



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 03:49 PM UTC

PDB ID : 2UUM / pdb_00002uum
Title : Crystal structure of C-phycoyanin from Phormidium, Lyngbya spp. (Marine) and Spirulina sp. (Fresh water) shows two different ways of energy transfer between two hexamers.
Authors : Satyanarayana, L.; Patel, A.; Mishra, S.; K Ghosh, P.; Suresh, C.G.
Deposited on : 2007-03-04
Resolution : 3.00 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtrriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

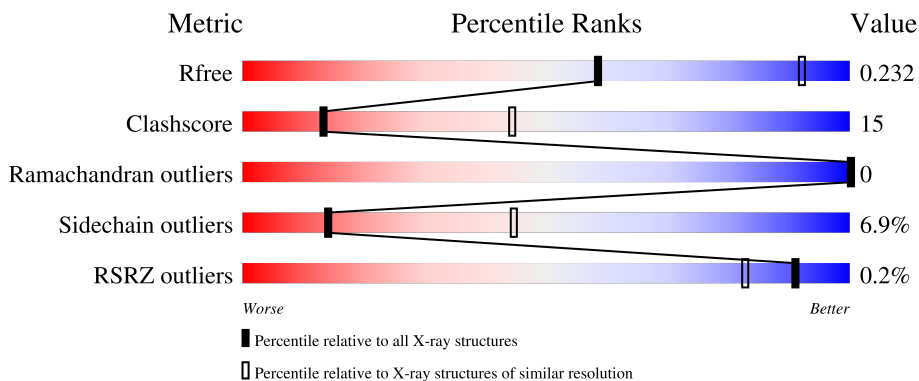
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

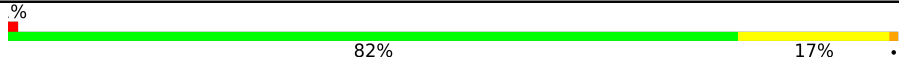
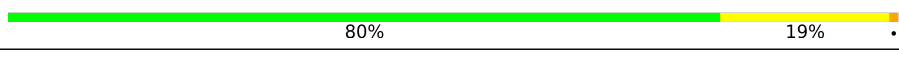
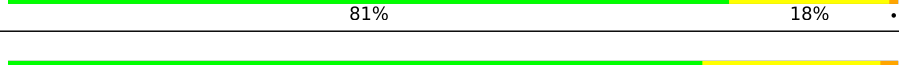

The reported resolution of this entry is 3.00 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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

Mol	Chain	Length	Quality of chain
1	A	162	 82% 17% .
1	C	162	 80% 19% .
1	E	162	 81% 18% .
1	G	162	 78% 20% .

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Mol	Chain	Length	Quality of chain	
1	I	162	82%	15%
1	K	162	81%	16%
1	M	162	77%	22%
1	O	162	82%	17%
1	Q	162	83%	16%
1	S	162	86%	13%
1	U	162	81%	19%
1	W	162	81%	17%
2	B	172	83%	17%
2	D	172	79%	20%
2	F	172	86%	13%
2	H	172	82%	17%
2	J	172	81%	17%
2	L	172	77%	22%
2	N	172	81%	17%
2	P	172	81%	17%
2	R	172	83%	16%
2	T	172	83%	15%
2	V	172	80%	19%
3	X	172	78%	20%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	CYC	D	1153	-	-	X	-
4	CYC	H	1153	-	-	X	-
4	CYC	J	1153	-	-	X	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	CYC	L	1153	-	-	X	-
4	CYC	M	1084	-	-	X	-
4	CYC	N	1153	-	-	X	-
4	CYC	O	1084	-	-	X	-
4	CYC	T	1153	-	-	X	-
4	CYC	X	1153	-	-	X	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 31610 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called C-PHYCOCYANIN ALPHA CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	162	1235	780	207	242	6	0	1	0
1	C	162	1233	778	207	242	6	0	0	0
1	E	162	1233	778	207	242	6	0	0	0
1	G	162	1233	778	207	242	6	0	0	0
1	I	162	1233	778	207	242	6	0	0	0
1	K	162	1233	778	207	242	6	0	0	0
1	M	162	1233	778	207	242	6	0	0	0
1	O	162	1233	778	207	242	6	0	0	0
1	Q	162	1233	778	207	242	6	0	0	0
1	S	162	1233	778	207	242	6	0	0	0
1	U	162	1233	778	207	242	6	0	0	0
1	W	162	1233	778	207	242	6	0	0	0

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	11	VAL	ILE	conflict	UNP P72509
A	148	VAL	THR	conflict	UNP P72509
C	11	VAL	ILE	conflict	UNP P72509
C	148	VAL	THR	conflict	UNP P72509
E	11	VAL	ILE	conflict	UNP P72509

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Chain	Residue	Modelled	Actual	Comment	Reference
E	148	VAL	THR	conflict	UNP P72509
G	11	VAL	ILE	conflict	UNP P72509
G	148	VAL	THR	conflict	UNP P72509
I	11	VAL	ILE	conflict	UNP P72509
I	148	VAL	THR	conflict	UNP P72509
K	11	VAL	ILE	conflict	UNP P72509
K	148	VAL	THR	conflict	UNP P72509
M	11	VAL	ILE	conflict	UNP P72509
M	148	VAL	THR	conflict	UNP P72509
O	11	VAL	ILE	conflict	UNP P72509
O	148	VAL	THR	conflict	UNP P72509
Q	11	VAL	ILE	conflict	UNP P72509
Q	148	VAL	THR	conflict	UNP P72509
S	11	VAL	ILE	conflict	UNP P72509
S	148	VAL	THR	conflict	UNP P72509
U	11	VAL	ILE	conflict	UNP P72509
U	148	VAL	THR	conflict	UNP P72509
W	11	VAL	ILE	conflict	UNP P72509
W	148	VAL	THR	conflict	UNP P72509

- Molecule 2 is a protein called C-PHYCOCYANIN BETA CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	172	1258	779	218	252	9	0	0	0
2	D	172	1258	779	218	252	9	0	0	0
2	F	172	1258	779	218	252	9	0	0	0
2	H	172	1258	779	218	252	9	0	0	0
2	J	172	1258	779	218	252	9	0	0	0
2	L	172	1258	779	218	252	9	0	0	0
2	N	172	1258	779	218	252	9	0	0	0
2	P	172	1258	779	218	252	9	0	0	0
2	R	172	1258	779	218	252	9	0	0	0
2	T	172	1258	779	218	252	9	0	0	0

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	V	172	1258	779	218	252	9	0	0	0

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	40	VAL	ALA	conflict	UNP P72508
B	77	THR	ARG	conflict	UNP P72508
D	40	VAL	ALA	conflict	UNP P72508
D	77	THR	ARG	conflict	UNP P72508
F	40	VAL	ALA	conflict	UNP P72508
F	77	THR	ARG	conflict	UNP P72508
H	40	VAL	ALA	conflict	UNP P72508
H	77	THR	ARG	conflict	UNP P72508
J	40	VAL	ALA	conflict	UNP P72508
J	77	THR	ARG	conflict	UNP P72508
L	40	VAL	ALA	conflict	UNP P72508
L	77	THR	ARG	conflict	UNP P72508
N	40	VAL	ALA	conflict	UNP P72508
N	77	THR	ARG	conflict	UNP P72508
P	40	VAL	ALA	conflict	UNP P72508
P	77	THR	ARG	conflict	UNP P72508
R	40	VAL	ALA	conflict	UNP P72508
R	77	THR	ARG	conflict	UNP P72508
T	40	VAL	ALA	conflict	UNP P72508
T	77	THR	ARG	conflict	UNP P72508
V	40	VAL	ALA	conflict	UNP P72508
V	77	THR	ARG	conflict	UNP P72508

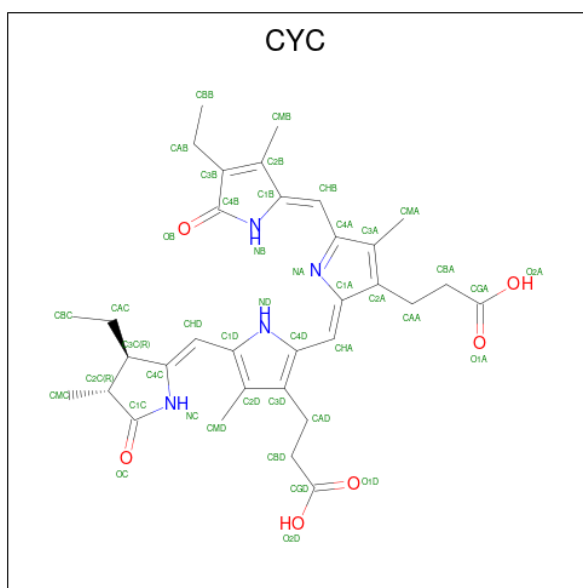
- Molecule 3 is a protein called C-PHYCOCYANIN BETA CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	X	172	1263	783	219	252	9	0	0	0

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
X	40	VAL	ALA	conflict	UNP P72508
X	77	THR	ARG	conflict	UNP P72508
X	132	LYS	GLY	conflict	UNP P72508

- Molecule 4 is PHYCOCYANOBILIN (CCD ID: CYC) (formula: $C_{33}H_{40}N_4O_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			43	33	4	6		
4	B	1	Total	C	N	O	0	0
			43	33	4	6		
4	B	1	Total	C	N	O	0	0
			43	33	4	6		
4	C	1	Total	C	N	O	0	0
			43	33	4	6		
4	D	1	Total	C	N	O	0	0
			43	33	4	6		
4	D	1	Total	C	N	O	0	0
			43	33	4	6		
4	E	1	Total	C	N	O	0	0
			43	33	4	6		
4	F	1	Total	C	N	O	0	0
			43	33	4	6		
4	F	1	Total	C	N	O	0	0
			43	33	4	6		
4	G	1	Total	C	N	O	0	0
			43	33	4	6		
4	H	1	Total	C	N	O	0	0
			43	33	4	6		
4	H	1	Total	C	N	O	0	0
			43	33	4	6		
4	I	1	Total	C	N	O	0	0
			43	33	4	6		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
4	J	1	Total 43	C 33	N 4	O 6	0	0
4	J	1	Total 43	C 33	N 4	O 6	0	0
4	K	1	Total 43	C 33	N 4	O 6	0	0
4	L	1	Total 43	C 33	N 4	O 6	0	0
4	L	1	Total 43	C 33	N 4	O 6	0	0
4	M	1	Total 43	C 33	N 4	O 6	0	0
4	N	1	Total 43	C 33	N 4	O 6	0	0
4	N	1	Total 43	C 33	N 4	O 6	0	0
4	O	1	Total 43	C 33	N 4	O 6	0	0
4	P	1	Total 43	C 33	N 4	O 6	0	0
4	P	1	Total 43	C 33	N 4	O 6	0	0
4	Q	1	Total 43	C 33	N 4	O 6	0	0
4	R	1	Total 43	C 33	N 4	O 6	0	0
4	R	1	Total 43	C 33	N 4	O 6	0	0
4	S	1	Total 43	C 33	N 4	O 6	0	0
4	T	1	Total 43	C 33	N 4	O 6	0	0
4	T	1	Total 43	C 33	N 4	O 6	0	0
4	U	1	Total 43	C 33	N 4	O 6	0	0
4	V	1	Total 43	C 33	N 4	O 6	0	0
4	V	1	Total 43	C 33	N 4	O 6	0	0
4	W	1	Total 43	C 33	N 4	O 6	0	0

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	X	1	Total	C	N	O	0	0
			43	33	4	6		
4	X	1	Total	C	N	O	0	0
			43	33	4	6		

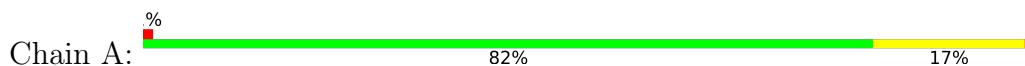
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	8	Total	O	0	0
			8	8		
5	C	10	Total	O	0	0
			10	10		
5	D	21	Total	O	0	0
			21	21		
5	F	6	Total	O	0	0
			6	6		
5	G	5	Total	O	0	0
			5	5		
5	I	13	Total	O	0	0
			13	13		
5	J	17	Total	O	0	0
			17	17		
5	L	5	Total	O	0	0
			5	5		
5	M	12	Total	O	0	0
			12	12		
5	N	16	Total	O	0	0
			16	16		
5	P	7	Total	O	0	0
			7	7		
5	Q	6	Total	O	0	0
			6	6		
5	S	2	Total	O	0	0
			2	2		
5	T	2	Total	O	0	0
			2	2		
5	U	5	Total	O	0	0
			5	5		
5	W	13	Total	O	0	0
			13	13		
5	X	15	Total	O	0	0
			15	15		

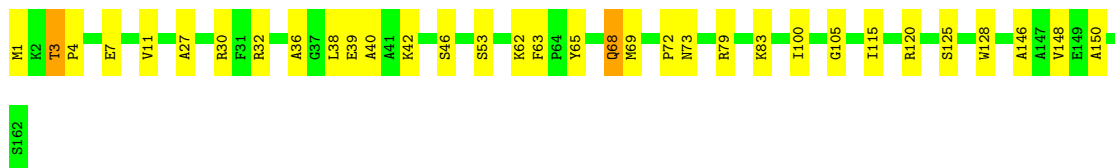
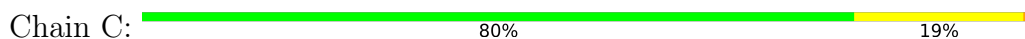
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

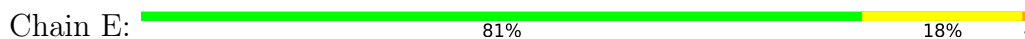
- Molecule 1: C-PHYCOCYANIN ALPHA CHAIN



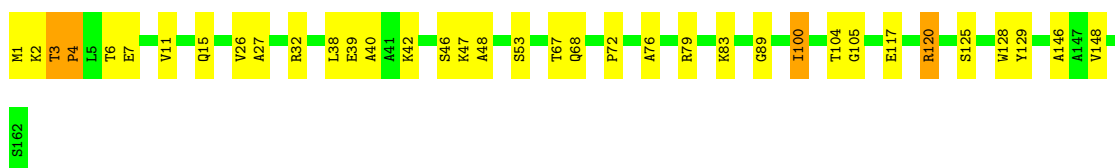
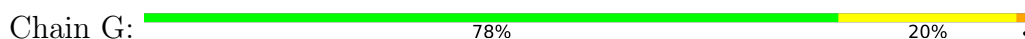
- Molecule 1: C-PHYCOCYANIN ALPHA CHAIN



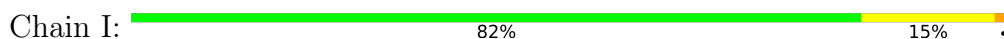
- Molecule 1: C-PHYCOCYANIN ALPHA CHAIN

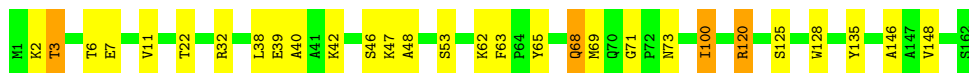


- Molecule 1: C-PHYCOCYANIN ALPHA CHAIN

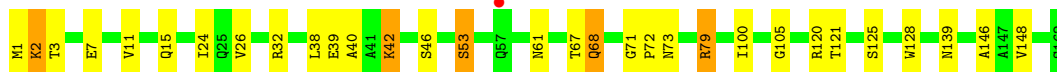
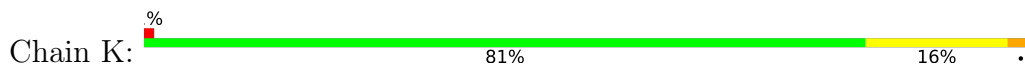


- Molecule 1: C-PHYCOCYANIN ALPHA CHAIN

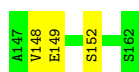
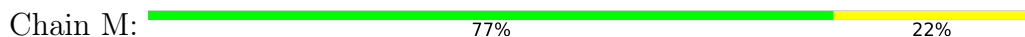




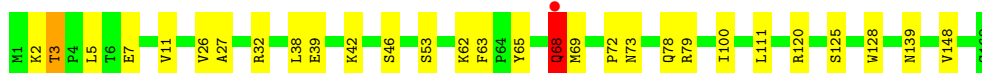
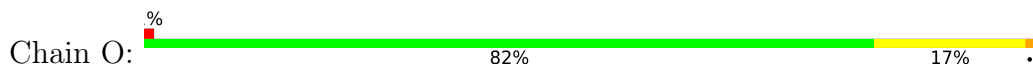
- Molecule 1: C-PHYCOCYANIN ALPHA CHAIN



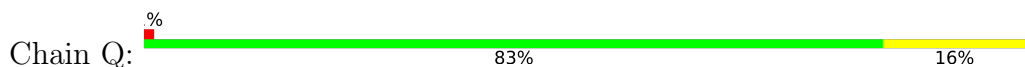
- Molecule 1: C-PHYCOCYANIN ALPHA CHAIN



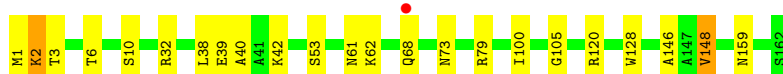
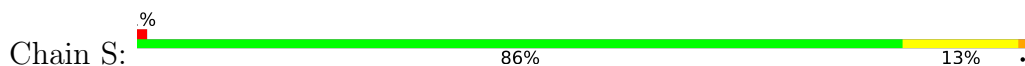
- Molecule 1: C-PHYCOCYANIN ALPHA CHAIN



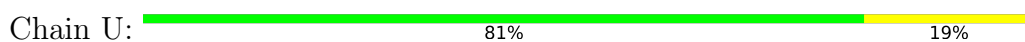
- Molecule 1: C-PHYCOCYANIN ALPHA CHAIN




- Molecule 1: C-PHYCOCYANIN ALPHA CHAIN

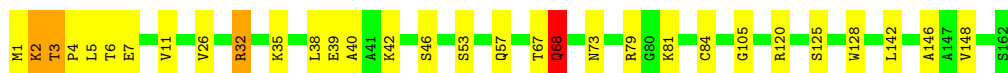


- Molecule 1: C-PHYCOCYANIN ALPHA CHAIN




- Molecule 1: C-PHYCOCYANIN ALPHA CHAIN

Chain W:  81% 17% ..




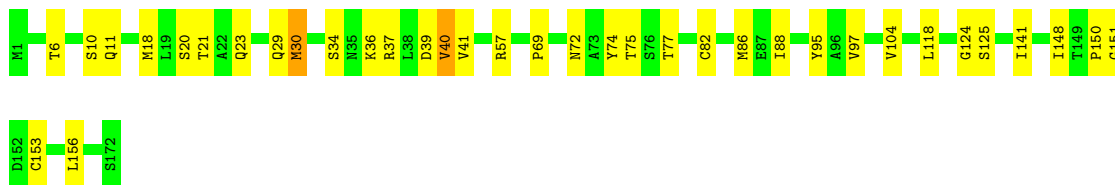
• Molecule 2: C-PHYCOCYANIN BETA CHAIN

Chain B:  83% 17% .




• Molecule 2: C-PHYCOCYANIN BETA CHAIN

Chain D:  79% 20% .




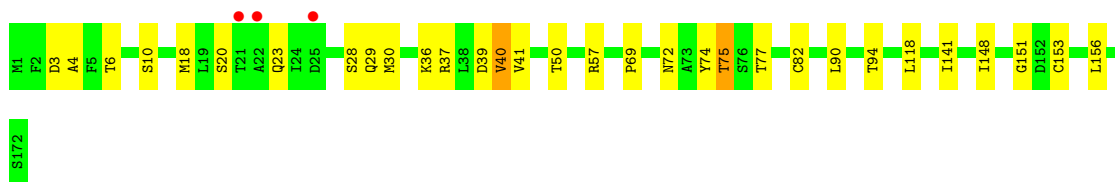
• Molecule 2: C-PHYCOCYANIN BETA CHAIN

Chain F:  86% 13% .




• Molecule 2: C-PHYCOCYANIN BETA CHAIN

Chain H:  2% 82% 17% .



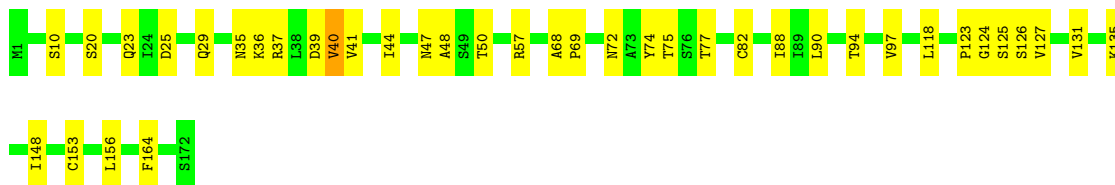
• Molecule 2: C-PHYCOCYANIN BETA CHAIN

Chain J:  81% 17% .



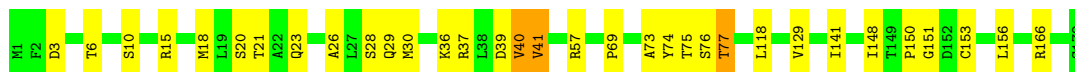
• Molecule 2: C-PHYCOCYANIN BETA CHAIN

Chain L:  77% 22% .



- Molecule 2: C-PHYCOCYANIN BETA CHAIN

Chain N: 81% 17%



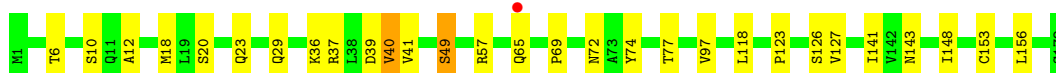
- Molecule 2: C-PHYCOCYANIN BETA CHAIN

Chain P: 81% 17%



- Molecule 2: C-PHYCOCYANIN BETA CHAIN

Chain R: 83% 16%



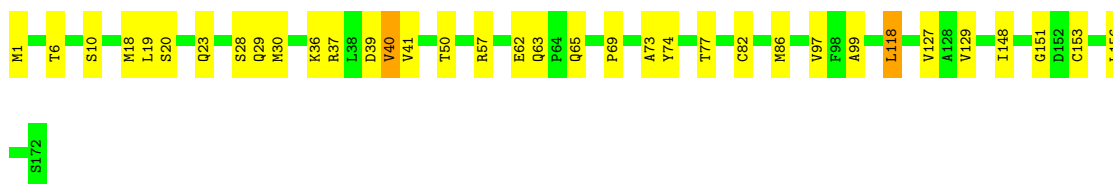
- Molecule 2: C-PHYCOCYANIN BETA CHAIN

Chain T: 83% 15%



- Molecule 2: C-PHYCOCYANIN BETA CHAIN

Chain V: 80% 19%



- Molecule 3: C-PHYCOCYANIN BETA CHAIN

Chain X: 78% 20%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	107.33Å 115.64Å 183.26Å 90.00° 90.03° 90.00°	Depositor
Resolution (Å)	182.57 – 3.00 182.57 – 3.00	Depositor EDS
% Data completeness (in resolution range)	96.5 (182.57-3.00) 96.6 (182.57-3.00)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.66 (at 3.01Å)	Xtrriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.201 , 0.242 0.193 , 0.232	Depositor DCC
R_{free} test set	4331 reflections (4.81%)	wwPDB-VP
Wilson B-factor (Å ²)	39.3	Xtrriage
Anisotropy	0.098	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 18.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.018 for h,-k,-l	Xtrriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	31610	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.26% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	1.18	0/1264	1.14	1/1710 (0.1%)
1	C	1.20	3/1256 (0.2%)	1.15	2/1699 (0.1%)
1	E	1.22	3/1256 (0.2%)	1.08	2/1699 (0.1%)
1	G	1.21	1/1256 (0.1%)	1.15	3/1699 (0.2%)
1	I	1.19	0/1256	1.16	3/1699 (0.2%)
1	K	1.18	1/1256 (0.1%)	1.19	7/1699 (0.4%)
1	M	1.14	1/1256 (0.1%)	1.16	0/1699
1	O	1.20	1/1256 (0.1%)	1.14	2/1699 (0.1%)
1	Q	1.22	3/1256 (0.2%)	1.15	1/1699 (0.1%)
1	S	1.19	1/1256 (0.1%)	1.11	1/1699 (0.1%)
1	U	1.17	1/1256 (0.1%)	1.12	2/1699 (0.1%)
1	W	1.16	2/1256 (0.2%)	1.15	2/1699 (0.1%)
2	B	1.16	1/1272 (0.1%)	1.19	1/1724 (0.1%)
2	D	1.18	1/1272 (0.1%)	1.21	1/1724 (0.1%)
2	F	1.16	1/1272 (0.1%)	1.18	2/1724 (0.1%)
2	H	1.19	2/1272 (0.2%)	1.21	1/1724 (0.1%)
2	J	1.22	2/1272 (0.2%)	1.23	3/1724 (0.2%)
2	L	1.19	2/1272 (0.2%)	1.14	2/1724 (0.1%)
2	N	1.22	4/1272 (0.3%)	1.20	1/1724 (0.1%)
2	P	1.18	3/1272 (0.2%)	1.20	2/1724 (0.1%)
2	R	1.18	2/1272 (0.2%)	1.20	2/1724 (0.1%)
2	T	1.17	2/1272 (0.2%)	1.18	0/1724
2	V	1.22	3/1272 (0.2%)	1.20	3/1724 (0.2%)
3	X	1.23	1/1277 (0.1%)	1.22	3/1730 (0.2%)
All	All	1.19	41/30349 (0.1%)	1.17	47/41093 (0.1%)

All (41) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	X	77	THR	CA-CB	11.89	1.72	1.53
2	R	77	THR	CA-CB	10.44	1.69	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	77	THR	CA-CB	9.62	1.68	1.53
2	V	77	THR	CA-CB	9.36	1.68	1.53
2	J	77	THR	CA-CB	9.30	1.68	1.53
2	L	77	THR	CA-CB	8.77	1.67	1.53
2	P	77	THR	CA-CB	8.75	1.67	1.53
2	B	77	THR	CA-CB	8.71	1.67	1.53
2	N	77	THR	CA-CB	7.67	1.65	1.53
2	T	77	THR	CA-CB	7.48	1.65	1.53
2	H	77	THR	CA-CB	7.38	1.65	1.53
2	D	77	THR	CA-CB	6.67	1.63	1.53
2	L	41	VAL	CA-CB	-6.48	1.46	1.54
1	U	27	ALA	CA-CB	-6.27	1.43	1.53
1	G	27	ALA	CA-CB	-6.24	1.43	1.53
1	Q	24	ILE	CA-CB	-6.05	1.47	1.54
1	M	35	LYS	CA-CB	5.99	1.63	1.53
1	Q	150	ALA	CA-CB	-5.83	1.44	1.53
1	E	88	ILE	CA-CB	-5.68	1.47	1.54
1	W	84	CYS	N-CA	-5.62	1.39	1.46
2	J	67	ILE	CA-CB	-5.50	1.48	1.54
1	C	27	ALA	CA-CB	-5.46	1.44	1.53
1	E	118	ILE	CA-CB	-5.39	1.48	1.54
2	V	99	ALA	CA-CB	-5.38	1.44	1.53
2	T	93	VAL	N-CA	-5.37	1.40	1.46
2	P	14	THR	CA-CB	5.37	1.62	1.53
2	N	26	ALA	CA-CB	5.30	1.61	1.53
1	C	150	ALA	CA-CB	-5.23	1.45	1.53
2	N	41	VAL	CA-CB	-5.21	1.48	1.54
1	O	27	ALA	CA-CB	-5.21	1.45	1.53
1	E	34	ALA	CA-CB	-5.16	1.45	1.53
2	H	75	THR	CA-CB	5.16	1.62	1.53
2	V	73	ALA	CA-CB	-5.15	1.46	1.53
2	P	93	VAL	CA-CB	-5.15	1.48	1.54
1	Q	27	ALA	CA-CB	-5.13	1.45	1.53
2	R	12	ALA	CA-CB	-5.12	1.45	1.53
1	W	35	LYS	CA-CB	5.10	1.61	1.53
1	C	30	ARG	CZ-NH1	5.07	1.39	1.32
1	S	148	VAL	N-CA	-5.03	1.40	1.46
2	N	75	THR	CA-CB	5.02	1.61	1.53
1	K	68	GLN	CB-CG	-5.01	1.37	1.52

All (47) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	O	139	ASN	N-CA-C	7.17	122.90	113.30
2	V	129	VAL	N-CA-C	-6.48	104.17	110.72
2	J	77	THR	N-CA-CB	6.32	119.54	110.06
2	V	86	MET	N-CA-C	-6.30	104.33	111.07
1	G	83	LYS	N-CA-C	-6.12	104.52	111.07
1	G	100	ILE	CB-CA-C	-6.03	104.25	111.97
2	N	129	VAL	N-CA-C	-5.98	104.68	110.72
2	L	50	THR	N-CA-CB	5.75	118.66	110.16
1	A	10	SER	N-CA-C	5.74	117.54	111.28
1	I	100	ILE	CB-CA-C	-5.71	104.66	111.97
1	U	121	THR	N-CA-C	5.66	117.90	111.11
2	F	131	VAL	CA-C-N	5.66	126.37	120.03
2	F	131	VAL	C-N-CA	5.66	126.37	120.03
1	I	71	GLY	CA-C-N	5.59	125.21	119.56
1	I	71	GLY	C-N-CA	5.59	125.21	119.56
2	L	44	ILE	N-CA-C	-5.58	105.28	110.53
1	K	121	THR	N-CA-C	5.58	117.80	111.11
3	X	77	THR	N-CA-CB	5.57	118.41	110.06
1	K	71	GLY	CA-C-N	5.54	125.16	119.56
1	K	71	GLY	C-N-CA	5.54	125.16	119.56
2	D	86	MET	N-CA-C	-5.54	105.24	111.28
1	E	121	THR	N-CA-C	5.46	117.66	111.11
1	C	68	GLN	N-CA-C	5.45	122.40	110.80
1	K	79	ARG	CA-C-N	5.39	125.93	120.00
1	K	79	ARG	C-N-CA	5.39	125.93	120.00
1	U	59	VAL	N-CA-C	-5.36	105.50	110.53
2	R	143	ASN	N-CA-C	5.33	119.41	113.01
2	B	77	THR	N-CA-CB	5.33	117.88	109.94
2	P	129	VAL	N-CA-C	-5.28	105.39	110.72
1	E	68	GLN	N-CA-C	5.28	122.04	110.80
2	J	68	ALA	CA-C-N	-5.26	114.53	119.90
2	J	68	ALA	C-N-CA	-5.26	114.53	119.90
3	X	18	MET	N-CA-C	-5.25	102.75	110.52
2	P	76	SER	N-CA-C	5.23	116.98	111.28
1	C	83	LYS	N-CA-C	-5.22	105.48	111.07
1	S	10	SER	N-CA-C	5.21	116.64	111.07
3	X	136	GLU	N-CA-C	-5.20	105.52	111.14
1	O	68	GLN	N-CA-C	5.17	121.81	114.16
1	K	139	ASN	N-CA-C	5.14	121.31	113.61
2	H	50	THR	N-CA-CB	5.10	117.81	110.20
1	W	68	GLN	N-CA-C	5.09	121.64	110.80
2	V	50	THR	N-CA-CB	5.08	117.76	110.20
2	R	65	GLN	N-CA-C	-5.06	106.21	112.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	W	32	ARG	N-CA-C	-5.03	105.70	111.14
1	G	89	GLY	N-CA-C	-5.02	106.67	112.50
1	K	53	SER	CB-CA-C	-5.00	102.80	110.81
1	Q	112	ILE	CB-CA-C	-5.00	105.47	112.02

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1235	0	1210	18	1
1	C	1233	0	1205	27	0
1	E	1233	0	1205	22	0
1	G	1233	0	1205	21	0
1	I	1233	0	1205	18	1
1	K	1233	0	1205	18	2
1	M	1233	0	1205	31	0
1	O	1233	0	1205	18	1
1	Q	1233	0	1205	18	1
1	S	1233	0	1205	12	1
1	U	1233	0	1205	16	0
1	W	1233	0	1205	22	1
2	B	1258	0	1255	32	0
2	D	1258	0	1254	44	1
2	F	1258	0	1254	24	0
2	H	1258	0	1255	40	0
2	J	1258	0	1255	33	0
2	L	1258	0	1255	44	0
2	N	1258	0	1254	32	0
2	P	1258	0	1255	30	0
2	R	1258	0	1254	30	0
2	T	1258	0	1254	31	0
2	V	1258	0	1254	35	1
3	X	1263	0	1265	47	0
4	A	43	0	36	15	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	86	0	76	29	0
4	C	43	0	37	14	0
4	D	86	0	75	37	0
4	E	43	0	37	16	0
4	F	86	0	75	27	0
4	G	43	0	35	11	0
4	H	86	0	76	39	0
4	I	43	0	36	13	0
4	J	86	0	76	38	0
4	K	43	0	37	19	0
4	L	86	0	76	46	0
4	M	43	0	36	22	0
4	N	86	0	75	35	0
4	O	43	0	37	21	0
4	P	86	0	76	30	0
4	Q	43	0	37	18	0
4	R	86	0	75	30	0
4	S	43	0	36	10	0
4	T	86	0	75	30	0
4	U	43	0	35	14	0
4	V	86	0	75	34	0
4	W	43	0	37	14	0
4	X	86	0	76	38	0
5	A	8	0	0	1	0
5	C	10	0	0	4	0
5	D	21	0	0	2	0
5	F	6	0	0	0	0
5	G	5	0	0	1	0
5	I	13	0	0	1	0
5	J	17	0	0	3	0
5	L	5	0	0	0	0
5	M	12	0	0	3	0
5	N	16	0	0	5	0
5	P	7	0	0	1	0
5	Q	6	0	0	1	0
5	S	2	0	0	0	0
5	T	2	0	0	0	0
5	U	5	0	0	1	0
5	W	13	0	0	7	0
5	X	15	0	0	9	0
All	All	31610	0	30871	917	5

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including

hydrogen atoms). The all-atom clashscore for this structure is 15.

All (917) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:82:CYS:SG	4:B:1082:CYC:HAC2	1.24	1.75
2:L:82:CYS:SG	4:L:1082:CYC:HAC2	1.22	1.72
2:P:82:CYS:SG	4:P:1082:CYC:HAC2	1.22	1.72
2:H:82:CYS:SG	4:H:1082:CYC:HAC2	1.30	1.71
2:F:82:CYS:SG	4:F:1082:CYC:HAC2	1.14	1.70
2:J:82:CYS:SG	4:J:1082:CYC:HAC2	1.13	1.70
2:D:153:CYS:SG	4:D:1153:CYC:HBC2	1.19	1.67
2:N:153:CYS:SG	4:N:1153:CYC:HAC1	1.36	1.66
3:X:82:CYS:SG	4:X:1082:CYC:HAC2	1.32	1.64
2:B:153:CYS:SG	4:B:1153:CYC:HAC1	1.32	1.64
2:R:153:CYS:SG	4:R:1153:CYC:HBC2	1.35	1.61
2:L:153:CYS:SG	4:L:1153:CYC:HAC2	1.45	1.55
2:V:153:CYS:SG	4:V:1153:CYC:HBC3	1.54	1.45
2:P:153:CYS:SG	4:P:1153:CYC:HAC1	1.58	1.42
2:D:153:CYS:SG	4:D:1153:CYC:CBC	2.13	1.36
2:B:82:CYS:SG	4:B:1082:CYC:CAC	2.13	1.35
2:H:82:CYS:SG	4:H:1082:CYC:CAC	2.16	1.34
2:P:82:CYS:SG	4:P:1082:CYC:CAC	2.15	1.34
4:A:1084:CYC:HC	4:A:1084:CYC:CMD	1.41	1.32
4:A:1084:CYC:NC	4:A:1084:CYC:HMD1	1.44	1.32
4:R:1082:CYC:HB	4:R:1082:CYC:CMA	1.43	1.30
3:X:82:CYS:SG	4:X:1082:CYC:CAC	2.20	1.29
2:H:153:CYS:SG	4:H:1153:CYC:HAC2	1.73	1.27
2:L:153:CYS:SG	4:L:1153:CYC:CAC	2.24	1.25
4:H:1082:CYC:HC	4:H:1082:CYC:CMD	1.49	1.25
4:F:1082:CYC:HB	4:F:1082:CYC:CMA	1.47	1.25
4:O:1084:CYC:HB	4:O:1084:CYC:CMA	1.50	1.24
2:R:153:CYS:SG	4:R:1153:CYC:CBC	2.25	1.24
2:J:153:CYS:SG	4:J:1153:CYC:HAC1	1.77	1.23
4:R:1082:CYC:HC	4:R:1082:CYC:CMD	1.50	1.22
4:R:1082:CYC:HMA1	4:R:1082:CYC:NB	1.53	1.22
4:N:1082:CYC:HB	4:N:1082:CYC:CMA	1.51	1.22
3:X:153:CYS:SG	4:X:1153:CYC:HAC2	1.81	1.21
2:B:153:CYS:SG	4:B:1153:CYC:CAC	2.28	1.19
4:W:1084:CYC:HB	4:W:1084:CYC:CMA	1.56	1.19
2:N:153:CYS:SG	4:N:1153:CYC:CAC	2.29	1.19
4:G:1084:CYC:HMD1	4:G:1084:CYC:HC	1.08	1.17
4:L:1082:CYC:HC	4:L:1082:CYC:CMD	1.55	1.17

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:G:1084:CYC:HB	4:G:1084:CYC:HMA3	1.10	1.16
4:O:1084:CYC:HC	4:O:1084:CYC:CMD	1.59	1.15
4:Q:1084:CYC:HMD1	4:Q:1084:CYC:HC	1.01	1.15
4:V:1153:CYC:HB	4:V:1153:CYC:HMA3	1.10	1.15
2:V:153:CYS:SG	4:V:1153:CYC:CBC	2.33	1.15
4:R:1082:CYC:NC	4:R:1082:CYC:HMD1	1.62	1.15
4:M:1084:CYC:HB	4:M:1084:CYC:HMA3	1.07	1.14
4:Q:1084:CYC:CMA	4:Q:1084:CYC:HB	1.59	1.14
4:V:1082:CYC:HC	4:V:1082:CYC:CMD	1.58	1.14
4:N:1082:CYC:HC	4:N:1082:CYC:CMD	1.60	1.14
4:O:1084:CYC:HMA3	4:O:1084:CYC:NB	1.64	1.13
4:U:1084:CYC:HC	4:U:1084:CYC:CMD	1.60	1.13
4:V:1082:CYC:HB	4:V:1082:CYC:HMA3	1.11	1.13
4:K:1084:CYC:HB	4:K:1084:CYC:CMA	1.61	1.13
4:U:1084:CYC:HMA3	4:U:1084:CYC:HB	1.05	1.12
4:D:1153:CYC:HMA3	4:D:1153:CYC:HB	0.95	1.12
4:E:1084:CYC:HB	4:E:1084:CYC:HMA3	0.96	1.12
4:F:1082:CYC:HMA3	4:F:1082:CYC:NB	1.64	1.11
4:W:1084:CYC:HB	4:W:1084:CYC:HMA3	1.01	1.11
4:E:1084:CYC:HB	4:E:1084:CYC:CMA	1.63	1.11
4:T:1082:CYC:HC	4:T:1082:CYC:HMD1	1.11	1.11
4:F:1153:CYC:HB	4:F:1153:CYC:HMA3	1.15	1.11
4:H:1082:CYC:NC	4:H:1082:CYC:HMD1	1.63	1.11
4:I:1084:CYC:HC	4:I:1084:CYC:HMD1	0.94	1.10
4:I:1084:CYC:HMA1	4:I:1084:CYC:HB	1.03	1.10
4:P:1082:CYC:HMA3	4:P:1082:CYC:HB	1.05	1.10
4:T:1153:CYC:HMD1	4:T:1153:CYC:HC	1.14	1.10
4:M:1084:CYC:HC	4:M:1084:CYC:CMD	1.65	1.10
4:N:1082:CYC:HMA1	4:N:1082:CYC:NB	1.66	1.10
4:C:1084:CYC:HB	4:C:1084:CYC:HMA3	1.04	1.09
4:Q:1084:CYC:HB	4:Q:1084:CYC:HMA3	0.96	1.09
4:K:1084:CYC:HC	4:K:1084:CYC:HMD1	1.03	1.09
4:D:1082:CYC:HMA3	4:D:1082:CYC:HB	1.13	1.09
4:I:1084:CYC:HC	4:I:1084:CYC:CMD	1.66	1.09
2:P:153:CYS:SG	4:P:1153:CYC:CAC	2.40	1.08
4:J:1153:CYC:HB	4:J:1153:CYC:HMA3	1.17	1.08
4:K:1084:CYC:HB	4:K:1084:CYC:HMA3	1.01	1.08
4:N:1082:CYC:HB	4:N:1082:CYC:HMA1	0.94	1.08
4:D:1082:CYC:HC	4:D:1082:CYC:HMD1	1.18	1.07
4:O:1084:CYC:NC	4:O:1084:CYC:HMD1	1.68	1.07
4:V:1082:CYC:NC	4:V:1082:CYC:HMD1	1.67	1.07

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:1082:CYC:HB	4:D:1082:CYC:CMA	1.68	1.07
4:P:1153:CYC:HC	4:P:1153:CYC:HMD1	1.14	1.07
4:N:1082:CYC:HMD1	4:N:1082:CYC:NC	1.69	1.07
4:A:1084:CYC:HB	4:A:1084:CYC:HMA1	0.93	1.06
4:H:1082:CYC:HB	4:H:1082:CYC:HMA3	1.14	1.06
4:U:1084:CYC:HC	4:U:1084:CYC:HMD1	0.92	1.06
1:W:142:LEU:HA	5:W:2008:HOH:O	1.52	1.06
4:B:1082:CYC:HMD1	4:B:1082:CYC:HC	1.18	1.05
4:L:1082:CYC:HC	4:L:1082:CYC:HMD1	0.91	1.05
4:A:1084:CYC:HB	4:A:1084:CYC:CMA	1.69	1.05
4:J:1082:CYC:HC	4:J:1082:CYC:HMD1	0.95	1.05
4:L:1082:CYC:HMD1	4:L:1082:CYC:NC	1.70	1.05
4:V:1153:CYC:HMD1	4:V:1153:CYC:HC	1.10	1.05
2:D:150:PRO:HD3	1:M:39:GLU:HB2	1.36	1.05
4:M:1084:CYC:HC	4:M:1084:CYC:HMD1	0.92	1.04
4:O:1084:CYC:HB	4:O:1084:CYC:HMA3	0.89	1.04
4:L:1153:CYC:HMD1	4:L:1153:CYC:HC	1.22	1.04
2:T:153:CYS:SG	4:T:1153:CYC:HAC1	1.97	1.04
4:M:1084:CYC:HMD1	4:M:1084:CYC:NC	1.73	1.04
4:U:1084:CYC:HMD1	4:U:1084:CYC:NC	1.73	1.04
4:Q:1084:CYC:HMA3	4:Q:1084:CYC:NB	1.73	1.03
4:T:1082:CYC:HB	4:T:1082:CYC:HMA3	1.19	1.03
4:U:1084:CYC:HB	4:U:1084:CYC:CMA	1.71	1.03
4:J:1082:CYC:HC	4:J:1082:CYC:CMD	1.71	1.03
4:J:1082:CYC:HMA1	4:J:1082:CYC:HB	1.23	1.03
4:W:1084:CYC:HC	4:W:1084:CYC:HMD1	1.21	1.03
4:F:1082:CYC:HB	4:F:1082:CYC:HMA3	0.89	1.02
4:N:1082:CYC:HC	4:N:1082:CYC:HMD1	0.88	1.02
4:S:1084:CYC:HB	4:S:1084:CYC:HMA3	1.20	1.02
4:X:1082:CYC:HB	4:X:1082:CYC:HMA3	1.23	1.02
2:D:82:CYS:SG	4:D:1082:CYC:CBC	2.45	1.02
4:C:1084:CYC:HB	4:C:1084:CYC:CMA	1.73	1.02
4:X:1082:CYC:HMD1	4:X:1082:CYC:HC	1.22	1.02
4:O:1084:CYC:HC	4:O:1084:CYC:HMD1	0.86	1.01
4:D:1153:CYC:HB	4:D:1153:CYC:CMA	1.72	1.01
4:V:1082:CYC:HC	4:V:1082:CYC:HMD1	0.86	1.01
3:X:153:CYS:SG	4:X:1153:CYC:C1C	2.48	1.01
3:X:153:CYS:SG	4:X:1153:CYC:C2C	2.49	1.01
4:H:1082:CYC:HC	4:H:1082:CYC:HMD1	0.84	1.01
4:J:1082:CYC:HMD1	4:J:1082:CYC:NC	1.76	1.00
4:P:1153:CYC:HB	4:P:1153:CYC:HMA3	1.21	1.00

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:U:1084:CYC:HBD2	4:U:1084:CYC:HHA	1.42	1.00
4:W:1084:CYC:HMA3	4:W:1084:CYC:NB	1.75	1.00
4:P:1082:CYC:HB	4:P:1082:CYC:CMA	1.74	1.00
4:R:1082:CYC:HC	4:R:1082:CYC:HMD1	0.83	1.00
4:V:1082:CYC:HB	4:V:1082:CYC:CMA	1.75	0.99
4:I:1084:CYC:HMD1	4:I:1084:CYC:NC	1.77	0.99
4:E:1084:CYC:HMA3	4:E:1084:CYC:NB	1.76	0.99
2:H:153:CYS:SG	4:H:1153:CYC:CAC	2.50	0.98
4:A:1084:CYC:HMA1	4:A:1084:CYC:NB	1.76	0.98
4:M:1084:CYC:HB	4:M:1084:CYC:CMA	1.77	0.98
4:N:1153:CYC:HMD3	4:N:1153:CYC:HC	1.29	0.98
4:K:1084:CYC:HMA3	4:K:1084:CYC:NB	1.78	0.97
4:X:1153:CYC:HMD1	4:X:1153:CYC:HC	1.30	0.97
4:Q:1084:CYC:HC	4:Q:1084:CYC:CMD	1.76	0.97
4:K:1084:CYC:HC	4:K:1084:CYC:CMD	1.77	0.96
4:G:1084:CYC:HC	4:G:1084:CYC:CMD	1.79	0.96
4:D:1153:CYC:HMA3	4:D:1153:CYC:NB	1.79	0.96
4:X:1153:CYC:HB	4:X:1153:CYC:HMA3	1.29	0.96
4:E:1084:CYC:HMD1	4:E:1084:CYC:HC	1.30	0.95
4:G:1084:CYC:HB	4:G:1084:CYC:CMA	1.78	0.95
4:I:1084:CYC:HB	4:I:1084:CYC:CMA	1.80	0.95
4:F:1153:CYC:HC	4:F:1153:CYC:HMD1	1.30	0.95
2:B:20:SER:H	2:B:23:GLN:HE21	1.14	0.95
4:H:1082:CYC:HB	4:H:1082:CYC:CMA	1.79	0.95
2:H:20:SER:H	2:H:23:GLN:HE21	1.07	0.94
4:T:1082:CYC:HC	4:T:1082:CYC:CMD	1.80	0.94
4:B:1153:CYC:HC	4:B:1153:CYC:HMD3	1.30	0.94
4:R:1153:CYC:HC	4:R:1153:CYC:HMD1	1.28	0.94
2:T:153:CYS:SG	4:T:1153:CYC:CAC	2.56	0.94
2:T:20:SER:H	2:T:23:GLN:HE21	1.15	0.94
4:B:1153:CYC:HC	4:B:1153:CYC:CMD	1.81	0.94
4:C:1084:CYC:HMA3	4:C:1084:CYC:NB	1.83	0.93
3:X:153:CYS:SG	4:X:1153:CYC:CAC	2.55	0.93
2:J:153:CYS:SG	4:J:1153:CYC:CAC	2.56	0.93
4:U:1084:CYC:HMA3	4:U:1084:CYC:NB	1.84	0.93
4:L:1153:CYC:HB	4:L:1153:CYC:HMA3	1.34	0.92
4:T:1082:CYC:HB	4:T:1082:CYC:CMA	1.83	0.92
4:R:1082:CYC:HB	4:R:1082:CYC:HMA1	0.78	0.92
4:P:1082:CYC:HC	4:P:1082:CYC:HMD1	1.33	0.92
4:V:1153:CYC:HC	4:V:1153:CYC:CMD	1.81	0.92
2:T:153:CYS:SG	4:T:1153:CYC:CBC	2.57	0.91

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:Q:1084:CYC:HMD1	4:Q:1084:CYC:NC	1.84	0.91
4:H:1153:CYC:HC	4:H:1153:CYC:HMD1	1.35	0.90
4:K:1084:CYC:HMD1	4:K:1084:CYC:NC	1.86	0.90
4:M:1084:CYC:HMA3	4:M:1084:CYC:NB	1.87	0.90
4:P:1153:CYC:HC	4:P:1153:CYC:CMD	1.84	0.90
4:M:1084:CYC:HBD2	4:M:1084:CYC:HHA	1.52	0.90
2:T:153:CYS:SG	4:T:1153:CYC:HBC3	2.11	0.90
4:B:1082:CYC:HB	4:B:1082:CYC:HMA3	1.35	0.90
2:F:20:SER:H	2:F:23:GLN:HE21	1.12	0.90
4:T:1153:CYC:HB	4:T:1153:CYC:HMA1	1.37	0.90
4:P:1082:CYC:HMA3	4:P:1082:CYC:NB	1.86	0.89
4:I:1084:CYC:HMA1	4:I:1084:CYC:NB	1.87	0.89
3:X:153:CYS:SG	4:X:1153:CYC:H2C	2.11	0.89
4:N:1153:CYC:HC	4:N:1153:CYC:CMD	1.86	0.89
2:R:20:SER:H	2:R:23:GLN:HE21	1.13	0.89
4:T:1153:CYC:HC	4:T:1153:CYC:CMD	1.86	0.89
2:D:20:SER:H	2:D:23:GLN:HE21	1.22	0.88
4:V:1082:CYC:HBD2	4:V:1082:CYC:HHA	1.53	0.88
4:L:1082:CYC:HB	4:L:1082:CYC:HMA3	1.35	0.88
1:C:148:VAL:HG11	4:J:1153:CYC:CMB	2.02	0.88
4:Q:1084:CYC:HBD2	4:Q:1084:CYC:HHA	1.54	0.88
4:V:1153:CYC:HB	4:V:1153:CYC:CMA	1.85	0.88
4:G:1084:CYC:HMA3	4:G:1084:CYC:NB	1.89	0.88
4:D:1082:CYC:HMA3	4:D:1082:CYC:NB	1.89	0.87
4:H:1153:CYC:HB	4:H:1153:CYC:HMA3	1.40	0.87
4:F:1153:CYC:HB	4:F:1153:CYC:CMA	1.87	0.87
2:L:20:SER:H	2:L:23:GLN:HE21	1.20	0.87
4:D:1153:CYC:HC	4:D:1153:CYC:HMD1	1.40	0.87
1:Q:148:VAL:HG11	4:T:1153:CYC:CMB	2.05	0.87
4:F:1082:CYC:HC	4:F:1082:CYC:HMD1	1.37	0.87
2:P:20:SER:H	2:P:23:GLN:HE21	1.20	0.87
4:R:1153:CYC:HB	4:R:1153:CYC:HMA3	1.37	0.87
4:V:1082:CYC:HMA3	4:V:1082:CYC:NB	1.90	0.87
4:B:1082:CYC:HC	4:B:1082:CYC:CMD	1.88	0.86
1:W:81:LYS:HE3	5:W:2006:HOH:O	1.73	0.86
2:L:82:CYS:SG	4:L:1082:CYC:CBC	2.63	0.86
2:N:21:THR:HB	5:N:2004:HOH:O	1.73	0.86
4:C:1084:CYC:HMD3	4:C:1084:CYC:HC	1.39	0.86
1:E:148:VAL:HG11	4:H:1153:CYC:CMB	2.06	0.85
4:S:1084:CYC:HB	4:S:1084:CYC:CMA	1.88	0.85
2:H:20:SER:H	2:H:23:GLN:NE2	1.75	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:V:1153:CYC:HMD1	4:V:1153:CYC:NC	1.92	0.85
4:G:1084:CYC:HMD1	4:G:1084:CYC:NC	1.91	0.84
4:S:1084:CYC:HC	4:S:1084:CYC:HMD1	1.40	0.84
4:S:1084:CYC:O1D	2:V:57:ARG:NH1	2.10	0.84
4:D:1082:CYC:HC	4:D:1082:CYC:CMD	1.89	0.84
3:X:20:SER:H	3:X:23:GLN:HE21	1.25	0.83
2:D:150:PRO:CD	1:M:39:GLU:HB2	2.08	0.83
2:H:82:CYS:SG	4:H:1082:CYC:CBC	2.67	0.83
4:X:1082:CYC:HC	4:X:1082:CYC:CMD	1.92	0.83
4:H:1082:CYC:HMA3	4:H:1082:CYC:NB	1.94	0.83
4:V:1153:CYC:HMA3	4:V:1153:CYC:NB	1.91	0.83
4:D:1153:CYC:CMB	1:I:148:VAL:HG11	2.09	0.83
2:T:20:SER:H	2:T:23:GLN:NE2	1.76	0.83
2:V:20:SER:H	2:V:23:GLN:HE21	1.25	0.82
4:L:1153:CYC:HC	4:L:1153:CYC:CMD	1.91	0.82
4:A:1084:CYC:HC	4:A:1084:CYC:HMD1	0.68	0.82
4:X:1082:CYC:HB	4:X:1082:CYC:CMA	1.92	0.82
2:H:57:ARG:HD3	4:K:1084:CYC:O2D	1.80	0.82
4:M:1084:CYC:HMA1	4:M:1084:CYC:HBA1	1.60	0.82
4:T:1082:CYC:HMD1	4:T:1082:CYC:NC	1.93	0.82
1:C:42:LYS:HZ2	2:N:150:PRO:HB3	1.45	0.81
2:J:23:GLN:HA	5:J:2003:HOH:O	1.80	0.81
4:I:1084:CYC:HBA1	4:I:1084:CYC:HMA3	1.61	0.81
4:U:1084:CYC:HHA	4:U:1084:CYC:CBD	2.09	0.81
2:B:20:SER:H	2:B:23:GLN:NE2	1.79	0.80
1:C:148:VAL:HG11	4:J:1153:CYC:HMB1	1.62	0.80
4:J:1153:CYC:HB	4:J:1153:CYC:CMA	1.94	0.80
1:Q:148:VAL:HG11	4:T:1153:CYC:HMB1	1.60	0.80
2:F:20:SER:H	2:F:23:GLN:NE2	1.79	0.80
1:Q:73:ASN:C	4:Q:1084:CYC:HMD3	2.07	0.80
4:J:1153:CYC:HMA3	4:J:1153:CYC:NB	1.96	0.79
2:R:20:SER:H	2:R:23:GLN:NE2	1.79	0.79
4:O:1084:CYC:HBD1	4:O:1084:CYC:HHA	1.65	0.78
4:O:1084:CYC:CMA	4:O:1084:CYC:NB	2.35	0.78
4:R:1153:CYC:HC	4:R:1153:CYC:CMD	1.96	0.78
4:T:1082:CYC:HMA3	4:T:1082:CYC:NB	1.99	0.78
1:C:148:VAL:CG1	4:J:1153:CYC:HMB1	2.15	0.77
2:D:150:PRO:HG2	1:M:39:GLU:HA	1.67	0.77
4:J:1082:CYC:HB	4:J:1082:CYC:CMA	1.97	0.77
2:J:20:SER:H	2:J:23:GLN:HE21	1.33	0.77
2:N:15:ARG:HD3	5:N:2002:HOH:O	1.83	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:L:1153:CYC:HBA1	4:L:1153:CYC:HHA	1.65	0.77
4:J:1153:CYC:HBD1	4:J:1153:CYC:HHA	1.65	0.77
2:L:20:SER:H	2:L:23:GLN:NE2	1.83	0.77
2:H:37:ARG:O	2:H:40:VAL:HG22	1.84	0.76
2:P:20:SER:H	2:P:23:GLN:NE2	1.84	0.76
4:S:1084:CYC:HMA3	4:S:1084:CYC:NB	2.00	0.76
4:N:1153:CYC:CMB	1:W:148:VAL:HG11	2.15	0.76
4:F:1153:CYC:HMA3	4:F:1153:CYC:NB	1.98	0.76
2:V:153:CYS:SG	4:V:1153:CYC:C1C	2.73	0.76
2:V:37:ARG:O	2:V:40:VAL:HG22	1.85	0.76
2:J:111:ASN:HB2	5:J:2014:HOH:O	1.86	0.76
2:B:82:CYS:SG	4:B:1082:CYC:CBC	2.74	0.76
1:E:148:VAL:HG11	4:H:1153:CYC:HMB1	1.68	0.75
1:U:3:THR:HG21	2:V:6:THR:HG21	1.66	0.75
2:F:82:CYS:SG	4:F:1082:CYC:CBC	2.75	0.75
2:H:148:ILE:HD13	4:H:1153:CYC:H3C	1.67	0.75
2:F:39:ASP:OD1	4:F:1153:CYC:HHB	1.87	0.75
4:R:1082:CYC:CMA	4:R:1082:CYC:NB	2.28	0.74
4:F:1082:CYC:CMA	4:F:1082:CYC:NB	2.33	0.74
4:L:1082:CYC:HB	4:L:1082:CYC:CMA	1.99	0.74
2:P:36:LYS:O	2:P:40:VAL:HG13	1.87	0.74
1:E:22:THR:HG23	1:G:4:PRO:HG3	1.69	0.74
2:H:20:SER:N	2:H:23:GLN:HE21	1.84	0.74
2:N:20:SER:H	2:N:23:GLN:HE21	1.33	0.73
2:L:153:CYS:SG	4:L:1153:CYC:HAC1	2.28	0.73
2:V:20:SER:H	2:V:23:GLN:NE2	1.85	0.73
4:W:1084:CYC:HC	4:W:1084:CYC:CMD	2.00	0.73
4:T:1153:CYC:HMD1	4:T:1153:CYC:NC	1.98	0.73
4:P:1153:CYC:HMD1	4:P:1153:CYC:NC	1.98	0.73
1:C:42:LYS:NZ	2:N:150:PRO:HB3	2.04	0.73
4:M:1084:CYC:HHA	4:M:1084:CYC:CBD	2.17	0.73
1:W:57:GLN:HG2	5:W:2004:HOH:O	1.88	0.73
4:W:1084:CYC:CMA	4:W:1084:CYC:NB	2.42	0.73
4:F:1153:CYC:CMB	1:G:148:VAL:HG11	2.18	0.73
4:P:1153:CYC:CMB	1:U:148:VAL:HG11	2.19	0.73
1:Q:7:GLU:O	1:Q:11:VAL:HG23	1.88	0.72
2:H:36:LYS:O	2:H:40:VAL:HG13	1.90	0.72
3:X:11:GLN:HA	5:X:2003:HOH:O	1.88	0.72
2:J:82:CYS:SG	4:J:1082:CYC:CBC	2.77	0.72
4:P:1153:CYC:HB	4:P:1153:CYC:CMA	2.00	0.72
2:P:82:CYS:SG	4:P:1082:CYC:CBC	2.77	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:153:CYS:SG	4:F:1153:CYC:HAC1	2.23	0.72
2:N:39:ASP:OD1	4:N:1153:CYC:HHB	1.90	0.72
2:D:153:CYS:HG	4:D:1153:CYC:HBC2	1.50	0.72
4:N:1082:CYC:HB	4:N:1082:CYC:C3A	2.00	0.71
2:R:39:ASP:OD1	4:R:1153:CYC:HMB3	1.89	0.71
1:G:7:GLU:O	1:G:11:VAL:HG23	1.90	0.71
4:P:1153:CYC:HMA3	4:P:1153:CYC:NB	2.01	0.71
1:A:148:VAL:HG11	4:L:1153:CYC:CMB	2.21	0.71
4:N:1082:CYC:CMA	4:N:1082:CYC:NB	2.36	0.71
4:O:1084:CYC:O1D	2:R:57:ARG:NH1	2.24	0.71
1:S:73:ASN:HA	4:S:1084:CYC:HBD2	1.72	0.70
2:H:153:CYS:SG	4:H:1153:CYC:CBC	2.80	0.70
2:D:11:GLN:HG3	5:D:2002:HOH:O	1.91	0.70
1:M:73:ASN:C	4:M:1084:CYC:HMD3	2.17	0.70
1:K:73:ASN:C	4:K:1084:CYC:HMD3	2.16	0.70
4:X:1153:CYC:HC	4:X:1153:CYC:CMD	2.05	0.70
4:B:1153:CYC:HMD3	4:B:1153:CYC:NC	2.06	0.69
1:Q:73:ASN:O	4:Q:1084:CYC:CMD	2.40	0.69
4:B:1082:CYC:HMD1	4:B:1082:CYC:NC	2.00	0.69
2:D:36:LYS:O	2:D:40:VAL:HG13	1.93	0.69
4:B:1082:CYC:HB	4:B:1082:CYC:CMA	2.05	0.69
1:C:73:ASN:C	4:C:1084:CYC:HMD2	2.17	0.69
3:X:36:LYS:O	3:X:40:VAL:HG13	1.92	0.69
2:D:39:ASP:OD1	4:D:1153:CYC:HHB	1.93	0.68
1:Q:148:VAL:CG1	4:T:1153:CYC:HMB1	2.22	0.68
2:H:39:ASP:OD1	4:H:1153:CYC:HHB	1.93	0.68
2:T:57:ARG:NH1	4:W:1084:CYC:O1D	2.24	0.68
4:H:1153:CYC:HC	4:H:1153:CYC:CMD	2.05	0.68
2:R:36:LYS:O	2:R:40:VAL:HG13	1.94	0.68
2:P:153:CYS:SG	4:P:1153:CYC:CBC	2.81	0.68
2:B:36:LYS:O	2:B:40:VAL:HG13	1.92	0.68
4:J:1153:CYC:HMD1	4:J:1153:CYC:HC	1.58	0.68
4:Q:1084:CYC:CMA	4:Q:1084:CYC:NB	2.44	0.68
4:E:1084:CYC:HC	4:E:1084:CYC:CMD	2.06	0.68
2:N:20:SER:H	2:N:23:GLN:NE2	1.90	0.68
4:J:1153:CYC:CBC	4:J:1153:CYC:CHD	2.71	0.68
3:X:40:VAL:HG12	4:X:1153:CYC:HBC1	1.75	0.68
2:D:20:SER:H	2:D:23:GLN:NE2	1.89	0.68
4:N:1153:CYC:HMD3	4:N:1153:CYC:NC	2.08	0.68
4:H:1082:CYC:CMD	4:H:1082:CYC:NC	2.35	0.67
3:X:20:SER:H	3:X:23:GLN:NE2	1.92	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:N:153:CYS:SG	4:N:1153:CYC:C1C	2.83	0.67
4:V:1082:CYC:HHA	4:V:1082:CYC:CBD	2.23	0.67
4:A:1084:CYC:O2D	2:D:57:ARG:HD3	1.93	0.67
4:X:1153:CYC:HB	4:X:1153:CYC:CMA	2.04	0.67
4:B:1153:CYC:CMB	1:K:148:VAL:HG11	2.24	0.67
3:X:82:CYS:SG	4:X:1082:CYC:CBC	2.83	0.67
4:F:1153:CYC:HC	4:F:1153:CYC:CMD	2.07	0.67
2:R:37:ARG:O	2:R:40:VAL:HG22	1.95	0.67
4:X:1082:CYC:HMD1	4:X:1082:CYC:NC	2.05	0.67
4:H:1082:CYC:HBD2	4:H:1082:CYC:HHA	1.75	0.67
4:R:1082:CYC:CMD	4:R:1082:CYC:NC	2.36	0.67
2:D:153:CYS:SG	4:D:1153:CYC:CAC	2.82	0.67
3:X:129:VAL:O	3:X:132:LYS:HG2	1.95	0.67
4:A:1084:CYC:CMD	4:A:1084:CYC:NC	2.23	0.66
2:V:36:LYS:O	2:V:40:VAL:HG13	1.96	0.66
4:L:1082:CYC:CMD	4:L:1082:CYC:NC	2.41	0.66
1:C:7:GLU:O	1:C:11:VAL:HG23	1.95	0.66
1:I:128:TRP:CE3	4:I:1084:CYC:HMC3	2.31	0.66
2:R:20:SER:N	2:R:23:GLN:HE21	1.92	0.66
2:V:40:VAL:HG11	2:V:156:LEU:HD21	1.78	0.66
3:X:54:ASN:HA	5:X:2006:HOH:O	1.93	0.66
1:C:73:ASN:O	4:C:1084:CYC:HMD2	1.96	0.66
4:U:1084:CYC:O2D	3:X:57:ARG:HD3	1.95	0.66
4:F:1082:CYC:HC	4:F:1082:CYC:CMD	2.09	0.66
1:O:73:ASN:C	4:O:1084:CYC:HMD3	2.20	0.66
1:W:67:THR:HG22	5:W:2006:HOH:O	1.96	0.66
1:C:3:THR:HG21	2:D:6:THR:HG21	1.77	0.66
2:N:21:THR:HG23	5:N:2003:HOH:O	1.95	0.66
2:R:153:CYS:SG	4:R:1153:CYC:CAC	2.84	0.65
4:W:1084:CYC:HMD1	4:W:1084:CYC:NC	2.02	0.65
1:Q:4:PRO:HG2	1:Q:30:ARG:HD3	1.76	0.65
1:C:39:GLU:HB2	2:N:150:PRO:HD3	1.76	0.65
1:C:72:PRO:O	4:C:1084:CYC:HBD1	1.96	0.65
4:T:1153:CYC:HB	4:T:1153:CYC:CMA	2.07	0.65
1:W:3:THR:HG21	3:X:6:THR:HG21	1.79	0.65
2:R:153:CYS:SG	4:R:1153:CYC:C2C	2.84	0.65
2:T:36:LYS:O	2:T:40:VAL:HG13	1.96	0.65
2:N:37:ARG:HA	2:N:156:LEU:HD21	1.77	0.65
1:Q:73:ASN:O	4:Q:1084:CYC:HMD1	1.96	0.65
2:N:37:ARG:O	2:N:40:VAL:HG22	1.96	0.65
1:O:148:VAL:HG11	4:V:1153:CYC:CMB	2.27	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:128:TRP:CE3	4:C:1084:CYC:HMC2	2.32	0.64
2:H:39:ASP:OD1	4:H:1153:CYC:HMB3	1.96	0.64
2:H:153:CYS:SG	4:H:1153:CYC:C1C	2.85	0.64
5:U:2002:HOH:O	3:X:76:SER:HB2	1.97	0.64
2:F:153:CYS:SG	4:F:1153:CYC:CBC	2.84	0.64
4:U:1084:CYC:HBD2	4:U:1084:CYC:CHA	2.21	0.64
2:V:153:CYS:SG	4:V:1153:CYC:C2C	2.85	0.64
2:D:153:CYS:CB	4:D:1153:CYC:HBC2	2.21	0.64
1:E:148:VAL:CG1	4:H:1153:CYC:HMB1	2.26	0.64
4:L:1153:CYC:HMD1	4:L:1153:CYC:NC	2.05	0.64
2:V:40:VAL:HG11	2:V:156:LEU:CD2	2.27	0.64
3:X:37:ARG:HA	3:X:156:LEU:HD21	1.79	0.64
4:X:1082:CYC:HMA3	4:X:1082:CYC:NB	2.04	0.64
2:L:36:LYS:O	2:L:40:VAL:HG13	1.96	0.64
2:R:153:CYS:CB	4:R:1153:CYC:HBC2	2.26	0.64
2:V:148:ILE:HD13	4:V:1153:CYC:H3C	1.79	0.64
2:L:148:ILE:CD1	4:L:1153:CYC:HHD	2.28	0.63
4:J:1082:CYC:HMA1	4:J:1082:CYC:NB	2.06	0.63
3:X:82:CYS:SG	4:X:1082:CYC:H2C	2.38	0.63
4:E:1084:CYC:CMA	4:E:1084:CYC:NB	2.47	0.63
4:D:1082:CYC:HMD1	4:D:1082:CYC:NC	2.02	0.63
4:G:1084:CYC:O1D	2:J:57:ARG:NH1	2.32	0.63
2:D:150:PRO:CG	1:M:39:GLU:HA	2.28	0.63
2:J:20:SER:H	2:J:23:GLN:NE2	1.96	0.62
3:X:148:ILE:HG12	5:X:2014:HOH:O	1.99	0.62
2:P:153:CYS:SG	4:P:1153:CYC:C1C	2.87	0.62
3:X:82:CYS:SG	4:X:1082:CYC:C2C	2.87	0.62
2:F:20:SER:N	2:F:23:GLN:HE21	1.90	0.62
1:W:2:LYS:HB3	5:W:2001:HOH:O	1.99	0.62
1:O:3:THR:HG21	2:P:6:THR:HG21	1.80	0.62
2:N:20:SER:N	2:N:23:GLN:HE21	1.97	0.62
4:N:1153:CYC:HB	4:N:1153:CYC:HMA1	1.64	0.62
2:R:153:CYS:HG	4:R:1153:CYC:HBC2	1.59	0.62
4:D:1153:CYC:HMB1	1:I:148:VAL:CG1	2.30	0.62
4:U:1084:CYC:CMD	4:U:1084:CYC:NC	2.46	0.62
1:C:148:VAL:HG11	4:J:1153:CYC:HMB3	1.82	0.62
2:J:37:ARG:HA	2:J:156:LEU:HD21	1.81	0.62
2:J:153:CYS:SG	4:J:1153:CYC:CBC	2.88	0.62
2:D:88:ILE:HG21	4:D:1082:CYC:HMB2	1.82	0.61
2:D:153:CYS:SG	4:D:1153:CYC:C1C	2.88	0.61
4:M:1084:CYC:CMD	4:M:1084:CYC:NC	2.48	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:3:THR:HG21	2:F:6:THR:HG21	1.81	0.61
2:F:148:ILE:HD13	4:F:1153:CYC:H3C	1.83	0.61
1:U:83:LYS:NZ	4:U:1084:CYC:O1A	2.28	0.61
3:X:37:ARG:O	3:X:40:VAL:HG22	2.00	0.61
2:J:35:ASN:HB2	4:J:1153:CYC:O1D	2.01	0.61
4:P:1153:CYC:HMB1	1:U:148:VAL:HG11	1.83	0.61
2:V:20:SER:N	2:V:23:GLN:HE21	1.98	0.61
2:D:82:CYS:SG	4:D:1082:CYC:HBC2	2.36	0.61
4:P:1082:CYC:HC	4:P:1082:CYC:CMD	2.08	0.61
2:D:37:ARG:HA	2:D:156:LEU:HD21	1.82	0.61
4:D:1153:CYC:HMB1	1:I:148:VAL:HG11	1.80	0.61
4:X:1153:CYC:C1C	5:X:2015:HOH:O	2.49	0.61
2:V:39:ASP:OD1	4:V:1153:CYC:HBB	2.00	0.60
3:X:77:THR:HG22	5:X:2007:HOH:O	2.01	0.60
2:B:37:ARG:HA	2:B:156:LEU:HD21	1.83	0.60
2:L:37:ARG:O	2:L:40:VAL:HG22	2.01	0.60
2:N:153:CYS:SG	4:N:1153:CYC:C2C	2.89	0.60
2:B:39:ASP:OD1	4:B:1153:CYC:HMB3	2.02	0.60
2:R:153:CYS:SG	4:R:1153:CYC:H2C	2.41	0.60
2:H:82:CYS:CB	4:H:1082:CYC:HAC2	2.29	0.60
2:L:37:ARG:HA	2:L:156:LEU:HD21	1.83	0.60
4:B:1153:CYC:HMB1	1:K:148:VAL:HG11	1.83	0.60
2:B:20:SER:N	2:B:23:GLN:HE21	1.94	0.60
2:F:37:ARG:HA	2:F:156:LEU:HD21	1.84	0.60
4:O:1084:CYC:CMD	4:O:1084:CYC:NC	2.43	0.60
4:R:1153:CYC:HB	4:R:1153:CYC:CMA	2.11	0.60
1:S:40:ALA:HB2	1:S:146:ALA:HB1	1.84	0.59
2:B:82:CYS:CB	4:B:1082:CYC:HAC2	2.28	0.59
2:F:36:LYS:O	2:F:40:VAL:HG13	2.02	0.59
2:R:153:CYS:HG	4:R:1153:CYC:CBC	2.13	0.59
4:C:1084:CYC:HMD3	4:C:1084:CYC:NC	2.15	0.59
4:D:1153:CYC:HMB3	1:I:148:VAL:HG11	1.85	0.59
4:M:1084:CYC:CMA	4:M:1084:CYC:HBA1	2.32	0.59
1:A:3:THR:HG21	2:B:6:THR:HG21	1.84	0.58
1:I:7:GLU:O	1:I:11:VAL:HG23	2.02	0.58
3:X:20:SER:N	3:X:23:GLN:HE21	1.98	0.58
4:J:1153:CYC:CHD	4:J:1153:CYC:HBC3	2.33	0.58
1:O:72:PRO:O	4:O:1084:CYC:HAD1	2.03	0.58
2:R:37:ARG:HA	2:R:156:LEU:HD21	1.85	0.58
4:T:1082:CYC:HMA1	4:T:1082:CYC:HBA1	1.86	0.58
2:N:36:LYS:O	2:N:40:VAL:HG13	2.03	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:P:37:ARG:O	2:P:40:VAL:HG22	2.02	0.58
2:R:148:ILE:HD13	4:R:1153:CYC:H3C	1.85	0.58
2:B:153:CYS:SG	4:B:1153:CYC:C2C	2.91	0.58
2:H:69:PRO:HA	2:H:74:TYR:CG	2.39	0.58
4:N:1153:CYC:HMB1	1:W:148:VAL:HG11	1.84	0.58
3:X:82:CYS:SG	4:X:1082:CYC:C3C	2.92	0.58
1:E:40:ALA:HB2	1:E:146:ALA:HB1	1.86	0.58
2:R:39:ASP:OD1	4:R:1153:CYC:HBB	2.04	0.58
5:C:2003:HOH:O	2:D:21:THR:HB	2.04	0.57
4:K:1084:CYC:CMA	4:K:1084:CYC:NB	2.46	0.57
1:I:3:THR:HG21	2:J:6:THR:HG21	1.85	0.57
2:P:37:ARG:HA	2:P:156:LEU:HD21	1.86	0.57
4:X:1153:CYC:HMD1	4:X:1153:CYC:NC	2.12	0.57
1:S:128:TRP:CE3	4:S:1084:CYC:HMC2	2.38	0.57
2:P:120:LEU:HD22	4:P:1082:CYC:HBD1	1.86	0.57
2:T:37:ARG:HA	2:T:156:LEU:HD21	1.87	0.57
1:C:36:ALA:HA	5:C:2002:HOH:O	2.03	0.57
4:E:1084:CYC:HMA1	4:E:1084:CYC:HBA1	1.86	0.57
1:I:135:TYR:HB2	5:I:2012:HOH:O	2.04	0.57
2:J:20:SER:N	2:J:23:GLN:HE21	2.01	0.57
1:M:7:GLU:O	1:M:11:VAL:HG23	2.04	0.57
1:G:72:PRO:O	4:G:1084:CYC:HAD1	2.05	0.57
1:W:7:GLU:O	1:W:11:VAL:HG23	2.05	0.57
1:M:22:THR:HG23	1:W:4:PRO:HG3	1.86	0.56
1:Q:73:ASN:C	4:Q:1084:CYC:CMD	2.78	0.56
4:C:1084:CYC:HC	4:C:1084:CYC:CMD	2.15	0.56
4:D:1082:CYC:HB	4:D:1082:CYC:C3A	2.18	0.56
1:K:40:ALA:HB2	1:K:146:ALA:HB1	1.86	0.56
2:D:151:GLY:HA3	4:D:1153:CYC:CMD	2.35	0.56
2:N:76:SER:HB2	5:Q:2001:HOH:O	2.05	0.56
2:P:20:SER:N	2:P:23:GLN:HE21	1.99	0.56
2:R:153:CYS:SG	4:R:1153:CYC:C1C	2.93	0.56
4:T:1153:CYC:HMA1	4:T:1153:CYC:NB	2.15	0.56
2:V:37:ARG:HA	2:V:156:LEU:HD21	1.87	0.56
1:A:125:SER:HB3	1:A:128:TRP:CE2	2.41	0.56
1:Q:72:PRO:O	4:Q:1084:CYC:HAD2	2.04	0.56
1:S:3:THR:HG21	2:T:6:THR:HG21	1.87	0.56
1:U:125:SER:HB3	1:U:128:TRP:CE2	2.40	0.56
1:A:73:ASN:C	4:A:1084:CYC:HMD3	2.30	0.56
4:L:1153:CYC:HMA3	4:L:1153:CYC:NB	2.13	0.56
2:T:40:VAL:HG11	2:T:156:LEU:CD2	2.36	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:72:PRO:O	4:C:1084:CYC:CBD	2.53	0.56
4:V:1082:CYC:CMD	4:V:1082:CYC:NC	2.43	0.56
1:A:7:GLU:O	1:A:11:VAL:HG23	2.05	0.56
2:B:37:ARG:O	2:B:40:VAL:HG22	2.05	0.56
4:F:1153:CYC:HMD1	4:F:1153:CYC:NC	2.12	0.56
2:H:37:ARG:HA	2:H:156:LEU:HD21	1.86	0.56
4:I:1084:CYC:CMD	4:I:1084:CYC:NC	2.51	0.56
2:T:148:ILE:CD1	4:T:1153:CYC:HHD	2.36	0.56
4:U:1084:CYC:CMA	4:U:1084:CYC:NB	2.55	0.56
4:E:1084:CYC:HMD1	4:E:1084:CYC:NC	2.13	0.56
1:K:73:ASN:O	4:K:1084:CYC:HMD1	2.06	0.56
1:G:125:SER:HB3	1:G:128:TRP:CE2	2.41	0.55
4:W:1084:CYC:HB	4:W:1084:CYC:C3A	2.19	0.55
2:H:57:ARG:NH1	4:K:1084:CYC:O1D	2.39	0.55
2:P:82:CYS:CB	4:P:1082:CYC:HAC2	2.28	0.55
2:J:36:LYS:O	2:J:40:VAL:HG13	2.07	0.55
4:J:1082:CYC:CMD	4:J:1082:CYC:NC	2.53	0.55
1:U:40:ALA:HB2	1:U:146:ALA:HB1	1.87	0.55
2:H:153:CYS:SG	4:H:1153:CYC:C2C	2.95	0.55
4:M:1084:CYC:HBD2	4:M:1084:CYC:CHA	2.32	0.55
1:K:73:ASN:O	4:K:1084:CYC:CMD	2.55	0.55
4:K:1084:CYC:HHA	4:K:1084:CYC:HBD1	1.89	0.55
4:B:1153:CYC:HC	4:B:1153:CYC:C2D	2.20	0.55
2:L:82:CYS:CB	4:L:1082:CYC:HAC2	2.28	0.54
4:L:1153:CYC:HB	4:L:1153:CYC:CMA	2.13	0.54
4:N:1153:CYC:HMB1	1:W:148:VAL:CG1	2.38	0.54
4:B:1153:CYC:HMA1	4:B:1153:CYC:HB	1.72	0.54
1:M:149:GLU:HG3	4:X:1153:CYC:HBB1	1.90	0.54
2:P:69:PRO:HA	2:P:74:TYR:CG	2.43	0.54
2:L:88:ILE:HG21	4:L:1082:CYC:HMB2	1.88	0.54
2:P:153:CYS:SG	4:P:1153:CYC:C2C	2.95	0.54
4:A:1084:CYC:HBB2	2:D:75:THR:HA	1.89	0.54
2:P:36:LYS:O	2:P:40:VAL:CG1	2.55	0.54
2:T:148:ILE:HD13	4:T:1153:CYC:H3C	1.90	0.54
1:E:1:MET:HG2	1:E:105:GLY:HA3	1.91	0.53
2:L:39:ASP:OD1	4:L:1153:CYC:HHB	2.07	0.53
2:N:151:GLY:HA3	4:N:1153:CYC:CMD	2.38	0.53
4:O:1084:CYC:CMA	4:O:1084:CYC:HBA1	2.39	0.53
4:P:1153:CYC:HMB1	1:U:148:VAL:CG1	2.38	0.53
2:V:82:CYS:SG	4:V:1082:CYC:HAC1	2.31	0.53
4:D:1153:CYC:CMA	4:D:1153:CYC:NB	2.55	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:H:1153:CYC:HB	4:H:1153:CYC:CMA	2.15	0.53
4:K:1084:CYC:HHA	4:K:1084:CYC:CBD	2.38	0.53
1:U:3:THR:HG21	2:V:6:THR:CG2	2.38	0.53
4:G:1084:CYC:HAA2	5:G:2005:HOH:O	2.08	0.53
2:B:57:ARG:NH1	4:E:1084:CYC:O1D	2.41	0.53
1:M:128:TRP:CE3	4:M:1084:CYC:HMC2	2.44	0.53
3:X:40:VAL:CG2	3:X:97:VAL:HG11	2.39	0.53
4:J:1082:CYC:CMA	5:J:2017:HOH:O	2.56	0.53
4:S:1084:CYC:HC	4:S:1084:CYC:CMD	2.16	0.53
1:E:3:THR:HG21	2:F:6:THR:CG2	2.39	0.53
4:L:1153:CYC:CBB	4:L:1153:CYC:OB	2.56	0.53
1:M:87:ASP:OD2	4:M:1084:CYC:NA	2.42	0.53
1:W:5:LEU:HB2	3:X:3:ASP:OD2	2.09	0.53
3:X:148:ILE:HD13	4:X:1153:CYC:H3C	1.90	0.53
1:S:1:MET:HG2	1:S:105:GLY:HA3	1.91	0.52
2:T:151:GLY:O	4:T:1153:CYC:OC	2.27	0.52
2:D:20:SER:N	2:D:23:GLN:HE21	1.98	0.52
1:I:40:ALA:HB2	1:I:146:ALA:HB1	1.92	0.52
1:M:40:ALA:HB2	1:M:146:ALA:HB1	1.92	0.52
2:N:151:GLY:O	4:N:1153:CYC:OC	2.28	0.52
4:I:1084:CYC:CMA	4:I:1084:CYC:HBA1	2.37	0.52
2:L:148:ILE:HD13	4:L:1153:CYC:H3C	1.91	0.52
2:P:90:LEU:O	2:P:94:THR:HG23	2.09	0.52
3:X:153:CYS:HG	4:X:1153:CYC:C1C	2.22	0.52
2:D:37:ARG:O	2:D:40:VAL:HG22	2.09	0.52
2:D:153:CYS:SG	4:D:1153:CYC:C2C	2.97	0.52
2:L:148:ILE:HD11	4:L:1153:CYC:HHD	1.91	0.52
1:O:65:TYR:HB2	1:O:69:MET:HE3	1.91	0.52
1:W:40:ALA:HB2	1:W:146:ALA:HB1	1.92	0.52
1:M:59:VAL:HG11	4:M:1084:CYC:HMC3	1.90	0.52
3:X:153:CYS:SG	4:X:1153:CYC:C3C	2.97	0.52
4:A:1084:CYC:HBD2	4:A:1084:CYC:HHA	1.91	0.52
1:E:73:ASN:C	4:E:1084:CYC:HMD3	2.34	0.52
2:H:82:CYS:SG	4:H:1082:CYC:H2C	2.50	0.52
2:L:39:ASP:OD1	4:L:1153:CYC:HMB3	2.08	0.52
2:L:82:CYS:SG	4:L:1082:CYC:C3C	2.96	0.52
4:D:1153:CYC:HC	4:D:1153:CYC:CMD	2.15	0.51
4:I:1084:CYC:O1D	2:L:57:ARG:NH1	2.37	0.51
2:T:77:THR:HG22	5:W:2007:HOH:O	2.10	0.51
4:L:1082:CYC:HMA3	4:L:1082:CYC:NB	2.17	0.51
2:B:151:GLY:O	4:B:1153:CYC:OC	2.28	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:B:1082:CYC:HMA3	4:B:1082:CYC:NB	2.15	0.51
4:F:1153:CYC:HMB1	1:G:148:VAL:HG11	1.93	0.51
2:N:39:ASP:HA	5:N:2006:HOH:O	2.11	0.51
1:O:148:VAL:HG11	4:V:1153:CYC:HMB1	1.92	0.51
1:Q:1:MET:HG2	1:Q:105:GLY:HA3	1.92	0.51
2:J:69:PRO:HA	2:J:74:TYR:CG	2.46	0.51
4:N:1153:CYC:OB	4:N:1153:CYC:CBB	2.58	0.51
4:X:1153:CYC:CMC	5:X:2015:HOH:O	2.58	0.51
1:O:7:GLU:O	1:O:11:VAL:HG23	2.09	0.51
2:B:82:CYS:SG	4:B:1082:CYC:H2C	2.50	0.51
4:F:1153:CYC:HMB1	1:G:148:VAL:CG1	2.41	0.51
2:J:68:ALA:O	2:J:69:PRO:C	2.51	0.51
1:W:125:SER:HB3	1:W:128:TRP:CE2	2.46	0.51
1:G:3:THR:HG21	2:H:6:THR:HG21	1.93	0.51
1:U:73:ASN:C	4:U:1084:CYC:HMD3	2.36	0.51
1:G:1:MET:HG2	1:G:105:GLY:HA3	1.92	0.51
1:M:152:SER:HA	5:M:2011:HOH:O	2.11	0.51
2:B:82:CYS:SG	4:B:1082:CYC:C3C	2.95	0.50
1:E:125:SER:HB3	1:E:128:TRP:CE2	2.46	0.50
2:J:148:ILE:HG21	4:J:1153:CYC:H3C	1.92	0.50
1:C:40:ALA:HB2	1:C:146:ALA:HB1	1.92	0.50
2:H:40:VAL:CG2	2:H:41:VAL:N	2.75	0.50
1:M:73:ASN:O	4:M:1084:CYC:HMD3	2.11	0.50
2:T:127:VAL:HG22	4:T:1082:CYC:H3C	1.93	0.50
1:Q:128:TRP:CE3	4:Q:1084:CYC:HMC2	2.46	0.50
1:A:148:VAL:CG1	4:L:1153:CYC:HMB1	2.41	0.50
4:C:1084:CYC:CMA	4:C:1084:CYC:NB	2.56	0.50
4:X:1153:CYC:HMA3	4:X:1153:CYC:NB	2.11	0.50
1:C:4:PRO:HG3	1:I:22:THR:HG23	1.92	0.50
2:J:37:ARG:O	2:J:40:VAL:HG22	2.11	0.50
4:N:1153:CYC:HC	4:N:1153:CYC:C2D	2.24	0.50
1:Q:73:ASN:O	4:Q:1084:CYC:HMD3	2.04	0.50
1:I:73:ASN:C	4:I:1084:CYC:HMD3	2.36	0.50
2:J:151:GLY:HA3	4:J:1153:CYC:CMD	2.42	0.50
1:M:125:SER:HB3	1:M:128:TRP:CE2	2.46	0.50
4:P:1082:CYC:HMD1	4:P:1082:CYC:NC	2.15	0.50
1:U:65:TYR:HB2	1:U:69:MET:HE3	1.94	0.50
2:B:69:PRO:HA	2:B:74:TYR:CG	2.47	0.50
2:D:39:ASP:OD2	4:D:1153:CYC:NA	2.44	0.50
4:N:1082:CYC:CMD	4:N:1082:CYC:NC	2.44	0.50
4:P:1082:CYC:CMA	4:P:1082:CYC:NB	2.58	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:W:73:ASN:HA	4:W:1084:CYC:HBD2	1.92	0.50
2:V:153:CYS:SG	4:V:1153:CYC:H2C	2.52	0.50
1:O:62:LYS:HG2	1:O:63:PHE:CE2	2.47	0.50
2:F:69:PRO:HA	2:F:74:TYR:CG	2.47	0.49
4:R:1153:CYC:HMD1	4:R:1153:CYC:NC	2.11	0.49
1:A:47:LYS:O	1:A:48:ALA:C	2.54	0.49
4:B:1153:CYC:HMB1	1:K:148:VAL:CG1	2.43	0.49
2:V:39:ASP:OD1	4:V:1153:CYC:HMB3	2.13	0.49
2:B:153:CYS:SG	4:B:1153:CYC:C1C	3.01	0.49
2:R:40:VAL:HG11	2:R:156:LEU:CD2	2.42	0.49
2:D:82:CYS:SG	4:D:1082:CYC:C3C	2.93	0.49
2:L:40:VAL:HG11	2:L:156:LEU:CD2	2.43	0.49
2:T:57:ARG:HD3	4:W:1084:CYC:O1D	2.13	0.49
2:H:153:CYS:SG	4:H:1153:CYC:HBC2	2.52	0.49
1:K:72:PRO:O	4:K:1084:CYC:HAD1	2.11	0.49
4:K:1084:CYC:CMD	4:K:1084:CYC:NC	2.60	0.49
1:M:3:THR:HG21	2:N:6:THR:HG21	1.93	0.49
4:P:1153:CYC:HMB3	1:U:148:VAL:HG11	1.94	0.49
1:A:73:ASN:O	4:A:1084:CYC:HMD3	2.12	0.49
2:B:82:CYS:SG	4:B:1082:CYC:C2C	3.00	0.49
1:E:148:VAL:HG11	4:H:1153:CYC:C2B	2.42	0.49
4:R:1082:CYC:HBD2	4:R:1082:CYC:HHA	1.95	0.49
2:T:20:SER:N	2:T:23:GLN:HE21	1.96	0.49
1:W:6:THR:HG21	5:X:2001:HOH:O	2.13	0.49
2:D:69:PRO:HA	2:D:74:TYR:CG	2.48	0.48
2:H:74:TYR:O	2:H:75:THR:OG1	2.29	0.48
1:A:148:VAL:HG11	4:L:1153:CYC:HMB1	1.93	0.48
2:V:63:GLN:HG2	2:V:65:GLN:HE22	1.79	0.48
4:A:1084:CYC:HMD1	4:A:1084:CYC:C4C	2.32	0.48
1:C:148:VAL:HG12	4:J:1153:CYC:HMB1	1.93	0.48
2:H:82:CYS:SG	4:H:1082:CYC:C3C	2.98	0.48
2:J:40:VAL:HG11	2:J:156:LEU:CD2	2.43	0.48
2:L:74:TYR:O	2:L:75:THR:OG1	2.25	0.48
1:S:73:ASN:HA	4:S:1084:CYC:HMD3	1.95	0.48
1:K:7:GLU:O	1:K:11:VAL:HG23	2.14	0.48
3:X:40:VAL:CG2	3:X:41:VAL:N	2.75	0.48
4:J:1153:CYC:HHA	4:J:1153:CYC:CBD	2.37	0.48
2:N:40:VAL:CG2	2:N:41:VAL:N	2.77	0.48
4:N:1082:CYC:NB	4:N:1082:CYC:C3A	2.74	0.48
2:T:153:CYS:SG	4:T:1153:CYC:C1C	3.02	0.48
2:F:37:ARG:O	2:F:40:VAL:HG22	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:22:THR:HG23	1:G:4:PRO:CG	2.40	0.47
1:K:73:ASN:C	4:K:1084:CYC:CMD	2.85	0.47
1:M:5:LEU:HB2	2:N:3:ASP:OD2	2.14	0.47
2:N:40:VAL:HG12	4:N:1153:CYC:HBC1	1.96	0.47
3:X:72:ASN:OD1	4:X:1082:CYC:CMD	2.62	0.47
1:C:1:MET:HG2	1:C:105:GLY:HA3	1.96	0.47
2:L:35:ASN:OD1	4:L:1153:CYC:HAA1	2.14	0.47
4:L:1153:CYC:HHA	4:L:1153:CYC:CBA	2.35	0.47
4:N:1153:CYC:OB	4:N:1153:CYC:HBB2	2.14	0.47
1:O:125:SER:HB3	1:O:128:TRP:CE2	2.49	0.47
2:T:148:ILE:HD13	4:T:1153:CYC:HHD	1.94	0.47
1:W:3:THR:HG21	3:X:6:THR:CG2	2.44	0.47
4:X:1153:CYC:HMC1	5:X:2015:HOH:O	2.13	0.47
2:B:40:VAL:CG2	2:B:41:VAL:N	2.78	0.47
2:V:40:VAL:CG1	2:V:156:LEU:HD21	2.44	0.47
3:X:149:THR:HG21	4:X:1153:CYC:HAD1	1.96	0.47
2:F:153:CYS:SG	4:F:1153:CYC:C2C	3.03	0.47
4:F:1153:CYC:HMB3	1:G:148:VAL:HG11	1.95	0.47
4:J:1153:CYC:HBC2	4:J:1153:CYC:HHD	1.96	0.47
2:L:20:SER:N	2:L:23:GLN:HE21	2.00	0.47
2:L:153:CYS:SG	4:L:1153:CYC:C1C	3.02	0.47
2:V:69:PRO:HA	2:V:74:TYR:CG	2.50	0.47
2:D:82:CYS:SG	4:D:1082:CYC:C2C	3.02	0.47
2:H:72:ASN:OD1	4:H:1082:CYC:HMD3	2.15	0.47
2:L:47:ASN:O	2:L:48:ALA:C	2.58	0.47
2:T:77:THR:CG2	5:W:2007:HOH:O	2.62	0.47
1:W:1:MET:HG2	1:W:105:GLY:HA3	1.97	0.47
1:K:125:SER:HB3	1:K:128:TRP:CE2	2.49	0.47
1:Q:3:THR:HG21	2:R:6:THR:HG21	1.96	0.47
2:F:40:VAL:CG2	2:F:97:VAL:HG11	2.45	0.47
1:A:1:MET:HG2	1:A:105:GLY:HA3	1.95	0.46
1:I:125:SER:HB3	1:I:128:TRP:CE2	2.50	0.46
1:E:73:ASN:O	4:E:1084:CYC:HMD1	2.16	0.46
1:M:59:VAL:CG1	4:M:1084:CYC:HMC3	2.44	0.46
1:M:73:ASN:O	4:M:1084:CYC:CMD	2.62	0.46
1:O:148:VAL:CG1	4:V:1153:CYC:HMB1	2.45	0.46
2:T:40:VAL:HG11	2:T:156:LEU:HD21	1.97	0.46
1:G:47:LYS:O	1:G:48:ALA:C	2.59	0.46
1:G:120:ARG:HG3	1:G:120:ARG:HH11	1.79	0.46
1:I:3:THR:HG21	2:J:6:THR:CG2	2.43	0.46
4:L:1153:CYC:OB	4:L:1153:CYC:HBB2	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Q:125:SER:HB3	1:Q:128:TRP:CE2	2.50	0.46
2:B:153:CYS:SG	4:B:1153:CYC:H2C	2.56	0.46
1:C:42:LYS:NZ	5:C:2003:HOH:O	2.44	0.46
1:E:72:PRO:HB2	4:E:1084:CYC:O1D	2.16	0.46
2:H:82:CYS:SG	4:H:1082:CYC:C2C	3.04	0.46
2:P:39:ASP:OD1	4:P:1153:CYC:HHB	2.16	0.46
2:R:123:PRO:O	2:R:126:SER:HB2	2.15	0.46
2:P:148:ILE:CD1	4:P:1153:CYC:HHD	2.46	0.46
3:X:40:VAL:HG11	3:X:156:LEU:CD2	2.45	0.46
1:O:3:THR:HG21	2:P:6:THR:CG2	2.46	0.46
2:T:149:THR:OG1	4:T:1153:CYC:O1A	2.30	0.46
2:T:153:CYS:SG	4:T:1153:CYC:C2C	3.03	0.46
2:V:127:VAL:HG22	4:V:1082:CYC:H3C	1.98	0.46
3:X:69:PRO:HA	3:X:74:TYR:CG	2.51	0.46
1:G:79:ARG:HE	1:G:79:ARG:HB3	1.71	0.46
2:N:69:PRO:HA	2:N:74:TYR:CG	2.51	0.46
2:N:148:ILE:CD1	4:N:1153:CYC:HHD	2.46	0.46
4:O:1084:CYC:HBA1	4:O:1084:CYC:HMA1	1.97	0.46
2:T:37:ARG:O	2:T:40:VAL:HG22	2.16	0.46
1:U:1:MET:HG2	1:U:105:GLY:HA3	1.97	0.46
2:V:40:VAL:CG2	2:V:41:VAL:N	2.79	0.46
2:H:148:ILE:HD13	4:H:1153:CYC:C3C	2.41	0.46
1:K:11:VAL:HG12	1:K:15:GLN:NE2	2.31	0.46
4:W:1084:CYC:HMA1	4:W:1084:CYC:HBA1	1.98	0.46
1:E:7:GLU:O	1:E:11:VAL:HG23	2.16	0.45
2:L:90:LEU:O	2:L:94:THR:HG23	2.16	0.45
2:L:148:ILE:CD1	4:L:1153:CYC:CHD	2.94	0.45
1:C:79:ARG:HH12	4:C:1084:CYC:HBD2	1.80	0.45
4:F:1082:CYC:HHA	4:F:1082:CYC:HAD2	1.64	0.45
2:J:39:ASP:OD1	4:J:1153:CYC:HHB	2.17	0.45
2:J:82:CYS:SG	4:J:1082:CYC:C3C	2.98	0.45
2:J:124:GLY:O	2:J:125:SER:C	2.59	0.45
4:M:1084:CYC:CMA	4:M:1084:CYC:NB	2.60	0.45
3:X:39:ASP:CG	3:X:148:ILE:HD11	2.42	0.45
2:J:148:ILE:HD13	4:J:1153:CYC:H3C	1.98	0.45
4:O:1084:CYC:HBB3	4:O:1084:CYC:HMB1	1.98	0.45
2:T:13:ASP:OD1	2:V:74:TYR:OH	2.26	0.45
1:I:120:ARG:HG3	1:I:120:ARG:HH11	1.81	0.45
1:M:65:TYR:HB2	1:M:69:MET:HE3	1.97	0.45
1:O:5:LEU:HB2	2:P:3:ASP:OD2	2.17	0.45
2:B:40:VAL:CG2	2:B:97:VAL:HG11	2.46	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:65:TYR:HB2	1:I:69:MET:HE3	1.98	0.45
1:M:21:SER:HB3	5:M:2003:HOH:O	2.15	0.45
1:O:111:LEU:C	1:O:111:LEU:HD23	2.42	0.45
4:H:1153:CYC:HBC2	4:H:1153:CYC:H2C	1.90	0.45
4:L:1153:CYC:HBA1	4:L:1153:CYC:CHA	2.42	0.45
2:N:73:ALA:HB2	4:N:1082:CYC:OC	2.17	0.45
2:V:82:CYS:SG	4:V:1082:CYC:C2C	3.04	0.45
2:V:82:CYS:SG	4:V:1082:CYC:C3C	2.97	0.45
2:F:153:CYS:SG	4:F:1153:CYC:C1C	3.05	0.45
3:X:69:PRO:HA	3:X:74:TYR:CD2	2.52	0.45
1:A:26:VAL:HG22	1:K:26:VAL:HG13	1.99	0.44
1:E:79:ARG:HE	1:E:79:ARG:HB3	1.68	0.44
1:M:1:MET:HG2	1:M:105:GLY:HA3	1.99	0.44
1:M:128:TRP:CD2	4:M:1084:CYC:HMC2	2.52	0.44
1:S:79:ARG:HE	1:S:79:ARG:HB3	1.72	0.44
2:B:144:ASP:HA	2:B:145:PRO:HD3	1.85	0.44
2:L:88:ILE:CG2	4:L:1082:CYC:HMB2	2.47	0.44
1:Q:148:VAL:HG11	4:T:1153:CYC:HMB3	1.92	0.44
2:D:72:ASN:OD1	4:D:1082:CYC:HMD1	2.16	0.44
1:K:79:ARG:HE	1:K:79:ARG:HB3	1.68	0.44
1:M:115:ILE:HD12	1:M:115:ILE:HA	1.88	0.44
2:N:77:THR:HG23	5:N:2014:HOH:O	2.16	0.44
4:Q:1084:CYC:CMD	4:Q:1084:CYC:NC	2.59	0.44
2:R:40:VAL:CG2	2:R:41:VAL:N	2.80	0.44
2:R:49:SER:OG	1:S:159:ASN:OD1	2.35	0.44
4:R:1153:CYC:CMB	1:S:148:VAL:HG11	2.48	0.44
3:X:72:ASN:OD1	4:X:1082:CYC:HMD3	2.17	0.44
1:A:64:PRO:HD2	5:A:2002:HOH:O	2.17	0.44
2:L:127:VAL:O	2:L:131:VAL:HG23	2.17	0.44
2:B:57:ARG:HD3	4:E:1084:CYC:O1D	2.17	0.44
3:X:57:ARG:NE	5:X:2006:HOH:O	2.34	0.44
1:I:47:LYS:O	1:I:48:ALA:C	2.61	0.44
2:J:82:CYS:SG	4:J:1082:CYC:C2C	3.05	0.44
4:N:1082:CYC:HMA1	4:N:1082:CYC:C4B	2.42	0.44
2:P:151:GLY:O	4:P:1153:CYC:OC	2.35	0.44
3:X:124:GLY:O	3:X:125:SER:C	2.59	0.44
2:D:148:ILE:HD13	4:D:1153:CYC:H3C	2.00	0.44
2:H:82:CYS:SG	4:H:1082:CYC:HBC2	2.55	0.44
2:L:69:PRO:HA	2:L:74:TYR:CG	2.53	0.44
2:L:148:ILE:HD13	4:L:1153:CYC:HHD	1.99	0.44
2:B:135:LYS:HG3	2:B:164:PHE:HB2	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:73:ASN:C	4:E:1084:CYC:CMD	2.91	0.44
2:L:148:ILE:HD13	4:L:1153:CYC:CHD	2.48	0.44
1:O:26:VAL:HG13	1:U:26:VAL:HG22	1.99	0.44
1:K:1:MET:HG2	1:K:105:GLY:HA3	2.00	0.43
4:L:1153:CYC:CBA	4:L:1153:CYC:CHA	2.96	0.43
1:G:11:VAL:HG12	1:G:15:GLN:NE2	2.32	0.43
2:J:153:CYS:SG	4:J:1153:CYC:C1C	3.06	0.43
2:B:153:CYS:SG	4:B:1153:CYC:CBC	3.02	0.43
1:C:65:TYR:HB2	1:C:69:MET:HE3	2.00	0.43
2:H:57:ARG:HH11	4:K:1084:CYC:CGD	2.30	0.43
2:D:124:GLY:O	2:D:125:SER:C	2.61	0.43
2:J:40:VAL:CG2	2:J:41:VAL:N	2.80	0.43
2:L:68:ALA:O	2:L:69:PRO:C	2.59	0.43
1:U:6:THR:HB	2:V:1:MET:HE2	2.00	0.43
2:D:72:ASN:OD1	4:D:1082:CYC:CMD	2.67	0.43
4:O:1084:CYC:HBB3	4:O:1084:CYC:CMB	2.49	0.43
2:P:40:VAL:CG2	2:P:97:VAL:HG11	2.48	0.43
2:R:40:VAL:HG11	2:R:156:LEU:HD21	1.99	0.43
4:T:1153:CYC:CBC	4:T:1153:CYC:CHD	2.96	0.43
2:D:82:CYS:SG	4:D:1082:CYC:H2C	2.58	0.43
4:G:1084:CYC:CMA	4:G:1084:CYC:NB	2.61	0.43
2:H:69:PRO:HA	2:H:74:TYR:CD2	2.54	0.43
4:S:1084:CYC:HHA	4:S:1084:CYC:HAD2	1.75	0.43
1:A:62:LYS:HG2	1:A:63:PHE:CE2	2.53	0.43
1:G:129:TYR:CE1	4:G:1084:CYC:HBC3	2.54	0.43
1:W:79:ARG:HE	1:W:79:ARG:HB3	1.73	0.43
2:B:135:LYS:HG3	2:B:164:PHE:CB	2.49	0.43
2:T:90:LEU:O	2:T:94:THR:HG23	2.19	0.43
2:N:57:ARG:HD3	4:Q:1084:CYC:O2D	2.18	0.43
4:A:1084:CYC:CMA	4:A:1084:CYC:NB	2.53	0.43
5:C:2009:HOH:O	2:J:47:ASN:HA	2.18	0.43
2:H:72:ASN:OD1	4:H:1082:CYC:CMD	2.67	0.42
1:Q:73:ASN:HA	4:Q:1084:CYC:HMD3	2.01	0.42
1:A:3:THR:HG21	2:B:6:THR:CG2	2.48	0.42
1:C:79:ARG:HE	1:C:79:ARG:HB3	1.63	0.42
1:M:148:VAL:HG11	4:X:1153:CYC:CMB	2.48	0.42
2:D:40:VAL:CG2	2:D:41:VAL:N	2.82	0.42
2:R:69:PRO:HA	2:R:74:TYR:CG	2.53	0.42
2:T:40:VAL:CG2	2:T:97:VAL:HG11	2.49	0.42
4:D:1082:CYC:HBC2	4:D:1082:CYC:H2C	1.73	0.42
2:R:40:VAL:CG2	2:R:97:VAL:HG11	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:T:30:MET:HE2	2:T:30:MET:HB3	1.91	0.42
1:W:73:ASN:C	4:W:1084:CYC:HMD3	2.44	0.42
3:X:153:CYS:SG	4:X:1153:CYC:CBC	3.08	0.42
1:G:40:ALA:HB2	1:G:146:ALA:HB1	2.00	0.42
1:O:128:TRP:CE3	4:O:1084:CYC:HMC2	2.54	0.42
2:B:102:ALA:O	2:B:103:SER:C	2.63	0.42
2:D:34:SER:HA	5:D:2006:HOH:O	2.19	0.42
2:H:39:ASP:OD1	4:H:1153:CYC:CMB	2.66	0.42
2:D:39:ASP:CG	2:D:148:ILE:HD11	2.44	0.42
2:D:40:VAL:CG2	2:D:97:VAL:HG11	2.50	0.42
2:F:40:VAL:HG21	2:F:160:ILE:HD11	2.02	0.42
1:G:67:THR:O	1:G:76:ALA:HA	2.19	0.42
2:H:90:LEU:O	2:H:94:THR:HG23	2.19	0.42
2:L:124:GLY:O	2:L:125:SER:C	2.62	0.42
1:Q:15:GLN:OE1	1:S:2:LYS:NZ	2.52	0.42
1:S:62:LYS:O	1:S:62:LYS:HG3	2.19	0.42
2:T:69:PRO:HA	2:T:74:TYR:CG	2.55	0.42
2:L:72:ASN:OD1	4:L:1082:CYC:CMD	2.68	0.42
1:M:2:LYS:HB2	5:M:2001:HOH:O	2.20	0.42
1:M:26:VAL:HG22	1:W:26:VAL:HG13	2.02	0.42
2:N:153:CYS:SG	4:N:1153:CYC:H2C	2.60	0.42
4:W:1084:CYC:HMC1	4:W:1084:CYC:HBC2	2.01	0.42
2:N:148:ILE:HD13	4:N:1153:CYC:H3C	2.02	0.42
2:V:151:GLY:O	4:V:1153:CYC:OC	2.37	0.42
2:F:40:VAL:HG11	2:F:156:LEU:CD2	2.50	0.41
1:M:79:ARG:HE	1:M:79:ARG:HB3	1.68	0.41
1:A:120:ARG:HG3	1:A:120:ARG:HH11	1.84	0.41
1:C:62:LYS:HG2	1:C:63:PHE:CE2	2.55	0.41
1:G:104:THR:O	1:G:105:GLY:C	2.63	0.41
2:L:36:LYS:O	2:L:40:VAL:CG1	2.67	0.41
2:L:123:PRO:O	2:L:126:SER:HB2	2.20	0.41
4:N:1153:CYC:HB	4:N:1153:CYC:CMA	2.33	0.41
2:R:72:ASN:OD1	4:R:1082:CYC:HMD3	2.20	0.41
1:C:3:THR:HG21	2:D:6:THR:CG2	2.49	0.41
1:E:111:LEU:C	1:E:111:LEU:HD23	2.45	0.41
2:P:30:MET:HE2	2:P:30:MET:HB3	1.94	0.41
3:X:103:SER:O	3:X:104:VAL:C	2.62	0.41
3:X:127:VAL:HG22	4:X:1082:CYC:H3C	2.03	0.41
2:L:82:CYS:SG	4:L:1082:CYC:HBC2	2.55	0.41
2:P:118:LEU:HD13	2:P:118:LEU:HA	1.80	0.41
1:C:125:SER:HB3	1:C:128:TRP:CE2	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:151:GLY:HA3	4:H:1153:CYC:CMD	2.51	0.41
1:K:42:LYS:NZ	2:L:25:ASP:OD1	2.54	0.41
1:U:47:LYS:O	1:U:48:ALA:C	2.62	0.41
2:J:73:ALA:HB2	4:J:1082:CYC:OC	2.20	0.41
1:M:47:LYS:O	1:M:48:ALA:C	2.64	0.41
4:N:1153:CYC:HMB3	1:W:148:VAL:HG11	2.01	0.41
1:U:55:ALA:O	1:U:56:ALA:C	2.63	0.41
1:O:73:ASN:O	4:O:1084:CYC:CMD	2.68	0.41
2:V:40:VAL:CG2	2:V:97:VAL:HG11	2.50	0.41
2:V:82:CYS:SG	4:V:1082:CYC:H2C	2.60	0.41
2:D:40:VAL:HG13	2:D:40:VAL:H	1.59	0.41
2:F:35:ASN:HD22	2:F:35:ASN:HA	1.70	0.41
1:I:73:ASN:HA	4:I:1084:CYC:HBD2	2.03	0.41
1:K:73:ASN:HA	4:K:1084:CYC:HAD1	2.02	0.41
2:L:40:VAL:CG2	2:L:97:VAL:HG11	2.51	0.41
2:L:82:CYS:SG	4:L:1082:CYC:H2C	2.60	0.41
2:L:135:LYS:HG3	2:L:164:PHE:HB2	2.03	0.41
2:R:36:LYS:O	2:R:40:VAL:CG1	2.67	0.41
2:V:153:CYS:SG	4:V:1153:CYC:HBC2	2.46	0.41
2:D:30:MET:HE2	2:D:30:MET:HB3	1.93	0.41
2:F:148:ILE:CD1	4:F:1153:CYC:HHD	2.51	0.41
2:H:3:ASP:O	2:H:4:ALA:C	2.63	0.41
4:J:1153:CYC:CBC	4:J:1153:CYC:HHD	2.48	0.41
1:O:78:GLN:O	1:O:79:ARG:C	2.64	0.41
2:P:40:VAL:CG2	2:P:41:VAL:N	2.84	0.41
2:R:127:VAL:HG22	4:R:1082:CYC:H3C	2.03	0.41
3:X:40:VAL:HG13	3:X:40:VAL:H	1.60	0.41
3:X:151:GLY:HA3	4:X:1153:CYC:CMD	2.50	0.41
1:M:72:PRO:HB2	4:M:1084:CYC:O1D	2.22	0.41
4:R:1153:CYC:HMA3	4:R:1153:CYC:NB	2.20	0.41
1:A:15:GLN:OE1	1:K:2:LYS:NZ	2.53	0.40
4:C:1084:CYC:CMA	4:C:1084:CYC:HBA1	2.51	0.40
2:F:39:ASP:CG	2:F:148:ILE:HD11	2.46	0.40
2:F:72:ASN:OD1	4:F:1082:CYC:HMD1	2.21	0.40
1:O:73:ASN:O	4:O:1084:CYC:HMD3	2.20	0.40
4:V:1153:CYC:CMA	4:V:1153:CYC:NB	2.66	0.40
1:C:115:ILE:HD12	1:C:115:ILE:HA	1.87	0.40
1:E:73:ASN:HA	4:E:1084:CYC:HMD3	2.04	0.40
2:J:10:SER:O	2:J:14:THR:HG23	2.21	0.40
4:O:1084:CYC:HB	4:O:1084:CYC:HMA2	1.66	0.40
2:T:72:ASN:OD1	2:T:122:THR:HA	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:73:ASN:O	4:A:1084:CYC:CMD	2.69	0.40
2:D:95:TYR:HB3	2:D:104:VAL:HG11	2.03	0.40
1:E:26:VAL:HG22	1:G:26:VAL:HG13	2.04	0.40
1:G:6:THR:HG23	2:H:3:ASP:OD2	2.21	0.40
2:H:36:LYS:O	2:H:40:VAL:CG1	2.66	0.40
2:J:82:CYS:SG	4:J:1082:CYC:H2C	2.62	0.40
4:O:1084:CYC:HBD1	4:O:1084:CYC:CHA	2.42	0.40
2:V:36:LYS:O	2:V:40:VAL:CG1	2.68	0.40
1:A:40:ALA:HB2	1:A:146:ALA:HB1	2.02	0.40
1:E:115:ILE:HD12	1:E:115:ILE:HA	1.84	0.40
1:I:62:LYS:HG2	1:I:63:PHE:CE2	2.56	0.40
1:M:43:ALA:HB3	1:M:142:LEU:HD21	2.03	0.40
1:S:3:THR:OG1	1:S:6:THR:HG23	2.22	0.40
1:E:87:ASP:OD2	4:E:1084:CYC:NA	2.55	0.40
2:L:72:ASN:OD1	4:L:1082:CYC:HMD3	2.22	0.40
2:L:82:CYS:SG	4:L:1082:CYC:C2C	3.10	0.40
2:P:108:ARG:HB3	5:P:2007:HOH:O	2.20	0.40
2:V:118:LEU:HD13	2:V:118:LEU:HA	1.99	0.40

All (5) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:125:SER:OG	2:V:62:GLU:OE1[2_656]	2.03	0.17
1:I:68:GLN:NE2	1:W:68:GLN:NE2[1_655]	2.06	0.14
1:K:67:THR:O	1:S:61:ASN:CG[1_656]	2.08	0.12
1:A:60:TYR:O	1:Q:68:GLN:OE1[1_655]	2.11	0.09
1:K:61:ASN:CG	1:O:68:GLN:NE2[1_656]	2.17	0.03

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	161/162 (99%)	158 (98%)	3 (2%)	0	100	100
1	C	160/162 (99%)	153 (96%)	7 (4%)	0	100	100
1	E	160/162 (99%)	153 (96%)	7 (4%)	0	100	100
1	G	160/162 (99%)	155 (97%)	5 (3%)	0	100	100
1	I	160/162 (99%)	153 (96%)	7 (4%)	0	100	100
1	K	160/162 (99%)	155 (97%)	5 (3%)	0	100	100
1	M	160/162 (99%)	152 (95%)	8 (5%)	0	100	100
1	O	160/162 (99%)	155 (97%)	5 (3%)	0	100	100
1	Q	160/162 (99%)	157 (98%)	3 (2%)	0	100	100
1	S	160/162 (99%)	156 (98%)	4 (2%)	0	100	100
1	U	160/162 (99%)	156 (98%)	4 (2%)	0	100	100
1	W	160/162 (99%)	155 (97%)	5 (3%)	0	100	100
2	B	170/172 (99%)	166 (98%)	4 (2%)	0	100	100
2	D	170/172 (99%)	166 (98%)	4 (2%)	0	100	100
2	F	170/172 (99%)	166 (98%)	4 (2%)	0	100	100
2	H	170/172 (99%)	166 (98%)	4 (2%)	0	100	100
2	J	170/172 (99%)	166 (98%)	4 (2%)	0	100	100
2	L	170/172 (99%)	166 (98%)	4 (2%)	0	100	100
2	N	170/172 (99%)	165 (97%)	5 (3%)	0	100	100
2	P	170/172 (99%)	167 (98%)	3 (2%)	0	100	100
2	R	170/172 (99%)	168 (99%)	2 (1%)	0	100	100
2	T	170/172 (99%)	165 (97%)	5 (3%)	0	100	100
2	V	170/172 (99%)	166 (98%)	4 (2%)	0	100	100
3	X	170/172 (99%)	166 (98%)	4 (2%)	0	100	100
All	All	3961/4008 (99%)	3851 (97%)	110 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	125/125 (100%)	115 (92%)	10 (8%)	11	38
1	C	124/125 (99%)	116 (94%)	8 (6%)	15	47
1	E	124/125 (99%)	114 (92%)	10 (8%)	11	38
1	G	124/125 (99%)	111 (90%)	13 (10%)	6	27
1	I	124/125 (99%)	112 (90%)	12 (10%)	8	30
1	K	124/125 (99%)	112 (90%)	12 (10%)	8	30
1	M	124/125 (99%)	115 (93%)	9 (7%)	13	42
1	O	124/125 (99%)	113 (91%)	11 (9%)	9	34
1	Q	124/125 (99%)	112 (90%)	12 (10%)	8	30
1	S	124/125 (99%)	115 (93%)	9 (7%)	13	42
1	U	124/125 (99%)	113 (91%)	11 (9%)	9	34
1	W	124/125 (99%)	114 (92%)	10 (8%)	11	38
2	B	131/133 (98%)	123 (94%)	8 (6%)	17	49
2	D	131/133 (98%)	124 (95%)	7 (5%)	20	54
2	F	131/133 (98%)	125 (95%)	6 (5%)	24	58
2	H	131/133 (98%)	123 (94%)	8 (6%)	17	49
2	J	131/133 (98%)	123 (94%)	8 (6%)	17	49
2	L	131/133 (98%)	127 (97%)	4 (3%)	35	68
2	N	131/133 (98%)	122 (93%)	9 (7%)	14	45
2	P	131/133 (98%)	123 (94%)	8 (6%)	17	49
2	R	131/133 (98%)	124 (95%)	7 (5%)	20	54
2	T	131/133 (98%)	124 (95%)	7 (5%)	20	54
2	V	131/133 (98%)	123 (94%)	8 (6%)	17	49
3	X	132/134 (98%)	127 (96%)	5 (4%)	29	63
All	All	3062/3097 (99%)	2850 (93%)	212 (7%)	14	45

All (212) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	LYS
1	A	3	THR
1	A	32	ARG
1	A	38	LEU

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Mol	Chain	Res	Type
1	A	39	GLU
1	A	42	LYS
1	A	46	SER
1	A	53	SER
1	A	68	GLN
1	A	120	ARG
2	B	10	SER
2	B	18	MET
2	B	29	GLN
2	B	30	MET
2	B	40	VAL
2	B	49	SER
2	B	118	LEU
2	B	141	ILE
1	C	3	THR
1	C	32	ARG
1	C	38	LEU
1	C	46	SER
1	C	53	SER
1	C	68	GLN
1	C	100	ILE
1	C	120	ARG
2	D	10	SER
2	D	18	MET
2	D	29	GLN
2	D	30	MET
2	D	40	VAL
2	D	118	LEU
2	D	141	ILE
1	E	2	LYS
1	E	3	THR
1	E	32	ARG
1	E	38	LEU
1	E	39	GLU
1	E	46	SER
1	E	53	SER
1	E	68	GLN
1	E	100	ILE
1	E	120	ARG
2	F	10	SER
2	F	18	MET
2	F	30	MET

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Mol	Chain	Res	Type
2	F	40	VAL
2	F	118	LEU
2	F	141	ILE
1	G	2	LYS
1	G	3	THR
1	G	4	PRO
1	G	32	ARG
1	G	38	LEU
1	G	39	GLU
1	G	42	LYS
1	G	46	SER
1	G	53	SER
1	G	68	GLN
1	G	100	ILE
1	G	117	GLU
1	G	120	ARG
2	H	10	SER
2	H	18	MET
2	H	28	SER
2	H	29	GLN
2	H	30	MET
2	H	40	VAL
2	H	118	LEU
2	H	141	ILE
1	I	2	LYS
1	I	3	THR
1	I	6	THR
1	I	32	ARG
1	I	38	LEU
1	I	39	GLU
1	I	42	LYS
1	I	46	SER
1	I	53	SER
1	I	68	GLN
1	I	100	ILE
1	I	120	ARG
2	J	10	SER
2	J	18	MET
2	J	29	GLN
2	J	30	MET
2	J	40	VAL
2	J	72	ASN

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Mol	Chain	Res	Type
2	J	118	LEU
2	J	141	ILE
1	K	2	LYS
1	K	3	THR
1	K	24	ILE
1	K	32	ARG
1	K	38	LEU
1	K	39	GLU
1	K	42	LYS
1	K	46	SER
1	K	53	SER
1	K	68	GLN
1	K	100	ILE
1	K	120	ARG
2	L	10	SER
2	L	29	GLN
2	L	40	VAL
2	L	118	LEU
1	M	2	LYS
1	M	3	THR
1	M	32	ARG
1	M	38	LEU
1	M	53	SER
1	M	68	GLN
1	M	100	ILE
1	M	120	ARG
1	M	121	THR
2	N	10	SER
2	N	18	MET
2	N	28	SER
2	N	29	GLN
2	N	30	MET
2	N	40	VAL
2	N	118	LEU
2	N	141	ILE
2	N	166	ARG
1	O	2	LYS
1	O	3	THR
1	O	32	ARG
1	O	38	LEU
1	O	39	GLU
1	O	42	LYS

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Mol	Chain	Res	Type
1	O	46	SER
1	O	53	SER
1	O	68	GLN
1	O	100	ILE
1	O	120	ARG
2	P	10	SER
2	P	18	MET
2	P	29	GLN
2	P	30	MET
2	P	40	VAL
2	P	72	ASN
2	P	118	LEU
2	P	141	ILE
1	Q	2	LYS
1	Q	3	THR
1	Q	32	ARG
1	Q	38	LEU
1	Q	39	GLU
1	Q	42	LYS
1	Q	46	SER
1	Q	53	SER
1	Q	68	GLN
1	Q	100	ILE
1	Q	115	ILE
1	Q	120	ARG
2	R	10	SER
2	R	18	MET
2	R	29	GLN
2	R	40	VAL
2	R	49	SER
2	R	118	LEU
2	R	141	ILE
1	S	2	LYS
1	S	32	ARG
1	S	38	LEU
1	S	39	GLU
1	S	42	LYS
1	S	53	SER
1	S	68	GLN
1	S	100	ILE
1	S	120	ARG
2	T	10	SER

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Mol	Chain	Res	Type
2	T	18	MET
2	T	29	GLN
2	T	30	MET
2	T	40	VAL
2	T	118	LEU
2	T	141	ILE
1	U	2	LYS
1	U	3	THR
1	U	32	ARG
1	U	38	LEU
1	U	39	GLU
1	U	42	LYS
1	U	46	SER
1	U	53	SER
1	U	68	GLN
1	U	100	ILE
1	U	120	ARG
2	V	10	SER
2	V	18	MET
2	V	19	LEU
2	V	28	SER
2	V	29	GLN
2	V	30	MET
2	V	40	VAL
2	V	118	LEU
1	W	2	LYS
1	W	3	THR
1	W	32	ARG
1	W	38	LEU
1	W	39	GLU
1	W	42	LYS
1	W	46	SER
1	W	53	SER
1	W	68	GLN
1	W	120	ARG
3	X	10	SER
3	X	29	GLN
3	X	40	VAL
3	X	118	LEU
3	X	141	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (47) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	119	ASN
2	B	23	GLN
2	B	35	ASN
2	B	47	ASN
2	D	23	GLN
2	D	35	ASN
2	D	47	ASN
2	F	23	GLN
2	F	42	ASN
2	F	47	ASN
1	G	70	GLN
2	H	23	GLN
2	H	35	ASN
2	H	42	ASN
2	H	47	ASN
2	J	23	GLN
2	J	47	ASN
1	K	70	GLN
2	L	23	GLN
2	L	35	ASN
2	L	42	ASN
2	L	47	ASN
1	M	33	GLN
1	M	119	ASN
2	N	23	GLN
2	N	42	ASN
2	N	47	ASN
1	O	68	GLN
2	P	23	GLN
2	P	35	ASN
2	P	47	ASN
1	Q	70	GLN
2	R	23	GLN
2	R	35	ASN
2	R	42	ASN
2	R	47	ASN
2	R	143	ASN
2	T	23	GLN
2	T	47	ASN
1	U	70	GLN
2	V	23	GLN
2	V	42	ASN
2	V	47	ASN

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Mol	Chain	Res	Type
1	W	70	GLN
3	X	23	GLN
3	X	35	ASN
3	X	47	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

36 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	CYC	Q	1084	1	46,46,46	2.41	13 (28%)	63,67,67	2.72	29 (46%)
4	CYC	D	1153	-	46,46,46	2.54	19 (41%)	63,67,67	3.76	29 (46%)
4	CYC	V	1153	-	46,46,46	2.41	17 (36%)	63,67,67	3.75	29 (46%)
4	CYC	K	1084	1	46,46,46	2.47	16 (34%)	63,67,67	2.97	26 (41%)
4	CYC	O	1084	1	46,46,46	2.37	19 (41%)	63,67,67	3.03	27 (42%)
4	CYC	L	1153	-	46,46,46	2.63	20 (43%)	63,67,67	3.84	31 (49%)
4	CYC	A	1084	1	46,46,46	2.13	15 (32%)	63,67,67	3.20	29 (46%)
4	CYC	H	1082	-	46,46,46	2.23	18 (39%)	63,67,67	3.16	32 (50%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	CYC	D	1082	2	46,46,46	2.67	17 (36%)	63,67,67	3.55	33 (52%)
4	CYC	J	1082	2	46,46,46	2.48	15 (32%)	63,67,67	3.11	27 (42%)
4	CYC	B	1153	-	46,46,46	2.66	18 (39%)	63,67,67	4.16	33 (52%)
4	CYC	N	1082	2	46,46,46	2.60	15 (32%)	63,67,67	3.72	38 (60%)
4	CYC	P	1082	-	46,46,46	2.40	15 (32%)	63,67,67	3.47	34 (53%)
4	CYC	U	1084	1	46,46,46	2.37	15 (32%)	63,67,67	2.76	27 (42%)
4	CYC	E	1084	1	46,46,46	2.37	17 (36%)	63,67,67	3.54	32 (50%)
4	CYC	V	1082	2	46,46,46	2.31	16 (34%)	63,67,67	3.40	31 (49%)
4	CYC	F	1082	2	46,46,46	2.39	16 (34%)	63,67,67	3.59	39 (61%)
4	CYC	R	1153	-	46,46,46	2.61	18 (39%)	63,67,67	3.90	35 (55%)
4	CYC	F	1153	2	46,46,46	2.46	16 (34%)	63,67,67	3.61	31 (49%)
4	CYC	X	1153	-	46,46,46	2.53	16 (34%)	63,67,67	2.86	25 (39%)
4	CYC	W	1084	1	46,46,46	2.27	12 (26%)	63,67,67	3.18	33 (52%)
4	CYC	L	1082	2	46,46,46	2.51	15 (32%)	63,67,67	3.28	27 (42%)
4	CYC	H	1153	-	46,46,46	2.70	19 (41%)	63,67,67	3.96	37 (58%)
4	CYC	X	1082	-	46,46,46	2.51	16 (34%)	63,67,67	3.13	26 (41%)
4	CYC	T	1153	-	46,46,46	2.46	19 (41%)	63,67,67	3.98	34 (53%)
4	CYC	N	1153	-	46,46,46	2.76	21 (45%)	63,67,67	3.75	35 (55%)
4	CYC	S	1084	1	46,46,46	2.42	15 (32%)	63,67,67	3.69	33 (52%)
4	CYC	J	1153	-	46,46,46	2.57	17 (36%)	63,67,67	3.88	33 (52%)
4	CYC	T	1082	2	46,46,46	2.25	15 (32%)	63,67,67	3.70	37 (58%)
4	CYC	C	1084	1	46,46,46	2.11	15 (32%)	63,67,67	3.86	38 (60%)
4	CYC	P	1153	-	46,46,46	2.43	17 (36%)	63,67,67	3.53	27 (42%)
4	CYC	I	1084	1	46,46,46	2.58	17 (36%)	63,67,67	3.81	36 (57%)
4	CYC	M	1084	1	46,46,46	2.42	18 (39%)	63,67,67	3.03	28 (44%)
4	CYC	G	1084	1	46,46,46	2.04	11 (23%)	63,67,67	3.02	36 (57%)
4	CYC	B	1082	-	46,46,46	2.66	18 (39%)	63,67,67	3.55	33 (52%)
4	CYC	R	1082	2	46,46,46	2.77	16 (34%)	63,67,67	3.53	30 (47%)

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
4	CYC	Q	1084	1	-	12/26/74/74	0/4/4/4

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	CYC	D	1153	-	-	11/26/74/74	0/4/4/4
4	CYC	V	1153	-	-	9/26/74/74	0/4/4/4
4	CYC	K	1084	1	-	12/26/74/74	0/4/4/4
4	CYC	O	1084	1	-	16/26/74/74	0/4/4/4
4	CYC	L	1153	-	-	14/26/74/74	0/4/4/4
4	CYC	A	1084	1	-	15/26/74/74	0/4/4/4
4	CYC	H	1082	-	-	10/26/74/74	0/4/4/4
4	CYC	D	1082	2	-	9/26/74/74	0/4/4/4
4	CYC	J	1082	2	-	11/26/74/74	0/4/4/4
4	CYC	B	1153	-	-	8/26/74/74	0/4/4/4
4	CYC	N	1082	2	-	10/26/74/74	0/4/4/4
4	CYC	P	1082	-	-	9/26/74/74	0/4/4/4
4	CYC	U	1084	1	-	12/26/74/74	0/4/4/4
4	CYC	E	1084	1	-	13/26/74/74	0/4/4/4
4	CYC	V	1082	2	-	13/26/74/74	0/4/4/4
4	CYC	F	1082	2	-	11/26/74/74	0/4/4/4
4	CYC	R	1153	-	-	11/26/74/74	0/4/4/4
4	CYC	F	1153	2	-	10/26/74/74	0/4/4/4
4	CYC	X	1153	-	-	12/26/74/74	0/4/4/4
4	CYC	W	1084	1	-	9/26/74/74	0/4/4/4
4	CYC	L	1082	2	-	10/26/74/74	0/4/4/4
4	CYC	H	1153	-	-	9/26/74/74	0/4/4/4
4	CYC	X	1082	-	-	7/26/74/74	0/4/4/4
4	CYC	T	1153	-	-	12/26/74/74	0/4/4/4
4	CYC	N	1153	-	-	10/26/74/74	0/4/4/4
4	CYC	S	1084	1	-	9/26/74/74	0/4/4/4
4	CYC	J	1153	-	-	13/26/74/74	0/4/4/4
4	CYC	T	1082	2	-	11/26/74/74	0/4/4/4
4	CYC	C	1084	1	-	15/26/74/74	0/4/4/4
4	CYC	P	1153	-	-	11/26/74/74	0/4/4/4
4	CYC	I	1084	1	-	9/26/74/74	0/4/4/4
4	CYC	M	1084	1	-	12/26/74/74	0/4/4/4
4	CYC	G	1084	1	-	14/26/74/74	0/4/4/4
4	CYC	B	1082	-	-	14/26/74/74	0/4/4/4

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	CYC	R	1082	2	-	11/26/74/74	0/4/4/4

All (592) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	N	1153	CYC	C4B-C3B	-7.37	1.34	1.48
4	H	1153	CYC	C2C-C1C	-7.29	1.45	1.52
4	J	1153	CYC	C2C-C1C	-7.19	1.45	1.52
4	N	1082	CYC	C2C-C1C	-7.17	1.45	1.52
4	I	1084	CYC	C4B-C3B	-7.05	1.35	1.48
4	J	1153	CYC	C4B-C3B	-7.04	1.35	1.48
4	M	1084	CYC	C4B-C3B	-6.88	1.35	1.48
4	L	1153	CYC	C4B-C3B	-6.87	1.35	1.48
4	R	1082	CYC	CHA-C1A	6.84	1.51	1.38
4	R	1153	CYC	C1A-C2A	-6.68	1.35	1.45
4	R	1082	CYC	C2C-C1C	-6.67	1.46	1.52
4	B	1153	CYC	C4C-NC	-6.63	1.24	1.37
4	F	1153	CYC	C1A-C2A	-6.52	1.35	1.45
4	X	1153	CYC	CHB-C1B	6.49	1.53	1.37
4	V	1153	CYC	C4B-C3B	-6.42	1.36	1.48
4	D	1082	CYC	C2C-C1C	-6.41	1.46	1.52
4	I	1084	CYC	C1C-NC	-6.37	1.29	1.37
4	R	1082	CYC	C4B-C3B	-6.37	1.36	1.48
4	M	1084	CYC	CHA-C1A	6.36	1.50	1.38
4	H	1153	CYC	C1A-C2A	-6.34	1.35	1.45
4	B	1082	CYC	C1C-NC	-6.33	1.29	1.37
4	D	1082	CYC	C1C-NC	-6.32	1.29	1.37
4	Q	1084	CYC	C4B-C3B	-6.27	1.36	1.48
4	E	1084	CYC	C1A-C2A	-6.20	1.36	1.45
4	J	1082	CYC	CHB-C1B	6.20	1.52	1.37
4	B	1153	CYC	C2C-C1C	-6.19	1.46	1.52
4	S	1084	CYC	C4B-C3B	-6.16	1.36	1.48
4	R	1153	CYC	C4B-C3B	-6.15	1.36	1.48
4	P	1153	CYC	C2C-C1C	-6.15	1.46	1.52
4	D	1082	CYC	C4B-C3B	-6.05	1.37	1.48
4	D	1153	CYC	C4B-C3B	-6.05	1.37	1.48
4	B	1153	CYC	C4B-C3B	-6.05	1.37	1.48
4	F	1153	CYC	C2C-C1C	-6.02	1.46	1.52
4	P	1082	CYC	C4B-C3B	-6.01	1.37	1.48
4	N	1082	CYC	C4B-C3B	-5.93	1.37	1.48
4	D	1082	CYC	CHA-C1A	5.87	1.49	1.38
4	D	1153	CYC	CHA-C1A	5.82	1.49	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	U	1084	CYC	CHD-C1D	5.81	1.53	1.40
4	X	1153	CYC	CHA-C1A	5.79	1.49	1.38
4	S	1084	CYC	C1A-C2A	-5.76	1.36	1.45
4	L	1153	CYC	C4C-NC	-5.75	1.26	1.37
4	K	1084	CYC	C1A-C2A	-5.73	1.36	1.45
4	N	1153	CYC	C4C-NC	-5.72	1.26	1.37
4	L	1082	CYC	CHB-C1B	5.71	1.51	1.37
4	B	1082	CYC	C4B-C3B	-5.65	1.37	1.48
4	T	1153	CYC	C4B-C3B	-5.61	1.37	1.48
4	J	1082	CYC	C4B-C3B	-5.57	1.37	1.48
4	L	1082	CYC	CHA-C1A	5.57	1.49	1.38
4	X	1082	CYC	C4B-C3B	-5.56	1.38	1.48
4	H	1082	CYC	C4B-C3B	-5.54	1.38	1.48
4	P	1153	CYC	C1A-C2A	-5.53	1.37	1.45
4	W	1084	CYC	CHD-C1D	5.52	1.52	1.40
4	R	1082	CYC	C1C-NC	-5.52	1.30	1.37
4	R	1082	CYC	CHA-C4D	5.50	1.52	1.40
4	F	1153	CYC	C4B-C3B	-5.49	1.38	1.48
4	U	1084	CYC	CHB-C1B	5.48	1.51	1.37
4	X	1153	CYC	C2C-C1C	-5.47	1.47	1.52
4	J	1082	CYC	CHA-C1A	5.45	1.49	1.38
4	H	1153	CYC	C4B-C3B	-5.45	1.38	1.48
4	N	1082	CYC	CHA-C1A	5.42	1.48	1.38
4	L	1082	CYC	C4B-C3B	-5.42	1.38	1.48
4	V	1082	CYC	CHB-C1B	5.42	1.50	1.37
4	H	1153	CYC	C4C-NC	-5.39	1.26	1.37
4	F	1082	CYC	C2C-C1C	-5.38	1.47	1.52
4	M	1084	CYC	CHB-C1B	5.35	1.50	1.37
4	N	1153	CYC	C2C-C1C	-5.33	1.47	1.52
4	L	1153	CYC	C1C-NC	-5.33	1.30	1.37
4	N	1153	CYC	C1A-C2A	-5.31	1.37	1.45
4	J	1082	CYC	CHB-C4A	5.29	1.52	1.40
4	L	1082	CYC	C1C-NC	-5.28	1.30	1.37
4	H	1153	CYC	CHB-C1B	5.27	1.50	1.37
4	J	1153	CYC	C1A-C2A	-5.26	1.37	1.45
4	X	1082	CYC	C2C-C1C	-5.24	1.47	1.52
4	D	1153	CYC	C4C-NC	-5.22	1.27	1.37
4	K	1084	CYC	C4B-C3B	-5.22	1.38	1.48
4	A	1084	CYC	CHD-C1D	5.22	1.52	1.40
4	P	1082	CYC	CHB-C1B	5.22	1.50	1.37
4	N	1153	CYC	CHB-C1B	5.21	1.50	1.37
4	O	1084	CYC	C4B-C3B	-5.20	1.38	1.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	X	1082	CYC	CHB-C1B	5.17	1.50	1.37
4	H	1082	CYC	CHA-C1A	5.17	1.48	1.38
4	V	1153	CYC	C1A-C2A	-5.16	1.37	1.45
4	P	1082	CYC	C2C-C1C	-5.13	1.47	1.52
4	T	1082	CYC	C4B-C3B	-5.12	1.38	1.48
4	B	1082	CYC	C1A-C2A	-5.12	1.37	1.45
4	P	1153	CYC	CHB-C1B	5.11	1.50	1.37
4	B	1153	CYC	CHB-C1B	5.09	1.50	1.37
4	A	1084	CYC	C4B-C3B	-5.08	1.38	1.48
4	K	1084	CYC	CHA-C1A	5.07	1.48	1.38
4	E	1084	CYC	CHB-C1B	5.06	1.50	1.37
4	T	1082	CYC	CHD-C1D	5.06	1.51	1.40
4	Q	1084	CYC	CHA-C1A	5.05	1.48	1.38
4	K	1084	CYC	CHB-C1B	5.05	1.50	1.37
4	T	1082	CYC	C1C-NC	-5.04	1.31	1.37
4	A	1084	CYC	CHA-C1A	5.04	1.48	1.38
4	M	1084	CYC	C1A-C2A	-5.04	1.37	1.45
4	X	1082	CYC	CHD-C1D	5.03	1.51	1.40
4	J	1153	CYC	CHB-C1B	5.03	1.49	1.37
4	B	1082	CYC	CHD-C1D	5.02	1.51	1.40
4	D	1153	CYC	C2C-C1C	-5.02	1.47	1.52
4	B	1153	CYC	C1A-C2A	-5.01	1.37	1.45
4	W	1084	CYC	CHB-C1B	4.98	1.49	1.37
4	O	1084	CYC	CHB-C1B	4.97	1.49	1.37
4	Q	1084	CYC	C1A-C2A	-4.96	1.38	1.45
4	G	1084	CYC	CHB-C1B	4.96	1.49	1.37
4	P	1153	CYC	C4B-C3B	-4.92	1.39	1.48
4	B	1082	CYC	CHB-C1B	4.92	1.49	1.37
4	K	1084	CYC	CHA-C4D	4.91	1.51	1.40
4	N	1082	CYC	CHD-C1D	4.88	1.51	1.40
4	R	1082	CYC	CHB-C1B	4.88	1.49	1.37
4	T	1082	CYC	C1A-C2A	-4.85	1.38	1.45
4	C	1084	CYC	C4B-C3B	-4.85	1.39	1.48
4	F	1082	CYC	C4B-C3B	-4.83	1.39	1.48
4	X	1153	CYC	CHB-C4A	4.83	1.51	1.40
4	L	1082	CYC	CHB-C4A	4.82	1.51	1.40
4	N	1082	CYC	CHB-C4A	4.80	1.51	1.40
4	F	1153	CYC	CHB-C1B	4.78	1.49	1.37
4	V	1082	CYC	C2C-C1C	-4.77	1.47	1.52
4	G	1084	CYC	CHD-C1D	4.76	1.51	1.40
4	B	1082	CYC	C2C-C1C	-4.75	1.47	1.52
4	V	1082	CYC	C4B-C3B	-4.74	1.39	1.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	S	1084	CYC	CHB-C1B	4.72	1.49	1.37
4	B	1082	CYC	C4C-NC	-4.72	1.28	1.37
4	I	1084	CYC	CHB-C1B	4.72	1.49	1.37
4	S	1084	CYC	CHD-C1D	4.72	1.51	1.40
4	W	1084	CYC	CHA-C1A	4.71	1.47	1.38
4	R	1153	CYC	C1C-NC	-4.71	1.31	1.37
4	K	1084	CYC	CHD-C1D	4.70	1.51	1.40
4	R	1153	CYC	C2C-C1C	-4.69	1.47	1.52
4	G	1084	CYC	CHA-C1A	4.68	1.47	1.38
4	U	1084	CYC	C4B-C3B	-4.68	1.39	1.48
4	F	1082	CYC	C1C-NC	-4.67	1.31	1.37
4	E	1084	CYC	CHA-C1A	4.65	1.47	1.38
4	T	1153	CYC	C1C-NC	-4.64	1.31	1.37
4	R	1082	CYC	CHD-C1D	4.64	1.50	1.40
4	H	1153	CYC	C1C-NC	-4.64	1.31	1.37
4	R	1082	CYC	CHB-C4A	4.63	1.51	1.40
4	N	1082	CYC	CHB-C1B	4.62	1.48	1.37
4	X	1082	CYC	C1A-C2A	-4.62	1.38	1.45
4	L	1153	CYC	CHB-C1B	4.60	1.48	1.37
4	I	1084	CYC	C2C-C1C	-4.59	1.48	1.52
4	O	1084	CYC	CHD-C1D	4.58	1.50	1.40
4	H	1082	CYC	CHD-C1D	4.58	1.50	1.40
4	E	1084	CYC	C4B-C3B	-4.58	1.39	1.48
4	V	1082	CYC	CHA-C1A	4.57	1.47	1.38
4	X	1082	CYC	C4C-NC	-4.56	1.28	1.37
4	X	1082	CYC	CHA-C1A	4.53	1.47	1.38
4	L	1153	CYC	CHA-C1A	4.53	1.47	1.38
4	T	1153	CYC	CHA-C1A	4.52	1.47	1.38
4	L	1153	CYC	C2C-C1C	-4.51	1.48	1.52
4	U	1084	CYC	CHA-C1A	4.51	1.47	1.38
4	D	1153	CYC	C1C-NC	-4.51	1.31	1.37
4	W	1084	CYC	C4B-C3B	-4.51	1.39	1.48
4	X	1153	CYC	C1A-C2A	-4.51	1.38	1.45
4	I	1084	CYC	CHB-C4A	4.50	1.51	1.40
4	F	1082	CYC	C4C-NC	-4.50	1.28	1.37
4	Q	1084	CYC	CHB-C1B	4.50	1.48	1.37
4	I	1084	CYC	CHD-C1D	4.50	1.50	1.40
4	P	1082	CYC	CHB-C4A	4.49	1.51	1.40
4	P	1153	CYC	CHA-C1A	4.48	1.47	1.38
4	J	1153	CYC	CHA-C1A	4.47	1.47	1.38
4	U	1084	CYC	C1C-NC	-4.46	1.31	1.37
4	T	1153	CYC	CHD-C1D	4.46	1.50	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	T	1082	CYC	C1D-C2D	4.44	1.52	1.41
4	S	1084	CYC	CHB-C4A	4.44	1.50	1.40
4	D	1082	CYC	CHB-C1B	4.44	1.48	1.37
4	E	1084	CYC	CHD-C1D	4.42	1.50	1.40
4	F	1082	CYC	CHA-C1A	4.41	1.47	1.38
4	O	1084	CYC	C1C-NC	-4.40	1.31	1.37
4	P	1082	CYC	CHD-C1D	4.39	1.50	1.40
4	X	1082	CYC	CHA-C4D	4.39	1.50	1.40
4	W	1084	CYC	CHB-C4A	4.37	1.50	1.40
4	V	1082	CYC	CHB-C4A	4.37	1.50	1.40
4	I	1084	CYC	C4C-NC	-4.37	1.28	1.37
4	P	1082	CYC	C1A-C2A	-4.37	1.38	1.45
4	R	1153	CYC	C1D-C2D	4.37	1.51	1.41
4	X	1153	CYC	C4B-C3B	-4.36	1.40	1.48
4	V	1153	CYC	CHA-C1A	4.36	1.46	1.38
4	T	1153	CYC	C1A-C2A	-4.35	1.38	1.45
4	R	1153	CYC	C4C-NC	-4.32	1.29	1.37
4	C	1084	CYC	CHA-C1A	4.32	1.46	1.38
4	C	1084	CYC	C1A-C2A	-4.29	1.39	1.45
4	S	1084	CYC	CHA-C1A	4.29	1.46	1.38
4	D	1082	CYC	CHD-C1D	4.29	1.50	1.40
4	O	1084	CYC	C1B-C2B	-4.28	1.37	1.45
4	Q	1084	CYC	C4C-NC	-4.27	1.29	1.37
4	C	1084	CYC	CHD-C1D	4.25	1.50	1.40
4	N	1153	CYC	CHA-C1A	4.25	1.46	1.38
4	F	1082	CYC	CHB-C1B	4.25	1.48	1.37
4	V	1082	CYC	CHD-C1D	4.24	1.49	1.40
4	L	1082	CYC	C4C-NC	-4.24	1.29	1.37
4	L	1153	CYC	C1B-C2B	-4.22	1.37	1.45
4	F	1153	CYC	CHA-C1A	4.21	1.46	1.38
4	P	1082	CYC	CHA-C1A	4.21	1.46	1.38
4	Q	1084	CYC	C1C-NC	-4.20	1.32	1.37
4	D	1082	CYC	CHA-C4D	4.20	1.49	1.40
4	M	1084	CYC	CHD-C1D	4.17	1.49	1.40
4	Q	1084	CYC	C1B-C2B	-4.16	1.37	1.45
4	U	1084	CYC	C1A-C2A	-4.14	1.39	1.45
4	L	1082	CYC	C2C-C1C	-4.13	1.48	1.52
4	V	1153	CYC	CHB-C1B	4.13	1.47	1.37
4	N	1153	CYC	CHD-C4C	-4.11	1.29	1.36
4	T	1153	CYC	C4C-NC	-4.11	1.29	1.37
4	L	1153	CYC	C1A-C2A	-4.11	1.39	1.45
4	B	1153	CYC	CHA-C1A	4.10	1.46	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	J	1082	CYC	CHA-C4D	4.10	1.49	1.40
4	D	1082	CYC	C4C-NC	-4.10	1.29	1.37
4	B	1153	CYC	C1D-C2D	4.10	1.51	1.41
4	D	1153	CYC	CHB-C1B	4.08	1.47	1.37
4	A	1084	CYC	CHB-C1B	4.08	1.47	1.37
4	H	1153	CYC	C1D-C2D	4.08	1.51	1.41
4	X	1082	CYC	CHB-C4A	4.05	1.50	1.40
4	F	1153	CYC	C4C-NC	-4.04	1.29	1.37
4	T	1153	CYC	C1D-C2D	4.04	1.51	1.41
4	T	1153	CYC	C2C-C1C	-4.03	1.48	1.52
4	V	1153	CYC	C4C-NC	-4.03	1.29	1.37
4	L	1082	CYC	C1A-C2A	-4.02	1.39	1.45
4	W	1084	CYC	C1A-C2A	-4.01	1.39	1.45
4	M	1084	CYC	CHB-C4A	4.01	1.49	1.40
4	V	1153	CYC	CHD-C1D	4.00	1.49	1.40
4	Q	1084	CYC	CHB-C4A	3.99	1.49	1.40
4	O	1084	CYC	CHA-C1A	3.99	1.46	1.38
4	X	1153	CYC	OB-C4B	3.98	1.31	1.23
4	V	1082	CYC	C1C-NC	-3.97	1.32	1.37
4	O	1084	CYC	CHA-C4D	3.97	1.49	1.40
4	N	1082	CYC	CHA-C4D	3.96	1.49	1.40
4	F	1082	CYC	C1A-C2A	-3.96	1.39	1.45
4	R	1153	CYC	CHB-C1B	3.95	1.47	1.37
4	H	1082	CYC	CHB-C1B	3.94	1.47	1.37
4	B	1082	CYC	OB-C4B	3.93	1.31	1.23
4	T	1082	CYC	CHA-C1A	3.91	1.46	1.38
4	O	1084	CYC	CHB-C4A	3.91	1.49	1.40
4	H	1082	CYC	CHB-C4A	3.89	1.49	1.40
4	G	1084	CYC	C4B-C3B	-3.88	1.41	1.48
4	J	1153	CYC	CHB-C4A	3.88	1.49	1.40
4	N	1082	CYC	C1C-NC	-3.88	1.32	1.37
4	D	1153	CYC	C4D-ND	-3.87	1.31	1.37
4	E	1084	CYC	CHB-C4A	3.87	1.49	1.40
4	F	1082	CYC	CHD-C1D	3.87	1.49	1.40
4	B	1082	CYC	C1D-C2D	3.86	1.50	1.41
4	L	1082	CYC	CHD-C1D	3.86	1.49	1.40
4	X	1082	CYC	C1D-C2D	3.84	1.50	1.41
4	D	1153	CYC	C1B-NB	-3.83	1.31	1.37
4	G	1084	CYC	C1D-C2D	3.82	1.50	1.41
4	P	1153	CYC	C4C-NC	-3.82	1.30	1.37
4	K	1084	CYC	CHB-C4A	3.81	1.49	1.40
4	V	1153	CYC	C1C-NC	-3.81	1.32	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	P	1153	CYC	CHD-C1D	3.81	1.48	1.40
4	D	1153	CYC	C1A-C2A	-3.79	1.39	1.45
4	B	1082	CYC	C1B-C2B	-3.78	1.38	1.45
4	J	1082	CYC	CHD-C1D	3.78	1.48	1.40
4	U	1084	CYC	CHB-C4A	3.78	1.49	1.40
4	H	1153	CYC	CHA-C1A	3.77	1.45	1.38
4	N	1153	CYC	CHB-C4A	3.76	1.49	1.40
4	P	1082	CYC	C1D-C2D	3.75	1.50	1.41
4	J	1082	CYC	OB-C4B	3.74	1.30	1.23
4	Q	1084	CYC	CHD-C1D	3.74	1.48	1.40
4	J	1153	CYC	CHD-C1D	3.73	1.48	1.40
4	I	1084	CYC	CHA-C1A	3.72	1.45	1.38
4	R	1153	CYC	CBA-CGA	3.72	1.59	1.50
4	T	1082	CYC	CHB-C1B	3.71	1.46	1.37
4	G	1084	CYC	CHB-C4A	3.70	1.49	1.40
4	V	1153	CYC	C1B-NB	-3.70	1.31	1.37
4	T	1153	CYC	CBA-CGA	3.70	1.59	1.50
4	E	1084	CYC	C1C-NC	-3.69	1.32	1.37
4	L	1082	CYC	CHA-C4D	3.69	1.48	1.40
4	S	1084	CYC	C1D-C2D	3.69	1.50	1.41
4	R	1082	CYC	C4C-NC	-3.68	1.30	1.37
4	V	1153	CYC	C1D-C2D	3.68	1.50	1.41
4	F	1082	CYC	CHA-C4D	3.68	1.48	1.40
4	E	1084	CYC	C1D-C2D	3.68	1.50	1.41
4	F	1082	CYC	OB-C4B	3.68	1.30	1.23
4	X	1082	CYC	C1C-NC	-3.67	1.32	1.37
4	H	1153	CYC	CHD-C1D	3.67	1.48	1.40
4	R	1153	CYC	CHD-C1D	3.67	1.48	1.40
4	V	1153	CYC	C2C-C1C	-3.66	1.48	1.52
4	N	1153	CYC	C1C-NC	-3.65	1.32	1.37
4	R	1153	CYC	C1B-NB	-3.65	1.31	1.37
4	D	1082	CYC	CHB-C4A	3.63	1.49	1.40
4	T	1153	CYC	CHB-C1B	3.63	1.46	1.37
4	N	1153	CYC	CHD-C1D	3.63	1.48	1.40
4	N	1153	CYC	C4D-C3D	3.62	1.50	1.42
4	U	1084	CYC	OB-C4B	3.62	1.30	1.23
4	S	1084	CYC	C2C-C1C	-3.61	1.48	1.52
4	H	1082	CYC	CHA-C4D	3.61	1.48	1.40
4	W	1084	CYC	CHA-C4D	3.60	1.48	1.40
4	J	1153	CYC	C1D-C2D	3.60	1.50	1.41
4	L	1153	CYC	CHD-C1D	3.59	1.48	1.40
4	P	1082	CYC	C1C-NC	-3.59	1.33	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	1084	CYC	CHB-C1B	3.58	1.46	1.37
4	I	1084	CYC	C1A-C2A	-3.58	1.40	1.45
4	W	1084	CYC	OB-C4B	3.57	1.30	1.23
4	F	1153	CYC	CHD-C1D	3.57	1.48	1.40
4	A	1084	CYC	CHB-C4A	3.56	1.48	1.40
4	B	1153	CYC	C1B-C2B	-3.55	1.38	1.45
4	C	1084	CYC	C1B-C2B	-3.54	1.38	1.45
4	N	1153	CYC	C1B-NB	-3.54	1.31	1.37
4	J	1082	CYC	C1A-C2A	-3.54	1.40	1.45
4	B	1082	CYC	CHB-C4A	3.53	1.48	1.40
4	N	1153	CYC	C1D-C2D	3.53	1.49	1.41
4	A	1084	CYC	CHA-C4D	3.53	1.48	1.40
4	D	1082	CYC	C1D-C2D	3.52	1.49	1.41
4	P	1153	CYC	CHB-C4A	3.52	1.48	1.40
4	M	1084	CYC	CHA-C4D	3.52	1.48	1.40
4	B	1153	CYC	CHB-C4A	3.51	1.48	1.40
4	J	1082	CYC	C1C-NC	-3.51	1.33	1.37
4	X	1153	CYC	C4C-NC	-3.51	1.30	1.37
4	F	1082	CYC	C1D-C2D	3.48	1.49	1.41
4	L	1082	CYC	C4A-C3A	3.47	1.53	1.45
4	B	1153	CYC	CHD-C1D	3.45	1.48	1.40
4	C	1084	CYC	C4C-NC	-3.45	1.30	1.37
4	X	1153	CYC	CHD-C1D	3.44	1.48	1.40
4	R	1082	CYC	C1B-NB	-3.44	1.31	1.37
4	L	1153	CYC	C4D-ND	-3.44	1.32	1.37
4	B	1153	CYC	C1C-NC	-3.43	1.33	1.37
4	B	1082	CYC	CHA-C1A	3.43	1.45	1.38
4	T	1153	CYC	C1B-C2B	-3.43	1.39	1.45
4	P	1153	CYC	C1D-C2D	3.42	1.49	1.41
4	R	1153	CYC	CHA-C1A	3.41	1.45	1.38
4	W	1084	CYC	C1D-C2D	3.41	1.49	1.41
4	B	1082	CYC	CHA-C4D	3.41	1.48	1.40
4	H	1082	CYC	C1D-C2D	3.40	1.49	1.41
4	O	1084	CYC	C4D-C3D	3.40	1.50	1.42
4	V	1082	CYC	C1A-C2A	-3.39	1.40	1.45
4	D	1082	CYC	C1B-NB	-3.38	1.32	1.37
4	X	1153	CYC	CHA-C4D	3.37	1.48	1.40
4	I	1084	CYC	C1D-C2D	3.37	1.49	1.41
4	J	1153	CYC	C1B-C2B	-3.36	1.39	1.45
4	O	1084	CYC	C4C-NC	-3.34	1.30	1.37
4	X	1153	CYC	C4D-C3D	3.33	1.50	1.42
4	E	1084	CYC	C1B-C2B	-3.32	1.39	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	N	1082	CYC	OB-C4B	3.32	1.29	1.23
4	S	1084	CYC	CHA-C4D	3.31	1.47	1.40
4	B	1153	CYC	C4D-C3D	3.31	1.50	1.42
4	K	1084	CYC	C4D-C3D	3.30	1.50	1.42
4	V	1082	CYC	CHA-C4D	3.30	1.47	1.40
4	U	1084	CYC	C4C-NC	-3.29	1.31	1.37
4	H	1153	CYC	CHB-C4A	3.28	1.48	1.40
4	K	1084	CYC	C1C-NC	-3.28	1.33	1.37
4	N	1082	CYC	C4C-NC	-3.28	1.31	1.37
4	W	1084	CYC	C4A-C3A	3.27	1.52	1.45
4	N	1082	CYC	C1D-C2D	3.26	1.49	1.41
4	T	1082	CYC	CHB-C4A	3.25	1.48	1.40
4	L	1153	CYC	C1D-C2D	3.24	1.49	1.41
4	E	1084	CYC	CHA-C4D	3.24	1.47	1.40
4	F	1153	CYC	CHB-C4A	3.23	1.48	1.40
4	I	1084	CYC	C1B-C2B	-3.22	1.39	1.45
4	E	1084	CYC	C4B-NB	-3.22	1.30	1.38
4	K	1084	CYC	C1D-C2D	3.21	1.49	1.41
4	V	1153	CYC	CHA-C4D	3.21	1.47	1.40
4	S	1084	CYC	C1B-C2B	-3.20	1.39	1.45
4	R	1153	CYC	OB-C4B	3.20	1.29	1.23
4	H	1082	CYC	C1C-NC	-3.19	1.33	1.37
4	H	1082	CYC	OB-C4B	3.17	1.29	1.23
4	J	1082	CYC	C4A-C3A	3.17	1.52	1.45
4	V	1082	CYC	C4C-NC	-3.16	1.31	1.37
4	R	1153	CYC	C1B-C2B	-3.15	1.39	1.45
4	O	1084	CYC	C1A-C2A	-3.15	1.40	1.45
4	D	1153	CYC	C1D-ND	-3.15	1.32	1.37
4	G	1084	CYC	C1C-NC	-3.14	1.33	1.37
4	C	1084	CYC	C4D-ND	-3.13	1.32	1.37
4	F	1153	CYC	OB-C4B	3.12	1.29	1.23
4	S	1084	CYC	C4C-NC	-3.12	1.31	1.37
4	C	1084	CYC	C1C-NC	-3.12	1.33	1.37
4	S	1084	CYC	C1B-NB	-3.11	1.32	1.37
4	W	1084	CYC	C1B-C2B	-3.11	1.39	1.45
4	R	1153	CYC	C1A-NA	-3.10	1.32	1.38
4	J	1082	CYC	C2C-C1C	-3.10	1.49	1.52
4	J	1082	CYC	C4D-C3D	3.10	1.49	1.42
4	U	1084	CYC	C1B-C2B	-3.10	1.39	1.45
4	R	1153	CYC	C4D-ND	-3.09	1.32	1.37
4	J	1082	CYC	C4C-NC	-3.09	1.31	1.37
4	D	1153	CYC	CHD-C1D	3.08	1.47	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	V	1082	CYC	C1D-C2D	3.08	1.48	1.41
4	U	1084	CYC	C2C-C1C	-3.08	1.49	1.52
4	F	1153	CYC	C1D-C2D	3.08	1.48	1.41
4	D	1082	CYC	C1A-C2A	-3.08	1.40	1.45
4	E	1084	CYC	C4C-NC	-3.06	1.31	1.37
4	A	1084	CYC	OB-C4B	3.04	1.29	1.23
4	J	1082	CYC	C1D-C2D	3.03	1.48	1.41
4	N	1082	CYC	C1A-C2A	-3.03	1.41	1.45
4	Q	1084	CYC	CHA-C4D	3.02	1.47	1.40
4	K	1084	CYC	OB-C4B	3.02	1.29	1.23
4	D	1153	CYC	CHA-C4D	3.02	1.47	1.40
4	M	1084	CYC	C1B-C2B	-3.01	1.39	1.45
4	P	1153	CYC	C4D-ND	-3.01	1.32	1.37
4	Q	1084	CYC	OB-C4B	3.00	1.29	1.23
4	T	1082	CYC	C2C-C1C	-3.00	1.49	1.52
4	A	1084	CYC	C1A-C2A	-3.00	1.41	1.45
4	T	1082	CYC	C4C-NC	-2.99	1.31	1.37
4	Q	1084	CYC	C1B-NB	-2.99	1.32	1.37
4	G	1084	CYC	OB-C4B	2.98	1.29	1.23
4	J	1153	CYC	C1B-NB	-2.97	1.32	1.37
4	N	1082	CYC	C1B-NB	-2.97	1.32	1.37
4	O	1084	CYC	C1D-C2D	2.96	1.48	1.41
4	L	1153	CYC	C1B-NB	-2.96	1.32	1.37
4	U	1084	CYC	CHA-C4D	2.95	1.47	1.40
4	J	1153	CYC	C4C-NC	-2.95	1.31	1.37
4	K	1084	CYC	C1A-NA	-2.93	1.32	1.38
4	K	1084	CYC	C1B-NB	-2.92	1.32	1.37
4	F	1082	CYC	CHB-C4A	2.92	1.47	1.40
4	X	1082	CYC	OB-C4B	2.91	1.29	1.23
4	I	1084	CYC	C4A-C3A	2.90	1.52	1.45
4	R	1082	CYC	C4D-C3D	2.90	1.49	1.42
4	V	1153	CYC	C1A-NA	-2.90	1.32	1.38
4	J	1153	CYC	C1C-NC	-2.90	1.33	1.37
4	H	1153	CYC	C1A-NA	-2.89	1.32	1.38
4	U	1084	CYC	C1D-C2D	2.87	1.48	1.41
4	M	1084	CYC	C4C-NC	-2.86	1.31	1.37
4	M	1084	CYC	C1C-NC	-2.84	1.34	1.37
4	P	1153	CYC	C1C-NC	-2.84	1.34	1.37
4	D	1153	CYC	CHB-C4A	2.84	1.47	1.40
4	E	1084	CYC	C1B-NB	-2.84	1.33	1.37
4	T	1153	CYC	C4D-ND	-2.83	1.33	1.37
4	B	1153	CYC	CHA-C4D	2.83	1.46	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	T	1153	CYC	C1B-NB	-2.82	1.33	1.37
4	A	1084	CYC	O1A-CGA	2.82	1.31	1.22
4	G	1084	CYC	C1A-C2A	-2.81	1.41	1.45
4	C	1084	CYC	C1D-C2D	2.81	1.48	1.41
4	P	1082	CYC	C1B-C2B	-2.80	1.40	1.45
4	H	1082	CYC	C2C-C1C	-2.79	1.49	1.52
4	R	1082	CYC	C1A-C2A	-2.78	1.41	1.45
4	N	1082	CYC	C4A-C3A	2.78	1.51	1.45
4	N	1153	CYC	C4D-ND	-2.78	1.33	1.37
4	A	1084	CYC	C1C-NC	-2.77	1.34	1.37
4	C	1084	CYC	C1B-NB	-2.76	1.33	1.37
4	T	1153	CYC	C1A-NA	-2.75	1.32	1.38
4	X	1153	CYC	C4D-ND	-2.75	1.33	1.37
4	O	1084	CYC	OB-C4B	2.75	1.28	1.23
4	D	1082	CYC	C4A-C3A	2.74	1.51	1.45
4	X	1153	CYC	C1D-C2D	2.74	1.48	1.41
4	X	1153	CYC	C1C-NC	-2.74	1.34	1.37
4	N	1153	CYC	C1B-C2B	-2.73	1.40	1.45
4	T	1153	CYC	CHB-C4A	2.73	1.46	1.40
4	T	1082	CYC	CHA-C4D	2.73	1.46	1.40
4	R	1153	CYC	CHB-C4A	2.72	1.46	1.40
4	J	1153	CYC	OB-C4B	2.72	1.28	1.23
4	N	1082	CYC	C4D-C3D	2.72	1.48	1.42
4	P	1082	CYC	CHA-C4D	2.72	1.46	1.40
4	H	1082	CYC	C4C-NC	-2.72	1.32	1.37
4	K	1084	CYC	C1B-C2B	-2.71	1.40	1.45
4	M	1084	CYC	C1B-NB	-2.69	1.33	1.37
4	S	1084	CYC	OB-C4B	2.69	1.28	1.23
4	P	1082	CYC	C1B-NB	-2.69	1.33	1.37
4	B	1153	CYC	C1B-NB	-2.68	1.33	1.37
4	I	1084	CYC	C1B-NB	-2.68	1.33	1.37
4	V	1153	CYC	O1A-CGA	2.66	1.30	1.22
4	A	1084	CYC	C1B-C2B	-2.64	1.40	1.45
4	R	1082	CYC	C1D-C2D	2.64	1.47	1.41
4	P	1153	CYC	CHA-C4D	2.62	1.46	1.40
4	H	1082	CYC	C1B-NB	-2.62	1.33	1.37
4	C	1084	CYC	C1A-NA	-2.60	1.33	1.38
4	C	1084	CYC	CHA-C4D	2.60	1.46	1.40
4	K	1084	CYC	C4C-NC	-2.59	1.32	1.37
4	L	1153	CYC	OB-C4B	2.58	1.28	1.23
4	H	1082	CYC	C1A-C2A	-2.58	1.41	1.45
4	S	1084	CYC	C4A-C3A	2.58	1.51	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	1082	CYC	C1D-ND	-2.57	1.33	1.37
4	B	1153	CYC	CHD-C4C	-2.55	1.32	1.36
4	K	1084	CYC	C4B-NB	-2.55	1.32	1.38
4	B	1082	CYC	C1B-NB	-2.54	1.33	1.37
4	C	1084	CYC	CHB-C4A	2.54	1.46	1.40
4	P	1082	CYC	C4A-C3A	2.53	1.51	1.45
4	G	1084	CYC	C4C-NC	-2.52	1.32	1.37
4	J	1153	CYC	C4D-C3D	2.52	1.48	1.42
4	T	1082	CYC	C4A-C3A	2.52	1.51	1.45
4	F	1153	CYC	C1B-C2B	-2.51	1.40	1.45
4	I	1084	CYC	C4D-C3D	2.51	1.48	1.42
4	F	1082	CYC	C1B-C2B	-2.50	1.40	1.45
4	H	1153	CYC	C1B-C2B	-2.50	1.40	1.45
4	F	1153	CYC	C1C-NC	-2.49	1.34	1.37
4	D	1153	CYC	CHD-C4C	-2.49	1.32	1.36
4	H	1082	CYC	C1B-C2B	-2.48	1.40	1.45
4	P	1153	CYC	C1D-ND	-2.47	1.33	1.37
4	V	1153	CYC	CBA-CGA	2.47	1.56	1.50
4	V	1082	CYC	C4A-C3A	2.47	1.51	1.45
4	D	1153	CYC	C1D-C2D	2.46	1.47	1.41
4	N	1153	CYC	C4A-C3A	2.46	1.51	1.45
4	N	1153	CYC	CBA-CGA	2.46	1.56	1.50
4	C	1084	CYC	OB-C4B	2.45	1.28	1.23
4	T	1153	CYC	C4D-C3D	2.44	1.48	1.42
4	M	1084	CYC	C1D-C2D	2.44	1.47	1.41
4	P	1082	CYC	C4C-NC	-2.43	1.32	1.37
4	V	1153	CYC	C1B-C2B	-2.43	1.40	1.45
4	N	1153	CYC	CHA-C4D	2.42	1.45	1.40
4	H	1153	CYC	CBA-CAA	2.42	1.60	1.51
4	L	1082	CYC	C4B-NB	-2.42	1.32	1.38
4	D	1082	CYC	C4D-C3D	2.41	1.48	1.42
4	H	1153	CYC	CHA-C4D	2.41	1.45	1.40
4	B	1153	CYC	C1A-NA	-2.41	1.33	1.38
4	A	1084	CYC	C1B-NB	-2.41	1.33	1.37
4	P	1153	CYC	C1A-NA	-2.41	1.33	1.38
4	T	1153	CYC	CBA-CAA	2.40	1.60	1.51
4	H	1153	CYC	C1D-ND	-2.40	1.33	1.37
4	F	1153	CYC	C1A-NA	-2.39	1.33	1.38
4	F	1153	CYC	C4D-ND	-2.39	1.33	1.37
4	P	1153	CYC	OB-C4B	2.39	1.28	1.23
4	R	1153	CYC	OC-C1C	-2.39	1.18	1.23
4	E	1084	CYC	OB-C4B	2.39	1.28	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	H	1153	CYC	CBA-CGA	2.38	1.56	1.50
4	V	1153	CYC	C4D-C3D	2.38	1.48	1.42
4	E	1084	CYC	C4D-C3D	2.38	1.48	1.42
4	J	1153	CYC	C4D-ND	-2.37	1.33	1.37
4	N	1153	CYC	OB-C4B	2.36	1.28	1.23
4	A	1084	CYC	C1A-NA	-2.36	1.33	1.38
4	R	1082	CYC	C4B-NB	-2.36	1.32	1.38
4	P	1153	CYC	C1B-C2B	-2.36	1.40	1.45
4	O	1084	CYC	O1D-CGD	2.35	1.29	1.22
4	L	1153	CYC	C1D-ND	-2.35	1.33	1.37
4	T	1082	CYC	C1A-NA	-2.35	1.33	1.38
4	L	1153	CYC	O2A-CGA	-2.35	1.23	1.30
4	L	1153	CYC	CBA-CGA	2.34	1.56	1.50
4	U	1084	CYC	C4A-C3A	2.33	1.50	1.45
4	L	1153	CYC	CHB-C4A	2.33	1.46	1.40
4	H	1082	CYC	C4D-C3D	2.33	1.48	1.42
4	X	1082	CYC	C1B-NB	-2.33	1.33	1.37
4	R	1082	CYC	C1B-C2B	-2.33	1.40	1.45
4	R	1153	CYC	CBA-CAA	2.32	1.60	1.51
4	B	1082	CYC	C1A-NA	-2.31	1.33	1.38
4	J	1153	CYC	CHA-C4D	2.31	1.45	1.40
4	H	1082	CYC	C2A-C3A	2.31	1.41	1.36
4	R	1082	CYC	OB-C4B	2.31	1.28	1.23
4	L	1153	CYC	CHA-C4D	2.31	1.45	1.40
4	H	1153	CYC	C4D-ND	-2.31	1.33	1.37
4	V	1153	CYC	CHB-C4A	2.31	1.45	1.40
4	D	1153	CYC	C4D-C3D	2.30	1.47	1.42
4	B	1153	CYC	OB-C4B	2.30	1.27	1.23
4	N	1153	CYC	C1A-NA	-2.29	1.33	1.38
4	D	1153	CYC	C1B-C2B	-2.29	1.41	1.45
4	V	1082	CYC	C2A-C3A	2.29	1.41	1.36
4	G	1084	CYC	C4A-C3A	2.29	1.50	1.45
4	X	1082	CYC	C1A-NA	-2.28	1.33	1.38
4	F	1153	CYC	C1B-NB	-2.28	1.33	1.37
4	O	1084	CYC	C1B-NB	-2.28	1.33	1.37
4	F	1082	CYC	C1D-ND	-2.27	1.33	1.37
4	T	1153	CYC	CHA-C4D	2.27	1.45	1.40
4	U	1084	CYC	O1D-CGD	2.27	1.29	1.22
4	D	1153	CYC	O1D-CGD	2.27	1.29	1.22
4	I	1084	CYC	CHA-C4D	2.26	1.45	1.40
4	M	1084	CYC	OB-C4B	2.26	1.27	1.23
4	D	1153	CYC	C1A-NA	-2.25	1.33	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	E	1084	CYC	C1A-NA	-2.24	1.33	1.38
4	Q	1084	CYC	C1D-C2D	2.22	1.46	1.41
4	X	1082	CYC	C4A-C3A	2.22	1.50	1.45
4	X	1153	CYC	C1B-C2B	-2.21	1.41	1.45
4	T	1082	CYC	OB-C4B	2.21	1.27	1.23
4	H	1153	CYC	C4B-NB	-2.21	1.32	1.38
4	D	1082	CYC	OB-C4B	2.20	1.27	1.23
4	A	1084	CYC	O1D-CGD	2.19	1.29	1.22
4	M	1084	CYC	C4D-ND	-2.19	1.34	1.37
4	B	1082	CYC	C4A-NA	-2.18	1.31	1.36
4	L	1082	CYC	C1B-NB	-2.17	1.34	1.37
4	F	1082	CYC	C1B-NB	-2.16	1.34	1.37
4	I	1084	CYC	C1A-NA	-2.16	1.34	1.38
4	T	1153	CYC	OB-C4B	2.16	1.27	1.23
4	H	1153	CYC	C1B-NB	-2.15	1.34	1.37
4	A	1084	CYC	C4D-C3D	2.15	1.47	1.42
4	P	1153	CYC	O2A-CGA	-2.15	1.23	1.30
4	V	1082	CYC	C1B-NB	-2.15	1.34	1.37
4	X	1082	CYC	C1B-C2B	-2.14	1.41	1.45
4	F	1153	CYC	CHA-C4D	2.14	1.45	1.40
4	M	1084	CYC	C2A-C3A	2.14	1.41	1.36
4	M	1084	CYC	C4B-NB	-2.14	1.33	1.38
4	B	1153	CYC	O2A-CGA	-2.13	1.23	1.30
4	P	1082	CYC	OB-C4B	2.13	1.27	1.23
4	M	1084	CYC	C4A-C3A	2.12	1.50	1.45
4	F	1082	CYC	C4A-C3A	2.12	1.50	1.45
4	L	1153	CYC	CBA-CAA	2.12	1.59	1.51
4	V	1082	CYC	OB-C4B	2.12	1.27	1.23
4	V	1082	CYC	CAC-C3C	-2.12	1.50	1.54
4	S	1084	CYC	C1A-NA	-2.11	1.34	1.38
4	D	1082	CYC	C4B-NB	-2.10	1.33	1.38
4	J	1082	CYC	C2A-C3A	2.10	1.41	1.36
4	E	1084	CYC	C1D-ND	-2.09	1.34	1.37
4	B	1082	CYC	C4D-ND	-2.09	1.34	1.37
4	L	1153	CYC	C1A-NA	-2.08	1.34	1.38
4	T	1082	CYC	C1B-C2B	-2.08	1.41	1.45
4	H	1082	CYC	C4A-C3A	2.08	1.50	1.45
4	J	1153	CYC	C1A-NA	-2.07	1.34	1.38
4	O	1084	CYC	O1A-CGA	2.07	1.28	1.22
4	X	1153	CYC	C1B-NB	-2.06	1.34	1.37
4	D	1082	CYC	C1B-C2B	-2.04	1.41	1.45
4	W	1084	CYC	C4B-NB	-2.04	1.33	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	L	1082	CYC	C1D-ND	-2.04	1.34	1.37
4	I	1084	CYC	O1D-CGD	2.03	1.28	1.22
4	L	1082	CYC	C1D-C2D	2.03	1.46	1.41
4	O	1084	CYC	CMC-C2C	-2.03	1.48	1.53
4	N	1153	CYC	C4B-NB	-2.02	1.33	1.38
4	H	1082	CYC	C1A-NA	-2.02	1.34	1.38
4	O	1084	CYC	C4B-NB	-2.01	1.33	1.38
4	M	1084	CYC	C4D-C3D	2.01	1.47	1.42
4	O	1084	CYC	C4A-C3A	2.00	1.50	1.45

All (1140) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	V	1153	CYC	C3B-C4B-NB	16.83	120.19	106.77
4	D	1153	CYC	C3B-C4B-NB	14.95	118.69	106.77
4	V	1082	CYC	C3B-C4B-NB	14.60	118.41	106.77
4	T	1153	CYC	C3B-C4B-NB	14.47	118.31	106.77
4	R	1082	CYC	C3B-C4B-NB	14.30	118.17	106.77
4	I	1084	CYC	C3B-C4B-NB	14.23	118.12	106.77
4	J	1153	CYC	C3B-C4B-NB	13.88	117.84	106.77
4	N	1153	CYC	C3B-C4B-NB	13.86	117.82	106.77
4	H	1153	CYC	C3B-C4B-NB	13.82	117.79	106.77
4	P	1082	CYC	C3B-C4B-NB	13.79	117.77	106.77
4	D	1082	CYC	C3B-C4B-NB	13.55	117.57	106.77
4	P	1153	CYC	C3B-C4B-NB	13.43	117.48	106.77
4	L	1153	CYC	C3B-C4B-NB	13.39	117.44	106.77
4	B	1153	CYC	C3B-C4B-NB	13.34	117.41	106.77
4	L	1082	CYC	C3B-C4B-NB	13.12	117.23	106.77
4	S	1084	CYC	C3B-C4B-NB	13.06	117.18	106.77
4	R	1153	CYC	C3B-C4B-NB	12.86	117.02	106.77
4	X	1082	CYC	C3B-C4B-NB	12.72	116.91	106.77
4	N	1082	CYC	C3B-C4B-NB	12.54	116.77	106.77
4	T	1082	CYC	C3B-C4B-NB	12.16	116.46	106.77
4	E	1084	CYC	CAA-CBA-CGA	-12.11	81.56	113.67
4	C	1084	CYC	C3B-C4B-NB	12.04	116.37	106.77
4	F	1153	CYC	C3B-C4B-NB	11.95	116.30	106.77
4	A	1084	CYC	C3B-C4B-NB	11.86	116.23	106.77
4	J	1153	CYC	OB-C4B-C3B	-11.74	115.70	128.03
4	M	1084	CYC	C3B-C4B-NB	11.62	116.03	106.77
4	L	1153	CYC	C2C-C1C-NC	11.60	117.95	108.29
4	H	1082	CYC	C3B-C4B-NB	11.56	115.99	106.77
4	J	1082	CYC	C3B-C4B-NB	11.48	115.92	106.77

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	E	1084	CYC	C3B-C4B-NB	11.47	115.91	106.77
4	B	1082	CYC	C3B-C4B-NB	11.46	115.91	106.77
4	O	1084	CYC	CAA-CBA-CGA	-11.33	83.61	113.67
4	T	1153	CYC	OB-C4B-C3B	-11.20	116.26	128.03
4	U	1084	CYC	C3B-C4B-NB	11.01	115.55	106.77
4	K	1084	CYC	C3B-C4B-NB	10.99	115.53	106.77
4	F	1153	CYC	OC-C1C-C2C	-10.91	117.50	126.17
4	F	1082	CYC	C3B-C4B-NB	10.51	115.15	106.77
4	R	1153	CYC	C2C-C1C-NC	10.47	117.00	108.29
4	J	1153	CYC	OC-C1C-C2C	-10.35	117.94	126.17
4	L	1153	CYC	OB-C4B-C3B	-10.27	117.24	128.03
4	G	1084	CYC	C3B-C4B-NB	10.15	114.86	106.77
4	V	1153	CYC	OB-C4B-C3B	-10.11	117.40	128.03
4	O	1084	CYC	C3B-C4B-NB	10.07	114.80	106.77
4	B	1153	CYC	OC-C1C-C2C	-10.07	118.17	126.17
4	B	1153	CYC	C2C-C1C-NC	9.98	116.59	108.29
4	W	1084	CYC	C3B-C4B-NB	9.95	114.70	106.77
4	N	1153	CYC	OB-C4B-C3B	-9.83	117.70	128.03
4	C	1084	CYC	CAA-CBA-CGA	-9.76	87.78	113.67
4	D	1153	CYC	C2C-C1C-NC	9.56	116.24	108.29
4	X	1153	CYC	C3B-C4B-NB	9.50	114.34	106.77
4	V	1153	CYC	C1B-NB-C4B	-9.26	99.30	110.66
4	P	1153	CYC	OC-C1C-C2C	-9.22	118.84	126.17
4	Q	1084	CYC	C3B-C4B-NB	9.19	114.10	106.77
4	H	1153	CYC	CMC-C2C-C1C	-9.12	92.74	112.40
4	T	1153	CYC	C1B-NB-C4B	-9.09	99.50	110.66
4	K	1084	CYC	CAA-CBA-CGA	-8.88	90.11	113.67
4	B	1153	CYC	CMC-C2C-C1C	-8.80	93.43	112.40
4	R	1153	CYC	OB-C4B-C3B	-8.79	118.80	128.03
4	H	1153	CYC	OC-C1C-C2C	-8.73	119.23	126.17
4	B	1153	CYC	OB-C4B-C3B	-8.70	118.89	128.03
4	B	1082	CYC	OB-C4B-C3B	-8.61	118.99	128.03
4	I	1084	CYC	C1B-NB-C4B	-8.59	100.12	110.66
4	C	1084	CYC	C1B-NB-C4B	-8.53	100.19	110.66
4	H	1082	CYC	OB-C4B-C3B	-8.47	119.13	128.03
4	R	1082	CYC	C1B-NB-C4B	-8.44	100.30	110.66
4	S	1084	CYC	CAA-CBA-CGA	-8.36	91.48	113.67
4	I	1084	CYC	C3D-C4D-ND	8.33	118.01	107.57
4	T	1082	CYC	C1B-NB-C4B	-8.33	100.44	110.66
4	X	1153	CYC	OC-C1C-C2C	-8.30	119.57	126.17
4	V	1082	CYC	C1B-NB-C4B	-8.27	100.51	110.66
4	H	1153	CYC	C2C-C1C-NC	8.22	115.12	108.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1153	CYC	C1B-NB-C4B	-8.15	100.66	110.66
4	L	1153	CYC	C1B-NB-C4B	-8.12	100.69	110.66
4	I	1084	CYC	CAA-CBA-CGA	-7.98	92.49	113.67
4	N	1082	CYC	OC-C1C-C2C	-7.97	119.84	126.17
4	J	1153	CYC	C1B-NB-C4B	-7.94	100.92	110.66
4	S	1084	CYC	OC-C1C-C2C	-7.94	119.86	126.17
4	D	1153	CYC	C1B-NB-C4B	-7.91	100.95	110.66
4	R	1153	CYC	C1B-NB-C4B	-7.89	100.98	110.66
4	F	1153	CYC	C2C-C1C-NC	7.88	114.84	108.29
4	I	1084	CYC	OB-C4B-C3B	-7.83	119.80	128.03
4	N	1082	CYC	CMB-C2B-C1B	7.81	133.66	124.16
4	D	1153	CYC	OB-C4B-C3B	-7.76	119.87	128.03
4	P	1153	CYC	OB-C4B-C3B	-7.76	119.88	128.03
4	X	1082	CYC	OB-C4B-C3B	-7.75	119.89	128.03
4	T	1082	CYC	C3D-C4D-ND	7.74	117.28	107.57
4	E	1084	CYC	CAB-C3B-C4B	7.69	133.26	121.37
4	A	1084	CYC	C1B-NB-C4B	-7.65	101.27	110.66
4	F	1082	CYC	OC-C1C-C2C	-7.62	120.11	126.17
4	D	1082	CYC	C1B-NB-C4B	-7.58	101.36	110.66
4	R	1153	CYC	OC-C1C-C2C	-7.58	120.15	126.17
4	N	1082	CYC	OB-C4B-C3B	-7.55	120.09	128.03
4	H	1082	CYC	C1B-NB-C4B	-7.54	101.40	110.66
4	P	1153	CYC	C1B-NB-C4B	-7.53	101.42	110.66
4	H	1153	CYC	C1B-NB-C4B	-7.49	101.47	110.66
4	B	1153	CYC	CHD-C4C-NC	-7.47	116.31	125.63
4	L	1153	CYC	CAC-C3C-C4C	-7.46	93.51	112.67
4	W	1084	CYC	CAA-CBA-CGA	-7.44	93.94	113.67
4	J	1082	CYC	OB-C4B-C3B	-7.42	120.23	128.03
4	N	1082	CYC	CHB-C1B-NB	-7.37	110.35	126.06
4	B	1082	CYC	C3D-C4D-ND	7.31	116.73	107.57
4	T	1153	CYC	C3D-C4D-ND	7.28	116.70	107.57
4	I	1084	CYC	C2A-C1A-NA	7.27	120.33	110.04
4	P	1153	CYC	CMC-C2C-C1C	-7.26	96.76	112.40
4	F	1082	CYC	C2C-C1C-NC	7.25	114.32	108.29
4	N	1153	CYC	CMC-C2C-C1C	-7.21	96.86	112.40
4	M	1084	CYC	CAA-CBA-CGA	-7.17	94.65	113.67
4	F	1082	CYC	C1B-NB-C4B	-7.16	101.87	110.66
4	E	1084	CYC	C1B-NB-C4B	-7.13	101.91	110.66
4	N	1153	CYC	C1B-NB-C4B	-7.12	101.92	110.66
4	D	1153	CYC	CAB-C3B-C4B	7.10	132.35	121.37
4	R	1082	CYC	OB-C4B-C3B	-7.10	120.57	128.03
4	B	1082	CYC	C1B-NB-C4B	-7.04	102.02	110.66

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	R	1082	CYC	CHB-C1B-NB	-7.03	111.06	126.06
4	L	1082	CYC	C1B-NB-C4B	-7.01	102.05	110.66
4	P	1082	CYC	OC-C1C-C2C	-7.01	120.60	126.17
4	P	1082	CYC	C1B-NB-C4B	-7.01	102.06	110.66
4	I	1084	CYC	C2C-C1C-NC	6.99	114.10	108.29
4	P	1082	CYC	C3D-C4D-ND	6.99	116.33	107.57
4	Q	1084	CYC	CAA-CBA-CGA	-6.96	95.19	113.67
4	F	1153	CYC	C1B-NB-C4B	-6.91	102.18	110.66
4	B	1082	CYC	C2A-C1A-NA	6.87	119.77	110.04
4	F	1153	CYC	C3D-C4D-ND	6.85	116.16	107.57
4	J	1153	CYC	CMC-C2C-C1C	-6.85	97.65	112.40
4	N	1082	CYC	C1B-NB-C4B	-6.83	102.28	110.66
4	S	1084	CYC	C1B-NB-C4B	-6.79	102.32	110.66
4	C	1084	CYC	C3D-C4D-ND	6.79	116.08	107.57
4	D	1153	CYC	OC-C1C-C2C	-6.72	120.83	126.17
4	U	1084	CYC	C1B-NB-C4B	-6.70	102.44	110.66
4	R	1153	CYC	C3D-C4D-ND	6.68	115.95	107.57
4	M	1084	CYC	OB-C4B-C3B	-6.67	121.03	128.03
4	T	1153	CYC	C2C-C1C-NC	6.66	113.83	108.29
4	D	1153	CYC	CHD-C1D-ND	-6.65	110.27	125.29
4	X	1153	CYC	C2C-C1C-NC	6.62	113.79	108.29
4	P	1153	CYC	C2C-C1C-NC	6.61	113.79	108.29
4	L	1082	CYC	CAA-CBA-CGA	-6.60	96.17	113.67
4	X	1082	CYC	OC-C1C-C2C	-6.58	120.94	126.17
4	M	1084	CYC	C1B-NB-C4B	-6.58	102.59	110.66
4	C	1084	CYC	C2C-C1C-NC	6.57	113.75	108.29
4	T	1153	CYC	OC-C1C-C2C	-6.56	120.96	126.17
4	T	1082	CYC	OB-C4B-C3B	-6.53	121.16	128.03
4	H	1153	CYC	C3D-C4D-ND	6.47	115.68	107.57
4	J	1153	CYC	C3D-C4D-ND	6.44	115.64	107.57
4	W	1084	CYC	C3D-C4D-ND	6.40	115.60	107.57
4	X	1153	CYC	CAB-C3B-C4B	6.40	131.27	121.37
4	P	1153	CYC	CAB-C3B-C4B	6.40	131.26	121.37
4	S	1084	CYC	OB-C4B-C3B	-6.39	121.31	128.03
4	K	1084	CYC	C1B-NB-C4B	-6.33	102.89	110.66
4	P	1082	CYC	OB-C4B-C3B	-6.33	121.37	128.03
4	H	1153	CYC	CHD-C1D-ND	-6.33	110.98	125.29
4	X	1082	CYC	C1B-NB-C4B	-6.33	102.90	110.66
4	R	1082	CYC	CMC-C2C-C1C	-6.29	98.84	112.40
4	W	1084	CYC	C1B-NB-C4B	-6.28	102.95	110.66
4	N	1153	CYC	CMA-C3A-C4A	6.26	134.82	125.10
4	B	1153	CYC	CMA-C3A-C4A	6.26	134.82	125.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	N	1153	CYC	CMD-C2D-C1D	6.25	134.70	125.62
4	L	1082	CYC	OB-C4B-C3B	-6.24	121.47	128.03
4	C	1084	CYC	CAD-C3D-C4D	-6.23	115.88	125.77
4	E	1084	CYC	C3D-C4D-ND	6.23	115.38	107.57
4	T	1082	CYC	CHB-C4A-NA	-6.23	111.50	124.95
4	D	1082	CYC	CAA-CBA-CGA	-6.21	97.19	113.67
4	B	1153	CYC	C3D-C4D-ND	6.19	115.33	107.57
4	D	1082	CYC	CHB-C1B-NB	-6.17	112.91	126.06
4	S	1084	CYC	CAC-C3C-C4C	6.14	128.43	112.67
4	D	1082	CYC	C2C-C1C-NC	6.13	113.39	108.29
4	G	1084	CYC	CAA-CBA-CGA	-6.12	97.42	113.67
4	J	1153	CYC	C2C-C1C-NC	6.10	113.37	108.29
4	A	1084	CYC	CAB-C3B-C4B	6.10	130.81	121.37
4	U	1084	CYC	CAA-CBA-CGA	-6.09	97.52	113.67
4	D	1153	CYC	C3D-C4D-ND	6.08	115.19	107.57
4	G	1084	CYC	C3D-C4D-ND	6.07	115.18	107.57
4	C	1084	CYC	C2B-C1B-NB	6.06	115.81	106.97
4	W	1084	CYC	OC-C1C-C2C	-6.05	121.36	126.17
4	T	1082	CYC	C2A-C1A-NA	6.01	118.55	110.04
4	N	1082	CYC	CAA-CBA-CGA	-5.99	97.78	113.67
4	B	1082	CYC	CAD-C3D-C2D	5.98	140.49	127.07
4	O	1084	CYC	C2C-C1C-NC	5.97	113.25	108.29
4	R	1153	CYC	CAB-C3B-C4B	5.95	130.57	121.37
4	F	1153	CYC	OB-C4B-C3B	-5.95	121.78	128.03
4	D	1082	CYC	OB-C4B-C3B	-5.93	121.80	128.03
4	L	1082	CYC	CMB-C2B-C1B	5.90	131.34	124.16
4	F	1082	CYC	C3D-C4D-ND	5.89	114.96	107.57
4	P	1153	CYC	C3D-C4D-ND	5.89	114.96	107.57
4	B	1153	CYC	CMD-C2D-C1D	5.89	134.18	125.62
4	G	1084	CYC	C1B-NB-C4B	-5.86	103.47	110.66
4	A	1084	CYC	OB-C4B-C3B	-5.86	121.87	128.03
4	P	1082	CYC	C2A-C1A-NA	5.84	118.31	110.04
4	T	1153	CYC	CAB-C3B-C4B	5.83	130.39	121.37
4	V	1082	CYC	CHB-C1B-NB	-5.83	113.62	126.06
4	A	1084	CYC	CAA-CBA-CGA	-5.82	98.24	113.67
4	F	1153	CYC	CAC-C3C-C4C	-5.81	97.77	112.67
4	F	1082	CYC	CAA-CBA-CGA	-5.81	98.27	113.67
4	N	1153	CYC	OC-C1C-C2C	-5.80	121.56	126.17
4	H	1082	CYC	CHB-C4A-NA	-5.76	112.52	124.95
4	N	1153	CYC	CHD-C4C-NC	-5.74	118.47	125.63
4	J	1082	CYC	CMB-C2B-C1B	5.73	131.13	124.16
4	V	1082	CYC	CMB-C2B-C1B	5.72	131.12	124.16

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	V	1082	CYC	C3D-C4D-ND	5.72	114.75	107.57
4	B	1082	CYC	CHA-C4D-C3D	-5.69	114.98	127.22
4	H	1153	CYC	CAC-C3C-C4C	-5.69	98.06	112.67
4	V	1153	CYC	CAB-C3B-C4B	5.68	130.15	121.37
4	T	1153	CYC	C4A-C3A-C2A	-5.67	100.05	106.48
4	K	1084	CYC	CHB-C1B-NB	-5.66	113.98	126.06
4	O	1084	CYC	C1B-NB-C4B	-5.66	103.71	110.66
4	G	1084	CYC	C2A-C1A-NA	5.65	118.05	110.04
4	H	1153	CYC	CMA-C3A-C4A	5.65	133.88	125.10
4	R	1082	CYC	C2C-C1C-NC	5.65	112.99	108.29
4	H	1153	CYC	CHB-C4A-NA	-5.64	112.78	124.95
4	J	1082	CYC	C2C-C1C-NC	5.62	112.97	108.29
4	F	1082	CYC	C2A-C1A-NA	5.62	118.00	110.04
4	D	1082	CYC	CMB-C2B-C1B	5.62	130.99	124.16
4	S	1084	CYC	C3D-C4D-ND	5.61	114.61	107.57
4	B	1082	CYC	CAD-C3D-C4D	-5.61	116.87	125.77
4	C	1084	CYC	CAD-C3D-C2D	5.60	139.63	127.07
4	L	1153	CYC	CHD-C1D-ND	-5.59	112.65	125.29
4	F	1082	CYC	CHD-C1D-ND	-5.58	112.68	125.29
4	N	1153	CYC	CHD-C1D-ND	-5.57	112.71	125.29
4	H	1153	CYC	CAB-C3B-C4B	5.55	129.96	121.37
4	V	1082	CYC	C2A-C1A-NA	5.53	117.88	110.04
4	V	1153	CYC	C3D-C4D-ND	5.53	114.50	107.57
4	N	1082	CYC	C2C-C1C-NC	5.52	112.88	108.29
4	J	1082	CYC	C1B-NB-C4B	-5.52	103.89	110.66
4	V	1153	CYC	C2C-C1C-NC	5.51	112.88	108.29
4	R	1153	CYC	CHB-C4A-NA	-5.48	113.12	124.95
4	J	1082	CYC	CHB-C1B-NB	-5.47	114.39	126.06
4	N	1082	CYC	CAD-C3D-C4D	5.46	134.44	125.77
4	T	1082	CYC	CMB-C2B-C1B	5.46	130.80	124.16
4	S	1084	CYC	C2A-C1A-NA	5.46	117.78	110.04
4	N	1153	CYC	C3D-C4D-ND	5.45	114.40	107.57
4	S	1084	CYC	C2C-C1C-NC	5.43	112.81	108.29
4	L	1082	CYC	CHB-C1B-NB	-5.43	114.48	126.06
4	R	1082	CYC	CMB-C2B-C1B	5.42	130.75	124.16
4	F	1082	CYC	OB-C4B-C3B	-5.40	122.36	128.03
4	V	1153	CYC	CHD-C1D-ND	-5.39	113.11	125.29
4	R	1153	CYC	CMD-C2D-C1D	5.39	133.45	125.62
4	B	1082	CYC	CAA-CBA-CGA	-5.37	99.41	113.67
4	L	1153	CYC	C2C-C3C-C4C	5.37	109.39	101.34
4	I	1084	CYC	C4D-C3D-C2D	-5.36	101.38	107.62
4	F	1082	CYC	CAD-C3D-C2D	5.33	139.03	127.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	N	1082	CYC	CHB-C1B-C2B	5.33	137.59	126.97
4	A	1084	CYC	C2C-C1C-NC	5.31	112.70	108.29
4	T	1082	CYC	C2B-C1B-NB	5.30	114.70	106.97
4	B	1153	CYC	CHD-C1D-ND	-5.29	113.33	125.29
4	W	1084	CYC	CHB-C1B-NB	-5.28	114.80	126.06
4	R	1153	CYC	CHD-C1D-ND	-5.28	113.36	125.29
4	B	1153	CYC	CBA-CAA-C2A	-5.26	97.98	112.53
4	N	1082	CYC	CMC-C2C-C1C	-5.24	101.10	112.40
4	R	1082	CYC	CAA-CBA-CGA	-5.24	99.76	113.67
4	Q	1084	CYC	C2C-C1C-NC	5.24	112.64	108.29
4	J	1153	CYC	CHD-C1D-ND	-5.21	113.51	125.29
4	T	1153	CYC	CHB-C4A-NA	-5.20	113.72	124.95
4	T	1082	CYC	CAD-C3D-C2D	5.20	138.73	127.07
4	T	1153	CYC	C2B-C1B-NB	5.20	114.54	106.97
4	D	1082	CYC	CBD-CAD-C3D	5.19	126.89	112.53
4	F	1082	CYC	CAD-C3D-C4D	-5.18	117.54	125.77
4	H	1082	CYC	CHB-C1B-NB	-5.17	115.02	126.06
4	L	1153	CYC	CHB-C4A-NA	-5.17	113.78	124.95
4	X	1082	CYC	C2C-C1C-NC	5.17	112.59	108.29
4	C	1084	CYC	OC-C1C-C2C	-5.16	122.07	126.17
4	P	1082	CYC	CHD-C1D-ND	-5.16	113.64	125.29
4	T	1082	CYC	CAD-C3D-C4D	-5.15	117.59	125.77
4	L	1082	CYC	C2A-C1A-NA	5.15	117.34	110.04
4	W	1084	CYC	C2A-C1A-NA	5.14	117.33	110.04
4	B	1082	CYC	CHD-C4C-NC	-5.14	119.21	125.63
4	A	1084	CYC	OC-C1C-C2C	-5.13	122.09	126.17
4	M	1084	CYC	C2C-C1C-NC	5.11	112.54	108.29
4	D	1082	CYC	CMA-C3A-C4A	5.10	133.03	125.10
4	I	1084	CYC	C4D-ND-C1D	-5.10	100.83	109.78
4	C	1084	CYC	CHA-C4D-C3D	-5.08	116.30	127.22
4	C	1084	CYC	CAB-C3B-C4B	5.08	129.22	121.37
4	N	1153	CYC	CHB-C4A-NA	-5.06	114.02	124.95
4	W	1084	CYC	C4D-C3D-C2D	-5.05	101.74	107.62
4	J	1082	CYC	C3D-C4D-ND	5.04	113.88	107.57
4	B	1153	CYC	CHB-C4A-NA	-5.03	114.08	124.95
4	H	1082	CYC	CMB-C2B-C1B	5.03	130.28	124.16
4	C	1084	CYC	C2A-C1A-NA	5.03	117.16	110.04
4	V	1082	CYC	CAB-C3B-C4B	5.03	129.15	121.37
4	R	1082	CYC	C2B-C1B-NB	5.02	114.29	106.97
4	D	1082	CYC	CHB-C4A-NA	-5.02	114.10	124.95
4	C	1084	CYC	OB-C4B-C3B	-5.01	122.77	128.03
4	L	1082	CYC	CMC-C2C-C1C	-4.99	101.65	112.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	G	1084	CYC	CAB-C3B-C4B	4.99	129.08	121.37
4	E	1084	CYC	C4D-C3D-C2D	-4.98	101.83	107.62
4	R	1082	CYC	OC-C1C-C2C	-4.97	122.22	126.17
4	X	1082	CYC	C2A-C1A-NA	4.96	117.07	110.04
4	S	1084	CYC	CAB-C3B-C4B	4.96	129.04	121.37
4	H	1153	CYC	OB-C4B-C3B	-4.95	122.83	128.03
4	T	1153	CYC	CMA-C3A-C4A	4.94	132.77	125.10
4	T	1153	CYC	CAA-CBA-CGA	4.94	126.75	113.67
4	T	1153	CYC	O2A-CGA-O1A	-4.91	110.69	123.33
4	T	1082	CYC	CHA-C4D-C3D	-4.91	116.67	127.22
4	X	1153	CYC	CHD-C1D-ND	-4.90	114.21	125.29
4	D	1153	CYC	C2D-C1D-ND	4.89	114.09	107.43
4	F	1153	CYC	CMA-C3A-C4A	4.89	132.69	125.10
4	C	1084	CYC	CHB-C4A-NA	-4.88	114.42	124.95
4	P	1153	CYC	CHD-C1D-ND	-4.86	114.32	125.29
4	O	1084	CYC	C3D-C4D-ND	4.85	113.66	107.57
4	S	1084	CYC	CHB-C4A-NA	-4.85	114.48	124.95
4	B	1082	CYC	C2C-C1C-NC	4.85	112.32	108.29
4	V	1153	CYC	C4A-C3A-C2A	-4.85	100.97	106.48
4	I	1084	CYC	C2D-C1D-ND	4.83	114.02	107.43
4	V	1153	CYC	C2B-C1B-NB	4.83	114.00	106.97
4	X	1082	CYC	CAA-CBA-CGA	-4.82	100.88	113.67
4	W	1084	CYC	CHB-C4A-NA	-4.82	114.55	124.95
4	L	1082	CYC	CMA-C3A-C4A	4.82	132.58	125.10
4	N	1082	CYC	C3D-C4D-ND	4.81	113.61	107.57
4	T	1153	CYC	C2A-C1A-NA	4.80	116.84	110.04
4	P	1153	CYC	CMD-C2D-C1D	4.79	132.59	125.62
4	O	1084	CYC	CAB-C3B-C4B	4.79	128.77	121.37
4	X	1153	CYC	C3D-C4D-ND	4.78	113.56	107.57
4	L	1082	CYC	CHB-C4A-NA	-4.78	114.64	124.95
4	D	1153	CYC	CAC-C3C-C4C	-4.78	100.41	112.67
4	J	1153	CYC	CAD-CBD-CGD	-4.76	101.03	113.67
4	A	1084	CYC	CMB-C2B-C1B	4.75	129.94	124.16
4	B	1082	CYC	CHB-C4A-NA	-4.75	114.70	124.95
4	H	1082	CYC	C3D-C4D-ND	4.74	113.51	107.57
4	E	1084	CYC	C2B-C1B-NB	4.74	113.88	106.97
4	D	1082	CYC	C3D-C4D-ND	4.72	113.49	107.57
4	C	1084	CYC	C2D-C1D-ND	4.72	113.87	107.43
4	L	1153	CYC	OC-C1C-NC	-4.71	119.37	124.93
4	N	1153	CYC	C2C-C1C-NC	4.71	112.21	108.29
4	Q	1084	CYC	C1B-NB-C4B	-4.71	104.88	110.66
4	B	1082	CYC	C1A-C2A-C3A	-4.71	101.63	106.73

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	U	1084	CYC	CAB-C3B-C4B	4.70	128.65	121.37
4	J	1082	CYC	C2A-C1A-NA	4.70	116.70	110.04
4	L	1082	CYC	C3D-C4D-ND	4.69	113.45	107.57
4	Q	1084	CYC	C3D-C4D-ND	4.68	113.44	107.57
4	V	1153	CYC	CMC-C2C-C1C	-4.68	102.32	112.40
4	F	1082	CYC	CHB-C1B-NB	-4.68	116.09	126.06
4	B	1082	CYC	CHD-C1D-ND	-4.67	114.74	125.29
4	K	1084	CYC	CAB-C3B-C4B	4.66	128.58	121.37
4	B	1153	CYC	C2D-C1D-ND	4.66	113.79	107.43
4	J	1082	CYC	OC-C1C-C2C	-4.66	122.47	126.17
4	K	1084	CYC	C3D-C4D-ND	4.63	113.38	107.57
4	I	1084	CYC	C2B-C1B-NB	4.63	113.72	106.97
4	H	1082	CYC	C2A-C1A-NA	4.63	116.60	110.04
4	N	1082	CYC	CMA-C3A-C4A	4.63	132.28	125.10
4	T	1153	CYC	CHD-C1D-ND	-4.62	114.85	125.29
4	N	1153	CYC	C4A-C3A-C2A	-4.61	101.25	106.48
4	R	1082	CYC	CAD-C3D-C4D	4.57	133.02	125.77
4	V	1082	CYC	CHB-C4A-NA	-4.56	115.10	124.95
4	B	1153	CYC	CAB-C3B-C4B	4.55	128.40	121.37
4	D	1153	CYC	CMA-C3A-C4A	4.54	132.16	125.10
4	S	1084	CYC	CAA-C2A-C3A	4.54	136.38	127.87
4	A	1084	CYC	CAD-C3D-C4D	4.54	132.97	125.77
4	H	1153	CYC	C2D-C1D-ND	4.53	113.61	107.43
4	R	1153	CYC	CMA-C3A-C4A	4.53	132.13	125.10
4	T	1082	CYC	CHB-C1B-NB	-4.53	116.41	126.06
4	T	1153	CYC	C2D-C1D-ND	4.52	113.60	107.43
4	V	1082	CYC	C2B-C1B-NB	4.52	113.56	106.97
4	F	1082	CYC	CMB-C2B-C1B	4.52	129.65	124.16
4	T	1082	CYC	CAA-C2A-C3A	4.51	136.31	127.87
4	E	1084	CYC	OB-C4B-NB	-4.51	114.51	125.08
4	F	1153	CYC	C2C-C3C-C4C	4.51	108.09	101.34
4	P	1082	CYC	CAD-C3D-C2D	4.50	137.17	127.07
4	T	1082	CYC	CHD-C1D-ND	-4.50	115.13	125.29
4	E	1084	CYC	C2A-C1A-NA	4.50	116.41	110.04
4	F	1153	CYC	C4D-C3D-C2D	-4.49	102.39	107.62
4	R	1153	CYC	CAA-CBA-CGA	4.49	125.57	113.67
4	S	1084	CYC	CHA-C4D-C3D	-4.48	117.58	127.22
4	F	1082	CYC	CHB-C4A-NA	-4.48	115.27	124.95
4	V	1153	CYC	CBA-CAA-C2A	-4.48	100.15	112.53
4	O	1084	CYC	C2A-C1A-NA	4.47	116.38	110.04
4	W	1084	CYC	C2B-C1B-NB	4.47	113.48	106.97
4	H	1153	CYC	C4A-C3A-C2A	-4.46	101.41	106.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1084	CYC	CHB-C4A-NA	-4.46	115.32	124.95
4	T	1082	CYC	C4D-ND-C1D	-4.46	101.95	109.78
4	X	1082	CYC	CHB-C4A-NA	-4.46	115.33	124.95
4	R	1153	CYC	CMC-C2C-C1C	-4.45	102.80	112.40
4	X	1153	CYC	C1B-NB-C4B	-4.44	105.21	110.66
4	H	1082	CYC	C2B-C1B-NB	4.44	113.44	106.97
4	P	1082	CYC	CAB-C3B-C4B	4.43	128.23	121.37
4	F	1082	CYC	CHA-C4D-C3D	-4.43	117.69	127.22
4	X	1082	CYC	CMC-C2C-C1C	-4.43	102.85	112.40
4	V	1082	CYC	CAA-CBA-CGA	-4.43	101.92	113.67
4	B	1153	CYC	C4A-C3A-C2A	-4.43	101.45	106.48
4	F	1153	CYC	CHB-C4A-NA	-4.42	115.40	124.95
4	L	1153	CYC	OC-C1C-C2C	-4.40	122.68	126.17
4	G	1084	CYC	OC-C1C-C2C	-4.39	122.68	126.17
4	E	1084	CYC	CHB-C1B-NB	-4.39	116.70	126.06
4	Q	1084	CYC	C2A-C1A-NA	4.38	116.25	110.04
4	T	1153	CYC	C4D-ND-C1D	-4.38	102.09	109.78
4	C	1084	CYC	CHD-C1D-ND	-4.38	115.40	125.29
4	O	1084	CYC	C4D-C3D-C2D	-4.38	102.53	107.62
4	M	1084	CYC	CAC-C3C-C4C	4.37	123.90	112.67
4	B	1082	CYC	C4D-C3D-C2D	-4.35	102.55	107.62
4	U	1084	CYC	CHB-C4A-NA	-4.35	115.55	124.95
4	H	1153	CYC	C2A-C1A-NA	4.35	116.20	110.04
4	F	1082	CYC	C2B-C1B-NB	4.35	113.31	106.97
4	U	1084	CYC	C2C-C1C-NC	4.34	111.90	108.29
4	I	1084	CYC	CHB-C4A-NA	-4.31	115.65	124.95
4	U	1084	CYC	C2A-C1A-NA	4.31	116.14	110.04
4	H	1153	CYC	C2B-C1B-NB	4.30	113.24	106.97
4	X	1082	CYC	C3D-C4D-ND	4.30	112.96	107.57
4	R	1153	CYC	C2B-C1B-NB	4.29	113.22	106.97
4	F	1082	CYC	CAB-C3B-C4B	4.29	128.00	121.37
4	A	1084	CYC	C2B-C1B-NB	4.28	113.22	106.97
4	B	1153	CYC	C2B-C1B-NB	4.28	113.21	106.97
4	L	1153	CYC	C3D-C4D-ND	4.27	112.93	107.57
4	G	1084	CYC	CHB-C4A-NA	-4.27	115.72	124.95
4	F	1153	CYC	CHD-C1D-ND	-4.25	115.68	125.29
4	J	1082	CYC	C2D-C1D-ND	4.25	113.23	107.43
4	R	1153	CYC	C2D-C1D-ND	4.25	113.23	107.43
4	R	1153	CYC	C4A-C3A-C2A	-4.24	101.66	106.48
4	T	1082	CYC	CAB-C3B-C4B	4.23	127.92	121.37
4	L	1153	CYC	C2B-C1B-NB	4.23	113.13	106.97
4	V	1153	CYC	CHD-C4C-NC	-4.23	120.35	125.63

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	V	1153	CYC	CBC-CAC-C3C	4.22	122.37	113.41
4	H	1082	CYC	CHB-C4A-C3A	4.21	135.70	124.87
4	C	1084	CYC	C1D-CHD-C4C	4.21	134.96	127.76
4	U	1084	CYC	C2B-C1B-NB	4.21	113.10	106.97
4	R	1153	CYC	CAC-C3C-C4C	-4.20	101.88	112.67
4	E	1084	CYC	CMB-C2B-C1B	4.20	129.27	124.16
4	G	1084	CYC	CAC-C3C-C4C	4.20	123.45	112.67
4	C	1084	CYC	C4D-ND-C1D	-4.19	102.42	109.78
4	S	1084	CYC	O1A-CGA-CBA	-4.19	109.79	123.09
4	F	1153	CYC	C2A-C1A-NA	4.18	115.96	110.04
4	M	1084	CYC	C3D-C4D-ND	4.17	112.80	107.57
4	E	1084	CYC	CAA-C2A-C3A	4.16	135.66	127.87
4	H	1082	CYC	CHA-C1A-NA	-4.16	115.86	124.60
4	D	1082	CYC	C2A-C1A-NA	4.15	115.92	110.04
4	T	1153	CYC	CMC-C2C-C1C	-4.15	103.46	112.40
4	V	1153	CYC	CHB-C4A-NA	-4.14	116.00	124.95
4	M	1084	CYC	CBB-CAB-C3B	-4.14	101.21	112.42
4	D	1082	CYC	CMC-C2C-C1C	-4.13	103.50	112.40
4	P	1153	CYC	C2D-C1D-ND	4.10	113.03	107.43
4	D	1153	CYC	C4D-C3D-C2D	-4.08	102.88	107.62
4	T	1082	CYC	CHB-C4A-C3A	4.07	135.34	124.87
4	L	1082	CYC	CHB-C4A-C3A	4.07	135.34	124.87
4	V	1153	CYC	C4D-C3D-C2D	-4.07	102.88	107.62
4	F	1153	CYC	CMC-C2C-C1C	-4.07	103.63	112.40
4	F	1082	CYC	CBD-CAD-C3D	4.07	123.79	112.53
4	L	1153	CYC	C1D-C2D-C3D	-4.06	102.68	107.76
4	J	1153	CYC	C2D-C1D-ND	4.04	112.94	107.43
4	A	1084	CYC	CAC-C3C-C4C	4.04	123.05	112.67
4	D	1082	CYC	CAD-C3D-C4D	4.04	132.18	125.77
4	F	1153	CYC	C2B-C1B-NB	4.03	112.85	106.97
4	K	1084	CYC	CHB-C4A-NA	-4.03	116.24	124.95
4	J	1153	CYC	C4D-C3D-C2D	-4.03	102.93	107.62
4	J	1082	CYC	CHB-C1B-C2B	4.03	135.00	126.97
4	X	1153	CYC	C2D-C1D-ND	4.01	112.90	107.43
4	T	1082	CYC	C4A-C3A-C2A	-4.00	101.93	106.48
4	B	1153	CYC	C4D-ND-C1D	-4.00	102.76	109.78
4	P	1082	CYC	CAD-C3D-C4D	-4.00	119.42	125.77
4	D	1082	CYC	CHD-C1D-ND	-3.98	116.30	125.29
4	D	1082	CYC	CAB-C3B-C4B	3.96	127.49	121.37
4	N	1153	CYC	C2D-C1D-ND	3.95	112.82	107.43
4	F	1153	CYC	CAB-C3B-C4B	3.95	127.48	121.37
4	D	1082	CYC	C2B-C1B-NB	3.94	112.71	106.97

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	J	1082	CYC	CBA-CAA-C2A	3.93	123.41	112.53
4	I	1084	CYC	CHA-C4D-C3D	-3.93	118.77	127.22
4	R	1153	CYC	CHB-C4A-C3A	3.93	134.96	124.87
4	T	1082	CYC	C2D-C1D-ND	3.93	112.79	107.43
4	W	1084	CYC	CAC-C3C-C4C	3.92	122.75	112.67
4	K	1084	CYC	CMB-C2B-C1B	3.90	128.91	124.16
4	W	1084	CYC	CHA-C4D-C3D	-3.90	118.83	127.22
4	E	1084	CYC	O1D-CGD-CBD	-3.89	110.77	123.09
4	L	1153	CYC	C2D-C1D-ND	3.88	112.73	107.43
4	P	1082	CYC	CHA-C4D-C3D	-3.88	118.87	127.22
4	C	1084	CYC	C4A-C3A-C2A	-3.88	102.08	106.48
4	L	1082	CYC	CHD-C4C-NC	-3.87	120.80	125.63
4	G	1084	CYC	C2B-C1B-NB	3.86	112.60	106.97
4	V	1082	CYC	OC-C1C-C2C	-3.86	123.10	126.17
4	W	1084	CYC	O1A-CGA-CBA	-3.86	110.84	123.09
4	K	1084	CYC	C2B-C1B-NB	3.86	112.60	106.97
4	P	1082	CYC	CHB-C1B-NB	-3.86	117.83	126.06
4	L	1153	CYC	C2A-C1A-NA	3.86	115.51	110.04
4	X	1082	CYC	CHA-C1A-NA	-3.86	116.50	124.60
4	G	1084	CYC	C2C-C1C-NC	3.85	111.49	108.29
4	R	1082	CYC	CHB-C1B-C2B	3.85	134.65	126.97
4	S	1084	CYC	C2C-C3C-C4C	3.85	107.10	101.34
4	U	1084	CYC	CAC-C3C-C4C	3.85	122.55	112.67
4	N	1082	CYC	C2A-C1A-NA	3.84	115.47	110.04
4	P	1082	CYC	C2D-C1D-ND	3.84	112.66	107.43
4	L	1153	CYC	CMD-C2D-C1D	3.83	131.19	125.62
4	P	1082	CYC	C4D-ND-C1D	-3.81	103.09	109.78
4	X	1153	CYC	C4D-C3D-C2D	-3.80	103.19	107.62
4	I	1084	CYC	CAC-C3C-C4C	3.79	122.40	112.67
4	T	1082	CYC	CHD-C4C-NC	-3.78	120.91	125.63
4	R	1153	CYC	C2A-C1A-NA	3.78	115.40	110.04
4	F	1153	CYC	C4D-ND-C1D	-3.78	103.14	109.78
4	G	1084	CYC	O1A-CGA-CBA	-3.77	111.12	123.09
4	I	1084	CYC	CHB-C1B-NB	-3.77	118.01	126.06
4	S	1084	CYC	O1D-CGD-CBD	-3.77	111.14	123.09
4	D	1082	CYC	CHB-C4A-C3A	3.76	134.54	124.87
4	P	1082	CYC	CAD-CBD-CGD	-3.76	103.69	113.67
4	A	1084	CYC	CAD-CBD-CGD	-3.76	103.69	113.67
4	V	1082	CYC	OB-C4B-C3B	-3.76	124.08	128.03
4	P	1153	CYC	C4D-ND-C1D	-3.76	103.19	109.78
4	Q	1084	CYC	CAB-C3B-C4B	3.75	127.17	121.37
4	S	1084	CYC	CHB-C4A-C3A	3.75	134.50	124.87

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	T	1082	CYC	CAA-C2A-C1A	-3.75	118.44	125.02
4	I	1084	CYC	C1A-NA-C4A	-3.74	99.64	106.52
4	L	1082	CYC	C2B-C1B-NB	3.74	112.43	106.97
4	P	1082	CYC	C2C-C1C-NC	3.73	111.39	108.29
4	H	1153	CYC	CHB-C4A-C3A	3.73	134.44	124.87
4	I	1084	CYC	CHD-C1D-ND	-3.73	116.87	125.29
4	T	1153	CYC	CMD-C2D-C1D	3.72	131.03	125.62
4	D	1082	CYC	CHB-C1B-C2B	3.72	134.38	126.97
4	P	1153	CYC	C2B-C1B-NB	3.71	112.38	106.97
4	R	1153	CYC	C4D-ND-C1D	-3.71	103.27	109.78
4	N	1082	CYC	C2D-C1D-ND	3.70	112.47	107.43
4	Q	1084	CYC	OB-C4B-C3B	-3.69	124.15	128.03
4	F	1153	CYC	C2D-C1D-ND	3.69	112.46	107.43
4	F	1153	CYC	C4A-C3A-C2A	-3.68	102.30	106.48
4	R	1153	CYC	C1D-C2D-C3D	-3.67	103.16	107.76
4	X	1153	CYC	CMC-C2C-C1C	-3.67	104.49	112.40
4	J	1082	CYC	CHB-C4A-NA	-3.67	117.03	124.95
4	H	1153	CYC	C4D-ND-C1D	-3.67	103.35	109.78
4	J	1153	CYC	C4D-ND-C1D	-3.66	103.36	109.78
4	J	1153	CYC	C2B-C1B-NB	3.66	112.30	106.97
4	R	1082	CYC	CAA-C2A-C1A	3.66	131.44	125.02
4	F	1082	CYC	CBA-CAA-C2A	3.65	122.63	112.53
4	P	1082	CYC	C4D-C3D-C2D	-3.65	103.38	107.62
4	S	1084	CYC	CAA-C2A-C1A	-3.64	118.62	125.02
4	T	1153	CYC	C4D-C3D-C2D	-3.64	103.38	107.62
4	N	1082	CYC	CAA-C2A-C1A	3.64	131.41	125.02
4	S	1084	CYC	CHD-C1D-ND	-3.64	117.07	125.29
4	C	1084	CYC	C1D-C2D-C3D	-3.64	103.21	107.76
4	J	1082	CYC	CHB-C4A-C3A	3.63	134.21	124.87
4	W	1084	CYC	C4A-C3A-C2A	-3.63	102.35	106.48
4	N	1153	CYC	C2A-C1A-NA	3.62	115.18	110.04
4	L	1153	CYC	CAA-CBA-CGA	3.62	123.26	113.67
4	F	1082	CYC	C4D-C3D-C2D	-3.62	103.41	107.62
4	K	1084	CYC	C4D-C3D-C2D	-3.61	103.42	107.62
4	U	1084	CYC	CHD-C4C-NC	-3.61	121.12	125.63
4	X	1082	CYC	CHB-C4A-C3A	3.61	134.14	124.87
4	H	1082	CYC	CAA-CBA-CGA	-3.61	104.10	113.67
4	A	1084	CYC	C2A-C1A-NA	3.60	115.15	110.04
4	H	1082	CYC	C2C-C1C-NC	3.60	111.28	108.29
4	S	1084	CYC	O2A-CGA-O1A	3.60	132.59	123.33
4	D	1153	CYC	C4D-ND-C1D	-3.60	103.46	109.78
4	T	1153	CYC	O2A-CGA-CBA	3.60	125.37	114.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	J	1153	CYC	C2A-C1A-NA	3.60	115.14	110.04
4	A	1084	CYC	CHA-C1A-NA	-3.60	117.05	124.60
4	N	1153	CYC	CHB-C4A-C3A	3.60	134.11	124.87
4	O	1084	CYC	CAC-C3C-C4C	3.59	121.90	112.67
4	W	1084	CYC	CAB-C3B-C4B	3.59	126.92	121.37
4	J	1082	CYC	C4D-ND-C1D	-3.59	103.48	109.78
4	M	1084	CYC	CMB-C2B-C1B	3.58	128.51	124.16
4	J	1153	CYC	CHA-C1A-C2A	-3.58	117.14	125.40
4	K	1084	CYC	CHB-C4A-C3A	3.58	134.06	124.87
4	D	1153	CYC	C4A-C3A-C2A	-3.58	102.42	106.48
4	N	1082	CYC	CHA-C1A-NA	-3.58	117.09	124.60
4	E	1084	CYC	CHD-C1D-ND	-3.57	117.23	125.29
4	V	1082	CYC	C2D-C1D-ND	3.56	112.29	107.43
4	N	1153	CYC	CAB-C3B-C4B	3.56	126.88	121.37
4	K	1084	CYC	C2D-C1D-ND	3.56	112.29	107.43
4	R	1082	CYC	C2D-C1D-ND	3.56	112.28	107.43
4	E	1084	CYC	CHA-C4D-C3D	-3.56	119.57	127.22
4	D	1082	CYC	CMD-C2D-C1D	3.54	130.77	125.62
4	A	1084	CYC	C3D-C4D-ND	3.54	112.01	107.57
4	A	1084	CYC	O2A-CGA-O1A	3.54	132.44	123.33
4	Q	1084	CYC	CAD-C3D-C4D	3.54	131.39	125.77
4	A	1084	CYC	CHB-C1B-NB	-3.53	118.53	126.06
4	U	1084	CYC	CHB-C1B-NB	-3.53	118.53	126.06
4	G	1084	CYC	C4D-ND-C1D	-3.53	103.58	109.78
4	J	1153	CYC	C2C-C3C-C4C	3.53	106.62	101.34
4	M	1084	CYC	CAB-C3B-C4B	3.53	126.83	121.37
4	L	1153	CYC	CAB-C3B-C4B	3.52	126.82	121.37
4	S	1084	CYC	CHB-C1B-NB	-3.52	118.55	126.06
4	V	1082	CYC	C4D-C3D-C2D	-3.52	103.53	107.62
4	D	1082	CYC	C4D-C3D-C2D	-3.52	103.53	107.62
4	B	1082	CYC	C2B-C1B-NB	3.51	112.09	106.97
4	F	1153	CYC	CBA-CAA-C2A	-3.51	102.83	112.53
4	P	1082	CYC	CBD-CAD-C3D	3.50	122.20	112.53
4	J	1153	CYC	C4A-C3A-C2A	-3.49	102.52	106.48
4	B	1082	CYC	CBD-CAD-C3D	3.49	122.18	112.53
4	R	1082	CYC	C3D-C4D-ND	3.49	111.94	107.57
4	X	1082	CYC	CHB-C1B-NB	-3.49	118.62	126.06
4	J	1082	CYC	C4D-C3D-C2D	-3.49	103.56	107.62
4	K	1084	CYC	CHD-C1D-ND	-3.48	117.42	125.29
4	N	1082	CYC	C2B-C1B-NB	3.48	112.05	106.97
4	A	1084	CYC	C2C-C3C-C4C	3.47	106.54	101.34
4	V	1082	CYC	C4D-ND-C1D	-3.47	103.70	109.78

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	E	1084	CYC	CHB-C4A-NA	-3.46	117.48	124.95
4	R	1082	CYC	CAB-C3B-C2B	3.46	133.91	127.56
4	D	1153	CYC	C2C-C3C-C4C	3.46	106.52	101.34
4	R	1082	CYC	CBB-CAB-C3B	-3.45	103.07	112.42
4	P	1082	CYC	CHB-C4A-NA	-3.45	117.51	124.95
4	O	1084	CYC	C2D-C1D-ND	3.45	112.13	107.43
4	K	1084	CYC	CBB-CAB-C3B	-3.45	103.08	112.42
4	D	1082	CYC	CAD-CBD-CGD	-3.45	104.53	113.67
4	H	1153	CYC	CMD-C2D-C1D	3.44	130.63	125.62
4	N	1153	CYC	C4D-ND-C1D	-3.44	103.75	109.78
4	O	1084	CYC	CHD-C4C-NC	-3.44	121.34	125.63
4	T	1082	CYC	C4D-C3D-C2D	-3.44	103.62	107.62
4	M	1084	CYC	O1A-CGA-CBA	-3.43	112.21	123.09
4	D	1153	CYC	C2B-C1B-NB	3.43	111.97	106.97
4	W	1084	CYC	CBD-CAD-C3D	3.42	122.00	112.53
4	D	1153	CYC	CHA-C4D-ND	-3.42	117.56	125.29
4	I	1084	CYC	CHD-C4C-NC	-3.42	121.36	125.63
4	X	1153	CYC	CMA-C3A-C4A	3.42	130.41	125.10
4	J	1153	CYC	CMD-C2D-C1D	3.41	130.58	125.62
4	H	1153	CYC	C1D-C2D-C3D	-3.40	103.50	107.76
4	J	1082	CYC	C1A-C2A-C3A	-3.40	103.05	106.73
4	B	1082	CYC	C4D-ND-C1D	-3.40	103.81	109.78
4	X	1082	CYC	CHD-C4C-NC	-3.39	121.40	125.63
4	B	1082	CYC	CHB-C4A-C3A	3.39	133.58	124.87
4	G	1084	CYC	CHB-C1B-NB	-3.39	118.83	126.06
4	T	1153	CYC	C1A-NA-C4A	-3.39	100.30	106.52
4	E	1084	CYC	CAA-C2A-C1A	-3.38	119.08	125.02
4	Q	1084	CYC	C3C-C4C-NC	3.37	112.29	107.94
4	L	1153	CYC	C4A-C3A-C2A	-3.37	102.65	106.48
4	N	1153	CYC	C2B-C1B-NB	3.37	111.88	106.97
4	V	1153	CYC	CMA-C3A-C4A	3.37	130.33	125.10
4	N	1082	CYC	C4D-ND-C1D	-3.37	103.87	109.78
4	N	1082	CYC	CHB-C4A-NA	-3.36	117.69	124.95
4	K	1084	CYC	OB-C4B-NB	-3.36	117.20	125.08
4	X	1153	CYC	C2C-C3C-C4C	3.36	106.37	101.34
4	X	1082	CYC	CAB-C3B-C2B	3.36	133.72	127.56
4	P	1153	CYC	C2A-C1A-NA	3.35	114.79	110.04
4	Q	1084	CYC	C2D-C1D-ND	3.35	111.99	107.43
4	B	1153	CYC	C2C-C3C-C4C	3.34	106.35	101.34
4	X	1153	CYC	CAD-C3D-C4D	3.34	131.06	125.77
4	T	1082	CYC	CMC-C2C-C1C	-3.33	105.22	112.40
4	V	1082	CYC	C2C-C1C-NC	3.33	111.06	108.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	W	1084	CYC	CMA-C3A-C4A	3.32	130.26	125.10
4	T	1082	CYC	CBB-CAB-C3B	-3.32	103.42	112.42
4	W	1084	CYC	C4D-ND-C1D	-3.31	103.98	109.78
4	J	1082	CYC	CAB-C3B-C2B	3.30	133.63	127.56
4	R	1082	CYC	CHA-C1A-NA	-3.30	117.66	124.60
4	V	1082	CYC	C1A-C2A-C3A	-3.30	103.15	106.73
4	I	1084	CYC	C4A-C3A-C2A	-3.30	102.73	106.48
4	O	1084	CYC	C2B-C1B-NB	3.30	111.78	106.97
4	T	1082	CYC	CMA-C3A-C4A	3.30	130.22	125.10
4	G	1084	CYC	CAA-C2A-C3A	3.30	134.04	127.87
4	D	1153	CYC	CHD-C1D-C2D	3.30	135.61	127.53
4	J	1153	CYC	CHD-C4C-NC	3.30	129.75	125.63
4	X	1082	CYC	C4D-C3D-C2D	-3.29	103.79	107.62
4	V	1082	CYC	CHB-C4A-C3A	3.29	133.32	124.87
4	X	1082	CYC	CMA-C3A-C4A	3.29	130.21	125.10
4	X	1082	CYC	CHD-C1D-ND	-3.29	117.86	125.29
4	P	1153	CYC	C1D-C2D-C3D	-3.28	103.66	107.76
4	L	1153	CYC	CMA-C3A-C4A	3.27	130.19	125.10
4	P	1153	CYC	C4A-C3A-C2A	-3.27	102.76	106.48
4	X	1153	CYC	CHA-C4D-ND	-3.27	117.90	125.29
4	V	1153	CYC	C2A-C1A-NA	3.26	114.66	110.04
4	D	1153	CYC	C4D-CHA-C1A	3.26	135.34	128.22
4	W	1084	CYC	CHB-C4A-C3A	3.26	133.25	124.87
4	M	1084	CYC	C2B-C1B-NB	3.26	111.72	106.97
4	E	1084	CYC	C2D-C1D-ND	3.25	111.87	107.43
4	N	1082	CYC	CBD-CAD-C3D	3.25	121.52	112.53
4	F	1153	CYC	CHB-C1B-NB	-3.25	119.13	126.06
4	V	1153	CYC	C2D-C1D-ND	3.25	111.86	107.43
4	E	1084	CYC	C4D-ND-C1D	-3.25	104.08	109.78
4	E	1084	CYC	O1A-CGA-CBA	-3.24	112.81	123.09
4	H	1153	CYC	C4D-C3D-C2D	-3.24	103.85	107.62
4	F	1082	CYC	CHD-C1D-C2D	3.24	135.47	127.53
4	B	1153	CYC	CHB-C4A-C3A	3.24	133.19	124.87
4	V	1082	CYC	OB-C4B-NB	-3.24	117.49	125.08
4	A	1084	CYC	C4A-C3A-C2A	-3.23	102.81	106.48
4	D	1153	CYC	CBC-CAC-C3C	3.23	120.27	113.41
4	P	1153	CYC	CHB-C4A-NA	-3.23	117.97	124.95
4	H	1082	CYC	CAD-C3D-C4D	3.23	130.90	125.77
4	U	1084	CYC	C4A-C3A-C2A	-3.23	102.81	106.48
4	Q	1084	CYC	CAD-CBD-CGD	-3.21	105.14	113.67
4	J	1153	CYC	CAD-C3D-C4D	3.21	130.87	125.77
4	H	1153	CYC	CHD-C1D-C2D	3.21	135.40	127.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	G	1084	CYC	CMD-C2D-C1D	3.21	130.29	125.62
4	G	1084	CYC	C1A-C2A-C3A	-3.21	103.25	106.73
4	X	1153	CYC	CAC-C3C-C4C	-3.21	104.43	112.67
4	K	1084	CYC	CHB-C1B-C2B	3.21	133.37	126.97
4	S	1084	CYC	C4D-ND-C1D	-3.21	104.15	109.78
4	T	1082	CYC	C1B-C2B-C3B	-3.20	104.57	107.86
4	H	1153	CYC	CHD-C4C-NC	-3.20	121.63	125.63
4	B	1082	CYC	C2D-C1D-ND	3.20	111.79	107.43
4	Q	1084	CYC	C4D-C3D-C2D	-3.19	103.90	107.62
4	C	1084	CYC	C1B-C2B-C3B	-3.19	104.58	107.86
4	L	1082	CYC	C2D-C1D-ND	3.19	111.78	107.43
4	F	1082	CYC	C2D-C1D-ND	3.18	111.77	107.43
4	L	1082	CYC	C4D-ND-C1D	-3.17	104.21	109.78
4	N	1153	CYC	C4D-C3D-C2D	-3.17	103.93	107.62
4	W	1084	CYC	C2C-C1C-NC	3.16	110.91	108.29
4	R	1153	CYC	CBA-CAA-C2A	-3.16	103.81	112.53
4	G	1084	CYC	CBA-CAA-C2A	3.15	121.26	112.53
4	G	1084	CYC	CHD-C4C-NC	-3.15	121.69	125.63
4	Q	1084	CYC	CHB-C1B-NB	-3.15	119.34	126.06
4	J	1153	CYC	OC-C1C-NC	3.15	128.64	124.93
4	P	1153	CYC	CMA-C3A-C4A	3.14	129.98	125.10
4	Q	1084	CYC	CHB-C4A-NA	-3.14	118.17	124.95
4	V	1082	CYC	CHA-C1A-NA	-3.14	118.01	124.60
4	L	1082	CYC	CMD-C2D-C1D	3.13	130.17	125.62
4	B	1153	CYC	C4D-C3D-C2D	-3.12	103.98	107.62
4	C	1084	CYC	CHB-C1B-NB	-3.12	119.41	126.06
4	P	1082	CYC	C1A-C2A-C3A	-3.12	103.36	106.73
4	N	1153	CYC	CBA-CAA-C2A	-3.11	103.94	112.53
4	N	1082	CYC	CHB-C4A-C3A	3.10	132.84	124.87
4	M	1084	CYC	CHB-C1B-NB	-3.10	119.44	126.06
4	M	1084	CYC	CHA-C1A-C2A	-3.10	118.24	125.40
4	P	1082	CYC	CAA-C2A-C3A	3.09	133.66	127.87
4	C	1084	CYC	O1A-CGA-CBA	-3.09	113.29	123.09
4	B	1082	CYC	CAB-C3B-C4B	3.09	126.15	121.37
4	S	1084	CYC	CAD-CBD-CGD	-3.08	105.49	113.67
4	O	1084	CYC	CHB-C1B-NB	-3.08	119.49	126.06
4	V	1153	CYC	CAD-C3D-C2D	3.08	133.98	127.07
4	M	1084	CYC	C4D-C3D-C2D	-3.08	104.04	107.62
4	L	1153	CYC	O2A-CGA-O1A	-3.07	115.43	123.33
4	L	1082	CYC	CHB-C1B-C2B	3.07	133.09	126.97
4	M	1084	CYC	C4A-C3A-C2A	-3.07	102.99	106.48
4	N	1153	CYC	CHA-C1A-C2A	-3.07	118.31	125.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	G	1084	CYC	CHD-C1D-ND	-3.06	118.37	125.29
4	V	1153	CYC	CHD-C1D-C2D	3.06	135.03	127.53
4	U	1084	CYC	OB-C4B-NB	-3.06	117.91	125.08
4	R	1082	CYC	CBD-CAD-C3D	3.06	120.98	112.53
4	T	1082	CYC	CAD-CBD-CGD	-3.05	105.58	113.67
4	B	1153	CYC	C2A-C1A-NA	3.05	114.36	110.04
4	F	1082	CYC	O1D-CGD-CBD	-3.04	113.45	123.09
4	F	1082	CYC	C4D-ND-C1D	-3.03	104.46	109.78
4	N	1082	CYC	C1B-CHB-C4A	-3.03	120.62	128.06
4	H	1153	CYC	O2A-CGA-O1A	-3.02	115.56	123.33
4	M	1084	CYC	CAA-C2A-C1A	-3.02	119.72	125.02
4	U	1084	CYC	C3D-C4D-ND	3.02	111.35	107.57
4	X	1082	CYC	CHA-C4D-C3D	-3.02	120.74	127.22
4	N	1082	CYC	C4D-C3D-C2D	-3.01	104.11	107.62
4	E	1084	CYC	CBD-CAD-C3D	3.01	120.85	112.53
4	J	1153	CYC	C1D-CHD-C4C	3.01	132.90	127.76
4	H	1153	CYC	C2C-C3C-C4C	3.00	105.84	101.34
4	F	1082	CYC	CBB-CAB-C3B	-3.00	104.29	112.42
4	P	1153	CYC	C2C-C3C-C4C	2.99	105.82	101.34
4	G	1084	CYC	O2A-CGA-CBA	2.99	123.45	114.00
4	L	1082	CYC	CHD-C1D-ND	-2.99	118.54	125.29
4	J	1082	CYC	CHD-C1D-ND	-2.99	118.54	125.29
4	K	1084	CYC	CBA-CAA-C2A	2.98	120.78	112.53
4	O	1084	CYC	OB-C4B-NB	-2.98	118.09	125.08
4	M	1084	CYC	C2D-C1D-ND	2.97	111.49	107.43
4	R	1153	CYC	O2A-CGA-CBA	2.97	123.37	114.00
4	T	1082	CYC	C2C-C1C-NC	2.96	110.75	108.29
4	A	1084	CYC	CHB-C4A-C3A	2.96	132.48	124.87
4	M	1084	CYC	CHD-C1D-ND	-2.96	118.60	125.29
4	R	1082	CYC	C1B-C2B-C3B	-2.96	104.82	107.86
4	Q	1084	CYC	CMC-C2C-C1C	2.96	118.78	112.40
4	U	1084	CYC	CHB-C4A-C3A	2.96	132.47	124.87
4	P	1082	CYC	C2B-C1B-NB	2.95	111.28	106.97
4	J	1153	CYC	CAB-C3B-C4B	2.95	125.94	121.37
4	L	1153	CYC	CBC-CAC-C3C	2.94	119.66	113.41
4	E	1084	CYC	O2D-CGD-CBD	2.94	123.29	114.00
4	T	1082	CYC	C1A-NA-C4A	-2.94	101.12	106.52
4	Q	1084	CYC	CAA-C2A-C3A	2.94	133.37	127.87
4	P	1153	CYC	CBD-CAD-C3D	-2.94	104.41	112.53
4	V	1082	CYC	CHB-C1B-C2B	2.94	132.82	126.97
4	H	1082	CYC	CMD-C2D-C1D	2.93	129.87	125.62
4	B	1153	CYC	C1D-C2D-C3D	-2.92	104.10	107.76

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	I	1084	CYC	CAB-C3B-C4B	2.92	125.89	121.37
4	Q	1084	CYC	OC-C1C-C2C	-2.92	123.85	126.17
4	C	1084	CYC	CMB-C2B-C1B	2.92	127.71	124.16
4	E	1084	CYC	CMC-C2C-C1C	2.92	118.69	112.40
4	J	1153	CYC	CAB-C3B-C2B	2.91	132.91	127.56
4	X	1082	CYC	C2B-C1B-NB	2.91	111.21	106.97
4	W	1084	CYC	CHA-C1A-NA	-2.91	118.49	124.60
4	D	1082	CYC	C2D-C1D-ND	2.91	111.40	107.43
4	Q	1084	CYC	C4D-ND-C1D	-2.91	104.67	109.78
4	V	1082	CYC	CHD-C1D-ND	-2.91	118.72	125.29
4	L	1153	CYC	CHD-C4C-NC	-2.90	122.01	125.63
4	S	1084	CYC	C4A-C3A-C2A	-2.90	103.19	106.48
4	L	1082	CYC	C4D-C3D-C2D	-2.90	104.25	107.62
4	T	1153	CYC	CHB-C1B-C2B	-2.90	121.19	126.97
4	L	1153	CYC	CHD-C1D-C2D	2.90	134.62	127.53
4	X	1153	CYC	C2A-C1A-NA	2.89	114.14	110.04
4	B	1153	CYC	CHA-C4D-C3D	-2.89	121.01	127.22
4	L	1153	CYC	CHB-C4A-C3A	2.88	132.28	124.87
4	V	1082	CYC	CBC-CAC-C3C	-2.88	107.30	113.41
4	M	1084	CYC	C2A-C1A-NA	2.88	114.12	110.04
4	J	1153	CYC	C4D-CHA-C1A	2.88	134.51	128.22
4	L	1082	CYC	CAD-CBD-CGD	-2.88	106.04	113.67
4	L	1082	CYC	CHA-C4D-C3D	-2.87	121.04	127.22
4	E	1084	CYC	C1D-CHD-C4C	2.87	132.67	127.76
4	K	1084	CYC	C4D-ND-C1D	-2.87	104.75	109.78
4	O	1084	CYC	O2A-CGA-O1A	2.87	130.70	123.33
4	D	1153	CYC	CHB-C1B-NB	-2.86	119.95	126.06
4	F	1082	CYC	CMA-C3A-C4A	2.86	129.54	125.10
4	H	1153	CYC	O2A-CGA-CBA	2.86	123.03	114.00
4	B	1082	CYC	CBA-CAA-C2A	2.86	120.44	112.53
4	P	1082	CYC	CMC-C2C-C1C	-2.86	106.25	112.40
4	H	1082	CYC	C4D-C3D-C2D	-2.85	104.30	107.62
4	H	1082	CYC	C4D-ND-C1D	-2.85	104.78	109.78
4	X	1153	CYC	CHB-C4A-NA	-2.84	118.81	124.95
4	Q	1084	CYC	CHD-C1D-ND	-2.84	118.88	125.29
4	G	1084	CYC	C2D-C1D-ND	2.83	111.30	107.43
4	B	1082	CYC	CHA-C1A-NA	-2.83	118.65	124.60
4	F	1153	CYC	CHB-C4A-C3A	2.83	132.15	124.87
4	I	1084	CYC	OC-C1C-NC	-2.83	121.59	124.93
4	N	1082	CYC	C2C-C3C-C4C	2.83	105.58	101.34
4	O	1084	CYC	O1A-CGA-CBA	-2.82	114.14	123.09
4	Q	1084	CYC	C2B-C1B-NB	2.82	111.08	106.97

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	T	1153	CYC	CHA-C1A-C2A	-2.82	118.88	125.40
4	T	1153	CYC	C1D-C2D-C3D	-2.82	104.23	107.76
4	C	1084	CYC	C2C-C3C-C4C	2.82	105.56	101.34
4	J	1082	CYC	CAA-CBA-CGA	-2.82	106.19	113.67
4	W	1084	CYC	OB-C4B-NB	-2.81	118.48	125.08
4	O	1084	CYC	C4A-C3A-C2A	-2.81	103.28	106.48
4	D	1153	CYC	CAA-CBA-CGA	2.81	121.13	113.67
4	N	1153	CYC	CHD-C1D-C2D	2.81	134.42	127.53
4	O	1084	CYC	CHD-C1D-ND	-2.81	118.94	125.29
4	A	1084	CYC	C4D-C3D-C2D	-2.81	104.35	107.62
4	T	1082	CYC	CBD-CAD-C3D	2.81	120.29	112.53
4	S	1084	CYC	C2B-C1B-NB	2.80	111.05	106.97
4	D	1082	CYC	C4D-ND-C1D	-2.80	104.86	109.78
4	L	1082	CYC	CAB-C3B-C4B	2.80	125.70	121.37
4	E	1084	CYC	C2C-C1C-NC	2.79	110.61	108.29
4	I	1084	CYC	C1A-C2A-C3A	-2.79	103.71	106.73
4	H	1082	CYC	C1A-C2A-C3A	-2.78	103.72	106.73
4	M	1084	CYC	CHB-C4A-NA	-2.78	118.94	124.95
4	H	1153	CYC	CAA-CBA-CGA	2.78	121.04	113.67
4	R	1153	CYC	C4D-C3D-C2D	-2.77	104.39	107.62
4	X	1153	CYC	CHB-C1B-NB	-2.77	120.15	126.06
4	P	1153	CYC	CHD-C4C-NC	-2.77	122.17	125.63
4	I	1084	CYC	CMA-C3A-C4A	2.76	129.39	125.10
4	K	1084	CYC	C2A-C1A-NA	2.76	113.95	110.04
4	C	1084	CYC	CMC-C2C-C1C	2.75	118.34	112.40
4	U	1084	CYC	CBC-CAC-C3C	-2.75	107.57	113.41
4	C	1084	CYC	C4D-C3D-C2D	-2.75	104.42	107.62
4	F	1082	CYC	C1A-NA-C4A	-2.75	101.47	106.52
4	S	1084	CYC	CMB-C2B-C1B	2.75	127.50	124.16
4	U	1084	CYC	CAD-C3D-C4D	2.74	130.12	125.77
4	H	1082	CYC	C2D-C1D-ND	2.74	111.16	107.43
4	F	1082	CYC	C4A-C3A-C2A	-2.73	103.37	106.48
4	H	1153	CYC	CAD-C3D-C2D	2.73	133.20	127.07
4	H	1082	CYC	CHA-C4D-C3D	-2.73	121.35	127.22
4	R	1082	CYC	C4D-C3D-C2D	-2.73	104.44	107.62
4	R	1082	CYC	CHB-C4A-NA	-2.73	119.06	124.95
4	D	1153	CYC	C1D-C2D-C3D	-2.73	104.35	107.76
4	M	1084	CYC	CAA-C2A-C3A	2.73	132.97	127.87
4	P	1082	CYC	CHA-C1A-C2A	-2.73	119.10	125.40
4	J	1153	CYC	CAC-C3C-C4C	-2.73	105.67	112.67
4	T	1082	CYC	C1D-C2D-C3D	-2.72	104.35	107.76
4	W	1084	CYC	C1B-CHB-C4A	-2.72	121.37	128.06

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	M	1084	CYC	OC-C1C-C2C	-2.72	124.01	126.17
4	O	1084	CYC	OC-C1C-C2C	-2.72	124.01	126.17
4	B	1153	CYC	CBC-CAC-C3C	2.71	119.16	113.41
4	B	1082	CYC	C1A-NA-C4A	-2.71	101.55	106.52
4	P	1153	CYC	CHA-C4D-C3D	-2.71	121.41	127.22
4	G	1084	CYC	CHB-C4A-C3A	2.70	131.82	124.87
4	S	1084	CYC	C1D-CHD-C4C	2.70	132.37	127.76
4	R	1153	CYC	CBB-CAB-C3B	2.69	119.72	112.42
4	G	1084	CYC	CHA-C4D-C3D	-2.69	121.44	127.22
4	L	1153	CYC	C1A-NA-C4A	-2.69	101.58	106.52
4	S	1084	CYC	CMD-C2D-C1D	2.69	129.52	125.62
4	I	1084	CYC	CBB-CAB-C3B	-2.69	105.14	112.42
4	V	1082	CYC	CAD-CBD-CGD	-2.68	106.55	113.67
4	R	1153	CYC	C1C-NC-C4C	-2.68	110.04	113.41
4	K	1084	CYC	CAD-C3D-C4D	2.67	130.01	125.77
4	D	1153	CYC	CMD-C2D-C1D	2.67	129.50	125.62
4	D	1082	CYC	CBA-CAA-C2A	2.66	119.90	112.53
4	R	1153	CYC	C2C-C3C-C4C	2.66	105.32	101.34
4	F	1153	CYC	CBC-CAC-C3C	2.66	119.05	113.41
4	D	1153	CYC	C2A-C1A-NA	2.65	113.80	110.04
4	D	1153	CYC	C1C-NC-C4C	-2.65	110.08	113.41
4	K	1084	CYC	CHA-C1A-NA	-2.65	119.04	124.60
4	C	1084	CYC	O2A-CGA-O1A	2.65	130.14	123.33
4	H	1082	CYC	CAD-CBD-CGD	-2.65	106.65	113.67
4	X	1153	CYC	C4D-ND-C1D	-2.65	105.14	109.78
4	I	1084	CYC	CAA-C2A-C3A	2.65	132.82	127.87
4	O	1084	CYC	C4D-ND-C1D	-2.64	105.14	109.78
4	W	1084	CYC	C1B-C2B-C3B	-2.64	105.15	107.86
4	F	1082	CYC	CAD-CBD-CGD	-2.64	106.67	113.67
4	H	1153	CYC	CHA-C4D-C3D	-2.64	121.55	127.22
4	H	1153	CYC	CBD-CAD-C3D	-2.64	105.24	112.53
4	F	1082	CYC	C1A-C2A-C3A	-2.64	103.88	106.73
4	S	1084	CYC	C2D-C1D-ND	2.64	111.03	107.43
4	B	1082	CYC	CMA-C3A-C4A	2.64	129.19	125.10
4	R	1082	CYC	C4D-ND-C1D	-2.63	105.16	109.78
4	O	1084	CYC	CMA-C3A-C4A	2.63	129.18	125.10
4	B	1153	CYC	C1C-NC-C4C	-2.63	110.11	113.41
4	C	1084	CYC	C1A-NA-C4A	-2.63	101.70	106.52
4	V	1153	CYC	C2C-C3C-C4C	2.61	105.25	101.34
4	L	1082	CYC	C1A-C2A-C3A	-2.61	103.90	106.73
4	J	1153	CYC	CBA-CAA-C2A	-2.61	105.31	112.53
4	N	1082	CYC	CMD-C2D-C1D	2.61	129.41	125.62

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	1084	CYC	CMD-C2D-C3D	2.61	131.16	125.62
4	N	1153	CYC	CBC-CAC-C3C	2.61	118.94	113.41
4	C	1084	CYC	CHB-C4A-C3A	2.60	131.56	124.87
4	A	1084	CYC	C4D-ND-C1D	-2.60	105.21	109.78
4	L	1153	CYC	CHB-C1B-C2B	-2.60	121.78	126.97
4	W	1084	CYC	C2C-C3C-C4C	2.60	105.23	101.34
4	P	1082	CYC	CHB-C4A-C3A	2.59	131.53	124.87
4	R	1153	CYC	CAA-C2A-C1A	-2.59	120.47	125.02
4	O	1084	CYC	CHA-C1A-NA	-2.59	119.16	124.60
4	X	1153	CYC	OB-C4B-NB	-2.59	119.01	125.08
4	Q	1084	CYC	O1A-CGA-CBA	-2.59	114.89	123.09
4	G	1084	CYC	C1A-NA-C4A	-2.58	101.78	106.52
4	U	1084	CYC	CAD-CBD-CGD	-2.58	106.83	113.67
4	P	1082	CYC	C1D-C2D-C3D	-2.58	104.53	107.76
4	H	1153	CYC	CHB-C1B-NB	-2.58	120.56	126.06
4	G	1084	CYC	O1D-CGD-CBD	-2.57	114.93	123.09
4	V	1082	CYC	CMC-C2C-C1C	-2.57	106.86	112.40
4	V	1082	CYC	CAA-C2A-C3A	2.57	132.68	127.87
4	R	1082	CYC	C2A-C1A-NA	2.57	113.68	110.04
4	X	1082	CYC	CAA-C2A-C1A	2.57	129.53	125.02
4	K	1084	CYC	OC-C1C-C2C	-2.57	124.13	126.17
4	J	1082	CYC	CMA-C3A-C4A	2.57	129.09	125.10
4	N	1153	CYC	CMD-C2D-C3D	-2.57	120.18	125.62
4	F	1082	CYC	CHB-C4A-C3A	2.56	131.46	124.87
4	B	1082	CYC	CMC-C2C-C1C	-2.56	106.88	112.40
4	C	1084	CYC	CAA-C2A-C3A	2.56	132.67	127.87
4	B	1082	CYC	CAB-C3B-C2B	2.56	132.26	127.56
4	P	1082	CYC	OC-C1C-NC	2.56	127.94	124.93
4	N	1153	CYC	O1D-CGD-CBD	-2.55	114.99	123.09
4	W	1084	CYC	C1D-CHD-C4C	2.55	132.12	127.76
4	G	1084	CYC	C4D-C3D-C2D	-2.54	104.66	107.62
4	N	1082	CYC	CAB-C3B-C2B	2.54	132.22	127.56
4	P	1082	CYC	CMB-C2B-C1B	2.53	127.23	124.16
4	N	1153	CYC	CAD-C3D-C4D	2.52	129.78	125.77
4	B	1153	CYC	C1B-CHB-C4A	2.52	134.25	128.06
4	X	1082	CYC	CAB-C3B-C4B	2.52	125.27	121.37
4	J	1082	CYC	CMC-C2C-C3C	-2.52	104.23	113.98
4	N	1153	CYC	CBD-CAD-C3D	-2.52	105.56	112.53
4	S	1084	CYC	C4D-C3D-C2D	-2.52	104.69	107.62
4	P	1082	CYC	CHD-C1D-C2D	2.52	133.70	127.53
4	C	1084	CYC	C4D-CHA-C1A	2.52	133.71	128.22
4	N	1082	CYC	CAD-C3D-C2D	-2.51	121.44	127.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	F	1153	CYC	C1A-NA-C4A	-2.51	101.91	106.52
4	J	1082	CYC	C2B-C1B-NB	2.49	110.60	106.97
4	Q	1084	CYC	CBA-CAA-C2A	2.49	119.42	112.53
4	E	1084	CYC	CHA-C1A-C2A	-2.49	119.65	125.40
4	F	1153	CYC	O2A-CGA-CBA	2.49	121.86	114.00
4	E	1084	CYC	C1B-C2B-C3B	-2.49	105.31	107.86
4	H	1153	CYC	CHA-C1A-C2A	-2.49	119.66	125.40
4	V	1082	CYC	CHA-C4D-C3D	-2.49	121.88	127.22
4	X	1153	CYC	CHB-C4A-C3A	2.48	131.25	124.87
4	K	1084	CYC	C4D-CHA-C1A	2.48	133.64	128.22
4	C	1084	CYC	CBB-CAB-C3B	-2.48	105.70	112.42
4	T	1153	CYC	CHB-C4A-C3A	2.48	131.24	124.87
4	T	1153	CYC	C1C-NC-C4C	-2.47	110.31	113.41
4	S	1084	CYC	CAD-C3D-C2D	2.47	132.61	127.07
4	E	1084	CYC	OC-C1C-C2C	-2.47	124.21	126.17
4	A	1084	CYC	C2D-C1D-ND	2.47	110.80	107.43
4	X	1153	CYC	CMD-C2D-C1D	2.46	129.19	125.62
4	H	1082	CYC	OC-C1C-C2C	-2.46	124.22	126.17
4	H	1153	CYC	OB-C4B-NB	-2.46	119.32	125.08
4	J	1153	CYC	CHD-C1D-C2D	2.45	133.54	127.53
4	T	1082	CYC	CMC-C2C-C3C	-2.45	104.51	113.98
4	H	1153	CYC	O2D-CGD-CBD	2.44	121.72	114.00
4	I	1084	CYC	CAD-C3D-C2D	2.44	132.55	127.07
4	D	1082	CYC	C2C-C3C-C4C	2.44	105.00	101.34
4	R	1153	CYC	O2A-CGA-O1A	-2.44	117.06	123.33
4	N	1082	CYC	CBA-CAA-C2A	2.44	119.28	112.53
4	I	1084	CYC	CHA-C1A-NA	-2.44	119.48	124.60
4	H	1082	CYC	CAB-C3B-C4B	2.43	125.13	121.37
4	O	1084	CYC	CBC-CAC-C3C	-2.43	108.25	113.41
4	C	1084	CYC	CMA-C3A-C4A	2.43	128.87	125.10
4	V	1153	CYC	C4D-ND-C1D	-2.43	105.52	109.78
4	R	1082	CYC	CAA-C2A-C3A	-2.43	123.32	127.87
4	N	1082	CYC	O1D-CGD-CBD	-2.43	115.39	123.09
4	D	1082	CYC	OC-C1C-NC	-2.42	122.07	124.93
4	F	1153	CYC	C1D-CHD-C4C	2.42	131.90	127.76
4	N	1082	CYC	O1A-CGA-CBA	-2.42	115.43	123.09
4	G	1084	CYC	CMB-C2B-C1B	2.41	127.09	124.16
4	L	1082	CYC	C1B-C2B-C3B	-2.41	105.38	107.86
4	V	1153	CYC	CBB-CAB-C3B	2.41	118.95	112.42
4	B	1082	CYC	CHD-C1D-C2D	2.41	133.43	127.53
4	E	1084	CYC	CHB-C4A-C3A	2.41	131.06	124.87
4	J	1153	CYC	CMA-C3A-C4A	2.41	128.84	125.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	V	1082	CYC	CBA-CAA-C2A	2.41	119.19	112.53
4	U	1084	CYC	C2C-C3C-C4C	2.40	104.94	101.34
4	I	1084	CYC	OC-C1C-C2C	-2.40	124.27	126.17
4	R	1153	CYC	CHD-C1D-C2D	2.40	133.40	127.53
4	C	1084	CYC	CAC-C3C-C4C	2.39	118.81	112.67
4	R	1153	CYC	CHB-C1B-NB	-2.39	120.96	126.06
4	G	1084	CYC	OB-C4B-C3B	-2.38	125.53	128.03
4	D	1153	CYC	CHA-C1A-C2A	-2.38	119.89	125.40
4	E	1084	CYC	O2A-CGA-O1A	2.38	129.46	123.33
4	A	1084	CYC	CMA-C3A-C4A	2.38	128.80	125.10
4	N	1153	CYC	CHB-C1B-NB	-2.38	120.98	126.06
4	H	1082	CYC	CAB-C3B-C2B	2.38	131.93	127.56
4	W	1084	CYC	CHB-C1B-C2B	2.38	131.72	126.97
4	J	1082	CYC	C1C-NC-C4C	-2.38	110.42	113.41
4	U	1084	CYC	CHA-C1A-NA	-2.38	119.61	124.60
4	T	1082	CYC	CAA-CBA-CGA	-2.37	107.37	113.67
4	Q	1084	CYC	C1C-NC-C4C	-2.37	110.43	113.41
4	Q	1084	CYC	CHB-C4A-C3A	2.37	130.97	124.87
4	O	1084	CYC	O1D-CGD-CBD	-2.37	115.56	123.09
4	J	1082	CYC	C2C-C3C-C4C	2.37	104.89	101.34
4	X	1082	CYC	C1A-C2A-C3A	-2.37	104.16	106.73
4	J	1153	CYC	CBD-CAD-C3D	2.37	119.08	112.53
4	M	1084	CYC	CMC-C2C-C1C	2.36	117.50	112.40
4	H	1082	CYC	C2C-C3C-C4C	2.36	104.88	101.34
4	N	1153	CYC	CHA-C4D-ND	-2.36	119.96	125.29
4	T	1082	CYC	C3C-C4C-NC	2.36	110.98	107.94
4	H	1082	CYC	CBD-CAD-C3D	2.35	119.04	112.53
4	E	1084	CYC	C4A-C3A-C2A	-2.35	103.81	106.48
4	B	1082	CYC	CHB-C1B-NB	-2.35	121.04	126.06
4	F	1153	CYC	CMD-C2D-C1D	2.35	129.04	125.62
4	W	1084	CYC	CHD-C1D-ND	-2.35	119.98	125.29
4	N	1153	CYC	O2A-CGA-CBA	2.35	121.42	114.00
4	K	1084	CYC	C3C-C4C-NC	2.34	110.96	107.94
4	R	1153	CYC	CHD-C4C-NC	-2.34	122.72	125.63
4	Q	1084	CYC	C4A-C3A-C2A	-2.34	103.83	106.48
4	B	1082	CYC	C3C-C4C-NC	2.33	110.95	107.94
4	X	1082	CYC	C2C-C3C-C4C	2.33	104.83	101.34
4	F	1082	CYC	CHA-C1A-NA	-2.33	119.70	124.60
4	G	1084	CYC	OB-C4B-NB	-2.33	119.61	125.08
4	W	1084	CYC	OC-C1C-NC	2.33	127.67	124.93
4	D	1082	CYC	C1B-CHB-C4A	-2.33	122.34	128.06
4	J	1153	CYC	CHB-C4A-NA	-2.33	119.93	124.95

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	T	1082	CYC	CMD-C2D-C1D	2.33	129.00	125.62
4	D	1082	CYC	CHA-C1A-NA	-2.33	119.72	124.60
4	I	1084	CYC	CAB-C3B-C2B	2.32	131.83	127.56
4	R	1082	CYC	CMA-C3A-C4A	2.32	128.71	125.10
4	L	1153	CYC	C1C-NC-C4C	-2.32	110.50	113.41
4	T	1153	CYC	CHD-C4C-NC	-2.32	122.74	125.63
4	H	1082	CYC	CBB-CAB-C3B	-2.32	106.14	112.42
4	V	1082	CYC	C1B-C2B-C3B	-2.32	105.48	107.86
4	F	1082	CYC	C1C-NC-C4C	-2.32	110.50	113.41
4	F	1153	CYC	OC-C1C-NC	2.31	127.65	124.93
4	F	1082	CYC	O2D-CGD-CBD	2.31	121.30	114.00
4	H	1153	CYC	O1D-CGD-CBD	-2.31	115.78	123.09
4	W	1084	CYC	CMB-C2B-C1B	2.30	126.95	124.16
4	I	1084	CYC	CHB-C4A-C3A	2.30	130.77	124.87
4	F	1082	CYC	C1B-CHB-C4A	-2.30	122.42	128.06
4	H	1082	CYC	C1D-CHD-C4C	-2.29	123.85	127.76
4	O	1084	CYC	C1A-NA-C4A	-2.29	102.31	106.52
4	H	1082	CYC	CHB-C1B-C2B	2.29	131.53	126.97
4	F	1153	CYC	CHA-C4D-C3D	-2.28	122.31	127.22
4	Q	1084	CYC	CAC-C3C-C4C	2.28	118.53	112.67
4	L	1153	CYC	CAB-C3B-C2B	2.28	131.74	127.56
4	I	1084	CYC	CHA-C1A-C2A	-2.27	120.15	125.40
4	B	1153	CYC	CBD-CAD-C3D	2.27	118.82	112.53
4	N	1082	CYC	CAB-C3B-C4B	2.27	124.89	121.37
4	R	1082	CYC	CHD-C1D-ND	-2.27	120.16	125.29
4	U	1084	CYC	O1A-CGA-CBA	-2.27	115.90	123.09
4	T	1153	CYC	CHA-C4D-C3D	-2.27	122.35	127.22
4	F	1082	CYC	C2C-C3C-C4C	2.25	104.71	101.34
4	V	1153	CYC	CHB-C1B-NB	-2.25	121.25	126.06
4	J	1082	CYC	CBD-CAD-C3D	2.25	118.75	112.53
4	P	1082	CYC	CMA-C3A-C4A	2.24	128.58	125.10
4	J	1082	CYC	CAD-C3D-C4D	2.24	129.32	125.77
4	L	1082	CYC	C3C-C4C-NC	2.24	110.82	107.94
4	K	1084	CYC	O2A-CGA-O1A	2.24	129.08	123.33
4	W	1084	CYC	O2A-CGA-CBA	2.23	121.06	114.00
4	T	1153	CYC	C3C-C4C-NC	2.23	110.81	107.94
4	T	1082	CYC	CBC-CAC-C3C	2.23	118.14	113.41
4	S	1084	CYC	CBB-CAB-C3B	-2.23	106.38	112.42
4	I	1084	CYC	O2A-CGA-O1A	2.23	129.07	123.33
4	N	1153	CYC	CMB-C2B-C1B	2.23	126.87	124.16
4	I	1084	CYC	CMB-C2B-C1B	2.23	126.87	124.16
4	H	1082	CYC	CBA-CAA-C2A	2.23	118.69	112.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	K	1084	CYC	CAA-C2A-C3A	2.23	132.04	127.87
4	I	1084	CYC	CBD-CAD-C3D	2.22	118.68	112.53
4	F	1153	CYC	CHA-C1A-C2A	-2.22	120.27	125.40
4	H	1082	CYC	C4A-C3A-C2A	-2.22	103.96	106.48
4	B	1153	CYC	C1A-NA-C4A	-2.21	102.46	106.52
4	M	1084	CYC	O2A-CGA-O1A	2.20	129.00	123.33
4	F	1082	CYC	CMC-C2C-C1C	-2.20	107.65	112.40
4	G	1084	CYC	CMC-C2C-C3C	2.20	122.50	113.98
4	A	1084	CYC	CHD-C1D-C2D	-2.20	122.14	127.53
4	I	1084	CYC	C2C-C3C-C4C	2.19	104.62	101.34
4	R	1153	CYC	CHA-C1A-C2A	-2.19	120.34	125.40
4	D	1153	CYC	CHB-C4A-NA	-2.19	120.23	124.95
4	X	1153	CYC	CHD-C1D-C2D	2.19	132.88	127.53
4	F	1082	CYC	CHD-C4C-NC	-2.18	122.90	125.63
4	O	1084	CYC	CHB-C4A-NA	-2.18	120.24	124.95
4	R	1153	CYC	CHA-C4D-ND	-2.18	120.37	125.29
4	P	1153	CYC	CHB-C4A-C3A	2.18	130.46	124.87
4	B	1153	CYC	CHD-C1D-C2D	2.17	132.84	127.53
4	O	1084	CYC	CHA-C4D-C3D	-2.17	122.56	127.22
4	X	1082	CYC	CHD-C1D-C2D	2.16	132.83	127.53
4	B	1153	CYC	O2A-CGA-O1A	-2.16	117.77	123.33
4	D	1082	CYC	CHD-C4C-NC	-2.16	122.93	125.63
4	N	1082	CYC	CMB-C2B-C3B	-2.16	120.30	126.15
4	N	1082	CYC	C1A-C2A-C3A	-2.16	104.39	106.73
4	C	1084	CYC	C1C-NC-C4C	-2.16	110.70	113.41
4	L	1153	CYC	C4D-ND-C1D	-2.16	106.00	109.78
4	T	1153	CYC	CBA-CAA-C2A	-2.15	106.58	112.53
4	M	1084	CYC	CBC-CAC-C3C	-2.15	108.85	113.41
4	E	1084	CYC	CAB-C3B-C2B	-2.15	123.61	127.56
4	N	1153	CYC	CAB-C3B-C2B	2.14	131.50	127.56
4	G	1084	CYC	CHA-C1A-NA	-2.14	120.10	124.60
4	S	1084	CYC	O2D-CGD-CBD	2.14	120.76	114.00
4	M	1084	CYC	C2C-C3C-C4C	2.14	104.54	101.34
4	V	1082	CYC	CBD-CAD-C3D	2.14	118.44	112.53
4	V	1153	CYC	CHB-C4A-C3A	2.14	130.36	124.87
4	N	1082	CYC	CBB-CAB-C3B	-2.14	106.63	112.42
4	B	1082	CYC	C1D-C2D-C3D	-2.13	105.09	107.76
4	U	1084	CYC	CMD-C2D-C1D	2.13	128.72	125.62
4	N	1153	CYC	C1D-C2D-C3D	-2.13	105.10	107.76
4	R	1153	CYC	CMB-C2B-C1B	2.13	126.74	124.16
4	D	1082	CYC	OC-C1C-C2C	-2.12	124.49	126.17
4	D	1082	CYC	C4A-C3A-C2A	-2.12	104.07	106.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	V	1082	CYC	CMC-C2C-C3C	-2.12	105.80	113.98
4	T	1153	CYC	C2C-C3C-C4C	2.12	104.51	101.34
4	W	1084	CYC	CBB-CAB-C3B	-2.12	106.69	112.42
4	P	1082	CYC	C3C-C4C-NC	2.11	110.66	107.94
4	F	1082	CYC	CMD-C2D-C1D	2.11	128.69	125.62
4	J	1153	CYC	C1D-C2D-C3D	-2.11	105.12	107.76
4	P	1153	CYC	C4D-C3D-C2D	-2.11	105.17	107.62
4	R	1082	CYC	CHD-C4C-NC	-2.11	123.00	125.63
4	P	1082	CYC	CAA-CBA-CGA	-2.11	108.08	113.67
4	D	1082	CYC	CHA-C4D-C3D	-2.10	122.70	127.22
4	U	1084	CYC	CHA-C4D-C3D	-2.10	122.70	127.22
4	P	1082	CYC	C1A-NA-C4A	-2.10	102.66	106.52
4	H	1153	CYC	C1A-NA-C4A	-2.09	102.68	106.52
4	P	1153	CYC	CBA-CAA-C2A	-2.09	106.76	112.53
4	G	1084	CYC	CAD-CBD-CGD	-2.09	108.12	113.67
4	F	1153	CYC	C1C-NC-C4C	-2.09	110.79	113.41
4	L	1082	CYC	CAB-C3B-C2B	2.09	131.39	127.56
4	U	1084	CYC	CMC-C2C-C3C	2.09	122.05	113.98
4	P	1153	CYC	CHD-C1D-C2D	2.09	132.64	127.53
4	C	1084	CYC	CHD-C4C-NC	2.08	128.23	125.63
4	H	1082	CYC	CHD-C4C-NC	-2.08	123.03	125.63
4	B	1153	CYC	C1A-C2A-C3A	2.08	108.98	106.73
4	P	1082	CYC	CMC-C2C-C3C	-2.08	105.95	113.98
4	U	1084	CYC	O2D-CGD-O1D	2.07	128.66	123.33
4	L	1153	CYC	C4D-CHA-C1A	2.07	132.74	128.22
4	R	1082	CYC	CHB-C4A-C3A	2.07	130.19	124.87
4	U	1084	CYC	CMA-C3A-C4A	2.06	128.31	125.10
4	I	1084	CYC	C1B-C2B-C3B	-2.05	105.75	107.86
4	A	1084	CYC	CAD-C3D-C2D	-2.05	122.47	127.07
4	B	1153	CYC	O2A-CGA-CBA	2.05	120.48	114.00
4	J	1153	CYC	C1A-NA-C4A	-2.05	102.75	106.52
4	N	1082	CYC	CBC-CAC-C3C	-2.05	109.07	113.41
4	R	1082	CYC	CAD-C3D-C2D	-2.05	122.48	127.07
4	X	1153	CYC	C1D-C2D-C3D	-2.05	105.20	107.76
4	N	1082	CYC	CAA-C2A-C3A	-2.04	124.04	127.87
4	A	1084	CYC	CBD-CAD-C3D	2.04	118.18	112.53
4	G	1084	CYC	C1B-C2B-C3B	-2.04	105.76	107.86
4	V	1153	CYC	C1A-NA-C4A	-2.04	102.78	106.52
4	D	1153	CYC	C1A-NA-C4A	-2.03	102.78	106.52
4	T	1153	CYC	C3A-C4A-NA	2.03	115.02	110.58
4	V	1153	CYC	OC-C1C-C2C	-2.03	124.56	126.17
4	B	1082	CYC	OC-C1C-C2C	-2.03	124.56	126.17

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	N	1153	CYC	O2D-CGD-CBD	2.03	120.40	114.00
4	B	1082	CYC	O1A-CGA-CBA	-2.02	116.67	123.09
4	M	1084	CYC	CHB-C4A-C3A	2.02	130.07	124.87
4	U	1084	CYC	OC-C1C-C2C	-2.02	124.57	126.17
4	V	1153	CYC	C1D-C2D-C3D	-2.02	105.23	107.76
4	P	1153	CYC	OC-C1C-NC	2.01	127.30	124.93
4	A	1084	CYC	CHA-C4D-C3D	-2.01	122.89	127.22
4	H	1153	CYC	C1B-C2B-C3B	-2.01	105.79	107.86
4	Q	1084	CYC	CAA-C2A-C1A	-2.01	121.49	125.02
4	G	1084	CYC	CMC-C2C-C1C	2.01	116.74	112.40
4	G	1084	CYC	C1D-C2D-C3D	-2.01	105.24	107.76
4	Q	1084	CYC	CBC-CAC-C3C	-2.01	109.15	113.41
4	V	1153	CYC	OC-C1C-NC	-2.01	122.56	124.93
4	T	1153	CYC	C4D-CHA-C1A	2.00	132.60	128.22
4	W	1084	CYC	C2D-C1D-ND	2.00	110.16	107.43
4	T	1082	CYC	CHA-C1A-C2A	-2.00	120.78	125.40
4	X	1153	CYC	CHA-C1A-C2A	-2.00	120.78	125.40
4	N	1082	CYC	CHD-C1D-ND	-2.00	120.77	125.29
4	X	1082	CYC	CMD-C2D-C1D	2.00	128.53	125.62

There are no chirality outliers.

All (404) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	1084	CYC	NA-C4A-CHB-C1B
4	A	1084	CYC	C3A-C4A-CHB-C1B
4	A	1084	CYC	C4C-C3C-CAC-CBC
4	A	1084	CYC	ND-C1D-CHD-C4C
4	A	1084	CYC	C2D-C1D-CHD-C4C
4	A	1084	CYC	C4D-C3D-CAD-CBD
4	B	1082	CYC	NA-C4A-CHB-C1B
4	B	1082	CYC	C3A-C4A-CHB-C1B
4	B	1082	CYC	ND-C1D-CHD-C4C
4	B	1082	CYC	C2D-C1D-CHD-C4C
4	B	1082	CYC	C2D-C3D-CAD-CBD
4	B	1153	CYC	NA-C4A-CHB-C1B
4	B	1153	CYC	C3A-C4A-CHB-C1B
4	B	1153	CYC	C2D-C1D-CHD-C4C
4	C	1084	CYC	C3A-C2A-CAA-CBA
4	C	1084	CYC	NA-C4A-CHB-C1B
4	C	1084	CYC	C3A-C4A-CHB-C1B
4	C	1084	CYC	ND-C1D-CHD-C4C

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Mol	Chain	Res	Type	Atoms
4	C	1084	CYC	C2D-C1D-CHD-C4C
4	D	1082	CYC	NA-C4A-CHB-C1B
4	D	1082	CYC	C3A-C4A-CHB-C1B
4	D	1082	CYC	ND-C1D-CHD-C4C
4	D	1082	CYC	C2D-C1D-CHD-C4C
4	D	1082	CYC	C4D-C3D-CAD-CBD
4	D	1153	CYC	NA-C4A-CHB-C1B
4	D	1153	CYC	C3A-C4A-CHB-C1B
4	D	1153	CYC	C2C-C3C-CAC-CBC
4	D	1153	CYC	C4C-C3C-CAC-CBC
4	D	1153	CYC	ND-C1D-CHD-C4C
4	D	1153	CYC	C2D-C1D-CHD-C4C
4	E	1084	CYC	C3A-C4A-CHB-C1B
4	E	1084	CYC	C4C-C3C-CAC-CBC
4	E	1084	CYC	ND-C1D-CHD-C4C
4	E	1084	CYC	C2D-C1D-CHD-C4C
4	F	1082	CYC	NA-C4A-CHB-C1B
4	F	1082	CYC	C3A-C4A-CHB-C1B
4	F	1082	CYC	ND-C1D-CHD-C4C
4	F	1082	CYC	C2D-C1D-CHD-C4C
4	F	1082	CYC	C2D-C3D-CAD-CBD
4	F	1082	CYC	C4D-C3D-CAD-CBD
4	F	1153	CYC	NA-C4A-CHB-C1B
4	F	1153	CYC	C3A-C4A-CHB-C1B
4	F	1153	CYC	C2C-C3C-CAC-CBC
4	F	1153	CYC	C4C-C3C-CAC-CBC
4	F	1153	CYC	ND-C1D-CHD-C4C
4	F	1153	CYC	C2D-C1D-CHD-C4C
4	G	1084	CYC	C1A-C2A-CAA-CBA
4	G	1084	CYC	C3A-C2A-CAA-CBA
4	G	1084	CYC	NA-C4A-CHB-C1B
4	G	1084	CYC	C3A-C4A-CHB-C1B
4	G	1084	CYC	C2C-C3C-CAC-CBC
4	G	1084	CYC	C4C-C3C-CAC-CBC
4	G	1084	CYC	ND-C1D-CHD-C4C
4	G	1084	CYC	C2D-C1D-CHD-C4C
4	H	1082	CYC	NA-C4A-CHB-C1B
4	H	1082	CYC	C3A-C4A-CHB-C1B
4	H	1082	CYC	ND-C1D-CHD-C4C
4	H	1082	CYC	C2D-C1D-CHD-C4C
4	H	1082	CYC	C4D-C3D-CAD-CBD
4	H	1153	CYC	NA-C4A-CHB-C1B

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Mol	Chain	Res	Type	Atoms
4	H	1153	CYC	C3A-C4A-CHB-C1B
4	H	1153	CYC	C2A-CAA-CBA-CGA
4	H	1153	CYC	C2B-C3B-CAB-CBB
4	H	1153	CYC	ND-C1D-CHD-C4C
4	H	1153	CYC	C2D-C1D-CHD-C4C
4	I	1084	CYC	C3A-C2A-CAA-CBA
4	I	1084	CYC	NA-C4A-CHB-C1B
4	I	1084	CYC	C3A-C4A-CHB-C1B
4	I	1084	CYC	C4C-C3C-CAC-CBC
4	I	1084	CYC	ND-C1D-CHD-C4C
4	I	1084	CYC	C2D-C1D-CHD-C4C
4	J	1082	CYC	NA-C4A-CHB-C1B
4	J	1082	CYC	C3A-C4A-CHB-C1B
4	J	1082	CYC	ND-C1D-CHD-C4C
4	J	1082	CYC	C2D-C1D-CHD-C4C
4	J	1153	CYC	NA-C4A-CHB-C1B
4	J	1153	CYC	C3A-C4A-CHB-C1B
4	J	1153	CYC	C2C-C3C-CAC-CBC
4	J	1153	CYC	C4C-C3C-CAC-CBC
4	J	1153	CYC	ND-C1D-CHD-C4C
4	J	1153	CYC	C2D-C1D-CHD-C4C
4	K	1084	CYC	C3A-C2A-CAA-CBA
4	K	1084	CYC	NA-C4A-CHB-C1B
4	K	1084	CYC	C3A-C4A-CHB-C1B
4	K	1084	CYC	ND-C1D-CHD-C4C
4	K	1084	CYC	C2D-C1D-CHD-C4C
4	L	1082	CYC	NA-C4A-CHB-C1B
4	L	1082	CYC	C3A-C4A-CHB-C1B
4	L	1082	CYC	ND-C1D-CHD-C4C
4	L	1082	CYC	C2D-C1D-CHD-C4C
4	L	1153	CYC	NA-C4A-CHB-C1B
4	L	1153	CYC	C3A-C4A-CHB-C1B
4	L	1153	CYC	C2B-C3B-CAB-CBB
4	L	1153	CYC	C2C-C3C-CAC-CBC
4	L	1153	CYC	C4C-C3C-CAC-CBC
4	L	1153	CYC	C2D-C1D-CHD-C4C
4	M	1084	CYC	C1A-C2A-CAA-CBA
4	M	1084	CYC	NA-C4A-CHB-C1B
4	M	1084	CYC	C3A-C4A-CHB-C1B
4	M	1084	CYC	C4C-C3C-CAC-CBC
4	M	1084	CYC	ND-C1D-CHD-C4C
4	M	1084	CYC	C2D-C1D-CHD-C4C

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Mol	Chain	Res	Type	Atoms
4	M	1084	CYC	C4D-C3D-CAD-CBD
4	N	1082	CYC	NA-C4A-CHB-C1B
4	N	1082	CYC	C3A-C4A-CHB-C1B
4	N	1082	CYC	ND-C1D-CHD-C4C
4	N	1082	CYC	C2D-C1D-CHD-C4C
4	N	1082	CYC	C4D-C3D-CAD-CBD
4	N	1153	CYC	NA-C4A-CHB-C1B
4	N	1153	CYC	C3A-C4A-CHB-C1B
4	N	1153	CYC	C2B-C3B-CAB-CBB
4	N	1153	CYC	ND-C1D-CHD-C4C
4	N	1153	CYC	C2D-C1D-CHD-C4C
4	O	1084	CYC	C1A-C2A-CAA-CBA
4	O	1084	CYC	C3A-C2A-CAA-CBA
4	O	1084	CYC	NA-C4A-CHB-C1B
4	O	1084	CYC	C3A-C4A-CHB-C1B
4	O	1084	CYC	C4B-C3B-CAB-CBB
4	O	1084	CYC	C4C-C3C-CAC-CBC
4	O	1084	CYC	ND-C1D-CHD-C4C
4	O	1084	CYC	C2D-C1D-CHD-C4C
4	O	1084	CYC	C4D-C3D-CAD-CBD
4	P	1082	CYC	C3A-C2A-CAA-CBA
4	P	1082	CYC	NA-C4A-CHB-C1B
4	P	1082	CYC	C3A-C4A-CHB-C1B
4	P	1082	CYC	ND-C1D-CHD-C4C
4	P	1082	CYC	C2D-C1D-CHD-C4C
4	P	1153	CYC	NA-C4A-CHB-C1B
4	P	1153	CYC	C3A-C4A-CHB-C1B
4	P	1153	CYC	C2C-C3C-CAC-CBC
4	P	1153	CYC	C4C-C3C-CAC-CBC
4	P	1153	CYC	ND-C1D-CHD-C4C
4	P	1153	CYC	C2D-C1D-CHD-C4C
4	Q	1084	CYC	C3A-C2A-CAA-CBA
4	Q	1084	CYC	NA-C4A-CHB-C1B
4	Q	1084	CYC	C3A-C4A-CHB-C1B
4	Q	1084	CYC	C4C-C3C-CAC-CBC
4	Q	1084	CYC	ND-C1D-CHD-C4C
4	Q	1084	CYC	C2D-C1D-CHD-C4C
4	R	1082	CYC	NA-C4A-CHB-C1B
4	R	1082	CYC	C3A-C4A-CHB-C1B
4	R	1082	CYC	ND-C1D-CHD-C4C
4	R	1082	CYC	C2D-C1D-CHD-C4C
4	R	1153	CYC	NA-C4A-CHB-C1B

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Mol	Chain	Res	Type	Atoms
4	R	1153	CYC	C3A-C4A-CHB-C1B
4	R	1153	CYC	C2B-C3B-CAB-CBB
4	R	1153	CYC	C4B-C3B-CAB-CBB
4	R	1153	CYC	C2C-C3C-CAC-CBC
4	R	1153	CYC	C4C-C3C-CAC-CBC
4	R	1153	CYC	ND-C1D-CHD-C4C
4	R	1153	CYC	C2D-C1D-CHD-C4C
4	S	1084	CYC	C3A-C2A-CAA-CBA
4	S	1084	CYC	NA-C4A-CHB-C1B
4	S	1084	CYC	C3A-C4A-CHB-C1B
4	S	1084	CYC	C4C-C3C-CAC-CBC
4	S	1084	CYC	ND-C1D-CHD-C4C
4	S	1084	CYC	C2D-C1D-CHD-C4C
4	T	1082	CYC	C3A-C2A-CAA-CBA
4	T	1082	CYC	NA-C4A-CHB-C1B
4	T	1082	CYC	C3A-C4A-CHB-C1B
4	T	1082	CYC	ND-C1D-CHD-C4C
4	T	1082	CYC	C2D-C1D-CHD-C4C
4	T	1082	CYC	C2D-C3D-CAD-CBD
4	T	1153	CYC	NA-C4A-CHB-C1B
4	T	1153	CYC	C3A-C4A-CHB-C1B
4	T	1153	CYC	C2A-CAA-CBA-CGA
4	T	1153	CYC	C4C-C3C-CAC-CBC
4	T	1153	CYC	ND-C1D-CHD-C4C
4	T	1153	CYC	C2D-C1D-CHD-C4C
4	U	1084	CYC	C1A-C2A-CAA-CBA
4	U	1084	CYC	C3A-C2A-CAA-CBA
4	U	1084	CYC	NA-C4A-CHB-C1B
4	U	1084	CYC	C3A-C4A-CHB-C1B
4	U	1084	CYC	C2C-C3C-CAC-CBC
4	U	1084	CYC	C4C-C3C-CAC-CBC
4	U	1084	CYC	ND-C1D-CHD-C4C
4	U	1084	CYC	C2D-C1D-CHD-C4C
4	U	1084	CYC	C4D-C3D-CAD-CBD
4	V	1082	CYC	C1A-C2A-CAA-CBA
4	V	1082	CYC	C3A-C2A-CAA-CBA
4	V	1082	CYC	NA-C4A-CHB-C1B
4	V	1082	CYC	C3A-C4A-CHB-C1B
4	V	1082	CYC	ND-C1D-CHD-C4C
4	V	1082	CYC	C2D-C1D-CHD-C4C
4	V	1153	CYC	NA-C4A-CHB-C1B
4	V	1153	CYC	C3A-C4A-CHB-C1B

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Mol	Chain	Res	Type	Atoms
4	V	1153	CYC	ND-C1D-CHD-C4C
4	V	1153	CYC	C2D-C1D-CHD-C4C
4	W	1084	CYC	C3A-C2A-CAA-CBA
4	W	1084	CYC	NA-C4A-CHB-C1B
4	W	1084	CYC	C3A-C4A-CHB-C1B
4	W	1084	CYC	ND-C1D-CHD-C4C
4	W	1084	CYC	C2D-C1D-CHD-C4C
4	X	1082	CYC	NA-C4A-CHB-C1B
4	X	1082	CYC	C3A-C4A-CHB-C1B
4	X	1082	CYC	ND-C1D-CHD-C4C
4	X	1082	CYC	C2D-C1D-CHD-C4C
4	X	1153	CYC	NA-C4A-CHB-C1B
4	X	1153	CYC	C3A-C4A-CHB-C1B
4	X	1153	CYC	ND-C1D-CHD-C4C
4	X	1153	CYC	C2D-C1D-CHD-C4C
4	X	1153	CYC	C4D-C3D-CAD-CBD
4	D	1153	CYC	C2B-C3B-CAB-CBB
4	T	1153	CYC	C2B-C3B-CAB-CBB
4	V	1153	CYC	C2B-C3B-CAB-CBB
4	P	1153	CYC	C2B-C3B-CAB-CBB
4	X	1153	CYC	C2B-C3B-CAB-CBB
4	O	1084	CYC	C2B-C3B-CAB-CBB
4	F	1153	CYC	C2B-C3B-CAB-CBB
4	E	1084	CYC	C3A-C2A-CAA-CBA
4	M	1084	CYC	C3A-C2A-CAA-CBA
4	C	1084	CYC	C2D-C3D-CAD-CBD
4	H	1082	CYC	C2D-C3D-CAD-CBD
4	B	1153	CYC	ND-C1D-CHD-C4C
4	L	1153	CYC	ND-C1D-CHD-C4C
4	C	1084	CYC	C1A-C2A-CAA-CBA
4	E	1084	CYC	C1A-C2A-CAA-CBA
4	I	1084	CYC	C1A-C2A-CAA-CBA
4	K	1084	CYC	C1A-C2A-CAA-CBA
4	P	1082	CYC	C1A-C2A-CAA-CBA
4	Q	1084	CYC	C1A-C2A-CAA-CBA
4	S	1084	CYC	C1A-C2A-CAA-CBA
4	T	1082	CYC	C1A-C2A-CAA-CBA
4	W	1084	CYC	C1A-C2A-CAA-CBA
4	A	1084	CYC	C2B-C3B-CAB-CBB
4	B	1082	CYC	C2B-C3B-CAB-CBB
4	C	1084	CYC	C4D-C3D-CAD-CBD
4	J	1153	CYC	C4D-C3D-CAD-CBD

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Mol	Chain	Res	Type	Atoms
4	K	1084	CYC	C4D-C3D-CAD-CBD
4	Q	1084	CYC	C4D-C3D-CAD-CBD
4	R	1082	CYC	C4D-C3D-CAD-CBD
4	V	1082	CYC	C4D-C3D-CAD-CBD
4	K	1084	CYC	C2D-C3D-CAD-CBD
4	R	1082	CYC	C2D-C3D-CAD-CBD
4	X	1153	CYC	C2D-C3D-CAD-CBD
4	B	1153	CYC	C2B-C3B-CAB-CBB
4	L	1082	CYC	C1A-C2A-CAA-CBA
4	H	1082	CYC	C3A-C2A-CAA-CBA
4	A	1084	CYC	C2D-C3D-CAD-CBD
4	D	1082	CYC	C2D-C3D-CAD-CBD
4	J	1153	CYC	C2D-C3D-CAD-CBD
4	Q	1084	CYC	C2D-C3D-CAD-CBD
4	V	1082	CYC	C2D-C3D-CAD-CBD
4	D	1153	CYC	C4B-C3B-CAB-CBB
4	H	1153	CYC	C4B-C3B-CAB-CBB
4	L	1153	CYC	C4B-C3B-CAB-CBB
4	N	1153	CYC	C4B-C3B-CAB-CBB
4	P	1153	CYC	C4B-C3B-CAB-CBB
4	V	1153	CYC	C4B-C3B-CAB-CBB
4	X	1153	CYC	C4B-C3B-CAB-CBB
4	L	1082	CYC	C2B-C3B-CAB-CBB
4	B	1082	CYC	C4D-C3D-CAD-CBD
4	T	1082	CYC	C4D-C3D-CAD-CBD
4	E	1084	CYC	NA-C4A-CHB-C1B
4	L	1082	CYC	C3A-C2A-CAA-CBA
4	R	1153	CYC	C2A-CAA-CBA-CGA
4	M	1084	CYC	C2D-C3D-CAD-CBD
4	N	1082	CYC	C2D-C3D-CAD-CBD
4	O	1084	CYC	C2D-C3D-CAD-CBD
4	U	1084	CYC	C2D-C3D-CAD-CBD
4	J	1082	CYC	C2B-C3B-CAB-CBB
4	H	1082	CYC	C1A-C2A-CAA-CBA
4	A	1084	CYC	C4B-C3B-CAB-CBB
4	F	1153	CYC	C4B-C3B-CAB-CBB
4	T	1153	CYC	C4B-C3B-CAB-CBB
4	J	1082	CYC	C4D-C3D-CAD-CBD
4	L	1153	CYC	C2A-CAA-CBA-CGA
4	N	1153	CYC	C2A-CAA-CBA-CGA
4	E	1084	CYC	C2C-C3C-CAC-CBC
4	O	1084	CYC	C2C-C3C-CAC-CBC

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Mol	Chain	Res	Type	Atoms
4	T	1153	CYC	C2C-C3C-CAC-CBC
4	L	1153	CYC	C1A-C2A-CAA-CBA
4	B	1082	CYC	C1A-C2A-CAA-CBA
4	F	1082	CYC	C1A-C2A-CAA-CBA
4	J	1082	CYC	C2D-C3D-CAD-CBD
4	A	1084	CYC	C1A-C2A-CAA-CBA
4	D	1153	CYC	C2A-CAA-CBA-CGA
4	R	1082	CYC	C2A-CAA-CBA-CGA
4	B	1082	CYC	C4B-C3B-CAB-CBB
4	C	1084	CYC	C2A-CAA-CBA-CGA
4	G	1084	CYC	C2A-CAA-CBA-CGA
4	P	1153	CYC	C3D-CAD-CBD-CGD
4	X	1153	CYC	C2A-CAA-CBA-CGA
4	C	1084	CYC	C4C-C3C-CAC-CBC
4	X	1082	CYC	C2B-C3B-CAB-CBB
4	B	1153	CYC	C4B-C3B-CAB-CBB
4	L	1082	CYC	C4B-C3B-CAB-CBB
4	E	1084	CYC	C2A-CAA-CBA-CGA
4	X	1153	CYC	C3D-CAD-CBD-CGD
4	F	1082	CYC	C3A-C2A-CAA-CBA
4	T	1153	CYC	C3D-CAD-CBD-CGD
4	L	1153	CYC	C3A-C2A-CAA-CBA
4	B	1082	CYC	C3A-C2A-CAA-CBA
4	K	1084	CYC	CAD-CBD-CGD-O2D
4	A	1084	CYC	CAA-CBA-CGA-O2A
4	F	1153	CYC	CAA-CBA-CGA-O2A
4	J	1082	CYC	C4B-C3B-CAB-CBB
4	P	1153	CYC	CAA-CBA-CGA-O2A
4	F	1082	CYC	CAA-CBA-CGA-O2A
4	C	1084	CYC	CAA-CBA-CGA-O1A
4	D	1082	CYC	CAA-CBA-CGA-O2A
4	O	1084	CYC	CAA-CBA-CGA-O1A
4	S	1084	CYC	CAA-CBA-CGA-O1A
4	V	1153	CYC	C3D-CAD-CBD-CGD
4	E	1084	CYC	CAA-CBA-CGA-O1A
4	P	1153	CYC	CAA-CBA-CGA-O1A
4	R	1153	CYC	CAA-CBA-CGA-O2A
4	W	1084	CYC	CAA-CBA-CGA-O2A
4	K	1084	CYC	CAA-CBA-CGA-O1A
4	K	1084	CYC	CAD-CBD-CGD-O1D
4	O	1084	CYC	CAD-CBD-CGD-O2D
4	B	1082	CYC	CAD-CBD-CGD-O2D

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Mol	Chain	Res	Type	Atoms
4	F	1153	CYC	CAA-CBA-CGA-O1A
4	G	1084	CYC	CAA-CBA-CGA-O1A
4	I	1084	CYC	CAA-CBA-CGA-O1A
4	C	1084	CYC	CAD-CBD-CGD-O2D
4	J	1082	CYC	CAA-CBA-CGA-O2A
4	O	1084	CYC	CAD-CBD-CGD-O1D
4	R	1153	CYC	CAA-CBA-CGA-O1A
4	T	1082	CYC	CAD-CBD-CGD-O2D
4	D	1153	CYC	CAA-CBA-CGA-O1A
4	H	1082	CYC	CAA-CBA-CGA-O1A
4	N	1082	CYC	CAA-CBA-CGA-O2A
4	C	1084	CYC	CAD-CBD-CGD-O1D
4	U	1084	CYC	CAA-CBA-CGA-O2A
4	A	1084	CYC	CAA-CBA-CGA-O1A
4	H	1153	CYC	CAA-CBA-CGA-O2A
4	M	1084	CYC	CAA-CBA-CGA-O1A
4	M	1084	CYC	CAA-CBA-CGA-O2A
4	D	1153	CYC	CAA-CBA-CGA-O2A
4	L	1153	CYC	CAA-CBA-CGA-O1A
4	Q	1084	CYC	CAA-CBA-CGA-O2A
4	X	1082	CYC	CAA-CBA-CGA-O1A
4	C	1084	CYC	C3D-CAD-CBD-CGD
4	L	1153	CYC	C3D-CAD-CBD-CGD
4	N	1153	CYC	C3D-CAD-CBD-CGD
4	K	1084	CYC	CAA-CBA-CGA-O2A
4	T	1153	CYC	CAA-CBA-CGA-O1A
4	U	1084	CYC	CAA-CBA-CGA-O1A
4	C	1084	CYC	CAA-CBA-CGA-O2A
4	N	1082	CYC	CAD-CBD-CGD-O1D
4	X	1082	CYC	CAA-CBA-CGA-O2A
4	X	1153	CYC	CAD-CBD-CGD-O1D
4	M	1084	CYC	C2C-C3C-CAC-CBC
4	I	1084	CYC	CAA-CBA-CGA-O2A
4	V	1082	CYC	CAA-CBA-CGA-O2A
4	V	1082	CYC	CAD-CBD-CGD-O2D
4	H	1082	CYC	CAA-CBA-CGA-O2A
4	O	1084	CYC	CAA-CBA-CGA-O2A
4	P	1082	CYC	CAA-CBA-CGA-O2A
4	Q	1084	CYC	CAA-CBA-CGA-O1A
4	E	1084	CYC	CAA-CBA-CGA-O2A
4	R	1082	CYC	CAA-CBA-CGA-O1A
4	A	1084	CYC	C3A-C2A-CAA-CBA

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Mol	Chain	Res	Type	Atoms
4	B	1082	CYC	CAD-CBD-CGD-O1D
4	G	1084	CYC	CAA-CBA-CGA-O2A
4	H	1153	CYC	CAA-CBA-CGA-O1A
4	S	1084	CYC	CAA-CBA-CGA-O2A
4	V	1082	CYC	CAA-CBA-CGA-O1A
4	B	1082	CYC	CAA-CBA-CGA-O1A
4	D	1082	CYC	CAA-CBA-CGA-O1A
4	F	1082	CYC	CAA-CBA-CGA-O1A
4	R	1082	CYC	CAA-CBA-CGA-O2A
4	J	1082	CYC	CAA-CBA-CGA-O1A
4	T	1082	CYC	CAD-CBD-CGD-O1D
4	V	1153	CYC	CAA-CBA-CGA-O2A
4	L	1153	CYC	CAA-CBA-CGA-O2A
4	Q	1084	CYC	C2B-C3B-CAB-CBB
4	B	1082	CYC	CAA-CBA-CGA-O2A
4	N	1082	CYC	CAD-CBD-CGD-O2D
4	P	1082	CYC	CAA-CBA-CGA-O1A
4	X	1153	CYC	CAD-CBD-CGD-O2D
4	J	1153	CYC	CAA-CBA-CGA-O2A
4	N	1082	CYC	CAA-CBA-CGA-O1A
4	W	1084	CYC	CAA-CBA-CGA-O1A
4	G	1084	CYC	C2B-C3B-CAB-CBB
4	A	1084	CYC	C3D-CAD-CBD-CGD
4	T	1153	CYC	CAA-CBA-CGA-O2A
4	J	1153	CYC	CAA-CBA-CGA-O1A
4	V	1153	CYC	CAA-CBA-CGA-O1A
4	L	1082	CYC	CAA-CBA-CGA-O2A
4	B	1153	CYC	CAD-CBD-CGD-O2D
4	L	1082	CYC	CAA-CBA-CGA-O1A
4	N	1153	CYC	CAA-CBA-CGA-O1A
4	V	1082	CYC	CAD-CBD-CGD-O1D
4	N	1153	CYC	CAA-CBA-CGA-O2A
4	D	1082	CYC	C2A-CAA-CBA-CGA
4	B	1153	CYC	CAD-CBD-CGD-O1D
4	R	1082	CYC	CAD-CBD-CGD-O2D
4	J	1153	CYC	C2A-CAA-CBA-CGA
4	J	1153	CYC	CAD-CBD-CGD-O1D
4	F	1082	CYC	C2B-C3B-CAB-CBB
4	E	1084	CYC	CAD-CBD-CGD-O1D
4	J	1153	CYC	CAD-CBD-CGD-O2D
4	A	1084	CYC	C2C-C3C-CAC-CBC
4	R	1082	CYC	CAD-CBD-CGD-O1D

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Mol	Chain	Res	Type	Atoms
4	V	1082	CYC	C2B-C3B-CAB-CBB
4	E	1084	CYC	CAD-CBD-CGD-O2D
4	G	1084	CYC	CAD-CBD-CGD-O1D
4	P	1082	CYC	CAD-CBD-CGD-O2D
4	J	1082	CYC	C1A-C2A-CAA-CBA
4	W	1084	CYC	CAD-CBD-CGD-O2D
4	T	1082	CYC	C2A-CAA-CBA-CGA
4	G	1084	CYC	CAD-CBD-CGD-O2D

There are no ring outliers.

36 monomers are involved in 600 short contacts:

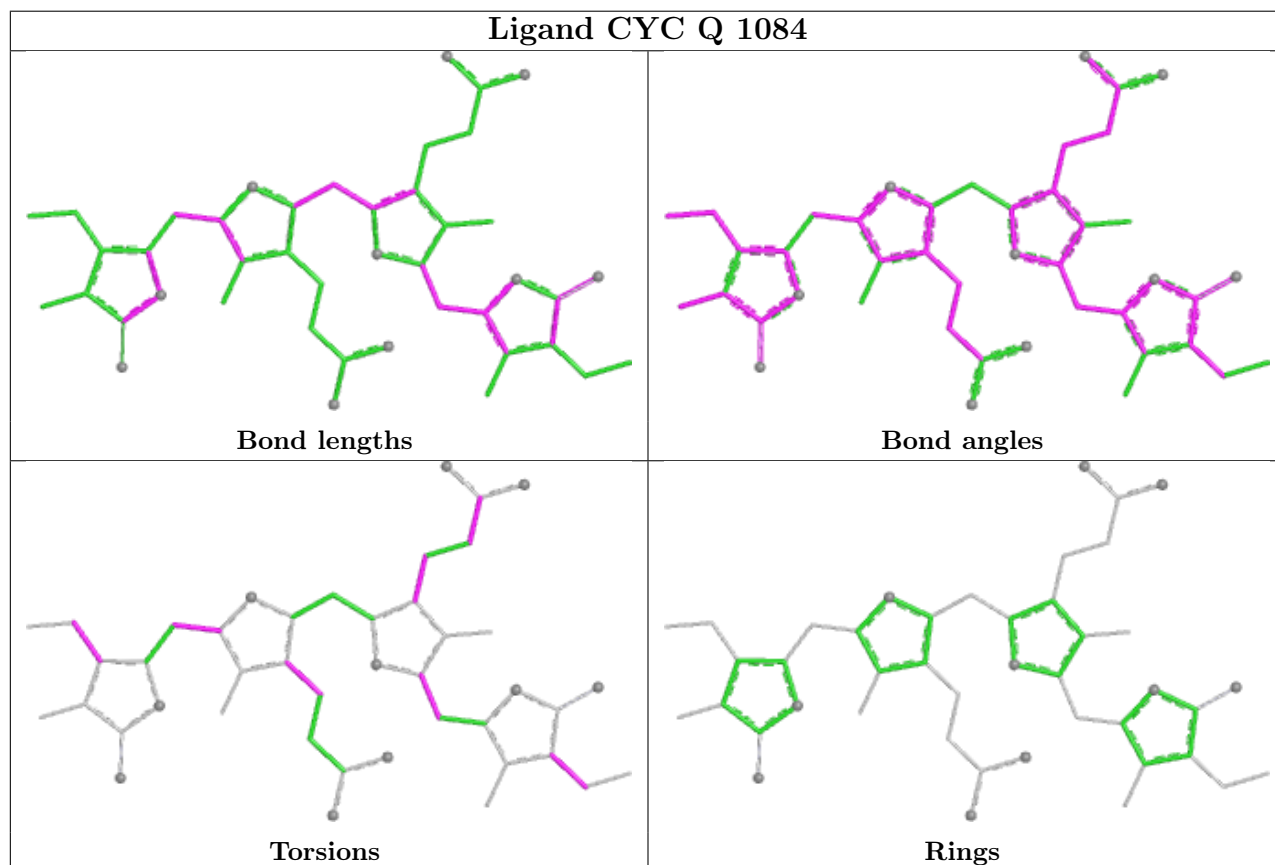
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	Q	1084	CYC	18	0
4	D	1153	CYC	21	0
4	V	1153	CYC	20	0
4	K	1084	CYC	19	0
4	O	1084	CYC	21	0
4	L	1153	CYC	28	0
4	A	1084	CYC	15	0
4	H	1082	CYC	18	0
4	D	1082	CYC	16	0
4	J	1082	CYC	14	0
4	B	1153	CYC	16	0
4	N	1082	CYC	12	0
4	P	1082	CYC	12	0
4	U	1084	CYC	14	0
4	E	1084	CYC	16	0
4	V	1082	CYC	14	0
4	F	1082	CYC	10	0
4	R	1153	CYC	19	0
4	F	1153	CYC	17	0
4	X	1153	CYC	23	0
4	W	1084	CYC	14	0
4	L	1082	CYC	18	0
4	H	1153	CYC	21	0
4	X	1082	CYC	15	0
4	T	1153	CYC	22	0
4	N	1153	CYC	23	0
4	S	1084	CYC	10	0
4	J	1153	CYC	24	0
4	T	1082	CYC	8	0

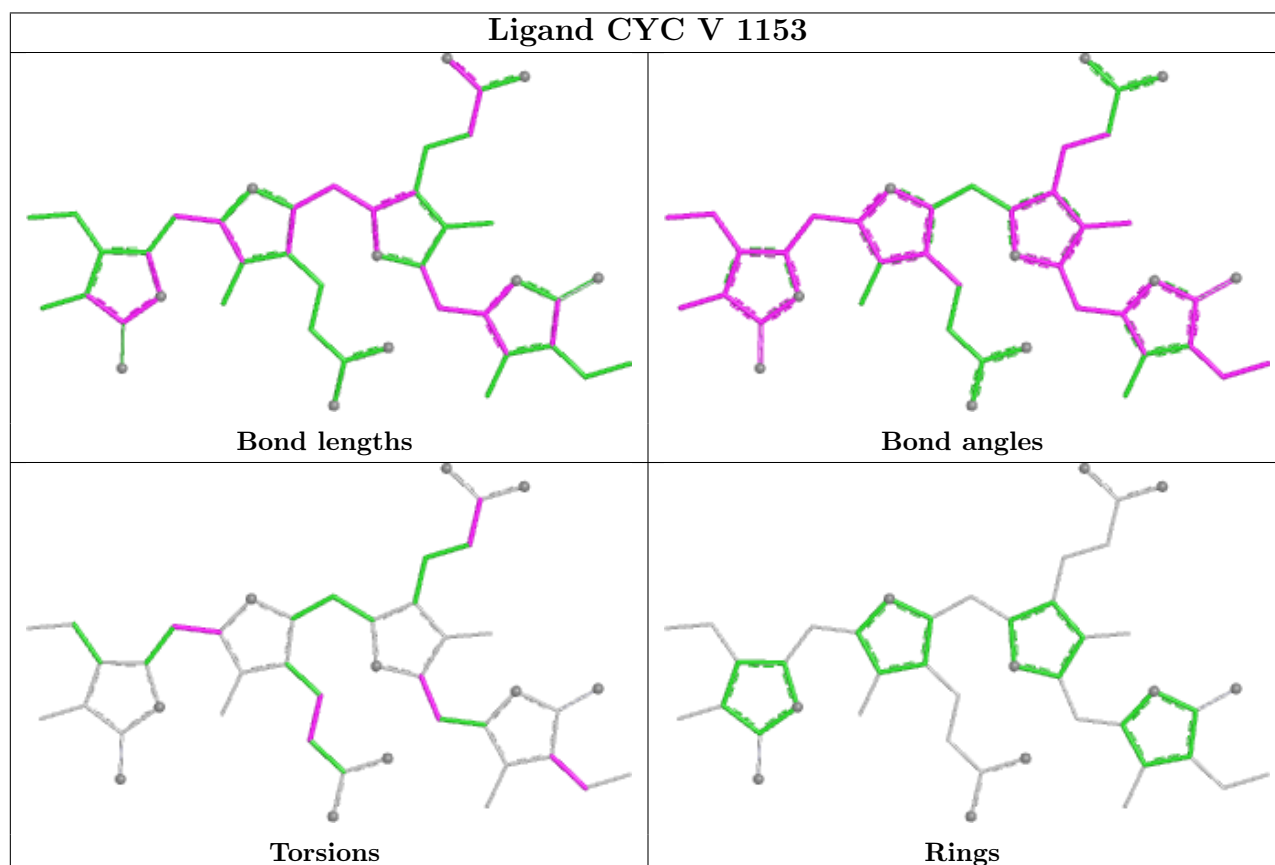
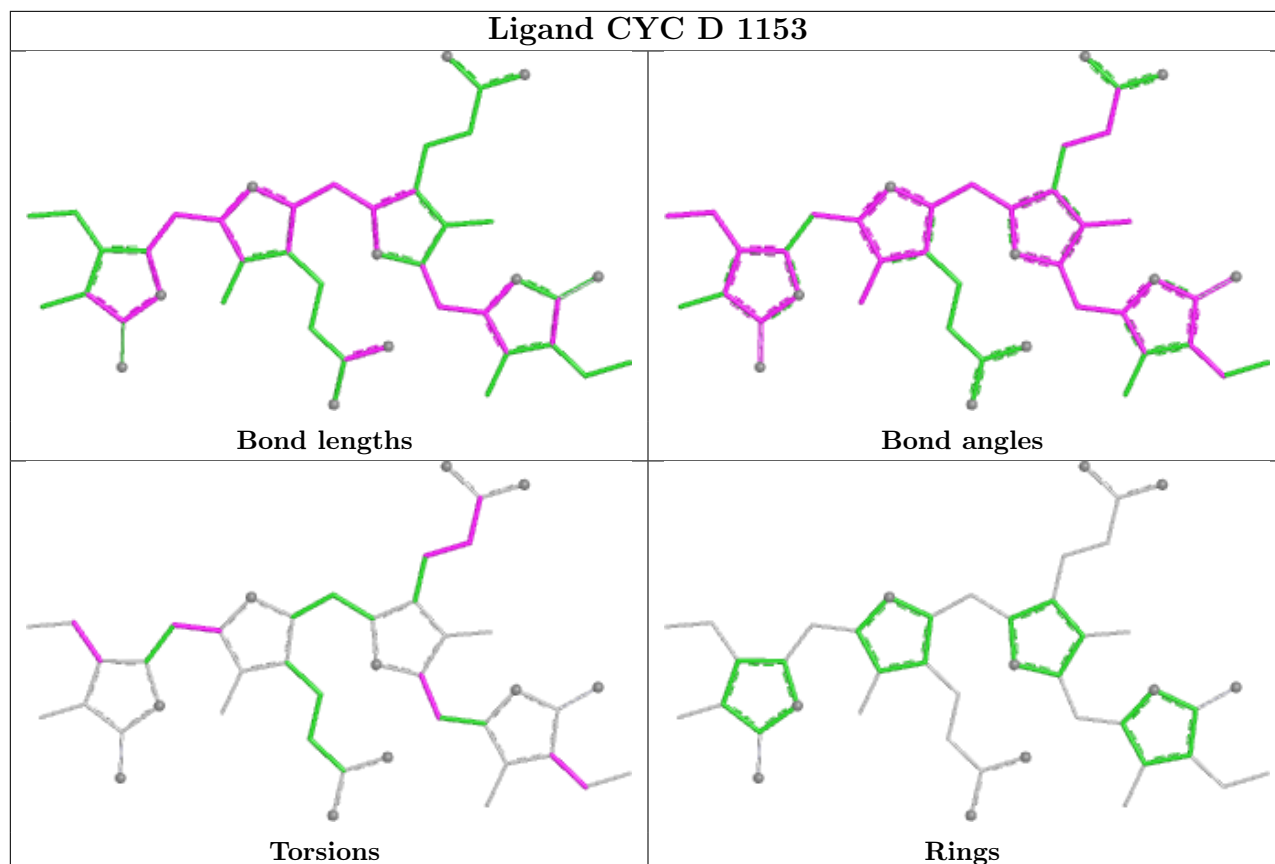
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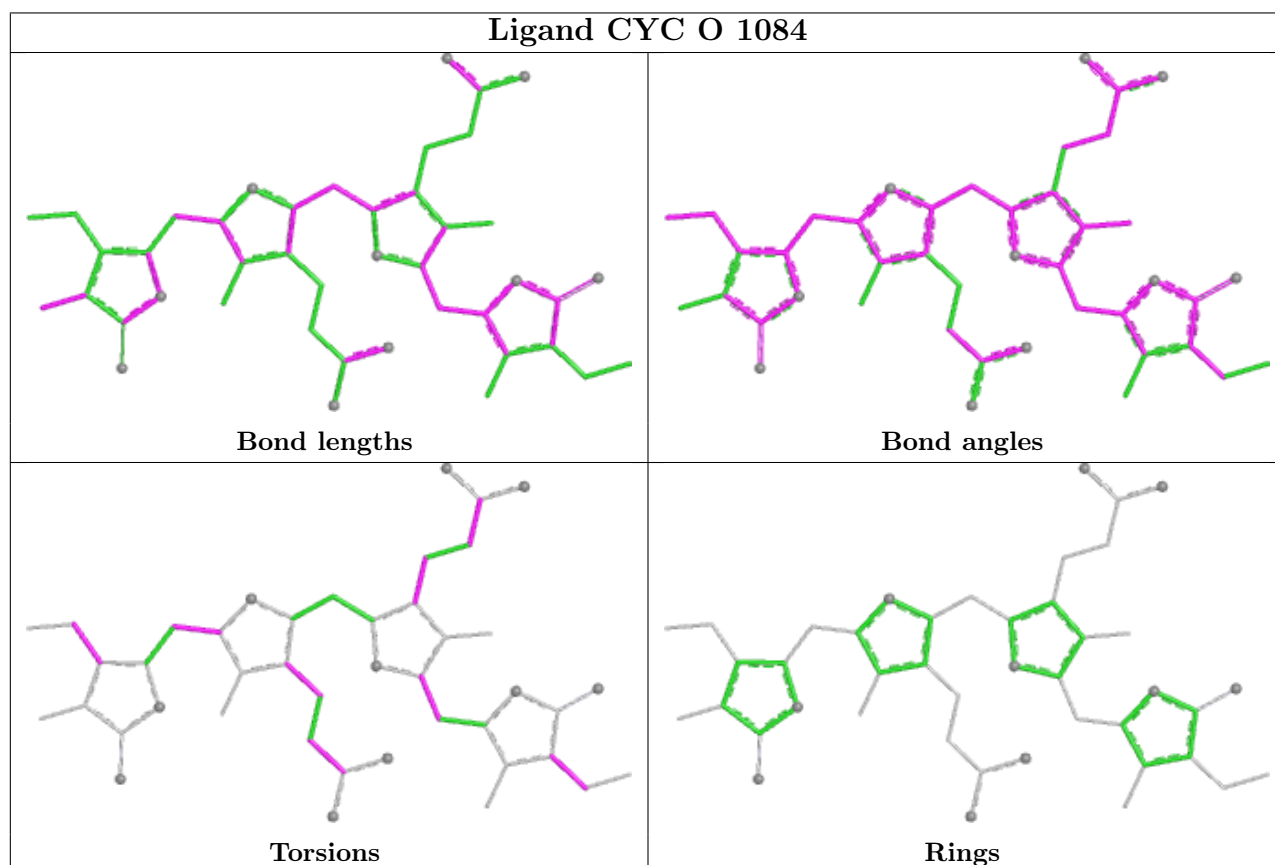
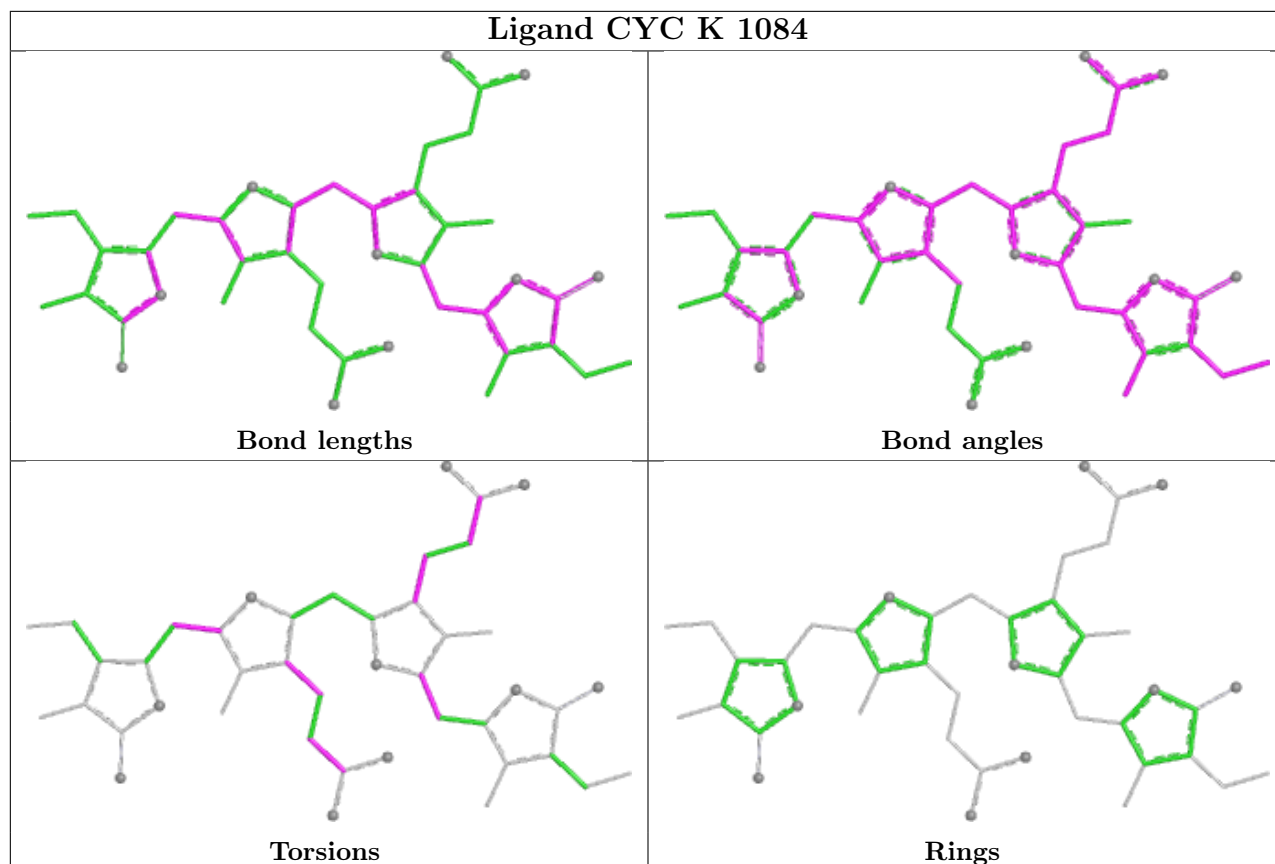
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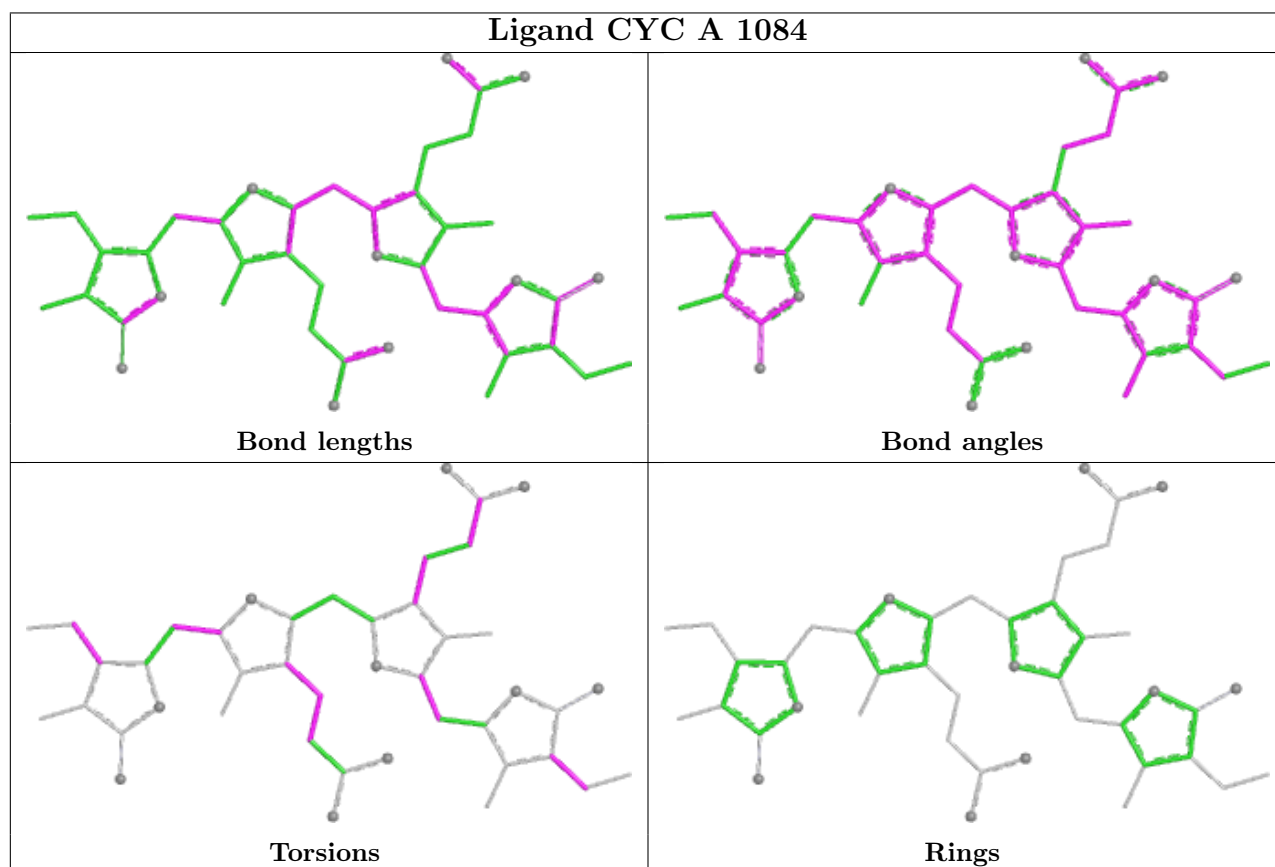
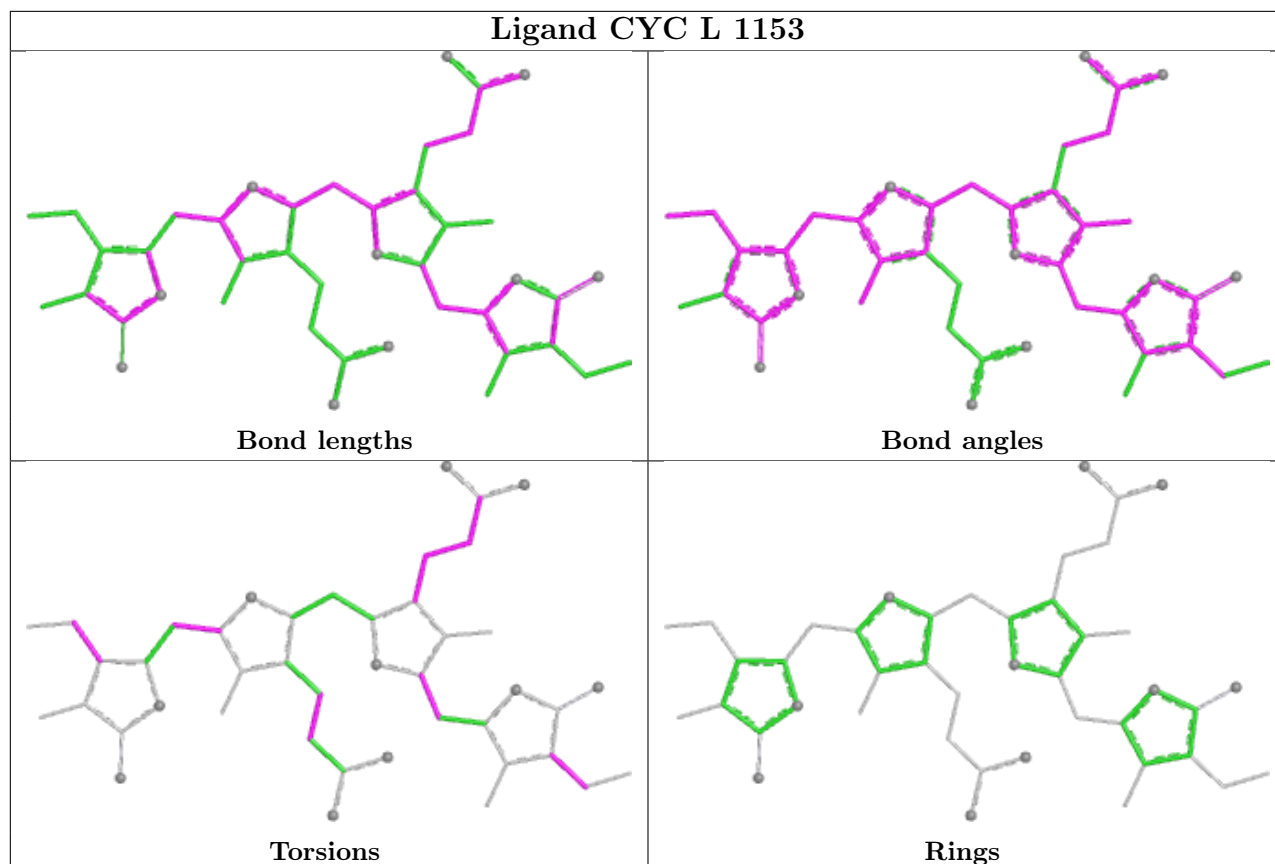
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	1084	CYC	14	0
4	P	1153	CYC	18	0
4	I	1084	CYC	13	0
4	M	1084	CYC	22	0
4	G	1084	CYC	11	0
4	B	1082	CYC	13	0
4	R	1082	CYC	11	0

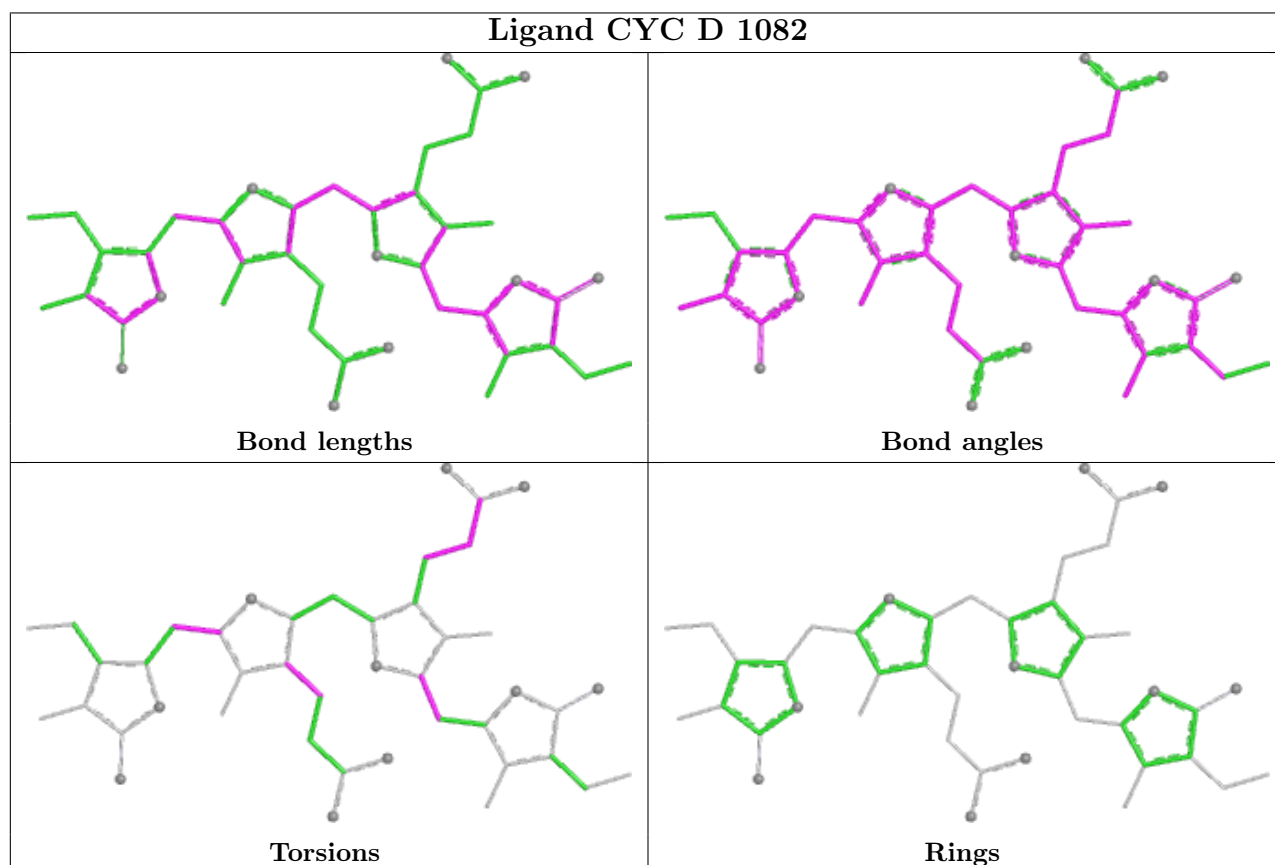
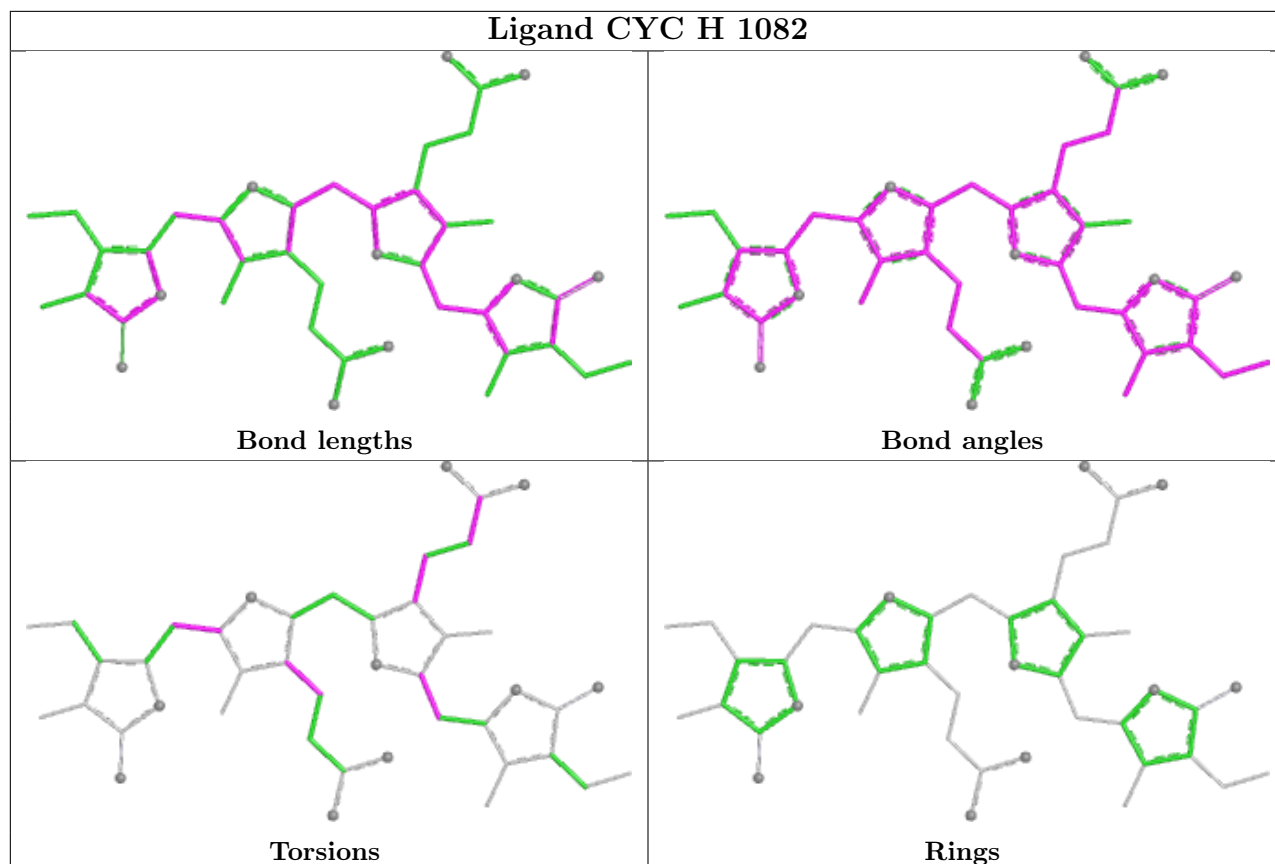
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

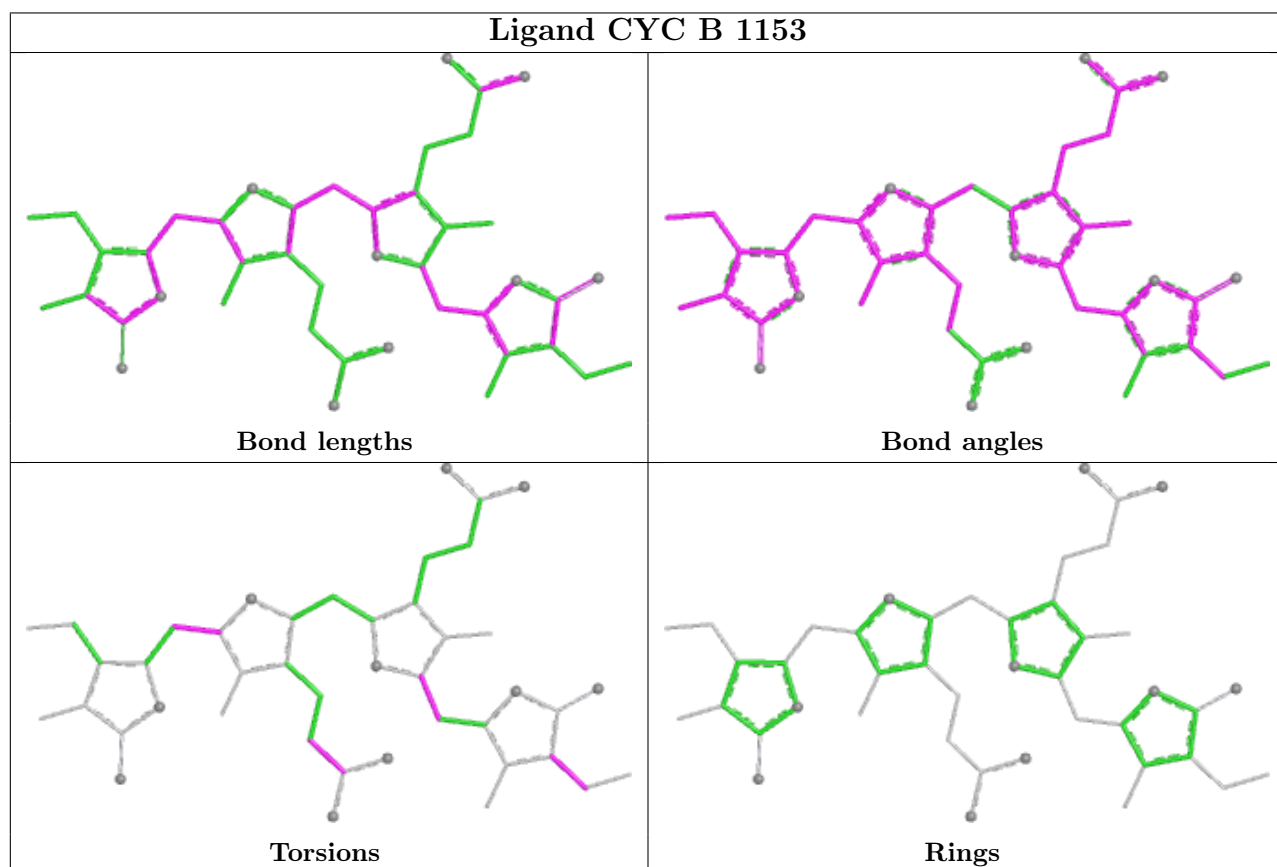
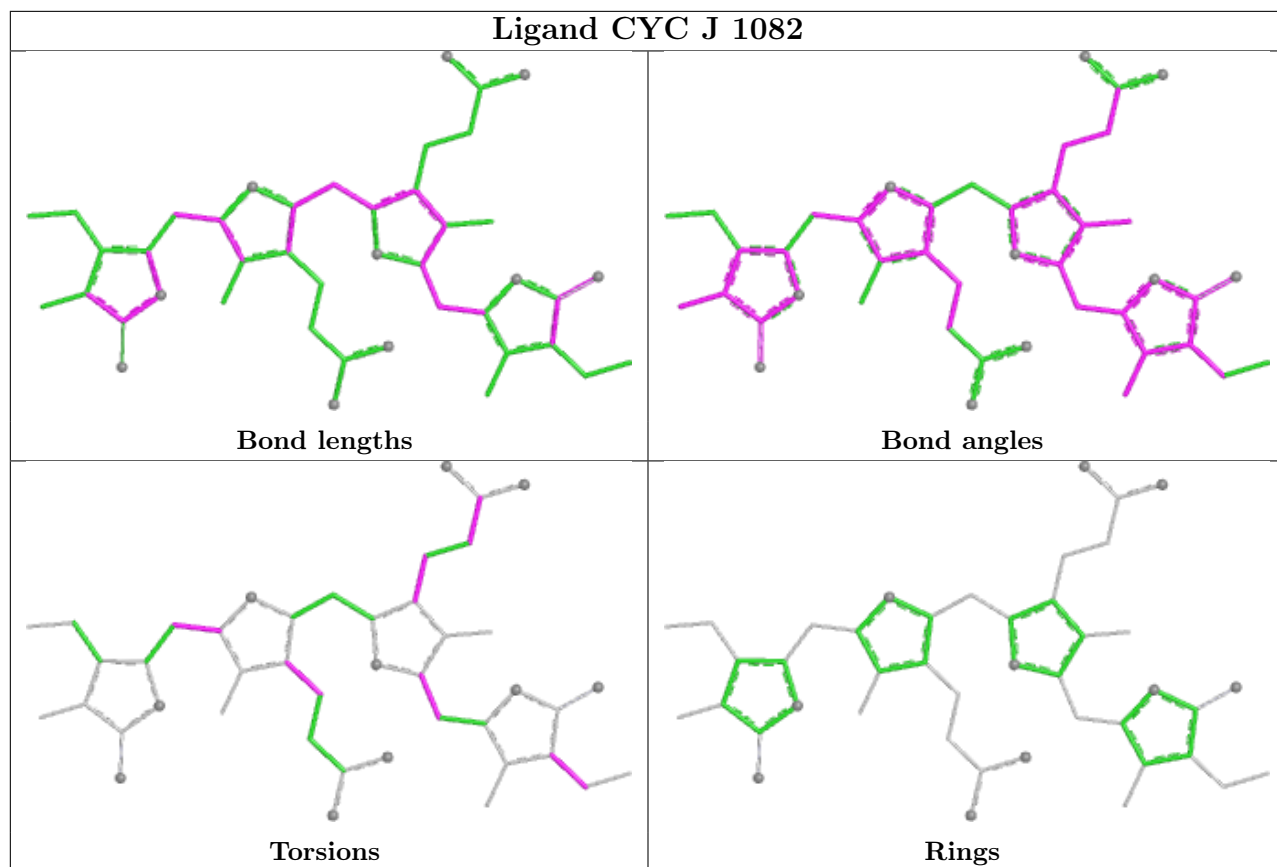


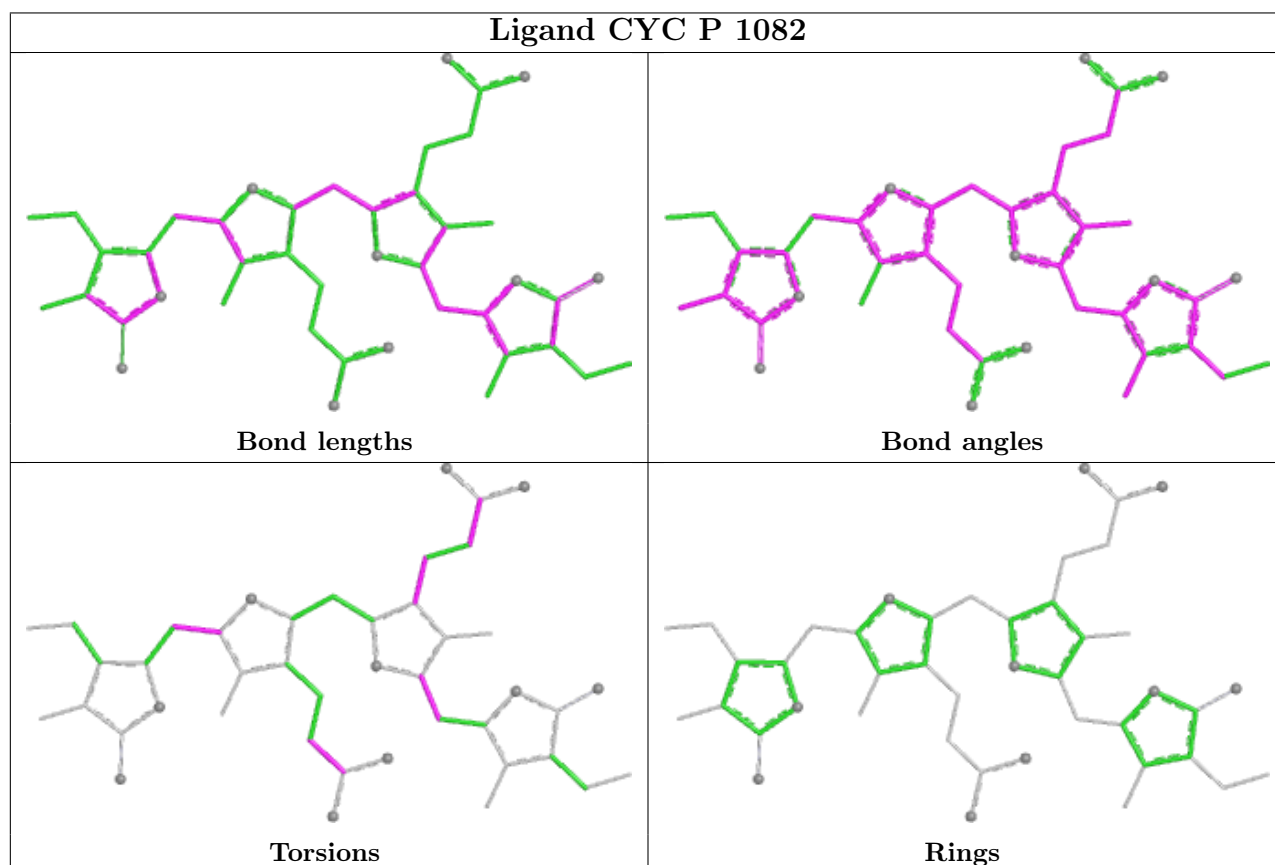
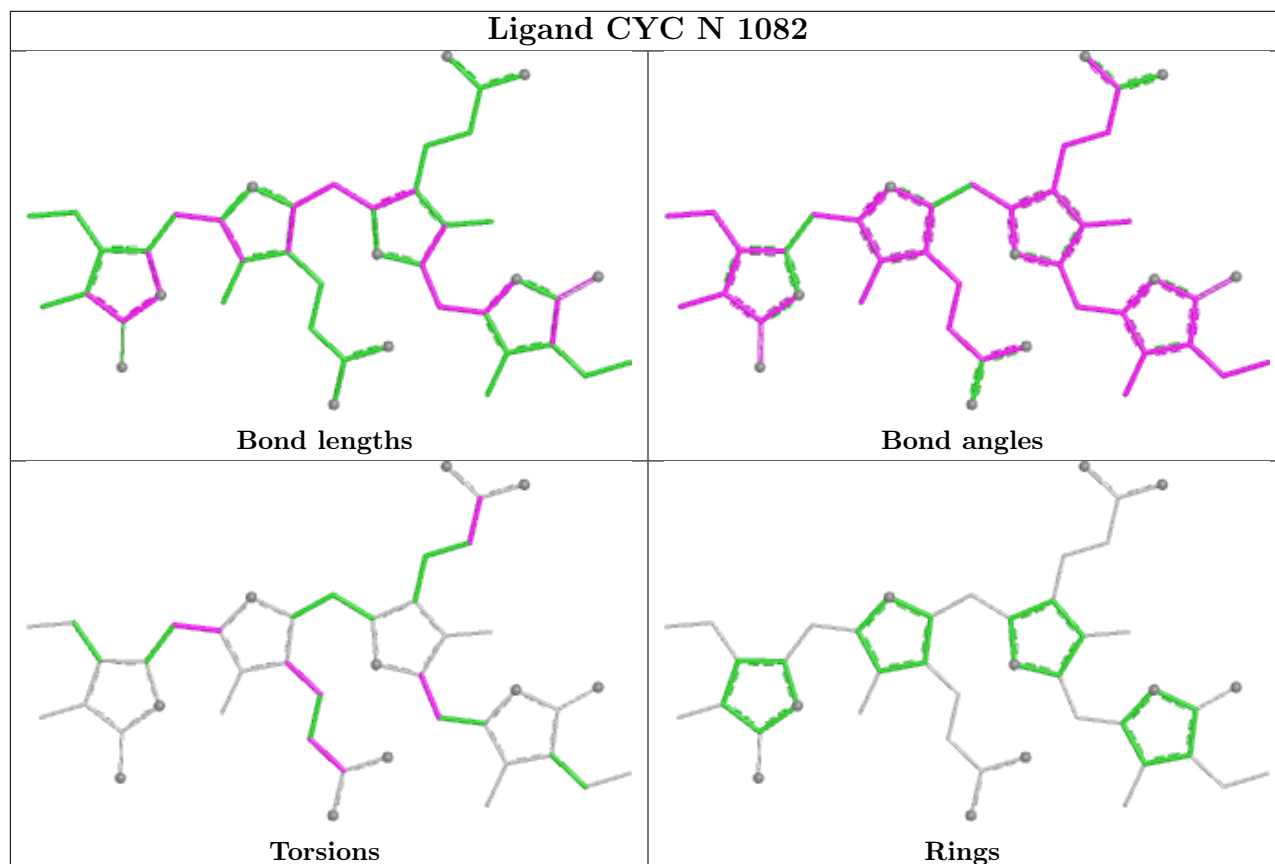


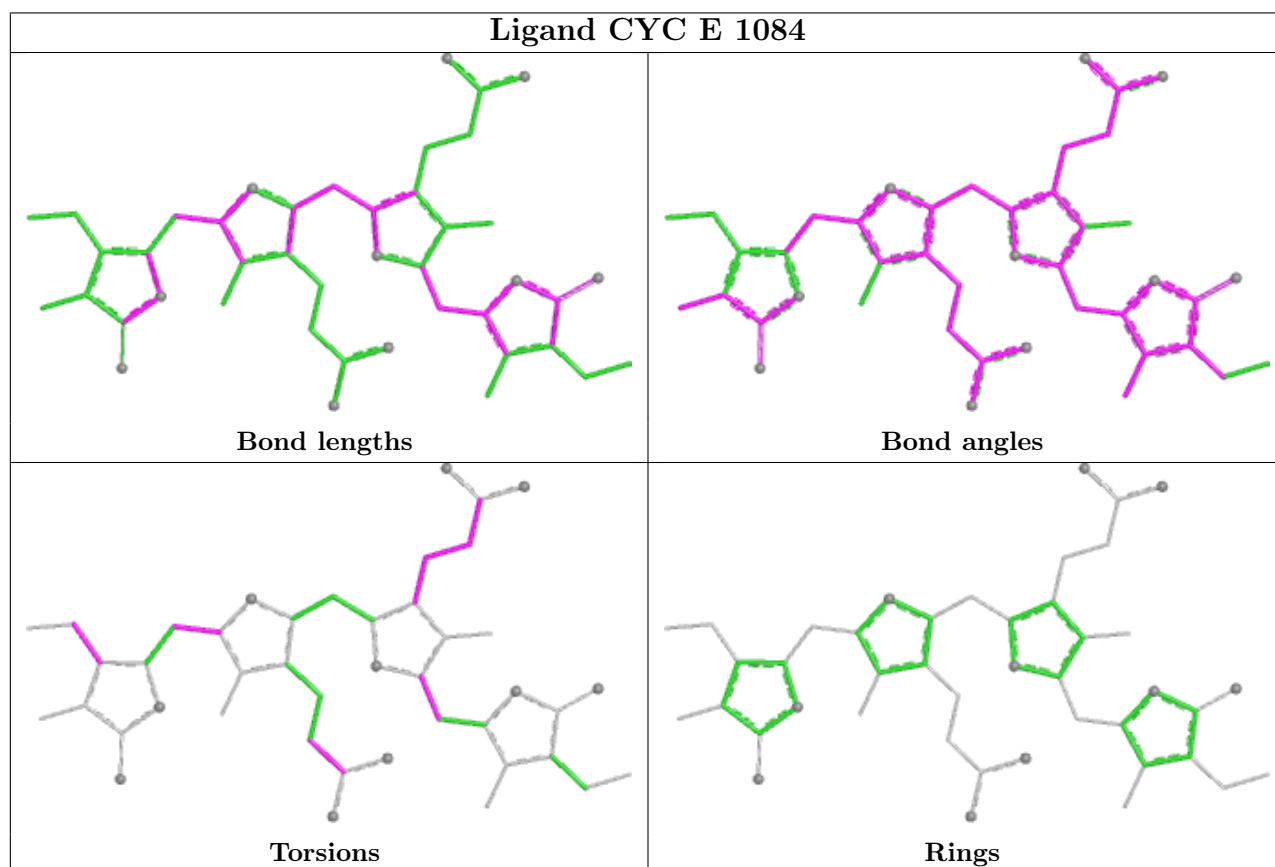
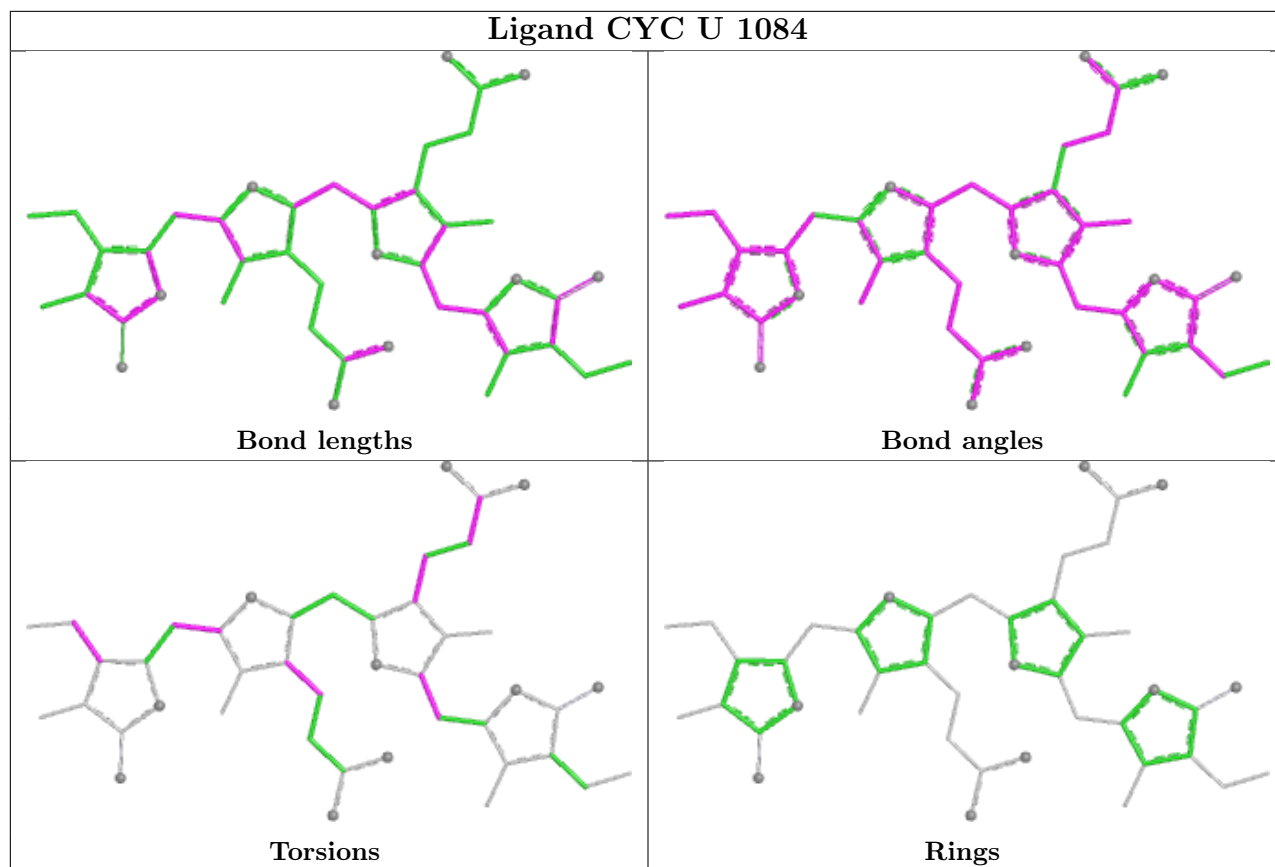


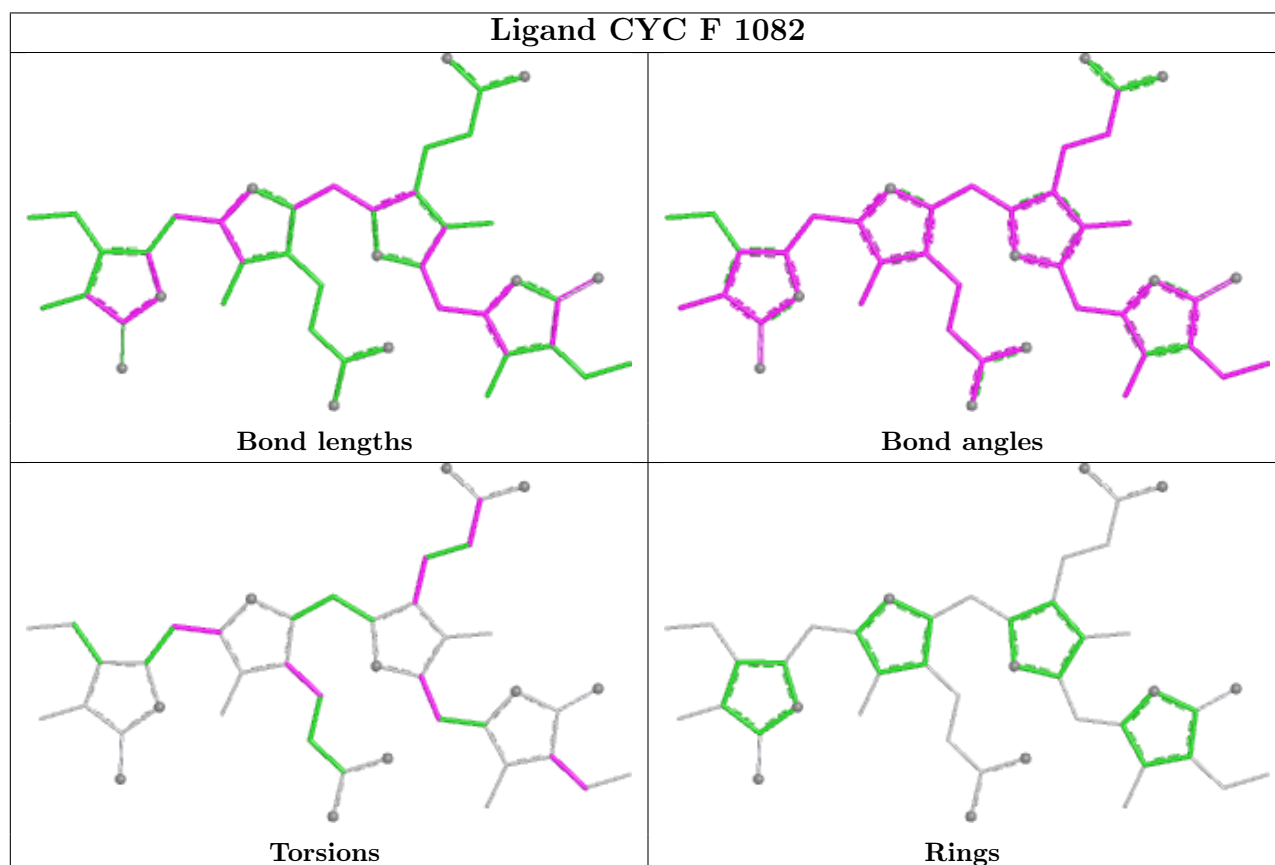
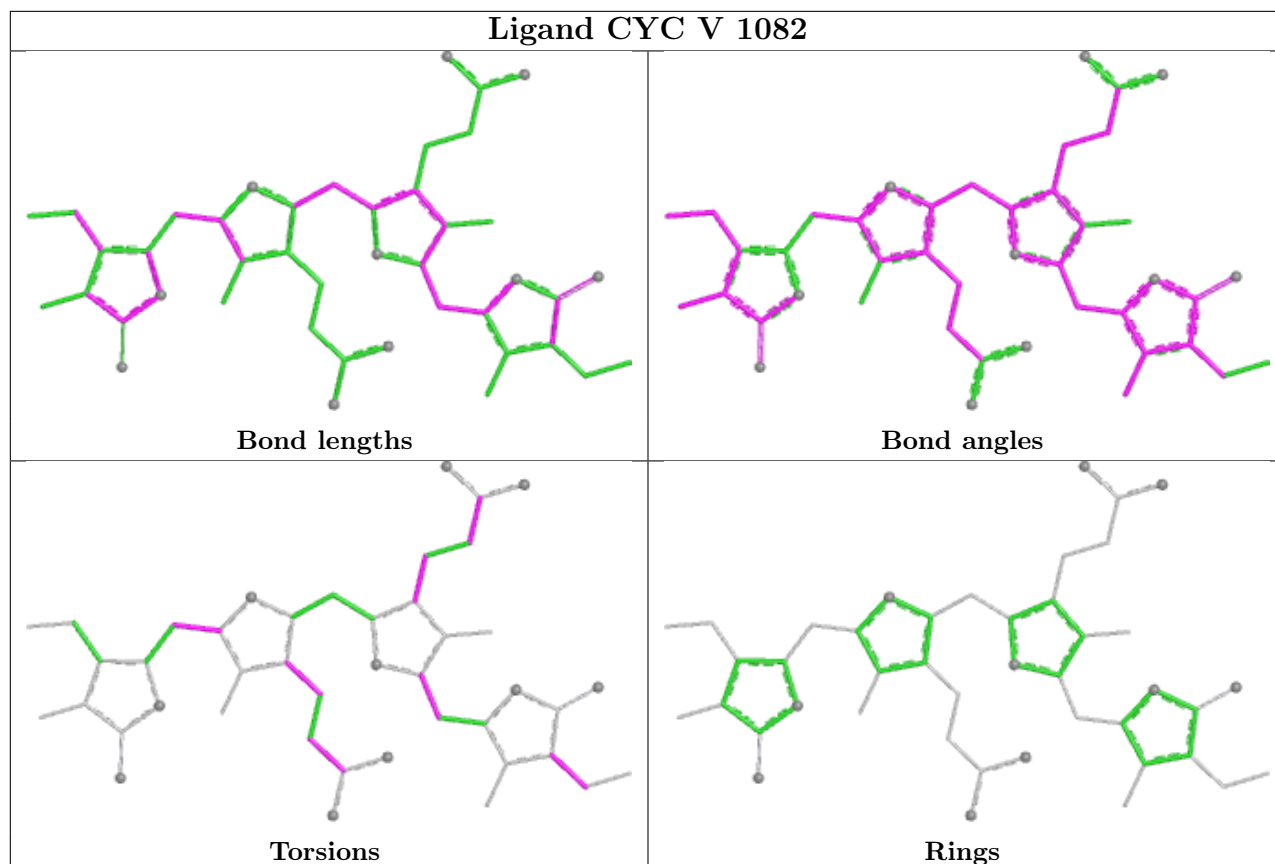


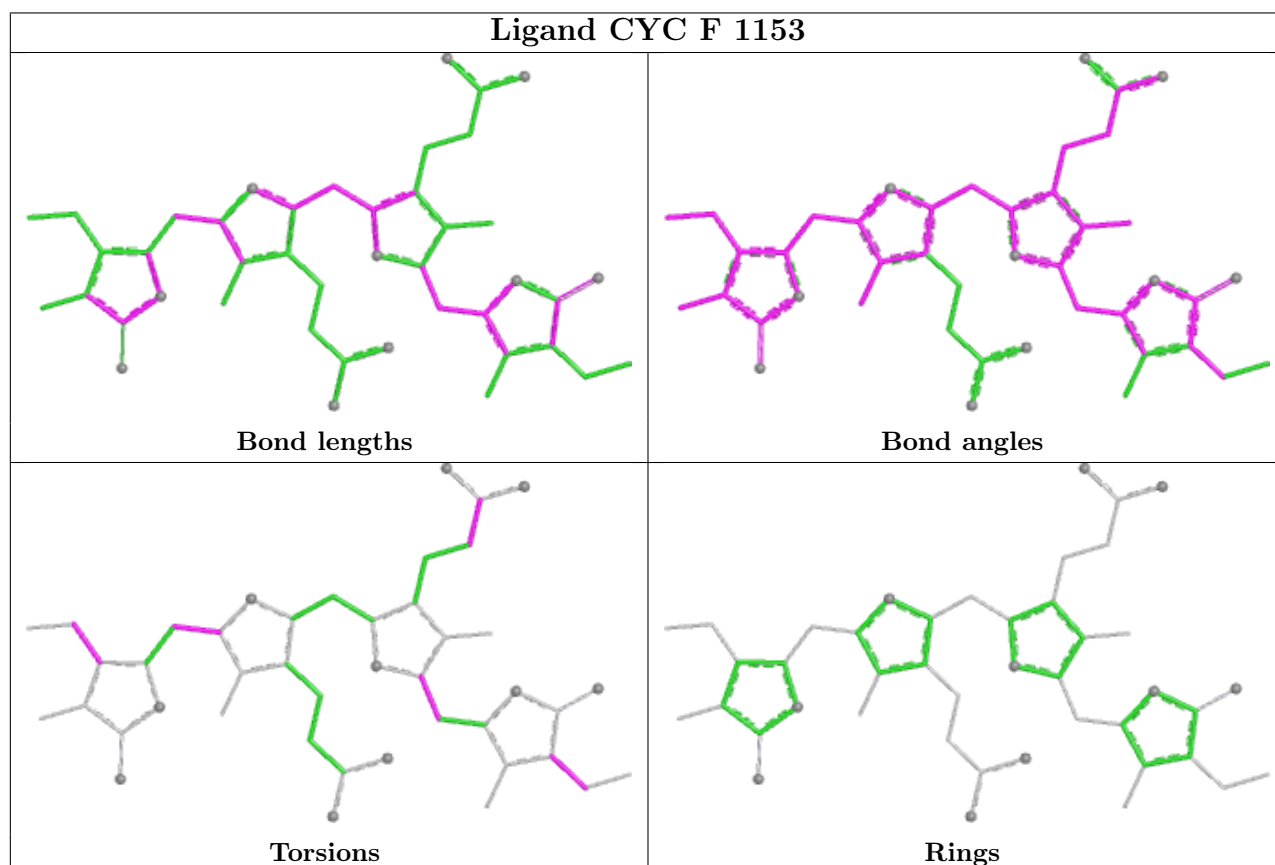
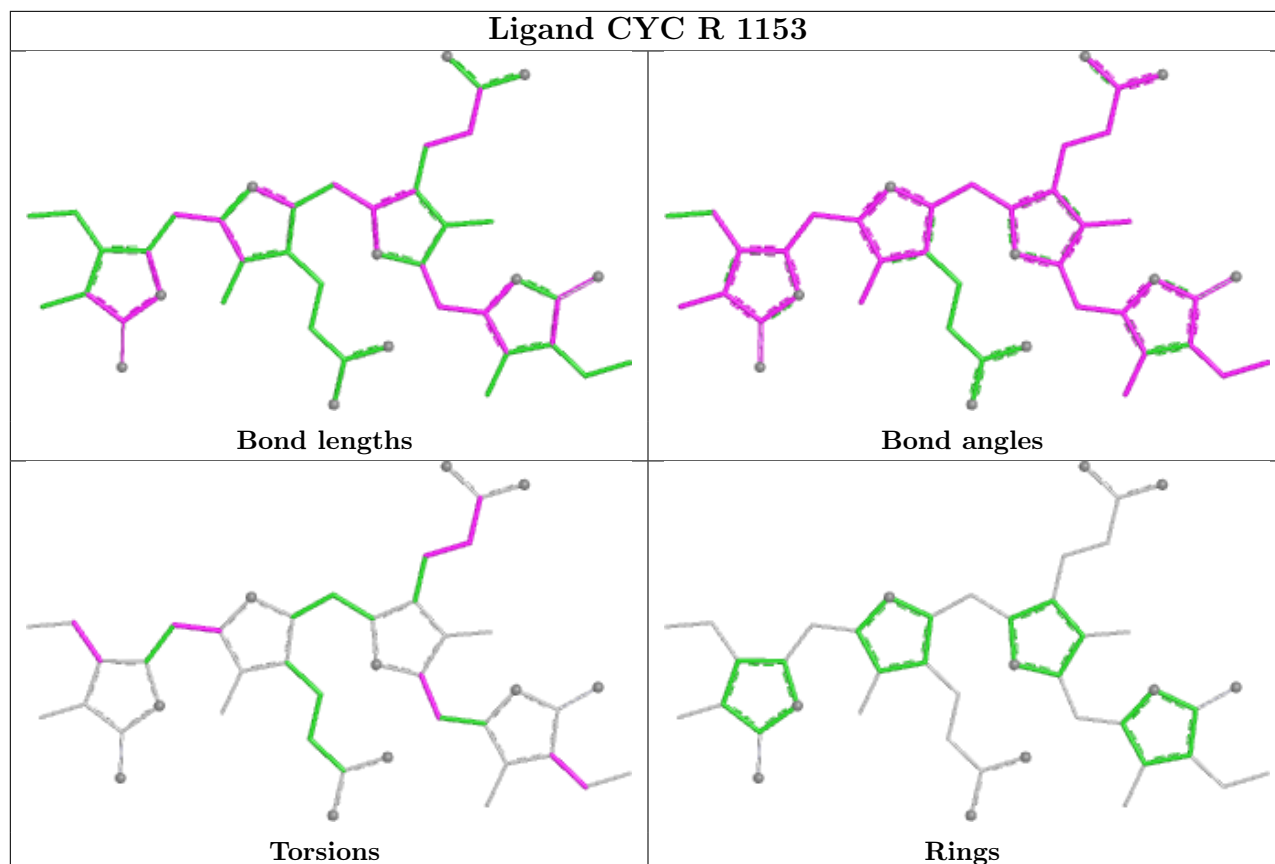


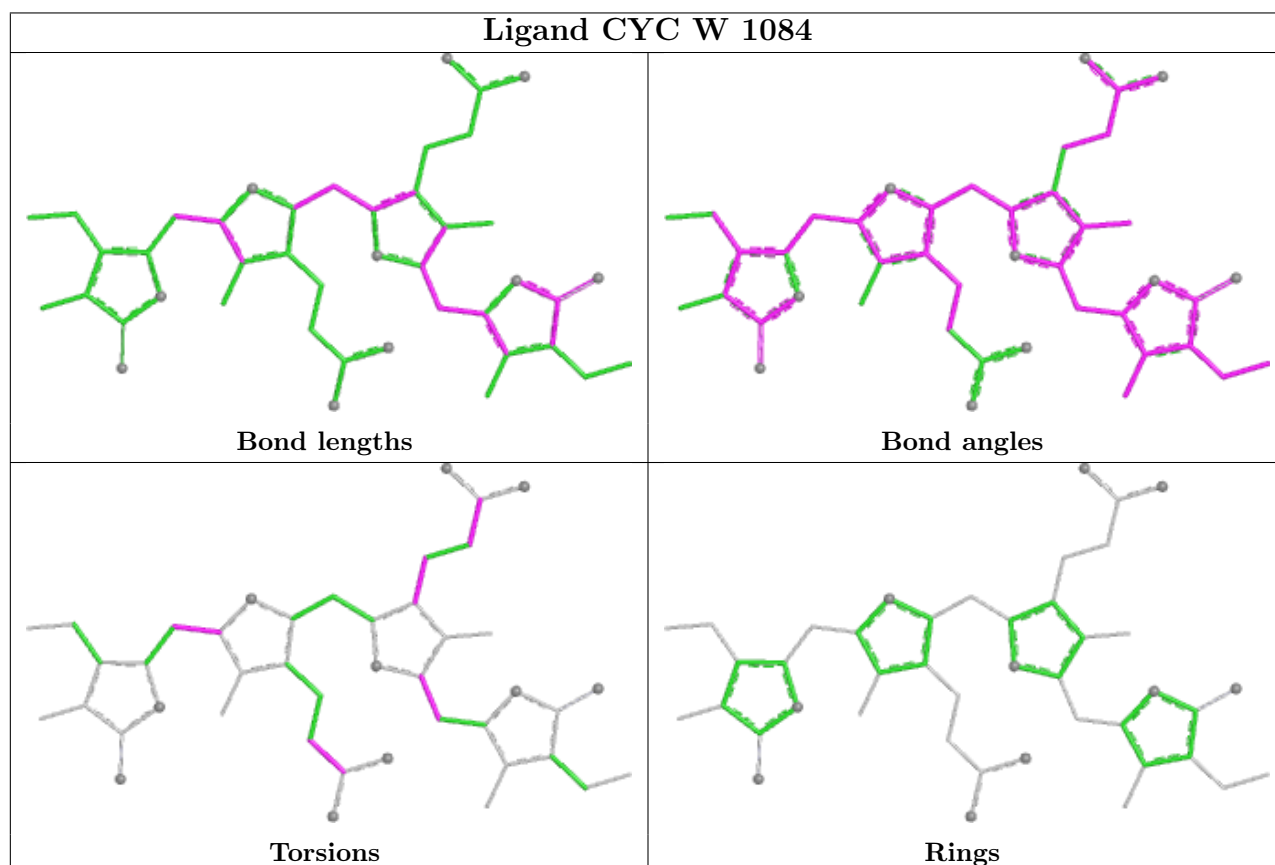
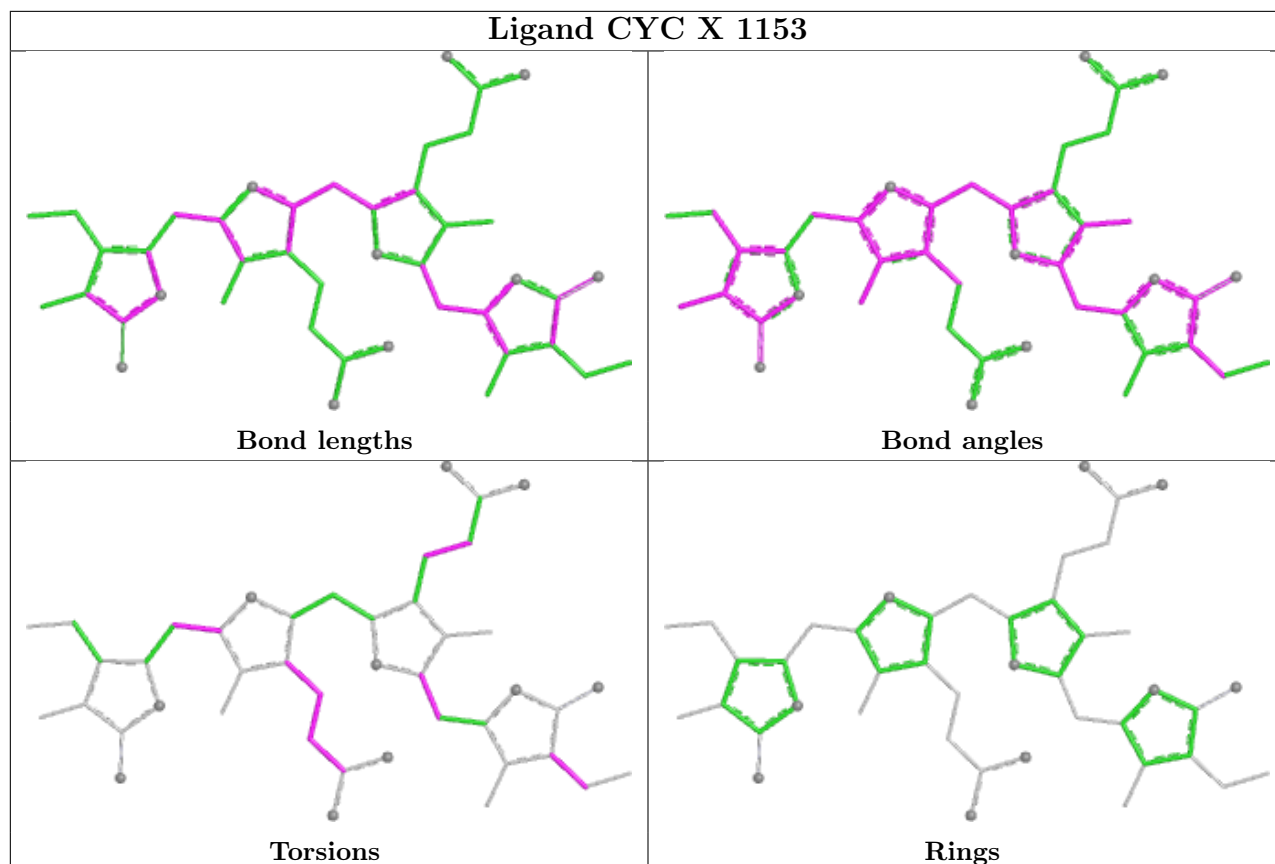


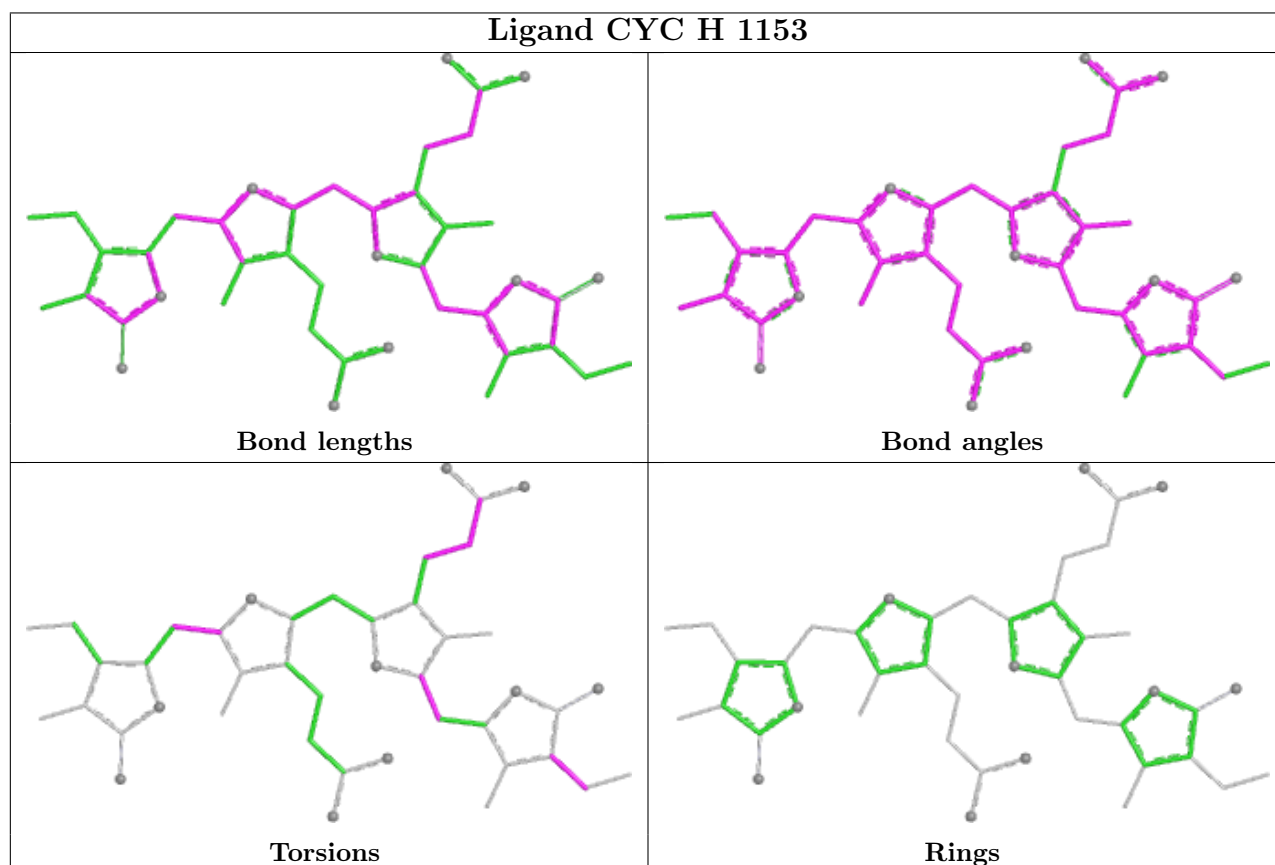
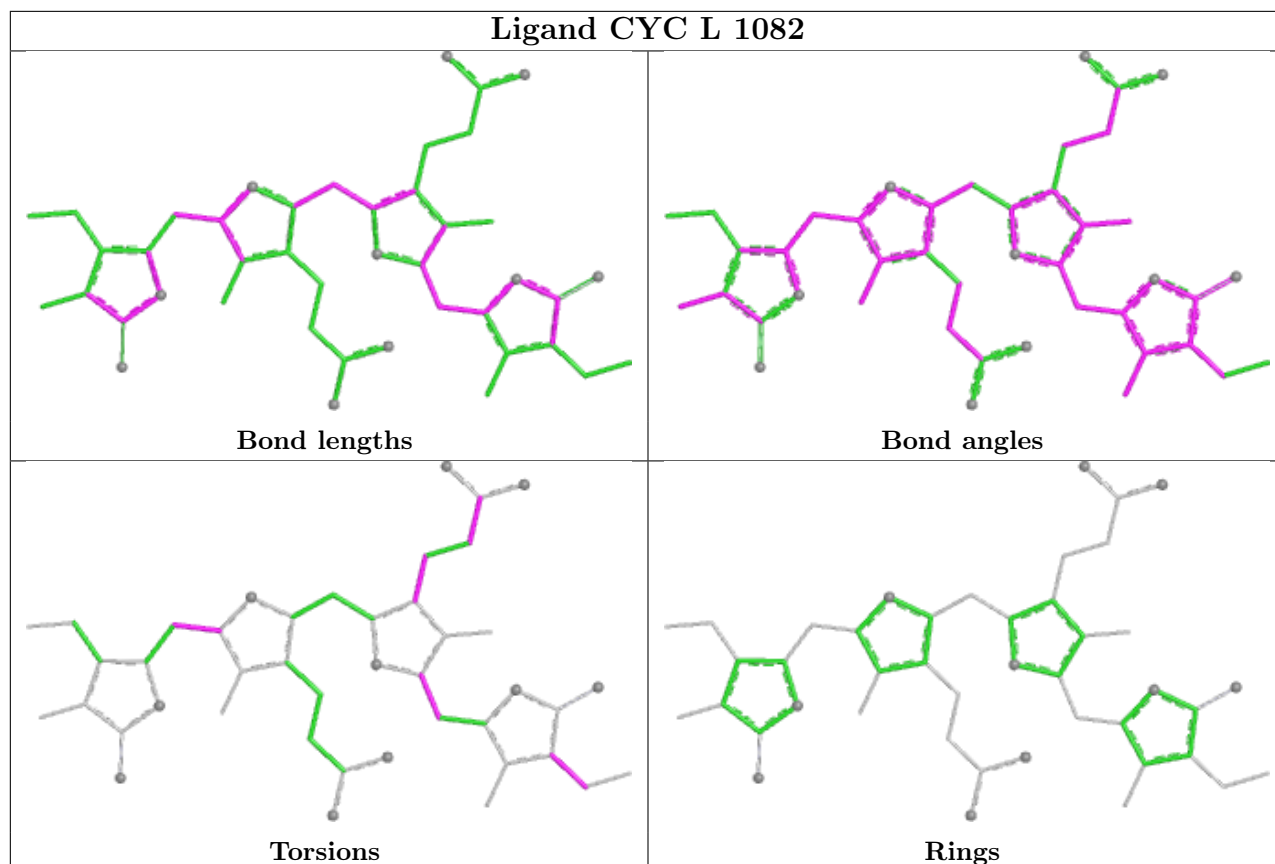


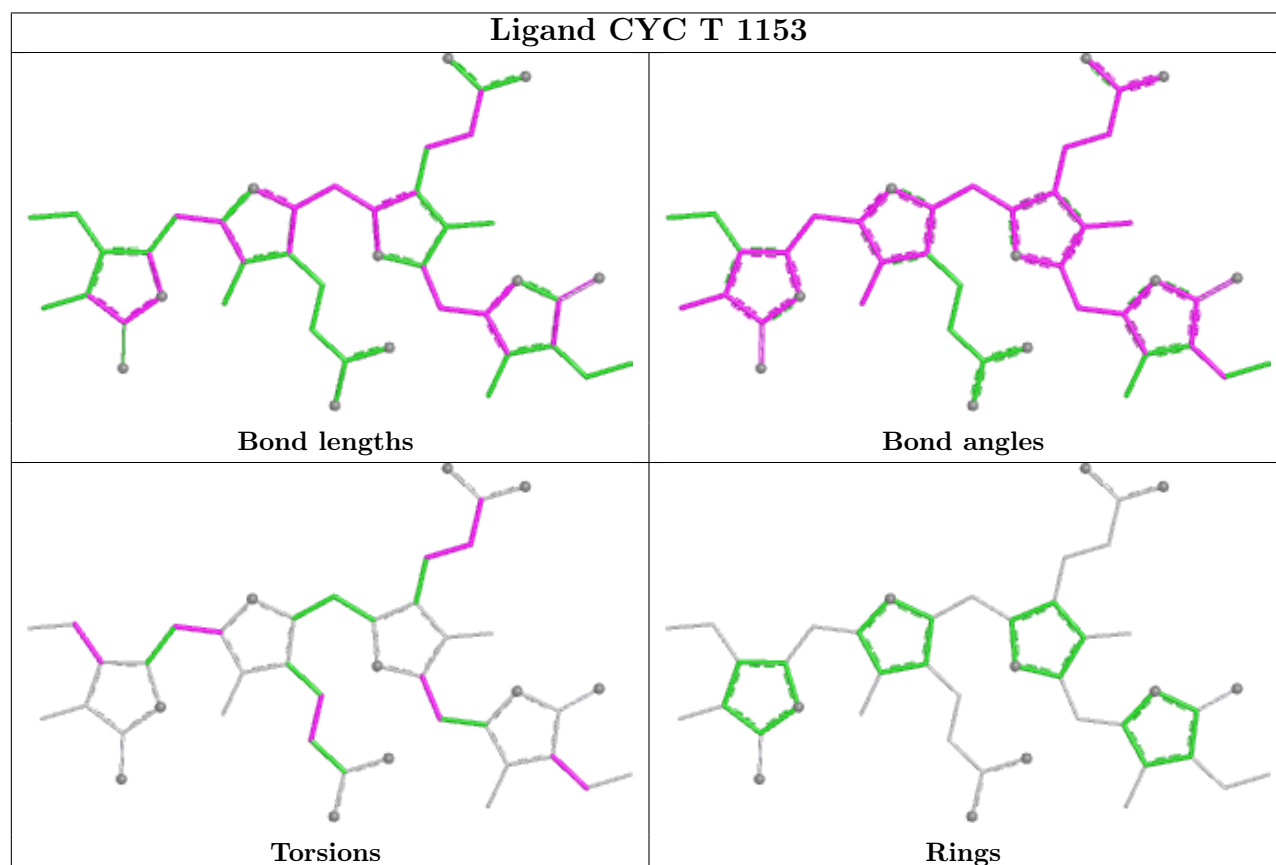
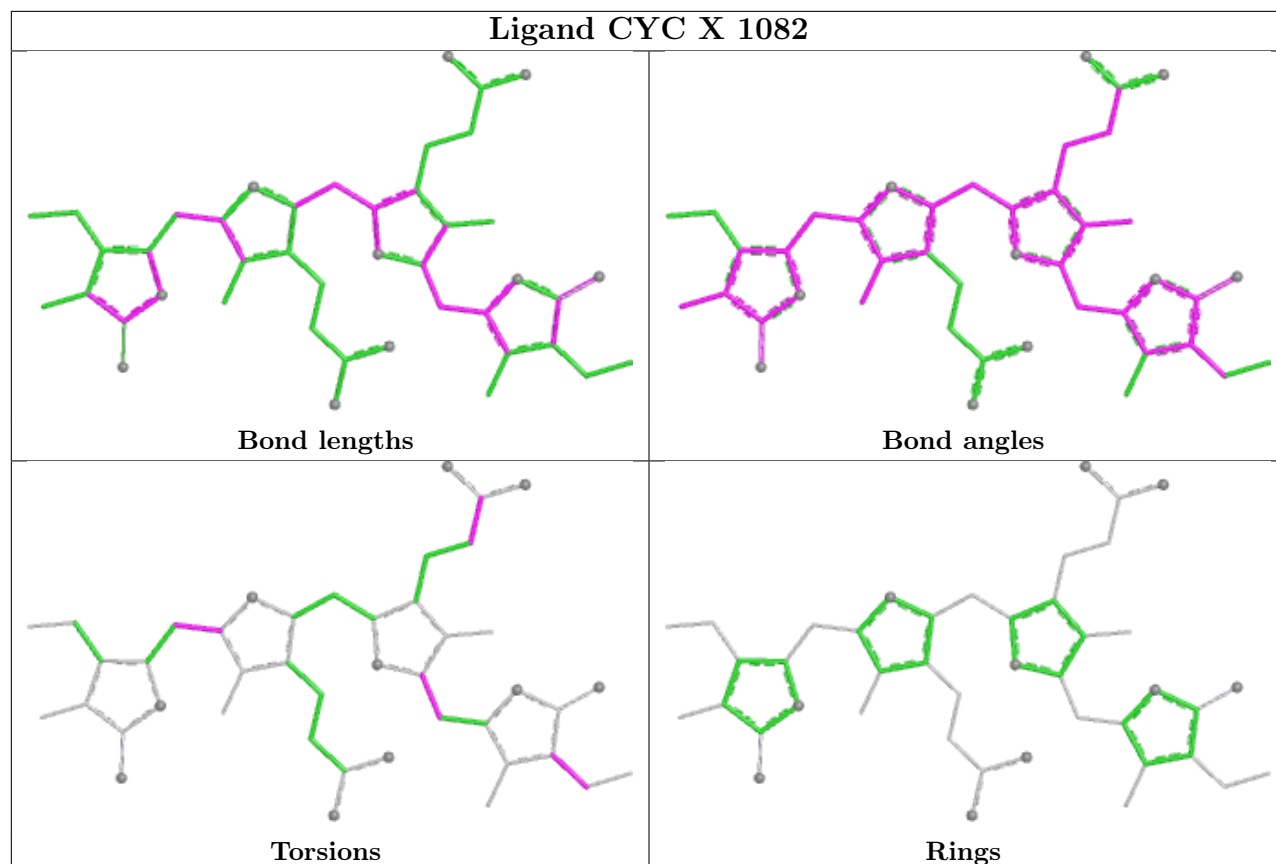


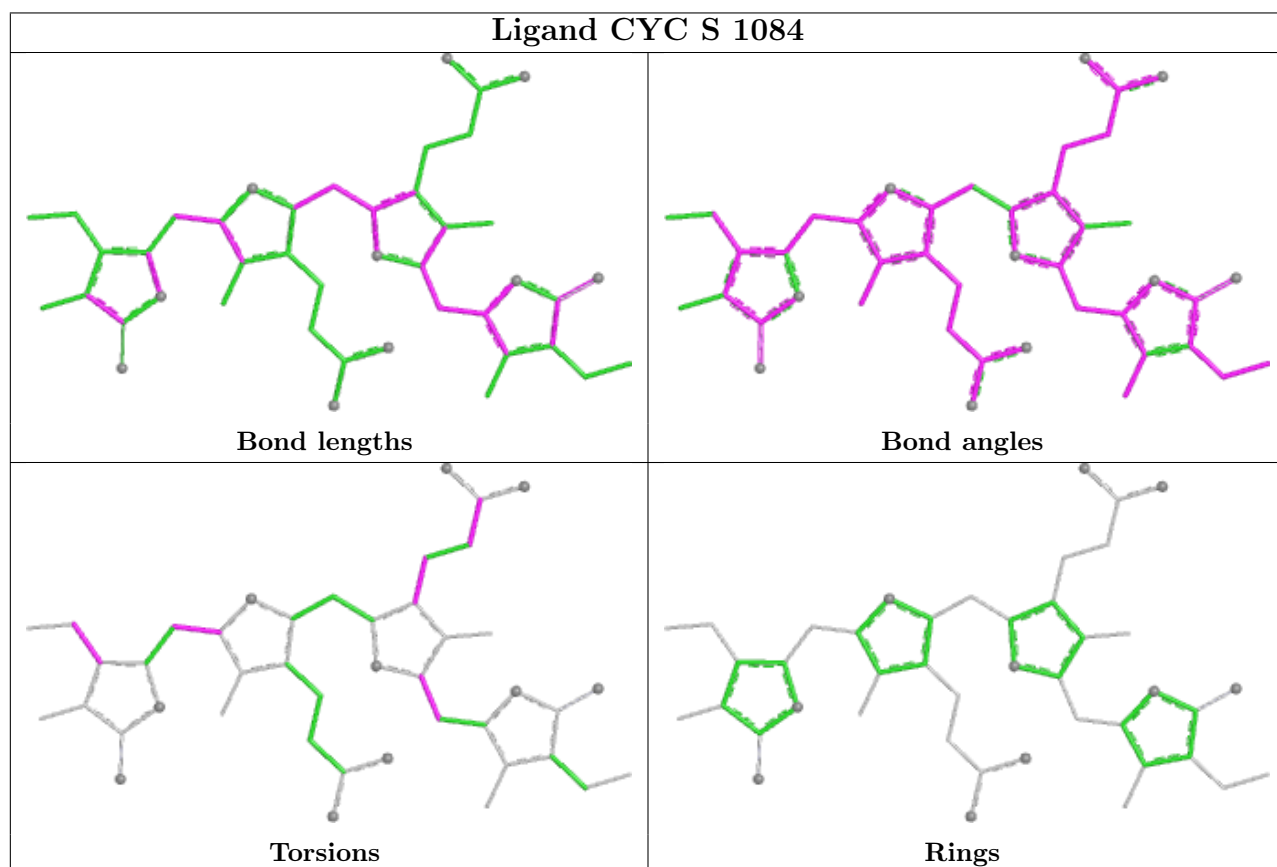
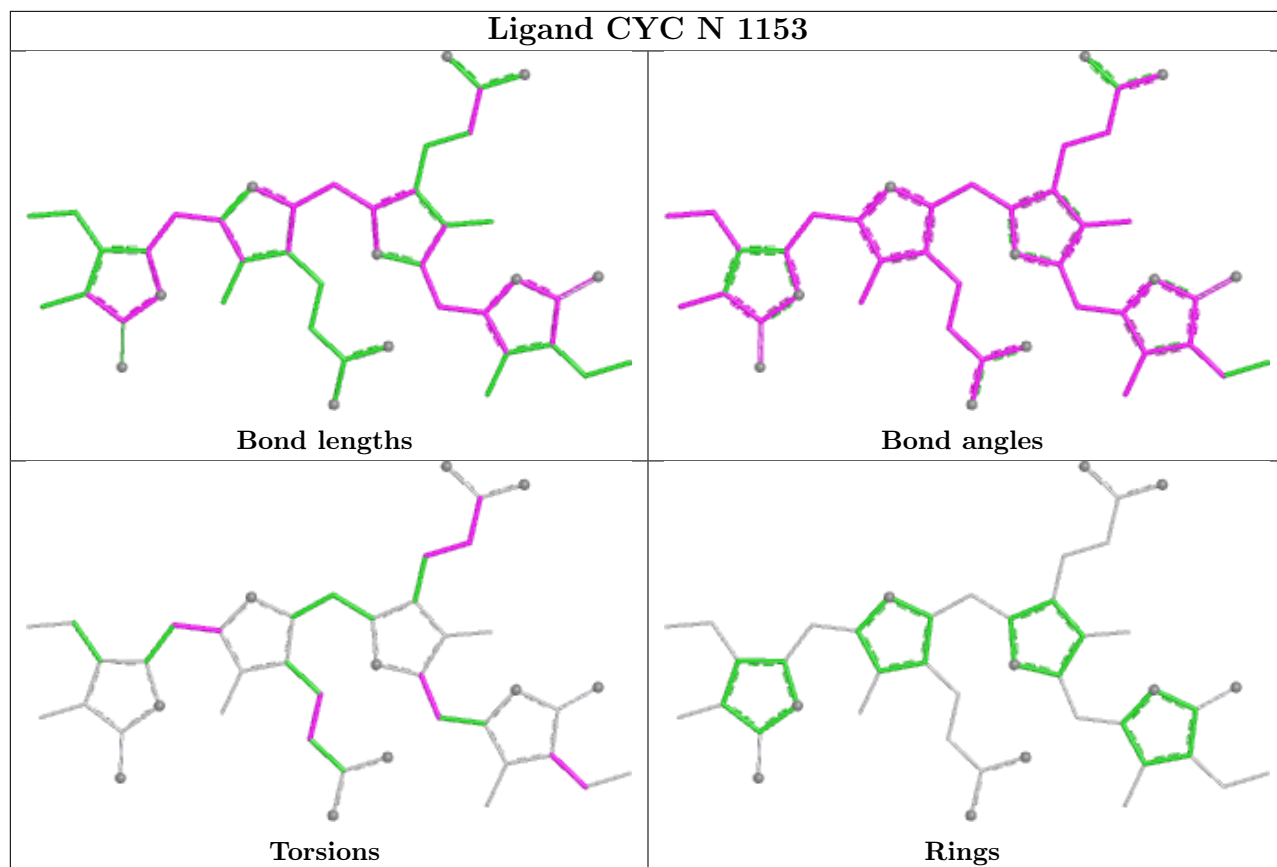


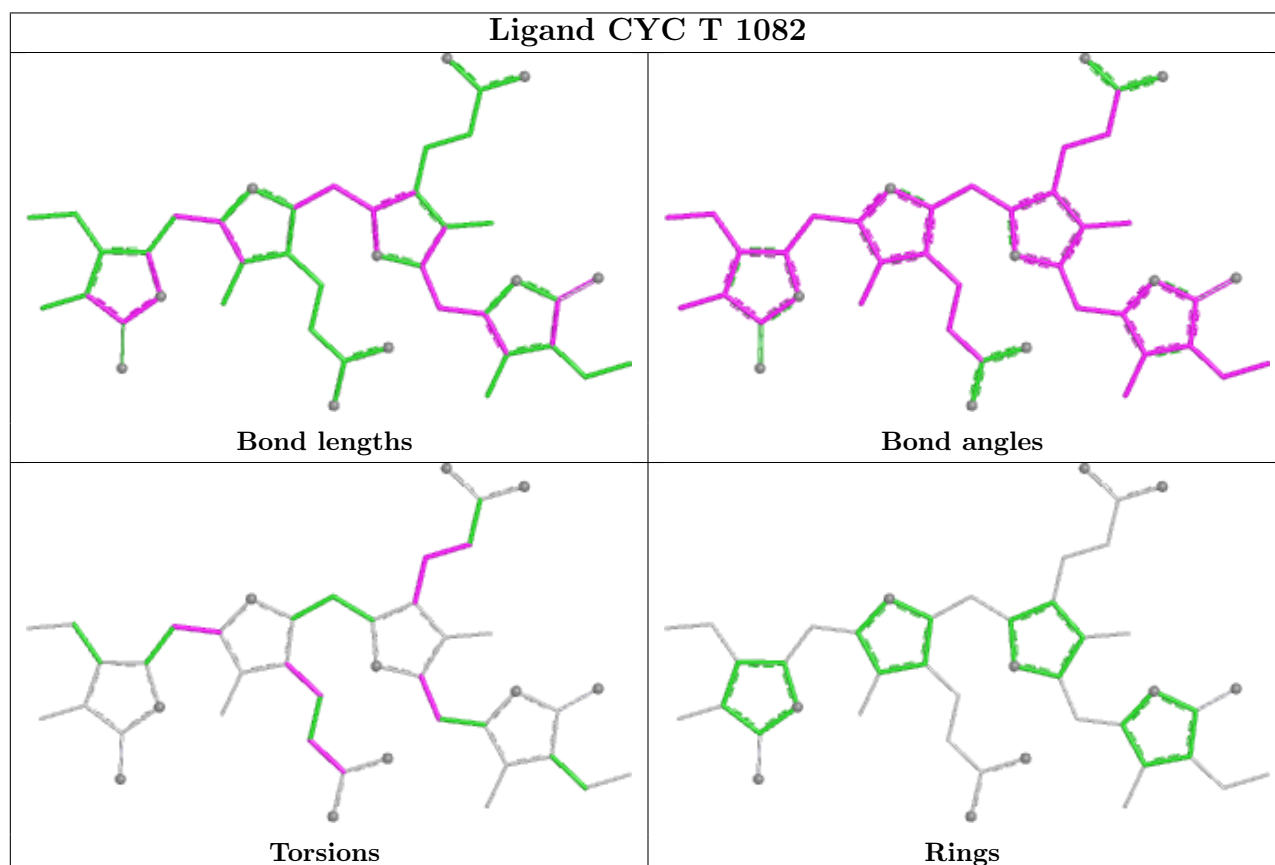
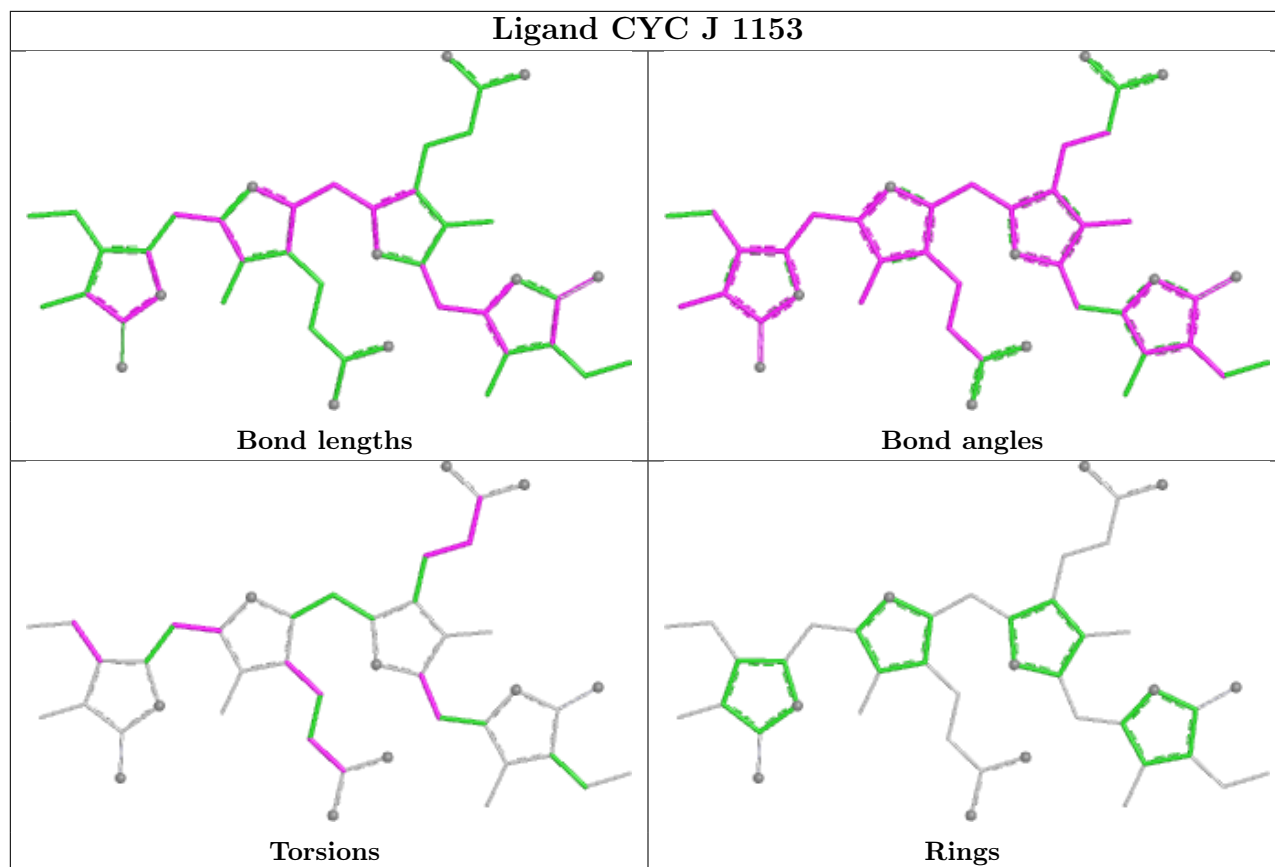


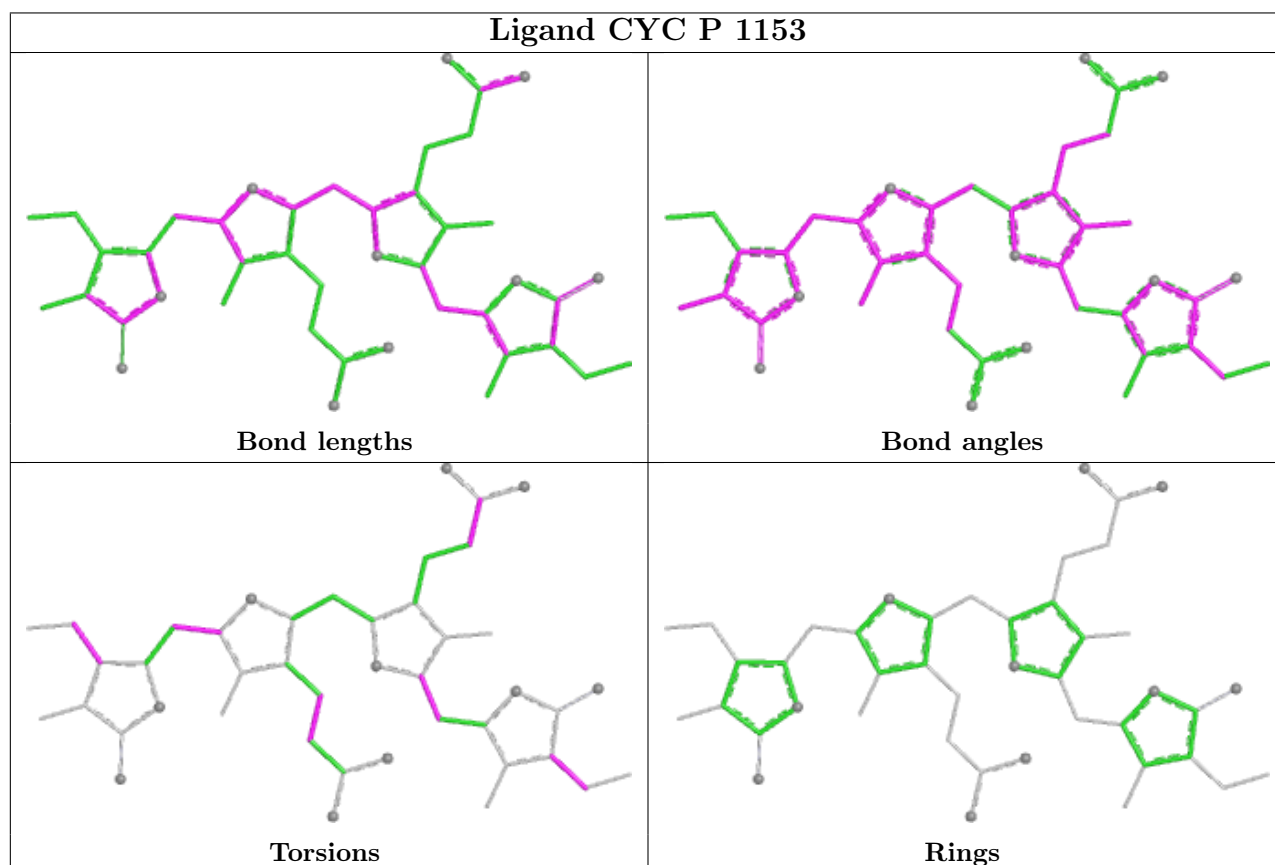
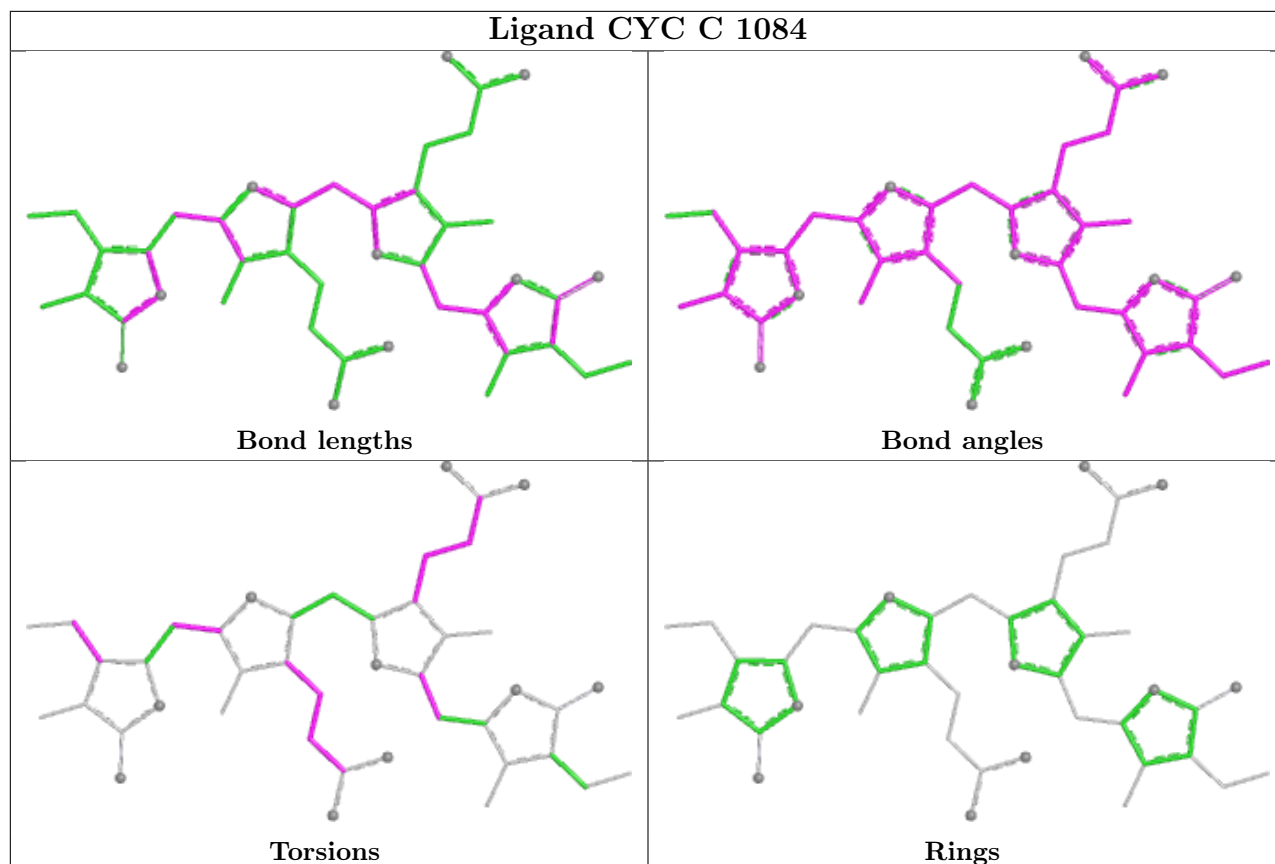


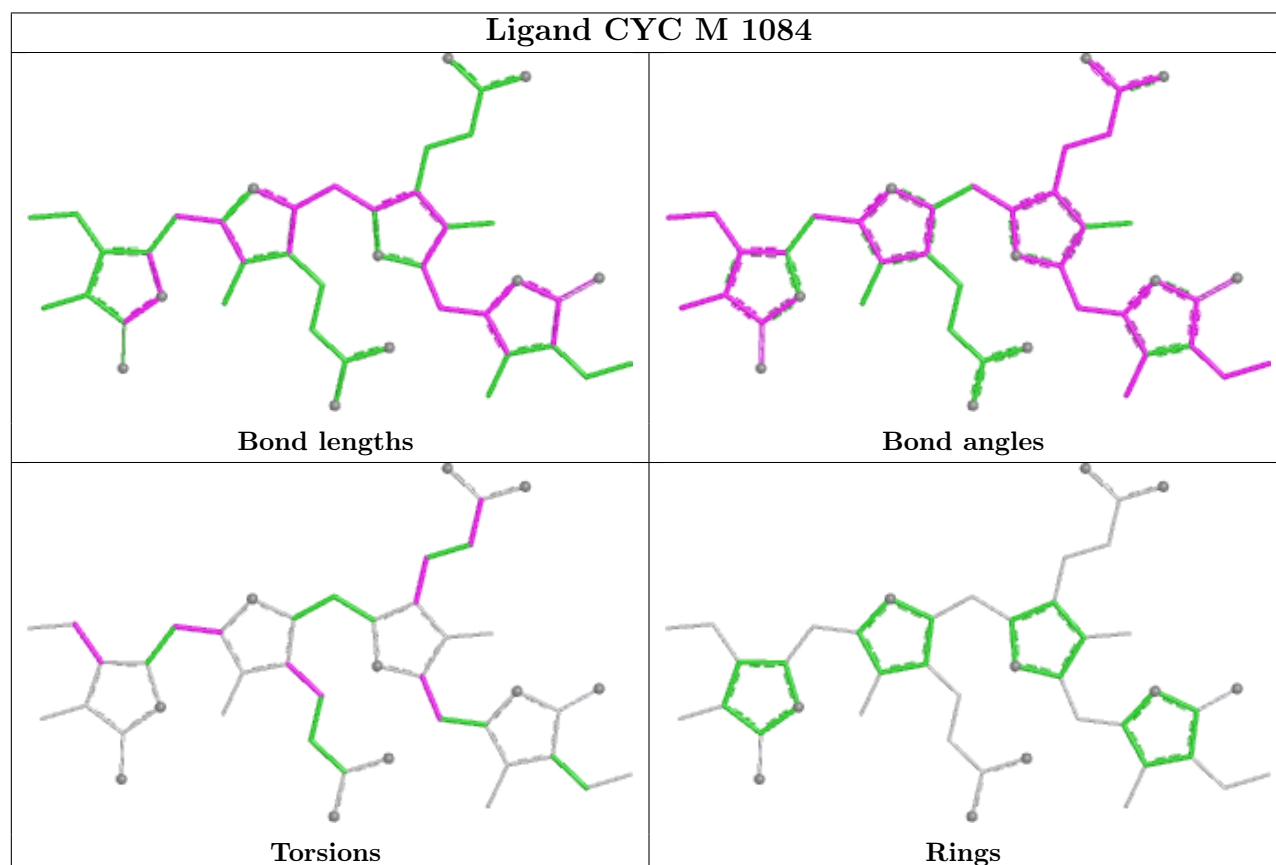
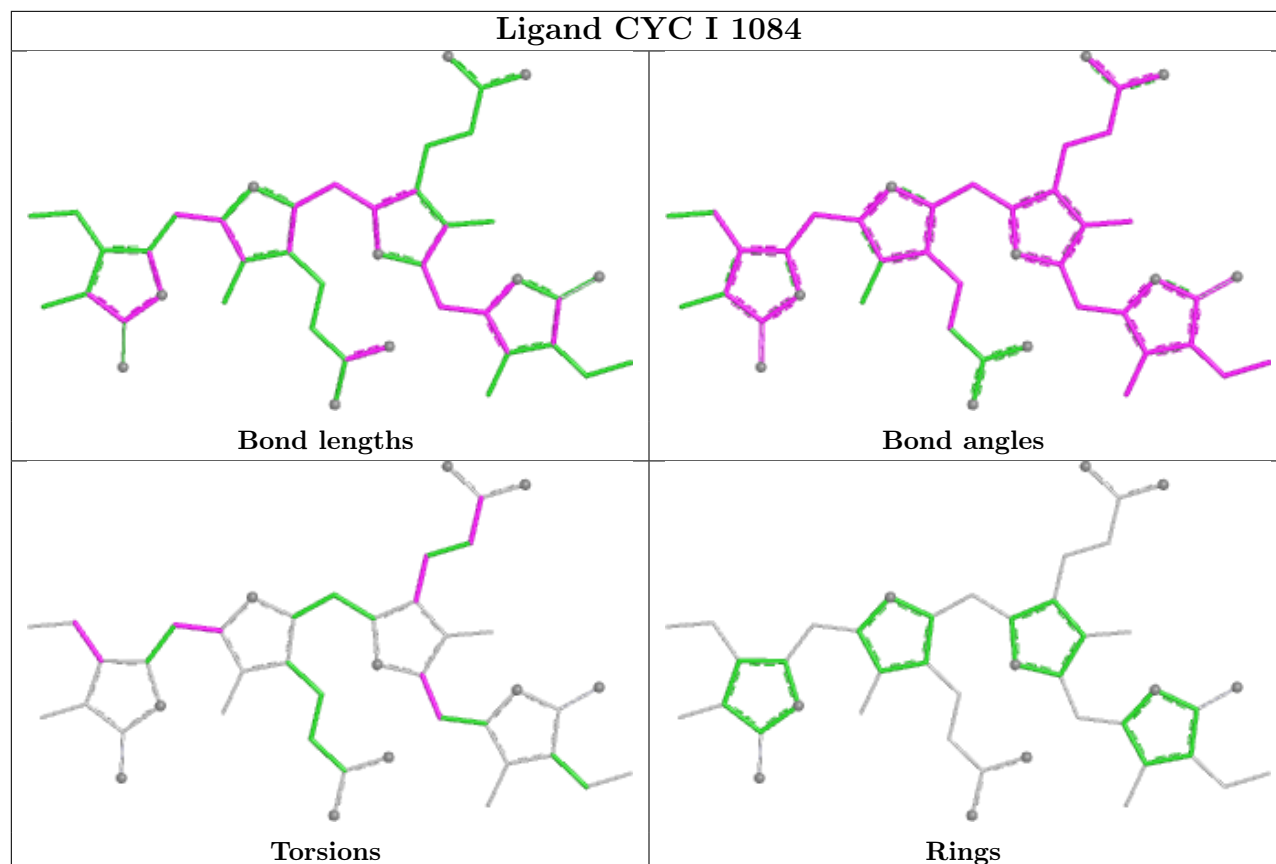


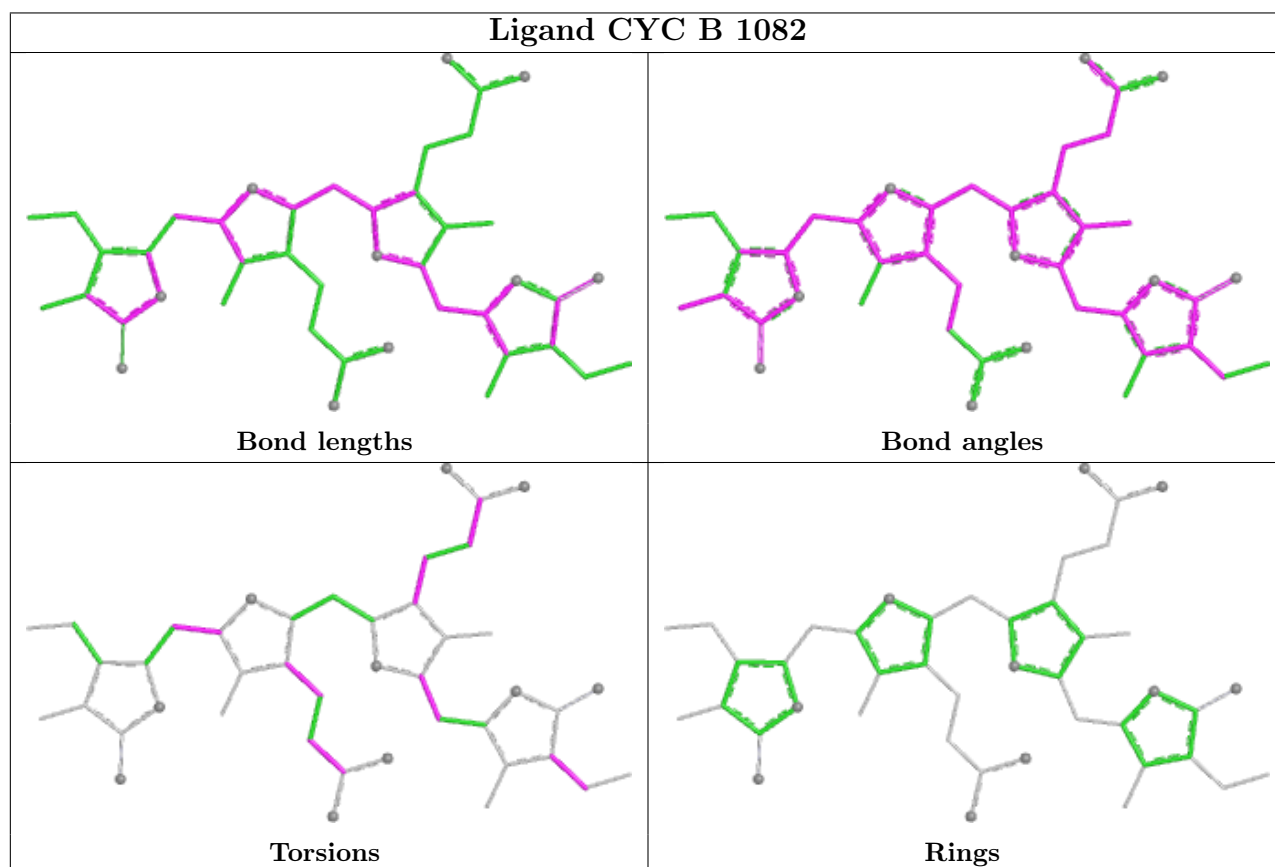
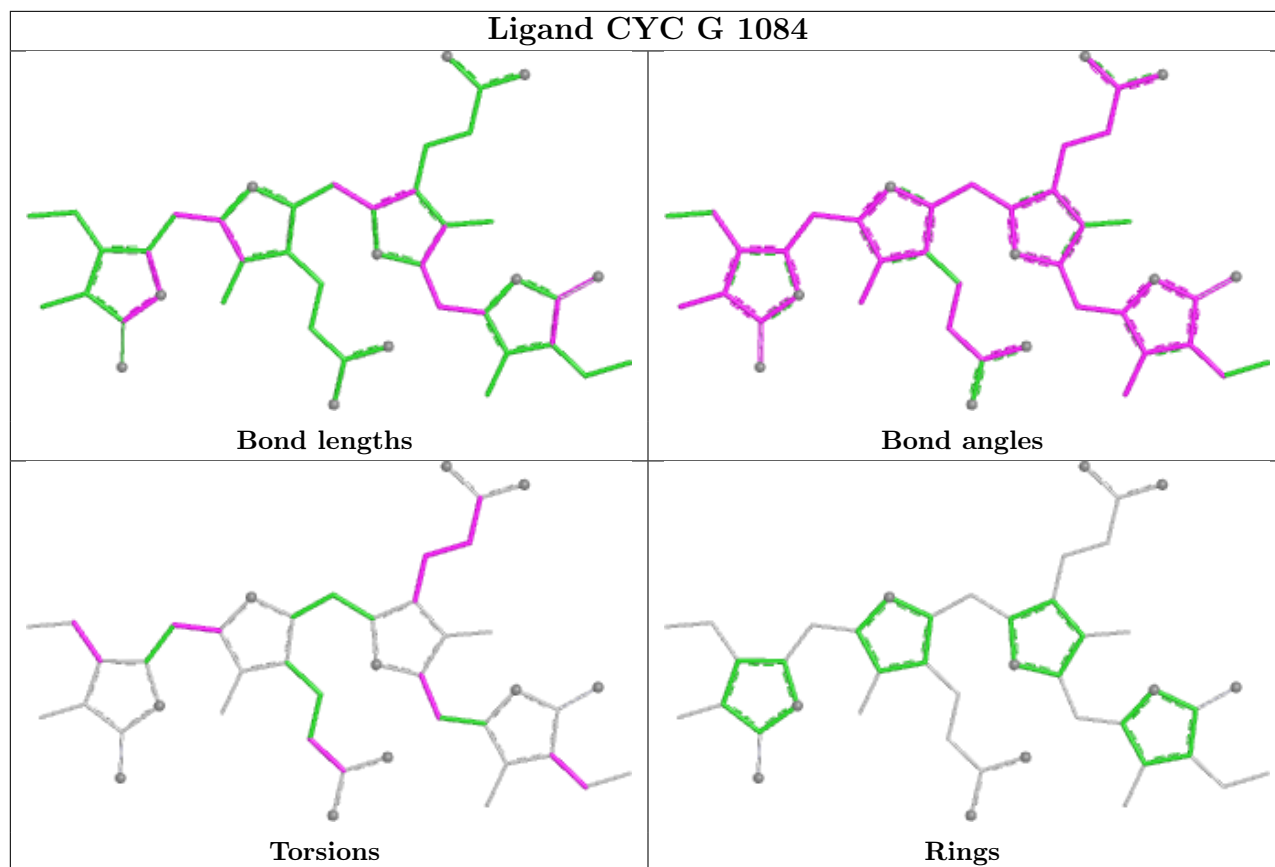


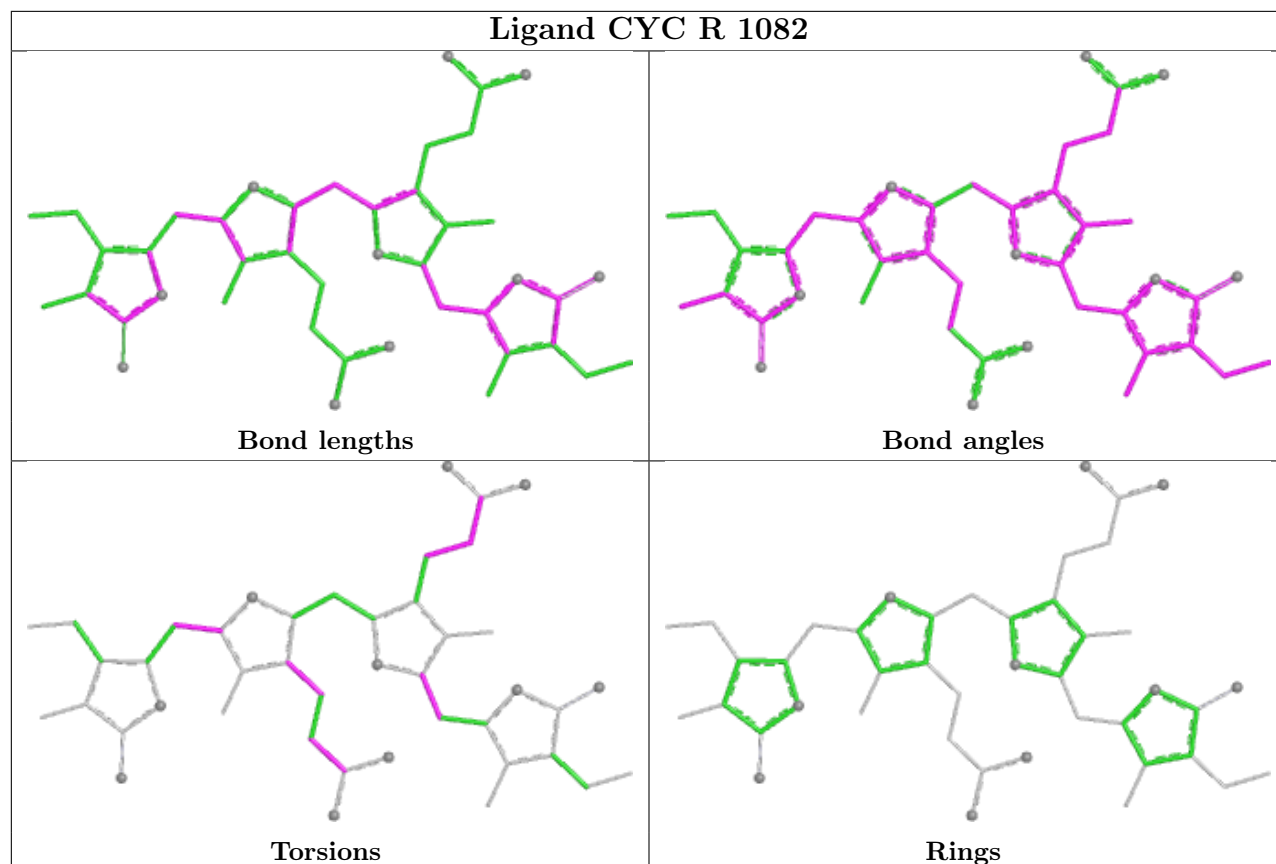












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data i

6.1 Protein, DNA and RNA chains i

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	162/162 (100%)	-0.48	1 (0%) 85 69	16, 24, 32, 42	1 (0%)
1	C	162/162 (100%)	-0.53	0 100 100	16, 24, 32, 42	0
1	E	162/162 (100%)	-0.50	0 100 100	16, 23, 32, 42	0
1	G	162/162 (100%)	-0.58	0 100 100	16, 24, 32, 42	0
1	I	162/162 (100%)	-0.55	0 100 100	16, 24, 32, 42	0
1	K	162/162 (100%)	-0.51	1 (0%) 85 69	16, 24, 32, 42	0
1	M	162/162 (100%)	-0.44	0 100 100	16, 24, 32, 42	0
1	O	162/162 (100%)	-0.49	1 (0%) 85 69	16, 24, 32, 42	0
1	Q	162/162 (100%)	-0.54	1 (0%) 85 69	16, 24, 32, 42	0
1	S	162/162 (100%)	-0.52	1 (0%) 85 69	16, 24, 32, 42	0
1	U	162/162 (100%)	-0.53	0 100 100	16, 24, 32, 42	0
1	W	162/162 (100%)	-0.59	0 100 100	16, 24, 32, 42	0
2	B	172/172 (100%)	-0.62	0 100 100	15, 24, 35, 41	0
2	D	172/172 (100%)	-0.55	0 100 100	15, 24, 35, 41	0
2	F	172/172 (100%)	-0.62	0 100 100	15, 24, 35, 41	0
2	H	172/172 (100%)	-0.49	3 (1%) 69 45	15, 24, 35, 41	0
2	J	172/172 (100%)	-0.50	0 100 100	15, 24, 35, 41	0
2	L	172/172 (100%)	-0.58	0 100 100	15, 24, 35, 41	0
2	N	172/172 (100%)	-0.49	0 100 100	15, 24, 35, 41	0
2	P	172/172 (100%)	-0.59	0 100 100	15, 24, 35, 41	0
2	R	172/172 (100%)	-0.58	1 (0%) 85 69	15, 24, 35, 41	0
2	T	172/172 (100%)	-0.57	0 100 100	15, 24, 35, 41	0
2	V	172/172 (100%)	-0.46	0 100 100	15, 24, 35, 41	0
3	X	172/172 (100%)	-0.54	0 100 100	15, 24, 35, 41	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
All	All	4008/4008 (100%)	-0.54	9 (0%) 91 83	15, 24, 35, 42	1 (0%)

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	68	GLN	4.1
1	O	68	GLN	2.9
2	H	25	ASP	2.7
1	Q	68	GLN	2.5
1	S	68	GLN	2.5
2	R	65	GLN	2.5
2	H	21	THR	2.4
1	K	57	GLN	2.3
2	H	22	ALA	2.3

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	CYC	L	1082	43/43	0.84	0.14	23,30,36,41	0
4	CYC	J	1082	43/43	0.85	0.16	38,45,49,50	0
4	CYC	H	1082	43/43	0.85	0.14	23,32,38,44	0
4	CYC	P	1082	43/43	0.85	0.15	45,52,54,57	0
4	CYC	R	1082	43/43	0.85	0.15	31,38,42,50	0
4	CYC	D	1082	43/43	0.86	0.15	32,41,43,44	0
4	CYC	F	1082	43/43	0.86	0.14	35,40,46,51	0
4	CYC	B	1082	43/43	0.86	0.14	25,32,39,41	0

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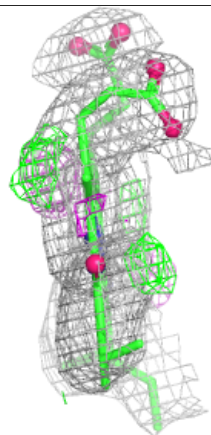
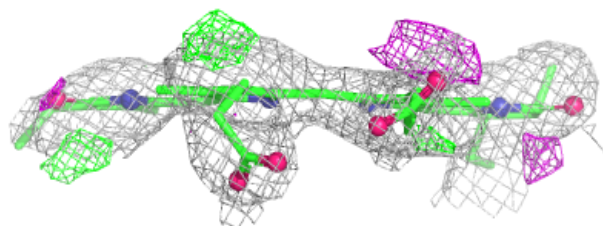
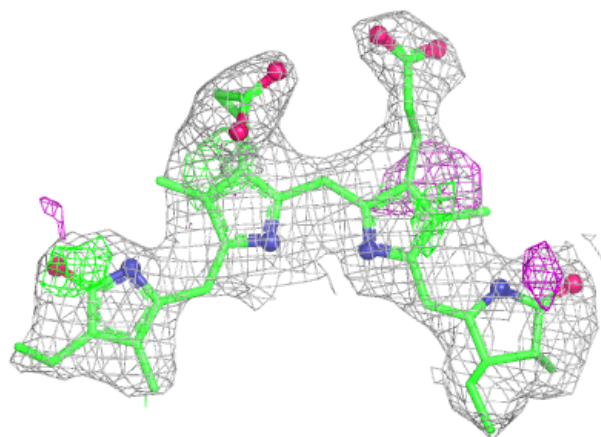
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	CYC	P	1153	43/43	0.86	0.15	29,43,44,45	0
4	CYC	H	1153	43/43	0.86	0.14	24,33,35,37	0
4	CYC	V	1082	43/43	0.86	0.15	33,39,47,54	0
4	CYC	X	1082	43/43	0.86	0.15	36,45,49,52	0
4	CYC	X	1153	43/43	0.86	0.14	34,45,45,46	0
4	CYC	T	1082	43/43	0.87	0.14	30,38,43,47	0
4	CYC	J	1153	43/43	0.87	0.16	34,51,51,54	0
4	CYC	B	1153	43/43	0.87	0.13	19,26,27,29	0
4	CYC	L	1153	43/43	0.87	0.15	17,31,31,33	0
4	CYC	T	1153	43/43	0.88	0.14	19,42,43,43	0
4	CYC	F	1153	43/43	0.88	0.13	18,34,34,35	0
4	CYC	V	1153	43/43	0.88	0.13	18,31,32,32	0
4	CYC	W	1084	43/43	0.88	0.13	26,33,33,34	0
4	CYC	M	1084	43/43	0.88	0.13	36,42,44,48	0
4	CYC	D	1153	43/43	0.88	0.14	26,34,35,38	0
4	CYC	E	1084	43/43	0.89	0.12	12,21,22,23	0
4	CYC	N	1082	43/43	0.89	0.12	28,31,36,39	0
4	CYC	R	1153	43/43	0.89	0.12	18,31,32,32	0
4	CYC	N	1153	43/43	0.89	0.12	16,31,31,32	0
4	CYC	G	1084	43/43	0.89	0.12	21,27,31,33	0
4	CYC	U	1084	43/43	0.90	0.12	27,29,30,31	0
4	CYC	I	1084	43/43	0.90	0.12	16,23,24,29	0
4	CYC	S	1084	43/43	0.90	0.12	13,23,24,28	0
4	CYC	O	1084	43/43	0.91	0.12	13,23,24,25	0
4	CYC	K	1084	43/43	0.91	0.11	10,22,22,23	0
4	CYC	C	1084	43/43	0.91	0.11	23,29,29,31	0
4	CYC	Q	1084	43/43	0.91	0.12	14,22,24,25	0
4	CYC	A	1084	43/43	0.91	0.11	15,21,21,26	0

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

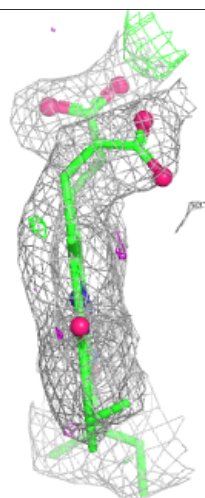
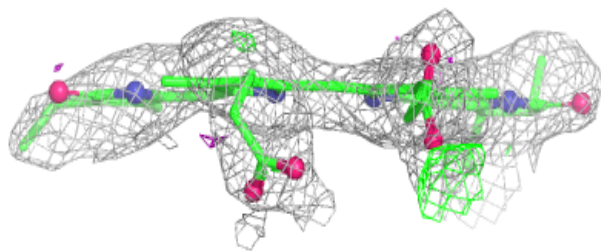
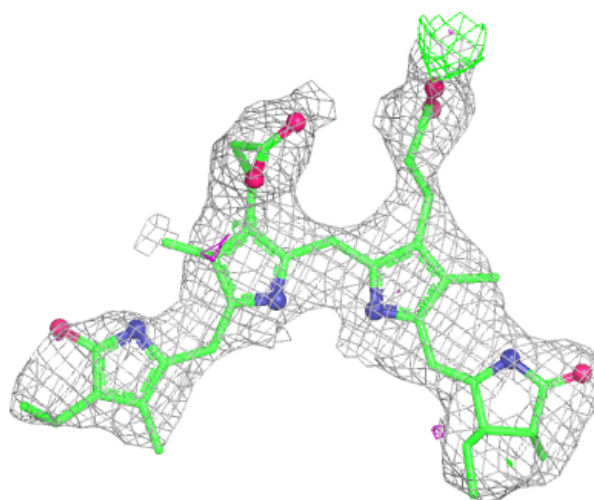
Electron density around CYC L 1082:

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



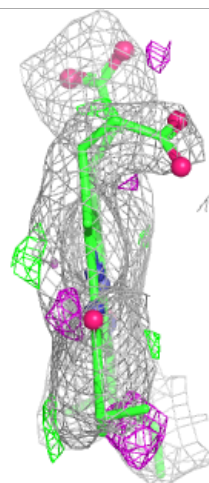
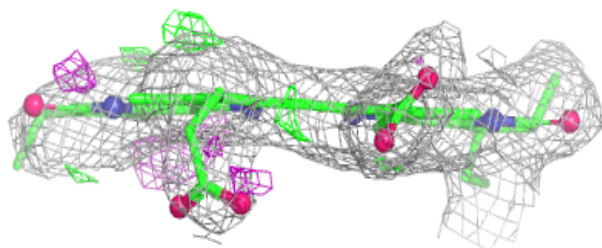
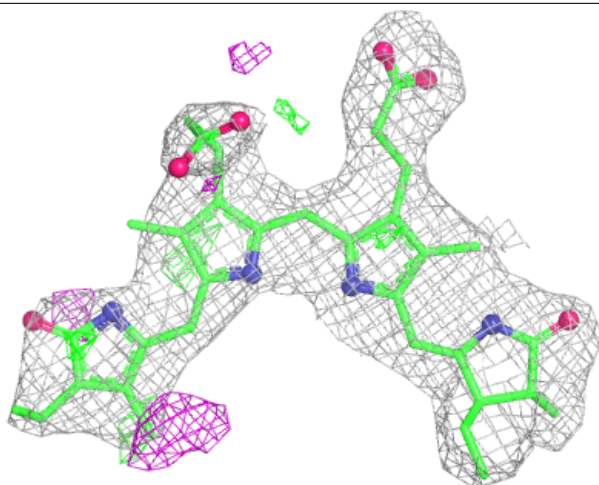
Electron density around CYC J 1082:

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



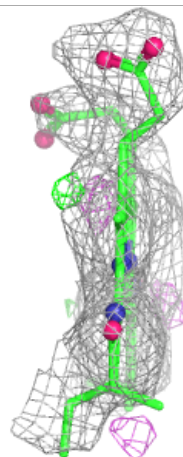
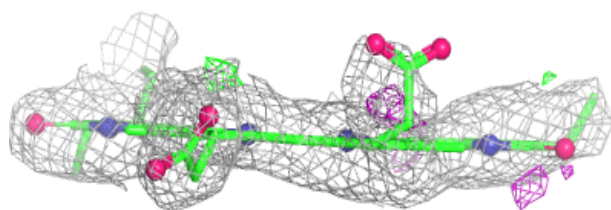
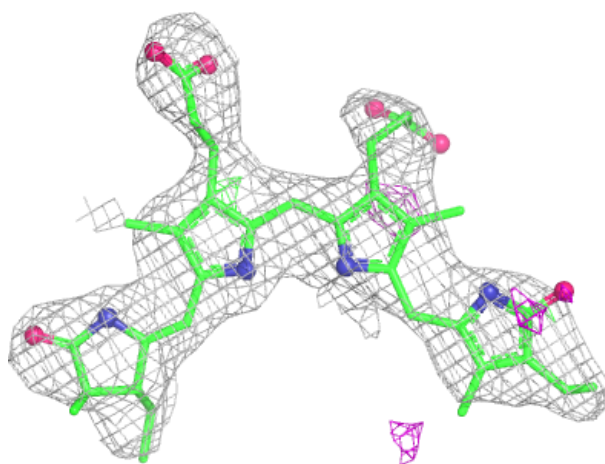
Electron density around CYC H 1082:

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



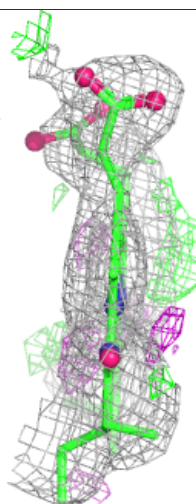
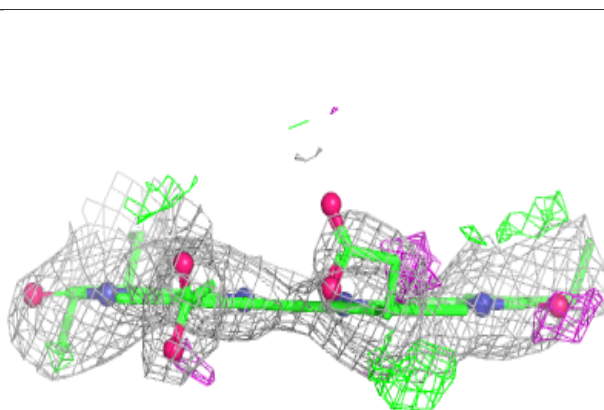
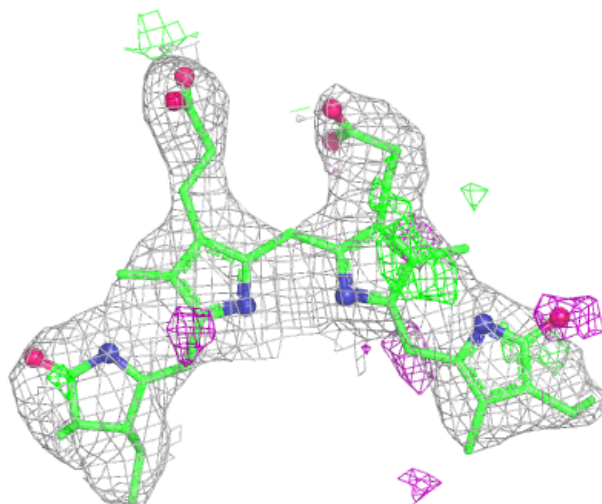
Electron density around CYC P 1082:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



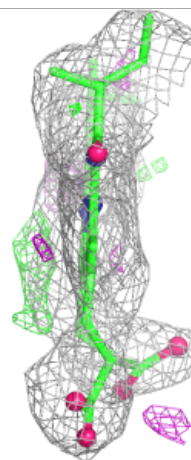
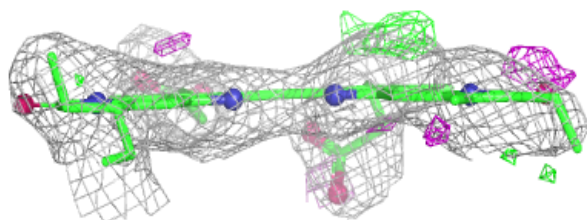
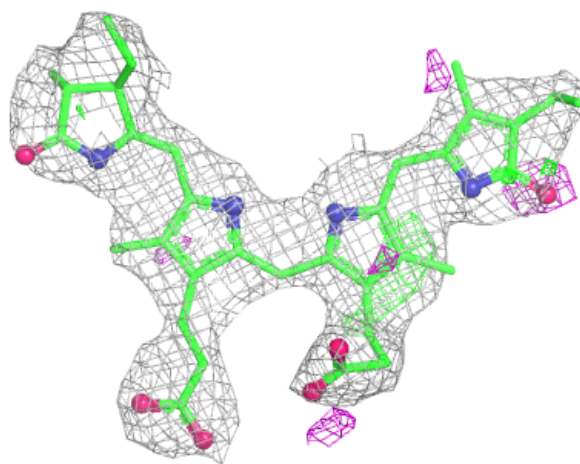
Electron density around CYC R 1082:

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



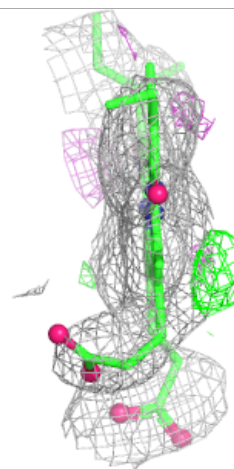
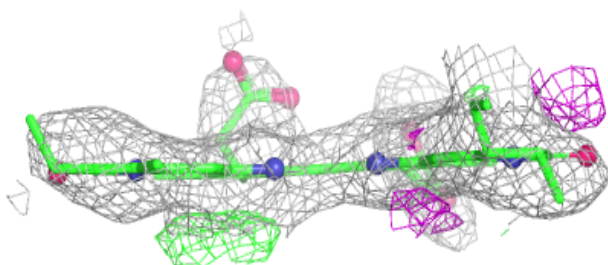
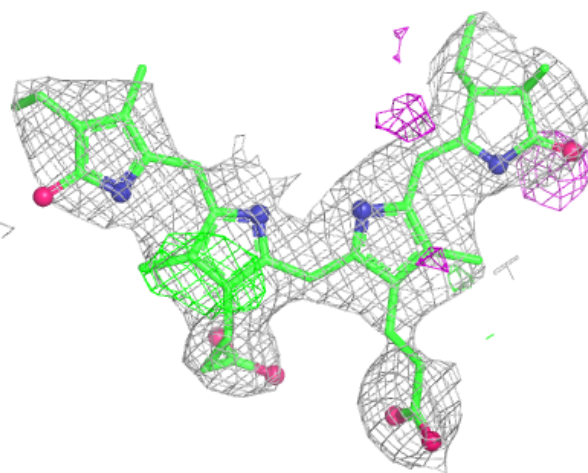
Electron density around CYC D 1082:

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



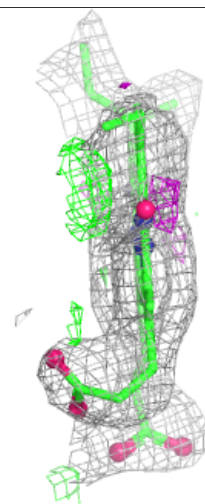
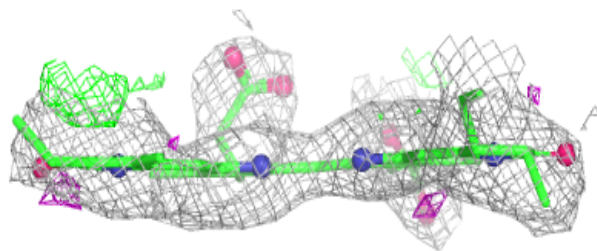
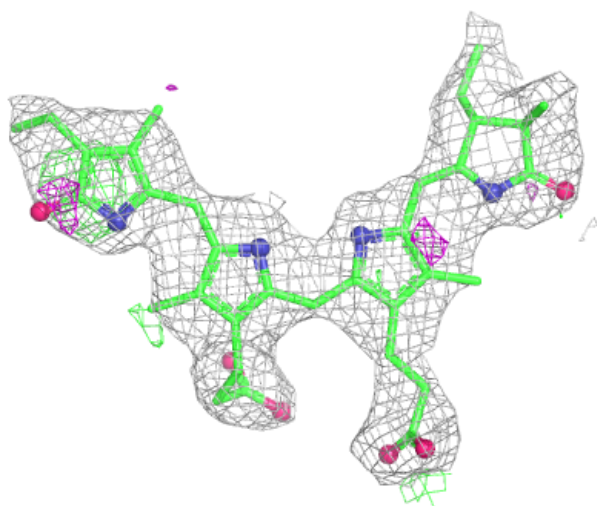
Electron density around CYC F 1082:

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



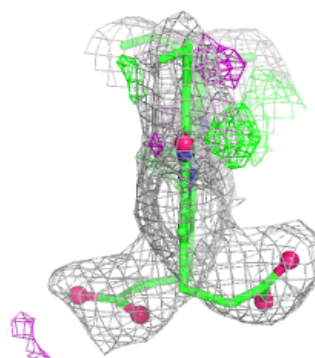
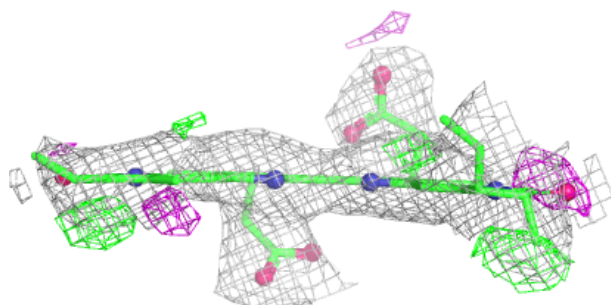
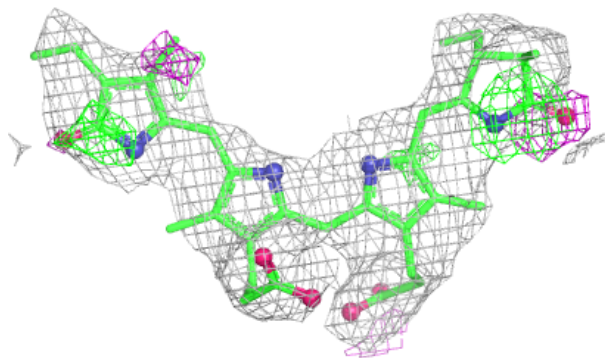
Electron density around CYC B 1082:

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

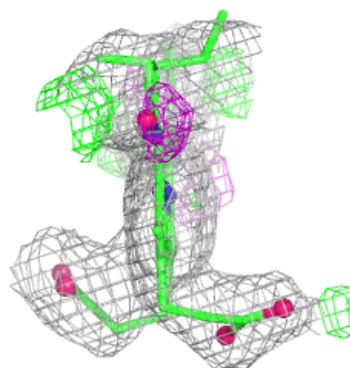
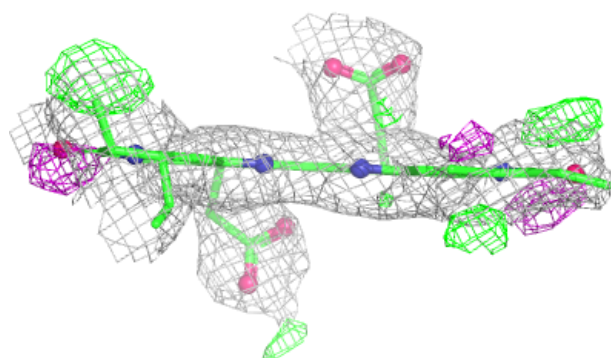
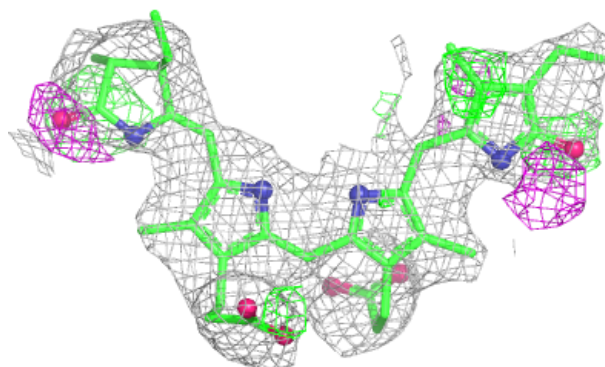


Electron density around CYC P 1153:

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

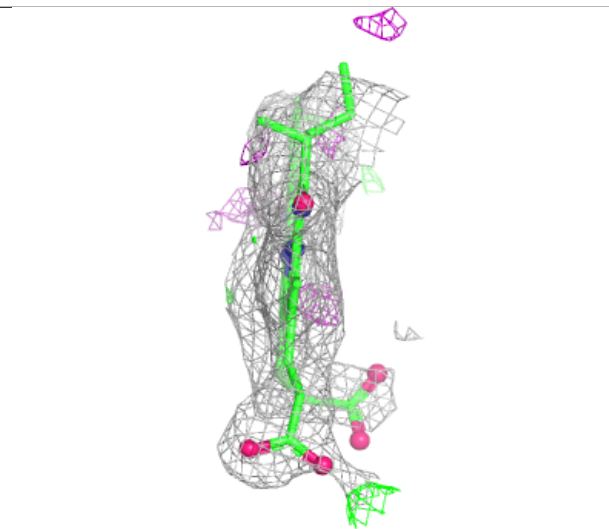
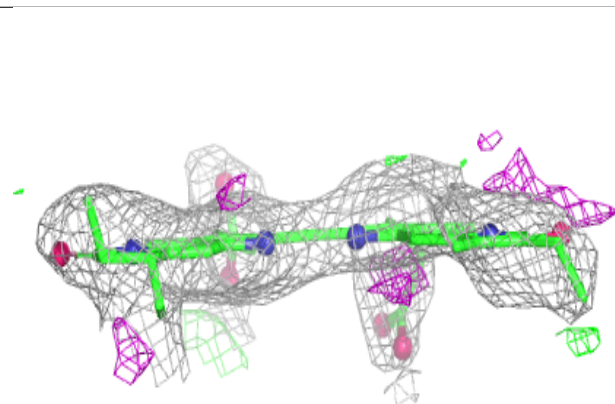
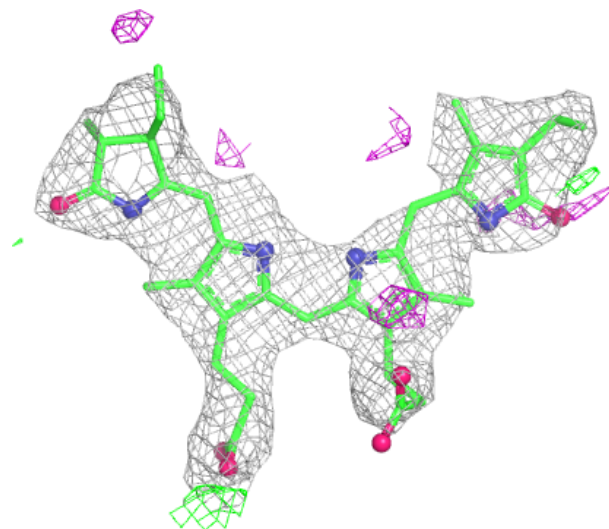
**Electron density around CYC H 1153:**

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



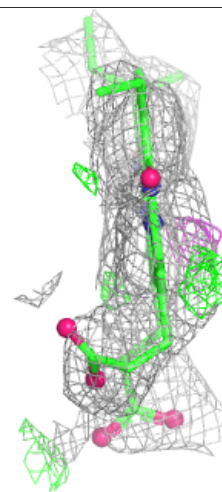
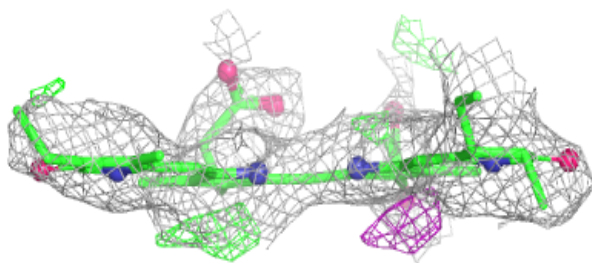
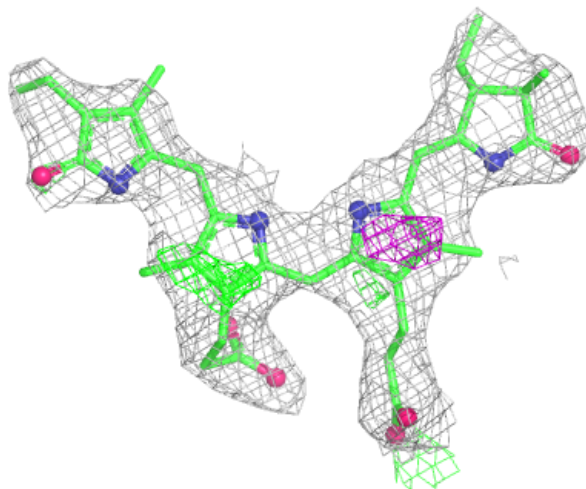
Electron density around CYC V 1082:

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



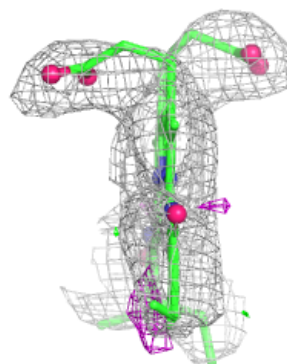
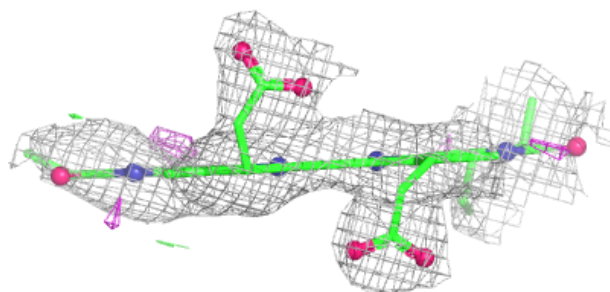
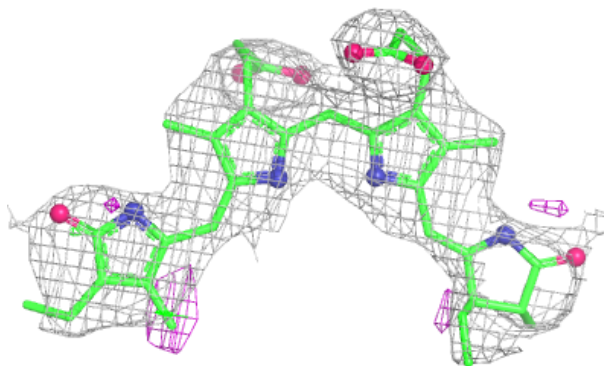
Electron density around CYC X 1082:

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



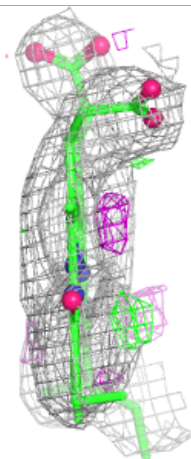
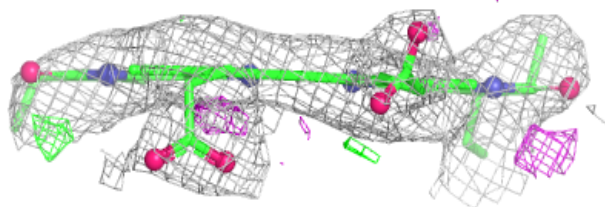
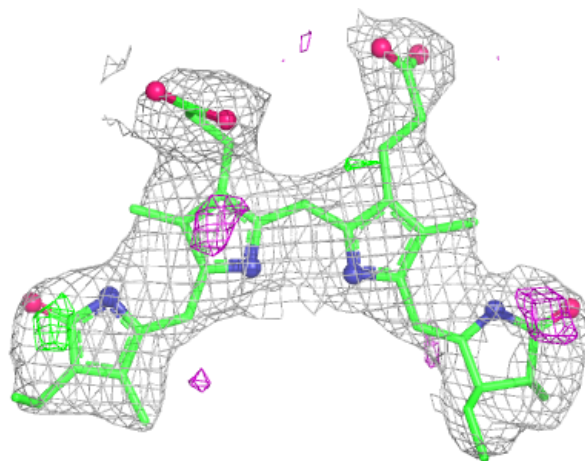
Electron density around CYC X 1153:

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



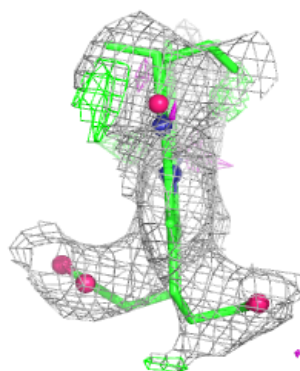
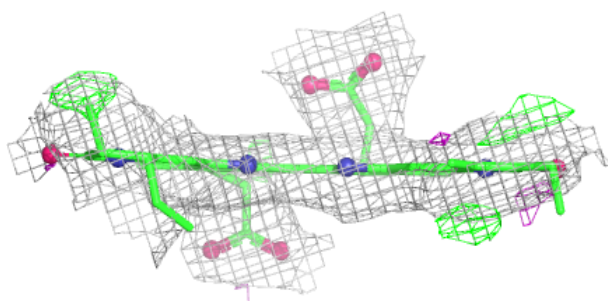
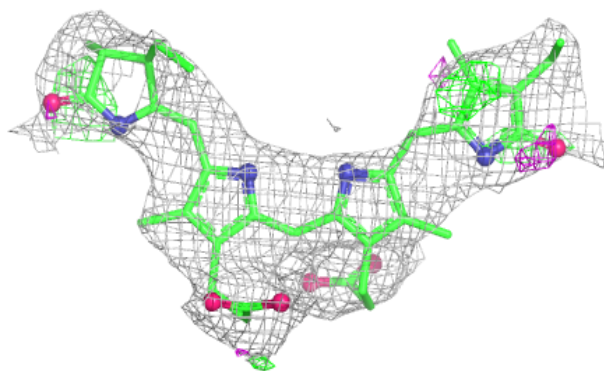
Electron density around CYC T 1082:

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

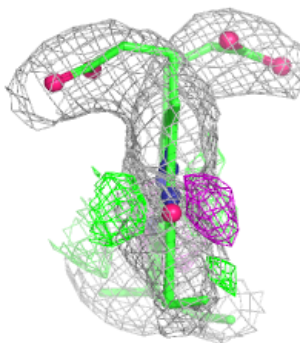
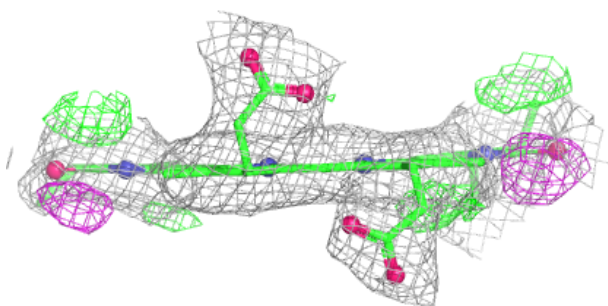
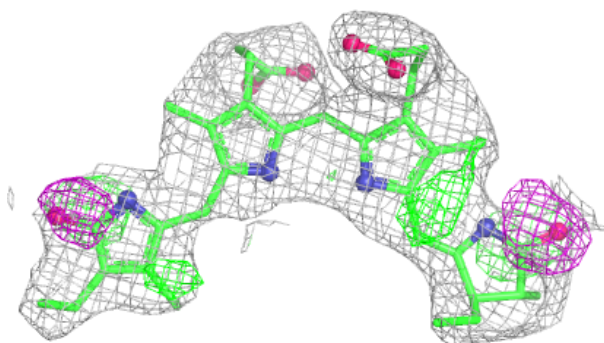


Electron density around CYC J 1153:

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

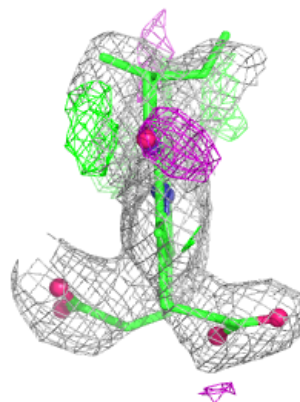
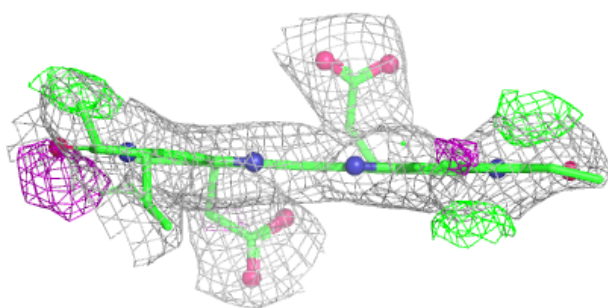
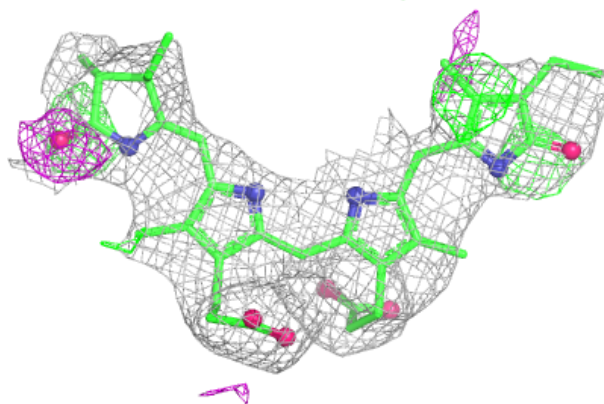
**Electron density around CYC B 1153:**

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

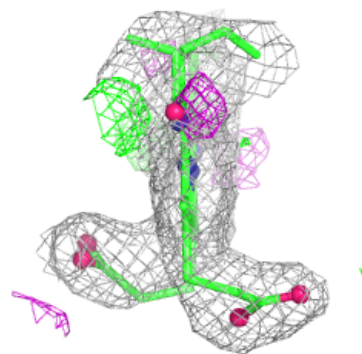
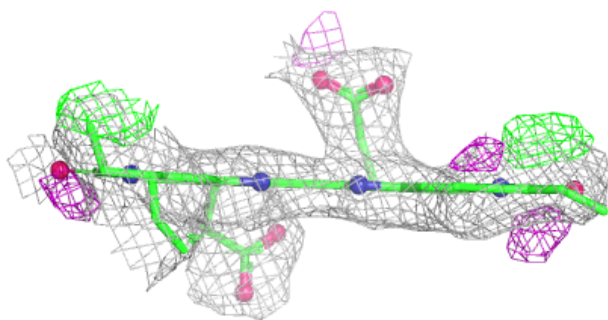
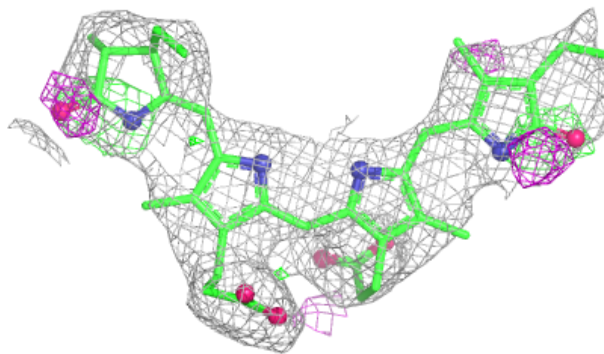


Electron density around CYC L 1153:

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

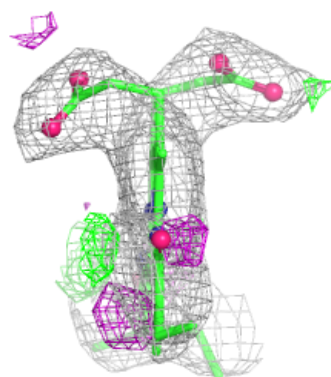
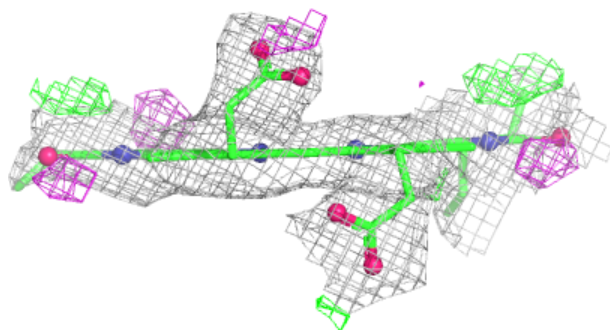
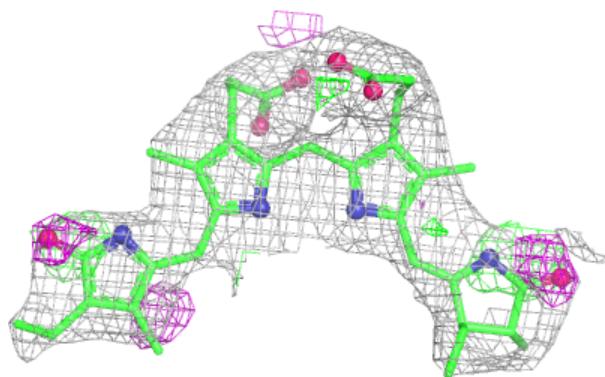
**Electron density around CYC T 1153:**

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

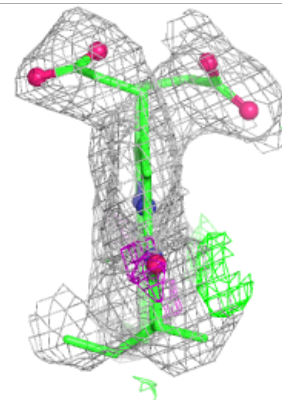
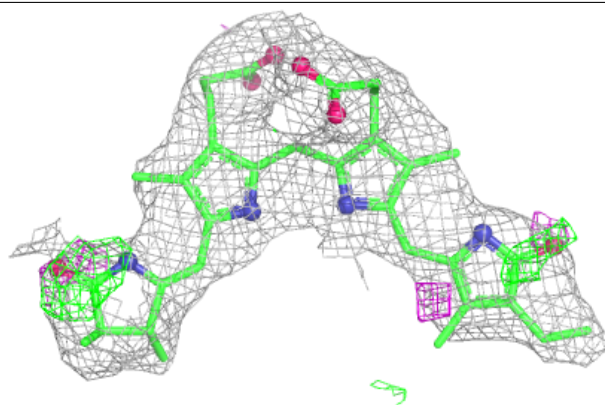


Electron density around CYC F 1153:

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

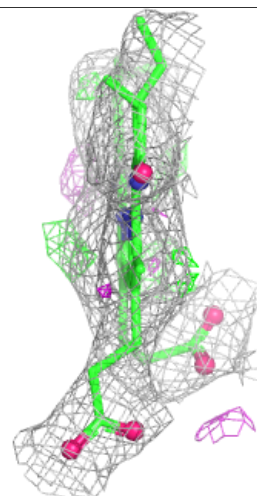
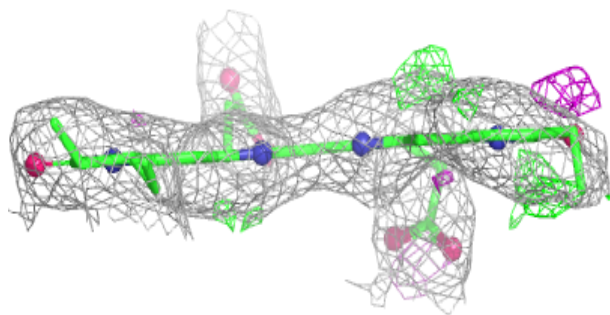
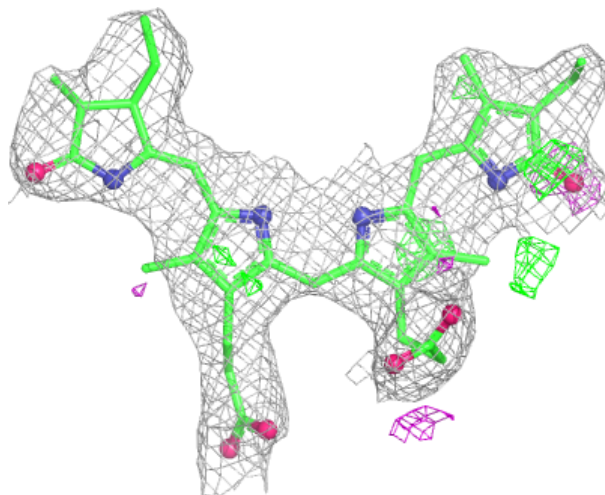
**Electron density around CYC V 1153:**

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



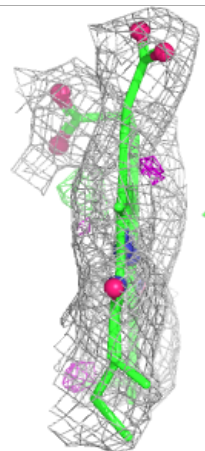
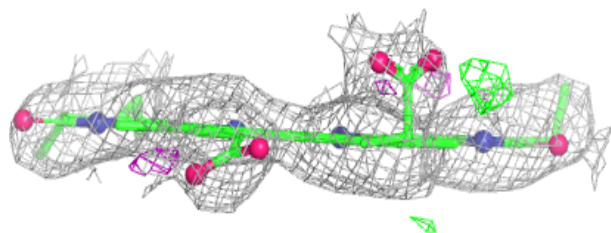
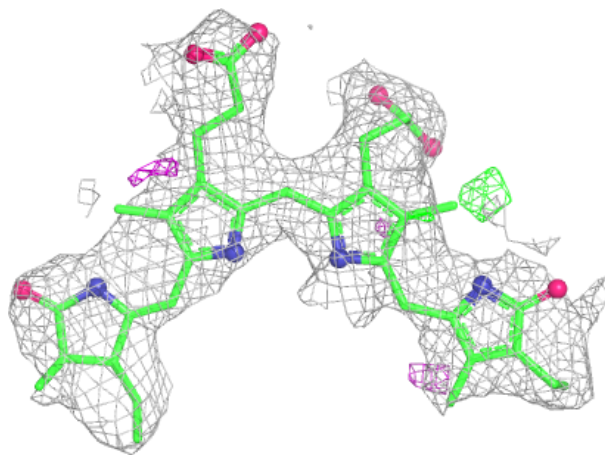
Electron density around CYC W 1084:

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



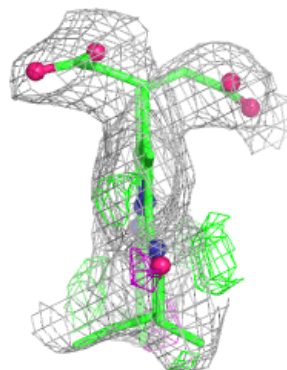
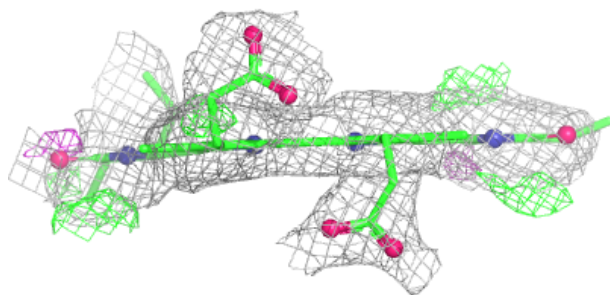
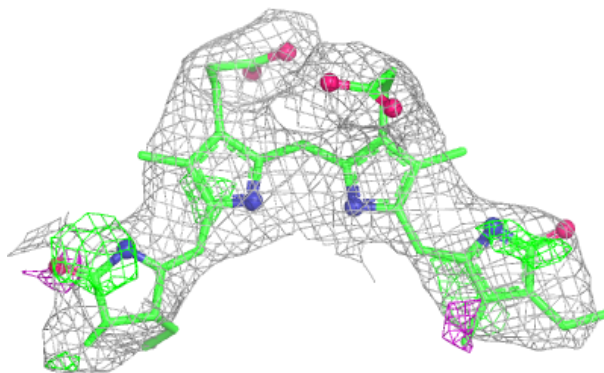
Electron density around CYC M 1084:

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



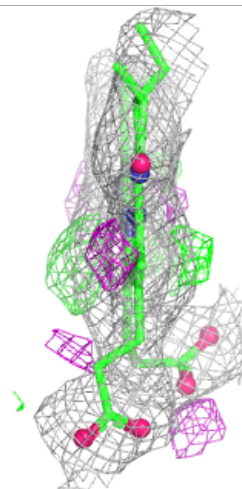
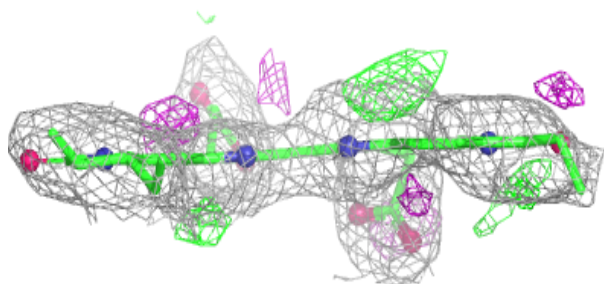
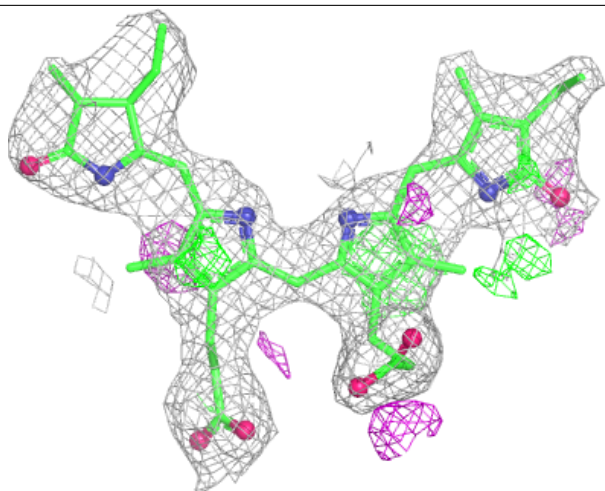
Electron density around CYC D 1153:

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



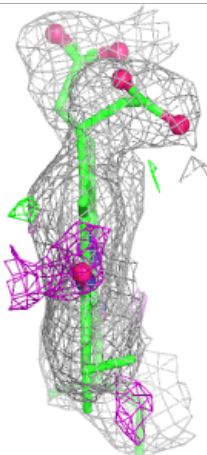
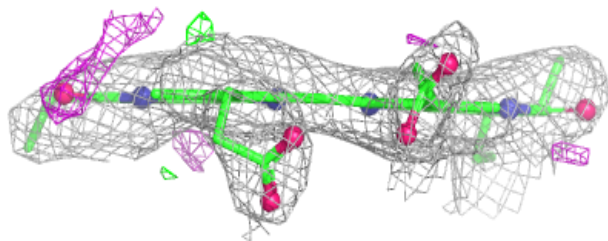
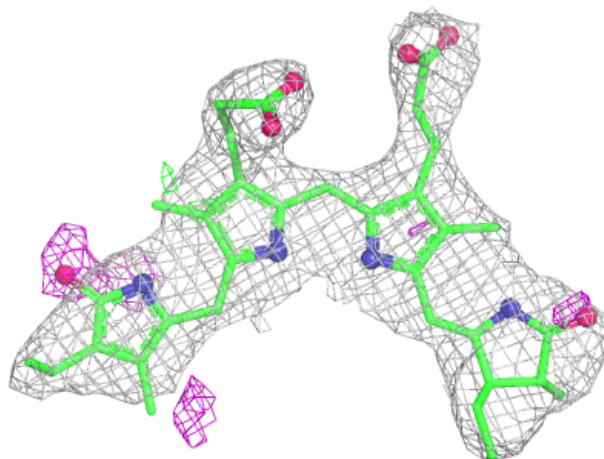
Electron density around CYC E 1084:

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



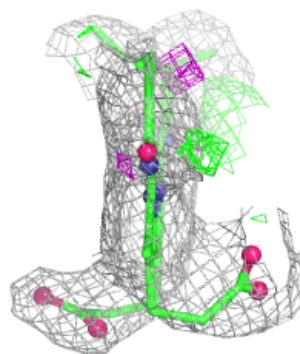
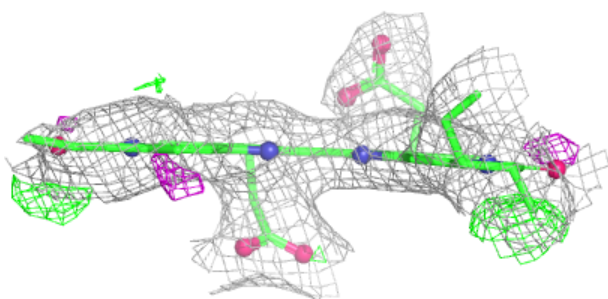
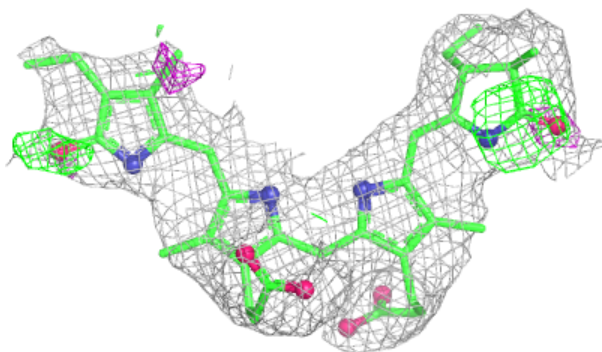
Electron density around CYC N 1082:

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

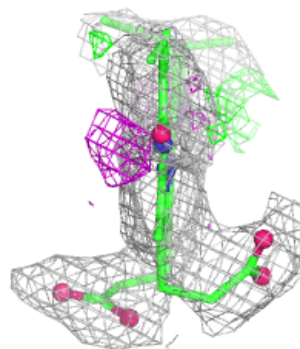
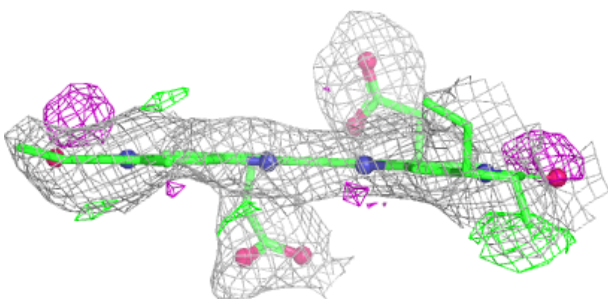
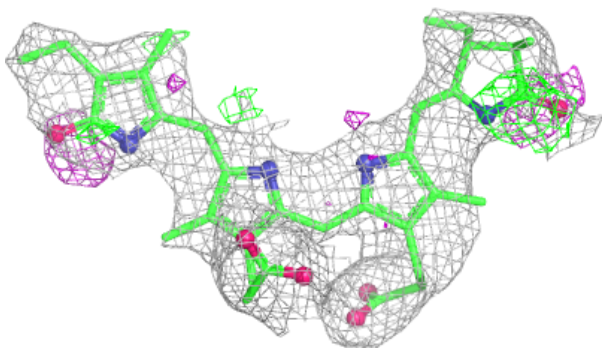


Electron density around CYC R 1153:

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

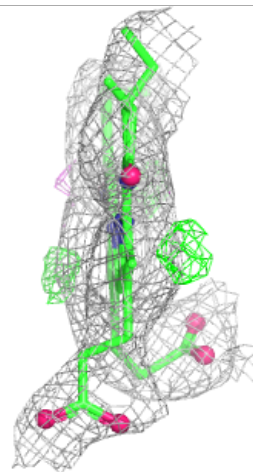
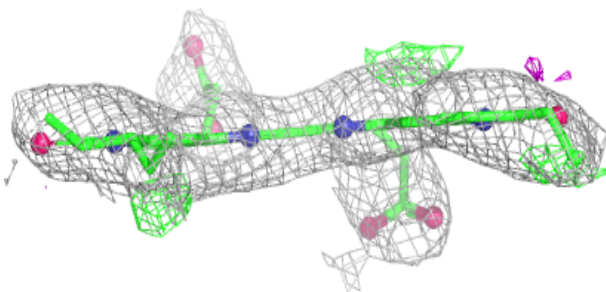
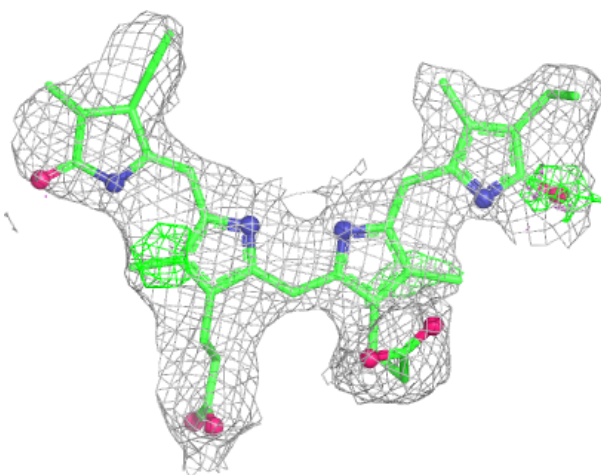
**Electron density around CYC N 1153:**

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



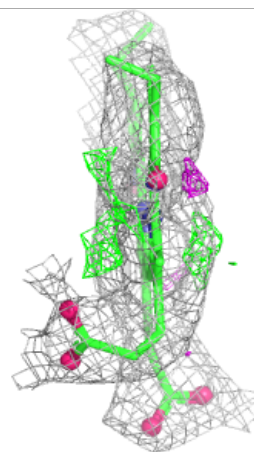
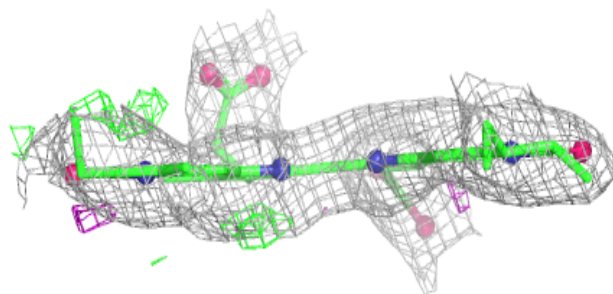
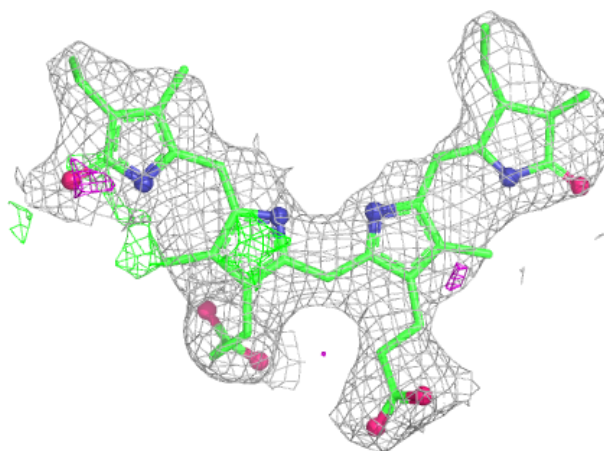
Electron density around CYC G 1084:

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



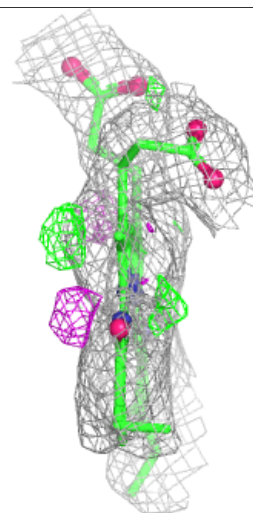
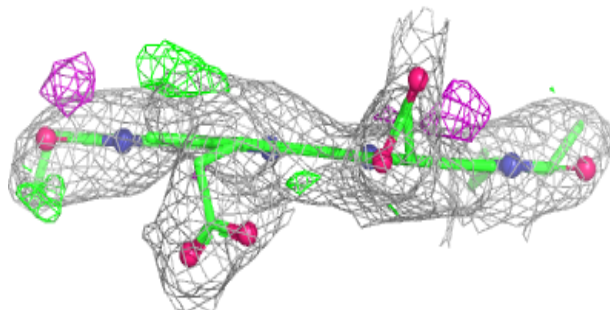
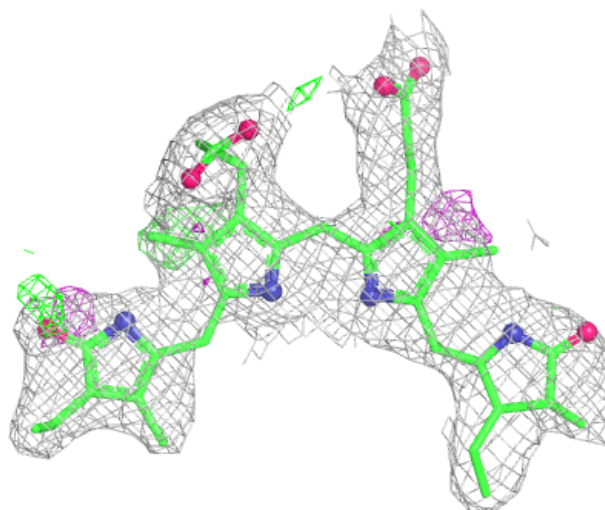
Electron density around CYC U 1084:

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



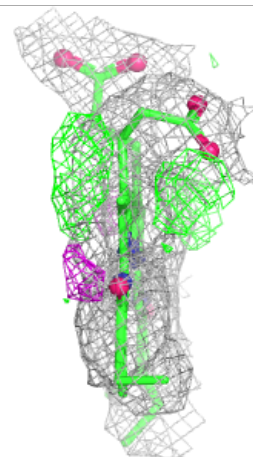
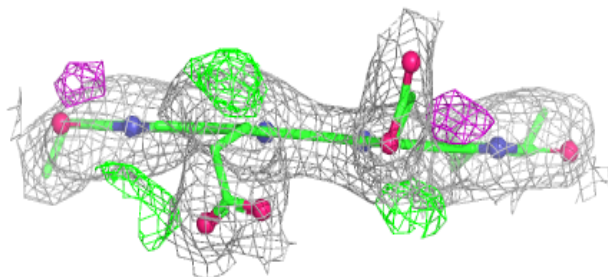
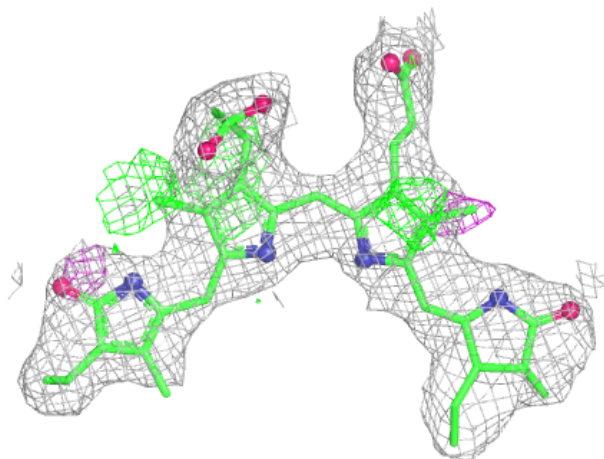
Electron density around CYC I 1084:

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



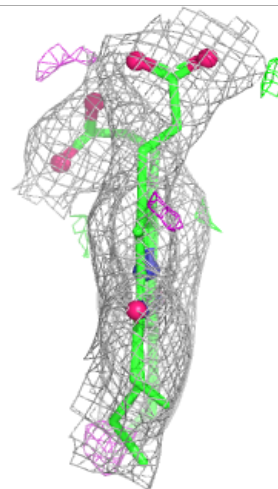
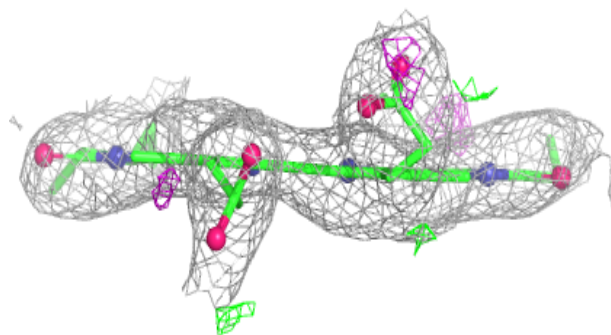
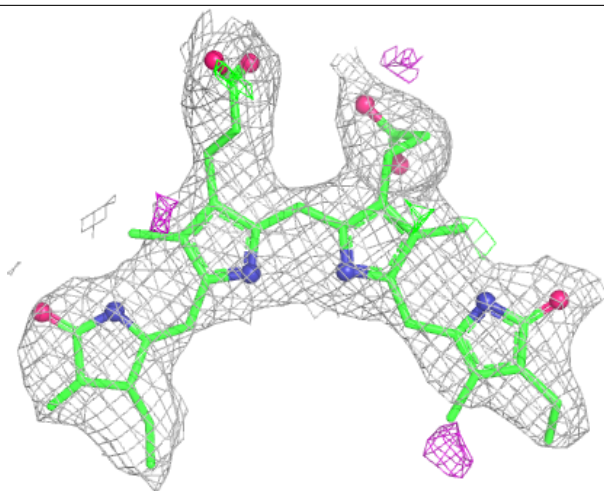
Electron density around CYC S 1084:

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



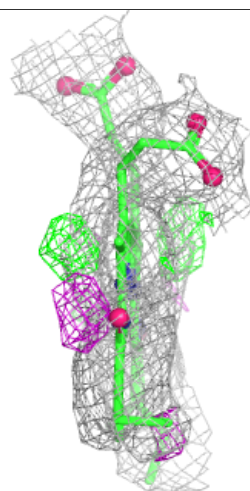
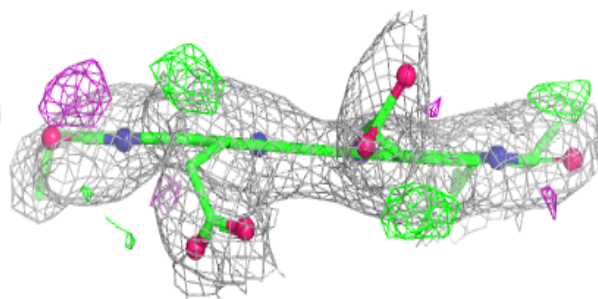
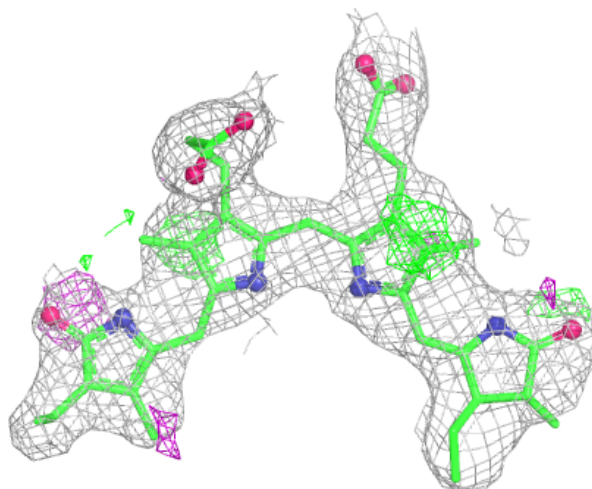
Electron density around CYC O 1084:

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



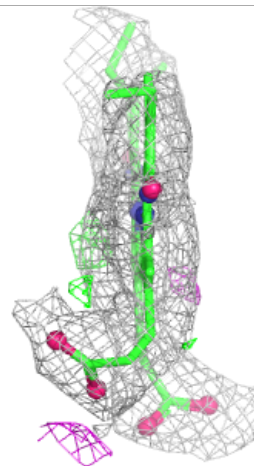
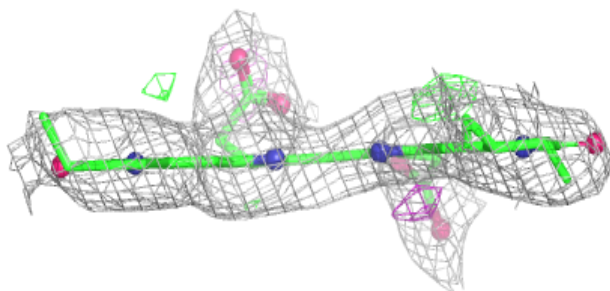
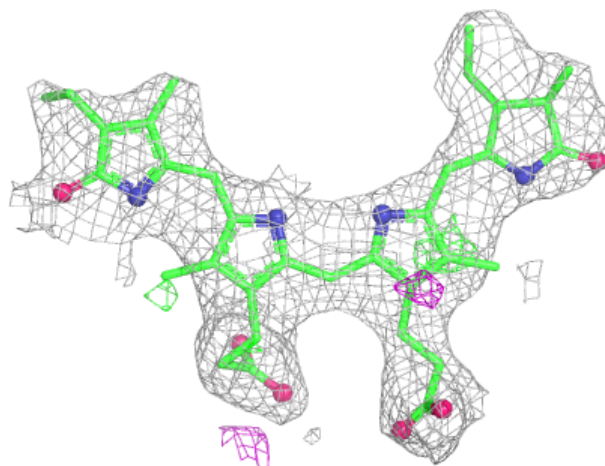
Electron density around CYC K 1084:

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



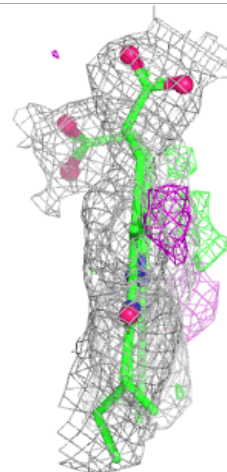
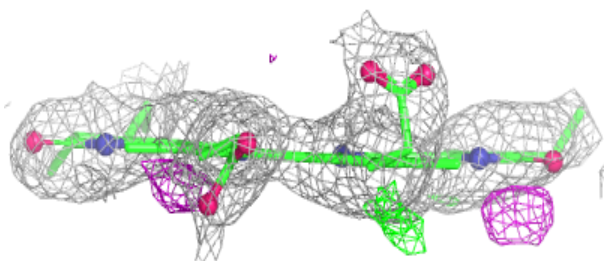
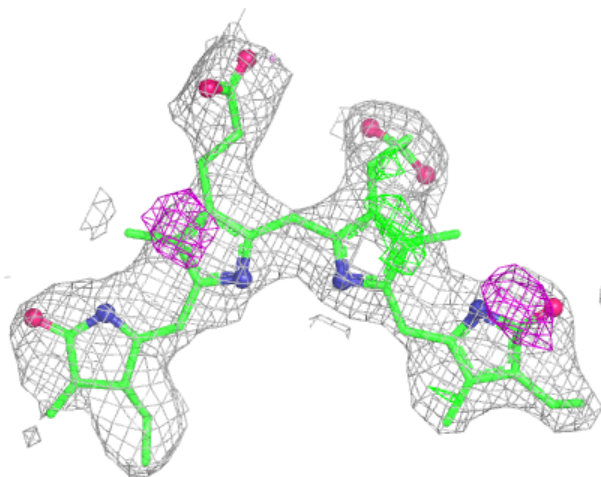
Electron density around CYC C 1084:

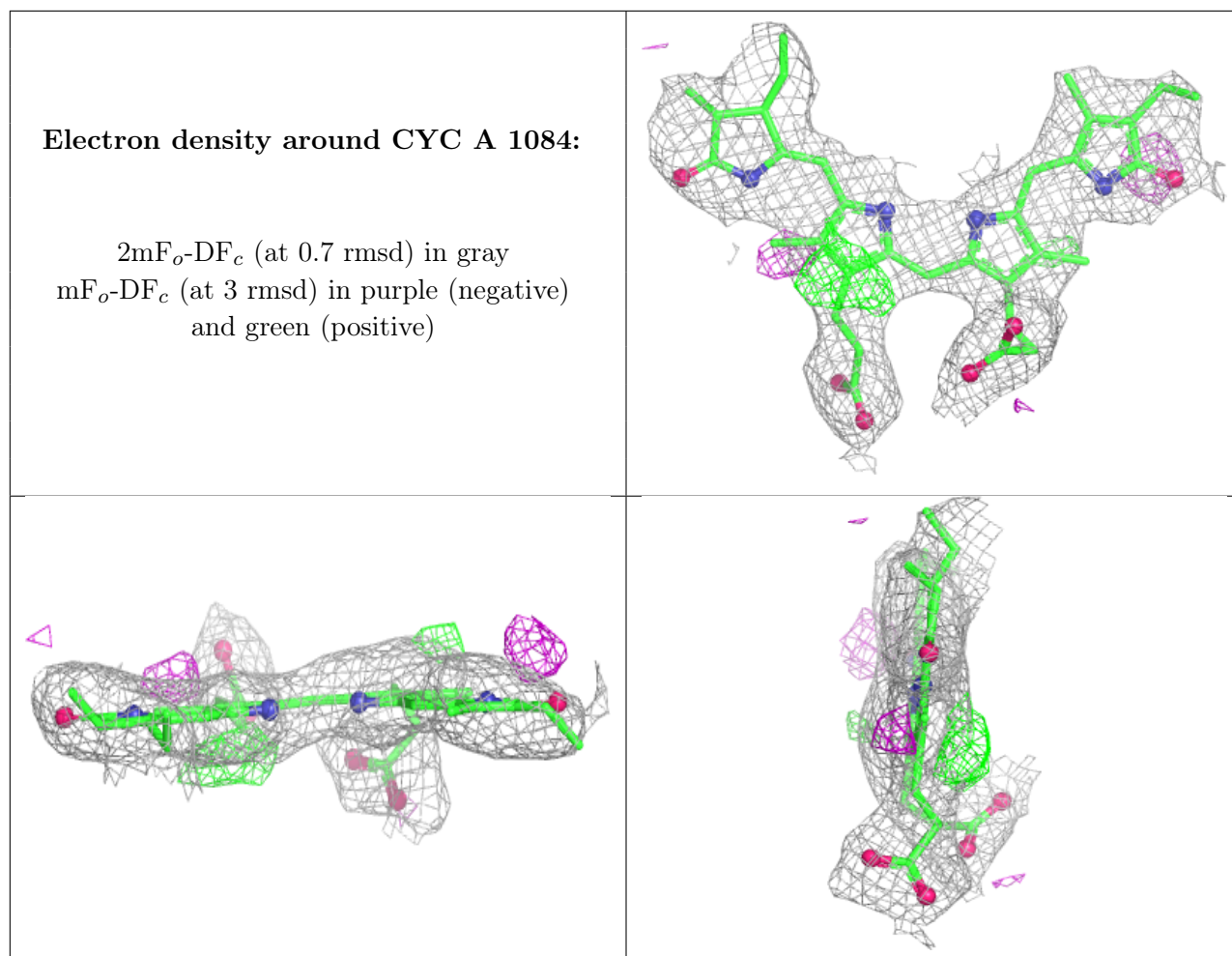
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around CYC Q 1084:

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





6.5 Other polymers [\(i\)](#)

There are no such residues in this entry.