



Full wwPDB EM Validation Report ⓘ

Mar 28, 2026 – 12:38 AM UTC

PDB ID : 7QSL / pdb_00007qsl
EMDB ID : EMD-14133
Title : Bovine complex I in lipid nanodisc, Active-apo
Authors : Chung, I.; Bridges, H.R.; Hirst, J.
Deposited on : 2022-01-13
Resolution : 2.76 Å(reported)

This is a Full wwPDB EM 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/EMValidationReportHelp>

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:

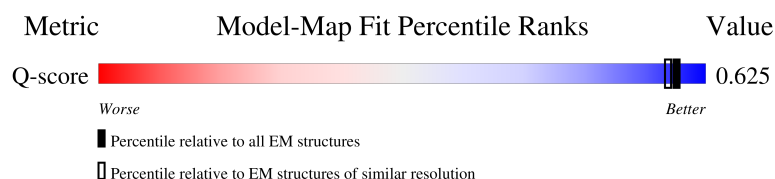
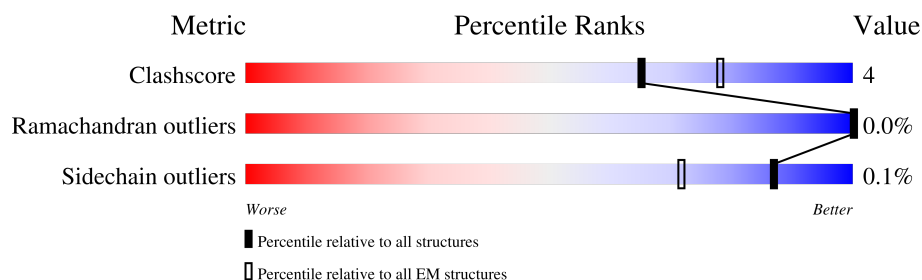
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
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:
ELECTRON MICROSCOPY

The reported resolution of this entry is 2.76 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.



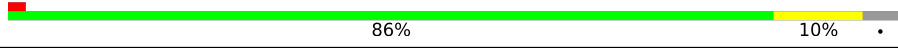
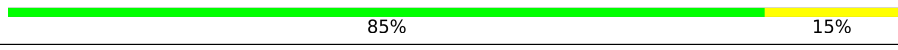
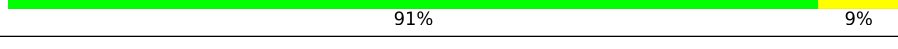
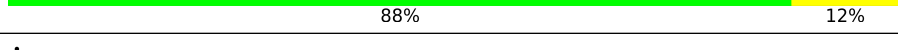
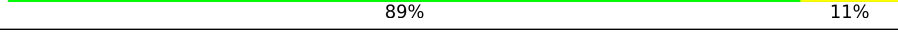
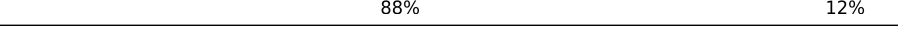
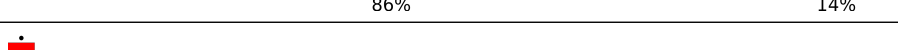
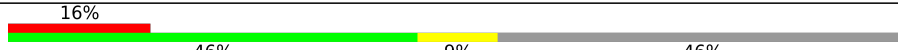
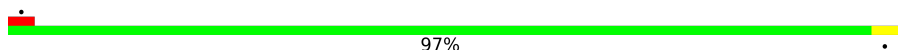
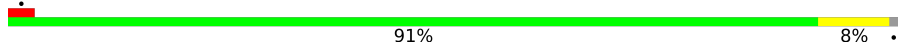
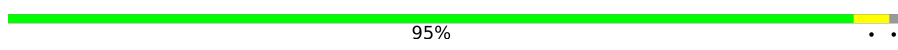
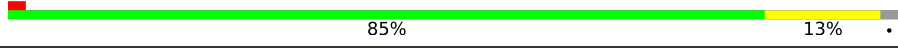
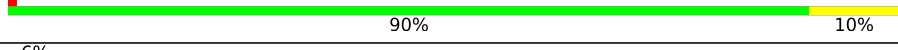
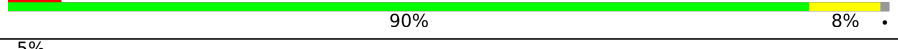
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	10642 (2.26 - 3.26)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the 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 EM map (all-atom inclusion $< 40\%$). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	115	 81% 19%
2	B	216	 65% 7% 28%
3	C	266	 73% 5% 21%
4	D	463	 84% 9% 7%

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Mol	Chain	Length	Quality of chain
5	E	249	
6	F	464	
7	G	727	
8	H	318	
9	I	212	
10	J	175	
11	K	98	
12	L	606	
13	M	459	
14	N	347	
15	O	343	
16	P	380	
17	Q	175	
18	R	124	
19	S	99	
20	T	156	
20	U	156	
21	V	116	
22	W	128	
23	X	172	
24	Y	141	
25	Z	144	
26	a	70	
27	b	84	
28	c	76	

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Mol	Chain	Length	Quality of chain
29	d	120	
30	e	106	
31	f	57	
32	g	154	
33	h	189	
34	i	127	
35	j	108	
36	k	98	
37	l	186	
38	m	129	
39	n	179	
40	o	137	
41	p	176	
42	q	145	
43	r	113	
44	s	109	

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
29	AME	d	1	-	-	X	-

2 Entry composition [i](#)

There are 59 unique types of molecules in this entry. The entry contains 69733 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, 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 NADH-ubiquinone oxidoreductase chain 3.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	115	Total	C	N	O	S	0	0
			921	622	133	159	7		

- Molecule 2 is a protein called NADH dehydrogenase [ubiquinone] iron-sulfur protein 7, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	155	Total	C	N	O	S	0	0
			1241	792	224	211	14		

- Molecule 3 is a protein called NADH dehydrogenase [ubiquinone] iron-sulfur protein 3, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	209	Total	C	N	O	S	0	0
			1738	1120	298	317	3		

- Molecule 4 is a protein called NADH dehydrogenase [ubiquinone] iron-sulfur protein 2, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	D	430	Total	C	N	O	S	0	0
			3459	2209	596	629	25		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	129	ARG	GLN	variant	UNP P17694

- Molecule 5 is a protein called NADH dehydrogenase [ubiquinone] flavoprotein 2, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	E	214	Total	C	N	O	S	0	0
			1659	1059	278	312	10		

- Molecule 6 is a protein called NADH dehydrogenase [ubiquinone] flavoprotein 1, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	F	432	Total	C	N	O	S	1	0
			3336	2102	597	617	20		

- Molecule 7 is a protein called NADH-ubiquinone oxidoreductase 75 kDa subunit, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	G	699	Total	C	N	O	S	1	0
			5366	3362	934	1030	40		

- Molecule 8 is a protein called NADH-ubiquinone oxidoreductase chain 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	H	318	Total	C	N	O	S	1	0
			2517	1687	386	421	23		

- Molecule 9 is a protein called NADH dehydrogenase [ubiquinone] iron-sulfur protein 8, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	I	176	Total	C	N	O	S	0	0
			1414	889	243	270	12		

- Molecule 10 is a protein called NADH-ubiquinone oxidoreductase chain 6.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	J	175	Total	C	N	O	S	0	0
			1345	906	191	236	12		

- Molecule 11 is a protein called NADH-ubiquinone oxidoreductase chain 4L.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	K	98	Total	C	N	O	S	0	0
			745	486	112	131	16		

- Molecule 12 is a protein called NADH-ubiquinone oxidoreductase chain 5.

Mol	Chain	Residues	Atoms					AltConf	Trace
12	L	606	Total	C	N	O	S	0	0
			4802	3195	737	827	43		

- Molecule 13 is a protein called NADH-ubiquinone oxidoreductase chain 4.

Mol	Chain	Residues	Atoms					AltConf	Trace
13	M	459	Total	C	N	O	S	0	0
			3654	2436	570	609	39		

- Molecule 14 is a protein called NADH-ubiquinone oxidoreductase chain 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	N	347	Total	C	N	O	S	0	0
			2733	1817	416	457	43		

- Molecule 15 is a protein called NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 10, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
15	O	320	Total	C	N	O	S	0	0
			2589	1662	429	488	10		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
O	255	LYS	ASN	variant	UNP P34942

- Molecule 16 is a protein called NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 9, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	P	342	Total	C	N	O	S	1	0
			2768	1792	489	482	5		

- Molecule 17 is a protein called NADH dehydrogenase [ubiquinone] iron-sulfur protein 4, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
17	Q	129	Total	C	N	O	S	0	0
			1049	659	188	199	3		

- Molecule 18 is a protein called NADH dehydrogenase [ubiquinone] iron-sulfur protein 6, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
18	R	96	Total	C	N	O	S	0	0
			740	454	140	143	3		

- Molecule 19 is a protein called NADH dehydrogenase [ubiquinone] 1 alpha subcomplex sub-unit 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
19	S	87	Total	C	N	O	S	0	0
			700	440	131	127	2		

- Molecule 20 is a protein called Acyl carrier protein, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
20	T	85	Total	C	N	O	S	0	0
			688	444	101	138	5		
20	U	86	Total	C	N	O	S	0	0
			693	447	102	139	5		

- Molecule 21 is a protein called NADH dehydrogenase [ubiquinone] 1 alpha subcomplex sub-unit 5.

Mol	Chain	Residues	Atoms					AltConf	Trace
21	V	115	Total	C	N	O	S	0	0
			928	600	157	168	3		

- Molecule 22 is a protein called NADH dehydrogenase [ubiquinone] 1 alpha subcomplex sub-unit 6.

Mol	Chain	Residues	Atoms					AltConf	Trace
22	W	115	Total	C	N	O	S	0	0
			976	625	181	166	4		

- Molecule 23 is a protein called NADH dehydrogenase [ubiquinone] 1 alpha subcomplex sub-unit 8.

Mol	Chain	Residues	Atoms					AltConf	Trace
23	X	171	Total	C	N	O	S	0	0
			1402	887	253	252	10		

- Molecule 24 is a protein called NADH dehydrogenase [ubiquinone] 1 alpha subcomplex sub-unit 11.

Mol	Chain	Residues	Atoms					AltConf	Trace
24	Y	140	Total	C	N	O	S	0	0
			1030	657	176	191	6		

- Molecule 25 is a protein called NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 13.

Mol	Chain	Residues	Atoms					AltConf	Trace
25	Z	141	Total	C	N	O	S	0	0
			1152	740	201	202	9		

- Molecule 26 is a protein called NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
26	a	70	Total	C	N	O	S	0	0
			569	365	104	95	5		

- Molecule 27 is a protein called NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 3.

Mol	Chain	Residues	Atoms					AltConf	Trace
27	b	83	Total	C	N	O	S	0	0
			651	425	109	115	2		

- Molecule 28 is a protein called NADH dehydrogenase [ubiquinone] 1 subunit C1, mitochondrial.

Mol	Chain	Residues	Atoms				AltConf	Trace
28	c	49	Total	C	N	O	0	0
			414	273	70	71		

- Molecule 29 is a protein called NADH dehydrogenase [ubiquinone] 1 subunit C2.

Mol	Chain	Residues	Atoms					AltConf	Trace
29	d	120	Total	C	N	O	S	0	0
			999	650	172	172	5		

- Molecule 30 is a protein called NADH dehydrogenase [ubiquinone] iron-sulfur protein 5.

Mol	Chain	Residues	Atoms					AltConf	Trace
30	e	99	Total	C	N	O	S	0	0
			829	523	158	142	6		

- Molecule 31 is a protein called NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
31	f	57	Total	C	N	O	S	0	0
			492	322	86	82	2		

- Molecule 32 is a protein called NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 11, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
32	g	100	Total	C	N	O	S	0	0
			839	539	139	157	4		

- Molecule 33 is a protein called NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 5, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
33	h	138	Total	C	N	O	S	0	0
			1154	759	196	197	2		

- Molecule 34 is a protein called NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 6.

Mol	Chain	Residues	Atoms					AltConf	Trace
34	i	127	Total	C	N	O	S	0	0
			1097	722	191	183	1		

- Molecule 35 is a protein called NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 2, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
35	j	67	Total	C	N	O	S	0	0
			580	381	95	103	1		

- Molecule 36 is a protein called NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 3.

Mol	Chain	Residues	Atoms					AltConf	Trace
36	k	81	Total	C	N	O	S	0	0
			653	427	110	114	2		

- Molecule 37 is a protein called NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 8, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
37	l	156	Total	C	N	O	S	0	0
			1314	850	216	240	8		

- Molecule 38 is a protein called NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 4.

Mol	Chain	Residues	Atoms					AltConf	Trace
38	m	128	Total	C	N	O	S	0	0
			1067	684	188	195			

- Molecule 39 is a protein called NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 9.

Mol	Chain	Residues	Atoms					AltConf	Trace
39	n	172	Total	C	N	O	S	0	0
			1492	955	273	257	7		

- Molecule 40 is a protein called NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 7.

Mol	Chain	Residues	Atoms					AltConf	Trace
40	o	122	Total	C	N	O	S	0	0
			1048	653	201	185	9		

- Molecule 41 is a protein called NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 10.

Mol	Chain	Residues	Atoms					AltConf	Trace
41	p	173	Total	C	N	O	S	0	0
			1453	910	268	267	8		

- Molecule 42 is a protein called NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 12.

Mol	Chain	Residues	Atoms					AltConf	Trace
42	q	145	Total	C	N	O	S	0	0
			1209	778	216	210	5		

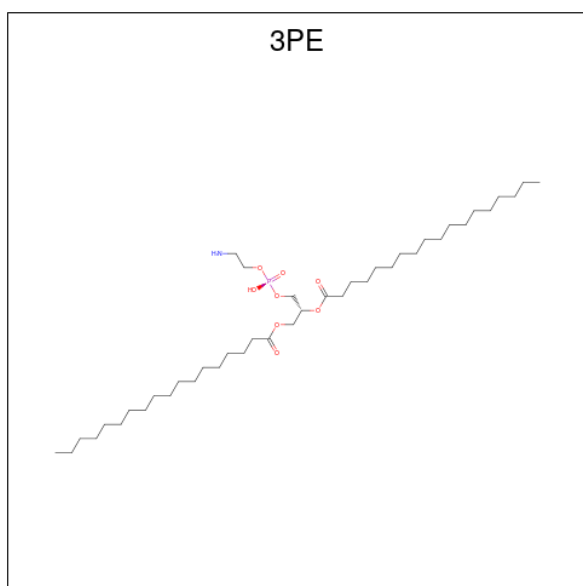
- Molecule 43 is a protein called NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 7.

Mol	Chain	Residues	Atoms					AltConf	Trace
43	r	95	Total	C	N	O	S	0	0
			776	490	144	139	3		

- Molecule 44 is a protein called NADH dehydrogenase [ubiquinone] flavoprotein 3, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
44	s	45	Total	C	N	O	S	1	0
			391	244	71	75	1		

- Molecule 45 is 1,2-Distearoyl-sn-glycerophosphoethanolamine (CCD ID: 3PE) (formula: C₄₁H₈₂NO₈P).



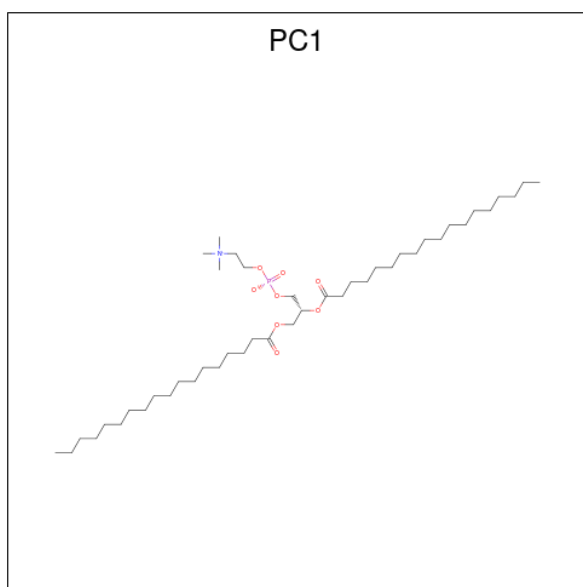
Mol	Chain	Residues	Atoms					AltConf
45	A	1	Total	C	N	O	P	0
			42	32	1	8	1	
45	A	1	Total	C	N	O	P	0
			37	27	1	8	1	
45	H	1	Total	C	N	O	P	0
			29	19	1	8	1	
45	I	1	Total	C	N	O	P	0
			38	28	1	8	1	
45	L	1	Total	C	N	O	P	0
			39	29	1	8	1	
45	L	1	Total	C	N	O	P	0
			45	35	1	8	1	
45	M	1	Total	C	N	O	P	0
			45	35	1	8	1	

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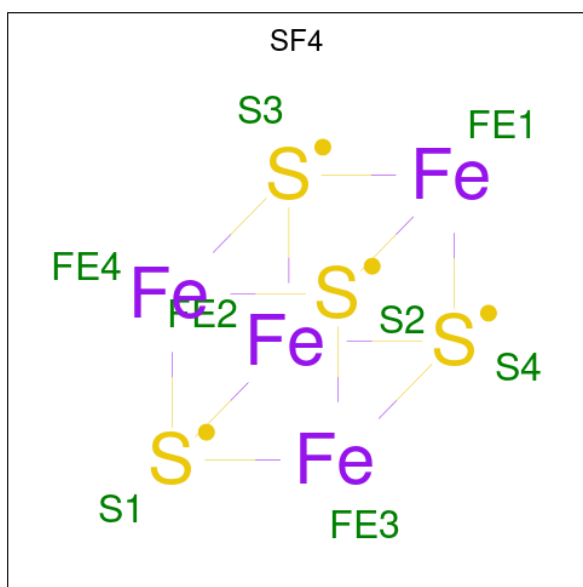
Mol	Chain	Residues	Atoms					AltConf
45	M	1	Total	C	N	O	P	0
			48	38	1	8	1	
45	M	1	Total	C	N	O	P	0
			51	41	1	8	1	
45	M	1	Total	C	N	O	P	0
			50	40	1	8	1	
45	N	1	Total	C	N	O	P	0
			48	38	1	8	1	
45	P	1	Total	C	N	O	P	0
			35	25	1	8	1	
45	Y	1	Total	C	N	O	P	0
			35	25	1	8	1	
45	Y	1	Total	C	N	O	P	0
			40	30	1	8	1	
45	Y	1	Total	C	N	O	P	0
			31	21	1	8	1	
45	Y	1	Total	C	N	O	P	0
			51	41	1	8	1	
45	Z	1	Total	C	N	O	P	0
			35	25	1	8	1	
45	b	1	Total	C	N	O	P	0
			47	37	1	8	1	
45	d	1	Total	C	N	O	P	0
			37	27	1	8	1	
45	m	1	Total	C	N	O	P	0
			40	30	1	8	1	

- Molecule 46 is 1,2-DIACYL-SN-GLYCERO-3-PHOSPHOCHOLINE (CCD ID: PC1) (formula: C₄₄H₈₈NO₈P).



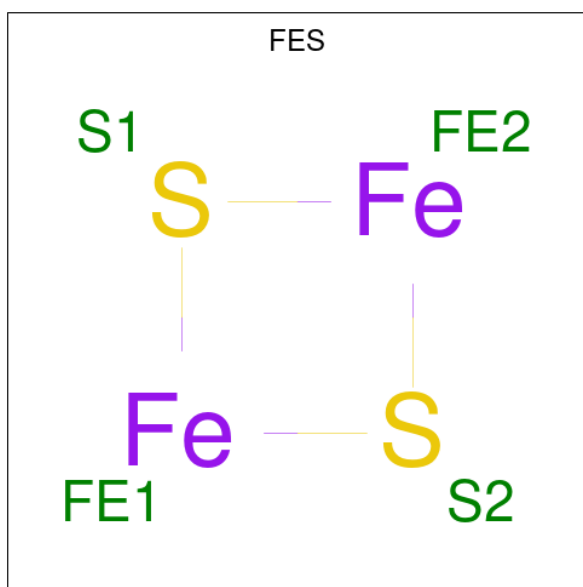
Mol	Chain	Residues	Atoms					AltConf
46	A	1	Total	C	N	O	P	0
			35	25	1	8	1	
46	A	1	Total	C	N	O	P	0
			33	23	1	8	1	
46	B	1	Total	C	N	O	P	0
			41	31	1	8	1	
46	M	1	Total	C	N	O	P	0
			39	29	1	8	1	
46	M	1	Total	C	N	O	P	0
			42	32	1	8	1	
46	P	1	Total	C	N	O	P	0
			54	44	1	8	1	
46	Z	1	Total	C	N	O	P	0
			48	38	1	8	1	
46	d	1	Total	C	N	O	P	0
			33	23	1	8	1	
46	h	1	Total	C	N	O	P	0
			44	34	1	8	1	
46	m	1	Total	C	N	O	P	0
			40	30	1	8	1	

- Molecule 47 is IRON/SULFUR CLUSTER (CCD ID: SF4) (formula: Fe₄S₄).



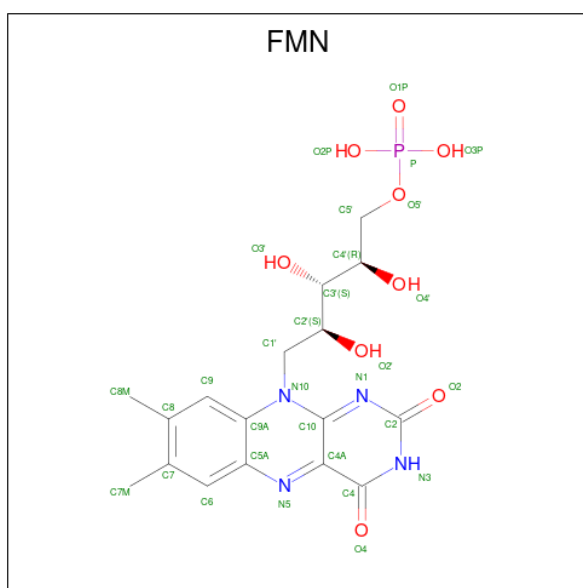
Mol	Chain	Residues	Atoms			AltConf
47	B	1	Total	Fe	S	0
			8	4	4	
47	F	1	Total	Fe	S	0
			8	4	4	
47	G	1	Total	Fe	S	0
			8	4	4	
47	G	1	Total	Fe	S	0
			8	4	4	
47	I	1	Total	Fe	S	0
			8	4	4	
47	I	1	Total	Fe	S	0
			8	4	4	

- Molecule 48 is FE2/S2 (INORGANIC) CLUSTER (CCD ID: FES) (formula: Fe₂S₂).



Mol	Chain	Residues	Atoms			AltConf
48	E	1	Total	Fe	S	0
			4	2	2	
48	G	1	Total	Fe	S	0
			4	2	2	

- Molecule 49 is FLAVIN MONONUCLEOTIDE (CCD ID: FMN) (formula: $C_{17}H_{21}N_4O_9P$).

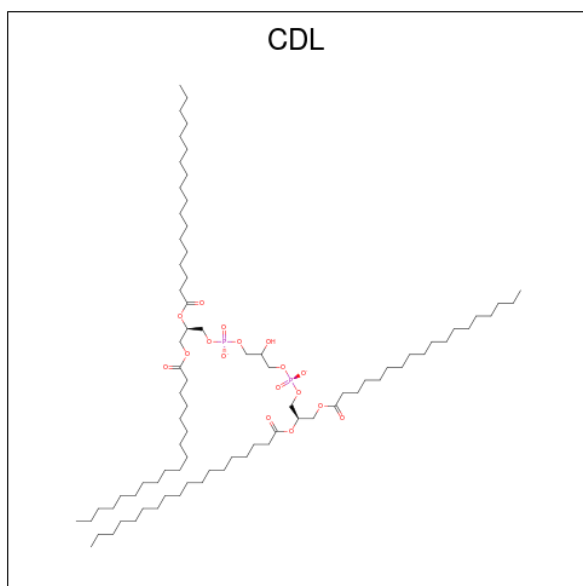


Mol	Chain	Residues	Atoms					AltConf
49	F	1	Total	C	N	O	P	0
			31	17	4	9	1	

- Molecule 50 is POTASSIUM ION (CCD ID: K) (formula: K).

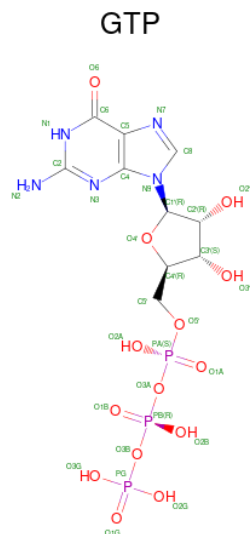
Mol	Chain	Residues	Atoms		AltConf
50	G	1	Total	K	0
			1	1	

- Molecule 51 is CARDIOLIPIN (CCD ID: CDL) (formula: $C_{81}H_{156}O_{17}P_2$).



Mol	Chain	Residues	Atoms				AltConf
51	H	1	Total	C	O	P	0
			50	31	17	2	
51	K	1	Total	C	O	P	0
			69	50	17	2	
51	L	1	Total	C	O	P	0
			60	41	17	2	
51	N	1	Total	C	O	P	0
			60	41	17	2	
51	X	1	Total	C	O	P	0
			73	54	17	2	
51	d	1	Total	C	O	P	0
			65	46	17	2	
51	h	1	Total	C	O	P	0
			78	59	17	2	
51	r	1	Total	C	O	P	0
			62	43	17	2	

- Molecule 52 is GUANOSINE-5'-TRIPHOSPHATE (CCD ID: GTP) (formula: $C_{10}H_{16}N_5O_{14}P_3$).

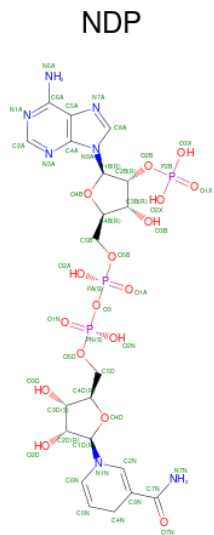


Mol	Chain	Residues	Atoms					AltConf
52	O	1	Total 32	C 10	N 5	O 14	P 3	0

- Molecule 53 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	AltConf
53	O	1	Total Mg 1 1	0

- Molecule 54 is NADPH DIHYDRO-NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (CCD ID: NDP) (formula: $C_{21}H_{30}N_7O_{17}P_3$).

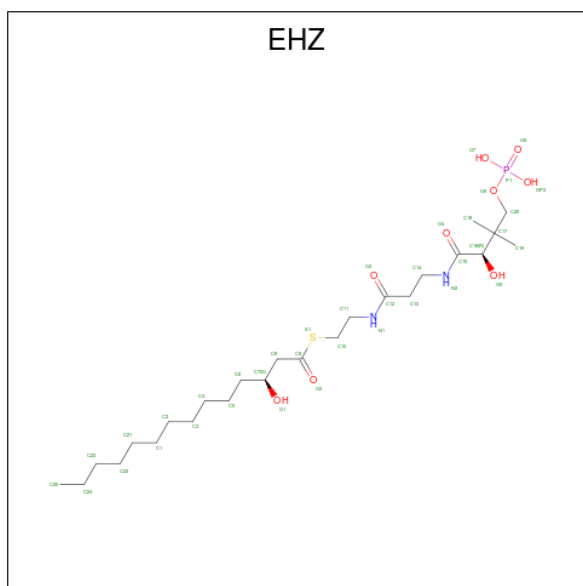


Mol	Chain	Residues	Atoms					AltConf
54	P	1	Total	C	N	O	P	0
			48	21	7	17	3	

- Molecule 55 is ZINC ION (CCD ID: ZN) (formula: Zn).

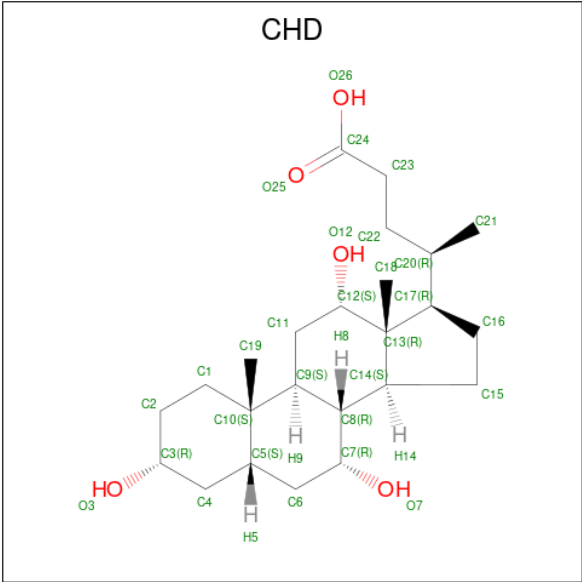
Mol	Chain	Residues	Atoms		AltConf
55	R	1	Total	Zn	0
			1	1	

- Molecule 56 is {S}-[2-[3-[(2 {R})-3,3-dimethyl-2-oxidanyl-4-phosphonooxy-butanoyl]amino]propanoylamino]ethyl] (3 {S})-3-oxidanyltetradecanethioate (CCD ID: EHZ) (formula: C₂₅H₄₉N₂O₉PS).



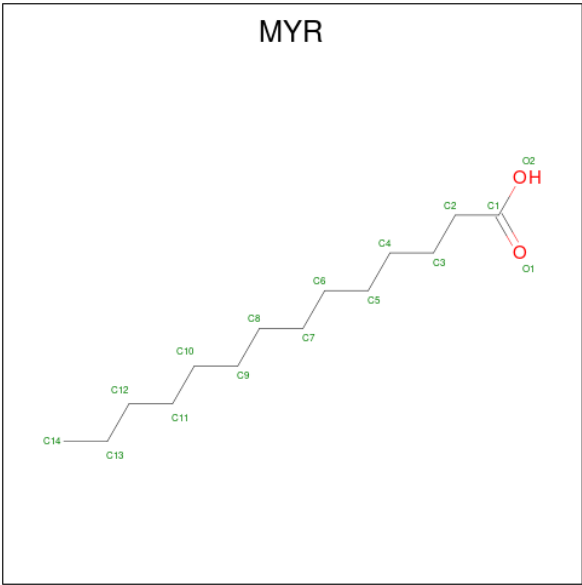
Mol	Chain	Residues	Atoms						AltConf
56	T	1	Total	C	N	O	P	S	0
			37	25	2	8	1	1	
56	U	1	Total	C	N	O	P	S	0
			37	25	2	8	1	1	

- Molecule 57 is CHOLIC ACID (CCD ID: CHD) (formula: C₂₄H₄₀O₅).



Mol	Chain	Residues	Atoms			AltConf
57	i	1	Total	C	O	0
			29	24	5	

- Molecule 58 is MYRISTIC ACID (CCD ID: MYR) (formula: C₁₄H₂₈O₂).



Mol	Chain	Residues	Atoms			AltConf
58	o	1	Total	C	O	0
			15	14	1	

- Molecule 59 is water.

Mol	Chain	Residues	Atoms	AltConf
59	A	17	Total O 17 17	0
59	B	59	Total O 59 59	0
59	C	54	Total O 54 54	0
59	D	123	Total O 123 123	0
59	E	9	Total O 9 9	0
59	F	30	Total O 30 30	0
59	G	120	Total O 120 120	0
59	H	64	Total O 64 64	0
59	I	64	Total O 64 64	0
59	J	17	Total O 17 17	0
59	K	10	Total O 10 10	0
59	L	40	Total O 40 40	0
59	M	62	Total O 62 62	0
59	N	60	Total O 60 60	0
59	O	10	Total O 10 10	0
59	P	44	Total O 44 44	0
59	Q	51	Total O 51 51	0
59	R	14	Total O 14 14	0
59	U	1	Total O 1 1	0
59	W	6	Total O 6 6	0
59	X	22	Total O 22 22	0
59	Y	1	Total O 1 1	0

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Mol	Chain	Residues	Atoms		AltConf
59	Z	23	Total 23	O 23	0
59	a	15	Total 15	O 15	0
59	b	2	Total 2	O 2	0
59	d	10	Total 10	O 10	0
59	e	13	Total 13	O 13	0
59	g	5	Total 5	O 5	0
59	h	11	Total 11	O 11	0
59	i	1	Total 1	O 1	0
59	k	1	Total 1	O 1	0
59	l	7	Total 7	O 7	0
59	m	5	Total 5	O 5	0
59	n	10	Total 10	O 10	0
59	p	12	Total 12	O 12	0
59	q	18	Total 18	O 18	0
59	r	11	Total 11	O 11	0
59	s	2	Total 2	O 2	0

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion in map 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 diamond above a residue indicates a poor fit to the EM map for this residue (all-atom inclusion < 40%). 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.

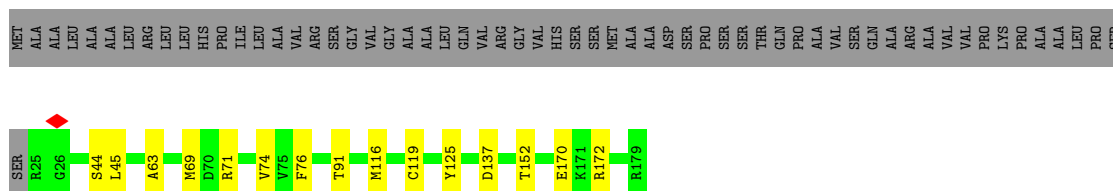
- Molecule 1: NADH-ubiquinone oxidoreductase chain 3

Chain A: 



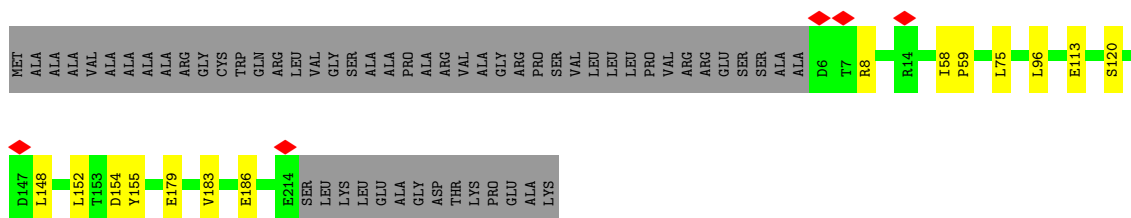
- Molecule 2: NADH dehydrogenase [ubiquinone] iron-sulfur protein 7, mitochondrial

Chain B: 



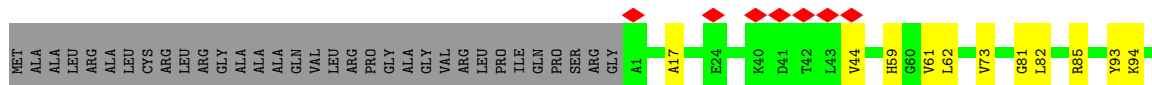
- Molecule 3: NADH dehydrogenase [ubiquinone] iron-sulfur protein 3, mitochondrial

Chain C: 



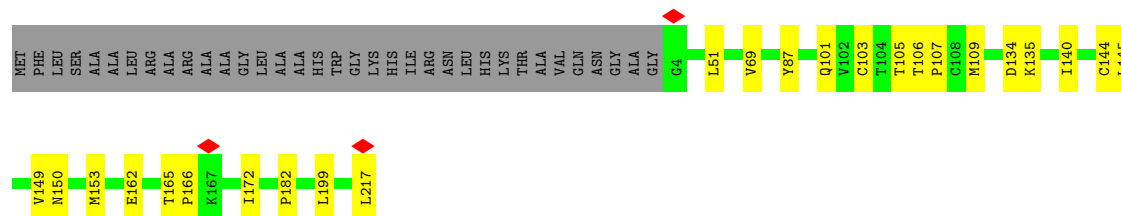
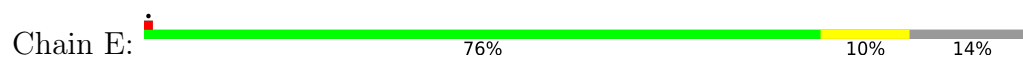
- Molecule 4: NADH dehydrogenase [ubiquinone] iron-sulfur protein 2, mitochondrial

Chain D: 

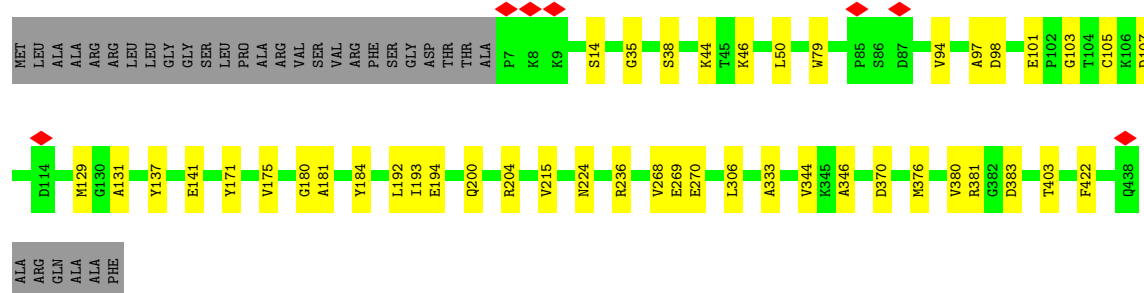
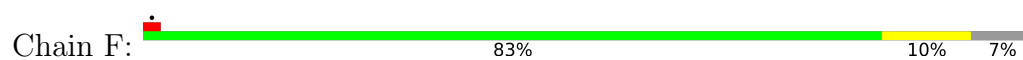




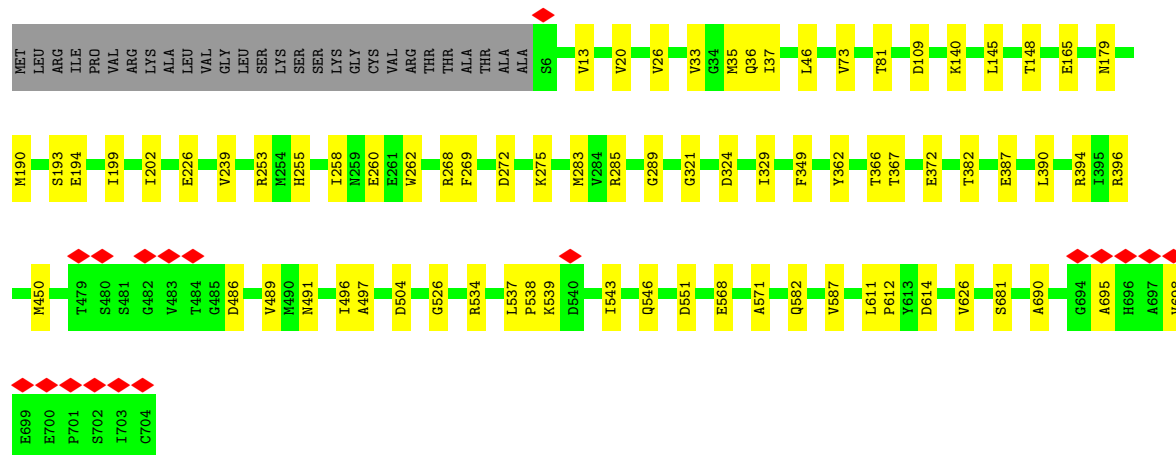
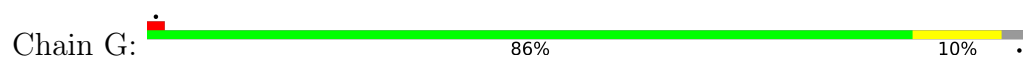
- Molecule 5: NADH dehydrogenase [ubiquinone] flavoprotein 2, mitochondrial



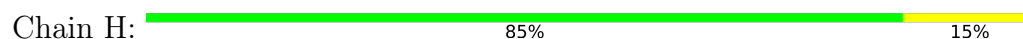
- Molecule 6: NADH dehydrogenase [ubiquinone] flavoprotein 1, mitochondrial

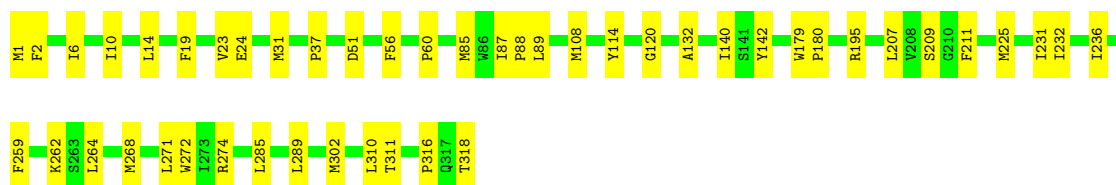


- Molecule 7: NADH-ubiquinone oxidoreductase 75 kDa subunit, mitochondrial



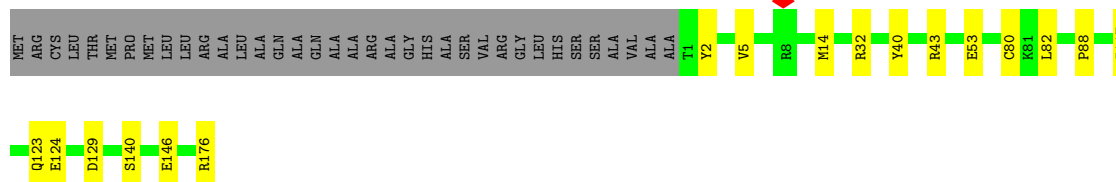
- Molecule 8: NADH-ubiquinone oxidoreductase chain 1





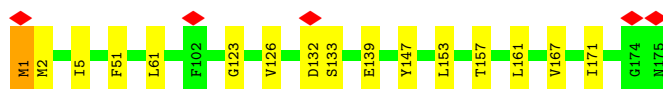
- Molecule 9: NADH dehydrogenase [ubiquinone] iron-sulfur protein 8, mitochondrial

Chain I: 75% 8% 17%



- Molecule 10: NADH-ubiquinone oxidoreductase chain 6

Chain J: 91% 9%



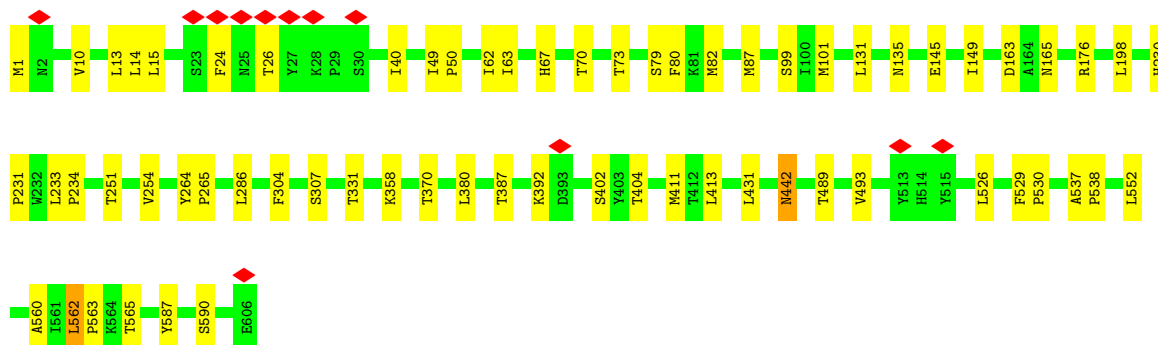
- Molecule 11: NADH-ubiquinone oxidoreductase chain 4L

Chain K: 88% 12%



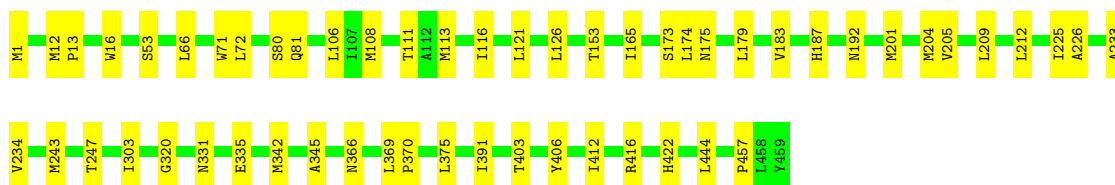
- Molecule 12: NADH-ubiquinone oxidoreductase chain 5

Chain L: 89% 11%



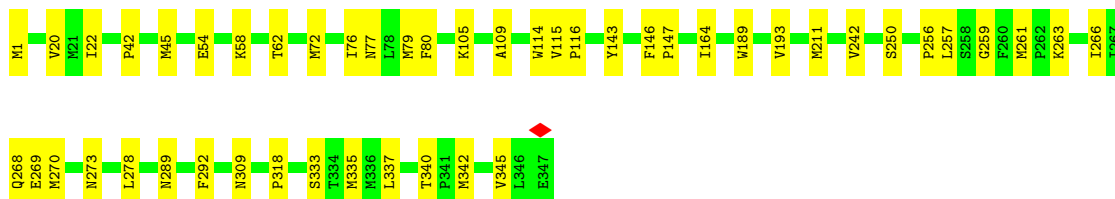
- Molecule 13: NADH-ubiquinone oxidoreductase chain 4

Chain M: 88% 12%



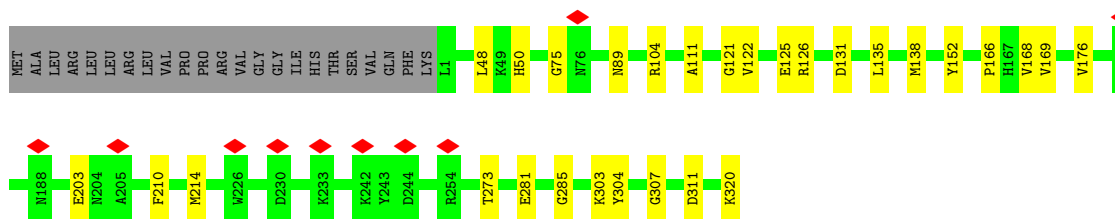
- Molecule 14: NADH-ubiquinone oxidoreductase chain 2

Chain N: 86% 14%



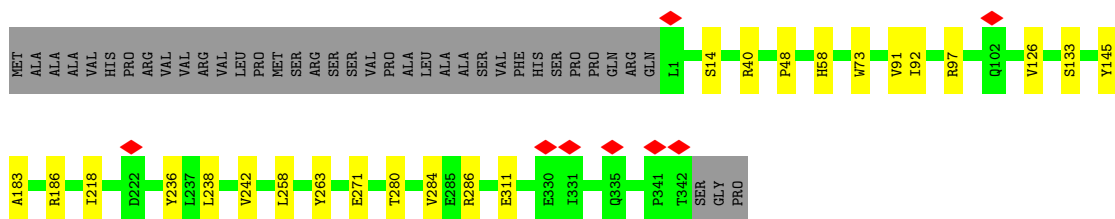
- Molecule 15: NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 10, mitochondrial

Chain O: 85% 8% 7%



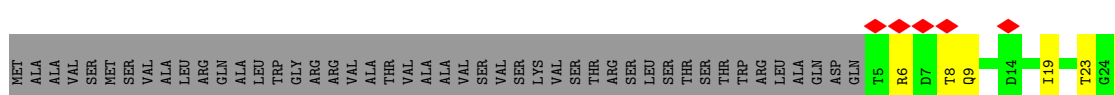
- Molecule 16: NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 9, mitochondrial

Chain P: 84% 6% 10%



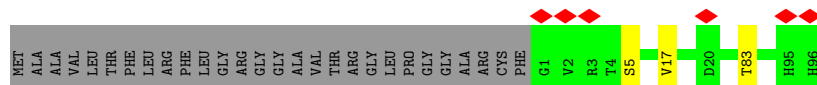
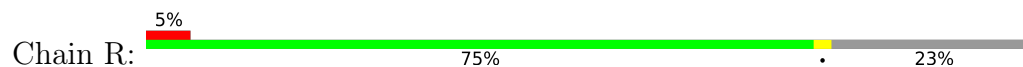
- Molecule 17: NADH dehydrogenase [ubiquinone] iron-sulfur protein 4, mitochondrial

Chain Q: 65% 9% 26%

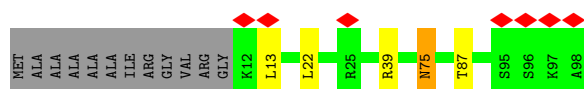
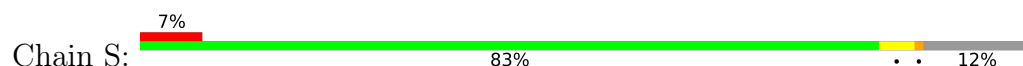




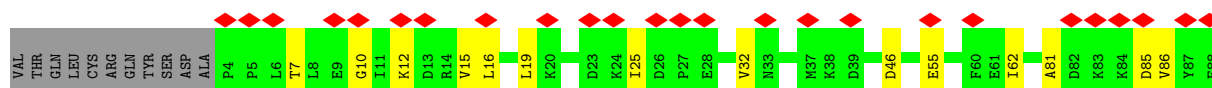
- Molecule 18: NADH dehydrogenase [ubiquinone] iron-sulfur protein 6, mitochondrial



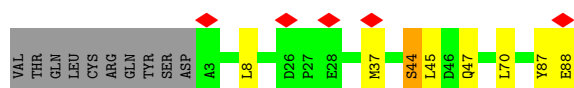
- Molecule 19: NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 2



- Molecule 20: Acyl carrier protein, mitochondrial



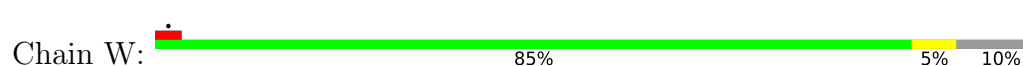
- Molecule 20: Acyl carrier protein, mitochondrial

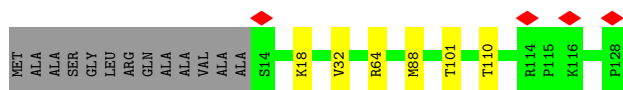


- Molecule 21: NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 5



- Molecule 22: NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 6





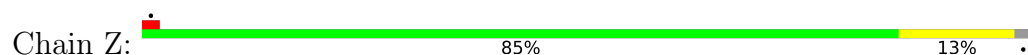
- Molecule 23: NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 8



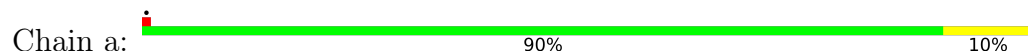
- Molecule 24: NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 11



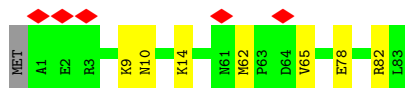
- Molecule 25: NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 13



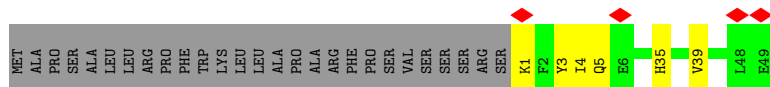
- Molecule 26: NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 1



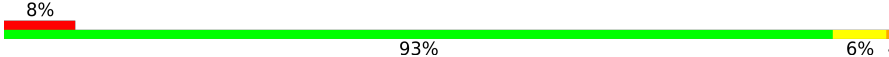
- Molecule 27: NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 3



- Molecule 28: NADH dehydrogenase [ubiquinone] 1 subunit C1, mitochondrial



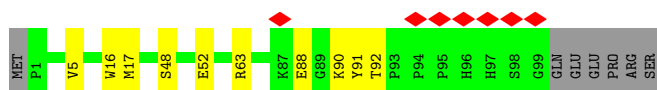
- Molecule 29: NADH dehydrogenase [ubiquinone] 1 subunit C2

Chain d: 

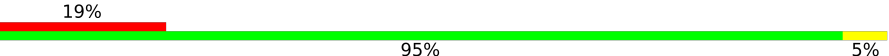


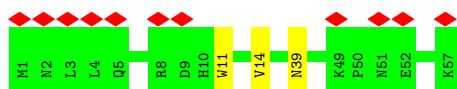
- Molecule 30: NADH dehydrogenase [ubiquinone] iron-sulfur protein 5

Chain e: 



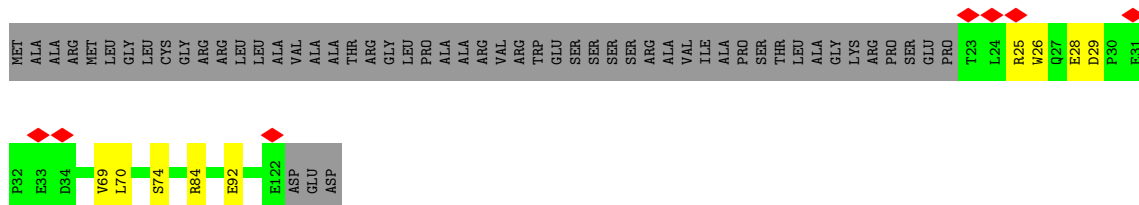
- Molecule 31: NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 1

Chain f: 



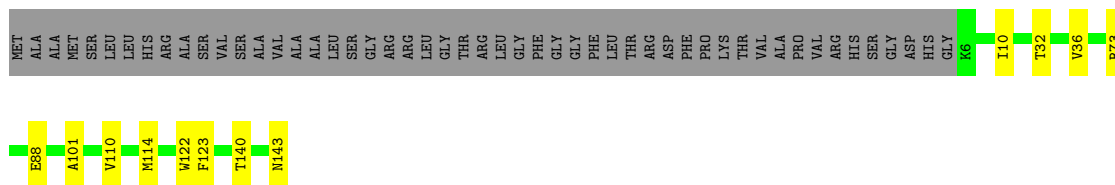
- Molecule 32: NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 11, mitochondrial

Chain g: 



- Molecule 33: NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 5, mitochondrial

Chain h: 

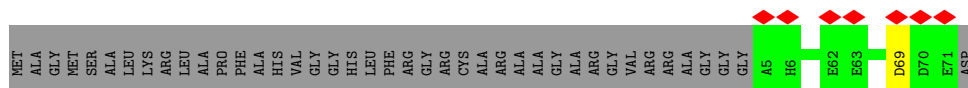


- Molecule 34: NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 6

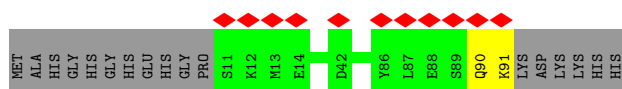
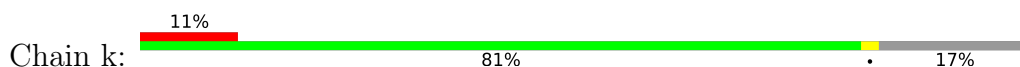
Chain i: 



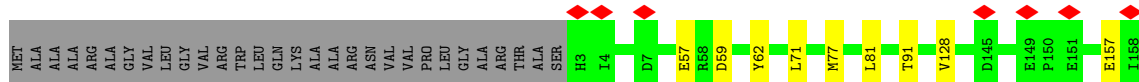
- Molecule 35: NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 2, mitochondrial



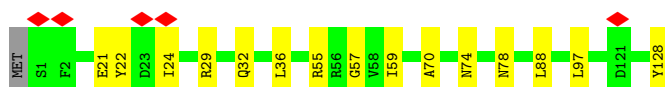
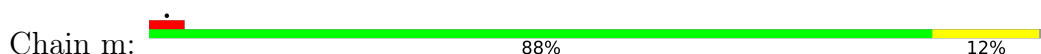
- Molecule 36: NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 3



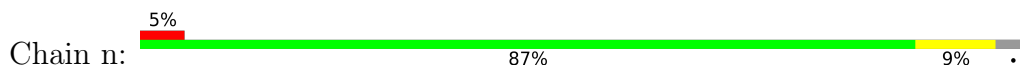
- Molecule 37: NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 8, mitochondrial



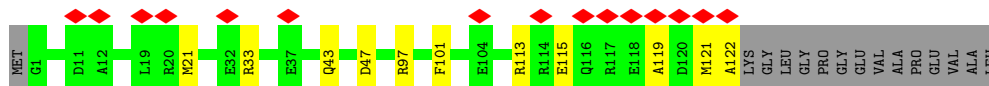
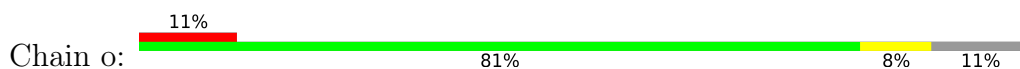
- Molecule 38: NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 4



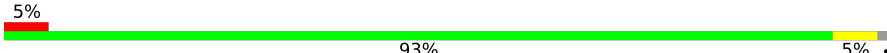
- Molecule 39: NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 9



- Molecule 40: NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 7

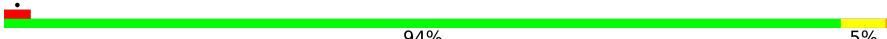


- Molecule 41: NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 10

Chain p:  5% 93% 5%



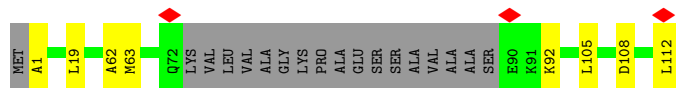
- Molecule 42: NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 12

Chain q:  5% 94% 5%



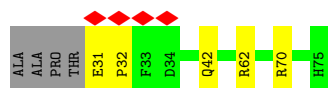
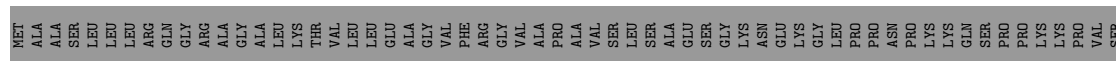
- Molecule 43: NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 7

Chain r:  5% 77% 7% 16%



- Molecule 44: NADH dehydrogenase [ubiquinone] flavoprotein 3, mitochondrial

Chain s:  5% 37% 5% 59%



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	38205	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	40.5	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2400	Depositor
Magnification	81000	Depositor
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	38.366	Depositor
Minimum map value	-15.492	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.938	Depositor
Recommended contour level	6.5	Depositor
Map size (Å)	479.744, 479.744, 479.744	wwPDB
Map dimensions	640, 640, 640	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.7496, 0.7496, 0.7496	Depositor

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: GTP, MYR, EHZ, AYA, NDP, SAC, FES, PC1, FME, AME, SF4, K, 3PE, ZN, CDL, 2MR, FMN, MG, CHD

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	0.39	0/936	0.40	0/1281
2	B	0.44	0/1272	0.49	0/1720
3	C	0.39	0/1789	0.44	2/2436 (0.1%)
4	D	0.40	0/3537	0.41	0/4794
5	E	0.35	0/1699	0.41	0/2312
6	F	0.36	0/3412	0.38	0/4610
7	G	0.39	0/5457	0.42	0/7397
8	H	0.38	0/2579	0.40	0/3524
9	I	0.43	0/1445	0.46	0/1956
10	J	0.35	0/1370	0.37	0/1859
11	K	0.36	0/745	0.43	0/1008
12	L	0.36	0/4920	0.36	0/6694
13	M	0.38	0/3738	0.38	0/5097
14	N	0.39	0/2792	0.40	0/3800
15	O	0.36	0/2651	0.35	0/3587
16	P	0.36	0/2847	0.42	0/3864
17	Q	0.39	0/1072	0.40	0/1449
18	R	0.37	0/753	0.34	0/1014
19	S	0.35	0/711	0.36	0/956
20	T	0.27	0/700	0.45	1/944 (0.1%)
20	U	0.37	0/705	0.35	0/952
21	V	0.34	0/948	0.33	0/1284
22	W	0.35	0/1000	0.39	0/1344
23	X	0.35	0/1439	0.41	0/1942
24	Y	0.34	0/1042	0.37	0/1414
25	Z	0.38	0/1181	0.37	0/1592
26	a	0.37	0/584	0.35	0/786
27	b	0.34	0/672	0.37	0/923
28	c	0.35	0/427	0.34	0/579
29	d	0.38	0/1018	0.39	0/1375
30	e	0.38	0/850	0.45	0/1136

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
31	f	0.31	0/505	0.42	0/681
32	g	0.36	0/865	0.46	0/1175
33	h	0.37	0/1188	0.37	0/1607
34	i	0.36	0/1127	0.50	0/1534
35	j	0.35	0/607	0.35	0/833
36	k	0.33	0/672	0.36	0/906
37	l	0.38	0/1369	0.37	0/1873
38	m	0.37	0/1094	0.38	0/1480
39	n	0.38	0/1545	0.37	0/2092
40	o	0.33	0/1073	0.40	0/1437
41	p	0.36	0/1486	0.38	0/2004
42	q	0.38	0/1250	0.37	0/1698
43	r	0.36	0/789	0.42	0/1068
44	s	0.36	0/403	0.46	0/545
All	All	0.37	0/68264	0.40	3/92562 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	120	SER	CA-C-N	7.07	124.70	120.24
3	C	120	SER	C-N-CA	7.07	124.70	120.24
20	T	86	VAL	CG1-CB-CG2	5.13	122.09	110.80

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	921	0	952	17	0
2	B	1241	0	1251	14	0
3	C	1738	0	1685	10	0
4	D	3459	0	3404	30	0
5	E	1659	0	1664	17	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	F	3336	0	3288	34	0
7	G	5366	0	5378	56	0
8	H	2517	0	2631	36	0
9	I	1414	0	1370	18	0
10	J	1345	0	1352	17	0
11	K	745	0	785	11	0
12	L	4802	0	4960	46	0
13	M	3654	0	3852	38	0
14	N	2733	0	2912	36	0
15	O	2589	0	2565	23	0
16	P	2768	0	2782	19	0
17	Q	1049	0	1045	12	0
18	R	740	0	714	2	0
19	S	700	0	719	5	0
20	T	688	0	684	10	0
20	U	693	0	688	5	0
21	V	928	0	972	3	0
22	W	976	0	991	6	0
23	X	1402	0	1381	14	0
24	Y	1030	0	1041	3	0
25	Z	1152	0	1151	19	0
26	a	569	0	568	8	0
27	b	651	0	662	5	0
28	c	414	0	415	3	0
29	d	999	0	988	12	0
30	e	829	0	829	10	0
31	f	492	0	501	2	0
32	g	839	0	790	7	0
33	h	1154	0	1168	11	0
34	i	1097	0	1108	14	0
35	j	580	0	519	1	0
36	k	653	0	639	2	0
37	l	1314	0	1210	6	0
38	m	1067	0	1067	14	0
39	n	1492	0	1438	15	0
40	o	1048	0	1016	8	0
41	p	1453	0	1425	6	0
42	q	1209	0	1182	15	0
43	r	776	0	782	7	0
44	s	391	0	361	4	0
45	A	79	0	109	0	0
45	H	29	0	32	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
45	I	38	0	50	0	0
45	L	84	0	119	0	0
45	M	194	0	296	6	0
45	N	48	0	70	0	0
45	P	35	0	44	0	0
45	Y	157	0	219	0	0
45	Z	35	0	44	0	0
45	b	47	0	71	0	0
45	d	37	0	48	0	0
45	m	40	0	57	2	0
46	A	68	0	84	0	0
46	B	41	0	56	1	0
46	M	81	0	113	1	0
46	P	54	0	88	1	0
46	Z	48	0	73	5	0
46	d	33	0	40	1	0
46	h	44	0	65	3	0
46	m	40	0	54	0	0
47	B	8	0	0	0	0
47	F	8	0	0	1	0
47	G	16	0	0	0	0
47	I	16	0	0	0	0
48	E	4	0	0	0	0
48	G	4	0	0	0	0
49	F	31	0	19	2	0
50	G	1	0	0	0	0
51	H	50	0	44	0	0
51	K	69	0	82	0	0
51	L	60	0	64	0	0
51	N	60	0	64	0	0
51	X	73	0	93	4	0
51	d	65	0	77	0	0
51	h	78	0	100	0	0
51	r	62	0	68	0	0
52	O	32	0	12	2	0
53	O	1	0	0	0	0
54	P	48	0	26	3	0
55	R	1	0	0	0	0
56	T	37	0	0	1	0
56	U	37	0	0	0	0
57	i	29	0	35	2	0
58	o	15	0	27	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
59	A	17	0	0	2	0
59	B	59	0	0	4	0
59	C	54	0	0	1	0
59	D	123	0	0	5	0
59	E	9	0	0	0	0
59	F	30	0	0	5	0
59	G	120	0	0	8	0
59	H	64	0	0	2	0
59	I	64	0	0	6	0
59	J	17	0	0	0	0
59	K	10	0	0	0	0
59	L	40	0	0	6	0
59	M	62	0	0	6	0
59	N	60	0	0	10	0
59	O	10	0	0	1	0
59	P	44	0	0	3	0
59	Q	51	0	0	2	0
59	R	14	0	0	1	0
59	U	1	0	0	0	0
59	W	6	0	0	0	0
59	X	22	0	0	2	0
59	Y	1	0	0	0	0
59	Z	23	0	0	1	0
59	a	15	0	0	2	0
59	b	2	0	0	0	0
59	d	10	0	0	0	0
59	e	13	0	0	0	0
59	g	5	0	0	0	0
59	h	11	0	0	0	0
59	i	1	0	0	0	0
59	k	1	0	0	0	0
59	l	7	0	0	0	0
59	m	5	0	0	2	0
59	n	10	0	0	0	0
59	p	12	0	0	0	0
59	q	18	0	0	1	0
59	r	11	0	0	0	0
59	s	2	0	0	0	0
All	All	69733	0	69328	549	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

All (549) 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)
25:Z:93:GLU:OE1	30:e:91:TYR:OH	1.71	1.05
1:A:68:GLU:OE1	1:A:98:LEU:HD13	1.65	0.96
7:G:260:GLU:OE1	59:G:901:HOH:O	1.84	0.96
23:X:91:SER:O	59:X:301:HOH:O	1.82	0.95
14:N:289:ASN:ND2	59:N:1002:HOH:O	2.06	0.87
13:M:81:GLN:OE1	59:M:701:HOH:O	1.93	0.86
8:H:311:THR:OG1	25:Z:51:MET:SD	2.33	0.86
6:F:200:GLN:NE2	59:F:605:HOH:O	2.10	0.85
4:D:145:THR:OG1	4:D:181:TYR:OH	1.95	0.85
23:X:84:TYR:OH	59:X:302:HOH:O	1.94	0.85
28:c:1:LYS:NZ	28:c:3:TYR:O	2.10	0.85
38:m:70:ALA:O	59:m:301:HOH:O	1.94	0.85
7:G:497:ALA:O	59:G:902:HOH:O	1.94	0.84
12:L:163:ASP:OD1	59:L:801:HOH:O	1.95	0.84
13:M:366:ASN:OD1	59:M:702:HOH:O	1.96	0.84
16:P:58:HIS:O	59:P:501:HOH:O	1.98	0.81
3:C:179:GLU:OE1	16:P:40:ARG:NH1	2.13	0.81
7:G:109:ASP:OD2	59:G:903:HOH:O	1.98	0.80
1:A:71:LEU:O	10:J:147:TYR:OH	2.00	0.80
7:G:366:THR:O	7:G:367:THR:OG1	1.99	0.80
13:M:303:ILE:O	59:M:703:HOH:O	2.00	0.79
26:a:46:TYR:OH	59:a:101:HOH:O	2.00	0.79
4:D:368:GLU:OE2	59:D:501:HOH:O	2.00	0.78
17:Q:132:THR:O	59:Q:201:HOH:O	2.02	0.78
1:A:26:GLN:O	59:A:301:HOH:O	2.00	0.78
14:N:143:TYR:O	59:N:1001:HOH:O	2.02	0.78
2:B:137:ASP:OD1	59:B:301:HOH:O	2.02	0.77
7:G:165:GLU:O	7:G:396:ARG:NH1	2.19	0.76
26:a:28:ARG:NH1	59:a:102:HOH:O	2.18	0.75
6:F:107:ASP:OD2	59:F:601:HOH:O	2.02	0.75
29:d:8:ARG:NH2	46:d:202:PC1:O14	2.19	0.75
6:F:98:ASP:OD2	59:F:603:HOH:O	2.05	0.74
49:F:501:FMN:O3P	59:F:602:HOH:O	2.04	0.74
7:G:194:GLU:OE2	59:G:904:HOH:O	2.04	0.74
2:B:170:GLU:OE1	2:B:172:ARG:NH2	2.21	0.74
3:C:152:LEU:O	59:C:301:HOH:O	2.05	0.74
27:b:10:ASN:OD1	27:b:14:LYS:NZ	2.21	0.73
8:H:24:GLU:HG3	8:H:271:LEU:HD22	1.72	0.72
15:O:273:THR:OG1	59:O:501:HOH:O	2.07	0.72
13:M:243:MET:O	13:M:247:THR:HG23	1.90	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:63:ALA:O	59:B:302:HOH:O	2.08	0.71
12:L:99:SER:OG	59:L:802:HOH:O	2.08	0.71
16:P:183:ALA:O	16:P:186:ARG:NH1	2.24	0.71
8:H:140:ILE:O	59:H:702:HOH:O	2.07	0.71
42:q:32:ASP:OD2	59:q:201:HOH:O	2.08	0.71
1:A:42:ASP:OD1	59:A:302:HOH:O	2.08	0.70
4:D:93:TYR:O	59:D:503:HOH:O	2.07	0.70
40:o:43:GLN:NE2	40:o:47:ASP:OD2	2.24	0.70
25:Z:21:TYR:OH	43:r:108:ASP:OD2	2.09	0.70
6:F:180:GLY:O	59:F:604:HOH:O	2.09	0.69
8:H:318:THR:HG23	26:a:41:TYR:CE2	2.27	0.69
12:L:145:GLU:OE1	59:L:803:HOH:O	2.10	0.69
16:P:14:SER:O	59:P:502:HOH:O	2.10	0.69
37:l:57:GLU:OE2	37:l:91:THR:OG1	2.11	0.69
46:B:202:PC1:O14	59:B:303:HOH:O	2.10	0.69
9:I:146:GLU:OE2	59:I:302:HOH:O	2.11	0.69
15:O:48:LEU:HD22	15:O:121:GLY:HA3	1.75	0.69
38:m:32:GLN:O	39:n:8:TYR:OH	2.10	0.68
17:Q:41:ALA:O	59:Q:202:HOH:O	2.12	0.68
14:N:256:PRO:O	59:N:1003:HOH:O	2.12	0.67
57:i:201:CHD:H212	57:i:201:CHD:O12	1.93	0.67
12:L:442:ASN:O	12:L:442:ASN:ND2	2.24	0.67
12:L:176:ARG:NH1	59:L:807:HOH:O	2.27	0.67
29:d:1:AME:HT23	29:d:1:AME:HB2	1.75	0.67
20:U:8:LEU:N	20:U:88:GLU:OE2	2.27	0.67
13:M:444:LEU:O	59:M:704:HOH:O	2.12	0.66
12:L:560:ALA:O	12:L:565:THR:OG1	2.14	0.66
18:R:83:THR:O	59:R:301:HOH:O	2.12	0.66
6:F:101:GLU:OE1	6:F:184:TYR:OH	2.07	0.66
9:I:117:ILE:O	59:I:303:HOH:O	2.14	0.66
13:M:345:ALA:O	59:M:705:HOH:O	2.12	0.66
16:P:97:ARG:NH1	54:P:402:NDP:O1A	2.29	0.66
23:X:44:LEU:HD22	23:X:130:VAL:CG1	2.26	0.65
9:I:124:GLU:OE2	59:I:304:HOH:O	2.15	0.65
34:i:6:GLU:OE2	39:n:157:ARG:NH2	2.30	0.64
2:B:116:MET:O	59:B:304:HOH:O	2.14	0.64
51:X:201:CDL:C33	51:X:201:CDL:H273	2.28	0.64
20:T:7:THR:O	20:T:10:GLY:N	2.31	0.63
42:q:144:TYR:HD2	42:q:145:LYS:HG3	1.63	0.63
8:H:195:ARG:HD3	8:H:231:ILE:HD11	1.80	0.63
14:N:79:MET:O	59:N:1004:HOH:O	2.15	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:G:534:ARG:CZ	42:q:145:LYS:HE2	2.29	0.63
20:T:81:ALA:O	20:T:85:ASP:N	2.32	0.63
15:O:138:MET:CE	52:O:401:GTP:HN21	2.11	0.62
12:L:370:THR:HG23	12:L:431:LEU:HD13	1.80	0.62
34:i:6:GLU:OE2	39:n:155:PRO:HD2	2.00	0.62
8:H:24:GLU:CG	8:H:271:LEU:HD22	2.29	0.61
16:P:48:PRO:HB2	16:P:73:TRP:CD1	2.34	0.61
12:L:358:LYS:NZ	39:n:27:GLU:OE1	2.33	0.61
12:L:251:THR:O	12:L:254:VAL:HG22	2.01	0.61
30:e:88:GLU:OE1	30:e:90:LYS:NZ	2.34	0.61
10:J:157:THR:HG22	11:K:66:PHE:HE1	1.66	0.60
14:N:318:PRO:O	59:N:1005:HOH:O	2.15	0.60
38:m:128:TYR:OH	59:m:302:HOH:O	2.15	0.60
9:I:5:VAL:HA	43:r:112:LEU:HD23	1.83	0.60
6:F:35:GLY:O	6:F:38:SER:OG	2.15	0.60
7:G:268:ARG:O	59:G:905:HOH:O	2.16	0.60
12:L:135:ASN:O	12:L:198:LEU:HD13	2.02	0.60
4:D:82:LEU:O	59:D:504:HOH:O	2.16	0.60
7:G:46:LEU:O	17:Q:116:LYS:NZ	2.22	0.59
12:L:380:LEU:O	12:L:392:LYS:NZ	2.21	0.59
3:C:113:GLU:N	3:C:113:GLU:OE1	2.36	0.59
8:H:236:ILE:HG23	8:H:259:PHE:HZ	1.68	0.59
12:L:14:LEU:CD2	34:i:73:HIS:O	2.51	0.59
13:M:106:LEU:HD13	13:M:234:VAL:HG11	1.84	0.59
14:N:22:ILE:O	59:N:1006:HOH:O	2.17	0.59
7:G:324:ASP:CB	7:G:571:ALA:HB1	2.33	0.58
16:P:236:TYR:OH	16:P:311:GLU:OE2	2.18	0.58
38:m:24:ILE:HD12	38:m:29:ARG:HD3	1.85	0.58
38:m:97:LEU:HD11	45:m:202:3PE:H2I3	1.86	0.58
19:S:39:ARG:NH1	19:S:87:THR:OG1	2.36	0.58
12:L:79:SER:OG	59:L:805:HOH:O	2.17	0.58
23:X:39:ASN:OD1	25:Z:73:PRO:CB	2.52	0.57
13:M:205:VAL:HG22	13:M:212:LEU:HD13	1.84	0.57
34:i:50:LEU:O	34:i:51:GLN:C	2.47	0.57
25:Z:125:TYR:HB3	25:Z:133:VAL:HG22	1.86	0.57
41:p:98:ASP:OD2	41:p:141:ARG:NH1	2.37	0.57
41:p:100:GLU:OE1	41:p:103:ASN:ND2	2.37	0.57
15:O:104:ARG:NE	15:O:131:ASP:OD2	2.38	0.57
12:L:73:THR:OG1	59:L:804:HOH:O	2.17	0.57
1:A:68:GLU:HG2	10:J:161:LEU:HD13	1.87	0.57
16:P:92:ILE:HD11	16:P:218:ILE:HD11	1.85	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
16:P:263:TYR:CD2	16:P:284:VAL:HG13	2.39	0.57
4:D:81:GLY:N	4:D:429:ASP:OD2	2.30	0.57
12:L:587:TYR:O	12:L:590:SER:OG	2.21	0.57
12:L:26:THR:O	12:L:26:THR:HG22	2.05	0.56
20:T:32:VAL:O	20:T:32:VAL:HG12	2.04	0.56
8:H:310:LEU:HD12	8:H:311:THR:HG23	1.87	0.56
32:g:70:LEU:O	32:g:74:SER:OG	2.18	0.56
8:H:264:LEU:HD11	26:a:13:GLY:HA2	1.86	0.56
13:M:225:ILE:HD13	13:M:331:ASN:HB2	1.87	0.56
5:E:145:LEU:HD12	5:E:153:MET:SD	2.45	0.56
12:L:402:SER:O	12:L:404:THR:HG22	2.05	0.56
18:R:5:SER:OG	18:R:17:VAL:HG21	2.06	0.55
9:I:140:SER:OG	59:I:301:HOH:O	2.06	0.55
7:G:35:MET:HE2	7:G:81:THR:HB	1.89	0.55
46:M:604:PC1:O13	46:M:604:PC1:H132	2.06	0.55
23:X:39:ASN:OD1	25:Z:73:PRO:HB2	2.06	0.55
5:E:105:THR:HG22	5:E:106:THR:H	1.72	0.55
41:p:2:ASP:OD1	41:p:3:SER:N	2.38	0.55
1:A:63:LEU:O	1:A:67:LEU:HD23	2.06	0.54
2:B:125:TYR:HB2	9:I:117:ILE:CG2	2.37	0.54
3:C:183:VAL:O	22:W:101:THR:OG1	2.18	0.54
7:G:538:PRO:O	7:G:539:LYS:HB3	2.08	0.54
4:D:349:PHE:CE2	9:I:82:LEU:HD11	2.43	0.54
7:G:140:LYS:O	7:G:148:THR:OG1	2.18	0.54
38:m:36:LEU:HB2	39:n:8:TYR:OH	2.07	0.54
44:s:31:GLU:N	44:s:32:PRO:HD2	2.22	0.54
14:N:80:PHE:HB3	30:e:63:ARG:HH22	1.73	0.54
14:N:20:VAL:O	59:N:1007:HOH:O	2.19	0.54
7:G:193:SER:OG	59:G:906:HOH:O	2.18	0.53
20:T:19:LEU:HD12	20:T:25:ILE:HD13	1.89	0.53
7:G:190:MET:HE3	7:G:695:ALA:HB2	1.89	0.53
37:l:77:MET:HE3	37:l:81:LEU:CD2	2.38	0.53
4:D:44:VAL:HG22	4:D:44:VAL:O	2.08	0.53
6:F:376:MET:O	6:F:380:VAL:HG23	2.09	0.53
15:O:135:LEU:HD22	15:O:152:TYR:CD1	2.44	0.53
23:X:35:CYS:O	23:X:39:ASN:HB2	2.09	0.53
33:h:73:ARG:NH2	46:h:202:PC1:O12	2.41	0.53
8:H:262:LYS:NZ	45:H:601:3PE:O14	2.37	0.53
25:Z:140:PHE:CZ	46:Z:201:PC1:H151	2.43	0.53
30:e:17:MET:O	30:e:17:MET:HG2	2.09	0.53
13:M:66:LEU:HD11	13:M:111:THR:CG2	2.39	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
45:M:602:3PE:H3A1	14:N:335:MET:HE1	1.90	0.52
5:E:217:LEU:CD1	6:F:236:ARG:HA	2.39	0.52
9:I:32:ARG:NH2	25:Z:26:PRO:O	2.40	0.52
15:O:138:MET:HE3	52:O:401:GTP:HN21	1.74	0.52
5:E:51:LEU:HD23	5:E:69:VAL:HG21	1.91	0.52
45:M:605:3PE:H3I2	45:M:605:3PE:C2H	2.39	0.52
4:D:322:GLU:OE1	4:D:322:GLU:N	2.30	0.52
29:d:1:AME:HE1	29:d:2:MET:CG	2.40	0.52
15:O:135:LEU:HD22	15:O:152:TYR:CG	2.45	0.52
29:d:1:AME:HT23	29:d:1:AME:CB	2.39	0.52
11:K:43:MET:HE1	14:N:72:MET:HE1	1.92	0.52
14:N:72:MET:HE3	14:N:76:ILE:HD11	1.92	0.52
16:P:258:LEU:HD23	16:P:263:TYR:HD1	1.74	0.52
15:O:48:LEU:HD22	15:O:121:GLY:CA	2.38	0.52
20:U:37:MET:HE1	20:U:44:SER:HA	1.92	0.52
7:G:366:THR:HB	7:G:450:MET:HE3	1.92	0.51
13:M:126:LEU:HD21	13:M:153:THR:HG21	1.91	0.51
34:i:24:ASP:OD1	39:n:120:LYS:CE	2.58	0.51
3:C:8:ARG:O	4:D:129:ARG:NH2	2.41	0.51
29:d:1:AME:HE2	33:h:122:TRP:CD2	2.46	0.51
1:A:68:GLU:CG	10:J:161:LEU:HD13	2.41	0.51
40:o:121:MET:HE2	40:o:121:MET:HA	1.91	0.51
23:X:58:GLU:OE1	23:X:58:GLU:N	2.41	0.51
8:H:2:PHE:CE2	8:H:6:ILE:HD11	2.45	0.51
13:M:342:MET:O	59:M:705:HOH:O	2.18	0.51
56:T:101:EHZ:O5	22:W:32:VAL:HG11	2.10	0.51
37:l:157:GLU:O	40:o:33:ARG:NH2	2.39	0.51
16:P:258:LEU:HD23	16:P:263:TYR:CD1	2.46	0.51
23:X:70:PHE:HB3	26:a:69:ILE:CD1	2.41	0.51
12:L:13:LEU:HD22	34:i:81:ILE:HD11	1.93	0.51
14:N:337:LEU:HD23	51:X:201:CDL:H181	1.92	0.51
7:G:283:MET:HE1	42:q:139:PRO:HG3	1.92	0.50
17:Q:25:VAL:HB	17:Q:30:ILE:HD11	1.94	0.50
6:F:224:ASN:ND2	49:F:501:FMN:O2	2.34	0.50
16:P:263:TYR:CE2	16:P:284:VAL:HG13	2.46	0.50
7:G:13:VAL:HG11	7:G:33:VAL:HG21	1.92	0.50
14:N:342:MET:O	14:N:345:VAL:HG22	2.11	0.50
39:n:175:GLU:HG3	39:n:176:ARG:HG2	1.92	0.50
13:M:201:MET:O	13:M:205:VAL:HG23	2.11	0.50
15:O:303:LYS:NZ	15:O:304:TYR:OH	2.45	0.50
30:e:16:TRP:CH2	30:e:17:MET:HE3	2.47	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:343:GLU:OE2	4:D:343:GLU:N	2.45	0.50
25:Z:68:ARG:O	25:Z:72:MET:HG3	2.12	0.50
7:G:329:ILE:HD11	7:G:626:VAL:HG21	1.93	0.50
12:L:331:THR:HB	12:L:387:THR:HG22	1.93	0.50
16:P:14:SER:OG	59:P:503:HOH:O	2.19	0.50
6:F:175:VAL:O	44:s:62:ARG:NH1	2.39	0.49
4:D:354:GLU:OE2	4:D:357:GLN:NE2	2.45	0.49
13:M:331:ASN:O	13:M:335:GLU:HG2	2.12	0.49
4:D:296:ARG:NH1	59:D:512:HOH:O	2.40	0.49
12:L:80:PHE:HB3	12:L:82:MET:HE2	1.93	0.49
7:G:289:GLY:C	42:q:144:TYR:HE1	2.21	0.49
8:H:114:TYR:OH	10:J:61:LEU:O	2.30	0.49
15:O:303:LYS:NZ	15:O:304:TYR:CZ	2.80	0.49
1:A:13:LEU:HD13	8:H:10:ILE:HD12	1.95	0.49
4:D:61:VAL:O	59:D:505:HOH:O	2.20	0.49
37:l:59:ASP:OD2	37:l:62:TYR:N	2.46	0.49
14:N:62:THR:HG21	14:N:114:TRP:CD1	2.47	0.49
6:F:370:ASP:OD1	7:G:179:ASN:ND2	2.39	0.48
7:G:226:GLU:OE1	59:G:907:HOH:O	2.20	0.48
9:I:80:CYS:SG	9:I:82:LEU:HD13	2.53	0.48
13:M:370:PRO:HB3	13:M:375:LEU:CD2	2.42	0.48
34:i:51:GLN:O	34:i:52:ASP:OD1	2.31	0.48
7:G:690:ALA:HA	7:G:698:VAL:HG21	1.95	0.48
13:M:187:HIS:O	13:M:192:ASN:ND2	2.38	0.48
16:P:238:LEU:O	16:P:242:VAL:HG23	2.13	0.48
7:G:285:ARG:NH2	42:q:144:TYR:CD1	2.81	0.48
20:T:19:LEU:HD12	20:T:25:ILE:CD1	2.44	0.48
1:A:66:ASP:O	1:A:69:ILE:HG12	2.13	0.48
7:G:537:LEU:CD1	7:G:543:ILE:HD11	2.43	0.48
8:H:89:LEU:HD13	8:H:236:ILE:HG21	1.95	0.48
9:I:53:GLU:OE2	42:q:34:ARG:NH2	2.47	0.48
9:I:123:GLN:OE1	59:I:306:HOH:O	2.20	0.48
42:q:144:TYR:O	42:q:145:LYS:C	2.56	0.48
7:G:35:MET:HE3	7:G:37:ILE:HD13	1.96	0.48
13:M:16:TRP:CZ2	45:M:602:3PE:H271	2.48	0.48
13:M:12:MET:HB2	13:M:13:PRO:HD3	1.95	0.48
37:l:71:LEU:HD11	37:l:81:LEU:HD21	1.96	0.48
6:F:306:LEU:C	6:F:306:LEU:HD12	2.39	0.48
8:H:211:PHE:O	59:H:703:HOH:O	2.19	0.48
6:F:103:GLY:O	6:F:333:ALA:HB1	2.14	0.48
7:G:194:GLU:OE1	7:G:194:GLU:N	2.44	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
30:e:92:THR:HG23	30:e:92:THR:O	2.14	0.48
38:m:22:TYR:OH	39:n:67:GLU:HG3	2.13	0.48
3:C:148:LEU:O	22:W:88:MET:HE2	2.14	0.47
12:L:304:PHE:CZ	12:L:526:LEU:HD22	2.49	0.47
14:N:77:ASN:ND2	59:N:1011:HOH:O	2.40	0.47
33:h:140:THR:O	33:h:143:ASN:ND2	2.47	0.47
36:k:90:GLN:O	36:k:91:LYS:C	2.57	0.47
34:i:56:TRP:HD1	34:i:57:LYS:HG3	1.79	0.47
13:M:403:THR:HA	13:M:406:TYR:CE1	2.50	0.47
13:M:457:PRO:O	32:g:84:ARG:NH2	2.48	0.47
1:A:95:ILE:HG21	8:H:302:MET:HG3	1.96	0.47
7:G:272:ASP:OD1	7:G:681:SER:OG	2.25	0.47
8:H:14:LEU:HD22	8:H:225:MET:HE3	1.97	0.47
12:L:489:THR:O	12:L:493:VAL:HG22	2.15	0.47
16:P:271:GLU:HG2	16:P:280:THR:HG22	1.96	0.47
17:Q:8:THR:O	17:Q:9:GLN:HB3	2.15	0.47
20:T:46:ASP:OD1	22:W:64:ARG:NH2	2.47	0.47
23:X:129:LYS:NZ	27:b:62:MET:O	2.38	0.47
29:d:101:PHE:CZ	33:h:110:VAL:HG22	2.49	0.47
10:J:157:THR:HG21	11:K:62:ILE:HD12	1.96	0.47
11:K:42:VAL:O	11:K:46:LEU:HD13	2.14	0.47
24:Y:81:GLU:O	24:Y:81:GLU:HG3	2.14	0.47
25:Z:47:TYR:OH	59:Z:401:HOH:O	2.05	0.47
13:M:165:ILE:HG21	14:N:268:GLN:HA	1.96	0.47
17:Q:19:ILE:O	17:Q:23:THR:HG23	2.14	0.47
1:A:73:LEU:N	1:A:74:PRO:CD	2.78	0.47
4:D:73:VAL:HG21	4:D:414:VAL:HG21	1.96	0.47
7:G:255:HIS:CD2	17:Q:109:LYS:CE	2.98	0.47
8:H:23:VAL:HG11	8:H:268:MET:HE1	1.97	0.47
40:o:21:MET:HE1	40:o:101:PHE:CD2	2.50	0.47
13:M:71:TRP:CZ3	32:g:69:VAL:HG11	2.50	0.47
17:Q:90:GLU:OE1	17:Q:90:GLU:N	2.43	0.47
34:i:24:ASP:OD1	39:n:120:LYS:HE2	2.15	0.47
4:D:335:ARG:NH2	9:I:129:ASP:OD1	2.49	0.46
12:L:230:HIS:N	12:L:231:PRO:CD	2.78	0.46
14:N:292:PHE:O	59:N:1008:HOH:O	2.20	0.46
9:I:176:ARG:NH2	59:I:309:HOH:O	2.45	0.46
12:L:442:ASN:HD22	12:L:442:ASN:C	2.19	0.46
14:N:115:VAL:HB	14:N:116:PRO:HD3	1.97	0.46
4:D:302:GLU:OE1	43:r:112:LEU:HD11	2.16	0.46
5:E:149:VAL:HG13	6:F:105:CYS:SG	2.55	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:F:141:GLU:OE1	6:F:141:GLU:N	2.46	0.46
21:V:57:MET:HE1	21:V:70:ARG:O	2.15	0.46
32:g:92:GLU:HG2	41:p:97:VAL:HG21	1.96	0.46
33:h:32:THR:O	33:h:36:VAL:HG23	2.14	0.46
2:B:45:LEU:O	2:B:74:VAL:HA	2.15	0.46
5:E:172:ILE:HG23	5:E:182:PRO:HG3	1.97	0.46
9:I:2:TYR:CE2	43:r:105:LEU:HD23	2.50	0.46
6:F:79:TRP:CD2	6:F:129:MET:HE2	2.51	0.46
29:d:39:TYR:CE2	29:d:43:LEU:HD11	2.51	0.46
10:J:132:ASP:OD1	10:J:133:SER:N	2.49	0.46
14:N:105:LYS:NZ	59:N:1009:HOH:O	2.23	0.46
20:T:55:GLU:HG2	20:T:62:ILE:HD12	1.98	0.46
46:h:202:PC1:H291	46:h:202:PC1:H2D1	1.98	0.46
16:P:91:VAL:HG23	16:P:126:VAL:HG11	1.98	0.46
7:G:26:VAL:HG22	7:G:73:VAL:HG12	1.97	0.46
10:J:5:ILE:HG21	25:Z:142:TRP:HZ2	1.81	0.46
7:G:324:ASP:HB2	7:G:571:ALA:HB1	1.98	0.45
7:G:382:THR:OG1	7:G:387:GLU:OE1	2.34	0.45
3:C:75:LEU:HD12	3:C:96:LEU:HD23	1.98	0.45
3:C:154:ASP:OD1	3:C:155:TYR:N	2.49	0.45
13:M:204:MET:SD	13:M:209:LEU:HD13	2.56	0.45
3:C:58:ILE:HB	3:C:59:PRO:HD3	1.98	0.45
4:D:269:LEU:HB2	4:D:368:GLU:HB2	1.98	0.45
5:E:217:LEU:HD23	6:F:44:LYS:HB3	1.97	0.45
12:L:24:PHE:O	12:L:24:PHE:CD1	2.70	0.45
12:L:233:LEU:HD23	12:L:307:SER:OG	2.17	0.45
34:i:55:LEU:HD23	34:i:56:TRP:N	2.31	0.45
7:G:36:GLN:OE1	17:Q:49:VAL:HG22	2.15	0.45
10:J:157:THR:HG22	11:K:66:PHE:CE1	2.49	0.45
14:N:80:PHE:O	30:e:63:ARG:NH2	2.50	0.45
14:N:189:TRP:CZ2	14:N:263:LYS:HE2	2.51	0.45
21:V:57:MET:CE	21:V:70:ARG:HB3	2.46	0.45
29:d:1:AME:HE1	29:d:2:MET:HG3	1.98	0.45
32:g:25:ARG:O	32:g:25:ARG:NE	2.48	0.45
6:F:46:LYS:O	6:F:50:LEU:HD23	2.17	0.45
8:H:232:ILE:HG21	26:a:16:LEU:HD21	1.97	0.45
10:J:139:GLU:HG3	11:K:49:LEU:HD21	1.98	0.45
15:O:311:ASP:OD1	15:O:311:ASP:N	2.48	0.45
29:d:1:AME:HB2	29:d:1:AME:CT2	2.45	0.45
40:o:121:MET:O	40:o:122:ALA:HB2	2.15	0.45
7:G:362:TYR:OH	7:G:504:ASP:OD1	2.31	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
15:O:111:ALA:HB1	15:O:122:VAL:HG21	1.97	0.45
25:Z:96:THR:O	25:Z:99:LYS:NZ	2.34	0.45
2:B:125:TYR:CE2	4:D:102:TYR:CD1	3.05	0.45
7:G:262:TRP:HB2	7:G:390:LEU:HD11	1.98	0.45
7:G:489:VAL:O	7:G:491:ASN:ND2	2.50	0.45
12:L:165:ASN:OD1	13:M:416:ARG:NH1	2.49	0.45
12:L:233:LEU:HB3	12:L:234:PRO:HD3	1.98	0.45
13:M:108:MET:CB	13:M:121:LEU:HD13	2.47	0.45
15:O:210:PHE:CE1	15:O:214:MET:HG3	2.51	0.45
5:E:106:THR:HB	5:E:107:PRO:HD3	1.97	0.45
8:H:19:PHE:O	8:H:23:VAL:HG23	2.16	0.45
51:X:201:CDL:H273	51:X:201:CDL:H331	1.97	0.45
27:b:78:GLU:OE1	27:b:82:ARG:NE	2.50	0.45
14:N:211:MET:HG2	14:N:333:SER:HB2	1.99	0.44
25:Z:140:PHE:O	46:Z:201:PC1:H142	2.18	0.44
29:d:1:AME:HG2	33:h:123:PHE:CZ	2.51	0.44
39:n:162:LEU:HD12	39:n:162:LEU:N	2.31	0.44
1:A:93:PHE:O	1:A:97:LEU:HD13	2.18	0.44
12:L:149:ILE:HG13	13:M:369:LEU:HD13	1.99	0.44
2:B:69:MET:HE2	2:B:74:VAL:HG12	1.98	0.44
12:L:562:LEU:CB	12:L:563:PRO:CD	2.95	0.44
13:M:108:MET:HE1	51:X:201:CDL:H222	2.00	0.44
28:c:35:HIS:O	28:c:39:VAL:HG23	2.17	0.44
1:A:25:PRO:HG2	8:H:60:PRO:HD3	1.99	0.44
6:F:97:ALA:HB3	6:F:137:TYR:O	2.18	0.44
7:G:534:ARG:NH1	42:q:145:LYS:HE2	2.31	0.44
9:I:43:ARG:HH21	43:r:19:LEU:CD1	2.30	0.44
14:N:109:ALA:HB1	14:N:164:ILE:HD12	2.00	0.44
23:X:129:LYS:HE3	27:b:65:VAL:O	2.18	0.44
2:B:44:SER:OG	8:H:51:ASP:OD1	2.27	0.44
2:B:91:THR:HG21	4:D:85:2MR:HG3	1.99	0.44
6:F:344:VAL:HG12	6:F:380:VAL:HG22	2.00	0.44
7:G:568:GLU:HA	7:G:587:VAL:O	2.18	0.44
12:L:40:ILE:HG13	12:L:101:MET:SD	2.58	0.44
15:O:166:PRO:HD2	15:O:214:MET:HE1	1.99	0.44
23:X:35:CYS:O	23:X:39:ASN:CB	2.66	0.44
7:G:372:GLU:OE2	7:G:394:ARG:NH1	2.46	0.44
11:K:19:LEU:HD22	11:K:36:MET:CE	2.48	0.44
12:L:286:LEU:HD22	12:L:411:MET:SD	2.58	0.44
38:m:74:ASN:OD1	38:m:78:ASN:ND2	2.51	0.44
42:q:144:TYR:HD2	42:q:145:LYS:CG	2.30	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
43:r:62:ALA:O	43:r:63:MET:HE2	2.17	0.44
4:D:218:PHE:CE2	4:D:222:ILE:HD11	2.52	0.44
5:E:165:THR:CG2	5:E:166:PRO:HD2	2.47	0.44
6:F:193:ILE:HG23	6:F:215:VAL:HA	2.00	0.44
10:J:153:LEU:O	10:J:157:THR:HG23	2.18	0.44
11:K:55:LEU:O	30:e:17:MET:HE2	2.18	0.44
15:O:307:GLY:O	15:O:320:LYS:NZ	2.46	0.44
40:o:115:GLU:O	40:o:119:ALA:N	2.45	0.44
1:A:24:LEU:N	1:A:25:PRO:CD	2.81	0.43
6:F:306:LEU:HD13	6:F:422:PHE:HE2	1.83	0.43
7:G:285:ARG:HD2	42:q:144:TYR:CE1	2.53	0.43
7:G:571:ALA:O	7:G:582:GLN:HA	2.18	0.43
8:H:142:TYR:CD1	8:H:142:TYR:C	2.96	0.43
10:J:171:ILE:HG23	14:N:45:MET:HE2	2.00	0.43
20:T:32:VAL:O	20:T:32:VAL:CG1	2.66	0.43
45:M:603:3PE:H2I1	14:N:257:LEU:HD21	1.99	0.43
31:f:39:ASN:N	33:h:88:GLU:OE1	2.51	0.43
2:B:71:ARG:HA	8:H:37:PRO:HA	2.00	0.43
10:J:1:FME:SD	10:J:2:MET:N	2.89	0.43
13:M:173:SER:HB2	33:h:101:ALA:HB2	2.00	0.43
13:M:422:HIS:HB2	38:m:59:ILE:HD12	2.00	0.43
31:f:11:TRP:O	31:f:14:VAL:HG22	2.17	0.43
7:G:255:HIS:CE1	7:G:258:ILE:HG13	2.53	0.43
9:I:14:MET:SD	27:b:9:LYS:NZ	2.89	0.43
4:D:106:LEU:HD11	4:D:391:ILE:HD13	2.00	0.43
6:F:194:GLU:CD	6:F:204:ARG:HE	2.27	0.43
5:E:217:LEU:HD23	6:F:44:LYS:CB	2.48	0.43
6:F:131:ALA:O	6:F:171:TYR:OH	2.16	0.43
7:G:199:ILE:HA	7:G:202:ILE:HG12	2.00	0.43
8:H:285:LEU:HD12	8:H:289:LEU:HD23	2.01	0.43
10:J:123:GLY:O	10:J:126:VAL:HG22	2.18	0.43
35:j:69:ASP:OD2	40:o:113:ARG:NH1	2.52	0.43
39:n:154:PRO:O	39:n:164:PRO:HD2	2.19	0.43
4:D:94:LYS:HE3	9:I:88:PRO:O	2.18	0.43
25:Z:47:TYR:O	25:Z:51:MET:HG2	2.18	0.43
33:h:36:VAL:HG22	46:h:202:PC1:H2D2	2.00	0.43
5:E:103:CYS:SG	5:E:144:CYS:HA	2.59	0.43
7:G:285:ARG:HD2	42:q:144:TYR:CZ	2.53	0.43
7:G:349:PHE:CZ	7:G:362:TYR:HB3	2.54	0.43
10:J:1:FME:SD	10:J:1:FME:N	2.92	0.43
12:L:82:MET:HB3	12:L:87:MET:HE3	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:N:269:GLU:O	14:N:273:ASN:ND2	2.45	0.43
23:X:79:GLU:HB3	23:X:80:PRO:HD3	2.01	0.43
8:H:316:PRO:HG3	25:Z:57:ARG:HB3	2.00	0.43
11:K:56:ALA:O	11:K:60:PRO:HD3	2.19	0.43
45:M:603:3PE:C33	14:N:242:VAL:HG11	2.49	0.43
25:Z:86:MET:HE1	25:Z:125:TYR:CE2	2.53	0.43
38:m:21:GLU:O	39:n:14:LYS:NZ	2.52	0.43
6:F:94:VAL:HG11	6:F:192:LEU:HD22	2.00	0.43
8:H:31:MET:HE1	8:H:272:TRP:CD1	2.54	0.43
12:L:62:ILE:HG21	12:L:135:ASN:ND2	2.33	0.43
13:M:53:SER:O	13:M:113:MET:HE1	2.19	0.43
19:S:75:ASN:O	19:S:75:ASN:ND2	2.40	0.43
20:U:45:LEU:HD23	39:n:16:LEU:CD2	2.49	0.43
4:D:141:PHE:HD2	4:D:181:TYR:HE1	1.67	0.42
7:G:272:ASP:O	7:G:275:LYS:HG2	2.18	0.42
9:I:40:TYR:CZ	9:I:43:ARG:NH1	2.87	0.42
15:O:50:HIS:NE2	15:O:125:GLU:OE2	2.44	0.42
15:O:285:GLY:HA2	32:g:26:TRP:CZ3	2.53	0.42
2:B:152:THR:HG22	4:D:188:ARG:HD3	2.01	0.42
4:D:17:ALA:HB1	11:K:98:CYS:SG	2.59	0.42
12:L:49:ILE:HB	12:L:50:PRO:HD3	2.02	0.42
12:L:537:ALA:HB3	12:L:538:PRO:HD3	2.01	0.42
15:O:125:GLU:O	15:O:126:ARG:HB2	2.20	0.42
16:P:97:ARG:HE	54:P:402:NDP:H8A	1.84	0.42
3:C:186:GLU:OE2	22:W:110:THR:OG1	2.29	0.42
7:G:20:VAL:HG21	7:G:73:VAL:HG21	2.00	0.42
7:G:255:HIS:ND1	7:G:258:ILE:HG13	2.35	0.42
7:G:551:ASP:N	59:G:916:HOH:O	2.44	0.42
8:H:85:MET:SD	8:H:108:MET:HB2	2.59	0.42
20:T:12:LYS:O	20:T:16:LEU:HD23	2.19	0.42
38:m:24:ILE:O	38:m:24:ILE:HG13	2.19	0.42
38:m:55:ARG:NH1	38:m:57:GLY:O	2.52	0.42
12:L:87:MET:HA	12:L:87:MET:HE2	2.01	0.42
13:M:113:MET:HG2	13:M:175:ASN:OD1	2.20	0.42
14:N:309:ASN:HD21	15:O:75:GLY:HA3	1.83	0.42
19:S:22:LEU:HD23	19:S:22:LEU:N	2.35	0.42
20:U:47:GLN:NE2	20:U:70:LEU:O	2.51	0.42
4:D:62:LEU:HB2	4:D:425:PHE:CZ	2.54	0.42
5:E:101:GLN:HG2	5:E:140:ILE:HD11	2.01	0.42
14:N:193:VAL:HB	14:N:266:ILE:HG23	2.01	0.42
57:i:201:CHD:H232	57:i:201:CHD:H211	1.81	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:59:HIS:CD2	4:D:59:HIS:N	2.86	0.42
8:H:56:PHE:CD2	46:P:401:PC1:C3H	3.02	0.42
29:d:106:LYS:HG2	41:p:77:GLN:HB3	2.01	0.42
37:l:128:VAL:HG22	40:o:97:ARG:HB3	2.01	0.42
7:G:289:GLY:O	42:q:144:TYR:CE1	2.73	0.42
7:G:526:GLY:N	7:G:546:GLN:O	2.49	0.42
10:J:167:VAL:HG22	14:N:42:PRO:HG3	2.00	0.42
41:p:16:THR:O	41:p:16:THR:HG23	2.18	0.42
5:E:87:TYR:CG	6:F:181:ALA:HB3	2.55	0.42
11:K:19:LEU:CD2	11:K:36:MET:HE1	2.50	0.42
25:Z:141:ILE:HA	46:Z:201:PC1:H153	2.02	0.42
4:D:143:GLU:HG3	4:D:310:ILE:HG21	2.02	0.42
4:D:155:THR:HB	4:D:167:PHE:HA	2.01	0.42
7:G:239:VAL:HG23	7:G:253:ARG:HB2	2.01	0.42
14:N:250:SER:O	14:N:259:GLY:HA3	2.18	0.42
25:Z:141:ILE:O	46:Z:201:PC1:H121	2.20	0.42
5:E:109:MET:HE1	6:F:346:ALA:HA	2.00	0.42
6:F:192:LEU:C	6:F:192:LEU:HD23	2.44	0.42
8:H:87:ILE:CG1	8:H:88:PRO:HD3	2.49	0.42
8:H:87:ILE:N	8:H:88:PRO:CD	2.83	0.42
17:Q:60:ASP:OD1	17:Q:60:ASP:N	2.52	0.42
2:B:69:MET:HE1	2:B:76:PHE:CZ	2.54	0.41
8:H:24:GLU:OE2	8:H:274:ARG:NH1	2.52	0.41
12:L:15:LEU:HD22	12:L:40:ILE:HD13	2.02	0.41
14:N:261:MET:HG3	14:N:340:THR:HG23	2.00	0.41
16:P:145:TYR:CD2	16:P:286:ARG:HD3	2.54	0.41
24:Y:28:GLY:O	24:Y:62:ALA:HB1	2.20	0.41
30:e:48:SER:O	30:e:52:GLU:HG3	2.19	0.41
33:h:110:VAL:CG1	33:h:114:MET:HE3	2.50	0.41
1:A:78:ALA:O	1:A:87:MET:SD	2.78	0.41
12:L:413:LEU:HD22	12:L:489:THR:HG22	2.02	0.41
13:M:80:SER:OG	13:M:226:ALA:HB2	2.20	0.41
14:N:146:PHE:N	14:N:147:PRO:CD	2.83	0.41
29:d:11:PHE:CE2	30:e:5:VAL:HG12	2.55	0.41
34:i:75:LEU:O	34:i:78:VAL:HG12	2.19	0.41
34:i:85:LEU:HD23	34:i:89:VAL:HG21	2.02	0.41
39:n:99:PRO:HB2	39:n:101:TRP:CD1	2.55	0.41
42:q:143:PRO:C	42:q:144:TYR:CD1	2.98	0.41
5:E:217:LEU:HD23	6:F:44:LYS:HD3	2.02	0.41
13:M:412:ILE:HA	13:M:416:ARG:HD2	2.02	0.41
8:H:264:LEU:HD11	26:a:13:GLY:CA	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:L:529:PHE:HB3	12:L:530:PRO:HD3	2.01	0.41
45:M:603:3PE:H331	14:N:242:VAL:HG11	2.02	0.41
12:L:552:LEU:HD21	38:m:88:LEU:C	2.46	0.41
15:O:89:ASN:O	15:O:89:ASN:OD1	2.39	0.41
33:h:10:ILE:HG23	39:n:178:MET:HG2	2.02	0.41
7:G:145:LEU:HD22	7:G:269:PHE:CG	2.56	0.41
17:Q:6:ARG:NH2	22:W:18:LYS:HD2	2.35	0.41
25:Z:13:VAL:HG13	25:Z:14:GLY:N	2.36	0.41
28:c:4:ILE:HG23	28:c:5:GLN:HG3	2.02	0.41
4:D:201:GLN:C	4:D:323:ILE:CD1	2.93	0.41
5:E:134:ASP:O	5:E:135:LYS:HB2	2.20	0.41
12:L:10:VAL:O	12:L:14:LEU:HG	2.21	0.41
15:O:176:VAL:HG23	15:O:203:GLU:OE1	2.20	0.41
23:X:7:PRO:HG2	23:X:56:LEU:HD11	2.03	0.41
46:Z:201:PC1:H292	46:Z:201:PC1:H331	2.02	0.41
1:A:65:PHE:CD2	1:A:98:LEU:HD11	2.55	0.41
12:L:63:ILE:O	12:L:79:SER:HA	2.21	0.41
13:M:179:LEU:O	13:M:183:VAL:HB	2.21	0.41
36:k:90:GLN:O	36:k:90:GLN:CD	2.64	0.41
2:B:91:THR:HA	2:B:119:CYS:HB3	2.02	0.41
6:F:14:SER:N	6:F:270:GLU:OE2	2.54	0.41
7:G:486:ASP:OD1	7:G:486:ASP:O	2.39	0.41
8:H:207:LEU:O	8:H:209:SER:N	2.53	0.41
10:J:51:PHE:HB3	10:J:139:GLU:OE2	2.21	0.41
13:M:72:LEU:HD11	13:M:233:ALA:HB1	2.02	0.41
13:M:233:ALA:HA	13:M:320:GLY:HA2	2.03	0.41
13:M:391:ILE:O	13:M:391:ILE:HG22	2.20	0.41
14:N:54:GLU:OE1	14:N:58:LYS:NZ	2.51	0.41
15:O:168:VAL:HG12	15:O:169:VAL:N	2.36	0.41
20:U:87:TYR:OH	34:i:26:GLU:OE1	2.29	0.41
32:g:28:GLU:HG3	32:g:29:ASP:N	2.36	0.41
38:m:97:LEU:CD1	45:m:202:3PE:H2I3	2.51	0.41
6:F:268:VAL:HG22	6:F:269:GLU:N	2.36	0.41
6:F:403:THR:HB	47:F:502:SF4:S4	2.60	0.41
16:P:133:SER:O	54:P:402:NDP:H6N	2.21	0.41
34:i:49:PHE:O	34:i:57:LYS:HD3	2.20	0.41
12:L:264:TYR:N	12:L:265:PRO:CD	2.85	0.40
5:E:150:ASN:HB3	5:E:162:GLU:HB3	2.02	0.40
19:S:13:LEU:HG	19:S:13:LEU:O	2.21	0.40
19:S:75:ASN:HD22	19:S:75:ASN:C	2.27	0.40
1:A:62:PHE:O	1:A:66:ASP:HB2	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:H:179:TRP:CG	8:H:180:PRO:HD3	2.56	0.40
17:Q:127:ARG:HE	44:s:70:ARG:HH12	1.69	0.40
20:T:15:VAL:O	20:T:19:LEU:HD23	2.22	0.40
6:F:381:ARG:NH2	6:F:383:ASP:OD2	2.55	0.40
8:H:120:GLY:HA3	8:H:132:ALA:HB2	2.03	0.40
12:L:131:LEU:HD23	12:L:131:LEU:C	2.46	0.40
13:M:116:ILE:HG12	13:M:174:LEU:CD1	2.52	0.40
14:N:270:MET:CE	14:N:278:LEU:HD23	2.52	0.40
15:O:281:GLU:OE1	15:O:281:GLU:N	2.53	0.40
23:X:23:VAL:HG21	26:a:65:GLY:HA2	2.02	0.40
42:q:37:THR:O	42:q:49:TYR:HA	2.21	0.40
7:G:321:GLY:HA2	7:G:496:ILE:HG23	2.04	0.40
7:G:611:LEU:HB3	7:G:612:PRO:HD2	2.04	0.40
12:L:67:HIS:NE2	12:L:70:THR:OG1	2.51	0.40
21:V:49:GLN:NE2	43:r:92:LYS:HE2	2.37	0.40
24:Y:48:PHE:O	24:Y:52:VAL:HG23	2.22	0.40
44:s:42:GLN:OE1	44:s:42:GLN:N	2.52	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

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	113/115 (98%)	108 (96%)	5 (4%)	0	100	100
2	B	153/216 (71%)	147 (96%)	6 (4%)	0	100	100
3	C	207/266 (78%)	202 (98%)	5 (2%)	0	100	100
4	D	427/463 (92%)	411 (96%)	16 (4%)	0	100	100
5	E	212/249 (85%)	200 (94%)	11 (5%)	1 (0%)	24	43
6	F	431/464 (93%)	414 (96%)	17 (4%)	0	100	100
7	G	698/727 (96%)	676 (97%)	22 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
8	H	317/318 (100%)	305 (96%)	12 (4%)	0	100	100
9	I	174/212 (82%)	170 (98%)	4 (2%)	0	100	100
10	J	173/175 (99%)	161 (93%)	12 (7%)	0	100	100
11	K	96/98 (98%)	94 (98%)	2 (2%)	0	100	100
12	L	604/606 (100%)	580 (96%)	23 (4%)	1 (0%)	43	64
13	M	457/459 (100%)	451 (99%)	6 (1%)	0	100	100
14	N	345/347 (99%)	338 (98%)	7 (2%)	0	100	100
15	O	318/343 (93%)	311 (98%)	7 (2%)	0	100	100
16	P	341/380 (90%)	330 (97%)	11 (3%)	0	100	100
17	Q	127/175 (73%)	124 (98%)	3 (2%)	0	100	100
18	R	94/124 (76%)	92 (98%)	2 (2%)	0	100	100
19	S	85/99 (86%)	82 (96%)	3 (4%)	0	100	100
20	T	83/156 (53%)	79 (95%)	4 (5%)	0	100	100
20	U	84/156 (54%)	82 (98%)	2 (2%)	0	100	100
21	V	113/116 (97%)	108 (96%)	5 (4%)	0	100	100
22	W	113/128 (88%)	110 (97%)	3 (3%)	0	100	100
23	X	169/172 (98%)	167 (99%)	2 (1%)	0	100	100
24	Y	138/141 (98%)	136 (99%)	2 (1%)	0	100	100
25	Z	139/144 (96%)	137 (99%)	2 (1%)	0	100	100
26	a	68/70 (97%)	66 (97%)	2 (3%)	0	100	100
27	b	81/84 (96%)	79 (98%)	2 (2%)	0	100	100
28	c	47/76 (62%)	46 (98%)	1 (2%)	0	100	100
29	d	118/120 (98%)	117 (99%)	1 (1%)	0	100	100
30	e	97/106 (92%)	95 (98%)	2 (2%)	0	100	100
31	f	55/57 (96%)	53 (96%)	2 (4%)	0	100	100
32	g	98/154 (64%)	92 (94%)	6 (6%)	0	100	100
33	h	136/189 (72%)	135 (99%)	1 (1%)	0	100	100
34	i	125/127 (98%)	118 (94%)	7 (6%)	0	100	100
35	j	65/108 (60%)	64 (98%)	1 (2%)	0	100	100
36	k	79/98 (81%)	75 (95%)	4 (5%)	0	100	100
37	l	154/186 (83%)	148 (96%)	6 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
38	m	126/129 (98%)	119 (94%)	7 (6%)	0	100	100
39	n	170/179 (95%)	164 (96%)	6 (4%)	0	100	100
40	o	120/137 (88%)	115 (96%)	5 (4%)	0	100	100
41	p	171/176 (97%)	168 (98%)	3 (2%)	0	100	100
42	q	143/145 (99%)	141 (99%)	1 (1%)	1 (1%)	18	34
43	r	91/113 (80%)	87 (96%)	4 (4%)	0	100	100
44	s	44/109 (40%)	41 (93%)	3 (7%)	0	100	100
All	All	8199/9212 (89%)	7938 (97%)	258 (3%)	3 (0%)	100	100

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
12	L	562	LEU
42	q	144	TYR
5	E	199	LEU

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

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	100/100 (100%)	100 (100%)	0	100	100
2	B	131/175 (75%)	131 (100%)	0	100	100
3	C	190/228 (83%)	190 (100%)	0	100	100
4	D	370/392 (94%)	370 (100%)	0	100	100
5	E	183/205 (89%)	183 (100%)	0	100	100
6	F	347/368 (94%)	347 (100%)	0	100	100
7	G	587/608 (96%)	586 (100%)	1 (0%)	87	95
8	H	275/274 (100%)	275 (100%)	0	100	100
9	I	151/175 (86%)	151 (100%)	0	100	100
10	J	141/141 (100%)	141 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
11	K	85/85 (100%)	85 (100%)	0	100	100
12	L	533/533 (100%)	532 (100%)	1 (0%)	87	95
13	M	412/412 (100%)	412 (100%)	0	100	100
14	N	315/315 (100%)	315 (100%)	0	100	100
15	O	283/303 (93%)	283 (100%)	0	100	100
16	P	297/327 (91%)	297 (100%)	0	100	100
17	Q	116/153 (76%)	116 (100%)	0	100	100
18	R	79/97 (81%)	79 (100%)	0	100	100
19	S	77/82 (94%)	76 (99%)	1 (1%)	61	76
20	T	79/135 (58%)	79 (100%)	0	100	100
20	U	79/135 (58%)	78 (99%)	1 (1%)	61	76
21	V	101/102 (99%)	101 (100%)	0	100	100
22	W	107/114 (94%)	107 (100%)	0	100	100
23	X	154/155 (99%)	154 (100%)	0	100	100
24	Y	101/102 (99%)	101 (100%)	0	100	100
25	Z	120/121 (99%)	120 (100%)	0	100	100
26	a	59/59 (100%)	59 (100%)	0	100	100
27	b	71/72 (99%)	71 (100%)	0	100	100
28	c	45/68 (66%)	45 (100%)	0	100	100
29	d	105/105 (100%)	105 (100%)	0	100	100
30	e	89/96 (93%)	89 (100%)	0	100	100
31	f	54/54 (100%)	54 (100%)	0	100	100
32	g	91/131 (70%)	91 (100%)	0	100	100
33	h	121/158 (77%)	121 (100%)	0	100	100
34	i	120/120 (100%)	120 (100%)	0	100	100
35	j	61/84 (73%)	61 (100%)	0	100	100
36	k	63/76 (83%)	63 (100%)	0	100	100
37	l	140/159 (88%)	140 (100%)	0	100	100
38	m	114/115 (99%)	114 (100%)	0	100	100
39	n	156/161 (97%)	156 (100%)	0	100	100
40	o	110/120 (92%)	110 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
41	p	155/157 (99%)	155 (100%)	0	100	100
42	q	131/131 (100%)	131 (100%)	0	100	100
43	r	85/97 (88%)	85 (100%)	0	100	100
44	s	45/92 (49%)	45 (100%)	0	100	100
All	All	7228/7892 (92%)	7224 (100%)	4 (0%)	87	95

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
7	G	614	ASP
12	L	442	ASN
19	S	75	ASN
20	U	44	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (48) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	162	GLN
4	D	55	HIS
4	D	59	HIS
4	D	280	GLN
5	E	42	HIS
5	E	91	ASN
5	E	99	HIS
6	F	257	ASN
6	F	398	GLN
7	G	155	GLN
8	H	235	ASN
8	H	287	HIS
12	L	56	HIS
12	L	170	GLN
12	L	175	ASN
12	L	269	ASN
12	L	296	ASN
12	L	354	GLN
12	L	400	ASN
12	L	541	ASN
13	M	81	GLN
13	M	82	HIS
13	M	88	ASN

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Mol	Chain	Res	Type
13	M	366	ASN
14	N	77	ASN
14	N	316	GLN
15	O	114	HIS
15	O	141	GLN
15	O	154	GLN
15	O	294	GLN
16	P	103	ASN
16	P	260	HIS
17	Q	9	GLN
20	T	74	GLN
23	X	30	HIS
23	X	34	GLN
23	X	76	HIS
23	X	162	HIS
24	Y	18	HIS
29	d	97	HIS
29	d	117	HIS
30	e	6	GLN
30	e	26	HIS
30	e	69	GLN
33	h	143	ASN
34	i	25	GLN
40	o	75	ASN
42	q	135	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

12 non-standard protein/DNA/RNA residues are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
10	FME	J	1	10	8,9,10	0.96	0	8,9,11	1.19	1 (12%)
4	2MR	D	85	4	10,12,13	0.57	0	5,13,15	1.07	0
8	FME	H	1	8	8,9,10	1.03	1 (12%)	8,9,11	1.01	0
12	FME	L	1	12	8,9,10	1.12	1 (12%)	8,9,11	0.95	0
13	FME	M	1	13	8,9,10	1.05	1 (12%)	8,9,11	0.90	0
24	AYA	Y	1	24	6,7,8	1.31	1 (16%)	6,8,10	1.38	1 (16%)
29	AME	d	1	29	9,10,11	1.75	1 (11%)	9,11,13	3.24	5 (55%)
34	SAC	i	1	34	7,8,9	1.12	1 (14%)	7,9,11	1.08	1 (14%)
1	FME	A	1	1	8,9,10	1.08	1 (12%)	8,9,11	0.80	0
43	AYA	r	1	43	6,7,8	1.21	1 (16%)	6,8,10	1.06	1 (16%)
11	FME	K	1	11	8,9,10	0.94	0	8,9,11	0.83	0
14	FME	N	1	14	8,9,10	1.06	1 (12%)	8,9,11	1.02	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
10	FME	J	1	10	-	3/7/9/11	-
4	2MR	D	85	4	-	2/10/13/15	-
8	FME	H	1	8	-	2/7/9/11	-
12	FME	L	1	12	-	2/7/9/11	-
13	FME	M	1	13	-	2/7/9/11	-
24	AYA	Y	1	24	-	0/5/6/8	-
29	AME	d	1	29	-	6/9/10/12	-
34	SAC	i	1	34	-	1/7/8/10	-
1	FME	A	1	1	-	4/7/9/11	-
43	AYA	r	1	43	-	3/5/6/8	-
11	FME	K	1	11	-	2/7/9/11	-
14	FME	N	1	14	-	3/7/9/11	-

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
29	d	1	AME	CT1-N	4.27	1.48	1.34
24	Y	1	AYA	CA-N	-2.79	1.43	1.46
12	L	1	FME	CA-N	-2.57	1.42	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1	FME	CA-N	-2.52	1.42	1.46
43	r	1	AYA	CA-N	-2.47	1.43	1.46
34	i	1	SAC	CA-N	-2.33	1.43	1.46
13	M	1	FME	CA-N	-2.26	1.43	1.46
8	H	1	FME	CA-N	-2.23	1.43	1.46
14	N	1	FME	CA-N	-2.13	1.43	1.46

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
29	d	1	AME	CT2-CT1-N	7.28	128.19	116.12
29	d	1	AME	OT-CT1-N	-3.71	115.42	121.98
29	d	1	AME	OT-CT1-CT2	-3.18	116.39	122.05
24	Y	1	AYA	CB-CA-N	3.07	113.14	109.68
10	J	1	FME	CB-CA-N	2.94	115.88	110.52
29	d	1	AME	CB-CA-N	2.52	115.11	110.52
29	d	1	AME	O-C-CA	-2.44	118.48	124.77
34	i	1	SAC	OG-CB-CA	-2.27	104.86	111.13
43	r	1	AYA	CA-N-CT	2.19	124.33	121.50

There are no chirality outliers.

All (30) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	1	FME	C-CA-CB-CG
1	A	1	FME	O-C-CA-CB
4	D	85	2MR	NE-CZ-NH2-CQ2
4	D	85	2MR	NH1-CZ-NH2-CQ2
10	J	1	FME	O1-CN-N-CA
10	J	1	FME	N-CA-CB-CG
10	J	1	FME	C-CA-CB-CG
11	K	1	FME	O1-CN-N-CA
12	L	1	FME	O1-CN-N-CA
13	M	1	FME	C-CA-CB-CG
14	N	1	FME	O1-CN-N-CA
14	N	1	FME	N-CA-CB-CG
14	N	1	FME	C-CA-CB-CG
29	d	1	AME	C-CA-CB-CG
29	d	1	AME	O-C-CA-CB
29	d	1	AME	CB-CA-N-CT1
34	i	1	SAC	O-C-CA-CB
43	r	1	AYA	O-C-CA-CB

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Mol	Chain	Res	Type	Atoms
8	H	1	FME	CA-CB-CG-SD
29	d	1	AME	CT2-CT1-N-CA
29	d	1	AME	OT-CT1-N-CA
1	A	1	FME	N-CA-CB-CG
29	d	1	AME	N-CA-CB-CG
8	H	1	FME	C-CA-CB-CG
12	L	1	FME	CA-CB-CG-SD
1	A	1	FME	CB-CA-N-CN
43	r	1	AYA	CB-CA-N-CT
13	M	1	FME	N-CA-CB-CG
11	K	1	FME	N-CA-CB-CG
43	r	1	AYA	C-CA-N-CT

There are no ring outliers.

3 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
10	J	1	FME	2	0
4	D	85	2MR	1	0
29	d	1	AME	7	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 56 ligands modelled in this entry, 3 are monoatomic - leaving 53 for Mogul analysis.

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	$\# Z > 2$
45	3PE	I	201	-	37,37,50	1.02	4 (10%)	40,42,55	1.10	3 (7%)
47	SF4	G	801	7	0,12,12	-	-	-		

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
45	3PE	d	201	-	36,36,50	1.02	3 (8%)	39,41,55	1.08	2 (5%)
45	3PE	M	601	-	44,44,50	0.92	3 (6%)	47,49,55	1.14	2 (4%)
45	3PE	M	603	-	50,50,50	0.87	4 (8%)	53,55,55	1.19	3 (5%)
46	PC1	h	202	-	43,43,53	1.06	4 (9%)	49,51,61	1.07	2 (4%)
45	3PE	P	403	-	34,34,50	1.06	4 (11%)	37,39,55	1.11	2 (5%)
45	3PE	L	703	-	44,44,50	0.92	3 (6%)	47,49,55	1.06	2 (4%)
46	PC1	A	202	-	34,34,53	1.21	3 (8%)	40,42,61	1.13	2 (5%)
46	PC1	m	201	-	39,39,53	1.11	4 (10%)	45,47,61	1.12	2 (4%)
47	SF4	I	203	9	0,12,12	-	-	-	-	-
51	CDL	L	702	-	59,59,99	1.14	6 (10%)	65,71,111	1.16	4 (6%)
45	3PE	Y	801	-	34,34,50	1.05	3 (8%)	37,39,55	1.23	3 (8%)
46	PC1	A	204	-	32,32,53	1.26	5 (15%)	38,40,61	1.15	2 (5%)
45	3PE	b	101	-	46,46,50	0.90	3 (6%)	49,51,55	1.03	2 (4%)
54	NDP	P	402	-	51,52,52	2.22	8 (15%)	71,80,80	1.44	12 (16%)
47	SF4	I	202	9	0,12,12	-	-	-	-	-
47	SF4	F	502	6	0,12,12	-	-	-	-	-
45	3PE	M	602	-	47,47,50	0.89	4 (8%)	50,52,55	1.14	2 (4%)
51	CDL	r	201	-	61,61,99	1.11	5 (8%)	67,73,111	1.10	4 (5%)
45	3PE	N	901	-	47,47,50	0.90	3 (6%)	50,52,55	1.05	2 (4%)
45	3PE	Y	803	-	30,30,50	1.07	3 (10%)	33,35,55	1.17	2 (6%)
51	CDL	H	602	-	49,49,99	1.23	8 (16%)	55,61,111	1.29	5 (9%)
46	PC1	M	604	-	38,38,53	1.14	3 (7%)	44,46,61	1.03	2 (4%)
58	MYR	o	201	40	13,14,15	0.71	0	12,13,15	0.50	0
45	3PE	H	601	-	28,28,50	1.07	2 (7%)	31,33,55	1.11	1 (3%)
56	EHZ	U	101	20	31,36,37	1.54	5 (16%)	36,44,47	1.41	3 (8%)
47	SF4	B	201	2	0,12,12	-	-	-	-	-
45	3PE	Z	202	-	34,34,50	1.05	4 (11%)	37,39,55	1.06	2 (5%)
49	FMN	F	501	-	33,33,33	1.12	2 (6%)	48,50,50	1.29	6 (12%)
51	CDL	d	203	-	64,64,99	1.08	8 (12%)	70,76,111	1.10	4 (5%)
45	3PE	Y	804	-	50,50,50	0.87	4 (8%)	53,55,55	1.00	2 (3%)
48	FES	G	803	7	0,4,4	-	-	-	-	-
45	3PE	M	605	-	49,49,50	0.86	4 (8%)	52,54,55	1.06	2 (3%)
51	CDL	K	101	-	68,68,99	1.06	7 (10%)	74,80,111	1.16	4 (5%)
52	GTP	O	401	53	33,34,34	3.22	15 (45%)	50,54,54	1.88	14 (28%)
46	PC1	Z	201	-	47,47,53	0.42	0	53,55,61	0.74	2 (3%)
45	3PE	L	701	-	38,38,50	1.02	3 (7%)	41,43,55	1.13	2 (4%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
46	PC1	P	401	-	53,53,53	1.05	3 (5%)	59,61,61	1.13	3 (5%)
45	3PE	Y	802	-	39,39,50	0.97	3 (7%)	42,44,55	1.12	2 (4%)
45	3PE	A	203	-	36,36,50	1.02	4 (11%)	39,41,55	1.10	3 (7%)
48	FES	E	301	5	0,4,4	-	-	-	-	-
46	PC1	d	202	-	32,32,53	1.27	6 (18%)	38,40,61	1.24	2 (5%)
51	CDL	X	201	-	72,72,99	1.03	7 (9%)	78,84,111	1.09	4 (5%)
51	CDL	N	902	-	59,59,99	1.12	7 (11%)	65,71,111	1.19	4 (6%)
51	CDL	h	201	-	77,77,99	0.99	8 (10%)	83,89,111	1.19	5 (6%)
56	EHZ	T	101	20	31,36,37	1.57	5 (16%)	36,44,47	1.82	8 (22%)
57	CHD	i	201	-	32,32,32	3.32	10 (31%)	51,51,51	3.86	22 (43%)
46	PC1	M	606	-	41,41,53	1.10	4 (9%)	47,49,61	1.17	3 (6%)
46	PC1	B	202	-	40,40,53	1.10	3 (7%)	46,48,61	1.14	2 (4%)
45	3PE	m	202	-	39,39,50	0.99	4 (10%)	42,44,55	1.17	2 (4%)
45	3PE	A	201	-	41,41,50	0.96	4 (9%)	44,46,55	1.08	2 (4%)
47	SF4	G	802	7	0,12,12	-	-	-	-	-

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
45	3PE	I	201	-	-	12/41/41/54	-
47	SF4	G	801	7	-	-	0/6/5/5
45	3PE	d	201	-	-	19/40/40/54	-
45	3PE	M	601	-	-	26/48/48/54	-
45	3PE	M	603	-	-	33/54/54/54	-
46	PC1	h	202	-	-	17/47/47/57	-
45	3PE	P	403	-	-	14/38/38/54	-
45	3PE	L	703	-	-	13/48/48/54	-
46	PC1	A	202	-	-	14/38/38/57	-
46	PC1	m	201	-	-	12/43/43/57	-
47	SF4	I	203	9	-	-	0/6/5/5
51	CDL	L	702	-	-	27/70/70/110	-
45	3PE	Y	801	-	-	14/38/38/54	-
46	PC1	A	204	-	-	16/36/36/57	-
45	3PE	b	101	-	-	22/50/50/54	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
54	NDP	P	402	-	-	6/34/77/77	0/5/5/5
47	SF4	I	202	9	-	-	0/6/5/5
47	SF4	F	502	6	-	-	0/6/5/5
45	3PE	M	602	-	-	25/51/51/54	-
51	CDL	r	201	-	-	22/72/72/110	-
45	3PE	N	901	-	-	19/51/51/54	-
45	3PE	Y	803	-	-	12/34/34/54	-
51	CDL	H	602	-	-	25/60/60/110	-
46	PC1	M	604	-	-	20/42/42/57	-
58	MYR	o	201	40	-	5/12/12/13	-
45	3PE	H	601	-	-	17/31/31/54	-
56	EHZ	U	101	20	-	7/42/44/45	-
47	SF4	B	201	2	-	-	0/6/5/5
45	3PE	Z	202	-	-	11/38/38/54	-
49	FMN	F	501	-	-	2/18/18/18	0/3/3/3
51	CDL	d	203	-	-	33/75/75/110	-
45	3PE	Y	804	-	-	20/54/54/54	-
48	FES	G	803	7	-	-	0/1/1/1
45	3PE	M	605	-	-	18/53/53/54	-
51	CDL	K	101	-	-	32/79/79/110	-
52	GTP	O	401	53	-	7/22/38/38	0/3/3/3
46	PC1	Z	201	-	-	18/51/51/57	-
45	3PE	L	701	-	-	18/42/42/54	-
46	PC1	P	401	-	-	19/57/57/57	-
45	3PE	Y	802	-	-	19/43/43/54	-
45	3PE	A	203	-	-	12/40/40/54	-
48	FES	E	301	5	-	-	0/1/1/1
46	PC1	d	202	-	-	18/36/36/57	-
51	CDL	X	201	-	-	41/83/83/110	-
51	CDL	N	902	-	-	25/69/69/110	-
51	CDL	h	201	-	-	38/88/88/110	-
56	EHZ	T	101	20	-	12/42/44/45	-
57	CHD	i	201	-	-	6/9/74/74	0/4/4/4
46	PC1	M	606	-	-	17/45/45/57	-
46	PC1	B	202	-	-	11/44/44/57	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
45	3PE	m	202	-	-	14/43/43/54	-
45	3PE	A	201	-	-	14/45/45/54	-
47	SF4	G	802	7	-	-	0/6/5/5

All (205) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
54	P	402	NDP	P2B-O2B	12.53	1.81	1.59
57	i	201	CHD	C11-C12	9.48	1.68	1.53
52	O	401	GTP	O6-C6	8.54	1.39	1.23
57	i	201	CHD	C16-C15	7.18	1.73	1.54
52	O	401	GTP	C5-N7	6.60	1.52	1.39
57	i	201	CHD	C13-C17	5.95	1.65	1.55
57	i	201	CHD	C6-C5	5.51	1.62	1.53
57	i	201	CHD	C8-C9	5.45	1.64	1.53
57	i	201	CHD	C20-C17	-5.33	1.45	1.54
57	i	201	CHD	O12-C12	-5.30	1.34	1.43
56	T	101	EHZ	C12-N1	4.88	1.44	1.33
52	O	401	GTP	C2-N1	4.83	1.49	1.37
56	U	101	EHZ	C15-N2	4.81	1.44	1.33
56	U	101	EHZ	C12-N1	4.80	1.44	1.33
56	T	101	EHZ	C15-N2	4.73	1.44	1.33
52	O	401	GTP	PA-O3A	4.64	1.64	1.59
52	O	401	GTP	C2-N3	4.61	1.44	1.33
52	O	401	GTP	C2-N2	4.59	1.44	1.34
52	O	401	GTP	PB-O3A	4.52	1.64	1.59
52	O	401	GTP	C8-N9	-4.51	1.27	1.37
52	O	401	GTP	PB-O3B	4.44	1.64	1.59
57	i	201	CHD	C15-C14	4.12	1.62	1.54
52	O	401	GTP	C4-N9	-3.86	1.28	1.38
46	P	401	PC1	O21-C2	-3.69	1.37	1.46
54	P	402	NDP	O2B-C2B	-3.66	1.31	1.44
54	P	402	NDP	PN-O5D	3.48	1.73	1.59
57	i	201	CHD	C6-C7	3.38	1.59	1.52
54	P	402	NDP	O4B-C4B	-3.28	1.37	1.45
51	L	702	CDL	OA6-CA4	-3.20	1.39	1.46
51	r	201	CDL	OA6-CA4	-3.19	1.39	1.46
49	F	501	FMN	C4A-N5	3.13	1.37	1.30
45	L	703	3PE	O21-C2	-3.05	1.39	1.46
45	b	101	3PE	O21-C2	-3.04	1.39	1.46
51	r	201	CDL	OB6-CB4	-2.98	1.39	1.46
46	P	401	PC1	O31-C3	-2.97	1.38	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
46	A	202	PC1	O21-C2	-2.96	1.39	1.46
45	L	701	3PE	O21-C2	-2.96	1.39	1.46
51	K	101	CDL	OA6-CA4	-2.93	1.39	1.46
51	H	602	CDL	OB6-CB4	-2.93	1.39	1.46
45	Y	802	3PE	O21-C2	-2.91	1.39	1.46
46	M	604	PC1	O21-C2	-2.90	1.39	1.46
46	h	202	PC1	O21-C2	-2.90	1.39	1.46
51	X	201	CDL	OA6-CA4	-2.87	1.39	1.46
51	d	203	CDL	OA6-CA4	-2.87	1.39	1.46
45	d	201	3PE	O21-C2	-2.86	1.39	1.46
45	H	601	3PE	O21-C2	-2.86	1.39	1.46
46	M	606	PC1	O21-C2	-2.85	1.39	1.46
51	H	602	CDL	OA6-CA4	-2.84	1.39	1.46
51	h	201	CDL	OA6-CA4	-2.84	1.39	1.46
45	A	201	3PE	O21-C2	-2.82	1.40	1.46
51	N	902	CDL	OA6-CA4	-2.82	1.40	1.46
46	d	202	PC1	O21-C2	-2.81	1.40	1.46
56	T	101	EHZ	O4-C15	-2.81	1.18	1.23
45	M	601	3PE	O21-C2	-2.79	1.40	1.46
51	d	203	CDL	OB6-CB4	-2.79	1.40	1.46
45	Z	202	3PE	O21-C2	-2.79	1.40	1.46
45	Y	801	3PE	O21-C2	-2.79	1.40	1.46
51	L	702	CDL	OB6-CB4	-2.79	1.40	1.46
52	O	401	GTP	O4'-C1'	2.77	1.48	1.42
51	K	101	CDL	OB6-CB4	-2.76	1.40	1.46
46	B	202	PC1	O21-C2	-2.74	1.40	1.46
46	A	204	PC1	O21-C2	-2.73	1.40	1.46
45	L	701	3PE	O31-C3	-2.72	1.39	1.45
45	M	603	3PE	O21-C2	-2.71	1.40	1.46
45	M	602	3PE	O21-C2	-2.71	1.40	1.46
45	I	201	3PE	O21-C2	-2.70	1.40	1.46
45	m	202	3PE	O21-C2	-2.69	1.40	1.46
52	O	401	GTP	C2'-C3'	-2.68	1.46	1.53
45	P	403	3PE	O21-C2	-2.68	1.40	1.46
56	U	101	EHZ	O4-C15	-2.67	1.18	1.23
51	N	902	CDL	OB6-CB4	-2.67	1.40	1.46
45	M	601	3PE	O31-C3	-2.65	1.39	1.45
51	K	101	CDL	OA8-CA6	-2.64	1.39	1.45
56	T	101	EHZ	O3-C12	-2.64	1.18	1.23
45	N	901	3PE	O21-C2	-2.63	1.40	1.46
51	L	702	CDL	OA8-CA7	2.61	1.41	1.33
51	L	702	CDL	OB8-CB7	2.59	1.40	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
46	A	202	PC1	O31-C3	-2.59	1.39	1.45
45	Y	803	3PE	O21-C2	-2.58	1.40	1.46
45	A	203	3PE	O21-C2	-2.54	1.40	1.46
51	r	201	CDL	OB8-CB7	2.54	1.40	1.33
45	I	201	3PE	O31-C3	-2.54	1.39	1.45
51	X	201	CDL	OB6-CB4	-2.54	1.40	1.46
51	N	902	CDL	OA8-CA7	2.53	1.40	1.33
51	h	201	CDL	OB6-CB4	-2.52	1.40	1.46
46	m	201	PC1	O31-C3	-2.52	1.39	1.45
45	d	201	3PE	O31-C31	2.51	1.40	1.33
46	m	201	PC1	O21-C2	-2.50	1.40	1.46
51	h	201	CDL	OA8-CA7	2.49	1.40	1.33
45	Y	802	3PE	O31-C31	2.49	1.40	1.33
45	M	603	3PE	O31-C3	-2.49	1.39	1.45
45	M	602	3PE	O31-C3	-2.48	1.39	1.45
46	M	604	PC1	O31-C3	-2.47	1.39	1.45
51	X	201	CDL	OA8-CA7	2.46	1.40	1.33
52	O	401	GTP	PG-O3G	-2.46	1.45	1.54
45	M	605	3PE	O21-C2	-2.45	1.40	1.46
46	M	606	PC1	O31-C3	-2.45	1.39	1.45
45	N	901	3PE	O31-C31	2.44	1.40	1.33
45	m	202	3PE	O31-C31	2.44	1.40	1.33
45	Y	801	3PE	O31-C3	-2.44	1.39	1.45
56	U	101	EHZ	O3-C12	-2.44	1.18	1.23
51	X	201	CDL	OB8-CB7	2.44	1.40	1.33
46	d	202	PC1	O31-C31	2.43	1.40	1.33
51	N	902	CDL	OB8-CB7	2.43	1.40	1.33
45	Y	804	3PE	O31-C31	2.43	1.40	1.33
51	d	203	CDL	OA8-CA7	2.42	1.40	1.33
46	M	604	PC1	O31-C31	2.42	1.40	1.33
46	h	202	PC1	O31-C3	-2.41	1.39	1.45
45	L	703	3PE	O31-C3	-2.40	1.39	1.45
45	P	403	3PE	O31-C31	2.40	1.40	1.33
51	r	201	CDL	OA8-CA6	-2.39	1.39	1.45
51	H	602	CDL	OB8-CB6	-2.38	1.39	1.45
46	A	204	PC1	O31-C3	-2.38	1.39	1.45
52	O	401	GTP	PG-O2G	-2.37	1.46	1.54
54	P	402	NDP	PA-O3	2.36	1.62	1.59
45	P	403	3PE	O31-C3	-2.35	1.39	1.45
46	B	202	PC1	O31-C31	2.34	1.40	1.33
45	N	901	3PE	O31-C3	-2.33	1.40	1.45
51	H	602	CDL	OA8-CA7	2.33	1.40	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
45	A	203	3PE	O31-C31	2.33	1.40	1.33
51	K	101	CDL	OB8-CB7	2.32	1.40	1.33
46	P	401	PC1	C12-N	-2.32	1.44	1.51
54	P	402	NDP	O3B-C3B	-2.32	1.37	1.43
45	A	201	3PE	O31-C31	2.32	1.40	1.33
45	Y	804	3PE	O21-C21	2.31	1.40	1.34
51	d	203	CDL	OB8-CB6	-2.31	1.40	1.45
45	b	101	3PE	O31-C31	2.31	1.40	1.33
51	h	201	CDL	OB8-CB6	-2.31	1.40	1.45
45	Z	202	3PE	O31-C31	2.31	1.40	1.33
51	K	101	CDL	OB6-CB5	2.30	1.40	1.34
45	M	605	3PE	O31-C31	2.30	1.40	1.33
46	A	204	PC1	O31-C31	2.29	1.40	1.33
45	b	101	3PE	O31-C3	-2.29	1.40	1.45
51	X	201	CDL	OB6-CB5	2.29	1.40	1.34
51	d	203	CDL	OB8-CB7	2.28	1.40	1.33
45	A	201	3PE	O31-C3	-2.28	1.40	1.45
46	m	201	PC1	O21-C21	2.27	1.40	1.34
45	Y	803	3PE	O31-C31	2.25	1.39	1.33
45	Y	804	3PE	O21-C2	-2.24	1.41	1.46
45	Z	202	3PE	O21-C21	2.24	1.40	1.34
45	Y	801	3PE	O31-C31	2.24	1.39	1.33
54	P	402	NDP	O2D-C2D	-2.24	1.37	1.43
45	P	403	3PE	O21-C21	2.23	1.40	1.34
57	i	201	CHD	O7-C7	-2.21	1.38	1.43
45	M	605	3PE	O31-C3	-2.20	1.40	1.45
45	I	201	3PE	O21-C21	2.20	1.40	1.34
45	M	605	3PE	O21-C21	2.20	1.40	1.34
46	A	204	PC1	O21-C21	2.20	1.40	1.34
45	Y	803	3PE	O31-C3	-2.20	1.40	1.45
51	h	201	CDL	OA6-CA5	2.19	1.40	1.34
51	d	203	CDL	OA8-CA6	-2.19	1.40	1.45
51	h	201	CDL	OB8-CB7	2.19	1.39	1.33
45	A	203	3PE	O21-C21	2.19	1.40	1.34
46	B	202	PC1	O31-C3	-2.18	1.40	1.45
45	M	601	3PE	O31-C31	2.17	1.39	1.33
51	H	602	CDL	OA6-CA5	2.17	1.40	1.34
45	H	601	3PE	O31-C3	-2.17	1.40	1.45
51	h	201	CDL	OA8-CA6	-2.17	1.40	1.45
45	A	203	3PE	O31-C3	-2.17	1.40	1.45
46	h	202	PC1	O31-C31	2.16	1.39	1.33
51	L	702	CDL	OB8-CB6	-2.16	1.40	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
46	M	606	PC1	O31-C31	2.14	1.39	1.33
51	N	902	CDL	OB8-CB6	-2.14	1.40	1.45
51	K	101	CDL	OB8-CB6	-2.13	1.40	1.45
45	m	202	3PE	O21-C21	2.13	1.40	1.34
45	M	603	3PE	O21-C21	2.12	1.40	1.34
51	X	201	CDL	OB8-CB6	-2.12	1.40	1.45
51	r	201	CDL	OA8-CA7	2.12	1.39	1.33
46	M	606	PC1	O21-C21	2.11	1.40	1.34
46	A	202	PC1	O31-C31	2.11	1.39	1.33
51	H	602	CDL	OA8-CA6	-2.11	1.40	1.45
45	d	201	3PE	O31-C3	-2.11	1.40	1.45
46	m	201	PC1	O31-C31	2.10	1.39	1.33
45	L	703	3PE	O31-C31	2.10	1.39	1.33
45	m	202	3PE	O31-C3	-2.10	1.40	1.45
51	N	902	CDL	OA6-CA5	2.09	1.40	1.34
45	M	602	3PE	O31-C31	2.09	1.39	1.33
51	N	902	CDL	OB6-CB5	2.09	1.39	1.35
45	Z	202	3PE	O31-C3	-2.09	1.40	1.45
56	T	101	EHZ	O2-C9	-2.09	1.17	1.21
46	d	202	PC1	O31-C3	-2.08	1.40	1.45
49	F	501	FMN	C4A-C10	-2.08	1.38	1.44
51	h	201	CDL	OB6-CB5	2.08	1.40	1.34
46	A	204	PC1	C12-N	-2.08	1.45	1.51
45	Y	804	3PE	O31-C3	-2.07	1.40	1.45
51	d	203	CDL	OB6-CB5	2.07	1.40	1.34
51	K	101	CDL	OA6-CA5	2.06	1.40	1.34
51	H	602	CDL	OB8-CB7	2.06	1.39	1.33
51	d	203	CDL	OA6-CA5	2.06	1.40	1.34
45	M	602	3PE	O21-C21	2.05	1.40	1.34
46	h	202	PC1	O21-C21	2.05	1.40	1.34
56	U	101	EHZ	C9-S1	2.05	1.81	1.76
46	d	202	PC1	O21-C21	2.04	1.40	1.34
51	L	702	CDL	OA8-CA6	-2.04	1.40	1.45
51	H	602	CDL	OB6-CB5	2.02	1.40	1.34
54	P	402	NDP	O5B-C5B	-2.02	1.36	1.44
46	d	202	PC1	C12-N	-2.02	1.45	1.51
46	d	202	PC1	C15-N	-2.02	1.44	1.50
45	L	701	3PE	O31-C31	2.01	1.39	1.33
45	A	201	3PE	O21-C21	2.01	1.39	1.34
45	M	603	3PE	O31-C31	2.01	1.39	1.33
52	O	401	GTP	C1'-N9	-2.01	1.41	1.47
51	X	201	CDL	OA8-CA6	-2.00	1.40	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
45	Y	802	3PE	O31-C3	-2.00	1.40	1.45
45	I	201	3PE	O31-C31	2.00	1.39	1.33

All (164) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
57	i	201	CHD	C11-C9-C10	11.22	125.09	113.70
57	i	201	CHD	C4-C5-C10	9.20	122.45	112.66
57	i	201	CHD	C18-C13-C12	-8.48	100.57	109.06
57	i	201	CHD	C14-C13-C12	8.43	115.12	107.42
57	i	201	CHD	C17-C13-C12	8.30	125.13	117.67
57	i	201	CHD	C14-C8-C7	8.24	122.79	111.85
52	O	401	GTP	C8-N9-C4	7.53	120.14	106.03
57	i	201	CHD	C6-C5-C10	7.18	120.30	112.66
56	T	101	EHZ	C8-C9-S1	6.66	121.97	113.56
56	U	101	EHZ	C8-C9-S1	5.22	120.15	113.56
57	i	201	CHD	C9-C8-C7	5.06	118.22	111.86
51	N	902	CDL	OB6-CB5-C51	4.80	119.65	111.09
51	H	602	CDL	OB6-CB5-C51	4.76	121.78	111.48
57	i	201	CHD	C11-C12-C13	4.71	116.05	111.26
51	h	201	CDL	OA6-CA5-C11	4.61	121.46	111.48
46	d	202	PC1	O21-C21-C22	4.57	121.37	111.48
57	i	201	CHD	C1-C10-C5	4.57	114.30	107.75
57	i	201	CHD	C15-C14-C8	4.49	124.52	118.36
45	M	603	3PE	O21-C21-C22	4.45	121.11	111.48
45	L	701	3PE	O21-C21-C22	4.44	121.09	111.48
51	h	201	CDL	OB6-CB5-C51	4.42	121.04	111.48
45	Y	801	3PE	O21-C21-C22	4.37	120.94	111.48
57	i	201	CHD	C10-C9-C8	4.33	116.67	111.84
45	m	202	3PE	O21-C21-C22	4.27	120.72	111.48
51	X	201	CDL	OB6-CB5-C51	4.26	120.69	111.48
45	M	602	3PE	O21-C21-C22	4.19	120.55	111.48
45	M	605	3PE	O21-C21-C22	4.19	120.54	111.48
46	B	202	PC1	O21-C21-C22	4.16	120.49	111.48
51	L	702	CDL	OB6-CB5-C51	4.15	120.45	111.48
45	H	601	3PE	O21-C21-C22	4.15	120.45	111.48
45	A	201	3PE	O21-C21-C22	4.11	120.38	111.48
46	M	606	PC1	O21-C21-C22	4.08	120.30	111.48
45	Y	803	3PE	O21-C21-C22	4.05	120.23	111.48
45	N	901	3PE	O21-C21-C22	4.04	120.23	111.48
46	A	202	PC1	O21-C21-C22	4.04	120.22	111.48
46	A	204	PC1	O21-C21-C22	4.04	120.21	111.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
51	d	203	CDL	OA6-CA5-C11	3.97	120.07	111.48
46	Z	201	PC1	O21-C21-C22	3.96	120.05	111.48
57	i	201	CHD	C13-C17-C20	-3.96	114.69	119.48
45	Y	802	3PE	O21-C21-C22	3.96	120.04	111.48
54	P	402	NDP	O2B-P2B-O1X	-3.90	95.44	109.33
45	Y	804	3PE	O21-C21-C22	3.83	119.77	111.48
46	m	201	PC1	O21-C21-C22	3.82	119.75	111.48
46	h	202	PC1	O21-C21-C22	3.81	119.72	111.48
45	P	403	3PE	O21-C21-C22	3.77	119.63	111.48
51	K	101	CDL	OB6-CB5-C51	3.75	119.59	111.48
57	i	201	CHD	C14-C8-C9	3.71	114.92	109.75
51	K	101	CDL	OA6-CA5-C11	3.70	119.49	111.48
45	d	201	3PE	O21-C21-C22	3.67	119.43	111.48
45	M	601	3PE	O21-C21-C22	3.66	119.39	111.48
45	I	201	3PE	O21-C21-C22	3.65	119.38	111.48
45	Z	202	3PE	O21-C21-C22	3.58	119.24	111.48
51	H	602	CDL	OA6-CA5-C11	3.55	119.17	111.48
45	L	703	3PE	O21-C21-C22	3.54	119.14	111.48
49	F	501	FMN	C4-N3-C2	-3.52	119.39	125.64
51	L	702	CDL	OA6-CA5-C11	3.49	119.03	111.48
46	P	401	PC1	O21-C21-C22	3.47	118.98	111.48
45	A	203	3PE	O21-C21-C22	3.46	118.97	111.48
45	m	202	3PE	O31-C31-C32	3.39	122.17	111.83
49	F	501	FMN	C4A-C10-N10	3.38	121.32	116.48
51	r	201	CDL	OA6-CA5-C11	3.35	118.73	111.48
52	O	401	GTP	C2-N1-C6	-3.34	119.05	125.11
51	N	902	CDL	OA6-CA5-C11	3.30	118.62	111.48
52	O	401	GTP	O2G-PG-O3B	3.29	115.66	104.64
51	r	201	CDL	OB6-CB5-C51	3.29	118.59	111.48
57	i	201	CHD	C21-C20-C17	-3.28	107.97	112.88
56	T	101	EHZ	C13-C12-N1	3.27	122.30	116.34
54	P	402	NDP	O3-PA-O1A	-3.25	100.94	110.70
51	X	201	CDL	OA6-CA5-C11	3.24	118.49	111.48
45	b	101	3PE	O21-C21-C22	3.24	118.49	111.48
56	T	101	EHZ	O2-C9-S1	-3.19	118.63	122.68
45	Y	801	3PE	O31-C31-C32	3.19	121.55	111.83
46	M	604	PC1	O31-C31-C32	3.17	120.79	111.15
46	M	604	PC1	O21-C21-C22	3.16	118.32	111.48
45	M	603	3PE	O31-C31-C32	3.12	121.36	111.83
51	N	902	CDL	OA8-CA7-C31	3.11	121.31	111.83
46	d	202	PC1	O31-C31-C32	3.11	121.31	111.83
46	B	202	PC1	O31-C31-C32	3.11	121.30	111.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
54	P	402	NDP	P2B-O2B-C2B	-3.06	115.25	123.43
45	d	201	3PE	O31-C31-C32	3.05	121.14	111.83
46	m	201	PC1	O31-C31-C32	3.04	121.10	111.83
51	r	201	CDL	OB8-CB7-C71	3.03	120.35	111.15
57	i	201	CHD	C11-C9-C8	3.01	115.35	110.89
46	M	606	PC1	O31-C31-C32	3.01	121.02	111.83
45	A	203	3PE	O31-C31-C32	2.98	120.93	111.83
45	M	601	3PE	O31-C31-C32	2.94	120.80	111.83
51	H	602	CDL	OA8-CA7-C31	2.90	120.68	111.83
51	K	101	CDL	OB8-CB7-C71	2.89	120.66	111.83
56	T	101	EHZ	C10-S1-C9	2.89	110.39	101.84
51	H	602	CDL	OB8-CB7-C71	2.89	120.65	111.83
51	L	702	CDL	OA8-CA7-C31	2.88	120.62	111.83
51	N	902	CDL	OB8-CB7-C71	2.87	120.58	111.83
46	A	204	PC1	O31-C31-C32	2.85	120.53	111.83
54	P	402	NDP	PA-O5B-C5B	-2.84	105.06	121.35
51	X	201	CDL	OA8-CA7-C31	2.81	120.39	111.83
54	P	402	NDP	O2N-PN-O3	2.80	114.84	107.27
45	Y	804	3PE	O31-C31-C32	2.79	120.33	111.83
45	I	201	3PE	O31-C31-C32	2.78	120.31	111.83
45	Z	202	3PE	O31-C31-C32	2.77	120.29	111.83
56	T	101	EHZ	O2-C9-C8	-2.76	119.41	123.74
57	i	201	CHD	C21-C20-C22	-2.75	106.08	110.34
46	h	202	PC1	O31-C31-C32	2.71	120.10	111.83
52	O	401	GTP	C1'-N9-C8	-2.68	119.11	126.73
46	A	202	PC1	O31-C31-C32	2.66	119.95	111.83
51	X	201	CDL	OB8-CB7-C71	2.66	119.94	111.83
45	Y	802	3PE	O31-C31-C32	2.66	119.93	111.83
57	i	201	CHD	C9-C11-C12	-2.65	110.82	114.29
54	P	402	NDP	C2A-N1A-C6A	-2.65	114.38	118.73
45	M	602	3PE	O31-C31-C32	2.64	119.89	111.83
49	F	501	FMN	C4A-C4-N3	2.63	119.94	113.25
45	Y	803	3PE	O31-C31-C32	2.60	119.77	111.83
51	K	101	CDL	OA8-CA7-C31	2.60	119.75	111.83
52	O	401	GTP	O3G-PG-O3B	2.59	113.33	104.64
54	P	402	NDP	O3X-P2B-O2X	2.59	117.51	107.80
52	O	401	GTP	C2'-C3'-C4'	2.58	107.59	102.61
51	d	203	CDL	OA8-CA7-C31	2.57	119.67	111.83
57	i	201	CHD	C18-C13-C14	-2.55	107.22	111.20
51	r	201	CDL	OA8-CA7-C31	2.54	119.59	111.83
56	U	101	EHZ	O2-C9-C8	-2.54	119.74	123.74
45	b	101	3PE	O31-C31-C32	2.53	119.56	111.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
51	h	201	CDL	OB8-CB7-C71	2.51	119.50	111.83
45	P	403	3PE	O31-C31-C32	2.51	119.49	111.83
52	O	401	GTP	C5-C6-N1	2.51	119.64	113.25
51	d	203	CDL	OB8-CB7-C71	2.50	119.45	111.83
51	d	203	CDL	OB6-CB5-C51	2.48	120.05	110.93
56	T	101	EHZ	C13-C14-N2	-2.48	106.72	112.00
51	L	702	CDL	OB8-CB7-C71	2.47	119.37	111.83
49	F	501	FMN	C10-C4A-N5	-2.46	119.78	124.81
52	O	401	GTP	O6-C6-C5	-2.45	120.06	126.53
46	P	401	PC1	O31-C31-C32	2.39	119.13	111.83
46	P	401	PC1	C24-C23-C22	-2.39	104.34	113.13
49	F	501	FMN	O4-C4-C4A	-2.38	120.25	126.53
51	h	201	CDL	OA8-CA7-C31	2.38	119.08	111.83
52	O	401	GTP	C5-C4-N9	-2.36	101.43	105.66
45	M	605	3PE	O31-C31-C32	2.36	119.02	111.83
51	H	602	CDL	OB6-CB5-OB7	-2.34	118.23	123.70
57	i	201	CHD	C19-C10-C5	-2.32	106.55	110.44
56	T	101	EHZ	C11-N1-C12	-2.32	118.51	122.82
56	T	101	EHZ	O3-C12-N1	-2.31	118.49	123.03
52	O	401	GTP	O2A-PA-O1A	-2.31	101.72	112.44
45	A	201	3PE	O31-C31-C32	2.29	118.81	111.83
45	N	901	3PE	O31-C31-C32	2.28	118.78	111.83
57	i	201	CHD	C6-C5-C4	2.26	113.81	111.23
45	I	201	3PE	C3-C2-C1	-2.26	106.52	111.78
45	L	703	3PE	O31-C31-C32	2.26	118.71	111.83
52	O	401	GTP	O2B-PB-O1B	-2.23	102.07	112.44
46	Z	201	PC1	O21-C21-O22	-2.21	118.55	123.70
46	M	606	PC1	C2-O21-C21	-2.20	112.52	117.80
45	Y	801	3PE	O21-C21-O22	-2.20	118.56	123.70
54	P	402	NDP	O2N-PN-O1N	2.18	122.59	112.44
54	P	402	NDP	O7N-C7N-N7N	-2.16	118.06	122.89
45	L	701	3PE	O31-C31-C32	2.13	118.34	111.83
49	F	501	FMN	C4A-C10-N1	-2.13	119.37	124.59
45	A	203	3PE	C24-C23-C22	-2.13	105.31	113.13
45	M	603	3PE	O31-C31-O32	-2.09	118.40	123.63
52	O	401	GTP	C3'-C2'-C1'	2.08	105.39	101.46
52	O	401	GTP	N9-C8-N7	-2.08	109.55	113.40
54	P	402	NDP	C5D-C4D-C3D	-2.05	107.83	115.21
57	i	201	CHD	O26-C24-C23	2.05	120.46	114.00
52	O	401	GTP	C8-N7-C5	-2.04	100.63	104.26
56	U	101	EHZ	O2-C9-S1	-2.03	120.11	122.68
51	h	201	CDL	CB6-CB4-CB3	-2.02	107.07	111.78

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
54	P	402	NDP	O3X-P2B-O2B	-2.02	97.98	105.85
54	P	402	NDP	C5B-C4B-C3B	-2.01	107.97	115.21

There are no chirality outliers.

All (802) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
45	A	201	3PE	C11-O13-P-O11
45	A	201	3PE	C11-O13-P-O14
45	A	201	3PE	O13-C11-C12-N
45	H	601	3PE	C1-O11-P-O12
45	H	601	3PE	C1-O11-P-O13
45	H	601	3PE	C1-O11-P-O14
45	H	601	3PE	O13-C11-C12-N
45	I	201	3PE	C1-O11-P-O14
45	L	701	3PE	C11-O13-P-O11
45	L	701	3PE	C11-O13-P-O14
45	L	701	3PE	O13-C11-C12-N
45	L	701	3PE	C22-C21-O21-C2
45	L	703	3PE	O13-C11-C12-N
45	M	601	3PE	C1-O11-P-O12
45	M	601	3PE	C1-O11-P-O13
45	M	601	3PE	C1-O11-P-O14
45	M	601	3PE	C11-O13-P-O14
45	M	601	3PE	C12-C11-O13-P
45	M	601	3PE	O13-C11-C12-N
45	M	602	3PE	O13-C11-C12-N
45	M	602	3PE	C22-C21-O21-C2
45	M	603	3PE	C1-O11-P-O12
45	M	603	3PE	C1-O11-P-O13
45	M	603	3PE	C1-O11-P-O14
45	M	603	3PE	O13-C11-C12-N
45	M	605	3PE	O13-C11-C12-N
45	N	901	3PE	C1-O11-P-O12
45	N	901	3PE	C1-O11-P-O13
45	N	901	3PE	C1-O11-P-O14
45	N	901	3PE	C11-O13-P-O14
45	N	901	3PE	O13-C11-C12-N
45	N	901	3PE	O11-C1-C2-O21
45	P	403	3PE	C1-O11-P-O12
45	P	403	3PE	C1-O11-P-O13
45	P	403	3PE	C1-O11-P-O14

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Mol	Chain	Res	Type	Atoms
45	P	403	3PE	C11-O13-P-O11
45	P	403	3PE	C11-O13-P-O14
45	Y	801	3PE	C22-C21-O21-C2
45	Y	802	3PE	C1-O11-P-O12
45	Y	802	3PE	C1-O11-P-O13
45	Y	802	3PE	C1-O11-P-O14
45	Y	802	3PE	C11-O13-P-O11
45	Y	802	3PE	C11-O13-P-O12
45	Y	802	3PE	C2-C1-O11-P
45	Y	802	3PE	C12-C11-O13-P
45	Y	803	3PE	O13-C11-C12-N
45	Y	804	3PE	C11-O13-P-O14
45	b	101	3PE	C1-O11-P-O14
45	b	101	3PE	O13-C11-C12-N
45	d	201	3PE	C1-O11-P-O13
45	d	201	3PE	C1-O11-P-O14
45	d	201	3PE	C11-O13-P-O11
45	d	201	3PE	C11-O13-P-O14
45	m	202	3PE	C22-C21-O21-C2
46	A	202	PC1	C1-O11-P-O14
46	A	204	PC1	O13-C11-C12-N
46	A	204	PC1	O22-C21-O21-C2
46	B	202	PC1	C22-C21-O21-C2
46	M	604	PC1	C11-O13-P-O14
46	M	604	PC1	C11-O13-P-O11
46	M	604	PC1	C1-O11-P-O12
46	M	604	PC1	C1-O11-P-O14
46	M	604	PC1	C1-O11-P-O13
46	M	606	PC1	C11-O13-P-O14
46	M	606	PC1	C22-C21-O21-C2
46	P	401	PC1	C1-O11-P-O12
46	P	401	PC1	C1-O11-P-O13
46	Z	201	PC1	C11-O13-P-O12
46	Z	201	PC1	C1-O11-P-O14
46	Z	201	PC1	C1-O11-P-O13
46	Z	201	PC1	O11-C1-C2-O21
46	d	202	PC1	C1-O11-P-O12
46	d	202	PC1	C1-O11-P-O14
46	d	202	PC1	C1-O11-P-O13
46	d	202	PC1	O32-C31-O31-C3
46	h	202	PC1	C11-O13-P-O12
46	h	202	PC1	C11-O13-P-O11

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Mol	Chain	Res	Type	Atoms
46	h	202	PC1	C1-O11-P-O14
46	h	202	PC1	C1-O11-P-O13
46	m	201	PC1	C11-O13-P-O12
46	m	201	PC1	C11-O13-P-O11
46	m	201	PC1	C1-O11-P-O12
46	m	201	PC1	C1-O11-P-O13
46	m	201	PC1	O13-C11-C12-N
51	H	602	CDL	CB3-OB5-PB2-OB4
51	H	602	CDL	OB7-CB5-OB6-CB4
51	K	101	CDL	CA2-OA2-PA1-OA4
51	K	101	CDL	CA2-OA2-PA1-OA5
51	K	101	CDL	CB2-OB2-PB2-OB3
51	K	101	CDL	CB2-OB2-PB2-OB5
51	L	702	CDL	CB2-C1-CA2-OA2
51	L	702	CDL	CB2-OB2-PB2-OB3
51	L	702	CDL	CB2-OB2-PB2-OB5
51	N	902	CDL	CB2-OB2-PB2-OB3
51	N	902	CDL	CB2-OB2-PB2-OB5
51	X	201	CDL	CA3-OA5-PA1-OA2
51	X	201	CDL	CA3-OA5-PA1-OA3
51	X	201	CDL	CB2-OB2-PB2-OB3
51	X	201	CDL	CB2-OB2-PB2-OB4
51	X	201	CDL	CB2-OB2-PB2-OB5
51	X	201	CDL	CB3-OB5-PB2-OB2
51	X	201	CDL	C51-CB5-OB6-CB4
51	d	203	CDL	CA3-OA5-PA1-OA3
51	d	203	CDL	CB2-OB2-PB2-OB4
51	d	203	CDL	CB2-OB2-PB2-OB5
51	d	203	CDL	CB3-OB5-PB2-OB2
51	d	203	CDL	CB3-OB5-PB2-OB3
51	d	203	CDL	CB3-OB5-PB2-OB4
51	d	203	CDL	C51-CB5-OB6-CB4
51	h	201	CDL	CB2-C1-CA2-OA2
51	h	201	CDL	CA2-OA2-PA1-OA3
51	h	201	CDL	CA2-OA2-PA1-OA4
51	h	201	CDL	CA2-OA2-PA1-OA5
51	h	201	CDL	OA7-CA5-OA6-CA4
51	h	201	CDL	C11-CA5-OA6-CA4
51	h	201	CDL	CB3-OB5-PB2-OB2
51	h	201	CDL	CB3-OB5-PB2-OB3
51	r	201	CDL	C11-CA5-OA6-CA4
51	r	201	CDL	CB3-OB5-PB2-OB2

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Mol	Chain	Res	Type	Atoms
51	r	201	CDL	CB3-OB5-PB2-OB3
51	r	201	CDL	OB5-CB3-CB4-OB6
52	O	401	GTP	PB-O3B-PG-O2G
52	O	401	GTP	C5'-O5'-PA-O3A
52	O	401	GTP	C5'-O5'-PA-O1A
54	P	402	NDP	C5D-O5D-PN-O3
54	P	402	NDP	C5D-O5D-PN-O1N
54	P	402	NDP	C5D-O5D-PN-O2N
56	T	101	EHZ	S1-C10-C11-N1
56	T	101	EHZ	C11-C10-S1-C9
56	T	101	EHZ	C16-C17-C20-O6
56	T	101	EHZ	C18-C17-C20-O6
56	T	101	EHZ	C19-C17-C20-O6
56	T	101	EHZ	O2-C9-S1-C10
56	T	101	EHZ	C8-C9-S1-C10
56	U	101	EHZ	O2-C9-S1-C10
56	U	101	EHZ	C8-C9-S1-C10
51	N	902	CDL	C51-CB5-OB6-CB4
45	d	201	3PE	O32-C31-O31-C3
51	L	702	CDL	OA9-CA7-OA8-CA6
51	L	702	CDL	C31-CA7-OA8-CA6
45	M	601	3PE	O32-C31-O31-C3
45	Y	801	3PE	O32-C31-O31-C3
45	Y	803	3PE	O32-C31-O31-C3
45	Y	804	3PE	O32-C31-O31-C3
45	m	202	3PE	O32-C31-O31-C3
51	X	201	CDL	OB9-CB7-OB8-CB6
57	i	201	CHD	C13-C17-C20-C21
45	L	701	3PE	O22-C21-O21-C2
45	M	602	3PE	O22-C21-O21-C2
45	Y	801	3PE	O22-C21-O21-C2
45	m	202	3PE	O22-C21-O21-C2
46	B	202	PC1	O22-C21-O21-C2
46	M	606	PC1	O22-C21-O21-C2
51	X	201	CDL	OB7-CB5-OB6-CB4
51	d	203	CDL	OB7-CB5-OB6-CB4
51	h	201	CDL	OB7-CB5-OB6-CB4
45	M	601	3PE	C32-C31-O31-C3
45	M	605	3PE	C32-C31-O31-C3
45	Y	801	3PE	C32-C31-O31-C3
45	Y	803	3PE	C32-C31-O31-C3
45	Y	804	3PE	C32-C31-O31-C3

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Mol	Chain	Res	Type	Atoms
45	d	201	3PE	C32-C31-O31-C3
45	m	202	3PE	C32-C31-O31-C3
46	d	202	PC1	C32-C31-O31-C3
51	X	201	CDL	C71-CB7-OB8-CB6
51	h	201	CDL	C71-CB7-OB8-CB6
46	A	204	PC1	C22-C21-O21-C2
51	H	602	CDL	C51-CB5-OB6-CB4
51	h	201	CDL	C51-CB5-OB6-CB4
57	i	201	CHD	C16-C17-C20-C21
45	A	203	3PE	C32-C31-O31-C3
45	M	603	3PE	C32-C31-O31-C3
46	A	202	PC1	C32-C31-O31-C3
46	A	204	PC1	C32-C31-O31-C3
46	M	606	PC1	C32-C31-O31-C3
51	N	902	CDL	C31-CA7-OA8-CA6
45	M	602	3PE	O32-C31-O31-C3
45	M	605	3PE	O32-C31-O31-C3
46	M	604	PC1	O32-C31-O31-C3
46	M	606	PC1	O32-C31-O31-C3
51	N	902	CDL	OA9-CA7-OA8-CA6
51	N	902	CDL	OB9-CB7-OB8-CB6
51	h	201	CDL	OB9-CB7-OB8-CB6
45	A	201	3PE	O22-C21-O21-C2
51	r	201	CDL	OA7-CA5-OA6-CA4
45	M	602	3PE	C32-C31-O31-C3
46	M	604	PC1	C32-C31-O31-C3
51	N	902	CDL	C71-CB7-OB8-CB6
45	A	203	3PE	O32-C31-O31-C3
51	N	902	CDL	OB7-CB5-OB6-CB4
45	A	201	3PE	C22-C21-O21-C2
45	A	203	3PE	C22-C21-O21-C2
45	M	601	3PE	C22-C21-O21-C2
45	Y	803	3PE	C22-C21-O21-C2
46	A	202	PC1	C22-C21-O21-C2
51	H	602	CDL	C11-CA5-OA6-CA4
45	M	603	3PE	O32-C31-O31-C3
46	A	202	PC1	O32-C31-O31-C3
46	A	204	PC1	O32-C31-O31-C3
46	A	204	PC1	C2-C1-O11-P
46	m	201	PC1	C2-C1-O11-P
51	H	602	CDL	C1-CA2-OA2-PA1
51	h	201	CDL	CB4-CB3-OB5-PB2

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Mol	Chain	Res	Type	Atoms
45	L	701	3PE	C32-C31-O31-C3
51	r	201	CDL	C71-CB7-OB8-CB6
57	i	201	CHD	C13-C17-C20-C22
45	M	603	3PE	C22-C21-O21-C2
45	A	203	3PE	O22-C21-O21-C2
45	M	601	3PE	O22-C21-O21-C2
51	H	602	CDL	OA7-CA5-OA6-CA4
45	I	201	3PE	C32-C31-O31-C3
51	L	702	CDL	C71-CB7-OB8-CB6
46	Z	201	PC1	C11-C12-N-C13
51	L	702	CDL	O1-C1-CA2-OA2
51	h	201	CDL	O1-C1-CA2-OA2
45	I	201	3PE	O32-C31-O31-C3
46	A	202	PC1	O22-C21-O21-C2
51	L	702	CDL	OA5-CA3-CA4-OA6
45	P	403	3PE	C32-C31-O31-C3
45	Y	803	3PE	O22-C21-O21-C2
45	Y	804	3PE	C21-C22-C23-C24
51	r	201	CDL	CA5-C11-C12-C13
45	L	703	3PE	C22-C21-O21-C2
51	K	101	CDL	CA5-C11-C12-C13
51	X	201	CDL	CB7-C71-C72-C73
51	r	201	CDL	OB9-CB7-OB8-CB6
51	K	101	CDL	O1-C1-CB2-OB2
51	h	201	CDL	O1-C1-CB2-OB2
45	L	701	3PE	O32-C31-O31-C3
51	L	702	CDL	OB9-CB7-OB8-CB6
45	M	603	3PE	C31-C32-C33-C34
51	L	702	CDL	C11-CA5-OA6-CA4
51	X	201	CDL	C11-CA5-OA6-CA4
45	L	703	3PE	O22-C21-O21-C2
45	M	603	3PE	O22-C21-O21-C2
51	L	702	CDL	OA7-CA5-OA6-CA4
51	d	203	CDL	OA7-CA5-OA6-CA4
51	K	101	CDL	CA2-C1-CB2-OB2
51	h	201	CDL	CA2-C1-CB2-OB2
46	Z	201	PC1	C32-C31-O31-C3
46	Z	201	PC1	C11-C12-N-C14
46	Z	201	PC1	C11-C12-N-C15
45	b	101	3PE	C32-C31-O31-C3
46	P	401	PC1	C32-C31-O31-C3
46	M	604	PC1	C22-C21-O21-C2

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Mol	Chain	Res	Type	Atoms
51	d	203	CDL	C11-CA5-OA6-CA4
51	X	201	CDL	OA7-CA5-OA6-CA4
51	d	203	CDL	O1-C1-CB2-OB2
45	P	403	3PE	O32-C31-O31-C3
51	h	201	CDL	CB6-CB4-OB6-CB5
46	M	604	PC1	O22-C21-O21-C2
45	P	403	3PE	C22-C21-O21-C2
51	X	201	CDL	C36-C37-C38-C39
45	I	201	3PE	C38-C39-C3A-C3B
45	L	703	3PE	C38-C39-C3A-C3B
46	M	606	PC1	C38-C39-C3A-C3B
51	X	201	CDL	C33-C34-C35-C36
45	Y	804	3PE	C34-C35-C36-C37
45	Y	804	3PE	C29-C2A-C2B-C2C
45	b	101	3PE	C2E-C2F-C2G-C2H
46	d	202	PC1	C32-C33-C34-C35
51	L	702	CDL	C52-C53-C54-C55
51	N	902	CDL	C38-C39-C40-C41
51	X	201	CDL	C31-C32-C33-C34
58	o	201	MYR	C2-C3-C4-C5
45	I	201	3PE	C33-C34-C35-C36
51	X	201	CDL	C15-C16-C17-C18
45	H	601	3PE	C22-C21-O21-C2
45	L	701	3PE	C37-C38-C39-C3A
46	A	204	PC1	C32-C33-C34-C35
46	M	606	PC1	C24-C25-C26-C27
51	h	201	CDL	C16-C17-C18-C19
45	H	601	3PE	C21-C22-C23-C24
51	L	702	CDL	C15-C16-C17-C18
46	Z	201	PC1	O32-C31-O31-C3
51	h	201	CDL	C54-C55-C56-C57
46	M	604	PC1	C24-C25-C26-C27
45	b	101	3PE	O32-C31-O31-C3
46	P	401	PC1	O32-C31-O31-C3
45	N	901	3PE	C33-C34-C35-C36
45	Y	803	3PE	C25-C26-C27-C28
45	P	403	3PE	O22-C21-O21-C2
45	M	602	3PE	C32-C33-C34-C35
45	M	603	3PE	C35-C36-C37-C38
51	d	203	CDL	CA2-C1-CB2-OB2
45	M	602	3PE	C35-C36-C37-C38
51	d	203	CDL	CA5-C11-C12-C13

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Mol	Chain	Res	Type	Atoms
45	L	701	3PE	C23-C24-C25-C26
45	M	601	3PE	C39-C3A-C3B-C3C
46	A	202	PC1	C26-C27-C28-C29
51	X	201	CDL	C20-C21-C22-C23
51	r	201	CDL	C59-C60-C61-C62
51	N	902	CDL	C33-C34-C35-C36
56	U	101	EHZ	C21-C1-C2-C3
45	M	602	3PE	C39-C3A-C3B-C3C
45	M	603	3PE	C23-C24-C25-C26
51	h	201	CDL	C12-C13-C14-C15
51	K	101	CDL	C11-CA5-OA6-CA4
45	A	201	3PE	C32-C31-O31-C3
45	M	603	3PE	C2A-C2B-C2C-C2D
51	K	101	CDL	C71-C72-C73-C74
46	M	604	PC1	C28-C29-C2A-C2B
45	H	601	3PE	O22-C21-O21-C2
45	Y	804	3PE	C3B-C3C-C3D-C3E
46	M	604	PC1	C2B-C2C-C2D-C2E
46	Z	201	PC1	C38-C39-C3A-C3B
45	H	601	3PE	C23-C24-C25-C26
45	M	601	3PE	C2D-C2E-C2F-C2G
45	H	601	3PE	C26-C27-C28-C29
51	r	201	CDL	C51-C52-C53-C54
51	d	203	CDL	C35-C36-C37-C38
45	d	201	3PE	C21-C22-C23-C24
45	m	202	3PE	C29-C2A-C2B-C2C
45	M	601	3PE	C34-C35-C36-C37
45	Y	802	3PE	C22-C21-O21-C2
45	Y	804	3PE	C22-C21-O21-C2
45	Z	202	3PE	C22-C21-O21-C2
46	h	202	PC1	C22-C21-O21-C2
45	Y	804	3PE	O22-C21-O21-C2
51	H	602	CDL	C51-C52-C53-C54
45	Z	202	3PE	C22-C23-C24-C25
51	X	201	CDL	C39-C40-C41-C42
51	d	203	CDL	C42-C43-C44-C45
45	L	701	3PE	C36-C37-C38-C39
46	Z	201	PC1	C34-C35-C36-C37
46	h	202	PC1	C26-C27-C28-C29
45	N	901	3PE	C2C-C2D-C2E-C2F
45	Z	202	3PE	C31-C32-C33-C34
45	N	901	3PE	C39-C3A-C3B-C3C

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Mol	Chain	Res	Type	Atoms
51	d	203	CDL	OB5-CB3-CB4-OB6
45	Z	202	3PE	C32-C33-C34-C35
45	b	101	3PE	C31-C32-C33-C34
45	Y	802	3PE	O21-C2-C3-O31
51	K	101	CDL	C71-CB7-OB8-CB6
45	A	201	3PE	O32-C31-O31-C3
45	Y	802	3PE	C23-C24-C25-C26
51	N	902	CDL	C37-C38-C39-C40
51	d	203	CDL	C33-C34-C35-C36
58	o	201	MYR	C5-C6-C7-C8
45	N	901	3PE	C34-C35-C36-C37
45	N	901	3PE	C23-C24-C25-C26
45	Y	804	3PE	C27-C28-C29-C2A
45	M	602	3PE	C2-C1-O11-P
51	X	201	CDL	C1-CB2-OB2-PB2
45	b	101	3PE	C28-C29-C2A-C2B
45	Y	802	3PE	O22-C21-O21-C2
58	o	201	MYR	C3-C4-C5-C6
51	d	203	CDL	C31-C32-C33-C34
45	M	605	3PE	C22-C21-O21-C2
45	L	703	3PE	C33-C34-C35-C36
46	m	201	PC1	C39-C3A-C3B-C3C
46	P	401	PC1	C26-C27-C28-C29
45	M	603	3PE	C3E-C3F-C3G-C3H
45	b	101	3PE	C35-C36-C37-C38
45	H	601	3PE	O11-C1-C2-C3
45	N	901	3PE	O11-C1-C2-C3
45	Y	801	3PE	O11-C1-C2-C3
45	m	202	3PE	O11-C1-C2-C3
46	A	202	PC1	O11-C1-C2-C3
46	Z	201	PC1	O11-C1-C2-C3
51	L	702	CDL	OA5-CA3-CA4-CA6
51	X	201	CDL	OB5-CB3-CB4-CB6
51	r	201	CDL	OB5-CB3-CB4-CB6
45	Z	202	3PE	O22-C21-O21-C2
51	K	101	CDL	OA7-CA5-OA6-CA4
46	A	202	PC1	C22-C23-C24-C25
51	X	201	CDL	C12-C13-C14-C15
51	h	201	CDL	C39-C40-C41-C42
45	d	201	3PE	C22-C21-O21-C2
45	M	601	3PE	C26-C27-C28-C29
46	A	204	PC1	C21-C22-C23-C24

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Mol	Chain	Res	Type	Atoms
45	Y	801	3PE	C1-C2-C3-O31
51	r	201	CDL	CA3-CA4-CA6-OA8
51	N	902	CDL	C14-C15-C16-C17
45	Y	802	3PE	C21-C22-C23-C24
45	L	701	3PE	C26-C27-C28-C29
46	P	401	PC1	C3B-C3C-C3D-C3E
45	m	202	3PE	C2D-C2E-C2F-C2G
56	T	101	EHZ	C2-C1-C21-C22
45	M	602	3PE	C31-C32-C33-C34
51	K	101	CDL	OB9-CB7-OB8-CB6
46	h	202	PC1	C2-C1-O11-P
45	I	201	3PE	C36-C37-C38-C39
45	Y	803	3PE	C3-C2-O21-C21
51	H	602	CDL	CA3-CA4-OA6-CA5
51	K	101	CDL	C40-C41-C42-C43
51	L	702	CDL	C56-C57-C58-C59
45	M	601	3PE	O11-C1-C2-O21
46	A	202	PC1	O11-C1-C2-O21
45	A	201	3PE	C33-C34-C35-C36
45	L	703	3PE	C39-C3A-C3B-C3C
46	Z	201	PC1	C2C-C2D-C2E-C2F
51	h	201	CDL	C75-C76-C77-C78
45	M	605	3PE	C23-C24-C25-C26
57	i	201	CHD	C16-C17-C20-C22
45	M	601	3PE	C36-C37-C38-C39
46	M	606	PC1	C3A-C3B-C3C-C3D
45	N	901	3PE	C28-C29-C2A-C2B
51	N	902	CDL	OA6-CA4-CA6-OA8
51	h	201	CDL	C73-C74-C75-C76
46	P	401	PC1	C37-C38-C39-C3A
46	Z	201	PC1	O21-C21-C22-C23
56	T	101	EHZ	C21-C22-C23-C24
45	d	201	3PE	C3D-C3E-C3F-C3G
45	b	101	3PE	C2C-C2D-C2E-C2F
51	H	602	CDL	C31-CA7-OA8-CA6
51	r	201	CDL	C57-C58-C59-C60
45	M	601	3PE	C35-C36-C37-C38
46	h	202	PC1	C2A-C2B-C2C-C2D
45	Y	804	3PE	C2F-C2G-C2H-C2I
51	d	203	CDL	CA7-C31-C32-C33
51	d	203	CDL	C76-C77-C78-C79
45	M	605	3PE	C3C-C3D-C3E-C3F

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Mol	Chain	Res	Type	Atoms
45	M	605	3PE	C25-C26-C27-C28
46	h	202	PC1	O22-C21-O21-C2
51	X	201	CDL	CB4-CB3-OB5-PB2
51	h	201	CDL	C1-CA2-OA2-PA1
45	H	601	3PE	C22-C23-C24-C25
45	M	603	3PE	C2D-C2E-C2F-C2G
45	L	701	3PE	C25-C26-C27-C28
45	M	601	3PE	O11-C1-C2-C3
45	Y	803	3PE	O11-C1-C2-C3
51	K	101	CDL	OB5-CB3-CB4-CB6
51	L	702	CDL	OB5-CB3-CB4-CB6
51	d	203	CDL	OB5-CB3-CB4-CB6
46	h	202	PC1	C35-C36-C37-C38
45	m	202	3PE	C26-C27-C28-C29
46	M	604	PC1	C2C-C2D-C2E-C2F
51	K	101	CDL	C38-C39-C40-C41
45	A	203	3PE	C1-C2-C3-O31
46	M	604	PC1	C1-C2-C3-O31
46	P	401	PC1	C1-C2-C3-O31
51	H	602	CDL	CB3-CB4-CB6-OB8
51	N	902	CDL	CA3-CA4-CA6-OA8
45	M	603	3PE	C39-C3A-C3B-C3C
45	M	601	3PE	C33-C34-C35-C36
45	L	701	3PE	C3A-C3B-C3C-C3D
46	M	606	PC1	C3C-C3D-C3E-C3F
45	M	601	3PE	C22-C23-C24-C25
46	M	604	PC1	C2E-C2F-C2G-C2H
51	K	101	CDL	C52-C53-C54-C55
56	U	101	EHZ	C4-C5-C6-C7
45	b	101	3PE	C39-C3A-C3B-C3C
45	I	201	3PE	O11-C1-C2-O21
46	d	202	PC1	O11-C1-C2-O21
51	L	702	CDL	OB5-CB3-CB4-OB6
46	B	202	PC1	C37-C38-C39-C3A
45	M	603	3PE	C2-C1-O11-P
45	N	901	3PE	C2-C1-O11-P
51	K	101	CDL	C1-CB2-OB2-PB2
45	m	202	3PE	C21-C22-C23-C24
56	T	101	EHZ	O1-C7-C8-C9
45	Y	801	3PE	C34-C35-C36-C37
46	M	606	PC1	C33-C34-C35-C36
56	T	101	EHZ	C3-C4-C5-C6

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Mol	Chain	Res	Type	Atoms
45	M	601	3PE	C2E-C2F-C2G-C2H
51	h	201	CDL	C52-C51-CB5-OB6
45	A	203	3PE	O21-C2-C3-O31
45	H	601	3PE	O21-C2-C3-O31
45	Y	801	3PE	O21-C2-C3-O31
45	M	605	3PE	C34-C35-C36-C37
45	M	601	3PE	C2C-C2D-C2E-C2F
45	M	603	3PE	C28-C29-C2A-C2B
46	Z	201	PC1	C35-C36-C37-C38
46	d	202	PC1	C34-C35-C36-C37
58	o	201	MYR	C7-C8-C9-C10
45	L	701	3PE	C21-C22-C23-C24
45	M	601	3PE	C31-C32-C33-C34
45	M	605	3PE	C3A-C3B-C3C-C3D
45	P	403	3PE	C35-C36-C37-C38
56	T	101	EHZ	C6-C7-C8-C9
45	d	201	3PE	O22-C21-O21-C2
51	N	902	CDL	C11-C12-C13-C14
45	b	101	3PE	O21-C21-C22-C23
45	Z	202	3PE	C29-C2A-C2B-C2C
45	I	201	3PE	O11-C1-C2-C3
45	M	605	3PE	O11-C1-C2-C3
46	P	401	PC1	O11-C1-C2-C3
51	d	203	CDL	OA5-CA3-CA4-CA6
45	M	605	3PE	O22-C21-O21-C2
45	b	101	3PE	C38-C39-C3A-C3B
46	P	401	PC1	C32-C33-C34-C35
51	K	101	CDL	C15-C16-C17-C18
45	P	403	3PE	C1-C2-O21-C21
46	A	204	PC1	C3-C2-O21-C21
46	d	202	PC1	C21-C22-C23-C24
45	b	101	3PE	C33-C34-C35-C36
51	K	101	CDL	C36-C37-C38-C39
51	H	602	CDL	OA9-CA7-OA8-CA6
45	d	201	3PE	C37-C38-C39-C3A
45	A	201	3PE	C28-C29-C2A-C2B
45	H	601	3PE	O11-C1-C2-O21
45	L	701	3PE	O11-C1-C2-O21
45	M	605	3PE	O11-C1-C2-O21
45	m	202	3PE	O11-C1-C2-O21
46	P	401	PC1	O11-C1-C2-O21
51	K	101	CDL	OB5-CB3-CB4-OB6

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Mol	Chain	Res	Type	Atoms
51	X	201	CDL	OA5-CA3-CA4-OA6
51	L	702	CDL	C31-C32-C33-C34
46	M	606	PC1	C1-C2-C3-O31
51	r	201	CDL	CB3-CB4-CB6-OB8
45	L	703	3PE	C32-C33-C34-C35
45	A	203	3PE	C12-C11-O13-P
45	M	605	3PE	C12-C11-O13-P
45	Y	803	3PE	C12-C11-O13-P
46	m	201	PC1	C12-C11-O13-P
45	M	602	3PE	O21-C2-C3-O31
45	M	603	3PE	O21-C2-C3-O31
46	M	604	PC1	O21-C2-C3-O31
46	P	401	PC1	O21-C2-C3-O31
51	H	602	CDL	OB6-CB4-CB6-OB8
51	r	201	CDL	OA6-CA4-CA6-OA8
51	r	201	CDL	OB6-CB4-CB6-OB8
51	K	101	CDL	C53-C54-C55-C56
45	M	603	3PE	C33-C34-C35-C36
45	Y	802	3PE	C2C-C2D-C2E-C2F
45	I	201	3PE	C24-C25-C26-C27
45	b	101	3PE	C24-C25-C26-C27
46	A	202	PC1	O13-C11-C12-N
46	M	606	PC1	O13-C11-C12-N
46	h	202	PC1	O13-C11-C12-N
46	P	401	PC1	C28-C29-C2A-C2B
45	L	703	3PE	C26-C27-C28-C29
46	M	604	PC1	C23-C24-C25-C26
45	M	602	3PE	C38-C39-C3A-C3B
45	Y	802	3PE	C2E-C2F-C2G-C2H
51	L	702	CDL	CB4-CB6-OB8-CB7
45	b	101	3PE	C3A-C3B-C3C-C3D
51	K	101	CDL	C31-CA7-OA8-CA6
46	h	202	PC1	C22-C23-C24-C25
51	h	201	CDL	C11-C12-C13-C14
45	m	202	3PE	C24-C25-C26-C27
45	M	603	3PE	C32-C33-C34-C35
45	Y	801	3PE	C37-C38-C39-C3A
46	h	202	PC1	C32-C33-C34-C35
51	L	702	CDL	C57-C58-C59-C60
45	I	201	3PE	C25-C26-C27-C28
45	M	601	3PE	C2-C1-O11-P
51	d	203	CDL	CB4-CB3-OB5-PB2

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Mol	Chain	Res	Type	Atoms
45	Y	801	3PE	O11-C1-C2-O21
45	d	201	3PE	O11-C1-C2-O21
51	H	602	CDL	OA5-CA3-CA4-OA6
51	d	203	CDL	OA5-CA3-CA4-OA6
45	M	602	3PE	C2B-C2C-C2D-C2E
46	M	606	PC1	C34-C35-C36-C37
51	K	101	CDL	C51-CB5-OB6-CB4
51	K	101	CDL	OA9-CA7-OA8-CA6
45	d	201	3PE	C3A-C3B-C3C-C3D
54	P	402	NDP	O4D-C1D-N1N-C6N
46	d	202	PC1	O21-C2-C3-O31
45	L	701	3PE	C22-C23-C24-C25
45	M	603	3PE	C29-C2A-C2B-C2C
45	Y	802	3PE	C1-C2-C3-O31
45	A	203	3PE	C26-C27-C28-C29
45	d	201	3PE	C33-C34-C35-C36
45	Y	804	3PE	C39-C3A-C3B-C3C
45	A	203	3PE	C24-C25-C26-C27
45	A	201	3PE	C1-O11-P-O14
45	A	201	3PE	C11-O13-P-O12
45	A	203	3PE	C11-O13-P-O11
45	A	203	3PE	C11-O13-P-O12
45	L	701	3PE	C1-O11-P-O14
45	M	601	3PE	C11-O13-P-O11
45	M	602	3PE	C1-O11-P-O12
45	M	602	3PE	C1-O11-P-O13
45	M	602	3PE	C1-O11-P-O14
45	M	602	3PE	C11-O13-P-O12
45	M	605	3PE	C11-O13-P-O11
45	M	605	3PE	C11-O13-P-O12
45	M	605	3PE	C11-O13-P-O14
45	N	901	3PE	C11-O13-P-O11
45	Y	802	3PE	C11-O13-P-O14
45	Y	803	3PE	C1-O11-P-O13
45	Z	202	3PE	C1-O11-P-O13
45	Z	202	3PE	C1-O11-P-O14
45	d	201	3PE	C11-O13-P-O12
46	A	202	PC1	C11-O13-P-O14
46	A	202	PC1	C11-O13-P-O11
46	B	202	PC1	C11-O13-P-O14
46	M	606	PC1	C1-O11-P-O14
46	P	401	PC1	C11-O13-P-O14

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Mol	Chain	Res	Type	Atoms
46	P	401	PC1	C1-O11-P-O14
46	h	202	PC1	C1-O11-P-O12
46	m	201	PC1	C1-O11-P-O14
51	H	602	CDL	CA2-OA2-PA1-OA3
51	H	602	CDL	CB2-OB2-PB2-OB3
51	H	602	CDL	CB3-OB5-PB2-OB2
51	H	602	CDL	CB3-OB5-PB2-OB3
51	K	101	CDL	CB3-OB5-PB2-OB3
51	L	702	CDL	CA3-OA5-PA1-OA4
51	L	702	CDL	CB2-OB2-PB2-OB4
51	N	902	CDL	CA2-OA2-PA1-OA4
51	N	902	CDL	CA3-OA5-PA1-OA2
51	N	902	CDL	CA3-OA5-PA1-OA3
51	N	902	CDL	CA3-OA5-PA1-OA4
51	X	201	CDL	CA3-OA5-PA1-OA4
51	X	201	CDL	CB3-OB5-PB2-OB3
51	d	203	CDL	CA2-OA2-PA1-OA3
51	d	203	CDL	CA3-OA5-PA1-OA2
51	h	201	CDL	CB3-OB5-PB2-OB4
51	r	201	CDL	CB3-OB5-PB2-OB4
52	O	401	GTP	C5'-O5'-PA-O2A
45	M	603	3PE	C34-C35-C36-C37
51	h	201	CDL	CA7-C31-C32-C33
45	M	605	3PE	C2-C1-O11-P
45	Y	801	3PE	C2-C1-O11-P
45	b	101	3PE	C2-C1-O11-P
51	H	602	CDL	CB4-CB3-OB5-PB2
51	N	902	CDL	CB4-CB3-OB5-PB2
51	X	201	CDL	CA4-CA3-OA5-PA1
51	h	201	CDL	C1-CB2-OB2-PB2
45	b	101	3PE	C23-C24-C25-C26
51	X	201	CDL	OA9-CA7-OA8-CA6
46	A	202	PC1	C33-C34-C35-C36
51	N	902	CDL	CA5-C11-C12-C13
51	X	201	CDL	C31-CA7-OA8-CA6
51	L	702	CDL	C14-C15-C16-C17
51	N	902	CDL	CB3-CB4-OB6-CB5
51	X	201	CDL	CB6-CB4-OB6-CB5
46	M	604	PC1	C2D-C2E-C2F-C2G
45	L	701	3PE	O11-C1-C2-C3
51	H	602	CDL	OA5-CA3-CA4-CA6
45	Y	801	3PE	C21-C22-C23-C24

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Mol	Chain	Res	Type	Atoms
45	Y	804	3PE	C31-C32-C33-C34
45	Y	803	3PE	O11-C1-C2-O21
51	X	201	CDL	OB5-CB3-CB4-OB6
51	X	201	CDL	C34-C35-C36-C37
51	h	201	CDL	C17-C18-C19-C20
45	M	603	3PE	C3A-C3B-C3C-C3D
56	U	101	EHZ	C21-C22-C23-C24
51	K	101	CDL	OB7-CB5-OB6-CB4
46	M	604	PC1	C26-C27-C28-C29
46	d	202	PC1	C36-C37-C38-C39
46	A	204	PC1	O21-C2-C3-O31
46	P	401	PC1	O21-C21-C22-C23
45	L	703	3PE	C36-C37-C38-C39
46	P	401	PC1	C2A-C2B-C2C-C2D
51	d	203	CDL	C74-C75-C76-C77
45	Y	802	3PE	O31-C31-C32-C33
45	M	603	3PE	C2C-C2D-C2E-C2F
51	X	201	CDL	C24-C25-C26-C27
51	d	203	CDL	C72-C73-C74-C75
56	U	101	EHZ	C7-C8-C9-O2
45	Y	801	3PE	C39-C3A-C3B-C3C
51	H	602	CDL	CB2-C1-CA2-OA2
46	B	202	PC1	C28-C29-C2A-C2B
45	A	201	3PE	C31-C32-C33-C34
51	K	101	CDL	C11-C12-C13-C14
46	Z	201	PC1	O22-C21-C22-C23
45	Y	804	3PE	C28-C29-C2A-C2B
45	M	603	3PE	C37-C38-C39-C3A
51	N	902	CDL	C32-C31-CA7-OA8
46	m	201	PC1	C32-C33-C34-C35
51	K	101	CDL	C35-C36-C37-C38
45	Y	802	3PE	C24-C25-C26-C27
45	I	201	3PE	C26-C27-C28-C29
51	h	201	CDL	C15-C16-C17-C18
46	M	606	PC1	C23-C24-C25-C26
45	b	101	3PE	C2A-C2B-C2C-C2D
45	Y	802	3PE	C28-C29-C2A-C2B
45	M	602	3PE	C1-C2-C3-O31
45	A	201	3PE	C23-C24-C25-C26
45	Y	804	3PE	C24-C25-C26-C27
45	I	201	3PE	C21-C22-C23-C24
45	L	703	3PE	C28-C29-C2A-C2B

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Mol	Chain	Res	Type	Atoms
46	B	202	PC1	C34-C35-C36-C37
46	Z	201	PC1	C22-C23-C24-C25
45	Y	804	3PE	C3-C2-O21-C21
51	X	201	CDL	CA6-CA4-OA6-CA5
51	d	203	CDL	C44-C45-C46-C47
45	M	603	3PE	C2B-C2C-C2D-C2E
45	M	603	3PE	O11-C1-C2-C3
45	P	403	3PE	O11-C1-C2-C3
45	d	201	3PE	O11-C1-C2-C3
51	X	201	CDL	OA5-CA3-CA4-CA6
46	d	202	PC1	O22-C21-O21-C2
45	b	101	3PE	C32-C33-C34-C35
45	m	202	3PE	C25-C26-C27-C28
46	M	606	PC1	O21-C2-C3-O31
45	L	703	3PE	C34-C35-C36-C37
45	N	901	3PE	C2A-C2B-C2C-C2D
51	r	201	CDL	C52-C53-C54-C55
51	r	201	CDL	C19-C20-C21-C22
51	h	201	CDL	C32-C33-C34-C35
45	Y	803	3PE	C22-C23-C24-C25
46	d	202	PC1	C22-C21-O21-C2
51	H	602	CDL	C33-C34-C35-C36
51	K	101	CDL	C16-C17-C18-C19
45	M	602	3PE	C28-C29-C2A-C2B
51	N	902	CDL	C1-CA2-OA2-PA1
46	h	202	PC1	C29-C2A-C2B-C2C
51	K	101	CDL	C33-C34-C35-C36
45	Z	202	3PE	C34-C35-C36-C37
46	d	202	PC1	C1-C2-C3-O31
57	i	201	CHD	C22-C23-C24-O26
51	N	902	CDL	OB5-CB3-CB4-OB6
45	Z	202	3PE	C27-C28-C29-C2A
51	h	201	CDL	C76-C77-C78-C79
45	m	202	3PE	O21-C21-C22-C23
46	d	202	PC1	O11-C1-C2-C3
51	h	201	CDL	C52-C51-CB5-OB7
49	F	501	FMN	C4'-C5'-O5'-P
51	L	702	CDL	CA4-CA3-OA5-PA1
45	M	602	3PE	C24-C25-C26-C27
46	A	204	PC1	C24-C25-C26-C27
57	i	201	CHD	C22-C23-C24-O25
45	P	403	3PE	C39-C3A-C3B-C3C

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Mol	Chain	Res	Type	Atoms
45	A	201	3PE	C34-C35-C36-C37
46	M	606	PC1	C39-C3A-C3B-C3C
45	m	202	3PE	O31-C31-C32-C33
51	X	201	CDL	CA3-CA4-OA6-CA5
46	A	204	PC1	C25-C26-C27-C28
51	r	201	CDL	C17-C18-C19-C20
51	d	203	CDL	OA9-CA7-OA8-CA6
51	r	201	CDL	CB2-C1-CA2-OA2
51	H	602	CDL	OB5-CB3-CB4-OB6
45	H	601	3PE	C1-C2-C3-O31
45	L	703	3PE	C23-C24-C25-C26
46	A	204	PC1	C12-C11-O13-P
46	P	401	PC1	C35-C36-C37-C38
51	d	203	CDL	C31-CA7-OA8-CA6
45	b	101	3PE	C37-C38-C39-C3A
51	K	101	CDL	C13-C14-C15-C16
45	M	601	3PE	O21-C2-C3-O31
51	h	201	CDL	C58-C59-C60-C61
46	A	202	PC1	C34-C35-C36-C37
51	H	602	CDL	O1-C1-CA2-OA2
45	M	603	3PE	O31-C31-C32-C33
51	h	201	CDL	OB5-CB3-CB4-CB6
45	M	603	3PE	C22-C23-C24-C25
51	K	101	CDL	C18-C19-C20-C21
51	r	201	CDL	C72-C71-CB7-OB8
46	m	201	PC1	O21-C21-C22-C23
46	P	401	PC1	C33-C34-C35-C36
45	N	901	3PE	C29-C2A-C2B-C2C
51	H	602	CDL	C72-C71-CB7-OB8
45	Y	804	3PE	C35-C36-C37-C38
52	O	401	GTP	PB-O3A-PA-O2A
46	M	604	PC1	C2A-C2B-C2C-C2D
45	M	603	3PE	C3B-C3C-C3D-C3E
45	M	602	3PE	C2C-C2D-C2E-C2F
51	h	201	CDL	C31-C32-C33-C34
45	H	601	3PE	C28-C29-C2A-C2B
51	X	201	CDL	C51-C52-C53-C54
45	d	201	3PE	O31-C31-C32-C33
58	o	201	MYR	C1-C2-C3-C4
45	M	602	3PE	C3C-C3D-C3E-C3F
46	d	202	PC1	O21-C21-C22-C23
46	B	202	PC1	C22-C23-C24-C25

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Mol	Chain	Res	Type	Atoms
46	Z	201	PC1	C21-C22-C23-C24
46	A	204	PC1	O21-C21-C22-C23
54	P	402	NDP	O4B-C4B-C5B-O5B
45	b	101	3PE	O22-C21-C22-C23
54	P	402	NDP	C2B-O2B-P2B-O3X
45	A	203	3PE	C32-C33-C34-C35
45	Y	801	3PE	C36-C37-C38-C39
51	X	201	CDL	CA5-C11-C12-C13
45	Y	804	3PE	O21-C21-C22-C23
46	B	202	PC1	O21-C21-C22-C23
45	M	603	3PE	C1-C2-C3-O31
46	A	204	PC1	C1-C2-C3-O31
45	M	602	3PE	O21-C21-C22-C23
46	h	202	PC1	O31-C31-C32-C33
51	X	201	CDL	C12-C11-CA5-OA6
51	H	602	CDL	OA6-CA4-CA6-OA8
51	h	201	CDL	C51-C52-C53-C54
45	Z	202	3PE	C28-C29-C2A-C2B
51	L	702	CDL	C54-C55-C56-C57
49	F	501	FMN	N10-C1'-C2'-O2'
56	U	101	EHZ	C7-C8-C9-S1
45	M	602	3PE	C2E-C2F-C2G-C2H
45	M	605	3PE	C2B-C2C-C2D-C2E
45	Y	804	3PE	C1-C2-O21-C21
46	d	202	PC1	C35-C36-C37-C38
51	L	702	CDL	C32-C31-CA7-OA8
46	A	204	PC1	O22-C21-C22-C23
45	M	603	3PE	O32-C31-C32-C33
45	d	201	3PE	C3B-C3C-C3D-C3E
51	X	201	CDL	C12-C11-CA5-OA7
46	B	202	PC1	C32-C31-O31-C3
45	d	201	3PE	O32-C31-C32-C33
51	d	203	CDL	C38-C39-C40-C41
45	H	601	3PE	O21-C21-C22-C23
45	N	901	3PE	O21-C21-C22-C23
46	d	202	PC1	O22-C21-C22-C23
46	m	201	PC1	O22-C21-C22-C23
45	N	901	3PE	O22-C21-O21-C2
46	h	202	PC1	O32-C31-C32-C33
45	L	703	3PE	C1-C2-C3-O31
46	B	202	PC1	C1-C2-C3-O31
45	Y	804	3PE	O22-C21-C22-C23

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Mol	Chain	Res	Type	Atoms
46	B	202	PC1	O22-C21-C22-C23
51	d	203	CDL	O1-C1-CA2-OA2
51	L	702	CDL	C32-C31-CA7-OA9
45	b	101	3PE	C21-C22-C23-C24
45	M	603	3PE	C2E-C2F-C2G-C2H
45	P	403	3PE	O21-C21-C22-C23
51	X	201	CDL	C72-C71-CB7-OB8
52	O	401	GTP	PG-O3B-PB-O2B
52	O	401	GTP	PA-O3A-PB-O2B
45	H	601	3PE	O22-C21-C22-C23
45	M	602	3PE	O22-C21-C22-C23

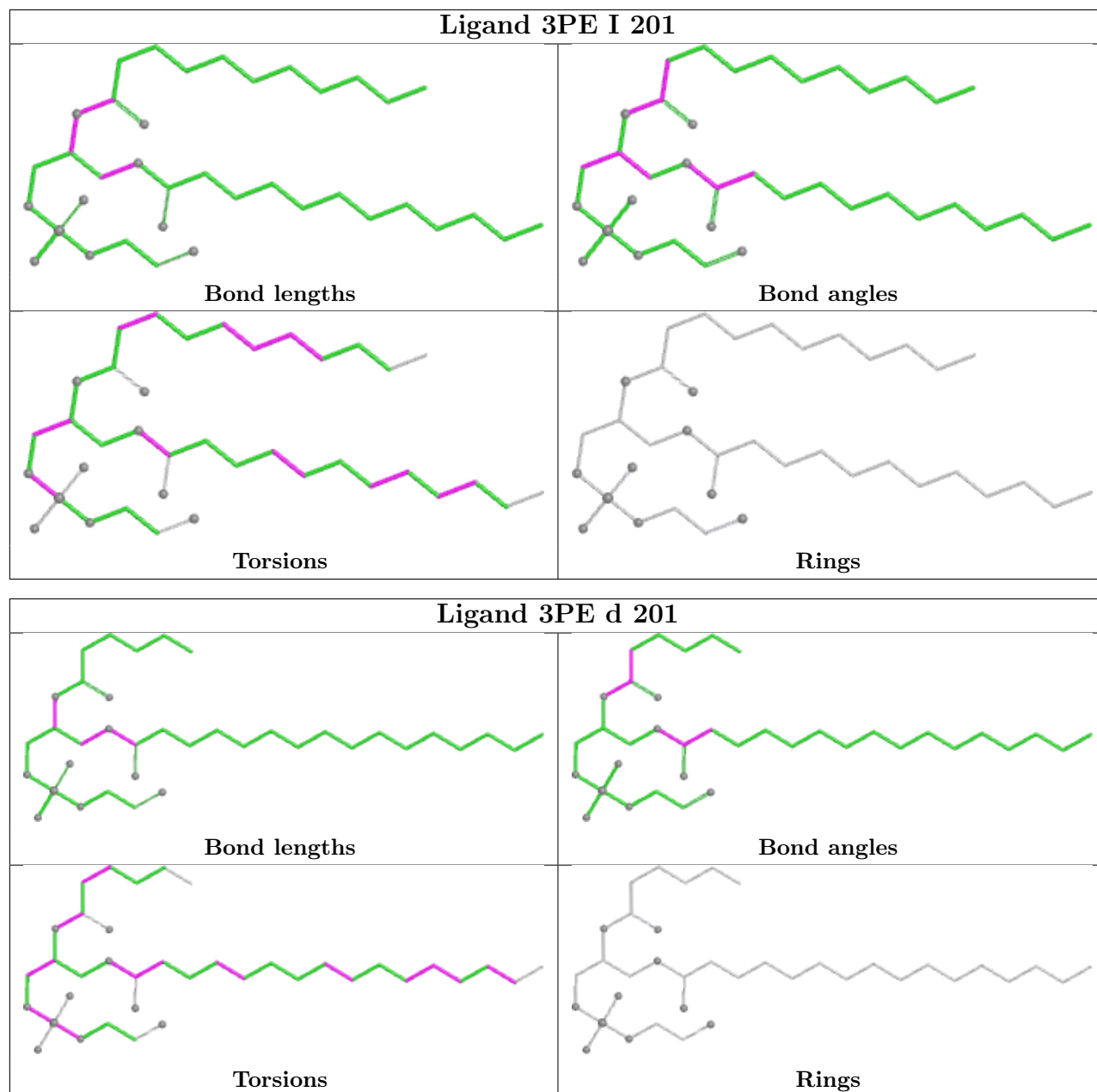
There are no ring outliers.

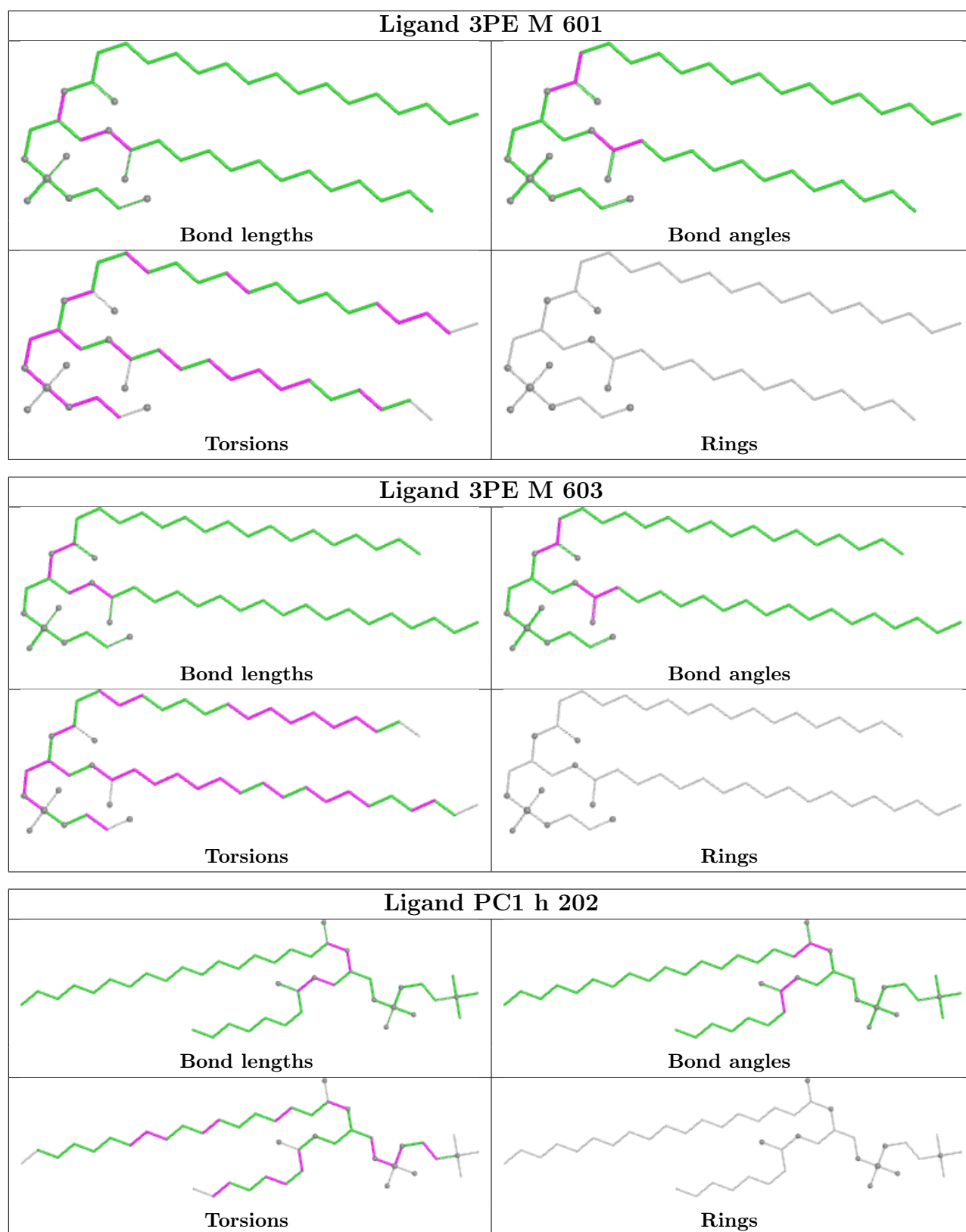
18 monomers are involved in 36 short contacts:

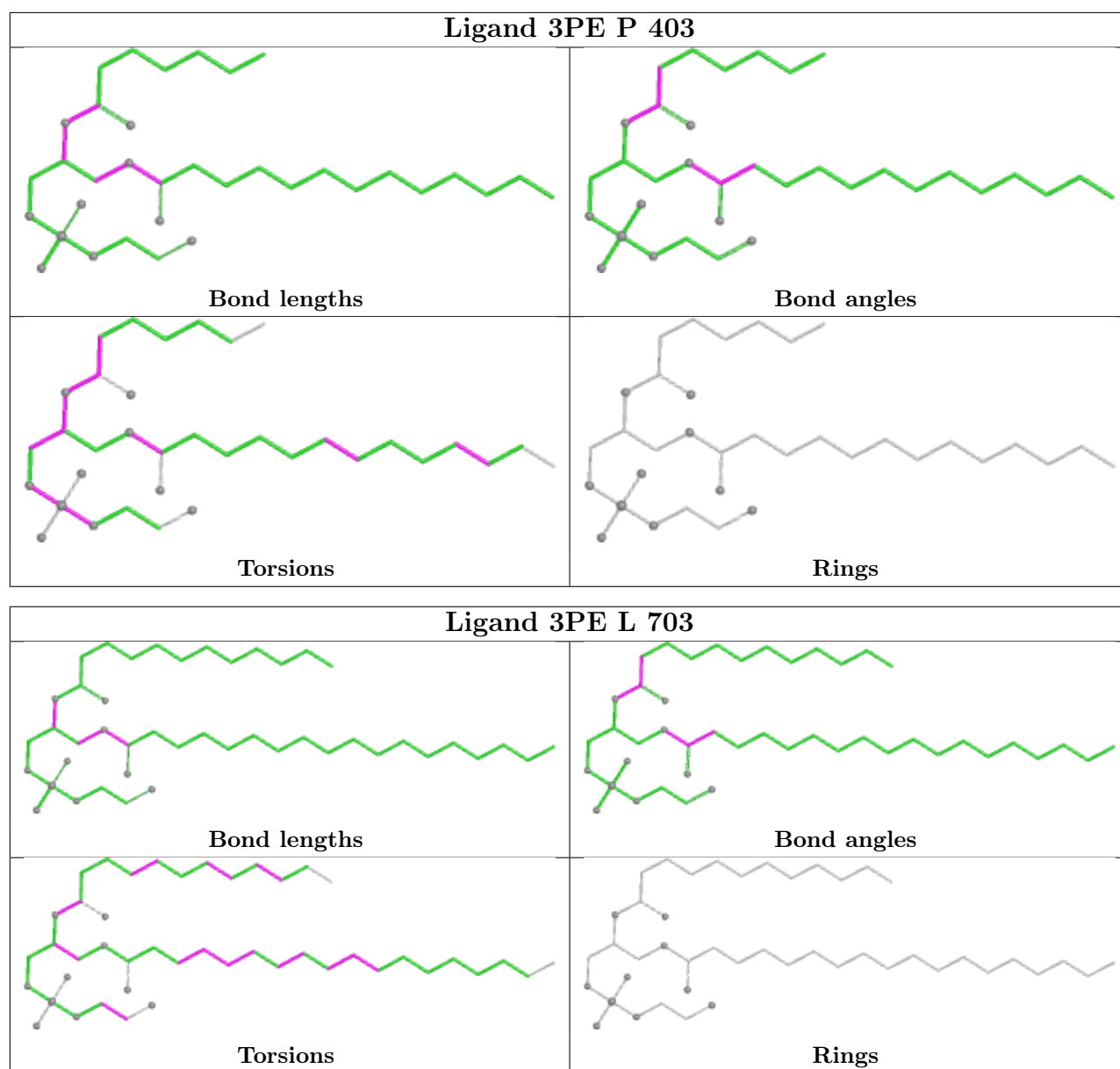
Mol	Chain	Res	Type	Clashes	Symm-Clashes
45	M	603	3PE	3	0
46	h	202	PC1	3	0
54	P	402	NDP	3	0
47	F	502	SF4	1	0
45	M	602	3PE	2	0
46	M	604	PC1	1	0
45	H	601	3PE	1	0
49	F	501	FMN	2	0
45	M	605	3PE	1	0
52	O	401	GTP	2	0
46	Z	201	PC1	5	0
46	P	401	PC1	1	0
46	d	202	PC1	1	0
51	X	201	CDL	4	0
56	T	101	EHZ	1	0
57	i	201	CHD	2	0
46	B	202	PC1	1	0
45	m	202	3PE	2	0

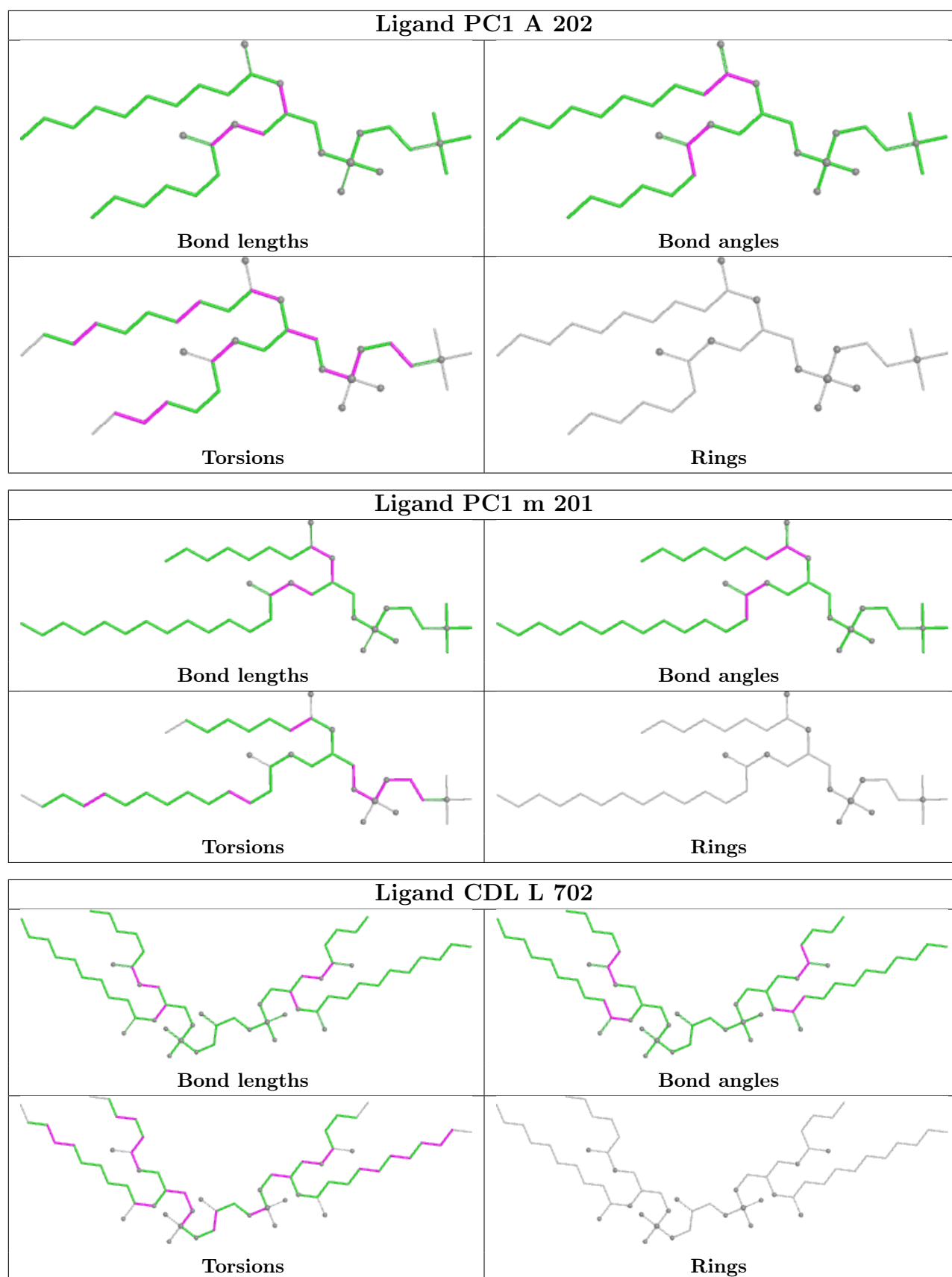
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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

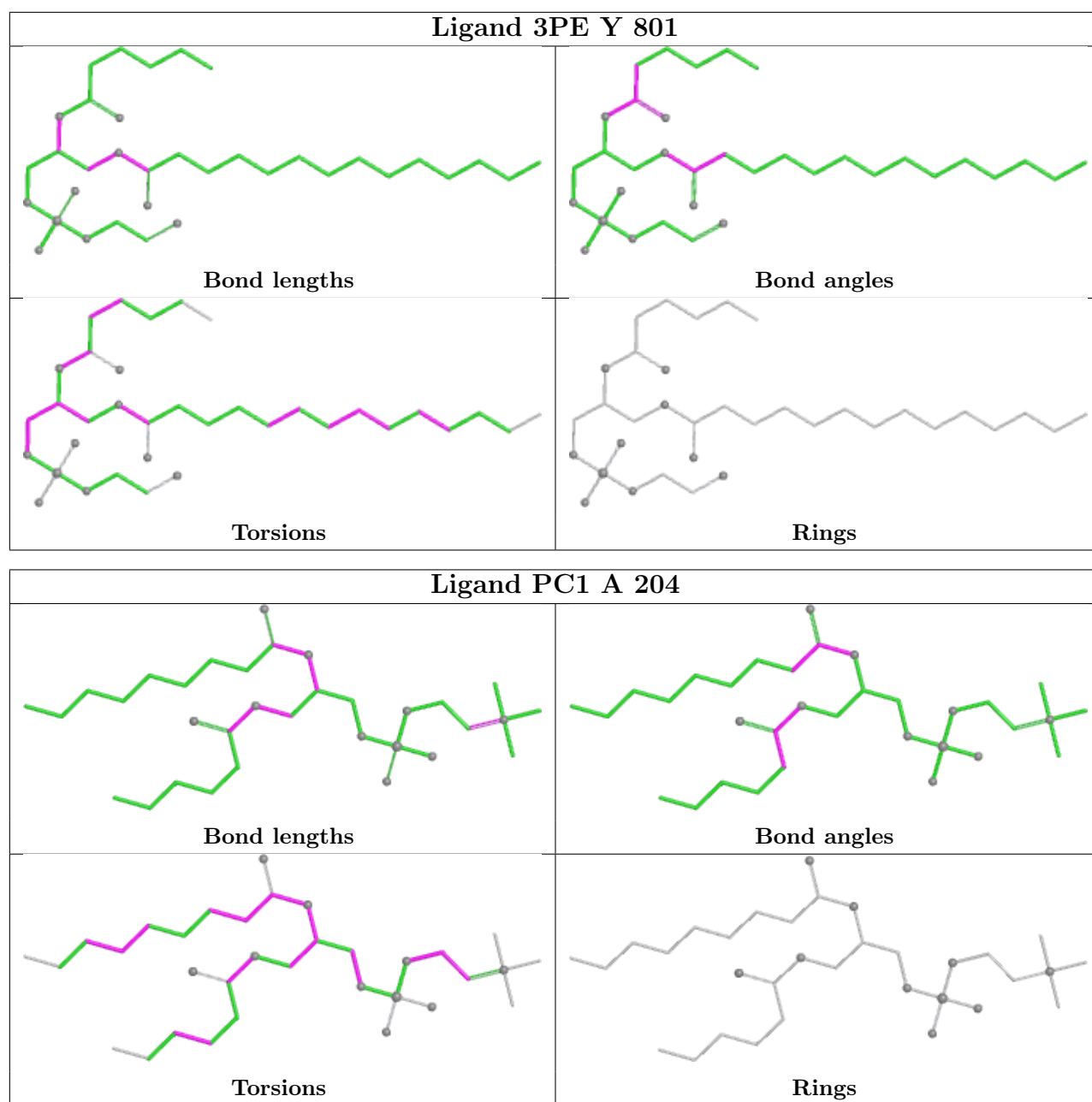
in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

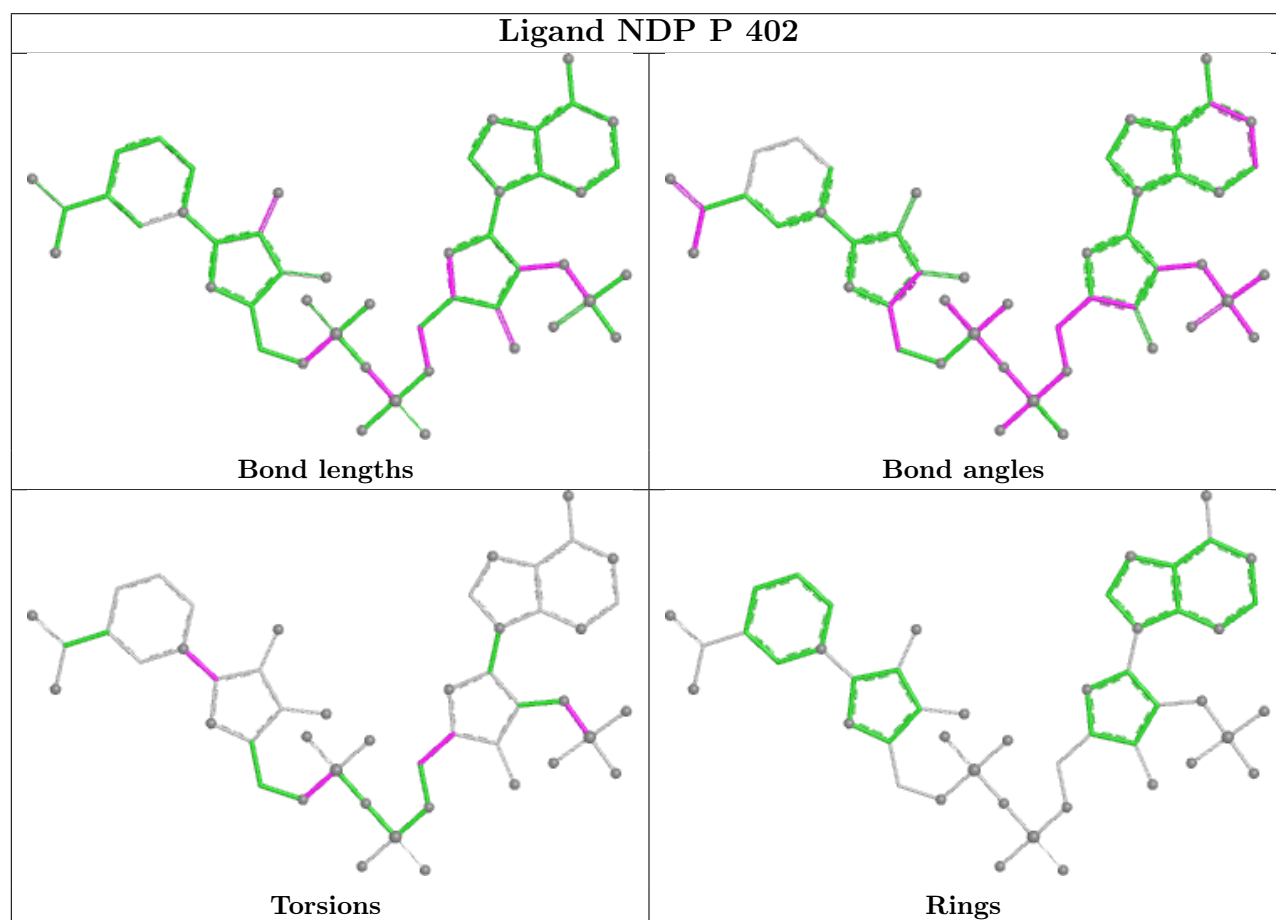
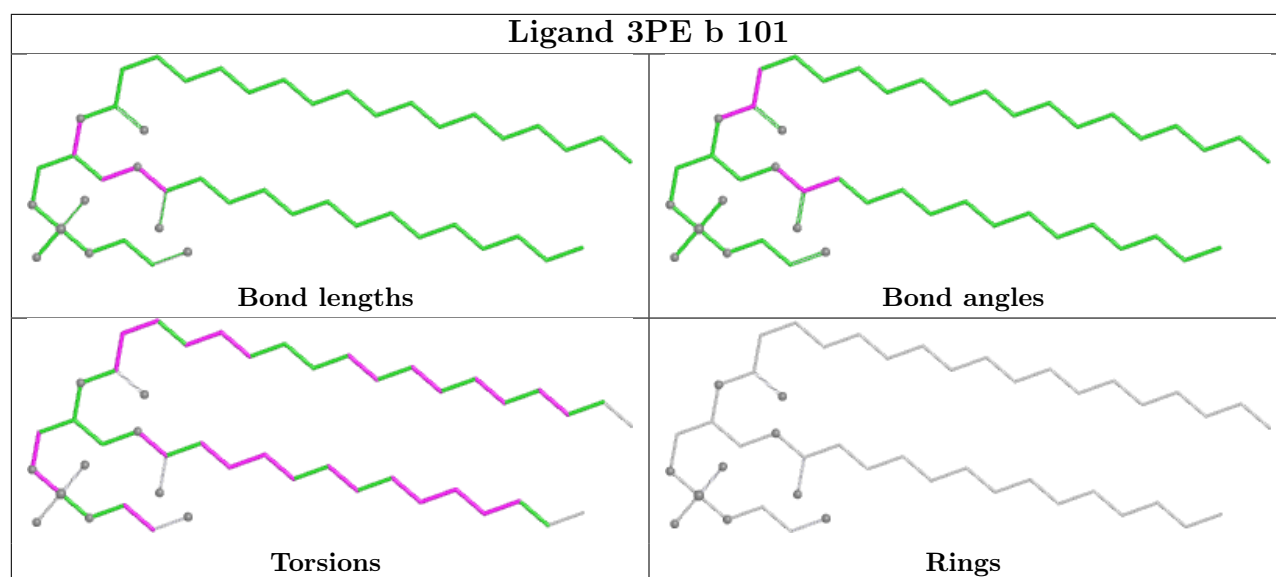


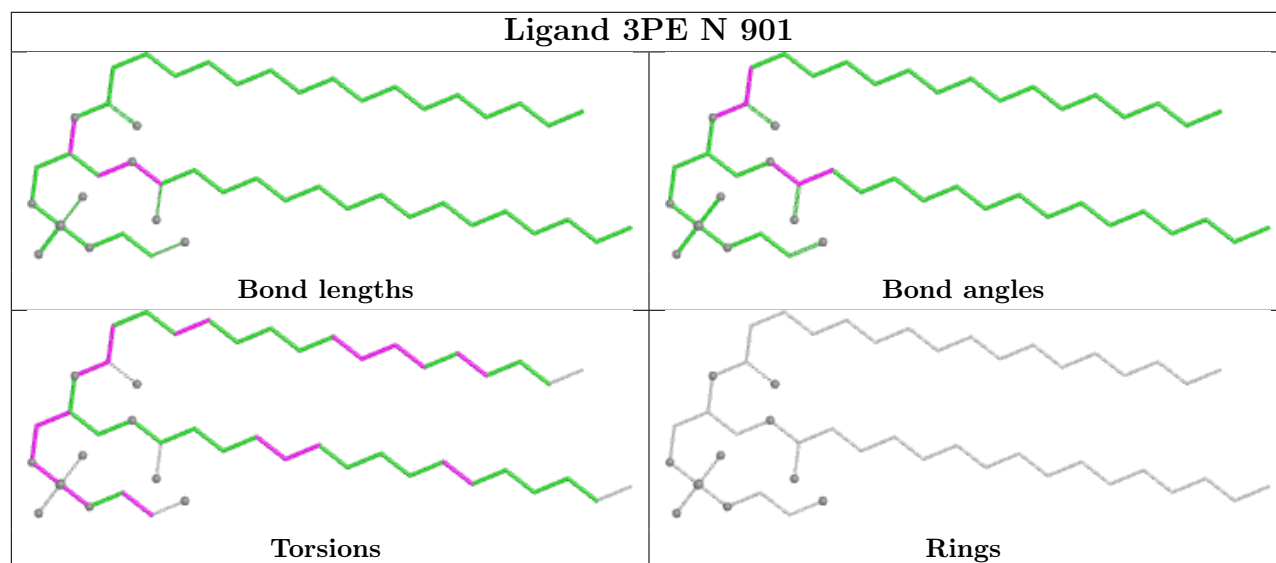
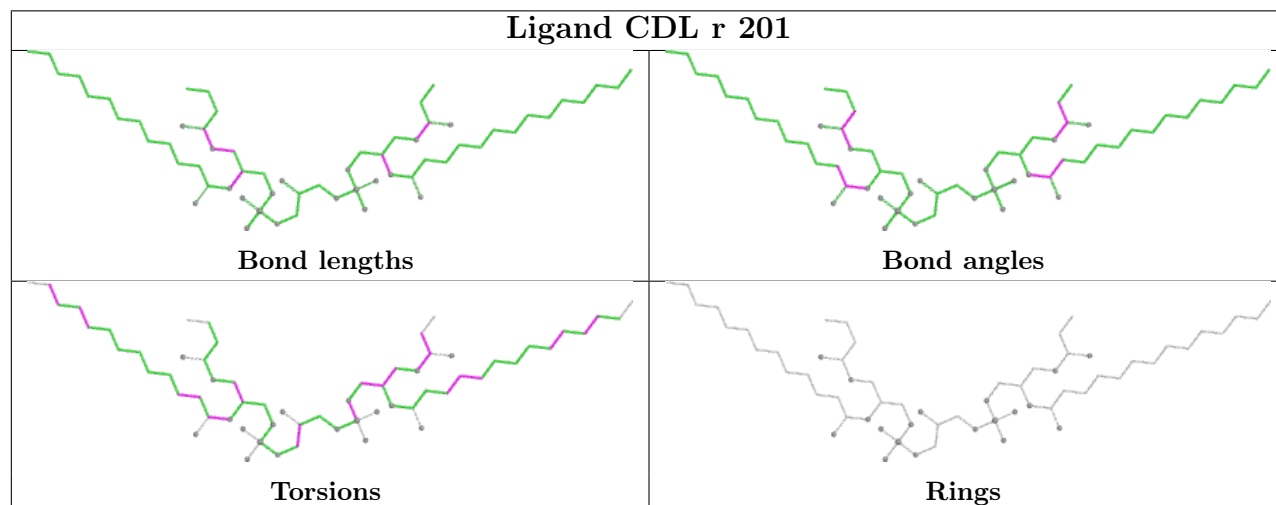
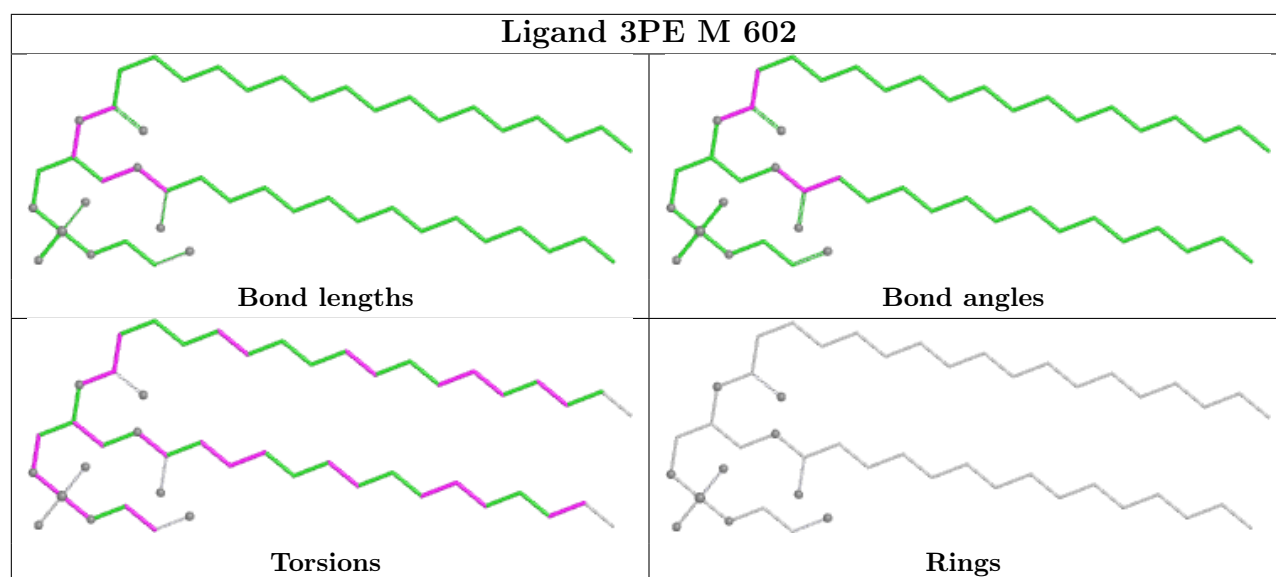


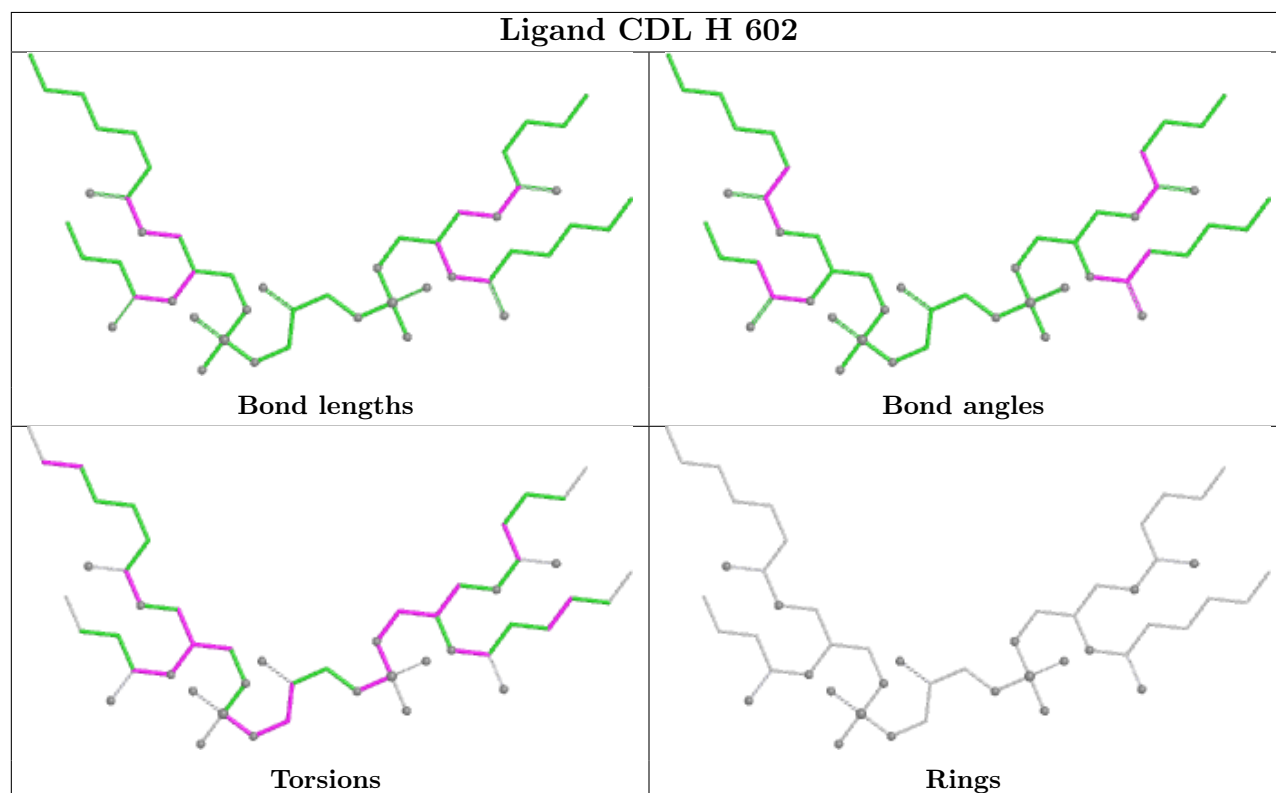
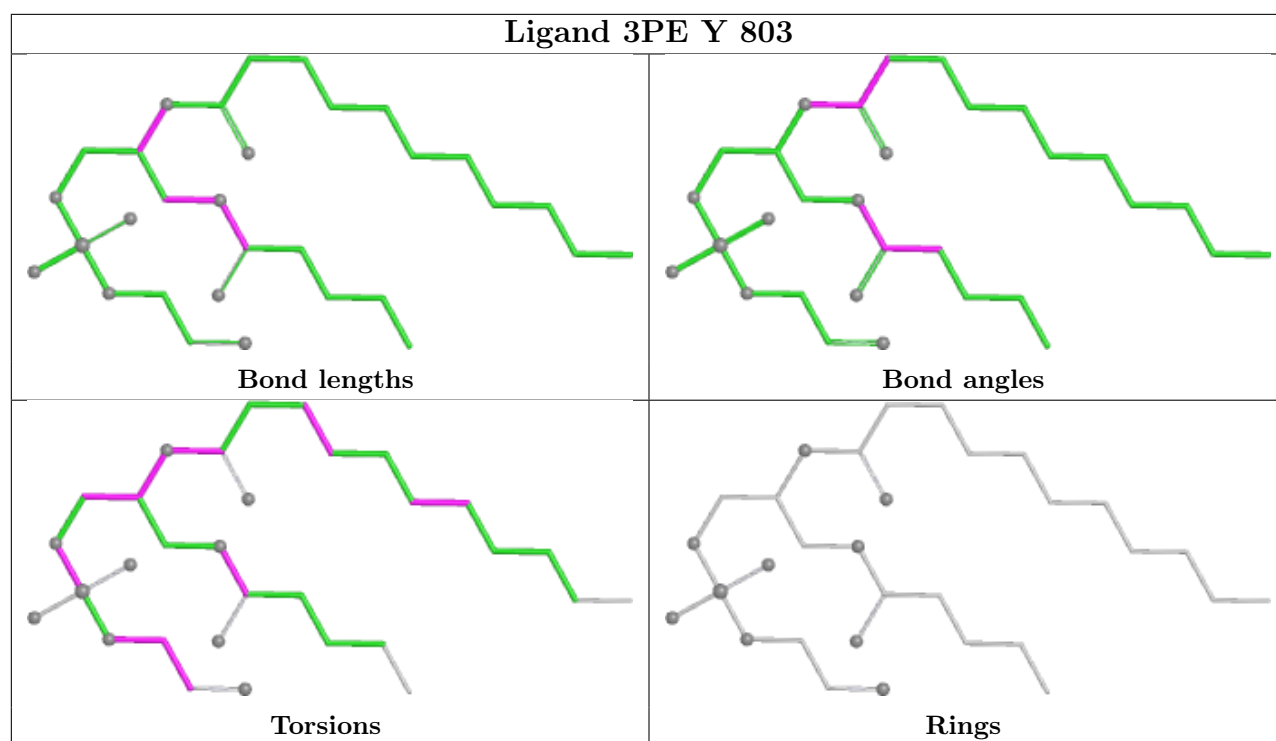


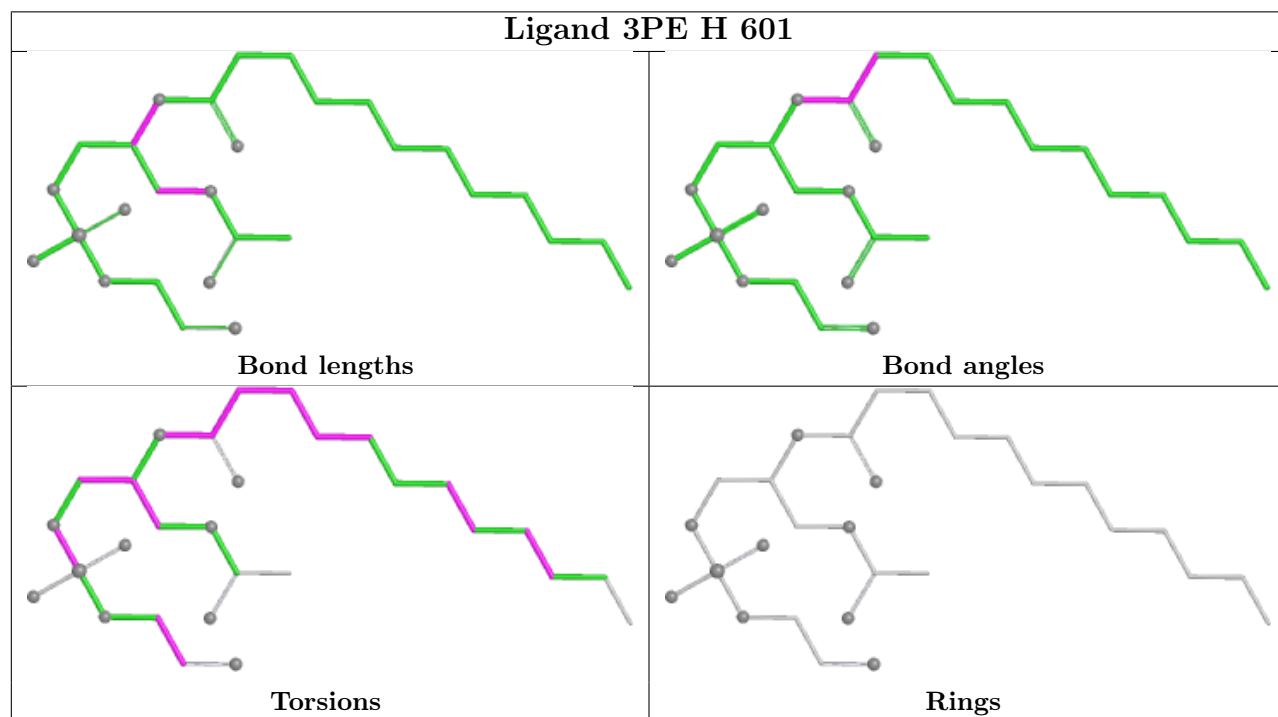
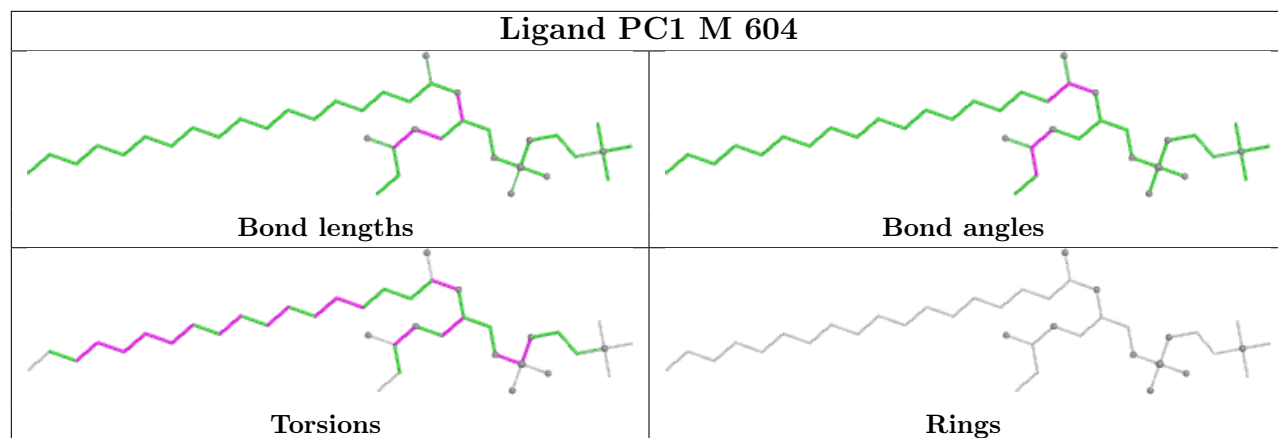


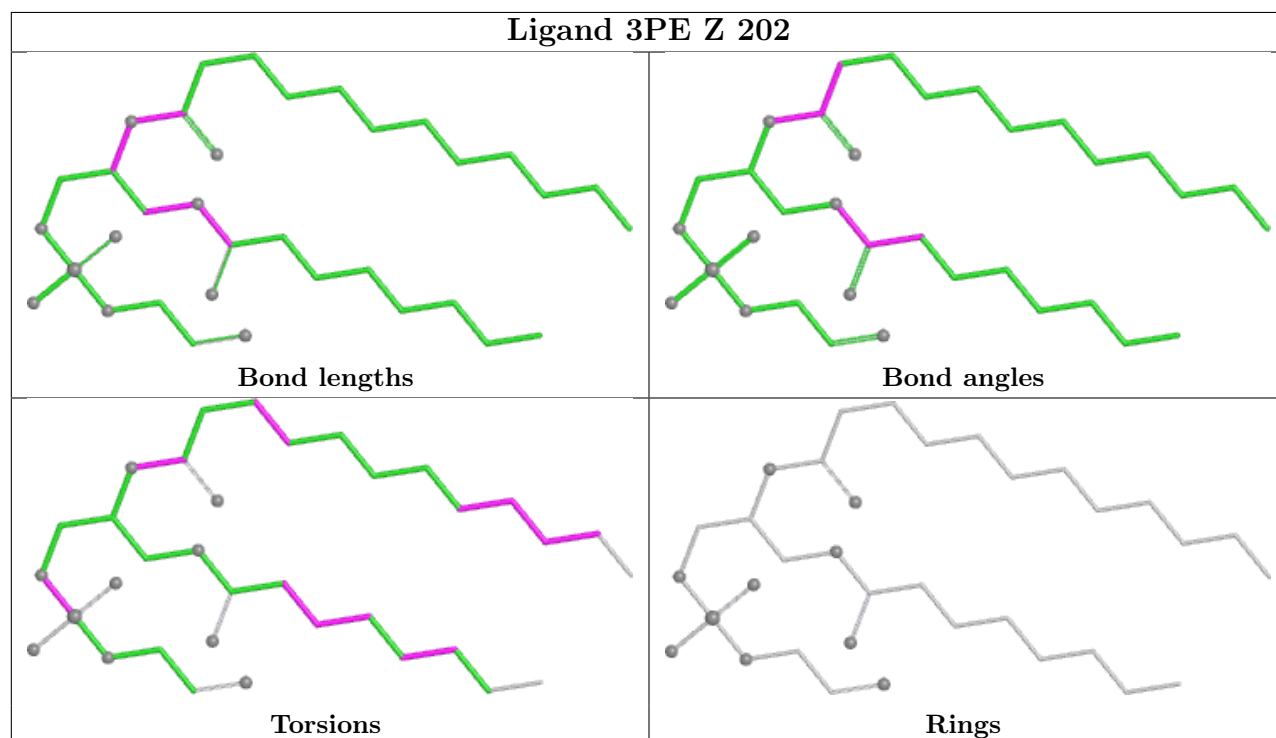
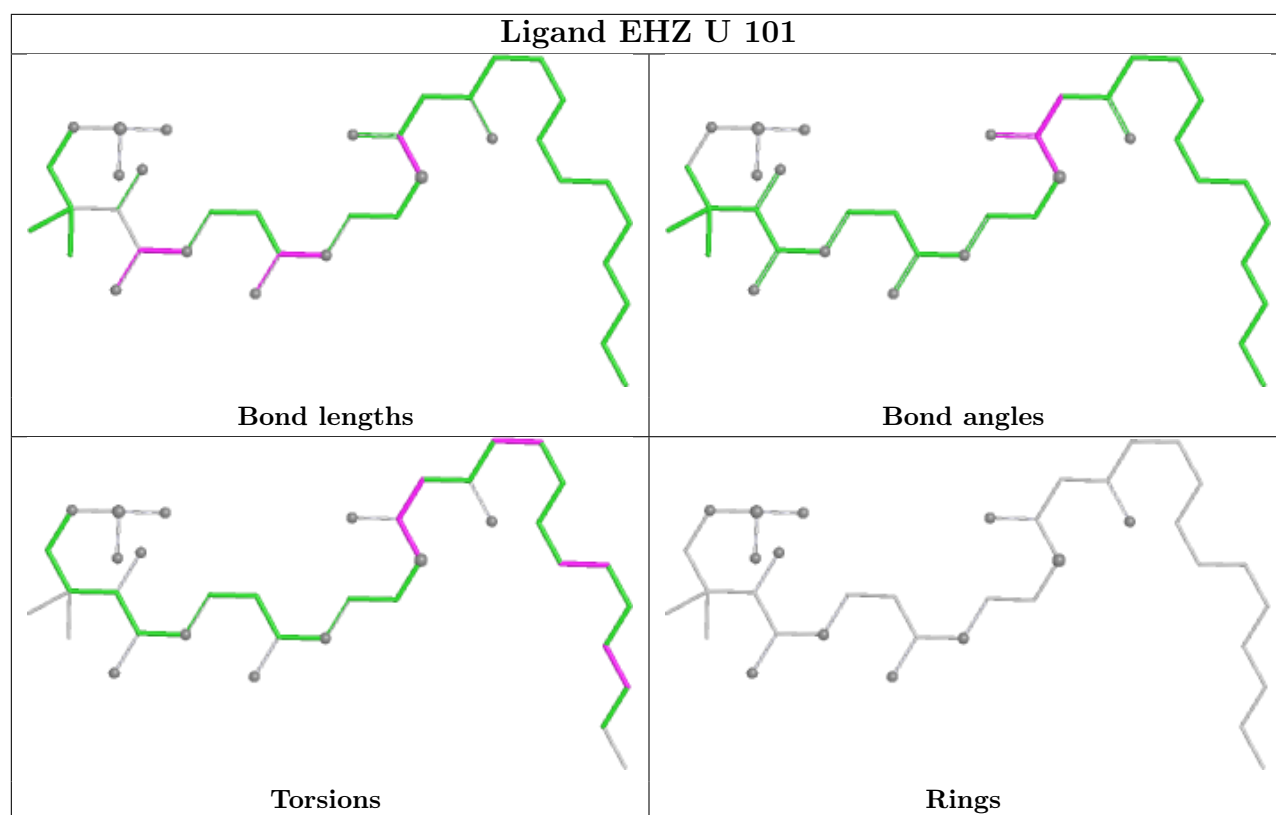


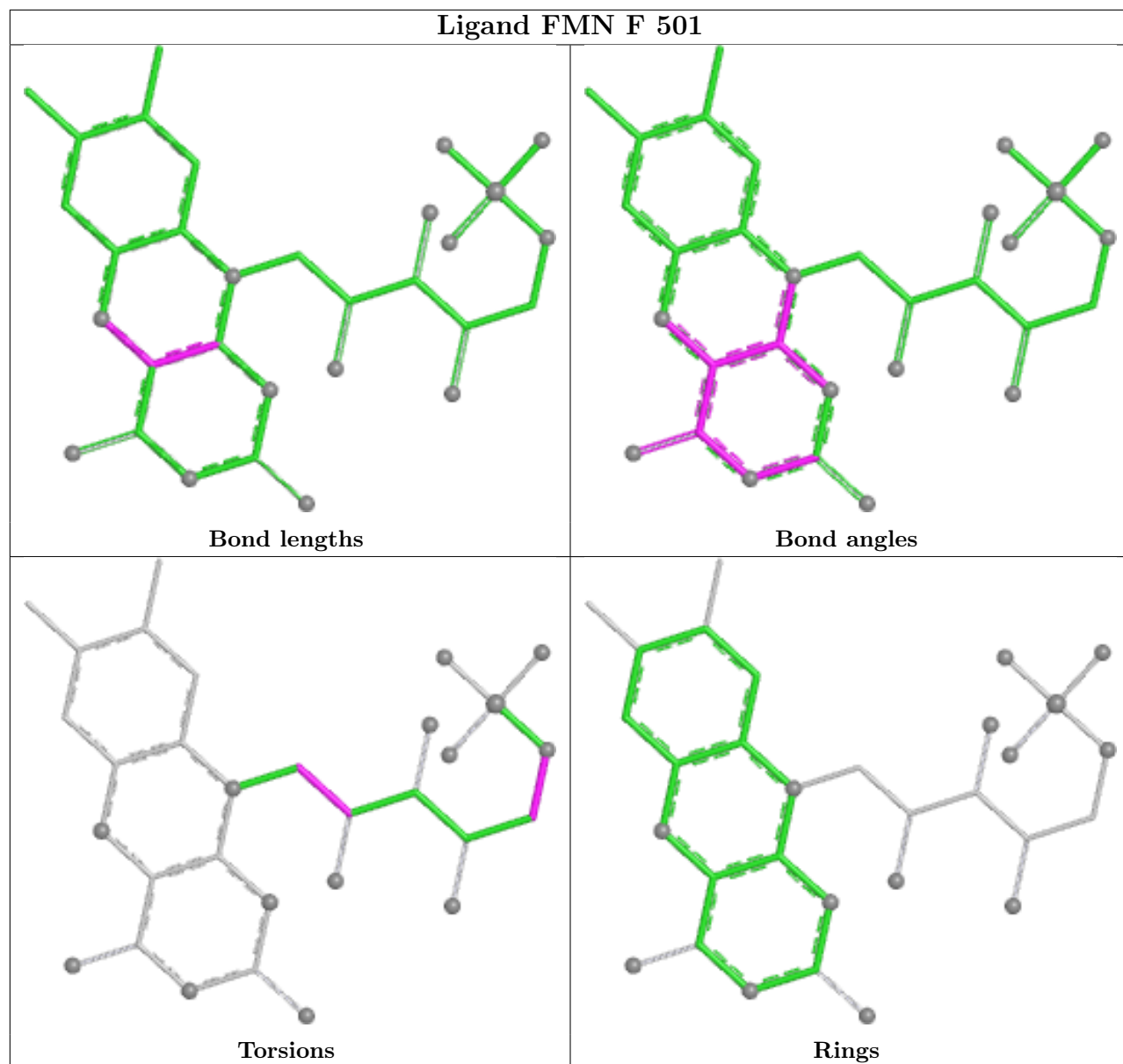


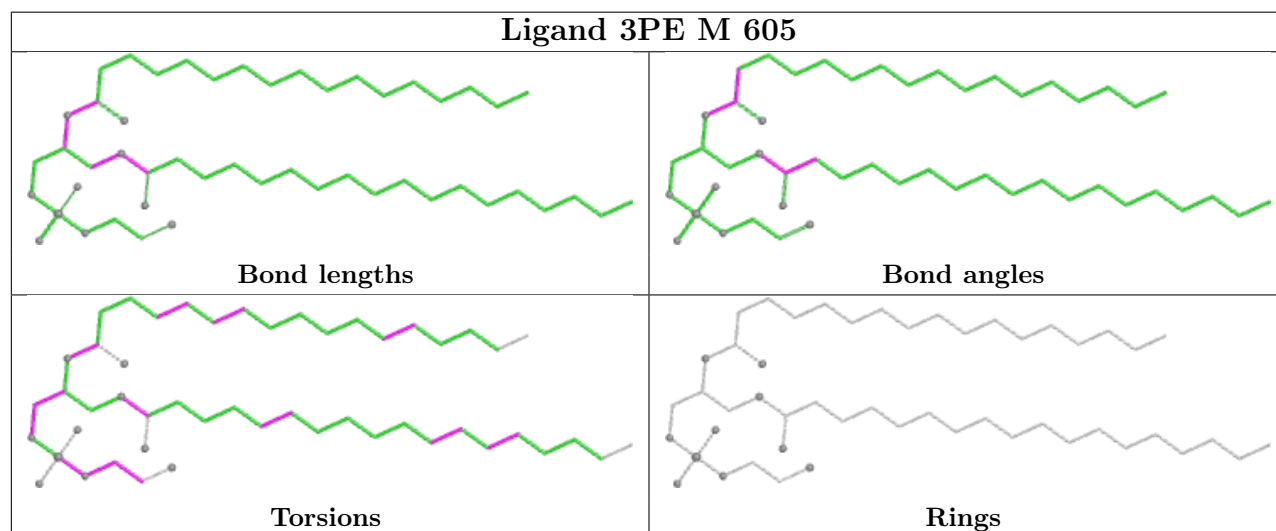
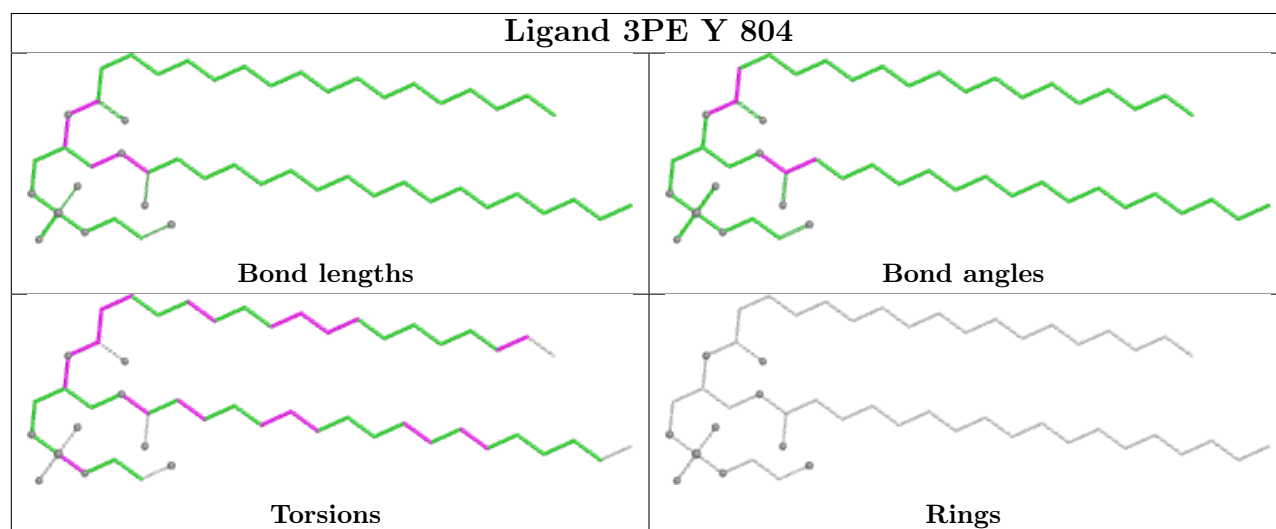
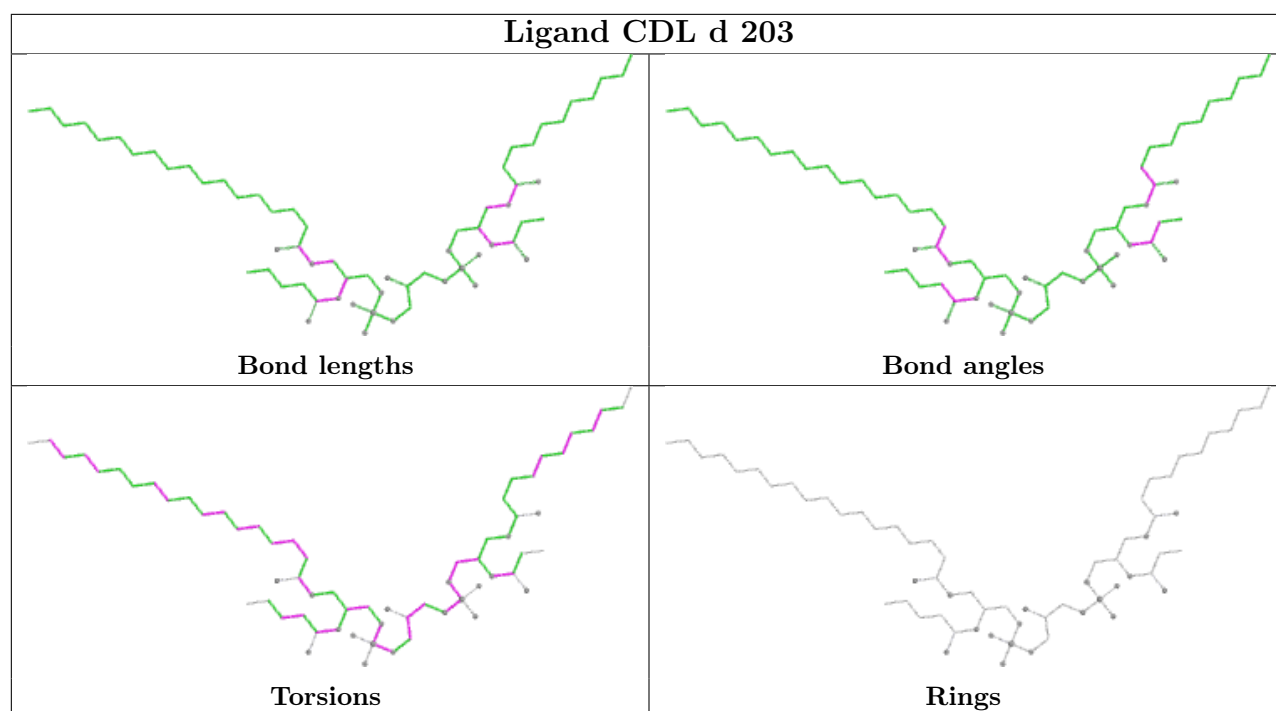


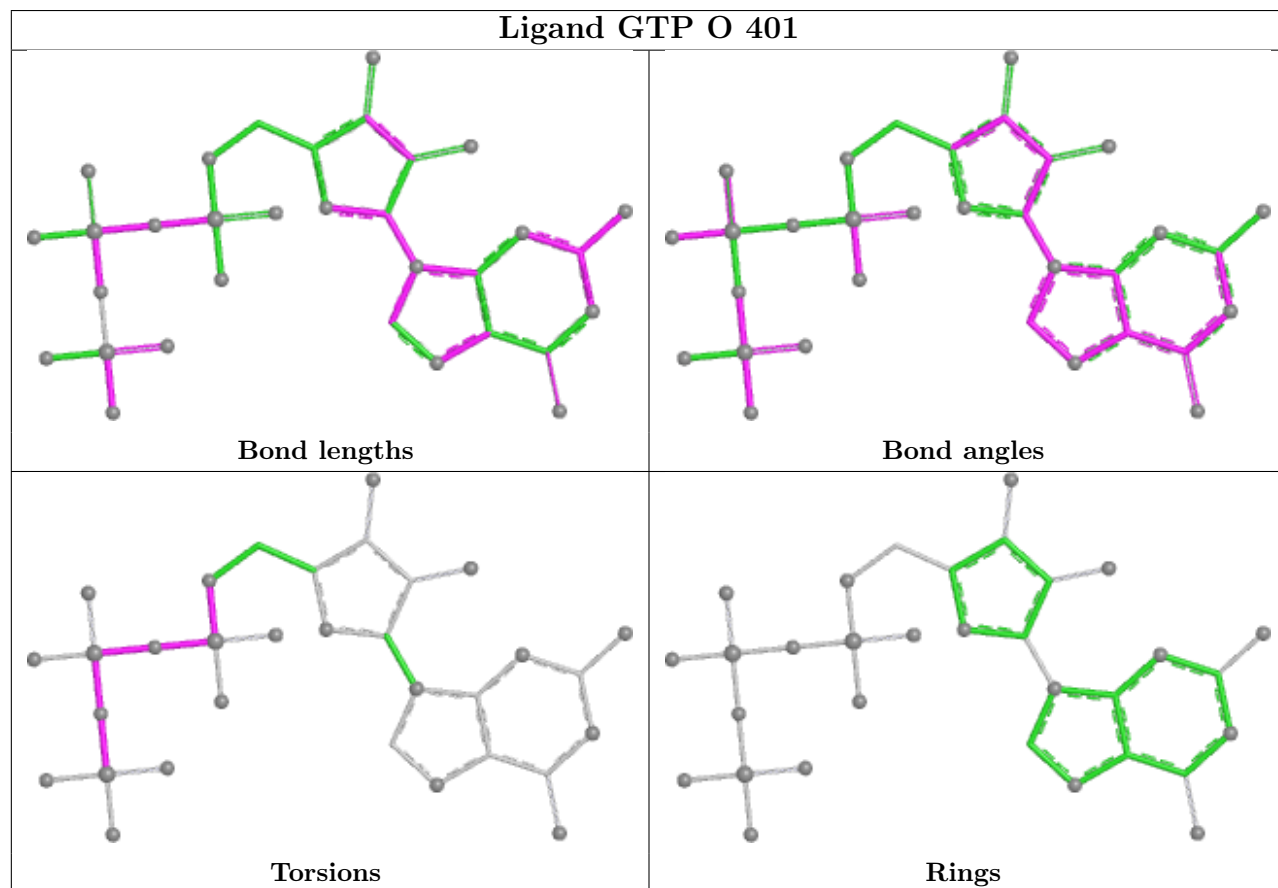
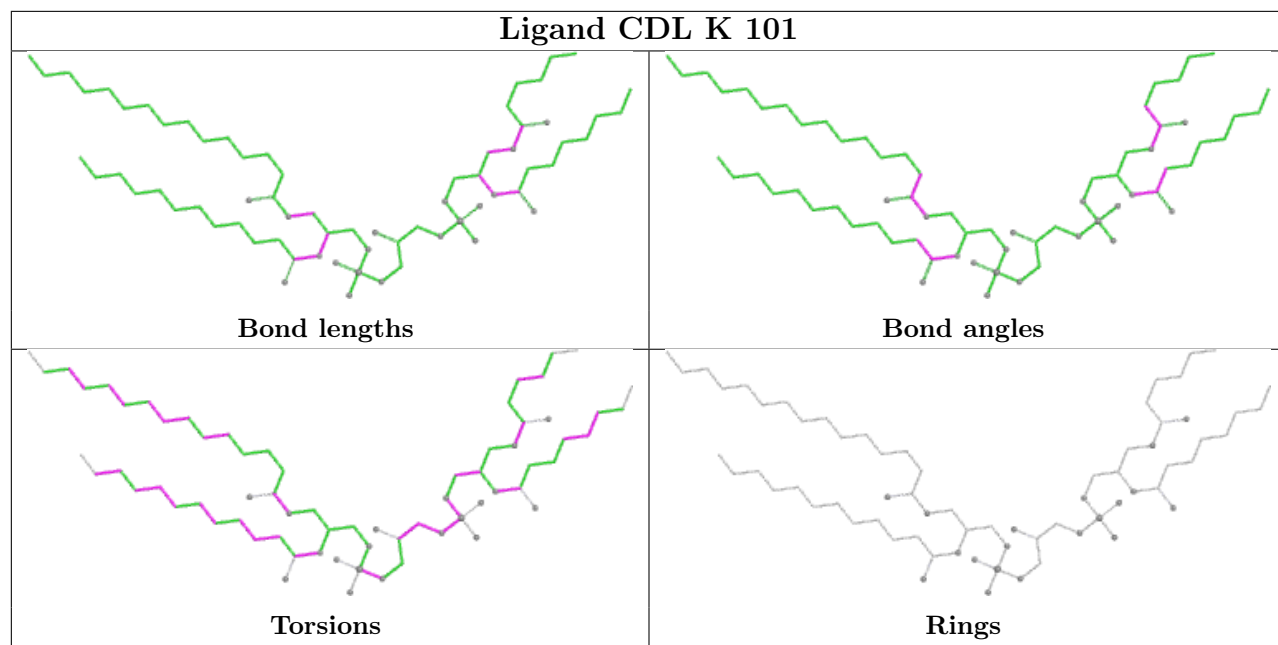


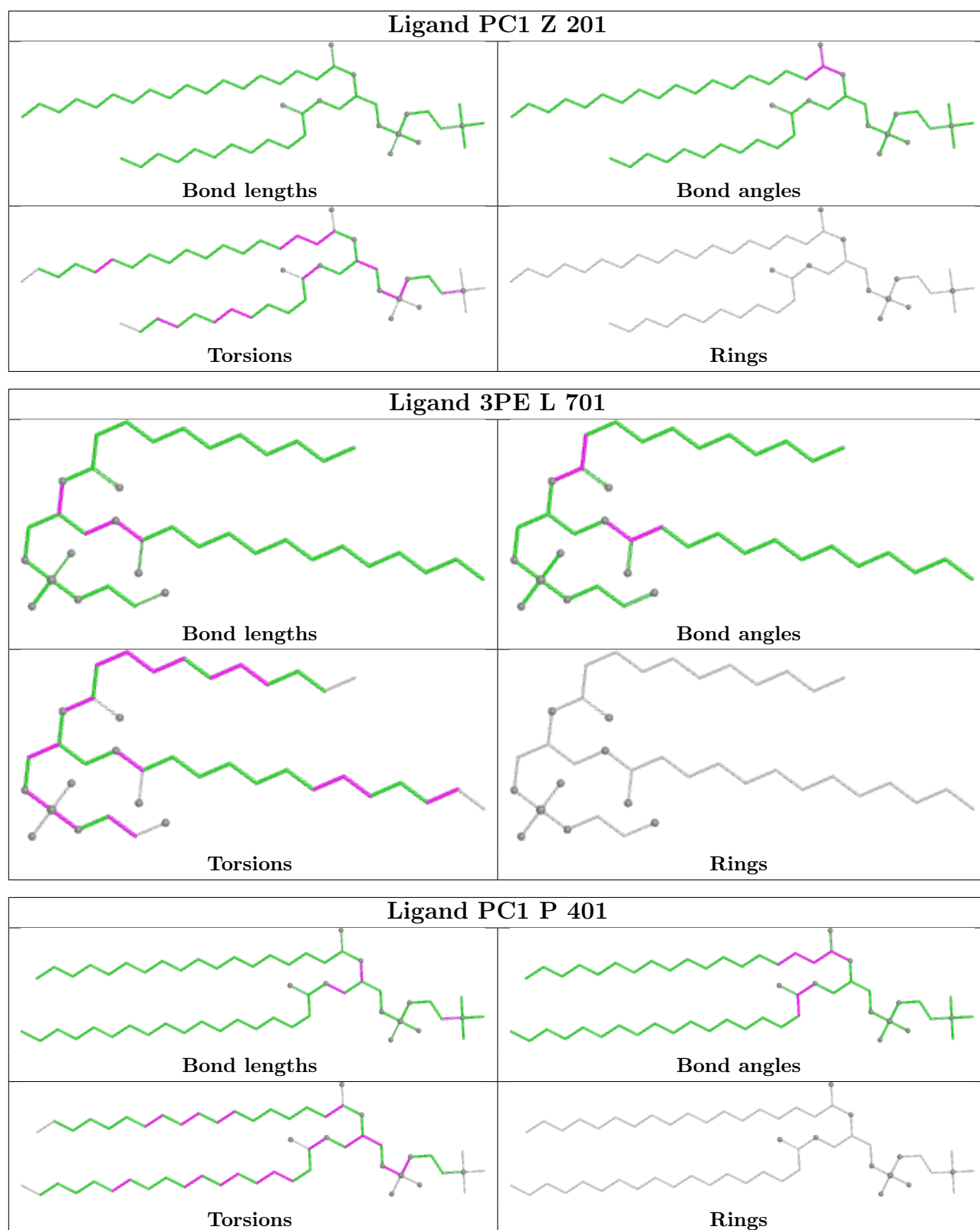


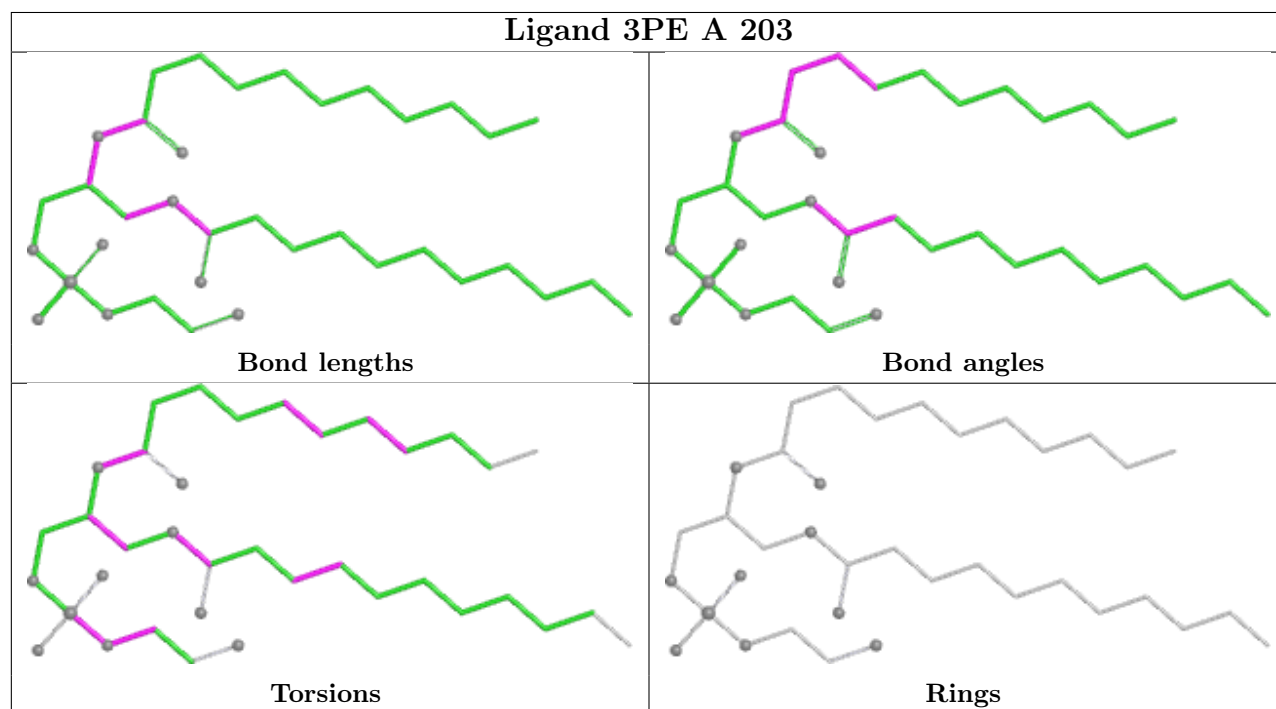
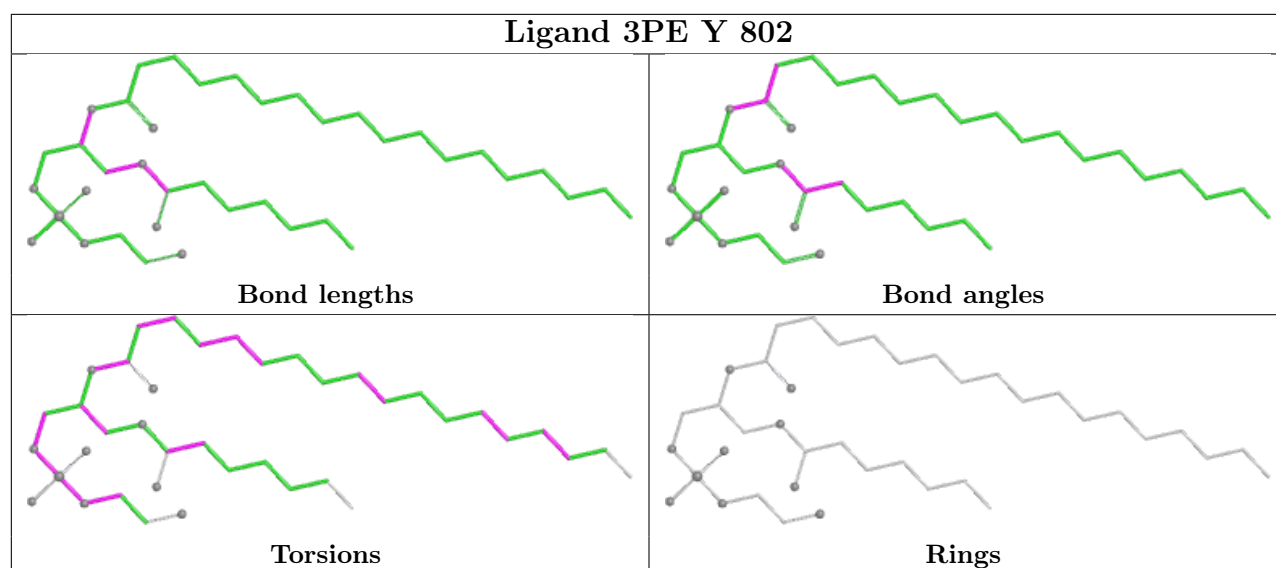


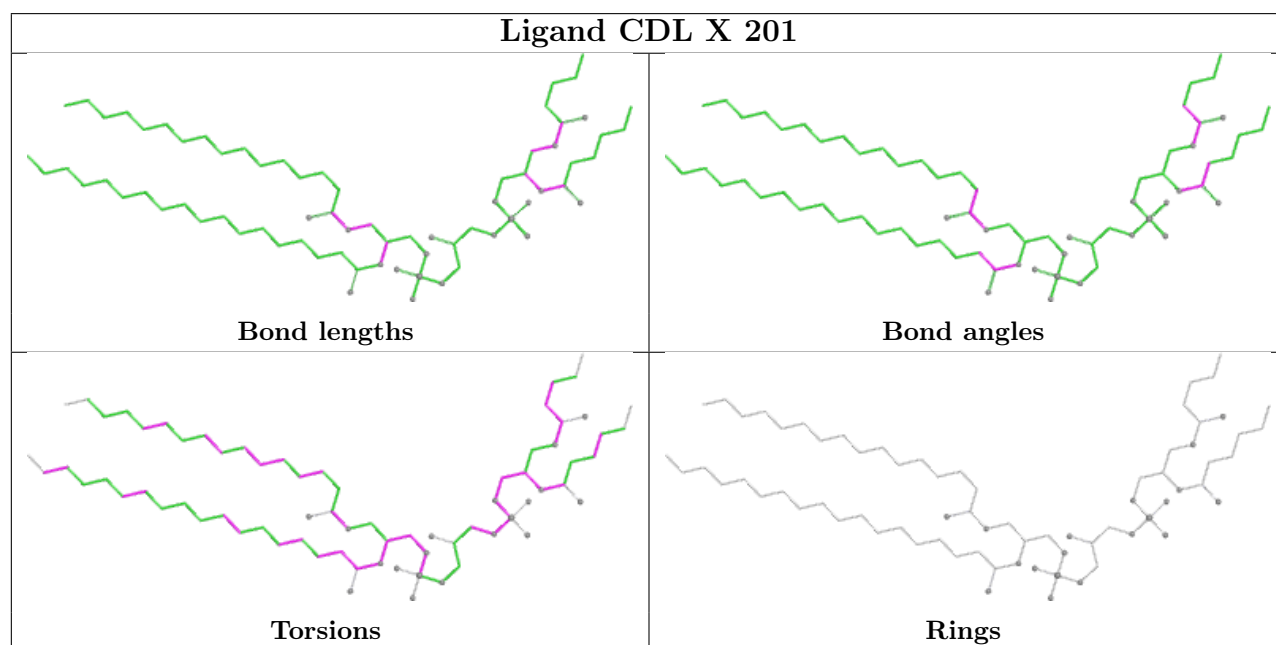
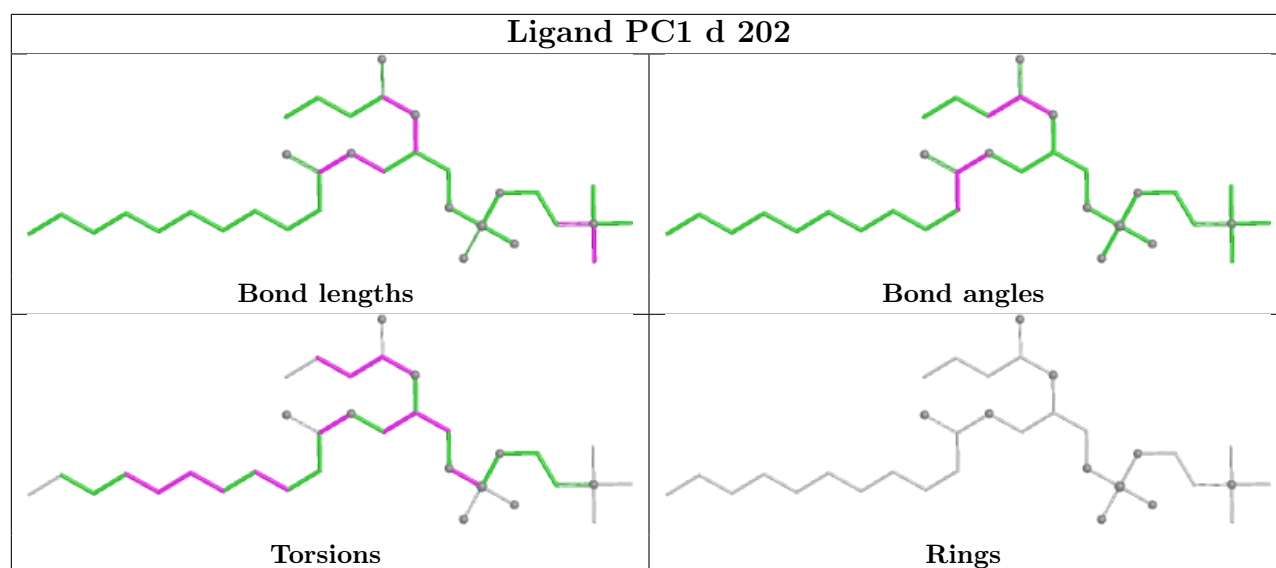


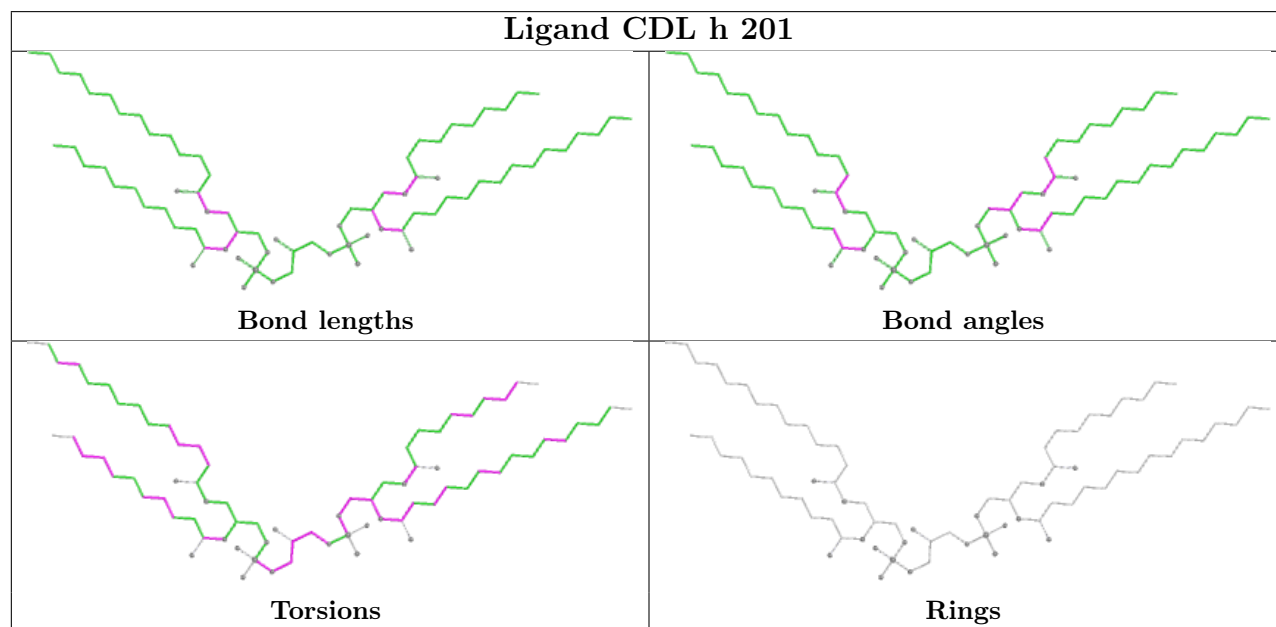
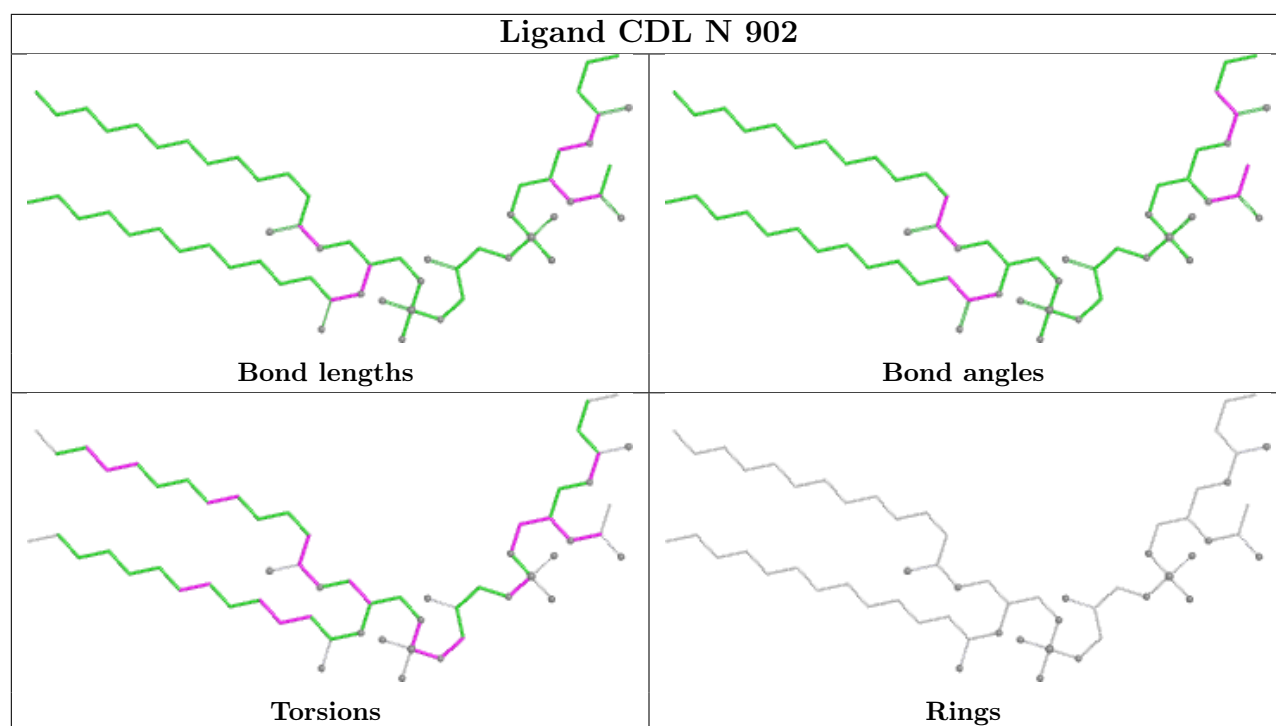


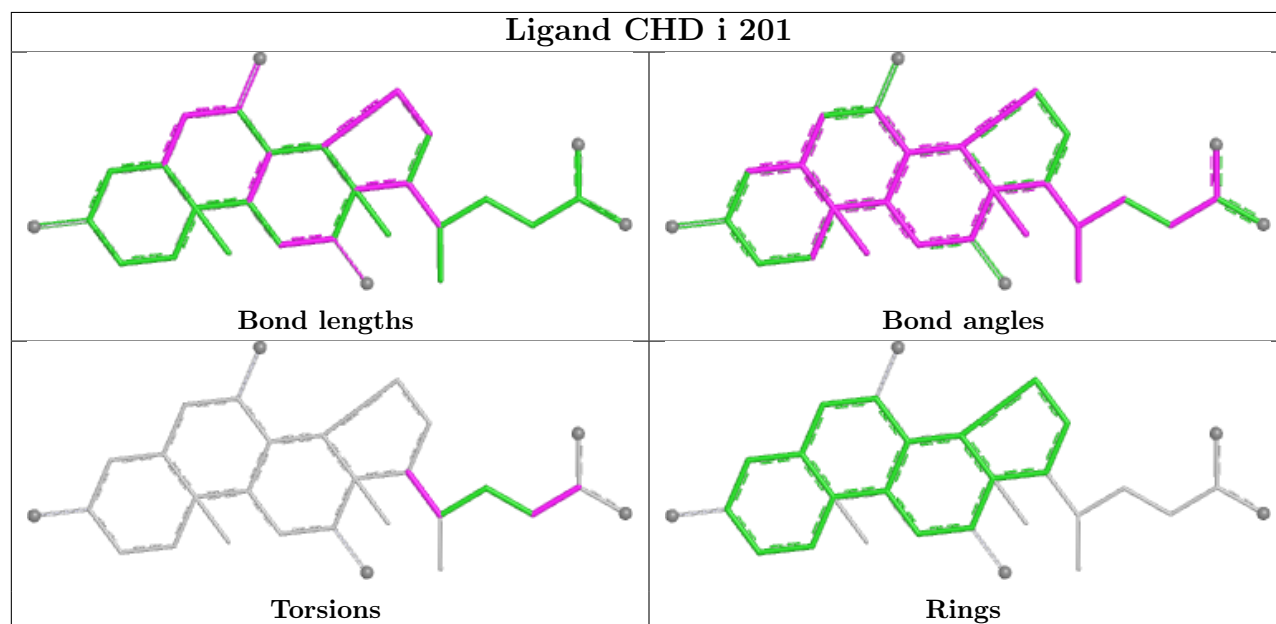
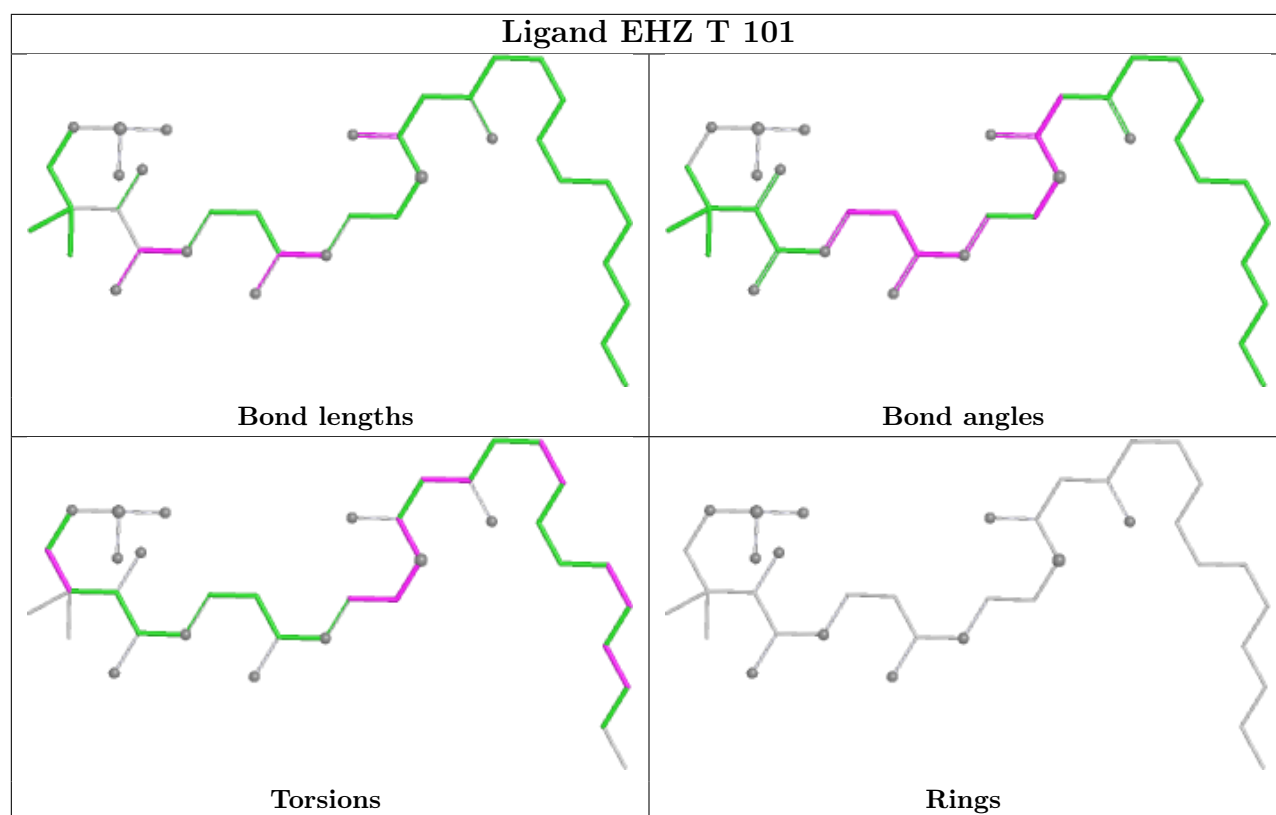


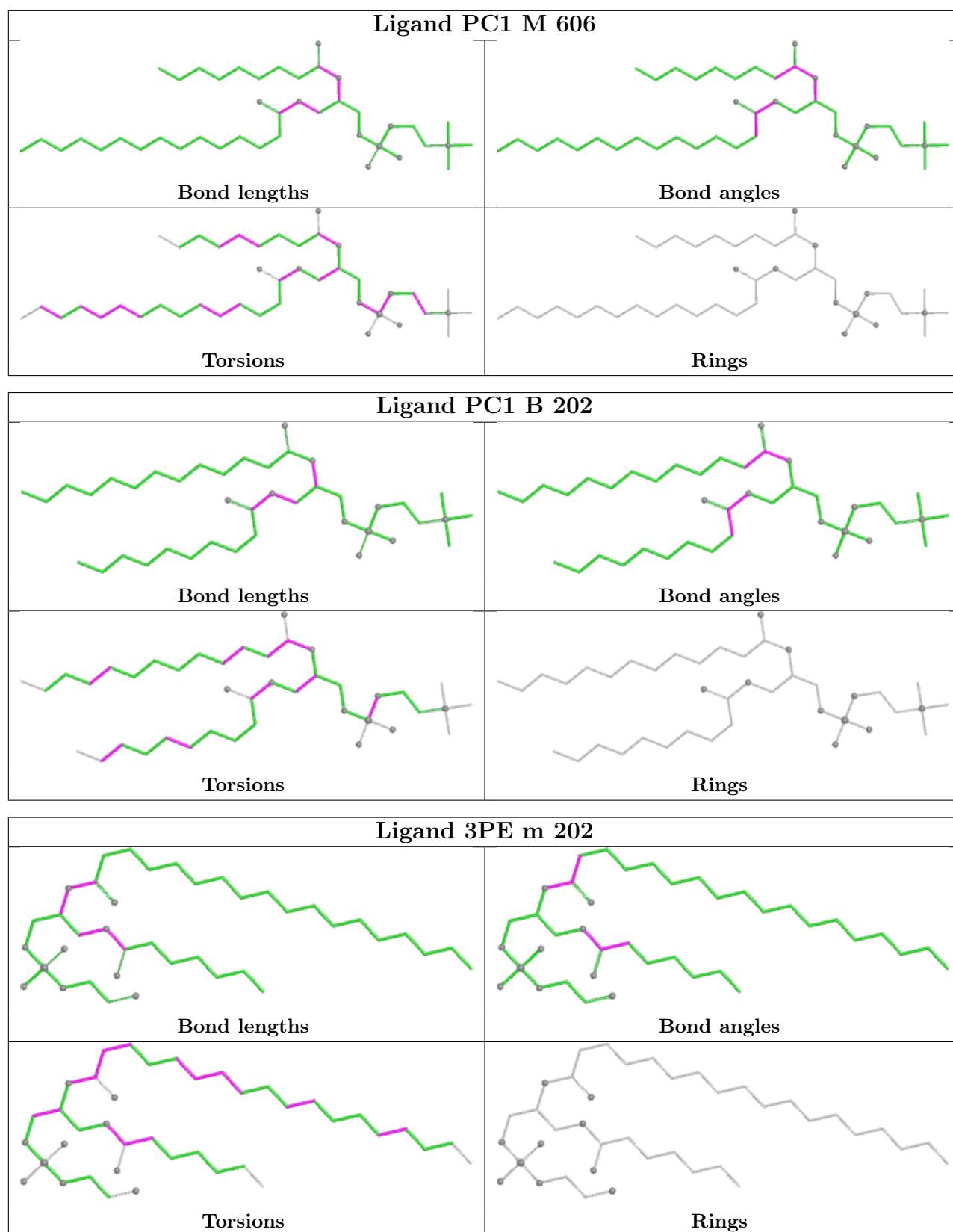


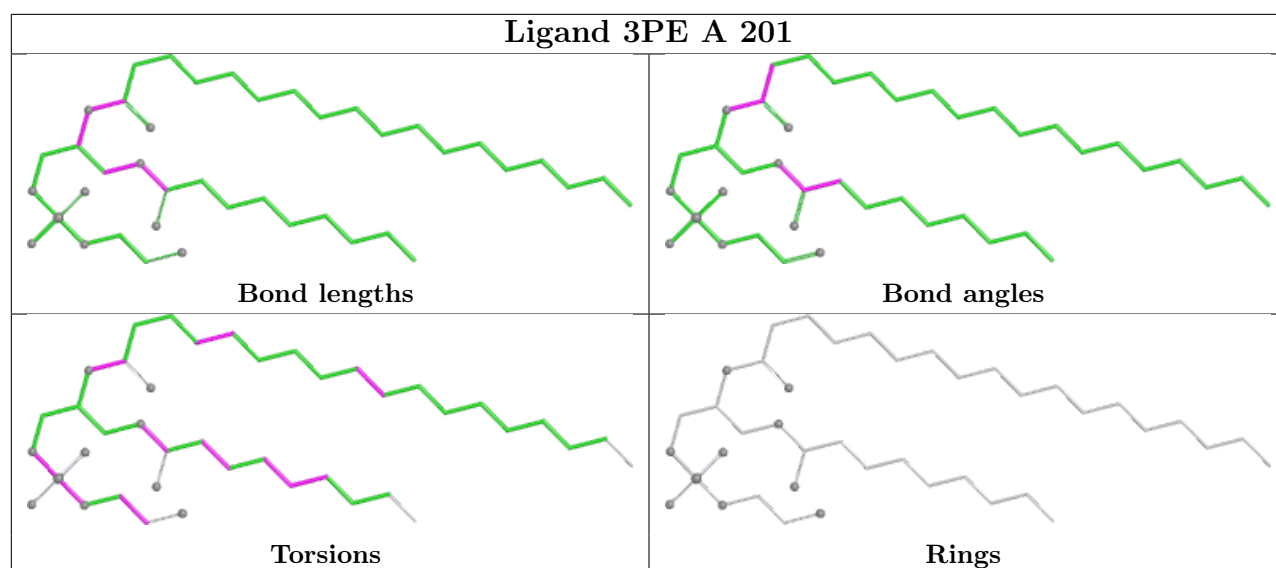












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

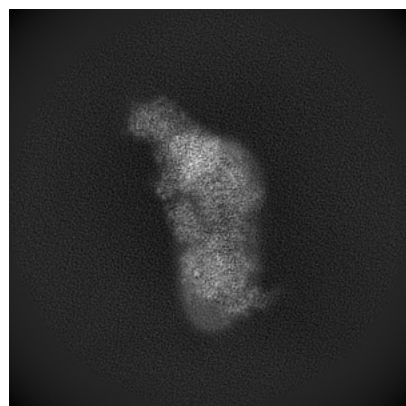
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-14133. These allow visual inspection of the internal detail of the map and identification of artifacts.

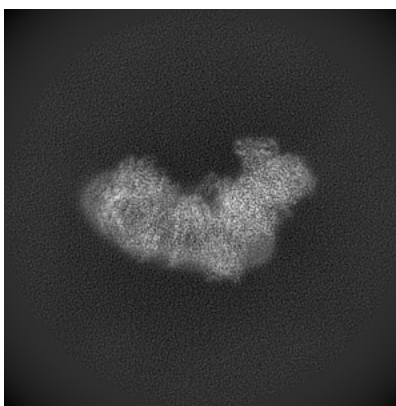
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

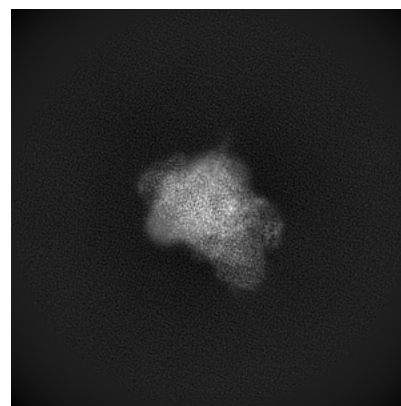
6.1.1 Primary map



X

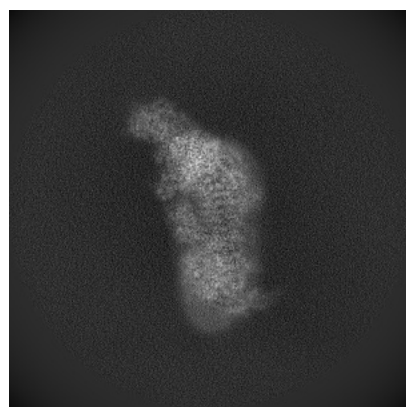


Y

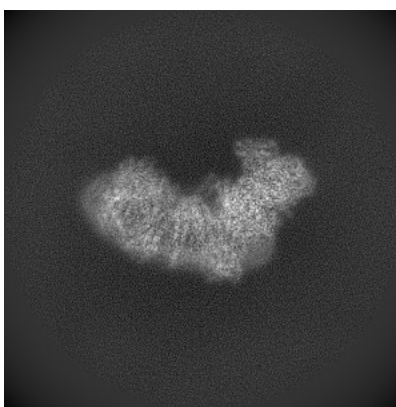


Z

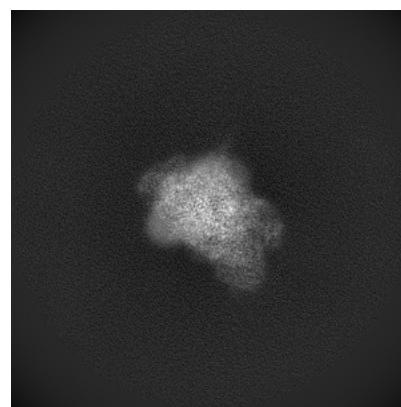
6.1.2 Raw map



X



Y

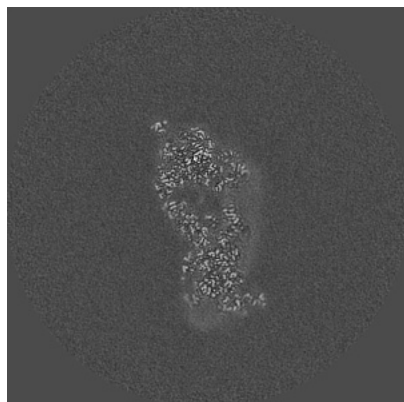


Z

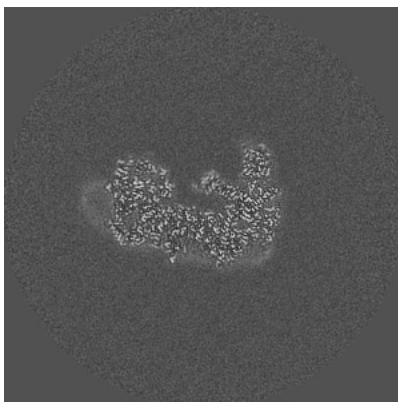
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

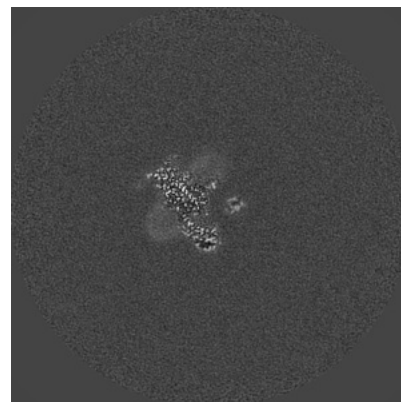
6.2.1 Primary map



X Index: 320

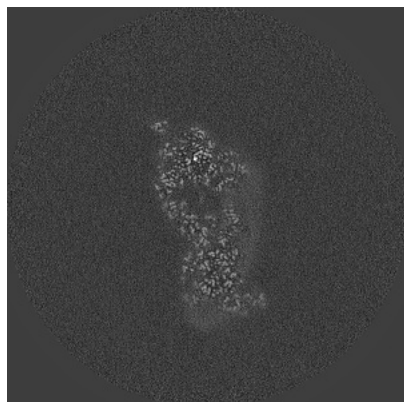


Y Index: 320

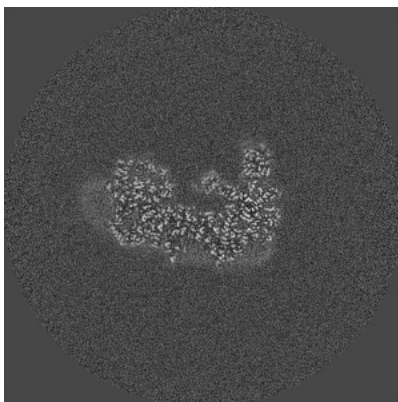


Z Index: 320

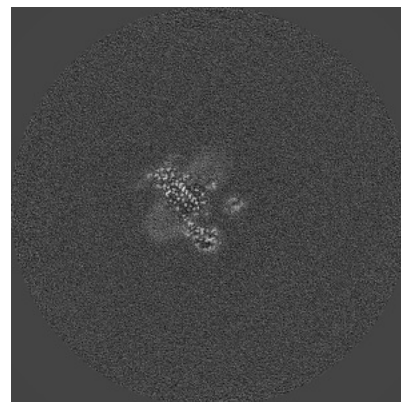
6.2.2 Raw map



X Index: 320



Y Index: 320

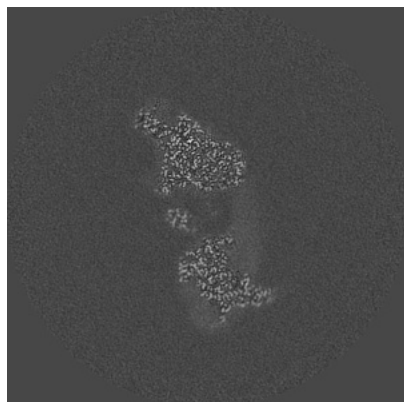


Z Index: 320

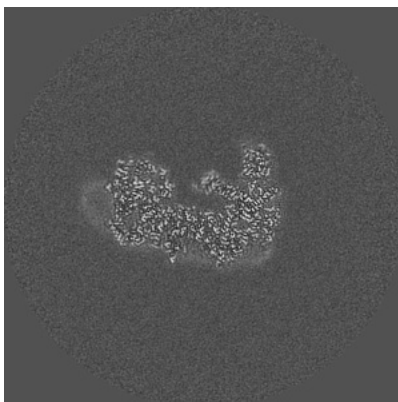
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

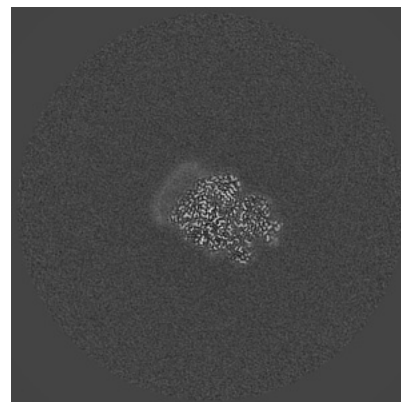
6.3.1 Primary map



X Index: 339

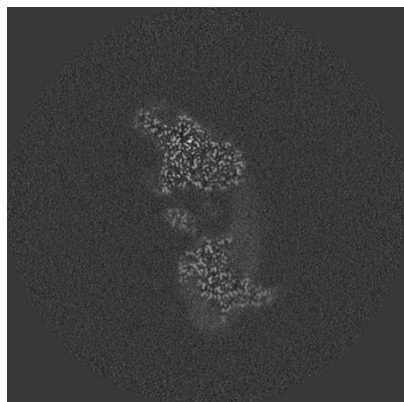


Y Index: 320

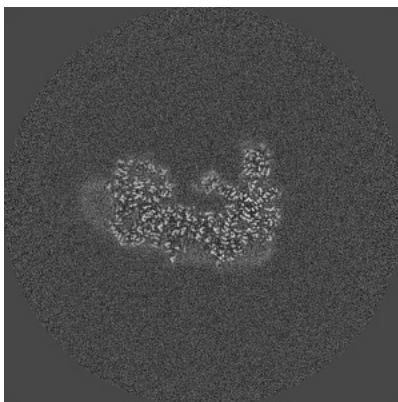


Z Index: 405

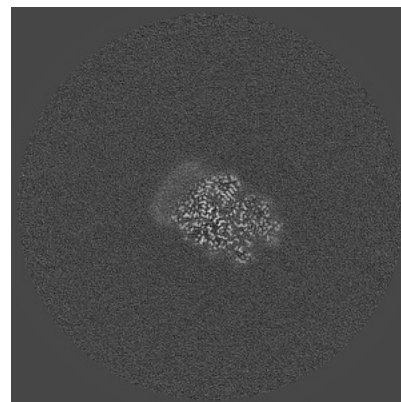
6.3.2 Raw map



X Index: 339



Y Index: 320

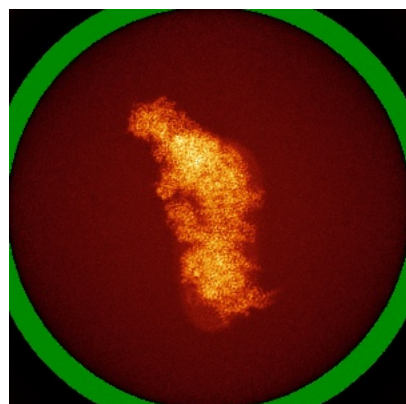


Z Index: 405

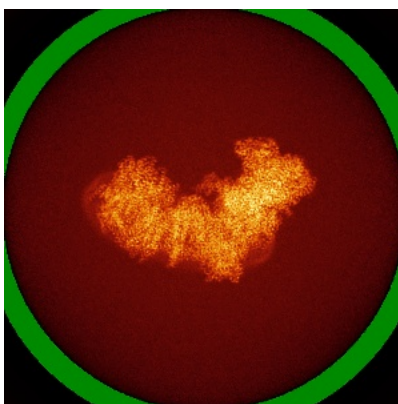
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) [i](#)

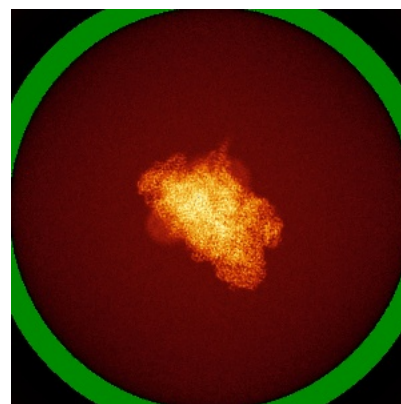
6.4.1 Primary map



X

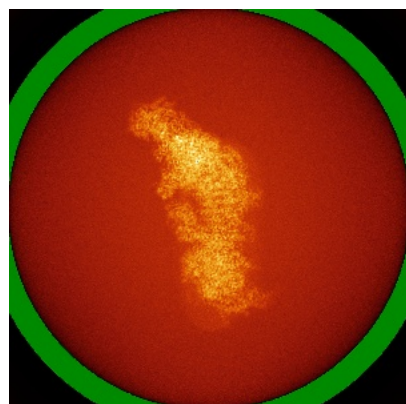


Y

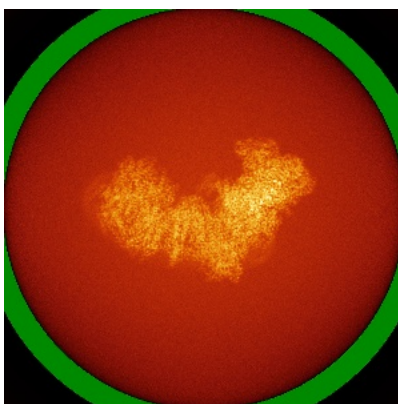


Z

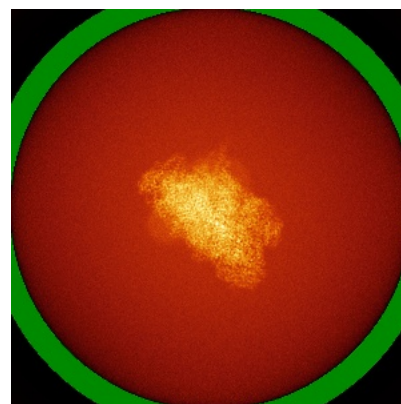
6.4.2 Raw map



X



Y



Z

The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



X



Y



Z

The images above show the 3D surface view of the map at the recommended contour level 6.5. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



X



Y



Z

These images show the 3D surface of the raw map. The raw map's contour level was selected so that its surface encloses the same volume as the primary map does at its recommended contour level.

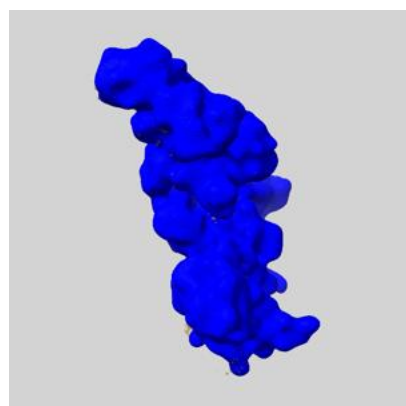
6.6 Mask visualisation [i](#)

This section shows the 3D surface view of the primary map at 50% transparency overlaid with the specified mask at 0% transparency

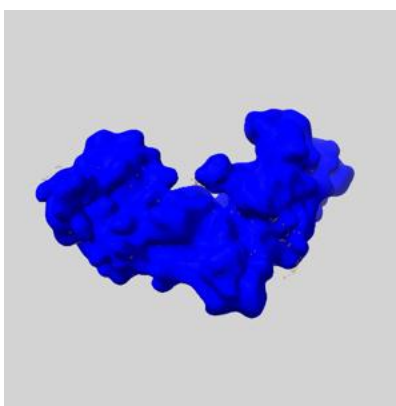
A mask typically either:

- Encompasses the whole structure
- Separates out a domain, a functional unit, a monomer or an area of interest from a larger structure

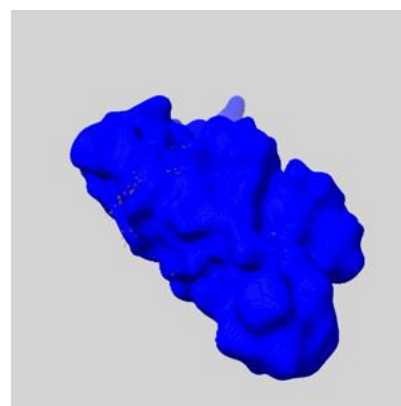
6.6.1 emd_14133_msk_1.map [i](#)



X



Y

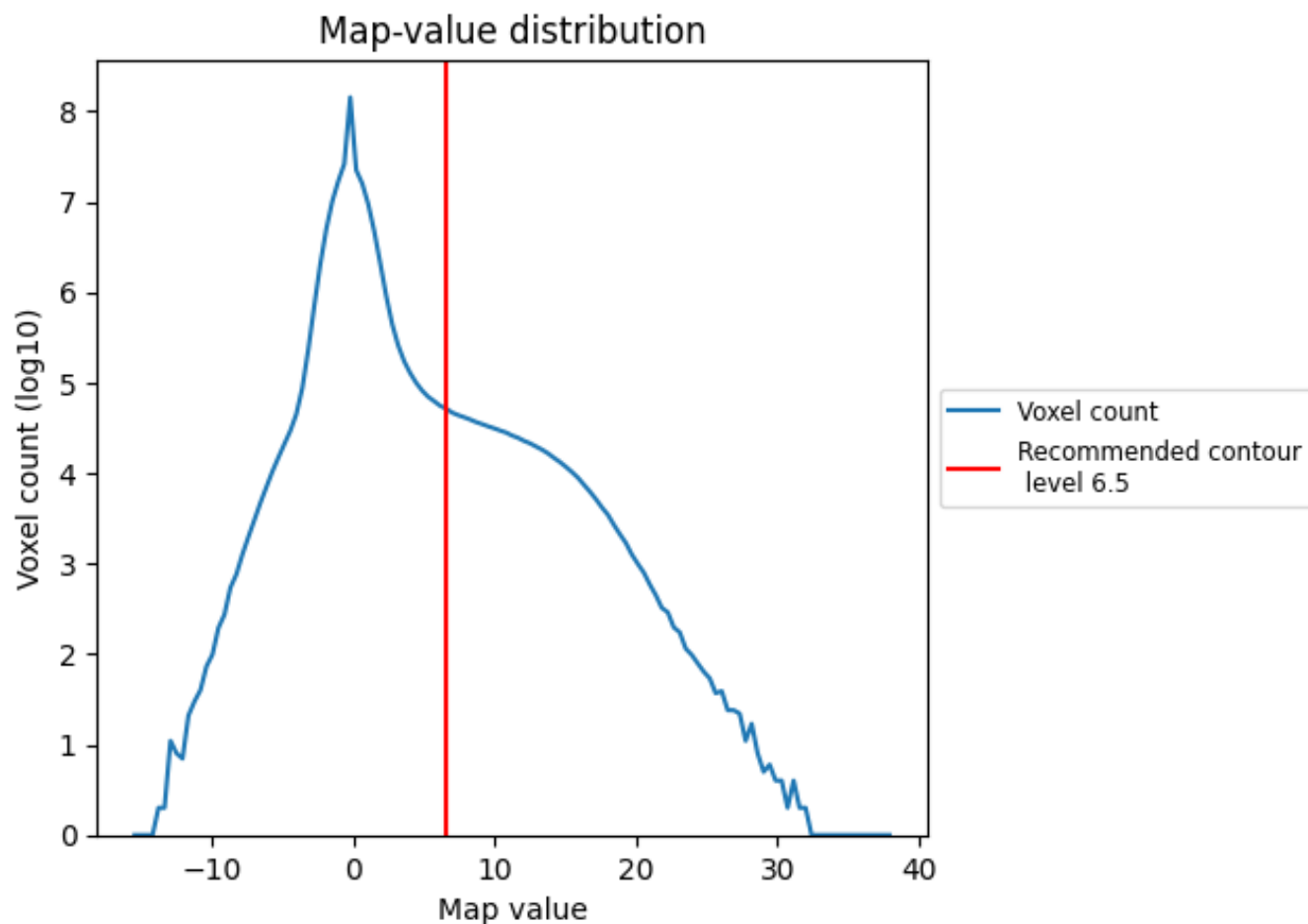


Z

7 Map analysis [i](#)

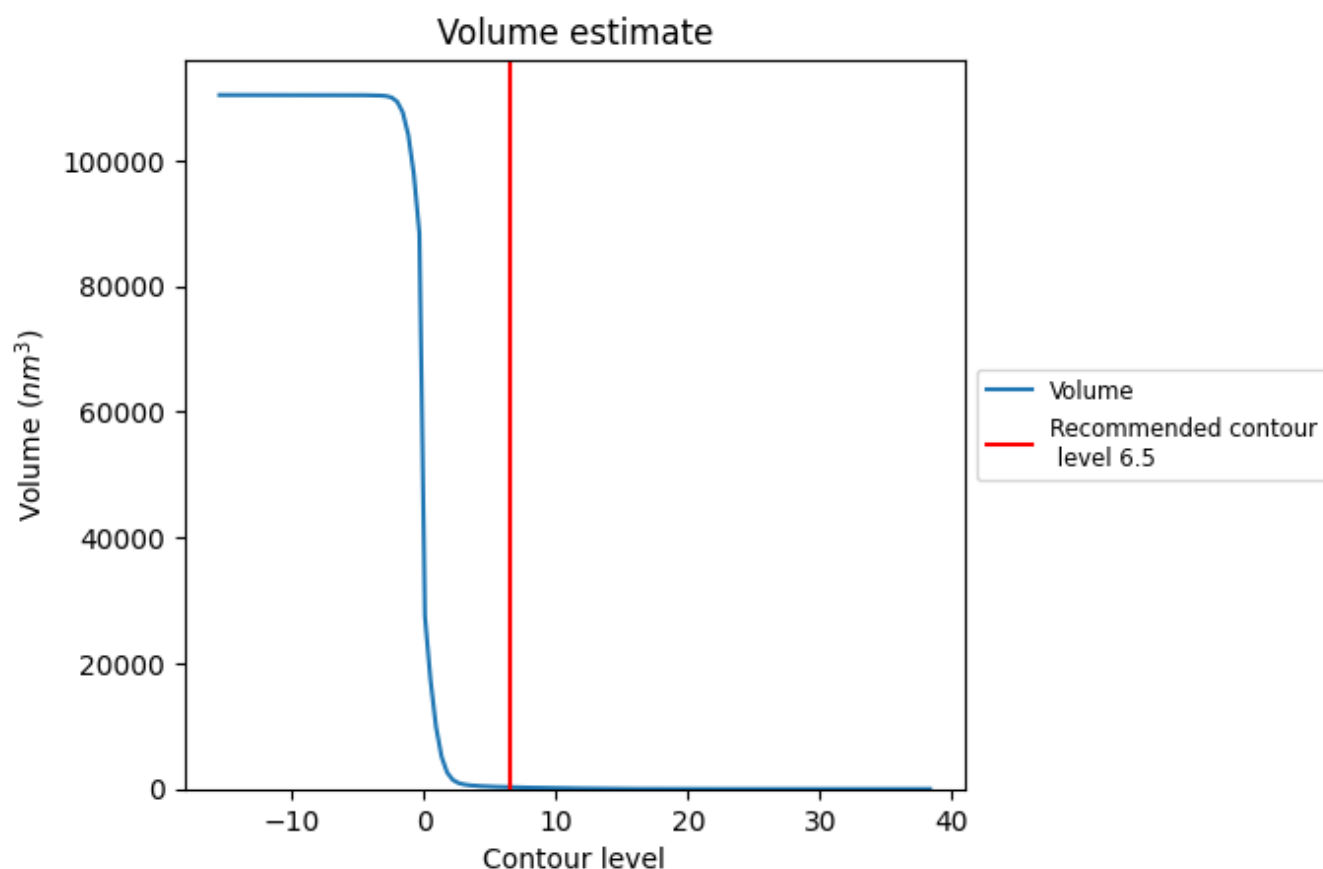
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

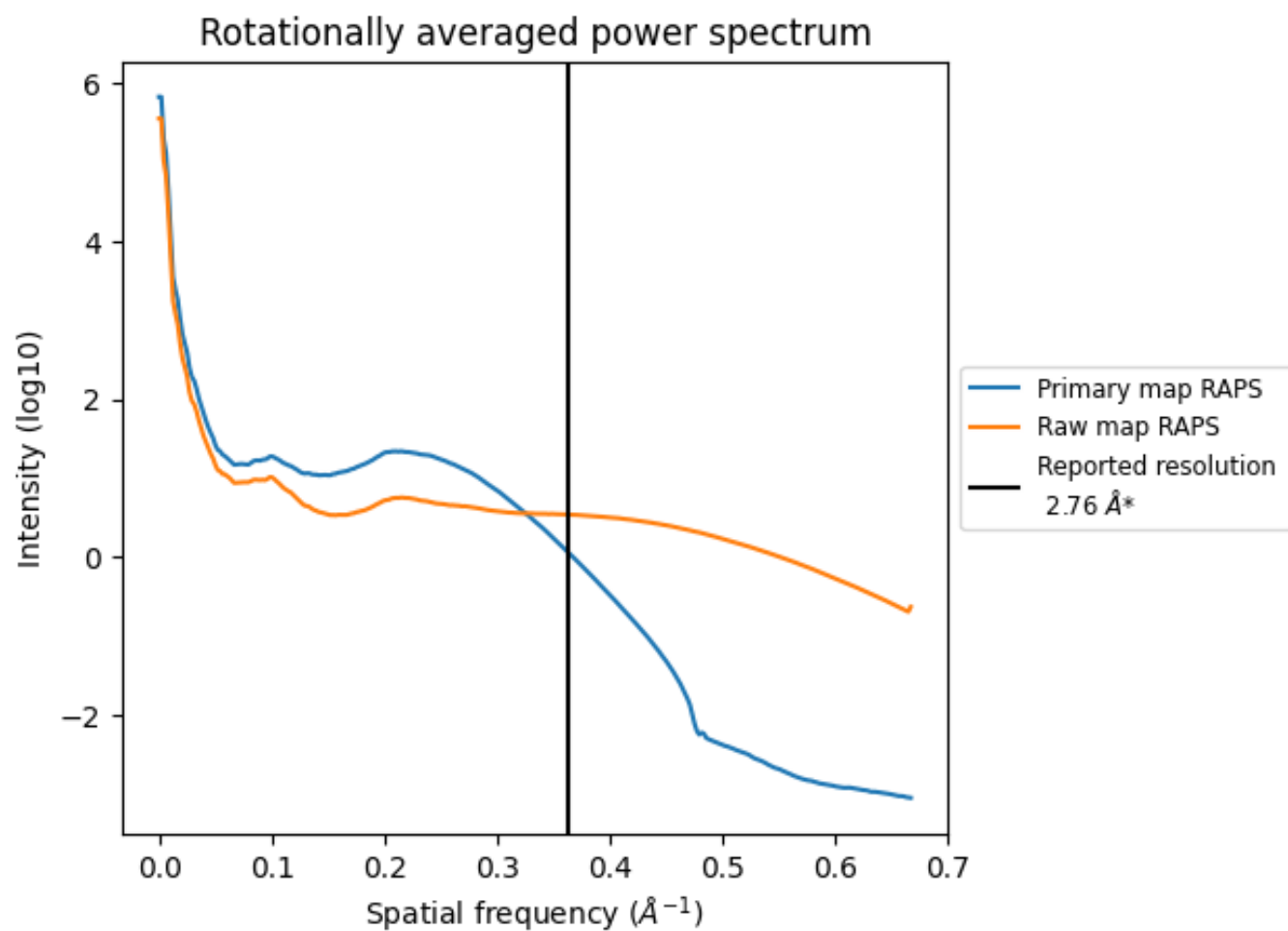
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 283 nm^3 ; this corresponds to an approximate mass of 256 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.

7.3 Rotationally averaged power spectrum ⓘ

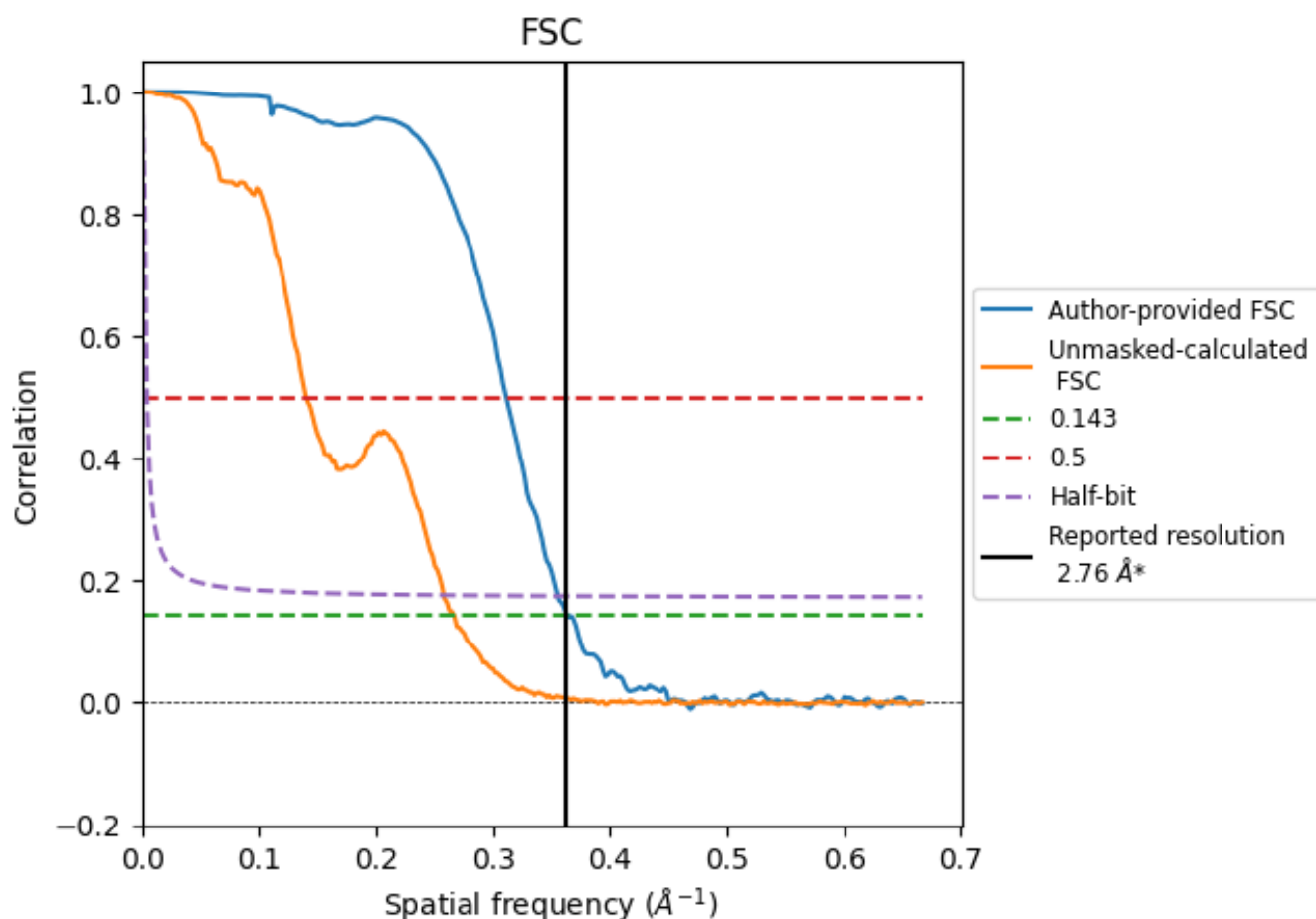


*Reported resolution corresponds to spatial frequency of 0.362 Å⁻¹

8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.362 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

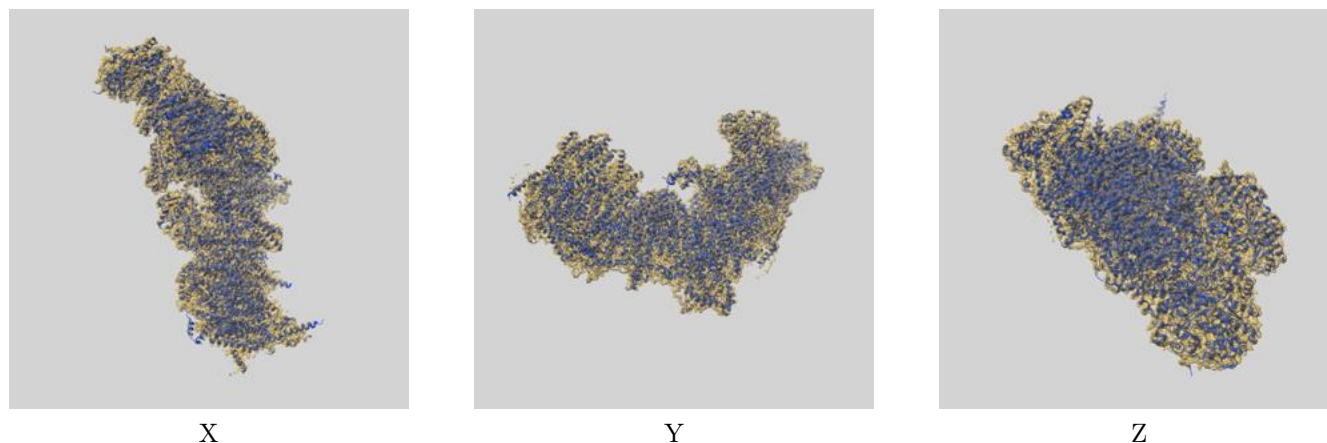
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.76	-	-
Author-provided FSC curve	2.75	3.21	2.81
Unmasked-calculated*	3.75	7.12	3.88

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.75 differs from the reported value 2.76 by more than 10 %

9 Map-model fit [i](#)

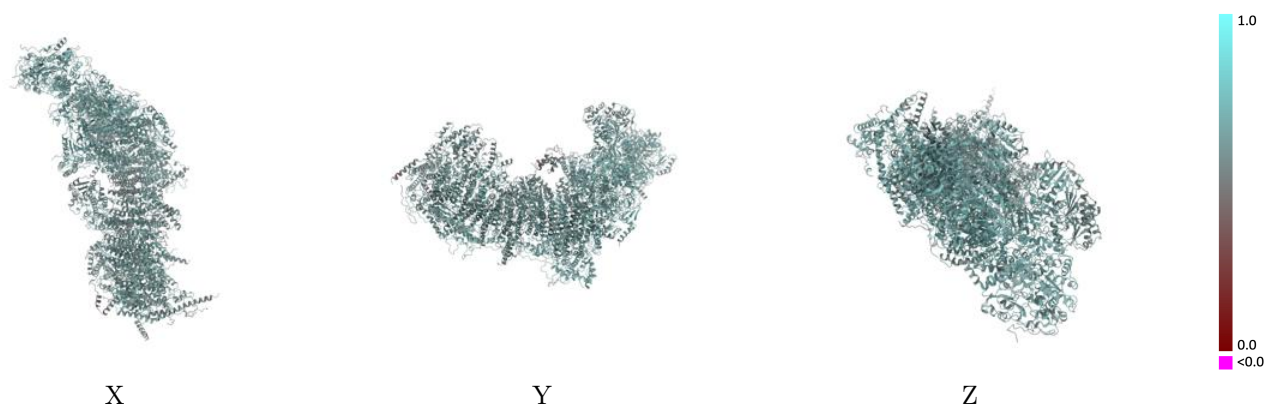
This section contains information regarding the fit between EMDB map EMD-14133 and PDB model 7QSL. Per-residue inclusion information can be found in section [3](#) on page [23](#).

9.1 Map-model overlay [i](#)



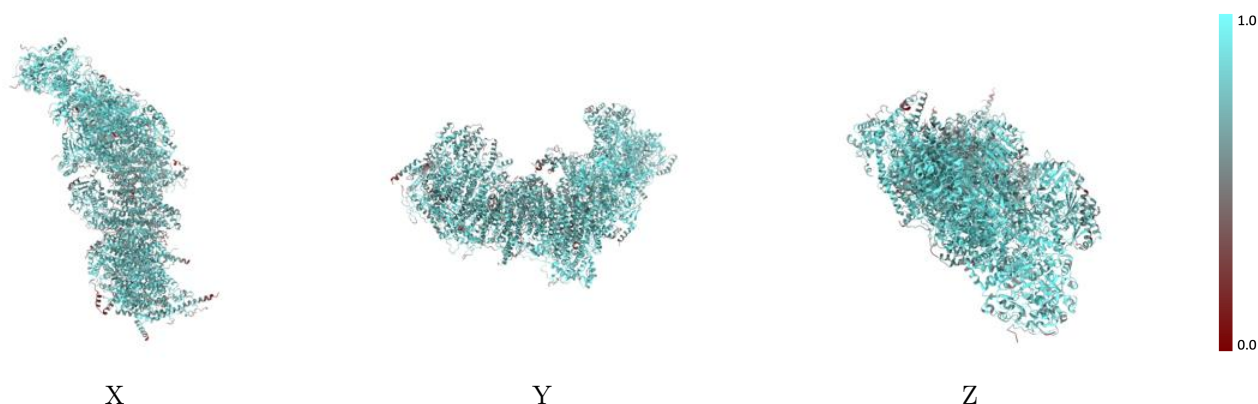
The images above show the 3D surface view of the map at the recommended contour level 6.5 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

9.2 Q-score mapped to coordinate model [i](#)



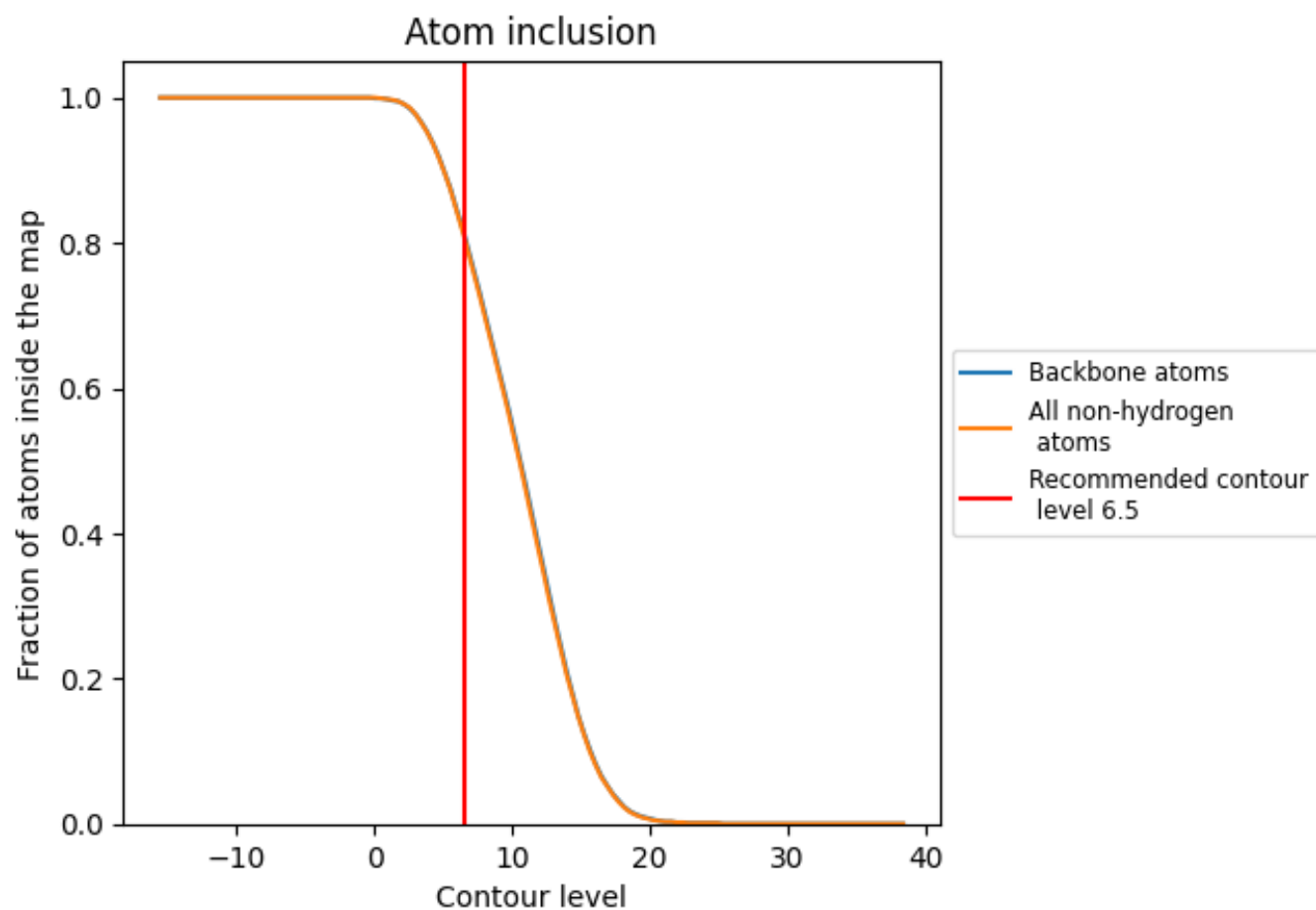
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

9.3 Atom inclusion mapped to coordinate model [i](#)



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (6.5).

9.4 Atom inclusion [i](#)



At the recommended contour level, 82% of all backbone atoms, 81% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary ⓘ

The table lists the average atom inclusion at the recommended contour level (6.5) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8110	 0.6250
A	 0.7800	 0.6350
B	 0.9330	 0.6740
C	 0.9040	 0.6680
D	 0.9050	 0.6650
E	 0.8130	 0.6160
F	 0.8450	 0.6230
G	 0.8360	 0.6300
H	 0.9060	 0.6630
I	 0.9270	 0.6750
J	 0.8270	 0.6330
K	 0.8490	 0.6500
L	 0.8180	 0.6240
M	 0.8750	 0.6510
N	 0.8960	 0.6600
O	 0.7740	 0.6010
P	 0.8290	 0.6300
Q	 0.8220	 0.6430
R	 0.8290	 0.6450
S	 0.6900	 0.5660
T	 0.5160	 0.4970
U	 0.6850	 0.5600
V	 0.7740	 0.6150
W	 0.7970	 0.6350
X	 0.7760	 0.6140
Y	 0.6900	 0.5790
Z	 0.8150	 0.6280
a	 0.8650	 0.6390
b	 0.7600	 0.6020
c	 0.7330	 0.5820
d	 0.7810	 0.6250
e	 0.7690	 0.6180
f	 0.6600	 0.5880
g	 0.7710	 0.6110
h	 0.8160	 0.6310



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Chain	Atom inclusion	Q-score
i	 0.6400	 0.5430
j	 0.7080	 0.5680
k	 0.6480	 0.5480
l	 0.7860	 0.5890
m	 0.7470	 0.6010
n	 0.7580	 0.5800
o	 0.6780	 0.5510
p	 0.7730	 0.6100
q	 0.8330	 0.6390
r	 0.7920	 0.6320
s	 0.7650	 0.6010