



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 08:19 PM UTC

PDB ID : 1KBO / pdb_00001kbo
Title : Complex of Human recombinant NAD(P)H:Quinone Oxide reductase type 1 with 5-methoxy-1,2-dimethyl-3-(phenoxyethyl)indole-4,7-dione (ES1340)
Authors : Faig, M.; Bianchet, M.A.; Amzel, L.M.
Deposited on : 2001-11-06
Resolution : 2.30 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

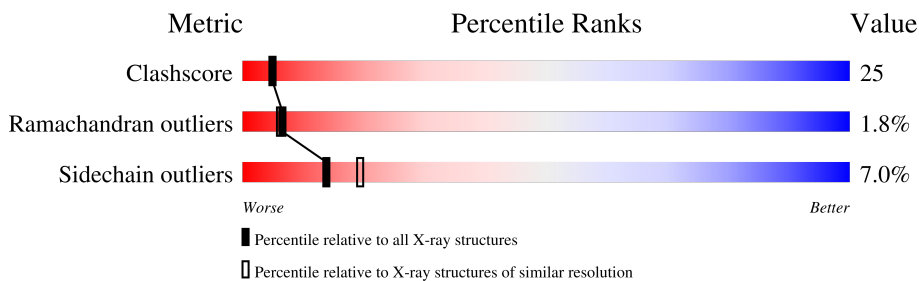
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)

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	273	55% 37% 7%
1	B	273	53% 39% 8%
1	C	273	51% 43% 7%
1	D	273	54% 41% 5%

2 Entry composition [i](#)

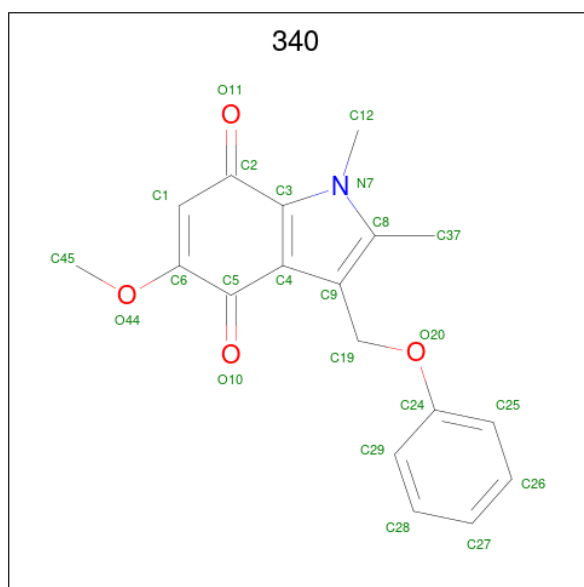
There are 4 unique types of molecules in this entry. The entry contains 9169 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 NAD(P)H dehydrogenase [quinone] 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	273	Total	C	N	O	S	0	0	0
			2175	1414	365	389	7			
1	B	273	Total	C	N	O	S	0	0	0
			2175	1414	365	389	7			
1	C	273	Total	C	N	O	S	0	0	0
			2175	1414	365	389	7			
1	D	273	Total	C	N	O	S	0	0	0
			2175	1414	365	389	7			

- Molecule 2 is 5-METHOXY-1,2-DIMETHYL-3-(PHENOXYMETHYL)INDOLE-4,7-DIONE (CCD ID: 340) (formula: C₁₈H₁₇NO₄).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			23	18	1	4		
2	B	1	Total	C	N	O	0	0
			23	18	1	4		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	C	1	Total	C	N	O	0	0
			23	18	1	4		
2	D	1	Total	C	N	O	0	0
			23	18	1	4		

- Molecule 3 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	40	Total	O	0	0
			40	40		
4	B	40	Total	O	0	0
			40	40		
4	C	42	Total	O	0	0
			42	42		

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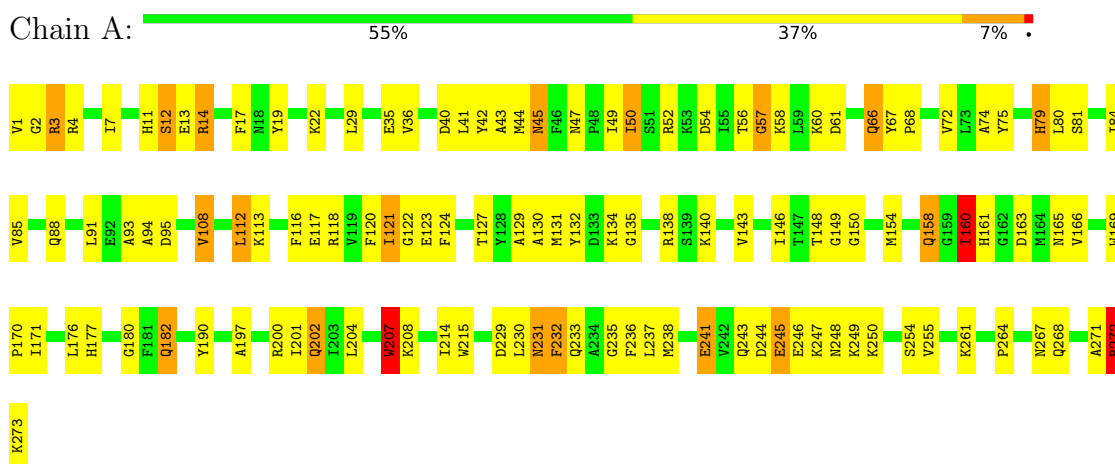
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	43	Total	O	0	0
			43	43		

3 Residue-property plots

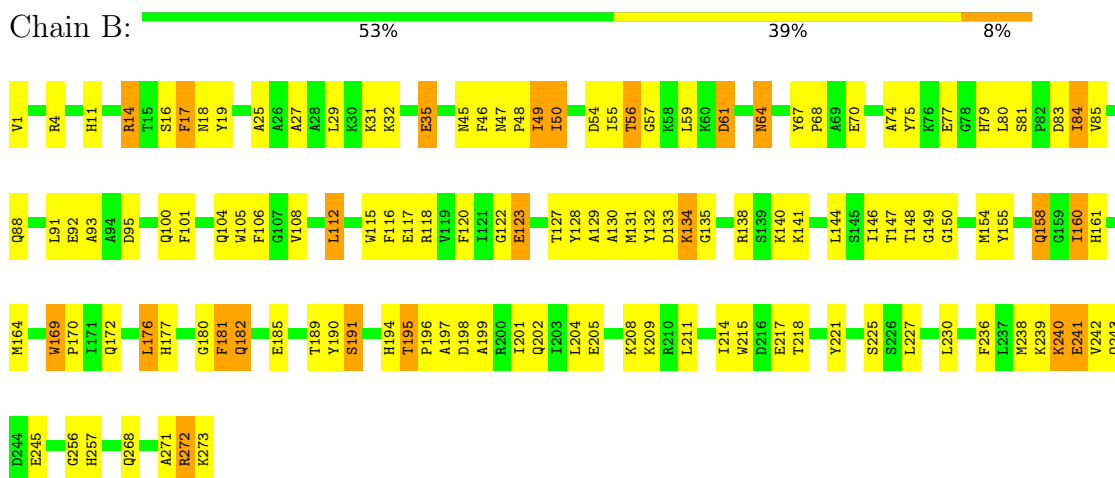
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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: NAD(P)H dehydrogenase [quinone] 1

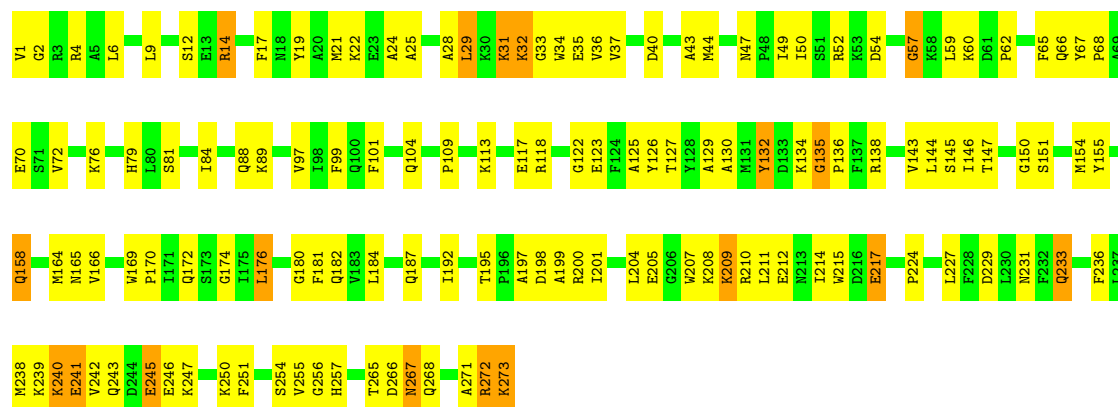


• Molecule 1: NAD(P)H dehydrogenase [quinone] 1



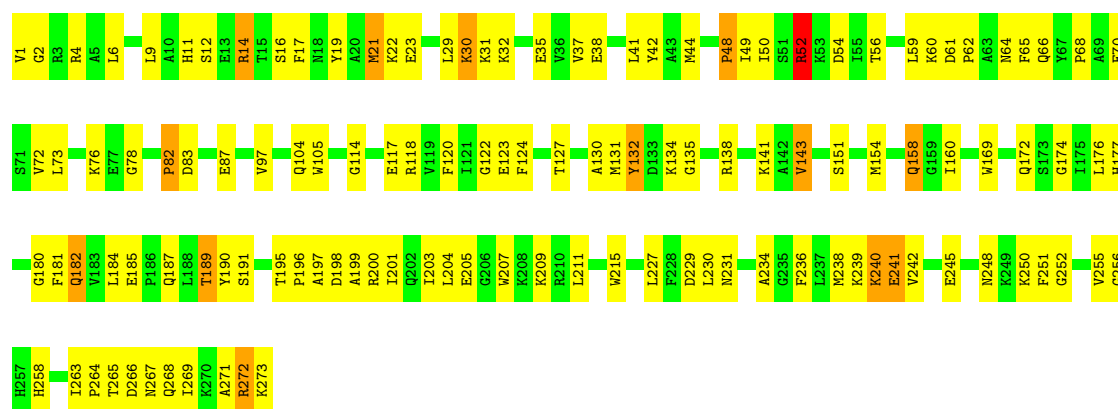
• Molecule 1: NAD(P)H dehydrogenase [quinone] 1





- Molecule 1: NAD(P)H dehydrogenase [quinone] 1

Chain D: 54% 41% 5%



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	55.64Å 57.03Å 96.88Å 76.80° 76.93° 86.28°	Depositor
Resolution (Å)	35.55 – 2.30	Depositor
% Data completeness (in resolution range)	74.4 (35.55-2.30)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.200 , 0.287	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	9169	wwPDB-VP
Average B, all atoms (Å ²)	32.0	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: FAD, 340

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	0.52	0/2233	1.07	15/3015 (0.5%)
1	B	0.51	0/2233	1.11	19/3015 (0.6%)
1	C	0.52	0/2233	1.00	7/3015 (0.2%)
1	D	0.53	0/2233	1.05	12/3015 (0.4%)
All	All	0.52	0/8932	1.06	53/12060 (0.4%)

There are no bond length outliers.

All (53) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	176	LEU	N-CA-C	8.52	120.26	110.97
1	C	101	PHE	N-CA-C	8.48	115.76	108.13
1	B	49	ILE	N-CA-C	8.12	119.84	107.99
1	D	135	GLY	CA-C-N	7.99	127.66	119.19
1	D	135	GLY	C-N-CA	7.99	127.66	119.19
1	B	61	ASP	CA-C-N	7.91	127.20	118.97
1	B	61	ASP	C-N-CA	7.91	127.20	118.97
1	C	135	GLY	CA-C-N	7.39	127.25	119.05
1	C	135	GLY	C-N-CA	7.39	127.25	119.05
1	A	79	HIS	N-CA-C	7.07	118.51	108.54
1	D	182	GLN	N-CA-C	-7.04	98.47	109.25
1	C	182	GLN	N-CA-C	-7.04	97.25	108.73
1	A	232	PHE	N-CA-C	6.45	118.31	111.28
1	B	181	PHE	N-CA-C	6.43	119.35	110.35
1	A	190	TYR	N-CA-C	6.38	118.96	110.53
1	B	19	TYR	N-CA-C	-6.31	104.48	111.36
1	A	49	ILE	N-CA-C	6.28	116.86	107.51
1	B	17	PHE	N-CA-C	-6.27	105.59	113.18
1	A	176	LEU	N-CA-C	6.24	117.77	110.97
1	A	12	SER	N-CA-C	6.22	119.04	111.82

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	195	THR	CA-C-N	6.19	126.16	119.78
1	B	195	THR	C-N-CA	6.19	126.16	119.78
1	D	61	ASP	CA-C-N	6.15	126.26	119.87
1	D	61	ASP	C-N-CA	6.15	126.26	119.87
1	D	174	GLY	CA-C-N	-6.11	116.98	122.97
1	D	174	GLY	C-N-CA	-6.11	116.98	122.97
1	A	182	GLN	N-CA-C	-6.07	97.81	108.20
1	A	36	VAL	N-CA-C	6.00	116.51	107.80
1	A	95	ASP	N-CA-C	-6.00	104.11	112.45
1	D	14	ARG	N-CA-C	-5.97	105.25	112.54
1	B	45	ASN	N-CA-C	-5.96	103.70	111.74
1	B	182	GLN	N-CA-C	-5.91	99.09	108.73
1	B	56	THR	N-CA-C	5.89	118.93	111.69
1	B	146	ILE	N-CA-C	5.88	117.06	108.53
1	D	38	GLU	N-CA-C	5.87	119.25	110.14
1	B	95	ASP	N-CA-C	-5.83	105.83	113.12
1	B	35	GLU	N-CA-C	-5.69	99.45	108.73
1	C	174	GLY	N-CA-C	5.63	119.49	112.73
1	B	169	TRP	CA-C-N	5.63	126.00	119.47
1	B	169	TRP	C-N-CA	5.63	126.00	119.47
1	D	190	TYR	N-CA-C	5.43	117.09	110.41
1	C	176	LEU	N-CA-C	5.43	117.06	111.03
1	A	171	ILE	N-CA-C	5.42	115.68	110.74
1	D	252	GLY	N-CA-C	-5.41	104.23	112.51
1	B	84	ILE	N-CA-C	-5.38	105.29	110.72
1	A	61	ASP	CA-C-N	5.36	125.02	119.56
1	A	61	ASP	C-N-CA	5.36	125.02	119.56
1	D	52	ARG	N-CA-C	-5.26	106.83	113.20
1	A	146	ILE	N-CA-C	5.24	115.45	108.11
1	A	207	TRP	N-CA-C	-5.15	105.35	110.97
1	A	160	ILE	CB-CA-C	-5.09	105.38	112.04
1	B	101	PHE	N-CA-C	5.07	112.69	108.13
1	C	12	SER	N-CA-C	5.05	119.44	113.28

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	2175	0	2179	111	0
1	B	2175	0	2179	135	0
1	C	2175	0	2179	121	0
1	D	2175	0	2179	109	0
2	A	23	0	17	0	0
2	B	23	0	17	0	0
2	C	23	0	17	0	0
2	D	23	0	17	1	0
3	A	53	0	31	5	0
3	B	53	0	31	5	0
3	C	53	0	31	4	0
3	D	53	0	31	4	0
4	A	40	0	0	3	0
4	B	40	0	0	3	0
4	C	42	0	0	6	0
4	D	43	0	0	5	0
All	All	9169	0	8908	452	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 25.

All (452) 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:14:ARG:HE	1:A:14:ARG:N	1.61	0.98
1:D:143:VAL:HG13	1:D:184:LEU:HB2	1.41	0.98
1:A:264:PRO:HG3	1:A:273:LYS:HE3	1.42	0.98
1:B:135:GLY:O	1:B:138:ARG:HG3	1.63	0.97
1:B:238:MET:HE2	1:D:158:GLN:HG3	1.46	0.97
1:A:14:ARG:H	1:A:14:ARG:NE	1.63	0.95
1:A:202:GLN:HA	1:A:202:GLN:HE21	1.31	0.94
1:B:154:MET:HE3	1:B:160:ILE:HD11	1.51	0.92
1:A:130:ALA:HB1	1:A:134:LYS:O	1.70	0.90
1:B:240:LYS:NZ	1:B:240:LYS:HB3	1.86	0.89
1:C:233:GLN:HA	1:C:233:GLN:HE21	1.36	0.89
1:D:17:PHE:HB2	3:D:604:FAD:H52A	1.53	0.89
1:A:66:GLN:HE21	1:A:66:GLN:HA	1.37	0.87
1:C:240:LYS:H	1:C:240:LYS:HD2	1.37	0.87
1:D:21:MET:HE3	1:D:204:LEU:HD23	1.55	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:68:PRO:O	1:A:72:VAL:HG23	1.75	0.86
1:A:127:THR:HG22	1:A:129:ALA:H	1.42	0.85
1:C:204:LEU:O	1:C:208:LYS:HG3	1.77	0.84
1:D:138:ARG:HA	1:D:180:GLY:O	1.77	0.83
1:C:209:LYS:HD3	1:C:210:ARG:N	1.94	0.82
1:A:163:ASP:OD2	1:A:165:ASN:HB2	1.80	0.81
1:A:17:PHE:HB2	3:A:601:FAD:H52A	1.62	0.81
1:C:238:MET:HE3	1:C:243:GLN:HG2	1.63	0.79
1:B:17:PHE:HB2	3:B:602:FAD:H52A	1.65	0.78
1:A:14:ARG:HE	1:A:14:ARG:H	0.83	0.77
1:C:151:SER:H	1:C:154:MET:HE3	1.49	0.77
1:B:176:LEU:O	1:B:181:PHE:HB2	1.84	0.77
1:D:1:VAL:HB	1:D:215:TRP:CD1	2.20	0.76
1:C:1:VAL:HG12	1:C:2:GLY:H	1.49	0.76
1:A:50:ILE:CD1	1:A:117:GLU:HB3	2.16	0.75
1:B:238:MET:HE3	1:B:243:GLN:HG2	1.67	0.75
1:C:197:ALA:O	1:C:201:ILE:HG13	1.87	0.75
1:A:40:ASP:HB3	1:A:43:ALA:HB3	1.70	0.74
1:A:118:ARG:O	1:A:121:ILE:HD12	1.88	0.73
1:A:241:GLU:O	1:A:245:GLU:HG3	1.88	0.73
1:C:76:LYS:HE3	1:C:123:GLU:HG3	1.71	0.73
1:B:48:PRO:HG3	1:D:49:ILE:HD11	1.70	0.73
1:A:255:VAL:HG23	1:A:267:ASN:HD22	1.54	0.72
1:B:88:GLN:O	1:B:92:GLU:HG3	1.89	0.72
1:C:169:TRP:HB3	1:C:170:PRO:HD3	1.72	0.72
1:A:271:ALA:O	1:A:272:ARG:HB3	1.91	0.71
1:D:201:ILE:O	1:D:205:GLU:HG2	1.91	0.71
1:A:202:GLN:HA	1:A:202:GLN:NE2	2.05	0.71
1:B:1:VAL:HA	1:B:215:TRP:CD1	2.27	0.70
1:D:199:ALA:O	1:D:203:ILE:HG13	1.90	0.70
1:B:240:LYS:HB3	1:B:240:LYS:HZ3	1.55	0.69
1:C:233:GLN:HA	1:C:233:GLN:NE2	2.06	0.69
1:B:1:VAL:HA	1:B:215:TRP:NE1	2.07	0.69
1:D:132:TYR:O	1:D:180:GLY:HA2	1.93	0.69
1:B:169:TRP:HB3	1:B:170:PRO:HD3	1.72	0.69
1:C:9:LEU:HD22	1:C:22:LYS:HD3	1.75	0.69
1:D:9:LEU:HD22	1:D:22:LYS:HD3	1.73	0.69
1:B:236:PHE:HD2	1:D:160:ILE:HD12	1.57	0.69
1:C:1:VAL:HA	1:C:215:TRP:NE1	2.07	0.69
1:C:143:VAL:HG22	1:C:184:LEU:HB2	1.73	0.69
1:D:127:THR:CG2	1:D:130:ALA:HB3	2.23	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:50:ILE:HD11	1:A:117:GLU:HB3	1.74	0.68
1:D:64:ASN:HD21	1:D:66:GLN:HE21	1.42	0.68
1:C:1:VAL:HA	1:C:215:TRP:CD1	2.29	0.68
1:A:148:THR:HG23	3:A:601:FAD:O2	1.93	0.68
1:D:132:TYR:HA	1:D:177:HIS:O	1.95	0.66
1:A:4:ARG:HG2	1:A:35:GLU:OE2	1.95	0.66
1:A:169:TRP:HB3	1:A:170:PRO:HD3	1.78	0.66
1:A:246:GLU:HA	1:A:249:LYS:HG3	1.77	0.66
1:B:160:ILE:HG12	1:D:236:PHE:HB3	1.77	0.65
1:A:108:VAL:HG13	1:A:112:LEU:HB3	1.77	0.65
1:C:29:LEU:O	1:C:34:TRP:HB2	1.97	0.65
1:D:1:VAL:HG23	1:D:2:GLY:N	2.10	0.65
1:C:138:ARG:HA	1:C:180:GLY:O	1.97	0.64
1:B:50:ILE:HD12	1:B:117:GLU:HB3	1.80	0.64
1:B:108:VAL:CG1	1:B:112:LEU:HB3	2.27	0.64
1:C:14:ARG:HH11	1:C:14:ARG:HB2	1.63	0.63
1:C:57:GLY:H	1:C:79:HIS:HB3	1.63	0.63
1:C:271:ALA:O	1:C:272:ARG:HB2	1.96	0.63
1:A:58:LYS:N	1:A:58:LYS:HD2	2.13	0.63
1:B:185:GLU:HG3	4:B:707:HOH:O	1.98	0.63
1:A:4:ARG:HG2	1:A:35:GLU:CD	2.24	0.62
1:B:83:ASP:OD2	1:B:118:ARG:NH2	2.31	0.62
1:B:214:ILE:HG23	1:B:215:TRP:N	2.14	0.62
1:A:238:MET:HE2	1:A:243:GLN:HG2	1.81	0.62
1:C:1:VAL:HG12	1:C:2:GLY:N	2.14	0.62
1:A:158:GLN:HB3	1:C:238:MET:HE2	1.82	0.62
1:D:1:VAL:HG23	1:D:2:GLY:H	1.63	0.62
1:C:14:ARG:HB2	1:C:14:ARG:NH1	2.14	0.62
1:B:221:TYR:CB	1:B:273:LYS:HA	2.30	0.62
1:C:231:ASN:OD1	1:C:233:GLN:HB3	2.00	0.62
1:D:1:VAL:HB	1:D:215:TRP:NE1	2.15	0.62
1:B:214:ILE:O	1:B:217:GLU:HG3	2.00	0.61
1:B:112:LEU:HD22	1:B:116:PHE:CE2	2.35	0.61
1:B:135:GLY:O	1:B:138:ARG:CG	2.45	0.61
1:D:52:ARG:HG2	4:D:712:HOH:O	2.00	0.61
1:D:197:ALA:O	1:D:201:ILE:HG13	2.00	0.61
1:B:84:ILE:O	1:B:88:GLN:HG3	2.00	0.61
1:C:50:ILE:HG22	1:C:118:ARG:HG2	1.82	0.60
1:C:122:GLY:O	1:C:123:GLU:HB2	2.01	0.60
1:A:1:VAL:HG12	1:A:2:GLY:H	1.65	0.60
1:C:240:LYS:HD2	1:C:240:LYS:N	2.13	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:127:THR:HG23	1:D:130:ALA:H	1.66	0.60
1:C:130:ALA:HB1	1:C:134:LYS:O	2.01	0.60
1:B:54:ASP:OD1	1:B:118:ARG:NH1	2.34	0.60
1:A:272:ARG:O	1:A:272:ARG:HG2	2.00	0.60
1:A:50:ILE:HD12	1:A:117:GLU:HB3	1.83	0.60
1:D:50:ILE:HG22	1:D:118:ARG:HG2	1.83	0.59
1:B:47:ASN:ND2	1:B:49:ILE:H	2.00	0.59
1:B:127:THR:HG22	1:B:129:ALA:H	1.66	0.59
1:B:198:ASP:O	1:B:202:GLN:HG2	2.02	0.59
1:D:241:GLU:O	1:D:245:GLU:HG3	2.02	0.59
1:B:134:LYS:HB2	1:B:134:LYS:NZ	2.17	0.59
1:C:151:SER:N	1:C:154:MET:HE3	2.17	0.59
1:C:240:LYS:H	1:C:240:LYS:CD	2.15	0.59
1:A:138:ARG:HA	1:A:180:GLY:O	2.02	0.59
1:D:60:LYS:HD2	1:D:73:LEU:HD22	1.85	0.59
1:D:17:PHE:HB2	3:D:604:FAD:C5B	2.31	0.59
1:A:41:LEU:HD23	1:A:44:MET:CE	2.33	0.58
1:C:57:GLY:HA3	1:C:79:HIS:ND1	2.18	0.58
1:B:158:GLN:HB3	1:D:238:MET:HE2	1.85	0.58
1:B:201:ILE:O	1:B:205:GLU:HG2	2.03	0.58
1:D:29:LEU:HD11	1:D:211:LEU:HD22	1.85	0.58
1:B:81:SER:O	1:B:85:VAL:HG23	2.04	0.58
1:B:50:ILE:HD11	1:B:117:GLU:O	2.03	0.58
1:C:243:GLN:O	1:C:247:LYS:HG3	2.04	0.58
1:D:248:ASN:HB3	4:D:731:HOH:O	2.03	0.58
1:B:147:THR:HA	1:B:189:THR:OG1	2.04	0.58
1:C:84:ILE:O	1:C:88:GLN:HG3	2.04	0.57
1:A:160:ILE:HG13	1:C:236:PHE:HB3	1.85	0.57
1:C:255:VAL:HG23	1:C:267:ASN:HD22	1.68	0.57
1:A:204:LEU:O	1:A:208:LYS:HG3	2.04	0.57
1:D:127:THR:HG23	1:D:130:ALA:HB3	1.87	0.57
1:A:7:ILE:HG21	1:A:22:LYS:HG2	1.85	0.57
1:D:6:LEU:HD12	1:D:37:VAL:O	2.03	0.57
1:A:200:ARG:NH1	3:A:601:FAD:H1B	2.20	0.57
1:A:238:MET:CE	1:A:243:GLN:HG2	2.34	0.57
1:C:17:PHE:HB2	3:C:603:FAD:H52A	1.87	0.56
1:D:143:VAL:CG1	1:D:184:LEU:HB2	2.26	0.56
1:B:50:ILE:HG22	1:B:50:ILE:O	2.03	0.56
1:B:236:PHE:CD2	1:D:160:ILE:HD12	2.39	0.56
1:B:134:LYS:HA	1:B:138:ARG:HD3	1.86	0.56
1:B:148:THR:HG23	3:B:602:FAD:O2	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:239:LYS:HB2	1:B:242:VAL:HG23	1.86	0.56
1:A:84:ILE:O	1:A:88:GLN:HG3	2.06	0.56
1:A:197:ALA:O	1:A:201:ILE:HG12	2.06	0.56
1:B:132:TYR:O	1:B:180:GLY:HA2	2.06	0.56
1:D:21:MET:CE	1:D:204:LEU:HA	2.36	0.56
1:D:72:VAL:O	1:D:76:LYS:HG3	2.06	0.56
1:C:176:LEU:O	1:C:181:PHE:HB2	2.06	0.56
1:D:4:ARG:HG3	1:D:35:GLU:HB2	1.88	0.56
1:D:122:GLY:C	1:D:124:PHE:H	2.14	0.55
1:C:24:ALA:O	1:C:208:LYS:HE3	2.06	0.55
1:D:240:LYS:CD	1:D:240:LYS:H	2.20	0.55
1:A:238:MET:HE3	1:A:243:GLN:HA	1.88	0.55
1:C:32:LYS:HE3	1:C:212:GLU:HB3	1.89	0.55
1:B:75:TYR:HB2	1:B:80:LEU:HD11	1.88	0.55
1:A:80:LEU:O	1:A:81:SER:C	2.50	0.55
1:C:146:ILE:HG22	1:C:147:THR:N	2.22	0.55
1:C:250:LYS:HE3	1:C:251:PHE:CZ	2.42	0.55
1:D:11:HIS:CE1	1:D:16:SER:HB3	2.42	0.55
1:B:54:ASP:HA	1:B:81:SER:HB2	1.88	0.54
1:C:198:ASP:OD1	1:C:199:ALA:N	2.40	0.54
1:A:47:ASN:HB3	1:A:118:ARG:NH2	2.22	0.54
1:B:230:LEU:HD23	1:D:160:ILE:CD1	2.38	0.54
1:D:240:LYS:H	1:D:240:LYS:HD2	1.71	0.54
1:C:209:LYS:HD3	1:C:210:ARG:H	1.72	0.54
1:A:50:ILE:HG13	1:A:118:ARG:HG2	1.89	0.54
1:B:104:GLN:HA	3:B:602:FAD:C5X	2.38	0.54
1:A:108:VAL:CG1	1:A:112:LEU:HB3	2.37	0.54
1:C:14:ARG:HH11	1:C:14:ARG:CB	2.21	0.54
1:C:169:TRP:CZ2	1:C:256:GLY:HA3	2.43	0.54
1:A:160:ILE:CG1	1:C:236:PHE:HB3	2.37	0.54
1:B:134:LYS:C	1:B:138:ARG:HD3	2.33	0.54
1:C:14:ARG:HG2	1:C:19:TYR:CZ	2.43	0.54
1:A:91:LEU:HD11	1:A:120:PHE:HE1	1.73	0.53
1:D:250:LYS:HD3	1:D:251:PHE:CE2	2.43	0.53
1:D:21:MET:CE	1:D:204:LEU:HD23	2.34	0.53
1:A:202:GLN:HE21	1:A:202:GLN:CA	2.04	0.53
1:D:255:VAL:HG23	1:D:267:ASN:HD22	1.72	0.53
1:A:1:VAL:CG1	1:A:3:ARG:HH21	2.22	0.53
1:A:231:ASN:HD21	1:A:233:GLN:HG2	1.74	0.53
1:A:14:ARG:N	1:A:14:ARG:NE	2.39	0.53
1:A:131:MET:HE1	1:A:236:PHE:CD2	2.44	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:238:MET:HE2	1:D:158:GLN:CG	2.29	0.52
1:D:239:LYS:HB2	1:D:242:VAL:HG23	1.92	0.52
1:A:1:VAL:HG12	1:A:2:GLY:N	2.23	0.52
1:A:1:VAL:HG13	1:A:215:TRP:CD1	2.44	0.52
1:C:134:LYS:HB3	4:C:728:HOH:O	2.10	0.52
1:C:255:VAL:H	1:C:267:ASN:ND2	2.07	0.52
1:D:158:GLN:HE21	1:D:158:GLN:N	2.08	0.52
1:B:169:TRP:CZ2	1:B:256:GLY:HA3	2.44	0.52
1:B:144:LEU:HD12	1:B:172:GLN:HE21	1.74	0.52
1:C:89:LYS:HD2	4:C:731:HOH:O	2.10	0.52
1:D:227:LEU:HB3	1:D:242:VAL:HG11	1.92	0.52
1:A:75:TYR:CZ	1:A:124:PHE:HB2	2.45	0.52
1:B:132:TYR:HA	1:B:177:HIS:O	2.09	0.52
1:D:130:ALA:HB1	1:D:134:LYS:O	2.09	0.52
1:A:113:LYS:O	1:A:117:GLU:HG3	2.10	0.52
1:C:47:ASN:ND2	1:C:49:ILE:H	2.08	0.52
1:B:130:ALA:O	1:B:135:GLY:HA2	2.10	0.51
1:D:44:MET:HE1	1:D:87:GLU:OE2	2.09	0.51
1:B:56:THR:HB	1:B:79:HIS:O	2.10	0.51
1:C:52:ARG:C	1:C:54:ASP:H	2.19	0.51
1:B:116:PHE:O	1:B:120:PHE:HB2	2.11	0.51
1:C:28:ALA:HA	1:C:31:LYS:HE2	1.92	0.51
1:D:30:LYS:C	1:D:32:LYS:H	2.18	0.51
1:A:271:ALA:C	1:A:272:ARG:HD3	2.36	0.51
1:B:18:ASN:OD1	1:B:100:GLN:NE2	2.40	0.51
1:B:67:TYR:HB3	1:B:68:PRO:HD3	1.93	0.51
1:A:50:ILE:HD11	1:A:117:GLU:O	2.10	0.51
1:A:135:GLY:O	1:A:138:ARG:NH1	2.37	0.51
1:B:197:ALA:O	1:B:201:ILE:HG12	2.11	0.51
1:B:108:VAL:HG13	1:B:112:LEU:HD13	1.92	0.51
1:A:2:GLY:O	1:A:3:ARG:CZ	2.59	0.50
1:B:134:LYS:HB2	1:B:134:LYS:HZ3	1.75	0.50
1:A:17:PHE:HB2	3:A:601:FAD:C5B	2.36	0.50
1:B:271:ALA:O	1:B:272:ARG:HB2	2.10	0.50
1:B:50:ILE:HG13	1:B:118:ARG:HG2	1.92	0.50
1:C:68:PRO:O	1:C:72:VAL:HG23	2.12	0.50
1:D:172:GLN:HB2	1:D:268:GLN:HE22	1.76	0.50
1:C:150:GLY:HA2	1:C:154:MET:HE3	1.93	0.50
1:A:50:ILE:CG1	1:A:118:ARG:HG2	2.42	0.50
1:B:127:THR:HG22	1:B:129:ALA:N	2.27	0.50
1:C:207:TRP:O	1:C:211:LEU:HG	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:108:VAL:HG13	1:B:112:LEU:HB3	1.92	0.50
1:B:230:LEU:CD2	1:D:160:ILE:HD13	2.42	0.50
1:B:236:PHE:HB3	1:D:160:ILE:HG13	1.94	0.50
1:C:273:LYS:H	1:C:273:LYS:HD2	1.77	0.50
1:D:104:GLN:HA	3:D:604:FAD:N5	2.27	0.50
1:C:31:LYS:C	1:C:33:GLY:H	2.20	0.50
1:D:141:LYS:HA	1:D:182:GLN:O	2.12	0.50
1:A:52:ARG:HG2	4:A:704:HOH:O	2.11	0.49
1:B:57:GLY:H	1:B:79:HIS:CD2	2.30	0.49
1:A:13:GLU:OE1	1:A:14:ARG:NH2	2.45	0.49
1:D:64:ASN:HD21	1:D:66:GLN:NE2	2.08	0.49
1:A:132:TYR:O	1:A:180:GLY:HA2	2.12	0.49
1:A:202:GLN:NE2	1:A:202:GLN:CA	2.67	0.49
1:A:244:ASP:HA	1:A:247:LYS:HE3	1.95	0.49
1:B:112:LEU:O	1:B:115:TRP:HB3	2.12	0.49
1:C:104:GLN:HA	3:C:603:FAD:C5X	2.42	0.49
1:A:14:ARG:HA	1:A:19:TYR:CG	2.47	0.49
1:D:169:TRP:CZ2	1:D:256:GLY:HA3	2.47	0.49
1:B:155:TYR:HB3	1:B:164:MET:HB2	1.94	0.49
1:B:241:GLU:O	1:B:245:GLU:HG3	2.12	0.49
1:C:241:GLU:O	1:C:245:GLU:HB2	2.13	0.49
1:B:148:THR:HB	1:B:190:TYR:HA	1.94	0.49
1:B:195:THR:HG22	1:B:199:ALA:HB3	1.95	0.49
1:B:239:LYS:HB2	1:B:242:VAL:CG2	2.43	0.49
1:D:54:ASP:OD1	1:D:118:ARG:NH1	2.46	0.48
1:B:169:TRP:HA	1:B:268:GLN:HE21	1.77	0.48
1:B:236:PHE:O	1:D:160:ILE:HG13	2.13	0.48
1:C:65:PHE:CE1	1:C:70:GLU:HG3	2.48	0.48
1:D:104:GLN:HA	3:D:604:FAD:C5X	2.44	0.48
1:B:29:LEU:CD1	1:B:211:LEU:HB3	2.43	0.48
1:B:104:GLN:HE21	1:D:117:GLU:CD	2.21	0.48
1:B:158:GLN:HB3	1:D:238:MET:CE	2.43	0.48
1:B:257:HIS:HA	4:D:729:HOH:O	2.13	0.48
1:D:264:PRO:HB2	1:D:273:LYS:HG2	1.96	0.47
1:B:91:LEU:O	1:B:140:LYS:HE3	2.14	0.47
1:B:32:LYS:HE3	4:B:731:HOH:O	2.14	0.47
1:C:271:ALA:O	1:C:272:ARG:CB	2.61	0.47
1:D:65:PHE:CD1	1:D:70:GLU:HG3	2.50	0.47
1:B:106:PHE:CD1	1:B:106:PHE:N	2.83	0.47
1:B:148:THR:CG2	1:B:149:GLY:N	2.78	0.47
1:A:50:ILE:HG23	1:A:67:TYR:CD1	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:50:ILE:HD11	1:B:117:GLU:C	2.40	0.47
1:C:57:GLY:HA3	1:C:79:HIS:CG	2.50	0.47
1:A:45:ASN:HD22	1:A:45:ASN:HA	1.47	0.47
1:C:172:GLN:HB2	1:C:268:GLN:NE2	2.30	0.47
1:D:59:LEU:HB2	1:D:62:PRO:HG3	1.97	0.47
1:D:198:ASP:HA	1:D:201:ILE:HD12	1.97	0.47
1:A:75:TYR:CE1	1:A:124:PHE:HB2	2.50	0.47
1:A:122:GLY:O	1:A:123:GLU:HB3	2.14	0.47
1:D:12:SER:HB3	1:D:42:TYR:CD1	2.50	0.47
1:A:243:GLN:O	1:A:247:LYS:HG3	2.15	0.47
1:D:42:TYR:HD2	4:D:703:HOH:O	1.97	0.47
1:B:47:ASN:ND2	1:B:47:ASN:C	2.71	0.47
1:B:240:LYS:HB3	1:B:240:LYS:HZ2	1.73	0.47
1:C:76:LYS:CE	1:C:123:GLU:HG3	2.44	0.47
1:D:189:THR:O	1:D:189:THR:OG1	2.33	0.47
1:B:27:ALA:O	1:B:31:LYS:HG3	2.15	0.46
1:C:59:LEU:HB2	1:C:62:PRO:HG3	1.98	0.46
1:D:60:LYS:HD2	1:D:73:LEU:CD2	2.45	0.46
1:B:55:ILE:HG23	1:B:74:ALA:HB2	1.97	0.46
1:D:83:ASP:OD2	1:D:118:ARG:NH2	2.48	0.46
1:B:236:PHE:HB3	1:D:160:ILE:CD1	2.44	0.46
1:C:113:LYS:O	1:C:117:GLU:HG3	2.15	0.46
1:D:239:LYS:HB2	1:D:242:VAL:CG2	2.46	0.46
1:A:154:MET:HE2	1:A:161:HIS:NE2	2.31	0.46
1:C:17:PHE:HB2	3:C:603:FAD:C5B	2.46	0.46
1:C:72:VAL:HG13	1:C:123:GLU:HG2	1.96	0.46
1:C:207:TRP:CH2	1:C:211:LEU:HD21	2.50	0.46
1:C:205:GLU:OE1	1:C:205:GLU:HA	2.15	0.46
1:A:67:TYR:HB3	1:A:68:PRO:HD3	1.96	0.46
1:C:195:THR:CG2	1:C:200:ARG:HG3	2.46	0.46
1:D:176:LEU:O	1:D:181:PHE:HB2	2.15	0.46
1:C:4:ARG:HG2	1:C:35:GLU:HB3	1.97	0.46
1:C:40:ASP:HB3	1:C:43:ALA:HB3	1.96	0.46
1:C:150:GLY:CA	1:C:154:MET:HE3	2.46	0.46
1:A:127:THR:HG22	1:A:129:ALA:N	2.21	0.46
1:B:214:ILE:CG2	1:B:215:TRP:N	2.79	0.46
1:C:127:THR:OG1	1:C:130:ALA:N	2.49	0.46
1:D:185:GLU:HB3	1:D:269:ILE:O	2.16	0.46
1:A:116:PHE:O	1:A:120:PHE:HB2	2.15	0.46
1:B:4:ARG:HH21	1:B:93:ALA:HB1	1.81	0.46
1:D:78:GLY:HA2	4:D:723:HOH:O	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:151:SER:H	1:D:154:MET:HE3	1.81	0.45
1:D:271:ALA:O	1:D:272:ARG:HB2	2.17	0.45
1:A:91:LEU:HD11	1:A:120:PHE:CE1	2.52	0.45
1:C:209:LYS:HA	1:C:212:GLU:OE2	2.15	0.45
1:D:229:ASP:OD1	1:D:239:LYS:HG2	2.15	0.45
1:C:14:ARG:HG2	1:C:19:TYR:CE1	2.51	0.45
1:C:9:LEU:CD2	1:C:22:LYS:HD3	2.46	0.45
1:A:66:GLN:HE21	1:A:66:GLN:CA	2.12	0.45
1:A:207:TRP:HA	1:A:207:TRP:CE3	2.52	0.45
1:C:214:ILE:O	1:C:217:GLU:HB2	2.17	0.45
1:A:94:ALA:O	1:A:140:LYS:NZ	2.38	0.45
1:D:172:GLN:HB2	1:D:268:GLN:NE2	2.31	0.45
1:A:11:HIS:NE2	3:A:601:FAD:O2P	2.48	0.45
1:B:122:GLY:O	1:B:123:GLU:CB	2.64	0.45
1:B:271:ALA:O	1:B:272:ARG:CB	2.65	0.45
1:C:99:PHE:HB2	1:C:144:LEU:HD23	1.98	0.45
1:C:165:ASN:ND2	1:C:266:ASP:HA	2.32	0.45
1:A:132:TYR:HA	1:A:177:HIS:O	2.16	0.45
1:C:36:VAL:HG12	1:C:37:VAL:N	2.31	0.45
1:C:207:TRP:HZ3	1:C:210:ARG:HD3	1.82	0.45
1:B:221:TYR:HB2	1:B:273:LYS:HA	1.97	0.45
1:D:19:TYR:O	1:D:22:LYS:HB3	2.16	0.45
1:D:68:PRO:HG2	2:D:702:340:H453	1.98	0.45
1:C:145:SER:HA	1:C:187:GLN:HB3	1.97	0.45
1:D:209:LYS:HD3	1:D:209:LYS:C	2.42	0.45
1:B:61:ASP:OD1	1:B:64:ASN:HB2	2.18	0.44
1:C:1:VAL:CG1	1:C:2:GLY:H	2.25	0.44
1:C:125:ALA:O	1:C:136:PRO:HD2	2.16	0.44
1:C:246:GLU:HG3	4:C:716:HOH:O	2.17	0.44
1:D:97:VAL:HG21	1:D:181:PHE:HZ	1.82	0.44
1:A:117:GLU:OE2	1:C:109:PRO:HA	2.17	0.44
1:B:190:TYR:O	1:B:191:SER:C	2.60	0.44
1:D:231:ASN:HD21	1:D:234:ALA:HB2	1.82	0.44
1:B:134:LYS:CA	1:B:138:ARG:HD3	2.47	0.44
1:A:237:LEU:HD22	1:C:158:GLN:HE21	1.82	0.44
1:B:4:ARG:HG2	1:B:35:GLU:CD	2.43	0.44
1:A:207:TRP:HA	1:A:207:TRP:HE3	1.82	0.44
1:B:4:ARG:NH2	1:B:93:ALA:HB1	2.33	0.44
1:A:148:THR:HG22	1:A:150:GLY:N	2.32	0.44
1:A:165:ASN:OD1	1:A:268:GLN:HB2	2.17	0.44
1:C:224:PRO:HD2	1:C:227:LEU:HD11	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:258:HIS:HB3	1:D:263:ILE:HG12	1.99	0.44
1:C:229:ASP:O	1:C:236:PHE:HA	2.17	0.44
1:B:134:LYS:HE3	1:B:225:SER:OG	2.18	0.44
1:B:202:GLN:HA	1:B:202:GLN:NE2	2.33	0.44
1:A:12:SER:HB3	1:A:42:TYR:CE1	2.52	0.43
1:B:122:GLY:O	1:B:123:GLU:HB2	2.17	0.43
1:B:230:LEU:HD23	1:D:160:ILE:HD11	2.00	0.43
1:A:50:ILE:HD11	1:A:117:GLU:C	2.43	0.43
1:B:64:ASN:ND2	4:B:716:HOH:O	2.50	0.43
1:C:52:ARG:C	1:C:54:ASP:N	2.75	0.43
1:D:196:PRO:O	1:D:200:ARG:HG3	2.18	0.43
1:B:31:LYS:HZ1	1:B:208:LYS:CE	2.32	0.43
1:C:25:ALA:HA	1:C:211:LEU:CD1	2.48	0.43
1:C:211:LEU:HA	1:C:214:ILE:HB	2.00	0.43
1:B:50:ILE:CD1	1:B:117:GLU:HB3	2.46	0.43
1:B:74:ALA:HA	1:B:79:HIS:NE2	2.33	0.43
1:B:191:SER:HB2	1:B:194:HIS:HB2	2.01	0.43
1:C:47:ASN:ND2	1:C:47:ASN:C	2.76	0.43
1:D:1:VAL:CG2	1:D:2:GLY:H	2.26	0.43
1:D:127:THR:HG23	1:D:130:ALA:N	2.33	0.43
1:B:25:ALA:O	1:B:29:LEU:HB2	2.19	0.43
1:B:141:LYS:HG2	1:B:182:GLN:HB2	2.01	0.43
1:C:143:VAL:HG12	1:C:144:LEU:N	2.33	0.43
1:B:169:TRP:CB	1:B:170:PRO:HD3	2.44	0.43
1:C:146:ILE:HG22	1:C:147:THR:H	1.83	0.43
1:A:81:SER:O	1:A:85:VAL:HG23	2.19	0.43
1:C:169:TRP:CB	1:C:170:PRO:HD3	2.45	0.43
1:D:56:THR:HG21	1:D:82:PRO:HD3	1.99	0.43
1:B:236:PHE:HD2	1:D:160:ILE:CD1	2.29	0.43
1:A:148:THR:HG22	1:A:149:GLY:N	2.33	0.43
1:C:6:LEU:O	1:C:97:VAL:HA	2.18	0.43
1:A:54:ASP:OD1	1:A:118:ARG:NH1	2.52	0.42
1:D:48:PRO:O	1:D:114:GLY:HA3	2.19	0.42
1:D:64:ASN:ND2	1:D:66:GLN:HE21	2.13	0.42
1:A:229:ASP:O	1:A:231:ASN:N	2.52	0.42
1:B:31:LYS:HZ1	1:B:208:LYS:NZ	2.16	0.42
1:B:128:TYR:O	1:B:131:MET:HG3	2.20	0.42
1:C:272:ARG:HA	4:C:702:HOH:O	2.19	0.42
1:D:68:PRO:O	1:D:72:VAL:HG23	2.19	0.42
1:A:42:TYR:HB2	4:A:706:HOH:O	2.19	0.42
1:B:105:TRP:CZ2	1:D:120:PHE:HB3	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:154:MET:O	1:B:161:HIS:HB2	2.18	0.42
1:C:150:GLY:HA2	1:C:154:MET:CE	2.49	0.42
1:A:4:ARG:CG	1:A:35:GLU:OE2	2.67	0.42
1:B:46:PHE:O	1:B:48:PRO:HD3	2.19	0.42
1:D:48:PRO:HB2	1:D:114:GLY:HA3	2.00	0.42
1:B:54:ASP:O	1:B:81:SER:N	2.49	0.42
1:B:148:THR:CG2	1:B:150:GLY:O	2.68	0.42
1:C:132:TYR:O	1:C:180:GLY:HA2	2.18	0.42
1:A:67:TYR:N	1:A:68:PRO:CD	2.83	0.42
1:B:138:ARG:HG3	1:B:138:ARG:HH11	1.83	0.42
1:C:130:ALA:O	1:C:135:GLY:HA2	2.20	0.42
1:D:131:MET:O	1:D:132:TYR:C	2.63	0.42
1:A:148:THR:CG2	1:A:149:GLY:N	2.82	0.42
1:C:67:TYR:HB3	1:C:68:PRO:HD3	2.02	0.42
1:C:239:LYS:HB3	1:C:241:GLU:OE2	2.20	0.42
1:D:12:SER:HB3	1:D:42:TYR:CE1	2.55	0.42
1:D:151:SER:OG	1:D:154:MET:HE3	2.20	0.42
1:B:27:ALA:C	1:B:31:LYS:HE3	2.44	0.41
1:B:204:LEU:O	1:B:208:LYS:HG3	2.19	0.41
1:B:211:LEU:HA	1:B:214:ILE:HB	2.02	0.41
1:C:122:GLY:HA2	1:C:126:TYR:CE1	2.55	0.41
1:C:197:ALA:HA	1:C:200:ARG:CZ	2.49	0.41
1:A:232:PHE:CD1	1:A:232:PHE:N	2.88	0.41
1:A:254:SER:HB2	1:A:267:ASN:HD21	1.85	0.41
1:B:104:GLN:HA	3:B:602:FAD:N5	2.34	0.41
1:A:235:GLY:O	1:A:236:PHE:HB2	2.20	0.41
1:B:11:HIS:CE1	1:B:16:SER:HB3	2.55	0.41
1:C:242:VAL:O	1:C:246:GLU:HB2	2.19	0.41
1:C:257:HIS:HA	4:C:709:HOH:O	2.20	0.41
1:A:166:VAL:HG13	1:C:166:VAL:HG13	2.03	0.41
1:B:158:GLN:HB3	1:B:158:GLN:HE21	1.68	0.41
1:A:4:ARG:HD3	1:A:93:ALA:O	2.20	0.41
1:A:47:ASN:ND2	4:A:738:HOH:O	2.54	0.41
1:B:31:LYS:NZ	1:B:208:LYS:NZ	2.69	0.41
1:B:77:GLU:HB2	1:B:79:HIS:CE1	2.56	0.41
1:C:155:TYR:HB3	1:C:164:MET:HB2	2.03	0.41
1:C:195:THR:HG22	1:C:200:ARG:HG3	2.01	0.41
1:C:254:SER:HB2	1:C:267:ASN:ND2	2.36	0.41
1:C:273:LYS:N	1:C:273:LYS:NZ	2.68	0.41
1:D:187:GLN:HG2	1:D:207:TRP:CE3	2.56	0.41
1:A:57:GLY:H	1:A:79:HIS:CD2	2.39	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:224:PRO:O	1:C:227:LEU:HG	2.21	0.41
1:B:195:THR:HA	1:B:196:PRO:HD3	1.85	0.41
1:C:89:LYS:HG3	4:C:730:HOH:O	2.21	0.41
1:C:192:ILE:CG2	3:C:603:FAD:H5'1	2.51	0.41
1:D:122:GLY:C	1:D:124:PHE:N	2.79	0.41
1:A:3:ARG:NH1	1:A:3:ARG:HA	2.35	0.41
1:B:14:ARG:HE	1:B:14:ARG:HA	1.85	0.41
1:B:132:TYR:O	1:B:134:LYS:N	2.53	0.41
1:D:30:LYS:C	1:D:32:LYS:N	2.77	0.41
1:D:195:THR:HG21	1:D:203:ILE:HD11	2.02	0.41
1:D:215:TRP:CD1	1:D:215:TRP:O	2.74	0.41
1:D:240:LYS:HD2	1:D:240:LYS:N	2.36	0.40
1:A:58:LYS:N	1:A:58:LYS:CD	2.83	0.40
1:B:214:ILE:HD12	1:B:217:GLU:CD	2.46	0.40
1:B:227:LEU:HB3	1:B:242:VAL:HG11	2.03	0.40
1:A:13:GLU:HA	1:A:14:ARG:HH21	1.85	0.40
1:A:214:ILE:HG23	1:A:215:TRP:N	2.36	0.40
1:B:59:LEU:HB3	1:B:70:GLU:OE2	2.22	0.40
1:C:272:ARG:O	1:C:273:LYS:O	2.39	0.40
1:B:17:PHE:HB2	3:B:602:FAD:C5B	2.43	0.40
1:D:265:THR:O	1:D:266:ASP:C	2.65	0.40
1:A:74:ALA:HB1	1:A:79:HIS:O	2.21	0.40
1:A:244:ASP:HA	1:A:247:LYS:CE	2.52	0.40
1:B:117:GLU:HG2	1:D:105:TRP:HD1	1.86	0.40
1:B:147:THR:HG22	1:B:189:THR:OG1	2.21	0.40
1:C:17:PHE:O	1:C:21:MET:HG2	2.22	0.40
1:C:31:LYS:HE2	1:C:31:LYS:HB2	1.87	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	271/273 (99%)	244 (90%)	23 (8%)	4 (2%)	8	8
1	B	271/273 (99%)	248 (92%)	19 (7%)	4 (2%)	8	8
1	C	271/273 (99%)	247 (91%)	19 (7%)	5 (2%)	6	6
1	D	271/273 (99%)	247 (91%)	18 (7%)	6 (2%)	5	4
All	All	1084/1092 (99%)	986 (91%)	79 (7%)	19 (2%)	6	6

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	272	ARG
1	C	272	ARG
1	B	123	GLU
1	A	245	GLU
1	B	272	ARG
1	D	41	LEU
1	D	191	SER
1	D	230	LEU
1	B	133	ASP
1	B	191	SER
1	C	132	TYR
1	A	230	LEU
1	C	32	LYS
1	D	31	LYS
1	C	57	GLY
1	C	129	ALA
1	D	132	TYR
1	D	272	ARG
1	A	57	GLY

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	227/227 (100%)	204 (90%)	23 (10%)	7	9
1	B	227/227 (100%)	216 (95%)	11 (5%)	23	35

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	227/227 (100%)	210 (92%)	17 (8%)	12	17
1	D	227/227 (100%)	214 (94%)	13 (6%)	18	27
All	All	908/908 (100%)	844 (93%)	64 (7%)	14	19

All (64) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	ARG
1	A	14	ARG
1	A	29	LEU
1	A	45	ASN
1	A	50	ILE
1	A	56	THR
1	A	60	LYS
1	A	66	GLN
1	A	108	VAL
1	A	112	LEU
1	A	121	ILE
1	A	143	VAL
1	A	158	GLN
1	A	160	ILE
1	A	182	GLN
1	A	202	GLN
1	A	207	TRP
1	A	231	ASN
1	A	241	GLU
1	A	248	ASN
1	A	250	LYS
1	A	261	LYS
1	A	272	ARG
1	B	14	ARG
1	B	50	ILE
1	B	64	ASN
1	B	112	LEU
1	B	134	LYS
1	B	158	GLN
1	B	160	ILE
1	B	209	LYS
1	B	218	THR
1	B	240	LYS
1	B	241	GLU

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Mol	Chain	Res	Type
1	C	14	ARG
1	C	29	LEU
1	C	31	LYS
1	C	44	MET
1	C	60	LYS
1	C	66	GLN
1	C	81	SER
1	C	158	GLN
1	C	209	LYS
1	C	217	GLU
1	C	233	GLN
1	C	240	LYS
1	C	241	GLU
1	C	245	GLU
1	C	265	THR
1	C	267	ASN
1	C	273	LYS
1	D	14	ARG
1	D	21	MET
1	D	23	GLU
1	D	30	LYS
1	D	48	PRO
1	D	52	ARG
1	D	82	PRO
1	D	123	GLU
1	D	143	VAL
1	D	158	GLN
1	D	189	THR
1	D	240	LYS
1	D	241	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (38) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	45	ASN
1	A	47	ASN
1	A	66	GLN
1	A	79	HIS
1	A	104	GLN
1	A	158	GLN
1	A	172	GLN
1	A	202	GLN

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Mol	Chain	Res	Type
1	A	267	ASN
1	A	268	GLN
1	B	47	ASN
1	B	66	GLN
1	B	79	HIS
1	B	104	GLN
1	B	158	GLN
1	B	172	GLN
1	B	202	GLN
1	B	231	ASN
1	B	248	ASN
1	B	267	ASN
1	B	268	GLN
1	C	47	ASN
1	C	64	ASN
1	C	158	GLN
1	C	172	GLN
1	C	233	GLN
1	C	267	ASN
1	C	268	GLN
1	D	45	ASN
1	D	64	ASN
1	D	66	GLN
1	D	79	HIS
1	D	104	GLN
1	D	158	GLN
1	D	172	GLN
1	D	243	GLN
1	D	267	ASN
1	D	268	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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

8 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	340	D	702	-	24,25,25	2.23	11 (45%)	27,36,36	2.43	4 (14%)
2	340	A	703	-	24,25,25	2.45	11 (45%)	27,36,36	4.80	6 (22%)
3	FAD	C	603	-	58,58,58	2.65	23 (39%)	85,89,89	1.13	9 (10%)
3	FAD	D	604	-	58,58,58	2.50	22 (37%)	85,89,89	1.12	6 (7%)
2	340	B	704	-	24,25,25	2.38	11 (45%)	27,36,36	2.87	5 (18%)
3	FAD	A	601	-	58,58,58	2.55	22 (37%)	85,89,89	1.14	6 (7%)
3	FAD	B	602	-	58,58,58	2.62	22 (37%)	85,89,89	1.08	5 (5%)
2	340	C	701	-	24,25,25	2.14	11 (45%)	27,36,36	5.56	3 (11%)

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
2	340	D	702	-	-	3/7/23/23	0/3/3/3
2	340	A	703	-	-	4/7/23/23	0/3/3/3
3	FAD	C	603	-	-	6/34/50/50	0/6/6/6
3	FAD	D	604	-	-	4/34/50/50	0/6/6/6
2	340	B	704	-	-	0/7/23/23	0/3/3/3
3	FAD	A	601	-	-	5/34/50/50	0/6/6/6
3	FAD	B	602	-	-	6/34/50/50	0/6/6/6
2	340	C	701	-	-	4/7/23/23	0/3/3/3

All (133) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	603	FAD	P-O3P	-12.93	1.45	1.59
3	B	602	FAD	P-O3P	-12.26	1.46	1.59
3	D	604	FAD	P-O3P	-11.92	1.46	1.59
3	A	601	FAD	P-O3P	-11.47	1.47	1.59
3	A	601	FAD	PA-O3P	6.22	1.66	1.59
3	B	602	FAD	PA-O3P	5.59	1.65	1.59
3	A	601	FAD	C9A-N10	4.88	1.49	1.41
3	B	602	FAD	C9A-N10	4.35	1.48	1.41
2	B	704	340	C25-C24	4.35	1.46	1.38
3	C	603	FAD	PA-O3P	4.30	1.64	1.59
3	C	603	FAD	PA-O2A	-4.30	1.35	1.55
3	C	603	FAD	C9A-N10	4.24	1.48	1.41
3	B	602	FAD	PA-O2A	-4.20	1.35	1.55
2	A	703	340	C29-C24	4.16	1.46	1.38
3	D	604	FAD	PA-O3P	4.15	1.64	1.59
2	A	703	340	C25-C24	4.15	1.46	1.38
3	D	604	FAD	C9A-N10	4.03	1.48	1.41
3	A	601	FAD	PA-O2A	-3.96	1.37	1.55
3	D	604	FAD	PA-O2A	-3.93	1.37	1.55
3	B	602	FAD	P-O2P	-3.91	1.37	1.55
3	A	601	FAD	P-O2P	-3.80	1.37	1.55
2	D	702	340	C25-C24	3.80	1.45	1.38
3	B	602	FAD	O5'-C5'	3.78	1.58	1.44
3	D	604	FAD	O5'-C5'	3.76	1.58	1.44
3	C	603	FAD	C4X-N5	3.74	1.38	1.30
3	D	604	FAD	C1'-C2'	3.74	1.57	1.52
3	C	603	FAD	C9A-C5X	3.71	1.47	1.41
2	D	702	340	C29-C24	3.63	1.45	1.38
3	C	603	FAD	O5'-C5'	3.63	1.58	1.44
3	D	604	FAD	P-O2P	-3.54	1.38	1.55
2	D	702	340	C8-C9	-3.51	1.32	1.40
3	C	603	FAD	P-O2P	-3.50	1.39	1.55
2	A	703	340	C26-C25	3.49	1.44	1.38
2	A	703	340	O20-C24	3.47	1.45	1.37
2	C	701	340	C25-C24	3.45	1.45	1.38
2	A	703	340	O44-C6	3.42	1.42	1.36
2	C	701	340	C8-C9	-3.39	1.32	1.40
2	D	702	340	C27-C26	3.37	1.45	1.38
3	B	602	FAD	C8-C7	3.36	1.49	1.40
2	D	702	340	C26-C25	3.31	1.44	1.38
3	A	601	FAD	O5'-C5'	3.30	1.57	1.44
2	B	704	340	C28-C27	3.30	1.45	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	704	340	C26-C25	3.29	1.44	1.38
2	B	704	340	C28-C29	3.29	1.44	1.38
2	B	704	340	C8-C9	-3.27	1.32	1.40
2	C	701	340	C29-C24	3.26	1.44	1.38
2	A	703	340	C27-C26	3.26	1.45	1.38
3	D	604	FAD	C8-C7	3.22	1.48	1.40
2	B	704	340	C29-C24	3.21	1.44	1.38
3	A	601	FAD	C9A-C5X	3.21	1.46	1.41
3	B	602	FAD	C6-C5X	3.21	1.44	1.40
3	C	603	FAD	C8-C7	3.18	1.48	1.40
3	A	601	FAD	O4B-C1B	3.16	1.49	1.42
3	B	602	FAD	C1'-C2'	3.14	1.57	1.52
2	A	703	340	C28-C29	3.14	1.44	1.38
2	C	701	340	C37-C8	3.12	1.54	1.49
3	A	601	FAD	C4X-N5	3.06	1.37	1.30
3	B	602	FAD	C9A-C5X	3.03	1.46	1.41
2	B	704	340	C37-C8	3.00	1.54	1.49
2	C	701	340	C26-C25	2.98	1.44	1.38
3	D	604	FAD	O4B-C1B	2.98	1.48	1.42
2	B	704	340	C27-C26	2.97	1.44	1.38
3	A	601	FAD	C10-N1	2.97	1.39	1.33
3	B	602	FAD	C4X-N5	2.94	1.37	1.30
3	A	601	FAD	C6-C5X	2.91	1.44	1.40
2	A	703	340	C28-C27	2.91	1.44	1.38
2	D	702	340	C28-C29	2.90	1.43	1.38
2	A	703	340	C37-C8	2.90	1.53	1.49
3	C	603	FAD	C4A-N3A	2.89	1.39	1.34
3	C	603	FAD	C1'-C2'	2.89	1.56	1.52
3	C	603	FAD	C6-C5X	2.89	1.44	1.40
3	D	604	FAD	C4A-N3A	2.86	1.39	1.34
3	B	602	FAD	C2A-N1A	2.85	1.39	1.33
2	B	704	340	O20-C24	2.84	1.44	1.37
3	C	603	FAD	C9-C9A	2.83	1.44	1.39
3	D	604	FAD	C5A-C6A	2.80	1.48	1.41
2	C	701	340	O44-C6	2.79	1.41	1.36
3	A	601	FAD	C9-C9A	2.79	1.44	1.39
3	A	601	FAD	C5A-C6A	2.79	1.48	1.41
3	B	602	FAD	C9-C9A	2.78	1.44	1.39
3	B	602	FAD	C10-N1	2.76	1.38	1.33
3	C	603	FAD	C10-N1	2.75	1.38	1.33
3	D	604	FAD	C9A-C5X	2.75	1.45	1.41
2	A	703	340	C8-C9	-2.75	1.33	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	702	340	C28-C27	2.74	1.44	1.38
2	C	701	340	C27-C26	2.73	1.44	1.38
2	C	701	340	C28-C29	2.69	1.43	1.38
3	D	604	FAD	C9-C9A	2.67	1.43	1.39
3	D	604	FAD	C4X-N5	2.65	1.36	1.30
2	C	701	340	C28-C27	2.65	1.44	1.38
3	C	603	FAD	C5A-C6A	2.64	1.48	1.41
3	B	602	FAD	C9-C8	2.63	1.43	1.39
2	C	701	340	C1-C2	2.61	1.52	1.44
2	D	702	340	C37-C8	2.61	1.53	1.49
2	B	704	340	O44-C6	2.61	1.41	1.36
2	B	704	340	C1-C2	2.57	1.51	1.44
2	D	702	340	O44-C6	2.57	1.41	1.36
3	C	603	FAD	O4B-C1B	2.56	1.47	1.42
3	B	602	FAD	C5A-C6A	2.53	1.48	1.41
3	D	604	FAD	C6-C5X	2.51	1.43	1.40
3	B	602	FAD	PA-O5B	-2.50	1.49	1.59
3	C	603	FAD	C5B-C4B	2.49	1.59	1.51
3	D	604	FAD	C5A-C4A	2.48	1.43	1.39
3	A	601	FAD	C5A-C4A	2.47	1.43	1.39
3	B	602	FAD	O4B-C1B	2.45	1.47	1.42
2	A	703	340	C1-C2	2.44	1.51	1.44
3	A	601	FAD	C5B-C4B	2.40	1.58	1.51
3	C	603	FAD	C9-C8	2.40	1.42	1.39
3	A	601	FAD	PA-O5B	-2.38	1.50	1.59
3	A	601	FAD	C8-C7	2.37	1.46	1.40
3	C	603	FAD	C2A-N1A	2.35	1.38	1.33
3	C	603	FAD	O4B-C4B	2.34	1.50	1.45
3	B	602	FAD	C4A-N3A	2.34	1.38	1.34
3	B	602	FAD	C2-N3	2.34	1.44	1.39
3	C	603	FAD	C5A-C4A	2.32	1.43	1.39
3	D	604	FAD	C2-N3	2.32	1.44	1.39
3	A	601	FAD	C2A-N1A	2.31	1.38	1.33
3	A	601	FAD	C4A-N3A	2.31	1.38	1.34
3	B	602	FAD	C5A-C4A	2.31	1.43	1.39
3	B	602	FAD	C5B-C4B	2.28	1.58	1.51
2	C	701	340	O20-C24	2.25	1.42	1.37
3	A	601	FAD	C1'-C2'	2.25	1.55	1.52
3	C	603	FAD	C2-N3	2.23	1.43	1.39
3	D	604	FAD	C10-N1	2.21	1.37	1.33
2	D	702	340	C1-C2	2.21	1.50	1.44
3	C	603	FAD	PA-O5B	-2.19	1.50	1.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	702	340	O20-C24	2.18	1.42	1.37
3	A	601	FAD	C9-C8	2.16	1.42	1.39
3	D	604	FAD	C2A-N1A	2.10	1.37	1.33
3	A	601	FAD	O2-C2	-2.10	1.20	1.24
3	D	604	FAD	PA-O5B	-2.09	1.51	1.59
3	D	604	FAD	C5B-C4B	2.02	1.57	1.51
3	D	604	FAD	O2-C2	-2.02	1.20	1.24

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	701	340	C19-O20-C24	27.75	167.34	117.63
2	A	703	340	C19-O20-C24	23.35	159.45	117.63
2	D	702	340	O20-C24-C25	7.15	155.28	120.01
2	B	704	340	O20-C19-C9	7.01	129.61	108.49
2	D	702	340	O20-C24-C29	-6.98	85.58	120.01
2	B	704	340	C19-O20-C24	6.87	129.93	117.63
2	B	704	340	O20-C24-C25	6.71	153.09	120.01
2	B	704	340	O20-C24-C29	-6.48	88.07	120.01
2	A	703	340	O20-C19-C9	5.09	123.82	108.49
2	C	701	340	O20-C19-C9	5.09	123.82	108.49
2	D	702	340	O20-C19-C9	4.48	121.98	108.49
2	C	701	340	C9-C4-C5	3.20	137.97	130.65
2	D	702	340	C9-C4-C5	3.15	137.85	130.65
3	D	604	FAD	O5B-PA-O1A	-2.99	97.09	108.94
3	A	601	FAD	O5B-PA-O1A	-2.94	97.29	108.94
2	B	704	340	C9-C4-C5	2.93	137.35	130.65
3	C	603	FAD	O5B-PA-O1A	-2.85	97.64	108.94
3	B	602	FAD	O5B-PA-O1A	-2.80	97.84	108.94
2	A	703	340	C9-C4-C5	2.73	136.89	130.65
3	C	603	FAD	C2A-N1A-C6A	2.69	123.16	118.73
3	D	604	FAD	C2A-N1A-C6A	2.63	123.06	118.73
2	A	703	340	O20-C24-C29	2.62	132.91	120.01
3	A	601	FAD	C2A-N1A-C6A	2.51	122.86	118.73
3	A	601	FAD	C5'-C4'-C3'	-2.49	107.53	112.22
3	C	603	FAD	C5'-C4'-C3'	-2.33	107.83	112.22
3	B	602	FAD	C5'-C4'-C3'	-2.31	107.86	112.22
3	B	602	FAD	C2A-N1A-C6A	2.31	122.52	118.73
3	D	604	FAD	C4-N3-C2	-2.28	121.60	125.64
2	A	703	340	O20-C24-C25	-2.27	108.81	120.01
3	A	601	FAD	O5B-C5B-C4B	-2.26	101.28	108.99
3	C	603	FAD	C5X-C9A-N10	-2.26	115.92	117.97

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	604	FAD	O2P-P-O3P	-2.24	101.21	107.27
3	B	602	FAD	C4-N3-C2	-2.23	121.69	125.64
3	C	603	FAD	O4B-C1B-N9A	-2.18	103.89	108.09
3	D	604	FAD	C5X-C9A-N10	-2.18	116.00	117.97
3	C	603	FAD	C4-N3-C2	-2.15	121.82	125.64
3	B	602	FAD	C5X-C9A-N10	-2.14	116.03	117.97
3	D	604	FAD	O4B-C1B-N9A	-2.12	104.02	108.09
3	A	601	FAD	O2P-P-O3P	-2.08	101.64	107.27
3	A	601	FAD	C4-N3-C2	-2.08	121.95	125.64
3	C	603	FAD	C5A-C4A-N9A	-2.08	103.55	105.81
2	A	703	340	C45-O44-C6	2.05	120.94	116.96
3	C	603	FAD	C10-N1-C2	2.04	121.26	116.85
3	C	603	FAD	O2P-P-O3P	-2.03	101.78	107.27

There are no chirality outliers.

All (32) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	703	340	C9-C19-O20-C24
2	D	702	340	O20-C19-C9-C4
2	D	702	340	O20-C19-C9-C8
2	A	703	340	C25-C24-O20-C19
2	A	703	340	C29-C24-O20-C19
2	D	702	340	C9-C19-O20-C24
3	B	602	FAD	C4'-C5'-O5'-P
3	C	603	FAD	C4'-C5'-O5'-P
3	D	604	FAD	C4'-C5'-O5'-P
3	A	601	FAD	P-O3P-PA-O5B
3	B	602	FAD	P-O3P-PA-O5B
3	C	603	FAD	P-O3P-PA-O5B
3	D	604	FAD	P-O3P-PA-O5B
3	C	603	FAD	C2B-C1B-N9A-C8A
3	A	601	FAD	C2B-C1B-N9A-C8A
3	B	602	FAD	C2B-C1B-N9A-C8A
3	D	604	FAD	C2B-C1B-N9A-C8A
2	C	701	340	C9-C19-O20-C24
3	A	601	FAD	C4'-C5'-O5'-P
3	C	603	FAD	C2B-C1B-N9A-C4A
2	C	701	340	C29-C24-O20-C19
2	C	701	340	C25-C24-O20-C19
3	A	601	FAD	PA-O3P-P-O5'
3	B	602	FAD	PA-O3P-P-O5'

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Mol	Chain	Res	Type	Atoms
3	C	603	FAD	PA-O3P-P-O5'
3	D	604	FAD	PA-O3P-P-O5'
3	A	601	FAD	O4B-C1B-N9A-C8A
3	B	602	FAD	O4B-C1B-N9A-C8A
3	C	603	FAD	O4B-C1B-N9A-C8A
2	A	703	340	O20-C19-C9-C4
2	C	701	340	O20-C19-C9-C4
3	B	602	FAD	P-O3P-PA-O2A

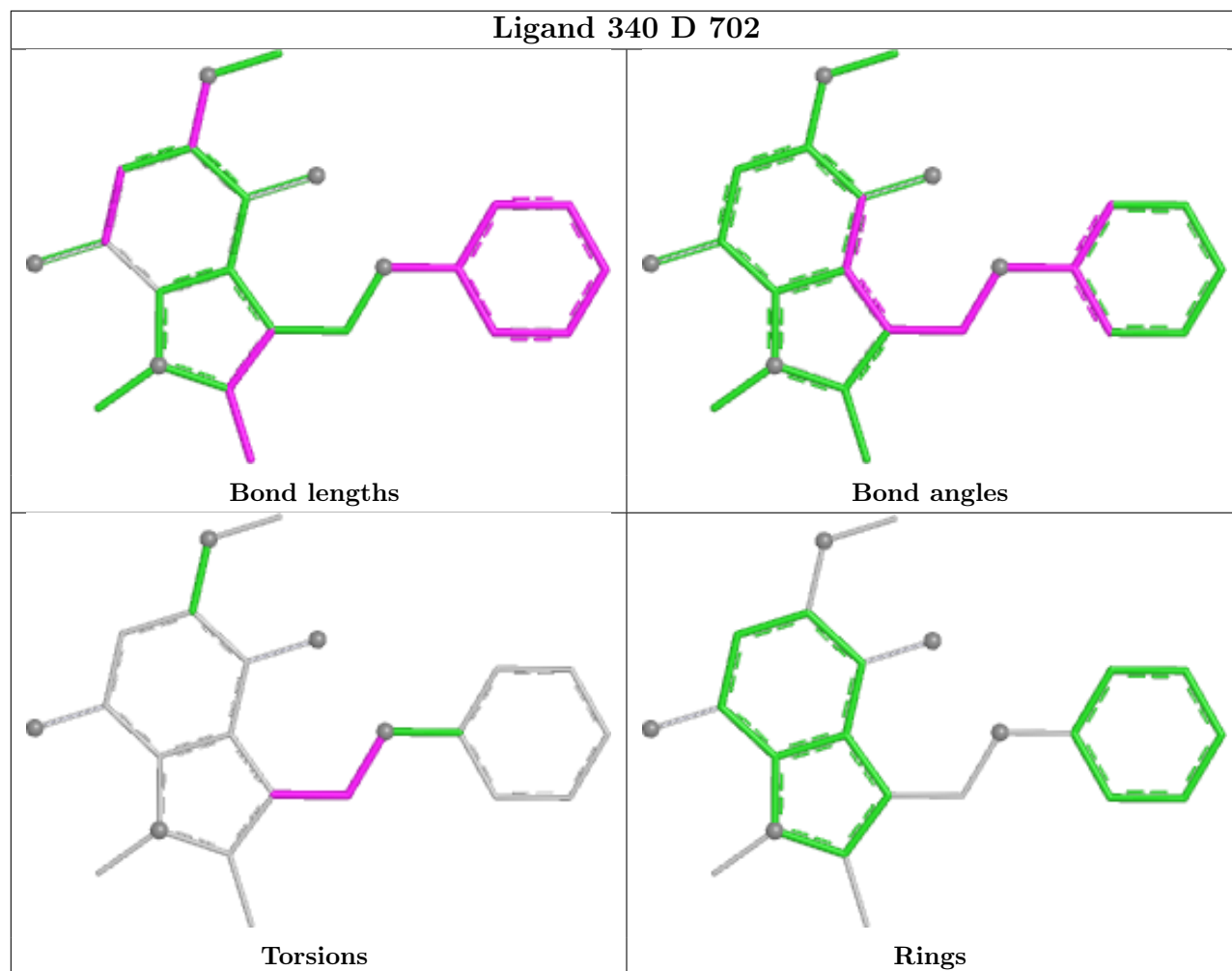
There are no ring outliers.

5 monomers are involved in 19 short contacts:

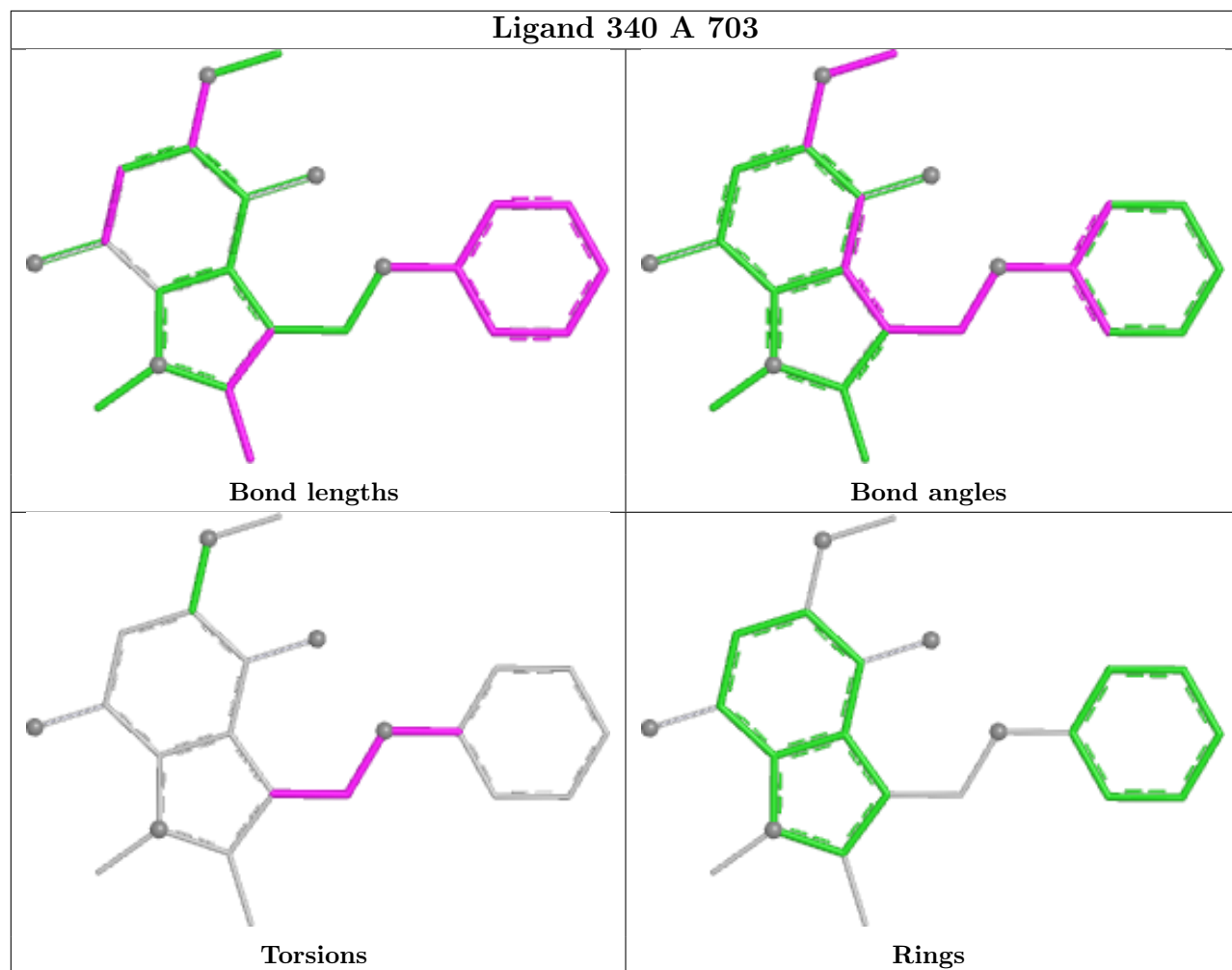
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	702	340	1	0
3	C	603	FAD	4	0
3	D	604	FAD	4	0
3	A	601	FAD	5	0
3	B	602	FAD	5	0

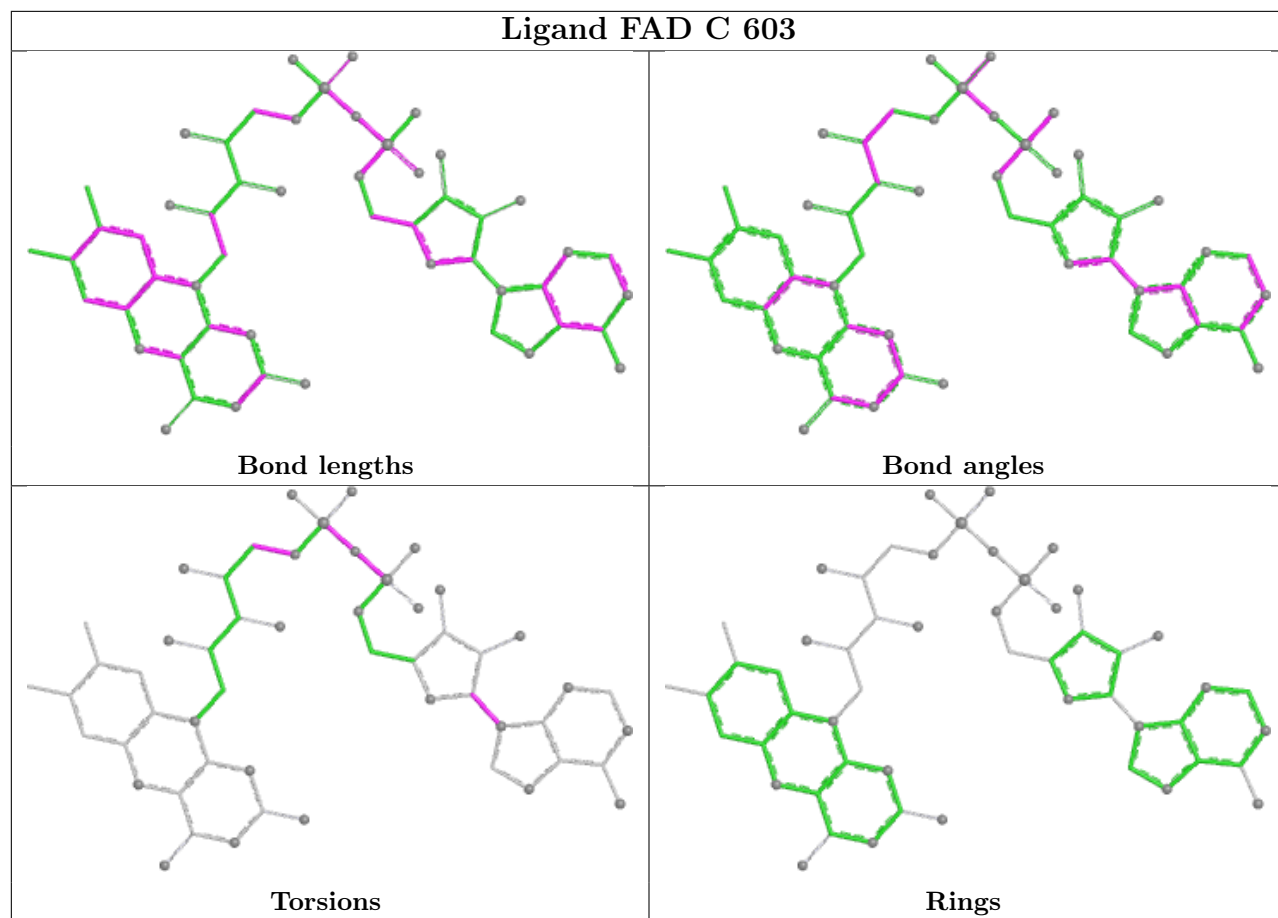
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

Ligand 340 D 702

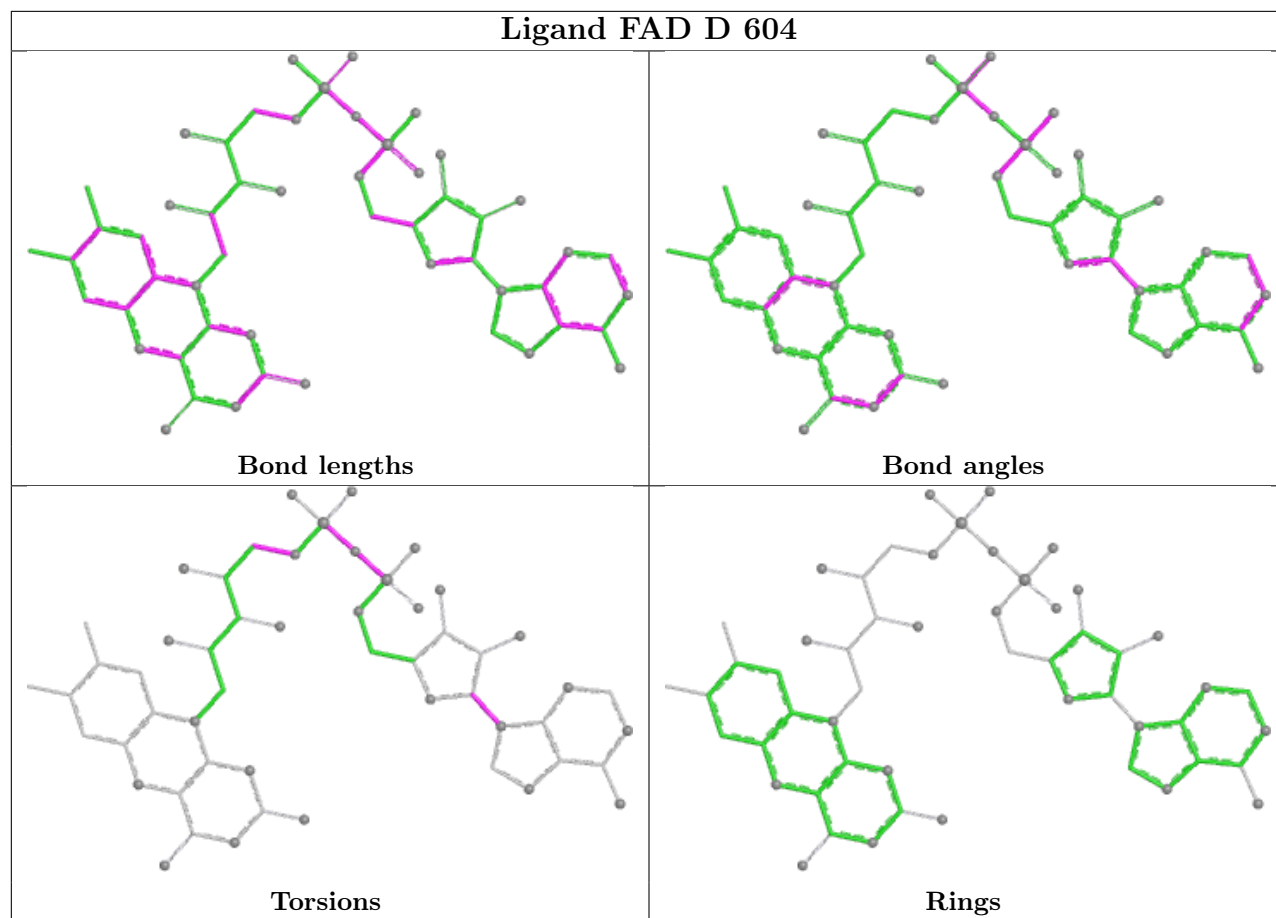


Ligand 340 A 703

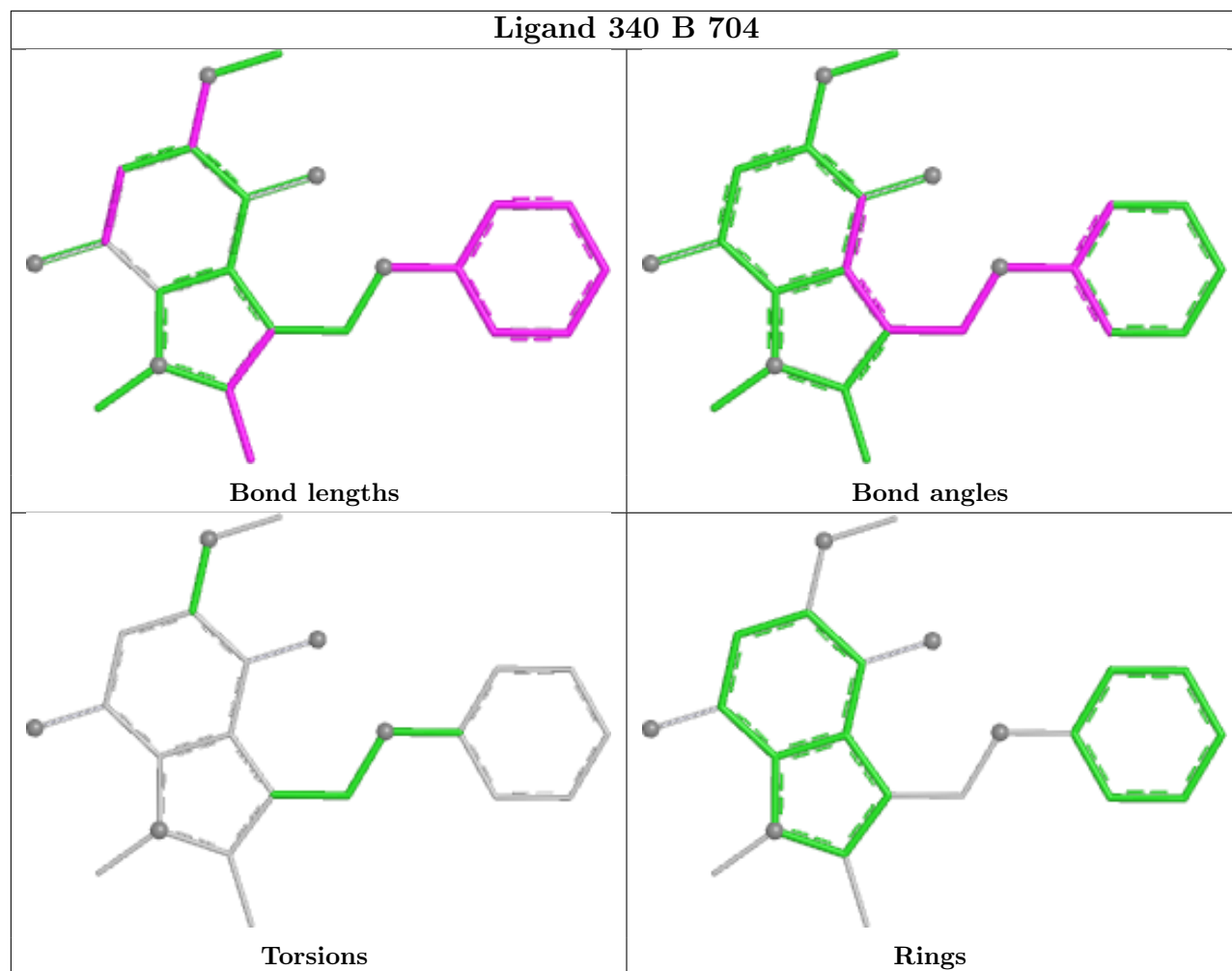


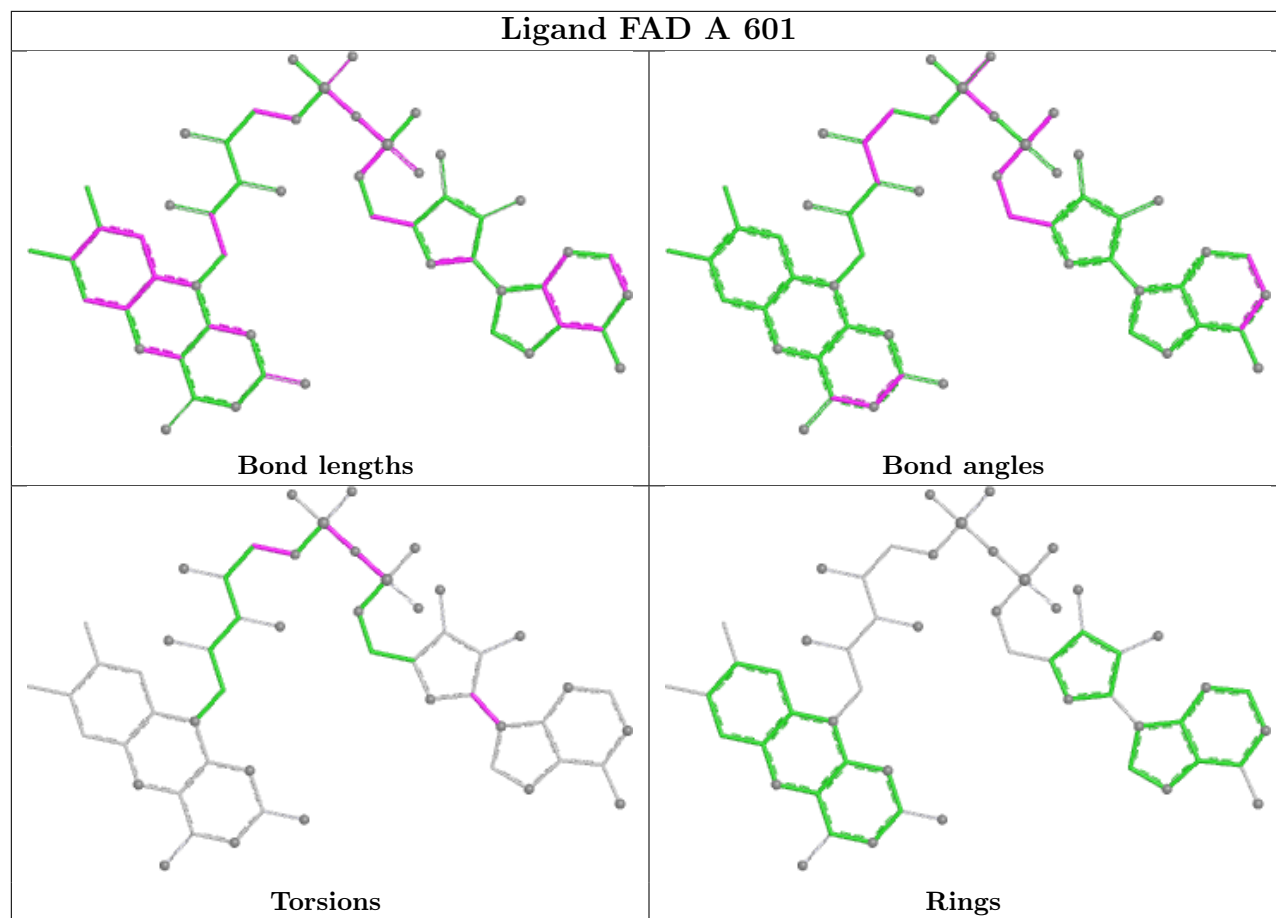


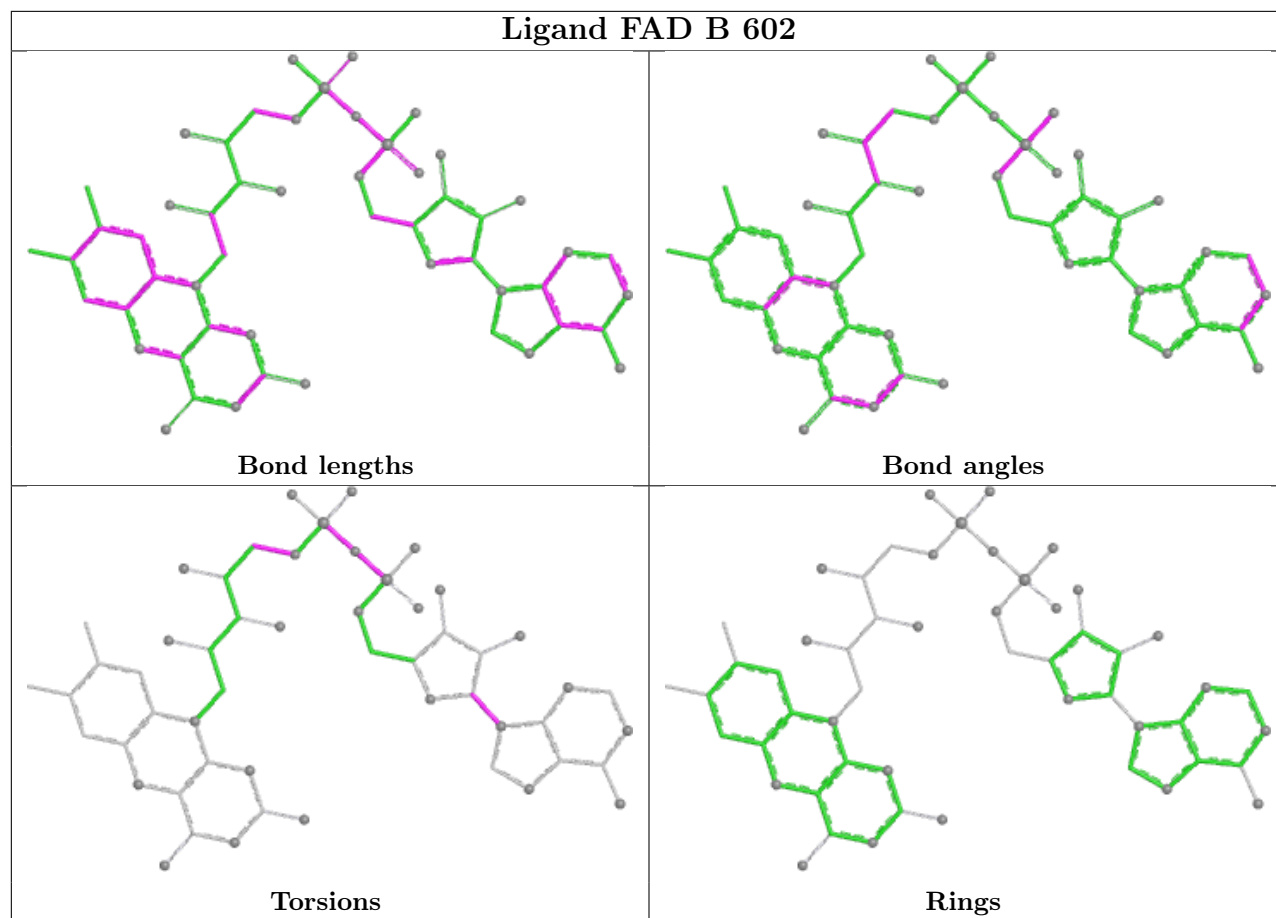
Ligand FAD D 604

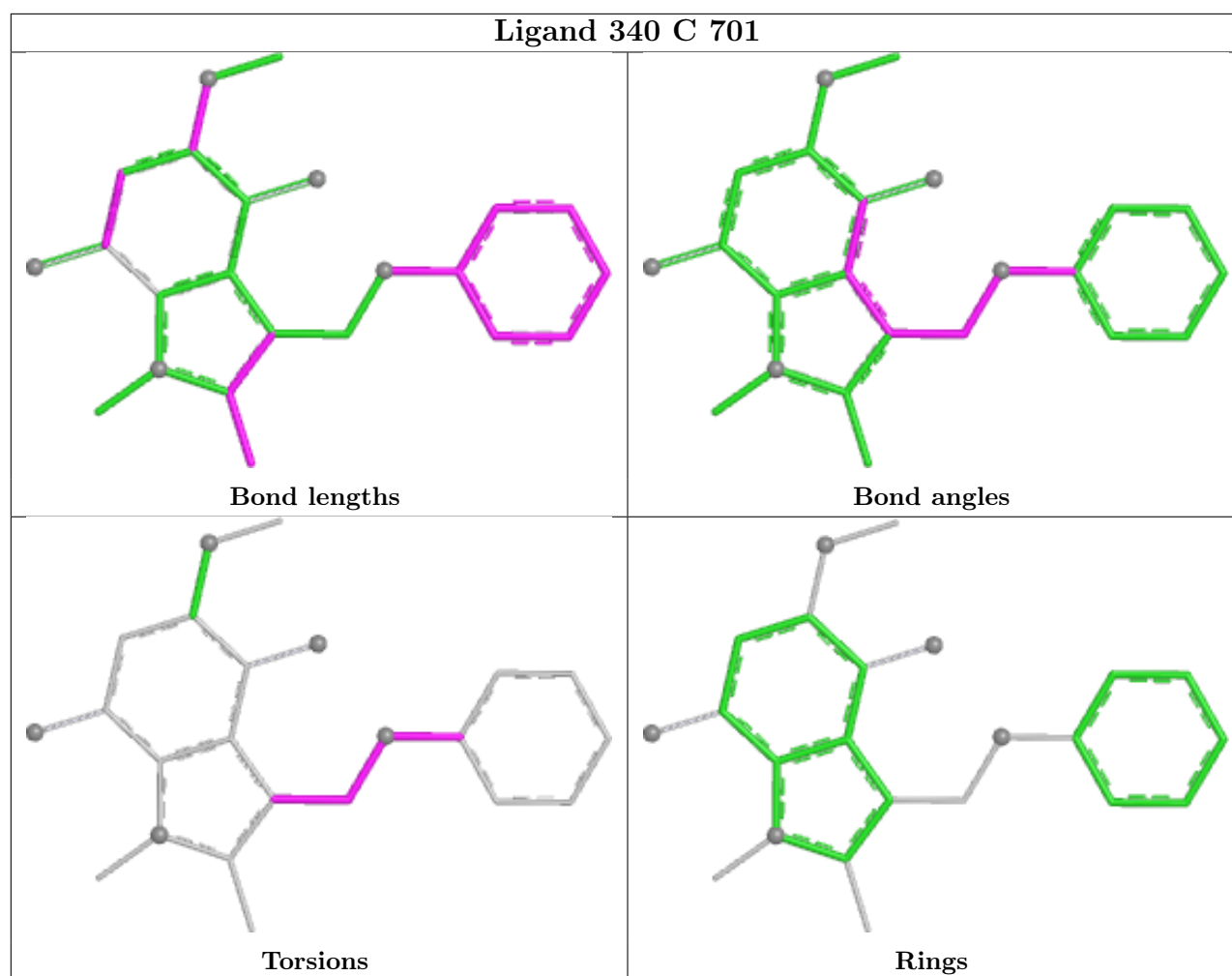


Ligand 340 B 704









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.