



Full wwPDB EM Validation Report ⓘ

Mar 30, 2026 – 03:28 AM UTC

PDB ID : 6ZKC / pdb_00006zkc
EMDB ID : EMD-11244
Title : Complex I during turnover, closed
Authors : Kampjut, D.; Sazanov, L.A.
Deposited on : 2020-06-30
Resolution : 3.10 Å(reported)
Based on initial model : 5LNK

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:

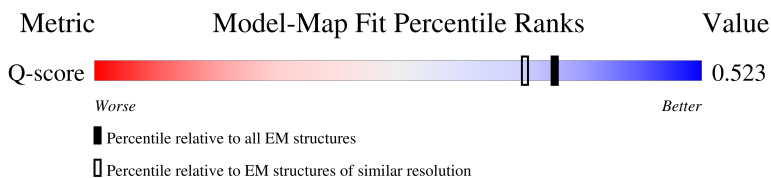
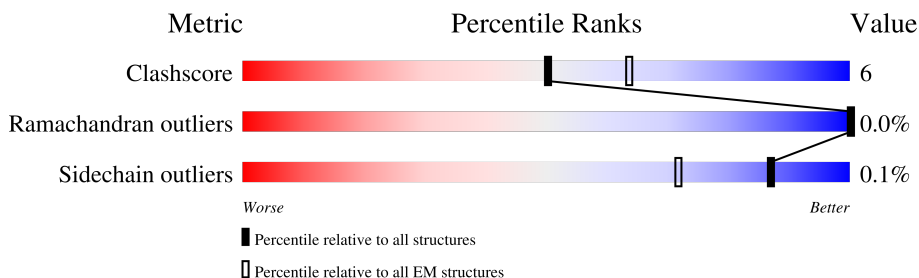
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data 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 3.10 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	14724 (2.60 - 3.60)

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	1	464	
2	2	246	
3	3	727	
4	4	463	

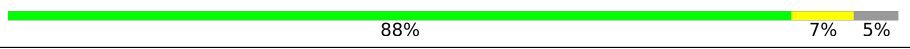
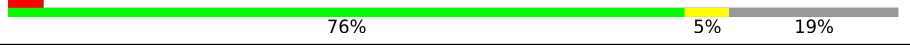
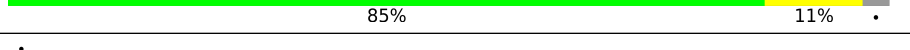
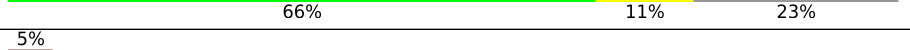
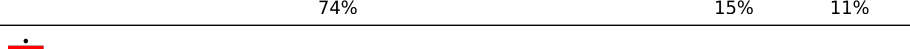
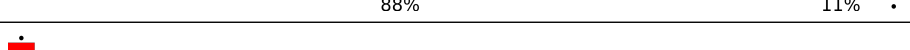
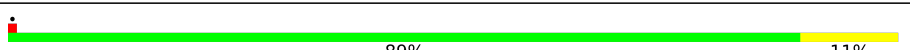
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Mol	Chain	Length	Quality of chain
5	5	266	
6	6	223	
7	9	217	
8	A	115	
9	H	318	
10	J	175	
11	K	98	
12	L	606	
13	M	459	
14	N	347	
15	V	141	
16	W	189	
17	X	157	
17	j	157	
18	Y	172	
19	Z	175	
20	a	109	
21	b	124	
22	c	170	
23	d	380	
24	e	99	
25	f	116	
26	g	140	
27	h	114	
28	i	145	

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Mol	Chain	Length	Quality of chain
29	k	355	
30	l	106	
31	m	84	
32	n	98	
33	o	122	
34	p	130	
35	q	144	
36	r	128	
37	s	137	
38	t	179	
39	u	108	
40	v	186	
41	w	154	
42	x	76	
43	y	58	
44	z	70	

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
45	SF4	1	501	-	-	X	-
53	CDL	L	1004	X	-	-	-
53	CDL	L	1005	X	-	-	-
53	CDL	M	503	X	-	-	-
53	CDL	x	101	X	-	-	-

2 Entry composition

There are 58 unique types of molecules in this entry. The entry contains 67967 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 dehydrogenase [ubiquinone] flavoprotein 1, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	1	430	Total	C	N	O	S	0	0
			3312	2086	593	613	20		

- Molecule 2 is a protein called Mitochondrial complex I, 24 kDa subunit.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	2	213	Total	C	N	O	S	0	0
			1655	1058	278	309	10		

- Molecule 3 is a protein called NADH:ubiquinone oxidoreductase core subunit S1.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	3	688	Total	C	N	O	S	0	0
			5275	3301	922	1011	41		

- Molecule 4 is a protein called Mitochondrial complex I, 49 kDa subunit.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	4	430	Total	C	N	O	S	0	0
			3457	2207	594	631	25		

- Molecule 5 is a protein called NADH:ubiquinone oxidoreductase core subunit S3.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	5	208	Total	C	N	O	S	0	0
			1726	1112	296	315	3		

- Molecule 6 is a protein called Mitochondrial complex I, PSST subunit.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	6	156	Total	C	N	O	S	0	0
			1247	795	225	213	14		

- Molecule 7 is a protein called Mitochondrial complex I, TYKY subunit.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	9	176	Total	C	N	O	S	0	0
			1414	889	243	270	12		

- Molecule 8 is a protein called NADH-ubiquinone oxidoreductase chain 3.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	A	115	Total	C	N	O	S	0	0
			922	621	133	161	7		

- Molecule 9 is a protein called NADH-ubiquinone oxidoreductase chain 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	H	318	Total	C	N	O	S	0	0
			2528	1704	384	421	19		

- 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
			1344	904	192	235	13		

- 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
			749	490	112	132	15		

- 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
			4807	3188	746	829	44		

- 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
			3647	2429	571	607	40		

- 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
			2723	1808	416	459	40		

- Molecule 15 is a protein called Mitochondrial complex I, B14.7 subunit.

Mol	Chain	Residues	Atoms					AltConf	Trace
15	V	140	Total	C	N	O	S	0	0
			1028	656	175	191	6		

- Molecule 16 is a protein called NADH:ubiquinone oxidoreductase subunit B5.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	W	139	Total	C	N	O	S	0	0
			1155	761	194	198	2		

- Molecule 17 is a protein called Acyl carrier protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
17	X	87	Total	C	N	O	S	0	0
			701	451	103	142	5		
17	j	82	Total	C	N	O	S	0	0
			660	425	98	132	5		

- Molecule 18 is a protein called NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 8.

Mol	Chain	Residues	Atoms					AltConf	Trace
18	Y	171	Total	C	N	O	S	0	0
			1403	889	253	251	10		

- Molecule 19 is a protein called Mitochondrial complex I, PDSW subunit.

Mol	Chain	Residues	Atoms					AltConf	Trace
19	Z	171	Total	C	N	O	S	0	0
			1441	905	266	262	8		

- Molecule 20 is a protein called Mitochondrial complex I, 10 kDa subunit.

Mol	Chain	Residues	Atoms					AltConf	Trace
20	a	44	Total	C	N	O	S	0	0
			371	233	66	71	1		

- Molecule 21 is a protein called Mitochondrial complex I, 13 kDa subunit.

Mol	Chain	Residues	Atoms					AltConf	Trace
21	b	95	Total	C	N	O	S	0	0
			737	451	139	144	3		

- Molecule 22 is a protein called NADH dehydrogenase [ubiquinone] iron-sulfur protein 4, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
22	c	126	Total	C	N	O	S	0	0
			1024	646	182	193	3		

- Molecule 23 is a protein called NADH:ubiquinone oxidoreductase subunit A9.

Mol	Chain	Residues	Atoms					AltConf	Trace
23	d	340	Total	C	N	O	S	0	0
			2748	1775	489	478	6		

- Molecule 24 is a protein called NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
24	e	86	Total	C	N	O	S	0	0
			691	434	129	126	2		

- Molecule 25 is a protein called Mitochondrial complex I, B13 subunit.

Mol	Chain	Residues	Atoms					AltConf	Trace
25	f	113	Total	C	N	O	S	0	0
			917	595	153	167	2		

- Molecule 26 is a protein called NADH:ubiquinone oxidoreductase subunit A6.

Mol	Chain	Residues	Atoms					AltConf	Trace
26	g	114	Total	C	N	O	S	0	0
			969	619	180	166	4		

- Molecule 27 is a protein called Mitochondrial complex I, B14.5a subunit.

Mol	Chain	Residues	Atoms					AltConf	Trace
27	h	96	Total	C	N	O	S	0	0
			769	480	146	140	3		

- Molecule 28 is a protein called NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 12.

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

- Molecule 29 is a protein called NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 10, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
29	k	320	Total	C	N	O	P S	0	0
			2596	1659	432	494	1 10		

- Molecule 30 is a protein called NADH:ubiquinone oxidoreductase subunit S5.

Mol	Chain	Residues	Atoms					AltConf	Trace
30	l	105	Total	C	N	O	S	0	0
			874	551	164	153	6		

- Molecule 31 is a protein called NADH:ubiquinone oxidoreductase subunit A3.

Mol	Chain	Residues	Atoms					AltConf	Trace
31	m	80	Total	C	N	O	S	0	0
			626	411	103	110	2		

- Molecule 32 is a protein called NADH:ubiquinone oxidoreductase subunit B3.

Mol	Chain	Residues	Atoms					AltConf	Trace
32	n	79	Total	C	N	O	S	0	0
			634	415	106	111	2		

- Molecule 33 is a protein called NADH dehydrogenase [ubiquinone] 1 subunit C2.

Mol	Chain	Residues	Atoms					AltConf	Trace
33	o	120	Total	C	N	O	S	0	0
			1004	652	175	172	5		

- Molecule 34 is a protein called NADH:ubiquinone oxidoreductase subunit B4.

Mol	Chain	Residues	Atoms					AltConf	Trace
34	p	128	Total	C	N	O	S	0	0
			1059	675	189	194	1		

- Molecule 35 is a protein called Mitochondrial complex I, B16.6 subunit.

Mol	Chain	Residues	Atoms					AltConf	Trace
35	q	139	Total	C	N	O	S	0	0
			1142	733	200	200	9		

- Molecule 36 is a protein called Mitochondrial complex I, B17 subunit.

Mol	Chain	Residues	Atoms					AltConf	Trace
36	r	99	Total	C	N	O	S	0	0
			846	554	149	142	1		

- Molecule 37 is a protein called NADH:ubiquinone oxidoreductase subunit B7.

Mol	Chain	Residues	Atoms					AltConf	Trace
37	s	122	Total	C	N	O	S	0	0
			1047	653	199	186	9		

- Molecule 38 is a protein called NADH:ubiquinone oxidoreductase subunit B9.

Mol	Chain	Residues	Atoms					AltConf	Trace
38	t	177	Total	C	N	O	S	0	0
			1520	973	279	262	6		

- Molecule 39 is a protein called NADH:ubiquinone oxidoreductase subunit B2.

Mol	Chain	Residues	Atoms					AltConf	Trace
39	u	65	Total	C	N	O	S	0	0
			563	372	93	97	1		

- Molecule 40 is a protein called NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 8, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
40	v	155	Total	C	N	O	S	0	0
			1307	846	213	239	9		

- Molecule 41 is a protein called Mitochondrial complex I, ESSS subunit.

Mol	Chain	Residues	Atoms					AltConf	Trace
41	w	101	Total	C	N	O	S	0	0
			846	542	140	160	4		

- Molecule 42 is a protein called Mitochondrial complex I, KFYI subunit.

Mol	Chain	Residues	Atoms				AltConf	Trace
42	x	49	Total	C	N	O	0	0
			412	271	70	71		

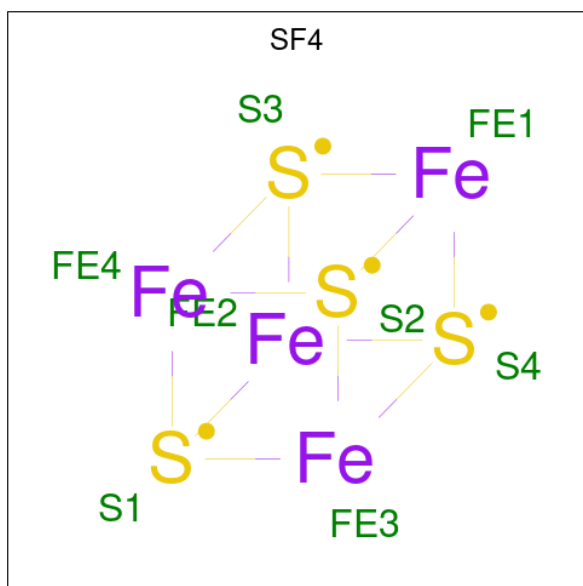
- Molecule 43 is a protein called Mitochondrial complex I, MNLL subunit.

Mol	Chain	Residues	Atoms				AltConf	Trace
43	y	50	Total	C	N	O	0	0
			436	287	77	72		

- Molecule 44 is a protein called Mitochondrial complex I, MWFE subunit.

Mol	Chain	Residues	Atoms					AltConf	Trace
44	z	70	Total	C	N	O	S	0	0
			576	369	106	96	5		

- Molecule 45 is IRON/SULFUR CLUSTER (CCD ID: SF4) (formula: Fe₄S₄).



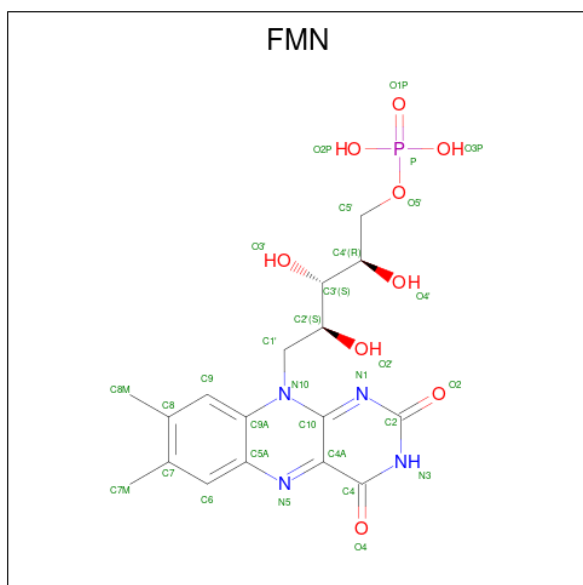
Mol	Chain	Residues	Atoms			AltConf
45	1	1	Total	Fe	S	0
			8	4	4	
45	3	1	Total	Fe	S	0
			8	4	4	
45	3	1	Total	Fe	S	0
			8	4	4	
45	6	1	Total	Fe	S	0
			8	4	4	

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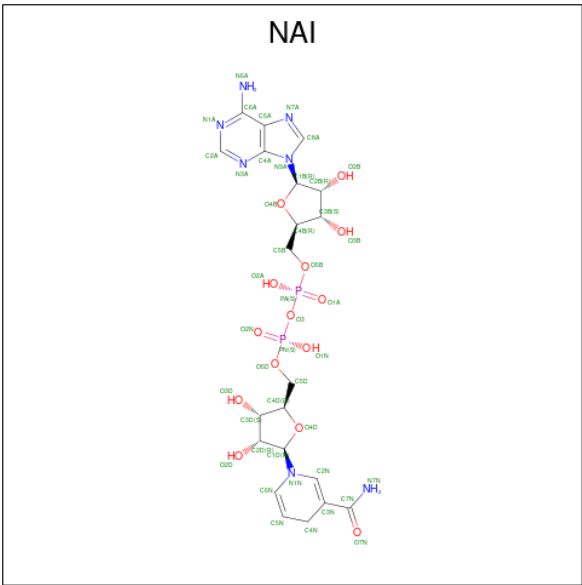
Mol	Chain	Residues	Atoms			AltConf
45	9	1	Total	Fe	S	0
			8	4	4	
45	9	1	Total	Fe	S	0
			8	4	4	

- Molecule 46 is FLAVIN MONONUCLEOTIDE (CCD ID: FMN) (formula: $C_{17}H_{21}N_4O_9P$).



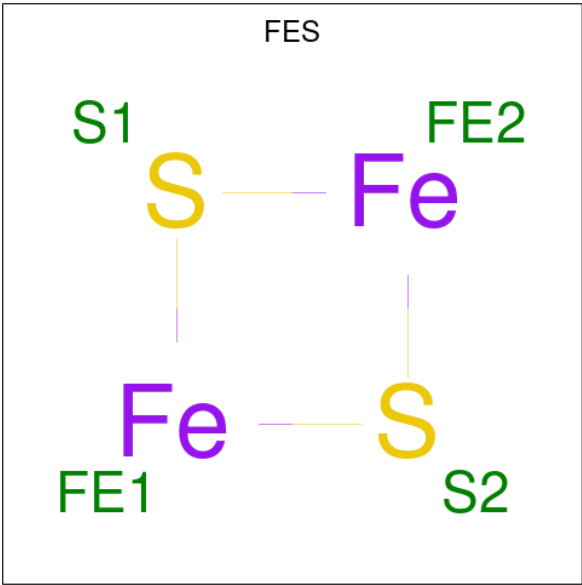
Mol	Chain	Residues	Atoms					AltConf
46	1	1	Total	C	N	O	P	0
			31	17	4	9	1	

- Molecule 47 is 1,4-DIHYDRONICOTINAMIDE ADENINE DINUCLEOTIDE (CCD ID: NAI) (formula: $C_{21}H_{29}N_7O_{14}P_2$).



Mol	Chain	Residues	Atoms					AltConf
47	1	1	Total	C	N	O	P	0
			44	21	7	14	2	

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

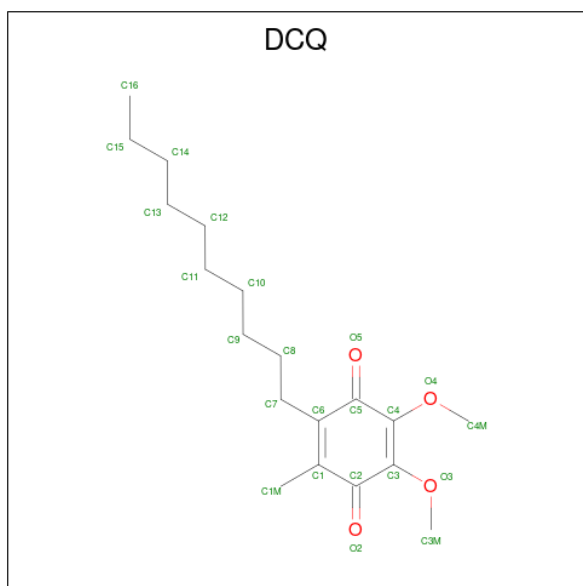


Mol	Chain	Residues	Atoms			AltConf
48	2	1	Total	Fe	S	0
			4	2	2	
48	3	1	Total	Fe	S	0
			4	2	2	

- Molecule 49 is POTASSIUM ION (CCD ID: K) (formula: K).

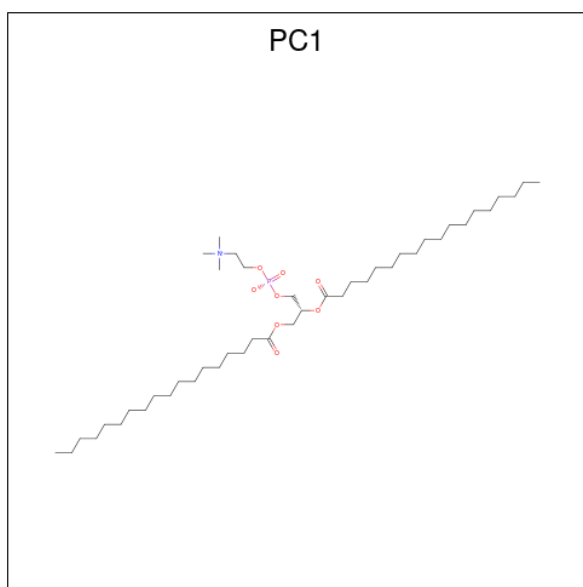
Mol	Chain	Residues	Atoms		AltConf
49	3	1	Total	K	0
			1	1	

- Molecule 50 is 2-decyl-5,6-dimethoxy-3-methylcyclohexa-2,5-diene-1,4-dione (CCD ID: DCQ) (formula: $C_{19}H_{30}O_4$).



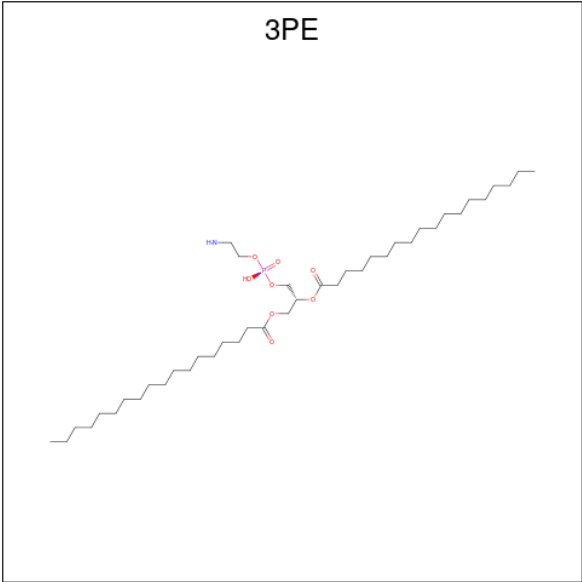
Mol	Chain	Residues	Atoms			AltConf
50	4	1	Total	C	O	0
			23	19	4	
50	H	1	Total	C	O	0
			23	19	4	

- Molecule 51 is 1,2-DIACYL-SN-GLYCERO-3-PHOSPHOCHOLINE (CCD ID: PC1) (formula: $C_{44}H_{88}NO_8P$).



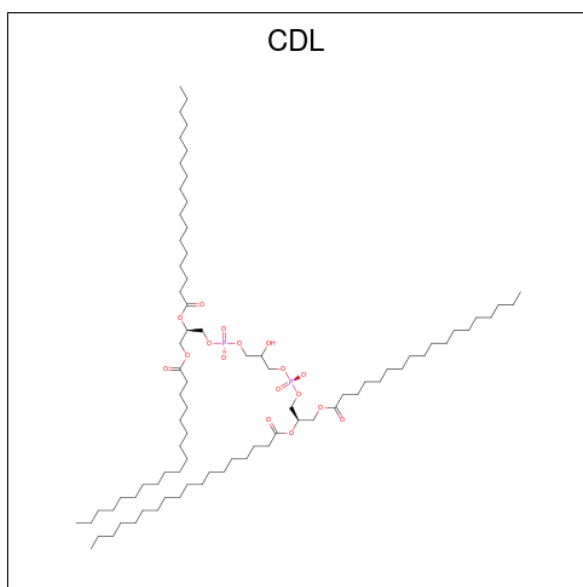
Mol	Chain	Residues	Atoms					AltConf
51	6	1	Total	C	N	O	P	0
			46	36	1	8	1	
51	9	1	Total	C	N	O	P	0
			54	44	1	8	1	
51	L	1	Total	C	N	O	P	0
			54	44	1	8	1	
51	M	1	Total	C	N	O	P	0
			54	44	1	8	1	
51	w	1	Total	C	N	O	P	0
			54	44	1	8	1	

- Molecule 52 is 1,2-Distearoyl-sn-glycerophosphoethanolamine (CCD ID: 3PE) (formula: $C_{41}H_{82}NO_8P$).



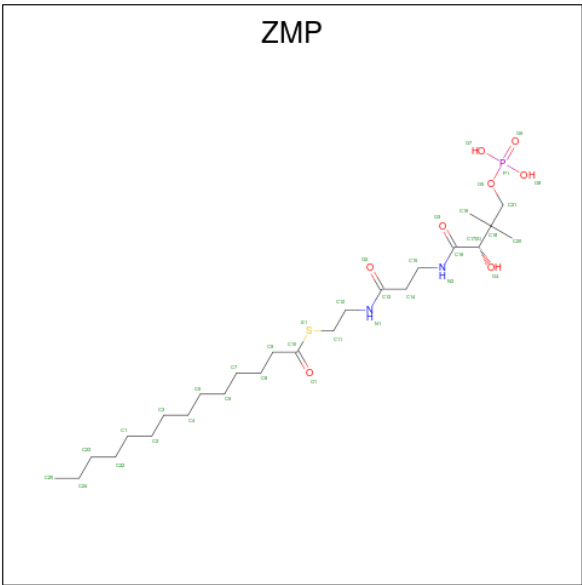
Mol	Chain	Residues	Atoms					AltConf
52	A	1	Total	C	N	O	P	0
			51	41	1	8	1	
52	J	1	Total	C	N	O	P	0
			51	41	1	8	1	
52	J	1	Total	C	N	O	P	0
			40	30	1	8	1	
52	L	1	Total	C	N	O	P	0
			51	41	1	8	1	
52	L	1	Total	C	N	O	P	0
			31	21	1	8	1	
52	M	1	Total	C	N	O	P	0
			44	34	1	8	1	
52	N	1	Total	C	N	O	P	0
			40	30	1	8	1	
52	N	1	Total	C	N	O	P	0
			51	41	1	8	1	
52	V	1	Total	C	N	O	P	0
			37	27	1	8	1	
52	i	1	Total	C	N	O	P	0
			51	41	1	8	1	
52	i	1	Total	C	N	O	P	0
			51	41	1	8	1	
52	o	1	Total	C	N	O	P	0
			31	21	1	8	1	
52	p	1	Total	C	O	P		0
			27	18	8	1		

- Molecule 53 is CARDIOLIPIN (CCD ID: CDL) (formula: C₈₁H₁₅₆O₁₇P₂).



Mol	Chain	Residues	Atoms				AltConf
53	L	1	Total	C	O	P	0
			85	66	17	2	
53	L	1	Total	C	O	P	0
			100	81	17	2	
53	M	1	Total	C	O	P	0
			100	81	17	2	
53	V	1	Total	C	O	P	0
			94	75	17	2	
53	W	1	Total	C	O	P	0
			100	81	17	2	
53	h	1	Total	C	O	P	0
			58	39	17	2	
53	o	1	Total	C	O	P	0
			90	71	17	2	
53	x	1	Total	C	O	P	0
			75	56	17	2	

- Molecule 54 is S-[2-({N-[(2S)-2-hydroxy-3,3-dimethyl-4-(phosphonooxy)butanoyl]-beta-alanyl}amino)ethyl] tetradecanethioate (CCD ID: ZMP) (formula: C₂₅H₄₉N₂O₈PS).

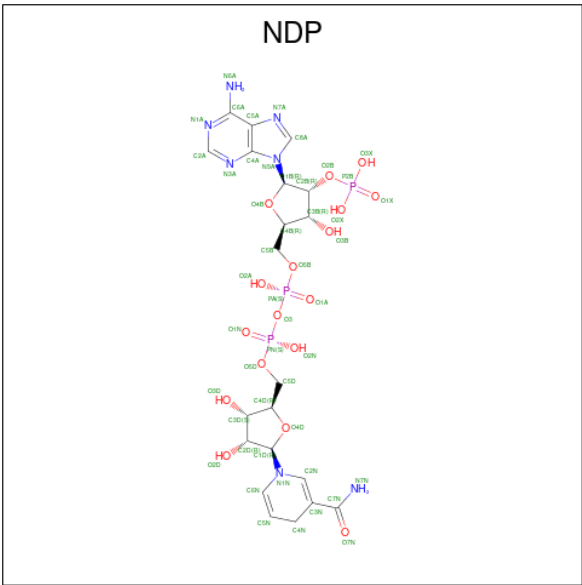


Mol	Chain	Residues	Atoms						AltConf
54	X	1	Total	C	N	O	P	S	0
			31	20	2	7	1	1	
54	g	1	Total	C	N	O	P	S	0
			34	23	2	7	1	1	

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

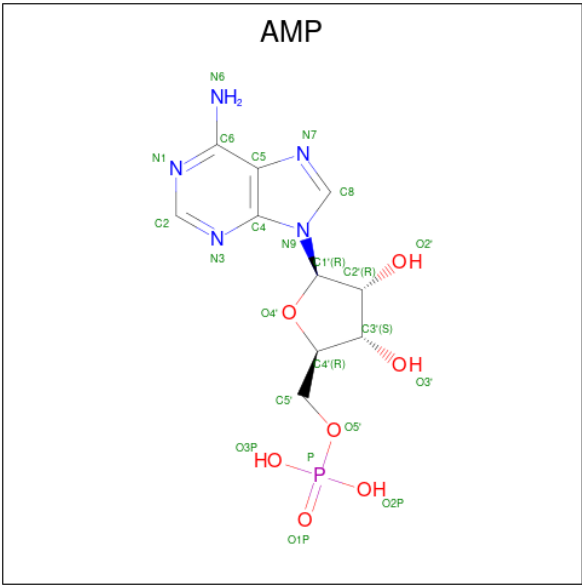
Mol	Chain	Residues	Atoms		AltConf
55	b	1	Total	Zn	0
			1	1	

- Molecule 56 is NADPH DIHYDRO-NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (CCD ID: NDP) (formula: C₂₁H₃₀N₇O₁₇P₃).



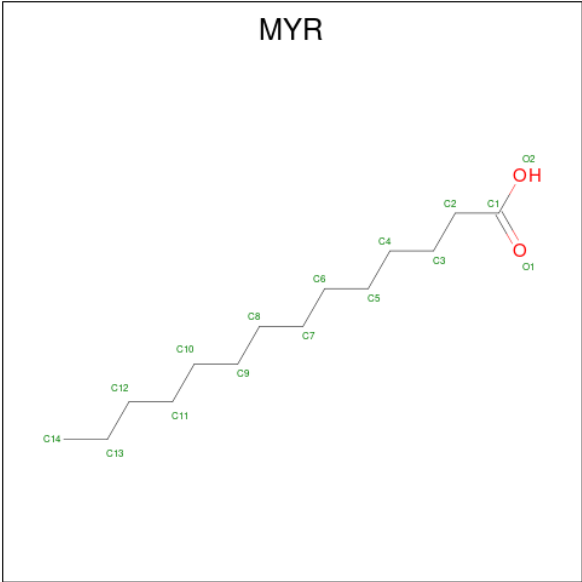
Mol	Chain	Residues	Atoms					AltConf
56	d	1	Total	C	N	O	P	0
			48	21	7	17	3	

- Molecule 57 is ADENOSINE MONOPHOSPHATE (CCD ID: AMP) (formula: $C_{10}H_{14}N_5O_7P$).



Mol	Chain	Residues	Atoms					AltConf
57	k	1	Total	C	N	O	P	0
			23	10	5	7	1	

- Molecule 58 is MYRISTIC ACID (CCD ID: MYR) (formula: $C_{14}H_{28}O_2$).



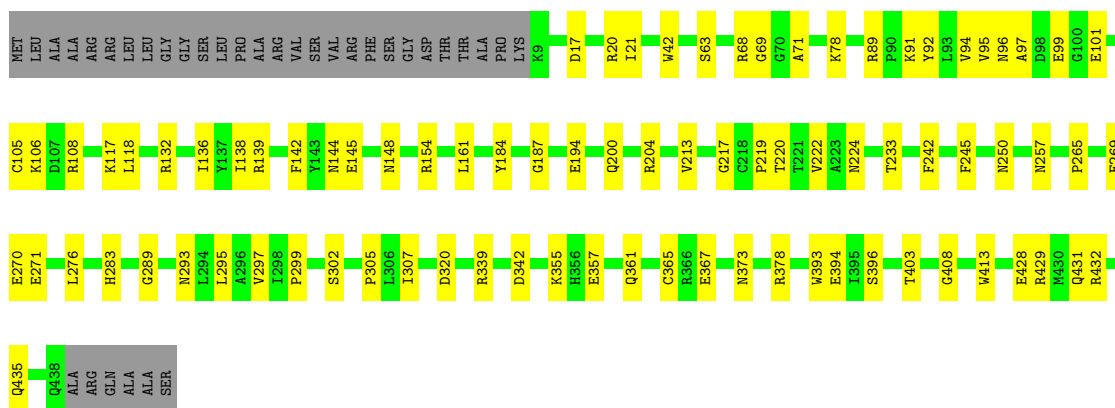
Mol	Chain	Residues	Atoms			AltConf
58	s	1	Total	C	O	0
			15	14	1	

3 Residue-property plots

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.

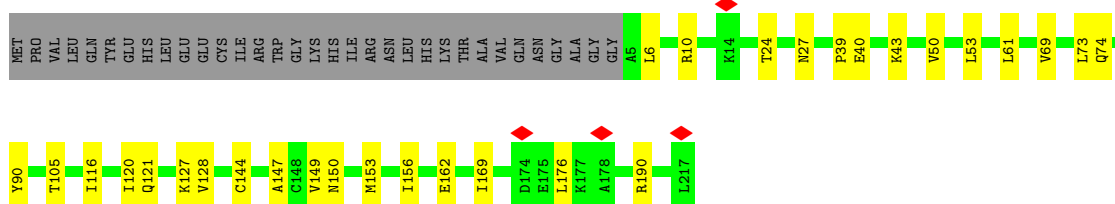
- Molecule 1: NADH dehydrogenase [ubiquinone] flavoprotein 1, mitochondrial

Chain 1: 



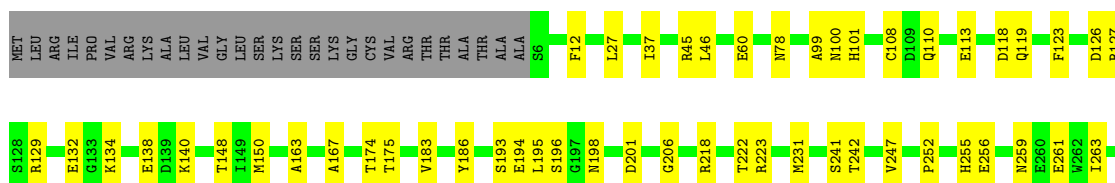
- Molecule 2: Mitochondrial complex I, 24 kDa subunit

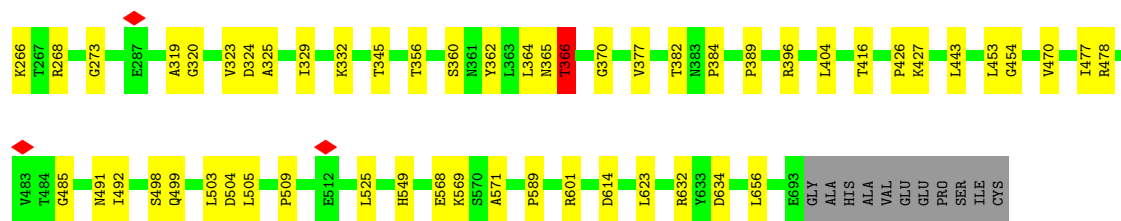
Chain 2: 



- Molecule 3: NADH:ubiquinone oxidoreductase core subunit S1

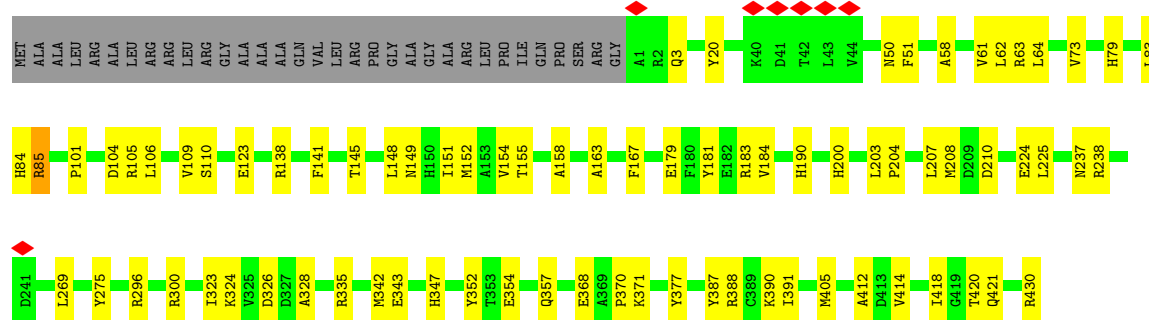
Chain 3: 





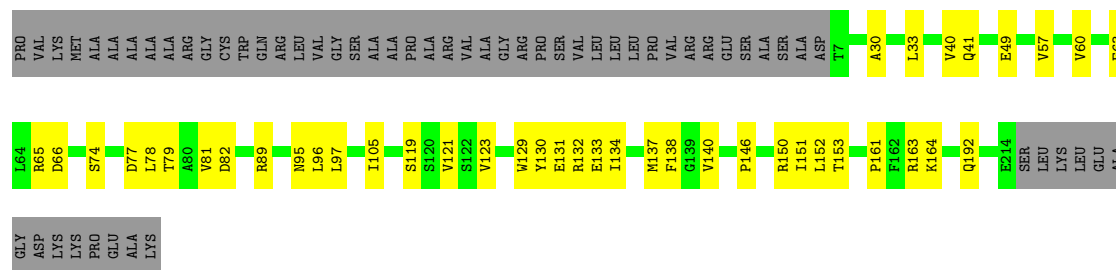
- Molecule 4: Mitochondrial complex I, 49 kDa subunit

Chain 4: 76% 17% 7%



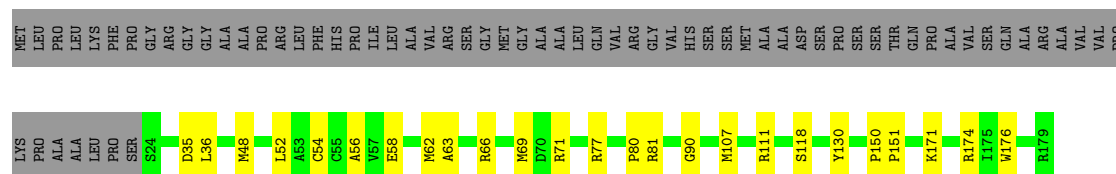
- Molecule 5: NADH:ubiquinone oxidoreductase core subunit S3

Chain 5: 62% 16% 22%



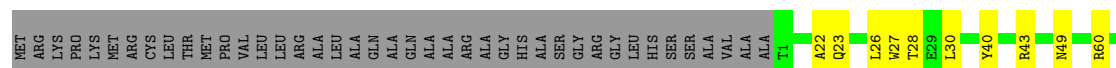
- Molecule 6: Mitochondrial complex I, PSST subunit

Chain 6: 59% 11% 30%



- Molecule 7: Mitochondrial complex I, TYKY subunit

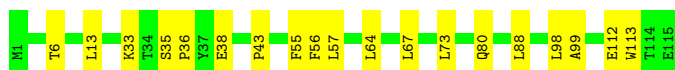
Chain 9: 63% 18% 19%





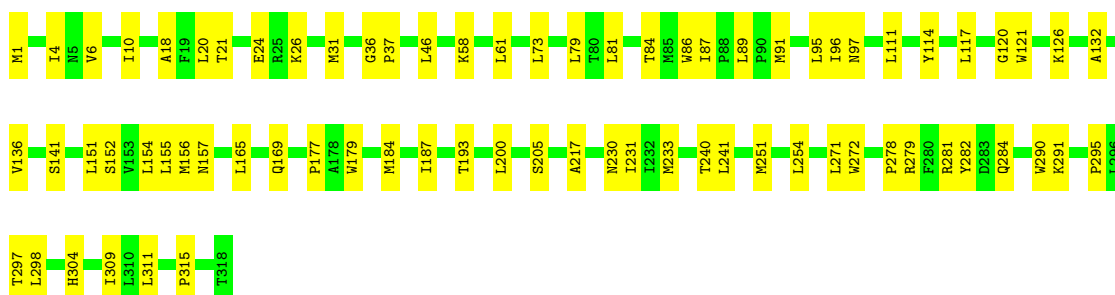
- Molecule 8: NADH-ubiquinone oxidoreductase chain 3

Chain A: 83% 17%



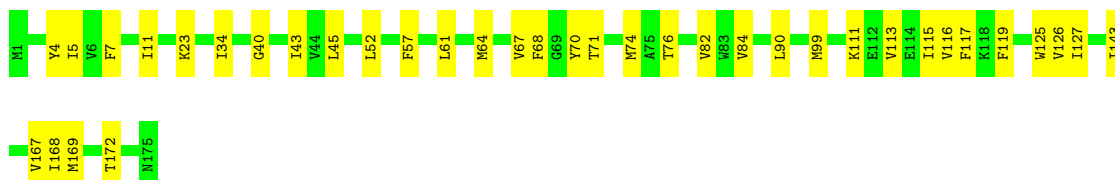
- Molecule 9: NADH-ubiquinone oxidoreductase chain 1

Chain H: 77% 23%



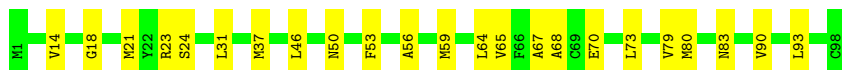
- Molecule 10: NADH-ubiquinone oxidoreductase chain 6

Chain J: 79% 21%



- Molecule 11: NADH-ubiquinone oxidoreductase chain 4L

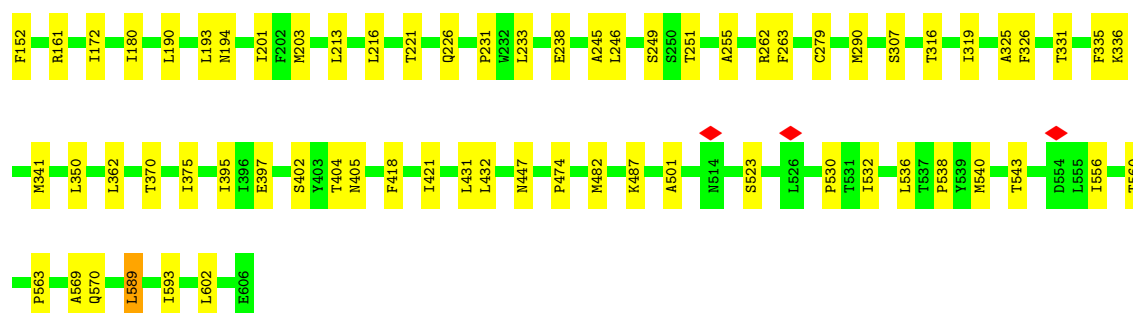
Chain K: 77% 23%



- Molecule 12: NADH-ubiquinone oxidoreductase chain 5

Chain L: 84% 16%





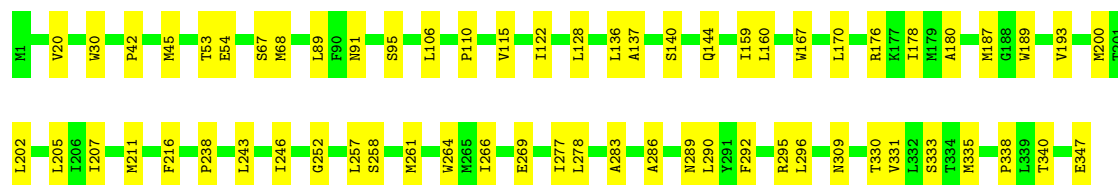
- Molecule 13: NADH-ubiquinone oxidoreductase chain 4

Chain M: 83% 17%



- Molecule 14: NADH-ubiquinone oxidoreductase chain 2

Chain N: 82% 18%



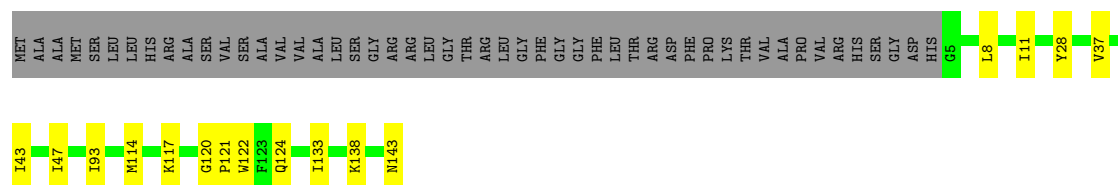
- Molecule 15: Mitochondrial complex I, B14.7 subunit

Chain V: 89% 10% ..

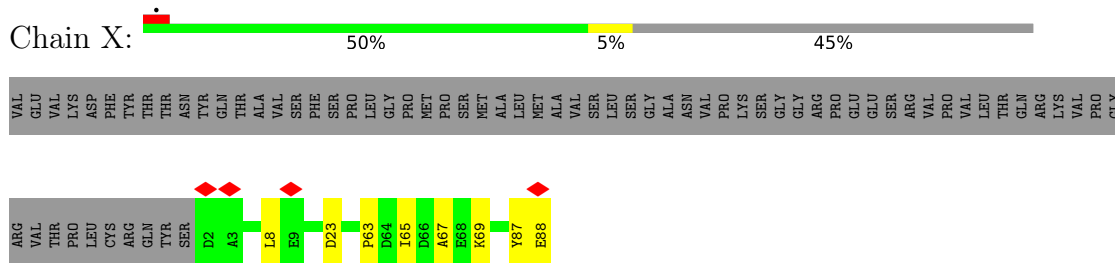


- Molecule 16: NADH:ubiquinone oxidoreductase subunit B5

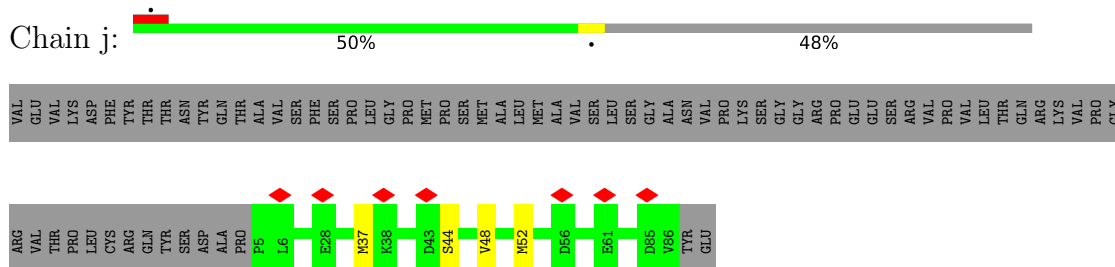
Chain W: 65% 8% 26%



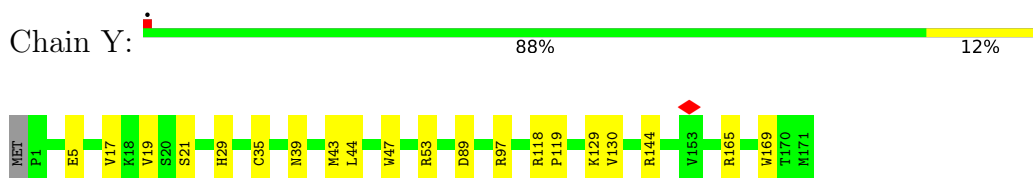
- Molecule 17: Acyl carrier protein



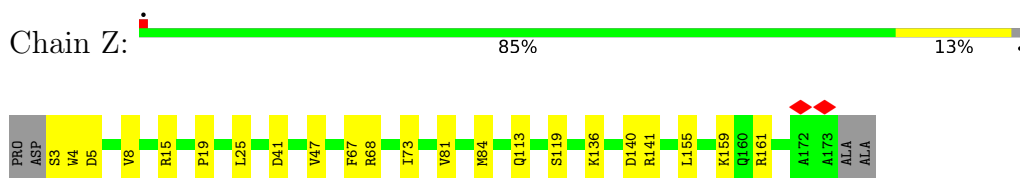
- Molecule 17: Acyl carrier protein



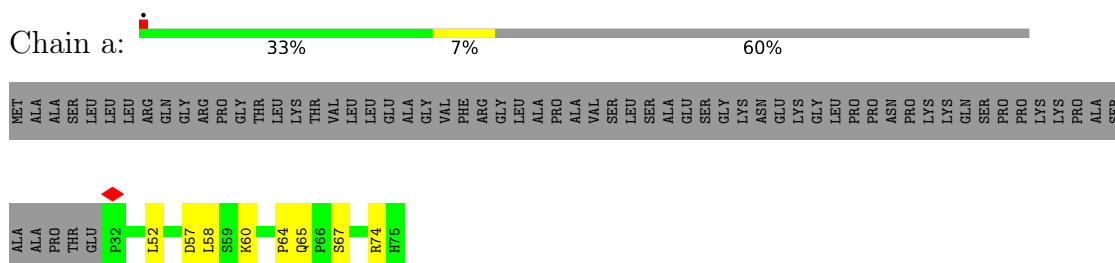
- Molecule 18: NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 8



- Molecule 19: Mitochondrial complex I, PDSW subunit

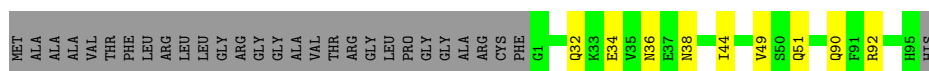


- Molecule 20: Mitochondrial complex I, 10 kDa subunit

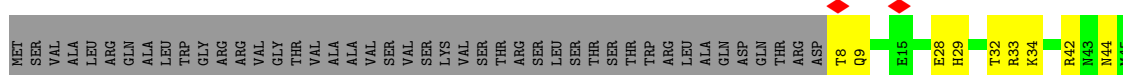


- Molecule 21: Mitochondrial complex I, 13 kDa subunit

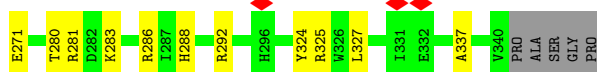
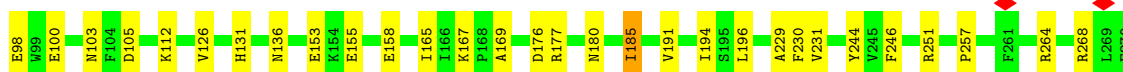
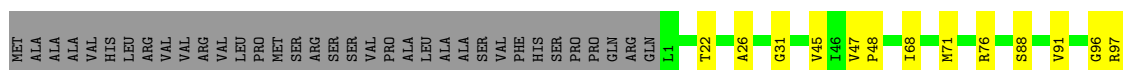
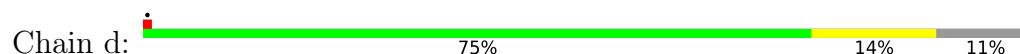




- Molecule 22: NADH dehydrogenase [ubiquinone] iron-sulfur protein 4, mitochondrial



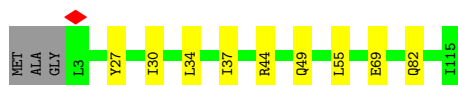
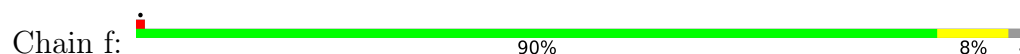
- Molecule 23: NADH:ubiquinone oxidoreductase subunit A9



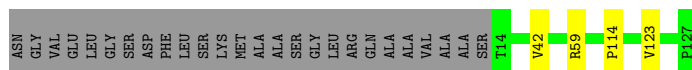
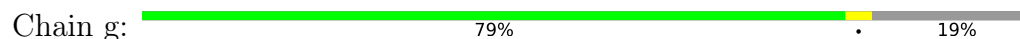
- Molecule 24: NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 2



- Molecule 25: Mitochondrial complex I, B13 subunit

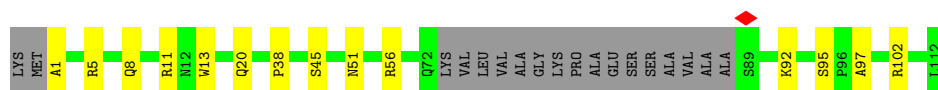


- Molecule 26: NADH:ubiquinone oxidoreductase subunit A6



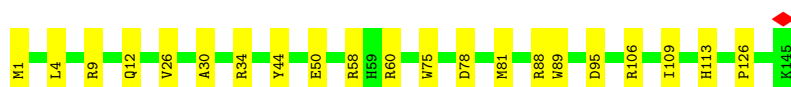
- Molecule 27: Mitochondrial complex I, B14.5a subunit

Chain h:  72% 12% 16%



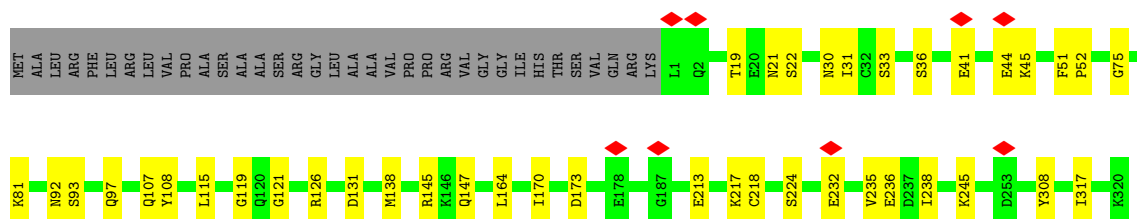
- Molecule 28: NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 12

Chain i:  86% 14%



- Molecule 29: NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 10, mitochondrial

Chain k:  79% 12% 10%



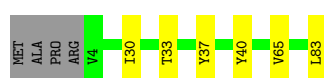
- Molecule 30: NADH:ubiquinone oxidoreductase subunit S5

Chain l:  91% 8%



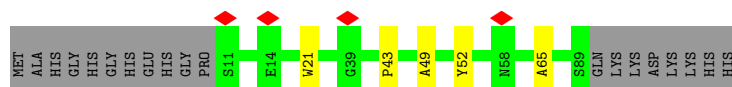
- Molecule 31: NADH:ubiquinone oxidoreductase subunit A3

Chain m:  88% 7% 5%

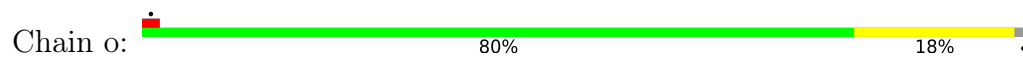


- Molecule 32: NADH:ubiquinone oxidoreductase subunit B3

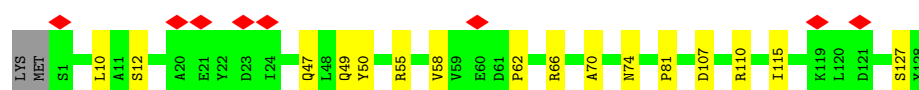
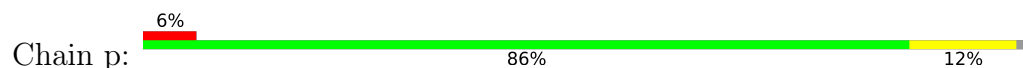
Chain n:  76% 5% 19%



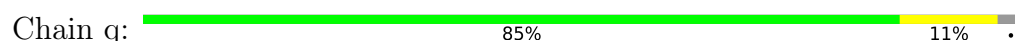
- Molecule 33: NADH dehydrogenase [ubiquinone] 1 subunit C2



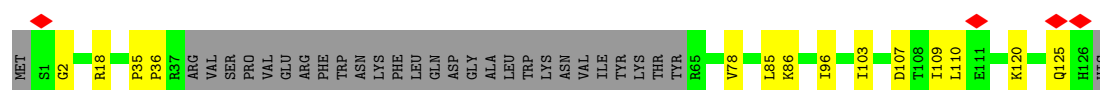
- Molecule 34: NADH:ubiquinone oxidoreductase subunit B4



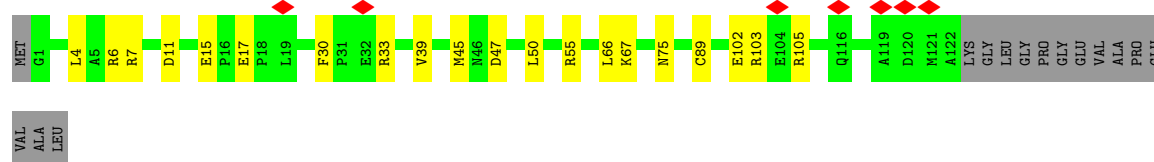
- Molecule 35: Mitochondrial complex I, B16.6 subunit



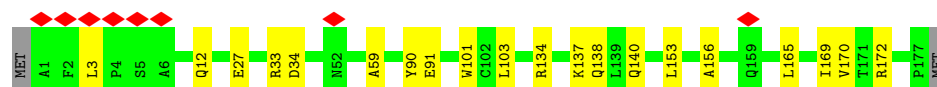
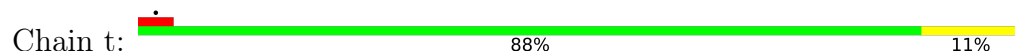
- Molecule 36: Mitochondrial complex I, B17 subunit



- Molecule 37: NADH:ubiquinone oxidoreductase subunit B7

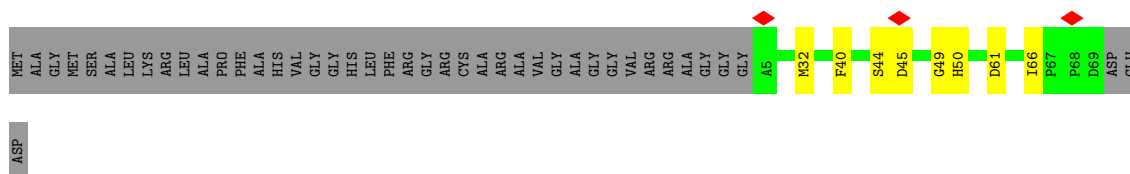


- Molecule 38: NADH:ubiquinone oxidoreductase subunit B9

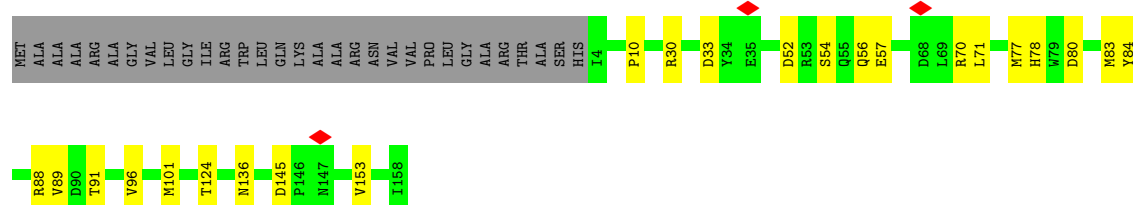


- Molecule 39: NADH:ubiquinone oxidoreductase subunit B2

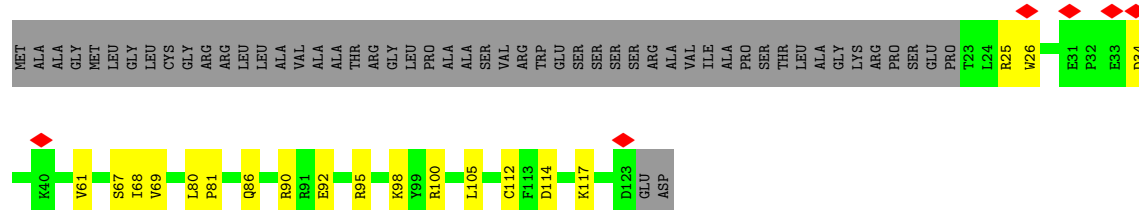




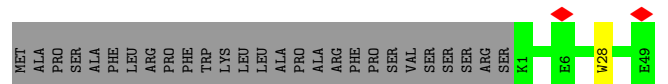
- Molecule 40: NADH dehydrogenase [ubiquinone] 1 beta subcomplex subunit 8, mitochondrial



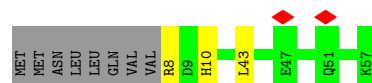
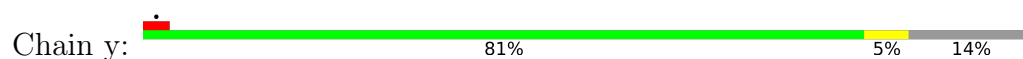
- Molecule 41: Mitochondrial complex I, ESSS subunit



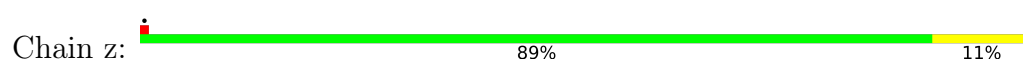
- Molecule 42: Mitochondrial complex I, KFYI subunit



- Molecule 43: Mitochondrial complex I, MNLL subunit



- Molecule 44: Mitochondrial complex I, MWFE subunit





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	15769	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 ($e^-/\text{\AA}^2$)	100	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	FEI FALCON III (4k x 4k)	Depositor
Maximum map value	0.447	Depositor
Minimum map value	-0.150	Depositor
Average map value	0.003	Depositor
Map value standard deviation	0.029	Depositor
Recommended contour level	0.045	Depositor
Map size (Å)	170.821, 193.102, 283.287	wwPDB
Map dimensions	267, 182, 161	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.061, 1.061, 1.061	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZMP, AMP, 3PE, NDP, FES, SF4, FMN, K, DCQ, 2MR, NAI, AYA, MYR, ZN, PC1, FME, SEP, CDL

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	1	0.29	0/3386	0.56	0/4575
2	2	0.27	0/1695	0.56	0/2306
3	3	0.29	0/5362	0.53	2/7266 (0.0%)
4	4	0.33	0/3535	0.55	0/4791
5	5	0.30	0/1776	0.56	2/2417 (0.1%)
6	6	0.36	0/1278	0.57	0/1728
7	9	0.35	0/1445	0.55	0/1956
8	A	0.31	0/947	0.57	0/1296
9	H	0.36	0/2603	0.67	3/3561 (0.1%)
10	J	0.31	0/1378	0.62	0/1868
11	K	0.33	0/749	0.58	0/1014
12	L	0.28	0/4925	0.56	0/6700
13	M	0.32	0/3731	0.57	0/5085
14	N	0.33	0/2787	0.57	0/3795
15	V	0.20	0/1041	0.43	0/1412
16	W	0.25	0/1188	0.48	0/1607
17	X	0.20	0/713	0.45	0/963
17	j	0.20	0/670	0.46	0/902
18	Y	0.27	0/1440	0.54	0/1942
19	Z	0.23	0/1475	0.47	0/1989
20	a	0.18	0/383	0.40	0/518
21	b	0.25	0/749	0.42	0/1009
22	c	0.28	0/1047	0.44	0/1415
23	d	0.27	0/2824	0.54	2/3830 (0.1%)
24	e	0.22	0/702	0.49	0/945
25	f	0.24	0/937	0.51	0/1271
26	g	0.26	0/993	0.52	0/1336
27	h	0.27	0/779	0.51	0/1053
28	i	0.26	0/1250	0.44	0/1698
29	k	0.24	0/2646	0.46	0/3579
30	l	0.27	0/896	0.51	0/1200

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
31	m	0.22	0/647	0.46	0/890
32	n	0.19	0/653	0.38	0/882
33	o	0.24	0/1035	0.46	0/1398
34	p	0.20	0/1085	0.46	0/1467
35	q	0.26	0/1171	0.48	0/1579
36	r	0.25	0/874	0.50	0/1188
37	s	0.20	0/1072	0.47	0/1436
38	t	0.23	0/1573	0.49	0/2130
39	u	0.20	0/590	0.47	0/810
40	v	0.21	0/1361	0.46	0/1861
41	w	0.23	0/872	0.53	0/1185
42	x	0.20	0/425	0.39	0/576
43	y	0.21	0/449	0.47	0/605
44	z	0.34	0/591	0.59	0/795
All	All	0.28	0/67728	0.53	9/91829 (0.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
3	3	0	2
4	4	0	1
10	J	0	2
All	All	0	5

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
23	d	112	LYS	CA-C-N	6.96	124.62	120.24
23	d	112	LYS	C-N-CA	6.96	124.62	120.24
3	3	366	THR	CA-C-N	6.03	133.06	121.54
3	3	366	THR	C-N-CA	6.03	133.06	121.54
9	H	91	MET	CA-C-N	5.77	140.86	127.00
9	H	91	MET	C-N-CA	5.77	140.86	127.00
9	H	91	MET	C-N-CD	-5.20	109.16	120.60
5	5	57	VAL	CA-C-N	5.01	123.40	120.24
5	5	57	VAL	C-N-CA	5.01	123.40	120.24

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	3	259	ASN	Peptide
3	3	366	THR	Peptide
4	4	275	TYR	Peptide
10	J	115	ILE	Peptide
10	J	82	VAL	Peptide

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	1	3312	0	3269	52	0
2	2	1655	0	1668	22	0
3	3	5275	0	5300	70	0
4	4	3457	0	3398	59	0
5	5	1726	0	1676	25	0
6	6	1247	0	1259	22	0
7	9	1414	0	1371	29	0
8	A	922	0	953	19	0
9	H	2528	0	2641	54	0
10	J	1344	0	1364	27	0
11	K	749	0	793	21	0
12	L	4807	0	4949	70	0
13	M	3647	0	3849	57	0
14	N	2723	0	2930	44	0
15	V	1028	0	1036	11	0
16	W	1155	0	1177	15	0
17	X	701	0	692	6	0
17	j	660	0	663	2	0
18	Y	1403	0	1392	15	0
19	Z	1441	0	1419	18	0
20	a	371	0	344	8	0
21	b	737	0	710	6	0
22	c	1024	0	1023	12	0
23	d	2748	0	2763	31	0
24	e	691	0	706	7	0
25	f	917	0	958	7	0
26	g	969	0	980	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
27	h	769	0	780	11	0
28	i	1209	0	1182	13	0
29	k	2596	0	2559	23	0
30	l	874	0	869	10	0
31	m	626	0	635	4	0
32	n	634	0	616	4	0
33	o	1004	0	995	16	0
34	p	1059	0	1062	13	0
35	q	1142	0	1137	14	0
36	r	846	0	864	13	0
37	s	1047	0	1013	16	0
38	t	1520	0	1477	16	0
39	u	563	0	509	5	0
40	v	1307	0	1207	18	0
41	w	846	0	792	14	0
42	x	412	0	411	1	0
43	y	436	0	437	3	0
44	z	576	0	570	7	0
45	1	8	0	0	2	0
45	3	16	0	0	1	0
45	6	8	0	0	1	0
45	9	16	0	0	0	0
46	1	31	0	19	0	0
47	1	44	0	27	3	0
48	2	4	0	0	1	0
48	3	4	0	0	0	0
49	3	1	0	0	0	0
50	4	23	0	30	1	0
50	H	23	0	30	2	0
51	6	46	0	69	2	0
51	9	54	0	88	1	0
51	L	54	0	88	3	0
51	M	54	0	88	3	0
51	w	54	0	88	8	0
52	A	51	0	82	3	0
52	J	91	0	136	8	0
52	L	82	0	118	2	0
52	M	44	0	65	0	0
52	N	91	0	136	1	0
52	V	37	0	48	2	0
52	i	102	0	164	5	0
52	o	31	0	36	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
52	p	27	0	27	1	0
53	L	185	0	276	11	0
53	M	100	0	156	13	0
53	V	94	0	141	6	0
53	W	100	0	156	4	0
53	h	58	0	60	0	0
53	o	90	0	133	9	0
53	x	75	0	97	3	0
54	X	31	0	34	3	0
54	g	34	0	40	0	0
55	b	1	0	0	0	0
56	d	48	0	26	2	0
57	k	23	0	12	1	0
58	s	15	0	27	0	0
All	All	67967	0	68895	767	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 6.

All (767) 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)
29:k:232:GLU:O	29:k:236:GLU:HB2	1.87	0.74
6:6:81:ARG:NH1	8:A:33:LYS:O	2.22	0.73
6:6:171:LYS:HG2	6:6:174:ARG:HH11	1.59	0.67
4:4:104:ASP:HB3	4:4:190:HIS:HD1	1.60	0.66
16:W:124:GLN:HB2	33:o:4:GLY:HA3	1.77	0.66
41:w:25:ARG:HG3	41:w:26:TRP:HD1	1.61	0.66
19:Z:119:SER:HB3	37:s:55:ARG:HH22	1.61	0.65
35:q:81:ARG:NH1	35:q:85:MET:SD	2.70	0.65
7:9:92:ILE:HG12	7:9:111:ILE:HG12	1.79	0.64
12:L:123:LEU:HD13	53:L:1005:CDL:H712	1.79	0.64
44:z:34:ARG:NH2	44:z:61:TYR:O	2.31	0.63
1:1:101:GLU:HB2	47:1:503:NAI:H42N	1.81	0.63
7:9:43:ARG:HG2	27:h:11:ARG:HD2	1.79	0.63
14:N:106:LEU:HD22	14:N:187:MET:HE2	1.79	0.63
3:3:175:THR:HG21	3:3:186:TYR:HB2	1.80	0.63
4:4:405:MET:SD	4:4:421:GLN:NE2	2.72	0.62
52:J:201:3PE:H2F2	52:J:201:3PE:H371	1.81	0.62
14:N:243:LEU:HD22	14:N:330:THR:HG21	1.80	0.62
18:Y:165:ARG:NH1	33:o:87:ASP:OD1	2.32	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:t:153:LEU:HD11	38:t:165:LEU:HD12	1.81	0.62
53:L:1005:CDL:H642	53:L:1005:CDL:H562	1.81	0.62
37:s:33:ARG:HH12	37:s:103:ARG:HH22	1.48	0.62
12:L:100:ILE:HG21	12:L:246:LEU:HB2	1.82	0.61
3:3:118:ASP:OD2	22:c:46:GLN:NE2	2.33	0.61
12:L:370:THR:HG23	12:L:431:LEU:HD13	1.82	0.61
5:5:129:TRP:O	5:5:132:ARG:HB3	2.00	0.61
14:N:140:SER:HA	30:l:1:PRO:HD2	1.82	0.61
3:3:360:SER:O	3:3:365:ASN:ND2	2.33	0.61
4:4:190:HIS:HD2	6:6:150:PRO:HD3	1.65	0.60
1:1:132:ARG:NH2	20:a:64:PRO:O	2.34	0.60
13:M:231:LEU:HA	13:M:235:LEU:HB2	1.81	0.60
1:1:355:LYS:HD3	1:1:373:ASN:HD22	1.65	0.60
9:H:97:ASN:H	35:q:143:THR:HG22	1.67	0.60
11:K:64:LEU:O	11:K:67:ALA:HB3	2.02	0.60
3:3:252:PRO:HG3	3:3:263:ILE:HG12	1.83	0.60
12:L:221:THR:HG23	12:L:226:GLN:HB2	1.81	0.59
4:4:51:PHE:HB3	4:4:64:LEU:HB2	1.84	0.59
9:H:169:GLN:NE2	9:H:241:LEU:O	2.35	0.59
4:4:58:ALA:HB1	4:4:62:LEU:HB3	1.83	0.59
19:Z:68:ARG:NH1	43:y:43:LEU:O	2.36	0.59
12:L:103:PHE:HB2	12:L:341:MET:HE3	1.85	0.59
4:4:101:PRO:O	4:4:105:ARG:NH1	2.36	0.59
5:5:151:ILE:HG23	5:5:152:LEU:HG	1.84	0.59
13:M:204:MET:HE1	13:M:302:LEU:HD11	1.85	0.59
1:1:21:ILE:O	1:1:117:LYS:NZ	2.36	0.59
12:L:161:ARG:HA	38:t:91:GLU:HG3	1.84	0.59
28:i:26:VAL:O	28:i:30:ALA:HB3	2.03	0.59
3:3:194:GLU:HG3	3:3:389:PRO:HB3	1.85	0.58
6:6:81:ARG:NH2	8:A:35:SER:O	2.36	0.58
15:V:39:SER:OG	15:V:54:ARG:NH2	2.36	0.58
12:L:73:THR:HB	12:L:194:ASN:HD21	1.68	0.58
43:y:8:ARG:HG3	43:y:10:HIS:H	1.68	0.58
4:4:63:ARG:HB3	4:4:79:HIS:HB2	1.86	0.58
4:4:343:GLU:O	4:4:347:HIS:ND1	2.33	0.58
12:L:106:TRP:HD1	12:L:447:ASN:HD22	1.52	0.58
23:d:48:PRO:HA	23:d:71:MET:O	2.03	0.58
9:H:20:LEU:HD21	9:H:231:ILE:HD11	1.85	0.58
12:L:161:ARG:NH2	38:t:90:TYR:O	2.36	0.58
12:L:233:LEU:HD23	12:L:307:SER:HB3	1.85	0.58
14:N:30:TRP:NE1	14:N:67:SER:OG	2.36	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:X:65:ILE:HD13	38:t:3:LEU:HD22	1.86	0.58
18:Y:29:HIS:HB3	18:Y:119:PRO:HD2	1.84	0.58
25:f:49:GLN:HE22	27:h:92:LYS:HA	1.69	0.58
28:i:34:ARG:NH2	28:i:58:ARG:O	2.36	0.58
2:2:150:ASN:HB3	2:2:162:GLU:HB3	1.84	0.58
4:4:269:LEU:HB2	4:4:368:GLU:HB2	1.86	0.58
6:6:36:LEU:HD13	52:i:501:3PE:H392	1.84	0.58
23:d:169:ALA:HB2	23:d:231:VAL:HG12	1.86	0.58
20:a:57:ASP:OD1	20:a:60:LYS:NZ	2.36	0.57
3:3:319:ALA:HB3	3:3:345:THR:HG22	1.86	0.57
8:A:67:LEU:HD11	11:K:68:ALA:HB3	1.87	0.57
10:J:117:PHE:HE1	14:N:91:ASN:HD21	1.52	0.57
4:4:179:GLU:OE2	6:6:66:ARG:NH1	2.34	0.57
52:J:202:3PE:H11	11:K:23:ARG:HG2	1.87	0.57
18:Y:35:CYS:O	18:Y:39:ASN:ND2	2.38	0.56
17:j:37:MET:HE1	17:j:44:SER:HA	1.87	0.56
17:j:48:VAL:HG12	17:j:52:MET:HE2	1.87	0.56
12:L:231:PRO:HB3	12:L:530:PRO:HG3	1.87	0.56
53:L:1004:CDL:H141	15:V:100:LEU:HD11	1.87	0.56
14:N:159:ILE:HG21	14:N:278:LEU:HD11	1.87	0.56
29:k:232:GLU:O	29:k:236:GLU:CB	2.53	0.56
37:s:7:ARG:HH21	37:s:17:GLU:HG3	1.70	0.56
19:Z:67:PHE:HD1	43:y:43:LEU:HD21	1.70	0.56
4:4:20:TYR:OH	14:N:295:ARG:NH2	2.38	0.56
4:4:184:VAL:O	7:9:60:ARG:NH1	2.38	0.56
6:6:77:ARG:HH21	50:H:501:DCQ:H4MA	1.71	0.56
3:3:362:TYR:HA	3:3:492:ILE:HD12	1.87	0.56
3:3:453:LEU:HD13	3:3:470:VAL:HG21	1.87	0.56
3:3:632:ARG:NH1	24:e:57:CYS:SG	2.79	0.56
1:1:342:ASP:OD1	1:1:429:ARG:NH2	2.37	0.56
52:A:201:3PE:H381	9:H:295:PRO:HB3	1.87	0.56
53:L:1005:CDL:H542	53:L:1005:CDL:H722	1.87	0.56
19:Z:5:ASP:OD2	41:w:98:LYS:NZ	2.39	0.56
1:1:200:GLN:NE2	3:3:174:THR:O	2.39	0.56
7:9:69:ARG:NH1	7:9:73:GLY:O	2.39	0.56
10:J:67:VAL:HG11	11:K:31:LEU:HD21	1.86	0.56
13:M:243:MET:HB3	13:M:301:ILE:HG21	1.88	0.56
5:5:150:ARG:NH1	5:5:153:THR:OG1	2.38	0.55
1:1:89:ARG:NH1	1:1:217:GLY:O	2.39	0.55
12:L:316:THR:HA	12:L:319:ILE:HG12	1.88	0.55
13:M:36:LEU:HB2	41:w:69:VAL:HG22	1.87	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
18:Y:118:ARG:NH2	35:q:62:GLU:OE2	2.40	0.55
25:f:69:GLU:HG3	27:h:102:ARG:HH12	1.71	0.55
10:J:5:ILE:HG21	35:q:141:TRP:HZ2	1.71	0.55
13:M:446:LEU:HB3	51:w:801:PC1:H332	1.88	0.55
23:d:167:LYS:HB2	23:d:229:ALA:HA	1.87	0.55
27:h:45:SER:O	27:h:51:ASN:ND2	2.37	0.55
9:H:114:TYR:OH	10:J:61:LEU:O	2.24	0.55
10:J:52:LEU:HD22	10:J:143:ILE:HD11	1.89	0.55
3:3:201:ASP:OD2	3:3:268:ARG:NH2	2.36	0.55
28:i:60:ARG:HH22	28:i:95:ASP:HA	1.71	0.55
34:p:74:ASN:HD22	40:v:10:PRO:HB2	1.72	0.54
3:3:46:LEU:O	22:c:116:LYS:NZ	2.37	0.54
6:6:107:MET:O	6:6:111:ARG:NH1	2.40	0.54
8:A:112:GLU:OE2	9:H:291:LYS:NZ	2.40	0.54
34:p:47:GLN:HG3	38:t:165:LEU:HD13	1.88	0.54
34:p:81:PRO:HB2	52:p:201:3PE:H222	1.90	0.54
3:3:332:LYS:NZ	3:3:505:LEU:O	2.38	0.54
4:4:352:TYR:HD1	7:9:86:VAL:HG21	1.71	0.54
5:5:74:SER:HB3	5:5:97:LEU:HB3	1.90	0.54
16:W:28:TYR:O	41:w:67:SER:OG	2.24	0.54
36:r:85:LEU:HD22	36:r:96:ILE:HD11	1.89	0.54
1:1:108:ARG:NH2	2:2:162:GLU:OE2	2.40	0.54
12:L:83:ASP:OD2	12:L:262:ARG:NH1	2.40	0.54
53:L:1005:CDL:H531	51:w:801:PC1:H2I1	1.90	0.54
53:x:101:CDL:H131	53:x:101:CDL:H522	1.90	0.54
5:5:65:ARG:NH1	5:5:123:VAL:O	2.41	0.54
1:1:378:ARG:NE	3:3:132:GLU:OE2	2.41	0.54
5:5:33:LEU:HD13	5:5:60:VAL:HG22	1.88	0.54
9:H:24:GLU:HG2	9:H:271:LEU:HD22	1.89	0.54
9:H:111:LEU:HD22	10:J:57:PHE:HZ	1.73	0.54
6:6:118:SER:N	45:6:201:SF4:S4	2.81	0.54
10:J:126:VAL:HG23	10:J:127:ILE:HG23	1.90	0.54
13:M:126:LEU:HD21	13:M:153:THR:HG21	1.90	0.54
5:5:121:VAL:HG21	5:5:146:PRO:HD3	1.90	0.54
9:H:81:LEU:HA	9:H:84:THR:HG22	1.90	0.54
12:L:69:LEU:HD13	13:M:451:PRO:HG2	1.90	0.54
33:o:45:ASP:OD2	33:o:51:ARG:NH2	2.41	0.54
3:3:261:GLU:OE1	22:c:42:ARG:NH2	2.40	0.53
13:M:447:LEU:HD11	51:w:801:PC1:H3F1	1.89	0.53
2:2:105:THR:OG1	48:2:300:FES:S2	2.67	0.53
5:5:119:SER:OG	5:5:131:GLU:OE2	2.25	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:9:145:GLU:HA	7:9:148:LEU:HD13	1.91	0.53
8:A:73:LEU:HD12	9:H:151:LEU:HD13	1.90	0.53
12:L:90:VAL:HG22	12:L:129:LEU:HD22	1.91	0.53
2:2:74:GLN:O	20:a:74:ARG:NH2	2.41	0.53
9:H:141:SER:HA	9:H:290:TRP:HE1	1.74	0.53
13:M:5:ILE:HG23	13:M:104:LEU:HD11	1.89	0.53
14:N:207:ILE:HG22	14:N:211:MET:HE2	1.90	0.53
4:4:388:ARG:HG2	5:5:81:VAL:HG22	1.90	0.53
9:H:87:ILE:HD12	9:H:96:ILE:HD12	1.89	0.53
10:J:84:VAL:HG12	10:J:90:LEU:HD13	1.89	0.53
1:1:428:GLU:HA	1:1:431:GLN:HG2	1.91	0.53
4:4:149:ASN:HD21	4:4:371:LYS:HG3	1.74	0.53
23:d:185:ILE:HD13	23:d:191:VAL:HG23	1.90	0.53
9:H:120:GLY:HA3	9:H:132:ALA:HB2	1.89	0.53
12:L:421:ILE:HG12	12:L:501:ALA:HB2	1.90	0.53
1:1:139:ARG:NH2	2:2:144:CYS:O	2.32	0.53
2:2:10:ARG:NH2	3:3:138:GLU:OE2	2.42	0.53
12:L:142:ILE:HG12	13:M:370:PRO:HB2	1.90	0.53
7:9:64:GLU:OE1	7:9:136:ASN:ND2	2.42	0.53
12:L:536:LEU:HG	12:L:540:MET:HE2	1.91	0.53
29:k:45:LYS:HB2	29:k:235:VAL:HG21	1.90	0.53
5:5:77:ASP:OD2	5:5:79:THR:OG1	2.27	0.52
2:2:147:ALA:HB3	2:2:153:MET:HG2	1.91	0.52
4:4:335:ARG:NH2	7:9:129:ASP:OD1	2.43	0.52
9:H:184:MET:SD	9:H:297:THR:OG1	2.68	0.52
10:J:23:LYS:NZ	11:K:18:GLY:O	2.40	0.52
10:J:84:VAL:HG23	23:d:325:ARG:HH22	1.75	0.52
19:Z:5:ASP:HB3	19:Z:8:VAL:HG12	1.90	0.52
1:1:270:GLU:OE1	1:1:283:HIS:NE2	2.43	0.52
13:M:64:PRO:HB3	13:M:454:ILE:HG22	1.90	0.52
25:f:34:LEU:HD12	25:f:37:ILE:HD12	1.90	0.52
1:1:365:CYS:HB3	45:1:501:SF4:S3	2.48	0.52
29:k:170:ILE:HD11	29:k:238:ILE:HD11	1.91	0.52
3:3:193:SER:H	3:3:196:SER:HB3	1.75	0.52
8:A:57:LEU:HD21	10:J:169:MET:HG2	1.92	0.52
3:3:126:ASP:HB2	4:4:328:ALA:HB3	1.92	0.52
14:N:45:MET:HE3	14:N:53:THR:HG22	1.91	0.52
39:u:40:PHE:O	39:u:44:SER:OG	2.27	0.52
14:N:257:LEU:HD21	53:o:502:CDL:H861	1.92	0.52
3:3:163:ALA:HA	3:3:167:ALA:HB3	1.91	0.52
29:k:145:ARG:NH1	29:k:147:GLN:OE1	2.43	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:3:632:ARG:HH12	24:e:59:ASP:H	1.58	0.52
29:k:81:LYS:HD2	29:k:92:ASN:HD22	1.75	0.52
1:1:213:VAL:HG13	1:1:217:GLY:HA2	1.92	0.51
1:1:289:GLY:HA3	1:1:293:ASN:HD22	1.75	0.51
29:k:30:ASN:N	29:k:33:SER:OG	2.43	0.51
2:2:53:LEU:HD12	20:a:52:LEU:HD13	1.93	0.51
3:3:443:LEU:HD13	3:3:477:ILE:HD11	1.91	0.51
4:4:3:GLN:NE2	13:M:135:ARG:O	2.42	0.51
14:N:160:LEU:HD21	53:V:202:CDL:H191	1.93	0.51
25:f:37:ILE:O	25:f:44:ARG:NH1	2.43	0.51
27:h:5:ARG:HA	27:h:8:GLN:HB3	1.92	0.51
2:2:149:VAL:O	2:2:190:ARG:NH2	2.42	0.51
3:3:255:HIS:NE2	3:3:634:ASP:OD1	2.43	0.51
4:4:224:GLU:OE2	7:9:40:TYR:OH	2.22	0.51
12:L:73:THR:OG1	34:p:127:SER:OG	2.28	0.51
12:L:487:LYS:HE3	39:u:49:GLY:HA2	1.93	0.51
53:M:503:CDL:H792	14:N:338:PRO:HG3	1.91	0.51
29:k:173:ASP:HB2	29:k:224:SER:HA	1.91	0.51
40:v:54:SER:HA	40:v:84:TYR:HB3	1.91	0.51
52:A:201:3PE:H2H1	52:A:201:3PE:H292	1.93	0.51
3:3:140:LYS:HD3	3:3:150:MET:HG3	1.93	0.51
4:4:50:ASN:HB3	8:A:38:GLU:HG3	1.92	0.51
41:w:114:ASP:HB3	41:w:117:LYS:HG2	1.92	0.51
4:4:145:THR:OG1	4:4:181:TYR:OH	2.28	0.51
51:9:401:PC1:H351	9:H:187:ILE:HD12	1.92	0.51
12:L:137:LEU:HD21	12:L:263:PHE:HZ	1.76	0.51
29:k:51:PHE:HB3	29:k:107:GLN:HE21	1.75	0.51
3:3:110:GLN:NE2	3:3:113:GLU:O	2.43	0.51
5:5:66:ASP:O	25:f:82:GLN:NE2	2.44	0.51
13:M:6:ILE:HG23	53:M:503:CDL:H202	1.92	0.51
28:i:44:TYR:OH	28:i:113:HIS:N	2.42	0.51
41:w:92:GLU:OE2	41:w:95:ARG:NH2	2.40	0.51
12:L:97:THR:HG21	12:L:125:LEU:HD22	1.92	0.51
53:M:503:CDL:H852	53:o:502:CDL:H673	1.92	0.51
14:N:238:PRO:HB2	53:o:502:CDL:HB4	1.92	0.51
16:W:120:GLY:O	16:W:124:GLN:NE2	2.42	0.51
5:5:96:LEU:HB2	5:5:105:ILE:HG22	1.91	0.50
7:9:74:GLU:HB2	21:b:44:ILE:HD13	1.93	0.50
9:H:46:LEU:HD21	52:i:501:3PE:H272	1.92	0.50
14:N:115:VAL:HG12	14:N:180:ALA:HB1	1.93	0.50
19:Z:41:ASP:OD1	36:r:86:LYS:NZ	2.41	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:d:100:GLU:HB2	23:d:105:ASP:HA	1.94	0.50
28:i:50:GLU:HG3	28:i:89:TRP:HZ2	1.76	0.50
3:3:499:GLN:HE21	3:3:503:LEU:HD11	1.76	0.50
9:H:165:LEU:HD21	9:H:241:LEU:HA	1.93	0.50
12:L:152:PHE:HB2	12:L:172:ILE:HD11	1.93	0.50
16:W:133:ILE:HG23	30:l:21:SER:HB3	1.93	0.50
34:p:47:GLN:HA	34:p:50:TYR:HB3	1.93	0.50
3:3:119:GLN:HB2	3:3:123:PHE:HD2	1.76	0.50
3:3:382:THR:HG23	3:3:384:PRO:HD3	1.94	0.50
5:5:82:ASP:OD2	5:5:89:ARG:NH2	2.43	0.50
3:3:222:THR:HA	3:3:242:THR:O	2.11	0.50
4:4:141:PHE:O	4:4:145:THR:OG1	2.29	0.50
8:A:98:LEU:HD22	9:H:298:LEU:HD11	1.94	0.50
29:k:52:PRO:O	29:k:107:GLN:NE2	2.45	0.50
31:m:83:LEU:HD13	35:q:54:ARG:HD3	1.94	0.50
40:v:56:GLN:O	40:v:70:ARG:NH2	2.45	0.50
10:J:40:GLY:HA2	10:J:43:ILE:HD12	1.94	0.50
19:Z:159:LYS:NZ	33:o:112:VAL:O	2.45	0.50
4:4:109:VAL:HG11	4:4:152:MET:HG2	1.94	0.50
15:V:65:ILE:HD11	15:V:100:LEU:HD23	1.93	0.50
18:Y:97:ARG:HD3	35:q:59:LEU:HD13	1.92	0.50
19:Z:140:ASP:O	19:Z:161:ARG:NH1	2.44	0.50
4:4:300:ARG:NH2	4:4:420:THR:O	2.44	0.50
12:L:570:GLN:OE1	14:N:167:TRP:NE1	2.42	0.50
13:M:155:ALA:HB1	51:M:502:PC1:H2H2	1.94	0.50
1:1:94:VAL:O	1:1:222:VAL:HA	2.12	0.50
12:L:67:HIS:NE2	12:L:70:THR:OG1	2.43	0.50
13:M:232:ALA:O	13:M:237:LYS:NZ	2.44	0.50
53:M:503:CDL:H822	53:o:502:CDL:H673	1.94	0.50
17:X:69:LYS:HA	36:r:2:GLY:HA2	1.94	0.50
23:d:91:VAL:HG23	23:d:126:VAL:HG11	1.93	0.50
23:d:280:THR:HG23	23:d:283:LYS:H	1.77	0.50
34:p:62:PRO:O	34:p:66:ARG:HB2	2.12	0.50
3:3:323:VAL:HG11	3:3:525:LEU:HD13	1.93	0.50
50:4:501:DCQ:H8	6:6:56:ALA:HB1	1.94	0.49
13:M:9:MET:HE2	53:M:503:CDL:H842	1.94	0.49
28:i:75:TRP:NE1	52:i:502:3PE:O12	2.44	0.49
1:1:394:GLU:OE1	3:3:129:ARG:NH1	2.45	0.49
4:4:200:HIS:NE2	7:9:124:GLU:OE1	2.45	0.49
7:9:162:GLU:HB3	28:i:109:ILE:HG12	1.93	0.49
8:A:67:LEU:HD22	11:K:65:VAL:HA	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:L:350:LEU:HD11	12:L:362:LEU:HD11	1.94	0.49
23:d:97:ARG:NH1	56:d:401:NDP:O1A	2.45	0.49
23:d:98:GLU:HG3	23:d:177:ARG:HE	1.77	0.49
23:d:136:ASN:HA	23:d:292:ARG:HE	1.77	0.49
6:6:52:LEU:HB2	6:6:90:GLY:HA3	1.93	0.49
13:M:180:GLN:HA	13:M:248:LEU:HD21	1.94	0.49
3:3:101:HIS:HD2	4:4:342:MET:HE2	1.77	0.49
3:3:273:GLY:O	3:3:549:HIS:NE2	2.36	0.49
4:4:190:HIS:CD2	6:6:150:PRO:HD3	2.47	0.49
4:4:354:GLU:OE2	4:4:357:GLN:NE2	2.45	0.49
6:6:35:ASP:HB3	52:i:501:3PE:H331	1.93	0.49
9:H:230:ASN:HA	9:H:233:MET:HE3	1.95	0.49
30:l:105:PRO:HG3	44:z:53:ARG:HE	1.78	0.49
1:1:184:TYR:HB3	1:1:357:GLU:HB3	1.95	0.49
12:L:532:ILE:HD11	40:v:101:MET:HB3	1.93	0.49
13:M:369:LEU:HD12	13:M:370:PRO:HD2	1.95	0.49
2:2:156:ILE:HD13	2:2:176:LEU:HD11	1.95	0.49
3:3:194:GLU:HG2	3:3:195:LEU:HG	1.94	0.49
7:9:69:ARG:NH2	21:b:38:ASN:O	2.40	0.49
33:o:66:THR:OG1	42:x:28:TRP:NE1	2.45	0.49
1:1:144:ASN:O	1:1:148:ASN:HB2	2.13	0.49
10:J:172:THR:HA	11:K:80:MET:HE3	1.95	0.49
13:M:18:SER:HB2	13:M:23:ILE:HG22	1.95	0.49
3:3:27:LEU:HA	3:3:37:ILE:HD12	1.95	0.49
3:3:198:ASN:OD1	3:3:268:ARG:NH2	2.38	0.49
12:L:316:THR:HG23	12:L:325:ALA:HB2	1.95	0.49
16:W:11:ILE:HG12	38:t:103:LEU:HD11	1.95	0.49
21:b:32:GLN:HE22	21:b:34:GLU:HB2	1.78	0.49
53:x:101:CDL:H541	53:x:101:CDL:H741	1.94	0.49
16:W:43:ILE:HG12	16:W:47:ILE:HD12	1.95	0.48
28:i:78:ASP:HB3	28:i:81:MET:HG3	1.94	0.48
4:4:183:ARG:NH1	4:4:210:ASP:OD2	2.38	0.48
5:5:138:PHE:O	5:5:163:ARG:NE	2.44	0.48
1:1:367:GLU:OE1	3:3:100:ASN:ND2	2.43	0.48
3:3:324:ASP:N	3:3:324:ASP:OD1	2.45	0.48
5:5:60:VAL:O	5:5:63:PHE:HB3	2.13	0.48
53:L:1005:CDL:H391	53:L:1005:CDL:H171	1.94	0.48
13:M:54:LEU:HD23	16:W:93:ILE:HG23	1.95	0.48
53:M:503:CDL:H622	53:M:503:CDL:H251	1.95	0.48
29:k:31:ILE:HG23	57:k:501:AMP:H3'	1.95	0.48
3:3:99:ALA:O	3:3:134:LYS:NZ	2.46	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:3:396:ARG:NH1	3:3:416:THR:O	2.46	0.48
19:Z:15:ARG:HB3	36:r:109:ILE:HD12	1.96	0.48
29:k:108:TYR:OH	29:k:164:LEU:O	2.23	0.48
37:s:6:ARG:NH1	37:s:15:GLU:OE2	2.46	0.48
3:3:323:VAL:HG21	3:3:525:LEU:HB3	1.95	0.48
13:M:68:LEU:HB2	51:w:801:PC1:H3G1	1.94	0.48
23:d:194:ILE:H	23:d:288:HIS:CE1	2.32	0.48
1:1:69:GLY:O	47:1:503:NAI:H2N	2.13	0.48
13:M:301:ILE:HA	13:M:309:TYR:HE1	1.78	0.48
15:V:1:AYA:HA	15:V:4:LEU:HD23	1.95	0.48
15:V:1:AYA:HM2	15:V:3:THR:HG22	1.96	0.48
19:Z:81:VAL:HA	19:Z:84:MET:HG2	1.94	0.48
4:4:149:ASN:HD22	4:4:370:PRO:HB2	1.79	0.48
9:H:154:LEU:HD22	9:H:157:ASN:HD22	1.79	0.48
10:J:125:TRP:HB2	35:q:136:THR:HG21	1.96	0.48
1:1:194:GLU:OE2	1:1:204:ARG:NE	2.35	0.48
1:1:295:LEU:HB2	1:1:339:ARG:HA	1.96	0.48
3:3:140:LYS:HB2	3:3:148:THR:HG21	1.95	0.48
12:L:203:MET:HB2	19:Z:113:GLN:HG3	1.96	0.48
14:N:309:ASN:ND2	29:k:75:GLY:O	2.39	0.48
38:t:137:LYS:HA	38:t:140:GLN:HG2	1.96	0.48
4:4:155:THR:HB	4:4:167:PHE:HA	1.96	0.48
12:L:63:ILE:O	12:L:79:SER:HA	2.14	0.48
19:Z:19:PRO:HD3	36:r:110:LEU:HD11	1.96	0.48
4:4:64:LEU:HD11	4:4:418:ILE:HD11	1.96	0.47
5:5:192:GLN:OE1	7:9:115:LYS:NZ	2.47	0.47
14:N:246:ILE:HG21	53:o:502:CDL:H591	1.97	0.47
16:W:138:LYS:HA	30:l:27:ARG:HB3	1.95	0.47
21:b:51:GLN:OE1	21:b:92:ARG:NH1	2.47	0.47
9:H:205:SER:H	9:H:279:ARG:HH22	1.61	0.47
13:M:158:LEU:HD23	14:N:283:ALA:HB1	1.96	0.47
3:3:601:ARG:NH2	3:3:614:ASP:OD1	2.47	0.47
9:H:156:MET:HE3	9:H:304:HIS:HD2	1.79	0.47
10:J:71:THR:O	10:J:76:THR:OG1	2.30	0.47
23:d:98:GLU:OE1	23:d:286:ARG:NH2	2.38	0.47
4:4:84:HIS:NE2	6:6:130:TYR:OH	2.37	0.47
18:Y:5:GLU:O	18:Y:53:ARG:NH1	2.47	0.47
40:v:52:ASP:OD1	40:v:78:HIS:NE2	2.35	0.47
3:3:45:ARG:NH2	3:3:256:GLU:OE2	2.39	0.47
9:H:251:MET:HB3	9:H:254:LEU:HD13	1.96	0.47
12:L:402:SER:HG	12:L:404:THR:HG1	1.58	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:4:61:VAL:HG21	4:4:83:LEU:HB2	1.96	0.47
12:L:593:ILE:HD12	15:V:41:ALA:HB2	1.97	0.47
13:M:229:MET:O	13:M:233:ALA:HB3	2.14	0.47
40:v:136:ASN:HA	40:v:153:VAL:HB	1.97	0.47
1:1:305:PRO:HG3	1:1:413:TRP:HB3	1.97	0.47
2:2:121:GLN:HE21	2:2:128:VAL:HG23	1.80	0.47
6:6:71:ARG:HA	9:H:37:PRO:HA	1.96	0.47
11:K:21:MET:HG2	12:L:589:LEU:HD23	1.96	0.47
11:K:46:LEU:O	11:K:50:ASN:HB2	2.15	0.47
12:L:72:GLN:NE2	13:M:458:LEU:O	2.47	0.47
13:M:51:ASN:HA	13:M:57:PHE:HB2	1.97	0.47
19:Z:136:LYS:NZ	19:Z:140:ASP:OD2	2.43	0.47
2:2:121:GLN:HG2	2:2:127:LYS:HA	1.97	0.47
4:4:73:VAL:HG21	4:4:414:VAL:HG21	1.95	0.47
23:d:26:ALA:HA	23:d:31:GLY:HA3	1.96	0.47
29:k:213:GLU:O	29:k:217:LYS:NZ	2.46	0.47
5:5:137:MET:HA	5:5:161:PRO:HD2	1.96	0.47
12:L:180:ILE:HD13	52:L:1001:3PE:H2H2	1.97	0.47
16:W:117:LYS:NZ	33:o:100:ASP:OD2	2.39	0.47
3:3:364:LEU:HD12	3:3:491:ASN:HB3	1.97	0.47
9:H:311:LEU:O	35:q:54:ARG:NH2	2.45	0.47
12:L:279:CYS:SG	12:L:405:ASN:ND2	2.88	0.47
38:t:134:ARG:HA	38:t:137:LYS:HE3	1.97	0.47
40:v:80:ASP:HB3	40:v:83:MET:HB3	1.96	0.47
2:2:40:GLU:HB2	20:a:67:SER:HA	1.97	0.46
23:d:176:ASP:OD1	23:d:180:ASN:ND2	2.48	0.46
29:k:97:GLN:NE2	29:k:131:ASP:OD2	2.48	0.46
1:1:361:GLN:N	45:1:501:SF4:S4	2.86	0.46
4:4:145:THR:HG1	4:4:181:TYR:HH	1.56	0.46
13:M:116:ILE:HD11	13:M:161:LEU:HD12	1.97	0.46
16:W:138:LYS:HB2	16:W:143:ASN:HD21	1.80	0.46
36:r:107:ASP:OD1	37:s:67:LYS:NZ	2.47	0.46
37:s:102:GLU:OE2	37:s:105:ARG:NH2	2.47	0.46
1:1:101:GLU:OE2	1:1:302:SER:N	2.48	0.46
3:3:478:ARG:NH2	3:3:485:GLY:O	2.48	0.46
10:J:99:MET:HE1	52:J:202:3PE:H2B2	1.98	0.46
33:o:45:ASP:OD1	33:o:49:ARG:NH2	2.48	0.46
13:M:352:LEU:HB3	13:M:355:MET:HB3	1.98	0.46
14:N:200:MET:HG3	14:N:269:GLU:HG3	1.97	0.46
15:V:34:VAL:HG11	53:V:202:CDL:H402	1.95	0.46
3:3:365:ASN:N	3:3:491:ASN:OD1	2.49	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:3:426:PRO:HB2	3:3:656:LEU:HD13	1.98	0.46
3:3:569:LYS:HG3	3:3:571:ALA:HB2	1.97	0.46
3:3:12:PHE:HB2	3:3:78:ASN:HA	1.97	0.46
3:3:356:THR:HG21	3:3:503:LEU:HD22	1.97	0.46
4:4:412:ALA:HB3	9:H:281:ARG:HG3	1.98	0.46
5:5:77:ASP:HB3	5:5:95:ASN:HD22	1.81	0.46
8:A:88:LEU:HD13	9:H:309:ILE:HD12	1.98	0.46
10:J:43:ILE:HG22	11:K:46:LEU:HD11	1.97	0.46
53:L:1005:CDL:H362	53:L:1005:CDL:H532	1.97	0.46
53:M:503:CDL:H741	18:Y:169:TRP:HE1	1.81	0.46
16:W:122:TRP:CD1	33:o:2:MET:HG2	2.50	0.46
22:c:34:LYS:NZ	22:c:102:SER:OG	2.43	0.46
24:e:41:VAL:HG12	24:e:45:LYS:HE2	1.98	0.46
10:J:167:VAL:HG22	14:N:42:PRO:HG2	1.97	0.46
12:L:10:VAL:HG21	36:r:78:VAL:HG22	1.98	0.46
38:t:169:ILE:O	38:t:172:ARG:NH1	2.45	0.46
3:3:127:ARG:NH2	4:4:326:ASP:O	2.49	0.46
16:W:114:MET:HE2	16:W:121:PRO:HD2	1.97	0.46
24:e:40:TYR:O	24:e:43:LEU:HB3	2.16	0.46
1:1:96:ASN:ND2	1:1:187:GLY:O	2.40	0.46
1:1:299:PRO:HG3	1:1:307:ILE:HG12	1.98	0.46
5:5:41:GLN:NE2	5:5:49:GLU:OE1	2.43	0.46
53:M:503:CDL:H872	53:M:503:CDL:H652	1.98	0.46
28:i:106:ARG:HB2	28:i:109:ILE:HG13	1.97	0.46
1:1:265:PRO:O	2:2:190:ARG:NH1	2.41	0.46
53:M:503:CDL:HA61	53:M:503:CDL:HA21	1.98	0.46
14:N:193:VAL:HB	14:N:266:ILE:HG23	1.98	0.46
23:d:244:TYR:HB2	23:d:337:ALA:HB2	1.98	0.46
29:k:19:THR:HG23	29:k:21:ASN:H	1.82	0.46
1:1:78:LYS:NZ	1:1:222:VAL:O	2.43	0.45
3:3:252:PRO:O	22:c:44:ASN:ND2	2.49	0.45
3:3:427:LYS:HD3	3:3:656:LEU:HD12	1.98	0.45
53:L:1005:CDL:H431	51:w:801:PC1:H261	1.97	0.45
14:N:331:VAL:HG13	14:N:335:MET:HE3	1.98	0.45
54:X:101:ZMP:H15	54:X:101:ZMP:H17	1.61	0.45
37:s:45:MET:HB2	37:s:50:LEU:HD12	1.98	0.45
38:t:138:GLN:HE22	38:t:156:ALA:HA	1.80	0.45
9:H:165:LEU:HD23	9:H:240:THR:HG22	1.98	0.45
11:K:93:LEU:HB3	14:N:54:GLU:HG3	1.97	0.45
14:N:246:ILE:HA	53:o:502:CDL:H873	1.97	0.45
17:X:23:ASP:HB2	32:n:43:PRO:HA	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
29:k:41:GLU:HA	29:k:44:GLU:HG2	1.98	0.45
6:6:63:ALA:HA	6:6:69:MET:HG2	1.98	0.45
11:K:56:ALA:O	11:K:59:MET:HB3	2.16	0.45
13:M:328:CYS:HB3	13:M:437:MET:HE1	1.98	0.45
23:d:268:ARG:HA	23:d:271:GLU:HG2	1.97	0.45
3:3:60:GLU:HG2	3:3:78:ASN:HB3	1.97	0.45
12:L:418:PHE:HD1	12:L:421:ILE:HD12	1.82	0.45
14:N:261:MET:HG3	14:N:340:THR:HG23	1.99	0.45
22:c:42:ARG:NH1	22:c:46:GLN:O	2.50	0.45
1:1:63:SER:HB2	1:1:242:PHE:HB3	1.98	0.45
12:L:432:LEU:HD21	32:n:65:ALA:HB2	1.99	0.45
13:M:23:ILE:HD11	13:M:92:LYS:HD2	1.99	0.45
41:w:34:ASP:OD1	41:w:34:ASP:N	2.50	0.45
1:1:21:ILE:HG12	1:1:233:THR:HG21	1.97	0.45
3:3:223:ARG:O	3:3:241:SER:HA	2.16	0.45
5:5:78:LEU:HB3	5:5:130:TYR:HB3	1.99	0.45
7:9:155:LEU:HB3	21:b:36:ASN:HD22	1.81	0.45
12:L:66:TRP:NE1	51:w:801:PC1:O12	2.39	0.45
13:M:105:PHE:O	13:M:109:THR:OG1	2.23	0.45
53:V:202:CDL:H351	53:V:202:CDL:H322	1.83	0.45
18:Y:43:MET:HE3	18:Y:47:TRP:HZ3	1.82	0.45
18:Y:129:LYS:NZ	31:m:65:VAL:O	2.39	0.45
22:c:28:GLU:O	22:c:32:THR:OG1	2.33	0.45
33:o:13:PHE:HE2	33:o:74:TYR:HB3	1.82	0.45
34:p:12:SER:OG	40:v:57:GLU:OE2	2.33	0.45
1:1:71:ALA:HB2	47:1:503:NAI:H4D	1.97	0.45
8:A:55:PHE:HB3	8:A:113:TRP:CE2	2.52	0.45
12:L:50:PRO:HA	12:L:53:MET:HG2	1.98	0.45
18:Y:89:ASP:OD2	44:z:34:ARG:NH1	2.49	0.45
22:c:42:ARG:NH2	22:c:44:ASN:OD1	2.49	0.45
4:4:110:SER:OG	4:4:149:ASN:ND2	2.50	0.45
9:H:86:TRP:CE2	9:H:233:MET:HB2	2.52	0.45
12:L:397:GLU:HB3	12:L:482:MET:HE1	1.98	0.45
53:L:1005:CDL:H621	13:M:372:SER:HB3	1.99	0.45
13:M:121:MET:HE1	53:M:503:CDL:H802	1.99	0.45
23:d:131:HIS:HB3	23:d:165:ILE:HG12	1.99	0.45
1:1:20:ARG:HD2	1:1:269:GLU:HB3	1.99	0.45
1:1:403:THR:HG21	1:1:408:GLY:HA3	1.98	0.45
7:9:27:TRP:HB3	7:9:30:LEU:HD12	1.99	0.45
53:W:201:CDL:H381	51:w:801:PC1:H251	1.99	0.45
22:c:8:THR:OG1	22:c:9:GLN:N	2.49	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:r:18:ARG:NH1	38:t:170:VAL:O	2.42	0.45
1:1:97:ALA:HB3	1:1:138:ILE:HA	1.99	0.45
11:K:37:MET:HG2	14:N:68:MET:HE1	1.98	0.45
51:M:502:PC1:H2C2	51:M:502:PC1:H3B1	1.99	0.45
37:s:4:LEU:HD11	40:v:124:THR:HA	1.99	0.45
40:v:71:LEU:HD21	40:v:77:MET:HG2	1.98	0.45
8:A:38:GLU:HB2	8:A:43:PRO:HG3	1.99	0.44
9:H:18:ALA:O	9:H:21:THR:OG1	2.29	0.44
52:J:202:3PE:H292	11:K:14:VAL:HG22	1.98	0.44
13:M:78:MET:HE2	41:w:61:VAL:HG22	1.99	0.44
13:M:290:SER:HA	13:M:319:HIS:HE2	1.81	0.44
13:M:343:ILE:O	13:M:346:ARG:NH1	2.48	0.44
17:X:63:PRO:O	17:X:67:ALA:N	2.50	0.44
29:k:115:LEU:HD13	29:k:121:GLY:HA2	1.99	0.44
36:r:103:ILE:HB	37:s:47:ASP:HB3	1.98	0.44
53:M:503:CDL:H432	33:o:27:THR:HG21	1.99	0.44
31:m:30:ILE:O	31:m:33:THR:OG1	2.33	0.44
7:9:175:TYR:CZ	27:h:38:PRO:HG3	2.52	0.44
9:H:200:LEU:HD22	9:H:282:TYR:HB3	1.98	0.44
11:K:24:SER:HA	11:K:90:VAL:HG22	2.00	0.44
51:L:1003:PC1:H32	51:w:801:PC1:H142	2.00	0.44
1:1:17:ASP:HA	1:1:20:ARG:HG2	1.98	0.44
3:3:377:VAL:HG23	3:3:404:LEU:HD11	2.00	0.44
8:A:64:LEU:HD11	10:J:168:ILE:HD12	1.99	0.44
53:L:1005:CDL:H741	13:M:369:LEU:HD21	1.99	0.44
13:M:94:LEU:HD13	53:o:502:CDL:H762	1.99	0.44
14:N:89:LEU:HD12	14:N:95:SER:HA	1.98	0.44
18:Y:144:ARG:HH12	30:l:57:ILE:HD12	1.82	0.44
19:Z:141:ARG:HG3	41:w:112:CYS:HB2	1.99	0.44
23:d:264:ARG:HH21	23:d:281:ARG:HD3	1.82	0.44
7:9:169:ILE:O	7:9:173:TYR:HB3	2.17	0.44
9:H:152:SER:HA	9:H:155:LEU:HD12	1.98	0.44
23:d:22:THR:OG1	23:d:88:SER:OG	2.26	0.44
34:p:70:ALA:HB2	40:v:10:PRO:HG2	2.00	0.44
39:u:45:ASP:HB3	39:u:50:HIS:HA	2.00	0.44
7:9:23:GLN:HG3	7:9:28:THR:HB	1.98	0.44
7:9:172:ASP:OD2	7:9:176:ARG:NE	2.50	0.44
11:K:70:GLU:O	11:K:73:LEU:HB3	2.18	0.44
13:M:1:FME:HE2	13:M:1:FME:HB2	1.81	0.44
14:N:110:PRO:HD3	14:N:160:LEU:HD23	2.00	0.44
14:N:211:MET:HG2	14:N:333:SER:HB2	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:d:45:VAL:O	23:d:68:ILE:HA	2.17	0.44
27:h:95:SER:OG	27:h:97:ALA:O	2.32	0.44
40:v:96:VAL:HB	40:v:101:MET:HE3	2.00	0.44
4:4:104:ASP:HB3	4:4:190:HIS:ND1	2.31	0.44
4:4:154:VAL:HG11	4:4:225:LEU:HD21	2.00	0.44
53:M:503:CDL:H311	53:M:503:CDL:H562	2.00	0.44
23:d:76:ARG:NH2	23:d:103:ASN:O	2.47	0.44
23:d:246:PHE:HD1	23:d:251:ARG:HB2	1.83	0.44
26:g:114:PRO:HG2	26:g:123:VAL:HG21	2.00	0.44
29:k:30:ASN:O	29:k:126:ARG:NH2	2.51	0.44
41:w:100:ARG:HB2	41:w:105:LEU:HB2	1.99	0.44
4:4:387:TYR:CZ	5:5:164:LYS:HE3	2.53	0.44
14:N:122:ILE:O	14:N:176:ARG:NH1	2.51	0.44
23:d:196:LEU:HD22	23:d:257:PRO:HB3	1.98	0.44
1:1:95:VAL:HB	1:1:136:ILE:HG12	1.99	0.43
4:4:106:LEU:HD13	4:4:391:ILE:HG21	2.00	0.43
4:4:237:ASN:O	9:H:284:GLN:NE2	2.39	0.43
9:H:6:VAL:HG22	9:H:95:LEU:HD21	1.99	0.43
23:d:153:GLU:HG3	23:d:165:ILE:HG21	1.99	0.43
29:k:218:CYS:O	29:k:245:LYS:NZ	2.49	0.43
1:1:91:LYS:HG2	1:1:219:PRO:HB2	1.98	0.43
3:3:382:THR:HB	3:3:454:GLY:HA3	1.99	0.43
4:4:123:GLU:OE2	4:4:138:ARG:NH2	2.51	0.43
8:A:6:THR:HG21	9:H:87:ILE:HG21	2.00	0.43
52:J:202:3PE:H222	11:K:21:MET:HB2	2.00	0.43
54:X:101:ZMP:H14	38:t:59:ALA:HB1	2.00	0.43
23:d:167:LYS:O	23:d:230:PHE:N	2.43	0.43
12:L:556:ILE:O	12:L:560:THR:N	2.51	0.43
12:L:563:PRO:HA	52:N:401:3PE:H3A1	2.00	0.43
33:o:34:ILE:HG22	33:o:68:PHE:HB3	1.99	0.43
9:H:121:TRP:HB2	10:J:68:PHE:HZ	1.83	0.43
12:L:474:PRO:HD2	37:s:66:LEU:HD21	2.01	0.43
38:t:27:GLU:OE2	38:t:33:ARG:NH1	2.51	0.43
3:3:242:THR:HG22	3:3:247:VAL:HA	2.01	0.43
51:6:202:PC1:H3H1	52:i:501:3PE:H2H1	2.00	0.43
12:L:128:MET:HE1	12:L:255:ALA:HB2	2.00	0.43
12:L:290:MET:O	12:L:523:SER:OG	2.35	0.43
13:M:423:ILE:HG12	34:p:58:VAL:HG22	2.00	0.43
3:3:231:MET:HB3	3:3:266:LYS:HD2	2.00	0.43
12:L:9:LEU:HD12	36:r:85:LEU:HD11	2.00	0.43
13:M:266:LEU:HB3	13:M:395:LEU:HD13	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
16:W:37:VAL:HG11	53:W:201:CDL:H862	2.01	0.43
54:X:101:ZMP:O6	38:t:12:GLN:NE2	2.51	0.43
2:2:116:ILE:HG23	2:2:169:ILE:HG21	2.01	0.43
7:9:114:THR:HG21	7:9:144:HIS:CD2	2.54	0.43
50:H:501:DCQ:H7A	50:H:501:DCQ:H1M	1.84	0.43
23:d:96:GLY:HA3	56:d:401:NDP:H3D	2.01	0.43
1:1:92:TYR:O	1:1:220:THR:HA	2.18	0.43
2:2:24:THR:OG1	2:2:27:ASN:OD1	2.32	0.43
6:6:176:TRP:HB2	51:6:202:PC1:H12	2.00	0.43
9:H:156:MET:HE1	9:H:177:PRO:HB2	1.99	0.43
13:M:207:MET:HE3	13:M:294:MET:HB3	2.01	0.43
27:h:11:ARG:NH2	27:h:20:GLN:OE1	2.52	0.43
32:n:52:TYR:HE2	38:t:34:ASP:HB3	1.83	0.43
2:2:39:PRO:HA	20:a:65:GLN:HB3	2.00	0.43
12:L:602:LEU:HD21	53:V:202:CDL:H422	2.01	0.43
33:o:43:LEU:HD22	33:o:53:VAL:HG12	2.01	0.43
1:1:393:TRP:O	1:1:396:SER:OG	2.32	0.43
1:1:432:ARG:HA	1:1:435:GLN:HG3	2.01	0.43
5:5:134:ILE:HG22	5:5:140:VAL:HB	2.01	0.43
8:A:80:GLN:NE2	9:H:315:PRO:O	2.51	0.43
9:H:24:GLU:HA	9:H:271:LEU:HD13	2.01	0.43
12:L:245:ALA:O	12:L:249:SER:OG	2.31	0.43
35:q:55:GLU:OE1	35:q:58:ARG:NH2	2.43	0.43
37:s:6:ARG:HH12	37:s:105:ARG:HD2	1.83	0.43
1:1:250:ASN:HD21	1:1:320:ASP:HB2	1.84	0.42
3:3:568:GLU:HB3	3:3:589:PRO:HG3	2.01	0.42
13:M:16:TRP:NE1	13:M:97:SER:OG	2.47	0.42
14:N:136:LEU:HD23	14:N:205:LEU:HD21	2.01	0.42
14:N:289:ASN:HA	14:N:292:PHE:CE1	2.54	0.42
3:3:329:ILE:HD12	3:3:329:ILE:HA	1.84	0.42
6:6:58:GLU:HG2	6:6:151:PRO:O	2.19	0.42
8:A:36:PRO:HB2	8:A:43:PRO:HG2	2.01	0.42
13:M:61:LEU:HB2	13:M:457:PRO:HG3	2.01	0.42
16:W:8:LEU:HD11	17:X:88:GLU:HG3	2.00	0.42
23:d:324:TYR:HA	23:d:327:LEU:HB2	2.01	0.42
34:p:107:ASP:OD1	34:p:110:ARG:NH2	2.52	0.42
1:1:68:ARG:C	1:1:224:ASN:HD22	2.28	0.42
7:9:22:ALA:O	7:9:26:LEU:HB2	2.19	0.42
9:H:117:LEU:HD11	9:H:136:VAL:HG23	2.01	0.42
13:M:329:LEU:HB3	13:M:359:TRP:CZ2	2.54	0.42
13:M:408:LEU:O	13:M:412:ILE:N	2.50	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
51:M:502:PC1:H251	51:M:502:PC1:H372	2.01	0.42
28:i:9:ARG:HA	28:i:12:GLN:HG2	2.00	0.42
4:4:430:ARG:NH2	5:5:133:GLU:OE2	2.53	0.42
6:6:62:MET:SD	6:6:69:MET:HB3	2.60	0.42
52:J:202:3PE:H31	11:K:21:MET:HE2	2.00	0.42
11:K:79:VAL:O	11:K:83:ASN:N	2.48	0.42
22:c:29:HIS:HD2	22:c:33:ARG:HH11	1.67	0.42
27:h:8:GLN:NE2	27:h:20:GLN:HG3	2.34	0.42
53:o:502:CDL:H552	53:o:502:CDL:H312	2.01	0.42
1:1:99:GLU:OE2	1:1:106:LYS:N	2.52	0.42
3:3:320:GLY:O	3:3:498:SER:OG	2.38	0.42
4:4:85:2MR:HE	4:4:85:2MR:HB3	1.69	0.42
12:L:128:MET:HG2	12:L:251:THR:HG22	2.02	0.42
13:M:379:LEU:O	13:M:383:MET:HG2	2.20	0.42
19:Z:25:LEU:HD12	37:s:75:ASN:HB2	2.01	0.42
25:f:30:ILE:O	25:f:34:LEU:HB2	2.20	0.42
33:o:120:ARG:HD3	34:p:115:ILE:HD13	2.02	0.42
39:u:61:ASP:HB3	39:u:66:ILE:HB	2.01	0.42
1:1:142:PHE:HB3	1:1:145:GLU:HB2	2.01	0.42
1:1:245:PHE:HB3	1:1:271:GLU:HG3	2.01	0.42
3:3:366:THR:HG22	3:3:370:GLY:HA3	2.02	0.42
4:4:203:LEU:HD22	4:4:207:LEU:HD23	2.00	0.42
4:4:208:MET:HE1	35:q:10:PRO:HD3	2.02	0.42
7:9:95:GLU:HB2	7:9:108:ARG:HB3	2.01	0.42
8:A:99:ALA:HB1	52:A:201:3PE:H391	2.02	0.42
9:H:179:TRP:HE1	35:q:42:LEU:HG	1.84	0.42
10:J:113:VAL:HG12	10:J:119:PHE:HB2	2.02	0.42
12:L:74:LEU:HD13	12:L:190:LEU:HD13	2.01	0.42
53:M:503:CDL:H552	53:M:503:CDL:H152	2.01	0.42
18:Y:43:MET:HE3	18:Y:47:TRP:CZ3	2.55	0.42
24:e:23:CYS:SG	24:e:24:GLN:N	2.93	0.42
24:e:64:LEU:HB3	24:e:76:VAL:HB	2.01	0.42
37:s:30:PHE:HB3	37:s:33:ARG:HB2	2.02	0.42
2:2:43:LYS:HE2	2:2:73:LEU:HA	2.02	0.42
2:2:120:ILE:HD11	2:2:169:ILE:HG12	2.01	0.42
9:H:89:LEU:HD11	9:H:233:MET:HG2	2.02	0.42
10:J:74:MET:HE2	11:K:79:VAL:HG12	2.01	0.42
12:L:129:LEU:HA	12:L:132:VAL:HG22	2.01	0.42
12:L:331:THR:HG22	12:L:335:PHE:HE1	1.84	0.42
13:M:297:VAL:HG13	13:M:312:ALA:HB1	2.02	0.42
14:N:128:LEU:HD13	14:N:216:PHE:HB2	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
18:Y:21:SER:N	44:z:51:ASP:OD1	2.49	0.42
25:f:27:TYR:HB2	25:f:55:LEU:HD13	2.01	0.42
2:2:61:LEU:HD12	2:2:90:TYR:HB3	2.01	0.42
3:3:362:TYR:OH	3:3:504:ASP:OD1	2.32	0.42
4:4:204:PRO:HD2	4:4:207:LEU:HD22	2.01	0.42
9:H:73:LEU:HD11	52:J:201:3PE:H3H2	2.01	0.42
12:L:249:SER:HB2	12:L:336:LYS:HG3	2.01	0.42
12:L:316:THR:HB	12:L:395:ILE:HG23	2.01	0.42
12:L:375:ILE:HG12	39:u:32:MET:SD	2.59	0.42
14:N:202:LEU:HD21	30:l:1:PRO:HD3	2.01	0.42
14:N:258:SER:O	14:N:261:MET:HB3	2.20	0.42
53:W:201:CDL:H192	19:Z:47:VAL:HG21	2.02	0.42
24:e:17:GLU:HG2	24:e:67:ARG:HB3	2.00	0.42
26:g:42:VAL:HG11	26:g:59:ARG:HG3	2.01	0.42
1:1:95:VAL:HG11	1:1:118:LEU:HD11	2.01	0.42
4:4:149:ASN:ND2	4:4:370:PRO:HB2	2.35	0.42
9:H:31:MET:HE1	9:H:272:TRP:HB2	2.02	0.42
10:J:4:TYR:OH	10:J:111:LYS:NZ	2.52	0.42
12:L:149:ILE:HD13	12:L:149:ILE:HA	1.91	0.42
12:L:569:ALA:HB2	53:L:1004:CDL:H541	2.01	0.42
13:M:282:LEU:HD21	13:M:359:TRP:HH2	1.84	0.42
23:d:26:ALA:HB3	23:d:47:VAL:HG13	2.02	0.42
28:i:1:MET:HB3	28:i:4:LEU:HD13	2.01	0.42
31:m:37:TYR:HA	31:m:40:TYR:HD2	1.85	0.42
35:q:86:LEU:HD11	35:q:118:PRO:HG3	2.01	0.42
22:c:69:LEU:HD11	28:i:126:PRO:HG2	2.02	0.42
27:h:13:TRP:O	35:q:27:ARG:NH1	2.40	0.42
4:4:323:ILE:HD12	4:4:324:LYS:HG3	2.02	0.41
5:5:30:ALA:HB2	5:5:40:VAL:HG21	2.02	0.41
11:K:53:PHE:CG	30:l:20:GLN:HG2	2.55	0.41
17:X:8:LEU:H	17:X:87:TYR:HE2	1.68	0.41
19:Z:3:SER:OG	19:Z:4:TRP:N	2.52	0.41
27:h:51:ASN:OD1	27:h:56:ARG:NE	2.52	0.41
3:3:206:GLY:N	45:3:801:SF4:S2	2.81	0.41
6:6:71:ARG:NH2	7:9:49:ASN:OD1	2.51	0.41
12:L:67:HIS:HE1	41:w:86:GLN:HE22	1.68	0.41
19:Z:73:ILE:HG23	19:Z:155:LEU:HD22	2.02	0.41
36:r:125:GLN:HE22	37:s:89:CYS:HB2	1.85	0.41
12:L:193:LEU:HD23	12:L:201:ILE:HA	2.02	0.41
3:3:324:ASP:HB3	3:3:571:ALA:HB1	2.02	0.41
7:9:132:VAL:HG11	7:9:169:ILE:HD11	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:L:75:LYS:HD3	41:w:90:ARG:HH22	1.85	0.41
51:L:1003:PC1:H372	53:W:201:CDL:H512	2.01	0.41
14:N:189:TRP:HZ2	14:N:286:ALA:HB2	1.84	0.41
14:N:347:GLU:HB2	33:o:82:MET:HE2	2.02	0.41
15:V:43:LYS:HD3	53:V:202:CDL:HA4	2.02	0.41
23:d:155:GLU:HA	23:d:158:GLU:HG2	2.03	0.41
33:o:52:PRO:HG2	33:o:56:ALA:HB2	2.01	0.41
1:1:105:CYS:H	1:1:257:ASN:ND2	2.18	0.41
1:1:276:LEU:HD21	1:1:297:VAL:HG11	2.03	0.41
2:2:6:LEU:HD13	2:2:10:ARG:HH12	1.85	0.41
4:4:238:ARG:NH1	9:H:279:ARG:O	2.46	0.41
7:9:164:GLU:HG3	28:i:88:ARG:HG3	2.02	0.41
9:H:58:LYS:HB3	9:H:217:ALA:HB3	2.02	0.41
10:J:34:ILE:HG13	10:J:64:MET:HG3	2.03	0.41
12:L:538:PRO:HB2	40:v:89:VAL:HG22	2.02	0.41
13:M:36:LEU:HG	41:w:68:ILE:HG22	2.01	0.41
13:M:161:LEU:HD11	14:N:264:TRP:CD1	2.55	0.41
13:M:229:MET:O	13:M:233:ALA:CB	2.68	0.41
18:Y:17:VAL:HG12	18:Y:19:VAL:HG22	2.01	0.41
4:4:163:ALA:HB2	9:H:278:PRO:HG3	2.03	0.41
12:L:543:THR:HG23	40:v:83:MET:HE1	2.02	0.41
13:M:426:ILE:HB	38:t:101:TRP:HH2	1.86	0.41
15:V:113:ALA:O	15:V:117:MET:HB2	2.20	0.41
36:r:120:LYS:HE3	37:s:39:VAL:HG13	2.03	0.41
4:4:296:ARG:HH21	4:4:420:THR:HG21	1.85	0.41
8:A:56:PHE:O	10:J:70:TYR:OH	2.37	0.41
10:J:45:LEU:HD21	52:J:201:3PE:H222	2.03	0.41
13:M:41:LEU:HD13	13:M:66:LEU:HD13	2.03	0.41
13:M:116:ILE:HD12	13:M:116:ILE:HA	1.97	0.41
15:V:124:VAL:HG11	53:V:202:CDL:H873	2.03	0.41
53:o:502:CDL:H592	53:o:502:CDL:H802	2.03	0.41
40:v:30:ARG:HB2	40:v:33:ASP:HB2	2.03	0.41
10:J:7:PHE:O	10:J:11:ILE:HG12	2.21	0.41
12:L:88:MET:HB2	12:L:326:PHE:HE2	1.85	0.41
29:k:93:SER:HB3	29:k:138:MET:HE3	2.02	0.41
30:l:94:PRO:HG3	35:q:92:GLU:HG3	2.02	0.41
3:3:503:LEU:HD21	3:3:509:PRO:HB3	2.01	0.41
5:5:137:MET:HE3	5:5:152:LEU:HB2	2.03	0.41
8:A:13:LEU:HD21	9:H:10:ILE:HD11	2.01	0.41
9:H:1:MET:HG3	9:H:4:ILE:HD12	2.03	0.41
9:H:26:LYS:HB3	44:z:5:VAL:HG12	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:L:136:ASN:HD21	51:L:1003:PC1:H2I1	1.86	0.41
12:L:213:LEU:HD23	12:L:216:LEU:HD12	2.02	0.41
52:L:1001:3PE:H221	40:v:88:ARG:HA	2.03	0.41
14:N:277:ILE:HD13	52:V:201:3PE:H372	2.03	0.41
23:d:271:GLU:OE1	23:d:280:THR:OG1	2.29	0.41
29:k:22:SER:OG	29:k:119:GLY:O	2.30	0.41
33:o:5:ARG:NH1	33:o:17:GLU:OE2	2.50	0.41
34:p:49:GLN:O	34:p:55:ARG:NE	2.49	0.41
37:s:11:ASP:OD1	37:s:11:ASP:N	2.54	0.41
40:v:145:ASP:N	40:v:145:ASP:OD1	2.52	0.41
3:3:108:CYS:O	3:3:218:ARG:NH1	2.50	0.41
4:4:151:ILE:O	4:4:155:THR:OG1	2.28	0.41
6:6:48:MET:HE1	6:6:80:PRO:HG3	2.02	0.41
7:9:158:GLY:O	7:9:162:GLU:HB2	2.21	0.41
9:H:20:LEU:HD13	44:z:12:MET:HE1	2.02	0.41
12:L:106:TRP:CD1	12:L:447:ASN:HD22	2.36	0.40
14:N:170:LEU:O	14:N:295:ARG:NH1	2.40	0.40
14:N:178:ILE:HD13	14:N:296:LEU:HD11	2.02	0.40
29:k:308:TYR:HD1	29:k:317:ILE:HG23	1.86	0.40
41:w:80:LEU:HA	41:w:81:PRO:HD3	1.95	0.40
3:3:325:ALA:HB1	3:3:623:LEU:HD11	2.03	0.40
4:4:145:THR:HA	4:4:148:LEU:HD12	2.03	0.40
9:H:26:LYS:HG2	9:H:36:GLY:HA3	2.03	0.40
52:V:201:3PE:H322	52:V:201:3PE:H261	2.03	0.40
7:9:92:ILE:HA	7:9:110:ASP:O	2.21	0.40
12:L:60:GLU:HG2	12:L:83:ASP:HA	2.03	0.40
16:W:143:ASN:HD22	30:l:29:PRO:HB3	1.86	0.40
34:p:10:LEU:HD23	40:v:91:THR:HG22	2.03	0.40
36:r:35:PRO:HA	36:r:36:PRO:HD3	1.95	0.40
1:1:42:TRP:CD2	1:1:161:LEU:HD13	2.56	0.40
1:1:154:ARG:HA	20:a:58:LEU:HD21	2.04	0.40
2:2:50:VAL:HG12	2:2:69:VAL:HG22	2.04	0.40
3:3:174:THR:HG22	3:3:183:VAL:HG22	2.03	0.40
4:4:158:ALA:HB1	4:4:163:ALA:HB3	2.04	0.40
4:4:377:TYR:HB3	4:4:390:LYS:HB3	2.03	0.40
7:9:108:ARG:NH1	22:c:70:MET:O	2.37	0.40
9:H:281:ARG:HH11	9:H:281:ARG:HD3	1.78	0.40
13:M:14:LEU:O	13:M:18:SER:OG	2.37	0.40
14:N:252:GLY:HA3	14:N:290:LEU:HD13	2.02	0.40
15:V:12:PRO:HB2	15:V:15:THR:HG21	2.03	0.40
18:Y:44:LEU:HD22	18:Y:130:VAL:HB	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
21:b:49:VAL:HG22	21:b:90:GLN:HG3	2.03	0.40
53:x:101:CDL:H582	53:x:101:CDL:H781	2.03	0.40
6:6:81:ARG:NH1	9:H:61:LEU:HD21	2.36	0.40
12:L:161:ARG:NE	12:L:238:GLU:OE2	2.55	0.40
13:M:118:PHE:O	13:M:122:PHE:CB	2.69	0.40
14:N:20:VAL:HG11	14:N:137:ALA:HB1	2.04	0.40
14:N:144:GLN:HE21	30:l:2:PHE:HB2	1.86	0.40
20:a:58:LEU:HD23	20:a:58:LEU:HA	1.90	0.40
32:n:21:TRP:HB3	32:n:49:ALA:HB3	2.03	0.40
44:z:3:PHE:O	44:z:6:LEU:HB2	2.22	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	1	428/464 (92%)	405 (95%)	23 (5%)	0	100	100
2	2	211/246 (86%)	195 (92%)	16 (8%)	0	100	100
3	3	686/727 (94%)	655 (96%)	31 (4%)	0	100	100
4	4	427/463 (92%)	403 (94%)	24 (6%)	0	100	100
5	5	206/266 (77%)	195 (95%)	11 (5%)	0	100	100
6	6	154/223 (69%)	147 (96%)	7 (4%)	0	100	100
7	9	174/217 (80%)	164 (94%)	10 (6%)	0	100	100
8	A	113/115 (98%)	106 (94%)	7 (6%)	0	100	100
9	H	316/318 (99%)	299 (95%)	17 (5%)	0	100	100
10	J	173/175 (99%)	164 (95%)	8 (5%)	1 (1%)	21	52
11	K	96/98 (98%)	93 (97%)	3 (3%)	0	100	100
12	L	604/606 (100%)	575 (95%)	29 (5%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
13	M	457/459 (100%)	446 (98%)	11 (2%)	0	100	100
14	N	345/347 (99%)	335 (97%)	10 (3%)	0	100	100
15	V	138/141 (98%)	137 (99%)	1 (1%)	0	100	100
16	W	137/189 (72%)	134 (98%)	3 (2%)	0	100	100
17	X	85/157 (54%)	80 (94%)	5 (6%)	0	100	100
17	j	80/157 (51%)	75 (94%)	5 (6%)	0	100	100
18	Y	169/172 (98%)	164 (97%)	5 (3%)	0	100	100
19	Z	169/175 (97%)	167 (99%)	2 (1%)	0	100	100
20	a	42/109 (38%)	42 (100%)	0	0	100	100
21	b	93/124 (75%)	90 (97%)	3 (3%)	0	100	100
22	c	124/170 (73%)	122 (98%)	2 (2%)	0	100	100
23	d	338/380 (89%)	325 (96%)	13 (4%)	0	100	100
24	e	84/99 (85%)	81 (96%)	3 (4%)	0	100	100
25	f	111/116 (96%)	110 (99%)	1 (1%)	0	100	100
26	g	112/140 (80%)	108 (96%)	4 (4%)	0	100	100
27	h	92/114 (81%)	87 (95%)	5 (5%)	0	100	100
28	i	143/145 (99%)	137 (96%)	6 (4%)	0	100	100
29	k	317/355 (89%)	301 (95%)	16 (5%)	0	100	100
30	l	103/106 (97%)	97 (94%)	6 (6%)	0	100	100
31	m	78/84 (93%)	72 (92%)	6 (8%)	0	100	100
32	n	77/98 (79%)	76 (99%)	1 (1%)	0	100	100
33	o	118/122 (97%)	115 (98%)	3 (2%)	0	100	100
34	p	126/130 (97%)	123 (98%)	3 (2%)	0	100	100
35	q	137/144 (95%)	135 (98%)	2 (2%)	0	100	100
36	r	95/128 (74%)	89 (94%)	6 (6%)	0	100	100
37	s	120/137 (88%)	116 (97%)	4 (3%)	0	100	100
38	t	175/179 (98%)	170 (97%)	5 (3%)	0	100	100
39	u	63/108 (58%)	59 (94%)	4 (6%)	0	100	100
40	v	153/186 (82%)	146 (95%)	7 (5%)	0	100	100
41	w	99/154 (64%)	91 (92%)	8 (8%)	0	100	100
42	x	47/76 (62%)	46 (98%)	1 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
43	y	48/58 (83%)	46 (96%)	2 (4%)	0	100	100
44	z	68/70 (97%)	67 (98%)	1 (2%)	0	100	100
All	All	8131/9247 (88%)	7790 (96%)	340 (4%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
10	J	116	VAL

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all 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	1	344/368 (94%)	344 (100%)	0	100	100
2	2	183/210 (87%)	183 (100%)	0	100	100
3	3	578/608 (95%)	578 (100%)	0	100	100
4	4	370/391 (95%)	370 (100%)	0	100	100
5	5	189/230 (82%)	189 (100%)	0	100	100
6	6	132/181 (73%)	131 (99%)	1 (1%)	73	81
7	9	151/179 (84%)	150 (99%)	1 (1%)	76	82
8	A	103/103 (100%)	103 (100%)	0	100	100
9	H	278/278 (100%)	275 (99%)	3 (1%)	65	78
10	J	144/144 (100%)	144 (100%)	0	100	100
11	K	86/86 (100%)	86 (100%)	0	100	100
12	L	538/538 (100%)	537 (100%)	1 (0%)	87	89
13	M	411/411 (100%)	411 (100%)	0	100	100
14	N	315/315 (100%)	315 (100%)	0	100	100
15	V	101/102 (99%)	101 (100%)	0	100	100
16	W	122/160 (76%)	122 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
17	X	80/141 (57%)	80 (100%)	0	100	100
17	j	76/141 (54%)	76 (100%)	0	100	100
18	Y	154/155 (99%)	154 (100%)	0	100	100
19	Z	155/157 (99%)	155 (100%)	0	100	100
20	a	43/93 (46%)	43 (100%)	0	100	100
21	b	79/97 (81%)	79 (100%)	0	100	100
22	c	113/150 (75%)	113 (100%)	0	100	100
23	d	294/326 (90%)	293 (100%)	1 (0%)	86	87
24	e	76/82 (93%)	76 (100%)	0	100	100
25	f	101/102 (99%)	101 (100%)	0	100	100
26	g	107/124 (86%)	107 (100%)	0	100	100
27	h	84/96 (88%)	84 (100%)	0	100	100
28	i	131/131 (100%)	131 (100%)	0	100	100
29	k	283/309 (92%)	283 (100%)	0	100	100
30	l	94/95 (99%)	94 (100%)	0	100	100
31	m	69/72 (96%)	69 (100%)	0	100	100
32	n	61/76 (80%)	61 (100%)	0	100	100
33	o	107/109 (98%)	107 (100%)	0	100	100
34	p	114/116 (98%)	114 (100%)	0	100	100
35	q	119/122 (98%)	119 (100%)	0	100	100
36	r	95/122 (78%)	95 (100%)	0	100	100
37	s	110/120 (92%)	110 (100%)	0	100	100
38	t	159/161 (99%)	159 (100%)	0	100	100
39	u	59/84 (70%)	59 (100%)	0	100	100
40	v	140/160 (88%)	140 (100%)	0	100	100
41	w	92/130 (71%)	92 (100%)	0	100	100
42	x	44/67 (66%)	44 (100%)	0	100	100
43	y	46/54 (85%)	46 (100%)	0	100	100
44	z	59/59 (100%)	59 (100%)	0	100	100
All	All	7189/7955 (90%)	7182 (100%)	7 (0%)	87	90

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
6	6	54	CYS
7	9	78	ILE
9	H	79	LEU
9	H	126	LYS
9	H	193	THR
12	L	589	LEU
23	d	185	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (74) such sidechains are listed below:

Mol	Chain	Res	Type
1	1	148	ASN
1	1	150	GLN
1	1	293	ASN
1	1	373	ASN
2	2	55	GLN
2	2	101	GLN
2	2	121	GLN
3	3	100	ASN
3	3	179	ASN
3	3	401	HIS
3	3	499	GLN
4	4	98	GLN
4	4	135	GLN
4	4	149	ASN
4	4	252	ASN
5	5	19	HIS
5	5	95	ASN
5	5	144	ASN
5	5	145	HIS
5	5	160	HIS
6	6	94	ASN
8	A	80	GLN
9	H	194	ASN
10	J	46	ASN
12	L	135	ASN
12	L	199	GLN
12	L	354	GLN
12	L	405	ASN
12	L	434	GLN
12	L	506	ASN
12	L	541	ASN
13	M	144	ASN

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Mol	Chain	Res	Type
13	M	184	GLN
13	M	374	ASN
13	M	390	ASN
13	M	415	GLN
14	N	48	HIS
15	V	7	GLN
18	Y	29	HIS
19	Z	123	ASN
20	a	40	ASN
21	b	36	ASN
23	d	44	GLN
23	d	87	HIS
23	d	180	ASN
23	d	250	HIS
23	d	288	HIS
23	d	296	HIS
24	e	85	GLN
25	f	49	GLN
26	g	57	GLN
28	i	31	ASN
28	i	69	ASN
28	i	72	ASN
28	i	112	ASN
29	k	107	GLN
29	k	141	GLN
29	k	239	GLN
29	k	271	ASN
29	k	288	GLN
33	o	61	GLN
34	p	78	ASN
36	r	25	GLN
37	s	46	ASN
38	t	50	HIS
38	t	138	GLN
40	v	56	GLN
40	v	87	ASN
40	v	126	GLN
41	w	119	GLN
42	x	9	HIS
42	x	34	GLN
42	x	46	ASN
44	z	27	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

7 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$
11	FME	K	1	11	8,9,10	1.02	0	8,9,11	1.03	0
27	AYA	h	1	27	6,7,8	1.26	1 (16%)	6,8,10	1.10	1 (16%)
29	SEP	k	36	29	8,9,10	1.61	1 (12%)	7,12,14	1.12	1 (14%)
13	FME	M	1	13	8,9,10	0.98	1 (12%)	8,9,11	0.84	0
4	2MR	4	85	4	10,12,13	2.41	3 (30%)	5,13,15	1.19	1 (20%)
15	AYA	V	1	15	6,7,8	1.18	0	6,8,10	1.65	2 (33%)
12	FME	L	1	12	8,9,10	0.96	0	8,9,11	0.97	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
11	FME	K	1	11	-	2/7/9/11	-
27	AYA	h	1	27	-	1/5/6/8	-
29	SEP	k	36	29	-	4/6/8/10	-
13	FME	M	1	13	-	3/7/9/11	-
4	2MR	4	85	4	-	3/10/13/15	-
15	AYA	V	1	15	-	2/5/6/8	-
12	FME	L	1	12	-	1/7/9/11	-

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	4	85	2MR	CZ-NE	4.87	1.44	1.34
4	4	85	2MR	CZ-NH2	4.84	1.43	1.33
29	k	36	SEP	P-O1P	3.53	1.61	1.50
27	h	1	AYA	CA-N	-2.43	1.43	1.46
4	4	85	2MR	CQ1-NH1	-2.20	1.42	1.46
13	M	1	FME	CA-N	-2.06	1.43	1.46

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
15	V	1	AYA	CA-N-CT	2.53	124.76	121.50
27	h	1	AYA	CB-CA-N	2.39	112.38	109.68
15	V	1	AYA	CB-CA-N	2.30	112.28	109.68
4	4	85	2MR	CQ2-NH2-CZ	-2.30	118.71	123.65
29	k	36	SEP	OG-CB-CA	2.27	110.36	108.14

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
13	M	1	FME	C-CA-CB-CG
13	M	1	FME	O-C-CA-CB
27	h	1	AYA	O-C-CA-CB
29	k	36	SEP	CB-OG-P-O2P
29	k	36	SEP	CB-OG-P-O3P
4	4	85	2MR	C-CA-CB-CG
12	L	1	FME	CA-CB-CG-SD
15	V	1	AYA	OT-CT-N-CA
15	V	1	AYA	CM-CT-N-CA
11	K	1	FME	CA-CB-CG-SD
4	4	85	2MR	CA-CB-CG-CD
29	k	36	SEP	CB-OG-P-O1P
4	4	85	2MR	NE-CD-CG-CB
29	k	36	SEP	CA-CB-OG-P
13	M	1	FME	N-CA-CB-CG
11	K	1	FME	CB-CA-N-CN

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
13	M	1	FME	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	4	85	2MR	1	0
15	V	1	AYA	2	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 45 ligands modelled in this entry, 2 are monoatomic - leaving 43 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	SF4	9	402	7	0,12,12	-	-	-		
45	SF4	3	801	3	0,12,12	-	-	-		
46	FMN	1	502	-	33,33,33	1.10	2 (6%)	48,50,50	1.25	7 (14%)
48	FES	2	300	2	0,4,4	-	-	-		
52	3PE	J	202	-	39,39,50	0.36	0	42,44,55	0.37	0
52	3PE	M	501	-	43,43,50	0.34	0	46,48,55	0.50	1 (2%)
53	CDL	x	101	-	74,74,99	0.31	0	80,86,111	0.45	1 (1%)
52	3PE	N	401	-	39,39,50	0.33	0	42,44,55	0.34	0
51	PC1	L	1003	-	53,53,53	0.32	0	59,61,61	0.61	2 (3%)
53	CDL	W	201	-	99,99,99	0.27	0	105,111,111	0.30	0
57	AMP	k	501	-	25,25,25	1.42	4 (16%)	37,38,38	1.89	9 (24%)
53	CDL	L	1005	-	99,99,99	0.25	0	105,111,111	0.31	0
53	CDL	V	202	-	93,93,99	0.26	0	99,105,111	0.27	0
45	SF4	6	201	6	0,12,12	-	-	-		
52	3PE	i	502	-	50,50,50	0.30	0	53,55,55	0.29	0
53	CDL	M	503	-	99,99,99	0.26	0	105,111,111	0.31	0
52	3PE	A	201	-	50,50,50	0.31	0	53,55,55	0.34	0
50	DCQ	4	501	-	23,23,23	0.19	0	27,29,29	0.50	0
51	PC1	9	401	-	53,53,53	0.30	0	59,61,61	0.52	1 (1%)
58	MYR	s	201	37	13,14,15	0.13	0	12,13,15	0.19	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
45	SF4	1	501	1	0,12,12	-	-	-		
52	3PE	o	501	-	30,30,50	0.38	0	33,35,55	0.42	0
51	PC1	M	502	-	53,53,53	0.31	0	59,61,61	0.38	0
54	ZMP	X	101	17	25,30,36	0.81	1 (4%)	29,37,45	1.03	2 (6%)
45	SF4	9	403	7	0,12,12	-	-	-		
48	FES	3	803	3	0,4,4	-	-	-		
52	3PE	L	1001	-	50,50,50	0.29	0	53,55,55	0.39	0
52	3PE	N	402	-	50,50,50	0.33	0	53,55,55	0.57	2 (3%)
52	3PE	J	201	-	50,50,50	0.31	0	53,55,55	0.31	0
52	3PE	V	201	-	36,36,50	0.35	0	39,41,55	0.32	0
50	DCQ	H	501	-	23,23,23	0.21	0	27,29,29	0.53	0
52	3PE	p	201	-	26,26,50	0.49	0	29,31,55	0.59	1 (3%)
53	CDL	o	502	-	89,89,99	0.28	0	95,101,111	0.41	0
54	ZMP	g	201	-	28,33,36	0.61	1 (3%)	32,40,45	1.20	4 (12%)
56	NDP	d	401	-	51,52,52	0.36	0	71,80,80	0.38	0
47	NAI	1	503	-	47,48,48	0.41	0	64,73,73	1.81	4 (6%)
53	CDL	L	1004	-	84,84,99	0.29	0	90,96,111	0.28	0
45	SF4	3	802	3	0,12,12	-	-	-		
52	3PE	i	501	-	50,50,50	0.30	0	53,55,55	0.30	0
51	PC1	w	801	-	53,53,53	0.29	0	59,61,61	0.38	0
53	CDL	h	201	-	57,57,99	0.33	0	63,69,111	0.28	0
51	PC1	6	202	-	45,45,53	0.31	0	51,53,61	0.34	0
52	3PE	L	1002	-	30,30,50	0.42	0	33,35,55	0.79	1 (3%)

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	SF4	3	801	3	-	-	0/6/5/5
45	SF4	9	402	7	-	-	0/6/5/5
46	FMN	1	502	-	-	7/18/18/18	0/3/3/3
48	FES	2	300	2	-	-	0/1/1/1
52	3PE	J	202	-	-	13/43/43/54	-
52	3PE	M	501	-	-	10/47/47/54	-
53	CDL	x	101	-	2/2/9/9	18/85/85/110	-
52	3PE	N	401	-	-	10/43/43/54	-
51	PC1	L	1003	-	-	19/57/57/57	-
53	CDL	W	201	-	-	23/110/110/110	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
57	AMP	k	501	-	-	4/10/26/26	0/3/3/3
53	CDL	L	1005	-	1/1/9/9	34/110/110/110	-
53	CDL	V	202	-	-	21/104/104/110	-
52	3PE	i	502	-	-	8/54/54/54	-
53	CDL	M	503	-	1/1/9/9	30/110/110/110	-
45	SF4	6	201	6	-	-	0/6/5/5
52	3PE	A	201	-	-	12/54/54/54	-
50	DCQ	4	501	-	-	5/14/38/38	0/1/1/1
51	PC1	9	401	-	-	14/57/57/57	-
58	MYR	s	201	37	-	2/12/12/13	-
52	3PE	o	501	-	-	7/34/34/54	-
45	SF4	1	501	1	-	-	0/6/5/5
51	PC1	M	502	-	-	19/57/57/57	-
54	ZMP	X	101	17	-	15/35/37/43	-
52	3PE	L	1001	-	-	14/54/54/54	-
45	SF4	9	403	7	-	-	0/6/5/5
48	FES	3	803	3	-	-	0/1/1/1
52	3PE	N	402	-	-	17/54/54/54	-
52	3PE	J	201	-	-	10/54/54/54	-
52	3PE	V	201	-	-	13/40/40/54	-
50	DCQ	H	501	-	-	5/14/38/38	0/1/1/1
52	3PE	p	201	-	-	4/27/27/54	-
53	CDL	o	502	-	-	29/100/100/110	-
54	ZMP	g	201	-	-	13/38/40/43	-
56	NDP	d	401	-	-	4/34/77/77	0/5/5/5
47	NAI	1	503	-	-	8/29/72/72	0/5/5/5
53	CDL	L	1004	-	1/1/9/9	30/95/95/110	-
45	SF4	3	802	3	-	-	0/6/5/5
52	3PE	i	501	-	-	7/54/54/54	-
51	PC1	w	801	-	-	13/57/57/57	-
53	CDL	h	201	-	-	15/68/68/110	-
51	PC1	6	202	-	-	8/49/49/57	-
52	3PE	L	1002	-	-	9/34/34/54	-

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
57	k	501	AMP	C5-C4	4.62	1.47	1.39
46	1	502	FMN	C4A-N5	3.08	1.37	1.30
57	k	501	AMP	C5-C6	2.81	1.48	1.41
54	X	101	ZMP	C9-C10	2.63	1.53	1.50
46	1	502	FMN	C10-N1	2.49	1.38	1.33
57	k	501	AMP	C8-N7	2.41	1.36	1.31
54	g	201	ZMP	C9-C10	2.28	1.53	1.50
57	k	501	AMP	C5-N7	-2.12	1.35	1.39

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
47	1	503	NAI	O5B-PA-O1A	-9.53	71.15	108.94
47	1	503	NAI	O2A-PA-O1A	-8.50	72.91	112.44
47	1	503	NAI	O3-PA-O1A	-5.76	93.39	110.70
57	k	501	AMP	C5-C4-N3	-5.70	118.88	126.72
57	k	501	AMP	N3-C4-N9	4.44	134.72	127.17
57	k	501	AMP	C2-N3-C4	3.71	120.89	111.83
57	k	501	AMP	C4-C5-N7	-3.61	106.45	110.58
46	1	502	FMN	C4-N3-C2	-3.39	119.61	125.64
57	k	501	AMP	N3-C2-N1	-3.26	123.64	128.58
54	g	201	ZMP	O1-C10-C9	-3.09	120.42	123.98
57	k	501	AMP	C4-N9-C8	2.84	108.72	105.74
46	1	502	FMN	C4A-C4-N3	2.70	120.13	113.25
54	g	201	ZMP	C14-C15-N2	-2.66	106.33	112.00
46	1	502	FMN	O4-C4-C4A	-2.66	119.51	126.53
54	X	101	ZMP	O1-C10-C9	-2.57	121.01	123.98
54	g	201	ZMP	C15-C14-C13	-2.54	108.16	112.39
57	k	501	AMP	C5-N7-C8	2.53	107.43	103.45
46	1	502	FMN	C4A-C10-N10	2.50	120.06	116.48
51	L	1003	PC1	C2-O21-C21	2.50	123.78	117.80
47	1	503	NAI	O2A-PA-O5B	2.48	118.80	107.57
52	L	1002	3PE	C2-O21-C21	2.44	123.63	117.80
52	N	402	3PE	C2-O21-C21	2.34	123.39	117.80
57	k	501	AMP	C6-C5-N7	2.32	136.57	132.09
46	1	502	FMN	C4A-C10-N1	-2.29	118.97	124.59
46	1	502	FMN	C10-C4A-N5	-2.21	120.30	124.81
51	L	1003	PC1	O21-C2-C1	2.18	116.16	108.34
54	X	101	ZMP	C9-C10-S1	2.11	115.92	113.40
53	x	101	CDL	CB4-OB6-CB5	2.11	122.84	117.80
51	9	401	PC1	C2-O21-C21	2.08	122.77	117.80
54	g	201	ZMP	C9-C10-S1	2.07	115.87	113.40
57	k	501	AMP	N9-C8-N7	-2.07	111.00	113.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
52	M	501	3PE	O21-C2-C3	-2.01	101.13	108.34
52	p	201	3PE	O12-P-O14	2.01	118.66	110.83
52	N	402	3PE	O21-C2-C3	2.00	115.53	108.34
46	1	502	FMN	C4-C4A-C10	2.00	120.36	116.93

All (5) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
53	L	1004	CDL	CB4
53	L	1005	CDL	CB4
53	M	503	CDL	CB4
53	x	101	CDL	CB4
53	x	101	CDL	CA4

All (470) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
47	1	503	NAI	C5B-O5B-PA-O1A
51	6	202	PC1	C1-O11-P-O12
51	6	202	PC1	C1-O11-P-O13
51	9	401	PC1	C11-O13-P-O14
51	9	401	PC1	C1-O11-P-O14
51	9	401	PC1	C1-O11-P-O13
51	L	1003	PC1	C11-O13-P-O14
51	L	1003	PC1	C11-O13-P-O11
51	L	1003	PC1	C1-O11-P-O12
51	M	502	PC1	C11-O13-P-O14
52	A	201	3PE	C11-O13-P-O11
52	A	201	3PE	C11-O13-P-O12
52	A	201	3PE	C11-O13-P-O14
52	J	201	3PE	C1-O11-P-O13
52	J	201	3PE	C1-O11-P-O14
52	J	201	3PE	C11-O13-P-O11
52	J	201	3PE	C11-O13-P-O12
52	J	201	3PE	C11-O13-P-O14
52	J	202	3PE	C1-O11-P-O13
52	J	202	3PE	C1-O11-P-O14
52	J	202	3PE	O13-C11-C12-N
52	L	1001	3PE	C11-O13-P-O11
52	L	1001	3PE	C11-O13-P-O14
52	L	1001	3PE	O13-C11-C12-N
52	L	1002	3PE	C1-O11-P-O12

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Mol	Chain	Res	Type	Atoms
52	L	1002	3PE	C1-O11-P-O13
52	L	1002	3PE	C1-O11-P-O14
52	L	1002	3PE	O13-C11-C12-N
52	M	501	3PE	C11-O13-P-O11
52	M	501	3PE	C11-O13-P-O12
52	M	501	3PE	C11-O13-P-O14
52	M	501	3PE	O13-C11-C12-N
52	N	401	3PE	C1-O11-P-O12
52	N	401	3PE	C1-O11-P-O13
52	N	401	3PE	O13-C11-C12-N
52	N	402	3PE	C1-O11-P-O12
52	N	402	3PE	C1-O11-P-O14
52	N	402	3PE	C11-O13-P-O14
52	N	402	3PE	O13-C11-C12-N
52	V	201	3PE	C11-O13-P-O11
52	V	201	3PE	O13-C11-C12-N
52	i	501	3PE	C11-O13-P-O11
52	i	501	3PE	C11-O13-P-O12
52	i	502	3PE	C1-O11-P-O12
52	i	502	3PE	C1-O11-P-O13
52	i	502	3PE	C1-O11-P-O14
52	i	502	3PE	O13-C11-C12-N
52	o	501	3PE	C11-O13-P-O11
52	o	501	3PE	C11-O13-P-O12
52	o	501	3PE	C11-O13-P-O14
52	o	501	3PE	O13-C11-C12-N
53	L	1004	CDL	O1-C1-CA2-OA2
53	L	1004	CDL	CA2-OA2-PA1-OA3
53	L	1004	CDL	CA2-OA2-PA1-OA4
53	L	1004	CDL	CA2-OA2-PA1-OA5
53	L	1004	CDL	CA3-OA5-PA1-OA2
53	L	1004	CDL	CA3-OA5-PA1-OA3
53	L	1004	CDL	CA3-OA5-PA1-OA4
53	L	1004	CDL	CB2-OB2-PB2-OB4
53	L	1004	CDL	CB3-OB5-PB2-OB3
53	L	1005	CDL	CA3-OA5-PA1-OA2
53	L	1005	CDL	CA3-OA5-PA1-OA3
53	L	1005	CDL	CA3-OA5-PA1-OA4
53	L	1005	CDL	CB3-OB5-PB2-OB2
53	L	1005	CDL	CB3-OB5-PB2-OB3
53	M	503	CDL	O1-C1-CB2-OB2
53	M	503	CDL	CA2-C1-CB2-OB2

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Mol	Chain	Res	Type	Atoms
53	M	503	CDL	CB2-OB2-PB2-OB3
53	M	503	CDL	CB2-OB2-PB2-OB5
53	V	202	CDL	CB2-OB2-PB2-OB3
53	V	202	CDL	CB2-OB2-PB2-OB5
53	W	201	CDL	CA2-OA2-PA1-OA4
53	W	201	CDL	CA2-OA2-PA1-OA5
53	W	201	CDL	CA3-OA5-PA1-OA2
53	W	201	CDL	CA3-OA5-PA1-OA3
53	W	201	CDL	CA3-OA5-PA1-OA4
53	h	201	CDL	CA2-OA2-PA1-OA3
53	h	201	CDL	CA2-OA2-PA1-OA4
53	h	201	CDL	CA2-OA2-PA1-OA5
53	h	201	CDL	CA3-OA5-PA1-OA3
53	h	201	CDL	CB3-OB5-PB2-OB2
53	h	201	CDL	CB3-OB5-PB2-OB3
53	o	502	CDL	CA2-C1-CB2-OB2
53	o	502	CDL	CA2-OA2-PA1-OA4
53	o	502	CDL	CB2-OB2-PB2-OB3
53	o	502	CDL	CB2-OB2-PB2-OB4
53	o	502	CDL	CB2-OB2-PB2-OB5
53	o	502	CDL	CB3-OB5-PB2-OB3
53	x	101	CDL	CA2-OA2-PA1-OA3
53	x	101	CDL	CA2-OA2-PA1-OA4
53	x	101	CDL	CA2-OA2-PA1-OA5
53	x	101	CDL	CA3-OA5-PA1-OA2
53	x	101	CDL	CA3-OA5-PA1-OA3
53	x	101	CDL	CA4-CA3-OA5-PA1
53	x	101	CDL	CB2-OB2-PB2-OB4
53	x	101	CDL	CB2-OB2-PB2-OB5
53	x	101	CDL	CB3-OB5-PB2-OB2
54	X	101	ZMP	O4-C17-C18-C21
54	X	101	ZMP	C16-C17-C18-C21
54	X	101	ZMP	O4-C17-C18-C19
54	X	101	ZMP	C16-C17-C18-C19
54	X	101	ZMP	C16-C17-C18-C20
54	X	101	ZMP	C17-C16-N2-C15
54	g	201	ZMP	S1-C11-C12-N1
54	g	201	ZMP	C12-C11-S1-C10
57	k	501	AMP	C5'-O5'-P-O1P
57	k	501	AMP	C5'-O5'-P-O2P
57	k	501	AMP	C5'-O5'-P-O3P
50	4	501	DCQ	C6-C7-C8-C9

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Mol	Chain	Res	Type	Atoms
53	h	201	CDL	O1-C1-CB2-OB2
53	o	502	CDL	O1-C1-CB2-OB2
54	X	101	ZMP	O3-C16-N2-C15
52	V	201	3PE	C2-C1-O11-P
53	L	1005	CDL	C1-CA2-OA2-PA1
53	L	1004	CDL	CB2-C1-CA2-OA2
53	h	201	CDL	CA2-C1-CB2-OB2
52	J	201	3PE	O21-C2-C3-O31
51	6	202	PC1	C11-C12-N-C14
51	M	502	PC1	C11-C12-N-C14
51	L	1003	PC1	C2-C1-O11-P
53	M	503	CDL	CB5-C51-C52-C53
51	M	502	PC1	C11-C12-N-C15
53	L	1004	CDL	C1-CA2-OA2-PA1
53	M	503	CDL	CA7-C31-C32-C33
51	6	202	PC1	C11-C12-N-C15
53	W	201	CDL	C35-C36-C37-C38
52	i	502	3PE	C29-C2A-C2B-C2C
53	o	502	CDL	C78-C79-C80-C81
51	9	401	PC1	C3B-C3C-C3D-C3E
53	L	1005	CDL	C77-C78-C79-C80
53	L	1005	CDL	C21-C22-C23-C24
53	W	201	CDL	C82-C83-C84-C85
50	H	501	DCQ	C11-C10-C9-C8
53	M	503	CDL	C81-C82-C83-C84
53	L	1005	CDL	C82-C83-C84-C85
53	M	503	CDL	C12-C13-C14-C15
52	N	402	3PE	C37-C38-C39-C3A
53	W	201	CDL	CB5-C51-C52-C53
52	i	501	3PE	C3C-C3D-C3E-C3F
53	M	503	CDL	C63-C64-C65-C66
51	L	1003	PC1	C38-C39-C3A-C3B
51	9	401	PC1	C3C-C3D-C3E-C3F
53	L	1005	CDL	C22-C23-C24-C25
53	L	1004	CDL	C13-C14-C15-C16
53	L	1004	CDL	C78-C79-C80-C81
53	o	502	CDL	CB5-C51-C52-C53
53	o	502	CDL	C81-C82-C83-C84
51	9	401	PC1	C2C-C2D-C2E-C2F
51	6	202	PC1	C11-C12-N-C13
52	J	202	3PE	C26-C27-C28-C29
46	1	502	FMN	C2'-C3'-C4'-O4'

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Mol	Chain	Res	Type	Atoms
53	V	202	CDL	C79-C80-C81-C82
52	J	201	3PE	C21-C22-C23-C24
53	M	503	CDL	C52-C53-C54-C55
52	J	202	3PE	C32-C33-C34-C35
53	o	502	CDL	C52-C53-C54-C55
52	J	202	3PE	C33-C34-C35-C36
53	L	1005	CDL	C56-C57-C58-C59
51	M	502	PC1	C11-C12-N-C13
51	L	1003	PC1	C31-C32-C33-C34
52	L	1002	3PE	C21-C22-C23-C24
52	N	402	3PE	C2B-C2C-C2D-C2E
51	M	502	PC1	C3A-C3B-C3C-C3D
52	L	1001	3PE	C34-C35-C36-C37
52	N	402	3PE	O21-C2-C3-O31
52	A	201	3PE	C38-C39-C3A-C3B
53	V	202	CDL	C39-C40-C41-C42
51	6	202	PC1	C3B-C3C-C3D-C3E
52	A	201	3PE	C2E-C2F-C2G-C2H
52	J	202	3PE	C34-C35-C36-C37
52	A	201	3PE	C2C-C2D-C2E-C2F
51	L	1003	PC1	O11-C1-C2-C3
51	M	502	PC1	O11-C1-C2-C3
53	W	201	CDL	OA5-CA3-CA4-CA6
52	A	201	3PE	C35-C36-C37-C38
53	M	503	CDL	C11-C12-C13-C14
52	J	201	3PE	C1-C2-C3-O31
52	J	202	3PE	C1-C2-C3-O31
53	M	503	CDL	CA3-CA4-CA6-OA8
52	N	402	3PE	C29-C2A-C2B-C2C
51	L	1003	PC1	C23-C24-C25-C26
53	M	503	CDL	C35-C36-C37-C38
53	L	1005	CDL	C41-C42-C43-C44
53	V	202	CDL	OB5-CB3-CB4-OB6
51	w	801	PC1	C3E-C3F-C3G-C3H
53	L	1005	CDL	OA6-CA4-CA6-OA8
53	o	502	CDL	C62-C63-C64-C65
54	X	101	ZMP	C2-C3-C4-C5
54	X	101	ZMP	O4-C17-C18-C20
54	g	201	ZMP	O4-C17-C18-C20
52	V	201	3PE	C33-C34-C35-C36
53	h	201	CDL	C57-C58-C59-C60
52	L	1002	3PE	C2-C1-O11-P

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Mol	Chain	Res	Type	Atoms
53	V	202	CDL	C34-C35-C36-C37
53	M	503	CDL	C36-C37-C38-C39
54	g	201	ZMP	C5-C6-C7-C8
53	o	502	CDL	OB5-CB3-CB4-CB6
53	L	1004	CDL	C72-C73-C74-C75
53	M	503	CDL	C14-C15-C16-C17
53	L	1005	CDL	CA3-CA4-CA6-OA8
51	L	1003	PC1	O11-C1-C2-O21
52	L	1001	3PE	O11-C1-C2-O21
53	V	202	CDL	OA5-CA3-CA4-OA6
53	W	201	CDL	OB5-CB3-CB4-OB6
58	s	201	MYR	C11-C10-C9-C8
54	g	201	ZMP	O1-C10-S1-C11
53	M	503	CDL	C59-C60-C61-C62
47	1	503	NAI	PN-O3-PA-O5B
53	V	202	CDL	C71-C72-C73-C74
46	1	502	FMN	O3'-C3'-C4'-C5'
53	W	201	CDL	C31-C32-C33-C34
53	L	1004	CDL	CB7-C71-C72-C73
53	L	1005	CDL	C17-C18-C19-C20
52	M	501	3PE	C32-C31-O31-C3
53	V	202	CDL	OA5-CA3-CA4-CA6
53	V	202	CDL	OB5-CB3-CB4-CB6
53	o	502	CDL	C12-C13-C14-C15
52	p	201	3PE	C22-C23-C24-C25
53	L	1004	CDL	CA5-C11-C12-C13
51	M	502	PC1	C25-C26-C27-C28
50	H	501	DCQ	C5-C4-O4-C4M
51	w	801	PC1	O11-C1-C2-O21
53	L	1004	CDL	OA5-CA3-CA4-OA6
53	W	201	CDL	OA5-CA3-CA4-OA6
53	M	503	CDL	OA6-CA4-CA6-OA8
52	i	501	3PE	C36-C37-C38-C39
46	1	502	FMN	O3'-C3'-C4'-O4'
52	M	501	3PE	C34-C35-C36-C37
53	o	502	CDL	C1-CA2-OA2-PA1
53	V	202	CDL	C77-C78-C79-C80
54	g	201	ZMP	C22-C1-C2-C3
51	9	401	PC1	O13-C11-C12-N
51	L	1003	PC1	O13-C11-C12-N
53	M	503	CDL	C57-C58-C59-C60
53	W	201	CDL	C16-C17-C18-C19

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Mol	Chain	Res	Type	Atoms
54	X	101	ZMP	O3-C16-C17-O4
50	4	501	DCQ	C2-C3-O3-C3M
53	L	1004	CDL	C73-C74-C75-C76
53	L	1004	CDL	OA5-CA3-CA4-CA6
46	1	502	FMN	N10-C1'-C2'-O2'
54	X	101	ZMP	C19-C18-C21-O5
53	M	503	CDL	C58-C59-C60-C61
51	9	401	PC1	C34-C35-C36-C37
51	M	502	PC1	O11-C1-C2-O21
52	N	402	3PE	C23-C24-C25-C26
53	L	1005	CDL	C37-C38-C39-C40
53	o	502	CDL	C76-C77-C78-C79
52	J	202	3PE	O21-C2-C3-O31
52	N	402	3PE	C1-C2-C3-O31
53	L	1005	CDL	C38-C39-C40-C41
47	1	503	NAI	C5D-O5D-PN-O3
51	9	401	PC1	C1-O11-P-O12
51	L	1003	PC1	C11-O13-P-O12
51	L	1003	PC1	C1-O11-P-O14
51	L	1003	PC1	C1-O11-P-O13
52	L	1001	3PE	C1-O11-P-O14
52	N	401	3PE	C1-O11-P-O14
52	N	402	3PE	C1-O11-P-O13
52	V	201	3PE	C1-O11-P-O12
52	V	201	3PE	C1-O11-P-O13
52	V	201	3PE	C1-O11-P-O14
52	V	201	3PE	C11-O13-P-O14
52	i	502	3PE	C11-O13-P-O14
53	L	1004	CDL	CB2-OB2-PB2-OB3
53	L	1004	CDL	CB2-OB2-PB2-OB5
53	L	1004	CDL	CB3-OB5-PB2-OB2
53	L	1004	CDL	CB3-OB5-PB2-OB4
53	L	1005	CDL	CA2-OA2-PA1-OA3
53	L	1005	CDL	CA2-OA2-PA1-OA4
53	L	1005	CDL	CA2-OA2-PA1-OA5
53	L	1005	CDL	CB3-OB5-PB2-OB4
53	M	503	CDL	CB2-OB2-PB2-OB4
53	M	503	CDL	CB3-OB5-PB2-OB3
53	V	202	CDL	CB2-OB2-PB2-OB4
53	W	201	CDL	CA2-OA2-PA1-OA3
53	W	201	CDL	CB3-OB5-PB2-OB2
53	h	201	CDL	CB3-OB5-PB2-OB4

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Mol	Chain	Res	Type	Atoms
53	o	502	CDL	CA2-OA2-PA1-OA3
53	o	502	CDL	CA2-OA2-PA1-OA5
53	o	502	CDL	CB3-OB5-PB2-OB2
53	x	101	CDL	CA3-OA5-PA1-OA4
54	g	201	ZMP	O4-C17-C18-C21
53	L	1005	CDL	CB7-C71-C72-C73
52	J	202	3PE	C25-C26-C27-C28
53	L	1005	CDL	C58-C59-C60-C61
52	o	501	3PE	C2-C1-O11-P
53	L	1005	CDL	CA4-CA3-OA5-PA1
53	L	1005	CDL	C1-CB2-OB2-PB2
53	V	202	CDL	CA4-CA3-OA5-PA1
53	h	201	CDL	C1-CB2-OB2-PB2
53	h	201	CDL	CB4-CB3-OB5-PB2
52	J	202	3PE	C21-C22-C23-C24
53	L	1005	CDL	C42-C43-C44-C45
52	M	501	3PE	C31-C32-C33-C34
53	W	201	CDL	CA5-C11-C12-C13
52	J	201	3PE	C27-C28-C29-C2A
51	9	401	PC1	C3-C2-O21-C21
51	w	801	PC1	C31-C32-C33-C34
52	o	501	3PE	C25-C26-C27-C28
52	L	1001	3PE	O11-C1-C2-C3
53	x	101	CDL	OB5-CB3-CB4-OB6
51	L	1003	PC1	C39-C3A-C3B-C3C
51	M	502	PC1	C2-C1-O11-P
52	A	201	3PE	C2D-C2E-C2F-C2G
51	w	801	PC1	C27-C28-C29-C2A
53	L	1004	CDL	C12-C13-C14-C15
51	9	401	PC1	C22-C21-O21-C2
46	1	502	FMN	C2'-C3'-C4'-C5'
53	x	101	CDL	C52-C51-CB5-OB6
53	o	502	CDL	CA5-C11-C12-C13
51	M	502	PC1	C39-C3A-C3B-C3C
53	V	202	CDL	C19-C20-C21-C22
52	M	501	3PE	O32-C31-O31-C3
53	W	201	CDL	O1-C1-CB2-OB2
52	o	501	3PE	C22-C23-C24-C25
47	1	503	NAI	C2D-C1D-N1N-C2N
47	1	503	NAI	O4D-C1D-N1N-C2N
52	J	202	3PE	C29-C2A-C2B-C2C
53	o	502	CDL	C19-C20-C21-C22

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Mol	Chain	Res	Type	Atoms
53	L	1004	CDL	CA4-CA3-OA5-PA1
51	w	801	PC1	C2B-C2C-C2D-C2E
52	i	501	3PE	C23-C24-C25-C26
53	o	502	CDL	C72-C73-C74-C75
52	N	402	3PE	C24-C25-C26-C27
51	L	1003	PC1	C1-C2-O21-C21
52	L	1002	3PE	C1-C2-O21-C21
53	W	201	CDL	CB7-C71-C72-C73
51	9	401	PC1	C22-C23-C24-C25
53	x	101	CDL	C16-C17-C18-C19
50	4	501	DCQ	C4-C3-O3-C3M
50	H	501	DCQ	C3-C4-O4-C4M
52	N	402	3PE	C27-C28-C29-C2A
53	o	502	CDL	C1-CB2-OB2-PB2
57	k	501	AMP	C4'-C5'-O5'-P
52	p	201	3PE	C26-C27-C28-C29
54	X	101	ZMP	O1-C10-S1-C11
53	L	1005	CDL	C33-C34-C35-C36
51	w	801	PC1	O11-C1-C2-C3
53	M	503	CDL	OB5-CB3-CB4-CB6
51	M	502	PC1	C2F-C2G-C2H-C2I
51	9	401	PC1	O22-C21-O21-C2
51	L	1003	PC1	C36-C37-C38-C39
47	l	503	NAI	C2N-C3N-C7N-N7N
56	d	401	NDP	O4D-C1D-N1N-C6N
54	g	201	ZMP	C1-C2-C3-C4
53	L	1005	CDL	C76-C77-C78-C79
53	W	201	CDL	CA2-C1-CB2-OB2
51	M	502	PC1	C2B-C2C-C2D-C2E
52	J	201	3PE	C2A-C2B-C2C-C2D
52	i	502	3PE	C32-C33-C34-C35
53	o	502	CDL	CB4-CB3-OB5-PB2
51	w	801	PC1	C39-C3A-C3B-C3C
52	i	501	3PE	C33-C34-C35-C36
52	L	1001	3PE	C3A-C3B-C3C-C3D
53	o	502	CDL	CB3-CB4-CB6-OB8
52	A	201	3PE	C3C-C3D-C3E-C3F
52	M	501	3PE	C37-C38-C39-C3A
53	M	503	CDL	OB5-CB3-CB4-OB6
52	N	401	3PE	C32-C33-C34-C35
53	V	202	CDL	C40-C41-C42-C43
54	g	201	ZMP	C9-C10-S1-C11

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Mol	Chain	Res	Type	Atoms
50	H	501	DCQ	C2-C3-O3-C3M
52	V	201	3PE	O21-C21-C22-C23
53	L	1004	CDL	C72-C71-CB7-OB8
51	M	502	PC1	C3B-C3C-C3D-C3E
52	N	402	3PE	C28-C29-C2A-C2B
53	W	201	CDL	OB5-CB3-CB4-CB6
51	L	1003	PC1	O31-C31-C32-C33
56	d	401	NDP	C2D-C1D-N1N-C6N
51	w	801	PC1	C28-C29-C2A-C2B
52	V	201	3PE	C32-C33-C34-C35
53	L	1005	CDL	C55-C56-C57-C58
51	M	502	PC1	C31-C32-C33-C34
53	x	101	CDL	CB6-CB4-OB6-CB5
52	p	201	3PE	C21-C22-C23-C24
51	6	202	PC1	O31-C31-C32-C33
53	W	201	CDL	C75-C76-C77-C78
52	J	202	3PE	C2-C1-O11-P
52	L	1001	3PE	C1-C2-C3-O31
52	i	502	3PE	C36-C37-C38-C39
53	h	201	CDL	CB7-C71-C72-C73
54	g	201	ZMP	C16-C17-C18-C20
51	L	1003	PC1	C24-C25-C26-C27
53	M	503	CDL	C12-C11-CA5-OA6
53	V	202	CDL	C44-C45-C46-C47
53	L	1004	CDL	OA6-CA4-CA6-OA8
53	L	1005	CDL	C11-C12-C13-C14
52	V	201	3PE	O11-C1-C2-C3
53	L	1004	CDL	C54-C55-C56-C57
51	w	801	PC1	C11-C12-N-C14
52	L	1002	3PE	C25-C26-C27-C28
51	L	1003	PC1	C3C-C3D-C3E-C3F
50	4	501	DCQ	C5-C4-O4-C4M
54	g	201	ZMP	O4-C17-C18-C19
53	M	503	CDL	C72-C71-CB7-OB8
53	L	1005	CDL	C34-C35-C36-C37
53	o	502	CDL	C72-C71-CB7-OB8
51	M	502	PC1	C28-C29-C2A-C2B
58	s	201	MYR	C7-C8-C9-C10
52	N	402	3PE	O21-C21-C22-C23
53	x	101	CDL	C72-C71-CB7-OB8
47	1	503	NAI	C2D-C1D-N1N-C6N
52	A	201	3PE	O21-C21-C22-C23

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Mol	Chain	Res	Type	Atoms
52	V	201	3PE	O31-C31-C32-C33
53	M	503	CDL	C52-C51-CB5-OB6
52	A	201	3PE	C24-C25-C26-C27
52	L	1001	3PE	O21-C21-C22-C23
52	N	401	3PE	O21-C21-C22-C23
53	o	502	CDL	C32-C31-CA7-OA8
51	w	801	PC1	C24-C25-C26-C27
53	V	202	CDL	OB6-CB4-CB6-OB8
53	L	1004	CDL	C33-C34-C35-C36
53	h	201	CDL	C52-C51-CB5-OB6
52	i	501	3PE	C3D-C3E-C3F-C3G
53	V	202	CDL	C31-C32-C33-C34
52	L	1001	3PE	O31-C31-C32-C33
53	L	1005	CDL	C32-C31-CA7-OA8
52	L	1001	3PE	C23-C24-C25-C26
46	1	502	FMN	N10-C1'-C2'-C3'
53	M	503	CDL	CB2-C1-CA2-OA2
54	g	201	ZMP	C16-C17-C18-C21
52	N	401	3PE	C31-C32-C33-C34
52	N	402	3PE	C1-C2-O21-C21
53	o	502	CDL	CB3-CB4-OB6-CB5
50	4	501	DCQ	C3-C4-O4-C4M
50	H	501	DCQ	C4-C3-O3-C3M
51	9	401	PC1	C31-C32-C33-C34
52	N	401	3PE	O31-C31-C32-C33
46	1	502	FMN	C4'-C5'-O5'-P
53	x	101	CDL	CB4-CB3-OB5-PB2
54	X	101	ZMP	C12-C11-S1-C10
52	M	501	3PE	O21-C21-C22-C23
53	x	101	CDL	C75-C76-C77-C78
53	M	503	CDL	C12-C11-CA5-OA7
53	o	502	CDL	C72-C71-CB7-OB9
53	x	101	CDL	C72-C71-CB7-OB9
52	V	201	3PE	O32-C31-C32-C33
53	M	503	CDL	C52-C51-CB5-OB7
53	W	201	CDL	C14-C15-C16-C17
53	L	1005	CDL	CB5-C51-C52-C53
52	A	201	3PE	O22-C21-C22-C23
52	N	402	3PE	O22-C21-C22-C23
53	M	503	CDL	C72-C71-CB7-OB9
53	h	201	CDL	C52-C51-CB5-OB7
51	w	801	PC1	C2E-C2F-C2G-C2H

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Mol	Chain	Res	Type	Atoms
51	M	502	PC1	C35-C36-C37-C38
51	w	801	PC1	C26-C27-C28-C29
53	W	201	CDL	C81-C82-C83-C84
52	p	201	3PE	C23-C24-C25-C26
52	L	1001	3PE	O32-C31-C32-C33
52	N	401	3PE	C27-C28-C29-C2A
53	V	202	CDL	C80-C81-C82-C83
54	X	101	ZMP	N2-C16-C17-O4
54	g	201	ZMP	N2-C16-C17-O4
53	L	1004	CDL	CB4-CB3-OB5-PB2
51	w	801	PC1	C11-C12-N-C13
52	L	1001	3PE	O22-C21-C22-C23
54	X	101	ZMP	C17-C18-C21-O5
53	L	1005	CDL	C32-C31-CA7-OA9
51	M	502	PC1	C3F-C3G-C3H-C3I
51	6	202	PC1	C35-C36-C37-C38
52	N	401	3PE	O22-C21-C22-C23
53	V	202	CDL	C13-C14-C15-C16
47	1	503	NAI	O4D-C1D-N1N-C6N
52	L	1002	3PE	O21-C21-C22-C23
56	d	401	NDP	PA-O3-PN-O2N
53	V	202	CDL	CA7-C31-C32-C33
51	M	502	PC1	C2D-C2E-C2F-C2G
56	d	401	NDP	O4B-C4B-C5B-O5B
51	M	502	PC1	O31-C31-C32-C33

There are no ring outliers.

31 monomers are involved in 94 short contacts:

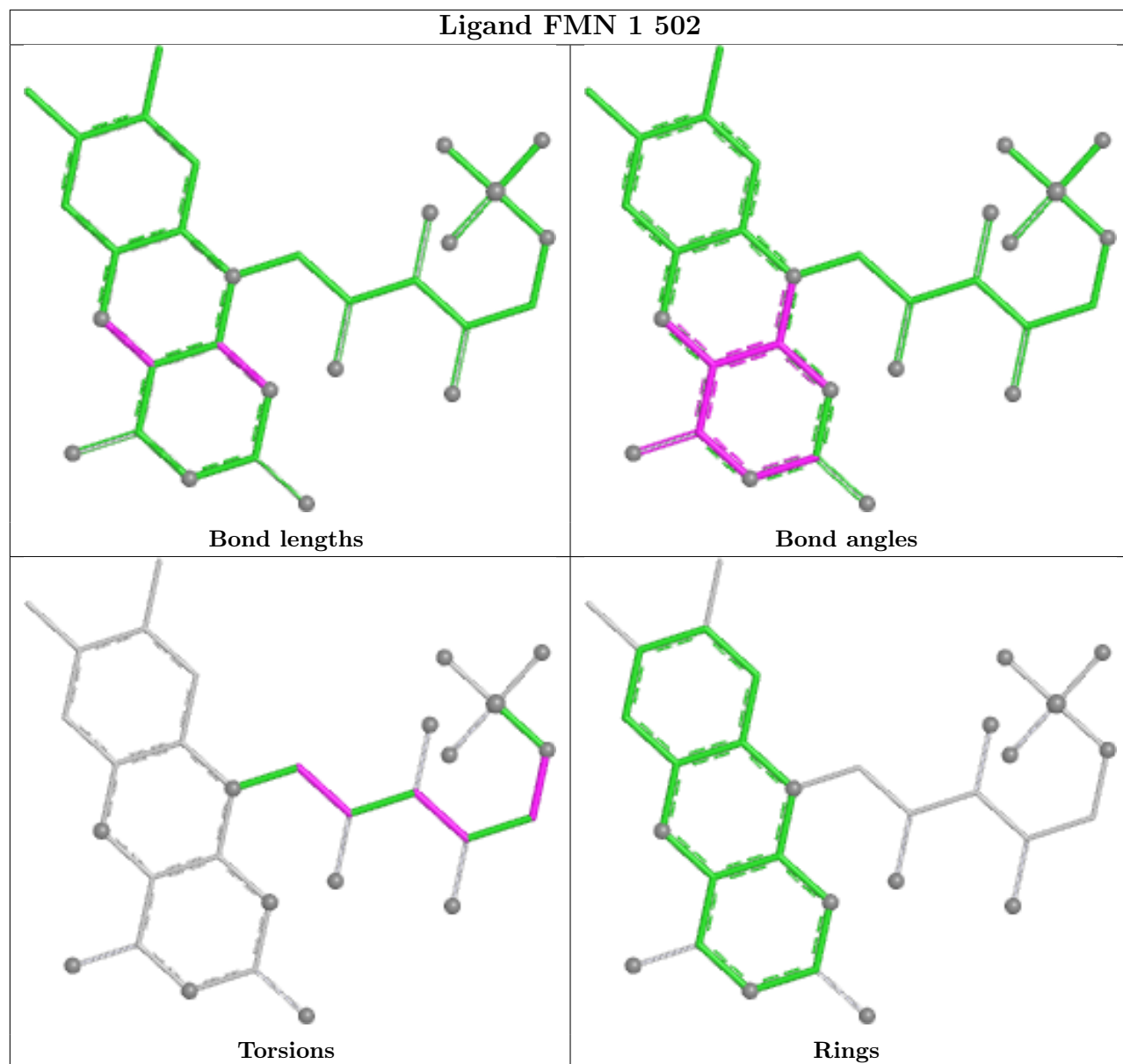
Mol	Chain	Res	Type	Clashes	Symm-Clashes
45	3	801	SF4	1	0
48	2	300	FES	1	0
52	J	202	3PE	5	0
53	x	101	CDL	3	0
52	N	401	3PE	1	0
51	L	1003	PC1	3	0
53	W	201	CDL	4	0
57	k	501	AMP	1	0
53	L	1005	CDL	9	0
53	V	202	CDL	6	0
45	6	201	SF4	1	0
52	i	502	3PE	1	0

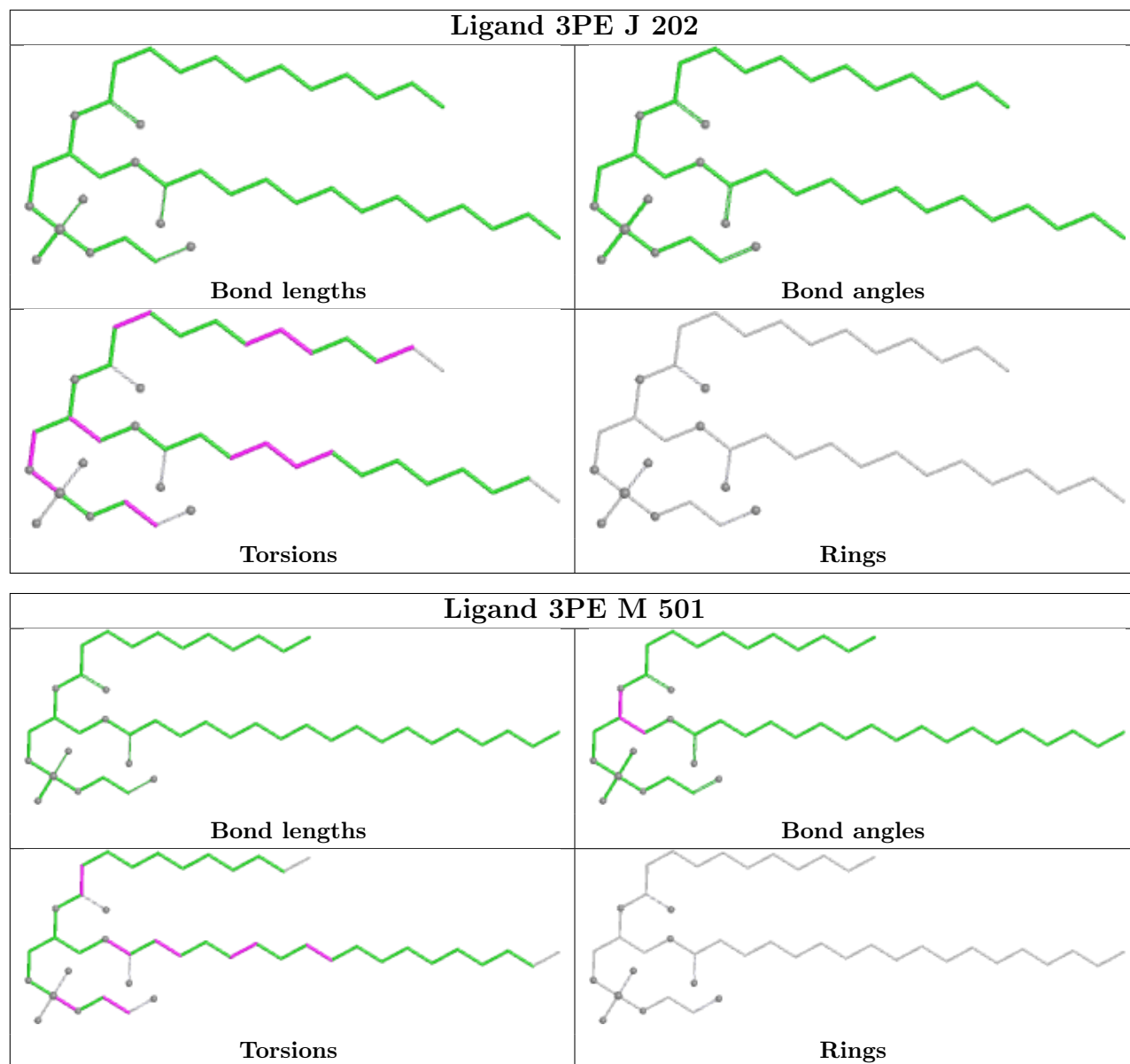
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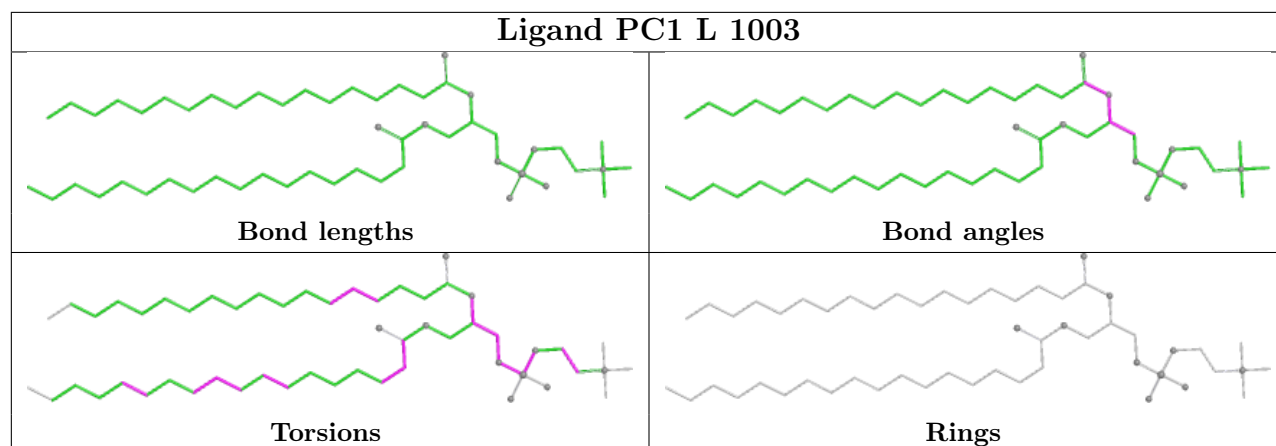
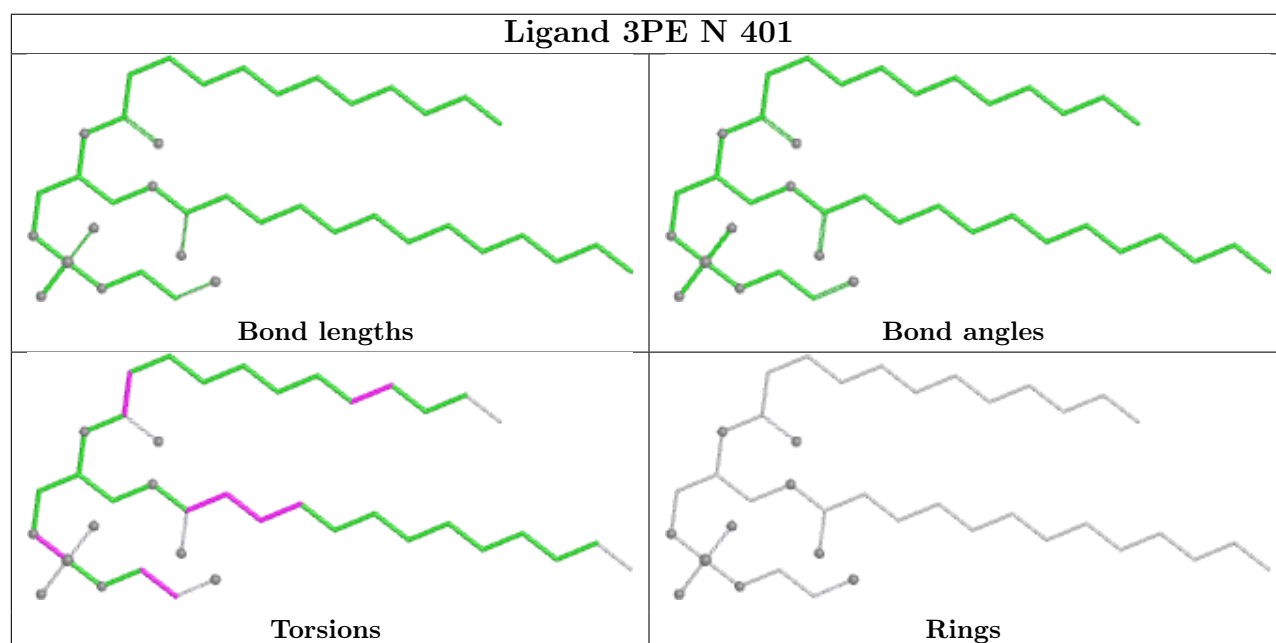
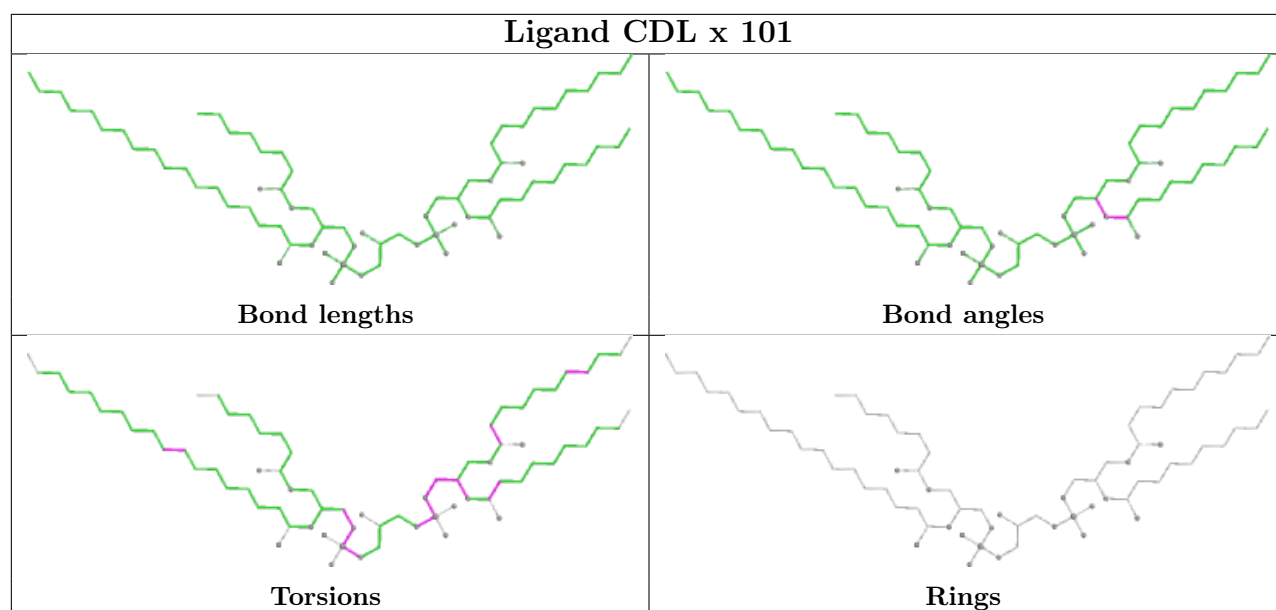
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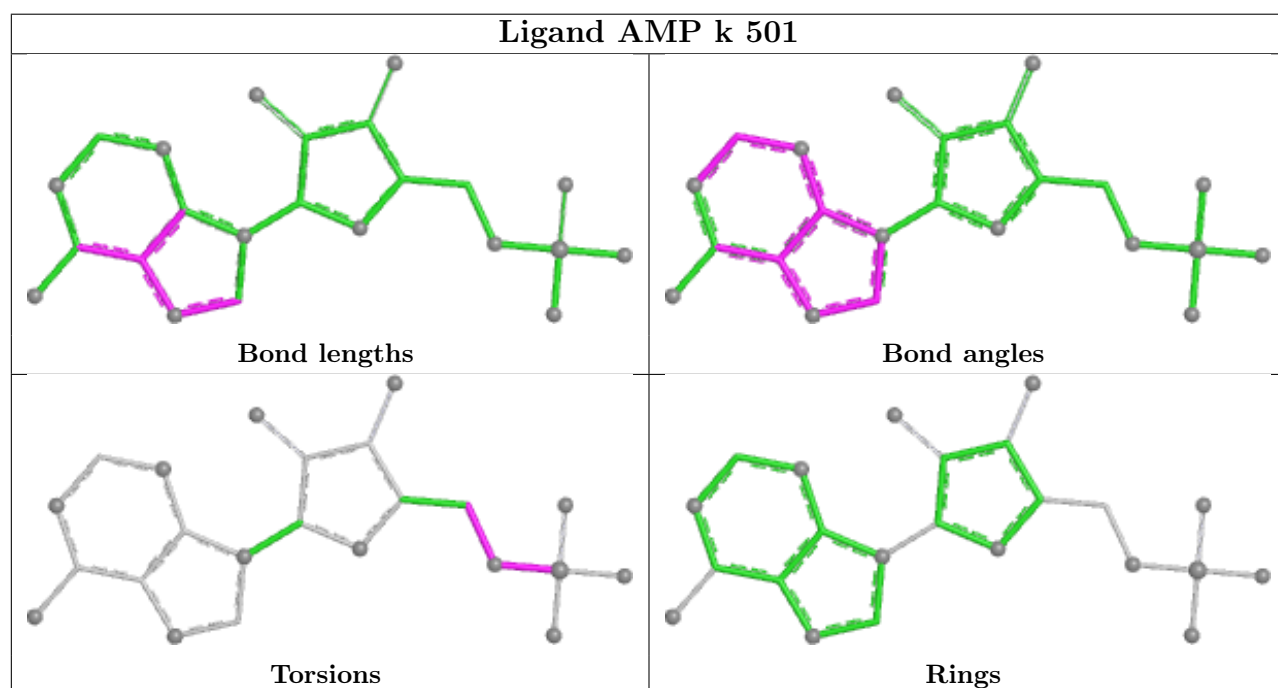
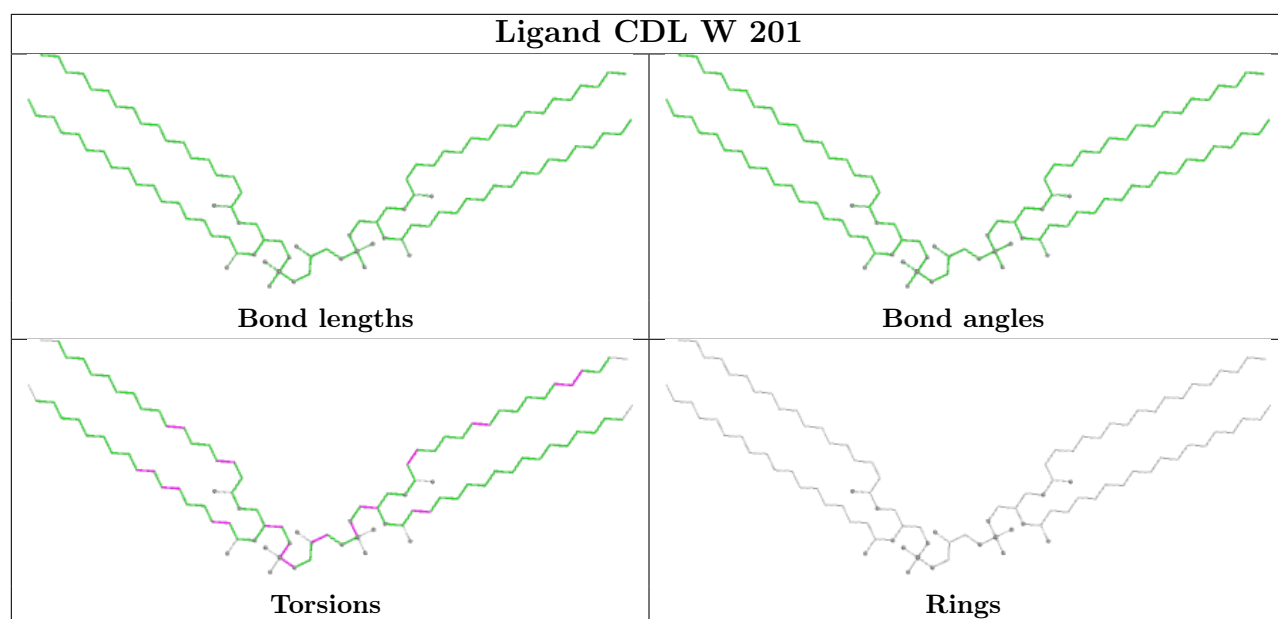
Mol	Chain	Res	Type	Clashes	Symm-Clashes
53	M	503	CDL	13	0
52	A	201	3PE	3	0
50	4	501	DCQ	1	0
51	9	401	PC1	1	0
45	1	501	SF4	2	0
51	M	502	PC1	3	0
54	X	101	ZMP	3	0
52	L	1001	3PE	2	0
52	J	201	3PE	3	0
52	V	201	3PE	2	0
50	H	501	DCQ	2	0
52	p	201	3PE	1	0
53	o	502	CDL	9	0
56	d	401	NDP	2	0
47	1	503	NAI	3	0
53	L	1004	CDL	2	0
52	i	501	3PE	4	0
51	w	801	PC1	8	0
51	6	202	PC1	2	0

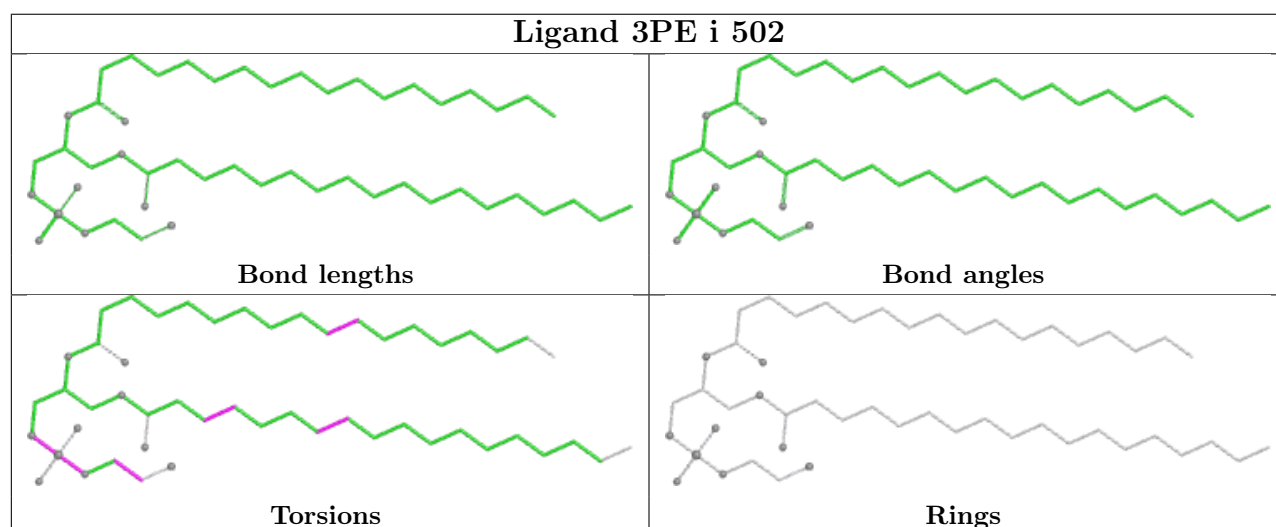
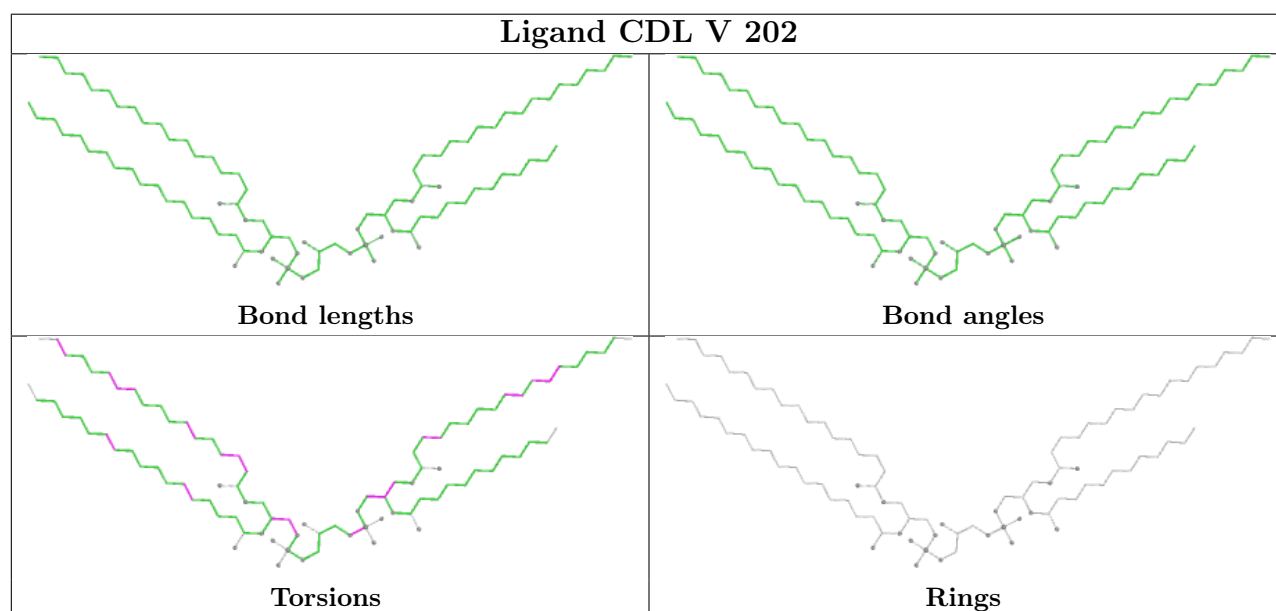
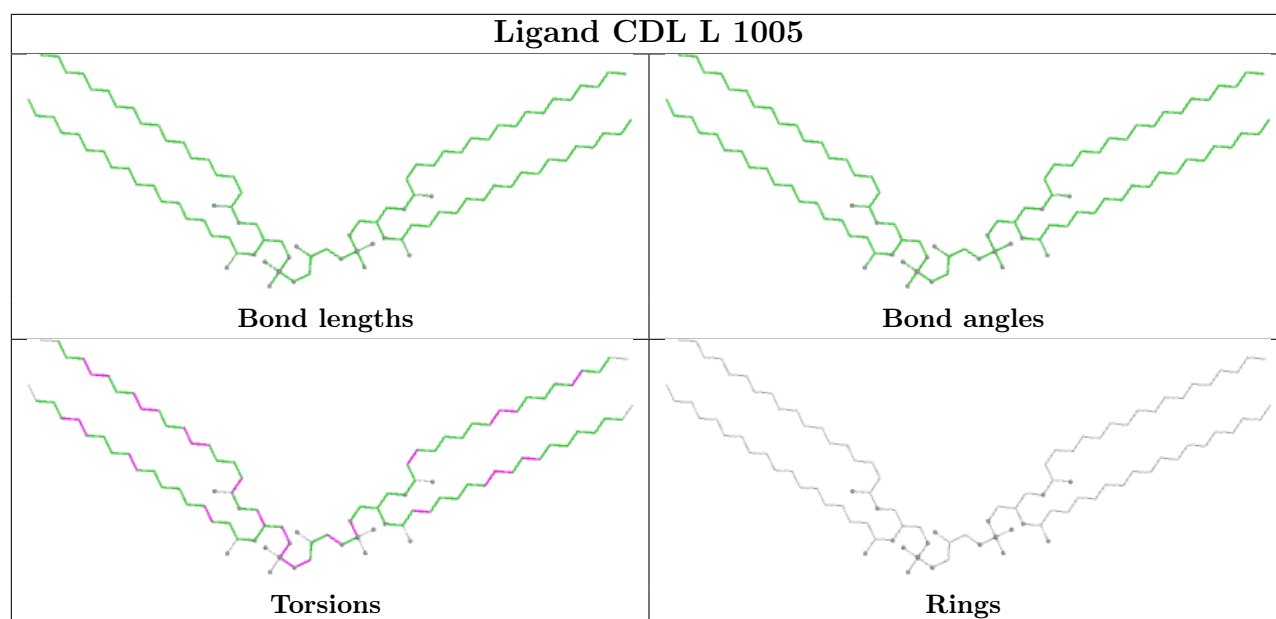
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

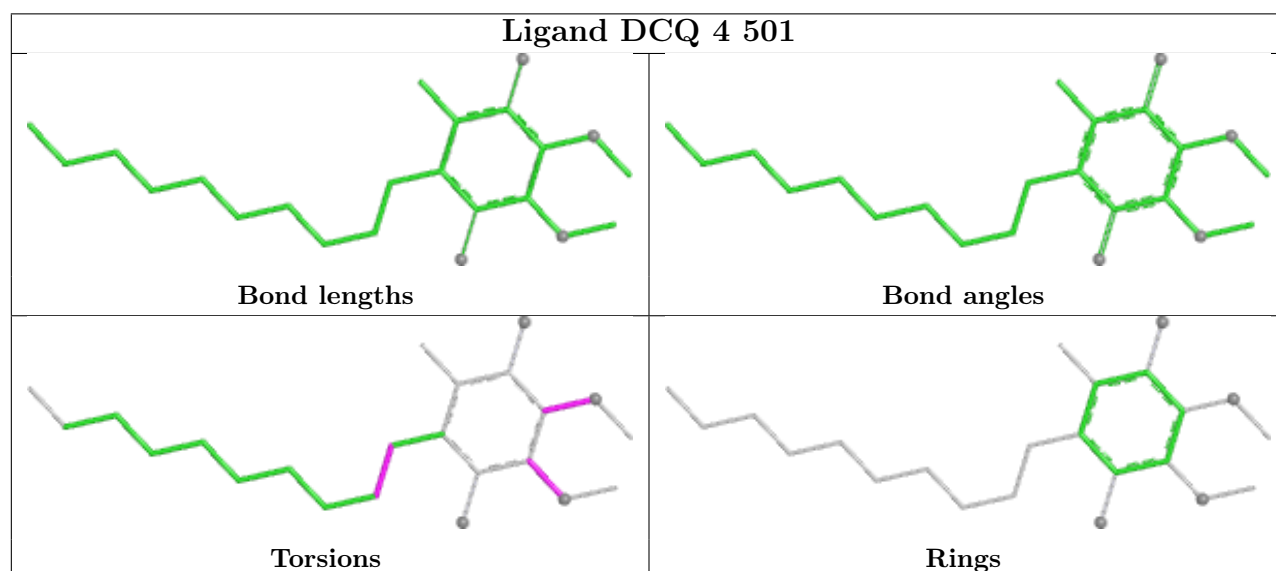
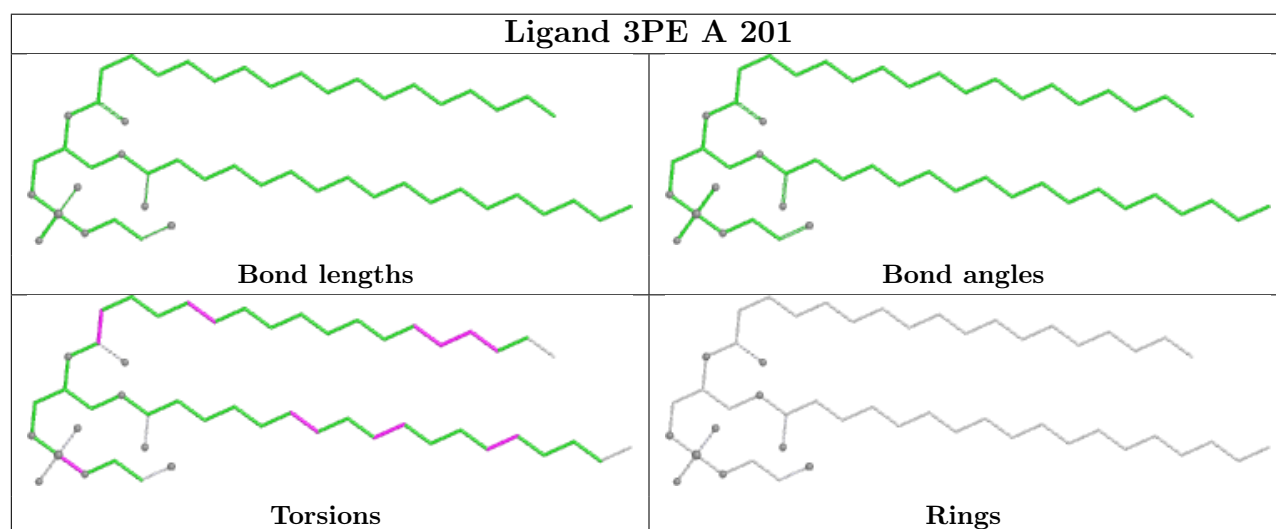
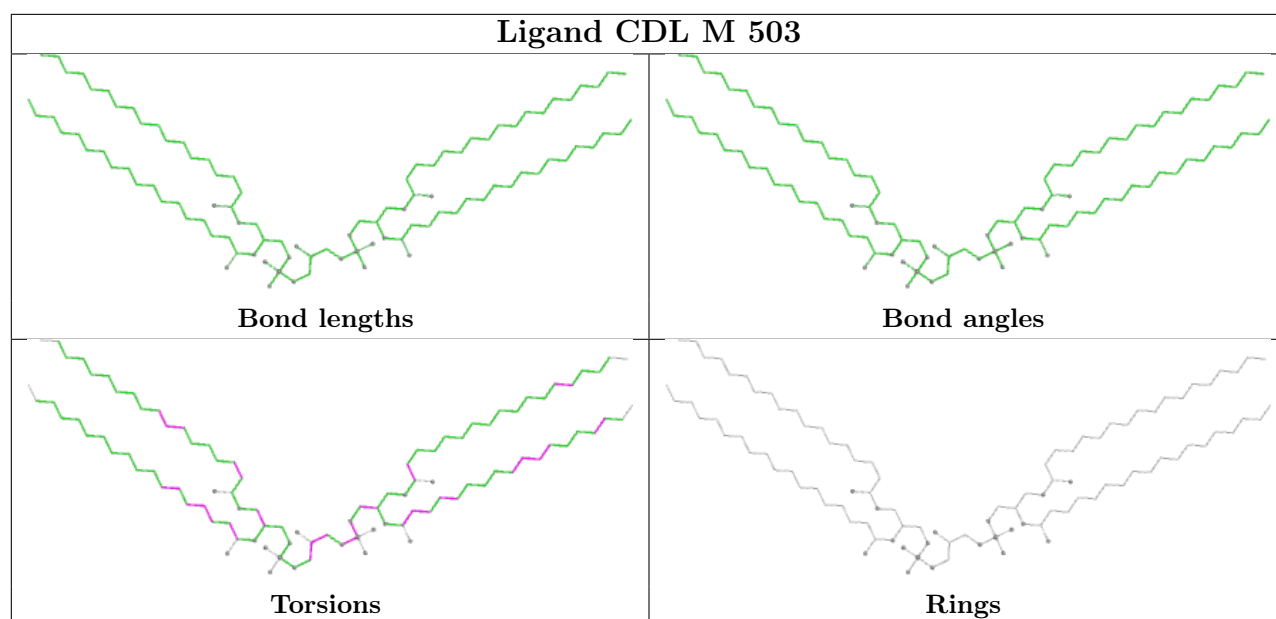


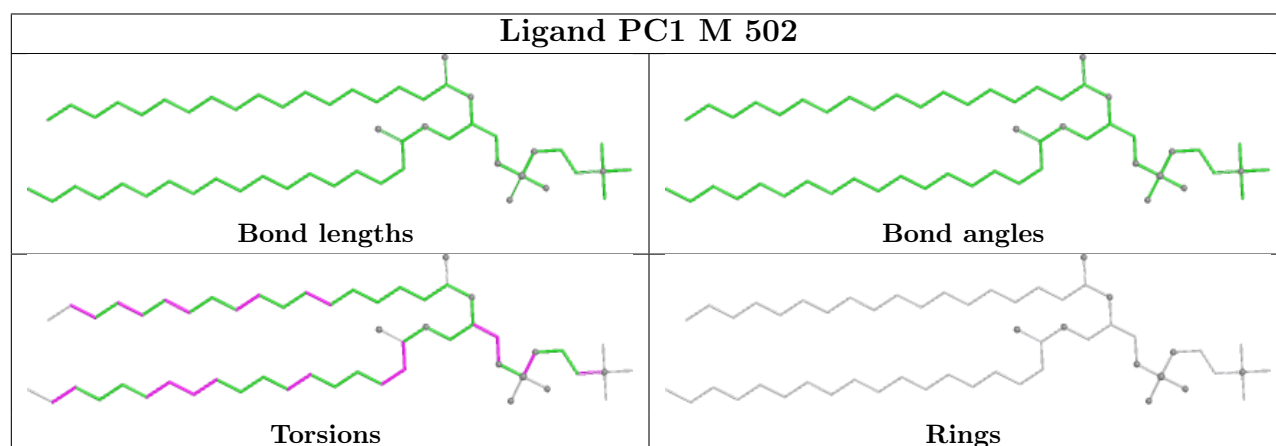
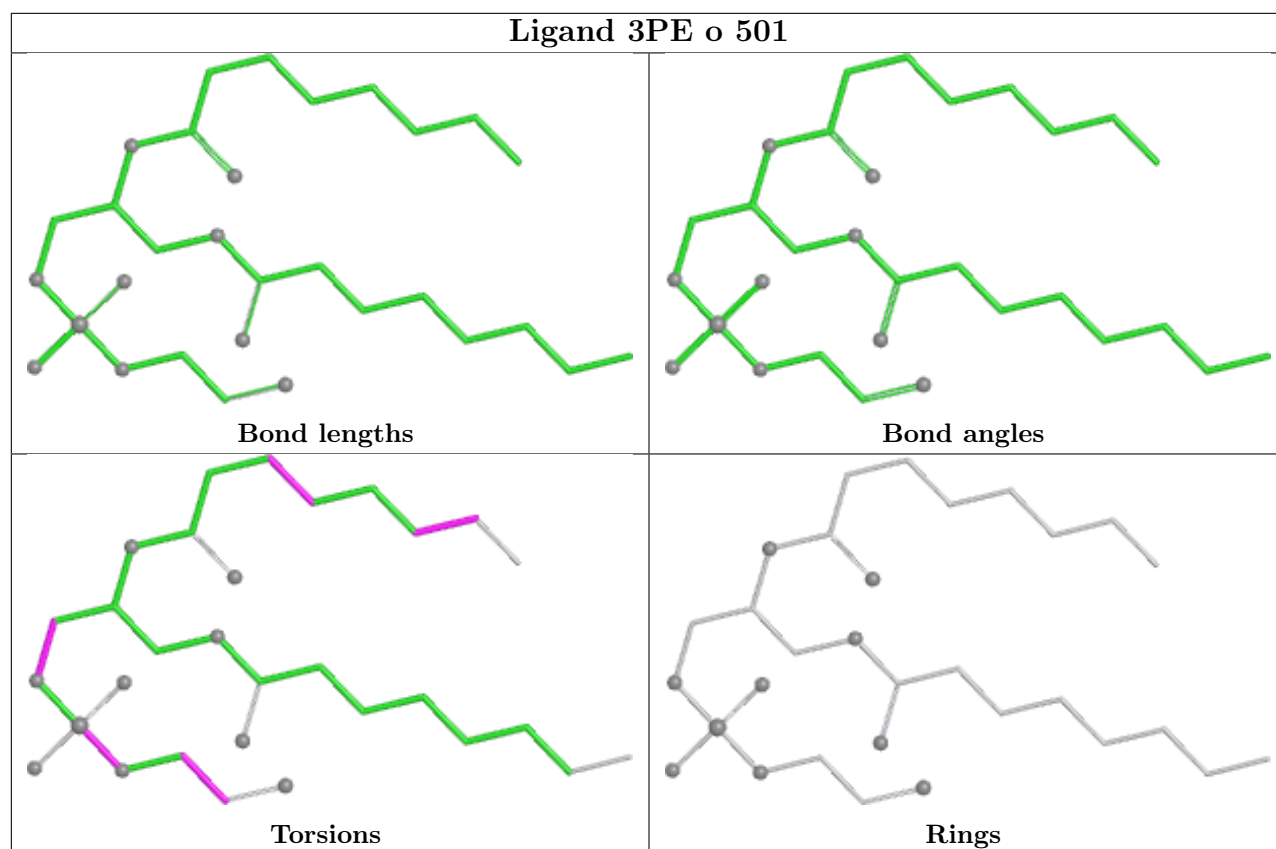
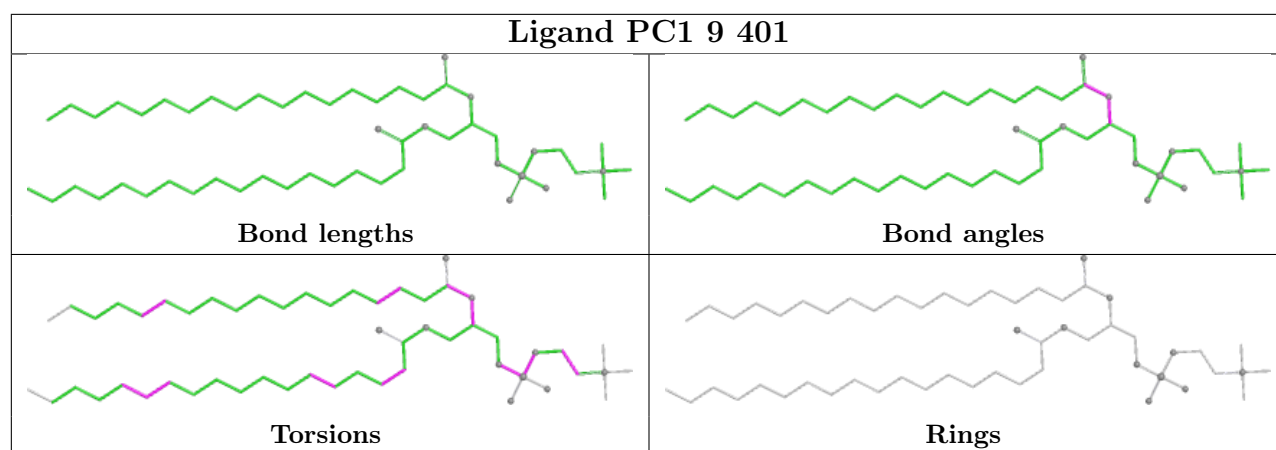


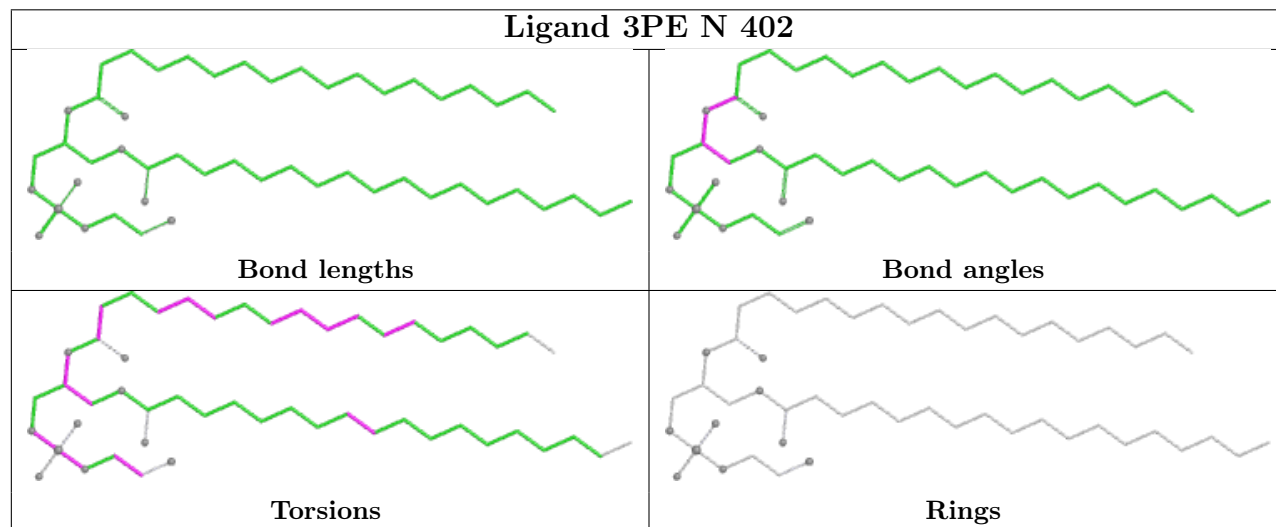
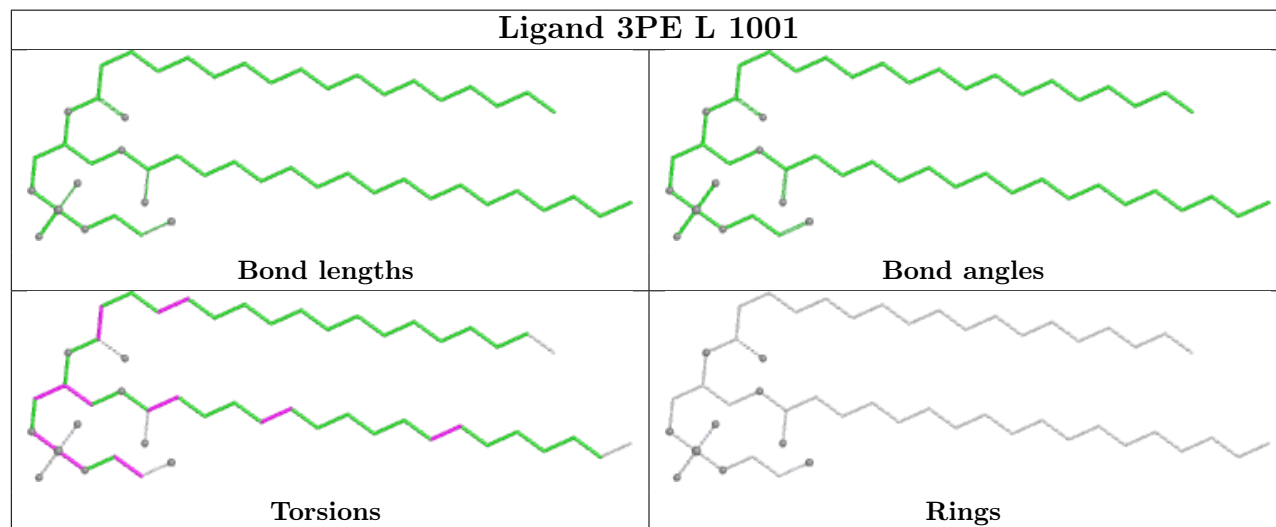
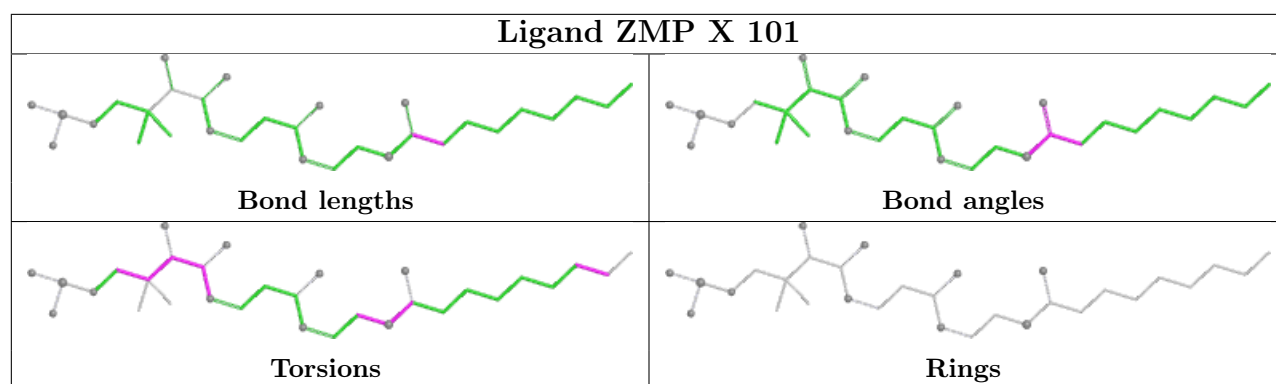


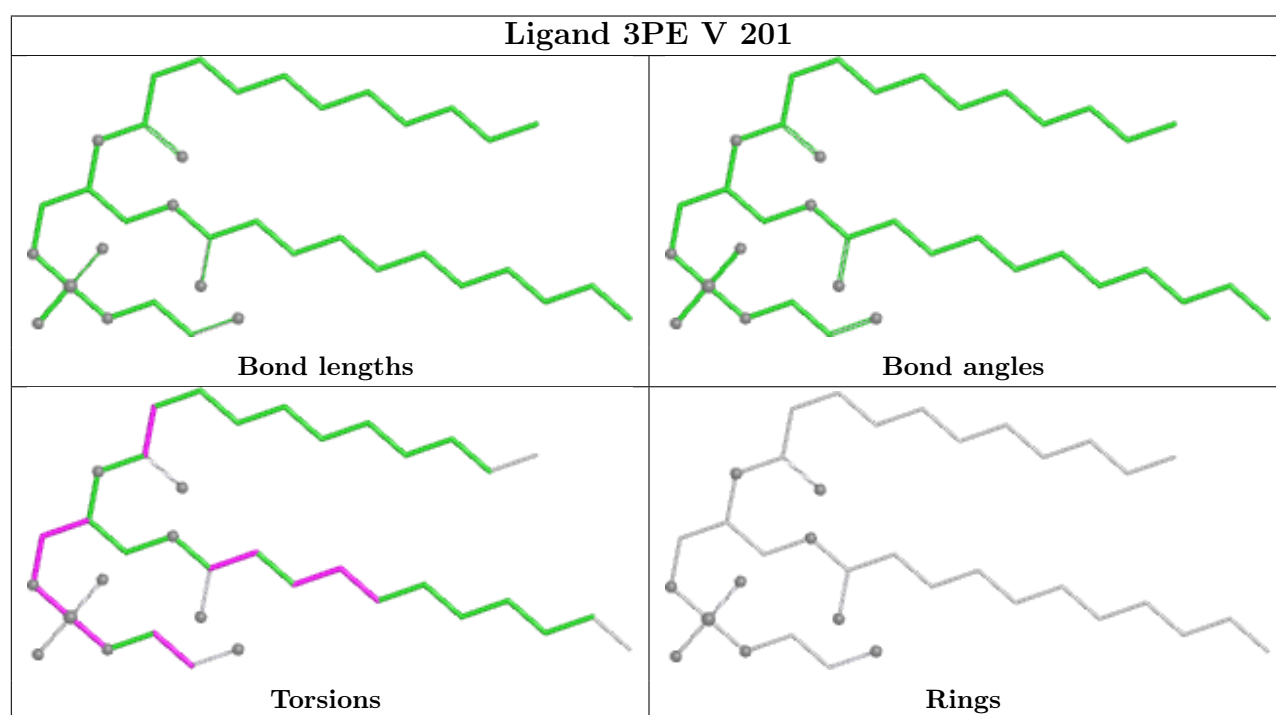
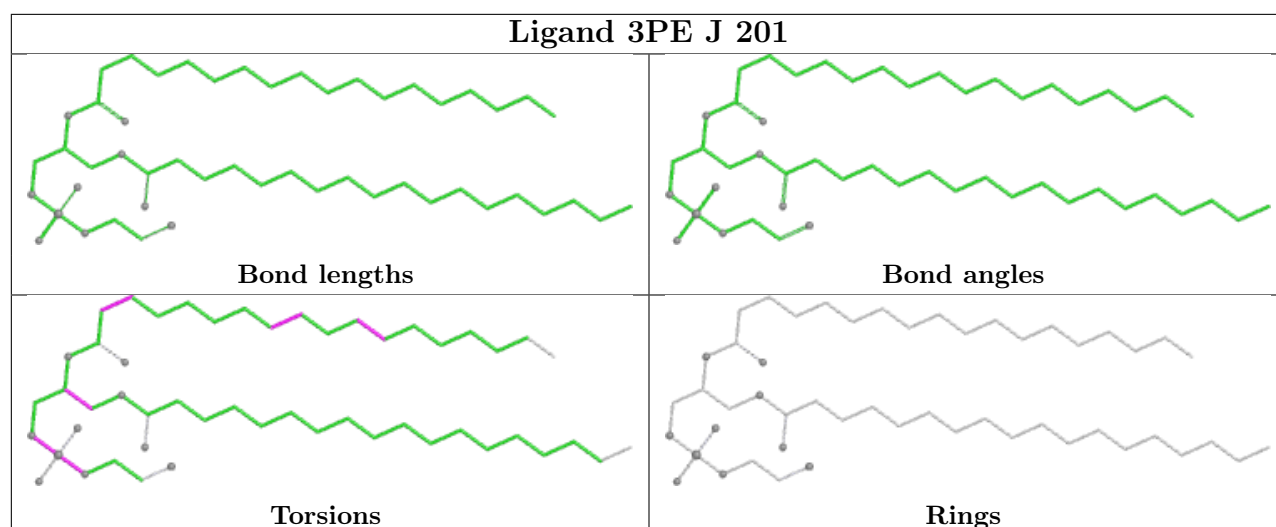


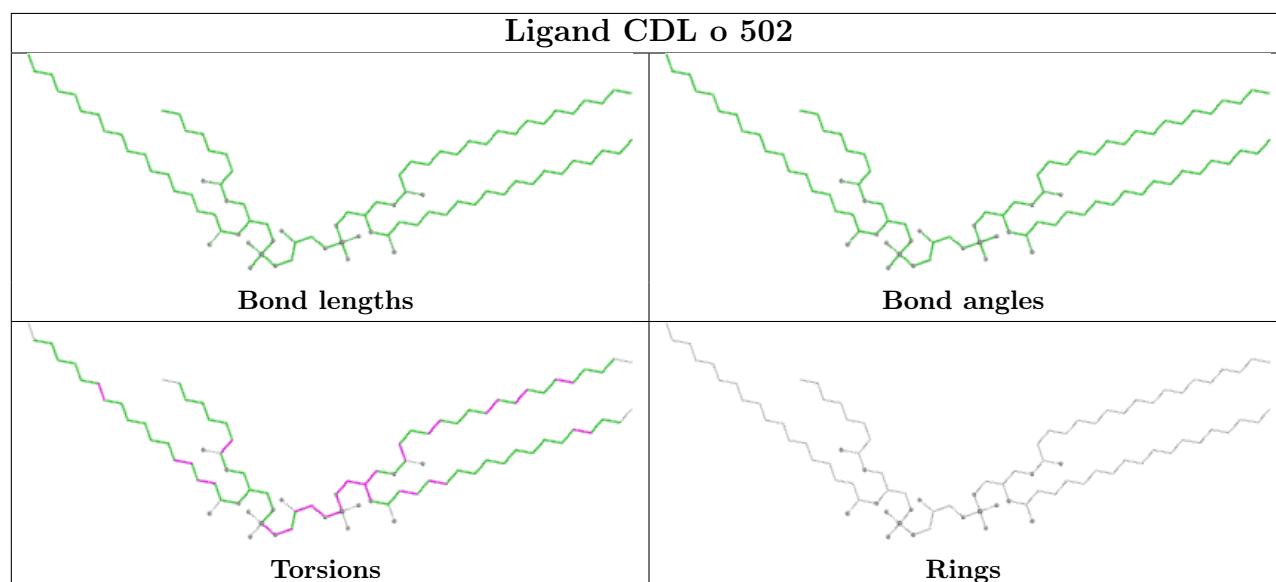
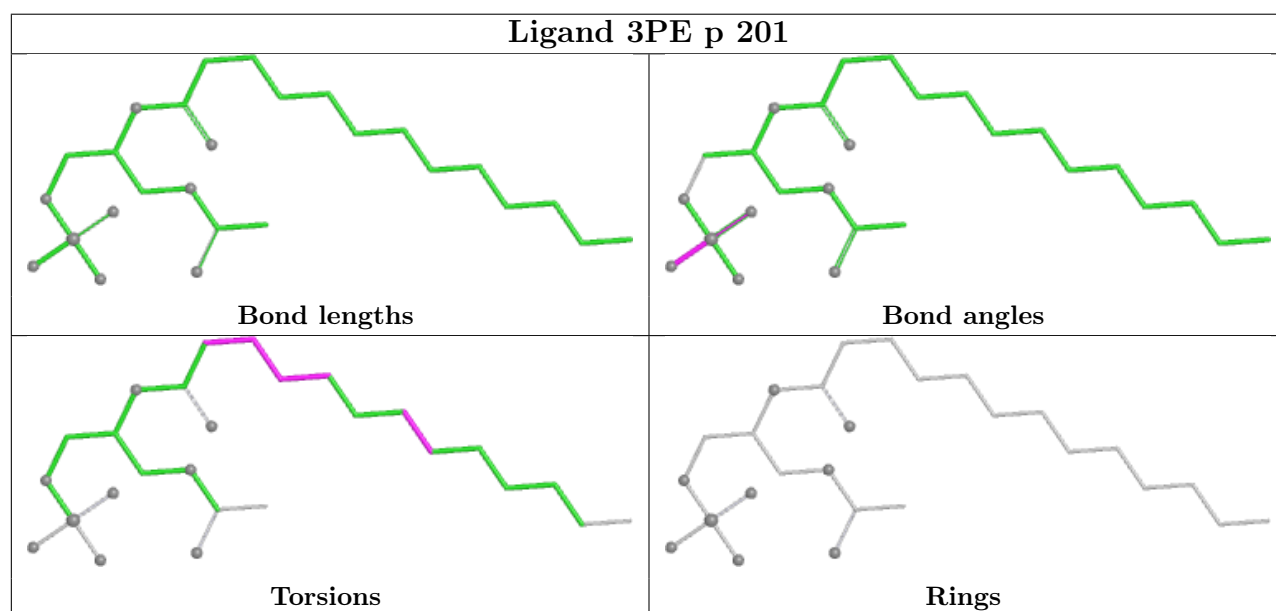
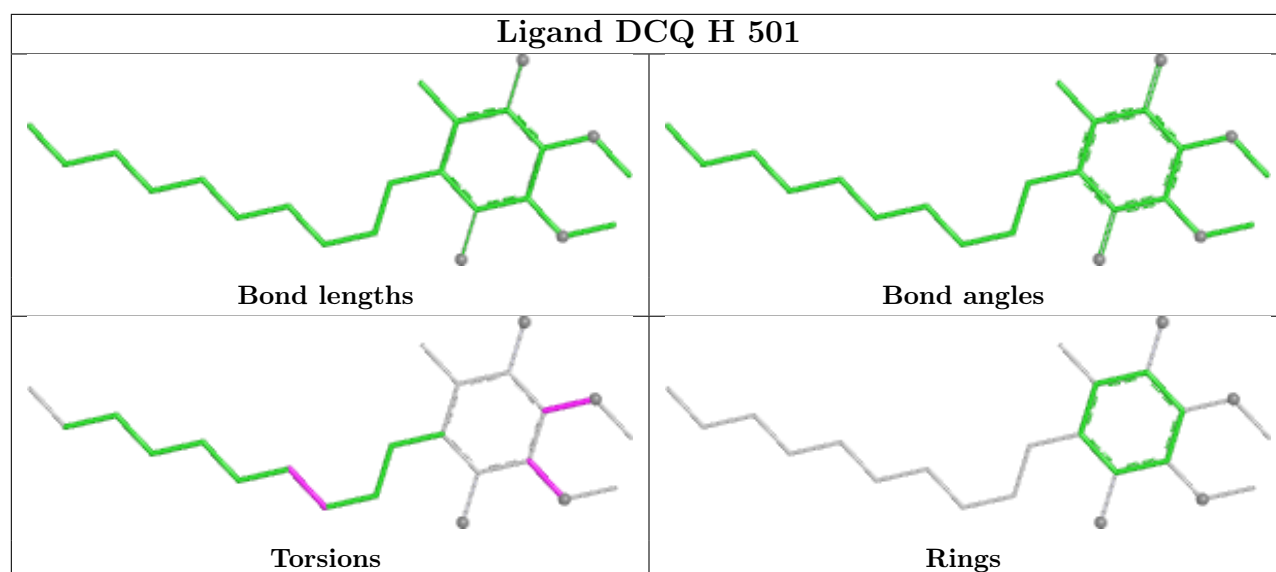


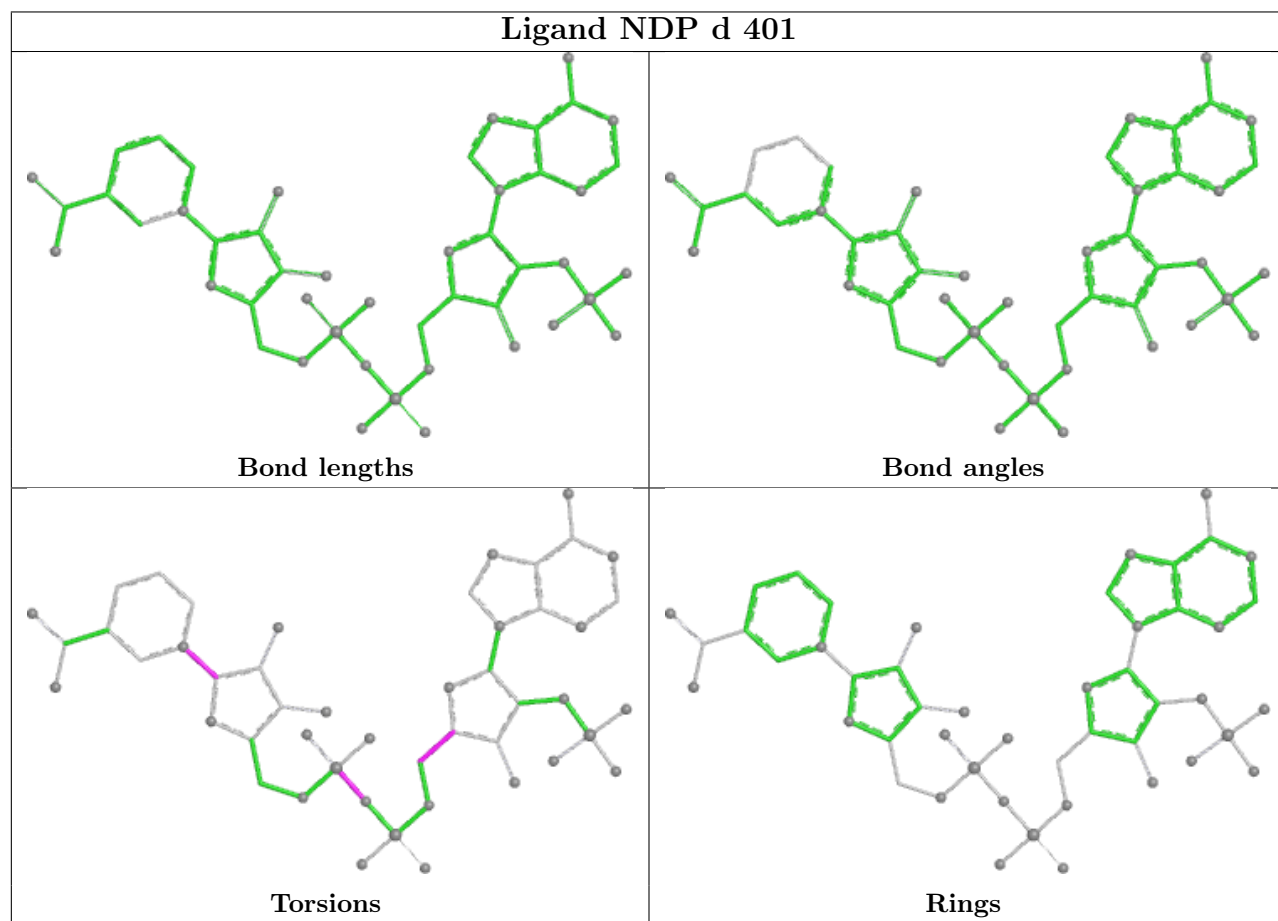
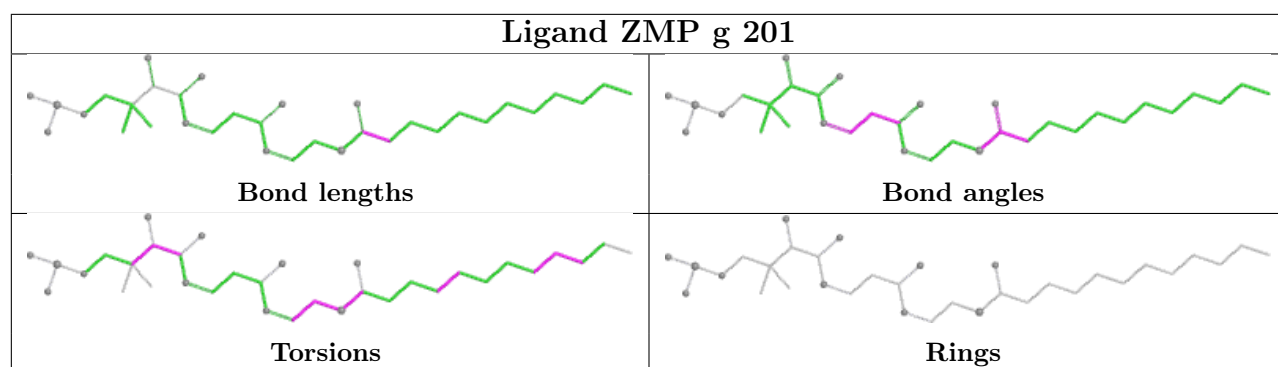


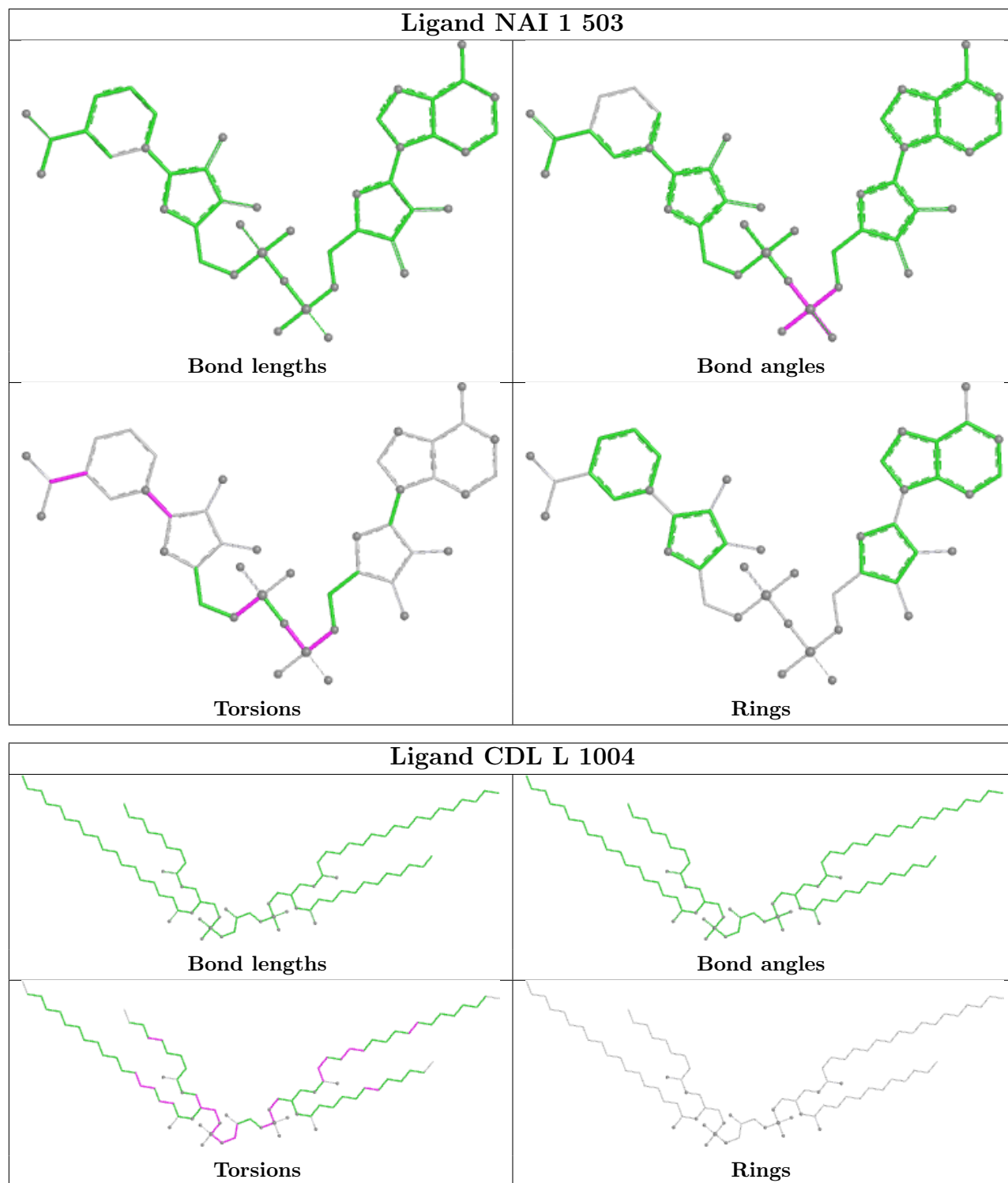


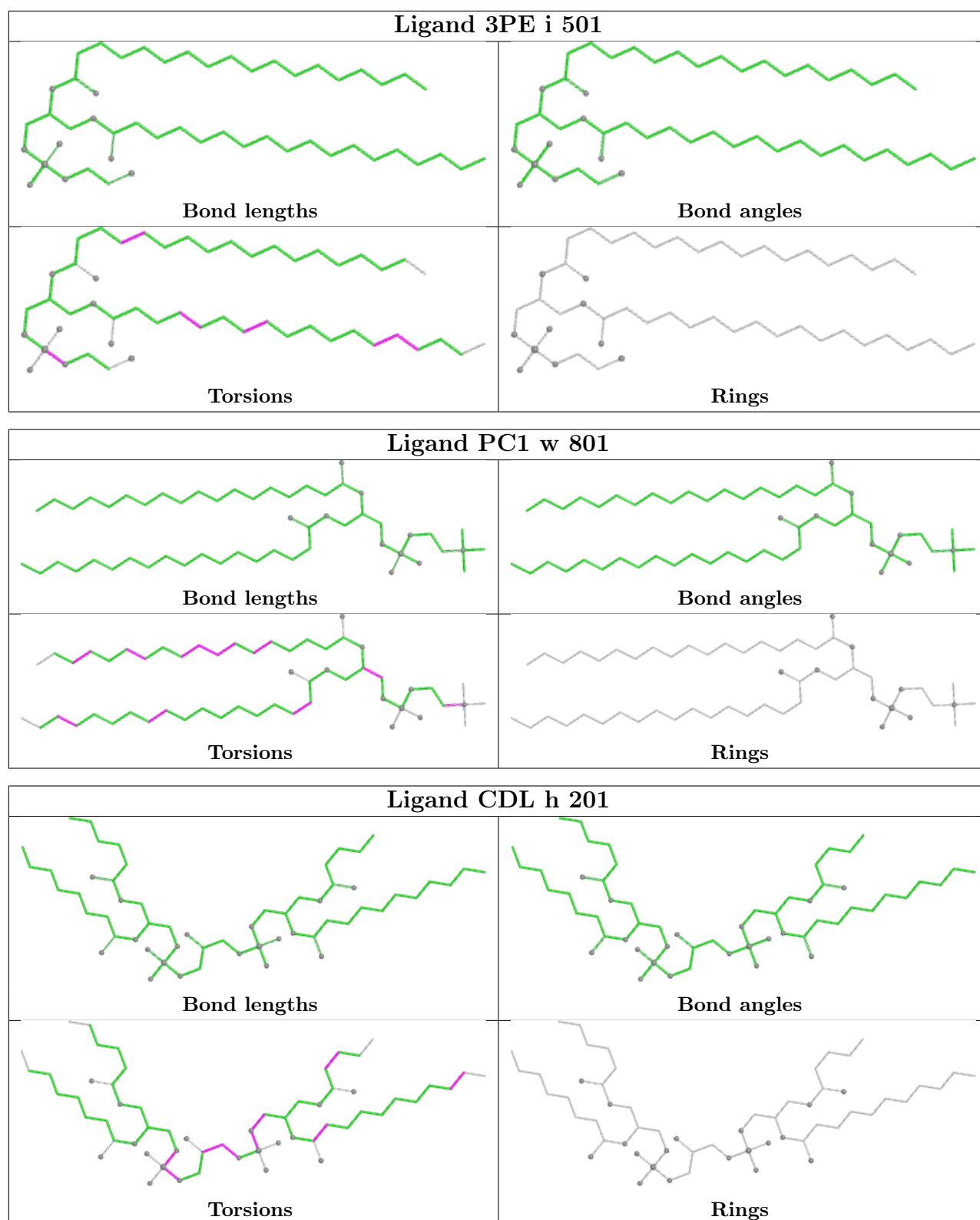


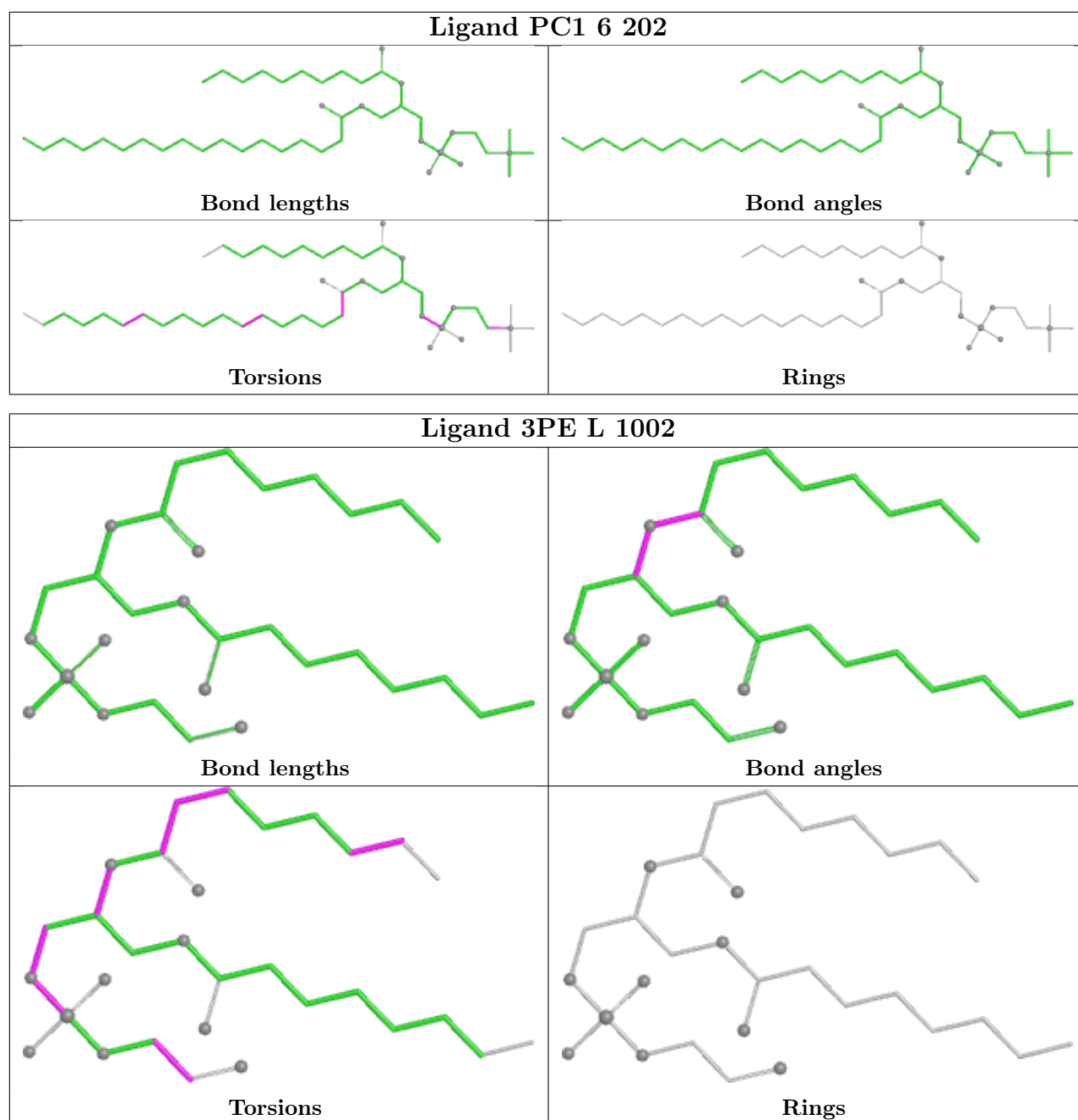












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

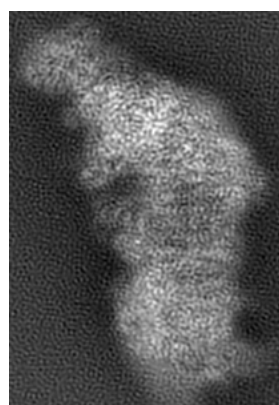
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-11244. These allow visual inspection of the internal detail of the map and identification of artifacts.

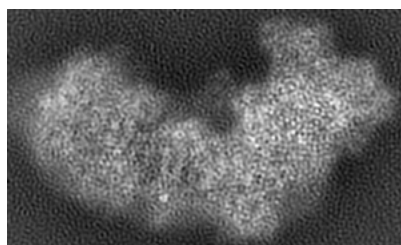
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

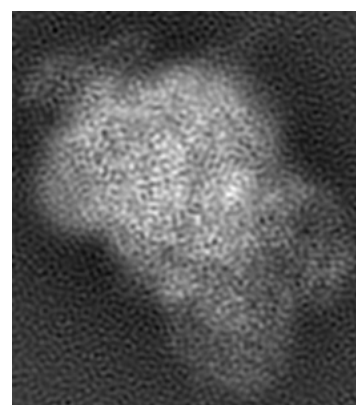
6.1.1 Primary map



X



Y

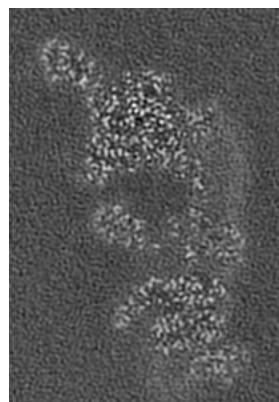


Z

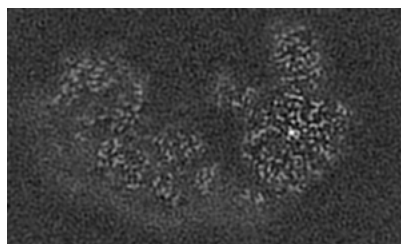
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

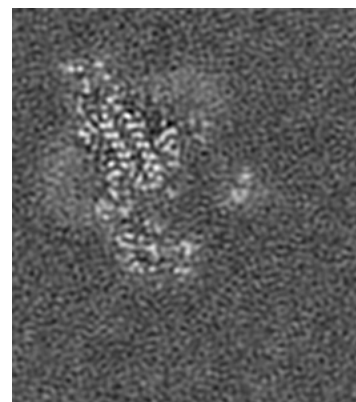
6.2.1 Primary map



X Index: 80



Y Index: 91

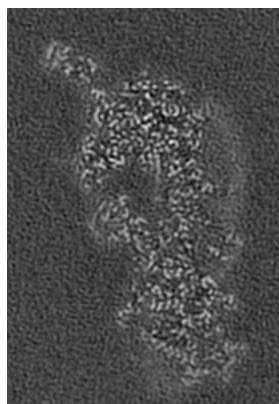


Z Index: 133

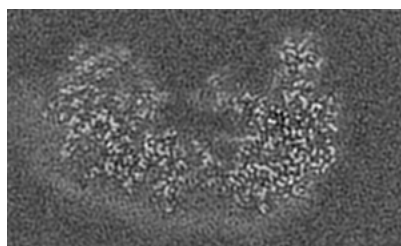
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

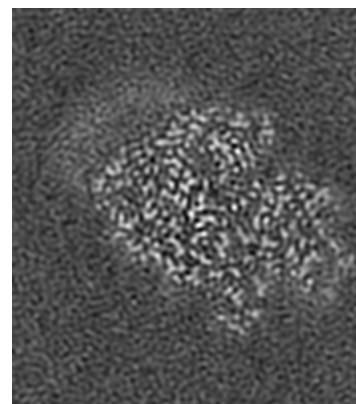
6.3.1 Primary map



X Index: 74



Y Index: 100

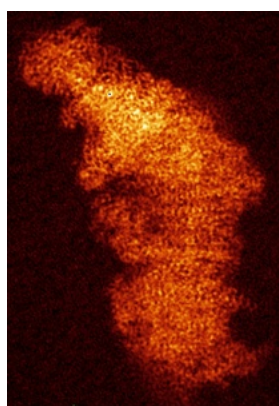


Z Index: 195

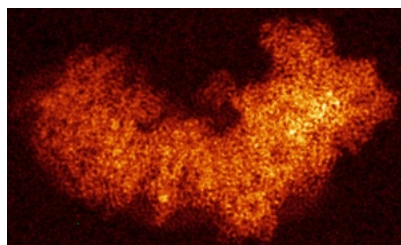
The images above show the largest variance slices of the map in three orthogonal directions.

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

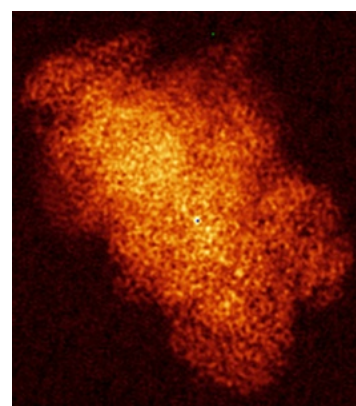
6.4.1 Primary map



X



Y



Z

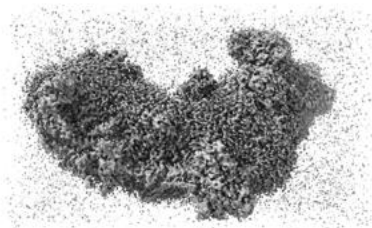
The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views [i](#)

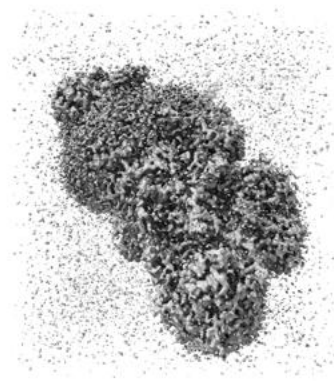
6.5.1 Primary map



X



Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.045. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

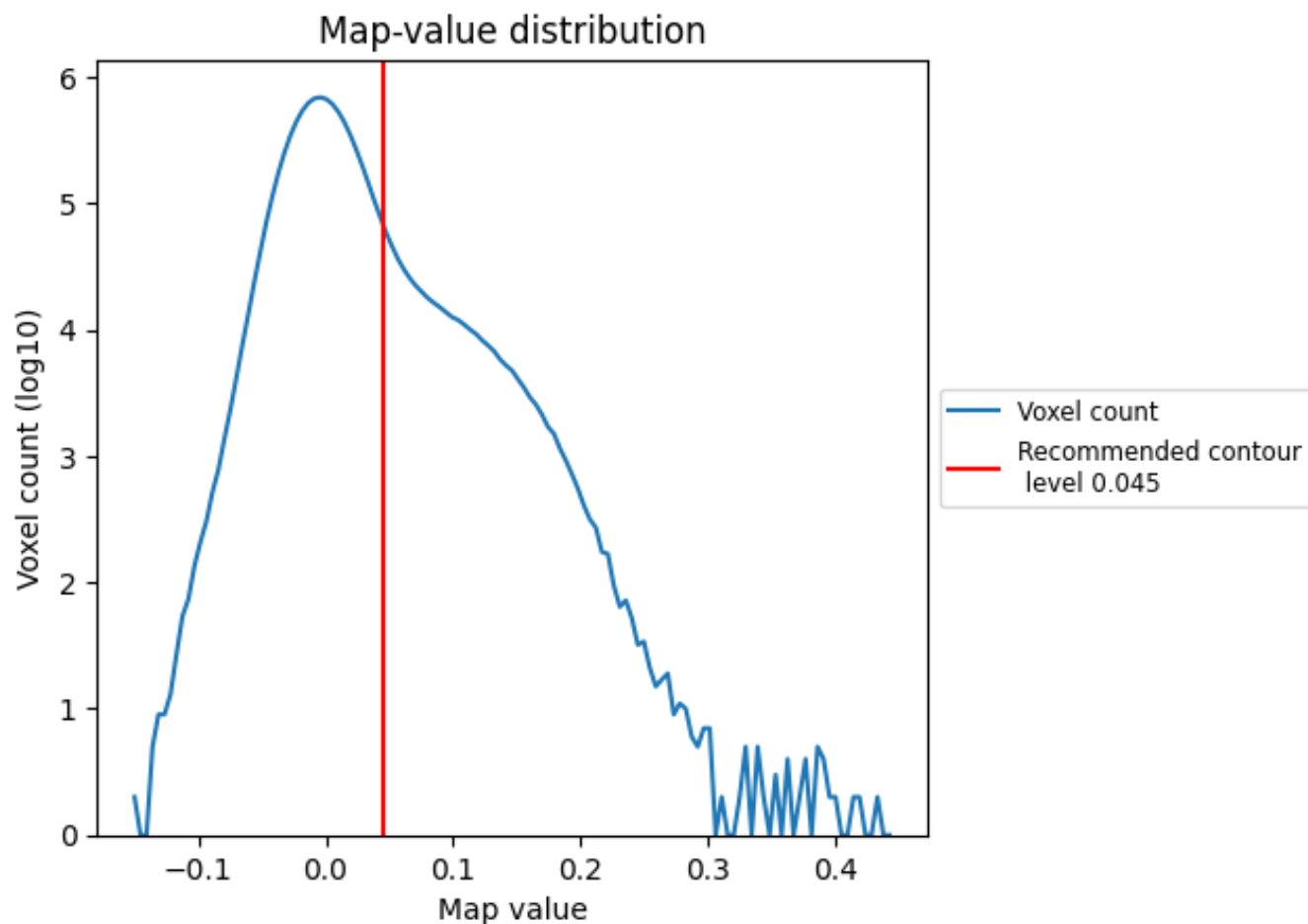
6.6 Mask visualisation [i](#)

This section was not generated. No masks/segmentation were deposited.

7 Map analysis [i](#)

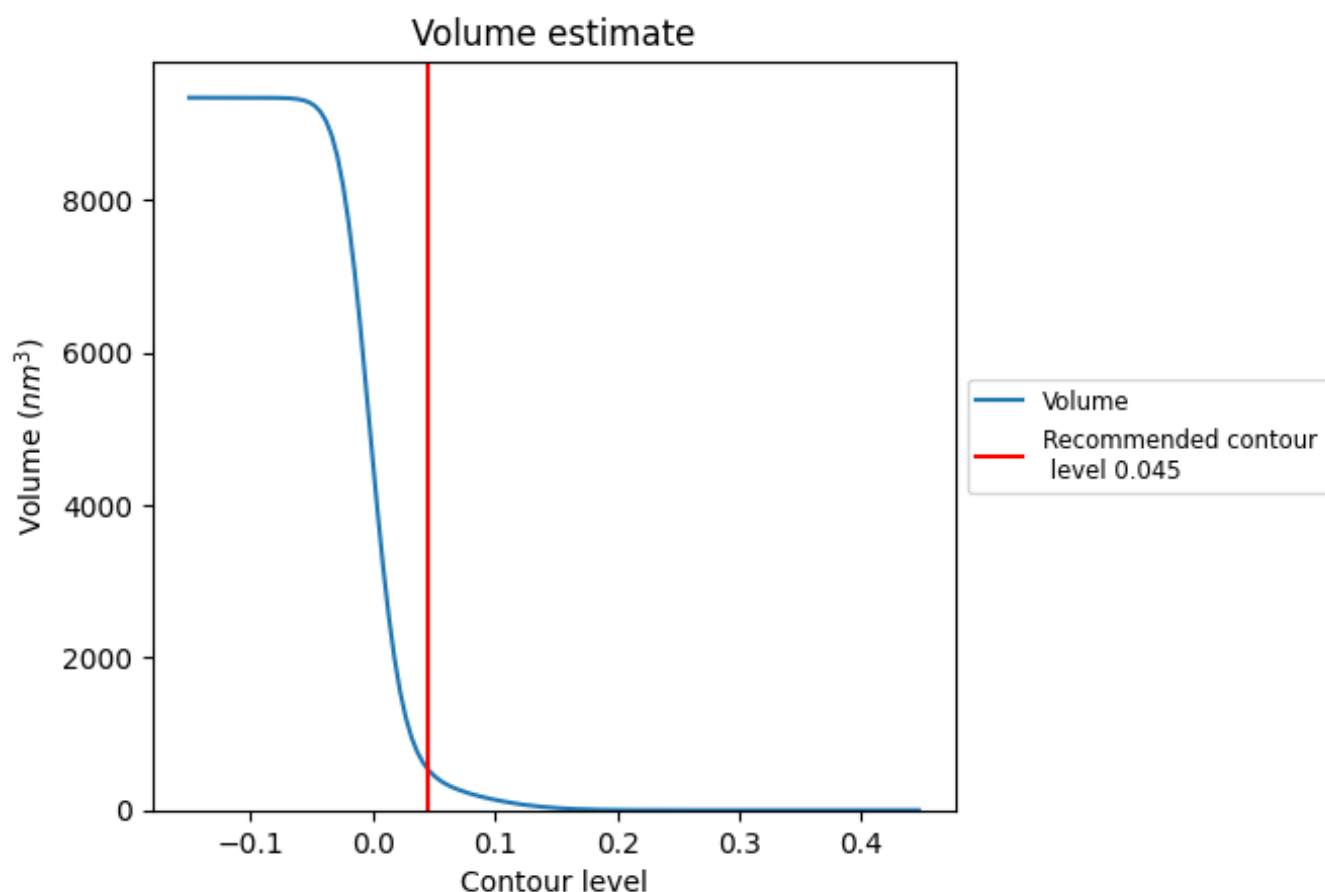
This section contains the results of statistical analysis of the map.

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 543 nm^3 ; this corresponds to an approximate mass of 491 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.

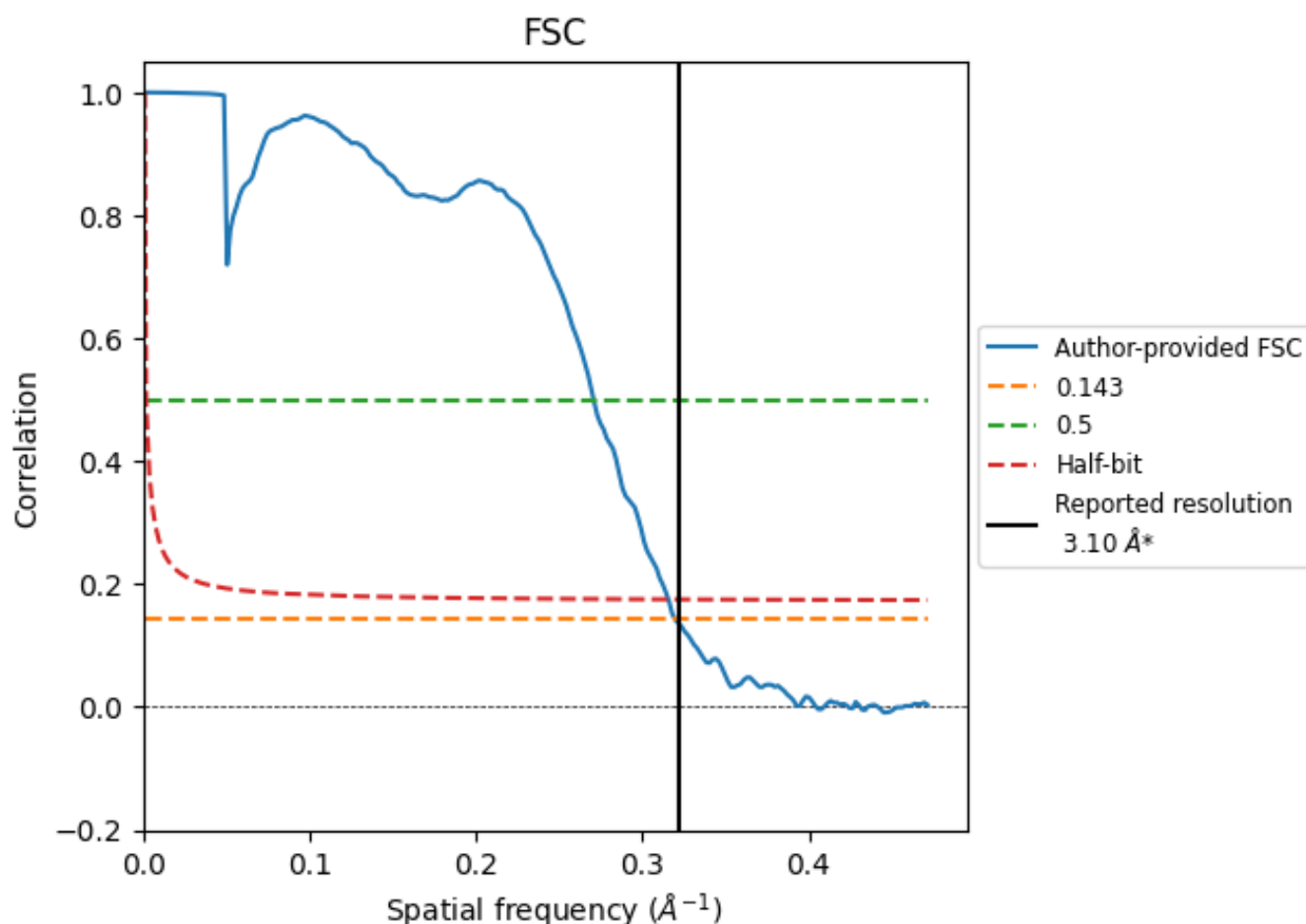
7.3 Rotationally averaged power spectrum [i](#)

This section was not generated. The rotationally averaged power spectrum is only generated for cubic maps.

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 0.323 Å⁻¹

8.2 Resolution estimates [i](#)

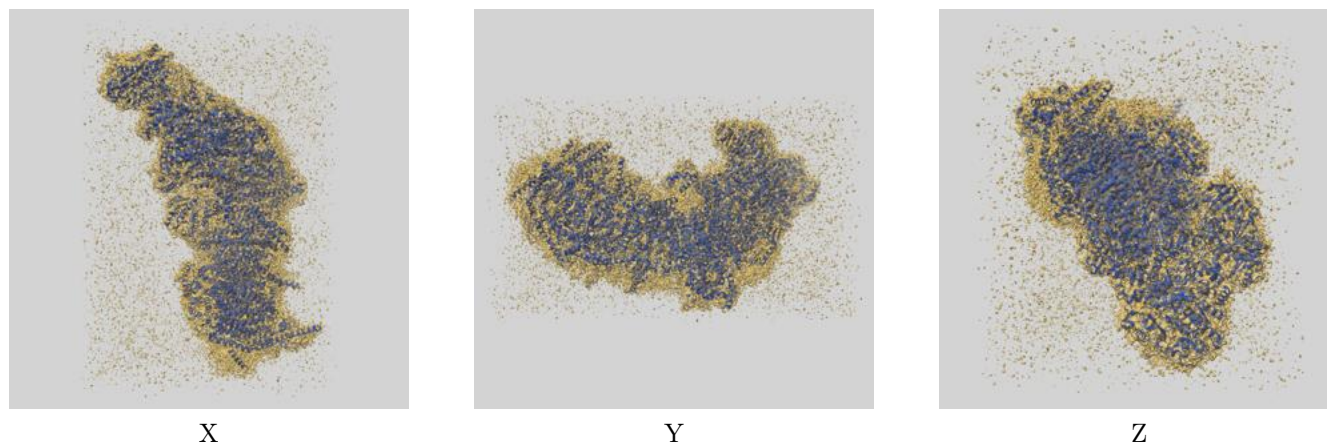
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.10	-	-
Author-provided FSC curve	3.13	3.69	3.17
Unmasked-calculated*	-	-	-

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps.

9 Map-model fit [i](#)

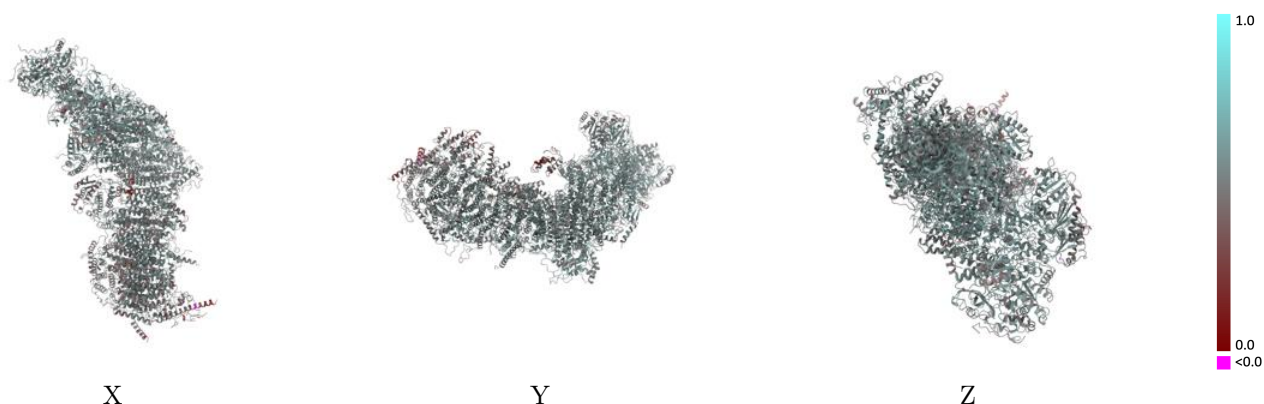
This section contains information regarding the fit between EMDB map EMD-11244 and PDB model 6ZKC. Per-residue inclusion information can be found in section 3 on page 21.

9.1 Map-model overlay [i](#)



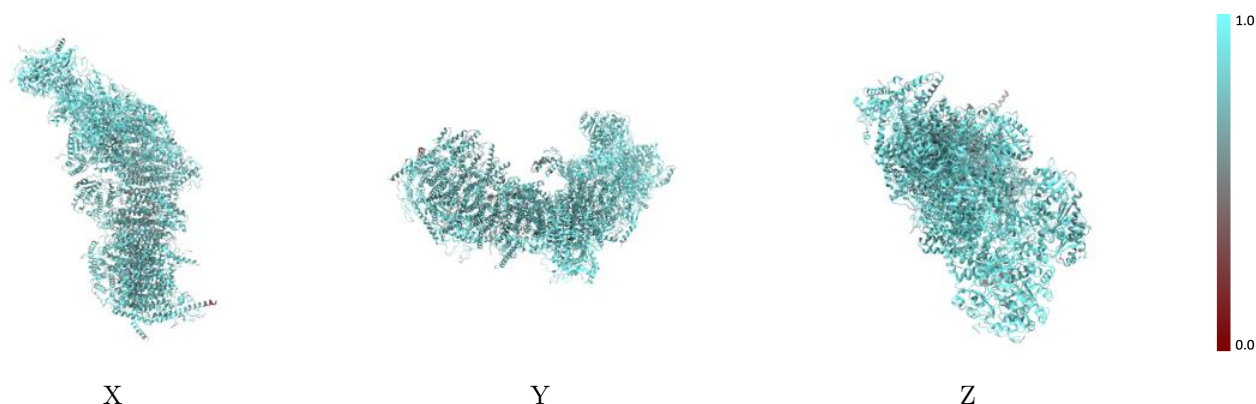
The images above show the 3D surface view of the map at the recommended contour level 0.045 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



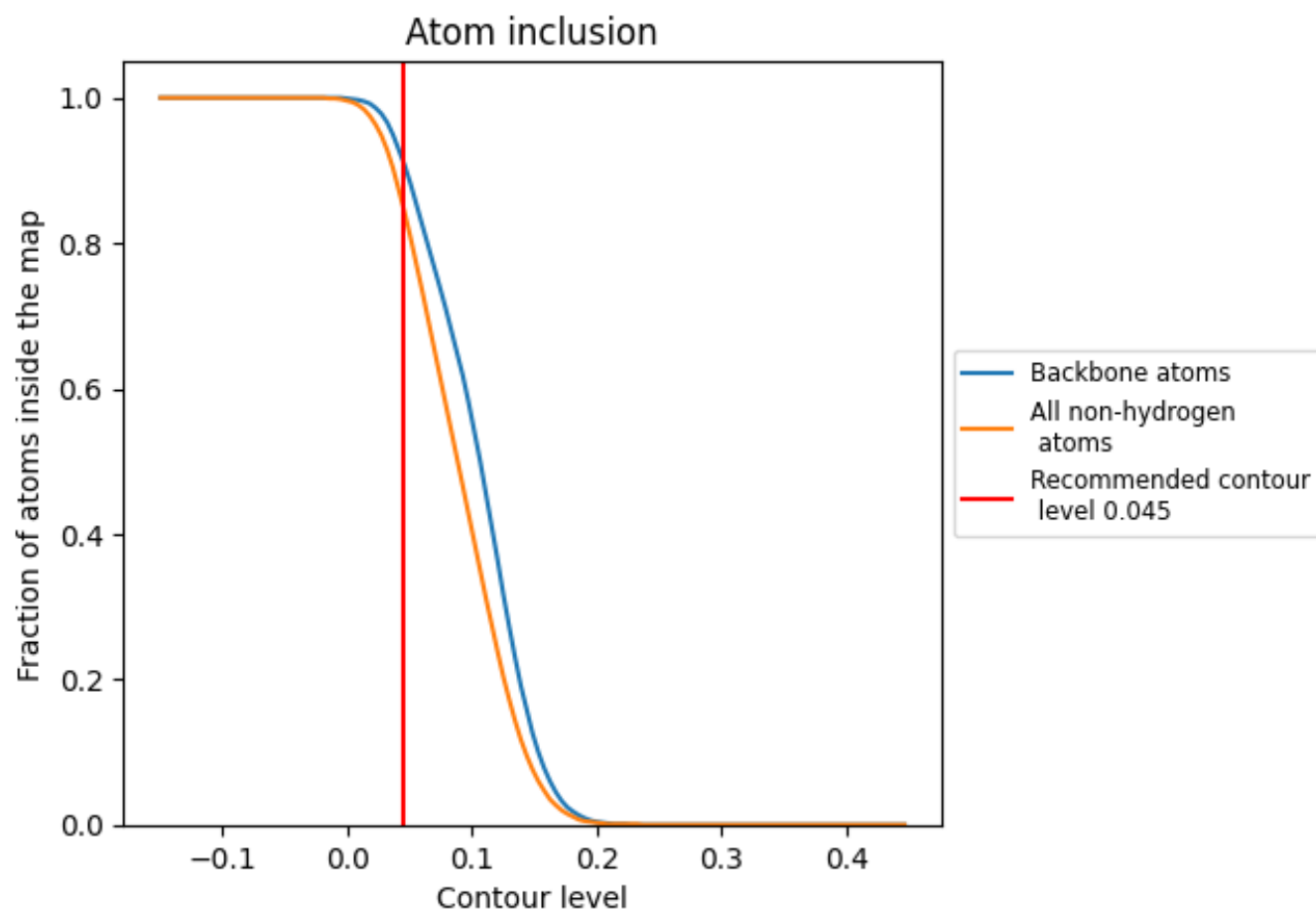
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.045).

9.4 Atom inclusion [i](#)



At the recommended contour level, 91% of all backbone atoms, 85% 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 (0.045) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8480	 0.5230
1	 0.8920	 0.5250
2	 0.8730	 0.5210
3	 0.8810	 0.5400
4	 0.8790	 0.5590
5	 0.9140	 0.5680
6	 0.9160	 0.5670
9	 0.9110	 0.5680
A	 0.8150	 0.5420
H	 0.8810	 0.5550
J	 0.8380	 0.5390
K	 0.8300	 0.5400
L	 0.8030	 0.5100
M	 0.8660	 0.5520
N	 0.8630	 0.5540
V	 0.7780	 0.4860
W	 0.8410	 0.5360
X	 0.7580	 0.4540
Y	 0.8690	 0.5190
Z	 0.8300	 0.5030
a	 0.8730	 0.5060
b	 0.8930	 0.5500
c	 0.8650	 0.5510
d	 0.8690	 0.5360
e	 0.8390	 0.4930
f	 0.8500	 0.5220
g	 0.8480	 0.5360
h	 0.8550	 0.5320
i	 0.8630	 0.5450
j	 0.6930	 0.3810
k	 0.8250	 0.5020
l	 0.8340	 0.5170
m	 0.8580	 0.5190
n	 0.7520	 0.4260
o	 0.8300	 0.5230



Continued on next page...

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Chain	Atom inclusion	Q-score
p	 0.7780	 0.4740
q	 0.8810	 0.5320
r	 0.8090	 0.4690
s	 0.7880	 0.4260
t	 0.7960	 0.4610
u	 0.8250	 0.4570
v	 0.7980	 0.4700
w	 0.7990	 0.5020
x	 0.7910	 0.4950
y	 0.7780	 0.4810
z	 0.8720	 0.5270