



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2014 – 11:26 AM GMT

PDB ID : 3B2C
Title : Crystal structure of the collagen triple helix model [{PRO-HYP(R)-GLY}4-{HYP(S)-Pro-GLY}2-{PRO-HYP(R)-GLY}4]3
Authors : Motooka, D.; Kawahara, K.; Nakamura, S.; Doi, M.; Nishi, Y.; Nishiuchi, Y.; Nakazawa, T.; Yoshida, T.; Ohkubo, T.; Kobayashi, Y.; Kang, Y.K.; Uchiyama, S.
Deposited on : 2011-07-26
Resolution : 1.36 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

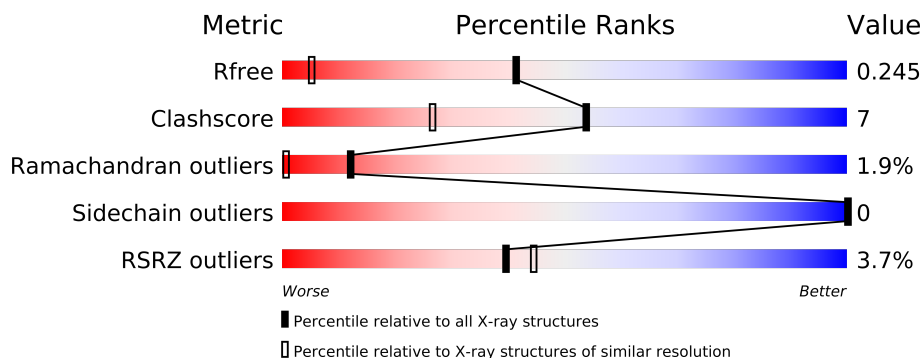
The following versions of software and data (see [references](#)) were used in the production of this report:

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 1.36 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	1519 (1.40-1.32)
Clashscore	79885	1707 (1.40-1.32)
Ramachandran outliers	78287	1662 (1.40-1.32)
Sidechain outliers	78261	1661 (1.40-1.32)
RSRZ outliers	66119	1519 (1.40-1.32)

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 ≥ 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	30	
1	B	30	
1	C	30	
1	D	30	
1	E	30	
1	F	30	
1	G	30	
1	H	30	
1	I	30	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 12591 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-like peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	A	30	Total	C	N	O	0	30	0
			1330	840	210	280			
1	B	30	Total	C	N	O	0	30	0
			1330	840	210	280			
1	C	30	Total	C	N	O	0	30	0
			1330	840	210	280			
1	D	30	Total	C	N	O	0	30	0
			1330	840	210	280			
1	E	30	Total	C	N	O	0	30	0
			1330	840	210	280			
1	F	30	Total	C	N	O	0	30	0
			1330	840	210	280			
1	G	30	Total	C	N	O	0	30	0
			1330	840	210	280			
1	H	30	Total	C	N	O	0	30	0
			1330	840	210	280			
1	I	30	Total	C	N	O	0	30	0
			1330	840	210	280			

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	56	Total	O	0	0
			56	56		
2	B	79	Total	O	0	0
			79	79		
2	C	63	Total	O	0	0
			63	63		
2	D	79	Total	O	0	0
			79	79		
2	E	69	Total	O	0	0
			69	69		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	F	81	Total 81	O 81	0	0
2	G	53	Total 53	O 53	0	0
2	H	76	Total 76	O 76	0	0
2	I	65	Total 65	O 65	0	0

3 Residue-property plots

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 ($RSRZ > 2$). 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-like peptide

Chain A: 



- Molecule 1: Collagen-like peptide

Chain B: 



- Molecule 1: Collagen-like peptide

Chain C: 



- Molecule 1: Collagen-like peptide

Chain D: 



- Molecule 1: Collagen-like peptide

Chain E: 



- Molecule 1: Collagen-like peptide

Chain F: 



- Molecule 1: Collagen-like peptide

Chain G: 



- Molecule 1: Collagen-like peptide

Chain H: 



- Molecule 1: Collagen-like peptide

Chain I: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	40.80Å 26.15Å 84.74Å 90.00° 92.48° 90.00°	Depositor
Resolution (Å)	10.00 – 1.36 9.97 – 1.36	Depositor EDS
% Data completeness (in resolution range)	98.8 (10.00-1.36) 98.8 (9.97-1.36)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.94 (at 1.36Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.203 , 0.248 0.206 , 0.245	Depositor DCC
R_{free} test set	1942 reflections (5.32%)	DCC
Wilson B-factor (Å ²)	6.9	Xtriage
Anisotropy	0.283	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 56.4	EDS
Estimated twinning fraction	0.000 for h,-k,-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.37$	Xtriage
Outliers	41 of 38474 reflections (0.107%)	Xtriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	12591	wwPDB-VP
Average B, all atoms (Å ²)	8.0	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 74.64 % 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 1.4721e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities 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, HZP

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	0.53	0/763	0.66	0/938
1	B	0.48	0/763	0.67	0/938
1	C	0.62	0/763	0.66	0/938
1	D	0.59	0/763	0.62	0/938
1	E	0.51	0/763	0.72	0/938
1	F	0.47	0/763	0.63	0/938
1	G	0.61	0/763	0.60	0/938
1	H	0.49	0/763	0.63	0/938
1	I	0.63	0/763	0.68	0/938
All	All	0.55	0/6867	0.65	0/8442

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3
1	B	0	1
1	C	0	4
1	D	0	5
1	E	0	4
1	F	0	4
1	G	0	7
1	H	0	6
1	I	0	4
All	All	0	38

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (38) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	27[E]	GLY	Mainchain
1	A	27[F]	GLY	Mainchain
1	A	27[G]	GLY	Mainchain
1	B	4[D]	PRO	Mainchain
1	C	1[B]	PRO	Mainchain
1	C	1[C]	PRO	Mainchain
1	C	1[D]	PRO	Mainchain
1	C	27[G]	GLY	Mainchain
1	D	27[E]	GLY	Mainchain
1	D	27[F]	GLY	Mainchain
1	D	29[E]	HYP	Mainchain
1	D	29[F]	HYP	Mainchain
1	D	29[G]	HYP	Mainchain
1	E	2[C]	HYP	Peptide
1	E	2[D]	HYP	Peptide
1	E	25[F]	PRO	Mainchain
1	E	25[G]	PRO	Mainchain
1	F	27[E]	GLY	Mainchain
1	F	27[F]	GLY	Mainchain
1	F	28[E]	PRO	Mainchain
1	F	28[F]	PRO	Mainchain
1	G	26[E]	HYP	Mainchain
1	G	26[G]	HYP	Mainchain
1	G	27[F]	GLY	Mainchain
1	G	28[E]	PRO	Mainchain
1	G	28[F]	PRO	Mainchain
1	G	28[G]	PRO	Mainchain
1	G	3[D]	GLY	Mainchain
1	H	29[F]	HYP	Mainchain
1	H	29[G]	HYP	Mainchain
1	H	3[C]	GLY	Mainchain
1	H	3[D]	GLY	Mainchain
1	H	4[C]	PRO	Mainchain
1	H	4[D]	PRO	Mainchain
1	I	28[G]	PRO	Mainchain
1	I	3[B]	GLY	Mainchain
1	I	3[C]	GLY	Mainchain
1	I	3[D]	GLY	Mainchain

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	1330	0	0	20	1
1	B	1330	0	0	10	0
1	C	1330	0	0	18	5
1	D	1330	0	0	13	7
1	E	1330	0	0	20	2
1	F	1330	0	0	18	4
1	G	1330	0	0	18	7
1	H	1330	0	0	20	2
1	I	1330	0	0	17	0
2	A	56	0	0	3	2
2	B	79	0	0	0	3
2	C	63	0	0	0	2
2	D	79	0	0	3	1
2	E	69	0	0	2	1
2	F	81	0	0	0	1
2	G	53	0	0	1	4
2	H	76	0	0	0	1
2	I	65	0	0	0	3
All	All	12591	0	0	83	23

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 7.

All (83) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:29[A]:HYP:CA	1:F:28[A]:PRO:CG	1.86	1.52
1:G:20[D]:HYP:OD1	2:G:578:HOH:O	1.66	1.12
1:E:29[A]:HYP:CB	1:F:28[A]:PRO:CG	2.47	0.93
1:E:29[A]:HYP:OD1	1:F:28[A]:PRO:CG	2.23	0.86
1:G:2[D]:HYP:OD1	1:H:1[D]:PRO:CG	2.25	0.85
1:D:27[F]:GLY:CA	1:F:27[F]:GLY:O	2.25	0.85
1:D:27[E]:GLY:CA	1:F:27[E]:GLY:O	2.25	0.85
1:H:23[G]:HYP:OD1	1:I:22[G]:PRO:CG	2.24	0.85
1:A:7[D]:PRO:CG	1:C:5[D]:HYP:OD1	2.28	0.82

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:4[D]:PRO:CA	2:D:424:HOH:O	2.26	0.82
1:E:29[A]:HYP:N	1:F:28[A]:PRO:CG	2.44	0.79
1:E:26[G]:HYP:OD1	1:F:25[G]:PRO:CG	2.31	0.78
1:D:4[D]:PRO:CB	2:D:424:HOH:O	2.33	0.75
1:G:2[B]:HYP:OD1	1:H:1[B]:PRO:CD	2.35	0.75
1:G:2[A]:HYP:OD1	1:H:1[A]:PRO:CD	2.35	0.75
1:D:4[D]:PRO:CB	1:D:5[D]:HYP:CD	2.66	0.73
1:G:2[B]:HYP:OD1	1:H:1[B]:PRO:CG	2.36	0.73
1:G:2[A]:HYP:OD1	1:H:1[A]:PRO:CG	2.36	0.72
1:A:11[G]:HYP:OD1	2:A:414:HOH:O	2.07	0.71
1:A:11[E]:HYP:OD1	2:A:414:HOH:O	2.07	0.71
1:A:11[F]:HYP:OD1	2:A:414:HOH:O	2.07	0.71
1:H:4[C]:PRO:CB	1:H:5[C]:HYP:CD	2.72	0.67
1:H:4[D]:PRO:CB	1:H:5[D]:HYP:CD	2.72	0.67
1:E:2[C]:HYP:OD1	2:E:279:HOH:O	2.13	0.67
1:E:2[D]:HYP:OD1	2:E:279:HOH:O	2.13	0.67
1:D:28[F]:PRO:O	1:F:30[F]:GLY:N	2.33	0.61
1:H:30[E]:GLY:C	1:I:30[E]:GLY:CA	2.70	0.60
1:H:30[A]:GLY:C	1:I:30[A]:GLY:CA	2.70	0.59
1:A:28[B]:PRO:CG	1:C:29[B]:HYP:OD1	2.51	0.58
1:A:28[C]:PRO:CG	1:C:29[C]:HYP:OD1	2.51	0.58
1:E:29[A]:HYP:CG	1:F:28[A]:PRO:CG	2.81	0.58
1:B:30[G]:GLY:N	1:C:28[G]:PRO:O	2.38	0.57
1:G:2[B]:HYP:CA	1:H:1[B]:PRO:CD	2.83	0.57
1:G:2[A]:HYP:CA	1:H:1[A]:PRO:CD	2.83	0.57
1:A:10[D]:PRO:CG	1:C:8[D]:HYP:OD1	2.53	0.56
1:D:27[E]:GLY:CA	1:F:27[E]:GLY:C	2.74	0.55
1:D:27[F]:GLY:CA	1:F:27[F]:GLY:C	2.74	0.55
1:A:1[B]:PRO:CG	1:C:2[B]:HYP:CB	2.85	0.55
1:A:1[C]:PRO:CG	1:C:2[C]:HYP:CB	2.85	0.55
1:G:28[A]:PRO:CG	1:I:26[A]:HYP:OD1	2.55	0.54
1:H:30[A]:GLY:N	1:I:28[A]:PRO:O	2.41	0.53
1:H:30[E]:GLY:N	1:I:28[E]:PRO:O	2.41	0.53
1:G:30[F]:GLY:CA	1:I:30[F]:GLY:O	2.56	0.53
1:G:30[E]:GLY:CA	1:I:30[E]:GLY:O	2.56	0.53
1:B:29[G]:HYP:OD1	1:C:28[G]:PRO:CG	2.57	0.53
1:E:29[A]:HYP:CA	1:F:28[A]:PRO:CD	2.81	0.52
1:B:29[G]:HYP:CA	1:C:28[G]:PRO:CD	2.88	0.51
1:E:28[A]:PRO:C	1:F:28[A]:PRO:CD	2.79	0.51
1:B:26[F]:HYP:OD1	1:E:26[F]:HYP:OD1	2.28	0.51
1:B:26[G]:HYP:OD1	1:E:26[G]:HYP:OD1	2.28	0.51
1:H:5[D]:HYP:OD1	1:I:4[D]:PRO:CG	2.59	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:28[E]:PRO:CG	1:C:29[E]:HYP:OD1	2.60	0.49
1:A:28[F]:PRO:CG	1:C:29[F]:HYP:OD1	2.60	0.49
1:G:28[E]:PRO:CG	1:I:29[E]:HYP:OD1	2.61	0.49
1:G:28[F]:PRO:O	1:I:30[F]:GLY:N	2.46	0.48
1:C:10[G]:PRO:CB	2:D:371:HOH:O	2.62	0.48
1:D:13[A]:HYP:OD1	1:F:11[A]:HYP:OD1	2.32	0.47
1:A:27[E]:GLY:CA	1:C:27[E]:GLY:O	2.64	0.46
1:B:11[F]:HYP:OD1	1:C:13[F]:HYP:OD1	2.34	0.46
1:A:27[F]:GLY:CA	1:C:27[F]:GLY:O	2.64	0.45
1:E:28[A]:PRO:O	1:F:28[A]:PRO:CD	2.64	0.45
1:E:30[A]:GLY:N	1:F:28[A]:PRO:O	2.50	0.45
1:A:13[A]:HYP:OD1	1:C:11[A]:HYP:OD1	2.36	0.44
1:H:29[A]:HYP:OD1	1:I:28[A]:PRO:CG	2.66	0.44
1:H:29[E]:HYP:OD1	1:I:28[E]:PRO:CG	2.66	0.44
1:A:13[D]:HYP:OD1	1:C:11[D]:HYP:OD1	2.36	0.43
1:G:3[B]:GLY:N	1:H:1[B]:PRO:O	2.50	0.43
1:G:28[E]:PRO:O	1:I:30[E]:GLY:N	2.51	0.43
1:G:3[A]:GLY:N	1:H:1[A]:PRO:O	2.50	0.43
1:A:28[G]:PRO:O	1:C:27[G]:GLY:N	2.52	0.43
1:D:2[D]:HYP:OD1	1:E:1[D]:PRO:CG	2.67	0.42
1:A:29[F]:HYP:OD1	1:B:28[F]:PRO:CG	2.67	0.42
1:A:29[G]:HYP:OD1	1:B:28[G]:PRO:CG	2.67	0.42
1:E:29[A]:HYP:N	1:F:28[A]:PRO:CD	2.83	0.41
1:A:30[B]:GLY:C	1:B:30[B]:GLY:CA	2.88	0.41
1:A:30[A]:GLY:C	1:B:30[A]:GLY:CA	2.88	0.41
1:G:3[A]:GLY:CA	1:I:1[A]:PRO:O	2.69	0.41
1:H:6[D]:GLY:N	1:I:4[D]:PRO:O	2.53	0.41
1:D:2[E]:HYP:OD1	1:E:4[E]:PRO:CG	2.68	0.41
1:D:5[B]:HYP:OD1	1:E:4[B]:PRO:CG	2.68	0.41
1:G:28[F]:PRO:CG	1:I:29[F]:HYP:OD1	2.69	0.41
1:D:5[A]:HYP:OD1	1:E:4[A]:PRO:CG	2.69	0.41
1:E:8[G]:HYP:OD1	1:F:7[G]:PRO:CG	2.70	0.40

All (23) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:G:502:HOH:O	2:I:319:HOH:O[2_655]	1.26	0.94
2:A:274:HOH:O	2:B:511:HOH:O[2_555]	1.34	0.86
2:G:523:HOH:O	2:I:561:HOH:O[2_645]	1.87	0.33
1:E:30[A]:GLY:C	1:H:1[A]:PRO:N[1_554]	1.92	0.28
1:E:30[B]:GLY:C	1:H:1[B]:PRO:N[1_554]	1.93	0.27

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:F:30[A]:GLY:N	1:G:1[A]:PRO:O[1_554]	1.94	0.26
2:B:585:HOH:O	2:D:613:HOH:O[2_555]	1.94	0.26
2:F:212:HOH:O	2:H:611:HOH:O[2_645]	2.02	0.18
2:A:476:HOH:O	2:B:618:HOH:O[2_545]	2.05	0.15
1:D:29[G]:HYP:O	2:G:612:HOH:O[1_554]	2.08	0.12
1:D:30[D]:GLY:O	1:G:1[D]:PRO:N[1_554]	2.08	0.12
1:C:30[A]:GLY:O	1:F:1[A]:PRO:N[1_554]	2.11	0.09
1:C:30[E]:GLY:O	1:F:1[E]:PRO:N[1_554]	2.11	0.09
1:C:30[F]:GLY:O	1:F:1[F]:PRO:N[1_554]	2.12	0.08
2:C:443:HOH:O	2:I:269:HOH:O[1_455]	2.12	0.08
2:C:443:HOH:O	2:E:584:HOH:O[2_545]	2.14	0.06
1:A:30[D]:GLY:O	1:D:1[D]:PRO:N[1_554]	2.17	0.03
1:C:1[B]:PRO:N	1:G:30[B]:GLY:O[1_456]	2.17	0.03
1:C:1[C]:PRO:N	1:G:30[C]:GLY:O[1_456]	2.17	0.03
1:D:29[F]:HYP:O	2:G:612:HOH:O[1_554]	2.18	0.02
1:D:30[A]:GLY:CA	1:G:1[A]:PRO:N[1_554]	2.18	0.02
1:D:30[B]:GLY:CA	1:G:1[B]:PRO:N[1_554]	2.18	0.02
1:D:30[C]:GLY:CA	1:G:1[C]:PRO:N[1_554]	2.18	0.02

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	126/30 (420%)	112 (89%)	0	14 (11%)	1	0
1	B	126/30 (420%)	126 (100%)	0	0	100	100
1	C	126/30 (420%)	119 (94%)	0	7 (6%)	3	0
1	D	126/30 (420%)	126 (100%)	0	0	100	100
1	E	126/30 (420%)	119 (94%)	7 (6%)	0	100	100
1	F	126/30 (420%)	119 (94%)	7 (6%)	0	100	100
1	G	126/30 (420%)	126 (100%)	0	0	100	100
1	H	126/30 (420%)	126 (100%)	0	0	100	100
1	I	126/30 (420%)	126 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	1134/270 (420%)	1099 (97%)	14 (1%)	21 (2%)	12 1

All (21) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	28[A]	PRO
1	A	28[B]	PRO
1	A	28[C]	PRO
1	A	28[D]	PRO
1	A	28[E]	PRO
1	A	28[F]	PRO
1	A	28[G]	PRO
1	C	27[A]	GLY
1	C	27[B]	GLY
1	C	27[C]	GLY
1	C	27[D]	GLY
1	C	27[E]	GLY
1	C	27[F]	GLY
1	C	27[G]	GLY
1	A	27[A]	GLY
1	A	27[B]	GLY
1	A	27[C]	GLY
1	A	27[D]	GLY
1	A	27[E]	GLY
1	A	27[F]	GLY
1	A	27[G]	GLY

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	70/10 (700%)	70 (100%)	0	100 100
1	B	70/10 (700%)	70 (100%)	0	100 100
1	C	70/10 (700%)	70 (100%)	0	100 100
1	D	70/10 (700%)	70 (100%)	0	100 100
1	E	70/10 (700%)	70 (100%)	0	100 100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	F	70/10 (700%)	70 (100%)	0	100	100
1	G	70/10 (700%)	70 (100%)	0	100	100
1	H	70/10 (700%)	70 (100%)	0	100	100
1	I	70/10 (700%)	70 (100%)	0	100	100
All	All	630/90 (700%)	630 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

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 ⓘ

630 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	11[A]	1	8,8,9	6.55	1 (12%)	8,10,12	1.45	2 (25%)
1	HYP	A	11[B]	1	8,8,9	6.51	1 (12%)	8,10,12	1.46	2 (25%)
1	HYP	A	11[C]	1	8,8,9	6.55	1 (12%)	8,10,12	1.46	2 (25%)
1	HYP	A	11[D]	1	8,8,9	6.81	1 (12%)	8,10,12	1.92	3 (37%)
1	HYP	A	11[E]	1	8,8,9	6.32	1 (12%)	8,10,12	2.39	4 (50%)
1	HYP	A	11[F]	1	8,8,9	6.36	1 (12%)	8,10,12	2.38	4 (50%)
1	HYP	A	11[G]	1	8,8,9	6.39	1 (12%)	8,10,12	2.40	4 (50%)
1	HZP	A	13[A]	1	8,8,9	6.37	1 (12%)	8,10,12	4.30	5 (62%)
1	HZP	A	13[B]	1	8,8,9	6.51	1 (12%)	8,10,12	4.33	5 (62%)
1	HZP	A	13[C]	1	8,8,9	6.43	1 (12%)	8,10,12	4.32	5 (62%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	HZP	A	13[D]	1	8,8,9	5.97	1 (12%)	8,10,12	3.74	5 (62%)
1	HZP	A	13[E]	1	8,8,9	6.52	2 (25%)	8,10,12	3.56	5 (62%)
1	HZP	A	13[F]	1	8,8,9	6.51	2 (25%)	8,10,12	3.57	5 (62%)
1	HZP	A	13[G]	1	8,8,9	6.49	2 (25%)	8,10,12	3.55	4 (50%)
1	HZP	A	16[A]	1	8,8,9	6.62	2 (25%)	8,10,12	3.45	4 (50%)
1	HZP	A	16[B]	1	8,8,9	6.57	2 (25%)	8,10,12	3.41	4 (50%)
1	HZP	A	16[C]	1	8,8,9	6.59	2 (25%)	8,10,12	3.44	4 (50%)
1	HZP	A	16[D]	1	8,8,9	6.33	1 (12%)	8,10,12	3.97	5 (62%)
1	HZP	A	16[E]	1	8,8,9	7.21	2 (25%)	8,10,12	5.05	5 (62%)
1	HZP	A	16[F]	1	8,8,9	7.29	2 (25%)	8,10,12	5.04	5 (62%)
1	HZP	A	16[G]	1	8,8,9	7.35	2 (25%)	8,10,12	5.05	5 (62%)
1	HYP	A	2[A]	1	8,8,9	6.45	1 (12%)	8,10,12	2.27	3 (37%)
1	HYP	A	2[B]	1	8,8,9	6.48	1 (12%)	8,10,12	2.27	3 (37%)
1	HYP	A	2[C]	1	8,8,9	6.39	1 (12%)	8,10,12	2.24	3 (37%)
1	HYP	A	2[D]	1	8,8,9	2.99	1 (12%)	8,10,12	0.86	0
1	HYP	A	2[E]	1	8,8,9	6.99	1 (12%)	8,10,12	2.17	3 (37%)
1	HYP	A	2[F]	1	8,8,9	7.01	1 (12%)	8,10,12	2.16	3 (37%)
1	HYP	A	2[G]	1	8,8,9	6.93	1 (12%)	8,10,12	2.16	3 (37%)
1	HYP	A	20[A]	1	8,8,9	6.79	1 (12%)	8,10,12	2.14	3 (37%)
1	HYP	A	20[B]	1	8,8,9	6.80	1 (12%)	8,10,12	2.15	3 (37%)
1	HYP	A	20[C]	1	8,8,9	6.86	1 (12%)	8,10,12	2.14	3 (37%)
1	HYP	A	20[D]	1	8,8,9	7.79	2 (25%)	8,10,12	2.62	4 (50%)
1	HYP	A	20[E]	1	8,8,9	6.96	1 (12%)	8,10,12	1.61	2 (25%)
1	HYP	A	20[F]	1	8,8,9	6.88	1 (12%)	8,10,12	1.60	2 (25%)
1	HYP	A	20[G]	1	8,8,9	6.86	1 (12%)	8,10,12	1.59	2 (25%)
1	HYP	A	23[A]	1	8,8,9	7.00	1 (12%)	8,10,12	1.56	2 (25%)
1	HYP	A	23[B]	1	8,8,9	7.02	1 (12%)	8,10,12	1.55	2 (25%)
1	HYP	A	23[C]	1	8,8,9	7.00	1 (12%)	8,10,12	1.55	2 (25%)
1	HYP	A	23[D]	1	8,8,9	6.82	1 (12%)	8,10,12	2.15	3 (37%)
1	HYP	A	23[E]	1	8,8,9	7.02	1 (12%)	8,10,12	1.46	2 (25%)
1	HYP	A	23[F]	1	8,8,9	6.95	1 (12%)	8,10,12	1.48	2 (25%)
1	HYP	A	23[G]	1	8,8,9	7.11	1 (12%)	8,10,12	1.46	2 (25%)
1	HYP	A	26[A]	1	8,8,9	7.04	1 (12%)	8,10,12	1.47	2 (25%)
1	HYP	A	26[B]	1	8,8,9	7.04	1 (12%)	8,10,12	1.48	2 (25%)
1	HYP	A	26[C]	1	8,8,9	7.07	1 (12%)	8,10,12	1.48	2 (25%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	HYP	A	26[D]	1	8,8,9	6.95	1 (12%)	8,10,12	1.54	2 (25%)
1	HYP	A	26[E]	1	8,8,9	7.02	1 (12%)	8,10,12	1.99	3 (37%)
1	HYP	A	26[F]	1	8,8,9	7.02	1 (12%)	8,10,12	1.95	3 (37%)
1	HYP	A	26[G]	1	8,8,9	7.09	1 (12%)	8,10,12	1.99	3 (37%)
1	HYP	A	29[A]	1	8,8,9	7.38	1 (12%)	8,10,12	2.15	2 (25%)
1	HYP	A	29[B]	1	8,8,9	7.38	1 (12%)	8,10,12	2.10	2 (25%)
1	HYP	A	29[C]	1	8,8,9	7.38	1 (12%)	8,10,12	2.12	2 (25%)
1	HYP	A	29[D]	1	8,8,9	7.08	1 (12%)	8,10,12	1.49	2 (25%)
1	HYP	A	29[E]	1	8,8,9	6.69	1 (12%)	8,10,12	2.50	2 (25%)
1	HYP	A	29[F]	1	8,8,9	6.67	1 (12%)	8,10,12	2.46	2 (25%)
1	HYP	A	29[G]	1	8,8,9	6.69	1 (12%)	8,10,12	2.47	2 (25%)
1	HYP	A	5[A]	1	8,8,9	7.01	1 (12%)	8,10,12	2.14	3 (37%)
1	HYP	A	5[B]	1	8,8,9	6.91	1 (12%)	8,10,12	2.18	3 (37%)
1	HYP	A	5[C]	1	8,8,9	7.11	1 (12%)	8,10,12	2.17	3 (37%)
1	HYP	A	5[D]	1	8,8,9	6.38	1 (12%)	8,10,12	2.26	3 (37%)
1	HYP	A	5[E]	1	8,8,9	6.79	1 (12%)	8,10,12	1.90	2 (25%)
1	HYP	A	5[F]	1	8,8,9	6.73	1 (12%)	8,10,12	1.89	2 (25%)
1	HYP	A	5[G]	1	8,8,9	6.77	1 (12%)	8,10,12	1.91	3 (37%)
1	HYP	A	8[A]	1	8,8,9	6.86	1 (12%)	8,10,12	1.92	2 (25%)
1	HYP	A	8[B]	1	8,8,9	6.77	1 (12%)	8,10,12	1.93	3 (37%)
1	HYP	A	8[C]	1	8,8,9	6.86	1 (12%)	8,10,12	1.91	2 (25%)
1	HYP	A	8[D]	1	8,8,9	6.81	1 (12%)	8,10,12	2.20	3 (37%)
1	HYP	A	8[E]	1	8,8,9	6.52	1 (12%)	8,10,12	1.47	2 (25%)
1	HYP	A	8[F]	1	8,8,9	6.51	1 (12%)	8,10,12	1.48	2 (25%)
1	HYP	A	8[G]	1	8,8,9	6.56	1 (12%)	8,10,12	1.47	2 (25%)
1	HYP	B	11[A]	1	8,8,9	7.01	1 (12%)	8,10,12	1.87	3 (37%)
1	HYP	B	11[B]	1	8,8,9	7.02	1 (12%)	8,10,12	1.85	3 (37%)
1	HYP	B	11[C]	1	8,8,9	7.28	1 (12%)	8,10,12	1.80	2 (25%)
1	HYP	B	11[D]	1	8,8,9	7.25	1 (12%)	8,10,12	1.78	2 (25%)
1	HYP	B	11[E]	1	8,8,9	7.01	1 (12%)	8,10,12	1.87	3 (37%)
1	HYP	B	11[F]	1	8,8,9	6.57	1 (12%)	8,10,12	2.14	4 (50%)
1	HYP	B	11[G]	1	8,8,9	6.62	1 (12%)	8,10,12	2.15	4 (50%)
1	HZP	B	13[A]	1	8,8,9	6.80	1 (12%)	8,10,12	5.82	5 (62%)
1	HZP	B	13[B]	1	8,8,9	6.73	1 (12%)	8,10,12	5.81	5 (62%)
1	HZP	B	13[C]	1	8,8,9	7.35	1 (12%)	8,10,12	3.78	5 (62%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	HZP	B	13[D]	1	8,8,9	7.25	1 (12%)	8,10,12	3.77	5 (62%)
1	HZP	B	13[E]	1	8,8,9	6.80	1 (12%)	8,10,12	5.81	5 (62%)
1	HZP	B	13[F]	1	8,8,9	6.96	2 (25%)	8,10,12	4.01	5 (62%)
1	HZP	B	13[G]	1	8,8,9	6.96	2 (25%)	8,10,12	4.01	5 (62%)
1	HZP	B	16[A]	1	8,8,9	6.86	2 (25%)	8,10,12	3.69	4 (50%)
1	HZP	B	16[B]	1	8,8,9	6.91	2 (25%)	8,10,12	3.67	4 (50%)
1	HZP	B	16[C]	1	8,8,9	5.05	2 (25%)	8,10,12	3.36	7 (87%)
1	HZP	B	16[D]	1	8,8,9	5.05	2 (25%)	8,10,12	3.37	7 (87%)
1	HZP	B	16[E]	1	8,8,9	6.84	2 (25%)	8,10,12	3.70	4 (50%)
1	HZP	B	16[F]	1	8,8,9	6.95	1 (12%)	8,10,12	5.86	5 (62%)
1	HZP	B	16[G]	1	8,8,9	6.89	1 (12%)	8,10,12	5.85	5 (62%)
1	HYP	B	2[A]	1	8,8,9	7.10	1 (12%)	8,10,12	1.87	3 (37%)
1	HYP	B	2[B]	1	8,8,9	7.05	1 (12%)	8,10,12	1.88	3 (37%)
1	HYP	B	2[C]	1	8,8,9	6.42	2 (25%)	8,10,12	2.14	3 (37%)
1	HYP	B	2[D]	1	8,8,9	6.47	2 (25%)	8,10,12	2.14	3 (37%)
1	HYP	B	2[E]	1	8,8,9	7.20	1 (12%)	8,10,12	1.86	3 (37%)
1	HYP	B	2[F]	1	8,8,9	6.75	1 (12%)	8,10,12	1.91	3 (37%)
1	HYP	B	2[G]	1	8,8,9	6.66	1 (12%)	8,10,12	1.89	2 (25%)
1	HYP	B	20[A]	1	8,8,9	6.42	1 (12%)	8,10,12	2.14	3 (37%)
1	HYP	B	20[B]	1	8,8,9	6.41	1 (12%)	8,10,12	2.14	3 (37%)
1	HYP	B	20[C]	1	8,8,9	6.18	1 (12%)	8,10,12	2.51	5 (62%)
1	HYP	B	20[D]	1	8,8,9	6.18	1 (12%)	8,10,12	2.52	5 (62%)
1	HYP	B	20[E]	1	8,8,9	6.41	1 (12%)	8,10,12	2.15	3 (37%)
1	HYP	B	20[F]	1	8,8,9	7.23	1 (12%)	8,10,12	1.37	1 (12%)
1	HYP	B	20[G]	1	8,8,9	7.22	1 (12%)	8,10,12	1.37	1 (12%)
1	HYP	B	23[A]	1	8,8,9	7.19	1 (12%)	8,10,12	1.36	1 (12%)
1	HYP	B	23[B]	1	8,8,9	7.10	1 (12%)	8,10,12	1.35	1 (12%)
1	HYP	B	23[C]	1	8,8,9	6.49	1 (12%)	8,10,12	2.13	3 (37%)
1	HYP	B	23[D]	1	8,8,9	6.42	1 (12%)	8,10,12	2.14	4 (50%)
1	HYP	B	23[E]	1	8,8,9	7.29	1 (12%)	8,10,12	1.36	1 (12%)
1	HYP	B	23[F]	1	8,8,9	7.03	1 (12%)	8,10,12	1.77	2 (25%)
1	HYP	B	23[G]	1	8,8,9	7.08	1 (12%)	8,10,12	1.78	2 (25%)
1	HYP	B	26[A]	1	8,8,9	7.10	1 (12%)	8,10,12	1.67	2 (25%)
1	HYP	B	26[B]	1	8,8,9	7.03	1 (12%)	8,10,12	1.68	2 (25%)
1	HYP	B	26[C]	1	8,8,9	7.15	1 (12%)	8,10,12	1.36	1 (12%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	HYP	B	26[D]	1	8,8,9	7.05	1 (12%)	8,10,12	1.35	1 (12%)
1	HYP	B	26[E]	1	8,8,9	7.12	1 (12%)	8,10,12	1.66	2 (25%)
1	HYP	B	26[F]	1	8,8,9	6.47	1 (12%)	8,10,12	2.33	5 (62%)
1	HYP	B	26[G]	1	8,8,9	6.41	1 (12%)	8,10,12	2.32	5 (62%)
1	HYP	B	29[A]	1	8,8,9	6.16	1 (12%)	8,10,12	2.08	3 (37%)
1	HYP	B	29[B]	1	8,8,9	6.16	1 (12%)	8,10,12	2.09	3 (37%)
1	HYP	B	29[C]	1	8,8,9	7.14	1 (12%)	8,10,12	1.63	2 (25%)
1	HYP	B	29[D]	1	8,8,9	7.14	1 (12%)	8,10,12	1.63	2 (25%)
1	HYP	B	29[E]	1	8,8,9	6.16	1 (12%)	8,10,12	2.08	3 (37%)
1	HYP	B	29[F]	1	8,8,9	5.67	1 (12%)	8,10,12	1.91	2 (25%)
1	HYP	B	29[G]	1	8,8,9	5.67	1 (12%)	8,10,12	1.91	2 (25%)
1	HYP	B	5[A]	1	8,8,9	6.72	1 (12%)	8,10,12	1.80	2 (25%)
1	HYP	B	5[B]	1	8,8,9	6.78	1 (12%)	8,10,12	1.83	2 (25%)
1	HYP	B	5[C]	1	8,8,9	7.02	1 (12%)	8,10,12	1.79	3 (37%)
1	HYP	B	5[D]	1	8,8,9	5.25	2 (25%)	8,10,12	3.02	3 (37%)
1	HYP	B	5[E]	1	8,8,9	6.72	1 (12%)	8,10,12	1.80	2 (25%)
1	HYP	B	5[F]	1	8,8,9	7.08	1 (12%)	8,10,12	1.90	2 (25%)
1	HYP	B	5[G]	1	8,8,9	7.11	1 (12%)	8,10,12	1.91	2 (25%)
1	HYP	B	8[A]	1	8,8,9	7.12	1 (12%)	8,10,12	1.88	2 (25%)
1	HYP	B	8[B]	1	8,8,9	7.15	1 (12%)	8,10,12	1.90	2 (25%)
1	HYP	B	8[C]	1	8,8,9	6.74	1 (12%)	8,10,12	1.98	3 (37%)
1	HYP	B	8[D]	1	8,8,9	6.81	1 (12%)	8,10,12	1.99	3 (37%)
1	HYP	B	8[E]	1	8,8,9	7.09	1 (12%)	8,10,12	1.86	2 (25%)
1	HYP	B	8[F]	1	8,8,9	6.92	1 (12%)	8,10,12	1.84	3 (37%)
1	HYP	B	8[G]	1	8,8,9	6.92	1 (12%)	8,10,12	1.85	3 (37%)
1	HYP	C	11[A]	1	8,8,9	6.80	1 (12%)	8,10,12	2.08	3 (37%)
1	HYP	C	11[B]	1	8,8,9	7.23	1 (12%)	8,10,12	1.63	2 (25%)
1	HYP	C	11[C]	1	8,8,9	7.23	1 (12%)	8,10,12	1.64	2 (25%)
1	HYP	C	11[D]	1	8,8,9	7.23	1 (12%)	8,10,12	1.64	2 (25%)
1	HYP	C	11[E]	1	8,8,9	6.87	1 (12%)	8,10,12	2.09	3 (37%)
1	HYP	C	11[F]	1	8,8,9	6.80	1 (12%)	8,10,12	2.08	3 (37%)
1	HYP	C	11[G]	1	8,8,9	6.66	1 (12%)	8,10,12	2.96	5 (62%)
1	HZP	C	13[A]	1	8,8,9	7.10	1 (12%)	8,10,12	3.28	5 (62%)
1	HZP	C	13[B]	1	8,8,9	6.74	1 (12%)	8,10,12	4.21	5 (62%)
1	HZP	C	13[C]	1	8,8,9	6.83	1 (12%)	8,10,12	4.20	5 (62%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	HZP	C	13[D]	1	8,8,9	6.65	1 (12%)	8,10,12	4.20	5 (62%)
1	HZP	C	13[E]	1	8,8,9	7.15	1 (12%)	8,10,12	3.26	5 (62%)
1	HZP	C	13[F]	1	8,8,9	7.11	1 (12%)	8,10,12	3.30	5 (62%)
1	HZP	C	13[G]	1	8,8,9	6.78	1 (12%)	8,10,12	3.18	5 (62%)
1	HZP	C	16[A]	1	8,8,9	6.92	1 (12%)	8,10,12	4.50	5 (62%)
1	HZP	C	16[B]	1	8,8,9	7.15	1 (12%)	8,10,12	3.14	5 (62%)
1	HZP	C	16[C]	1	8,8,9	7.23	1 (12%)	8,10,12	3.15	5 (62%)
1	HZP	C	16[D]	1	8,8,9	7.07	1 (12%)	8,10,12	3.14	5 (62%)
1	HZP	C	16[E]	1	8,8,9	6.92	1 (12%)	8,10,12	4.49	5 (62%)
1	HZP	C	16[F]	1	8,8,9	6.92	1 (12%)	8,10,12	4.50	5 (62%)
1	HZP	C	16[G]	1	8,8,9	7.52	2 (25%)	8,10,12	5.60	5 (62%)
1	HYP	C	2[A]	1	8,8,9	6.60	1 (12%)	8,10,12	1.88	2 (25%)
1	HYP	C	2[B]	1	8,8,9	6.59	1 (12%)	8,10,12	2.21	4 (50%)
1	HYP	C	2[C]	1	8,8,9	6.54	1 (12%)	8,10,12	2.21	4 (50%)
1	HYP	C	2[D]	1	8,8,9	6.52	1 (12%)	8,10,12	2.19	4 (50%)
1	HYP	C	2[E]	1	8,8,9	6.60	1 (12%)	8,10,12	1.88	2 (25%)
1	HYP	C	2[F]	1	8,8,9	6.61	1 (12%)	8,10,12	1.88	2 (25%)
1	HYP	C	2[G]	1	8,8,9	6.72	1 (12%)	8,10,12	1.76	3 (37%)
1	HYP	C	20[A]	1	8,8,9	7.22	1 (12%)	8,10,12	1.61	2 (25%)
1	HYP	C	20[B]	1	8,8,9	6.41	1 (12%)	8,10,12	2.45	4 (50%)
1	HYP	C	20[C]	1	8,8,9	6.47	1 (12%)	8,10,12	2.45	4 (50%)
1	HYP	C	20[D]	1	8,8,9	6.36	1 (12%)	8,10,12	2.45	4 (50%)
1	HYP	C	20[E]	1	8,8,9	7.29	1 (12%)	8,10,12	1.62	2 (25%)
1	HYP	C	20[F]	1	8,8,9	7.15	1 (12%)	8,10,12	1.60	2 (25%)
1	HYP	C	20[G]	1	8,8,9	6.16	1 (12%)	8,10,12	1.76	2 (25%)
1	HYP	C	23[A]	1	8,8,9	6.20	1 (12%)	8,10,12	1.73	2 (25%)
1	HYP	C	23[B]	1	8,8,9	7.24	1 (12%)	8,10,12	1.56	2 (25%)
1	HYP	C	23[C]	1	8,8,9	7.15	1 (12%)	8,10,12	1.57	2 (25%)
1	HYP	C	23[D]	1	8,8,9	7.26	1 (12%)	8,10,12	1.55	2 (25%)
1	HYP	C	23[E]	1	8,8,9	6.19	1 (12%)	8,10,12	1.71	2 (25%)
1	HYP	C	23[F]	1	8,8,9	6.19	1 (12%)	8,10,12	1.74	2 (25%)
1	HYP	C	23[G]	1	8,8,9	6.96	1 (12%)	8,10,12	2.50	4 (50%)
1	HYP	C	26[A]	1	8,8,9	6.99	1 (12%)	8,10,12	2.42	4 (50%)
1	HYP	C	26[B]	1	8,8,9	6.22	1 (12%)	8,10,12	1.73	2 (25%)
1	HYP	C	26[C]	1	8,8,9	6.23	1 (12%)	8,10,12	1.72	2 (25%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	HYP	C	26[D]	1	8,8,9	6.22	1 (12%)	8,10,12	1.72	2 (25%)
1	HYP	C	26[E]	1	8,8,9	6.99	1 (12%)	8,10,12	2.39	4 (50%)
1	HYP	C	26[F]	1	8,8,9	6.99	1 (12%)	8,10,12	2.43	4 (50%)
1	HYP	C	26[G]	1	8,8,9	6.93	1 (12%)	8,10,12	1.84	3 (37%)
1	HYP	C	29[A]	1	8,8,9	6.91	1 (12%)	8,10,12	2.05	4 (50%)
1	HYP	C	29[B]	1	8,8,9	6.94	1 (12%)	8,10,12	2.37	4 (50%)
1	HYP	C	29[C]	1	8,8,9	7.00	1 (12%)	8,10,12	2.37	4 (50%)
1	HYP	C	29[D]	1	8,8,9	6.96	1 (12%)	8,10,12	2.36	4 (50%)
1	HYP	C	29[E]	1	8,8,9	6.95	1 (12%)	8,10,12	2.04	4 (50%)
1	HYP	C	29[F]	1	8,8,9	6.91	1 (12%)	8,10,12	2.03	4 (50%)
1	HYP	C	29[G]	1	8,8,9	6.54	1 (12%)	8,10,12	1.82	2 (25%)
1	HYP	C	5[A]	1	8,8,9	6.72	1 (12%)	8,10,12	1.75	3 (37%)
1	HYP	C	5[B]	1	8,8,9	6.59	1 (12%)	8,10,12	1.84	2 (25%)
1	HYP	C	5[C]	1	8,8,9	6.49	1 (12%)	8,10,12	1.84	2 (25%)
1	HYP	C	5[D]	1	8,8,9	6.68	1 (12%)	8,10,12	1.83	2 (25%)
1	HYP	C	5[E]	1	8,8,9	6.72	1 (12%)	8,10,12	1.75	3 (37%)
1	HYP	C	5[F]	1	8,8,9	6.72	1 (12%)	8,10,12	1.75	3 (37%)
1	HYP	C	5[G]	1	8,8,9	7.15	1 (12%)	8,10,12	1.51	1 (12%)
1	HYP	C	8[A]	1	8,8,9	7.10	1 (12%)	8,10,12	1.53	1 (12%)
1	HYP	C	8[B]	1	8,8,9	6.72	1 (12%)	8,10,12	1.76	3 (37%)
1	HYP	C	8[C]	1	8,8,9	6.72	1 (12%)	8,10,12	1.75	3 (37%)
1	HYP	C	8[D]	1	8,8,9	6.72	1 (12%)	8,10,12	1.77	3 (37%)
1	HYP	C	8[E]	1	8,8,9	7.06	1 (12%)	8,10,12	1.53	1 (12%)
1	HYP	C	8[F]	1	8,8,9	7.15	1 (12%)	8,10,12	1.53	1 (12%)
1	HYP	C	8[G]	1	8,8,9	6.74	1 (12%)	8,10,12	2.12	3 (37%)
1	HYP	D	11[A]	1	8,8,9	7.18	1 (12%)	8,10,12	1.76	2 (25%)
1	HYP	D	11[B]	1	8,8,9	7.22	1 (12%)	8,10,12	1.75	2 (25%)
1	HYP	D	11[C]	1	8,8,9	7.11	1 (12%)	8,10,12	1.76	2 (25%)
1	HYP	D	11[D]	1	8,8,9	7.27	1 (12%)	8,10,12	1.55	2 (25%)
1	HYP	D	11[E]	1	8,8,9	7.20	1 (12%)	8,10,12	1.85	2 (25%)
1	HYP	D	11[F]	1	8,8,9	7.20	1 (12%)	8,10,12	1.84	2 (25%)
1	HYP	D	11[G]	1	8,8,9	7.20	1 (12%)	8,10,12	1.83	2 (25%)
1	HZP	D	13[A]	1	8,8,9	7.41	1 (12%)	8,10,12	3.35	5 (62%)
1	HZP	D	13[B]	1	8,8,9	7.41	1 (12%)	8,10,12	3.35	5 (62%)
1	HZP	D	13[C]	1	8,8,9	7.42	1 (12%)	8,10,12	3.36	5 (62%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	HZP	D	13[D]	1	8,8,9	7.09	2 (25%)	8,10,12	4.24	5 (62%)
1	HZP	D	13[E]	1	8,8,9	7.30	1 (12%)	8,10,12	3.62	5 (62%)
1	HZP	D	13[F]	1	8,8,9	7.20	1 (12%)	8,10,12	3.63	5 (62%)
1	HZP	D	13[G]	1	8,8,9	7.10	1 (12%)	8,10,12	3.63	5 (62%)
1	HZP	D	16[A]	1	8,8,9	7.06	1 (12%)	8,10,12	3.40	5 (62%)
1	HZP	D	16[B]	1	8,8,9	7.06	1 (12%)	8,10,12	3.42	5 (62%)
1	HZP	D	16[C]	1	8,8,9	7.06	1 (12%)	8,10,12	3.38	5 (62%)
1	HZP	D	16[D]	1	8,8,9	7.46	1 (12%)	8,10,12	3.35	5 (62%)
1	HZP	D	16[E]	1	8,8,9	7.35	2 (25%)	8,10,12	3.36	4 (50%)
1	HZP	D	16[F]	1	8,8,9	7.35	2 (25%)	8,10,12	3.38	4 (50%)
1	HZP	D	16[G]	1	8,8,9	7.35	2 (25%)	8,10,12	3.39	4 (50%)
1	HYP	D	2[A]	1	8,8,9	6.58	2 (25%)	8,10,12	2.21	3 (37%)
1	HYP	D	2[B]	1	8,8,9	6.58	2 (25%)	8,10,12	2.21	3 (37%)
1	HYP	D	2[C]	1	8,8,9	6.58	2 (25%)	8,10,12	2.23	3 (37%)
1	HYP	D	2[D]	1	8,8,9	6.98	1 (12%)	8,10,12	1.98	3 (37%)
1	HYP	D	2[E]	1	8,8,9	6.87	1 (12%)	8,10,12	1.96	3 (37%)
1	HYP	D	2[F]	1	8,8,9	6.87	1 (12%)	8,10,12	1.96	3 (37%)
1	HYP	D	2[G]	1	8,8,9	6.87	1 (12%)	8,10,12	1.96	3 (37%)
1	HYP	D	20[A]	1	8,8,9	7.40	1 (12%)	8,10,12	1.94	3 (37%)
1	HYP	D	20[B]	1	8,8,9	7.31	1 (12%)	8,10,12	1.95	3 (37%)
1	HYP	D	20[C]	1	8,8,9	7.42	1 (12%)	8,10,12	1.95	3 (37%)
1	HYP	D	20[D]	1	8,8,9	6.36	1 (12%)	8,10,12	2.45	5 (62%)
1	HYP	D	20[E]	1	8,8,9	6.18	1 (12%)	8,10,12	1.73	2 (25%)
1	HYP	D	20[F]	1	8,8,9	6.23	1 (12%)	8,10,12	1.76	2 (25%)
1	HYP	D	20[G]	1	8,8,9	6.24	1 (12%)	8,10,12	1.78	2 (25%)
1	HYP	D	23[A]	1	8,8,9	6.19	1 (12%)	8,10,12	1.80	2 (25%)
1	HYP	D	23[B]	1	8,8,9	6.19	1 (12%)	8,10,12	1.77	2 (25%)
1	HYP	D	23[C]	1	8,8,9	6.19	1 (12%)	8,10,12	1.76	2 (25%)
1	HYP	D	23[D]	1	8,8,9	7.34	1 (12%)	8,10,12	1.95	3 (37%)
1	HYP	D	23[E]	1	8,8,9	6.74	1 (12%)	8,10,12	1.89	4 (50%)
1	HYP	D	23[F]	1	8,8,9	6.80	1 (12%)	8,10,12	1.88	4 (50%)
1	HYP	D	23[G]	1	8,8,9	6.87	1 (12%)	8,10,12	1.88	4 (50%)
1	HYP	D	26[A]	1	8,8,9	6.80	1 (12%)	8,10,12	1.91	4 (50%)
1	HYP	D	26[B]	1	8,8,9	6.80	1 (12%)	8,10,12	1.91	4 (50%)
1	HYP	D	26[C]	1	8,8,9	6.73	1 (12%)	8,10,12	1.91	4 (50%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	HYP	D	26[D]	1	8,8,9	6.18	1 (12%)	8,10,12	1.78	2 (25%)
1	HYP	D	26[E]	1	8,8,9	6.63	1 (12%)	8,10,12	2.14	3 (37%)
1	HYP	D	26[F]	1	8,8,9	6.62	1 (12%)	8,10,12	2.15	4 (50%)
1	HYP	D	26[G]	1	8,8,9	6.63	1 (12%)	8,10,12	2.13	3 (37%)
1	HYP	D	29[A]	1	8,8,9	6.64	1 (12%)	8,10,12	1.78	3 (37%)
1	HYP	D	29[B]	1	8,8,9	6.68	1 (12%)	8,10,12	1.78	3 (37%)
1	HYP	D	29[C]	1	8,8,9	6.64	1 (12%)	8,10,12	1.78	3 (37%)
1	HYP	D	29[D]	1	8,8,9	6.78	1 (12%)	8,10,12	1.89	4 (50%)
1	HYP	D	29[E]	1	8,8,9	7.23	1 (12%)	8,10,12	1.70	3 (37%)
1	HYP	D	29[F]	1	8,8,9	7.20	1 (12%)	8,10,12	1.70	3 (37%)
1	HYP	D	29[G]	1	8,8,9	7.19	1 (12%)	8,10,12	1.71	3 (37%)
1	HYP	D	5[A]	1	8,8,9	6.86	1 (12%)	8,10,12	1.88	2 (25%)
1	HYP	D	5[B]	1	8,8,9	6.86	1 (12%)	8,10,12	1.86	2 (25%)
1	HYP	D	5[C]	1	8,8,9	6.79	1 (12%)	8,10,12	1.89	2 (25%)
1	HYP	D	5[D]	1	8,8,9	6.67	2 (25%)	8,10,12	2.67	4 (50%)
1	HYP	D	5[E]	1	8,8,9	7.27	1 (12%)	8,10,12	1.48	1 (12%)
1	HYP	D	5[F]	1	8,8,9	7.27	1 (12%)	8,10,12	1.49	1 (12%)
1	HYP	D	5[G]	1	8,8,9	7.23	1 (12%)	8,10,12	1.49	1 (12%)
1	HYP	D	8[A]	1	8,8,9	7.23	1 (12%)	8,10,12	1.53	1 (12%)
1	HYP	D	8[B]	1	8,8,9	7.23	1 (12%)	8,10,12	1.54	1 (12%)
1	HYP	D	8[C]	1	8,8,9	7.23	1 (12%)	8,10,12	1.52	1 (12%)
1	HYP	D	8[D]	1	8,8,9	6.82	1 (12%)	8,10,12	1.88	3 (37%)
1	HYP	D	8[E]	1	8,8,9	7.18	1 (12%)	8,10,12	1.67	2 (25%)
1	HYP	D	8[F]	1	8,8,9	7.18	1 (12%)	8,10,12	1.65	2 (25%)
1	HYP	D	8[G]	1	8,8,9	7.18	1 (12%)	8,10,12	1.63	2 (25%)
1	HYP	E	11[A]	1	8,8,9	6.07	1 (12%)	8,10,12	2.09	3 (37%)
1	HYP	E	11[B]	1	8,8,9	6.07	1 (12%)	8,10,12	2.09	3 (37%)
1	HYP	E	11[C]	1	8,8,9	6.35	1 (12%)	8,10,12	1.73	2 (25%)
1	HYP	E	11[D]	1	8,8,9	6.35	1 (12%)	8,10,12	1.72	2 (25%)
1	HYP	E	11[E]	1	8,8,9	6.07	1 (12%)	8,10,12	2.08	3 (37%)
1	HYP	E	11[F]	1	8,8,9	7.29	1 (12%)	8,10,12	2.12	3 (37%)
1	HYP	E	11[G]	1	8,8,9	7.29	1 (12%)	8,10,12	2.11	3 (37%)
1	HZP	E	13[A]	1	8,8,9	6.37	2 (25%)	8,10,12	4.09	5 (62%)
1	HZP	E	13[B]	1	8,8,9	6.37	2 (25%)	8,10,12	4.08	5 (62%)
1	HZP	E	13[C]	1	8,8,9	6.31	1 (12%)	8,10,12	5.15	5 (62%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	HZP	E	13[D]	1	8,8,9	6.31	1 (12%)	8,10,12	5.16	5 (62%)
1	HZP	E	13[E]	1	8,8,9	6.36	2 (25%)	8,10,12	4.09	5 (62%)
1	HZP	E	13[F]	1	8,8,9	6.28	1 (12%)	8,10,12	3.50	4 (50%)
1	HZP	E	13[G]	1	8,8,9	6.25	1 (12%)	8,10,12	3.48	4 (50%)
1	HZP	E	16[A]	1	8,8,9	6.38	1 (12%)	8,10,12	4.26	5 (62%)
1	HZP	E	16[B]	1	8,8,9	6.48	1 (12%)	8,10,12	4.28	5 (62%)
1	HZP	E	16[C]	1	8,8,9	6.37	2 (25%)	8,10,12	4.09	5 (62%)
1	HZP	E	16[D]	1	8,8,9	6.37	2 (25%)	8,10,12	4.09	5 (62%)
1	HZP	E	16[E]	1	8,8,9	6.43	1 (12%)	8,10,12	4.24	5 (62%)
1	HZP	E	16[F]	1	8,8,9	6.32	1 (12%)	8,10,12	5.14	7 (87%)
1	HZP	E	16[G]	1	8,8,9	6.40	1 (12%)	8,10,12	5.15	7 (87%)
1	HYP	E	2[A]	1	8,8,9	6.87	1 (12%)	8,10,12	2.08	3 (37%)
1	HYP	E	2[B]	1	8,8,9	7.04	1 (12%)	8,10,12	2.07	3 (37%)
1	HYP	E	2[C]	1	8,8,9	7.03	1 (12%)	8,10,12	2.33	4 (50%)
1	HYP	E	2[D]	1	8,8,9	7.03	1 (12%)	8,10,12	2.33	4 (50%)
1	HYP	E	2[E]	1	8,8,9	6.89	1 (12%)	8,10,12	2.08	3 (37%)
1	HYP	E	2[F]	1	8,8,9	6.69	1 (12%)	8,10,12	1.82	2 (25%)
1	HYP	E	2[G]	1	8,8,9	6.60	1 (12%)	8,10,12	1.83	2 (25%)
1	HYP	E	20[A]	1	8,8,9	6.82	1 (12%)	8,10,12	2.14	2 (25%)
1	HYP	E	20[B]	1	8,8,9	6.79	1 (12%)	8,10,12	2.12	2 (25%)
1	HYP	E	20[C]	1	8,8,9	6.80	1 (12%)	8,10,12	2.13	4 (50%)
1	HYP	E	20[D]	1	8,8,9	6.72	1 (12%)	8,10,12	2.11	4 (50%)
1	HYP	E	20[E]	1	8,8,9	6.89	1 (12%)	8,10,12	2.15	2 (25%)
1	HYP	E	20[F]	1	8,8,9	6.86	1 (12%)	8,10,12	1.57	2 (25%)
1	HYP	E	20[G]	1	8,8,9	6.86	1 (12%)	8,10,12	1.59	2 (25%)
1	HYP	E	23[A]	1	8,8,9	6.82	1 (12%)	8,10,12	1.56	2 (25%)
1	HYP	E	23[B]	1	8,8,9	6.82	1 (12%)	8,10,12	1.55	2 (25%)
1	HYP	E	23[C]	1	8,8,9	6.85	1 (12%)	8,10,12	2.08	2 (25%)
1	HYP	E	23[D]	1	8,8,9	6.85	1 (12%)	8,10,12	2.08	2 (25%)
1	HYP	E	23[E]	1	8,8,9	6.82	1 (12%)	8,10,12	1.59	2 (25%)
1	HYP	E	23[F]	1	8,8,9	7.19	1 (12%)	8,10,12	1.64	2 (25%)
1	HYP	E	23[G]	1	8,8,9	7.19	1 (12%)	8,10,12	1.66	2 (25%)
1	HYP	E	26[A]	1	8,8,9	7.24	1 (12%)	8,10,12	1.65	2 (25%)
1	HYP	E	26[B]	1	8,8,9	7.30	1 (12%)	8,10,12	1.66	2 (25%)
1	HYP	E	26[C]	1	8,8,9	6.82	1 (12%)	8,10,12	1.56	2 (25%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	HYP	E	26[D]	1	8,8,9	6.88	1 (12%)	8,10,12	1.55	2 (25%)
1	HYP	E	26[E]	1	8,8,9	7.17	1 (12%)	8,10,12	1.65	2 (25%)
1	HYP	E	26[F]	1	8,8,9	5.84	1 (12%)	8,10,12	2.08	3 (37%)
1	HYP	E	26[G]	1	8,8,9	5.84	1 (12%)	8,10,12	2.08	3 (37%)
1	HYP	E	29[A]	1	8,8,9	6.93	2 (25%)	8,10,12	2.61	5 (62%)
1	HYP	E	29[B]	1	8,8,9	6.93	2 (25%)	8,10,12	2.60	5 (62%)
1	HYP	E	29[C]	1	8,8,9	7.21	1 (12%)	8,10,12	1.62	2 (25%)
1	HYP	E	29[D]	1	8,8,9	7.21	1 (12%)	8,10,12	1.65	2 (25%)
1	HYP	E	29[E]	1	8,8,9	6.89	2 (25%)	8,10,12	2.62	5 (62%)
1	HYP	E	29[F]	1	8,8,9	6.99	1 (12%)	8,10,12	2.00	4 (50%)
1	HYP	E	29[G]	1	8,8,9	6.93	1 (12%)	8,10,12	1.99	4 (50%)
1	HYP	E	5[A]	1	8,8,9	6.82	1 (12%)	8,10,12	1.82	3 (37%)
1	HYP	E	5[B]	1	8,8,9	6.73	1 (12%)	8,10,12	1.86	3 (37%)
1	HYP	E	5[C]	1	8,8,9	5.67	1 (12%)	8,10,12	1.91	2 (25%)
1	HYP	E	5[D]	1	8,8,9	5.66	1 (12%)	8,10,12	1.92	2 (25%)
1	HYP	E	5[E]	1	8,8,9	6.91	1 (12%)	8,10,12	1.86	3 (37%)
1	HYP	E	5[F]	1	8,8,9	6.38	1 (12%)	8,10,12	1.71	2 (25%)
1	HYP	E	5[G]	1	8,8,9	6.37	1 (12%)	8,10,12	1.71	2 (25%)
1	HYP	E	8[A]	1	8,8,9	6.40	1 (12%)	8,10,12	1.72	2 (25%)
1	HYP	E	8[B]	1	8,8,9	6.40	1 (12%)	8,10,12	1.75	2 (25%)
1	HYP	E	8[C]	1	8,8,9	6.79	1 (12%)	8,10,12	1.77	2 (25%)
1	HYP	E	8[D]	1	8,8,9	6.74	1 (12%)	8,10,12	1.80	2 (25%)
1	HYP	E	8[E]	1	8,8,9	6.54	1 (12%)	8,10,12	1.73	2 (25%)
1	HYP	E	8[F]	1	8,8,9	6.28	1 (12%)	8,10,12	2.18	3 (37%)
1	HYP	E	8[G]	1	8,8,9	6.28	1 (12%)	8,10,12	2.18	3 (37%)
1	HYP	F	11[A]	1	8,8,9	6.80	1 (12%)	8,10,12	2.07	4 (50%)
1	HYP	F	11[B]	1	8,8,9	7.26	1 (12%)	8,10,12	1.81	2 (25%)
1	HYP	F	11[C]	1	8,8,9	7.26	1 (12%)	8,10,12	1.78	2 (25%)
1	HYP	F	11[D]	1	8,8,9	7.26	1 (12%)	8,10,12	1.83	2 (25%)
1	HYP	F	11[E]	1	8,8,9	6.80	1 (12%)	8,10,12	2.06	4 (50%)
1	HYP	F	11[F]	1	8,8,9	6.80	1 (12%)	8,10,12	2.09	4 (50%)
1	HYP	F	11[G]	1	8,8,9	6.39	2 (25%)	8,10,12	2.80	6 (75%)
1	HZP	F	13[A]	1	8,8,9	6.62	1 (12%)	8,10,12	3.94	4 (50%)
1	HZP	F	13[B]	1	8,8,9	7.16	1 (12%)	8,10,12	5.20	5 (62%)
1	HZP	F	13[C]	1	8,8,9	7.17	1 (12%)	8,10,12	5.20	5 (62%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	HZP	F	13[D]	1	8,8,9	7.17	1 (12%)	8,10,12	5.21	5 (62%)
1	HZP	F	13[E]	1	8,8,9	6.53	1 (12%)	8,10,12	3.95	4 (50%)
1	HZP	F	13[F]	1	8,8,9	6.72	1 (12%)	8,10,12	3.94	4 (50%)
1	HZP	F	13[G]	1	8,8,9	6.23	1 (12%)	8,10,12	4.52	5 (62%)
1	HZP	F	16[A]	1	8,8,9	6.17	1 (12%)	8,10,12	4.47	6 (75%)
1	HZP	F	16[B]	1	8,8,9	6.73	2 (25%)	8,10,12	3.98	5 (62%)
1	HZP	F	16[C]	1	8,8,9	6.67	2 (25%)	8,10,12	3.98	5 (62%)
1	HZP	F	16[D]	1	8,8,9	6.82	1 (12%)	8,10,12	3.98	5 (62%)
1	HZP	F	16[E]	1	8,8,9	6.17	1 (12%)	8,10,12	4.48	6 (75%)
1	HZP	F	16[F]	1	8,8,9	6.17	1 (12%)	8,10,12	4.46	6 (75%)
1	HZP	F	16[G]	1	8,8,9	6.78	2 (25%)	8,10,12	5.47	6 (75%)
1	HYP	F	2[A]	1	8,8,9	6.59	1 (12%)	8,10,12	1.80	3 (37%)
1	HYP	F	2[B]	1	8,8,9	6.76	1 (12%)	8,10,12	1.97	4 (50%)
1	HYP	F	2[C]	1	8,8,9	6.76	1 (12%)	8,10,12	1.96	4 (50%)
1	HYP	F	2[D]	1	8,8,9	6.76	1 (12%)	8,10,12	1.98	4 (50%)
1	HYP	F	2[E]	1	8,8,9	6.68	1 (12%)	8,10,12	1.77	2 (25%)
1	HYP	F	2[F]	1	8,8,9	6.53	1 (12%)	8,10,12	1.82	3 (37%)
1	HYP	F	2[G]	1	8,8,9	6.62	1 (12%)	8,10,12	1.96	2 (25%)
1	HYP	F	20[A]	1	8,8,9	7.14	1 (12%)	8,10,12	1.73	2 (25%)
1	HYP	F	20[B]	1	8,8,9	6.35	1 (12%)	8,10,12	1.99	3 (37%)
1	HYP	F	20[C]	1	8,8,9	6.36	1 (12%)	8,10,12	1.98	3 (37%)
1	HYP	F	20[D]	1	8,8,9	6.35	1 (12%)	8,10,12	1.98	3 (37%)
1	HYP	F	20[E]	1	8,8,9	7.04	1 (12%)	8,10,12	1.72	2 (25%)
1	HYP	F	20[F]	1	8,8,9	7.23	1 (12%)	8,10,12	1.73	2 (25%)
1	HYP	F	20[G]	1	8,8,9	7.03	1 (12%)	8,10,12	1.79	3 (37%)
1	HYP	F	23[A]	1	8,8,9	7.08	1 (12%)	8,10,12	1.58	2 (25%)
1	HYP	F	23[B]	1	8,8,9	7.11	1 (12%)	8,10,12	1.75	2 (25%)
1	HYP	F	23[C]	1	8,8,9	7.11	1 (12%)	8,10,12	1.74	2 (25%)
1	HYP	F	23[D]	1	8,8,9	7.06	1 (12%)	8,10,12	1.74	2 (25%)
1	HYP	F	23[E]	1	8,8,9	7.04	1 (12%)	8,10,12	1.60	2 (25%)
1	HYP	F	23[F]	1	8,8,9	7.13	1 (12%)	8,10,12	1.57	2 (25%)
1	HYP	F	23[G]	1	8,8,9	7.02	1 (12%)	8,10,12	1.90	2 (25%)
1	HYP	F	26[A]	1	8,8,9	6.92	1 (12%)	8,10,12	1.99	4 (50%)
1	HYP	F	26[B]	1	8,8,9	7.02	1 (12%)	8,10,12	1.66	2 (25%)
1	HYP	F	26[C]	1	8,8,9	7.02	1 (12%)	8,10,12	1.65	2 (25%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	HYP	F	26[D]	1	8,8,9	7.02	1 (12%)	8,10,12	1.67	2 (25%)
1	HYP	F	26[E]	1	8,8,9	6.90	1 (12%)	8,10,12	1.98	3 (37%)
1	HYP	F	26[F]	1	8,8,9	6.95	1 (12%)	8,10,12	2.01	4 (50%)
1	HYP	F	26[G]	1	8,8,9	6.40	1 (12%)	8,10,12	2.08	2 (25%)
1	HYP	F	29[A]	1	8,8,9	6.91	1 (12%)	8,10,12	1.98	2 (25%)
1	HYP	F	29[B]	1	8,8,9	7.01	1 (12%)	8,10,12	2.32	3 (37%)
1	HYP	F	29[C]	1	8,8,9	7.07	1 (12%)	8,10,12	2.34	3 (37%)
1	HYP	F	29[D]	1	8,8,9	6.95	1 (12%)	8,10,12	2.31	2 (25%)
1	HYP	F	29[E]	1	8,8,9	6.97	1 (12%)	8,10,12	1.96	2 (25%)
1	HYP	F	29[F]	1	8,8,9	6.82	1 (12%)	8,10,12	1.99	2 (25%)
1	HYP	F	29[G]	1	8,8,9	7.07	1 (12%)	8,10,12	1.85	3 (37%)
1	HYP	F	5[A]	1	8,8,9	6.62	1 (12%)	8,10,12	1.98	2 (25%)
1	HYP	F	5[B]	1	8,8,9	6.59	1 (12%)	8,10,12	1.82	3 (37%)
1	HYP	F	5[C]	1	8,8,9	6.50	1 (12%)	8,10,12	1.83	3 (37%)
1	HYP	F	5[D]	1	8,8,9	6.61	1 (12%)	8,10,12	1.79	3 (37%)
1	HYP	F	5[E]	1	8,8,9	6.73	1 (12%)	8,10,12	1.97	2 (25%)
1	HYP	F	5[F]	1	8,8,9	6.59	1 (12%)	8,10,12	1.98	2 (25%)
1	HYP	F	5[G]	1	8,8,9	7.22	1 (12%)	8,10,12	1.57	2 (25%)
1	HYP	F	8[A]	1	8,8,9	7.24	1 (12%)	8,10,12	1.58	2 (25%)
1	HYP	F	8[B]	1	8,8,9	6.81	1 (12%)	8,10,12	1.83	2 (25%)
1	HYP	F	8[C]	1	8,8,9	6.81	1 (12%)	8,10,12	1.82	2 (25%)
1	HYP	F	8[D]	1	8,8,9	6.81	1 (12%)	8,10,12	1.83	2 (25%)
1	HYP	F	8[E]	1	8,8,9	7.33	1 (12%)	8,10,12	1.59	2 (25%)
1	HYP	F	8[F]	1	8,8,9	7.14	1 (12%)	8,10,12	1.57	1 (12%)
1	HYP	F	8[G]	1	8,8,9	6.67	1 (12%)	8,10,12	2.13	4 (50%)
1	HYP	G	11[A]	1	8,8,9	7.05	1 (12%)	8,10,12	1.94	3 (37%)
1	HYP	G	11[B]	1	8,8,9	7.07	1 (12%)	8,10,12	1.94	3 (37%)
1	HYP	G	11[C]	1	8,8,9	7.09	1 (12%)	8,10,12	1.94	3 (37%)
1	HYP	G	11[D]	1	8,8,9	7.07	2 (25%)	8,10,12	2.07	4 (50%)
1	HYP	G	11[E]	1	8,8,9	6.87	1 (12%)	8,10,12	2.12	3 (37%)
1	HYP	G	11[F]	1	8,8,9	6.87	2 (25%)	8,10,12	2.11	3 (37%)
1	HYP	G	11[G]	1	8,8,9	6.87	1 (12%)	8,10,12	2.13	3 (37%)
1	HZP	G	13[A]	1	8,8,9	6.51	1 (12%)	8,10,12	4.15	5 (62%)
1	HZP	G	13[B]	1	8,8,9	6.46	1 (12%)	8,10,12	4.14	5 (62%)
1	HZP	G	13[C]	1	8,8,9	6.56	1 (12%)	8,10,12	4.16	5 (62%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	HZP	G	13[D]	1	8,8,9	7.49	1 (12%)	8,10,12	5.06	7 (87%)
1	HZP	G	13[E]	1	8,8,9	6.34	2 (25%)	8,10,12	4.06	5 (62%)
1	HZP	G	13[F]	1	8,8,9	6.34	2 (25%)	8,10,12	4.06	5 (62%)
1	HZP	G	13[G]	1	8,8,9	6.34	2 (25%)	8,10,12	4.05	5 (62%)
1	HZP	G	16[A]	1	8,8,9	6.06	2 (25%)	8,10,12	4.09	6 (75%)
1	HZP	G	16[B]	1	8,8,9	6.04	2 (25%)	8,10,12	4.11	6 (75%)
1	HZP	G	16[C]	1	8,8,9	6.01	2 (25%)	8,10,12	4.09	6 (75%)
1	HZP	G	16[D]	1	8,8,9	6.41	1 (12%)	8,10,12	4.12	5 (62%)
1	HZP	G	16[E]	1	8,8,9	6.86	2 (25%)	8,10,12	4.38	5 (62%)
1	HZP	G	16[F]	1	8,8,9	6.86	2 (25%)	8,10,12	4.40	5 (62%)
1	HZP	G	16[G]	1	8,8,9	6.86	2 (25%)	8,10,12	4.36	5 (62%)
1	HYP	G	2[A]	1	8,8,9	6.76	1 (12%)	8,10,12	1.96	2 (25%)
1	HYP	G	2[B]	1	8,8,9	6.64	1 (12%)	8,10,12	1.92	2 (25%)
1	HYP	G	2[C]	1	8,8,9	6.75	1 (12%)	8,10,12	1.96	2 (25%)
1	HYP	G	2[D]	1	8,8,9	6.73	1 (12%)	8,10,12	1.78	3 (37%)
1	HYP	G	2[E]	1	8,8,9	6.62	1 (12%)	8,10,12	2.34	2 (25%)
1	HYP	G	2[F]	1	8,8,9	6.56	1 (12%)	8,10,12	2.34	2 (25%)
1	HYP	G	2[G]	1	8,8,9	6.59	1 (12%)	8,10,12	2.34	2 (25%)
1	HYP	G	20[A]	1	8,8,9	7.44	1 (12%)	8,10,12	2.37	4 (50%)
1	HYP	G	20[B]	1	8,8,9	7.53	1 (12%)	8,10,12	2.37	4 (50%)
1	HYP	G	20[C]	1	8,8,9	7.39	1 (12%)	8,10,12	2.37	4 (50%)
1	HYP	G	20[D]	1	8,8,9	6.52	1 (12%)	8,10,12	2.38	3 (37%)
1	HYP	G	20[E]	1	8,8,9	6.82	1 (12%)	8,10,12	1.69	2 (25%)
1	HYP	G	20[F]	1	8,8,9	6.82	1 (12%)	8,10,12	1.69	2 (25%)
1	HYP	G	20[G]	1	8,8,9	6.83	1 (12%)	8,10,12	1.70	2 (25%)
1	HYP	G	23[A]	1	8,8,9	6.78	1 (12%)	8,10,12	1.83	2 (25%)
1	HYP	G	23[B]	1	8,8,9	6.80	1 (12%)	8,10,12	1.78	2 (25%)
1	HYP	G	23[C]	1	8,8,9	6.73	1 (12%)	8,10,12	1.81	2 (25%)
1	HYP	G	23[D]	1	8,8,9	7.40	1 (12%)	8,10,12	2.38	4 (50%)
1	HYP	G	23[E]	1	8,8,9	6.81	1 (12%)	8,10,12	1.71	2 (25%)
1	HYP	G	23[F]	1	8,8,9	6.85	1 (12%)	8,10,12	1.71	2 (25%)
1	HYP	G	23[G]	1	8,8,9	6.78	1 (12%)	8,10,12	1.71	2 (25%)
1	HYP	G	26[A]	1	8,8,9	6.80	1 (12%)	8,10,12	1.67	2 (25%)
1	HYP	G	26[B]	1	8,8,9	6.83	1 (12%)	8,10,12	1.67	2 (25%)
1	HYP	G	26[C]	1	8,8,9	6.77	1 (12%)	8,10,12	1.67	2 (25%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	HYP	G	26[D]	1	8,8,9	6.83	1 (12%)	8,10,12	1.80	2 (25%)
1	HYP	G	26[E]	1	8,8,9	7.18	1 (12%)	8,10,12	1.70	3 (37%)
1	HYP	G	26[F]	1	8,8,9	6.91	1 (12%)	8,10,12	1.91	2 (25%)
1	HYP	G	26[G]	1	8,8,9	7.21	1 (12%)	8,10,12	1.70	3 (37%)
1	HYP	G	29[A]	1	8,8,9	7.07	1 (12%)	8,10,12	1.73	2 (25%)
1	HYP	G	29[B]	1	8,8,9	7.04	1 (12%)	8,10,12	1.74	2 (25%)
1	HYP	G	29[C]	1	8,8,9	7.04	1 (12%)	8,10,12	1.72	2 (25%)
1	HYP	G	29[D]	1	8,8,9	6.83	1 (12%)	8,10,12	1.69	2 (25%)
1	HYP	G	29[E]	1	8,8,9	6.52	1 (12%)	8,10,12	2.20	4 (50%)
1	HYP	G	29[F]	1	8,8,9	6.52	1 (12%)	8,10,12	2.20	4 (50%)
1	HYP	G	29[G]	1	8,8,9	6.52	1 (12%)	8,10,12	2.19	4 (50%)
1	HYP	G	5[A]	1	8,8,9	6.73	1 (12%)	8,10,12	2.23	4 (50%)
1	HYP	G	5[B]	1	8,8,9	6.73	1 (12%)	8,10,12	2.23	4 (50%)
1	HYP	G	5[C]	1	8,8,9	6.73	1 (12%)	8,10,12	2.22	4 (50%)
1	HYP	G	5[D]	1	8,8,9	6.75	1 (12%)	8,10,12	1.96	2 (25%)
1	HYP	G	5[E]	1	8,8,9	6.99	2 (25%)	8,10,12	1.84	2 (25%)
1	HYP	G	5[F]	1	8,8,9	6.89	2 (25%)	8,10,12	1.84	2 (25%)
1	HYP	G	5[G]	1	8,8,9	7.09	2 (25%)	8,10,12	1.84	2 (25%)
1	HYP	G	8[A]	1	8,8,9	6.99	2 (25%)	8,10,12	1.86	2 (25%)
1	HYP	G	8[B]	1	8,8,9	7.02	2 (25%)	8,10,12	1.83	2 (25%)
1	HYP	G	8[C]	1	8,8,9	7.03	1 (12%)	8,10,12	1.87	2 (25%)
1	HYP	G	8[D]	1	8,8,9	3.83	1 (12%)	8,10,12	1.62	2 (25%)
1	HYP	G	8[E]	1	8,8,9	6.99	1 (12%)	8,10,12	1.95	3 (37%)
1	HYP	G	8[F]	1	8,8,9	7.01	1 (12%)	8,10,12	1.94	3 (37%)
1	HYP	G	8[G]	1	8,8,9	6.97	1 (12%)	8,10,12	1.97	3 (37%)
1	HYP	H	11[A]	1	8,8,9	7.35	1 (12%)	8,10,12	1.71	2 (25%)
1	HYP	H	11[B]	1	8,8,9	7.35	1 (12%)	8,10,12	1.67	2 (25%)
1	HYP	H	11[C]	1	8,8,9	6.91	1 (12%)	8,10,12	1.68	2 (25%)
1	HYP	H	11[D]	1	8,8,9	6.91	1 (12%)	8,10,12	1.70	3 (37%)
1	HYP	H	11[E]	1	8,8,9	7.35	1 (12%)	8,10,12	1.67	2 (25%)
1	HYP	H	11[F]	1	8,8,9	6.99	1 (12%)	8,10,12	2.33	4 (50%)
1	HYP	H	11[G]	1	8,8,9	6.89	1 (12%)	8,10,12	2.35	4 (50%)
1	HZP	H	13[A]	1	8,8,9	7.07	1 (12%)	8,10,12	4.37	5 (62%)
1	HZP	H	13[B]	1	8,8,9	7.06	1 (12%)	8,10,12	4.38	5 (62%)
1	HZP	H	13[C]	1	8,8,9	7.10	1 (12%)	8,10,12	4.19	6 (75%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	HZP	H	13[D]	1	8,8,9	7.10	1 (12%)	8,10,12	4.21	6 (75%)
1	HZP	H	13[E]	1	8,8,9	7.07	1 (12%)	8,10,12	4.37	5 (62%)
1	HZP	H	13[F]	1	8,8,9	6.63	1 (12%)	8,10,12	4.00	5 (62%)
1	HZP	H	13[G]	1	8,8,9	6.70	1 (12%)	8,10,12	4.01	5 (62%)
1	HZP	H	16[A]	1	8,8,9	6.55	1 (12%)	8,10,12	3.68	6 (75%)
1	HZP	H	16[B]	1	8,8,9	6.63	1 (12%)	8,10,12	3.70	6 (75%)
1	HZP	H	16[C]	1	8,8,9	7.06	1 (12%)	8,10,12	4.16	5 (62%)
1	HZP	H	16[D]	1	8,8,9	7.05	1 (12%)	8,10,12	4.16	5 (62%)
1	HZP	H	16[E]	1	8,8,9	6.55	1 (12%)	8,10,12	3.68	6 (75%)
1	HZP	H	16[F]	1	8,8,9	7.04	1 (12%)	8,10,12	4.56	5 (62%)
1	HZP	H	16[G]	1	8,8,9	6.94	1 (12%)	8,10,12	4.57	5 (62%)
1	HYP	H	2[A]	1	8,8,9	7.04	1 (12%)	8,10,12	2.07	4 (50%)
1	HYP	H	2[B]	1	8,8,9	7.04	1 (12%)	8,10,12	2.06	4 (50%)
1	HYP	H	2[C]	1	8,8,9	6.60	1 (12%)	8,10,12	2.05	3 (37%)
1	HYP	H	2[D]	1	8,8,9	6.60	1 (12%)	8,10,12	2.05	3 (37%)
1	HYP	H	2[E]	1	8,8,9	7.04	1 (12%)	8,10,12	2.07	4 (50%)
1	HYP	H	2[F]	1	8,8,9	7.32	1 (12%)	8,10,12	1.60	2 (25%)
1	HYP	H	2[G]	1	8,8,9	7.32	1 (12%)	8,10,12	1.60	2 (25%)
1	HYP	H	20[A]	1	8,8,9	6.09	1 (12%)	8,10,12	1.94	2 (25%)
1	HYP	H	20[B]	1	8,8,9	6.10	1 (12%)	8,10,12	1.96	2 (25%)
1	HYP	H	20[C]	1	8,8,9	7.06	1 (12%)	8,10,12	1.86	3 (37%)
1	HYP	H	20[D]	1	8,8,9	7.06	1 (12%)	8,10,12	1.84	3 (37%)
1	HYP	H	20[E]	1	8,8,9	6.09	1 (12%)	8,10,12	1.94	2 (25%)
1	HYP	H	20[F]	1	8,8,9	6.64	1 (12%)	8,10,12	1.72	2 (25%)
1	HYP	H	20[G]	1	8,8,9	6.64	1 (12%)	8,10,12	1.73	2 (25%)
1	HYP	H	23[A]	1	8,8,9	6.67	1 (12%)	8,10,12	1.81	2 (25%)
1	HYP	H	23[B]	1	8,8,9	6.66	1 (12%)	8,10,12	1.78	2 (25%)
1	HYP	H	23[C]	1	8,8,9	6.13	1 (12%)	8,10,12	1.91	3 (37%)
1	HYP	H	23[D]	1	8,8,9	6.13	1 (12%)	8,10,12	1.91	3 (37%)
1	HYP	H	23[E]	1	8,8,9	6.62	1 (12%)	8,10,12	1.77	2 (25%)
1	HYP	H	23[F]	1	8,8,9	7.28	1 (12%)	8,10,12	1.61	3 (37%)
1	HYP	H	23[G]	1	8,8,9	7.28	1 (12%)	8,10,12	1.61	3 (37%)
1	HYP	H	26[A]	1	8,8,9	7.31	1 (12%)	8,10,12	1.65	3 (37%)
1	HYP	H	26[B]	1	8,8,9	7.31	1 (12%)	8,10,12	1.64	3 (37%)
1	HYP	H	26[C]	1	8,8,9	6.60	1 (12%)	8,10,12	1.75	2 (25%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	HYP	H	26[D]	1	8,8,9	6.59	1 (12%)	8,10,12	1.81	2 (25%)
1	HYP	H	26[E]	1	8,8,9	7.31	1 (12%)	8,10,12	1.63	3 (37%)
1	HYP	H	26[F]	1	8,8,9	7.14	1 (12%)	8,10,12	1.96	2 (25%)
1	HYP	H	26[G]	1	8,8,9	7.08	1 (12%)	8,10,12	1.96	2 (25%)
1	HYP	H	29[A]	1	8,8,9	6.82	1 (12%)	8,10,12	2.19	3 (37%)
1	HYP	H	29[B]	1	8,8,9	6.77	1 (12%)	8,10,12	2.19	3 (37%)
1	HYP	H	29[C]	1	8,8,9	7.32	1 (12%)	8,10,12	1.69	3 (37%)
1	HYP	H	29[D]	1	8,8,9	7.39	1 (12%)	8,10,12	1.70	3 (37%)
1	HYP	H	29[E]	1	8,8,9	6.82	1 (12%)	8,10,12	2.19	2 (25%)
1	HYP	H	29[F]	1	8,8,9	7.26	1 (12%)	8,10,12	1.71	3 (37%)
1	HYP	H	29[G]	1	8,8,9	7.24	1 (12%)	8,10,12	1.69	3 (37%)
1	HYP	H	5[A]	1	8,8,9	7.32	1 (12%)	8,10,12	1.61	2 (25%)
1	HYP	H	5[B]	1	8,8,9	7.32	1 (12%)	8,10,12	1.60	2 (25%)
1	HYP	H	5[C]	1	8,8,9	4.96	1 (12%)	8,10,12	1.66	2 (25%)
1	HYP	H	5[D]	1	8,8,9	5.02	1 (12%)	8,10,12	1.66	2 (25%)
1	HYP	H	5[E]	1	8,8,9	7.32	1 (12%)	8,10,12	1.61	2 (25%)
1	HYP	H	5[F]	1	8,8,9	7.10	1 (12%)	8,10,12	1.81	3 (37%)
1	HYP	H	5[G]	1	8,8,9	7.10	1 (12%)	8,10,12	1.81	3 (37%)
1	HYP	H	8[A]	1	8,8,9	7.12	1 (12%)	8,10,12	1.83	3 (37%)
1	HYP	H	8[B]	1	8,8,9	7.16	1 (12%)	8,10,12	1.83	3 (37%)
1	HYP	H	8[C]	1	8,8,9	7.35	1 (12%)	8,10,12	1.60	2 (25%)
1	HYP	H	8[D]	1	8,8,9	7.35	1 (12%)	8,10,12	1.61	2 (25%)
1	HYP	H	8[E]	1	8,8,9	7.12	1 (12%)	8,10,12	1.83	3 (37%)
1	HYP	H	8[F]	1	8,8,9	7.20	1 (12%)	8,10,12	1.64	2 (25%)
1	HYP	H	8[G]	1	8,8,9	7.17	1 (12%)	8,10,12	1.64	2 (25%)
1	HYP	I	11[A]	1	8,8,9	7.33	1 (12%)	8,10,12	1.72	2 (25%)
1	HYP	I	11[B]	1	8,8,9	6.41	1 (12%)	8,10,12	1.92	4 (50%)
1	HYP	I	11[C]	1	8,8,9	6.37	1 (12%)	8,10,12	1.91	3 (37%)
1	HYP	I	11[D]	1	8,8,9	6.44	1 (12%)	8,10,12	1.90	3 (37%)
1	HYP	I	11[E]	1	8,8,9	7.24	1 (12%)	8,10,12	1.72	2 (25%)
1	HYP	I	11[F]	1	8,8,9	7.14	1 (12%)	8,10,12	1.72	2 (25%)
1	HYP	I	11[G]	1	8,8,9	6.69	1 (12%)	8,10,12	2.14	4 (50%)
1	HZP	I	13[A]	1	8,8,9	6.34	1 (12%)	8,10,12	3.95	5 (62%)
1	HZP	I	13[B]	1	8,8,9	6.70	1 (12%)	8,10,12	4.83	5 (62%)
1	HZP	I	13[C]	1	8,8,9	6.70	1 (12%)	8,10,12	4.81	5 (62%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	HZP	I	13[D]	1	8,8,9	6.70	1 (12%)	8,10,12	4.81	5 (62%)
1	HZP	I	13[E]	1	8,8,9	6.34	1 (12%)	8,10,12	3.96	5 (62%)
1	HZP	I	13[F]	1	8,8,9	6.34	1 (12%)	8,10,12	3.97	5 (62%)
1	HZP	I	13[G]	1	8,8,9	7.02	1 (12%)	8,10,12	4.32	6 (75%)
1	HZP	I	16[A]	1	8,8,9	6.97	1 (12%)	8,10,12	4.37	6 (75%)
1	HZP	I	16[B]	1	8,8,9	6.42	1 (12%)	8,10,12	4.15	5 (62%)
1	HZP	I	16[C]	1	8,8,9	6.42	1 (12%)	8,10,12	4.16	5 (62%)
1	HZP	I	16[D]	1	8,8,9	6.57	1 (12%)	8,10,12	4.16	5 (62%)
1	HZP	I	16[E]	1	8,8,9	6.97	1 (12%)	8,10,12	4.39	6 (75%)
1	HZP	I	16[F]	1	8,8,9	7.03	1 (12%)	8,10,12	4.39	7 (87%)
1	HZP	I	16[G]	1	8,8,9	6.91	2 (25%)	8,10,12	4.53	6 (75%)
1	HYP	I	2[A]	1	8,8,9	7.03	1 (12%)	8,10,12	1.89	4 (50%)
1	HYP	I	2[B]	1	8,8,9	6.24	1 (12%)	8,10,12	1.79	3 (37%)
1	HYP	I	2[C]	1	8,8,9	6.43	1 (12%)	8,10,12	1.85	3 (37%)
1	HYP	I	2[D]	1	8,8,9	6.43	1 (12%)	8,10,12	1.85	3 (37%)
1	HYP	I	2[E]	1	8,8,9	7.03	1 (12%)	8,10,12	1.90	4 (50%)
1	HYP	I	2[F]	1	8,8,9	7.03	1 (12%)	8,10,12	1.90	4 (50%)
1	HYP	I	2[G]	1	8,8,9	6.36	1 (12%)	8,10,12	1.80	2 (25%)
1	HYP	I	20[A]	1	8,8,9	6.36	1 (12%)	8,10,12	1.94	3 (37%)
1	HYP	I	20[B]	1	8,8,9	6.73	1 (12%)	8,10,12	2.33	3 (37%)
1	HYP	I	20[C]	1	8,8,9	6.73	1 (12%)	8,10,12	2.31	2 (25%)
1	HYP	I	20[D]	1	8,8,9	6.73	1 (12%)	8,10,12	2.29	2 (25%)
1	HYP	I	20[E]	1	8,8,9	6.36	1 (12%)	8,10,12	1.95	3 (37%)
1	HYP	I	20[F]	1	8,8,9	6.41	1 (12%)	8,10,12	1.96	3 (37%)
1	HYP	I	20[G]	1	8,8,9	7.31	1 (12%)	8,10,12	1.35	1 (12%)
1	HYP	I	23[A]	1	8,8,9	7.28	1 (12%)	8,10,12	1.40	1 (12%)
1	HYP	I	23[B]	1	8,8,9	6.38	1 (12%)	8,10,12	1.93	3 (37%)
1	HYP	I	23[C]	1	8,8,9	6.37	1 (12%)	8,10,12	1.96	3 (37%)
1	HYP	I	23[D]	1	8,8,9	6.29	1 (12%)	8,10,12	1.95	3 (37%)
1	HYP	I	23[E]	1	8,8,9	7.17	1 (12%)	8,10,12	1.42	1 (12%)
1	HYP	I	23[F]	1	8,8,9	7.17	1 (12%)	8,10,12	1.40	1 (12%)
1	HYP	I	23[G]	1	8,8,9	6.48	1 (12%)	8,10,12	1.89	2 (25%)
1	HYP	I	26[A]	1	8,8,9	6.59	1 (12%)	8,10,12	2.16	2 (25%)
1	HYP	I	26[B]	1	8,8,9	7.22	1 (12%)	8,10,12	1.40	1 (12%)
1	HYP	I	26[C]	1	8,8,9	7.24	1 (12%)	8,10,12	1.38	1 (12%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	HYP	I	26[D]	1	8,8,9	7.26	1 (12%)	8,10,12	1.36	1 (12%)
1	HYP	I	26[E]	1	8,8,9	6.59	1 (12%)	8,10,12	2.15	2 (25%)
1	HYP	I	26[F]	1	8,8,9	6.59	1 (12%)	8,10,12	2.15	2 (25%)
1	HYP	I	26[G]	1	8,8,9	6.70	1 (12%)	8,10,12	2.16	4 (50%)
1	HYP	I	29[A]	1	8,8,9	7.05	1 (12%)	8,10,12	2.10	3 (37%)
1	HYP	I	29[B]	1	8,8,9	6.37	1 (12%)	8,10,12	1.87	2 (25%)
1	HYP	I	29[C]	1	8,8,9	6.37	1 (12%)	8,10,12	1.87	2 (25%)
1	HYP	I	29[D]	1	8,8,9	6.37	1 (12%)	8,10,12	1.87	2 (25%)
1	HYP	I	29[E]	1	8,8,9	7.01	1 (12%)	8,10,12	2.08	3 (37%)
1	HYP	I	29[F]	1	8,8,9	7.06	1 (12%)	8,10,12	2.10	3 (37%)
1	HYP	I	29[G]	1	8,8,9	7.08	1 (12%)	8,10,12	2.01	3 (37%)
1	HYP	I	5[A]	1	8,8,9	6.36	1 (12%)	8,10,12	1.80	2 (25%)
1	HYP	I	5[B]	1	8,8,9	6.93	1 (12%)	8,10,12	2.10	4 (50%)
1	HYP	I	5[C]	1	8,8,9	6.88	1 (12%)	8,10,12	2.11	4 (50%)
1	HYP	I	5[D]	1	8,8,9	6.83	1 (12%)	8,10,12	2.11	4 (50%)
1	HYP	I	5[E]	1	8,8,9	6.37	1 (12%)	8,10,12	1.81	2 (25%)
1	HYP	I	5[F]	1	8,8,9	6.37	1 (12%)	8,10,12	1.82	2 (25%)
1	HYP	I	5[G]	1	8,8,9	6.43	1 (12%)	8,10,12	1.96	4 (50%)
1	HYP	I	8[A]	1	8,8,9	6.33	1 (12%)	8,10,12	1.97	4 (50%)
1	HYP	I	8[B]	1	8,8,9	6.37	1 (12%)	8,10,12	1.82	2 (25%)
1	HYP	I	8[C]	1	8,8,9	6.36	1 (12%)	8,10,12	1.83	2 (25%)
1	HYP	I	8[D]	1	8,8,9	6.35	1 (12%)	8,10,12	1.81	2 (25%)
1	HYP	I	8[E]	1	8,8,9	6.36	1 (12%)	8,10,12	1.97	4 (50%)
1	HYP	I	8[F]	1	8,8,9	6.43	1 (12%)	8,10,12	1.97	4 (50%)
1	HYP	I	8[G]	1	8,8,9	7.09	1 (12%)	8,10,12	1.85	2 (25%)

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	11[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	11[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	11[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	11[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	11[E]	1	-	0/0/11/13	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	HYP	A	11[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	11[G]	1	-	0/0/11/13	0/1/1/1
1	HZP	A	13[A]	1	-	0/0/11/13	0/1/1/1
1	HZP	A	13[B]	1	-	0/0/11/13	0/1/1/1
1	HZP	A	13[C]	1	-	0/0/11/13	0/1/1/1
1	HZP	A	13[D]	1	-	0/0/11/13	0/1/1/1
1	HZP	A	13[E]	1	-	0/0/11/13	0/1/1/1
1	HZP	A	13[F]	1	-	0/0/11/13	0/1/1/1
1	HZP	A	13[G]	1	-	0/0/11/13	0/1/1/1
1	HZP	A	16[A]	1	-	0/0/11/13	0/1/1/1
1	HZP	A	16[B]	1	-	0/0/11/13	0/1/1/1
1	HZP	A	16[C]	1	-	0/0/11/13	0/1/1/1
1	HZP	A	16[D]	1	-	0/0/11/13	0/1/1/1
1	HZP	A	16[E]	1	-	0/0/11/13	0/1/1/1
1	HZP	A	16[F]	1	-	0/0/11/13	0/1/1/1
1	HZP	A	16[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	2[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	2[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	2[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	2[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	2[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	2[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	2[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	20[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	20[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	20[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	20[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	20[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	20[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	20[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	23[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	23[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	23[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	23[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	23[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	23[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	23[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	26[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	26[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	26[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	26[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	26[E]	1	-	0/0/11/13	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	HYP	A	26[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	26[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	29[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	29[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	29[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	29[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	29[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	29[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	29[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	5[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	5[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	5[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	5[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	5[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	5[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	5[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	8[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	8[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	8[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	8[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	8[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	8[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	A	8[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	11[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	11[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	11[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	11[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	11[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	11[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	11[G]	1	-	0/0/11/13	0/1/1/1
1	HZP	B	13[A]	1	-	0/0/11/13	0/1/1/1
1	HZP	B	13[B]	1	-	0/0/11/13	0/1/1/1
1	HZP	B	13[C]	1	-	0/0/11/13	0/1/1/1
1	HZP	B	13[D]	1	-	0/0/11/13	0/1/1/1
1	HZP	B	13[E]	1	-	0/0/11/13	0/1/1/1
1	HZP	B	13[F]	1	-	0/0/11/13	0/1/1/1
1	HZP	B	13[G]	1	-	0/0/11/13	0/1/1/1
1	HZP	B	16[A]	1	-	0/0/11/13	0/1/1/1
1	HZP	B	16[B]	1	-	0/0/11/13	0/1/1/1
1	HZP	B	16[C]	1	-	0/0/11/13	0/1/1/1
1	HZP	B	16[D]	1	-	0/0/11/13	0/1/1/1
1	HZP	B	16[E]	1	-	0/0/11/13	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	HZP	B	16[F]	1	-	0/0/11/13	0/1/1/1
1	HZP	B	16[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	2[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	2[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	2[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	2[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	2[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	2[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	2[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	20[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	20[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	20[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	20[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	20[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	20[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	20[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	23[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	23[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	23[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	23[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	23[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	23[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	23[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	26[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	26[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	26[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	26[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	26[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	26[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	26[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	29[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	29[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	29[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	29[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	29[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	29[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	29[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	5[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	5[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	5[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	5[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	5[E]	1	-	0/0/11/13	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	HYP	B	5[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	5[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	8[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	8[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	8[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	8[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	8[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	8[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	B	8[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	11[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	11[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	11[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	11[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	11[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	11[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	11[G]	1	-	0/0/11/13	0/1/1/1
1	HZP	C	13[A]	1	-	0/0/11/13	0/1/1/1
1	HZP	C	13[B]	1	-	0/0/11/13	0/1/1/1
1	HZP	C	13[C]	1	-	0/0/11/13	0/1/1/1
1	HZP	C	13[D]	1	-	0/0/11/13	0/1/1/1
1	HZP	C	13[E]	1	-	0/0/11/13	0/1/1/1
1	HZP	C	13[F]	1	-	0/0/11/13	0/1/1/1
1	HZP	C	13[G]	1	-	0/0/11/13	0/1/1/1
1	HZP	C	16[A]	1	-	0/0/11/13	0/1/1/1
1	HZP	C	16[B]	1	-	0/0/11/13	0/1/1/1
1	HZP	C	16[C]	1	-	0/0/11/13	0/1/1/1
1	HZP	C	16[D]	1	-	0/0/11/13	0/1/1/1
1	HZP	C	16[E]	1	-	0/0/11/13	0/1/1/1
1	HZP	C	16[F]	1	-	0/0/11/13	0/1/1/1
1	HZP	C	16[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	2[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	2[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	2[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	2[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	2[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	2[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	2[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	20[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	20[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	20[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	20[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	20[E]	1	-	0/0/11/13	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	HYP	C	20[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	20[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	23[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	23[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	23[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	23[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	23[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	23[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	23[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	26[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	26[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	26[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	26[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	26[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	26[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	26[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	29[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	29[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	29[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	29[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	29[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	29[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	29[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	5[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	5[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	5[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	5[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	5[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	5[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	5[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	8[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	8[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	8[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	8[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	8[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	8[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	C	8[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	11[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	11[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	11[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	11[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	11[E]	1	-	0/0/11/13	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	HYP	D	11[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	11[G]	1	-	0/0/11/13	0/1/1/1
1	HZP	D	13[A]	1	-	0/0/11/13	0/1/1/1
1	HZP	D	13[B]	1	-	0/0/11/13	0/1/1/1
1	HZP	D	13[C]	1	-	0/0/11/13	0/1/1/1
1	HZP	D	13[D]	1	-	0/0/11/13	0/1/1/1
1	HZP	D	13[E]	1	-	0/0/11/13	0/1/1/1
1	HZP	D	13[F]	1	-	0/0/11/13	0/1/1/1
1	HZP	D	13[G]	1	-	0/0/11/13	0/1/1/1
1	HZP	D	16[A]	1	-	0/0/11/13	0/1/1/1
1	HZP	D	16[B]	1	-	0/0/11/13	0/1/1/1
1	HZP	D	16[C]	1	-	0/0/11/13	0/1/1/1
1	HZP	D	16[D]	1	-	0/0/11/13	0/1/1/1
1	HZP	D	16[E]	1	-	0/0/11/13	0/1/1/1
1	HZP	D	16[F]	1	-	0/0/11/13	0/1/1/1
1	HZP	D	16[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	2[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	2[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	2[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	2[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	2[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	2[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	2[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	20[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	20[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	20[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	20[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	20[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	20[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	20[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	23[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	23[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	23[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	23[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	23[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	23[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	23[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	26[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	26[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	26[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	26[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	26[E]	1	-	0/0/11/13	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	HYP	D	26[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	26[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	29[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	29[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	29[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	29[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	29[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	29[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	29[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	5[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	5[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	5[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	5[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	5[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	5[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	5[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	8[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	8[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	8[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	8[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	8[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	8[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	D	8[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	11[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	11[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	11[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	11[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	11[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	11[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	11[G]	1	-	0/0/11/13	0/1/1/1
1	HZP	E	13[A]	1	-	0/0/11/13	0/1/1/1
1	HZP	E	13[B]	1	-	0/0/11/13	0/1/1/1
1	HZP	E	13[C]	1	-	0/0/11/13	0/1/1/1
1	HZP	E	13[D]	1	-	0/0/11/13	0/1/1/1
1	HZP	E	13[E]	1	-	0/0/11/13	0/1/1/1
1	HZP	E	13[F]	1	-	0/0/11/13	0/1/1/1
1	HZP	E	13[G]	1	-	0/0/11/13	0/1/1/1
1	HZP	E	16[A]	1	-	0/0/11/13	0/1/1/1
1	HZP	E	16[B]	1	-	0/0/11/13	0/1/1/1
1	HZP	E	16[C]	1	-	0/0/11/13	0/1/1/1
1	HZP	E	16[D]	1	-	0/0/11/13	0/1/1/1
1	HZP	E	16[E]	1	-	0/0/11/13	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	HZP	E	16[F]	1	-	0/0/11/13	0/1/1/1
1	HZP	E	16[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	2[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	2[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	2[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	2[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	2[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	2[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	2[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	20[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	20[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	20[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	20[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	20[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	20[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	20[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	23[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	23[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	23[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	23[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	23[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	23[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	23[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	26[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	26[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	26[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	26[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	26[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	26[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	26[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	29[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	29[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	29[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	29[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	29[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	29[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	29[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	5[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	5[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	5[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	5[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	5[E]	1	-	0/0/11/13	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	HYP	E	5[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	5[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	8[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	8[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	8[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	8[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	8[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	8[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	E	8[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	11[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	11[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	11[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	11[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	11[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	11[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	11[G]	1	-	0/0/11/13	0/1/1/1
1	HZP	F	13[A]	1	-	0/0/11/13	0/1/1/1
1	HZP	F	13[B]	1	-	0/0/11/13	0/1/1/1
1	HZP	F	13[C]	1	-	0/0/11/13	0/1/1/1
1	HZP	F	13[D]	1	-	0/0/11/13	0/1/1/1
1	HZP	F	13[E]	1	-	0/0/11/13	0/1/1/1
1	HZP	F	13[F]	1	-	0/0/11/13	0/1/1/1
1	HZP	F	13[G]	1	-	0/0/11/13	0/1/1/1
1	HZP	F	16[A]	1	-	0/0/11/13	0/1/1/1
1	HZP	F	16[B]	1	-	0/0/11/13	0/1/1/1
1	HZP	F	16[C]	1	-	0/0/11/13	0/1/1/1
1	HZP	F	16[D]	1	-	0/0/11/13	0/1/1/1
1	HZP	F	16[E]	1	-	0/0/11/13	0/1/1/1
1	HZP	F	16[F]	1	-	0/0/11/13	0/1/1/1
1	HZP	F	16[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	2[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	2[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	2[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	2[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	2[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	2[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	2[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	20[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	20[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	20[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	20[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	20[E]	1	-	0/0/11/13	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	HYP	F	20[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	20[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	23[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	23[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	23[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	23[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	23[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	23[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	23[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	26[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	26[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	26[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	26[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	26[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	26[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	26[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	29[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	29[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	29[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	29[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	29[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	29[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	29[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	5[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	5[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	5[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	5[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	5[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	5[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	5[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	8[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	8[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	8[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	8[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	8[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	8[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	F	8[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	11[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	11[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	11[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	11[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	11[E]	1	-	0/0/11/13	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	HYP	G	11[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	11[G]	1	-	0/0/11/13	0/1/1/1
1	HZP	G	13[A]	1	-	0/0/11/13	0/1/1/1
1	HZP	G	13[B]	1	-	0/0/11/13	0/1/1/1
1	HZP	G	13[C]	1	-	0/0/11/13	0/1/1/1
1	HZP	G	13[D]	1	-	0/0/11/13	0/1/1/1
1	HZP	G	13[E]	1	-	0/0/11/13	0/1/1/1
1	HZP	G	13[F]	1	-	0/0/11/13	0/1/1/1
1	HZP	G	13[G]	1	-	0/0/11/13	0/1/1/1
1	HZP	G	16[A]	1	-	0/0/11/13	0/1/1/1
1	HZP	G	16[B]	1	-	0/0/11/13	0/1/1/1
1	HZP	G	16[C]	1	-	0/0/11/13	0/1/1/1
1	HZP	G	16[D]	1	-	0/0/11/13	0/1/1/1
1	HZP	G	16[E]	1	-	0/0/11/13	0/1/1/1
1	HZP	G	16[F]	1	-	0/0/11/13	0/1/1/1
1	HZP	G	16[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	2[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	2[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	2[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	2[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	2[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	2[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	2[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	20[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	20[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	20[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	20[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	20[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	20[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	20[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	23[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	23[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	23[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	23[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	23[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	23[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	23[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	26[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	26[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	26[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	26[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	26[E]	1	-	0/0/11/13	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	HYP	G	26[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	26[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	29[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	29[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	29[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	29[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	29[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	29[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	29[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	5[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	5[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	5[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	5[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	5[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	5[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	5[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	8[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	8[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	8[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	8[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	8[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	8[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	G	8[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	11[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	11[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	11[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	11[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	11[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	11[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	11[G]	1	-	0/0/11/13	0/1/1/1
1	HZP	H	13[A]	1	-	0/0/11/13	0/1/1/1
1	HZP	H	13[B]	1	-	0/0/11/13	0/1/1/1
1	HZP	H	13[C]	1	-	0/0/11/13	0/1/1/1
1	HZP	H	13[D]	1	-	0/0/11/13	0/1/1/1
1	HZP	H	13[E]	1	-	0/0/11/13	0/1/1/1
1	HZP	H	13[F]	1	-	0/0/11/13	0/1/1/1
1	HZP	H	13[G]	1	-	0/0/11/13	0/1/1/1
1	HZP	H	16[A]	1	-	0/0/11/13	0/1/1/1
1	HZP	H	16[B]	1	-	0/0/11/13	0/1/1/1
1	HZP	H	16[C]	1	-	0/0/11/13	0/1/1/1
1	HZP	H	16[D]	1	-	0/0/11/13	0/1/1/1
1	HZP	H	16[E]	1	-	0/0/11/13	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	HZP	H	16[F]	1	-	0/0/11/13	0/1/1/1
1	HZP	H	16[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	2[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	2[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	2[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	2[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	2[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	2[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	2[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	20[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	20[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	20[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	20[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	20[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	20[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	20[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	23[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	23[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	23[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	23[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	23[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	23[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	23[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	26[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	26[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	26[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	26[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	26[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	26[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	26[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	29[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	29[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	29[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	29[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	29[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	29[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	29[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	5[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	5[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	5[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	5[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	5[E]	1	-	0/0/11/13	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	HYP	H	5[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	5[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	8[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	8[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	8[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	8[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	8[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	8[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	H	8[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	11[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	11[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	11[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	11[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	11[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	11[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	11[G]	1	-	0/0/11/13	0/1/1/1
1	HZP	I	13[A]	1	-	0/0/11/13	0/1/1/1
1	HZP	I	13[B]	1	-	0/0/11/13	0/1/1/1
1	HZP	I	13[C]	1	-	0/0/11/13	0/1/1/1
1	HZP	I	13[D]	1	-	0/0/11/13	0/1/1/1
1	HZP	I	13[E]	1	-	0/0/11/13	0/1/1/1
1	HZP	I	13[F]	1	-	0/0/11/13	0/1/1/1
1	HZP	I	13[G]	1	-	0/0/11/13	0/1/1/1
1	HZP	I	16[A]	1	-	0/0/11/13	0/1/1/1
1	HZP	I	16[B]	1	-	0/0/11/13	0/1/1/1
1	HZP	I	16[C]	1	-	0/0/11/13	0/1/1/1
1	HZP	I	16[D]	1	-	0/0/11/13	0/1/1/1
1	HZP	I	16[E]	1	-	0/0/11/13	0/1/1/1
1	HZP	I	16[F]	1	-	0/0/11/13	0/1/1/1
1	HZP	I	16[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	2[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	2[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	2[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	2[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	2[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	2[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	2[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	20[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	20[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	20[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	20[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	20[E]	1	-	0/0/11/13	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	HYP	I	20[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	20[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	23[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	23[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	23[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	23[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	23[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	23[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	23[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	26[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	26[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	26[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	26[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	26[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	26[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	26[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	29[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	29[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	29[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	29[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	29[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	29[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	29[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	5[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	5[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	5[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	5[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	5[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	5[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	5[G]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	8[A]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	8[B]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	8[C]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	8[D]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	8[E]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	8[F]	1	-	0/0/11/13	0/1/1/1
1	HYP	I	8[G]	1	-	0/0/11/13	0/1/1/1

All (688) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	20[D]	HYP	O-C	21.88	1.26	1.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	G	20[B]	HYP	O-C	21.26	1.26	1.11
1	G	13[D]	HZP	O-C	21.08	1.26	1.11
1	C	16[G]	HZP	O-C	21.07	1.25	1.11
1	D	16[D]	HZP	O-C	21.04	1.25	1.11
1	G	20[A]	HYP	O-C	20.99	1.25	1.11
1	G	23[D]	HYP	O-C	20.90	1.25	1.11
1	D	13[B]	HZP	O-C	20.90	1.25	1.11
1	D	13[A]	HZP	O-C	20.90	1.25	1.11
1	D	13[C]	HZP	O-C	20.90	1.25	1.11
1	D	20[C]	HYP	O-C	20.90	1.25	1.11
1	G	20[C]	HYP	O-C	20.86	1.25	1.11
1	H	29[D]	HYP	O-C	20.86	1.25	1.11
1	D	20[A]	HYP	O-C	20.83	1.25	1.11
1	A	29[B]	HYP	O-C	20.80	1.25	1.11
1	A	29[C]	HYP	O-C	20.80	1.25	1.11
1	A	29[A]	HYP	O-C	20.80	1.25	1.11
1	H	8[D]	HYP	O-C	20.75	1.25	1.11
1	H	8[C]	HYP	O-C	20.75	1.25	1.11
1	H	11[B]	HYP	O-C	20.68	1.25	1.11
1	H	11[A]	HYP	O-C	20.68	1.25	1.11
1	H	11[E]	HYP	O-C	20.68	1.25	1.11
1	I	11[A]	HYP	O-C	20.68	1.25	1.11
1	B	13[C]	HZP	O-C	20.67	1.25	1.11
1	H	29[C]	HYP	O-C	20.66	1.25	1.11
1	H	2[G]	HYP	O-C	20.66	1.25	1.11
1	H	2[F]	HYP	O-C	20.66	1.25	1.11
1	H	5[E]	HYP	O-C	20.66	1.25	1.11
1	H	5[A]	HYP	O-C	20.66	1.25	1.11
1	H	5[B]	HYP	O-C	20.65	1.25	1.11
1	D	23[D]	HYP	O-C	20.65	1.25	1.11
1	F	8[E]	HYP	O-C	20.65	1.25	1.11
1	A	16[G]	HZP	O-C	20.64	1.25	1.11
1	H	26[B]	HYP	O-C	20.64	1.25	1.11
1	H	26[A]	HYP	O-C	20.64	1.25	1.11
1	H	26[E]	HYP	O-C	20.64	1.25	1.11
1	I	20[G]	HYP	O-C	20.63	1.25	1.11
1	E	26[B]	HYP	O-C	20.63	1.25	1.11
1	D	13[E]	HZP	O-C	20.58	1.25	1.11
1	D	20[B]	HYP	O-C	20.58	1.25	1.11
1	B	23[E]	HYP	O-C	20.57	1.25	1.11
1	C	20[E]	HYP	O-C	20.56	1.25	1.11
1	I	23[A]	HYP	O-C	20.55	1.25	1.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	H	23[F]	HYP	O-C	20.54	1.25	1.11
1	D	11[D]	HYP	O-C	20.54	1.25	1.11
1	B	11[C]	HYP	O-C	20.54	1.25	1.11
1	D	5[F]	HYP	O-C	20.54	1.25	1.11
1	D	5[E]	HYP	O-C	20.54	1.25	1.11
1	E	11[G]	HYP	O-C	20.54	1.25	1.11
1	E	11[F]	HYP	O-C	20.54	1.25	1.11
1	H	23[G]	HYP	O-C	20.53	1.25	1.11
1	D	16[F]	HZP	O-C	20.50	1.25	1.11
1	D	16[G]	HZP	O-C	20.50	1.25	1.11
1	D	16[E]	HZP	O-C	20.50	1.25	1.11
1	I	26[D]	HYP	O-C	20.48	1.25	1.11
1	H	29[F]	HYP	O-C	20.47	1.25	1.11
1	C	23[D]	HYP	O-C	20.47	1.25	1.11
1	A	16[F]	HZP	O-C	20.47	1.25	1.11
1	B	11[D]	HYP	O-C	20.45	1.25	1.11
1	F	11[B]	HYP	O-C	20.43	1.25	1.11
1	F	11[D]	HYP	O-C	20.43	1.25	1.11
1	F	11[C]	HYP	O-C	20.43	1.25	1.11
1	E	26[A]	HYP	O-C	20.43	1.25	1.11
1	I	26[C]	HYP	O-C	20.43	1.25	1.11
1	D	8[B]	HYP	O-C	20.43	1.25	1.11
1	C	11[C]	HYP	O-C	20.42	1.25	1.11
1	C	11[B]	HYP	O-C	20.42	1.25	1.11
1	D	8[A]	HYP	O-C	20.42	1.25	1.11
1	C	11[D]	HYP	O-C	20.42	1.25	1.11
1	F	20[F]	HYP	O-C	20.42	1.25	1.11
1	D	8[C]	HYP	O-C	20.42	1.25	1.11
1	H	29[G]	HYP	O-C	20.42	1.25	1.11
1	C	23[B]	HYP	O-C	20.41	1.25	1.11
1	I	11[E]	HYP	O-C	20.41	1.25	1.11
1	D	5[G]	HYP	O-C	20.40	1.25	1.11
1	D	29[E]	HYP	O-C	20.38	1.25	1.11
1	B	20[F]	HYP	O-C	20.38	1.25	1.11
1	B	13[D]	HZP	O-C	20.38	1.25	1.11
1	B	20[G]	HYP	O-C	20.38	1.25	1.11
1	I	26[B]	HYP	O-C	20.38	1.25	1.11
1	F	8[A]	HYP	O-C	20.38	1.25	1.11
1	C	20[A]	HYP	O-C	20.37	1.25	1.11
1	D	11[B]	HYP	O-C	20.35	1.25	1.11
1	E	29[C]	HYP	O-C	20.35	1.25	1.11
1	E	29[D]	HYP	O-C	20.35	1.25	1.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	16[C]	HZP	O-C	20.34	1.25	1.11
1	G	26[G]	HYP	O-C	20.32	1.25	1.11
1	F	5[G]	HYP	O-C	20.31	1.25	1.11
1	B	2[E]	HYP	O-C	20.31	1.25	1.11
1	B	23[A]	HYP	O-C	20.30	1.25	1.11
1	D	13[F]	HZP	O-C	20.30	1.25	1.11
1	D	29[F]	HYP	O-C	20.30	1.25	1.11
1	E	23[G]	HYP	O-C	20.30	1.25	1.11
1	E	23[F]	HYP	O-C	20.30	1.25	1.11
1	D	11[F]	HYP	O-C	20.29	1.25	1.11
1	D	11[E]	HYP	O-C	20.29	1.25	1.11
1	D	11[G]	HYP	O-C	20.29	1.25	1.11
1	D	29[G]	HYP	O-C	20.26	1.25	1.11
1	A	16[E]	HZP	O-C	20.25	1.25	1.11
1	H	8[F]	HYP	O-C	20.25	1.25	1.11
1	G	26[E]	HYP	O-C	20.24	1.25	1.11
1	E	26[E]	HYP	O-C	20.24	1.25	1.11
1	I	23[F]	HYP	O-C	20.24	1.25	1.11
1	I	23[E]	HYP	O-C	20.24	1.25	1.11
1	D	11[A]	HYP	O-C	20.23	1.25	1.11
1	D	8[F]	HYP	O-C	20.21	1.25	1.11
1	D	8[G]	HYP	O-C	20.21	1.25	1.11
1	D	8[E]	HYP	O-C	20.21	1.25	1.11
1	F	13[D]	HZP	O-C	20.20	1.25	1.11
1	C	8[F]	HYP	O-C	20.19	1.25	1.11
1	H	8[B]	HYP	O-C	20.19	1.25	1.11
1	F	13[B]	HZP	O-C	20.18	1.25	1.11
1	F	13[C]	HZP	O-C	20.18	1.25	1.11
1	C	5[G]	HYP	O-C	20.18	1.25	1.11
1	C	20[F]	HYP	O-C	20.17	1.25	1.11
1	B	26[C]	HYP	O-C	20.17	1.25	1.11
1	C	23[C]	HYP	O-C	20.16	1.25	1.11
1	H	8[G]	HYP	O-C	20.15	1.25	1.11
1	F	20[A]	HYP	O-C	20.15	1.25	1.11
1	B	8[B]	HYP	O-C	20.14	1.25	1.11
1	I	11[F]	HYP	O-C	20.14	1.25	1.11
1	C	13[E]	HZP	O-C	20.12	1.25	1.11
1	B	29[D]	HYP	O-C	20.11	1.25	1.11
1	C	16[B]	HZP	O-C	20.11	1.25	1.11
1	B	29[C]	HYP	O-C	20.11	1.25	1.11
1	F	8[F]	HYP	O-C	20.10	1.25	1.11
1	H	26[F]	HYP	O-C	20.09	1.25	1.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	H	8[E]	HYP	O-C	20.06	1.25	1.11
1	H	8[A]	HYP	O-C	20.06	1.25	1.11
1	F	23[B]	HYP	O-C	20.05	1.25	1.11
1	F	23[C]	HYP	O-C	20.05	1.25	1.11
1	C	8[A]	HYP	O-C	20.05	1.25	1.11
1	B	8[A]	HYP	O-C	20.05	1.25	1.11
1	B	26[E]	HYP	O-C	20.05	1.25	1.11
1	F	23[F]	HYP	O-C	20.05	1.25	1.11
1	D	11[C]	HYP	O-C	20.04	1.25	1.11
1	B	2[A]	HYP	O-C	20.04	1.25	1.11
1	B	23[B]	HYP	O-C	20.03	1.25	1.11
1	A	23[G]	HYP	O-C	20.03	1.25	1.11
1	A	5[C]	HYP	O-C	20.03	1.25	1.11
1	G	11[C]	HYP	O-C	20.02	1.25	1.11
1	B	5[G]	HYP	O-C	20.02	1.25	1.11
1	H	5[G]	HYP	O-C	20.02	1.25	1.11
1	H	5[F]	HYP	O-C	20.02	1.25	1.11
1	D	13[G]	HZP	O-C	20.01	1.25	1.11
1	C	13[F]	HZP	O-C	20.01	1.25	1.11
1	A	26[G]	HYP	O-C	20.01	1.25	1.11
1	I	8[G]	HYP	O-C	19.99	1.25	1.11
1	B	26[A]	HYP	O-C	19.98	1.25	1.11
1	H	13[C]	HZP	O-C	19.98	1.25	1.11
1	H	13[D]	HZP	O-C	19.98	1.25	1.11
1	C	13[A]	HZP	O-C	19.98	1.25	1.11
1	I	29[G]	HYP	O-C	19.96	1.25	1.11
1	G	11[B]	HYP	O-C	19.96	1.25	1.11
1	B	8[E]	HYP	O-C	19.96	1.25	1.11
1	B	23[G]	HYP	O-C	19.95	1.25	1.11
1	A	29[D]	HYP	O-C	19.95	1.25	1.11
1	D	13[D]	HZP	O-C	19.94	1.25	1.11
1	F	29[G]	HYP	O-C	19.94	1.25	1.11
1	G	29[A]	HYP	O-C	19.93	1.25	1.11
1	B	5[F]	HYP	O-C	19.93	1.25	1.11
1	H	13[E]	HZP	O-C	19.93	1.25	1.11
1	H	13[A]	HZP	O-C	19.93	1.25	1.11
1	C	8[E]	HYP	O-C	19.92	1.25	1.11
1	F	23[D]	HYP	O-C	19.92	1.25	1.11
1	A	26[C]	HYP	O-C	19.92	1.25	1.11
1	G	5[G]	HYP	O-C	19.92	1.25	1.11
1	H	13[B]	HZP	O-C	19.92	1.25	1.11
1	H	26[G]	HYP	O-C	19.92	1.25	1.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	23[A]	HYP	O-C	19.92	1.25	1.11
1	F	29[C]	HYP	O-C	19.91	1.25	1.11
1	B	26[D]	HYP	O-C	19.90	1.25	1.11
1	B	2[B]	HYP	O-C	19.90	1.25	1.11
1	H	16[C]	HZP	O-C	19.90	1.25	1.11
1	G	11[A]	HYP	O-C	19.90	1.25	1.11
1	H	16[D]	HZP	O-C	19.89	1.25	1.11
1	C	16[D]	HZP	O-C	19.89	1.25	1.11
1	D	16[A]	HZP	O-C	19.88	1.25	1.11
1	D	16[B]	HZP	O-C	19.88	1.25	1.11
1	D	16[C]	HZP	O-C	19.88	1.25	1.11
1	I	29[F]	HYP	O-C	19.88	1.25	1.11
1	F	20[E]	HYP	O-C	19.88	1.25	1.11
1	G	11[D]	HYP	O-C	19.87	1.25	1.11
1	H	20[D]	HYP	O-C	19.87	1.25	1.11
1	H	20[C]	HYP	O-C	19.87	1.25	1.11
1	E	2[B]	HYP	O-C	19.86	1.25	1.11
1	G	29[B]	HYP	O-C	19.84	1.25	1.11
1	G	29[C]	HYP	O-C	19.84	1.25	1.11
1	I	29[A]	HYP	O-C	19.84	1.25	1.11
1	A	26[B]	HYP	O-C	19.84	1.25	1.11
1	H	2[E]	HYP	O-C	19.84	1.25	1.11
1	H	2[A]	HYP	O-C	19.84	1.25	1.11
1	H	2[B]	HYP	O-C	19.84	1.25	1.11
1	I	2[F]	HYP	O-C	19.83	1.25	1.11
1	I	2[A]	HYP	O-C	19.83	1.25	1.11
1	I	2[E]	HYP	O-C	19.83	1.25	1.11
1	A	26[A]	HYP	O-C	19.82	1.25	1.11
1	B	23[F]	HYP	O-C	19.82	1.25	1.11
1	H	16[F]	HZP	O-C	19.81	1.25	1.11
1	I	16[F]	HZP	O-C	19.81	1.25	1.11
1	A	26[E]	HYP	O-C	19.80	1.25	1.11
1	B	26[B]	HYP	O-C	19.80	1.25	1.11
1	A	23[B]	HYP	O-C	19.79	1.25	1.11
1	A	26[F]	HYP	O-C	19.79	1.25	1.11
1	B	5[C]	HYP	O-C	19.79	1.25	1.11
1	F	23[E]	HYP	O-C	19.79	1.25	1.11
1	G	8[F]	HYP	O-C	19.78	1.25	1.11
1	F	23[G]	HYP	O-C	19.78	1.25	1.11
1	B	11[A]	HYP	O-C	19.77	1.25	1.11
1	B	11[B]	HYP	O-C	19.77	1.25	1.11
1	B	11[E]	HYP	O-C	19.77	1.25	1.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	2[C]	HYP	O-C	19.77	1.25	1.11
1	E	2[D]	HYP	O-C	19.77	1.25	1.11
1	I	13[G]	HZP	O-C	19.76	1.25	1.11
1	G	8[C]	HYP	O-C	19.76	1.25	1.11
1	A	2[F]	HYP	O-C	19.76	1.25	1.11
1	A	23[E]	HYP	O-C	19.76	1.25	1.11
1	A	5[A]	HYP	O-C	19.75	1.25	1.11
1	F	20[G]	HYP	O-C	19.74	1.25	1.11
1	F	29[B]	HYP	O-C	19.74	1.25	1.11
1	A	23[A]	HYP	O-C	19.74	1.25	1.11
1	A	23[C]	HYP	O-C	19.74	1.25	1.11
1	G	8[B]	HYP	O-C	19.74	1.25	1.11
1	F	26[B]	HYP	O-C	19.73	1.25	1.11
1	F	26[D]	HYP	O-C	19.73	1.25	1.11
1	F	26[C]	HYP	O-C	19.73	1.25	1.11
1	C	29[C]	HYP	O-C	19.72	1.25	1.11
1	E	29[F]	HYP	O-C	19.72	1.25	1.11
1	I	29[E]	HYP	O-C	19.72	1.25	1.11
1	G	8[E]	HYP	O-C	19.72	1.25	1.11
1	H	11[F]	HYP	O-C	19.70	1.25	1.11
1	C	26[A]	HYP	O-C	19.69	1.25	1.11
1	C	26[F]	HYP	O-C	19.69	1.25	1.11
1	D	2[D]	HYP	O-C	19.69	1.25	1.11
1	C	26[E]	HYP	O-C	19.69	1.25	1.11
1	A	2[E]	HYP	O-C	19.69	1.25	1.11
1	G	8[A]	HYP	O-C	19.66	1.25	1.11
1	G	8[G]	HYP	O-C	19.66	1.25	1.11
1	G	5[E]	HYP	O-C	19.64	1.24	1.11
1	F	29[E]	HYP	O-C	19.63	1.24	1.11
1	C	29[D]	HYP	O-C	19.62	1.24	1.11
1	A	20[E]	HYP	O-C	19.61	1.24	1.11
1	I	16[A]	HZP	O-C	19.61	1.24	1.11
1	I	16[E]	HZP	O-C	19.61	1.24	1.11
1	C	29[E]	HYP	O-C	19.61	1.24	1.11
1	C	23[G]	HYP	O-C	19.60	1.24	1.11
1	F	26[F]	HYP	O-C	19.59	1.24	1.11
1	A	26[D]	HYP	O-C	19.59	1.24	1.11
1	A	23[F]	HYP	O-C	19.58	1.24	1.11
1	F	29[D]	HYP	O-C	19.58	1.24	1.11
1	E	29[G]	HYP	O-C	19.54	1.24	1.11
1	C	29[B]	HYP	O-C	19.54	1.24	1.11
1	C	26[G]	HYP	O-C	19.54	1.24	1.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	I	5[B]	HYP	O-C	19.54	1.24	1.11
1	H	16[G]	HZP	O-C	19.53	1.24	1.11
1	B	13[G]	HZP	O-C	19.52	1.24	1.11
1	B	13[F]	HZP	O-C	19.52	1.24	1.11
1	F	26[A]	HYP	O-C	19.52	1.24	1.11
1	A	2[G]	HYP	O-C	19.52	1.24	1.11
1	B	16[F]	HZP	O-C	19.51	1.24	1.11
1	B	8[G]	HYP	O-C	19.50	1.24	1.11
1	B	8[F]	HYP	O-C	19.50	1.24	1.11
1	A	5[B]	HYP	O-C	19.48	1.24	1.11
1	C	29[F]	HYP	O-C	19.48	1.24	1.11
1	C	29[A]	HYP	O-C	19.47	1.24	1.11
1	E	5[E]	HYP	O-C	19.47	1.24	1.11
1	F	29[A]	HYP	O-C	19.47	1.24	1.11
1	G	26[F]	HYP	O-C	19.46	1.24	1.11
1	F	26[E]	HYP	O-C	19.46	1.24	1.11
1	H	11[D]	HYP	O-C	19.45	1.24	1.11
1	H	11[C]	HYP	O-C	19.45	1.24	1.11
1	C	16[E]	HZP	O-C	19.45	1.24	1.11
1	C	16[A]	HZP	O-C	19.45	1.24	1.11
1	C	16[F]	HZP	O-C	19.45	1.24	1.11
1	E	29[A]	HYP	O-C	19.44	1.24	1.11
1	E	29[B]	HYP	O-C	19.44	1.24	1.11
1	H	11[G]	HYP	O-C	19.43	1.24	1.11
1	E	2[E]	HYP	O-C	19.42	1.24	1.11
1	E	20[E]	HYP	O-C	19.41	1.24	1.11
1	D	23[G]	HYP	O-C	19.41	1.24	1.11
1	I	16[G]	HZP	O-C	19.41	1.24	1.11
1	I	5[C]	HYP	O-C	19.40	1.24	1.11
1	B	16[B]	HZP	O-C	19.40	1.24	1.11
1	A	20[F]	HYP	O-C	19.39	1.24	1.11
1	D	2[G]	HYP	O-C	19.39	1.24	1.11
1	D	2[E]	HYP	O-C	19.39	1.24	1.11
1	D	2[F]	HYP	O-C	19.39	1.24	1.11
1	E	2[A]	HYP	O-C	19.38	1.24	1.11
1	G	5[F]	HYP	O-C	19.37	1.24	1.11
1	D	5[A]	HYP	O-C	19.36	1.24	1.11
1	D	5[B]	HYP	O-C	19.36	1.24	1.11
1	E	26[D]	HYP	O-C	19.36	1.24	1.11
1	C	11[E]	HYP	O-C	19.36	1.24	1.11
1	B	16[G]	HZP	O-C	19.34	1.24	1.11
1	A	20[G]	HYP	O-C	19.34	1.24	1.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	8[C]	HYP	O-C	19.33	1.24	1.11
1	A	8[A]	HYP	O-C	19.33	1.24	1.11
1	E	29[E]	HYP	O-C	19.32	1.24	1.11
1	A	20[C]	HYP	O-C	19.31	1.24	1.11
1	G	11[G]	HYP	O-C	19.31	1.24	1.11
1	G	11[E]	HYP	O-C	19.31	1.24	1.11
1	G	11[F]	HYP	O-C	19.31	1.24	1.11
1	E	23[C]	HYP	O-C	19.31	1.24	1.11
1	E	23[D]	HYP	O-C	19.31	1.24	1.11
1	G	23[F]	HYP	O-C	19.30	1.24	1.11
1	E	20[F]	HYP	O-C	19.29	1.24	1.11
1	E	20[G]	HYP	O-C	19.29	1.24	1.11
1	I	5[D]	HYP	O-C	19.27	1.24	1.11
1	G	29[D]	HYP	O-C	19.27	1.24	1.11
1	B	16[A]	HZP	O-C	19.26	1.24	1.11
1	G	26[B]	HYP	O-C	19.26	1.24	1.11
1	D	8[D]	HYP	O-C	19.26	1.24	1.11
1	G	16[F]	HZP	O-C	19.25	1.24	1.11
1	G	16[E]	HZP	O-C	19.25	1.24	1.11
1	G	16[G]	HZP	O-C	19.25	1.24	1.11
1	E	20[A]	HYP	O-C	19.24	1.24	1.11
1	E	5[A]	HYP	O-C	19.22	1.24	1.11
1	C	13[C]	HZP	O-C	19.22	1.24	1.11
1	B	16[E]	HZP	O-C	19.22	1.24	1.11
1	D	23[F]	HYP	O-C	19.21	1.24	1.11
1	G	23[E]	HYP	O-C	19.21	1.24	1.11
1	D	26[B]	HYP	O-C	19.21	1.24	1.11
1	D	26[A]	HYP	O-C	19.21	1.24	1.11
1	G	26[D]	HYP	O-C	19.20	1.24	1.11
1	A	23[D]	HYP	O-C	19.20	1.24	1.11
1	F	8[B]	HYP	O-C	19.20	1.24	1.11
1	F	8[D]	HYP	O-C	19.20	1.24	1.11
1	F	8[C]	HYP	O-C	19.20	1.24	1.11
1	A	11[D]	HYP	O-C	19.20	1.24	1.11
1	H	29[E]	HYP	O-C	19.19	1.24	1.11
1	H	29[A]	HYP	O-C	19.19	1.24	1.11
1	F	29[F]	HYP	O-C	19.19	1.24	1.11
1	B	8[D]	HYP	O-C	19.19	1.24	1.11
1	E	23[E]	HYP	O-C	19.19	1.24	1.11
1	E	26[C]	HYP	O-C	19.19	1.24	1.11
1	E	23[A]	HYP	O-C	19.19	1.24	1.11
1	A	8[D]	HYP	O-C	19.18	1.24	1.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	16[D]	HZP	O-C	19.18	1.24	1.11
1	E	23[B]	HYP	O-C	19.17	1.24	1.11
1	G	20[G]	HYP	O-C	19.17	1.24	1.11
1	E	20[C]	HYP	O-C	19.17	1.24	1.11
1	G	20[E]	HYP	O-C	19.17	1.24	1.11
1	G	20[F]	HYP	O-C	19.17	1.24	1.11
1	D	5[C]	HYP	O-C	19.17	1.24	1.11
1	G	26[A]	HYP	O-C	19.17	1.24	1.11
1	C	11[A]	HYP	O-C	19.17	1.24	1.11
1	C	11[F]	HYP	O-C	19.17	1.24	1.11
1	F	11[F]	HYP	O-C	19.16	1.24	1.11
1	F	11[E]	HYP	O-C	19.16	1.24	1.11
1	F	11[A]	HYP	O-C	19.16	1.24	1.11
1	A	20[B]	HYP	O-C	19.15	1.24	1.11
1	A	5[E]	HYP	O-C	19.15	1.24	1.11
1	D	29[D]	HYP	O-C	19.15	1.24	1.11
1	E	8[C]	HYP	O-C	19.14	1.24	1.11
1	E	20[B]	HYP	O-C	19.14	1.24	1.11
1	A	20[A]	HYP	O-C	19.12	1.24	1.11
1	B	5[B]	HYP	O-C	19.12	1.24	1.11
1	G	23[G]	HYP	O-C	19.12	1.24	1.11
1	B	13[A]	HZP	O-C	19.12	1.24	1.11
1	B	13[E]	HZP	O-C	19.12	1.24	1.11
1	G	23[B]	HYP	O-C	19.12	1.24	1.11
1	A	5[G]	HYP	O-C	19.10	1.24	1.11
1	C	13[G]	HZP	O-C	19.09	1.24	1.11
1	A	8[B]	HYP	O-C	19.08	1.24	1.11
1	G	26[C]	HYP	O-C	19.08	1.24	1.11
1	F	2[C]	HYP	O-C	19.07	1.24	1.11
1	F	2[D]	HYP	O-C	19.07	1.24	1.11
1	F	2[B]	HYP	O-C	19.07	1.24	1.11
1	H	29[B]	HYP	O-C	19.05	1.24	1.11
1	G	23[A]	HYP	O-C	19.05	1.24	1.11
1	F	16[G]	HZP	O-C	19.05	1.24	1.11
1	B	2[F]	HYP	O-C	19.04	1.24	1.11
1	D	23[E]	HYP	O-C	19.03	1.24	1.11
1	G	2[A]	HYP	O-C	19.03	1.24	1.11
1	D	26[C]	HYP	O-C	19.01	1.24	1.11
1	G	2[C]	HYP	O-C	19.00	1.24	1.11
1	G	5[D]	HYP	O-C	19.00	1.24	1.11
1	B	8[C]	HYP	O-C	19.00	1.24	1.11
1	C	8[G]	HYP	O-C	19.00	1.24	1.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	8[D]	HYP	O-C	19.00	1.24	1.11
1	G	2[D]	HYP	O-C	18.99	1.24	1.11
1	G	5[B]	HYP	O-C	18.98	1.24	1.11
1	G	5[A]	HYP	O-C	18.98	1.24	1.11
1	G	5[C]	HYP	O-C	18.98	1.24	1.11
1	I	20[D]	HYP	O-C	18.97	1.24	1.11
1	I	20[C]	HYP	O-C	18.97	1.24	1.11
1	I	20[B]	HYP	O-C	18.97	1.24	1.11
1	E	5[B]	HYP	O-C	18.97	1.24	1.11
1	C	5[F]	HYP	O-C	18.96	1.24	1.11
1	C	5[A]	HYP	O-C	18.96	1.24	1.11
1	C	5[E]	HYP	O-C	18.96	1.24	1.11
1	C	2[G]	HYP	O-C	18.96	1.24	1.11
1	A	5[F]	HYP	O-C	18.95	1.24	1.11
1	C	8[C]	HYP	O-C	18.95	1.24	1.11
1	C	8[B]	HYP	O-C	18.95	1.24	1.11
1	C	8[D]	HYP	O-C	18.95	1.24	1.11
1	C	13[B]	HYP	O-C	18.95	1.24	1.11
1	E	20[D]	HYP	O-C	18.94	1.24	1.11
1	B	5[E]	HYP	O-C	18.93	1.24	1.11
1	B	5[A]	HYP	O-C	18.93	1.24	1.11
1	F	5[E]	HYP	O-C	18.93	1.24	1.11
1	G	23[C]	HYP	O-C	18.93	1.24	1.11
1	B	13[B]	HYP	O-C	18.92	1.24	1.11
1	F	16[B]	HYP	O-C	18.90	1.24	1.11
1	F	13[F]	HYP	O-C	18.89	1.24	1.11
1	D	29[B]	HYP	O-C	18.86	1.24	1.11
1	H	23[A]	HYP	O-C	18.85	1.24	1.11
1	E	2[F]	HYP	O-C	18.85	1.24	1.11
1	F	2[E]	HYP	O-C	18.84	1.24	1.11
1	I	11[G]	HYP	O-C	18.84	1.24	1.11
1	H	13[G]	HYP	O-C	18.83	1.24	1.11
1	I	26[G]	HYP	O-C	18.82	1.24	1.11
1	H	23[B]	HYP	O-C	18.82	1.24	1.11
1	F	8[G]	HYP	O-C	18.81	1.24	1.11
1	I	13[D]	HYP	O-C	18.80	1.24	1.11
1	I	13[C]	HYP	O-C	18.80	1.24	1.11
1	I	13[B]	HYP	O-C	18.80	1.24	1.11
1	C	11[G]	HYP	O-C	18.79	1.24	1.11
1	C	5[D]	HYP	O-C	18.79	1.24	1.11
1	B	2[G]	HYP	O-C	18.79	1.24	1.11
1	A	29[G]	HYP	O-C	18.78	1.24	1.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	29[E]	HYP	O-C	18.78	1.24	1.11
1	H	20[F]	HYP	O-C	18.75	1.24	1.11
1	H	20[G]	HYP	O-C	18.75	1.24	1.11
1	F	16[C]	HZP	O-C	18.74	1.24	1.11
1	A	29[F]	HYP	O-C	18.73	1.24	1.11
1	H	23[E]	HYP	O-C	18.72	1.24	1.11
1	D	29[A]	HYP	O-C	18.71	1.24	1.11
1	D	29[C]	HYP	O-C	18.71	1.24	1.11
1	D	5[D]	HYP	O-C	18.71	1.24	1.11
1	G	2[B]	HYP	O-C	18.69	1.24	1.11
1	C	13[D]	HZP	O-C	18.69	1.24	1.11
1	D	26[F]	HYP	O-C	18.69	1.24	1.11
1	D	26[G]	HYP	O-C	18.69	1.24	1.11
1	D	26[E]	HYP	O-C	18.69	1.24	1.11
1	G	2[E]	HYP	O-C	18.67	1.24	1.11
1	H	16[B]	HZP	O-C	18.65	1.24	1.11
1	F	5[D]	HYP	O-C	18.64	1.24	1.11
1	H	26[C]	HYP	O-C	18.64	1.24	1.11
1	F	2[G]	HYP	O-C	18.63	1.24	1.11
1	H	13[F]	HZP	O-C	18.62	1.24	1.11
1	F	5[A]	HYP	O-C	18.61	1.24	1.11
1	F	13[A]	HZP	O-C	18.61	1.24	1.11
1	B	11[G]	HYP	O-C	18.61	1.24	1.11
1	E	2[G]	HYP	O-C	18.60	1.24	1.11
1	H	26[D]	HYP	O-C	18.60	1.24	1.11
1	F	5[B]	HYP	O-C	18.60	1.24	1.11
1	F	2[A]	HYP	O-C	18.59	1.24	1.11
1	G	2[G]	HYP	O-C	18.58	1.24	1.11
1	C	2[E]	HYP	O-C	18.55	1.24	1.11
1	C	2[A]	HYP	O-C	18.55	1.24	1.11
1	C	2[F]	HYP	O-C	18.55	1.24	1.11
1	H	2[D]	HYP	O-C	18.55	1.24	1.11
1	H	2[C]	HYP	O-C	18.55	1.24	1.11
1	C	2[B]	HYP	O-C	18.53	1.24	1.11
1	F	5[F]	HYP	O-C	18.53	1.24	1.11
1	I	26[F]	HYP	O-C	18.52	1.24	1.11
1	I	26[E]	HYP	O-C	18.52	1.24	1.11
1	I	26[A]	HYP	O-C	18.52	1.24	1.11
1	A	16[A]	HZP	O-C	18.52	1.24	1.11
1	C	5[B]	HYP	O-C	18.50	1.24	1.11
1	I	16[D]	HZP	O-C	18.50	1.24	1.11
1	G	2[F]	HYP	O-C	18.50	1.24	1.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	8[G]	HYP	O-C	18.47	1.24	1.11
1	D	2[B]	HYP	O-C	18.46	1.24	1.11
1	D	2[A]	HYP	O-C	18.46	1.24	1.11
1	D	2[C]	HYP	O-C	18.46	1.24	1.11
1	G	13[C]	HZP	O-C	18.46	1.24	1.11
1	B	11[F]	HYP	O-C	18.46	1.24	1.11
1	C	29[G]	HYP	O-C	18.44	1.24	1.11
1	A	16[C]	HZP	O-C	18.43	1.24	1.11
1	A	11[C]	HYP	O-C	18.43	1.24	1.11
1	A	11[A]	HYP	O-C	18.43	1.24	1.11
1	H	16[E]	HZP	O-C	18.43	1.24	1.11
1	H	16[A]	HZP	O-C	18.43	1.24	1.11
1	E	8[E]	HYP	O-C	18.42	1.24	1.11
1	F	2[F]	HYP	O-C	18.41	1.24	1.11
1	C	2[C]	HYP	O-C	18.39	1.24	1.11
1	A	16[B]	HZP	O-C	18.35	1.24	1.11
1	G	29[F]	HYP	O-C	18.35	1.24	1.11
1	C	2[D]	HYP	O-C	18.35	1.24	1.11
1	G	29[G]	HYP	O-C	18.35	1.24	1.11
1	A	8[E]	HYP	O-C	18.34	1.24	1.11
1	A	8[F]	HYP	O-C	18.34	1.24	1.11
1	G	29[E]	HYP	O-C	18.34	1.24	1.11
1	F	13[E]	HZP	O-C	18.33	1.24	1.11
1	F	5[C]	HYP	O-C	18.33	1.24	1.11
1	G	20[D]	HYP	O-C	18.33	1.24	1.11
1	B	23[C]	HYP	O-C	18.32	1.24	1.11
1	A	11[B]	HYP	O-C	18.32	1.24	1.11
1	G	13[A]	HZP	O-C	18.31	1.24	1.11
1	A	13[B]	HZP	O-C	18.30	1.24	1.11
1	A	2[B]	HYP	O-C	18.26	1.24	1.11
1	C	5[C]	HYP	O-C	18.22	1.24	1.11
1	I	23[G]	HYP	O-C	18.22	1.24	1.11
1	E	16[B]	HZP	O-C	18.22	1.23	1.11
1	C	20[C]	HYP	O-C	18.21	1.23	1.11
1	A	13[E]	HZP	O-C	18.21	1.23	1.11
1	B	26[F]	HYP	O-C	18.20	1.23	1.11
1	A	2[A]	HYP	O-C	18.18	1.23	1.11
1	A	13[F]	HZP	O-C	18.18	1.23	1.11
1	G	13[B]	HZP	O-C	18.16	1.23	1.11
1	I	11[D]	HYP	O-C	18.15	1.23	1.11
1	B	2[D]	HYP	O-C	18.15	1.23	1.11
1	A	13[G]	HZP	O-C	18.14	1.23	1.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	23[D]	HYP	O-C	18.14	1.23	1.11
1	I	8[F]	HYP	O-C	18.13	1.23	1.11
1	I	5[G]	HYP	O-C	18.12	1.23	1.11
1	B	20[B]	HYP	O-C	18.11	1.23	1.11
1	B	20[E]	HYP	O-C	18.11	1.23	1.11
1	B	20[A]	HYP	O-C	18.11	1.23	1.11
1	I	2[D]	HYP	O-C	18.10	1.23	1.11
1	I	2[C]	HYP	O-C	18.10	1.23	1.11
1	I	16[C]	HZP	O-C	18.09	1.23	1.11
1	I	16[B]	HZP	O-C	18.09	1.23	1.11
1	E	16[E]	HZP	O-C	18.08	1.23	1.11
1	I	11[B]	HYP	O-C	18.08	1.23	1.11
1	A	13[C]	HZP	O-C	18.07	1.23	1.11
1	C	20[B]	HYP	O-C	18.06	1.23	1.11
1	G	16[D]	HZP	O-C	18.05	1.23	1.11
1	E	16[G]	HZP	O-C	18.03	1.23	1.11
1	I	20[F]	HYP	O-C	18.02	1.23	1.11
1	B	2[C]	HYP	O-C	18.01	1.23	1.11
1	B	26[G]	HYP	O-C	18.00	1.23	1.11
1	E	8[B]	HYP	O-C	17.99	1.23	1.11
1	E	8[A]	HYP	O-C	17.99	1.23	1.11
1	F	26[G]	HYP	O-C	17.99	1.23	1.11
1	A	2[C]	HYP	O-C	17.99	1.23	1.11
1	A	11[G]	HYP	O-C	17.98	1.23	1.11
1	I	11[C]	HYP	O-C	17.97	1.23	1.11
1	I	29[B]	HYP	O-C	17.95	1.23	1.11
1	I	29[D]	HYP	O-C	17.95	1.23	1.11
1	I	29[C]	HYP	O-C	17.95	1.23	1.11
1	E	16[A]	HZP	O-C	17.95	1.23	1.11
1	A	5[D]	HYP	O-C	17.94	1.23	1.11
1	I	8[E]	HYP	O-C	17.93	1.23	1.11
1	E	5[G]	HYP	O-C	17.93	1.23	1.11
1	E	5[F]	HYP	O-C	17.93	1.23	1.11
1	I	23[C]	HYP	O-C	17.92	1.23	1.11
1	I	23[B]	HYP	O-C	17.92	1.23	1.11
1	F	20[B]	HYP	O-C	17.92	1.23	1.11
1	F	20[C]	HYP	O-C	17.92	1.23	1.11
1	F	20[D]	HYP	O-C	17.92	1.23	1.11
1	D	20[D]	HYP	O-C	17.92	1.23	1.11
1	A	13[A]	HZP	O-C	17.91	1.23	1.11
1	C	20[D]	HYP	O-C	17.90	1.23	1.11
1	I	20[A]	HYP	O-C	17.89	1.23	1.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	11[G]	HYP	O-C	17.89	1.23	1.11
1	A	11[F]	HYP	O-C	17.89	1.23	1.11
1	I	20[E]	HYP	O-C	17.89	1.23	1.11
1	I	5[F]	HYP	O-C	17.88	1.23	1.11
1	I	5[A]	HYP	O-C	17.88	1.23	1.11
1	I	5[E]	HYP	O-C	17.88	1.23	1.11
1	I	8[B]	HYP	O-C	17.88	1.23	1.11
1	I	2[G]	HYP	O-C	17.86	1.23	1.11
1	I	8[C]	HYP	O-C	17.86	1.23	1.11
1	E	11[D]	HYP	O-C	17.86	1.23	1.11
1	E	11[C]	HYP	O-C	17.86	1.23	1.11
1	I	13[E]	HZP	O-C	17.85	1.23	1.11
1	I	13[A]	HZP	O-C	17.85	1.23	1.11
1	I	13[F]	HZP	O-C	17.85	1.23	1.11
1	I	8[D]	HYP	O-C	17.84	1.23	1.11
1	E	16[D]	HZP	O-C	17.84	1.23	1.11
1	E	16[C]	HZP	O-C	17.84	1.23	1.11
1	I	8[A]	HYP	O-C	17.83	1.23	1.11
1	A	16[D]	HZP	O-C	17.82	1.23	1.11
1	E	13[B]	HZP	O-C	17.82	1.23	1.11
1	E	13[A]	HZP	O-C	17.82	1.23	1.11
1	E	13[E]	HZP	O-C	17.82	1.23	1.11
1	E	16[F]	HZP	O-C	17.81	1.23	1.11
1	A	11[E]	HYP	O-C	17.79	1.23	1.11
1	E	13[C]	HZP	O-C	17.76	1.23	1.11
1	E	13[D]	HZP	O-C	17.76	1.23	1.11
1	G	13[E]	HZP	O-C	17.76	1.23	1.11
1	G	13[G]	HZP	O-C	17.76	1.23	1.11
1	G	13[F]	HZP	O-C	17.76	1.23	1.11
1	E	8[F]	HYP	O-C	17.70	1.23	1.11
1	E	8[G]	HYP	O-C	17.70	1.23	1.11
1	I	23[D]	HYP	O-C	17.69	1.23	1.11
1	E	13[F]	HZP	O-C	17.69	1.23	1.11
1	E	13[G]	HZP	O-C	17.61	1.23	1.11
1	D	20[G]	HYP	O-C	17.57	1.23	1.11
1	I	2[B]	HYP	O-C	17.55	1.23	1.11
1	D	20[F]	HYP	O-C	17.55	1.23	1.11
1	C	26[B]	HYP	O-C	17.54	1.23	1.11
1	C	26[D]	HYP	O-C	17.54	1.23	1.11
1	C	26[C]	HYP	O-C	17.54	1.23	1.11
1	F	13[G]	HZP	O-C	17.48	1.23	1.11
1	C	23[A]	HYP	O-C	17.47	1.23	1.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	23[E]	HYP	O-C	17.45	1.23	1.11
1	C	23[F]	HYP	O-C	17.45	1.23	1.11
1	D	23[B]	HYP	O-C	17.44	1.23	1.11
1	D	23[C]	HYP	O-C	17.44	1.23	1.11
1	D	23[A]	HYP	O-C	17.42	1.23	1.11
1	D	26[D]	HYP	O-C	17.41	1.23	1.11
1	D	20[E]	HYP	O-C	17.40	1.23	1.11
1	B	20[C]	HYP	O-C	17.37	1.23	1.11
1	B	20[D]	HYP	O-C	17.37	1.23	1.11
1	C	20[G]	HYP	O-C	17.35	1.23	1.11
1	B	29[B]	HYP	O-C	17.32	1.23	1.11
1	B	29[E]	HYP	O-C	17.32	1.23	1.11
1	B	29[A]	HYP	O-C	17.32	1.23	1.11
1	F	16[A]	HZP	O-C	17.30	1.23	1.11
1	F	16[F]	HZP	O-C	17.30	1.23	1.11
1	F	16[E]	HZP	O-C	17.30	1.23	1.11
1	H	23[C]	HYP	O-C	17.27	1.23	1.11
1	H	23[D]	HYP	O-C	17.26	1.23	1.11
1	H	20[B]	HYP	O-C	17.18	1.23	1.11
1	H	20[E]	HYP	O-C	17.15	1.23	1.11
1	H	20[A]	HYP	O-C	17.15	1.23	1.11
1	E	11[B]	HYP	O-C	17.11	1.23	1.11
1	E	11[A]	HYP	O-C	17.11	1.23	1.11
1	E	11[E]	HYP	O-C	17.10	1.23	1.11
1	G	16[A]	HZP	O-C	16.97	1.23	1.11
1	G	16[B]	HZP	O-C	16.91	1.23	1.11
1	G	16[C]	HZP	O-C	16.84	1.23	1.11
1	A	13[D]	HZP	O-C	16.78	1.22	1.11
1	E	26[F]	HYP	O-C	16.41	1.22	1.11
1	E	26[G]	HYP	O-C	16.41	1.22	1.11
1	B	29[G]	HYP	O-C	15.79	1.22	1.11
1	B	29[F]	HYP	O-C	15.79	1.22	1.11
1	E	5[D]	HYP	O-C	15.77	1.22	1.11
1	E	5[C]	HYP	O-C	15.77	1.22	1.11
1	B	5[D]	HYP	O-C	14.45	1.21	1.11
1	H	5[D]	HYP	O-C	14.02	1.21	1.11
1	B	16[D]	HZP	O-C	13.90	1.20	1.11
1	B	16[C]	HZP	O-C	13.90	1.20	1.11
1	H	5[C]	HYP	O-C	13.85	1.20	1.11
1	G	8[D]	HYP	O-C	10.70	1.18	1.11
1	A	2[D]	HYP	O-C	8.30	1.17	1.11
1	B	16[D]	HZP	CA-C	3.01	1.54	1.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	16[C]	HZP	CA-C	3.01	1.54	1.48
1	A	13[F]	HZP	CA-C	2.67	1.53	1.48
1	A	13[E]	HZP	CA-C	2.65	1.53	1.48
1	A	13[G]	HZP	CA-C	2.65	1.53	1.48
1	D	16[G]	HZP	CA-C	2.64	1.53	1.48
1	A	16[B]	HZP	CA-C	2.64	1.53	1.48
1	D	16[E]	HZP	CA-C	2.61	1.53	1.48
1	D	16[F]	HZP	CA-C	2.61	1.53	1.48
1	A	16[C]	HZP	CA-C	2.60	1.53	1.48
1	A	16[A]	HZP	CA-C	2.56	1.53	1.48
1	C	16[G]	HZP	CA-C	2.52	1.53	1.48
1	E	13[B]	HZP	CA-C	2.43	1.52	1.48
1	E	16[D]	HZP	CA-C	2.43	1.52	1.48
1	E	16[C]	HZP	CA-C	2.43	1.52	1.48
1	E	13[A]	HZP	CA-C	2.40	1.52	1.48
1	E	13[E]	HZP	CA-C	2.36	1.52	1.48
1	F	11[G]	HYP	CA-C	2.36	1.52	1.48
1	E	29[B]	HYP	CA-C	2.33	1.52	1.48
1	A	16[E]	HZP	CA-C	2.32	1.52	1.48
1	A	16[G]	HZP	CA-C	2.32	1.52	1.48
1	A	16[F]	HZP	CA-C	2.32	1.52	1.48
1	B	13[F]	HZP	CA-C	2.29	1.52	1.48
1	A	20[D]	HYP	CA-C	2.28	1.52	1.48
1	B	13[G]	HZP	CA-C	2.28	1.52	1.48
1	E	29[A]	HYP	CA-C	2.26	1.52	1.48
1	E	29[E]	HYP	CA-C	2.26	1.52	1.48
1	G	16[F]	HZP	CA-C	2.25	1.52	1.48
1	G	16[E]	HZP	CA-C	2.25	1.52	1.48
1	G	16[G]	HZP	CA-C	2.25	1.52	1.48
1	D	5[D]	HYP	CA-C	2.22	1.52	1.48
1	F	16[G]	HZP	CA-C	2.15	1.52	1.48
1	D	2[B]	HYP	CA-C	2.12	1.52	1.48
1	D	2[A]	HYP	CA-C	2.11	1.52	1.48
1	G	16[B]	HZP	CA-C	2.11	1.52	1.48
1	G	16[A]	HZP	CA-C	2.11	1.52	1.48
1	G	16[C]	HZP	CA-C	2.11	1.52	1.48
1	F	16[C]	HZP	CA-C	2.09	1.52	1.48
1	B	16[B]	HZP	CA-C	2.08	1.52	1.48
1	I	16[G]	HZP	CA-C	2.07	1.52	1.48
1	B	16[A]	HZP	CA-C	2.07	1.52	1.48
1	G	13[E]	HZP	CA-C	2.06	1.52	1.48
1	G	13[G]	HZP	CA-C	2.06	1.52	1.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	5[D]	HYP	CA-C	2.06	1.52	1.48
1	G	8[B]	HYP	CA-C	2.06	1.52	1.48
1	G	5[G]	HYP	CA-C	2.05	1.52	1.48
1	G	5[E]	HYP	CA-C	2.05	1.52	1.48
1	B	16[E]	HZP	CA-C	2.05	1.52	1.48
1	G	5[F]	HYP	CA-C	2.05	1.52	1.48
1	G	11[F]	HYP	CA-C	2.04	1.52	1.48
1	F	16[B]	HZP	CA-C	2.03	1.52	1.48
1	G	11[D]	HYP	CA-C	2.03	1.52	1.48
1	G	13[F]	HZP	CA-C	2.03	1.52	1.48
1	D	2[C]	HYP	CA-C	2.03	1.52	1.48
1	B	2[D]	HYP	CA-C	2.01	1.52	1.48
1	B	2[C]	HYP	CA-C	2.01	1.52	1.48
1	D	13[D]	HZP	CA-C	2.01	1.52	1.48
1	G	8[A]	HYP	CA-C	2.01	1.52	1.48

All (1962) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	13[B]	HZP	OD1-CG-CB	-10.46	81.81	110.16
1	B	13[A]	HZP	OD1-CG-CB	-10.45	81.85	110.16
1	B	13[E]	HZP	OD1-CG-CB	-10.44	81.88	110.16
1	B	16[F]	HZP	OD1-CG-CD	-10.15	86.47	110.31
1	B	16[G]	HZP	OD1-CG-CD	-10.15	86.47	110.31
1	C	16[G]	HZP	OD1-CG-CB	-9.49	84.45	110.16
1	B	16[F]	HZP	OD1-CG-CB	-9.41	84.66	110.16
1	B	16[G]	HZP	OD1-CG-CB	-9.41	84.67	110.16
1	F	16[G]	HZP	OD1-CG-CB	-9.26	85.07	110.16
1	A	16[F]	HZP	OD1-CG-CD	-8.90	89.40	110.31
1	A	16[G]	HZP	OD1-CG-CD	-8.87	89.48	110.31
1	A	16[E]	HZP	OD1-CG-CD	-8.84	89.54	110.31
1	G	13[D]	HZP	OD1-CG-CB	-8.76	86.43	110.16
1	C	16[G]	HZP	OD1-CG-CD	-8.72	89.81	110.31
1	E	13[C]	HZP	OD1-CG-CD	-8.35	90.68	110.31
1	E	13[D]	HZP	OD1-CG-CD	-8.35	90.68	110.31
1	A	16[E]	HZP	OD1-CG-CB	-8.25	87.82	110.16
1	A	16[F]	HZP	OD1-CG-CB	-8.24	87.84	110.16
1	A	16[G]	HZP	OD1-CG-CB	-8.23	87.86	110.16
1	F	13[C]	HZP	OD1-CG-CB	-8.20	87.94	110.16
1	F	13[B]	HZP	OD1-CG-CB	-8.19	87.97	110.16
1	F	13[D]	HZP	OD1-CG-CB	-8.18	88.00	110.16
1	I	13[D]	HZP	OD1-CG-CD	-7.99	91.53	110.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	I	13[B]	HZP	OD1-CG-CD	-7.99	91.54	110.31
1	I	13[C]	HZP	OD1-CG-CD	-7.96	91.61	110.31
1	E	16[B]	HZP	OD1-CG-CD	-7.91	91.72	110.31
1	E	16[A]	HZP	OD1-CG-CD	-7.89	91.78	110.31
1	E	16[E]	HZP	OD1-CG-CD	-7.86	91.83	110.31
1	H	16[G]	HZP	OD1-CG-CD	-7.69	92.25	110.31
1	G	16[B]	HZP	OD1-CG-CB	-7.67	89.38	110.16
1	G	16[A]	HZP	OD1-CG-CB	-7.64	89.45	110.16
1	H	16[F]	HZP	OD1-CG-CD	-7.64	92.35	110.31
1	B	13[A]	HZP	OD1-CG-CD	-7.64	92.36	110.31
1	B	13[B]	HZP	OD1-CG-CD	-7.63	92.37	110.31
1	B	13[E]	HZP	OD1-CG-CD	-7.63	92.39	110.31
1	G	16[C]	HZP	OD1-CG-CB	-7.62	89.52	110.16
1	F	16[E]	HZP	OD1-CG-CB	-7.58	89.62	110.16
1	E	13[C]	HZP	CB-CG-CD	7.58	114.23	103.46
1	E	13[D]	HZP	CB-CG-CD	7.58	114.23	103.46
1	F	16[F]	HZP	OD1-CG-CD	-7.56	92.55	110.31
1	F	16[E]	HZP	OD1-CG-CD	-7.56	92.55	110.31
1	F	16[A]	HZP	OD1-CG-CD	-7.56	92.56	110.31
1	F	16[A]	HZP	OD1-CG-CB	-7.55	89.70	110.16
1	E	16[F]	HZP	OD1-CG-CB	-7.54	89.73	110.16
1	E	16[G]	HZP	OD1-CG-CB	-7.54	89.74	110.16
1	F	16[F]	HZP	OD1-CG-CB	-7.53	89.75	110.16
1	C	16[E]	HZP	OD1-CG-CB	-7.51	89.80	110.16
1	G	16[G]	HZP	OD1-CG-CB	-7.51	89.81	110.16
1	G	16[E]	HZP	OD1-CG-CB	-7.51	89.83	110.16
1	G	16[F]	HZP	OD1-CG-CB	-7.50	89.84	110.16
1	C	16[A]	HZP	OD1-CG-CB	-7.49	89.86	110.16
1	C	16[F]	HZP	OD1-CG-CB	-7.48	89.90	110.16
1	A	13[B]	HZP	OD1-CG-CD	-7.47	92.76	110.31
1	F	13[C]	HZP	OD1-CG-CD	-7.47	92.77	110.31
1	F	13[B]	HZP	OD1-CG-CD	-7.45	92.80	110.31
1	F	13[D]	HZP	OD1-CG-CD	-7.45	92.80	110.31
1	H	13[G]	HZP	CG-CD-N	-7.45	93.47	105.47
1	A	13[C]	HZP	OD1-CG-CD	-7.45	92.80	110.31
1	A	13[A]	HZP	OD1-CG-CD	-7.43	92.84	110.31
1	H	13[F]	HZP	CG-CD-N	-7.43	93.50	105.47
1	F	16[C]	HZP	OD1-CG-CB	-7.42	90.06	110.16
1	F	16[B]	HZP	OD1-CG-CB	-7.42	90.07	110.16
1	F	16[D]	HZP	OD1-CG-CB	-7.41	90.08	110.16
1	C	16[E]	HZP	OD1-CG-CD	-7.40	92.93	110.31
1	C	16[A]	HZP	OD1-CG-CD	-7.39	92.96	110.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	16[F]	HZP	OD1-CG-CD	-7.36	93.02	110.31
1	I	16[G]	HZP	OD1-CG-CB	-7.33	90.30	110.16
1	G	13[C]	HZP	OD1-CG-CB	-7.30	90.40	110.16
1	G	13[B]	HZP	OD1-CG-CB	-7.29	90.42	110.16
1	G	13[A]	HZP	OD1-CG-CB	-7.28	90.43	110.16
1	B	13[A]	HZP	CB-CG-CD	7.27	113.80	103.46
1	F	13[G]	HZP	OD1-CG-CB	-7.26	90.50	110.16
1	H	16[D]	HZP	OD1-CG-CB	-7.25	90.51	110.16
1	H	16[C]	HZP	OD1-CG-CB	-7.25	90.51	110.16
1	F	13[E]	HZP	OD1-CG-CB	-7.24	90.54	110.16
1	F	13[F]	HZP	OD1-CG-CB	-7.23	90.57	110.16
1	F	13[A]	HZP	OD1-CG-CB	-7.23	90.57	110.16
1	D	13[D]	HZP	OD1-CG-CD	-7.23	93.32	110.31
1	B	13[B]	HZP	CB-CG-CD	7.23	113.74	103.46
1	H	13[E]	HZP	OD1-CG-CB	-7.22	90.61	110.16
1	H	13[A]	HZP	OD1-CG-CB	-7.22	90.61	110.16
1	B	13[E]	HZP	CB-CG-CD	7.21	113.71	103.46
1	H	13[B]	HZP	OD1-CG-CB	-7.19	90.67	110.16
1	I	16[G]	HZP	OD1-CG-CD	-7.17	93.45	110.31
1	I	16[D]	HZP	OD1-CG-CB	-7.17	90.73	110.16
1	I	16[C]	HZP	OD1-CG-CB	-7.17	90.75	110.16
1	I	16[B]	HZP	OD1-CG-CB	-7.15	90.78	110.16
1	G	16[D]	HZP	OD1-CG-CB	-7.13	90.84	110.16
1	B	16[F]	HZP	CB-CG-CD	7.09	113.55	103.46
1	F	16[G]	HZP	OD1-CG-CD	-7.05	93.75	110.31
1	E	13[C]	HZP	OD1-CG-CB	-7.05	91.07	110.16
1	E	13[D]	HZP	OD1-CG-CB	-7.05	91.07	110.16
1	B	16[G]	HZP	CB-CG-CD	7.04	113.47	103.46
1	D	13[D]	HZP	OD1-CG-CB	-7.03	91.13	110.16
1	B	13[G]	HZP	OD1-CG-CB	-6.94	91.35	110.16
1	B	13[F]	HZP	OD1-CG-CB	-6.94	91.35	110.16
1	C	13[B]	HZP	OD1-CG-CD	-6.88	94.15	110.31
1	C	13[D]	HZP	OD1-CG-CD	-6.84	94.24	110.31
1	C	13[C]	HZP	OD1-CG-CD	-6.84	94.25	110.31
1	F	13[G]	HZP	OD1-CG-CD	-6.78	94.37	110.31
1	B	13[D]	HZP	OD1-CG-CD	-6.76	94.43	110.31
1	B	13[C]	HZP	OD1-CG-CD	-6.75	94.46	110.31
1	C	16[G]	HZP	CB-CG-CD	6.74	113.04	103.46
1	H	13[E]	HZP	OD1-CG-CD	-6.73	94.49	110.31
1	H	13[A]	HZP	OD1-CG-CD	-6.73	94.49	110.31
1	H	13[B]	HZP	OD1-CG-CD	-6.72	94.53	110.31
1	C	11[G]	HYP	CG-CD-N	-6.71	94.66	105.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	16[G]	HZP	OD1-CG-CD	-6.70	94.57	110.31
1	F	13[D]	HZP	CB-CG-CD	6.68	112.96	103.46
1	E	16[F]	HZP	OD1-CG-CD	-6.65	94.68	110.31
1	F	13[B]	HZP	CB-CG-CD	6.60	112.85	103.46
1	I	13[E]	HZP	OD1-CG-CB	-6.58	92.33	110.16
1	I	13[A]	HZP	OD1-CG-CB	-6.58	92.34	110.16
1	C	13[C]	HZP	OD1-CG-CB	-6.56	92.38	110.16
1	I	13[F]	HZP	OD1-CG-CB	-6.55	92.41	110.16
1	C	13[D]	HZP	OD1-CG-CB	-6.55	92.41	110.16
1	C	13[B]	HZP	OD1-CG-CB	-6.55	92.41	110.16
1	G	16[F]	HZP	OD1-CG-CD	-6.55	94.92	110.31
1	F	13[C]	HZP	CB-CG-CD	6.53	112.75	103.46
1	D	16[G]	HZP	OD1-CG-CD	-6.51	95.01	110.31
1	G	16[E]	HZP	OD1-CG-CD	-6.51	95.02	110.31
1	H	16[F]	HZP	OD1-CG-CB	-6.49	92.58	110.16
1	F	16[G]	HZP	CB-CG-CD	6.48	112.68	103.46
1	D	16[F]	HZP	OD1-CG-CD	-6.48	95.09	110.31
1	H	16[G]	HZP	OD1-CG-CB	-6.47	92.62	110.16
1	G	16[G]	HZP	OD1-CG-CD	-6.47	95.11	110.31
1	E	16[E]	HZP	OD1-CG-CB	-6.46	92.65	110.16
1	E	16[A]	HZP	OD1-CG-CB	-6.46	92.67	110.16
1	E	13[E]	HZP	OD1-CG-CB	-6.45	92.68	110.16
1	E	16[B]	HZP	OD1-CG-CB	-6.45	92.69	110.16
1	D	16[E]	HZP	OD1-CG-CD	-6.45	95.16	110.31
1	E	13[A]	HZP	OD1-CG-CB	-6.44	92.72	110.16
1	H	16[B]	HZP	CG-CD-N	-6.43	95.11	105.47
1	E	13[B]	HZP	OD1-CG-CD	-6.42	95.23	110.31
1	E	13[A]	HZP	OD1-CG-CD	-6.42	95.23	110.31
1	E	13[E]	HZP	OD1-CG-CD	-6.42	95.23	110.31
1	E	13[B]	HZP	OD1-CG-CB	-6.42	92.77	110.16
1	G	13[F]	HZP	CB-CG-CD	6.42	112.59	103.46
1	E	16[D]	HZP	OD1-CG-CD	-6.41	95.24	110.31
1	E	16[C]	HZP	OD1-CG-CD	-6.41	95.25	110.31
1	H	16[E]	HZP	CG-CD-N	-6.39	95.17	105.47
1	H	16[A]	HZP	CG-CD-N	-6.39	95.17	105.47
1	E	16[D]	HZP	OD1-CG-CB	-6.38	92.87	110.16
1	E	16[C]	HZP	OD1-CG-CB	-6.38	92.87	110.16
1	I	13[D]	HZP	OD1-CG-CB	-6.38	92.87	110.16
1	G	13[E]	HZP	CB-CG-CD	6.38	112.53	103.46
1	I	13[C]	HZP	OD1-CG-CB	-6.37	92.91	110.16
1	I	13[B]	HZP	OD1-CG-CB	-6.36	92.95	110.16
1	H	16[D]	HZP	OD1-CG-CD	-6.35	95.39	110.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	16[C]	HZP	OD1-CG-CD	-6.35	95.39	110.31
1	G	13[G]	HZP	CB-CG-CD	6.34	112.47	103.46
1	G	13[D]	HZP	OD1-CG-CD	-6.32	95.46	110.31
1	E	16[G]	HZP	CG-CB-CA	-6.31	95.02	104.22
1	E	16[F]	HZP	CG-CB-CA	-6.31	95.02	104.22
1	I	16[B]	HZP	OD1-CG-CD	-6.29	95.53	110.31
1	I	16[E]	HZP	CG-CD-N	-6.28	95.35	105.47
1	I	16[F]	HZP	CG-CD-N	-6.27	95.37	105.47
1	I	16[C]	HZP	OD1-CG-CD	-6.27	95.58	110.31
1	I	16[A]	HZP	CG-CD-N	-6.27	95.37	105.47
1	A	16[D]	HZP	OD1-CG-CD	-6.26	95.59	110.31
1	G	13[B]	HZP	OD1-CG-CD	-6.25	95.62	110.31
1	A	13[B]	HZP	CB-CG-CD	6.25	112.34	103.46
1	I	16[D]	HZP	OD1-CG-CD	-6.24	95.65	110.31
1	G	13[A]	HZP	OD1-CG-CD	-6.23	95.66	110.31
1	G	13[C]	HZP	OD1-CG-CD	-6.21	95.71	110.31
1	A	13[C]	HZP	CB-CG-CD	6.21	112.29	103.46
1	G	16[D]	HZP	OD1-CG-CD	-6.19	95.76	110.31
1	B	5[D]	HYP	CD-N-CA	6.18	119.81	108.15
1	A	13[A]	HZP	CB-CG-CD	6.17	112.23	103.46
1	F	13[G]	HZP	CB-CG-CD	6.16	112.22	103.46
1	I	13[G]	HZP	CG-CD-N	-6.11	95.63	105.47
1	A	16[G]	HZP	CB-CG-CD	6.07	112.09	103.46
1	A	16[E]	HZP	CB-CG-CD	6.06	112.08	103.46
1	I	13[D]	HZP	CB-CG-CD	6.05	112.06	103.46
1	E	13[G]	HZP	OD1-CG-CD	-6.03	96.13	110.31
1	E	13[F]	HZP	OD1-CG-CD	-6.03	96.14	110.31
1	I	13[B]	HZP	CB-CG-CD	6.01	112.01	103.46
1	F	16[G]	HZP	CG-CD-N	-6.00	95.80	105.47
1	B	13[F]	HZP	OD1-CG-CD	-6.00	96.22	110.31
1	I	13[C]	HZP	CB-CG-CD	6.00	111.99	103.46
1	G	13[E]	HZP	OD1-CG-CD	-5.98	96.25	110.31
1	G	13[G]	HZP	OD1-CG-CD	-5.98	96.25	110.31
1	G	13[F]	HZP	OD1-CG-CD	-5.98	96.25	110.31
1	D	13[E]	HZP	OD1-CG-CD	-5.98	96.26	110.31
1	H	13[D]	HZP	OD1-CG-CB	-5.97	93.99	110.16
1	B	13[G]	HZP	OD1-CG-CD	-5.97	96.29	110.31
1	A	16[F]	HZP	CB-CG-CD	5.96	111.93	103.46
1	D	13[F]	HZP	OD1-CG-CD	-5.95	96.33	110.31
1	H	13[C]	HZP	OD1-CG-CB	-5.95	94.05	110.16
1	B	16[D]	HZP	CG-CD-N	-5.95	95.89	105.47
1	D	13[G]	HZP	OD1-CG-CD	-5.94	96.36	110.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	16[E]	HZP	OD1-CG-CD	-5.94	96.36	110.31
1	H	13[D]	HZP	OD1-CG-CD	-5.94	96.36	110.31
1	B	16[C]	HZP	CG-CD-N	-5.93	95.92	105.47
1	B	16[A]	HZP	OD1-CG-CB	-5.92	94.11	110.16
1	B	16[E]	HZP	OD1-CG-CB	-5.92	94.12	110.16
1	B	16[B]	HZP	OD1-CG-CB	-5.90	94.17	110.16
1	B	16[A]	HZP	OD1-CG-CD	-5.90	96.45	110.31
1	H	13[D]	HZP	CG-CB-CA	-5.89	95.62	104.22
1	H	13[C]	HZP	OD1-CG-CD	-5.89	96.47	110.31
1	B	16[B]	HZP	OD1-CG-CD	-5.87	96.52	110.31
1	C	13[G]	HZP	OD1-CG-CD	-5.87	96.52	110.31
1	H	13[C]	HZP	CG-CB-CA	-5.87	95.66	104.22
1	I	13[F]	HZP	OD1-CG-CD	-5.85	96.57	110.31
1	A	16[D]	HZP	OD1-CG-CB	-5.84	94.33	110.16
1	I	13[E]	HZP	OD1-CG-CD	-5.83	96.62	110.31
1	D	16[B]	HZP	OD1-CG-CD	-5.80	96.67	110.31
1	I	13[A]	HZP	OD1-CG-CD	-5.80	96.67	110.31
1	E	16[G]	HZP	CG-CD-N	-5.80	96.13	105.47
1	H	16[G]	HZP	CB-CG-CD	5.80	111.70	103.46
1	A	16[D]	HZP	CB-CG-CD	5.78	111.68	103.46
1	H	16[F]	HZP	CB-CG-CD	5.78	111.67	103.46
1	E	16[F]	HZP	CG-CD-N	-5.77	96.18	105.47
1	D	16[A]	HZP	OD1-CG-CD	-5.76	96.78	110.31
1	F	13[F]	HZP	OD1-CG-CD	-5.75	96.80	110.31
1	I	13[G]	HZP	CG-CB-CA	-5.75	95.83	104.22
1	A	13[G]	HZP	OD1-CG-CD	-5.74	96.83	110.31
1	I	16[F]	HZP	CG-CB-CA	-5.72	95.87	104.22
1	F	13[A]	HZP	OD1-CG-CD	-5.72	96.88	110.31
1	A	13[F]	HZP	OD1-CG-CD	-5.71	96.89	110.31
1	D	16[C]	HZP	OD1-CG-CD	-5.71	96.89	110.31
1	I	16[A]	HZP	CG-CB-CA	-5.70	95.90	104.22
1	I	16[E]	HZP	CG-CB-CA	-5.70	95.90	104.22
1	C	13[F]	HZP	CG-CD-N	-5.70	96.29	105.47
1	A	13[E]	HZP	OD1-CG-CD	-5.70	96.92	110.31
1	F	13[E]	HZP	OD1-CG-CD	-5.68	96.95	110.31
1	G	16[F]	HZP	CB-CG-CD	5.64	111.49	103.46
1	C	13[A]	HZP	CG-CD-N	-5.63	96.40	105.47
1	G	16[E]	HZP	CB-CG-CD	5.63	111.46	103.46
1	G	16[G]	HZP	CB-CG-CD	5.61	111.44	103.46
1	F	16[C]	HZP	OD1-CG-CD	-5.58	97.19	110.31
1	F	16[B]	HZP	OD1-CG-CD	-5.59	97.19	110.31
1	A	13[D]	HZP	OD1-CG-CB	-5.58	95.06	110.16

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	16[D]	HZP	OD1-CG-CD	-5.57	97.22	110.31
1	C	13[E]	HZP	CG-CD-N	-5.57	96.50	105.47
1	C	16[B]	HZP	CG-CD-N	-5.55	96.53	105.47
1	C	16[C]	HZP	CG-CD-N	-5.55	96.53	105.47
1	C	16[D]	HZP	CG-CD-N	-5.52	96.57	105.47
1	G	13[D]	HZP	CB-CG-CD	5.47	111.24	103.46
1	A	13[A]	HZP	OD1-CG-CB	-5.45	95.41	110.16
1	A	13[C]	HZP	OD1-CG-CB	-5.43	95.45	110.16
1	G	16[B]	HZP	OD1-CG-CD	-5.42	97.59	110.31
1	A	13[B]	HZP	OD1-CG-CB	-5.41	95.49	110.16
1	G	16[A]	HZP	OD1-CG-CD	-5.42	97.59	110.31
1	G	16[C]	HZP	OD1-CG-CD	-5.42	97.59	110.31
1	G	13[G]	HZP	OD1-CG-CB	-5.40	95.52	110.16
1	G	13[E]	HZP	OD1-CG-CB	-5.39	95.55	110.16
1	D	13[G]	HZP	OD1-CG-CB	-5.39	95.55	110.16
1	D	13[F]	HZP	OD1-CG-CB	-5.39	95.57	110.16
1	A	16[A]	HZP	OD1-CG-CD	-5.39	97.66	110.31
1	A	16[C]	HZP	OD1-CG-CD	-5.38	97.66	110.31
1	D	13[E]	HZP	OD1-CG-CB	-5.38	95.58	110.16
1	G	13[F]	HZP	OD1-CG-CB	-5.38	95.59	110.16
1	A	16[B]	HZP	OD1-CG-CD	-5.37	97.69	110.31
1	H	13[B]	HZP	CB-CG-CD	5.35	111.06	103.46
1	A	13[D]	HZP	CB-CG-CD	5.34	111.06	103.46
1	C	13[D]	HZP	CB-CG-CD	5.34	111.05	103.46
1	C	13[B]	HZP	CB-CG-CD	5.34	111.05	103.46
1	C	13[C]	HZP	CB-CG-CD	5.29	110.98	103.46
1	H	13[E]	HZP	CB-CG-CD	5.26	110.94	103.46
1	H	13[A]	HZP	CB-CG-CD	5.26	110.94	103.46
1	G	13[C]	HZP	CB-CG-CD	5.22	110.89	103.46
1	G	16[D]	HZP	CB-CG-CD	5.22	110.88	103.46
1	B	13[A]	HZP	CG-CD-N	-5.20	97.10	105.47
1	H	16[B]	HZP	CG-CB-CA	-5.19	96.65	104.22
1	A	29[E]	HYP	CG-CD-N	-5.19	97.11	105.47
1	B	13[E]	HZP	CG-CD-N	-5.17	97.14	105.47
1	G	13[A]	HZP	CB-CG-CD	5.17	110.81	103.46
1	H	16[E]	HZP	CG-CB-CA	-5.16	96.70	104.22
1	B	13[B]	HZP	CG-CD-N	-5.14	97.19	105.47
1	F	13[C]	HZP	CG-CD-N	-5.12	97.22	105.47
1	F	13[D]	HZP	CG-CD-N	-5.12	97.22	105.47
1	H	16[A]	HZP	CG-CB-CA	-5.12	96.75	104.22
1	G	13[B]	HZP	CB-CG-CD	5.12	110.74	103.46
1	F	13[B]	HZP	CG-CD-N	-5.10	97.25	105.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	29[G]	HYP	CG-CD-N	-5.08	97.28	105.47
1	I	16[D]	HZP	CB-CG-CD	5.08	110.68	103.46
1	I	16[F]	HZP	OD1-CG-CB	-5.08	96.41	110.16
1	I	13[B]	HZP	CD-N-CA	5.05	117.69	108.15
1	I	16[C]	HZP	CB-CG-CD	5.05	110.64	103.46
1	I	16[F]	HZP	OD1-CG-CD	-5.05	98.45	110.31
1	I	16[A]	HZP	OD1-CG-CD	-5.05	98.45	110.31
1	I	16[E]	HZP	OD1-CG-CD	-5.05	98.45	110.31
1	I	16[A]	HZP	OD1-CG-CB	-5.04	96.50	110.16
1	I	16[E]	HZP	OD1-CG-CB	-5.04	96.50	110.16
1	I	16[B]	HZP	CB-CG-CD	5.04	110.63	103.46
1	I	13[G]	HZP	OD1-CG-CB	-5.03	96.54	110.16
1	E	13[G]	HZP	OD1-CG-CB	-5.02	96.56	110.16
1	H	13[F]	HZP	OD1-CG-CB	-5.02	96.56	110.16
1	A	13[G]	HZP	OD1-CG-CB	-5.02	96.57	110.16
1	H	13[G]	HZP	OD1-CG-CB	-5.02	96.57	110.16
1	A	29[F]	HYP	CG-CD-N	-5.01	97.40	105.47
1	I	13[G]	HZP	OD1-CG-CD	-5.00	98.55	110.31
1	E	13[F]	HZP	OD1-CG-CB	-5.00	96.61	110.16
1	A	13[E]	HZP	OD1-CG-CB	-5.00	96.61	110.16
1	E	16[B]	HZP	CB-CG-CD	5.00	110.57	103.46
1	A	13[F]	HZP	OD1-CG-CB	-4.99	96.64	110.16
1	I	13[C]	HZP	CD-N-CA	4.99	117.57	108.15
1	B	13[D]	HZP	OD1-CG-CB	-4.94	96.77	110.16
1	E	16[A]	HZP	CB-CG-CD	4.94	110.48	103.46
1	B	13[C]	HZP	OD1-CG-CB	-4.93	96.81	110.16
1	I	13[D]	HZP	CD-N-CA	4.93	117.45	108.15
1	G	2[G]	HYP	CG-CD-N	-4.92	97.55	105.47
1	G	2[E]	HYP	CG-CD-N	-4.92	97.55	105.47
1	G	2[F]	HYP	CG-CD-N	-4.92	97.55	105.47
1	A	13[D]	HZP	OD1-CG-CD	-4.92	98.76	110.31
1	H	16[D]	HZP	CB-CG-CD	4.89	110.42	103.46
1	H	16[C]	HZP	CB-CG-CD	4.89	110.42	103.46
1	E	29[B]	HYP	CG-CD-N	-4.89	97.60	105.47
1	E	29[A]	HYP	CG-CD-N	-4.88	97.61	105.47
1	E	16[E]	HZP	CB-CG-CD	4.88	110.40	103.46
1	E	13[B]	HZP	CB-CG-CD	4.87	110.38	103.46
1	E	16[C]	HZP	CB-CG-CD	4.86	110.38	103.46
1	E	13[A]	HZP	CB-CG-CD	4.86	110.36	103.46
1	E	29[E]	HYP	CG-CD-N	-4.85	97.66	105.47
1	E	13[E]	HZP	CB-CG-CD	4.84	110.35	103.46
1	E	16[D]	HZP	CB-CG-CD	4.83	110.32	103.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	16[F]	HZP	CB-CG-CD	4.82	110.32	103.46
1	I	16[G]	HZP	CD-N-CA	4.79	117.19	108.15
1	C	16[A]	HZP	CB-CG-CD	4.77	110.25	103.46
1	A	16[C]	HZP	OD1-CG-CB	-4.75	97.28	110.16
1	A	16[A]	HZP	OD1-CG-CB	-4.75	97.28	110.16
1	I	13[F]	HZP	CB-CG-CD	4.75	110.22	103.46
1	B	13[G]	HZP	CB-CG-CD	4.75	110.21	103.46
1	A	16[B]	HZP	OD1-CG-CB	-4.75	97.30	110.16
1	D	13[B]	HZP	OD1-CG-CB	-4.74	97.31	110.16
1	D	13[C]	HZP	OD1-CG-CB	-4.74	97.32	110.16
1	D	13[A]	HZP	OD1-CG-CB	-4.74	97.33	110.16
1	B	13[E]	HZP	CD-N-CA	4.73	117.07	108.15
1	I	13[A]	HZP	CB-CG-CD	4.72	110.17	103.46
1	F	16[D]	HZP	CB-CG-CD	4.71	110.16	103.46
1	B	13[A]	HZP	CD-N-CA	4.71	117.05	108.15
1	B	13[B]	HZP	CD-N-CA	4.71	117.03	108.15
1	G	20[D]	HYP	CD-N-CA	4.70	117.02	108.15
1	D	16[D]	HZP	OD1-CG-CB	-4.70	97.44	110.16
1	I	13[E]	HZP	CB-CG-CD	4.70	110.14	103.46
1	B	13[F]	HZP	CB-CG-CD	4.69	110.14	103.46
1	C	16[E]	HZP	CB-CG-CD	4.69	110.13	103.46
1	F	16[B]	HZP	CB-CG-CD	4.69	110.13	103.46
1	D	16[G]	HZP	CG-CD-N	-4.68	97.93	105.47
1	G	13[D]	HZP	CG-CD-N	-4.68	97.93	105.47
1	F	16[C]	HZP	CB-CG-CD	4.67	110.11	103.46
1	C	16[G]	HZP	CD-N-CA	4.67	116.96	108.15
1	I	20[B]	HYP	CG-CD-N	-4.64	97.99	105.47
1	B	13[C]	HZP	CB-CG-CD	4.63	110.05	103.46
1	H	13[D]	HZP	CG-CD-N	-4.63	98.02	105.47
1	I	20[C]	HYP	CG-CD-N	-4.63	98.02	105.47
1	D	16[F]	HZP	CG-CD-N	-4.62	98.03	105.47
1	D	5[D]	HYP	C-CA-N	4.61	119.40	110.86
1	D	16[B]	HZP	OD1-CG-CB	-4.61	97.68	110.16
1	H	13[C]	HZP	CG-CD-N	-4.60	98.06	105.47
1	D	16[A]	HZP	OD1-CG-CB	-4.59	97.72	110.16
1	B	13[D]	HZP	CB-CG-CD	4.59	109.99	103.46
1	D	16[E]	HZP	CG-CD-N	-4.59	98.08	105.47
1	D	16[C]	HZP	OD1-CG-CB	-4.58	97.75	110.16
1	F	13[E]	HZP	CB-CG-CD	4.58	109.97	103.46
1	G	13[D]	HZP	CD-N-CA	4.57	116.78	108.15
1	E	13[F]	HZP	CB-CG-CD	4.57	109.96	103.46
1	D	13[D]	HZP	CB-CG-CD	4.56	109.95	103.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	13[A]	HZP	CB-CG-CD	4.55	109.93	103.46
1	A	29[A]	HYP	CD-N-CA	4.55	116.74	108.15
1	I	20[D]	HYP	CG-CD-N	-4.55	98.15	105.47
1	D	13[C]	HZP	CG-CD-N	-4.54	98.16	105.47
1	B	16[D]	HZP	CD-N-CA	4.53	116.71	108.15
1	D	13[B]	HZP	CG-CD-N	-4.51	98.20	105.47
1	D	13[A]	HZP	CG-CD-N	-4.51	98.21	105.47
1	B	16[C]	HZP	CD-N-CA	4.51	116.66	108.15
1	I	16[G]	HZP	CB-CG-CD	4.50	109.86	103.46
1	F	13[F]	HZP	CB-CG-CD	4.50	109.86	103.46
1	D	16[D]	HZP	CG-CD-N	-4.50	98.23	105.47
1	E	13[G]	HZP	CB-CG-CD	4.49	109.84	103.46
1	A	29[C]	HYP	CD-N-CA	4.48	116.61	108.15
1	F	13[C]	HZP	CD-N-CA	4.47	116.58	108.15
1	F	13[B]	HZP	CD-N-CA	4.45	116.56	108.15
1	A	13[E]	HZP	CB-CG-CD	4.44	109.77	103.46
1	E	13[A]	HZP	CD-N-CA	4.44	116.53	108.15
1	E	13[B]	HZP	CD-N-CA	4.42	116.50	108.15
1	F	13[D]	HZP	CD-N-CA	4.42	116.50	108.15
1	D	13[G]	HZP	CB-CG-CD	4.43	109.75	103.46
1	A	29[B]	HYP	CD-N-CA	4.42	116.49	108.15
1	D	13[B]	HZP	OD1-CG-CD	-4.42	99.93	110.31
1	A	13[F]	HZP	CD-N-CA	4.42	116.48	108.15
1	E	16[F]	HZP	C-CA-N	4.41	119.02	110.86
1	D	13[F]	HZP	CB-CG-CD	4.41	109.73	103.46
1	E	16[G]	HZP	C-CA-N	4.40	119.01	110.86
1	E	13[E]	HZP	CD-N-CA	4.39	116.44	108.15
1	D	16[D]	HZP	CG-CB-CA	-4.39	97.81	104.22
1	E	2[C]	HYP	CD-N-CA	4.39	116.44	108.15
1	E	2[D]	HYP	CD-N-CA	4.39	116.44	108.15
1	D	13[E]	HZP	CB-CG-CD	4.39	109.70	103.46
1	D	13[A]	HZP	OD1-CG-CD	-4.39	100.01	110.31
1	E	16[D]	HZP	CD-N-CA	4.38	116.42	108.15
1	E	26[F]	HYP	C-CA-N	4.38	118.96	110.86
1	E	26[G]	HYP	C-CA-N	4.38	118.96	110.86
1	A	13[F]	HZP	CB-CG-CD	4.38	109.68	103.46
1	D	13[C]	HZP	OD1-CG-CD	-4.37	100.03	110.31
1	C	29[C]	HYP	CG-CD-N	-4.37	98.44	105.47
1	E	16[C]	HZP	CD-N-CA	4.36	116.37	108.15
1	G	29[E]	HYP	CD-N-CA	4.35	116.36	108.15
1	A	13[G]	HZP	CD-N-CA	4.35	116.36	108.15
1	A	13[E]	HZP	CD-N-CA	4.35	116.36	108.15

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	29[B]	HYP	CG-CD-N	-4.35	98.47	105.47
1	C	2[C]	HYP	CD-N-CA	4.34	116.35	108.15
1	G	29[F]	HYP	CD-N-CA	4.34	116.34	108.15
1	A	16[C]	HZP	CB-CG-CD	4.34	109.63	103.46
1	C	2[B]	HYP	CD-N-CA	4.33	116.33	108.15
1	B	16[A]	HZP	CB-CG-CD	4.33	109.62	103.46
1	C	29[D]	HYP	CG-CD-N	-4.33	98.50	105.47
1	A	11[G]	HYP	CD-N-CA	4.33	116.32	108.15
1	A	11[E]	HYP	CD-N-CA	4.33	116.32	108.15
1	G	23[D]	HYP	CG-CD-N	-4.33	98.50	105.47
1	C	23[G]	HYP	CG-CD-N	-4.32	98.51	105.47
1	G	20[C]	HYP	CG-CD-N	-4.32	98.51	105.47
1	C	2[D]	HYP	CD-N-CA	4.32	116.30	108.15
1	G	29[G]	HYP	CD-N-CA	4.32	116.30	108.15
1	A	11[F]	HYP	CD-N-CA	4.32	116.30	108.15
1	B	16[B]	HZP	CB-CG-CD	4.31	109.60	103.46
1	A	13[G]	HZP	CB-CG-CD	4.31	109.59	103.46
1	B	16[E]	HZP	CB-CG-CD	4.31	109.59	103.46
1	H	13[G]	HZP	OD1-CG-CD	-4.31	100.19	110.31
1	G	20[A]	HYP	CG-CD-N	-4.30	98.54	105.47
1	A	16[A]	HZP	CB-CG-CD	4.30	109.58	103.46
1	A	8[D]	HYP	CD-N-CA	4.30	116.27	108.15
1	G	20[B]	HYP	CG-CD-N	-4.29	98.56	105.47
1	H	13[F]	HZP	OD1-CG-CD	-4.29	100.24	110.31
1	A	16[A]	HZP	CD-N-CA	4.28	116.23	108.15
1	I	26[A]	HYP	CG-CD-N	-4.27	98.59	105.47
1	A	20[D]	HYP	CD-N-CA	4.26	116.19	108.15
1	A	2[E]	HYP	CD-N-CA	4.26	116.19	108.15
1	A	2[F]	HYP	CD-N-CA	4.25	116.18	108.15
1	C	13[G]	HZP	OD1-CG-CB	-4.25	98.65	110.16
1	A	16[B]	HZP	CB-CG-CD	4.25	109.50	103.46
1	I	26[E]	HYP	CG-CD-N	-4.24	98.64	105.47
1	D	13[C]	HZP	CG-CB-CA	-4.24	98.04	104.22
1	D	16[D]	HZP	OD1-CG-CD	-4.23	100.36	110.31
1	A	2[G]	HYP	CD-N-CA	4.24	116.14	108.15
1	A	16[C]	HZP	CD-N-CA	4.22	116.11	108.15
1	H	13[B]	HZP	CD-N-CA	4.21	116.10	108.15
1	I	26[F]	HYP	CG-CD-N	-4.21	98.69	105.47
1	H	11[G]	HYP	CD-N-CA	4.20	116.08	108.15
1	D	13[A]	HZP	CG-CB-CA	-4.20	98.10	104.22
1	D	16[B]	HZP	CG-CD-N	-4.19	98.73	105.47
1	D	13[B]	HZP	CG-CB-CA	-4.19	98.12	104.22

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	13[E]	HZP	CD-N-CA	4.17	116.03	108.15
1	H	13[A]	HZP	CD-N-CA	4.17	116.03	108.15
1	D	16[F]	HZP	CD-N-CA	4.17	116.03	108.15
1	D	16[E]	HZP	CD-N-CA	4.16	116.01	108.15
1	A	16[B]	HZP	CD-N-CA	4.16	116.00	108.15
1	H	11[F]	HYP	CD-N-CA	4.15	115.99	108.15
1	D	16[G]	HZP	CD-N-CA	4.15	115.98	108.15
1	E	13[D]	HZP	CD-N-CA	4.15	115.98	108.15
1	C	16[E]	HZP	CD-N-CA	4.15	115.98	108.15
1	C	16[A]	HZP	CD-N-CA	4.15	115.98	108.15
1	C	16[F]	HZP	CD-N-CA	4.15	115.98	108.15
1	F	29[C]	HYP	CD-N-CA	4.14	115.97	108.15
1	D	16[A]	HZP	CG-CD-N	-4.14	98.81	105.47
1	B	16[G]	HZP	CD-N-CA	4.13	115.95	108.15
1	E	13[C]	HZP	CD-N-CA	4.13	115.94	108.15
1	B	16[F]	HZP	CD-N-CA	4.13	115.94	108.15
1	F	16[G]	HZP	CD-N-CA	4.12	115.93	108.15
1	C	26[F]	HYP	CD-N-CA	4.12	115.93	108.15
1	F	29[B]	HYP	CD-N-CA	4.11	115.92	108.15
1	I	29[A]	HYP	CD-N-CA	4.11	115.91	108.15
1	C	13[F]	HZP	OD1-CG-CD	-4.10	100.69	110.31
1	I	29[F]	HYP	CD-N-CA	4.10	115.88	108.15
1	F	29[C]	HYP	CG-CD-N	-4.09	98.88	105.47
1	I	29[E]	HYP	CD-N-CA	4.09	115.87	108.15
1	C	26[A]	HYP	CD-N-CA	4.09	115.87	108.15
1	E	8[G]	HYP	CD-N-CA	4.09	115.87	108.15
1	F	29[D]	HYP	CD-N-CA	4.09	115.87	108.15
1	I	13[F]	HZP	CD-N-CA	4.09	115.87	108.15
1	F	5[A]	HYP	CD-N-CA	4.08	115.86	108.15
1	D	16[C]	HZP	CG-CD-N	-4.09	98.89	105.47
1	I	13[E]	HZP	CD-N-CA	4.08	115.86	108.15
1	F	5[F]	HYP	CD-N-CA	4.08	115.86	108.15
1	H	2[E]	HYP	CD-N-CA	4.08	115.86	108.15
1	H	2[A]	HYP	CD-N-CA	4.08	115.86	108.15
1	H	2[B]	HYP	CD-N-CA	4.08	115.86	108.15
1	H	29[E]	HYP	CD-N-CA	4.08	115.84	108.15
1	E	8[F]	HYP	CD-N-CA	4.07	115.83	108.15
1	F	5[E]	HYP	CD-N-CA	4.07	115.83	108.15
1	A	2[A]	HYP	CD-N-CA	4.07	115.83	108.15
1	H	29[B]	HYP	CD-N-CA	4.07	115.83	108.15
1	C	13[A]	HZP	OD1-CG-CD	-4.07	100.75	110.31
1	A	5[D]	HYP	CD-N-CA	4.06	115.82	108.15

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	2[B]	HYP	CD-N-CA	4.06	115.82	108.15
1	F	29[B]	HYP	CG-CD-N	-4.06	98.93	105.47
1	C	26[E]	HYP	CD-N-CA	4.06	115.82	108.15
1	H	29[A]	HYP	CD-N-CA	4.06	115.81	108.15
1	I	13[B]	HZP	CG-CD-N	-4.06	98.94	105.47
1	I	13[D]	HZP	CG-CD-N	-4.05	98.95	105.47
1	F	2[G]	HYP	CD-N-CA	4.05	115.79	108.15
1	C	13[E]	HZP	OD1-CG-CD	-4.04	100.82	110.31
1	I	13[A]	HZP	CD-N-CA	4.04	115.78	108.15
1	I	13[C]	HZP	CG-CD-N	-4.03	98.98	105.47
1	F	29[D]	HYP	CG-CD-N	-4.03	98.98	105.47
1	A	2[C]	HYP	CD-N-CA	4.02	115.73	108.15
1	B	13[C]	HZP	CD-N-CA	4.01	115.71	108.15
1	B	5[G]	HYP	CD-N-CA	4.00	115.70	108.15
1	B	5[F]	HYP	CD-N-CA	4.00	115.70	108.15
1	B	13[D]	HZP	CD-N-CA	4.00	115.69	108.15
1	G	13[G]	HZP	CD-N-CA	3.99	115.69	108.15
1	A	5[B]	HYP	CD-N-CA	3.99	115.68	108.15
1	A	11[G]	HYP	CG-CD-N	-3.99	99.05	105.47
1	G	2[A]	HYP	CD-N-CA	3.98	115.67	108.15
1	B	8[B]	HYP	CD-N-CA	3.98	115.66	108.15
1	G	2[C]	HYP	CD-N-CA	3.98	115.66	108.15
1	G	5[D]	HYP	CD-N-CA	3.98	115.66	108.15
1	G	13[E]	HZP	CD-N-CA	3.98	115.66	108.15
1	A	5[C]	HYP	CD-N-CA	3.98	115.66	108.15
1	D	5[D]	HYP	CG-CD-N	-3.96	99.09	105.47
1	H	13[F]	HZP	CG-CB-CA	-3.96	98.45	104.22
1	E	29[E]	HYP	CD-N-CA	3.96	115.63	108.15
1	H	16[G]	HZP	CD-N-CA	3.96	115.62	108.15
1	H	16[F]	HZP	CD-N-CA	3.96	115.62	108.15
1	A	11[E]	HYP	CG-CD-N	-3.96	99.10	105.47
1	G	13[F]	HZP	CD-N-CA	3.95	115.61	108.15
1	E	11[B]	HYP	CD-N-CA	3.95	115.60	108.15
1	C	23[G]	HYP	CD-N-CA	3.94	115.59	108.15
1	C	16[G]	HZP	CG-CD-N	-3.94	99.12	105.47
1	E	29[A]	HYP	CD-N-CA	3.94	115.58	108.15
1	A	5[A]	HYP	CD-N-CA	3.93	115.58	108.15
1	E	11[A]	HYP	CD-N-CA	3.93	115.57	108.15
1	H	13[G]	HZP	CG-CB-CA	-3.93	98.50	104.22
1	B	8[A]	HYP	CD-N-CA	3.92	115.55	108.15
1	E	29[B]	HYP	CD-N-CA	3.92	115.55	108.15
1	G	2[B]	HYP	CD-N-CA	3.92	115.55	108.15

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	11[E]	HYP	CD-N-CA	3.91	115.54	108.15
1	B	16[E]	HZP	CD-N-CA	3.91	115.53	108.15
1	F	13[G]	HZP	CD-N-CA	3.90	115.52	108.15
1	B	16[A]	HZP	CD-N-CA	3.90	115.51	108.15
1	B	16[B]	HZP	CD-N-CA	3.89	115.50	108.15
1	E	13[D]	HZP	CG-CD-N	-3.89	99.21	105.47
1	A	11[F]	HYP	CG-CD-N	-3.89	99.21	105.47
1	B	8[E]	HYP	CD-N-CA	3.87	115.45	108.15
1	A	13[B]	HZP	CD-N-CA	3.87	115.45	108.15
1	C	13[D]	HZP	CD-N-CA	3.86	115.44	108.15
1	C	13[B]	HZP	CD-N-CA	3.86	115.44	108.15
1	A	13[D]	HZP	CG-CD-N	-3.86	99.26	105.47
1	H	16[G]	HZP	CG-CD-N	-3.85	99.27	105.47
1	C	13[C]	HZP	CD-N-CA	3.85	115.42	108.15
1	E	13[C]	HZP	CG-CD-N	-3.85	99.27	105.47
1	H	26[F]	HYP	CD-N-CA	3.84	115.40	108.15
1	H	16[B]	HZP	OD1-CG-CB	-3.83	99.78	110.16
1	A	13[C]	HZP	CD-N-CA	3.83	115.38	108.15
1	H	16[F]	HZP	CG-CD-N	-3.83	99.30	105.47
1	H	26[G]	HYP	CD-N-CA	3.82	115.37	108.15
1	H	16[E]	HZP	OD1-CG-CB	-3.82	99.81	110.16
1	H	16[A]	HZP	OD1-CG-CB	-3.82	99.82	110.16
1	A	20[B]	HYP	CD-N-CA	3.82	115.35	108.15
1	A	20[A]	HYP	CD-N-CA	3.81	115.35	108.15
1	G	16[B]	HZP	CB-CG-CD	3.81	108.88	103.46
1	A	20[D]	HYP	C-CA-N	3.80	117.90	110.86
1	A	13[A]	HZP	CD-N-CA	3.80	115.33	108.15
1	B	5[D]	HYP	CG-CD-N	-3.79	99.37	105.47
1	D	13[D]	HZP	CD-N-CA	3.79	115.31	108.15
1	G	16[A]	HZP	CB-CG-CD	3.79	108.85	103.46
1	A	20[C]	HYP	CD-N-CA	3.79	115.30	108.15
1	E	20[E]	HYP	CD-N-CA	3.78	115.29	108.15
1	B	29[E]	HYP	CD-N-CA	3.78	115.28	108.15
1	F	16[A]	HZP	CG-CD-N	-3.78	99.39	105.47
1	F	16[F]	HZP	CG-CD-N	-3.78	99.39	105.47
1	F	16[E]	HZP	CG-CD-N	-3.78	99.39	105.47
1	A	8[B]	HYP	CD-N-CA	3.77	115.27	108.15
1	G	16[C]	HZP	CB-CG-CD	3.77	108.83	103.46
1	A	20[A]	HYP	CG-CD-N	-3.77	99.40	105.47
1	E	20[A]	HYP	CD-N-CA	3.77	115.26	108.15
1	A	20[C]	HYP	CG-CD-N	-3.77	99.40	105.47
1	A	20[B]	HYP	CG-CD-N	-3.77	99.40	105.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	13[F]	HZP	CD-N-CA	3.76	115.25	108.15
1	D	2[G]	HYP	CD-N-CA	3.76	115.25	108.15
1	D	2[F]	HYP	CD-N-CA	3.76	115.25	108.15
1	A	11[D]	HYP	CD-N-CA	3.75	115.24	108.15
1	E	20[B]	HYP	CD-N-CA	3.75	115.23	108.15
1	B	29[A]	HYP	CD-N-CA	3.75	115.23	108.15
1	A	23[D]	HYP	CD-N-CA	3.75	115.22	108.15
1	I	16[F]	HZP	CD-N-CA	3.74	115.22	108.15
1	A	8[A]	HYP	CD-N-CA	3.74	115.22	108.15
1	I	16[E]	HZP	CD-N-CA	3.74	115.21	108.15
1	A	23[D]	HYP	CG-CD-N	-3.74	99.45	105.47
1	B	29[B]	HYP	CD-N-CA	3.74	115.20	108.15
1	F	16[A]	HZP	CD-N-CA	3.73	115.19	108.15
1	D	2[E]	HYP	CD-N-CA	3.73	115.19	108.15
1	F	16[F]	HZP	CD-N-CA	3.73	115.19	108.15
1	F	16[E]	HZP	CD-N-CA	3.73	115.19	108.15
1	H	20[C]	HYP	CG-CD-N	-3.73	99.46	105.47
1	A	8[C]	HYP	CD-N-CA	3.72	115.18	108.15
1	G	8[C]	HYP	CD-N-CA	3.72	115.17	108.15
1	B	20[B]	HYP	CD-N-CA	3.72	115.17	108.15
1	A	5[G]	HYP	CD-N-CA	3.71	115.16	108.15
1	I	16[A]	HZP	CD-N-CA	3.71	115.15	108.15
1	B	13[G]	HZP	CD-N-CA	3.71	115.15	108.15
1	C	13[E]	HZP	CG-CB-CA	-3.71	98.82	104.22
1	B	20[E]	HYP	CD-N-CA	3.70	115.14	108.15
1	D	13[G]	HZP	CD-N-CA	3.70	115.14	108.15
1	G	5[G]	HYP	CD-N-CA	3.70	115.13	108.15
1	D	13[F]	HZP	CD-N-CA	3.70	115.13	108.15
1	G	5[E]	HYP	CD-N-CA	3.70	115.13	108.15
1	G	5[F]	HYP	CD-N-CA	3.70	115.13	108.15
1	B	20[A]	HYP	CD-N-CA	3.70	115.13	108.15
1	C	20[C]	HYP	CD-N-CA	3.69	115.12	108.15
1	C	20[D]	HYP	CD-N-CA	3.69	115.12	108.15
1	C	20[B]	HYP	CD-N-CA	3.69	115.12	108.15
1	G	8[A]	HYP	CD-N-CA	3.69	115.12	108.15
1	A	5[E]	HYP	CD-N-CA	3.69	115.12	108.15
1	H	20[D]	HYP	CG-CD-N	-3.69	99.53	105.47
1	F	13[E]	HZP	CD-N-CA	3.69	115.11	108.15
1	C	13[A]	HZP	CG-CB-CA	-3.69	98.84	104.22
1	B	23[C]	HYP	CD-N-CA	3.68	115.10	108.15
1	D	2[C]	HYP	CD-N-CA	3.68	115.11	108.15
1	B	23[D]	HYP	CD-N-CA	3.68	115.10	108.15

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	16[C]	HZP	CG-CD-N	-3.68	99.54	105.47
1	D	13[E]	HZP	CD-N-CA	3.68	115.09	108.15
1	C	13[F]	HZP	CG-CB-CA	-3.68	98.86	104.22
1	F	13[A]	HZP	CD-N-CA	3.67	115.08	108.15
1	E	23[D]	HYP	CD-N-CA	3.67	115.08	108.15
1	G	8[G]	HYP	CD-N-CA	3.67	115.07	108.15
1	G	16[B]	HZP	CG-CD-N	-3.67	99.56	105.47
1	C	29[B]	HYP	CD-N-CA	3.67	115.07	108.15
1	A	5[F]	HYP	CD-N-CA	3.67	115.07	108.15
1	G	16[A]	HZP	CG-CD-N	-3.66	99.57	105.47
1	A	16[D]	HZP	CD-N-CA	3.66	115.06	108.15
1	C	29[D]	HYP	CD-N-CA	3.66	115.05	108.15
1	C	29[C]	HYP	CD-N-CA	3.66	115.05	108.15
1	F	13[F]	HZP	CD-N-CA	3.65	115.05	108.15
1	E	23[C]	HYP	CD-N-CA	3.65	115.05	108.15
1	D	2[A]	HYP	CD-N-CA	3.65	115.04	108.15
1	F	20[B]	HYP	CD-N-CA	3.65	115.04	108.15
1	F	20[D]	HYP	CD-N-CA	3.65	115.04	108.15
1	D	5[C]	HYP	CD-N-CA	3.65	115.04	108.15
1	F	20[C]	HYP	CD-N-CA	3.65	115.04	108.15
1	I	8[E]	HYP	CD-N-CA	3.65	115.03	108.15
1	G	8[B]	HYP	CD-N-CA	3.64	115.03	108.15
1	D	2[B]	HYP	CD-N-CA	3.64	115.03	108.15
1	G	11[C]	HYP	CD-N-CA	3.64	115.03	108.15
1	F	29[F]	HYP	CD-N-CA	3.64	115.02	108.15
1	G	8[E]	HYP	CD-N-CA	3.64	115.03	108.15
1	B	26[F]	HYP	CB-CG-CD	3.64	108.64	103.46
1	I	5[D]	HYP	CD-N-CA	3.64	115.02	108.15
1	G	11[B]	HYP	CD-N-CA	3.64	115.02	108.15
1	I	8[F]	HYP	CD-N-CA	3.64	115.02	108.15
1	G	11[A]	HYP	CD-N-CA	3.63	115.01	108.15
1	I	8[A]	HYP	CD-N-CA	3.63	115.00	108.15
1	B	26[G]	HYP	CB-CG-CD	3.63	108.62	103.46
1	H	26[D]	HYP	CD-N-CA	3.62	114.99	108.15
1	A	29[E]	HYP	CD-N-CA	3.62	114.99	108.15
1	I	5[B]	HYP	CD-N-CA	3.62	114.99	108.15
1	C	13[G]	HZP	CG-CD-N	-3.62	99.64	105.47
1	I	5[G]	HYP	CD-N-CA	3.62	114.98	108.15
1	D	5[A]	HYP	CD-N-CA	3.62	114.98	108.15
1	G	8[F]	HYP	CD-N-CA	3.62	114.98	108.15
1	G	5[B]	HYP	CD-N-CA	3.61	114.96	108.15
1	F	16[C]	HZP	CD-N-CA	3.61	114.96	108.15

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	16[B]	HZP	CD-N-CA	3.61	114.96	108.15
1	I	5[C]	HYP	CD-N-CA	3.61	114.96	108.15
1	I	16[C]	HZP	CD-N-CA	3.61	114.96	108.15
1	B	16[F]	HZP	CG-CD-N	-3.61	99.66	105.47
1	I	16[D]	HZP	CD-N-CA	3.60	114.95	108.15
1	C	16[B]	HZP	CG-CB-CA	-3.60	98.97	104.22
1	F	29[A]	HYP	CD-N-CA	3.60	114.94	108.15
1	C	16[D]	HZP	CG-CB-CA	-3.60	98.97	104.22
1	C	16[C]	HZP	CG-CB-CA	-3.60	98.97	104.22
1	G	5[A]	HYP	CD-N-CA	3.60	114.94	108.15
1	I	16[B]	HZP	CD-N-CA	3.60	114.94	108.15
1	F	16[D]	HZP	CD-N-CA	3.59	114.93	108.15
1	I	26[G]	HYP	CD-N-CA	3.59	114.93	108.15
1	H	23[A]	HYP	CD-N-CA	3.59	114.93	108.15
1	D	5[B]	HYP	CD-N-CA	3.59	114.92	108.15
1	I	2[G]	HYP	CD-N-CA	3.59	114.92	108.15
1	E	5[B]	HYP	CD-N-CA	3.58	114.92	108.15
1	G	5[C]	HYP	CD-N-CA	3.58	114.91	108.15
1	A	29[F]	HYP	CD-N-CA	3.58	114.91	108.15
1	I	13[G]	HZP	CD-N-CA	3.58	114.91	108.15
1	E	5[E]	HYP	CD-N-CA	3.58	114.90	108.15
1	F	11[G]	HYP	CG-CD-N	-3.58	99.71	105.47
1	A	29[G]	HYP	CD-N-CA	3.58	114.90	108.15
1	H	20[B]	HYP	CD-N-CA	3.58	114.90	108.15
1	E	11[C]	HYP	CD-N-CA	3.58	114.90	108.15
1	I	5[F]	HYP	CD-N-CA	3.57	114.89	108.15
1	E	2[E]	HYP	CD-N-CA	3.57	114.89	108.15
1	B	16[G]	HZP	CG-CD-N	-3.57	99.72	105.47
1	F	29[E]	HYP	CD-N-CA	3.57	114.89	108.15
1	F	26[F]	HYP	CD-N-CA	3.56	114.88	108.15
1	D	20[D]	HYP	OD1-CG-CD	3.56	118.69	110.31
1	E	2[A]	HYP	CD-N-CA	3.56	114.87	108.15
1	D	23[D]	HYP	CD-N-CA	3.56	114.87	108.15
1	I	5[A]	HYP	CD-N-CA	3.56	114.87	108.15
1	I	5[E]	HYP	CD-N-CA	3.55	114.86	108.15
1	I	20[B]	HYP	CD-N-CA	3.55	114.85	108.15
1	I	8[C]	HYP	CD-N-CA	3.55	114.85	108.15
1	I	8[B]	HYP	CD-N-CA	3.55	114.85	108.15
1	I	8[D]	HYP	CD-N-CA	3.55	114.85	108.15
1	E	2[B]	HYP	CD-N-CA	3.55	114.85	108.15
1	H	20[E]	HYP	CD-N-CA	3.55	114.85	108.15
1	H	23[B]	HYP	CD-N-CA	3.55	114.85	108.15

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	20[A]	HYP	CD-N-CA	3.55	114.85	108.15
1	B	8[D]	HYP	CD-N-CA	3.55	114.84	108.15
1	E	5[A]	HYP	CD-N-CA	3.54	114.84	108.15
1	E	11[D]	HYP	CD-N-CA	3.54	114.83	108.15
1	F	26[A]	HYP	CD-N-CA	3.54	114.83	108.15
1	A	26[G]	HYP	CD-N-CA	3.54	114.83	108.15
1	F	26[E]	HYP	CD-N-CA	3.54	114.83	108.15
1	I	20[C]	HYP	CD-N-CA	3.53	114.82	108.15
1	E	11[F]	HYP	CG-CD-N	-3.53	99.78	105.47
1	B	5[D]	HYP	CB-CA-N	-3.53	99.18	104.59
1	D	23[A]	HYP	CD-N-CA	3.53	114.81	108.15
1	E	8[B]	HYP	CD-N-CA	3.53	114.81	108.15
1	E	23[C]	HYP	CG-CD-N	-3.53	99.79	105.47
1	D	2[D]	HYP	CD-N-CA	3.53	114.81	108.15
1	I	20[D]	HYP	CD-N-CA	3.52	114.80	108.15
1	D	8[D]	HYP	CD-N-CA	3.52	114.80	108.15
1	E	29[G]	HYP	CD-N-CA	3.52	114.80	108.15
1	C	8[D]	HYP	CD-N-CA	3.52	114.80	108.15
1	E	29[F]	HYP	CD-N-CA	3.52	114.79	108.15
1	D	20[B]	HYP	CD-N-CA	3.51	114.78	108.15
1	A	26[E]	HYP	CD-N-CA	3.51	114.77	108.15
1	H	2[D]	HYP	CD-N-CA	3.51	114.77	108.15
1	H	2[C]	HYP	CD-N-CA	3.51	114.77	108.15
1	B	8[C]	HYP	CD-N-CA	3.51	114.77	108.15
1	F	26[G]	HYP	CD-N-CA	3.51	114.77	108.15
1	D	20[A]	HYP	CD-N-CA	3.50	114.76	108.15
1	B	5[B]	HYP	CD-N-CA	3.50	114.76	108.15
1	D	20[C]	HYP	CD-N-CA	3.50	114.76	108.15
1	C	5[A]	HYP	CD-N-CA	3.50	114.75	108.15
1	I	11[C]	HYP	CD-N-CA	3.50	114.75	108.15
1	C	5[E]	HYP	CD-N-CA	3.50	114.75	108.15
1	H	26[C]	HYP	CD-N-CA	3.50	114.75	108.15
1	C	8[G]	HYP	CD-N-CA	3.50	114.75	108.15
1	I	11[B]	HYP	CD-N-CA	3.50	114.75	108.15
1	C	5[F]	HYP	CD-N-CA	3.49	114.75	108.15
1	F	8[G]	HYP	CD-N-CA	3.50	114.75	108.15
1	I	11[D]	HYP	CD-N-CA	3.49	114.74	108.15
1	C	8[C]	HYP	CD-N-CA	3.49	114.74	108.15
1	C	8[B]	HYP	CD-N-CA	3.49	114.74	108.15
1	E	23[D]	HYP	CG-CD-N	-3.49	99.85	105.47
1	B	2[D]	HYP	CG-CD-N	-3.49	99.85	105.47
1	B	2[C]	HYP	CG-CD-N	-3.49	99.85	105.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	23[C]	HYP	CD-N-CA	3.49	114.74	108.15
1	H	23[D]	HYP	CD-N-CA	3.49	114.74	108.15
1	E	11[G]	HYP	CG-CD-N	-3.49	99.85	105.47
1	E	8[E]	HYP	CD-N-CA	3.49	114.73	108.15
1	G	29[B]	HYP	CD-N-CA	3.49	114.73	108.15
1	E	20[B]	HYP	CG-CD-N	-3.49	99.86	105.47
1	E	20[E]	HYP	CG-CD-N	-3.49	99.86	105.47
1	D	26[D]	HYP	CD-N-CA	3.49	114.73	108.15
1	E	20[A]	HYP	CG-CD-N	-3.49	99.86	105.47
1	E	20[C]	HYP	CD-N-CA	3.48	114.73	108.15
1	I	20[E]	HYP	CD-N-CA	3.48	114.73	108.15
1	C	26[F]	HYP	CB-CG-CD	3.48	108.41	103.46
1	I	26[A]	HYP	CD-N-CA	3.48	114.72	108.15
1	I	26[F]	HYP	CD-N-CA	3.48	114.72	108.15
1	I	26[E]	HYP	CD-N-CA	3.48	114.72	108.15
1	C	2[G]	HYP	CD-N-CA	3.48	114.72	108.15
1	C	16[B]	HZP	OD1-CG-CD	-3.48	102.14	110.31
1	C	16[C]	HZP	OD1-CG-CD	-3.48	102.14	110.31
1	C	26[A]	HYP	CB-CG-CD	3.48	108.40	103.46
1	I	20[F]	HYP	CD-N-CA	3.48	114.71	108.15
1	F	16[E]	HZP	CB-CG-CD	3.48	108.41	103.46
1	I	20[A]	HYP	CD-N-CA	3.47	114.70	108.15
1	B	20[D]	HYP	CD-N-CA	3.47	114.70	108.15
1	E	20[D]	HYP	CD-N-CA	3.47	114.70	108.15
1	B	20[C]	HYP	CD-N-CA	3.47	114.70	108.15
1	G	16[F]	HZP	CD-N-CA	3.47	114.70	108.15
1	C	11[E]	HYP	CD-N-CA	3.47	114.69	108.15
1	H	23[E]	HYP	CD-N-CA	3.47	114.69	108.15
1	G	16[B]	HZP	CD-N-CA	3.47	114.69	108.15
1	E	8[A]	HYP	CD-N-CA	3.47	114.69	108.15
1	B	5[E]	HYP	CD-N-CA	3.46	114.69	108.15
1	D	23[B]	HYP	CD-N-CA	3.46	114.69	108.15
1	F	16[A]	HZP	CB-CG-CD	3.46	108.39	103.46
1	C	11[F]	HYP	CD-N-CA	3.46	114.69	108.15
1	E	5[G]	HYP	CD-N-CA	3.46	114.69	108.15
1	C	11[A]	HYP	CD-N-CA	3.46	114.68	108.15
1	G	23[F]	HYP	CD-N-CA	3.46	114.68	108.15
1	G	23[G]	HYP	CD-N-CA	3.46	114.68	108.15
1	B	5[A]	HYP	CD-N-CA	3.46	114.68	108.15
1	E	5[D]	HYP	CD-N-CA	3.46	114.68	108.15
1	G	23[E]	HYP	CD-N-CA	3.46	114.68	108.15
1	B	29[G]	HYP	CD-N-CA	3.46	114.68	108.15

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	29[F]	HYP	CD-N-CA	3.46	114.68	108.15
1	C	16[D]	HZP	OD1-CG-CD	-3.46	102.19	110.31
1	A	26[F]	HYP	CD-N-CA	3.45	114.67	108.15
1	F	11[F]	HYP	CD-N-CA	3.45	114.67	108.15
1	G	29[A]	HYP	CD-N-CA	3.45	114.66	108.15
1	F	11[G]	HYP	CB-CA-N	-3.45	99.30	104.59
1	F	16[F]	HZP	CB-CG-CD	3.45	108.36	103.46
1	E	5[F]	HYP	CD-N-CA	3.44	114.65	108.15
1	E	5[C]	HYP	CD-N-CA	3.44	114.64	108.15
1	G	26[F]	HYP	CD-N-CA	3.44	114.64	108.15
1	F	11[A]	HYP	CD-N-CA	3.44	114.64	108.15
1	B	2[C]	HYP	CD-N-CA	3.44	114.64	108.15
1	G	16[E]	HZP	CD-N-CA	3.44	114.64	108.15
1	D	23[C]	HYP	CD-N-CA	3.43	114.63	108.15
1	G	16[G]	HZP	CD-N-CA	3.43	114.63	108.15
1	F	11[E]	HYP	CD-N-CA	3.43	114.62	108.15
1	G	20[G]	HYP	CD-N-CA	3.43	114.62	108.15
1	I	23[G]	HYP	CD-N-CA	3.43	114.62	108.15
1	B	20[D]	HYP	CG-CD-N	-3.42	99.96	105.47
1	B	2[D]	HYP	CD-N-CA	3.42	114.61	108.15
1	G	16[A]	HZP	CD-N-CA	3.42	114.61	108.15
1	G	16[C]	HZP	CD-N-CA	3.42	114.61	108.15
1	E	16[B]	HZP	CD-N-CA	3.42	114.60	108.15
1	G	20[E]	HYP	CD-N-CA	3.41	114.59	108.15
1	G	11[D]	HYP	CG-CD-N	-3.41	99.97	105.47
1	G	29[C]	HYP	CD-N-CA	3.41	114.59	108.15
1	B	20[C]	HYP	CG-CD-N	-3.41	99.98	105.47
1	C	26[E]	HYP	CB-CG-CD	3.41	108.31	103.46
1	G	29[D]	HYP	CD-N-CA	3.40	114.57	108.15
1	F	11[G]	HYP	CD-N-CA	3.40	114.57	108.15
1	G	20[F]	HYP	CD-N-CA	3.40	114.57	108.15
1	I	2[D]	HYP	CD-N-CA	3.40	114.56	108.15
1	I	23[D]	HYP	CD-N-CA	3.40	114.56	108.15
1	I	2[C]	HYP	CD-N-CA	3.40	114.56	108.15
1	B	2[F]	HYP	CD-N-CA	3.39	114.55	108.15
1	I	23[C]	HYP	CD-N-CA	3.39	114.54	108.15
1	D	20[G]	HYP	CD-N-CA	3.39	114.54	108.15
1	F	8[B]	HYP	CD-N-CA	3.39	114.54	108.15
1	F	8[D]	HYP	CD-N-CA	3.39	114.54	108.15
1	F	8[C]	HYP	CD-N-CA	3.39	114.54	108.15
1	G	23[D]	HYP	CD-N-CA	3.38	114.54	108.15
1	C	20[C]	HYP	OD1-CG-CB	-3.38	101.01	110.16

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	5[D]	HYP	CD-N-CA	3.38	114.52	108.15
1	C	2[A]	HYP	CD-N-CA	3.38	114.52	108.15
1	C	2[F]	HYP	CD-N-CA	3.38	114.52	108.15
1	E	16[A]	HZP	CD-N-CA	3.37	114.52	108.15
1	C	2[E]	HYP	CD-N-CA	3.37	114.52	108.15
1	C	20[D]	HYP	OD1-CG-CB	-3.37	101.03	110.16
1	C	20[B]	HYP	OD1-CG-CB	-3.37	101.03	110.16
1	F	23[G]	HYP	CD-N-CA	3.37	114.51	108.15
1	G	8[D]	HYP	CD-N-CA	3.36	114.50	108.15
1	D	20[F]	HYP	CD-N-CA	3.36	114.49	108.15
1	H	8[G]	HYP	CD-N-CA	3.36	114.49	108.15
1	B	2[G]	HYP	CD-N-CA	3.36	114.49	108.15
1	H	8[F]	HYP	CD-N-CA	3.36	114.49	108.15
1	C	29[A]	HYP	OD1-CG-CB	-3.36	101.07	110.16
1	I	29[G]	HYP	CD-N-CA	3.35	114.48	108.15
1	C	29[F]	HYP	OD1-CG-CB	-3.35	101.08	110.16
1	H	11[D]	HYP	CD-N-CA	3.35	114.47	108.15
1	E	8[F]	HYP	CG-CD-N	-3.35	100.08	105.47
1	G	26[B]	HYP	CD-N-CA	3.35	114.47	108.15
1	F	13[G]	HZP	CG-CD-N	-3.34	100.08	105.47
1	G	26[C]	HYP	CD-N-CA	3.35	114.47	108.15
1	G	26[A]	HYP	CD-N-CA	3.35	114.47	108.15
1	E	23[G]	HYP	CD-N-CA	3.34	114.46	108.15
1	A	20[D]	HYP	CG-CD-N	-3.34	100.09	105.47
1	F	20[G]	HYP	CG-CD-N	-3.34	100.09	105.47
1	A	5[B]	HYP	CG-CD-N	-3.34	100.09	105.47
1	I	23[B]	HYP	CD-N-CA	3.34	114.46	108.15
1	E	26[B]	HYP	CD-N-CA	3.34	114.45	108.15
1	G	20[C]	HYP	CD-N-CA	3.34	114.45	108.15
1	G	11[D]	HYP	CD-N-CA	3.34	114.45	108.15
1	G	20[A]	HYP	CD-N-CA	3.34	114.45	108.15
1	G	20[B]	HYP	CD-N-CA	3.34	114.45	108.15
1	H	16[D]	HZP	CD-N-CA	3.33	114.44	108.15
1	H	16[C]	HZP	CD-N-CA	3.33	114.44	108.15
1	A	20[E]	HYP	CD-N-CA	3.33	114.44	108.15
1	E	11[A]	HYP	CG-CD-N	-3.33	100.10	105.47
1	E	11[E]	HYP	CG-CD-N	-3.33	100.10	105.47
1	A	20[G]	HYP	CD-N-CA	3.33	114.44	108.15
1	E	8[G]	HYP	CG-CD-N	-3.33	100.11	105.47
1	E	13[F]	HZP	CD-N-CA	3.32	114.42	108.15
1	E	16[E]	HZP	CD-N-CA	3.32	114.42	108.15
1	D	20[E]	HYP	CD-N-CA	3.32	114.42	108.15

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	13[G]	HZP	CD-N-CA	3.32	114.42	108.15
1	F	8[A]	HYP	CD-N-CA	3.32	114.42	108.15
1	F	8[E]	HYP	CD-N-CA	3.32	114.42	108.15
1	A	5[C]	HYP	CG-CD-N	-3.32	100.12	105.47
1	E	11[B]	HYP	CG-CD-N	-3.32	100.13	105.47
1	E	26[A]	HYP	CD-N-CA	3.31	114.41	108.15
1	G	16[D]	HZP	CD-N-CA	3.31	114.41	108.15
1	H	11[C]	HYP	CD-N-CA	3.31	114.41	108.15
1	C	29[E]	HYP	OD1-CG-CB	-3.31	101.19	110.16
1	B	8[G]	HYP	CD-N-CA	3.31	114.40	108.15
1	D	29[G]	HYP	CD-N-CA	3.31	114.40	108.15
1	F	8[F]	HYP	CD-N-CA	3.31	114.40	108.15
1	D	26[G]	HYP	OD1-CG-CB	-3.31	101.20	110.16
1	A	20[F]	HYP	CD-N-CA	3.31	114.39	108.15
1	H	29[F]	HYP	CD-N-CA	3.31	114.39	108.15
1	C	5[C]	HYP	CD-N-CA	3.31	114.39	108.15
1	F	5[G]	HYP	CD-N-CA	3.31	114.39	108.15
1	E	26[E]	HYP	CD-N-CA	3.31	114.39	108.15
1	C	5[B]	HYP	CD-N-CA	3.30	114.39	108.15
1	I	16[E]	HZP	C-CA-N	3.30	116.98	110.86
1	E	8[D]	HYP	CD-N-CA	3.30	114.39	108.15
1	E	20[C]	HYP	CG-CD-N	-3.30	100.15	105.47
1	E	23[F]	HYP	CD-N-CA	3.30	114.38	108.15
1	C	13[F]	HZP	OD1-CG-CB	-3.30	101.23	110.16
1	D	26[E]	HYP	OD1-CG-CB	-3.30	101.23	110.16
1	B	23[F]	HYP	CD-N-CA	3.29	114.37	108.15
1	B	23[G]	HYP	CD-N-CA	3.29	114.37	108.15
1	G	26[G]	HYP	CD-N-CA	3.29	114.37	108.15
1	C	5[D]	HYP	CD-N-CA	3.29	114.37	108.15
1	C	13[A]	HZP	OD1-CG-CB	-3.30	101.23	110.16
1	E	20[D]	HYP	CG-CD-N	-3.29	100.17	105.47
1	E	2[B]	HYP	CG-CD-N	-3.29	100.17	105.47
1	C	13[E]	HZP	OD1-CG-CB	-3.29	101.25	110.16
1	D	29[F]	HYP	CD-N-CA	3.29	114.36	108.15
1	E	2[A]	HYP	CG-CD-N	-3.29	100.17	105.47
1	I	2[B]	HYP	CD-N-CA	3.29	114.36	108.15
1	G	26[E]	HYP	CD-N-CA	3.29	114.36	108.15
1	B	8[F]	HYP	CD-N-CA	3.29	114.36	108.15
1	I	16[F]	HZP	C-CA-N	3.29	116.94	110.86
1	F	8[D]	HYP	CG-CD-N	-3.29	100.18	105.47
1	B	2[B]	HYP	CD-N-CA	3.29	114.35	108.15
1	D	26[F]	HYP	OD1-CG-CB	-3.28	101.27	110.16

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	23[F]	HYP	CD-N-CA	3.28	114.34	108.15
1	E	29[D]	HYP	CD-N-CA	3.28	114.34	108.15
1	B	11[A]	HYP	CD-N-CA	3.28	114.34	108.15
1	C	11[C]	HYP	CG-CD-N	-3.28	100.19	105.47
1	B	2[A]	HYP	CD-N-CA	3.28	114.34	108.15
1	C	11[B]	HYP	CG-CD-N	-3.28	100.19	105.47
1	B	11[E]	HYP	CD-N-CA	3.28	114.34	108.15
1	C	26[F]	HYP	CG-CD-N	-3.28	100.19	105.47
1	B	2[E]	HYP	CD-N-CA	3.28	114.34	108.15
1	A	5[A]	HYP	CG-CD-N	-3.28	100.19	105.47
1	G	2[G]	HYP	CD-N-CA	3.27	114.33	108.15
1	G	2[E]	HYP	CD-N-CA	3.27	114.33	108.15
1	G	2[F]	HYP	CD-N-CA	3.27	114.33	108.15
1	A	13[B]	HZP	CG-CD-N	-3.27	100.20	105.47
1	C	11[D]	HYP	CG-CD-N	-3.27	100.21	105.47
1	E	2[E]	HYP	CG-CD-N	-3.27	100.21	105.47
1	B	26[B]	HYP	CD-N-CA	3.26	114.31	108.15
1	C	16[C]	HZP	OD1-CG-CB	-3.26	101.32	110.16
1	C	23[A]	HYP	CD-N-CA	3.26	114.31	108.15
1	D	29[E]	HYP	CD-N-CA	3.26	114.31	108.15
1	I	16[A]	HZP	C-CA-N	3.26	116.90	110.86
1	D	11[E]	HYP	CD-N-CA	3.26	114.30	108.15
1	E	8[C]	HYP	CD-N-CA	3.26	114.30	108.15
1	H	29[G]	HYP	CD-N-CA	3.26	114.30	108.15
1	F	8[B]	HYP	CG-CD-N	-3.26	100.23	105.47
1	C	16[D]	HZP	OD1-CG-CB	-3.25	101.35	110.16
1	E	20[G]	HYP	CD-N-CA	3.25	114.29	108.15
1	E	29[C]	HYP	CD-N-CA	3.25	114.28	108.15
1	B	11[B]	HYP	CD-N-CA	3.25	114.28	108.15
1	C	16[B]	HZP	OD1-CG-CB	-3.24	101.38	110.16
1	C	23[E]	HYP	CD-N-CA	3.24	114.27	108.15
1	B	26[A]	HYP	CD-N-CA	3.24	114.27	108.15
1	C	26[B]	HYP	CD-N-CA	3.24	114.27	108.15
1	C	26[A]	HYP	CG-CD-N	-3.24	100.25	105.47
1	I	29[F]	HYP	CG-CD-N	-3.24	100.25	105.47
1	I	23[G]	HYP	CG-CD-N	-3.24	100.25	105.47
1	C	26[C]	HYP	CD-N-CA	3.23	114.25	108.15
1	A	13[C]	HZP	CG-CD-N	-3.23	100.26	105.47
1	C	26[D]	HYP	CD-N-CA	3.23	114.25	108.15
1	I	11[G]	HYP	CB-CG-CD	3.23	108.05	103.46
1	E	2[F]	HYP	CD-N-CA	3.23	114.24	108.15
1	F	8[C]	HYP	CG-CD-N	-3.22	100.28	105.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	2[G]	HYP	CD-N-CA	3.22	114.23	108.15
1	D	11[F]	HYP	CD-N-CA	3.22	114.23	108.15
1	B	26[E]	HYP	CD-N-CA	3.22	114.23	108.15
1	F	2[D]	HYP	OD1-CG-CB	-3.22	101.45	110.16
1	C	20[G]	HYP	CD-N-CA	3.21	114.22	108.15
1	G	11[G]	HYP	CD-N-CA	3.21	114.22	108.15
1	F	26[G]	HYP	CB-CG-CD	3.21	108.03	103.46
1	D	11[G]	HYP	CD-N-CA	3.21	114.21	108.15
1	D	11[B]	HYP	CD-N-CA	3.21	114.21	108.15
1	B	29[C]	HYP	CD-N-CA	3.21	114.20	108.15
1	F	2[B]	HYP	OD1-CG-CB	-3.21	101.47	110.16
1	E	20[F]	HYP	CD-N-CA	3.21	114.20	108.15
1	C	26[E]	HYP	CG-CD-N	-3.21	100.30	105.47
1	F	2[C]	HYP	OD1-CG-CB	-3.21	101.48	110.16
1	D	11[A]	HYP	CD-N-CA	3.21	114.20	108.15
1	G	5[B]	HYP	CG-CD-N	-3.20	100.31	105.47
1	A	23[B]	HYP	CD-N-CA	3.20	114.20	108.15
1	D	11[C]	HYP	CD-N-CA	3.20	114.19	108.15
1	B	29[D]	HYP	CD-N-CA	3.20	114.19	108.15
1	G	11[E]	HYP	CD-N-CA	3.20	114.19	108.15
1	A	23[C]	HYP	CD-N-CA	3.19	114.17	108.15
1	H	20[F]	HYP	CD-N-CA	3.19	114.17	108.15
1	H	20[G]	HYP	CD-N-CA	3.19	114.17	108.15
1	H	11[A]	HYP	CD-N-CA	3.18	114.16	108.15
1	G	5[A]	HYP	CG-CD-N	-3.18	100.34	105.47
1	D	2[C]	HYP	C-CA-N	3.19	116.76	110.86
1	A	16[E]	HZP	CG-CD-N	-3.18	100.34	105.47
1	G	11[F]	HYP	CD-N-CA	3.18	114.16	108.15
1	D	2[B]	HYP	C-CA-N	3.18	116.75	110.86
1	I	29[E]	HYP	CG-CD-N	-3.18	100.35	105.47
1	I	29[A]	HYP	CG-CD-N	-3.18	100.35	105.47
1	I	11[G]	HYP	CD-N-CA	3.18	114.15	108.15
1	B	20[D]	HYP	CB-CA-N	-3.18	99.72	104.59
1	A	13[A]	HZP	CG-CD-N	-3.17	100.36	105.47
1	A	23[A]	HYP	CD-N-CA	3.17	114.14	108.15
1	H	13[G]	HZP	CD-N-CA	3.17	114.14	108.15
1	A	26[D]	HYP	CD-N-CA	3.17	114.13	108.15
1	G	5[C]	HYP	CG-CD-N	-3.17	100.37	105.47
1	D	2[A]	HYP	C-CA-N	3.16	116.72	110.86
1	G	23[A]	HYP	CD-N-CA	3.16	114.12	108.15
1	G	13[E]	HZP	CG-CD-N	-3.16	100.38	105.47
1	G	13[G]	HZP	CG-CD-N	-3.16	100.38	105.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	13[A]	HZP	CD-N-CA	3.16	114.12	108.15
1	G	13[C]	HZP	CD-N-CA	3.16	114.12	108.15
1	G	13[F]	HZP	CG-CD-N	-3.16	100.38	105.47
1	H	16[E]	HZP	OD1-CG-CD	-3.16	102.89	110.31
1	I	13[G]	HZP	C-CA-N	3.16	116.70	110.86
1	H	16[B]	HZP	OD1-CG-CD	-3.16	102.89	110.31
1	H	16[A]	HZP	OD1-CG-CD	-3.16	102.89	110.31
1	F	5[C]	HYP	CG-CD-N	-3.16	100.39	105.47
1	B	20[C]	HYP	CB-CA-N	-3.15	99.75	104.59
1	H	13[F]	HZP	CD-N-CA	3.15	114.10	108.15
1	A	16[G]	HZP	CG-CD-N	-3.15	100.40	105.47
1	G	23[A]	HYP	CG-CD-N	-3.14	100.41	105.47
1	D	11[G]	HYP	CG-CD-N	-3.14	100.41	105.47
1	G	13[B]	HZP	CD-N-CA	3.14	114.09	108.15
1	F	11[D]	HYP	CD-N-CA	3.14	114.08	108.15
1	D	11[F]	HYP	CG-CD-N	-3.14	100.41	105.47
1	B	8[C]	HYP	CG-CD-N	-3.14	100.42	105.47
1	H	8[A]	HYP	CD-N-CA	3.14	114.08	108.15
1	G	23[C]	HYP	CD-N-CA	3.14	114.07	108.15
1	B	8[D]	HYP	CG-CD-N	-3.14	100.42	105.47
1	B	2[F]	HYP	CG-CD-N	-3.14	100.42	105.47
1	A	2[A]	HYP	CB-CG-CD	3.13	107.92	103.46
1	A	2[B]	HYP	CB-CG-CD	3.13	107.92	103.46
1	D	11[E]	HYP	CG-CD-N	-3.13	100.43	105.47
1	H	29[E]	HYP	CG-CD-N	-3.12	100.44	105.47
1	H	29[A]	HYP	CG-CD-N	-3.12	100.44	105.47
1	D	8[E]	HYP	CD-N-CA	3.12	114.05	108.15
1	H	29[B]	HYP	CG-CD-N	-3.12	100.44	105.47
1	G	16[F]	HZP	CG-CD-N	-3.12	100.44	105.47
1	C	29[G]	HYP	CG-CD-N	-3.12	100.45	105.47
1	H	8[B]	HYP	CD-N-CA	3.12	114.04	108.15
1	H	8[E]	HYP	CD-N-CA	3.12	114.03	108.15
1	G	11[G]	HYP	CG-CD-N	-3.12	100.45	105.47
1	B	2[G]	HYP	CG-CD-N	-3.12	100.45	105.47
1	G	26[D]	HYP	CG-CD-N	-3.11	100.46	105.47
1	F	5[B]	HYP	CG-CD-N	-3.11	100.47	105.47
1	F	11[B]	HYP	CD-N-CA	3.10	114.01	108.15
1	B	5[C]	HYP	CD-N-CA	3.10	114.01	108.15
1	D	8[F]	HYP	CD-N-CA	3.10	114.01	108.15
1	H	11[B]	HYP	CD-N-CA	3.10	114.01	108.15
1	H	11[E]	HYP	CD-N-CA	3.10	114.00	108.15
1	G	23[C]	HYP	CG-CD-N	-3.10	100.48	105.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	11[C]	HYP	CG-CD-N	-3.10	100.48	105.47
1	G	11[E]	HYP	CG-CD-N	-3.10	100.48	105.47
1	C	20[G]	HYP	CG-CD-N	-3.09	100.49	105.47
1	G	26[D]	HYP	CD-N-CA	3.09	113.98	108.15
1	G	11[F]	HYP	CG-CD-N	-3.09	100.50	105.47
1	F	8[G]	HYP	OD1-CG-CB	-3.09	101.80	110.16
1	A	5[D]	HYP	CB-CG-CD	3.09	107.85	103.46
1	D	8[G]	HYP	CD-N-CA	3.09	113.97	108.15
1	A	16[F]	HYP	CG-CD-N	-3.08	100.50	105.47
1	H	5[D]	HYP	CD-N-CA	3.08	113.97	108.15
1	H	5[C]	HYP	CD-N-CA	3.08	113.97	108.15
1	C	20[E]	HYP	CD-N-CA	3.08	113.97	108.15
1	C	23[F]	HYP	CG-CD-N	-3.08	100.51	105.47
1	H	5[G]	HYP	CD-N-CA	3.08	113.96	108.15
1	H	5[F]	HYP	CD-N-CA	3.08	113.96	108.15
1	B	26[F]	HYP	CD-N-CA	3.07	113.95	108.15
1	G	23[B]	HYP	CG-CD-N	-3.07	100.52	105.47
1	F	23[E]	HYP	CD-N-CA	3.07	113.95	108.15
1	F	11[C]	HYP	CD-N-CA	3.07	113.94	108.15
1	B	11[D]	HYP	CG-CD-N	-3.07	100.53	105.47
1	C	20[A]	HYP	CD-N-CA	3.07	113.94	108.15
1	I	29[B]	HYP	CD-N-CA	3.06	113.93	108.15
1	I	29[D]	HYP	CD-N-CA	3.06	113.93	108.15
1	I	29[C]	HYP	CD-N-CA	3.06	113.93	108.15
1	D	26[C]	HYP	CD-N-CA	3.06	113.93	108.15
1	D	26[B]	HYP	CD-N-CA	3.06	113.93	108.15
1	D	26[A]	HYP	CD-N-CA	3.06	113.93	108.15
1	G	16[E]	HYP	CG-CD-N	-3.06	100.54	105.47
1	B	11[C]	HYP	CD-N-CA	3.06	113.92	108.15
1	A	2[C]	HYP	CB-CG-CD	3.06	107.81	103.46
1	F	26[B]	HYP	CD-N-CA	3.06	113.93	108.15
1	F	26[D]	HYP	CD-N-CA	3.06	113.93	108.15
1	F	26[C]	HYP	CD-N-CA	3.06	113.93	108.15
1	B	26[G]	HYP	CD-N-CA	3.06	113.92	108.15
1	C	23[A]	HYP	CG-CD-N	-3.06	100.55	105.47
1	C	13[F]	HYP	CD-N-CA	3.06	113.92	108.15
1	F	11[E]	HYP	OD1-CG-CB	-3.05	101.89	110.16
1	H	2[D]	HYP	CG-CD-N	-3.05	100.56	105.47
1	H	2[C]	HYP	CG-CD-N	-3.05	100.56	105.47
1	F	11[A]	HYP	OD1-CG-CB	-3.05	101.90	110.16
1	F	11[F]	HYP	OD1-CG-CB	-3.05	101.91	110.16
1	F	5[D]	HYP	CG-CD-N	-3.05	100.57	105.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	23[E]	HYP	CD-N-CA	3.04	113.89	108.15
1	C	26[D]	HYP	CG-CD-N	-3.04	100.57	105.47
1	G	23[B]	HYP	CD-N-CA	3.04	113.89	108.15
1	E	16[G]	HZP	CD-N-CA	3.04	113.89	108.15
1	C	26[B]	HYP	CG-CD-N	-3.04	100.58	105.47
1	A	29[D]	HYP	CD-N-CA	3.04	113.88	108.15
1	E	23[E]	HYP	CD-N-CA	3.04	113.88	108.15
1	D	23[F]	HYP	CD-N-CA	3.03	113.88	108.15
1	I	23[C]	HYP	CG-CD-N	-3.03	100.58	105.47
1	I	23[D]	HYP	CG-CD-N	-3.03	100.58	105.47
1	F	23[A]	HYP	CD-N-CA	3.03	113.87	108.15
1	C	23[E]	HYP	CG-CD-N	-3.03	100.59	105.47
1	D	29[D]	HYP	CD-N-CA	3.03	113.87	108.15
1	C	20[F]	HYP	CD-N-CA	3.03	113.87	108.15
1	A	13[D]	HZP	CD-N-CA	3.03	113.86	108.15
1	B	11[D]	HYP	CD-N-CA	3.02	113.86	108.15
1	E	16[F]	HZP	CD-N-CA	3.02	113.86	108.15
1	C	26[C]	HYP	CG-CD-N	-3.02	100.61	105.47
1	I	23[B]	HYP	CG-CD-N	-3.01	100.62	105.47
1	D	23[G]	HYP	CD-N-CA	3.01	113.83	108.15
1	C	2[E]	HYP	CG-CD-N	-3.01	100.62	105.47
1	C	2[A]	HYP	CG-CD-N	-3.01	100.62	105.47
1	C	13[A]	HZP	CD-N-CA	3.01	113.83	108.15
1	C	2[F]	HYP	CG-CD-N	-3.01	100.62	105.47
1	C	26[G]	HYP	OD1-CG-CB	-3.01	102.02	110.16
1	E	26[C]	HYP	CD-N-CA	3.01	113.83	108.15
1	C	23[C]	HYP	CD-N-CA	3.01	113.83	108.15
1	B	23[F]	HYP	CG-CD-N	-3.00	100.63	105.47
1	B	23[G]	HYP	CG-CD-N	-3.00	100.63	105.47
1	C	8[G]	HYP	CG-CD-N	-3.00	100.64	105.47
1	E	23[A]	HYP	CD-N-CA	3.00	113.81	108.15
1	G	16[G]	HZP	CG-CD-N	-3.00	100.64	105.47
1	B	2[B]	HYP	CG-CD-N	-3.00	100.64	105.47
1	C	13[E]	HZP	CD-N-CA	3.00	113.81	108.15
1	A	16[E]	HZP	CD-N-CA	2.99	113.80	108.15
1	E	26[D]	HYP	CD-N-CA	3.00	113.80	108.15
1	F	23[F]	HYP	CD-N-CA	2.99	113.80	108.15
1	A	26[C]	HYP	CD-N-CA	2.99	113.80	108.15
1	C	23[D]	HYP	CD-N-CA	2.99	113.79	108.15
1	A	26[B]	HYP	CD-N-CA	2.99	113.79	108.15
1	D	29[C]	HYP	CG-CD-N	-2.99	100.66	105.47
1	C	23[B]	HYP	CD-N-CA	2.98	113.78	108.15

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	26[A]	HYP	CD-N-CA	2.98	113.78	108.15
1	D	29[A]	HYP	CG-CD-N	-2.98	100.67	105.47
1	A	16[G]	HZP	CD-N-CA	2.98	113.77	108.15
1	A	16[F]	HZP	CD-N-CA	2.97	113.77	108.15
1	D	20[D]	HYP	OD1-CG-CB	-2.97	102.11	110.16
1	C	20[C]	HYP	CG-CD-N	-2.97	100.68	105.47
1	C	20[D]	HYP	CG-CD-N	-2.97	100.68	105.47
1	C	20[B]	HYP	CG-CD-N	-2.97	100.68	105.47
1	C	5[C]	HYP	CG-CD-N	-2.97	100.68	105.47
1	A	23[F]	HYP	CD-N-CA	2.97	113.76	108.15
1	D	29[B]	HYP	CG-CD-N	-2.97	100.69	105.47
1	G	2[D]	HYP	CG-CD-N	-2.97	100.69	105.47
1	I	11[G]	HYP	OD1-CG-CB	-2.97	102.12	110.16
1	G	13[C]	HZP	CG-CD-N	-2.96	100.70	105.47
1	C	5[D]	HYP	CG-CD-N	-2.96	100.70	105.47
1	D	2[C]	HYP	CG-CD-N	-2.96	100.70	105.47
1	E	23[B]	HYP	CD-N-CA	2.96	113.74	108.15
1	C	5[B]	HYP	CG-CD-N	-2.96	100.70	105.47
1	B	2[A]	HYP	CG-CD-N	-2.96	100.71	105.47
1	A	20[D]	HYP	CB-CA-N	-2.95	100.06	104.59
1	D	2[A]	HYP	CG-CD-N	-2.94	100.74	105.47
1	A	23[G]	HYP	CD-N-CA	2.93	113.68	108.15
1	G	13[A]	HZP	CG-CD-N	-2.93	100.75	105.47
1	A	23[E]	HYP	CD-N-CA	2.92	113.67	108.15
1	F	11[G]	HYP	OD1-CG-CD	2.92	117.18	110.31
1	C	16[D]	HZP	CD-N-CA	2.92	113.67	108.15
1	D	2[B]	HYP	CG-CD-N	-2.92	100.77	105.47
1	E	11[F]	HYP	CG-CB-CA	-2.92	99.96	104.22
1	C	16[B]	HZP	CD-N-CA	2.92	113.66	108.15
1	C	16[C]	HZP	CD-N-CA	2.92	113.66	108.15
1	D	20[D]	HYP	CD-N-CA	2.92	113.66	108.15
1	F	26[D]	HYP	CG-CD-N	-2.92	100.77	105.47
1	E	16[G]	HZP	CB-CA-N	-2.92	100.12	104.59
1	B	2[E]	HYP	CG-CD-N	-2.91	100.78	105.47
1	I	5[C]	HYP	CB-CG-CD	2.91	107.60	103.46
1	B	16[D]	HZP	OD1-CG-CB	2.91	118.05	110.16
1	F	23[D]	HYP	CG-CD-N	-2.91	100.79	105.47
1	B	16[C]	HZP	OD1-CG-CB	2.91	118.04	110.16
1	F	29[G]	HYP	CD-N-CA	2.91	113.63	108.15
1	C	8[A]	HYP	CD-N-CA	2.90	113.63	108.15
1	I	29[G]	HYP	CG-CD-N	-2.90	100.80	105.47
1	I	5[B]	HYP	CB-CG-CD	2.90	107.58	103.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	8[E]	HYP	CD-N-CA	2.90	113.62	108.15
1	C	29[G]	HYP	CD-N-CA	2.90	113.62	108.15
1	I	5[D]	HYP	CB-CG-CD	2.90	107.58	103.46
1	F	23[B]	HYP	CG-CD-N	-2.90	100.80	105.47
1	H	20[E]	HYP	CG-CD-N	-2.89	100.81	105.47
1	G	13[B]	HYP	CG-CD-N	-2.89	100.81	105.47
1	H	20[A]	HYP	CG-CD-N	-2.89	100.81	105.47
1	C	8[F]	HYP	CD-N-CA	2.89	113.61	108.15
1	H	20[B]	HYP	CG-CD-N	-2.89	100.82	105.47
1	E	11[G]	HYP	CG-CB-CA	-2.89	100.01	104.22
1	H	16[B]	HYP	CD-N-CA	2.89	113.60	108.15
1	I	29[D]	HYP	CG-CD-N	-2.89	100.82	105.47
1	I	29[C]	HYP	CG-CD-N	-2.89	100.82	105.47
1	E	16[F]	HYP	CB-CA-N	-2.88	100.17	104.59
1	A	11[C]	HYP	CD-N-CA	2.88	113.59	108.15
1	A	11[B]	HYP	CD-N-CA	2.87	113.58	108.15
1	F	26[B]	HYP	CG-CD-N	-2.87	100.84	105.47
1	F	23[C]	HYP	CG-CD-N	-2.87	100.85	105.47
1	A	11[A]	HYP	CD-N-CA	2.87	113.57	108.15
1	C	5[G]	HYP	CD-N-CA	2.87	113.57	108.15
1	H	26[F]	HYP	CG-CD-N	-2.87	100.85	105.47
1	H	29[D]	HYP	CG-CD-N	-2.87	100.85	105.47
1	H	29[C]	HYP	CG-CD-N	-2.87	100.85	105.47
1	H	16[E]	HYP	CD-N-CA	2.87	113.56	108.15
1	E	5[D]	HYP	OD1-CG-CB	-2.87	102.40	110.16
1	E	5[C]	HYP	OD1-CG-CB	-2.87	102.40	110.16
1	H	26[G]	HYP	CG-CD-N	-2.87	100.85	105.47
1	H	16[A]	HYP	CD-N-CA	2.87	113.56	108.15
1	I	16[G]	HYP	CG-CD-N	-2.86	100.86	105.47
1	B	29[G]	HYP	OD1-CG-CB	-2.86	102.41	110.16
1	B	29[F]	HYP	OD1-CG-CB	-2.86	102.41	110.16
1	A	8[F]	HYP	CD-N-CA	2.86	113.55	108.15
1	F	20[F]	HYP	CG-CD-N	-2.86	100.87	105.47
1	F	20[A]	HYP	CG-CD-N	-2.86	100.87	105.47
1	F	20[E]	HYP	CG-CD-N	-2.86	100.87	105.47
1	C	11[G]	HYP	OD1-CG-CB	-2.85	102.43	110.16
1	I	29[B]	HYP	CG-CD-N	-2.85	100.88	105.47
1	A	26[E]	HYP	CG-CD-N	-2.85	100.88	105.47
1	G	16[D]	HYP	CG-CD-N	-2.85	100.88	105.47
1	B	16[D]	HYP	C-CA-N	2.85	116.13	110.86
1	A	8[G]	HYP	CD-N-CA	2.85	113.53	108.15
1	G	13[D]	HYP	CB-CA-N	-2.85	100.22	104.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	23[A]	HYP	CG-CD-N	-2.85	100.89	105.47
1	A	8[E]	HYP	CD-N-CA	2.84	113.52	108.15
1	D	20[F]	HYP	CG-CD-N	-2.84	100.89	105.47
1	F	2[G]	HYP	CG-CD-N	-2.84	100.89	105.47
1	D	20[G]	HYP	CG-CD-N	-2.84	100.90	105.47
1	I	20[F]	HYP	CG-CD-N	-2.84	100.90	105.47
1	B	26[G]	HYP	OD1-CG-CB	-2.83	102.49	110.16
1	B	26[F]	HYP	OD1-CG-CB	-2.83	102.49	110.16
1	I	26[G]	HYP	CG-CD-N	-2.83	100.91	105.47
1	F	26[C]	HYP	CG-CD-N	-2.83	100.91	105.47
1	F	5[F]	HYP	CG-CD-N	-2.83	100.92	105.47
1	D	26[D]	HYP	CG-CD-N	-2.82	100.92	105.47
1	F	5[A]	HYP	CG-CD-N	-2.82	100.92	105.47
1	H	8[D]	HYP	CD-N-CA	2.82	113.47	108.15
1	B	11[F]	HYP	CD-N-CA	2.82	113.47	108.15
1	I	23[E]	HYP	CD-N-CA	2.82	113.47	108.15
1	I	20[A]	HYP	CG-CD-N	-2.82	100.93	105.47
1	D	20[D]	HYP	CG-CD-N	-2.82	100.94	105.47
1	I	20[E]	HYP	CG-CD-N	-2.81	100.94	105.47
1	E	20[C]	HYP	OD1-CG-CB	-2.81	102.54	110.16
1	H	8[C]	HYP	CD-N-CA	2.81	113.46	108.15
1	C	11[E]	HYP	CG-CD-N	-2.81	100.94	105.47
1	D	16[B]	HZP	CD-N-CA	2.81	113.45	108.15
1	D	23[B]	HYP	CG-CD-N	-2.81	100.94	105.47
1	B	16[C]	HZP	C-CA-N	2.81	116.06	110.86
1	D	2[D]	HYP	CG-CD-N	-2.81	100.95	105.47
1	B	11[G]	HYP	CD-N-CA	2.81	113.45	108.15
1	D	23[E]	HYP	CG-CD-N	-2.81	100.95	105.47
1	D	23[F]	HYP	CG-CD-N	-2.81	100.95	105.47
1	D	23[G]	HYP	CG-CD-N	-2.81	100.95	105.47
1	F	5[E]	HYP	CG-CD-N	-2.80	100.96	105.47
1	A	26[G]	HYP	CG-CD-N	-2.80	100.96	105.47
1	D	23[C]	HYP	CG-CD-N	-2.80	100.96	105.47
1	H	11[F]	HYP	OD1-CG-CB	-2.80	102.58	110.16
1	D	8[B]	HYP	CD-N-CA	2.80	113.43	108.15
1	I	23[A]	HYP	CD-N-CA	2.80	113.43	108.15
1	H	11[G]	HYP	OD1-CG-CB	-2.80	102.58	110.16
1	F	11[G]	HYP	C-CA-N	2.79	116.03	110.86
1	E	20[D]	HYP	OD1-CG-CB	-2.79	102.60	110.16
1	D	20[D]	HYP	C-CA-N	2.79	116.03	110.86
1	F	2[F]	HYP	CG-CD-N	-2.79	100.98	105.47
1	I	23[F]	HYP	CD-N-CA	2.79	113.42	108.15

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	16[A]	HZP	CD-N-CA	2.79	113.41	108.15
1	I	26[B]	HYP	CD-N-CA	2.78	113.40	108.15
1	C	11[A]	HYP	CG-CD-N	-2.78	100.99	105.47
1	C	23[G]	HYP	OD1-CG-CB	-2.78	102.64	110.16
1	B	5[B]	HYP	CG-CD-N	-2.78	101.00	105.47
1	F	2[F]	HYP	CD-N-CA	2.78	113.39	108.15
1	B	16[D]	HZP	CB-CG-CD	2.78	107.41	103.46
1	H	23[C]	HYP	CG-CD-N	-2.78	101.00	105.47
1	H	23[D]	HYP	CG-CD-N	-2.78	101.00	105.47
1	H	5[E]	HYP	CD-N-CA	2.77	113.39	108.15
1	D	20[E]	HYP	CG-CD-N	-2.77	101.00	105.47
1	H	5[A]	HYP	CD-N-CA	2.77	113.39	108.15
1	H	5[B]	HYP	CD-N-CA	2.77	113.39	108.15
1	E	2[E]	HYP	OD1-CG-CB	-2.77	102.66	110.16
1	G	8[F]	HYP	CG-CD-N	-2.77	101.01	105.47
1	G	8[E]	HYP	CG-CD-N	-2.77	101.01	105.47
1	G	8[G]	HYP	CG-CD-N	-2.77	101.01	105.47
1	I	2[F]	HYP	CD-N-CA	2.76	113.37	108.15
1	C	11[F]	HYP	CG-CD-N	-2.76	101.02	105.47
1	B	23[D]	HYP	OD1-CG-CB	-2.76	102.67	110.16
1	D	8[A]	HYP	CD-N-CA	2.76	113.37	108.15
1	I	2[E]	HYP	CD-N-CA	2.76	113.36	108.15
1	C	20[D]	HYP	OD1-CG-CD	2.76	116.80	110.31
1	C	20[B]	HYP	OD1-CG-CD	2.76	116.80	110.31
1	D	16[C]	HZP	CD-N-CA	2.76	113.36	108.15
1	F	23[C]	HYP	CD-N-CA	2.76	113.36	108.15
1	F	23[B]	HYP	CD-N-CA	2.76	113.36	108.15
1	I	5[E]	HYP	CG-CD-N	-2.76	101.03	105.47
1	I	8[C]	HYP	CG-CD-N	-2.75	101.04	105.47
1	E	2[A]	HYP	OD1-CG-CB	-2.75	102.71	110.16
1	I	2[A]	HYP	CD-N-CA	2.75	113.35	108.15
1	F	2[A]	HYP	CD-N-CA	2.75	113.34	108.15
1	F	20[G]	HYP	CD-N-CA	2.75	113.34	108.15
1	F	2[A]	HYP	CG-CD-N	-2.75	101.05	105.47
1	C	13[G]	HZP	CD-N-CA	2.75	113.34	108.15
1	B	5[A]	HYP	CG-CD-N	-2.75	101.05	105.47
1	A	8[B]	HYP	CG-CD-N	-2.75	101.04	105.47
1	F	5[C]	HYP	CD-N-CA	2.75	113.34	108.15
1	D	16[C]	HZP	CG-CB-CA	-2.75	100.22	104.22
1	B	16[C]	HZP	CB-CG-CD	2.75	107.37	103.46
1	I	26[C]	HYP	CD-N-CA	2.75	113.33	108.15
1	D	8[C]	HYP	CD-N-CA	2.74	113.33	108.15

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	23[C]	HYP	OD1-CG-CB	-2.74	102.73	110.16
1	B	20[E]	HYP	OD1-CG-CB	-2.74	102.73	110.16
1	A	26[F]	HYP	CG-CD-N	-2.74	101.05	105.47
1	C	20[C]	HYP	OD1-CG-CD	2.74	116.75	110.31
1	I	5[F]	HYP	CG-CD-N	-2.74	101.06	105.47
1	B	29[B]	HYP	OD1-CG-CB	-2.74	102.74	110.16
1	H	20[F]	HYP	CG-CD-N	-2.74	101.06	105.47
1	B	20[E]	HYP	CG-CD-N	-2.74	101.06	105.47
1	E	2[B]	HYP	OD1-CG-CB	-2.74	102.74	110.16
1	H	20[G]	HYP	CG-CD-N	-2.74	101.06	105.47
1	D	16[A]	HZP	CG-CB-CA	-2.74	100.23	104.22
1	F	23[D]	HYP	CD-N-CA	2.74	113.32	108.15
1	H	16[D]	HZP	CG-CD-N	-2.74	101.06	105.47
1	H	16[C]	HZP	CG-CD-N	-2.74	101.06	105.47
1	B	29[A]	HYP	OD1-CG-CB	-2.73	102.75	110.16
1	H	5[D]	HYP	OD1-CG-CB	-2.73	102.75	110.16
1	B	5[E]	HYP	CG-CD-N	-2.73	101.07	105.47
1	H	13[B]	HZP	CG-CD-N	-2.73	101.07	105.47
1	D	11[D]	HYP	CD-N-CA	2.73	113.31	108.15
1	I	8[B]	HYP	CG-CD-N	-2.73	101.08	105.47
1	G	11[C]	HYP	CG-CD-N	-2.73	101.08	105.47
1	B	29[E]	HYP	OD1-CG-CB	-2.73	102.77	110.16
1	D	26[F]	HYP	CG-CD-N	-2.73	101.08	105.47
1	F	2[D]	HYP	C-CA-N	2.73	115.91	110.86
1	G	26[F]	HYP	CG-CD-N	-2.73	101.08	105.47
1	D	16[B]	HZP	CG-CB-CA	-2.73	100.25	104.22
1	I	5[A]	HYP	CG-CD-N	-2.73	101.08	105.47
1	H	8[E]	HYP	CG-CD-N	-2.72	101.08	105.47
1	F	2[E]	HYP	CD-N-CA	2.72	113.29	108.15
1	A	8[A]	HYP	CG-CD-N	-2.72	101.08	105.47
1	B	20[B]	HYP	CG-CD-N	-2.72	101.09	105.47
1	B	20[A]	HYP	OD1-CG-CB	-2.72	102.80	110.16
1	B	20[A]	HYP	CG-CD-N	-2.72	101.09	105.47
1	F	20[F]	HYP	CD-N-CA	2.72	113.28	108.15
1	H	5[C]	HYP	OD1-CG-CB	-2.72	102.79	110.16
1	H	2[G]	HYP	CD-N-CA	2.72	113.28	108.15
1	H	2[F]	HYP	CD-N-CA	2.72	113.28	108.15
1	A	8[C]	HYP	CG-CD-N	-2.72	101.09	105.47
1	D	26[C]	HYP	CG-CD-N	-2.72	101.09	105.47
1	D	26[B]	HYP	CG-CD-N	-2.72	101.09	105.47
1	D	26[A]	HYP	CG-CD-N	-2.72	101.09	105.47
1	A	2[E]	HYP	CG-CD-N	-2.72	101.09	105.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	5[B]	HYP	CD-N-CA	2.72	113.28	108.15
1	D	5[G]	HYP	CD-N-CA	2.71	113.27	108.15
1	D	29[D]	HYP	CG-CD-N	-2.71	101.10	105.47
1	F	20[A]	HYP	CD-N-CA	2.71	113.27	108.15
1	D	5[F]	HYP	CD-N-CA	2.71	113.27	108.15
1	G	13[D]	HZP	C-CA-N	2.71	115.88	110.86
1	I	26[D]	HYP	CD-N-CA	2.71	113.26	108.15
1	F	20[E]	HYP	CD-N-CA	2.71	113.26	108.15
1	A	5[E]	HYP	CG-CD-N	-2.71	101.11	105.47
1	F	2[E]	HYP	CG-CD-N	-2.71	101.11	105.47
1	I	8[D]	HYP	CG-CD-N	-2.71	101.11	105.47
1	D	26[E]	HYP	CG-CD-N	-2.71	101.11	105.47
1	B	20[D]	HYP	CB-CG-CD	-2.70	99.62	103.46
1	G	11[B]	HYP	CG-CD-N	-2.70	101.12	105.47
1	F	2[B]	HYP	C-CA-N	2.70	115.86	110.86
1	I	11[E]	HYP	CG-CD-N	-2.70	101.12	105.47
1	A	2[F]	HYP	CG-CD-N	-2.70	101.12	105.47
1	G	11[A]	HYP	CG-CD-N	-2.70	101.13	105.47
1	B	20[B]	HYP	OD1-CG-CB	-2.70	102.86	110.16
1	H	5[G]	HYP	CG-CD-N	-2.70	101.13	105.47
1	H	5[F]	HYP	CG-CD-N	-2.70	101.13	105.47
1	F	5[D]	HYP	CD-N-CA	2.69	113.23	108.15
1	D	5[E]	HYP	CD-N-CA	2.69	113.23	108.15
1	H	13[E]	HZP	CG-CD-N	-2.69	101.14	105.47
1	D	20[C]	HYP	OD1-CG-CB	-2.69	102.88	110.16
1	H	13[A]	HZP	CG-CD-N	-2.69	101.14	105.47
1	A	8[D]	HYP	CG-CD-N	-2.69	101.14	105.47
1	I	26[G]	HYP	OD1-CG-CB	-2.69	102.88	110.16
1	D	2[E]	HYP	OD1-CG-CB	-2.69	102.88	110.16
1	I	11[F]	HYP	CG-CD-N	-2.69	101.14	105.47
1	D	26[G]	HYP	CG-CD-N	-2.69	101.14	105.47
1	F	29[A]	HYP	CG-CD-N	-2.69	101.15	105.47
1	D	2[F]	HYP	OD1-CG-CB	-2.68	102.89	110.16
1	H	8[A]	HYP	CG-CD-N	-2.68	101.15	105.47
1	F	2[C]	HYP	C-CA-N	2.68	115.82	110.86
1	I	20[G]	HYP	CD-N-CA	2.68	113.21	108.15
1	I	11[A]	HYP	CG-CD-N	-2.68	101.16	105.47
1	A	2[G]	HYP	CG-CD-N	-2.68	101.16	105.47
1	A	5[F]	HYP	CG-CD-N	-2.68	101.16	105.47
1	H	11[G]	HYP	CB-CG-CD	2.67	107.26	103.46
1	I	13[F]	HZP	CG-CD-N	-2.67	101.17	105.47
1	A	5[G]	HYP	CG-CD-N	-2.67	101.17	105.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	29[F]	HYP	CG-CD-N	-2.67	101.17	105.47
1	D	20[B]	HYP	OD1-CG-CB	-2.67	102.93	110.16
1	D	20[A]	HYP	OD1-CG-CB	-2.67	102.93	110.16
1	H	11[F]	HYP	CB-CG-CD	2.67	107.25	103.46
1	F	20[D]	HYP	CG-CD-N	-2.66	101.18	105.47
1	H	23[F]	HYP	CD-N-CA	2.66	113.18	108.15
1	H	23[G]	HYP	CD-N-CA	2.66	113.18	108.15
1	H	8[B]	HYP	CG-CD-N	-2.66	101.18	105.47
1	B	26[E]	HYP	CG-CD-N	-2.66	101.19	105.47
1	B	23[D]	HYP	CG-CD-N	-2.66	101.19	105.47
1	D	5[C]	HYP	OD1-CG-CB	-2.66	102.97	110.16
1	B	23[C]	HYP	CG-CD-N	-2.66	101.19	105.47
1	B	20[C]	HYP	CB-CG-CD	-2.65	99.69	103.46
1	I	2[G]	HYP	CG-CD-N	-2.65	101.20	105.47
1	F	20[B]	HYP	CG-CD-N	-2.65	101.20	105.47
1	D	2[G]	HYP	OD1-CG-CB	-2.65	102.98	110.16
1	F	11[B]	HYP	CG-CD-N	-2.64	101.21	105.47
1	D	5[A]	HYP	OD1-CG-CB	-2.65	103.00	110.16
1	F	11[D]	HYP	CG-CD-N	-2.64	101.21	105.47
1	F	11[C]	HYP	CG-CD-N	-2.64	101.21	105.47
1	B	26[A]	HYP	CG-CD-N	-2.64	101.22	105.47
1	I	13[A]	HZP	CG-CD-N	-2.64	101.22	105.47
1	F	29[E]	HYP	CG-CD-N	-2.63	101.23	105.47
1	C	29[E]	HYP	C-CA-N	2.64	115.74	110.86
1	E	5[E]	HYP	CG-CD-N	-2.63	101.23	105.47
1	I	8[F]	HYP	CG-CD-N	-2.63	101.23	105.47
1	D	5[B]	HYP	OD1-CG-CB	-2.63	103.03	110.16
1	F	29[G]	HYP	CG-CD-N	-2.63	101.24	105.47
1	I	13[E]	HZP	CG-CD-N	-2.63	101.24	105.47
1	D	8[D]	HYP	OD1-CG-CB	-2.63	103.05	110.16
1	D	26[F]	HYP	CD-N-CA	2.62	113.10	108.15
1	B	11[G]	HYP	CG-CD-N	-2.62	101.24	105.47
1	B	26[B]	HYP	CG-CD-N	-2.63	101.24	105.47
1	A	2[B]	HYP	OD1-CG-CB	-2.62	103.06	110.16
1	F	23[F]	HYP	CG-CD-N	-2.62	101.25	105.47
1	I	8[A]	HYP	CG-CD-N	-2.62	101.25	105.47
1	A	2[A]	HYP	OD1-CG-CB	-2.62	103.07	110.16
1	F	23[A]	HYP	CG-CD-N	-2.62	101.25	105.47
1	I	11[B]	HYP	CG-CD-N	-2.62	101.25	105.47
1	I	2[F]	HYP	OD1-CG-CB	-2.62	103.07	110.16
1	C	29[A]	HYP	C-CA-N	2.62	115.70	110.86
1	I	11[D]	HYP	CG-CD-N	-2.62	101.26	105.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	23[E]	HYP	CG-CD-N	-2.62	101.26	105.47
1	I	8[E]	HYP	CG-CD-N	-2.61	101.26	105.47
1	I	11[C]	HYP	CG-CD-N	-2.61	101.26	105.47
1	I	2[E]	HYP	OD1-CG-CB	-2.61	103.08	110.16
1	G	8[C]	HYP	CG-CD-N	-2.61	101.26	105.47
1	B	13[F]	HZP	CG-CD-N	-2.61	101.26	105.47
1	A	5[D]	HYP	OD1-CG-CB	-2.61	103.09	110.16
1	F	20[C]	HYP	CG-CD-N	-2.61	101.27	105.47
1	C	29[F]	HYP	C-CA-N	2.60	115.68	110.86
1	C	29[E]	HYP	CG-CD-N	-2.60	101.28	105.47
1	D	11[C]	HYP	CG-CD-N	-2.60	101.28	105.47
1	B	11[F]	HYP	CG-CD-N	-2.60	101.28	105.47
1	D	23[D]	HYP	OD1-CG-CB	-2.60	103.12	110.16
1	C	29[A]	HYP	CG-CD-N	-2.60	101.28	105.47
1	D	26[E]	HYP	CD-N-CA	2.60	113.06	108.15
1	C	11[E]	HYP	OD1-CG-CB	-2.60	103.12	110.16
1	C	11[A]	HYP	OD1-CG-CB	-2.60	103.12	110.16
1	F	29[G]	HYP	OD1-CG-CB	-2.60	103.12	110.16
1	C	11[F]	HYP	OD1-CG-CB	-2.60	103.12	110.16
1	B	13[G]	HZP	CG-CD-N	-2.60	101.28	105.47
1	D	26[G]	HYP	CD-N-CA	2.60	113.05	108.15
1	C	29[F]	HYP	CG-CD-N	-2.60	101.29	105.47
1	D	29[C]	HYP	CD-N-CA	2.59	113.05	108.15
1	D	29[A]	HYP	CD-N-CA	2.59	113.04	108.15
1	A	2[C]	HYP	OD1-CG-CB	-2.59	103.14	110.16
1	D	29[B]	HYP	CD-N-CA	2.59	113.04	108.15
1	G	2[D]	HYP	CD-N-CA	2.59	113.04	108.15
1	G	8[A]	HYP	CG-CD-N	-2.59	101.29	105.47
1	I	2[A]	HYP	OD1-CG-CB	-2.59	103.14	110.16
1	I	5[G]	HYP	CG-CD-N	-2.59	101.31	105.47
1	D	11[A]	HYP	CG-CD-N	-2.58	101.31	105.47
1	G	11[G]	HYP	CB-CA-N	-2.58	100.63	104.59
1	B	2[D]	HYP	CB-CG-CD	2.58	107.13	103.46
1	B	2[C]	HYP	CB-CG-CD	2.58	107.13	103.46
1	E	29[F]	HYP	OD1-CG-CB	-2.58	103.18	110.16
1	E	8[B]	HYP	CG-CD-N	-2.57	101.32	105.47
1	E	29[G]	HYP	OD1-CG-CB	-2.57	103.19	110.16
1	D	11[B]	HYP	CG-CD-N	-2.57	101.33	105.47
1	E	5[B]	HYP	CG-CD-N	-2.56	101.35	105.47
1	E	2[G]	HYP	CG-CD-N	-2.55	101.36	105.47
1	G	8[B]	HYP	CG-CD-N	-2.55	101.36	105.47
1	G	11[E]	HYP	CB-CA-N	-2.55	100.68	104.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	I	2[F]	HYP	CG-CD-N	-2.55	101.36	105.47
1	I	2[E]	HYP	CG-CD-N	-2.55	101.37	105.47
1	C	26[A]	HYP	OD1-CG-CB	-2.55	103.27	110.16
1	H	26[E]	HYP	CG-CD-N	-2.54	101.37	105.47
1	D	29[C]	HYP	OD1-CG-CB	-2.54	103.28	110.16
1	E	26[F]	HYP	CD-N-CA	2.54	112.94	108.15
1	E	26[G]	HYP	CD-N-CA	2.54	112.94	108.15
1	G	5[G]	HYP	CG-CD-N	-2.53	101.39	105.47
1	G	5[E]	HYP	CG-CD-N	-2.53	101.39	105.47
1	A	11[D]	HYP	CG-CD-N	-2.53	101.39	105.47
1	G	5[F]	HYP	CG-CD-N	-2.53	101.39	105.47
1	C	13[B]	HZP	CG-CD-N	-2.53	101.39	105.47
1	D	29[A]	HYP	OD1-CG-CB	-2.53	103.30	110.16
1	B	29[B]	HYP	CG-CD-N	-2.53	101.39	105.47
1	B	11[G]	HYP	CB-CG-CD	2.53	107.06	103.46
1	E	2[C]	HYP	OD1-CG-CB	-2.53	103.32	110.16
1	E	2[F]	HYP	CG-CD-N	-2.53	101.40	105.47
1	E	2[D]	HYP	OD1-CG-CB	-2.53	103.32	110.16
1	I	2[A]	HYP	CG-CD-N	-2.53	101.40	105.47
1	C	26[E]	HYP	OD1-CG-CB	-2.53	103.32	110.16
1	G	20[B]	HYP	OD1-CG-CB	-2.53	103.32	110.16
1	C	26[F]	HYP	OD1-CG-CB	-2.52	103.33	110.16
1	G	23[D]	HYP	OD1-CG-CB	-2.52	103.34	110.16
1	G	11[F]	HYP	CB-CA-N	-2.52	100.73	104.59
1	I	11[G]	HYP	CG-CD-N	-2.52	101.42	105.47
1	D	29[B]	HYP	OD1-CG-CB	-2.52	103.34	110.16
1	D	13[B]	HZP	CD-N-CA	2.52	112.90	108.15
1	D	13[A]	HZP	CD-N-CA	2.52	112.90	108.15
1	G	2[D]	HYP	OD1-CG-CB	-2.52	103.34	110.16
1	D	13[C]	HZP	CD-N-CA	2.52	112.90	108.15
1	E	5[F]	HYP	CG-CD-N	-2.52	101.42	105.47
1	H	26[B]	HYP	CG-CD-N	-2.51	101.42	105.47
1	E	5[G]	HYP	CG-CD-N	-2.51	101.42	105.47
1	E	8[E]	HYP	CG-CD-N	-2.51	101.43	105.47
1	H	11[A]	HYP	CG-CD-N	-2.50	101.44	105.47
1	B	29[C]	HYP	CG-CD-N	-2.50	101.44	105.47
1	G	29[C]	HYP	CG-CD-N	-2.50	101.44	105.47
1	B	29[A]	HYP	CG-CD-N	-2.50	101.44	105.47
1	C	13[D]	HZP	CG-CD-N	-2.50	101.45	105.47
1	C	11[G]	HYP	C-CA-N	-2.50	106.25	110.86
1	I	5[C]	HYP	CG-CD-N	-2.49	101.46	105.47
1	G	20[C]	HYP	OD1-CG-CD	2.49	116.17	110.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	20[A]	HYP	OD1-CG-CB	-2.49	103.41	110.16
1	H	26[A]	HYP	CG-CD-N	-2.49	101.46	105.47
1	I	5[D]	HYP	CG-CD-N	-2.49	101.46	105.47
1	I	5[B]	HYP	CG-CD-N	-2.49	101.46	105.47
1	B	29[D]	HYP	CG-CD-N	-2.49	101.46	105.47
1	E	8[A]	HYP	CG-CD-N	-2.49	101.46	105.47
1	C	13[C]	HYP	CG-CD-N	-2.49	101.47	105.47
1	B	11[F]	HYP	CB-CG-CD	2.48	106.99	103.46
1	H	29[D]	HYP	OD1-CG-CB	-2.48	103.45	110.16
1	H	11[B]	HYP	CG-CD-N	-2.48	101.48	105.47
1	H	29[C]	HYP	OD1-CG-CB	-2.48	103.45	110.16
1	D	29[G]	HYP	CB-CG-CD	2.48	106.98	103.46
1	B	11[F]	HYP	OD1-CG-CD	2.48	116.13	110.31
1	H	26[B]	HYP	CD-N-CA	2.47	112.81	108.15
1	B	29[E]	HYP	CG-CD-N	-2.47	101.49	105.47
1	E	11[C]	HYP	CG-CD-N	-2.47	101.49	105.47
1	G	20[A]	HYP	OD1-CG-CD	2.47	116.12	110.31
1	D	29[E]	HYP	CB-CG-CD	2.47	106.97	103.46
1	E	8[D]	HYP	CG-CD-N	-2.47	101.49	105.47
1	E	5[A]	HYP	CG-CD-N	-2.47	101.50	105.47
1	G	26[G]	HYP	CB-CG-CD	2.47	106.97	103.46
1	B	11[G]	HYP	OD1-CG-CD	2.46	116.10	110.31
1	D	29[F]	HYP	CB-CG-CD	2.46	106.97	103.46
1	G	8[D]	HYP	CG-CD-N	-2.46	101.51	105.47
1	G	20[C]	HYP	OD1-CG-CB	-2.46	103.50	110.16
1	I	26[G]	HYP	CB-CG-CD	2.46	106.96	103.46
1	G	20[G]	HYP	CG-CD-N	-2.45	101.52	105.47
1	H	29[F]	HYP	CB-CG-CD	2.45	106.95	103.46
1	D	16[D]	HYP	CD-N-CA	2.45	112.78	108.15
1	G	29[A]	HYP	CG-CD-N	-2.45	101.52	105.47
1	G	20[B]	HYP	OD1-CG-CD	2.45	116.06	110.31
1	E	11[D]	HYP	CG-CD-N	-2.44	101.53	105.47
1	F	11[F]	HYP	CG-CD-N	-2.44	101.54	105.47
1	G	26[E]	HYP	CB-CG-CD	2.44	106.93	103.46
1	G	20[E]	HYP	CG-CD-N	-2.44	101.55	105.47
1	F	11[G]	HYP	CB-CG-CD	-2.44	100.00	103.46
1	H	26[E]	HYP	CD-N-CA	2.43	112.75	108.15
1	I	8[G]	HYP	CG-CD-N	-2.43	101.56	105.47
1	C	26[G]	HYP	OD1-CG-CD	2.43	116.02	110.31
1	H	26[A]	HYP	CD-N-CA	2.43	112.73	108.15
1	G	23[D]	HYP	OD1-CG-CD	2.42	116.01	110.31
1	H	11[E]	HYP	CG-CD-N	-2.42	101.57	105.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	2[E]	HYP	CG-CD-N	-2.42	101.57	105.47
1	H	2[A]	HYP	CG-CD-N	-2.42	101.57	105.47
1	H	2[B]	HYP	CG-CD-N	-2.42	101.57	105.47
1	F	11[A]	HYP	CG-CD-N	-2.42	101.57	105.47
1	G	20[F]	HYP	CG-CD-N	-2.42	101.57	105.47
1	H	29[G]	HYP	CB-CG-CD	2.42	106.90	103.46
1	B	26[G]	HYP	C-CA-N	2.42	115.33	110.86
1	E	8[C]	HYP	CG-CD-N	-2.41	101.58	105.47
1	H	23[A]	HYP	CG-CD-N	-2.41	101.59	105.47
1	C	26[G]	HYP	CG-CD-N	-2.41	101.59	105.47
1	E	23[E]	HYP	CG-CD-N	-2.41	101.59	105.47
1	B	26[F]	HYP	C-CA-N	2.41	115.32	110.86
1	G	29[B]	HYP	CG-CD-N	-2.40	101.60	105.47
1	F	11[E]	HYP	CG-CD-N	-2.40	101.61	105.47
1	C	29[D]	HYP	OD1-CG-CB	-2.39	103.69	110.16
1	H	11[D]	HYP	CG-CD-N	-2.39	101.62	105.47
1	H	11[C]	HYP	CG-CD-N	-2.39	101.62	105.47
1	G	29[F]	HYP	OD1-CG-CB	-2.39	103.69	110.16
1	H	26[D]	HYP	CG-CD-N	-2.39	101.62	105.47
1	C	2[D]	HYP	OD1-CG-CB	-2.39	103.69	110.16
1	G	29[G]	HYP	OD1-CG-CB	-2.39	103.69	110.16
1	G	20[D]	HYP	C-CA-N	2.38	115.27	110.86
1	C	2[B]	HYP	OD1-CG-CB	-2.38	103.71	110.16
1	G	29[E]	HYP	OD1-CG-CB	-2.38	103.72	110.16
1	A	29[A]	HYP	CG-CD-N	-2.38	101.64	105.47
1	C	29[B]	HYP	OD1-CG-CB	-2.38	103.72	110.16
1	E	20[F]	HYP	CG-CD-N	-2.38	101.64	105.47
1	H	26[A]	HYP	OD1-CG-CB	-2.37	103.73	110.16
1	D	26[C]	HYP	CB-CG-CD	2.37	106.84	103.46
1	D	26[B]	HYP	CB-CG-CD	2.37	106.84	103.46
1	D	26[A]	HYP	CB-CG-CD	2.37	106.84	103.46
1	E	20[G]	HYP	CG-CD-N	-2.37	101.65	105.47
1	D	13[G]	HYP	CG-CD-N	-2.37	101.66	105.47
1	C	8[G]	HYP	OD1-CG-CB	-2.37	103.75	110.16
1	H	8[E]	HYP	C-CA-N	2.36	115.24	110.86
1	B	11[E]	HYP	CG-CD-N	-2.36	101.66	105.47
1	H	8[A]	HYP	C-CA-N	2.36	115.24	110.86
1	H	8[B]	HYP	C-CA-N	2.36	115.24	110.86
1	E	26[D]	HYP	CG-CD-N	-2.36	101.67	105.47
1	E	23[A]	HYP	CG-CD-N	-2.36	101.67	105.47
1	E	26[C]	HYP	CG-CD-N	-2.36	101.67	105.47
1	B	8[G]	HYP	CG-CD-N	-2.36	101.67	105.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	13[F]	HZP	CG-CD-N	-2.35	101.68	105.47
1	C	2[C]	HYP	OD1-CG-CB	-2.35	103.78	110.16
1	B	8[D]	HYP	C-CA-N	2.35	115.21	110.86
1	G	26[E]	HYP	CG-CB-CA	2.35	107.65	104.22
1	B	11[A]	HYP	CG-CD-N	-2.34	101.70	105.47
1	H	5[G]	HYP	C-CA-N	2.34	115.20	110.86
1	H	23[E]	HYP	CG-CD-N	-2.34	101.70	105.47
1	H	5[F]	HYP	C-CA-N	2.34	115.20	110.86
1	H	29[G]	HYP	CG-CB-CA	2.34	107.64	104.22
1	I	11[E]	HYP	CD-N-CA	2.34	112.57	108.15
1	D	13[E]	HZP	CG-CD-N	-2.34	101.70	105.47
1	I	11[F]	HYP	CD-N-CA	2.34	112.57	108.15
1	H	29[F]	HYP	CG-CB-CA	2.34	107.64	104.22
1	H	13[C]	HZP	CB-CA-N	-2.34	101.00	104.59
1	H	13[D]	HZP	CB-CA-N	-2.34	101.00	104.59
1	B	8[F]	HYP	CG-CD-N	-2.34	101.70	105.47
1	A	11[F]	HYP	OD1-CG-CB	-2.34	103.82	110.16
1	E	23[B]	HYP	CG-CD-N	-2.34	101.71	105.47
1	B	8[C]	HYP	C-CA-N	2.34	115.19	110.86
1	I	8[G]	HYP	CD-N-CA	2.34	112.56	108.15
1	C	29[C]	HYP	OD1-CG-CB	-2.34	103.83	110.16
1	H	26[E]	HYP	OD1-CG-CB	-2.34	103.83	110.16
1	C	20[F]	HYP	CG-CD-N	-2.34	101.71	105.47
1	H	2[E]	HYP	CB-CG-CD	2.34	106.78	103.46
1	H	2[A]	HYP	CB-CG-CD	2.34	106.78	103.46
1	H	2[B]	HYP	CB-CG-CD	2.34	106.78	103.46
1	I	5[C]	HYP	OD1-CG-CB	-2.33	103.84	110.16
1	G	26[G]	HYP	CG-CB-CA	2.33	107.63	104.22
1	I	5[G]	HYP	OD1-CG-CB	-2.33	103.84	110.16
1	I	8[A]	HYP	OD1-CG-CB	-2.33	103.85	110.16
1	H	26[B]	HYP	OD1-CG-CB	-2.33	103.86	110.16
1	C	20[A]	HYP	CG-CD-N	-2.33	101.72	105.47
1	E	29[F]	HYP	CG-CD-N	-2.33	101.72	105.47
1	E	16[C]	HZP	CG-CD-N	-2.33	101.72	105.47
1	C	2[C]	HYP	CB-CG-CD	2.33	106.77	103.46
1	I	8[F]	HYP	OD1-CG-CB	-2.33	103.86	110.16
1	I	29[G]	HYP	OD1-CG-CB	-2.32	103.87	110.16
1	H	8[D]	HYP	CG-CD-N	-2.32	101.73	105.47
1	E	16[D]	HZP	CG-CD-N	-2.32	101.73	105.47
1	I	8[E]	HYP	OD1-CG-CB	-2.32	103.87	110.16
1	E	29[G]	HYP	CG-CD-N	-2.32	101.73	105.47
1	A	29[B]	HYP	CG-CD-N	-2.32	101.73	105.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	29[F]	HYP	CG-CB-CA	2.32	107.61	104.22
1	I	11[A]	HYP	CD-N-CA	2.32	112.53	108.15
1	H	16[E]	HZP	C-CA-N	2.32	115.15	110.86
1	H	16[B]	HZP	C-CA-N	2.32	115.15	110.86
1	A	11[E]	HYP	OD1-CG-CB	-2.32	103.88	110.16
1	H	16[A]	HZP	C-CA-N	2.32	115.15	110.86
1	A	29[C]	HYP	CG-CD-N	-2.32	101.74	105.47
1	D	29[E]	HYP	CG-CB-CA	2.32	107.60	104.22
1	A	8[D]	HYP	CB-CG-CD	2.32	106.75	103.46
1	B	11[A]	HYP	OD1-CG-CB	-2.31	103.90	110.16
1	B	11[B]	HYP	OD1-CG-CB	-2.31	103.90	110.16
1	B	11[E]	HYP	OD1-CG-CB	-2.31	103.90	110.16
1	D	29[G]	HYP	CG-CB-CA	2.31	107.60	104.22
1	I	11[B]	HYP	OD1-CG-CB	-2.31	103.92	110.16
1	B	20[C]	HYP	C-CA-N	2.31	115.13	110.86
1	H	23[B]	HYP	CG-CD-N	-2.31	101.75	105.47
1	B	11[B]	HYP	CG-CD-N	-2.30	101.76	105.47
1	F	26[F]	HYP	CB-CG-CD	2.30	106.74	103.46
1	G	20[D]	HYP	CG-CD-N	-2.30	101.76	105.47
1	E	13[B]	HZP	CG-CD-N	-2.30	101.76	105.47
1	E	13[A]	HZP	CG-CD-N	-2.30	101.76	105.47
1	I	11[C]	HYP	OD1-CG-CB	-2.30	103.93	110.16
1	C	16[F]	HZP	CG-CD-N	-2.30	101.77	105.47
1	C	20[E]	HYP	CG-CD-N	-2.30	101.77	105.47
1	G	16[C]	HZP	C-CA-N	2.30	115.11	110.86
1	H	8[C]	HYP	CG-CD-N	-2.30	101.77	105.47
1	G	2[A]	HYP	CG-CD-N	-2.30	101.77	105.47
1	H	20[D]	HYP	OD1-CG-CB	-2.30	103.94	110.16
1	I	5[D]	HYP	OD1-CG-CB	-2.30	103.94	110.16
1	H	20[C]	HYP	OD1-CG-CB	-2.30	103.94	110.16
1	F	8[G]	HYP	CG-CD-N	-2.30	101.77	105.47
1	F	16[C]	HZP	CG-CD-N	-2.29	101.78	105.47
1	I	11[D]	HYP	OD1-CG-CB	-2.29	103.95	110.16
1	F	16[B]	HZP	CG-CD-N	-2.29	101.78	105.47
1	H	23[F]	HYP	CG-CD-N	-2.29	101.78	105.47
1	H	23[G]	HYP	CG-CD-N	-2.29	101.78	105.47
1	F	23[G]	HYP	CG-CD-N	-2.29	101.79	105.47
1	B	5[G]	HYP	CG-CD-N	-2.29	101.79	105.47
1	B	5[C]	HYP	OD1-CG-CB	-2.29	103.97	110.16
1	A	11[G]	HYP	OD1-CG-CB	-2.29	103.96	110.16
1	C	16[A]	HZP	CG-CD-N	-2.28	101.79	105.47
1	F	8[G]	HYP	CB-CG-CD	2.28	106.71	103.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	16[D]	HZP	CG-CD-N	-2.28	101.80	105.47
1	G	2[C]	HYP	CG-CD-N	-2.28	101.80	105.47
1	I	5[B]	HYP	OD1-CG-CB	-2.28	103.99	110.16
1	D	2[E]	HYP	CG-CD-N	-2.28	101.80	105.47
1	G	5[D]	HYP	CG-CD-N	-2.28	101.80	105.47
1	D	29[D]	HYP	CB-CG-CD	2.28	106.70	103.46
1	H	26[C]	HYP	CG-CD-N	-2.28	101.80	105.47
1	G	16[B]	HZP	C-CA-N	2.28	115.08	110.86
1	G	16[A]	HZP	C-CA-N	2.28	115.08	110.86
1	H	5[E]	HYP	CG-CD-N	-2.28	101.81	105.47
1	B	20[D]	HYP	C-CA-N	2.28	115.07	110.86
1	D	23[D]	HYP	CG-CD-N	-2.28	101.81	105.47
1	H	5[A]	HYP	CG-CD-N	-2.28	101.81	105.47
1	E	13[E]	HZP	CG-CD-N	-2.27	101.81	105.47
1	E	2[C]	HYP	CG-CD-N	-2.27	101.81	105.47
1	E	2[D]	HYP	CG-CD-N	-2.27	101.81	105.47
1	B	23[B]	HYP	CD-N-CA	2.27	112.44	108.15
1	E	29[E]	HYP	CB-CA-N	-2.27	101.11	104.59
1	C	29[C]	HYP	CB-CG-CD	2.27	106.69	103.46
1	B	8[B]	HYP	CG-CD-N	-2.26	101.82	105.47
1	G	29[F]	HYP	CB-CG-CD	2.26	106.68	103.46
1	G	26[B]	HYP	CG-CD-N	-2.26	101.83	105.47
1	C	2[D]	HYP	CB-CG-CD	2.26	106.68	103.46
1	D	8[E]	HYP	CG-CD-N	-2.26	101.83	105.47
1	G	29[G]	HYP	CB-CG-CD	2.26	106.68	103.46
1	G	26[C]	HYP	CG-CD-N	-2.26	101.83	105.47
1	G	26[A]	HYP	CG-CD-N	-2.26	101.83	105.47
1	B	13[C]	HZP	CG-CD-N	-2.26	101.83	105.47
1	H	23[F]	HYP	OD1-CG-CB	-2.26	104.04	110.16
1	H	23[G]	HYP	OD1-CG-CB	-2.26	104.04	110.16
1	H	2[G]	HYP	CG-CD-N	-2.26	101.83	105.47
1	H	2[F]	HYP	CG-CD-N	-2.26	101.83	105.47
1	B	5[F]	HYP	CG-CD-N	-2.26	101.83	105.47
1	F	26[A]	HYP	CB-CG-CD	2.26	106.67	103.46
1	C	2[B]	HYP	CB-CG-CD	2.26	106.67	103.46
1	B	23[A]	HYP	CD-N-CA	2.26	112.41	108.15
1	D	26[A]	HYP	OD1-CG-CB	-2.26	104.05	110.16
1	G	23[F]	HYP	CG-CD-N	-2.26	101.84	105.47
1	G	23[G]	HYP	CG-CD-N	-2.26	101.84	105.47
1	G	23[E]	HYP	CG-CD-N	-2.26	101.84	105.47
1	D	2[G]	HYP	CG-CD-N	-2.25	101.84	105.47
1	D	2[F]	HYP	CG-CD-N	-2.25	101.84	105.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	20[G]	HYP	CD-N-CA	2.25	112.40	108.15
1	B	20[F]	HYP	CD-N-CA	2.25	112.40	108.15
1	G	11[B]	HYP	OD1-CG-CB	-2.25	104.07	110.16
1	E	5[B]	HYP	OD1-CG-CB	-2.25	104.07	110.16
1	I	16[C]	HZP	C-CA-N	2.25	115.02	110.86
1	E	29[E]	HYP	OD1-CG-CB	-2.25	104.08	110.16
1	D	23[E]	HYP	OD1-CG-CB	-2.24	104.08	110.16
1	D	26[B]	HYP	OD1-CG-CB	-2.25	104.08	110.16
1	B	8[A]	HYP	CG-CD-N	-2.24	101.86	105.47
1	B	26[C]	HYP	CD-N-CA	2.24	112.39	108.15
1	D	29[D]	HYP	OD1-CG-CB	-2.24	104.09	110.16
1	B	23[E]	HYP	CD-N-CA	2.24	112.39	108.15
1	B	8[G]	HYP	OD1-CG-CB	-2.24	104.09	110.16
1	C	16[E]	HZP	CG-CD-N	-2.24	101.86	105.47
1	D	26[C]	HYP	OD1-CG-CB	-2.24	104.10	110.16
1	C	29[B]	HYP	CB-CG-CD	2.24	106.65	103.46
1	B	26[D]	HYP	CD-N-CA	2.24	112.38	108.15
1	A	26[G]	HYP	OD1-CG-CB	-2.24	104.10	110.16
1	G	11[A]	HYP	OD1-CG-CB	-2.23	104.11	110.16
1	D	23[G]	HYP	OD1-CG-CB	-2.24	104.11	110.16
1	B	5[C]	HYP	CG-CD-N	-2.23	101.87	105.47
1	H	29[D]	HYP	CD-N-CA	2.23	112.37	108.15
1	B	13[D]	HZP	CG-CD-N	-2.23	101.87	105.47
1	D	8[F]	HYP	CG-CD-N	-2.23	101.87	105.47
1	H	5[B]	HYP	CG-CD-N	-2.23	101.87	105.47
1	H	29[C]	HYP	CD-N-CA	2.23	112.37	108.15
1	A	26[E]	HYP	OD1-CG-CB	-2.23	104.12	110.16
1	D	16[F]	HZP	CB-CA-N	-2.23	101.17	104.59
1	G	29[D]	HYP	CG-CD-N	-2.23	101.88	105.47
1	C	29[D]	HYP	CB-CG-CD	2.23	106.63	103.46
1	B	26[F]	HYP	OD1-CG-CD	2.23	115.55	110.31
1	A	11[G]	HYP	CB-CA-N	-2.23	101.17	104.59
1	E	5[A]	HYP	OD1-CG-CB	-2.23	104.13	110.16
1	D	23[F]	HYP	OD1-CG-CB	-2.23	104.13	110.16
1	B	8[E]	HYP	CG-CD-N	-2.23	101.89	105.47
1	A	11[B]	HYP	CG-CD-N	-2.23	101.89	105.47
1	E	29[A]	HYP	CB-CA-N	-2.22	101.18	104.59
1	D	8[G]	HYP	CG-CD-N	-2.22	101.89	105.47
1	D	20[C]	HYP	CG-CD-N	-2.22	101.89	105.47
1	D	2[D]	HYP	OD1-CG-CB	-2.22	104.14	110.16
1	D	20[B]	HYP	CG-CD-N	-2.22	101.89	105.47
1	G	8[G]	HYP	OD1-CG-CB	-2.22	104.14	110.16

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	11[C]	HYP	OD1-CG-CB	-2.22	104.15	110.16
1	F	20[G]	HYP	C-CA-N	2.22	114.97	110.86
1	D	16[E]	HZP	CB-CA-N	-2.22	101.19	104.59
1	E	20[C]	HYP	CB-CG-CD	2.22	106.61	103.46
1	F	26[E]	HYP	CB-CG-CD	2.22	106.61	103.46
1	D	16[G]	HZP	CB-CA-N	-2.21	101.19	104.59
1	A	26[F]	HYP	OD1-CG-CB	-2.21	104.16	110.16
1	A	11[E]	HYP	CB-CA-N	-2.21	101.19	104.59
1	B	2[A]	HYP	OD1-CG-CB	-2.21	104.17	110.16
1	I	16[D]	HZP	C-CA-N	2.21	114.95	110.86
1	B	2[E]	HYP	OD1-CG-CB	-2.21	104.17	110.16
1	B	8[F]	HYP	OD1-CG-CB	-2.21	104.17	110.16
1	B	2[B]	HYP	OD1-CG-CB	-2.21	104.17	110.16
1	G	29[E]	HYP	CB-CG-CD	2.21	106.61	103.46
1	A	23[A]	HYP	CG-CD-N	-2.21	101.92	105.47
1	C	11[G]	HYP	CD-N-CA	2.21	112.31	108.15
1	E	5[E]	HYP	OD1-CG-CB	-2.21	104.19	110.16
1	E	29[A]	HYP	OD1-CG-CB	-2.21	104.19	110.16
1	B	26[G]	HYP	OD1-CG-CD	2.21	115.50	110.31
1	A	11[C]	HYP	CG-CD-N	-2.21	101.92	105.47
1	G	11[D]	HYP	C-CA-N	2.20	114.94	110.86
1	G	8[E]	HYP	OD1-CG-CB	-2.20	104.19	110.16
1	E	29[B]	HYP	OD1-CG-CB	-2.20	104.19	110.16
1	E	29[F]	HYP	CB-CG-CD	2.20	106.59	103.46
1	A	23[C]	HYP	CG-CD-N	-2.20	101.93	105.47
1	A	11[F]	HYP	CB-CA-N	-2.20	101.22	104.59
1	I	20[E]	HYP	OD1-CG-CB	-2.20	104.21	110.16
1	G	2[B]	HYP	CG-CD-N	-2.19	101.94	105.47
1	D	20[A]	HYP	CG-CD-N	-2.19	101.94	105.47
1	A	26[C]	HYP	CG-CD-N	-2.19	101.94	105.47
1	A	8[F]	HYP	CG-CD-N	-2.19	101.94	105.47
1	D	8[D]	HYP	CG-CD-N	-2.19	101.94	105.47
1	H	13[C]	HZP	CD-N-CA	2.19	112.29	108.15
1	H	13[D]	HZP	CD-N-CA	2.19	112.29	108.15
1	A	11[A]	HYP	CG-CD-N	-2.19	101.94	105.47
1	A	23[B]	HYP	CG-CD-N	-2.19	101.95	105.47
1	A	26[B]	HYP	CG-CD-N	-2.19	101.95	105.47
1	F	20[C]	HYP	OD1-CG-CB	-2.19	104.24	110.16
1	G	8[F]	HYP	OD1-CG-CB	-2.19	104.24	110.16
1	A	26[D]	HYP	CG-CD-N	-2.18	101.95	105.47
1	E	29[A]	HYP	C-CA-N	2.18	114.90	110.86
1	E	29[E]	HYP	C-CA-N	2.18	114.90	110.86

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	I	20[F]	HYP	OD1-CG-CB	-2.18	104.26	110.16
1	I	20[A]	HYP	OD1-CG-CB	-2.18	104.25	110.16
1	D	13[D]	HZP	CG-CD-N	-2.18	101.96	105.47
1	I	2[D]	HYP	OD1-CG-CB	-2.18	104.26	110.16
1	I	2[C]	HYP	OD1-CG-CB	-2.18	104.26	110.16
1	I	2[D]	HYP	CG-CD-N	-2.18	101.97	105.47
1	G	11[D]	HYP	CB-CA-N	-2.18	101.25	104.59
1	I	16[B]	HZP	C-CA-N	2.18	114.89	110.86
1	I	2[C]	HYP	CG-CD-N	-2.18	101.97	105.47
1	E	29[B]	HYP	CB-CA-N	-2.17	101.25	104.59
1	I	29[F]	HYP	CB-CG-CD	2.17	106.55	103.46
1	I	2[B]	HYP	OD1-CG-CB	-2.17	104.28	110.16
1	F	5[D]	HYP	OD1-CG-CB	-2.17	104.29	110.16
1	D	23[E]	HYP	CB-CG-CD	2.17	106.54	103.46
1	A	23[D]	HYP	OD1-CG-CB	-2.17	104.29	110.16
1	D	23[F]	HYP	CB-CG-CD	2.17	106.54	103.46
1	D	23[G]	HYP	CB-CG-CD	2.17	106.54	103.46
1	E	29[B]	HYP	C-CA-N	2.17	114.87	110.86
1	A	29[D]	HYP	CG-CD-N	-2.16	101.98	105.47
1	F	20[B]	HYP	OD1-CG-CB	-2.16	104.31	110.16
1	A	16[D]	HZP	CG-CD-N	-2.16	101.99	105.47
1	F	5[B]	HYP	OD1-CG-CB	-2.16	104.30	110.16
1	I	8[A]	HYP	CB-CG-CD	2.16	106.53	103.46
1	I	29[E]	HYP	CB-CG-CD	2.16	106.53	103.46
1	A	20[E]	HYP	CG-CD-N	-2.15	102.00	105.47
1	E	16[B]	HZP	CG-CD-N	-2.15	102.01	105.47
1	E	20[D]	HYP	CB-CG-CD	2.15	106.51	103.46
1	H	2[E]	HYP	OD1-CG-CB	-2.15	104.35	110.16
1	H	2[A]	HYP	OD1-CG-CB	-2.15	104.35	110.16
1	C	29[E]	HYP	CD-N-CA	2.15	112.20	108.15
1	I	29[A]	HYP	CB-CG-CD	2.15	106.51	103.46
1	A	20[G]	HYP	CG-CD-N	-2.14	102.02	105.47
1	A	2[G]	HYP	CB-CG-CD	2.14	106.51	103.46
1	E	11[E]	HYP	OD1-CG-CB	-2.14	104.36	110.16
1	A	8[G]	HYP	CG-CD-N	-2.14	102.03	105.47
1	A	8[E]	HYP	CG-CD-N	-2.14	102.03	105.47
1	E	23[F]	HYP	CG-CD-N	-2.14	102.02	105.47
1	A	20[B]	HYP	OD1-CG-CB	-2.14	104.37	110.16
1	F	2[D]	HYP	CD-N-CA	2.14	112.18	108.15
1	A	20[C]	HYP	OD1-CG-CB	-2.13	104.38	110.16
1	C	2[G]	HYP	OD1-CG-CB	-2.13	104.38	110.16
1	A	26[A]	HYP	CG-CD-N	-2.13	102.04	105.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	20[F]	HYP	CG-CD-N	-2.13	102.04	105.47
1	E	23[G]	HYP	CG-CD-N	-2.13	102.04	105.47
1	E	8[F]	HYP	OD1-CG-CB	-2.12	104.41	110.16
1	G	5[C]	HYP	OD1-CG-CB	-2.12	104.41	110.16
1	A	20[A]	HYP	OD1-CG-CB	-2.12	104.41	110.16
1	H	2[B]	HYP	OD1-CG-CB	-2.12	104.41	110.16
1	E	29[G]	HYP	CB-CG-CD	2.12	106.48	103.46
1	F	5[C]	HYP	OD1-CG-CB	-2.12	104.41	110.16
1	I	2[E]	HYP	C-CA-N	2.12	114.79	110.86
1	A	23[F]	HYP	CG-CD-N	-2.12	102.05	105.47
1	I	8[F]	HYP	CB-CG-CD	2.12	106.48	103.46
1	F	20[D]	HYP	OD1-CG-CB	-2.12	104.42	110.16
1	E	11[A]	HYP	OD1-CG-CB	-2.12	104.42	110.16
1	I	2[F]	HYP	C-CA-N	2.12	114.78	110.86
1	F	16[G]	HYP	CB-CA-N	-2.11	101.35	104.59
1	C	29[A]	HYP	CD-N-CA	2.11	112.14	108.15
1	C	8[B]	HYP	CG-CD-N	-2.11	102.07	105.47
1	G	5[A]	HYP	OD1-CG-CB	-2.11	104.45	110.16
1	I	2[A]	HYP	C-CA-N	2.11	114.77	110.86
1	C	8[D]	HYP	CG-CD-N	-2.11	102.07	105.47
1	H	2[D]	HYP	CB-CA-N	-2.11	101.36	104.59
1	H	2[C]	HYP	CB-CA-N	-2.11	101.36	104.59
1	A	2[E]	HYP	CB-CG-CD	2.11	106.46	103.46
1	C	5[F]	HYP	CG-CD-N	-2.10	102.08	105.47
1	A	23[G]	HYP	CG-CD-N	-2.10	102.08	105.47
1	E	8[G]	HYP	OD1-CG-CB	-2.10	104.47	110.16
1	G	5[C]	HYP	C-CA-N	2.10	114.75	110.86
1	I	23[B]	HYP	OD1-CG-CB	-2.10	104.48	110.16
1	C	23[C]	HYP	CG-CD-N	-2.10	102.09	105.47
1	F	2[B]	HYP	CD-N-CA	2.10	112.11	108.15
1	C	5[A]	HYP	CG-CD-N	-2.10	102.09	105.47
1	E	11[B]	HYP	OD1-CG-CB	-2.10	104.48	110.16
1	A	5[B]	HYP	OD1-CG-CB	-2.09	104.49	110.16
1	E	11[G]	HYP	C-CA-N	2.09	114.74	110.86
1	E	11[F]	HYP	C-CA-N	2.09	114.74	110.86
1	G	5[A]	HYP	C-CA-N	2.09	114.73	110.86
1	G	5[B]	HYP	OD1-CG-CB	-2.09	104.50	110.16
1	I	23[C]	HYP	OD1-CG-CB	-2.09	104.50	110.16
1	H	20[C]	HYP	CD-N-CA	2.09	112.09	108.15
1	F	11[F]	HYP	CB-CG-CD	2.09	106.43	103.46
1	A	23[E]	HYP	CG-CD-N	-2.08	102.11	105.47
1	C	8[C]	HYP	CG-CD-N	-2.08	102.11	105.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	I	8[E]	HYP	CB-CG-CD	2.08	106.43	103.46
1	A	5[A]	HYP	OD1-CG-CB	-2.08	104.52	110.16
1	G	5[B]	HYP	C-CA-N	2.08	114.71	110.86
1	D	5[D]	HYP	CB-CA-N	-2.08	101.40	104.59
1	B	16[D]	HZP	CG-CB-CA	-2.08	101.19	104.22
1	H	11[F]	HYP	OD1-CG-CD	2.08	115.21	110.31
1	E	29[D]	HYP	CG-CD-N	-2.08	102.12	105.47
1	B	16[C]	HZP	CG-CB-CA	-2.08	101.19	104.22
1	I	23[D]	HYP	OD1-CG-CB	-2.08	104.52	110.16
1	H	11[G]	HYP	OD1-CG-CD	2.08	115.20	110.31
1	I	5[G]	HYP	CB-CG-CD	2.08	106.42	103.46
1	F	29[C]	HYP	CB-CA-N	-2.08	101.40	104.59
1	E	16[A]	HZP	CG-CD-N	-2.08	102.12	105.47
1	C	29[F]	HYP	CD-N-CA	2.08	112.07	108.15
1	C	23[B]	HYP	CG-CD-N	-2.07	102.13	105.47
1	G	29[E]	HYP	CG-CD-N	-2.07	102.13	105.47
1	H	8[G]	HYP	CG-CD-N	-2.07	102.13	105.47
1	H	8[F]	HYP	CG-CD-N	-2.07	102.13	105.47
1	A	11[D]	HYP	OD1-CG-CB	-2.07	104.56	110.16
1	H	20[D]	HYP	CD-N-CA	2.07	112.05	108.15
1	C	5[E]	HYP	OD1-CG-CB	-2.07	104.56	110.16
1	C	2[C]	HYP	CG-CD-N	-2.07	102.14	105.47
1	E	26[F]	HYP	CG-CD-N	-2.07	102.14	105.47
1	A	5[C]	HYP	OD1-CG-CB	-2.07	104.56	110.16
1	E	26[G]	HYP	CG-CD-N	-2.07	102.14	105.47
1	A	2[F]	HYP	CB-CG-CD	2.06	106.40	103.46
1	C	5[E]	HYP	CG-CD-N	-2.06	102.15	105.47
1	F	16[A]	HZP	C-CA-N	2.06	114.68	110.86
1	B	2[F]	HYP	C-CA-N	2.06	114.68	110.86
1	F	16[F]	HZP	C-CA-N	2.06	114.68	110.86
1	F	16[E]	HZP	C-CA-N	2.06	114.68	110.86
1	F	2[C]	HYP	CD-N-CA	2.06	112.04	108.15
1	C	13[G]	HZP	C-CA-N	2.06	114.67	110.86
1	I	2[B]	HYP	CG-CD-N	-2.06	102.16	105.47
1	F	26[F]	HYP	CB-CA-N	-2.06	101.43	104.59
1	I	16[G]	HZP	C-CA-N	2.06	114.67	110.86
1	A	13[F]	HZP	CG-CD-N	-2.05	102.17	105.47
1	E	29[C]	HYP	CG-CD-N	-2.05	102.17	105.47
1	C	5[A]	HYP	OD1-CG-CB	-2.05	104.61	110.16
1	C	8[B]	HYP	OD1-CG-CB	-2.05	104.62	110.16
1	G	29[F]	HYP	CG-CD-N	-2.04	102.18	105.47
1	A	13[E]	HZP	CG-CD-N	-2.04	102.18	105.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	11[A]	HYP	CB-CG-CD	2.05	106.37	103.46
1	F	2[F]	HYP	C-CA-N	2.04	114.64	110.86
1	H	29[A]	HYP	OD1-CG-CB	-2.04	104.64	110.16
1	C	11[D]	HYP	OD1-CG-CB	-2.04	104.64	110.16
1	C	23[D]	HYP	CG-CD-N	-2.04	102.19	105.47
1	C	8[D]	HYP	OD1-CG-CB	-2.04	104.64	110.16
1	B	16[C]	HZP	OD1-CG-CD	-2.03	105.54	110.31
1	E	26[B]	HYP	CG-CD-N	-2.03	102.20	105.47
1	D	11[D]	HYP	CG-CD-N	-2.03	102.20	105.47
1	F	26[E]	HYP	CB-CA-N	-2.03	101.47	104.59
1	I	16[F]	HZP	CB-CA-N	-2.03	101.48	104.59
1	C	8[C]	HYP	OD1-CG-CB	-2.03	104.67	110.16
1	F	8[E]	HYP	CG-CD-N	-2.03	102.21	105.47
1	C	11[C]	HYP	OD1-CG-CB	-2.02	104.68	110.16
1	C	2[D]	HYP	CG-CD-N	-2.02	102.21	105.47
1	B	23[D]	HYP	CB-CG-CD	2.02	106.34	103.46
1	G	29[G]	HYP	CG-CD-N	-2.02	102.21	105.47
1	A	5[G]	HYP	CB-CA-N	-2.02	101.49	104.59
1	F	2[A]	HYP	C-CA-N	2.02	114.60	110.86
1	F	29[B]	HYP	CB-CA-N	-2.02	101.49	104.59
1	F	26[A]	HYP	CB-CA-N	-2.02	101.49	104.59
1	F	5[G]	HYP	CG-CD-N	-2.02	102.22	105.47
1	H	29[B]	HYP	OD1-CG-CB	-2.02	104.69	110.16
1	B	16[D]	HZP	OD1-CG-CD	-2.02	105.57	110.31
1	C	2[G]	HYP	CG-CD-N	-2.02	102.22	105.47
1	F	2[C]	HYP	OD1-CG-CD	2.01	115.05	110.31
1	F	2[B]	HYP	OD1-CG-CD	2.01	115.05	110.31
1	C	23[G]	HYP	CB-CG-CD	2.01	106.32	103.46
1	C	11[B]	HYP	OD1-CG-CB	-2.01	104.71	110.16
1	D	26[F]	HYP	C-CA-N	2.01	114.58	110.86
1	H	23[C]	HYP	OD1-CG-CB	-2.01	104.71	110.16
1	H	23[D]	HYP	OD1-CG-CB	-2.01	104.71	110.16
1	I	11[B]	HYP	CB-CG-CD	2.01	106.32	103.46
1	F	2[D]	HYP	OD1-CG-CD	2.01	115.03	110.31
1	I	20[B]	HYP	OD1-CG-CB	-2.01	104.72	110.16
1	C	5[F]	HYP	OD1-CG-CB	-2.01	104.73	110.16
1	E	16[E]	HZP	CG-CD-N	-2.01	102.24	105.47
1	E	26[A]	HYP	CG-CD-N	-2.01	102.24	105.47
1	E	26[E]	HYP	CG-CD-N	-2.01	102.24	105.47
1	H	11[D]	HYP	CB-CA-N	-2.00	101.52	104.59
1	E	2[C]	HYP	CB-CG-CD	2.00	106.31	103.46
1	A	8[B]	HYP	CB-CA-N	-2.00	101.52	104.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	26[F]	HYP	CG-CD-N	-2.00	102.24	105.47
1	E	2[D]	HYP	CB-CG-CD	2.00	106.31	103.46
1	F	11[E]	HYP	CB-CG-CD	2.00	106.31	103.46
1	F	26[A]	HYP	CG-CD-N	-2.00	102.25	105.47
1	F	8[A]	HYP	CG-CD-N	-2.00	102.25	105.47
1	C	2[B]	HYP	CG-CD-N	-2.00	102.25	105.47
1	C	11[G]	HYP	CB-CA-N	-2.00	101.52	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 ⓘ

There are no ligands in this entry.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å ²)	Q<0.9
1	A	30/30 (100%)	-0.06	0	100	100	4, 5, 15, 16	0
1	B	30/30 (100%)	-0.05	0	100	100	3, 5, 14, 15	0
1	C	30/30 (100%)	-0.06	1 (3%)	44	50	3, 5, 17, 18	0
1	D	30/30 (100%)	0.08	2 (6%)	17	19	3, 5, 15, 17	0
1	E	30/30 (100%)	0.06	2 (6%)	17	19	4, 6, 15, 17	0
1	F	30/30 (100%)	0.04	2 (6%)	17	19	4, 5, 16, 19	0
1	G	30/30 (100%)	0.40	3 (10%)	8	8	4, 6, 16, 20	0
1	H	30/30 (100%)	-0.06	0	100	100	3, 6, 15, 16	0
1	I	30/30 (100%)	0.00	0	100	100	3, 6, 17, 18	0
All	All	270/270 (100%)	0.04	10 (3%)	39	44	3, 5, 16, 20	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	2[A]	HYP	3.8
1	D	29[A]	HYP	2.8
1	G	1[A]	PRO	2.6
1	C	29[A]	HYP	2.4
1	D	30[A]	GLY	2.2
1	F	28[A]	PRO	2.2
1	G	3[A]	GLY	2.2
1	E	29[A]	HYP	2.2
1	E	1[A]	PRO	2.1
1	F	29[A]	HYP	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, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(Å ²)	Q<0.9
1	HYP	D	29[D]	8/9	0.18	3.00	5,6,7,10	8
1	HYP	C	29[F]	8/9	0.16	2.68	19,20,20,20	8
1	HYP	C	29[E]	8/9	0.16	2.68	19,20,20,20	8
1	HYP	C	29[A]	8/9	0.16	2.68	19,20,20,20	8
1	HYP	G	2[D]	8/9	0.22	2.63	14,15,16,17	8
1	HYP	C	29[B]	8/9	0.16	2.62	11,12,14,16	8
1	HYP	C	29[C]	8/9	0.16	2.62	11,12,14,16	8
1	HYP	C	29[D]	8/9	0.16	2.62	11,12,14,16	8
1	HYP	G	2[C]	8/9	0.22	2.45	21,22,23,24	8
1	HYP	G	2[B]	8/9	0.22	2.45	21,22,23,24	8
1	HYP	G	2[A]	8/9	0.22	2.45	21,22,23,24	8
1	HYP	G	2[E]	8/9	0.22	2.39	8,9,11,12	8
1	HYP	G	2[F]	8/9	0.22	2.39	8,9,11,12	8
1	HYP	G	2[G]	8/9	0.22	2.39	8,9,11,12	8
1	HYP	C	29[G]	8/9	0.16	2.33	12,14,14,16	8
1	HYP	E	29[D]	8/9	0.16	2.17	7,8,9,10	8
1	HYP	E	29[C]	8/9	0.16	2.17	7,8,9,10	8
1	HYP	H	2[G]	8/9	0.15	2.02	6,7,7,9	8
1	HYP	H	2[F]	8/9	0.15	2.02	6,7,7,9	8
1	HYP	F	2[G]	8/9	0.12	1.96	5,5,7,8	8
1	HYP	D	2[G]	8/9	0.14	1.73	7,8,9,11	8
1	HYP	D	2[E]	8/9	0.14	1.73	7,8,9,11	8
1	HYP	D	2[F]	8/9	0.14	1.73	7,8,9,11	8
1	HYP	F	5[G]	8/9	0.07	1.71	2,3,4,6	8
1	HYP	D	29[A]	8/9	0.18	1.66	14,15,16,17	8
1	HYP	D	29[C]	8/9	0.18	1.66	14,15,16,17	8
1	HYP	D	29[B]	8/9	0.18	1.66	14,15,16,17	8
1	HYP	G	5[E]	8/9	0.13	1.57	3,4,5,6	8
1	HYP	E	26[D]	8/9	0.09	1.57	3,3,4,5	8
1	HYP	G	5[G]	8/9	0.13	1.57	3,4,5,6	8
1	HYP	E	26[C]	8/9	0.09	1.57	3,3,4,5	8
1	HYP	G	5[F]	8/9	0.13	1.57	3,4,5,6	8
1	HYP	E	29[F]	8/9	0.16	1.53	15,16,17,18	8
1	HYP	E	29[G]	8/9	0.16	1.53	15,16,17,18	8

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
1	HYP	A	29[D]	8/9	0.12	1.52	5,5,6,7	8
1	HYP	E	29[A]	8/9	0.16	1.49	20,21,22,22	8
1	HYP	E	29[E]	8/9	0.16	1.49	20,21,22,22	8
1	HYP	E	29[B]	8/9	0.16	1.49	20,21,22,22	8
1	HYP	G	29[D]	8/9	0.11	1.47	6,7,8,11	8
1	HYP	A	26[D]	8/9	0.08	1.45	4,5,6,9	8
1	HYP	D	2[A]	8/9	0.14	1.43	16,17,18,18	8
1	HYP	E	5[G]	8/9	0.08	1.43	2,3,5,6	8
1	HYP	D	2[B]	8/9	0.14	1.43	16,17,18,18	8
1	HYP	D	2[C]	8/9	0.14	1.43	16,17,18,18	8
1	HYP	E	5[F]	8/9	0.08	1.42	2,3,5,6	8
1	HYP	C	2[G]	8/9	0.10	1.40	4,4,5,5	8
1	HYP	D	2[D]	8/9	0.14	1.38	14,14,14,15	8
1	HYP	D	29[E]	8/9	0.18	1.34	20,21,22,23	8
1	HYP	D	29[F]	8/9	0.18	1.34	20,21,22,23	8
1	HYP	D	29[G]	8/9	0.18	1.34	20,21,22,23	8
1	HYP	D	26[D]	8/9	0.08	1.30	3,3,4,6	8
1	HYP	C	26[C]	8/9	0.11	1.27	5,5,6,8	8
1	HYP	C	26[D]	8/9	0.11	1.27	5,5,6,8	8
1	HYP	C	26[B]	8/9	0.11	1.27	5,5,6,8	8
1	HYP	F	29[D]	8/9	0.15	1.24	10,12,12,13	8
1	HYP	F	29[C]	8/9	0.15	1.24	10,12,12,13	8
1	HYP	F	29[B]	8/9	0.15	1.24	10,12,12,13	8
1	HYP	H	2[A]	8/9	0.15	1.21	15,16,17,17	8
1	HYP	H	2[B]	8/9	0.15	1.21	15,16,17,17	8
1	HYP	H	2[E]	8/9	0.15	1.21	15,16,17,17	8
1	HYP	E	2[G]	8/9	0.12	1.15	6,6,7,9	8
1	HYP	E	2[F]	8/9	0.12	1.15	6,6,7,9	8
1	HYP	H	2[D]	8/9	0.15	1.04	18,20,22,22	8
1	HYP	H	2[C]	8/9	0.15	1.04	18,20,22,22	8
1	HYP	F	2[F]	8/9	0.12	0.97	12,14,15,18	8
1	HYP	B	29[D]	8/9	0.12	0.97	6,7,8,10	8
1	HYP	B	29[C]	8/9	0.12	0.97	6,7,8,10	8
1	HYP	F	2[E]	8/9	0.12	0.97	12,14,15,18	8
1	HYP	F	2[A]	8/9	0.12	0.97	12,14,15,18	8
1	HYP	F	29[G]	8/9	0.15	0.88	13,15,16,18	8
1	HZP	I	16[F]	8/9	0.06	0.82	3,4,5,7	8
1	HZP	I	16[A]	8/9	0.06	0.82	3,4,5,7	8
1	HZP	I	16[E]	8/9	0.06	0.82	3,4,5,7	8
1	HYP	E	20[D]	8/9	0.06	0.77	2,3,4,4	8
1	HYP	E	20[C]	8/9	0.06	0.77	2,3,4,4	8
1	HZP	I	16[G]	8/9	0.06	0.76	3,3,4,4	8

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
1	HYP	I	2[G]	8/9	0.11	0.75	3,4,6,7	8
1	HYP	A	29[B]	8/9	0.12	0.74	15,17,19,20	8
1	HYP	A	29[C]	8/9	0.12	0.74	15,17,19,20	8
1	HYP	A	29[A]	8/9	0.12	0.74	15,17,19,20	8
1	HZP	B	16[D]	8/9	0.06	0.74	3,4,5,11	8
1	HZP	B	16[C]	8/9	0.06	0.74	3,4,5,11	8
1	HYP	E	20[A]	8/9	0.06	0.72	3,4,6,7	8
1	HYP	E	20[E]	8/9	0.06	0.72	3,4,6,7	8
1	HYP	E	20[B]	8/9	0.06	0.72	3,4,6,7	8
1	HYP	D	5[E]	8/9	0.08	0.69	4,4,5,6	8
1	HYP	F	2[B]	8/9	0.12	0.69	19,20,20,22	8
1	HYP	D	5[F]	8/9	0.08	0.69	4,4,5,6	8
1	HYP	F	2[C]	8/9	0.12	0.69	19,20,20,22	8
1	HYP	F	2[D]	8/9	0.12	0.69	19,20,20,22	8
1	HYP	D	5[G]	8/9	0.08	0.69	4,4,5,6	8
1	HZP	B	16[B]	8/9	0.06	0.67	3,5,7,12	8
1	HZP	B	16[A]	8/9	0.06	0.67	3,5,7,12	8
1	HZP	B	16[E]	8/9	0.06	0.67	3,5,7,12	8
1	HYP	F	29[A]	8/9	0.15	0.63	24,24,25,27	8
1	HYP	F	29[F]	8/9	0.15	0.63	24,24,25,27	8
1	HYP	F	29[E]	8/9	0.15	0.63	24,24,25,27	8
1	HZP	B	16[G]	8/9	0.06	0.61	3,3,5,8	8
1	HZP	B	16[F]	8/9	0.06	0.61	3,3,5,8	8
1	HYP	I	5[G]	8/9	0.07	0.60	3,3,4,6	8
1	HZP	I	13[G]	8/9	0.06	0.60	3,4,5,7	8
1	HYP	A	5[G]	8/9	0.07	0.57	3,3,4,6	8
1	HYP	A	5[E]	8/9	0.07	0.57	3,3,4,6	8
1	HYP	H	29[C]	8/9	0.10	0.57	8,9,9,12	8
1	HYP	A	5[F]	8/9	0.07	0.57	3,3,4,6	8
1	HYP	A	20[D]	8/9	0.06	0.57	3,3,4,4	8
1	HYP	H	29[D]	8/9	0.10	0.57	8,9,9,12	8
1	HYP	H	5[F]	8/9	0.08	0.56	3,3,5,6	8
1	HYP	H	5[G]	8/9	0.08	0.56	3,3,5,6	8
1	HYP	G	5[A]	8/9	0.13	0.54	8,9,11,13	8
1	HYP	G	5[C]	8/9	0.13	0.54	8,9,11,13	8
1	HYP	G	5[B]	8/9	0.13	0.54	8,9,11,13	8
1	HYP	A	2[E]	8/9	0.10	0.53	6,7,8,10	8
1	HYP	A	2[G]	8/9	0.10	0.53	6,7,8,10	8
1	HYP	A	2[F]	8/9	0.10	0.53	6,7,8,10	8
1	HYP	A	29[F]	8/9	0.12	0.51	16,18,19,21	8
1	HYP	A	29[E]	8/9	0.12	0.51	16,18,19,21	8
1	HYP	A	29[G]	8/9	0.12	0.51	16,18,19,21	8

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(Å ²)	Q<0.9
1	HYP	I	29[G]	8/9	0.10	0.48	11,12,13,15	8
1	HYP	G	29[A]	8/9	0.11	0.47	12,13,14,16	8
1	HYP	G	29[C]	8/9	0.11	0.47	12,13,14,16	8
1	HYP	G	29[B]	8/9	0.11	0.47	12,13,14,16	8
1	HYP	A	20[B]	8/9	0.06	0.44	2,3,3,4	8
1	HYP	A	20[C]	8/9	0.06	0.44	2,3,3,4	8
1	HYP	A	20[A]	8/9	0.06	0.44	2,3,3,4	8
1	HZP	I	16[D]	8/9	0.06	0.43	5,5,7,12	8
1	HZP	I	16[B]	8/9	0.06	0.43	5,5,7,12	8
1	HZP	I	16[C]	8/9	0.06	0.43	5,5,7,12	8
1	HYP	B	29[G]	8/9	0.12	0.43	16,17,18,18	8
1	HZP	D	16[F]	8/9	0.06	0.43	2,3,5,8	8
1	HYP	B	29[F]	8/9	0.12	0.43	16,17,18,18	8
1	HZP	D	16[G]	8/9	0.06	0.43	2,3,5,8	8
1	HZP	D	16[E]	8/9	0.06	0.43	2,3,5,8	8
1	HZP	G	16[F]	8/9	0.06	0.41	3,4,6,6	8
1	HZP	G	16[G]	8/9	0.06	0.41	3,4,6,6	8
1	HZP	G	16[E]	8/9	0.06	0.41	3,4,6,6	8
1	HZP	F	16[E]	8/9	0.06	0.34	4,5,6,13	8
1	HZP	F	16[F]	8/9	0.06	0.34	4,5,6,13	8
1	HZP	F	16[A]	8/9	0.06	0.34	4,5,6,13	8
1	HYP	G	26[D]	8/9	0.06	0.33	3,3,4,6	8
1	HYP	F	26[C]	8/9	0.09	0.31	3,4,5,6	8
1	HYP	I	11[G]	8/9	0.06	0.31	2,3,4,4	8
1	HYP	F	26[B]	8/9	0.09	0.31	3,4,5,6	8
1	HYP	F	26[D]	8/9	0.09	0.31	3,4,5,6	8
1	HYP	B	29[B]	8/9	0.12	0.27	18,19,19,19	8
1	HYP	B	29[E]	8/9	0.12	0.27	18,19,19,19	8
1	HYP	B	29[A]	8/9	0.12	0.27	18,19,19,19	8
1	HZP	I	13[F]	8/9	0.06	0.26	5,5,6,8	8
1	HZP	I	13[A]	8/9	0.06	0.26	5,5,6,8	8
1	HZP	F	16[G]	8/9	0.06	0.26	2,3,4,4	8
1	HZP	I	13[E]	8/9	0.06	0.26	5,5,6,8	8
1	HYP	E	2[B]	8/9	0.12	0.24	14,15,16,16	8
1	HYP	E	2[E]	8/9	0.12	0.24	14,15,16,16	8
1	HYP	E	2[A]	8/9	0.12	0.24	14,15,16,16	8
1	HYP	G	5[D]	8/9	0.13	0.23	21,22,23,24	8
1	HYP	G	29[G]	8/9	0.11	0.22	25,26,29,51	8
1	HZP	G	13[F]	8/9	0.06	0.22	4,5,5,8	8
1	HYP	G	29[E]	8/9	0.11	0.22	25,26,29,51	8
1	HZP	G	13[G]	8/9	0.06	0.22	4,5,5,8	8
1	HYP	C	26[F]	8/9	0.11	0.22	11,12,14,14	8

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(Å ²)	Q<0.9
1	HYP	C	26[E]	8/9	0.11	0.22	11,12,14,14	8
1	HYP	G	29[F]	8/9	0.11	0.22	25,26,29,51	8
1	HZP	G	13[E]	8/9	0.06	0.22	4,5,5,8	8
1	HYP	C	26[A]	8/9	0.11	0.22	11,12,14,14	8
1	HYP	C	2[A]	8/9	0.10	0.21	7,8,10,13	8
1	HYP	C	2[E]	8/9	0.10	0.21	7,8,10,13	8
1	HYP	C	2[F]	8/9	0.10	0.21	7,8,10,13	8
1	HZP	D	16[C]	8/9	0.06	0.20	5,6,7,13	8
1	HZP	D	16[A]	8/9	0.06	0.20	5,6,7,13	8
1	HZP	D	16[B]	8/9	0.06	0.20	5,6,7,13	8
1	HYP	E	2[C]	8/9	0.12	0.19	18,19,21,21	8
1	HYP	E	2[D]	8/9	0.12	0.19	18,19,21,21	8
1	HYP	I	29[D]	8/9	0.10	0.18	10,10,11,14	8
1	HYP	I	29[B]	8/9	0.10	0.18	10,10,11,14	8
1	HYP	I	29[C]	8/9	0.10	0.18	10,10,11,14	8
1	HYP	I	11[E]	8/9	0.06	0.17	4,4,5,5	8
1	HYP	I	11[A]	8/9	0.06	0.17	4,4,5,5	8
1	HYP	I	11[F]	8/9	0.06	0.17	4,4,5,5	8
1	HZP	F	16[B]	8/9	0.06	0.14	5,6,8,12	8
1	HZP	F	16[D]	8/9	0.06	0.14	5,6,8,12	8
1	HZP	F	16[C]	8/9	0.06	0.14	5,6,8,12	8
1	HZP	D	16[D]	8/9	0.06	0.13	2,3,4,7	8
1	HZP	I	13[C]	8/9	0.06	0.11	4,5,6,9	8
1	HYP	E	26[B]	8/9	0.09	0.11	7,8,9,10	8
1	HZP	I	13[B]	8/9	0.06	0.11	4,5,6,9	8
1	HYP	E	26[A]	8/9	0.09	0.11	7,8,9,10	8
1	HZP	I	13[D]	8/9	0.06	0.11	4,5,6,9	8
1	HYP	E	26[E]	8/9	0.09	0.11	7,8,9,10	8
1	HYP	B	2[G]	8/9	0.08	0.07	5,6,7,8	8
1	HYP	B	2[F]	8/9	0.08	0.07	5,6,7,8	8
1	HYP	B	26[D]	8/9	0.07	0.06	4,5,5,7	8
1	HZP	G	16[B]	8/9	0.06	0.06	4,5,7,12	8
1	HYP	B	26[C]	8/9	0.07	0.06	4,5,5,7	8
1	HZP	G	16[A]	8/9	0.06	0.06	4,5,7,12	8
1	HYP	A	26[B]	8/9	0.08	0.06	5,5,6,7	8
1	HZP	G	16[C]	8/9	0.06	0.06	4,5,7,12	8
1	HYP	A	26[C]	8/9	0.08	0.06	5,5,6,7	8
1	HYP	I	26[B]	8/9	0.07	0.06	3,5,6,6	8
1	HYP	I	26[C]	8/9	0.07	0.06	3,5,6,6	8
1	HYP	A	26[A]	8/9	0.08	0.06	5,5,6,7	8
1	HYP	I	26[D]	8/9	0.07	0.06	3,5,6,6	8
1	HYP	E	5[B]	8/9	0.08	0.05	6,6,7,9	8

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
1	HYP	E	5[A]	8/9	0.08	0.05	6,6,7,9	8
1	HYP	E	5[E]	8/9	0.08	0.05	6,6,7,9	8
1	HYP	F	5[E]	8/9	0.07	0.03	5,5,7,8	8
1	HYP	F	5[A]	8/9	0.07	0.03	5,5,7,8	8
1	HYP	F	5[F]	8/9	0.07	0.03	5,5,7,8	8
1	HZP	G	13[A]	8/9	0.06	-0.01	4,5,7,11	8
1	HZP	G	13[B]	8/9	0.06	-0.01	4,5,7,11	8
1	HZP	G	13[C]	8/9	0.06	-0.01	4,5,7,11	8
1	HYP	F	23[C]	8/9	0.06	-0.02	3,4,5,5	8
1	HYP	F	23[D]	8/9	0.06	-0.02	3,4,5,5	8
1	HYP	F	23[B]	8/9	0.06	-0.02	3,4,5,5	8
1	HYP	E	8[F]	8/9	0.06	-0.04	3,3,5,8	8
1	HYP	E	8[G]	8/9	0.06	-0.04	3,3,5,8	8
1	HYP	D	26[C]	8/9	0.08	-0.05	5,6,7,10	8
1	HYP	D	26[A]	8/9	0.08	-0.05	5,6,7,10	8
1	HYP	D	26[B]	8/9	0.08	-0.05	5,6,7,10	8
1	HYP	F	11[G]	8/9	0.06	-0.10	3,3,3,4	8
1	HZP	G	16[D]	8/9	0.06	-0.14	4,5,7,11	8
1	HYP	C	23[D]	8/9	0.06	-0.14	3,3,4,5	8
1	HYP	C	23[C]	8/9	0.06	-0.14	3,3,4,5	8
1	HYP	C	23[B]	8/9	0.06	-0.14	3,3,4,5	8
1	HYP	F	11[B]	8/9	0.06	-0.15	2,3,3,5	8
1	HYP	F	11[D]	8/9	0.06	-0.15	2,3,3,5	8
1	HYP	C	26[G]	8/9	0.11	-0.15	20,20,21,22	8
1	HYP	H	5[A]	8/9	0.08	-0.15	6,7,7,9	8
1	HYP	I	2[A]	8/9	0.11	-0.15	13,15,16,19	8
1	HYP	H	5[B]	8/9	0.08	-0.15	6,7,7,9	8
1	HYP	H	5[E]	8/9	0.08	-0.15	6,7,7,9	8
1	HYP	I	2[F]	8/9	0.11	-0.15	13,15,16,19	8
1	HYP	I	2[E]	8/9	0.11	-0.15	13,15,16,19	8
1	HYP	F	11[C]	8/9	0.06	-0.15	2,3,3,5	8
1	HYP	G	8[E]	8/9	0.06	-0.16	2,3,4,5	8
1	HYP	G	8[F]	8/9	0.06	-0.16	2,3,4,5	8
1	HYP	G	8[G]	8/9	0.06	-0.16	2,3,4,5	8
1	HYP	E	11[B]	8/9	0.05	-0.18	3,4,5,8	8
1	HYP	E	11[E]	8/9	0.05	-0.18	3,4,5,8	8
1	HYP	E	11[A]	8/9	0.05	-0.18	3,4,5,8	8
1	HYP	I	29[A]	8/9	0.10	-0.21	20,22,23,24	8
1	HYP	F	11[F]	8/9	0.06	-0.21	3,4,5,6	8
1	HYP	H	29[B]	8/9	0.10	-0.21	14,15,17,17	8
1	HYP	H	29[A]	8/9	0.10	-0.21	14,15,17,17	8
1	HYP	I	29[F]	8/9	0.10	-0.21	20,22,23,24	8

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
1	HYP	I	29[E]	8/9	0.10	-0.21	20,22,23,24	8
1	HYP	F	11[E]	8/9	0.06	-0.21	3,4,5,6	8
1	HYP	H	29[E]	8/9	0.10	-0.21	14,15,17,17	8
1	HYP	F	11[A]	8/9	0.06	-0.21	3,4,5,6	8
1	HYP	A	2[D]	8/9	0.10	-0.21	14,15,17,17	8
1	HYP	I	5[A]	8/9	0.07	-0.24	3,4,6,7	8
1	HYP	I	5[F]	8/9	0.07	-0.24	3,4,6,7	8
1	HYP	I	5[E]	8/9	0.07	-0.24	3,4,6,7	8
1	HYP	A	20[F]	8/9	0.06	-0.26	4,5,7,9	8
1	HYP	A	20[E]	8/9	0.06	-0.26	4,5,7,9	8
1	HYP	A	23[D]	8/9	0.05	-0.26	2,3,3,4	8
1	HYP	A	20[G]	8/9	0.06	-0.26	4,5,7,9	8
1	HZP	G	13[D]	8/9	0.06	-0.27	2,3,3,4	8
1	HZP	C	13[G]	8/9	0.05	-0.28	3,3,4,7	8
1	HZP	B	13[G]	8/9	0.05	-0.30	3,4,7,12	8
1	HYP	A	2[B]	8/9	0.10	-0.30	18,20,20,21	8
1	HYP	A	2[C]	8/9	0.10	-0.30	18,20,20,21	8
1	HZP	B	13[F]	8/9	0.05	-0.30	3,4,7,12	8
1	HYP	A	2[A]	8/9	0.10	-0.30	18,20,20,21	8
1	HYP	I	11[C]	8/9	0.06	-0.32	3,4,4,6	8
1	HYP	I	11[B]	8/9	0.06	-0.32	3,4,4,6	8
1	HYP	I	11[D]	8/9	0.06	-0.32	3,4,4,6	8
1	HZP	H	13[G]	8/9	0.05	-0.36	4,5,7,11	8
1	HZP	H	13[F]	8/9	0.05	-0.36	4,5,7,11	8
1	HYP	D	5[A]	8/9	0.08	-0.38	7,8,9,11	8
1	HYP	H	29[F]	8/9	0.10	-0.38	16,17,19,21	8
1	HYP	D	5[B]	8/9	0.08	-0.38	7,8,9,11	8
1	HYP	H	29[G]	8/9	0.10	-0.38	16,17,19,21	8
1	HYP	D	5[C]	8/9	0.08	-0.38	7,8,9,11	8
1	HYP	E	8[E]	8/9	0.06	-0.39	2,3,5,6	8
1	HYP	E	8[B]	8/9	0.06	-0.39	2,3,5,6	8
1	HYP	E	8[A]	8/9	0.06	-0.39	2,3,5,6	8
1	HZP	H	13[E]	8/9	0.05	-0.41	4,4,6,12	8
1	HZP	H	13[B]	8/9	0.05	-0.41	4,4,6,12	8
1	HYP	A	5[C]	8/9	0.07	-0.41	6,7,8,9	8
1	HYP	A	5[A]	8/9	0.07	-0.41	6,7,8,9	8
1	HYP	A	5[B]	8/9	0.07	-0.41	6,7,8,9	8
1	HYP	G	20[D]	8/9	0.05	-0.41	2,3,4,4	8
1	HZP	H	13[A]	8/9	0.05	-0.41	4,4,6,12	8
1	HYP	E	11[F]	8/9	0.05	-0.41	3,3,4,6	8
1	HYP	E	11[G]	8/9	0.05	-0.41	3,3,4,6	8
1	HYP	E	20[G]	8/9	0.06	-0.42	3,3,4,5	8

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(Å ²)	Q<0.9
1	HYP	E	20[F]	8/9	0.06	-0.42	3,3,4,5	8
1	HZP	C	13[A]	8/9	0.05	-0.43	3,4,5,9	8
1	HZP	C	13[E]	8/9	0.05	-0.43	3,4,5,9	8
1	HZP	C	13[F]	8/9	0.05	-0.43	3,4,5,9	8
1	HYP	B	2[E]	8/9	0.08	-0.44	11,12,13,14	8
1	HYP	B	2[A]	8/9	0.08	-0.44	11,12,13,14	8
1	HYP	B	2[B]	8/9	0.08	-0.44	11,12,13,14	8
1	HZP	A	13[F]	8/9	0.05	-0.45	4,4,7,10	8
1	HYP	C	11[G]	8/9	0.05	-0.45	3,3,5,6	8
1	HYP	E	11[C]	8/9	0.05	-0.45	2,3,5,6	8
1	HZP	A	13[E]	8/9	0.05	-0.45	4,4,7,10	8
1	HYP	E	11[D]	8/9	0.05	-0.45	2,3,5,6	8
1	HZP	A	13[G]	8/9	0.05	-0.45	4,4,7,10	8
1	HZP	C	13[B]	8/9	0.05	-0.46	3,3,4,7	8
1	HYP	G	20[E]	8/9	0.05	-0.46	3,3,5,5	8
1	HZP	C	13[D]	8/9	0.05	-0.46	3,3,4,7	8
1	HZP	B	13[D]	8/9	0.05	-0.46	2,3,3,4	8
1	HYP	G	20[G]	8/9	0.05	-0.46	3,3,5,5	8
1	HZP	B	13[C]	8/9	0.05	-0.46	2,3,3,4	8
1	HZP	C	13[C]	8/9	0.05	-0.46	3,3,4,7	8
1	HYP	G	20[F]	8/9	0.05	-0.46	3,3,5,5	8
1	HZP	B	13[E]	8/9	0.05	-0.47	3,4,6,9	8
1	HZP	B	13[B]	8/9	0.05	-0.47	3,4,6,9	8
1	HYP	G	20[A]	8/9	0.05	-0.47	3,3,5,5	8
1	HYP	G	20[B]	8/9	0.05	-0.47	3,3,5,5	8
1	HYP	G	20[C]	8/9	0.05	-0.47	3,3,5,5	8
1	HZP	B	13[A]	8/9	0.05	-0.47	3,4,6,9	8
1	HYP	I	2[D]	8/9	0.11	-0.49	20,22,22,22	8
1	HYP	I	2[B]	8/9	0.11	-0.49	20,22,22,22	8
1	HYP	I	2[C]	8/9	0.11	-0.49	20,22,22,22	8
1	HYP	E	23[C]	8/9	0.05	-0.50	3,4,6,7	8
1	HYP	E	23[D]	8/9	0.05	-0.50	3,4,6,7	8
1	HZP	H	16[G]	8/9	0.05	-0.52	2,4,6,10	8
1	HZP	H	16[F]	8/9	0.05	-0.52	2,4,6,10	8
1	HZP	H	16[B]	8/9	0.05	-0.52	4,4,6,9	8
1	HZP	H	16[E]	8/9	0.05	-0.52	4,4,6,9	8
1	HZP	H	16[A]	8/9	0.05	-0.52	4,4,6,9	8
1	HYP	F	23[F]	8/9	0.06	-0.53	3,4,5,6	8
1	HYP	F	26[A]	8/9	0.09	-0.53	10,12,13,13	8
1	HZP	H	16[C]	8/9	0.05	-0.53	4,5,7,12	8
1	HYP	F	23[A]	8/9	0.06	-0.53	3,4,5,6	8
1	HZP	H	16[D]	8/9	0.05	-0.53	4,5,7,12	8

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(Å ²)	Q<0.9
1	HYP	G	26[B]	8/9	0.06	-0.53	6,7,8,11	8
1	HYP	G	26[C]	8/9	0.06	-0.53	6,7,8,11	8
1	HYP	F	23[E]	8/9	0.06	-0.53	3,4,5,6	8
1	HYP	G	26[A]	8/9	0.06	-0.53	6,7,8,11	8
1	HYP	F	26[E]	8/9	0.09	-0.53	10,12,13,13	8
1	HYP	F	26[F]	8/9	0.09	-0.53	10,12,13,13	8
1	HYP	C	2[D]	8/9	0.10	-0.56	25,26,29,51	8
1	HYP	C	2[C]	8/9	0.10	-0.56	25,26,29,51	8
1	HYP	C	2[B]	8/9	0.10	-0.56	25,26,29,51	8
1	HYP	H	26[D]	8/9	0.06	-0.58	3,4,4,6	8
1	HYP	H	26[C]	8/9	0.06	-0.58	3,4,4,6	8
1	HYP	F	20[A]	8/9	0.05	-0.59	3,4,5,5	8
1	HYP	C	16[E]	8/9	0.05	-0.59	3,4,6,11	8
1	HYP	F	20[E]	8/9	0.05	-0.59	3,4,5,5	8
1	HYP	C	16[F]	8/9	0.05	-0.59	3,4,6,11	8
1	HYP	F	20[F]	8/9	0.05	-0.59	3,4,5,5	8
1	HYP	C	16[A]	8/9	0.05	-0.59	3,4,6,11	8
1	HYP	E	26[F]	8/9	0.09	-0.60	20,21,22,22	8
1	HYP	E	26[G]	8/9	0.09	-0.60	20,21,22,22	8
1	HYP	G	8[C]	8/9	0.06	-0.60	3,4,5,6	8
1	HYP	G	8[B]	8/9	0.06	-0.60	3,4,5,6	8
1	HYP	G	8[A]	8/9	0.06	-0.60	3,4,5,6	8
1	HYP	A	13[D]	8/9	0.05	-0.61	3,3,5,6	8
1	HYP	H	13[C]	8/9	0.05	-0.61	2,2,3,6	8
1	HYP	B	5[F]	8/9	0.05	-0.61	3,4,6,7	8
1	HYP	B	5[G]	8/9	0.05	-0.61	3,4,6,7	8
1	HYP	H	13[D]	8/9	0.05	-0.61	2,2,3,6	8
1	HYP	A	23[C]	8/9	0.05	-0.62	4,5,6,9	8
1	HYP	A	23[B]	8/9	0.05	-0.62	4,5,6,9	8
1	HYP	A	23[A]	8/9	0.05	-0.62	4,5,6,9	8
1	HYP	C	11[A]	8/9	0.05	-0.66	3,3,5,7	8
1	HYP	C	11[F]	8/9	0.05	-0.66	3,3,5,7	8
1	HYP	C	11[E]	8/9	0.05	-0.66	3,3,5,7	8
1	HYP	F	20[C]	8/9	0.05	-0.67	2,3,3,4	8
1	HYP	F	20[D]	8/9	0.05	-0.67	2,3,3,4	8
1	HYP	F	20[B]	8/9	0.05	-0.67	2,3,3,4	8
1	HYP	A	26[G]	8/9	0.08	-0.68	14,14,14,15	8
1	HYP	A	26[E]	8/9	0.08	-0.68	14,14,14,15	8
1	HYP	A	26[F]	8/9	0.08	-0.68	14,14,14,15	8
1	HYP	G	11[E]	8/9	0.05	-0.69	2,3,4,4	8
1	HYP	G	11[F]	8/9	0.05	-0.69	2,3,4,4	8
1	HYP	G	11[G]	8/9	0.05	-0.69	2,3,4,4	8

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(Å ²)	Q<0.9
1	HYP	G	23[C]	8/9	0.05	-0.70	3,3,4,6	8
1	HZP	C	16[G]	8/9	0.05	-0.70	2,2,3,3	8
1	HYP	G	23[B]	8/9	0.05	-0.70	3,3,4,6	8
1	HYP	G	23[A]	8/9	0.05	-0.70	3,3,4,6	8
1	HZP	A	13[B]	8/9	0.05	-0.71	3,3,7,11	8
1	HZP	A	13[A]	8/9	0.05	-0.71	3,3,7,11	8
1	HZP	A	13[C]	8/9	0.05	-0.71	3,3,7,11	8
1	HYP	B	26[B]	8/9	0.07	-0.72	6,7,8,10	8
1	HYP	D	20[F]	8/9	0.05	-0.72	3,3,4,6	8
1	HYP	F	5[C]	8/9	0.07	-0.72	12,13,14,16	8
1	HYP	F	5[B]	8/9	0.07	-0.72	12,13,14,16	8
1	HYP	B	26[E]	8/9	0.07	-0.72	6,7,8,10	8
1	HYP	G	11[A]	8/9	0.05	-0.72	2,3,4,5	8
1	HYP	G	11[C]	8/9	0.05	-0.72	2,3,4,5	8
1	HYP	F	5[D]	8/9	0.07	-0.72	12,13,14,16	8
1	HYP	G	11[B]	8/9	0.05	-0.72	2,3,4,5	8
1	HYP	D	20[G]	8/9	0.05	-0.72	3,3,4,6	8
1	HYP	B	26[A]	8/9	0.07	-0.72	6,7,8,10	8
1	HYP	F	20[G]	8/9	0.05	-0.72	3,3,4,6	8
1	HYP	D	20[E]	8/9	0.05	-0.72	3,3,4,6	8
1	HZP	E	16[C]	8/9	0.05	-0.75	4,4,7,9	8
1	HZP	E	16[D]	8/9	0.05	-0.75	4,4,7,9	8
1	HZP	C	16[B]	8/9	0.05	-0.76	3,4,5,9	8
1	HZP	C	16[D]	8/9	0.05	-0.76	3,4,5,9	8
1	HZP	C	16[C]	8/9	0.05	-0.76	3,4,5,9	8
1	HYP	E	23[E]	8/9	0.05	-0.79	3,3,4,5	8
1	HYP	E	23[A]	8/9	0.05	-0.79	3,3,4,5	8
1	HYP	E	23[B]	8/9	0.05	-0.79	3,3,4,5	8
1	HYP	D	26[E]	8/9	0.08	-0.81	14,15,16,16	8
1	HYP	D	26[F]	8/9	0.08	-0.81	14,15,16,16	8
1	HYP	D	26[G]	8/9	0.08	-0.81	14,15,16,16	8
1	HZP	F	13[C]	8/9	0.05	-0.82	3,3,4,6	8
1	HZP	F	13[D]	8/9	0.05	-0.82	3,3,4,6	8
1	HYP	D	8[B]	8/9	0.05	-0.82	4,4,5,6	8
1	HYP	D	8[C]	8/9	0.05	-0.82	4,4,5,6	8
1	HZP	F	13[B]	8/9	0.05	-0.82	3,3,4,6	8
1	HYP	D	8[A]	8/9	0.05	-0.82	4,4,5,6	8
1	HYP	H	5[D]	8/9	0.08	-0.82	16,17,18,20	8
1	HYP	H	5[C]	8/9	0.08	-0.82	16,17,18,20	8
1	HYP	D	8[G]	8/9	0.05	-0.83	2,3,4,5	8
1	HYP	D	8[E]	8/9	0.05	-0.83	2,3,4,5	8
1	HYP	D	8[F]	8/9	0.05	-0.83	2,3,4,5	8

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(Å ²)	Q<0.9
1	HYP	C	23[F]	8/9	0.06	-0.84	5,5,6,8	8
1	HYP	I	26[E]	8/9	0.07	-0.84	10,11,11,13	8
1	HZP	E	13[D]	8/9	0.05	-0.84	2,3,6,10	8
1	HYP	C	23[A]	8/9	0.06	-0.84	5,5,6,8	8
1	HZP	E	13[C]	8/9	0.05	-0.84	2,3,6,10	8
1	HYP	C	23[E]	8/9	0.06	-0.84	5,5,6,8	8
1	HYP	I	26[F]	8/9	0.07	-0.84	10,11,11,13	8
1	HYP	G	11[D]	8/9	0.05	-0.84	3,3,4,5	8
1	HYP	I	26[A]	8/9	0.07	-0.84	10,11,11,13	8
1	HYP	E	5[D]	8/9	0.08	-0.85	15,15,16,16	8
1	HYP	B	20[C]	8/9	0.05	-0.85	2,2,2,3	8
1	HYP	B	20[D]	8/9	0.05	-0.85	2,2,2,3	8
1	HYP	E	5[C]	8/9	0.08	-0.85	15,15,16,16	8
1	HYP	E	8[D]	8/9	0.06	-0.86	6,6,7,9	8
1	HYP	E	8[C]	8/9	0.06	-0.86	6,6,7,9	8
1	HZP	D	13[D]	8/9	0.05	-0.87	2,2,3,4	8
1	HYP	C	11[C]	8/9	0.05	-0.88	3,4,5,7	8
1	HYP	C	11[D]	8/9	0.05	-0.88	3,4,5,7	8
1	HYP	C	11[B]	8/9	0.05	-0.88	3,4,5,7	8
1	HYP	B	11[C]	8/9	0.05	-0.89	3,4,6,6	8
1	HYP	B	11[D]	8/9	0.05	-0.89	3,4,6,6	8
1	HZP	E	16[G]	8/9	0.05	-0.91	3,4,4,7	8
1	HZP	E	16[F]	8/9	0.05	-0.91	3,4,4,7	8
1	HYP	B	5[E]	8/9	0.05	-0.92	5,6,7,8	8
1	HZP	A	16[E]	8/9	0.05	-0.92	3,4,5,9	8
1	HYP	B	5[B]	8/9	0.05	-0.92	5,6,7,8	8
1	HYP	B	5[A]	8/9	0.05	-0.92	5,6,7,8	8
1	HZP	A	16[G]	8/9	0.05	-0.92	3,4,5,9	8
1	HZP	A	16[F]	8/9	0.05	-0.92	3,4,5,9	8
1	HYP	I	5[C]	8/9	0.07	-0.93	14,16,16,18	8
1	HYP	I	5[D]	8/9	0.07	-0.93	14,16,16,18	8
1	HYP	I	5[B]	8/9	0.07	-0.93	14,16,16,18	8
1	HYP	H	11[D]	8/9	0.05	-0.94	3,3,5,5	8
1	HYP	H	11[C]	8/9	0.05	-0.94	3,3,5,5	8
1	HZP	D	13[A]	8/9	0.05	-0.95	2,3,4,7	8
1	HZP	D	13[C]	8/9	0.05	-0.95	2,3,4,7	8
1	HZP	D	13[B]	8/9	0.05	-0.95	2,3,4,7	8
1	HYP	B	20[A]	8/9	0.05	-0.96	3,3,4,4	8
1	HYP	B	23[C]	8/9	0.05	-0.96	3,3,4,4	8
1	HYP	B	23[D]	8/9	0.05	-0.96	3,3,4,4	8
1	HYP	B	20[B]	8/9	0.05	-0.96	3,3,4,4	8
1	HYP	B	20[E]	8/9	0.05	-0.96	3,3,4,4	8

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(Å ²)	Q<0.9
1	HYP	F	26[G]	8/9	0.09	-1.02	21,22,22,22	8
1	HYP	F	8[C]	8/9	0.04	-1.03	5,5,7,8	8
1	HYP	B	23[A]	8/9	0.05	-1.03	4,5,5,7	8
1	HYP	B	23[E]	8/9	0.05	-1.03	4,5,5,7	8
1	HYP	I	8[D]	8/9	0.04	-1.03	3,4,6,7	8
1	HYP	F	8[B]	8/9	0.04	-1.03	5,5,7,8	8
1	HYP	F	8[D]	8/9	0.04	-1.03	5,5,7,8	8
1	HYP	I	8[C]	8/9	0.04	-1.03	3,4,6,7	8
1	HYP	B	23[B]	8/9	0.05	-1.03	4,5,5,7	8
1	HYP	D	5[D]	8/9	0.08	-1.03	16,17,18,18	8
1	HYP	I	8[B]	8/9	0.04	-1.03	3,4,6,7	8
1	HYP	B	20[G]	8/9	0.05	-1.03	4,5,5,7	8
1	HYP	B	20[F]	8/9	0.05	-1.03	4,5,5,7	8
1	HYP	D	11[D]	8/9	0.05	-1.05	4,4,5,6	8
1	HYP	A	23[E]	8/9	0.05	-1.06	5,5,6,7	8
1	HYP	A	23[F]	8/9	0.05	-1.06	5,5,6,7	8
1	HYP	A	23[G]	8/9	0.05	-1.06	5,5,6,7	8
1	HYP	G	23[E]	8/9	0.05	-1.07	6,7,8,11	8
1	HYP	G	23[F]	8/9	0.05	-1.07	6,7,8,11	8
1	HYP	G	23[G]	8/9	0.05	-1.07	6,7,8,11	8
1	HZP	A	16[B]	8/9	0.05	-1.08	3,4,7,10	8
1	HZP	A	16[C]	8/9	0.05	-1.08	3,4,7,10	8
1	HZP	A	16[A]	8/9	0.05	-1.08	3,4,7,10	8
1	HYP	G	8[D]	8/9	0.06	-1.10	8,9,10,10	8
1	HZP	E	13[A]	8/9	0.05	-1.12	4,5,7,9	8
1	HZP	E	13[E]	8/9	0.05	-1.12	4,5,7,9	8
1	HZP	E	13[B]	8/9	0.05	-1.12	4,5,7,9	8
1	HYP	I	23[C]	8/9	0.05	-1.13	2,3,4,5	8
1	HYP	I	23[F]	8/9	0.05	-1.13	3,5,6,6	8
1	HYP	I	23[D]	8/9	0.05	-1.13	2,3,4,5	8
1	HYP	I	23[A]	8/9	0.05	-1.13	3,5,6,6	8
1	HYP	I	23[B]	8/9	0.05	-1.13	2,3,4,5	8
1	HYP	I	23[E]	8/9	0.05	-1.13	3,5,6,6	8
1	HYP	E	23[G]	8/9	0.05	-1.14	7,8,9,10	8
1	HYP	B	8[C]	8/9	0.04	-1.14	5,6,7,8	8
1	HYP	B	8[D]	8/9	0.04	-1.14	5,6,7,8	8
1	HYP	E	23[F]	8/9	0.05	-1.14	7,8,9,10	8
1	HYP	F	23[G]	8/9	0.06	-1.14	10,10,12,12	8
1	HYP	G	23[D]	8/9	0.05	-1.15	3,3,5,5	8
1	HYP	I	8[E]	8/9	0.04	-1.17	3,3,4,6	8
1	HYP	I	8[A]	8/9	0.04	-1.17	3,3,4,6	8
1	HYP	I	8[F]	8/9	0.04	-1.17	3,3,4,6	8

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
1	HYP	B	8[E]	8/9	0.04	-1.17	3,4,6,7	8
1	HYP	B	8[B]	8/9	0.04	-1.17	3,4,6,7	8
1	HYP	B	8[A]	8/9	0.04	-1.17	3,4,6,7	8
1	HYP	H	11[A]	8/9	0.05	-1.18	4,4,5,7	8
1	HYP	C	5[F]	8/9	0.05	-1.18	4,4,4,5	8
1	HYP	H	26[E]	8/9	0.06	-1.18	8,9,9,12	8
1	HYP	H	11[E]	8/9	0.05	-1.18	4,4,5,7	8
1	HYP	C	5[E]	8/9	0.05	-1.18	4,4,4,5	8
1	HYP	H	26[B]	8/9	0.06	-1.18	8,9,9,12	8
1	HYP	H	26[A]	8/9	0.06	-1.18	8,9,9,12	8
1	HYP	C	5[A]	8/9	0.05	-1.18	4,4,4,5	8
1	HYP	H	11[B]	8/9	0.05	-1.18	4,4,5,7	8
1	HYP	H	23[A]	8/9	0.05	-1.20	3,4,4,6	8
1	HYP	H	20[F]	8/9	0.05	-1.20	3,4,4,6	8
1	HYP	H	23[B]	8/9	0.05	-1.20	3,4,4,6	8
1	HYP	H	20[G]	8/9	0.05	-1.20	3,4,4,6	8
1	HYP	H	23[E]	8/9	0.05	-1.20	3,4,4,6	8
1	HZP	E	16[E]	8/9	0.05	-1.21	5,5,7,11	8
1	HZP	E	16[B]	8/9	0.05	-1.21	5,5,7,11	8
1	HZP	E	16[A]	8/9	0.05	-1.21	5,5,7,11	8
1	HYP	D	23[F]	8/9	0.04	-1.22	5,6,7,10	8
1	HYP	D	23[E]	8/9	0.04	-1.22	5,6,7,10	8
1	HYP	D	23[G]	8/9	0.04	-1.22	5,6,7,10	8
1	HYP	A	11[D]	8/9	0.04	-1.25	3,3,4,6	8
1	HYP	D	20[A]	8/9	0.05	-1.27	3,4,4,5	8
1	HZP	F	13[E]	8/9	0.05	-1.27	5,6,8,12	8
1	HZP	F	13[F]	8/9	0.05	-1.27	5,6,8,12	8
1	HYP	D	20[B]	8/9	0.05	-1.27	3,4,4,5	8
1	HYP	D	20[C]	8/9	0.05	-1.27	3,4,4,5	8
1	HZP	D	13[F]	8/9	0.05	-1.27	4,5,5,8	8
1	HZP	D	13[E]	8/9	0.05	-1.27	4,5,5,8	8
1	HZP	F	13[A]	8/9	0.05	-1.27	5,6,8,12	8
1	HZP	D	13[G]	8/9	0.05	-1.27	4,5,5,8	8
1	HYP	D	8[D]	8/9	0.05	-1.28	7,8,9,11	8
1	HYP	D	23[A]	8/9	0.04	-1.29	3,3,4,6	8
1	HZP	F	13[G]	8/9	0.05	-1.29	4,4,5,5	8
1	HYP	D	23[C]	8/9	0.04	-1.29	3,3,4,6	8
1	HYP	D	23[B]	8/9	0.04	-1.29	3,3,4,6	8
1	HYP	A	8[B]	8/9	0.04	-1.31	3,3,4,6	8
1	HYP	A	8[C]	8/9	0.04	-1.31	3,3,4,6	8
1	HYP	A	8[A]	8/9	0.04	-1.31	3,3,4,6	8
1	HYP	H	11[G]	8/9	0.05	-1.32	3,3,4,4	8

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
1	HYP	H	11[F]	8/9	0.05	-1.32	3,3,4,4	8
1	HYP	H	8[C]	8/9	0.04	-1.33	6,7,7,9	8
1	HYP	H	8[D]	8/9	0.04	-1.33	6,7,7,9	8
1	HYP	F	8[E]	8/9	0.04	-1.36	2,3,4,6	8
1	HYP	F	8[A]	8/9	0.04	-1.36	2,3,4,6	8
1	HYP	F	8[F]	8/9	0.04	-1.36	2,3,4,6	8
1	HYP	D	11[C]	8/9	0.05	-1.37	2,3,4,6	8
1	HYP	D	11[A]	8/9	0.05	-1.37	2,3,4,6	8
1	HYP	D	20[D]	8/9	0.05	-1.37	3,3,4,4	8
1	HYP	A	5[D]	8/9	0.07	-1.37	18,20,20,21	8
1	HYP	D	11[B]	8/9	0.05	-1.37	2,3,4,6	8
1	HYP	C	8[D]	8/9	0.04	-1.39	4,4,4,5	8
1	HYP	C	8[C]	8/9	0.04	-1.39	4,4,4,5	8
1	HYP	C	8[B]	8/9	0.04	-1.39	4,4,4,5	8
1	HYP	I	20[G]	8/9	0.04	-1.40	4,5,6,6	8
1	HYP	B	23[F]	8/9	0.05	-1.44	6,7,8,10	8
1	HYP	B	23[G]	8/9	0.05	-1.44	6,7,8,10	8
1	HYP	A	16[D]	8/9	0.05	-1.44	3,3,5,6	8
1	HYP	B	11[G]	8/9	0.05	-1.45	3,4,5,6	8
1	HYP	B	11[F]	8/9	0.05	-1.45	3,4,5,6	8
1	HYP	H	8[E]	8/9	0.04	-1.47	3,3,5,6	8
1	HYP	H	8[B]	8/9	0.04	-1.47	3,3,5,6	8
1	HYP	H	8[A]	8/9	0.04	-1.47	3,3,5,6	8
1	HYP	G	26[E]	8/9	0.06	-1.48	12,13,14,15	8
1	HYP	G	26[F]	8/9	0.06	-1.48	12,13,14,15	8
1	HYP	G	26[G]	8/9	0.06	-1.48	12,13,14,15	8
1	HYP	C	5[C]	8/9	0.05	-1.48	7,8,10,12	8
1	HYP	C	5[B]	8/9	0.05	-1.48	7,8,10,12	8
1	HYP	C	5[D]	8/9	0.05	-1.48	7,8,10,12	8
1	HYP	C	5[G]	8/9	0.05	-1.49	3,4,5,6	8
1	HYP	C	20[G]	8/9	0.04	-1.51	5,5,6,8	8
1	HYP	B	11[B]	8/9	0.05	-1.52	2,3,4,5	8
1	HYP	B	11[E]	8/9	0.05	-1.52	2,3,4,5	8
1	HYP	C	20[A]	8/9	0.04	-1.52	3,3,4,5	8
1	HYP	C	20[E]	8/9	0.04	-1.52	3,3,4,5	8
1	HYP	B	11[A]	8/9	0.05	-1.52	2,3,4,5	8
1	HYP	C	20[F]	8/9	0.04	-1.52	3,3,4,5	8
1	HYP	A	8[D]	8/9	0.04	-1.53	6,7,9,10	8
1	HYP	C	23[G]	8/9	0.06	-1.56	11,12,14,16	8
1	HYP	H	23[F]	8/9	0.05	-1.57	8,9,9,12	8
1	HYP	H	23[G]	8/9	0.05	-1.57	8,9,9,12	8
1	HYP	B	5[D]	8/9	0.05	-1.57	11,12,13,14	8

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
1	HYP	I	23[G]	8/9	0.05	-1.66	10,11,11,13	8
1	HYP	B	2[C]	8/9	0.08	-1.67	19,20,21,21	8
1	HYP	B	2[D]	8/9	0.08	-1.67	19,20,21,21	8
1	HYP	I	8[G]	8/9	0.04	-1.70	4,4,5,6	8
1	HYP	D	11[G]	8/9	0.05	-1.71	4,4,6,10	8
1	HYP	D	11[E]	8/9	0.05	-1.71	4,4,6,10	8
1	HYP	D	11[F]	8/9	0.05	-1.71	4,4,6,10	8
1	HZP	E	13[F]	8/9	0.05	-1.72	4,5,5,9	8
1	HZP	E	13[G]	8/9	0.05	-1.72	4,5,5,9	8
1	HYP	B	5[C]	8/9	0.05	-1.75	11,12,13,14	8
1	HYP	B	26[F]	8/9	0.07	-1.77	18,19,20,20	8
1	HYP	B	26[G]	8/9	0.07	-1.77	18,19,20,20	8
1	HYP	C	20[D]	8/9	0.04	-1.79	2,3,4,4	8
1	HYP	C	20[B]	8/9	0.04	-1.79	2,3,4,4	8
1	HYP	C	20[C]	8/9	0.04	-1.79	2,3,4,4	8
1	HYP	A	11[B]	8/9	0.04	-1.82	2,3,4,6	8
1	HYP	I	20[A]	8/9	0.04	-1.82	3,3,4,5	8
1	HYP	I	20[E]	8/9	0.04	-1.82	3,3,4,5	8
1	HYP	A	11[C]	8/9	0.04	-1.82	2,3,4,6	8
1	HYP	A	11[A]	8/9	0.04	-1.82	2,3,4,6	8
1	HYP	I	20[F]	8/9	0.04	-1.82	3,3,4,5	8
1	HYP	I	20[B]	8/9	0.04	-1.83	4,5,7,8	8
1	HYP	I	20[C]	8/9	0.04	-1.83	4,5,7,8	8
1	HYP	I	20[D]	8/9	0.04	-1.83	4,5,7,8	8
1	HYP	F	8[G]	8/9	0.04	-1.86	3,4,4,6	8
1	HYP	A	8[F]	8/9	0.04	-1.98	2,3,4,6	8
1	HYP	A	8[E]	8/9	0.04	-1.98	2,3,4,6	8
1	HYP	A	8[G]	8/9	0.04	-1.98	2,3,4,6	8
1	HYP	C	8[A]	8/9	0.04	-2.01	3,4,5,6	8
1	HYP	C	8[F]	8/9	0.04	-2.01	3,4,5,6	8
1	HYP	C	8[E]	8/9	0.04	-2.01	3,4,5,6	8
1	HYP	H	20[D]	8/9	0.05	-2.05	3,4,5,6	8
1	HYP	H	20[C]	8/9	0.05	-2.05	3,4,5,6	8
1	HYP	C	8[G]	8/9	0.04	-2.11	2,3,5,7	8
1	HYP	B	8[F]	8/9	0.04	-2.12	2,3,4,5	8
1	HYP	I	26[G]	8/9	0.07	-2.12	10,18,20,20	8
1	HYP	B	8[G]	8/9	0.04	-2.12	2,3,4,5	8
1	HYP	A	11[G]	8/9	0.04	-2.15	4,4,6,11	8
1	HYP	A	11[E]	8/9	0.04	-2.15	4,4,6,11	8
1	HYP	A	11[F]	8/9	0.04	-2.15	4,4,6,11	8
1	HYP	H	26[F]	8/9	0.06	-2.20	14,15,16,16	8
1	HYP	H	26[G]	8/9	0.06	-2.20	14,15,16,16	8

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
1	HYP	H	20[B]	8/9	0.05	-2.29	3,3,4,5	8
1	HYP	H	23[C]	8/9	0.05	-2.29	3,3,4,5	8
1	HYP	H	20[A]	8/9	0.05	-2.29	3,3,4,5	8
1	HYP	H	20[E]	8/9	0.05	-2.29	3,3,4,5	8
1	HYP	H	23[D]	8/9	0.05	-2.29	3,3,4,5	8
1	HYP	H	8[F]	8/9	0.04	-2.59	3,4,5,7	8
1	HYP	H	8[G]	8/9	0.04	-2.59	3,4,5,7	8
1	HYP	D	23[D]	8/9	0.04	-2.96	3,4,4,5	8

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.