



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 03:24 AM UTC

PDB ID : 1NEK / pdb_00001nek
Title : Complex II (Succinate Dehydrogenase) From E. Coli with ubiquinone bound
Authors : Yankovskaya, V.; Horsefield, R.; Tornroth, S.; Luna-Chavez, C.; Miyoshi, H.;
Leger, C.; Byrne, B.; Cecchini, G.; Iwata, S.
Deposited on : 2002-12-11
Resolution : 2.60 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

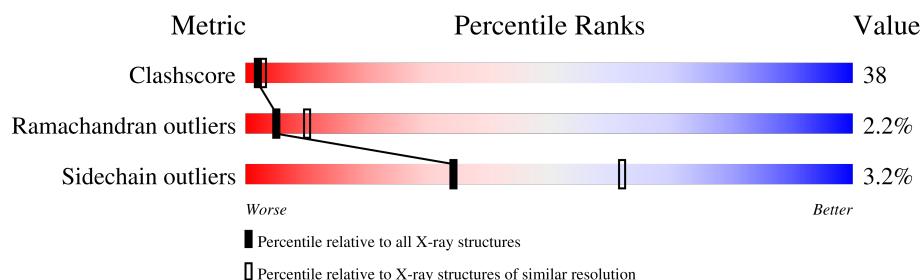
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)

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

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	588	36% 56% 7%
2	B	238	51% 47% .
3	C	129	66% 31% ..
4	D	115	61% 35% ..

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
10	F3S	B	304	-	-	X	-
12	CDN	C	308	X	-	-	-
8	FES	B	302	-	-	X	-

2 Entry composition

There are 15 unique types of molecules in this entry. The entry contains 8698 atoms, of which 0 are hydrogens and 0 are deuteriums.

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

- Molecule 1 is a protein called Succinate dehydrogenase flavoprotein subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	588	Total	C	N	O	S	89	0	0
			4522	2812	821	861	28			

- Molecule 2 is a protein called Succinate dehydrogenase iron-sulfur protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	238	Total	C	N	O	S	0	0	0
			1869	1172	329	348	20			

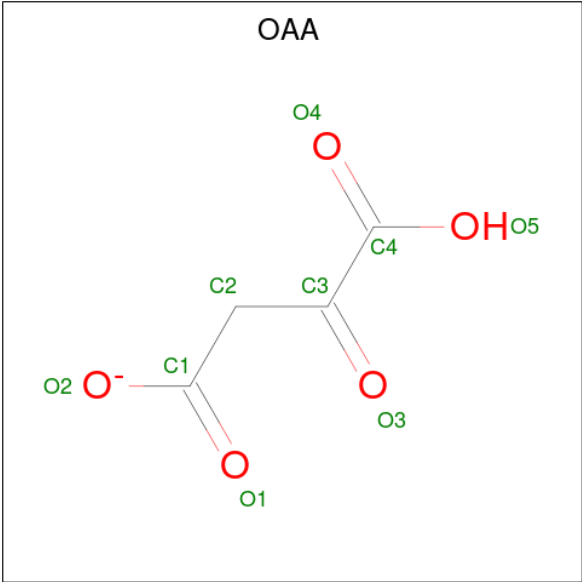
- Molecule 3 is a protein called Succinate dehydrogenase cytochrome b-556 subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	129	Total	C	N	O	S	0	0	0
			1008	668	166	168	6			

- Molecule 4 is a protein called Succinate dehydrogenase hydrophobic membrane anchor protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	113	Total	C	N	O	S	0	0	0
			898	615	136	144	3			

- Molecule 5 is OXALOACETATE ION (CCD ID: OAA) (formula: $C_4H_3O_5$).

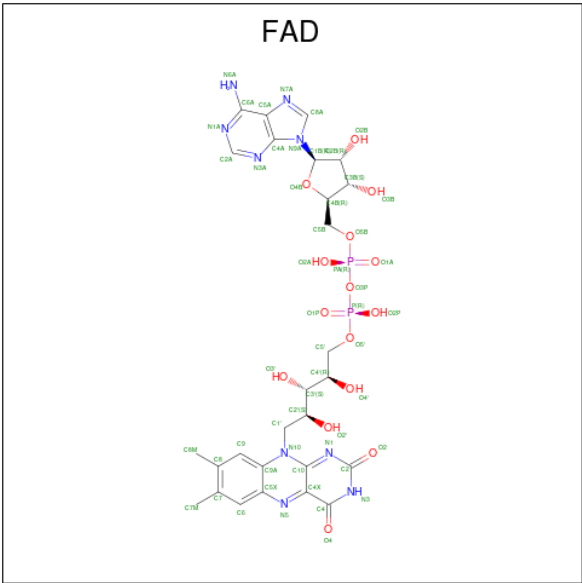


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			9	4	5		

- Molecule 6 is CALCIUM ION (CCD ID: CA) (formula: Ca).

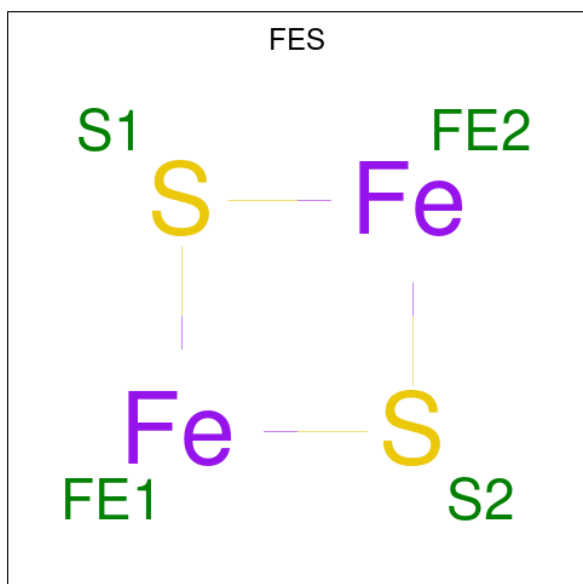
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total	Ca	0	0
			1	1		
6	B	1	Total	Ca	0	0
			1	1		

- Molecule 7 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: C₂₇H₃₃N₉O₁₅P₂).



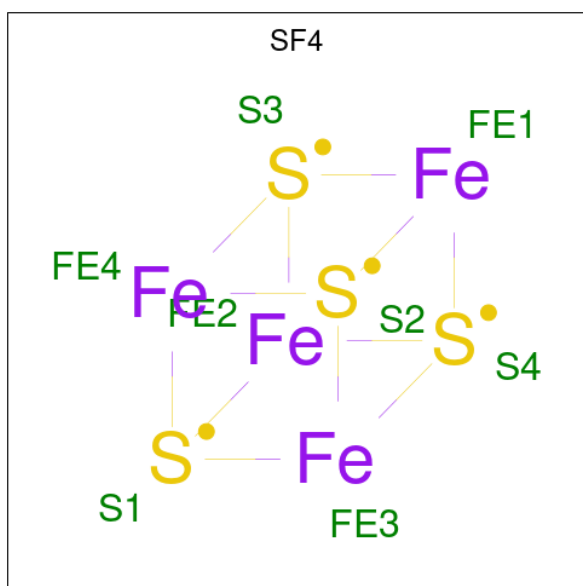
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 8 is FE2/S2 (INORGANIC) CLUSTER (CCD ID: FES) (formula: Fe_2S_2).



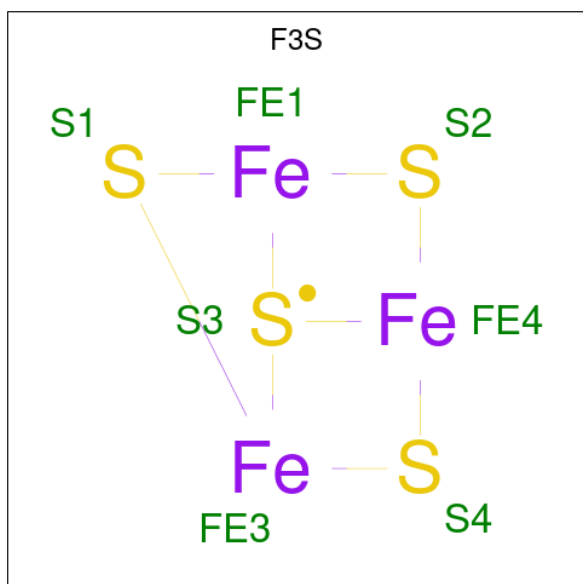
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	B	1	Total	Fe	S	0	0
			4	2	2		

- Molecule 9 is IRON/SULFUR CLUSTER (CCD ID: SF4) (formula: Fe_4S_4).



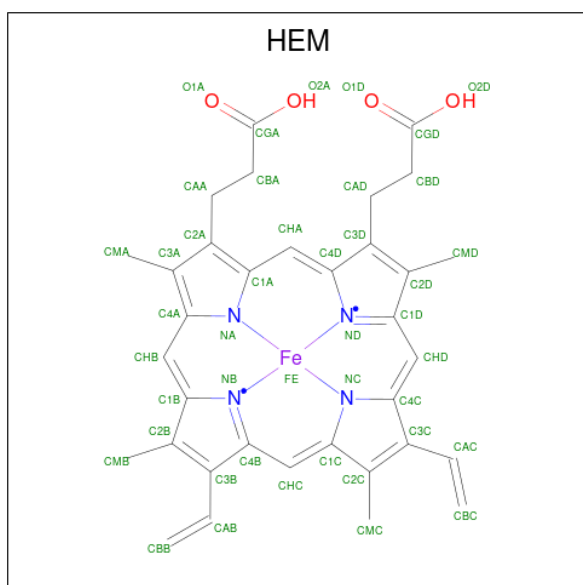
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	B	1	Total	Fe	S	0	0
			8	4	4		

- Molecule 10 is FE3-S4 CLUSTER (CCD ID: F3S) (formula: Fe_3S_4).



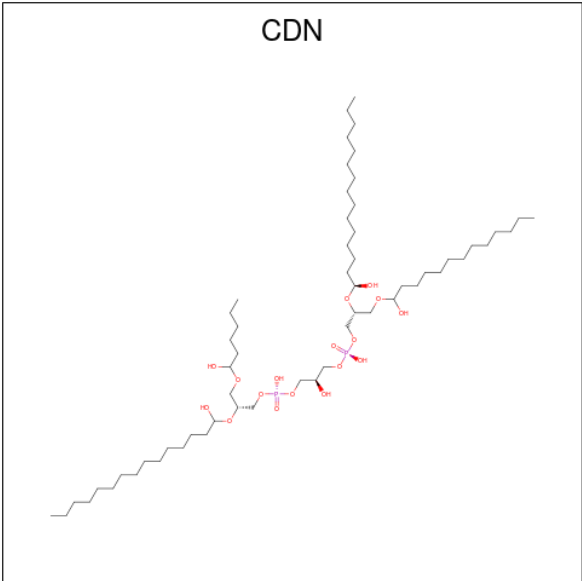
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
10	B	1	Total	Fe	S	0	0
			7	3	4		

- Molecule 11 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: $\text{C}_{34}\text{H}_{32}\text{FeN}_4\text{O}_4$).



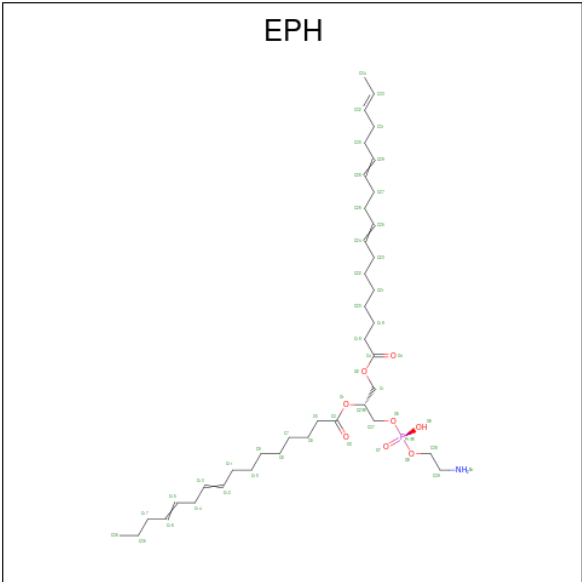
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
11	C	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- Molecule 12 is CARDIOLIPIN (CCD ID: CDN) (formula: C₅₈H₁₂₀O₁₇P₂).



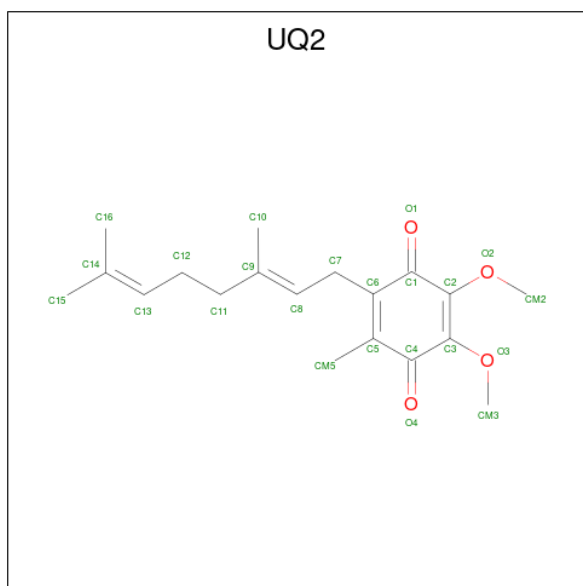
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
12	C	1	Total	C	O	P	0	0
			77	58	17	2		

- Molecule 13 is L-ALPHA-PHOSPHATIDYL-BETA-OLEOYL-GAMMA-PALMITOYL-PHOSPHATIDYLETHANOLAMINE (CCD ID: EPH) (formula: C₃₉H₆₈NO₈P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
13	C	1	Total	C	N	O	P	0	0
			35	25	1	8	1		

- Molecule 14 is UBIQUINONE-2 (CCD ID: UQ2) (formula: $C_{19}H_{26}O_4$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
14	D	1	Total	C	O	0	0
			23	19	4		

- Molecule 15 is water.

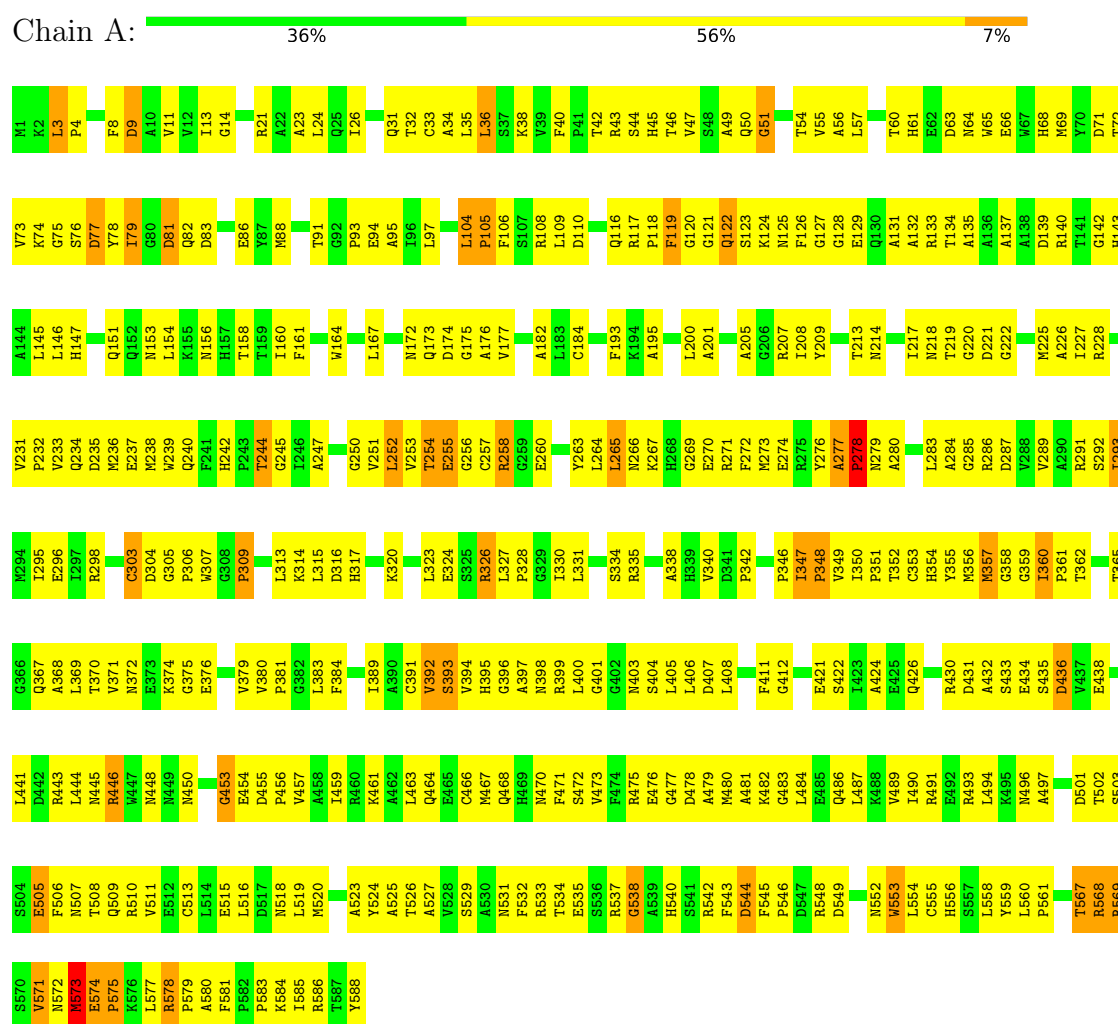
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
15	A	61	Total	O	0	0
			61	61		
15	B	50	Total	O	0	0
			50	50		
15	C	15	Total	O	0	0
			15	15		
15	D	14	Total	O	0	0
			14	14		

3 Residue-property plots

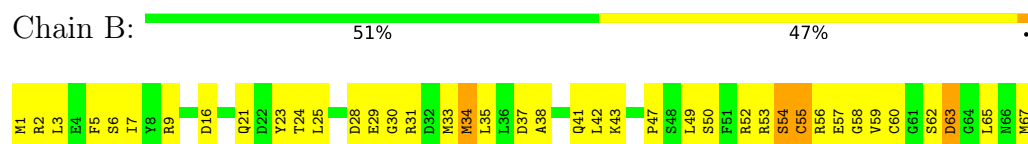
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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. 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.

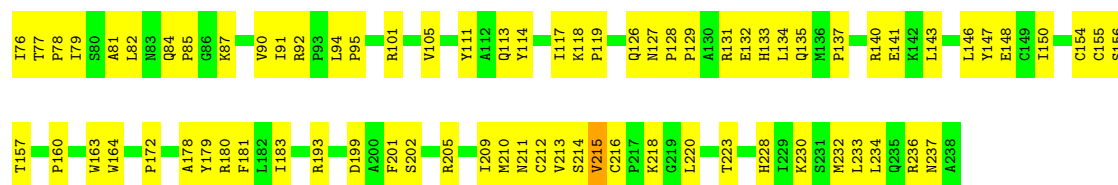
Note EDS was not executed.

- Molecule 1: Succinate dehydrogenase flavoprotein subunit



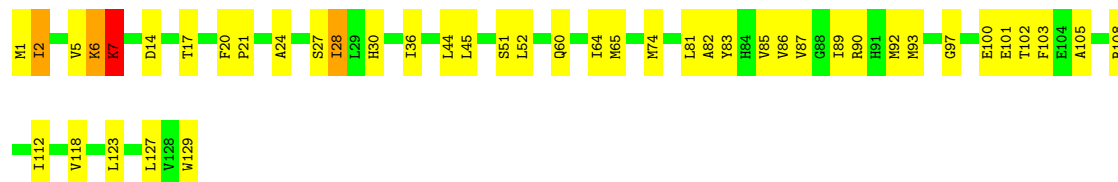
- Molecule 2: Succinate dehydrogenase iron-sulfur protein





• Molecule 3: Succinate dehydrogenase cytochrome b-556 subunit

Chain C: 66% 31% ..



• Molecule 4: Succinate dehydrogenase hydrophobic membrane anchor protein

Chain D: 61% 35% ..



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	138.80Å 138.80Å 521.90Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	40.00 – 2.60	Depositor
% Data completeness (in resolution range)	(Not available) (40.00-2.60)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.247 , 0.289	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	8698	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: HEM, F3S, CA, UQ2, OAA, FAD, EPH, CDN, FES, SF4

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.47	0/4611	1.07	31/6237 (0.5%)
2	B	0.51	0/1908	1.02	9/2578 (0.3%)
3	C	0.50	0/1030	1.03	5/1394 (0.4%)
4	D	0.54	0/923	0.99	2/1262 (0.2%)
All	All	0.49	0/8472	1.04	47/11471 (0.4%)

There are no bond length outliers.

All (47) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	571	VAL	N-CA-C	8.27	118.61	108.53
1	A	121	GLY	N-CA-C	-8.22	100.37	111.54
1	A	79	ILE	N-CA-C	-7.88	105.51	113.47
1	A	254	THR	N-CA-C	7.72	121.15	110.35
1	A	347	ILE	N-CA-C	7.56	116.14	107.89
1	A	252	LEU	N-CA-C	7.40	120.93	110.50
1	A	392	VAL	N-CA-C	-7.31	103.13	113.07
1	A	104	LEU	CA-C-N	7.13	128.76	119.84
1	A	104	LEU	C-N-CA	7.13	128.76	119.84
1	A	567	THR	N-CA-C	6.87	121.02	110.36
2	B	62	SER	N-CA-C	6.84	118.39	111.07
1	A	360	ILE	CA-C-N	6.60	126.96	119.83
1	A	360	ILE	C-N-CA	6.60	126.96	119.83
2	B	7	ILE	N-CA-C	6.56	117.93	108.48
1	A	255	GLU	N-CA-C	-6.39	104.74	112.54
4	D	112	VAL	N-CA-C	6.33	116.50	110.42
1	A	277	ALA	CA-C-N	6.18	127.57	119.84
1	A	277	ALA	C-N-CA	6.18	127.57	119.84
3	C	6	LYS	N-CA-C	6.13	119.64	109.96
2	B	178	ALA	N-CA-C	-6.13	104.68	111.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	63	ASP	N-CA-C	6.03	120.50	111.81
3	C	7	LYS	N-CA-C	6.00	123.58	110.80
1	A	575	PRO	N-CA-C	5.98	121.40	113.57
1	A	574	GLU	CA-C-N	5.95	125.49	119.19
1	A	574	GLU	C-N-CA	5.95	125.49	119.19
1	A	412	GLY	N-CA-C	-5.91	105.20	112.77
1	A	265	LEU	N-CA-C	5.90	118.84	109.81
3	C	90	ARG	N-CA-C	-5.78	104.18	111.11
2	B	215	VAL	N-CA-C	5.75	120.30	113.22
1	A	574	GLU	N-CA-C	5.74	122.49	109.81
1	A	244	THR	N-CA-C	5.73	120.04	112.25
1	A	573	MET	N-CA-C	5.71	116.94	108.60
2	B	16	ASP	N-CA-C	-5.66	106.26	114.12
1	A	3	LEU	CA-C-N	5.52	125.46	119.78
1	A	3	LEU	C-N-CA	5.52	125.46	119.78
3	C	65	MET	N-CA-C	-5.42	105.76	112.38
1	A	9	ASP	N-CA-C	-5.42	106.65	113.20
1	A	258	ARG	N-CA-C	-5.38	105.52	111.71
2	B	154	CYS	N-CA-C	-5.32	105.56	111.36
2	B	216	CYS	CA-C-N	5.31	124.94	119.05
2	B	216	CYS	C-N-CA	5.31	124.94	119.05
1	A	553	TRP	N-CA-C	5.26	119.40	112.25
4	D	83	TYR	N-CA-C	5.15	119.56	113.12
1	A	455	ASP	CA-C-N	5.12	124.78	119.56
1	A	455	ASP	C-N-CA	5.12	124.78	119.56
1	A	160	ILE	N-CA-C	5.10	115.25	108.11
3	C	27	SER	N-CA-C	5.07	117.70	111.82

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4522	0	4426	457	0
2	B	1869	0	1850	136	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	1008	0	1066	42	0
4	D	898	0	936	44	0
5	A	9	0	2	3	0
6	A	1	0	0	0	0
6	B	1	0	0	0	0
7	A	53	0	29	6	0
8	B	4	0	0	2	0
9	B	8	0	0	1	0
10	B	7	0	0	3	0
11	C	43	0	30	5	0
12	C	77	0	112	5	0
13	C	35	0	40	1	0
14	D	23	0	26	3	0
15	A	61	0	0	4	0
15	B	50	0	0	0	0
15	C	15	0	0	1	0
15	D	14	0	0	0	0
All	All	8698	0	8517	646	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 38.

All (646) 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)
1:A:559:TYR:HB2	1:A:569:ARG:NH2	1.66	1.09
3:C:6:LYS:HG2	3:C:7:LYS:H	1.19	1.06
1:A:584:LYS:HG2	1:A:585:ILE:H	1.19	1.05
1:A:534:THR:HB	1:A:553:TRP:HE1	1.21	1.04
2:B:95:PRO:HD2	2:B:157:THR:HG22	1.38	1.04
2:B:92:ARG:HH11	3:C:17:THR:HG21	1.13	1.02
1:A:79:ILE:HD13	1:A:397:ALA:HB2	1.40	0.99
1:A:534:THR:HB	1:A:553:TRP:NE1	1.78	0.97
2:B:94:LEU:HB3	2:B:157:THR:HG21	1.44	0.97
1:A:172:ASN:HD21	1:A:430:ARG:HG3	1.33	0.93
2:B:223:THR:HG22	10:B:304:F3S:S1	2.09	0.93
1:A:236:MET:HE3	1:A:236:MET:HA	1.50	0.93
4:D:6:SER:HB3	4:D:94:GLN:NE2	1.85	0.92
1:A:578:ARG:HB3	1:A:579:PRO:HD3	1.52	0.92
1:A:139:ASP:OD2	1:A:330:ILE:HG12	1.72	0.90
1:A:238:MET:HE3	1:A:397:ALA:HB1	1.55	0.89

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:9:ASP:H	1:A:32:THR:HG23	1.38	0.89
2:B:140:ARG:O	2:B:140:ARG:HD3	1.73	0.89
1:A:242:HIS:O	1:A:351:PRO:HA	1.72	0.89
1:A:555:CYS:HB2	1:A:571:VAL:HG13	1.52	0.89
1:A:516:LEU:HA	1:A:519:LEU:HD12	1.56	0.88
2:B:77:THR:HG22	2:B:82:LEU:HD11	1.55	0.88
2:B:92:ARG:NH1	3:C:17:THR:HG21	1.89	0.88
4:D:9:GLY:HA2	4:D:14:HIS:HD2	1.37	0.88
2:B:25:LEU:HB2	2:B:42:LEU:HD21	1.56	0.87
1:A:126:PHE:HE2	1:A:401:GLY:HA3	1.38	0.86
1:A:9:ASP:OD2	1:A:32:THR:HG22	1.76	0.86
1:A:126:PHE:O	1:A:283:LEU:HD11	1.76	0.86
4:D:6:SER:HB3	4:D:94:GLN:HE22	1.40	0.84
2:B:95:PRO:HD2	2:B:157:THR:CG2	2.08	0.84
1:A:233:VAL:HG11	1:A:236:MET:SD	2.18	0.83
3:C:89:ILE:HG22	3:C:93:MET:HE3	1.58	0.83
3:C:51:SER:HB3	4:D:48:TRP:HE1	1.43	0.83
1:A:36:LEU:HD22	1:A:161:PHE:HB2	1.61	0.82
1:A:324:GLU:O	1:A:328:PRO:HG3	1.80	0.82
1:A:205:ALA:HB2	1:A:220:GLY:N	1.94	0.82
1:A:559:TYR:HB2	1:A:569:ARG:HH21	1.42	0.81
2:B:34:MET:HE3	2:B:34:MET:HA	1.60	0.81
1:A:277:ALA:HB1	1:A:588:TYR:O	1.80	0.81
1:A:584:LYS:HG2	1:A:585:ILE:N	1.96	0.81
1:A:520:MET:HE2	1:A:520:MET:HA	1.63	0.80
1:A:104:LEU:HD12	1:A:105:PRO:HD2	1.63	0.80
1:A:265:LEU:HD22	1:A:271:ARG:HG2	1.63	0.80
1:A:128:GLY:HA3	1:A:400:LEU:HD11	1.64	0.80
1:A:47:VAL:HG13	1:A:146:LEU:HD23	1.64	0.80
1:A:272:PHE:CZ	1:A:293:ILE:HG22	2.17	0.79
1:A:238:MET:HE3	1:A:397:ALA:CB	2.12	0.79
3:C:6:LYS:HE3	3:C:7:LYS:HG3	1.63	0.79
1:A:49:ALA:HB3	1:A:142:GLY:HA3	1.64	0.79
1:A:273:MET:HE1	1:A:292:SER:CB	2.13	0.79
1:A:240:GLN:N	1:A:357:MET:HE1	1.98	0.79
1:A:273:MET:HE3	1:A:289:VAL:HG13	1.65	0.79
1:A:567:THR:O	1:A:568:ARG:HG3	1.82	0.78
1:A:254:THR:HA	5:A:589:OAA:O5	1.83	0.77
1:A:47:VAL:HG13	1:A:146:LEU:CD2	2.15	0.77
1:A:313:LEU:HD12	1:A:349:VAL:HG11	1.66	0.77
1:A:534:THR:CB	1:A:553:TRP:HE1	1.98	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:242:HIS:CD2	1:A:252:LEU:HD11	2.19	0.77
1:A:578:ARG:CB	1:A:579:PRO:HD3	2.14	0.76
1:A:126:PHE:CE2	1:A:401:GLY:HA3	2.21	0.76
1:A:42:THR:O	1:A:47:VAL:HG11	1.84	0.76
1:A:69:MET:O	1:A:73:VAL:HG23	1.86	0.76
1:A:245:GLY:HA3	1:A:352:THR:HG22	1.67	0.75
1:A:534:THR:HG22	1:A:534:THR:O	1.87	0.75
3:C:30:HIS:HD2	3:C:87:VAL:HB	1.52	0.75
3:C:89:ILE:CG2	3:C:93:MET:HE3	2.17	0.74
1:A:272:PHE:CZ	1:A:293:ILE:CG2	2.70	0.74
2:B:77:THR:CG2	2:B:82:LEU:HD11	2.17	0.74
3:C:2:ILE:HG12	3:C:5:VAL:HB	1.70	0.74
1:A:127:GLY:HA3	1:A:255:GLU:OE2	1.87	0.74
2:B:58:GLY:HA2	8:B:302:FES:S2	2.28	0.74
1:A:120:GLY:HA2	2:B:132:GLU:OE2	1.87	0.74
3:C:60:GLN:O	3:C:64:ILE:HG13	1.88	0.73
1:A:572:ASN:C	1:A:573:MET:HG3	2.12	0.73
2:B:9:ARG:HH11	2:B:49:LEU:HD13	1.54	0.73
1:A:556:HIS:HB2	1:A:572:ASN:HB3	1.69	0.73
4:D:9:GLY:HA2	4:D:14:HIS:CD2	2.22	0.72
1:A:55:VAL:HG13	1:A:57:LEU:HG	1.71	0.72
1:A:109:LEU:C	1:A:117:ARG:O	2.32	0.72
3:C:30:HIS:CD2	3:C:87:VAL:HB	2.24	0.72
2:B:172:PRO:HG3	10:B:304:F3S:S3	2.29	0.72
1:A:453:GLY:HA3	1:A:496:ASN:O	1.89	0.71
1:A:395:HIS:ND1	1:A:399:ARG:HG3	2.04	0.71
2:B:34:MET:HA	2:B:34:MET:CE	2.20	0.71
1:A:77:ASP:OD2	1:A:586:ARG:NH1	2.20	0.71
1:A:172:ASN:ND2	1:A:430:ARG:HG3	2.06	0.71
1:A:240:GLN:H	1:A:357:MET:HE1	1.54	0.71
1:A:392:VAL:HG13	1:A:394:VAL:HG13	1.73	0.71
1:A:276:TYR:C	1:A:278:PRO:HD3	2.15	0.71
1:A:463:LEU:O	1:A:467:MET:HG2	1.91	0.71
1:A:559:TYR:O	1:A:561:PRO:HD3	1.91	0.71
1:A:126:PHE:HE2	1:A:401:GLY:CA	2.04	0.70
1:A:463:LEU:HD13	1:A:520:MET:HE1	1.71	0.70
2:B:146:LEU:CD1	2:B:183:ILE:HD11	2.21	0.70
1:A:147:HIS:O	1:A:151:GLN:HG3	1.92	0.69
1:A:482:LYS:O	1:A:486:GLN:HG3	1.92	0.69
1:A:255:GLU:CD	1:A:286:ARG:HH22	1.99	0.69
1:A:454:GLU:OE1	1:A:493:ARG:HG3	1.92	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:34:MET:HE1	2:B:78:PRO:HG3	1.74	0.69
1:A:293:ILE:HD11	1:A:351:PRO:HG3	1.75	0.69
3:C:6:LYS:CG	3:C:7:LYS:H	1.99	0.69
1:A:221:ASP:HA	1:A:518:ASN:ND2	2.07	0.68
1:A:446:ARG:NH2	1:A:497:ALA:O	2.24	0.68
1:A:555:CYS:CB	1:A:571:VAL:HG13	2.23	0.68
1:A:245:GLY:HA3	1:A:352:THR:CG2	2.23	0.68
1:A:123:SER:HA	1:A:134:THR:O	1.91	0.68
1:A:139:ASP:OD1	1:A:139:ASP:O	2.10	0.68
1:A:255:GLU:HB2	1:A:286:ARG:HH22	1.57	0.68
1:A:291:ARG:NH2	1:A:538:GLY:O	2.27	0.68
2:B:164:TRP:HZ2	14:D:306:UQ2:H72	1.57	0.68
1:A:205:ALA:HB2	1:A:220:GLY:CA	2.22	0.68
2:B:94:LEU:HB3	2:B:157:THR:CG2	2.21	0.68
2:B:1:MET:H1	2:B:29:GLU:HB2	1.58	0.68
3:C:6:LYS:HG2	3:C:7:LYS:N	2.01	0.68
3:C:14:ASP:O	3:C:17:THR:HB	1.94	0.68
1:A:534:THR:HG23	1:A:545:PHE:CE2	2.29	0.67
1:A:47:VAL:CG1	1:A:146:LEU:HD23	2.23	0.67
1:A:537:ARG:HH21	1:A:548:ARG:CG	2.07	0.67
1:A:140:ARG:HH11	2:B:148:GLU:HG2	1.59	0.67
1:A:280:ALA:HB3	1:A:588:TYR:O	1.94	0.67
1:A:578:ARG:HB3	1:A:579:PRO:CD	2.25	0.67
1:A:285:GLY:O	1:A:289:VAL:HG23	1.95	0.67
1:A:323:LEU:HB3	1:A:331:LEU:HD21	1.76	0.67
1:A:454:GLU:OE2	1:A:493:ARG:NE	2.28	0.67
1:A:60:THR:CG2	1:A:123:SER:HB2	2.24	0.67
1:A:270:GLU:HG2	1:A:271:ARG:N	2.10	0.67
1:A:264:LEU:C	1:A:265:LEU:HD23	2.19	0.66
1:A:534:THR:HB	1:A:553:TRP:CD1	2.30	0.66
1:A:60:THR:HG21	1:A:123:SER:O	1.95	0.66
1:A:126:PHE:O	1:A:129:GLU:HG2	1.95	0.66
1:A:287:ASP:H	1:A:398:ASN:ND2	1.93	0.66
1:A:340:VAL:O	1:A:342:PRO:HD3	1.95	0.66
1:A:237:GLU:HB3	1:A:529:SER:HB2	1.77	0.66
1:A:571:VAL:HG12	1:A:572:ASN:N	2.11	0.66
2:B:65:LEU:HD12	2:B:74:ALA:HB2	1.77	0.65
1:A:537:ARG:HH21	1:A:548:ARG:HB2	1.61	0.65
1:A:237:GLU:OE1	1:A:529:SER:HB3	1.96	0.65
1:A:369:LEU:CD2	1:A:379:VAL:HG22	2.26	0.65
1:A:537:ARG:HA	1:A:586:ARG:HH12	1.60	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:574:GLU:HG3	1:A:574:GLU:O	1.97	0.65
1:A:9:ASP:CG	1:A:32:THR:HG22	2.21	0.65
2:B:37:ASP:O	2:B:41:GLN:HG2	1.97	0.65
1:A:78:TYR:CD2	1:A:583:PRO:HA	2.32	0.65
2:B:1:MET:HG3	2:B:29:GLU:OE2	1.97	0.65
2:B:77:THR:HG22	2:B:82:LEU:CD1	2.24	0.65
3:C:2:ILE:CG1	3:C:5:VAL:HB	2.26	0.65
1:A:75:GLY:O	1:A:398:ASN:HB3	1.96	0.64
1:A:404:SER:O	1:A:407:ASP:HB3	1.97	0.64
1:A:251:VAL:HG13	2:B:57:GLU:OE1	1.96	0.64
1:A:314:LYS:HZ1	1:A:317:HIS:HB3	1.62	0.64
1:A:537:ARG:HH12	1:A:554:LEU:HD13	1.62	0.64
1:A:129:GLU:O	1:A:283:LEU:HD21	1.98	0.64
4:D:92:MET:O	4:D:96:VAL:HG23	1.97	0.64
1:A:577:LEU:N	1:A:577:LEU:HD23	2.12	0.64
1:A:226:ALA:HB2	1:A:360:ILE:HD11	1.80	0.64
1:A:242:HIS:CE1	1:A:286:ARG:HH21	2.16	0.64
2:B:56:ARG:O	2:B:56:ARG:HG2	1.96	0.64
1:A:571:VAL:CG1	1:A:572:ASN:N	2.61	0.64
1:A:255:GLU:N	5:A:589:OAA:O4	2.25	0.64
2:B:129:PRO:HG2	2:B:133:HIS:CD2	2.34	0.63
1:A:63:ASP:OD1	1:A:64:ASN:N	2.31	0.63
4:D:98:VAL:O	4:D:102:VAL:HG23	1.99	0.63
1:A:357:MET:HE3	1:A:357:MET:H	1.62	0.63
1:A:529:SER:OG	1:A:569:ARG:NH2	2.31	0.63
1:A:239:TRP:CZ3	1:A:356:MET:HB2	2.33	0.63
1:A:252:LEU:HG	1:A:253:VAL:N	2.12	0.63
1:A:35:LEU:C	1:A:36:LEU:HD23	2.23	0.63
1:A:161:PHE:HB3	1:A:164:TRP:CD1	2.33	0.63
1:A:483:GLY:O	1:A:487:LEU:HD13	1.99	0.63
1:A:119:PHE:CD2	2:B:134:LEU:HD12	2.33	0.63
1:A:208:ILE:HG13	1:A:209:TYR:CD1	2.34	0.63
1:A:54:THR:OG1	1:A:403:ASN:ND2	2.32	0.63
1:A:227:ILE:HG23	1:A:561:PRO:HB3	1.81	0.62
2:B:58:GLY:CA	8:B:302:FES:S2	2.87	0.62
1:A:23:ALA:HB2	1:A:35:LEU:HD13	1.80	0.62
1:A:119:PHE:HB2	2:B:135:GLN:H	1.64	0.62
11:C:305:HEM:HAC	4:D:23:ALA:HB1	1.81	0.62
1:A:544:ASP:HB2	1:A:545:PHE:CD1	2.35	0.62
2:B:25:LEU:HB2	2:B:42:LEU:CD2	2.27	0.62
1:A:255:GLU:HB2	1:A:286:ARG:NH2	2.14	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:39:THR:O	4:D:39:THR:HG22	1.98	0.62
1:A:584:LYS:CG	1:A:585:ILE:H	2.00	0.62
2:B:94:LEU:CB	2:B:157:THR:HG21	2.26	0.62
1:A:369:LEU:HD21	1:A:379:VAL:HG22	1.82	0.62
2:B:1:MET:N	2:B:29:GLU:HB2	2.14	0.62
1:A:234:GLN:HB2	1:A:361:PRO:HG3	1.80	0.62
2:B:59:VAL:HG12	2:B:59:VAL:O	1.99	0.62
1:A:81:ASP:CG	1:A:393:SER:HB2	2.25	0.61
1:A:76:SER:O	1:A:79:ILE:HG22	2.00	0.61
1:A:559:TYR:CB	1:A:569:ARG:NH2	2.55	0.61
4:D:6:SER:HB2	4:D:77:TRP:CD1	2.35	0.61
1:A:265:LEU:CD2	1:A:271:ARG:HG2	2.30	0.61
1:A:272:PHE:HZ	1:A:293:ILE:CG2	2.11	0.61
1:A:38:LYS:CE	1:A:217:ILE:HG23	2.30	0.61
1:A:369:LEU:HD23	1:A:379:VAL:HA	1.82	0.61
1:A:572:ASN:ND2	1:A:573:MET:H	1.99	0.61
1:A:173:GLN:HB3	1:A:430:ARG:HE	1.66	0.61
1:A:362:THR:HG22	1:A:368:ALA:HA	1.81	0.61
1:A:79:ILE:HG21	1:A:397:ALA:CB	2.31	0.61
1:A:79:ILE:HG23	1:A:396:GLY:HA3	1.83	0.61
1:A:72:THR:HA	1:A:400:LEU:HD22	1.81	0.61
1:A:213:THR:HA	1:A:250:GLY:O	2.01	0.60
1:A:124:LYS:HE2	1:A:260:GLU:OE2	2.02	0.60
1:A:531:ASN:O	1:A:542:ARG:NH1	2.34	0.60
1:A:145:LEU:C	1:A:145:LEU:HD23	2.27	0.60
2:B:164:TRP:CZ2	14:D:306:UQ2:H72	2.36	0.60
1:A:434:GLU:O	1:A:438:GLU:HB2	2.01	0.60
1:A:560:LEU:HB3	1:A:568:ARG:HB2	1.83	0.60
3:C:44:LEU:HD21	3:C:74:MET:HE1	1.84	0.60
2:B:35:LEU:HD11	2:B:91:ILE:HD11	1.84	0.59
1:A:213:THR:HG23	1:A:353:CYS:O	2.02	0.59
1:A:511:VAL:O	1:A:515:GLU:HG3	2.03	0.59
1:A:293:ILE:HD11	1:A:351:PRO:CG	2.32	0.59
1:A:567:THR:HG23	15:A:606:HOH:O	2.01	0.59
4:D:94:GLN:O	4:D:98:VAL:HG23	2.02	0.59
1:A:537:ARG:HH12	1:A:554:LEU:CD1	2.15	0.59
1:A:295:ILE:HA	1:A:298:ARG:HG3	1.84	0.59
2:B:211:ASN:HD21	3:C:21:PRO:HD2	1.67	0.59
1:A:208:ILE:HG13	1:A:209:TYR:CE1	2.37	0.59
1:A:263:TYR:HB2	1:A:314:LYS:HB3	1.85	0.59
1:A:356:MET:HE3	1:A:358:GLY:O	2.03	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:405:LEU:HG	7:A:601:FAD:C2	2.33	0.58
2:B:146:LEU:HD12	2:B:183:ILE:HD11	1.84	0.58
4:D:69:LEU:O	4:D:73:TRP:HB2	2.02	0.58
2:B:209:ILE:HG23	3:C:24:ALA:HA	1.85	0.58
2:B:9:ARG:HH11	2:B:49:LEU:CD1	2.15	0.58
2:B:95:PRO:CD	2:B:157:THR:CG2	2.82	0.58
1:A:436:ASP:OD1	1:A:436:ASP:N	2.36	0.58
1:A:544:ASP:HB2	1:A:545:PHE:CE1	2.39	0.58
1:A:357:MET:H	1:A:357:MET:CE	2.16	0.58
2:B:67:MET:HE1	2:B:74:ALA:HA	1.86	0.58
1:A:221:ASP:HA	1:A:518:ASN:CG	2.29	0.58
2:B:33:MET:O	2:B:79:ILE:HG12	2.03	0.58
4:D:95:LEU:O	4:D:99:VAL:HG23	2.03	0.57
1:A:331:LEU:O	1:A:335:ARG:HG3	2.04	0.57
3:C:100:GLU:HB2	3:C:105:ALA:CB	2.33	0.57
2:B:25:LEU:HD13	2:B:42:LEU:HD23	1.86	0.57
1:A:81:ASP:HB3	1:A:83:ASP:OD1	2.04	0.57
1:A:140:ARG:NH1	2:B:148:GLU:HG2	2.18	0.57
1:A:579:PRO:O	1:A:580:ALA:HB3	2.04	0.57
2:B:2:ARG:HE	2:B:24:THR:HG21	1.68	0.57
1:A:177:VAL:HG21	1:A:383:LEU:HB2	1.86	0.57
1:A:236:MET:HA	1:A:236:MET:CE	2.31	0.57
1:A:104:LEU:HD12	1:A:105:PRO:CD	2.34	0.57
1:A:463:LEU:HD13	1:A:520:MET:CE	2.35	0.57
3:C:108:ARG:O	3:C:112:ILE:HG13	2.05	0.57
1:A:537:ARG:HH21	1:A:548:ARG:CB	2.18	0.56
1:A:534:THR:HG23	1:A:545:PHE:CD2	2.40	0.56
1:A:273:MET:HE1	1:A:292:SER:HB2	1.86	0.56
1:A:314:LYS:NZ	1:A:317:HIS:HB3	2.19	0.56
4:D:52:PHE:C	4:D:54:SER:H	2.14	0.56
1:A:60:THR:HG23	1:A:123:SER:HB2	1.86	0.56
1:A:459:ILE:HD13	1:A:494:LEU:HA	1.88	0.56
3:C:83:TYR:CZ	3:C:87:VAL:HG21	2.40	0.56
1:A:532:PHE:CD2	1:A:569:ARG:HB3	2.41	0.56
1:A:365:THR:OG1	1:A:367:GLN:HG3	2.06	0.56
1:A:139:ASP:OD2	1:A:330:ILE:CG1	2.52	0.56
1:A:255:GLU:CB	1:A:286:ARG:HH22	2.19	0.56
2:B:210:MET:HE1	3:C:102:THR:C	2.31	0.56
1:A:63:ASP:CG	1:A:133:ARG:HH12	2.13	0.55
1:A:355:TYR:CD2	1:A:399:ARG:HD3	2.41	0.55
2:B:114:TYR:O	2:B:117:ILE:HG13	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:205:ARG:HH22	4:D:81:THR:HB	1.70	0.55
2:B:95:PRO:CD	2:B:157:THR:HG22	2.24	0.55
1:A:327:LEU:N	1:A:328:PRO:HD3	2.22	0.55
1:A:361:PRO:HA	1:A:391:CYS:O	2.06	0.55
1:A:532:PHE:CE2	1:A:569:ARG:HB3	2.42	0.55
2:B:179:TYR:O	2:B:183:ILE:HG12	2.07	0.55
1:A:44:SER:O	1:A:47:VAL:HG22	2.06	0.55
1:A:486:GLN:O	1:A:489:VAL:HB	2.07	0.55
3:C:123:LEU:HB3	12:C:308:CDN:H521	1.89	0.55
1:A:38:LYS:HG3	7:A:601:FAD:C4A	2.37	0.54
1:A:45:HIS:CE1	1:A:214:ASN:HA	2.41	0.54
3:C:82:ALA:O	3:C:86:VAL:HG23	2.07	0.54
1:A:293:ILE:CD1	1:A:351:PRO:HD3	2.37	0.54
2:B:9:ARG:HH12	2:B:49:LEU:CA	2.20	0.54
1:A:255:GLU:HG3	1:A:258:ARG:NH1	2.22	0.54
3:C:83:TYR:CE2	3:C:87:VAL:HG21	2.43	0.54
1:A:14:GLY:HA3	1:A:201:ALA:O	2.08	0.54
1:A:242:HIS:HE1	1:A:286:ARG:HH21	1.54	0.54
1:A:263:TYR:CZ	1:A:271:ARG:CZ	2.90	0.54
1:A:287:ASP:HB2	1:A:538:GLY:O	2.08	0.54
1:A:164:TRP:CH2	1:A:184:CYS:HB2	2.42	0.54
1:A:54:THR:O	1:A:406:LEU:HD22	2.07	0.54
1:A:266:ASN:HD21	1:A:270:GLU:HB3	1.72	0.54
1:A:399:ARG:HH11	1:A:399:ARG:HG2	1.73	0.54
1:A:222:GLY:HA3	1:A:389:ILE:CD1	2.38	0.54
1:A:255:GLU:CG	1:A:286:ARG:HH22	2.21	0.54
3:C:51:SER:HB3	4:D:48:TRP:NE1	2.20	0.54
1:A:535:GLU:OE1	1:A:549:ASP:N	2.41	0.53
1:A:556:HIS:N	1:A:572:ASN:O	2.35	0.53
1:A:263:TYR:HB3	1:A:265:LEU:HD21	1.90	0.53
1:A:476:GLU:HG3	1:A:478:ASP:CG	2.33	0.53
1:A:571:VAL:CG1	1:A:572:ASN:H	2.21	0.53
1:A:586:ARG:O	1:A:586:ARG:HG3	2.09	0.53
2:B:210:MET:HE1	3:C:102:THR:O	2.07	0.53
4:D:45:TYR:O	4:D:49:ILE:HG22	2.08	0.53
2:B:180:ARG:HD3	2:B:181:PHE:CE1	2.44	0.53
3:C:1:MET:HG3	3:C:6:LYS:HA	1.90	0.53
1:A:466:CYS:SG	1:A:470:ASN:ND2	2.81	0.53
2:B:52:ARG:NH2	2:B:105:VAL:O	2.42	0.53
1:A:91:THR:O	1:A:91:THR:HG22	2.09	0.53
1:A:501:ASP:OD2	1:A:507:ASN:ND2	2.41	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:273:MET:HE1	1:A:292:SER:OG	2.08	0.53
2:B:55:CYS:HB3	2:B:60:CYS:HB3	1.91	0.53
2:B:111:TYR:O	2:B:114:TYR:HB3	2.08	0.53
1:A:61:HIS:NE2	1:A:131:ALA:HB3	2.24	0.52
1:A:244:THR:HG23	1:A:258:ARG:HH21	1.73	0.52
2:B:160:PRO:HA	2:B:163:TRP:CE3	2.43	0.52
1:A:350:ILE:HG13	1:A:350:ILE:O	2.08	0.52
1:A:444:LEU:HG	1:A:448:ASN:ND2	2.25	0.52
1:A:467:MET:SD	1:A:523:ALA:HB1	2.48	0.52
11:C:305:HEM:HBB1	12:C:308:CDN:C24	2.39	0.52
1:A:572:ASN:C	1:A:573:MET:CG	2.81	0.52
1:A:76:SER:O	1:A:77:ASP:CB	2.57	0.52
1:A:82:GLN:HB2	1:A:577:LEU:HB3	1.91	0.52
1:A:109:LEU:O	1:A:117:ARG:O	2.28	0.52
1:A:255:GLU:OE1	1:A:286:ARG:NH2	2.42	0.52
1:A:578:ARG:CG	1:A:579:PRO:HD3	2.40	0.52
1:A:399:ARG:HG2	1:A:399:ARG:NH1	2.24	0.52
1:A:272:PHE:CZ	1:A:293:ILE:HG23	2.44	0.52
2:B:67:MET:SD	2:B:77:THR:HG21	2.49	0.52
1:A:38:LYS:HE2	1:A:217:ILE:HG23	1.92	0.52
4:D:115:VAL:OXT	4:D:115:VAL:HG12	2.10	0.52
1:A:65:TRP:O	1:A:68:HIS:HB3	2.10	0.52
1:A:232:PRO:HB2	1:A:558:LEU:HD11	1.92	0.52
1:A:477:GLY:HA2	1:A:542:ARG:NH2	2.25	0.51
1:A:303:CYS:O	1:A:309:PRO:HA	2.11	0.51
1:A:76:SER:O	1:A:79:ILE:CG2	2.58	0.51
1:A:534:THR:CB	1:A:553:TRP:NE1	2.62	0.51
1:A:548:ARG:NH2	1:A:584:LYS:O	2.44	0.51
1:A:456:PRO:HG3	1:A:497:ALA:HB1	1.92	0.51
2:B:84:GLN:OE1	2:B:87:LYS:HD2	2.09	0.51
1:A:71:ASP:HB3	1:A:128:GLY:O	2.11	0.51
4:D:6:SER:HB2	4:D:77:TRP:NE1	2.26	0.51
1:A:81:ASP:N	1:A:81:ASP:OD1	2.44	0.51
1:A:97:LEU:HD22	2:B:132:GLU:HB3	1.93	0.51
1:A:72:THR:HG22	1:A:400:LEU:HD23	1.92	0.50
1:A:126:PHE:CE2	1:A:401:GLY:C	2.88	0.50
1:A:392:VAL:CG1	1:A:394:VAL:HG13	2.39	0.50
1:A:63:ASP:OD2	1:A:133:ARG:NH1	2.44	0.50
2:B:76:ILE:O	2:B:78:PRO:HD3	2.11	0.50
2:B:234:LEU:HD23	4:D:13:VAL:HG13	1.94	0.50
1:A:176:ALA:HB1	1:A:381:PRO:HB2	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:242:HIS:HB2	1:A:354:HIS:HB2	1.91	0.50
1:A:9:ASP:H	1:A:32:THR:CG2	2.18	0.50
1:A:40:PHE:HD2	1:A:42:THR:HG1	1.60	0.50
1:A:484:LEU:HD11	1:A:531:ASN:HB2	1.93	0.50
1:A:555:CYS:HB3	1:A:573:MET:HG2	1.93	0.50
1:A:559:TYR:CE2	1:A:561:PRO:HG3	2.45	0.50
2:B:25:LEU:HD21	2:B:41:GLN:HG3	1.93	0.50
2:B:54:SER:O	2:B:56:ARG:N	2.45	0.50
1:A:306:PRO:O	2:B:31:ARG:NH2	2.43	0.50
1:A:370:THR:HG23	1:A:380:VAL:CG2	2.42	0.50
1:A:46:THR:HB	1:A:146:LEU:HD13	1.94	0.50
1:A:234:GLN:CB	1:A:361:PRO:HG3	2.41	0.50
1:A:355:TYR:CE2	1:A:399:ARG:HD3	2.46	0.50
1:A:483:GLY:HA2	1:A:486:GLN:NE2	2.27	0.50
1:A:559:TYR:CB	1:A:569:ARG:HH21	2.18	0.50
1:A:209:TYR:HB3	1:A:468:GLN:OE1	2.12	0.50
1:A:389:ILE:HG12	1:A:389:ILE:O	2.11	0.50
2:B:201:PHE:CE2	4:D:81:THR:HG21	2.47	0.50
2:B:201:PHE:CZ	4:D:81:THR:HG21	2.46	0.50
1:A:79:ILE:CD1	1:A:397:ALA:HB2	2.28	0.50
1:A:108:ARG:HG3	2:B:135:GLN:O	2.11	0.50
1:A:372:ASN:HB3	1:A:376:GLU:H	1.76	0.50
1:A:483:GLY:HA2	1:A:486:GLN:HE21	1.77	0.50
1:A:513:CYS:O	1:A:516:LEU:N	2.44	0.49
2:B:35:LEU:HD21	2:B:91:ILE:CD1	2.42	0.49
2:B:70:LYS:HE3	2:B:215:VAL:HG12	1.94	0.49
1:A:8:PHE:CE2	1:A:34:ALA:HB2	2.48	0.49
1:A:560:LEU:HD23	1:A:568:ARG:HD3	1.94	0.49
12:C:308:CDN:HA21	4:D:41:GLY:O	2.12	0.49
1:A:221:ASP:O	1:A:225:MET:HG3	2.13	0.49
1:A:119:PHE:CG	2:B:134:LEU:HA	2.47	0.49
1:A:314:LYS:HA	1:A:346:PRO:HG3	1.94	0.49
1:A:545:PHE:N	1:A:546:PRO:HD3	2.27	0.49
1:A:50:GLN:O	1:A:51:GLY:C	2.55	0.49
1:A:125:ASN:CG	1:A:132:ALA:HB2	2.37	0.49
1:A:255:GLU:CD	1:A:286:ARG:NH2	2.69	0.49
1:A:293:ILE:HD12	1:A:351:PRO:HD3	1.95	0.49
1:A:574:GLU:O	1:A:574:GLU:CG	2.57	0.49
1:A:214:ASN:HD22	1:A:214:ASN:N	2.10	0.49
1:A:584:LYS:HG2	1:A:585:ILE:HG12	1.95	0.49
1:A:244:THR:CG2	1:A:258:ARG:HH21	2.25	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:450:ASN:HB3	15:A:633:HOH:O	2.13	0.49
1:A:457:VAL:O	1:A:461:LYS:HG3	2.12	0.49
2:B:35:LEU:HD21	2:B:91:ILE:HD11	1.94	0.49
2:B:9:ARG:NH1	2:B:49:LEU:HA	2.28	0.48
2:B:218:LYS:HB2	2:B:220:LEU:HG	1.94	0.48
1:A:49:ALA:HA	7:A:601:FAD:C5X	2.44	0.48
1:A:176:ALA:CB	1:A:381:PRO:HB2	2.44	0.48
1:A:443:ARG:NH1	15:A:602:HOH:O	2.42	0.48
1:A:537:ARG:HH21	1:A:548:ARG:CD	2.26	0.48
1:A:583:PRO:O	1:A:584:LYS:HB2	2.13	0.48
1:A:567:THR:C	1:A:568:ARG:HG3	2.38	0.48
1:A:314:LYS:HG2	1:A:316:ASP:OD1	2.13	0.48
1:A:502:THR:C	1:A:510:ARG:NH2	2.72	0.48
1:A:525:ALA:O	1:A:529:SER:OG	2.25	0.48
2:B:118:LYS:HB2	2:B:193:ARG:HH12	1.79	0.48
1:A:205:ALA:CB	1:A:220:GLY:N	2.71	0.48
1:A:320:LYS:HA	1:A:323:LEU:HD12	1.95	0.48
1:A:384:PHE:N	1:A:384:PHE:CD1	2.82	0.48
1:A:408:LEU:HD11	7:A:601:FAD:H4'	1.96	0.48
2:B:220:LEU:HD12	9:B:303:SF4:S4	2.54	0.48
1:A:49:ALA:HB3	1:A:142:GLY:CA	2.40	0.48
1:A:284:ALA:HB3	1:A:289:VAL:HG22	1.95	0.48
1:A:476:GLU:HG3	1:A:478:ASP:OD2	2.14	0.48
2:B:34:MET:HE3	2:B:34:MET:CA	2.37	0.48
4:D:83:TYR:OH	14:D:306:UQ2:O1	2.27	0.48
1:A:126:PHE:CE2	1:A:401:GLY:CA	2.90	0.47
1:A:236:MET:HE3	1:A:358:GLY:HA3	1.95	0.47
1:A:287:ASP:H	1:A:398:ASN:HD21	1.61	0.47
2:B:82:LEU:HD12	2:B:82:LEU:N	2.28	0.47
2:B:211:ASN:HD21	3:C:24:ALA:HB2	1.79	0.47
1:A:209:TYR:O	1:A:464:GLN:NE2	2.38	0.47
1:A:537:ARG:O	1:A:540:HIS:N	2.47	0.47
1:A:238:MET:CE	1:A:397:ALA:HB1	2.37	0.47
1:A:21:ARG:O	1:A:21:ARG:NH1	2.47	0.47
1:A:533:ARG:HB3	1:A:540:HIS:CE1	2.50	0.47
2:B:9:ARG:HH12	2:B:49:LEU:HA	1.80	0.47
1:A:182:ALA:HB3	1:A:193:PHE:HE1	1.79	0.47
1:A:255:GLU:CD	1:A:286:ARG:HH12	2.22	0.47
1:A:272:PHE:CE1	1:A:273:MET:HE2	2.49	0.47
3:C:127:LEU:HB2	12:C:308:CDN:H512	1.97	0.47
1:A:326:ARG:C	1:A:328:PRO:HD3	2.39	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:490:ILE:O	1:A:493:ARG:HB3	2.15	0.47
2:B:211:ASN:ND2	3:C:21:PRO:HD2	2.30	0.47
3:C:2:ILE:HD11	3:C:5:VAL:HG21	1.96	0.47
11:C:305:HEM:CAC	4:D:23:ALA:HB1	2.44	0.47
4:D:52:PHE:C	4:D:54:SER:N	2.70	0.47
2:B:37:ASP:O	2:B:38:ALA:C	2.56	0.47
1:A:124:LYS:N	1:A:134:THR:O	2.44	0.47
1:A:240:GLN:CB	1:A:357:MET:HE1	2.44	0.47
1:A:441:LEU:O	1:A:445:ASN:ND2	2.48	0.47
11:C:305:HEM:CBB	4:D:68:ILE:HG12	2.45	0.47
1:A:63:ASP:OD2	1:A:133:ARG:NH2	2.48	0.47
1:A:76:SER:O	1:A:77:ASP:HB2	2.14	0.47
4:D:70:ILE:O	4:D:74:ILE:HG13	2.14	0.46
1:A:233:VAL:CG1	1:A:236:MET:SD	2.99	0.46
1:A:234:GLN:CG	1:A:361:PRO:HG3	2.46	0.46
1:A:40:PHE:O	1:A:43:ARG:HG2	2.15	0.46
1:A:172:ASN:O	1:A:175:GLY:N	2.42	0.46
1:A:372:ASN:O	1:A:375:GLY:N	2.49	0.46
2:B:54:SER:O	2:B:55:CYS:C	2.57	0.46
1:A:356:MET:O	1:A:356:MET:HG2	2.15	0.46
2:B:9:ARG:NH1	2:B:49:LEU:CD1	2.78	0.46
2:B:95:PRO:O	2:B:157:THR:HG23	2.15	0.46
2:B:213:VAL:HG23	2:B:223:THR:HG23	1.98	0.46
1:A:404:SER:HB3	7:A:601:FAD:N1	2.31	0.46
1:A:472:SER:OG	1:A:473:VAL:N	2.45	0.46
1:A:559:TYR:OH	1:A:567:THR:HG21	2.16	0.46
2:B:211:ASN:O	2:B:214:SER:HB3	2.16	0.46
3:C:52:LEU:HD13	4:D:115:VAL:HG21	1.96	0.46
12:C:308:CDN:H152	4:D:37:PHE:CE2	2.49	0.46
1:A:470:ASN:HD21	1:A:486:GLN:HE22	1.63	0.46
2:B:29:GLU:O	2:B:29:GLU:HG3	2.16	0.46
11:C:305:HEM:HBB1	4:D:68:ILE:HG12	1.98	0.46
4:D:84:VAL:O	4:D:90:ARG:HD3	2.16	0.46
1:A:267:LYS:C	1:A:269:GLY:H	2.23	0.46
2:B:6:SER:HB3	2:B:90:VAL:HA	1.98	0.46
2:B:113:GLN:OE1	2:B:113:GLN:HA	2.16	0.46
2:B:150:ILE:HD13	2:B:218:LYS:HD3	1.97	0.46
1:A:372:ASN:N	1:A:376:GLU:O	2.49	0.46
1:A:432:ALA:HB1	1:A:436:ASP:HB2	1.98	0.46
3:C:118:VAL:HG21	13:C:309:EPH:H221	1.98	0.46
1:A:559:TYR:HD1	1:A:569:ARG:HE	1.64	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:94:GLU:HG3	1:A:95:ALA:N	2.30	0.45
1:A:182:ALA:HB3	1:A:193:PHE:CE1	2.51	0.45
1:A:47:VAL:HG13	1:A:146:LEU:HD22	1.94	0.45
1:A:236:MET:HG3	1:A:236:MET:O	2.15	0.45
1:A:471:PHE:CE2	1:A:527:ALA:HA	2.51	0.45
1:A:544:ASP:C	1:A:546:PRO:HD3	2.41	0.45
2:B:228:HIS:O	2:B:232:MET:HG3	2.17	0.45
1:A:13:ILE:HB	1:A:200:LEU:HD23	1.99	0.45
1:A:56:ALA:O	1:A:93:PRO:HG3	2.16	0.45
2:B:140:ARG:HD3	2:B:140:ARG:C	2.40	0.45
4:D:48:TRP:CH2	4:D:52:PHE:HE1	2.34	0.45
1:A:126:PHE:HB2	1:A:134:THR:OG1	2.17	0.45
1:A:559:TYR:CD2	1:A:561:PRO:HG3	2.52	0.45
1:A:244:THR:HG23	1:A:258:ARG:NH2	2.31	0.45
1:A:277:ALA:O	1:A:279:ASN:N	2.49	0.45
2:B:34:MET:HE1	2:B:78:PRO:CG	2.44	0.45
1:A:475:ARG:O	1:A:542:ARG:HA	2.17	0.45
1:A:505:GLU:O	2:B:101:ARG:HD2	2.16	0.45
2:B:81:ALA:C	2:B:82:LEU:HD12	2.41	0.45
1:A:122:GLN:O	1:A:135:ALA:HA	2.16	0.45
1:A:151:GLN:O	1:A:154:LEU:HB2	2.17	0.45
1:A:334:SER:O	1:A:338:ALA:HB3	2.17	0.45
2:B:34:MET:CE	2:B:34:MET:CA	2.91	0.45
2:B:126:GLN:O	2:B:127:ASN:HB2	2.16	0.45
1:A:106:PHE:HA	1:A:137:ALA:HB2	1.99	0.45
4:D:8:LEU:HG	4:D:73:TRP:CD2	2.52	0.45
2:B:21:GLN:NE2	2:B:23:TYR:OH	2.50	0.45
3:C:92:MET:HE1	4:D:24:ILE:HG13	1.98	0.45
4:D:62:LEU:HD23	4:D:65:LEU:HD12	1.99	0.45
1:A:104:LEU:HA	1:A:105:PRO:HD3	1.76	0.44
1:A:119:PHE:CD2	2:B:134:LEU:CD1	3.00	0.44
1:A:222:GLY:HA3	1:A:389:ILE:HD11	1.98	0.44
1:A:298:ARG:NH1	1:A:543:PHE:CZ	2.85	0.44
1:A:327:LEU:O	1:A:331:LEU:HG	2.17	0.44
1:A:61:HIS:NE2	1:A:131:ALA:CB	2.80	0.44
1:A:133:ARG:HB2	1:A:133:ARG:HH11	1.83	0.44
1:A:88:MET:HB2	1:A:411:PHE:CZ	2.52	0.44
1:A:304:ASP:OD1	1:A:305:GLY:N	2.51	0.44
1:A:478:ASP:O	1:A:481:ALA:HB3	2.17	0.44
2:B:236:ARG:O	2:B:236:ARG:HG2	2.17	0.44
1:A:68:HIS:O	1:A:72:THR:HG23	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:476:GLU:HG3	1:A:478:ASP:OD1	2.16	0.44
1:A:578:ARG:CB	1:A:579:PRO:CD	2.85	0.44
2:B:55:CYS:O	2:B:56:ARG:HB3	2.16	0.44
2:B:180:ARG:HG2	2:B:180:ARG:HH11	1.83	0.44
3:C:28:ILE:O	3:C:28:ILE:HG12	2.17	0.44
2:B:79:ILE:HG22	2:B:79:ILE:O	2.16	0.44
2:B:160:PRO:HA	2:B:163:TRP:CD2	2.52	0.44
3:C:81:LEU:O	3:C:85:VAL:HG23	2.18	0.44
1:A:36:LEU:HD23	1:A:36:LEU:N	2.33	0.44
1:A:60:THR:HG21	1:A:123:SER:HB2	1.97	0.44
1:A:97:LEU:CD1	2:B:131:ARG:HB3	2.48	0.44
4:D:111:VAL:O	4:D:115:VAL:HG23	2.17	0.44
1:A:172:ASN:N	1:A:176:ALA:O	2.51	0.44
1:A:234:GLN:HG3	1:A:361:PRO:HG3	1.99	0.44
1:A:3:LEU:HA	1:A:4:PRO:HD3	1.88	0.44
1:A:82:GLN:NE2	1:A:581:PHE:HB3	2.32	0.44
2:B:127:ASN:N	2:B:128:PRO:HD3	2.32	0.44
1:A:13:ILE:HD12	1:A:200:LEU:CD2	2.48	0.43
1:A:537:ARG:NH2	1:A:548:ARG:HB2	2.29	0.43
1:A:571:VAL:HG13	1:A:572:ASN:H	1.83	0.43
2:B:143:LEU:HD12	2:B:143:LEU:N	2.33	0.43
1:A:167:LEU:O	1:A:225:MET:HG2	2.18	0.43
1:A:272:PHE:CD1	1:A:272:PHE:C	2.95	0.43
1:A:222:GLY:HA3	1:A:389:ILE:HD12	2.00	0.43
1:A:257:CYS:HB3	1:A:315:LEU:HD21	2.00	0.43
4:D:82:ASP:O	4:D:85:LYS:HE3	2.19	0.43
1:A:26:ILE:O	1:A:31:GLN:HB2	2.18	0.43
1:A:532:PHE:HD2	1:A:569:ARG:HD3	1.83	0.43
2:B:210:MET:CE	3:C:103:PHE:HA	2.48	0.43
1:A:392:VAL:CG1	1:A:394:VAL:HG22	2.49	0.43
1:A:433:SER:C	1:A:435:SER:N	2.75	0.43
1:A:577:LEU:HD11	1:A:580:ALA:C	2.44	0.43
2:B:223:THR:CG2	10:B:304:F3S:S1	2.95	0.43
2:B:236:ARG:O	2:B:237:ASN:ND2	2.51	0.43
3:C:45:LEU:HD23	3:C:45:LEU:HA	1.83	0.43
1:A:74:LYS:O	1:A:74:LYS:HG2	2.18	0.43
1:A:164:TRP:CZ3	1:A:184:CYS:HB2	2.54	0.43
1:A:370:THR:HG23	1:A:380:VAL:HG22	2.01	0.43
1:A:395:HIS:CG	1:A:399:ARG:HG3	2.52	0.43
1:A:453:GLY:CA	1:A:496:ASN:O	2.63	0.43
1:A:79:ILE:HG21	1:A:397:ALA:HB3	1.98	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:292:SER:O	1:A:296:GLU:HG2	2.18	0.43
1:A:549:ASP:CG	1:A:552:ASN:HB2	2.44	0.43
2:B:179:TYR:CE2	2:B:183:ILE:HD13	2.54	0.43
1:A:79:ILE:HG21	1:A:397:ALA:N	2.33	0.43
1:A:129:GLU:HG3	1:A:131:ALA:O	2.18	0.43
1:A:529:SER:OG	1:A:569:ARG:CZ	2.67	0.43
2:B:155:CYS:SG	2:B:156:SER:N	2.92	0.43
1:A:11:VAL:HG23	1:A:195:ALA:HB2	2.01	0.42
1:A:133:ARG:HH11	1:A:133:ARG:CB	2.32	0.42
1:A:491:ARG:NH1	1:A:524:TYR:CE2	2.87	0.42
1:A:234:GLN:HG3	1:A:361:PRO:CG	2.48	0.42
1:A:253:VAL:HG13	1:A:330:ILE:HD12	2.00	0.42
1:A:503:SER:O	1:A:510:ARG:NH2	2.52	0.42
1:A:572:ASN:HD22	1:A:573:MET:H	1.65	0.42
2:B:210:MET:HG2	2:B:223:THR:HG21	2.01	0.42
2:B:137:PRO:O	2:B:141:GLU:HG3	2.19	0.42
2:B:212:CYS:SG	2:B:223:THR:HG23	2.59	0.42
1:A:228:ARG:HG3	15:A:602:HOH:O	2.18	0.42
1:A:537:ARG:O	1:A:538:GLY:C	2.60	0.42
2:B:21:GLN:HG2	2:B:23:TYR:CE1	2.55	0.42
2:B:54:SER:C	2:B:56:ARG:N	2.77	0.42
1:A:24:LEU:HD11	1:A:156:ASN:ND2	2.35	0.42
2:B:29:GLU:HA	2:B:29:GLU:OE1	2.20	0.42
1:A:219:THR:HG21	1:A:515:GLU:OE1	2.19	0.42
1:A:231:VAL:HA	1:A:232:PRO:HD3	1.81	0.42
2:B:67:MET:HB3	2:B:82:LEU:HD21	2.02	0.42
1:A:72:THR:HG22	1:A:400:LEU:CD2	2.49	0.42
1:A:480:MET:HE2	1:A:542:ARG:NH1	2.35	0.42
1:A:64:ASN:OD1	1:A:64:ASN:C	2.62	0.42
1:A:560:LEU:HD23	1:A:568:ARG:HB2	2.01	0.42
2:B:201:PHE:O	2:B:202:SER:C	2.63	0.42
3:C:20:PHE:HA	3:C:21:PRO:HD3	1.86	0.42
1:A:338:ALA:O	1:A:340:VAL:HG23	2.19	0.42
2:B:50:SER:OG	2:B:101:ARG:HD3	2.20	0.42
1:A:392:VAL:O	1:A:393:SER:CB	2.67	0.41
1:A:572:ASN:ND2	1:A:573:MET:N	2.66	0.41
1:A:264:LEU:O	1:A:265:LEU:HD23	2.20	0.41
1:A:463:LEU:HD11	1:A:523:ALA:CB	2.51	0.41
1:A:479:ALA:O	1:A:480:MET:C	2.62	0.41
1:A:502:THR:HA	1:A:510:ARG:HH21	1.85	0.41
1:A:508:THR:HB	2:B:43:LYS:HZ1	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:509:GLN:N	2:B:43:LYS:HZ1	2.18	0.41
1:A:77:ASP:C	1:A:78:TYR:CD1	2.98	0.41
7:A:601:FAD:H9	7:A:601:FAD:H1'1	1.80	0.41
1:A:421:GLU:O	1:A:424:ALA:HB3	2.20	0.41
1:A:476:GLU:O	1:A:479:ALA:HB3	2.20	0.41
1:A:42:THR:O	1:A:47:VAL:CG1	2.63	0.41
1:A:477:GLY:CA	1:A:542:ARG:NH2	2.83	0.41
4:D:85:LYS:HA	4:D:85:LYS:HD3	1.74	0.41
1:A:360:ILE:HA	1:A:361:PRO:HD3	1.85	0.41
2:B:210:MET:HE2	3:C:103:PHE:HA	2.02	0.41
4:D:44:THR:HG22	4:D:46:GLU:H	1.84	0.41
4:D:93:LEU:O	4:D:94:GLN:C	2.63	0.41
1:A:79:ILE:HG23	1:A:396:GLY:CA	2.49	0.41
1:A:277:ALA:CB	1:A:588:TYR:O	2.62	0.41
1:A:422:SER:O	1:A:426:GLN:HG2	2.20	0.41
4:D:73:TRP:CD1	4:D:73:TRP:C	2.99	0.41
2:B:3:LEU:HB3	2:B:5:PHE:CE1	2.56	0.41
1:A:79:ILE:CG2	1:A:397:ALA:N	2.84	0.41
1:A:133:ARG:NH1	1:A:133:ARG:CB	2.84	0.41
1:A:143:HIS:HE1	2:B:147:TYR:O	2.03	0.41
1:A:151:GLN:HB3	2:B:119:PRO:O	2.21	0.41
1:A:254:THR:CA	5:A:589:OAA:O5	2.63	0.41
1:A:264:LEU:HD12	1:A:264:LEU:HA	1.88	0.41
1:A:274:GLU:O	1:A:278:PRO:HA	2.20	0.41
1:A:463:LEU:HD22	1:A:520:MET:CE	2.50	0.41
1:A:509:GLN:O	1:A:510:ARG:C	2.62	0.41
1:A:545:PHE:CD1	1:A:545:PHE:N	2.89	0.41
1:A:574:GLU:N	1:A:575:PRO:HD3	2.35	0.41
2:B:43:LYS:HE2	2:B:47:PRO:O	2.20	0.41
4:D:48:TRP:HA	4:D:48:TRP:CE3	2.55	0.41
4:D:65:LEU:HA	4:D:68:ILE:HD12	2.02	0.41
1:A:205:ALA:HA	1:A:218:ASN:O	2.21	0.41
1:A:334:SER:O	1:A:338:ALA:N	2.53	0.41
1:A:359:GLY:O	1:A:391:CYS:HB3	2.21	0.41
1:A:374:LYS:HB2	1:A:376:GLU:HG3	2.03	0.41
2:B:67:MET:HE1	2:B:74:ALA:CA	2.51	0.41
2:B:180:ARG:HG2	2:B:180:ARG:NH1	2.36	0.41
2:B:230:LYS:HA	2:B:233:LEU:HD12	2.01	0.41
1:A:78:TYR:O	1:A:554:LEU:HD11	2.21	0.40
1:A:86:GLU:OE1	1:A:578:ARG:HB3	2.21	0.40
1:A:307:TRP:CE3	1:A:307:TRP:HA	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:307:TRP:HA	1:A:307:TRP:HE3	1.86	0.40
1:A:153:ASN:HD22	1:A:158:THR:CB	2.34	0.40
1:A:236:MET:O	1:A:526:THR:HA	2.21	0.40
1:A:247:ALA:HB2	1:A:350:ILE:HG23	2.03	0.40
1:A:256:GLY:O	1:A:260:GLU:HG2	2.21	0.40
3:C:101:GLU:HG2	15:C:312:HOH:O	2.20	0.40
1:A:174:ASP:HB2	1:A:430:ARG:NH2	2.36	0.40
1:A:347:ILE:HA	1:A:348:PRO:HD3	1.97	0.40
1:A:220:GLY:O	1:A:221:ASP:C	2.64	0.40
1:A:272:PHE:HZ	1:A:293:ILE:HG22	1.74	0.40
2:B:28:ASP:C	2:B:30:GLY:H	2.28	0.40
1:A:64:ASN:OD1	1:A:66:GLU:HB2	2.21	0.40
1:A:280:ALA:HB3	1:A:588:TYR:C	2.46	0.40
1:A:470:ASN:HB2	1:A:471:PHE:CE1	2.57	0.40
1:A:579:PRO:O	1:A:580:ALA:CB	2.70	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	586/588 (100%)	504 (86%)	65 (11%)	17 (3%)	3	6
2	B	236/238 (99%)	205 (87%)	28 (12%)	3 (1%)	9	21
3	C	127/129 (98%)	117 (92%)	7 (6%)	3 (2%)	4	9
4	D	111/115 (96%)	107 (96%)	4 (4%)	0	100	100
All	All	1060/1070 (99%)	933 (88%)	104 (10%)	23 (2%)	5	10

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	51	GLY
1	A	110	ASP
1	A	119	PHE
1	A	278	PRO
3	C	7	LYS
1	A	77	ASP
1	A	431	ASP
1	A	538	GLY
1	A	393	SER
1	A	506	PHE
2	B	54	SER
2	B	55	CYS
3	C	2	ILE
1	A	505	GLU
2	B	199	ASP
1	A	116	GLN
1	A	578	ARG
1	A	118	PRO
1	A	453	GLY
3	C	97	GLY
1	A	105	PRO
1	A	309	PRO
1	A	348	PRO

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	473/473 (100%)	455 (96%)	18 (4%)	29	56
2	B	208/208 (100%)	204 (98%)	4 (2%)	50	75
3	C	109/109 (100%)	106 (97%)	3 (3%)	38	66
4	D	94/96 (98%)	91 (97%)	3 (3%)	34	62
All	All	884/886 (100%)	856 (97%)	28 (3%)	34	62

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	33	CYS
1	A	36	LEU
1	A	81	ASP
1	A	122	GLN
1	A	207	ARG
1	A	235	ASP
1	A	278	PRO
1	A	293	ILE
1	A	303	CYS
1	A	326	ARG
1	A	357	MET
1	A	371	VAL
1	A	436	ASP
1	A	446	ARG
1	A	544	ASP
1	A	568	ARG
1	A	569	ARG
1	A	573	MET
2	B	34	MET
2	B	53	ARG
2	B	63	ASP
2	B	85	PRO
3	C	28	ILE
3	C	36	ILE
3	C	129	TRP
4	D	6	SER
4	D	42	GLU
4	D	112	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (31) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	25	GLN
1	A	122	GLN
1	A	147	HIS
1	A	152	GLN
1	A	153	ASN
1	A	172	ASN
1	A	210	GLN
1	A	218	ASN
1	A	242	HIS
1	A	367	GLN
1	A	398	ASN

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Mol	Chain	Res	Type
1	A	403	ASN
1	A	420	GLN
1	A	445	ASN
1	A	448	ASN
1	A	470	ASN
1	A	486	GLN
1	A	540	HIS
1	A	572	ASN
2	B	109	GLN
2	B	123	ASN
2	B	133	HIS
2	B	135	GLN
2	B	211	ASN
2	B	237	ASN
3	C	4	ASN
3	C	30	HIS
3	C	91	HIS
4	D	14	HIS
4	D	78	GLN
4	D	94	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 11 ligands modelled in this entry, 2 are monoatomic - leaving 9 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
10	F3S	B	304	2	0,9,9	-	-	-		
7	FAD	A	601	1	58,58,58	2.15	16 (27%)	85,89,89	1.52	9 (10%)
8	FES	B	302	2	0,4,4	-	-	-		
11	HEM	C	305	3,4	50,50,50	1.71	10 (20%)	67,82,82	1.89	18 (26%)
14	UQ2	D	306	-	23,23,23	2.74	9 (39%)	30,31,31	1.39	3 (10%)
5	OAA	A	589	-	8,8,8	4.68	3 (37%)	8,10,10	2.32	4 (50%)
9	SF4	B	303	2	0,12,12	-	-	-		
12	CDN	C	308	-	76,76,76	2.05	11 (14%)	74,88,88	1.65	9 (12%)
13	EPH	C	309	-	34,34,48	1.13	3 (8%)	37,39,53	0.96	2 (5%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
10	F3S	B	304	2	-	-	0/3/3/3
7	FAD	A	601	1	-	3/34/50/50	0/6/6/6
11	HEM	C	305	3,4	-	8/14/54/54	-
8	FES	B	302	2	-	-	0/1/1/1
14	UQ2	D	306	-	-	4/15/39/39	0/1/1/1
5	OAA	A	589	-	-	2/8/8/8	-
9	SF4	B	303	2	-	-	0/6/5/5
12	CDN	C	308	-	3/3/9/9	24/87/87/87	-
13	EPH	C	309	-	-	6/38/38/52	-

All (52) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	589	OAA	C3-C4	-12.51	1.34	1.53
14	D	306	UQ2	C7-C8	-10.03	1.34	1.50
7	A	601	FAD	C4X-N5	7.54	1.47	1.30
12	C	308	CDN	OA8-CA7	7.41	1.53	1.40
12	C	308	CDN	OA6-CA5	-7.12	1.21	1.41
12	C	308	CDN	OB8-CB7	6.94	1.52	1.40
12	C	308	CDN	OA7-CA5	4.99	1.56	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
12	C	308	CDN	OB7-CB5	-4.98	1.21	1.39
12	C	308	CDN	OA9-CA7	-4.90	1.22	1.39
12	C	308	CDN	OB9-CB7	-4.80	1.22	1.39
11	C	305	HEM	CBB-CAB	4.79	1.53	1.30
7	A	601	FAD	C4A-N3A	4.69	1.43	1.34
7	A	601	FAD	C9A-N10	4.68	1.49	1.41
7	A	601	FAD	C2A-N3A	4.56	1.42	1.33
7	A	601	FAD	C5'-C4'	-4.26	1.46	1.51
7	A	601	FAD	C5A-C4A	-4.03	1.31	1.39
12	C	308	CDN	OB6-CB5	3.99	1.53	1.41
11	C	305	HEM	FE-ND	3.95	2.07	1.94
11	C	305	HEM	FE-NA	3.90	2.08	1.95
13	C	309	EPH	C25-C24	3.76	1.53	1.29
14	D	306	UQ2	O4-C4	3.50	1.31	1.23
11	C	305	HEM	C4B-NB	-3.40	1.32	1.38
11	C	305	HEM	C1A-NA	-3.39	1.33	1.39
7	A	601	FAD	C9A-C5X	3.39	1.46	1.41
7	A	601	FAD	C6-C7	3.38	1.44	1.39
13	C	309	EPH	C12-C13	3.04	1.51	1.29
11	C	305	HEM	FE-NC	3.00	2.05	1.95
7	A	601	FAD	C10-N1	2.95	1.39	1.33
14	D	306	UQ2	O3-CM3	-2.85	1.38	1.45
14	D	306	UQ2	C13-C14	2.83	1.40	1.32
11	C	305	HEM	C4A-NA	-2.77	1.34	1.39
11	C	305	HEM	C1C-NC	-2.77	1.34	1.39
14	D	306	UQ2	C8-C9	2.74	1.39	1.33
7	A	601	FAD	C8A-N9A	2.68	1.42	1.37
11	C	305	HEM	C4D-ND	-2.67	1.35	1.40
14	D	306	UQ2	C16-C14	2.66	1.57	1.50
7	A	601	FAD	C2B-C3B	-2.63	1.46	1.53
12	C	308	CDN	OB6-CB4	-2.59	1.41	1.44
14	D	306	UQ2	CM5-C5	2.54	1.56	1.50
7	A	601	FAD	C5X-N5	2.44	1.43	1.39
12	C	308	CDN	OA6-CA4	-2.42	1.41	1.44
7	A	601	FAD	O4-C4	2.41	1.28	1.23
7	A	601	FAD	C2'-C3'	-2.38	1.49	1.53
13	C	309	EPH	P1-O7	2.37	1.59	1.50
14	D	306	UQ2	C6-C5	2.34	1.39	1.35
5	A	589	OAA	O4-C4	2.24	1.28	1.22
11	C	305	HEM	C4C-NC	-2.24	1.35	1.39
7	A	601	FAD	C1'-N10	-2.22	1.42	1.47
7	A	601	FAD	C9-C9A	2.22	1.43	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	589	OAA	O2-C1	-2.21	1.23	1.30
14	D	306	UQ2	O3-C3	2.14	1.41	1.36
12	C	308	CDN	C11-CA5	2.02	1.54	1.50

All (45) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
12	C	308	CDN	OB9-CB7-C71	5.81	125.22	109.37
12	C	308	CDN	OB7-CB5-C51	5.67	124.84	109.37
12	C	308	CDN	OA9-CA7-C31	5.65	124.80	109.37
7	A	601	FAD	C4A-N9A-C8A	-5.30	100.17	105.74
7	A	601	FAD	C5A-C4A-N9A	5.24	111.52	105.81
11	C	305	HEM	CHC-C4B-NB	4.86	129.65	124.42
7	A	601	FAD	N3A-C4A-N9A	-4.77	119.06	127.17
5	A	589	OAA	O3-C3-C2	4.57	127.44	120.41
11	C	305	HEM	C4D-ND-C1D	4.47	110.50	105.21
14	D	306	UQ2	CM3-O3-C3	4.43	132.05	116.47
11	C	305	HEM	C1B-NB-C4B	4.34	110.34	105.21
12	C	308	CDN	C19-C18-C17	-3.83	95.02	114.37
7	A	601	FAD	C1'-N10-C9A	-3.82	113.22	120.63
11	C	305	HEM	CHB-C1B-NB	3.79	129.06	124.37
11	C	305	HEM	CHD-C1D-ND	3.63	128.32	124.42
12	C	308	CDN	C15-C14-C13	-3.61	96.10	114.37
11	C	305	HEM	CBB-CAB-C3B	-3.58	109.64	127.53
11	C	305	HEM	CHA-C4D-ND	3.44	128.62	124.37
14	D	306	UQ2	C12-C13-C14	3.12	138.03	127.64
11	C	305	HEM	C4A-NA-C1A	3.11	110.89	105.82
11	C	305	HEM	CHA-C1A-NA	2.93	129.17	123.86
7	A	601	FAD	C1'-C2'-C3'	2.89	117.51	109.66
11	C	305	HEM	CHC-C1C-NC	2.77	127.47	124.45
7	A	601	FAD	C5X-C9A-N10	-2.73	115.50	117.97
5	A	589	OAA	O3-C3-C4	-2.70	115.81	119.67
11	C	305	HEM	C3B-C4B-NB	-2.69	107.53	109.47
11	C	305	HEM	CHD-C4C-NC	2.61	127.30	124.45
11	C	305	HEM	C3C-C2C-C1C	-2.60	104.58	107.05
11	C	305	HEM	C4C-C3C-C2C	2.59	109.06	106.81
11	C	305	HEM	CHB-C4A-NA	2.55	128.49	123.86
5	A	589	OAA	O4-C4-C3	-2.55	118.64	121.81
11	C	305	HEM	C2A-C1A-NA	-2.53	107.35	110.15
7	A	601	FAD	N3A-C2A-N1A	-2.53	124.76	128.58
12	C	308	CDN	C21-C20-C19	-2.48	101.83	114.37
12	C	308	CDN	OA7-CA5-C11	2.43	116.00	109.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	A	601	FAD	O2B-C2B-C3B	2.42	119.58	111.82
12	C	308	CDN	C22-C21-C20	-2.38	102.31	114.37
11	C	305	HEM	C3D-C4D-ND	-2.34	107.61	110.17
7	A	601	FAD	C6-C5X-C9A	-2.25	115.96	119.05
13	C	309	EPH	C6-C5-C3	-2.16	105.77	113.69
11	C	305	HEM	C4C-NC-C1C	2.14	109.31	105.82
12	C	308	CDN	C16-C15-C14	-2.10	103.74	114.37
13	C	309	EPH	C14-C13-C12	-2.04	110.51	126.43
14	D	306	UQ2	C6-C5-C4	2.03	120.78	119.17
5	A	589	OAA	O5-C4-C3	2.02	119.18	113.59

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
12	C	308	CDN	CA7
12	C	308	CDN	CB7
12	C	308	CDN	CB5

All (47) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
11	C	305	HEM	C2B-C3B-CAB-CBB
11	C	305	HEM	C4B-C3B-CAB-CBB
11	C	305	HEM	C2C-C3C-CAC-CBC
11	C	305	HEM	C4C-C3C-CAC-CBC
12	C	308	CDN	CB2-OB2-PB2-OB5
12	C	308	CDN	CB2-OB2-PB2-OB3
12	C	308	CDN	CB2-OB2-PB2-OB4
12	C	308	CDN	OB7-CB5-OB6-CB4
12	C	308	CDN	C52-C51-CB5-OB7
12	C	308	CDN	OA7-CA5-OA6-CA4
13	C	309	EPH	C37-O5-P1-O8
11	C	305	HEM	C2A-CAA-CBA-CGA
12	C	308	CDN	C58-C59-C60-C61
12	C	308	CDN	C60-C61-C62-C63
12	C	308	CDN	C55-C56-C57-C58
13	C	309	EPH	C21-C22-C23-C24
13	C	309	EPH	C9-C10-C11-C12
13	C	309	EPH	C18-C19-C20-C21
12	C	308	CDN	CB3-CB4-CB6-OB8
13	C	309	EPH	C11-C10-C9-C8
12	C	308	CDN	C54-C55-C56-C57

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Mol	Chain	Res	Type	Atoms
12	C	308	CDN	C14-C15-C16-C17
12	C	308	CDN	C72-C71-CB7-OB8
12	C	308	CDN	C61-C62-C63-C64
14	D	306	UQ2	C12-C11-C9-C10
5	A	589	OAA	O1-C1-C2-C3
12	C	308	CDN	C31-CA7-OA8-CA6
12	C	308	CDN	OB6-CB4-CB6-OB8
14	D	306	UQ2	C12-C11-C9-C8
12	C	308	CDN	C1-CB2-OB2-PB2
12	C	308	CDN	C32-C31-CA7-OA8
12	C	308	CDN	CB7-C71-C72-C73
14	D	306	UQ2	C4-C3-O3-CM3
11	C	305	HEM	CAA-CBA-CGA-O1A
7	A	601	FAD	P-O3P-PA-O1A
5	A	589	OAA	O2-C1-C2-C3
11	C	305	HEM	CAA-CBA-CGA-O2A
12	C	308	CDN	CB2-C1-CA2-OA2
12	C	308	CDN	C35-C36-C37-C38
12	C	308	CDN	C56-C57-C58-C59
12	C	308	CDN	C52-C51-CB5-OB6
13	C	309	EPH	C10-C11-C12-C13
7	A	601	FAD	P-O3P-PA-O2A
7	A	601	FAD	O4B-C4B-C5B-O5B
11	C	305	HEM	C4D-C3D-CAD-CBD
14	D	306	UQ2	C6-C7-C8-C9
12	C	308	CDN	OA9-CA7-OA8-CA6

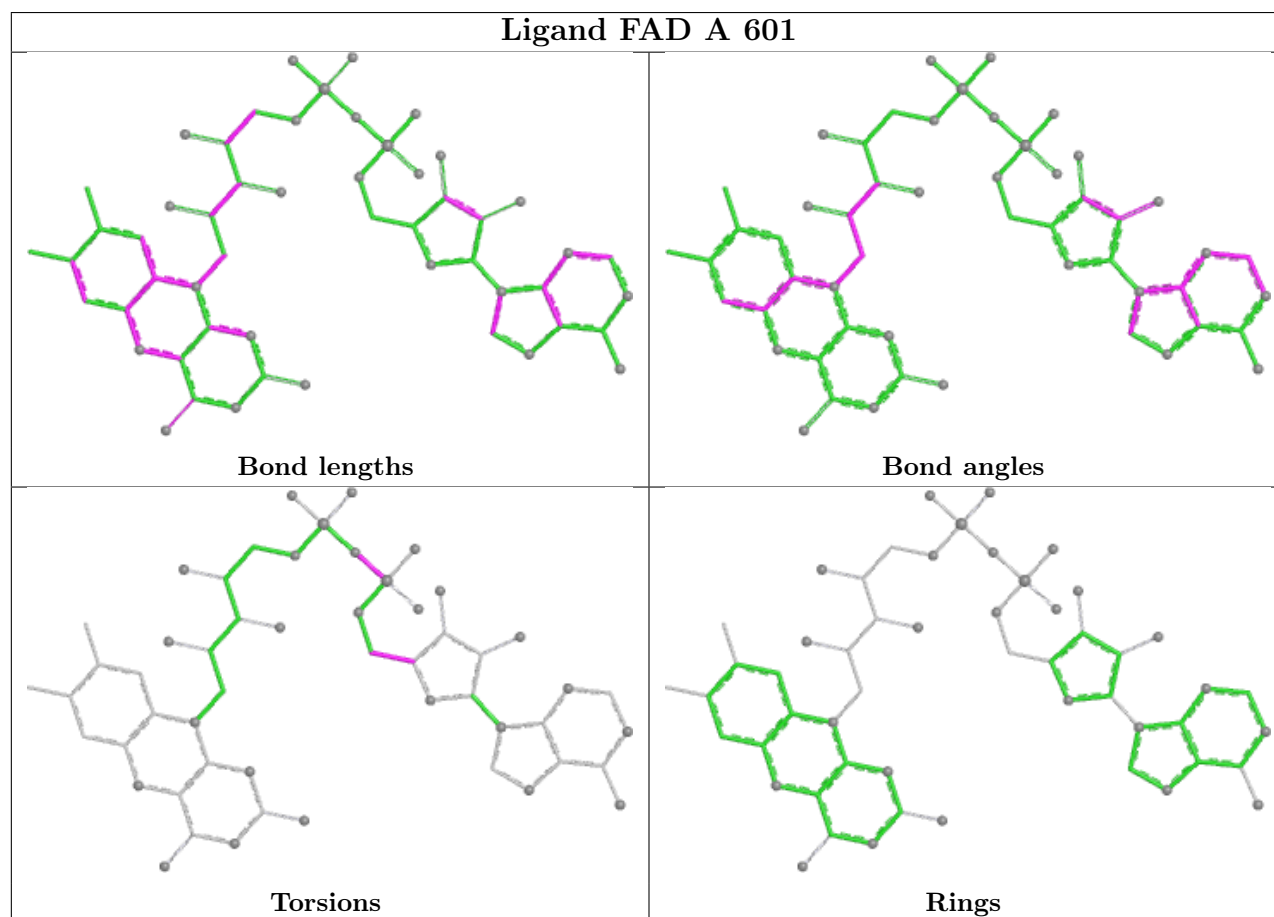
There are no ring outliers.

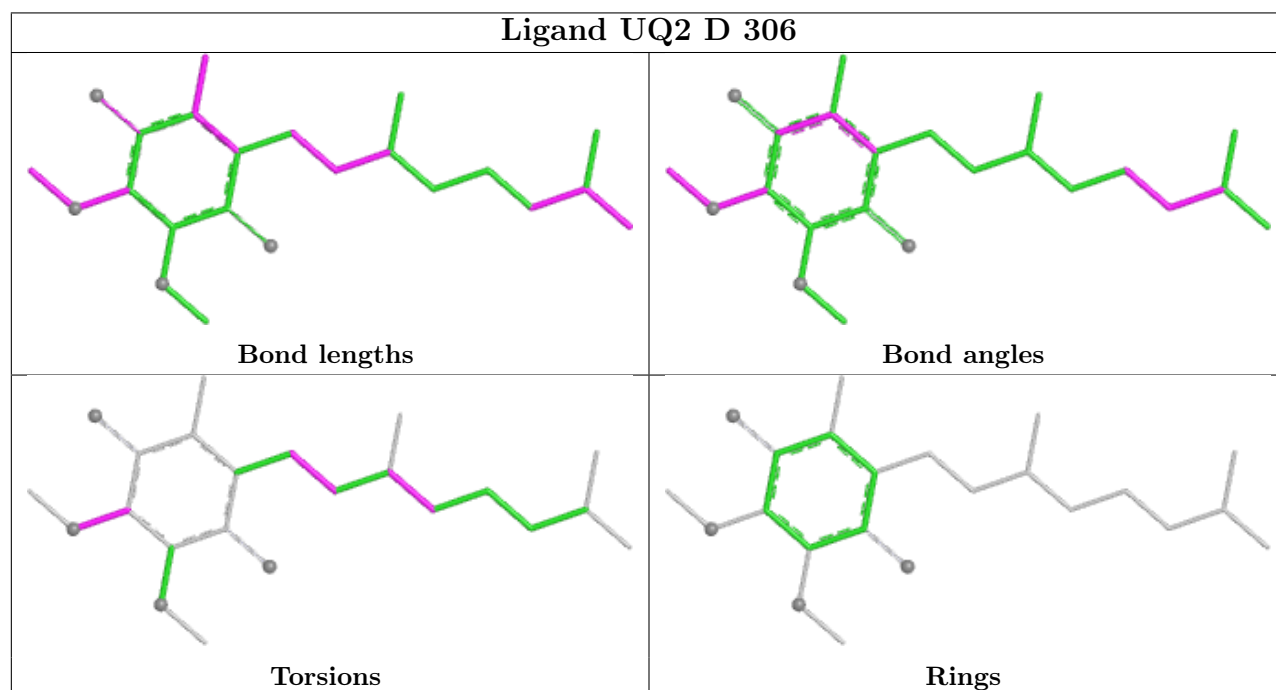
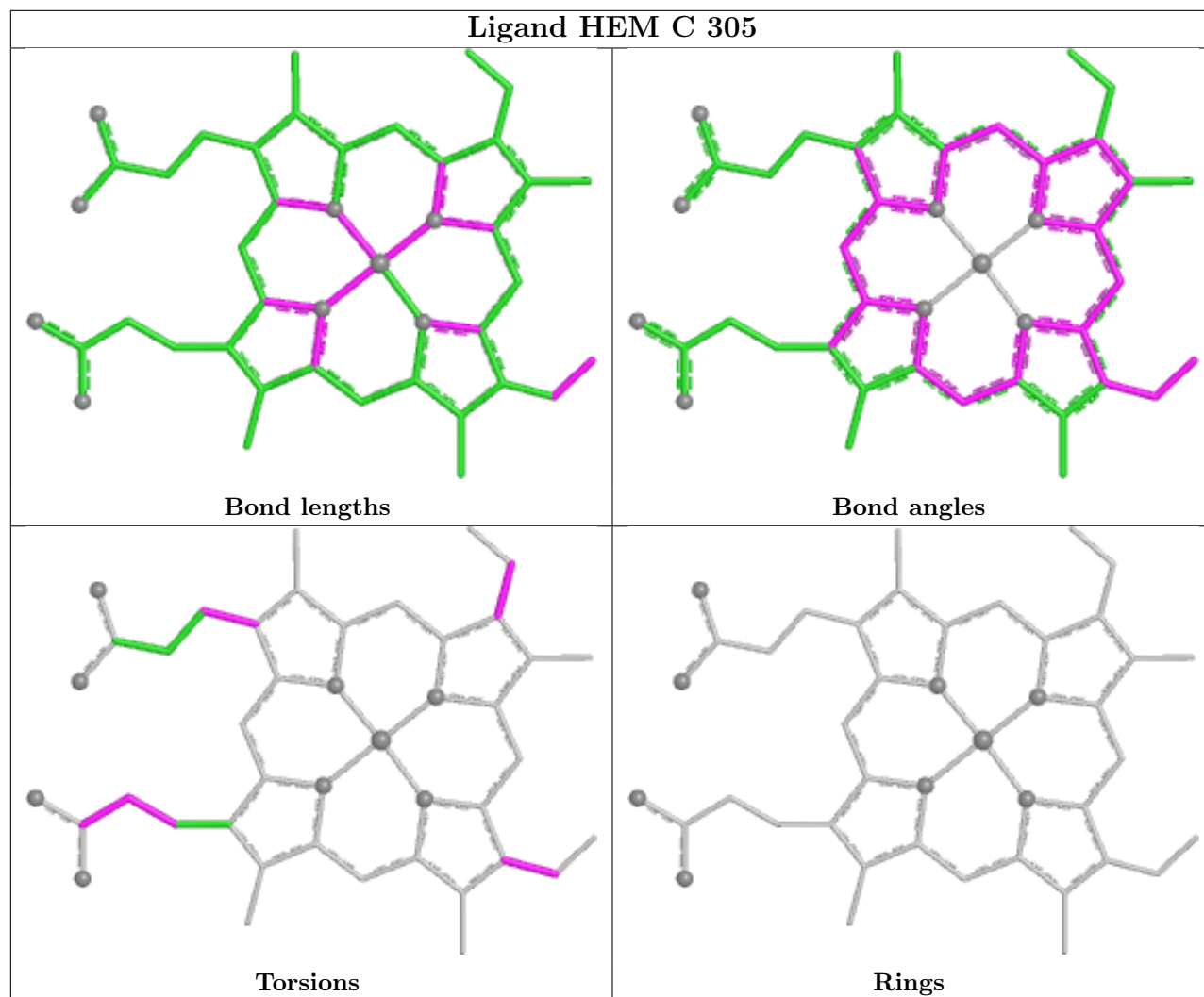
9 monomers are involved in 28 short contacts:

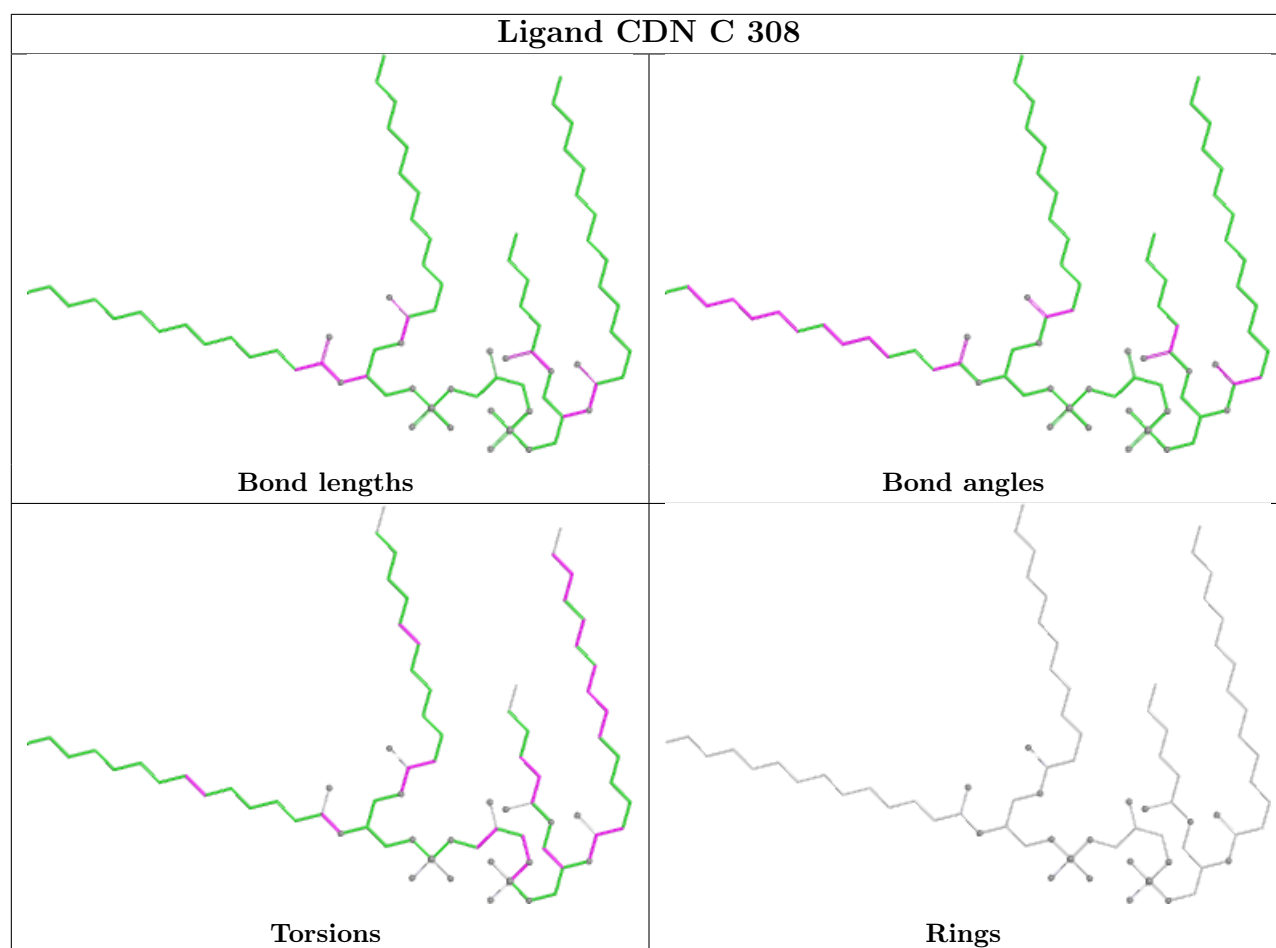
Mol	Chain	Res	Type	Clashes	Symm-Clashes
10	B	304	F3S	3	0
7	A	601	FAD	6	0
8	B	302	FES	2	0
11	C	305	HEM	5	0
14	D	306	UQ2	3	0
5	A	589	OAA	3	0
9	B	303	SF4	1	0
12	C	308	CDN	5	0
13	C	309	EPH	1	0

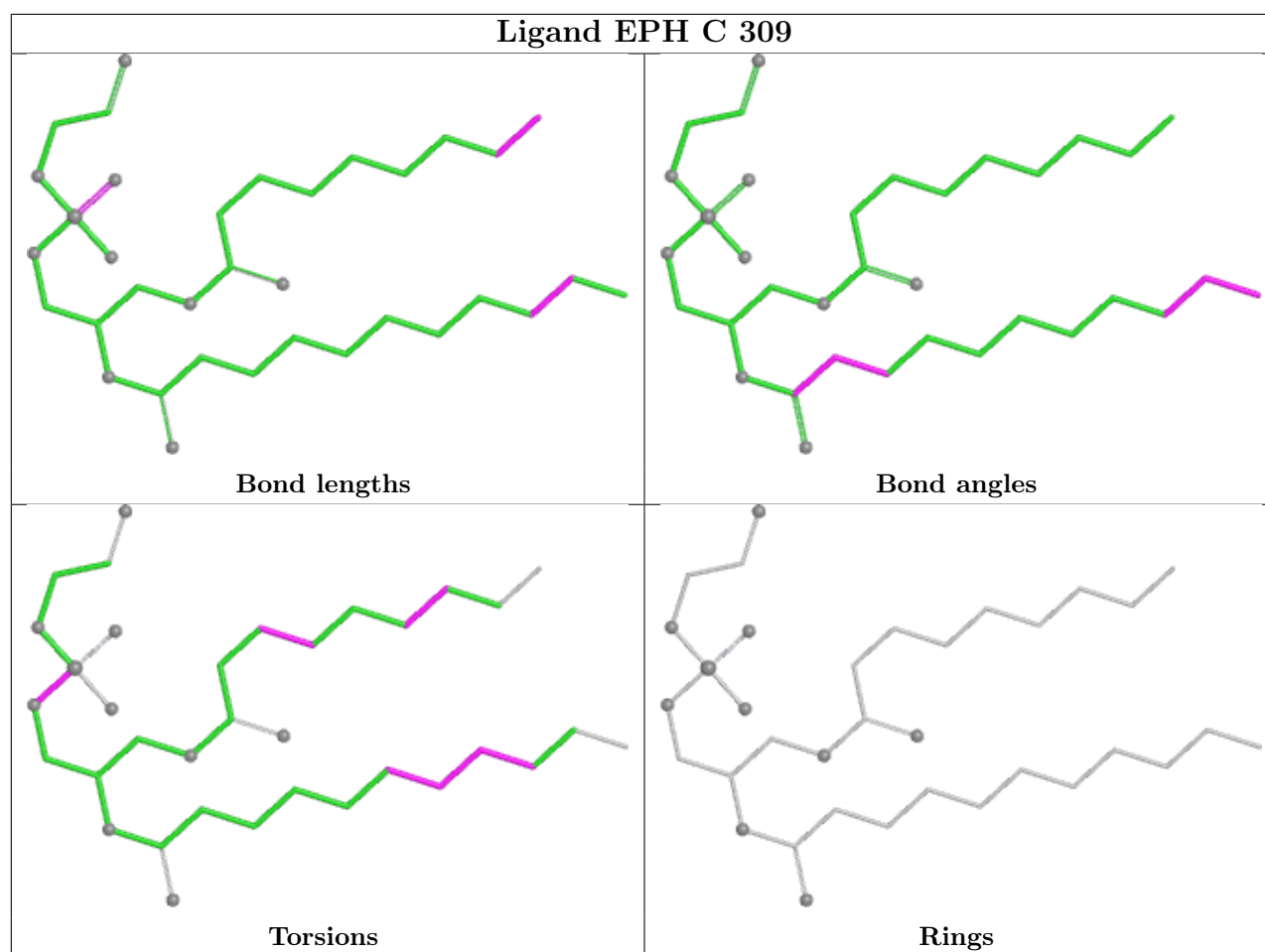
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths,

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









5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.