	107.70	104.22
1	A	221	HYP	CD-N-CA	2.38	112.65	108.15
1	G	6	HYP	CB-CG-CD	2.38	106.84	103.46
1	C	103	HYP	CD-N-CA	2.37	112.63	108.15
1	A	124	HYP	CB-CG-CD	2.36	106.82	103.46
1	C	209	HYP	C-CA-N	2.35	115.21	110.86
1	G	103	HYP	CG-CB-CA	2.35	107.65	104.22
1	E	124	HYP	CB-CG-CD	2.34	106.79	103.46
1	E	15	HYP	OD1-CG-CD	2.34	115.81	110.31
1	I	209	HYP	CB-CG-CD	2.34	106.78	103.46
1	E	6	HYP	CD-N-CA	2.33	112.56	108.15
1	E	206	HYP	CB-CG-CD	-2.33	100.15	103.46
1	C	115	HYP	C-CA-N	2.33	115.17	110.86
1	I	227	HYP	CG-CD-N	-2.33	101.72	105.47
1	K	209	HYP	CB-CG-CD	2.32	106.76	103.46
1	C	24	HYP	CB-CG-CD	2.31	106.75	103.46
1	K	9	HYP	CG-CD-N	-2.31	101.75	105.47
1	E	3	HYP	CD-N-CA	2.31	112.51	108.15
1	I	115	HYP	C-CA-N	2.31	115.13	110.86
1	K	15	HYP	CG-CD-N	-2.30	101.76	105.47
1	C	209	HYP	CB-CG-CD	2.30	106.73	103.46
1	A	124	HYP	CG-CB-CA	2.30	107.58	104.22
1	E	109	HYP	OD1-CG-CB	-2.30	103.94	110.16

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	127	HYP	CD-N-CA	2.30	112.48	108.15
1	A	221	HYP	CB-CG-CD	2.29	106.72	103.46
1	A	203	HYP	CB-CG-CD	2.30	106.73	103.46
1	C	21	HYP	CD-N-CA	2.29	112.48	108.15
1	G	227	HYP	CD-N-CA	2.29	112.47	108.15
1	E	115	HYP	CG-CB-CA	2.29	107.56	104.22
1	C	127	HYP	CD-N-CA	2.28	112.46	108.15
1	E	227	HYP	C-CA-N	-2.28	106.65	110.86
1	A	27	HYP	CG-CB-CA	2.28	107.54	104.22
1	C	3	HYP	CG-CD-N	-2.27	101.81	105.47
1	I	215	HYP	CG-CD-N	-2.27	101.81	105.47
1	A	24	HYP	CG-CB-CA	2.26	107.52	104.22
1	I	24	HYP	C-CA-N	2.25	115.02	110.86
1	I	103	HYP	CG-CB-CA	2.25	107.50	104.22
1	C	6	HYP	CB-CG-CD	2.24	106.65	103.46
1	E	127	HYP	CG-CB-CA	2.24	107.49	104.22
1	G	103	HYP	CB-CG-CD	2.24	106.64	103.46
1	G	224	HYP	CD-N-CA	2.23	112.37	108.15
1	A	24	HYP	CD-N-CA	2.22	112.34	108.15
1	I	124	HYP	CD-N-CA	2.21	112.32	108.15
1	K	27	HYP	CB-CG-CD	2.20	106.60	103.46
1	C	209	HYP	OD1-CG-CB	-2.19	104.24	110.16
1	G	27	HYP	CB-CG-CD	2.18	106.57	103.46
1	I	21	HYP	OD1-CG-CD	-2.18	105.19	110.31
1	G	209	HYP	CD-N-CA	2.18	112.26	108.15
1	I	109	HYP	CG-CD-N	-2.17	101.97	105.47
1	I	109	HYP	CD-N-CA	2.17	112.24	108.15
1	C	221	HYP	CG-CB-CA	2.16	107.38	104.22
1	G	3	HYP	CB-CG-CD	2.16	106.54	103.46
1	E	6	HYP	CB-CG-CD	2.15	106.52	103.46
1	G	209	HYP	CG-CD-N	-2.15	102.01	105.47
1	C	121	HYP	CG-CB-CA	2.13	107.33	104.22
1	I	103	HYP	CD-N-CA	2.13	112.17	108.15
1	K	3	HYP	CB-CG-CD	2.13	106.49	103.46
1	A	9	HYP	CD-N-CA	2.12	112.15	108.15
1	C	15	HYP	CD-N-CA	2.12	112.15	108.15
1	C	24	HYP	CG-CD-N	-2.12	102.06	105.47
1	I	203	HYP	CB-CG-CD	2.12	106.47	103.46
1	I	15	HYP	OD1-CG-CB	-2.11	104.45	110.16
1	C	109	HYP	OD1-CG-CD	-2.11	105.36	110.31
1	E	24	HYP	CD-N-CA	2.10	112.12	108.15
1	C	9	HYP	CB-CA-N	-2.10	101.36	104.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	9	HYP	C-CA-N	-2.10	106.98	110.86
1	E	106	HYP	C-CA-N	2.10	114.74	110.86
1	K	203	HYP	OD1-CG-CB	-2.09	104.51	110.16
1	A	15	HYP	C-CA-N	2.08	114.72	110.86
1	A	3	HYP	CD-N-CA	2.08	112.08	108.15
1	A	106	HYP	CB-CG-CD	2.08	106.42	103.46
1	E	3	HYP	OD1-CG-CB	2.08	115.80	110.16
1	E	124	HYP	CG-CB-CA	2.08	107.25	104.22
1	G	215	HYP	CG-CB-CA	2.07	107.25	104.22
1	A	24	HYP	CB-CG-CD	2.07	106.40	103.46
1	E	121	HYP	CB-CG-CD	2.07	106.40	103.46
1	A	6	HYP	CG-CB-CA	2.06	107.23	104.22
1	A	15	HYP	CD-N-CA	2.06	112.05	108.15
1	G	3	HYP	CB-CA-N	-2.06	101.43	104.59
1	K	106	HYP	CD-N-CA	2.06	112.03	108.15
1	E	203	HYP	CG-CD-N	-2.05	102.16	105.47
1	G	106	HYP	CD-N-CA	2.04	112.01	108.15
1	C	15	HYP	OD1-CG-CB	-2.04	104.65	110.16
1	I	209	HYP	C-CA-N	2.03	114.62	110.86
1	E	215	HYP	OD1-CG-CD	-2.02	105.55	110.31
1	K	206	HYP	OD1-CG-CB	-2.02	104.69	110.16
1	A	215	HYP	CB-CA-N	2.01	107.67	104.59
1	E	24	HYP	CB-CG-CD	2.01	106.31	103.46
1	A	227	HYP	CG-CB-CA	2.00	107.14	104.22
1	K	115	HYP	CB-CA-N	2.00	107.65	104.59

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

4 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	$\# Z  > 2$
3	SO4	C	301	-	4,4,4	0.34	0	6,6,6	0.65	0
3	SO4	F	1901	-	4,4,4	0.38	0	6,6,6	0.57	0
3	SO4	I	301	-	4,4,4	0.50	0	6,6,6	0.41	0
3	SO4	K	301	-	4,4,4	0.33	0	6,6,6	0.44	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
3	SO4	C	301	-	-	0/0/0/0	0/0/0/0
3	SO4	F	1901	-	-	0/0/0/0	0/0/0/0
3	SO4	I	301	-	-	0/0/0/0	0/0/0/0
3	SO4	K	301	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 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	79/87 (90%)	1.08	14 (17%) 2 2	42, 62, 137, 164	0
1	C	79/87 (90%)	0.16	4 (5%) 27 27	23, 45, 93, 114	0
1	E	79/87 (90%)	-0.03	2 (2%) 54 55	31, 41, 84, 95	0
1	G	79/87 (90%)	0.70	9 (11%) 6 4	53, 67, 125, 149	0
1	I	79/87 (90%)	0.04	2 (2%) 54 55	19, 39, 80, 101	0
1	K	80/87 (91%)	0.00	4 (5%) 28 28	25, 45, 93, 109	0
2	B	188/192 (97%)	-0.12	1 (0%) 88 90	36, 55, 84, 103	0
2	D	188/192 (97%)	-0.19	4 (2%) 60 61	27, 45, 64, 79	0
2	F	188/192 (97%)	-0.03	2 (1%) 77 78	35, 55, 78, 89	0
2	H	187/192 (97%)	0.28	7 (3%) 39 39	54, 67, 98, 116	0
2	J	191/192 (99%)	-0.18	3 (1%) 68 69	20, 34, 56, 73	0
2	L	188/192 (97%)	-0.16	2 (1%) 77 78	21, 42, 76, 99	0
All	All	1605/1674 (95%)	0.05	54 (3%) 43 44	19, 51, 89, 164	0

All (54) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	227	HYP	14.1
1	A	127	HYP	13.4
1	A	227	HYP	12.8
1	G	127	HYP	9.0
1	A	27	HYP	8.2
1	A	226	PRO	7.0
1	A	26	PRO	6.1
1	A	126	PRO	5.7
1	C	227	HYP	5.6
1	G	3	HYP	5.1
1	I	127	HYP	5.0

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Mol	Chain	Res	Type	RSRZ
1	A	24	HYP	4.7
2	L	1862	THR	4.0
1	G	224	HYP	3.9
1	A	221	HYP	3.9
1	A	123	PRO	3.8
1	E	227	HYP	3.6
2	H	1746	ASN	3.4
1	I	227	HYP	3.4
1	G	126	PRO	3.2
1	K	227	HYP	3.1
1	A	124	HYP	3.1
1	A	224	HYP	3.1
2	H	1823	PHE	3.1
1	C	224	HYP	3.0
2	F	1862	THR	2.9
2	H	1788	ALA	2.7
1	C	127	HYP	2.7
2	B	1862	THR	2.7
1	G	226	PRO	2.7
2	J	1686	CYS	2.7
2	J	1746	ASN	2.7
2	D	1746	ASN	2.6
1	C	3	HYP	2.6
1	G	27	HYP	2.5
2	H	1748	VAL	2.5
1	E	3	HYP	2.5
1	G	24	HYP	2.4
2	L	1746	ASN	2.3
2	F	1873	SER	2.3
1	A	115	HYP	2.2
2	D	1873	SER	2.2
1	A	217	VAL	2.2
1	K	3	HYP	2.1
1	K	124	HYP	2.1
2	H	1747	VAL	2.1
1	A	9	HYP	2.1
2	D	1869	HIS	2.1
2	J	1873	SER	2.1
1	G	215	HYP	2.0
2	D	1862	THR	2.0
2	H	1795	ALA	2.0
1	K	203	HYP	2.0

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Mol	Chain	Res	Type	RSRZ
2	H	1871	LEU	2.0

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

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
1	HYP	K	227	8/9	0.26	8.45	103,107,113,114	0
1	HYP	G	227	8/9	0.74	5.97	143,146,153,157	0
1	HYP	C	227	8/9	0.37	3.51	108,114,116,118	0
1	HYP	E	3	8/9	0.23	2.76	55,58,62,64	0
1	HYP	C	124	8/9	0.20	2.75	66,70,72,73	0
1	HYP	I	127	8/9	0.34	2.59	85,89,94,98	0
1	HYP	C	203	8/9	0.21	2.46	52,55,61,63	0
1	HYP	E	227	8/9	0.28	2.41	85,87,93,94	0
1	HYP	G	127	8/9	0.52	2.36	126,133,134,135	0
1	HYP	K	124	8/9	0.22	2.02	64,67,71,73	0
1	HYP	I	103	8/9	0.19	1.77	55,59,67,69	0
1	HYP	C	127	8/9	0.25	1.73	99,106,108,113	0
1	HYP	G	3	8/9	0.35	1.73	71,73,78,80	0
1	HYP	A	227	8/9	0.68	1.71	157,165,167,170	0
1	HYP	C	3	8/9	0.24	1.69	78,83,91,95	0
1	HYP	A	115	8/9	0.22	1.63	45,46,48,51	0
1	HYP	A	127	8/9	0.71	1.57	148,156,161,161	0
1	HYP	C	224	8/9	0.26	1.51	80,84,90,93	0
1	HYP	C	206	8/9	0.19	1.49	35,36,39,41	0
1	HYP	E	124	8/9	0.18	1.37	65,67,69,72	0
1	HYP	K	109	8/9	0.16	1.31	31,35,39,42	0
1	HYP	A	106	8/9	0.18	1.26	55,56,59,61	0
1	HYP	G	9	8/9	0.19	1.24	55,58,62,63	0
1	HYP	A	221	8/9	0.29	1.15	86,91,92,93	0
1	HYP	A	9	8/9	0.22	1.13	45,47,50,53	0
1	HYP	I	227	8/9	0.27	1.03	94,100,104,104	0
1	HYP	E	27	8/9	0.21	1.02	78,81,89,90	0
1	HYP	K	121	8/9	0.19	1.01	48,49,55,56	0
1	HYP	E	115	8/9	0.18	0.95	32,33,36,38	0
1	HYP	G	224	8/9	0.30	0.91	119,122,129,130	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
1	HYP	E	15	8/9	0.18	0.85	31,32,33,34	0
1	HYP	E	103	8/9	0.16	0.82	44,45,47,48	0
1	HYP	C	24	8/9	0.19	0.80	67,70,73,74	0
1	HYP	C	221	8/9	0.20	0.78	61,62,70,72	0
1	HYP	E	121	8/9	0.18	0.73	46,47,50,51	0
1	HYP	K	203	8/9	0.21	0.71	60,63,67,70	0
1	HYP	K	224	8/9	0.15	0.69	77,80,87,88	0
1	HYP	K	106	8/9	0.17	0.69	46,51,58,59	0
1	HYP	A	109	8/9	0.15	0.69	45,48,51,54	0
1	HYP	I	121	8/9	0.20	0.63	40,41,46,48	0
1	HYP	E	127	8/9	0.19	0.62	91,95,97,97	0
1	HYP	K	103	8/9	0.17	0.57	69,75,84,86	0
1	HYP	C	15	8/9	0.16	0.56	21,22,23,25	0
1	HYP	G	15	8/9	0.20	0.54	57,58,62,63	0
1	HYP	K	3	8/9	0.22	0.53	74,79,85,86	0
1	HYP	E	221	8/9	0.17	0.53	51,53,56,57	0
1	HYP	G	215	8/9	0.22	0.52	61,63,65,65	0
1	HYP	K	127	8/9	0.17	0.50	90,95,97,102	0
1	HYP	K	215	8/9	0.17	0.49	28,29,32,34	0
1	HYP	K	9	8/9	0.15	0.41	36,38,40,41	0
1	HYP	E	24	8/9	0.15	0.39	66,68,73,74	0
1	HYP	I	109	8/9	0.15	0.35	26,28,31,32	0
1	HYP	G	24	8/9	0.23	0.34	97,100,102,103	0
1	HYP	C	106	8/9	0.16	0.34	46,49,51,54	0
1	HYP	A	206	8/9	0.15	0.33	49,50,52,53	0
1	HYP	K	206	8/9	0.16	0.33	45,48,55,55	0
1	HYP	C	209	8/9	0.15	0.30	25,27,29,29	0
1	HYP	I	221	8/9	0.17	0.29	53,55,62,63	0
1	HYP	A	209	8/9	0.15	0.27	43,44,46,47	0
1	HYP	I	24	8/9	0.14	0.27	55,59,61,62	0
1	HYP	G	206	8/9	0.14	0.26	55,57,60,62	0
1	HYP	K	6	8/9	0.15	0.24	49,51,55,56	0
1	HYP	K	27	8/9	0.15	0.21	88,90,99,102	0
1	HYP	K	209	8/9	0.15	0.19	32,36,42,45	0
1	HYP	I	21	8/9	0.16	0.18	34,36,37,39	0
1	HYP	A	15	8/9	0.16	0.17	44,46,46,47	0
1	HYP	C	9	8/9	0.15	0.14	34,41,46,49	0
1	HYP	K	221	8/9	0.15	0.14	54,57,62,62	0
1	HYP	G	106	8/9	0.14	0.12	57,59,62,63	0
1	HYP	I	27	8/9	0.19	0.08	84,88,92,95	0
1	HYP	I	9	8/9	0.15	0.05	31,34,40,40	0
1	HYP	I	6	8/9	0.14	0.05	44,47,50,52	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
1	HYP	E	6	8/9	0.15	0.01	43,45,49,50	0
1	HYP	I	206	8/9	0.14	-0.01	33,37,42,47	0
1	HYP	A	21	8/9	0.21	-0.02	75,78,81,81	0
1	HYP	E	215	8/9	0.14	-0.02	37,38,43,43	0
1	HYP	K	24	8/9	0.12	-0.02	59,63,64,64	0
1	HYP	G	103	8/9	0.14	-0.02	65,66,71,72	0
1	HYP	G	209	8/9	0.16	-0.03	51,53,56,60	0
1	HYP	C	21	8/9	0.17	-0.04	42,44,46,49	0
1	HYP	G	124	8/9	0.20	-0.04	105,107,113,116	0
1	HYP	A	24	8/9	0.33	-0.06	101,105,110,110	0
1	HYP	E	9	8/9	0.15	-0.08	35,38,42,43	0
1	HYP	K	21	8/9	0.17	-0.09	41,44,45,45	0
1	HYP	E	203	8/9	0.15	-0.11	44,47,52,54	0
1	HYP	I	124	8/9	0.15	-0.11	56,61,62,62	0
1	HYP	E	21	8/9	0.16	-0.13	49,52,53,54	0
1	HYP	A	27	8/9	0.48	-0.13	132,137,143,145	0
1	HYP	I	209	8/9	0.14	-0.16	22,25,28,33	0
1	HYP	I	15	8/9	0.15	-0.19	19,21,25,29	0
1	HYP	A	215	8/9	0.17	-0.20	51,53,56,58	0
1	HYP	G	203	8/9	0.16	-0.21	62,65,70,71	0
1	HYP	C	215	8/9	0.15	-0.22	32,35,38,38	0
1	HYP	K	15	8/9	0.15	-0.22	28,30,35,36	0
1	HYP	E	209	8/9	0.14	-0.23	30,31,33,34	0
1	HYP	G	115	8/9	0.17	-0.23	62,64,68,69	0
1	HYP	I	3	8/9	0.17	-0.23	66,70,75,77	0
1	HYP	A	124	8/9	0.26	-0.27	111,117,119,119	0
1	HYP	G	121	8/9	0.19	-0.27	88,90,97,100	0
1	HYP	I	106	8/9	0.13	-0.33	36,38,43,44	0
1	HYP	C	109	8/9	0.13	-0.36	36,39,47,47	0
1	HYP	C	27	8/9	0.16	-0.36	98,102,107,109	0
1	HYP	G	221	8/9	0.16	-0.38	96,100,104,104	0
1	HYP	A	3	8/9	0.19	-0.41	81,87,93,97	0
1	HYP	I	203	8/9	0.13	-0.41	52,57,66,67	0
1	HYP	K	115	8/9	0.14	-0.41	27,29,31,33	0
1	HYP	A	6	8/9	0.13	-0.42	58,61,67,67	0
1	HYP	G	109	8/9	0.14	-0.43	54,55,58,59	0
1	HYP	C	121	8/9	0.15	-0.46	45,47,50,50	0
1	HYP	G	6	8/9	0.12	-0.46	60,63,67,68	0
1	HYP	I	224	8/9	0.14	-0.47	70,73,82,86	0
1	HYP	I	215	8/9	0.15	-0.49	22,25,27,28	0
1	HYP	A	224	8/9	0.26	-0.52	116,121,124,125	0
1	HYP	C	115	8/9	0.14	-0.57	27,30,37,39	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
1	HYP	E	224	8/9	0.13	-0.58	63,66,67,68	0
1	HYP	I	115	8/9	0.14	-0.63	22,24,31,34	0
1	HYP	E	109	8/9	0.13	-0.64	35,37,39,41	0
1	HYP	C	103	8/9	0.14	-0.65	59,63,66,66	0
1	HYP	G	27	8/9	0.24	-0.66	126,130,132,133	0
1	HYP	A	203	8/9	0.13	-0.66	61,63,67,67	0
1	HYP	E	206	8/9	0.12	-0.73	34,36,38,39	0
1	HYP	A	121	8/9	0.17	-0.76	80,82,86,86	0
1	HYP	G	21	8/9	0.16	-0.81	78,81,82,83	0
1	HYP	A	103	8/9	0.12	-0.90	70,72,75,76	0
1	HYP	E	106	8/9	0.12	-0.95	36,37,38,39	0
1	HYP	C	6	8/9	0.12	-1.15	53,58,66,69	0

### 6.3 Carbohydrates ⓘ

There are no carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
3	SO4	F	1901	5/5	0.18	-0.17	70,73,74,74	0
3	SO4	I	301	5/5	0.13	-0.94	44,50,50,52	0
3	SO4	C	301	5/5	0.12	-1.05	43,49,49,49	0
3	SO4	K	301	5/5	0.12	-1.53	53,56,59,59	0

### 6.5 Other polymers ⓘ

There are no such residues in this entry.