



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 06:53 AM UTC

PDB ID : 1DXO / pdb_00001dxo
Title : Crystal structure of human NAD[P]H-QUINONE oxidoreductase CO with 2, 3,5,6,tetramethyl-P-benzoquinone (duroquinone) at 2.5 Angstrom resolution
Authors : Faig, M.; Bianchet, M.A.; Chen, S.; Winski, S.; Ross, D.; Amzel, L.M.
Deposited on : 2000-01-12
Resolution : 2.50 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

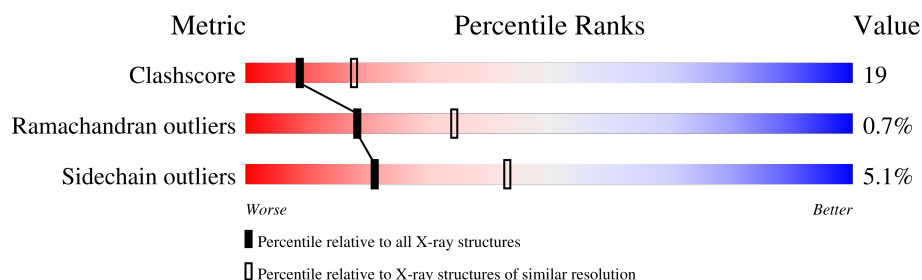
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)

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	 61% 36% •
1	B	273	 62% 33% 5%
1	C	273	 60% 37% •
1	D	273	 67% 29% •

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
3	DQN	B	302	-	-	X	-

2 Entry composition [i](#)

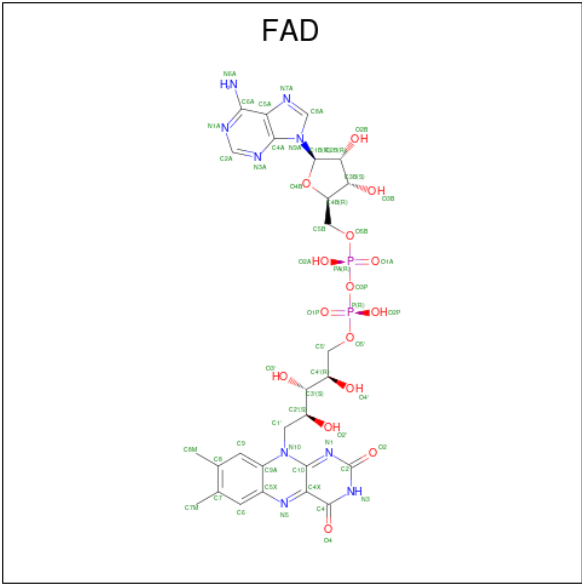
There are 4 unique types of molecules in this entry. The entry contains 9057 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 QUINONE REDUCTASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	273	Total	C	N	O	S	0	0	0
			2173	1412	365	389	7			
1	B	273	Total	C	N	O	S	0	0	0
			2173	1412	365	389	7			
1	C	273	Total	C	N	O	S	0	0	0
			2173	1412	365	389	7			
1	D	273	Total	C	N	O	S	0	0	0
			2173	1412	365	389	7			

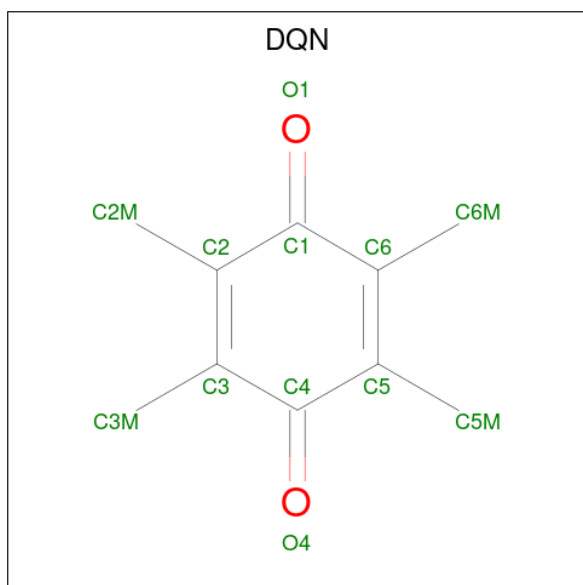
- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: C₂₇H₃₃N₉O₁₅P₂).



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

- Molecule 3 is DUROQUINONE (CCD ID: DQN) (formula: $C_{10}H_{12}O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			12	10	2		
3	B	1	Total	C	O	0	0
			12	10	2		
3	C	1	Total	C	O	0	0
			12	10	2		
3	D	1	Total	C	O	0	0
			12	10	2		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	22	Total	O	0	0
			22	22		
4	B	28	Total	O	0	0
			28	28		
4	C	23	Total	O	0	0
			23	23		

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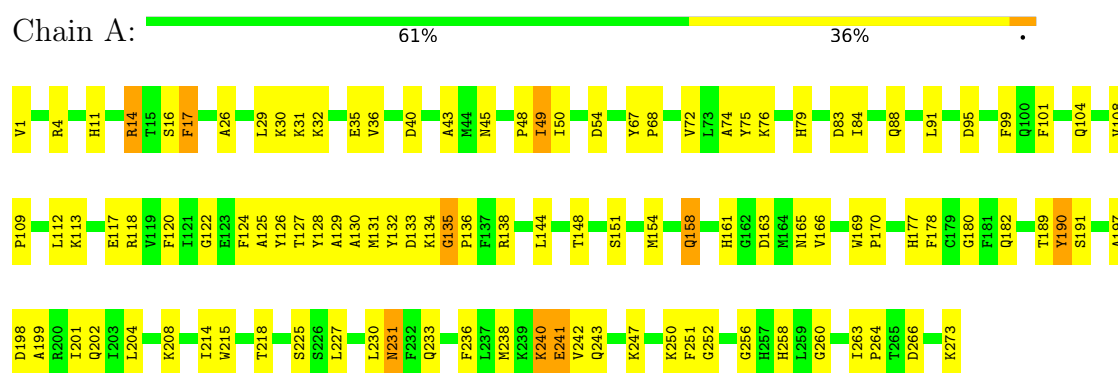
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	32	Total	O	0	0
			32	32		

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.

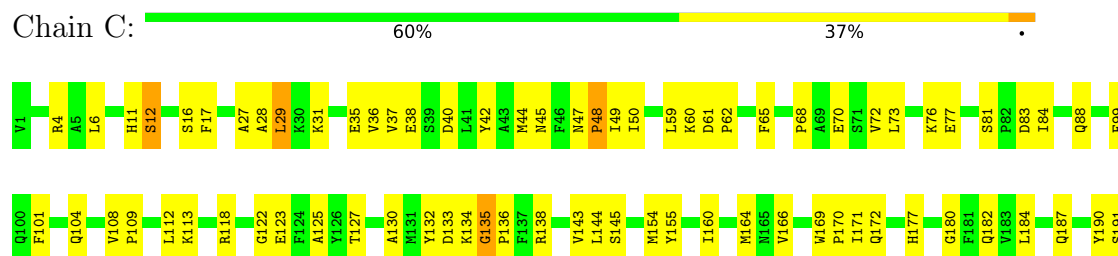
• Molecule 1: QUINONE REDUCTASE

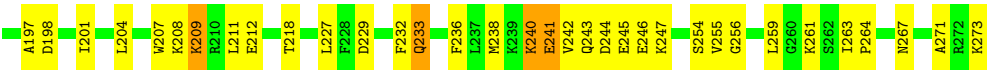


• Molecule 1: QUINONE REDUCTASE

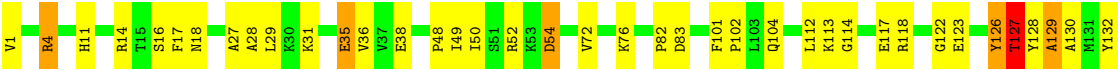


• Molecule 1: QUINONE REDUCTASE





● Molecule 1: QUINONE REDUCTASE



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	55.58Å 56.91Å 97.75Å 76.24° 76.73° 86.33°	Depositor
Resolution (Å)	27.66 – 2.50	Depositor
% Data completeness (in resolution range)	70.8 (27.66-2.50)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
Refinement program	CNS 0.9	Depositor
R, R_{free}	0.202 , 0.264	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	9057	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	0.54	0/2231	0.99	10/3012 (0.3%)
1	B	0.54	0/2231	1.00	8/3012 (0.3%)
1	C	0.56	0/2231	1.02	12/3012 (0.4%)
1	D	0.57	0/2231	1.02	11/3012 (0.4%)
All	All	0.55	0/8924	1.01	41/12048 (0.3%)

There are no bond length outliers.

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	182	GLN	N-CA-C	-7.43	96.62	108.73
1	B	182	GLN	N-CA-C	-7.42	97.16	108.96
1	B	49	ILE	N-CA-C	6.76	118.28	107.73
1	C	135	GLY	CA-C-N	6.63	126.57	119.28
1	C	135	GLY	C-N-CA	6.63	126.57	119.28
1	A	95	ASP	N-CA-C	-6.54	102.63	112.04
1	D	129	ALA	N-CA-C	-6.41	105.42	113.18
1	C	182	GLN	N-CA-C	-6.39	98.31	108.73
1	A	182	GLN	N-CA-C	-6.29	98.48	108.73
1	B	95	ASP	N-CA-C	-6.22	105.35	113.12
1	B	35	GLU	N-CA-C	-6.05	99.38	109.24
1	B	135	GLY	CA-C-N	6.04	126.21	119.32
1	B	135	GLY	C-N-CA	6.04	126.21	119.32
1	A	49	ILE	N-CA-C	6.00	117.08	107.73
1	C	101	PHE	N-CA-C	5.92	113.45	108.13
1	B	176	LEU	N-CA-C	5.85	117.35	110.97
1	A	36	VAL	N-CA-C	5.80	115.95	107.37
1	D	195	THR	CA-C-N	5.61	125.61	119.89
1	D	195	THR	C-N-CA	5.61	125.61	119.89
1	A	252	GLY	N-CA-C	-5.60	103.58	112.66

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	135	GLY	CA-C-N	5.59	125.43	119.28
1	A	135	GLY	C-N-CA	5.59	125.43	119.28
1	D	38	GLU	N-CA-C	5.52	118.57	109.85
1	D	135	GLY	CA-C-N	5.38	125.20	119.28
1	D	135	GLY	C-N-CA	5.38	125.20	119.28
1	D	35	GLU	N-CA-C	-5.26	100.67	109.24
1	D	36	VAL	N-CA-C	5.21	115.35	107.80
1	D	54	ASP	N-CA-C	-5.20	106.77	113.01
1	C	172	GLN	N-CA-C	5.16	116.90	111.28
1	C	36	VAL	N-CA-C	5.15	115.39	108.17
1	C	81	SER	CA-C-N	5.14	125.18	119.32
1	C	81	SER	C-N-CA	5.14	125.18	119.32
1	A	35	GLU	N-CA-C	-5.13	100.16	108.52
1	D	101	PHE	N-CA-C	5.12	112.83	108.22
1	A	190	TYR	N-CA-C	5.11	117.28	110.53
1	B	160	ILE	N-CA-C	5.09	116.42	110.62
1	C	38	GLU	N-CA-C	5.07	118.00	110.14
1	C	12	SER	N-CA-C	5.06	119.17	112.89
1	C	61	ASP	CA-C-N	5.04	124.83	119.28
1	C	61	ASP	C-N-CA	5.04	124.83	119.28
1	A	17	PHE	N-CA-C	-5.04	105.97	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	2173	0	2172	83	0
1	B	2173	0	2172	82	0
1	C	2173	0	2172	88	0
1	D	2173	0	2172	92	0
2	A	53	0	31	4	0
2	B	53	0	31	5	0
2	C	53	0	31	4	0
2	D	53	0	31	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	12	0	12	0	0
3	B	12	0	12	7	0
3	C	12	0	12	1	0
3	D	12	0	12	3	0
4	A	22	0	0	7	0
4	B	28	0	0	4	0
4	C	23	0	0	0	0
4	D	32	0	0	3	0
All	All	9057	0	8860	330	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 19.

All (330) 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:D:72:VAL:HG12	1:D:76:LYS:HE2	1.34	1.05
1:C:17:PHE:HB2	2:C:301:FAD:H52A	1.46	0.98
1:C:143:VAL:HG22	1:C:184:LEU:HB2	1.47	0.95
1:B:17:PHE:HB2	2:B:301:FAD:H52A	1.48	0.95
1:B:151:SER:H	1:B:154:MET:HE3	1.28	0.94
1:A:151:SER:H	1:A:154:MET:HE3	1.33	0.93
1:A:17:PHE:HB2	2:A:301:FAD:H52A	1.49	0.92
1:D:127:THR:OG1	1:D:130:ALA:HB3	1.68	0.91
1:C:233:GLN:HA	1:C:233:GLN:HE21	1.37	0.90
1:C:238:MET:HE3	1:C:243:GLN:HG2	1.53	0.89
1:D:17:PHE:HB2	2:D:301:FAD:H52A	1.52	0.88
3:B:302:DQN:H5M2	3:D:302:DQN:H2M3	1.55	0.85
1:A:48:PRO:HG3	1:C:49:ILE:HD11	1.62	0.81
1:D:126:TYR:CG	1:D:127:THR:N	2.50	0.80
1:A:50:ILE:HG22	1:A:118:ARG:HG2	1.65	0.79
1:B:48:PRO:HG3	1:D:49:ILE:HD11	1.64	0.78
1:B:60:LYS:HE3	1:B:73:LEU:HD21	1.67	0.76
1:C:169:TRP:HB3	1:C:170:PRO:HD3	1.68	0.75
1:A:14:ARG:HB2	1:A:14:ARG:HH11	1.53	0.74
1:D:130:ALA:HB1	1:D:134:LYS:O	1.88	0.74
3:B:302:DQN:H5M1	4:B:2020:HOH:O	1.88	0.73
1:D:255:VAL:HG23	1:D:267:ASN:HD22	1.54	0.73
1:A:166:VAL:HG13	1:C:166:VAL:HG13	1.72	0.72
1:B:255:VAL:HG23	1:B:267:ASN:HD22	1.56	0.71
1:D:127:THR:CB	1:D:130:ALA:HB3	2.20	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:49:ILE:HD11	1:C:48:PRO:HG3	1.73	0.71
1:A:158:GLN:HB3	1:C:238:MET:HE2	1.73	0.70
1:B:128:TYR:O	1:B:131:MET:HG3	1.91	0.69
1:B:72:VAL:O	1:B:76:LYS:HD3	1.92	0.69
1:B:4:ARG:HD2	1:B:35:GLU:OE1	1.93	0.68
1:C:138:ARG:HA	1:C:180:GLY:O	1.93	0.68
1:B:17:PHE:HB2	2:B:301:FAD:C5B	2.23	0.68
1:C:17:PHE:HB2	2:C:301:FAD:C5B	2.21	0.68
1:B:151:SER:N	1:B:154:MET:HE3	2.08	0.66
1:B:1:VAL:HA	1:B:215:TRP:NE1	2.10	0.66
1:D:28:ALA:HA	1:D:31:LYS:HE3	1.78	0.66
1:D:227:LEU:HB3	1:D:242:VAL:HG11	1.78	0.66
1:A:138:ARG:HA	1:A:180:GLY:O	1.96	0.65
1:B:160:ILE:HD12	1:D:230:LEU:CD2	2.26	0.65
1:D:14:ARG:HH11	1:D:14:ARG:HB2	1.60	0.65
1:D:83:ASP:OD2	1:D:118:ARG:NH2	2.28	0.65
1:A:130:ALA:HB1	1:A:134:LYS:O	1.97	0.65
1:C:99:PHE:HB2	1:C:144:LEU:HD23	1.78	0.65
1:C:122:GLY:O	1:C:123:GLU:HB2	1.96	0.65
1:B:169:TRP:CZ2	1:B:256:GLY:HA3	2.32	0.65
1:D:123:GLU:H	1:D:126:TYR:HD1	1.40	0.64
1:B:49:ILE:HD11	1:D:48:PRO:HG3	1.80	0.64
1:A:225:SER:HA	4:A:2019:HOH:O	1.97	0.64
1:B:14:ARG:HH11	1:B:14:ARG:CG	2.10	0.64
1:C:59:LEU:HB2	1:C:62:PRO:HG3	1.79	0.64
1:A:127:THR:HG22	1:A:129:ALA:H	1.62	0.64
1:B:60:LYS:CE	1:B:73:LEU:HD21	2.29	0.63
3:B:302:DQN:C5M	3:D:302:DQN:H2M3	2.28	0.63
1:A:240:LYS:H	1:A:240:LYS:HD2	1.63	0.63
1:B:235:GLY:O	1:B:237:LEU:HG	1.99	0.63
1:C:72:VAL:HG12	1:C:76:LYS:HE3	1.81	0.63
1:C:4:ARG:HH11	1:C:4:ARG:HG3	1.63	0.63
1:A:17:PHE:HB2	2:A:301:FAD:C5B	2.24	0.63
1:A:72:VAL:HG12	1:A:76:LYS:HE3	1.81	0.63
1:B:243:GLN:O	1:B:247:LYS:HG2	1.99	0.62
1:B:131:MET:HE3	1:D:160:ILE:HD13	1.81	0.62
1:C:204:LEU:O	1:C:208:LYS:HG3	2.00	0.62
1:C:50:ILE:HG22	1:C:118:ARG:HG2	1.80	0.62
1:B:54:ASP:OD1	1:B:118:ARG:NH1	2.33	0.62
1:D:17:PHE:HB2	2:D:301:FAD:C5B	2.26	0.62
1:B:50:ILE:HG22	1:B:118:ARG:HG2	1.81	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:14:ARG:HB2	1:D:14:ARG:NH1	2.15	0.61
1:C:255:VAL:HG23	1:C:267:ASN:HD22	1.64	0.61
1:A:83:ASP:OD2	1:A:118:ARG:NH2	2.31	0.61
1:A:151:SER:N	1:A:154:MET:HE3	2.08	0.61
1:C:47:ASN:ND2	1:C:49:ILE:H	1.99	0.61
3:B:302:DQN:H5M2	3:D:302:DQN:C2M	2.27	0.61
1:D:72:VAL:CG1	1:D:76:LYS:HE2	2.20	0.60
1:D:132:TYR:O	1:D:180:GLY:HA2	2.00	0.60
1:B:166:VAL:HG13	1:D:166:VAL:HG13	1.83	0.60
1:A:165:ASN:HD21	1:A:266:ASP:HA	1.66	0.60
1:D:27:ALA:O	1:D:31:LYS:HG3	2.00	0.60
1:B:138:ARG:HA	1:B:180:GLY:O	2.01	0.60
1:A:113:LYS:NZ	1:C:108:VAL:O	2.35	0.59
1:A:75:TYR:CE1	1:A:124:PHE:HB2	2.37	0.59
1:B:113:LYS:O	1:B:117:GLU:HG3	2.01	0.59
1:B:161:HIS:HE1	1:D:178:PHE:HE1	1.51	0.59
1:D:126:TYR:CD1	1:D:126:TYR:C	2.80	0.59
1:C:246:GLU:O	1:C:261:LYS:NZ	2.35	0.59
1:C:11:HIS:CE1	1:C:16:SER:HB3	2.38	0.59
1:D:197:ALA:O	1:D:201:ILE:HG13	2.03	0.58
1:D:122:GLY:O	1:D:123:GLU:HB2	2.01	0.58
1:A:1:VAL:HA	1:A:215:TRP:NE1	2.18	0.58
1:A:236:PHE:CD2	1:C:154:MET:HE3	2.38	0.58
1:C:127:THR:OG1	1:C:130:ALA:HB3	2.02	0.58
1:A:40:ASP:HB3	1:A:43:ALA:HB3	1.85	0.58
1:A:197:ALA:O	1:A:201:ILE:HG13	2.03	0.58
1:C:4:ARG:HD3	1:C:35:GLU:OE1	2.04	0.58
1:A:243:GLN:O	1:A:247:LYS:HG3	2.04	0.57
1:C:27:ALA:O	1:C:31:LYS:HG3	2.04	0.57
1:D:14:ARG:NH1	4:D:2005:HOH:O	2.37	0.57
1:D:1:VAL:HA	1:D:215:TRP:NE1	2.19	0.57
1:D:31:LYS:HZ1	1:D:208:LYS:HE2	1.69	0.57
1:D:155:TYR:HB3	1:D:164:MET:HB2	1.84	0.57
1:B:160:ILE:HD12	1:D:230:LEU:HD22	1.86	0.57
1:D:14:ARG:HH11	1:D:14:ARG:CB	2.17	0.57
1:A:101:PHE:CD2	1:A:109:PRO:HD3	2.40	0.57
1:D:126:TYR:HE2	1:D:128:TYR:HB2	1.68	0.57
1:D:123:GLU:N	1:D:126:TYR:CE1	2.70	0.56
1:B:147:THR:HG22	1:B:189:THR:OG1	2.06	0.56
1:D:122:GLY:HA2	1:D:126:TYR:CD1	2.41	0.56
1:B:83:ASP:OD2	1:B:118:ARG:NH2	2.38	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:40:ASP:O	1:C:44:MET:HG3	2.06	0.56
1:D:204:LEU:O	1:D:208:LYS:HG3	2.06	0.56
1:D:273:LYS:HD3	1:D:273:LYS:C	2.31	0.56
1:B:163:ASP:OD2	1:B:165:ASN:HB2	2.06	0.56
1:A:250:LYS:HD3	1:A:251:PHE:CE2	2.41	0.56
1:B:50:ILE:O	1:B:50:ILE:HG13	2.06	0.56
1:B:67:TYR:HB3	1:B:68:PRO:HD3	1.87	0.56
1:A:72:VAL:CG1	1:A:76:LYS:HE3	2.36	0.55
1:C:50:ILE:O	1:C:50:ILE:HG13	2.07	0.55
3:B:302:DQN:O1	1:D:126:TYR:CG	2.59	0.55
1:D:138:ARG:HA	1:D:180:GLY:O	2.07	0.55
1:B:230:LEU:CD2	1:D:160:ILE:HD12	2.36	0.55
1:B:132:TYR:HA	1:B:177:HIS:O	2.07	0.55
1:B:28:ALA:HA	1:B:31:LYS:HE3	1.89	0.54
1:A:32:LYS:HG2	4:A:2005:HOH:O	2.08	0.54
1:A:132:TYR:HA	1:A:177:HIS:O	2.08	0.54
1:D:169:TRP:CZ2	1:D:256:GLY:HA3	2.43	0.54
1:B:238:MET:HE3	1:B:243:GLN:HG2	1.89	0.54
1:A:54:ASP:OD1	1:A:118:ARG:NH1	2.39	0.54
1:A:227:LEU:HB3	1:A:242:VAL:HG11	1.90	0.54
1:C:143:VAL:CG2	1:C:184:LEU:HD12	2.38	0.54
1:D:126:TYR:CD2	1:D:127:THR:N	2.76	0.53
1:B:130:ALA:HB1	1:B:134:LYS:O	2.08	0.53
1:C:254:SER:HB2	1:C:267:ASN:HD21	1.73	0.53
1:D:112:LEU:HD12	1:D:112:LEU:O	2.08	0.53
1:D:127:THR:O	1:D:129:ALA:N	2.34	0.53
1:D:132:TYR:CD1	1:D:178:PHE:HA	2.44	0.53
1:B:238:MET:CE	1:B:243:GLN:HG2	2.38	0.53
1:C:72:VAL:CG1	1:C:76:LYS:HE3	2.39	0.53
1:D:127:THR:HG1	1:D:130:ALA:HB3	1.74	0.53
1:C:83:ASP:OD2	1:C:118:ARG:NH2	2.36	0.52
1:A:1:VAL:HA	1:A:215:TRP:CD1	2.45	0.52
1:C:209:LYS:HD3	1:C:209:LYS:O	2.09	0.52
1:A:108:VAL:O	1:C:113:LYS:NZ	2.42	0.52
1:D:132:TYR:HA	1:D:177:HIS:O	2.10	0.52
1:C:209:LYS:HE3	1:C:212:GLU:OE1	2.08	0.52
1:D:4:ARG:HD3	1:D:35:GLU:OE1	2.10	0.52
1:A:91:LEU:HD11	1:A:120:PHE:HE1	1.75	0.52
1:D:18:ASN:ND2	1:D:102:PRO:HG3	2.24	0.52
1:C:209:LYS:HD3	1:C:209:LYS:C	2.35	0.52
1:A:132:TYR:CD1	1:A:178:PHE:HA	2.45	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:148:THR:OG1	1:B:190:TYR:HA	2.11	0.51
1:C:130:ALA:HB1	1:C:134:LYS:O	2.11	0.51
1:A:166:VAL:O	1:A:170:PRO:HD3	2.11	0.51
1:A:99:PHE:HB2	1:A:144:LEU:HD23	1.92	0.51
1:C:227:LEU:HB3	1:C:242:VAL:HG11	1.91	0.51
1:D:127:THR:HB	1:D:130:ALA:HB3	1.93	0.51
1:D:190:TYR:O	1:D:191:SER:C	2.53	0.51
1:B:132:TYR:CD1	1:B:178:PHE:HA	2.46	0.51
1:C:4:ARG:HG3	1:C:4:ARG:NH1	2.22	0.51
1:A:31:LYS:HG3	4:A:2004:HOH:O	2.11	0.51
1:A:50:ILE:HG23	4:A:2007:HOH:O	2.11	0.50
1:B:263:ILE:HD11	1:D:263:ILE:HD11	1.94	0.50
1:D:126:TYR:HE2	1:D:128:TYR:CB	2.24	0.50
1:B:224:PRO:HD2	1:B:227:LEU:HD11	1.94	0.50
1:B:4:ARG:HB2	1:B:4:ARG:HH11	1.77	0.50
1:D:54:ASP:OD1	1:D:118:ARG:NH1	2.44	0.50
1:A:50:ILE:HG13	1:A:50:ILE:O	2.11	0.50
1:B:25:ALA:O	1:B:29:LEU:HB2	2.12	0.50
1:A:14:ARG:HH11	1:A:14:ARG:CB	2.23	0.49
1:A:258:HIS:HB3	1:A:263:ILE:HG12	1.94	0.49
1:A:131:MET:HE1	1:A:236:PHE:CD1	2.47	0.49
1:D:128:TYR:CE1	1:D:236:PHE:HZ	2.30	0.49
3:B:302:DQN:C5M	4:B:2020:HOH:O	2.53	0.49
1:D:130:ALA:O	1:D:135:GLY:HA2	2.12	0.49
1:D:200:ARG:HA	1:D:203:ILE:HD12	1.93	0.49
4:A:2019:HOH:O	1:C:160:ILE:CD1	2.61	0.49
1:C:255:VAL:H	1:C:267:ASN:ND2	2.11	0.49
1:C:241:GLU:O	1:C:245:GLU:HG3	2.13	0.49
1:C:12:SER:HB3	1:C:42:TYR:CD1	2.47	0.49
1:D:31:LYS:NZ	1:D:208:LYS:NZ	2.61	0.48
1:A:26:ALA:O	1:A:30:LYS:HG3	2.13	0.48
1:C:145:SER:HA	1:C:187:GLN:HB3	1.95	0.48
1:A:169:TRP:HB3	1:A:170:PRO:HD3	1.95	0.48
1:C:125:ALA:O	1:C:136:PRO:HD2	2.13	0.48
1:C:240:LYS:H	1:C:240:LYS:HD2	1.78	0.48
1:D:141:LYS:HG2	1:D:182:GLN:HB2	1.94	0.48
1:B:104:GLN:HA	2:B:301:FAD:C5X	2.44	0.48
1:C:65:PHE:CE1	1:C:70:GLU:HG3	2.49	0.48
1:B:161:HIS:HE1	1:D:178:PHE:CE1	2.30	0.48
1:B:1:VAL:HA	1:B:215:TRP:HE1	1.76	0.48
1:D:188:LEU:HD13	1:D:190:TYR:OH	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:233:GLN:HA	1:C:233:GLN:NE2	2.18	0.47
1:B:75:TYR:HB2	1:B:80:LEU:HD11	1.95	0.47
1:C:73:LEU:HD11	1:C:77:GLU:OE2	2.14	0.47
1:C:169:TRP:CZ2	1:C:256:GLY:HA3	2.49	0.47
1:C:197:ALA:O	1:C:201:ILE:HG13	2.14	0.47
1:D:31:LYS:HZ1	1:D:208:LYS:CE	2.26	0.47
1:D:113:LYS:O	1:D:117:GLU:HG3	2.14	0.47
1:C:47:ASN:ND2	1:C:47:ASN:C	2.72	0.47
1:D:239:LYS:HB2	1:D:242:VAL:HG23	1.96	0.47
1:A:68:PRO:O	1:A:72:VAL:HG23	2.15	0.47
1:A:263:ILE:HD11	1:C:263:ILE:HD11	1.96	0.47
1:B:14:ARG:CG	1:B:14:ARG:NH1	2.72	0.47
1:D:123:GLU:N	1:D:126:TYR:CD1	2.69	0.47
1:A:165:ASN:ND2	1:A:266:ASP:HA	2.29	0.47
1:C:238:MET:HE1	1:C:259:LEU:HD22	1.96	0.47
1:C:244:ASP:HA	1:C:247:LYS:HG3	1.97	0.46
1:D:27:ALA:C	1:D:31:LYS:HE2	2.40	0.46
1:D:126:TYR:CD2	1:D:128:TYR:N	2.83	0.46
1:A:31:LYS:HZ1	1:A:208:LYS:HE2	1.80	0.46
1:D:187:GLN:HG2	1:D:207:TRP:CE3	2.51	0.46
1:A:11:HIS:CE1	1:A:16:SER:HB3	2.50	0.46
1:A:154:MET:O	1:A:161:HIS:HB2	2.15	0.46
1:B:54:ASP:HA	1:B:81:SER:HB2	1.96	0.46
1:C:68:PRO:O	1:C:72:VAL:HG23	2.15	0.46
1:D:127:THR:C	1:D:129:ALA:H	2.21	0.46
1:D:145:SER:HA	1:D:187:GLN:HB3	1.98	0.46
1:A:238:MET:HE3	1:A:243:GLN:HG2	1.96	0.46
1:B:246:GLU:OE1	1:B:249:LYS:HG3	2.16	0.46
1:D:14:ARG:O	1:D:14:ARG:HG2	2.15	0.46
1:B:227:LEU:HB3	1:B:242:VAL:HG11	1.97	0.46
1:D:243:GLN:O	1:D:247:LYS:HG2	2.15	0.46
1:B:104:GLN:HA	2:B:301:FAD:N5	2.30	0.46
1:C:218:THR:HG23	1:C:271:ALA:HB3	1.96	0.46
1:B:14:ARG:HH11	1:B:14:ARG:HG2	1.79	0.46
1:B:190:TYR:O	1:B:191:SER:C	2.59	0.46
1:B:250:LYS:HD3	1:B:251:PHE:CE2	2.50	0.46
1:A:204:LEU:O	1:A:208:LYS:HG3	2.16	0.46
1:B:127:THR:OG1	1:B:130:ALA:HB3	2.15	0.46
1:C:243:GLN:O	1:C:247:LYS:HG3	2.16	0.46
1:D:11:HIS:CE1	1:D:16:SER:HB3	2.51	0.46
1:D:104:GLN:HA	2:D:301:FAD:C5X	2.46	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:198:ASP:OD1	1:A:198:ASP:N	2.43	0.45
1:C:132:TYR:HA	1:C:177:HIS:O	2.16	0.45
1:A:31:LYS:HZ1	1:A:208:LYS:CE	2.29	0.45
1:A:190:TYR:O	1:A:191:SER:C	2.59	0.45
1:B:31:LYS:HZ1	1:B:208:LYS:HE2	1.81	0.45
1:C:84:ILE:O	1:C:88:GLN:HG3	2.17	0.45
1:D:229:ASP:O	1:D:236:PHE:HA	2.16	0.45
1:A:125:ALA:O	1:A:136:PRO:HD2	2.16	0.45
1:A:263:ILE:O	1:A:264:PRO:C	2.59	0.45
1:B:60:LYS:NZ	1:B:73:LEU:HD21	2.32	0.45
1:A:258:HIS:CD2	1:A:260:GLY:H	2.35	0.45
1:B:189:THR:HG23	4:B:2022:HOH:O	2.17	0.45
1:C:47:ASN:C	1:C:47:ASN:HD22	2.24	0.45
1:B:1:VAL:HA	1:B:215:TRP:CD1	2.52	0.45
1:D:29:LEU:CD1	1:D:211:LEU:HB3	2.47	0.45
1:D:218:THR:O	1:D:218:THR:HG23	2.17	0.45
1:A:132:TYR:O	1:A:180:GLY:HA2	2.17	0.44
1:C:104:GLN:HA	2:C:301:FAD:C5X	2.47	0.44
1:C:238:MET:HE1	1:C:259:LEU:CD2	2.48	0.44
1:B:84:ILE:O	1:B:88:GLN:HG3	2.17	0.44
1:D:255:VAL:H	1:D:267:ASN:ND2	2.14	0.44
1:C:60:LYS:HE3	1:C:73:LEU:CD2	2.48	0.44
1:B:8:VAL:HG22	1:B:39:SER:HB2	2.00	0.44
1:A:122:GLY:HA2	1:A:126:TYR:CE1	2.53	0.43
1:A:241:GLU:CD	1:A:241:GLU:H	2.25	0.43
1:B:31:LYS:HG3	4:B:2004:HOH:O	2.17	0.43
1:B:50:ILE:CG2	1:B:118:ARG:HG2	2.47	0.43
1:C:6:LEU:HA	1:C:37:VAL:O	2.17	0.43
1:C:45:ASN:CG	1:C:45:ASN:O	2.60	0.43
4:A:2014:HOH:O	3:C:302:DQN:H5M1	2.18	0.43
1:A:54:ASP:CG	1:A:118:ARG:HH11	2.26	0.43
1:B:211:LEU:HD23	1:B:214:ILE:HD13	2.01	0.43
1:C:130:ALA:O	1:C:135:GLY:HA2	2.18	0.43
1:D:151:SER:H	1:D:154:MET:HE3	1.82	0.43
1:A:138:ARG:HG2	1:A:138:ARG:HH11	1.84	0.43
1:B:233:GLN:C	1:B:235:GLY:H	2.27	0.43
1:B:258:HIS:HB3	1:B:263:ILE:HG12	1.99	0.43
1:C:241:GLU:CD	1:C:241:GLU:H	2.26	0.43
1:D:28:ALA:HA	1:D:31:LYS:CE	2.47	0.43
1:C:28:ALA:HA	1:C:31:LYS:HE3	2.01	0.43
1:C:229:ASP:O	1:C:236:PHE:HA	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:130:ALA:O	1:A:135:GLY:HA2	2.19	0.43
1:C:47:ASN:HD22	1:C:48:PRO:N	2.17	0.43
1:B:125:ALA:O	1:B:179:CYS:HB3	2.19	0.43
1:D:123:GLU:N	1:D:126:TYR:HE1	2.17	0.43
1:D:192:ILE:O	1:D:192:ILE:HG13	2.18	0.43
1:A:169:TRP:CD1	1:C:166:VAL:HG11	2.54	0.43
1:A:169:TRP:CZ2	1:A:256:GLY:HA3	2.54	0.43
1:B:108:VAL:HB	1:B:112:LEU:HD13	2.01	0.43
1:C:263:ILE:O	1:C:264:PRO:C	2.60	0.43
1:D:48:PRO:HB2	1:D:114:GLY:HA3	2.01	0.43
1:D:126:TYR:CE2	1:D:128:TYR:CB	3.02	0.43
1:D:218:THR:HG23	1:D:271:ALA:HB3	2.00	0.43
1:D:52:ARG:HG2	4:D:2008:HOH:O	2.19	0.42
1:C:44:MET:O	1:C:45:ASN:HB3	2.19	0.42
1:C:104:GLN:HA	2:C:301:FAD:N5	2.34	0.42
1:C:232:PHE:O	1:C:233:GLN:C	2.63	0.42
1:D:50:ILE:O	1:D:50:ILE:HG13	2.18	0.42
1:A:68:PRO:HB2	4:A:2010:HOH:O	2.20	0.42
1:A:231:ASN:OD1	1:A:233:GLN:HB2	2.20	0.42
1:A:240:LYS:HD2	1:A:240:LYS:N	2.33	0.42
1:B:130:ALA:O	1:B:135:GLY:HA2	2.19	0.42
1:C:155:TYR:HB3	1:C:164:MET:HB2	2.01	0.42
1:B:154:MET:O	1:B:161:HIS:HB2	2.20	0.42
1:B:7:ILE:HG21	1:B:22:LYS:HG2	2.01	0.41
3:B:302:DQN:H3M1	4:D:2007:HOH:O	2.21	0.41
1:A:104:GLN:HA	2:A:301:FAD:C5X	2.51	0.41
1:B:106:PHE:CD1	1:B:106:PHE:N	2.89	0.41
1:C:144:LEU:HD11	1:C:171:ILE:CG2	2.50	0.41
1:B:176:LEU:O	1:B:181:PHE:HB2	2.19	0.41
1:A:104:GLN:HA	2:A:301:FAD:N5	2.35	0.41
1:A:117:GLU:OE2	1:C:109:PRO:HA	2.21	0.41
1:C:144:LEU:HD11	1:C:171:ILE:HG21	2.03	0.41
1:A:163:ASP:OD2	1:A:165:ASN:HB2	2.20	0.41
1:B:204:LEU:O	1:B:208:LYS:HG3	2.21	0.41
1:B:195:THR:HA	1:B:196:PRO:HD3	1.92	0.41
1:C:65:PHE:CD1	1:C:70:GLU:HG3	2.55	0.41
1:A:67:TYR:HB3	1:A:68:PRO:HD3	2.02	0.41
1:B:155:TYR:HB3	1:B:164:MET:HB2	2.02	0.41
1:C:12:SER:HB3	1:C:42:TYR:CE1	2.56	0.41
1:D:127:THR:C	1:D:129:ALA:N	2.77	0.41
1:A:84:ILE:O	1:A:88:GLN:HG3	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:199:ALA:O	1:A:202:GLN:N	2.54	0.41
1:B:11:HIS:NE2	2:B:301:FAD:O2P	2.52	0.41
1:B:132:TYR:O	1:B:134:LYS:N	2.54	0.41
1:C:112:LEU:HD12	1:C:112:LEU:O	2.21	0.41
1:A:1:VAL:HA	1:A:215:TRP:HE1	1.86	0.41
1:A:74:ALA:O	1:A:79:HIS:N	2.45	0.41
1:A:247:LYS:HE2	1:A:247:LYS:HB3	1.85	0.41
1:B:238:MET:HE3	1:B:243:GLN:HA	2.03	0.41
1:D:195:THR:HA	1:D:196:PRO:HD3	1.86	0.41
1:D:218:THR:HG23	1:D:271:ALA:CB	2.51	0.41
1:C:190:TYR:O	1:C:191:SER:C	2.64	0.40
1:D:127:THR:O	1:D:128:TYR:HB3	2.21	0.40
1:B:122:GLY:HA2	1:B:126:TYR:CE1	2.57	0.40
1:A:148:THR:OG1	1:A:190:TYR:HA	2.22	0.40
1:A:230:LEU:CD2	1:C:160:ILE:HD12	2.51	0.40
1:B:230:LEU:HD21	1:D:160:ILE:HD12	2.04	0.40
1:C:29:LEU:CD1	1:C:211:LEU:HB3	2.51	0.40
1:C:60:LYS:HE3	1:C:60:LYS:HB2	1.94	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%)	254 (94%)	16 (6%)	1 (0%)	30	49
1	B	271/273 (99%)	250 (92%)	17 (6%)	4 (2%)	8	16
1	C	271/273 (99%)	252 (93%)	18 (7%)	1 (0%)	30	49
1	D	271/273 (99%)	249 (92%)	20 (7%)	2 (1%)	18	34
All	All	1084/1092 (99%)	1005 (93%)	71 (6%)	8 (1%)	18	34

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	133	ASP
1	D	127	THR
1	A	133	ASP
1	B	234	ALA
1	D	191	SER
1	B	230	LEU
1	C	133	ASP
1	B	233	GLN

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	226/227 (100%)	212 (94%)	14 (6%)	16	34
1	B	226/227 (100%)	212 (94%)	14 (6%)	16	34
1	C	226/227 (100%)	217 (96%)	9 (4%)	28	54
1	D	226/227 (100%)	217 (96%)	9 (4%)	28	54
All	All	904/908 (100%)	858 (95%)	46 (5%)	21	43

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	ARG
1	A	14	ARG
1	A	29	LEU
1	A	45	ASN
1	A	112	LEU
1	A	128	TYR
1	A	158	GLN
1	A	189	THR
1	A	214	ILE
1	A	218	THR
1	A	231	ASN
1	A	240	LYS
1	A	241	GLU
1	A	273	LYS

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Mol	Chain	Res	Type
1	B	4	ARG
1	B	14	ARG
1	B	29	LEU
1	B	32	LYS
1	B	76	LYS
1	B	112	LEU
1	B	127	THR
1	B	143	VAL
1	B	158	GLN
1	B	198	ASP
1	B	205	GLU
1	B	214	ILE
1	B	233	GLN
1	B	273	LYS
1	C	29	LEU
1	C	48	PRO
1	C	198	ASP
1	C	207	TRP
1	C	209	LYS
1	C	233	GLN
1	C	240	LYS
1	C	241	GLU
1	C	273	LYS
1	D	4	ARG
1	D	82	PRO
1	D	126	TYR
1	D	127	THR
1	D	143	VAL
1	D	158	GLN
1	D	189	THR
1	D	198	ASP
1	D	273	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	45	ASN
1	A	104	GLN
1	A	158	GLN
1	A	268	GLN
1	B	104	GLN
1	B	158	GLN

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Mol	Chain	Res	Type
1	B	213	ASN
1	B	267	ASN
1	B	268	GLN
1	C	47	ASN
1	C	182	GLN
1	C	194	HIS
1	C	233	GLN
1	C	267	ASN
1	D	64	ASN
1	D	66	GLN
1	D	104	GLN
1	D	158	GLN
1	D	172	GLN
1	D	233	GLN
1	D	267	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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	FAD	B	301	-	58,58,58	2.55	21 (36%)	85,89,89	1.13	5 (5%)
2	FAD	D	301	-	58,58,58	2.58	21 (36%)	85,89,89	1.15	8 (9%)
3	DQN	B	302	-	12,12,12	1.29	1 (8%)	18,18,18	1.08	2 (11%)
2	FAD	A	301	-	58,58,58	2.44	21 (36%)	85,89,89	1.13	5 (5%)
3	DQN	D	302	-	12,12,12	1.11	0	18,18,18	0.67	0
3	DQN	C	302	-	12,12,12	1.06	0	18,18,18	0.66	0
3	DQN	A	302	-	12,12,12	1.14	0	18,18,18	0.58	0
2	FAD	C	301	-	58,58,58	2.60	22 (37%)	85,89,89	1.16	7 (8%)

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	FAD	B	301	-	-	5/34/50/50	0/6/6/6
2	FAD	D	301	-	-	5/34/50/50	0/6/6/6
3	DQN	B	302	-	-	-	0/1/1/1
2	FAD	A	301	-	-	5/34/50/50	0/6/6/6
3	DQN	D	302	-	-	-	0/1/1/1
3	DQN	C	302	-	-	-	0/1/1/1
3	DQN	A	302	-	-	-	0/1/1/1
2	FAD	C	301	-	-	5/34/50/50	0/6/6/6

All (86) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	301	FAD	P-O3P	-12.32	1.46	1.59
2	D	301	FAD	P-O3P	-11.95	1.46	1.59
2	C	301	FAD	P-O3P	-11.50	1.47	1.59
2	A	301	FAD	P-O3P	-10.59	1.48	1.59
2	A	301	FAD	PA-O3P	6.99	1.67	1.59
2	C	301	FAD	PA-O3P	5.68	1.65	1.59
2	C	301	FAD	C9A-N10	4.77	1.49	1.41
2	A	301	FAD	C9A-N10	4.46	1.48	1.41
2	B	301	FAD	PA-O2A	-4.45	1.34	1.55
2	D	301	FAD	C9A-N10	4.36	1.48	1.41
2	C	301	FAD	PA-O2A	-4.25	1.35	1.55
2	D	301	FAD	PA-O2A	-4.25	1.35	1.55
2	B	301	FAD	C9A-N10	4.15	1.48	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	301	FAD	PA-O3P	4.10	1.63	1.59
2	D	301	FAD	C1'-C2'	4.10	1.58	1.52
2	B	301	FAD	PA-O3P	4.10	1.63	1.59
2	C	301	FAD	C9A-C5X	4.08	1.47	1.41
2	A	301	FAD	PA-O2A	-4.03	1.36	1.55
2	C	301	FAD	C4X-N5	3.85	1.39	1.30
2	B	301	FAD	P-O2P	-3.80	1.37	1.55
2	D	301	FAD	C4X-N5	3.76	1.38	1.30
2	C	301	FAD	P-O2P	-3.70	1.38	1.55
2	D	301	FAD	P-O2P	-3.67	1.38	1.55
2	A	301	FAD	P-O2P	-3.57	1.38	1.55
2	B	301	FAD	O5'-C5'	3.53	1.58	1.44
2	C	301	FAD	O5'-C5'	3.37	1.57	1.44
2	B	301	FAD	C1'-C2'	3.31	1.57	1.52
2	B	301	FAD	C2A-N1A	3.24	1.39	1.33
2	B	301	FAD	C4X-N5	3.19	1.37	1.30
2	A	301	FAD	C8-C7	3.15	1.48	1.40
2	C	301	FAD	C9-C9A	3.14	1.44	1.39
2	D	301	FAD	O5'-C5'	3.11	1.56	1.44
2	A	301	FAD	C4X-N5	3.08	1.37	1.30
2	C	301	FAD	C8-C7	3.08	1.48	1.40
2	D	301	FAD	C9A-C5X	3.05	1.46	1.41
2	C	301	FAD	C4A-N3A	3.03	1.40	1.34
2	D	301	FAD	C5A-C4A	3.02	1.44	1.39
2	A	301	FAD	O5'-C5'	3.02	1.56	1.44
2	C	301	FAD	C2-N3	3.01	1.45	1.39
2	D	301	FAD	C5A-C6A	3.00	1.49	1.41
2	B	301	FAD	C10-N1	2.99	1.39	1.33
2	D	301	FAD	C9-C9A	2.96	1.44	1.39
2	B	301	FAD	C4A-N3A	2.90	1.39	1.34
2	B	301	FAD	C5A-C4A	2.87	1.44	1.39
2	B	301	FAD	PA-O5B	-2.86	1.48	1.59
2	C	301	FAD	PA-O5B	-2.84	1.48	1.59
2	D	301	FAD	O4B-C1B	2.83	1.48	1.42
2	D	301	FAD	C8-C7	2.81	1.47	1.40
2	D	301	FAD	C2-N3	2.80	1.45	1.39
2	C	301	FAD	C5A-C6A	2.76	1.48	1.41
2	C	301	FAD	O4B-C1B	2.73	1.48	1.42
2	B	301	FAD	C5A-C6A	2.70	1.48	1.41
2	B	301	FAD	C6-C5X	2.69	1.44	1.40
2	A	301	FAD	C10-N1	2.66	1.38	1.33
2	D	301	FAD	C10-N1	2.65	1.38	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	301	FAD	O4B-C1B	2.63	1.48	1.42
2	A	301	FAD	C5A-C6A	2.59	1.48	1.41
2	A	301	FAD	C9A-C5X	2.59	1.45	1.41
2	D	301	FAD	C2A-N1A	2.58	1.38	1.33
3	B	302	DQN	C6-C1	2.57	1.56	1.47
2	C	301	FAD	C10-N1	2.55	1.38	1.33
2	B	301	FAD	C8-C7	2.51	1.47	1.40
2	A	301	FAD	PA-O5B	-2.50	1.49	1.59
2	B	301	FAD	C2-N3	2.49	1.44	1.39
2	A	301	FAD	C4A-N3A	2.46	1.39	1.34
2	D	301	FAD	PA-O5B	-2.44	1.49	1.59
2	C	301	FAD	C5B-C4B	2.43	1.58	1.51
2	D	301	FAD	C6-C5X	2.41	1.43	1.40
2	A	301	FAD	C2-N3	2.37	1.44	1.39
2	A	301	FAD	C5B-C4B	2.34	1.58	1.51
2	C	301	FAD	C1'-C2'	2.34	1.55	1.52
2	A	301	FAD	C9-C9A	2.34	1.43	1.39
2	C	301	FAD	C4X-C10	2.31	1.51	1.44
2	B	301	FAD	O4B-C1B	2.23	1.47	1.42
2	B	301	FAD	C9-C9A	2.19	1.43	1.39
2	B	301	FAD	C9A-C5X	2.19	1.44	1.41
2	A	301	FAD	O2-C2	-2.18	1.20	1.24
2	A	301	FAD	C5A-C4A	2.17	1.43	1.39
2	B	301	FAD	C5B-C4B	2.17	1.58	1.51
2	C	301	FAD	C9-C8	2.14	1.42	1.39
2	D	301	FAD	C4A-N3A	2.13	1.38	1.34
2	A	301	FAD	C6-C5X	2.12	1.43	1.40
2	C	301	FAD	C2A-N1A	2.11	1.37	1.33
2	A	301	FAD	C2A-N1A	2.07	1.37	1.33
2	D	301	FAD	C3B-C4B	2.03	1.58	1.53
2	C	301	FAD	C6-C5X	2.00	1.43	1.40

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	301	FAD	O5B-PA-O1A	-3.25	96.04	108.94
2	A	301	FAD	O5B-PA-O1A	-3.23	96.13	108.94
2	D	301	FAD	O5B-PA-O1A	-3.21	96.23	108.94
2	B	301	FAD	O5B-PA-O1A	-3.11	96.61	108.94
2	C	301	FAD	O4B-C1B-N9A	-2.70	102.90	108.09
2	A	301	FAD	C2A-N1A-C6A	2.64	123.07	118.73
2	C	301	FAD	C2A-N1A-C6A	2.62	123.04	118.73

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	301	FAD	O4B-C1B-N9A	-2.58	103.13	108.09
2	B	301	FAD	O4B-C1B-N9A	-2.56	103.18	108.09
2	D	301	FAD	C4-N3-C2	-2.47	121.25	125.64
2	A	301	FAD	C4-N3-C2	-2.40	121.37	125.64
2	D	301	FAD	C2A-N1A-C6A	2.38	122.64	118.73
2	B	301	FAD	C2A-N1A-C6A	2.33	122.56	118.73
2	C	301	FAD	C5'-C4'-C3'	-2.32	107.83	112.22
2	A	301	FAD	C5'-C4'-C3'	-2.30	107.87	112.22
2	B	301	FAD	C4-N3-C2	-2.27	121.61	125.64
2	D	301	FAD	O2P-P-O3P	-2.25	101.18	107.27
2	C	301	FAD	C5X-C9A-N10	-2.25	115.94	117.97
2	D	301	FAD	C5'-C4'-C3'	-2.24	107.99	112.22
2	D	301	FAD	C5X-C9A-N10	-2.20	115.98	117.97
3	B	302	DQN	O1-C1-C6	2.17	126.79	120.73
2	D	301	FAD	C10-N1-C2	2.10	121.39	116.85
2	C	301	FAD	C4-N3-C2	-2.10	121.92	125.64
2	C	301	FAD	O2P-P-O3P	-2.05	101.74	107.27
2	A	301	FAD	C5X-C9A-N10	-2.01	116.15	117.97
3	B	302	DQN	O1-C1-C2	-2.01	115.11	120.73
2	B	301	FAD	C5X-C9A-N10	-2.00	116.16	117.97

There are no chirality outliers.

All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	301	FAD	C4'-C5'-O5'-P
2	B	301	FAD	C4'-C5'-O5'-P
2	C	301	FAD	C4'-C5'-O5'-P
2	D	301	FAD	C4'-C5'-O5'-P
2	A	301	FAD	P-O3P-PA-O5B
2	B	301	FAD	P-O3P-PA-O5B
2	C	301	FAD	P-O3P-PA-O5B
2	D	301	FAD	P-O3P-PA-O5B
2	A	301	FAD	C2B-C1B-N9A-C8A
2	C	301	FAD	C2B-C1B-N9A-C8A
2	D	301	FAD	C2B-C1B-N9A-C8A
2	B	301	FAD	C2B-C1B-N9A-C8A
2	A	301	FAD	PA-O3P-P-O5'
2	B	301	FAD	PA-O3P-P-O5'
2	C	301	FAD	PA-O3P-P-O5'
2	D	301	FAD	PA-O3P-P-O5'
2	D	301	FAD	O4B-C1B-N9A-C8A

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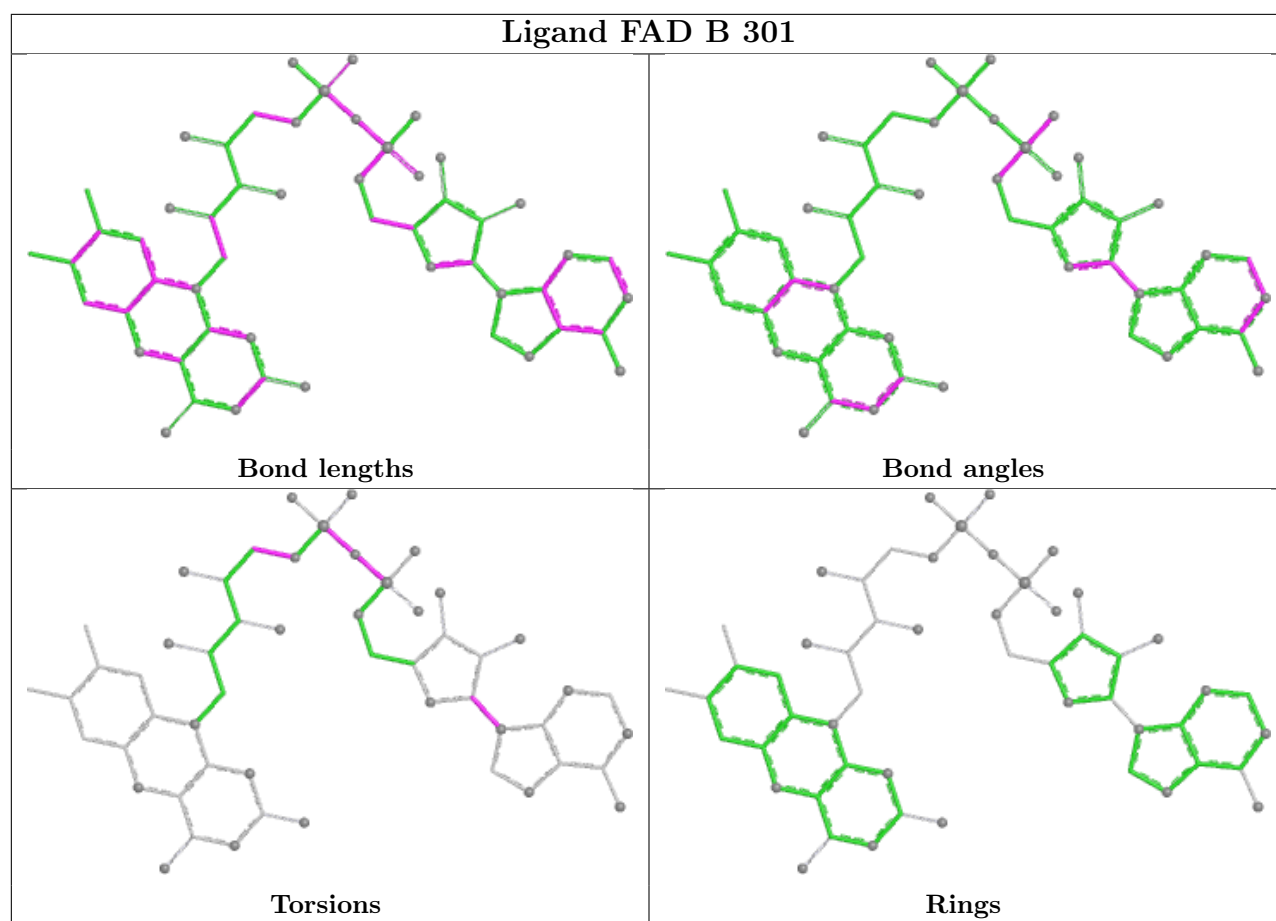
Mol	Chain	Res	Type	Atoms
2	A	301	FAD	O4B-C1B-N9A-C8A
2	B	301	FAD	O4B-C1B-N9A-C8A
2	C	301	FAD	O4B-C1B-N9A-C8A

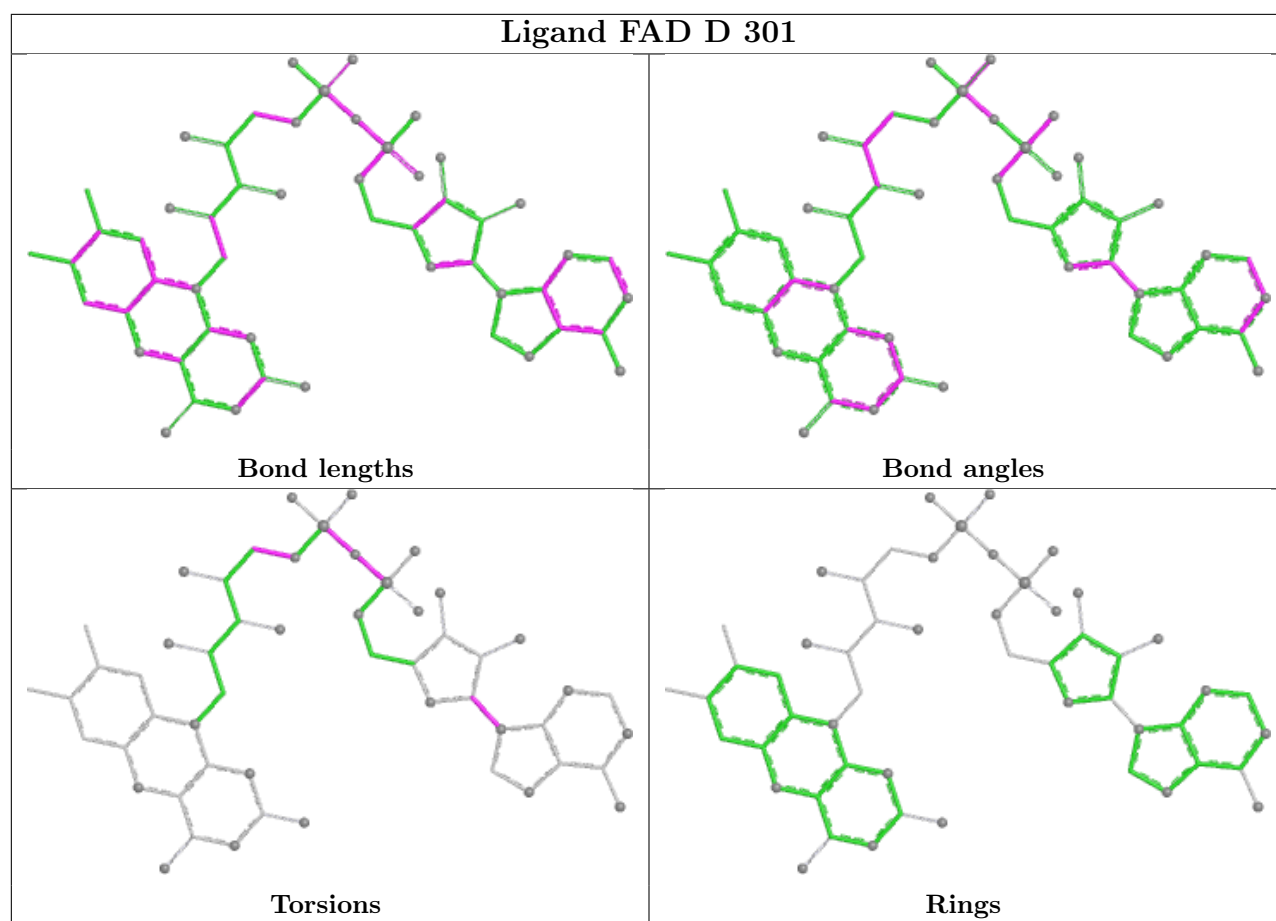
There are no ring outliers.

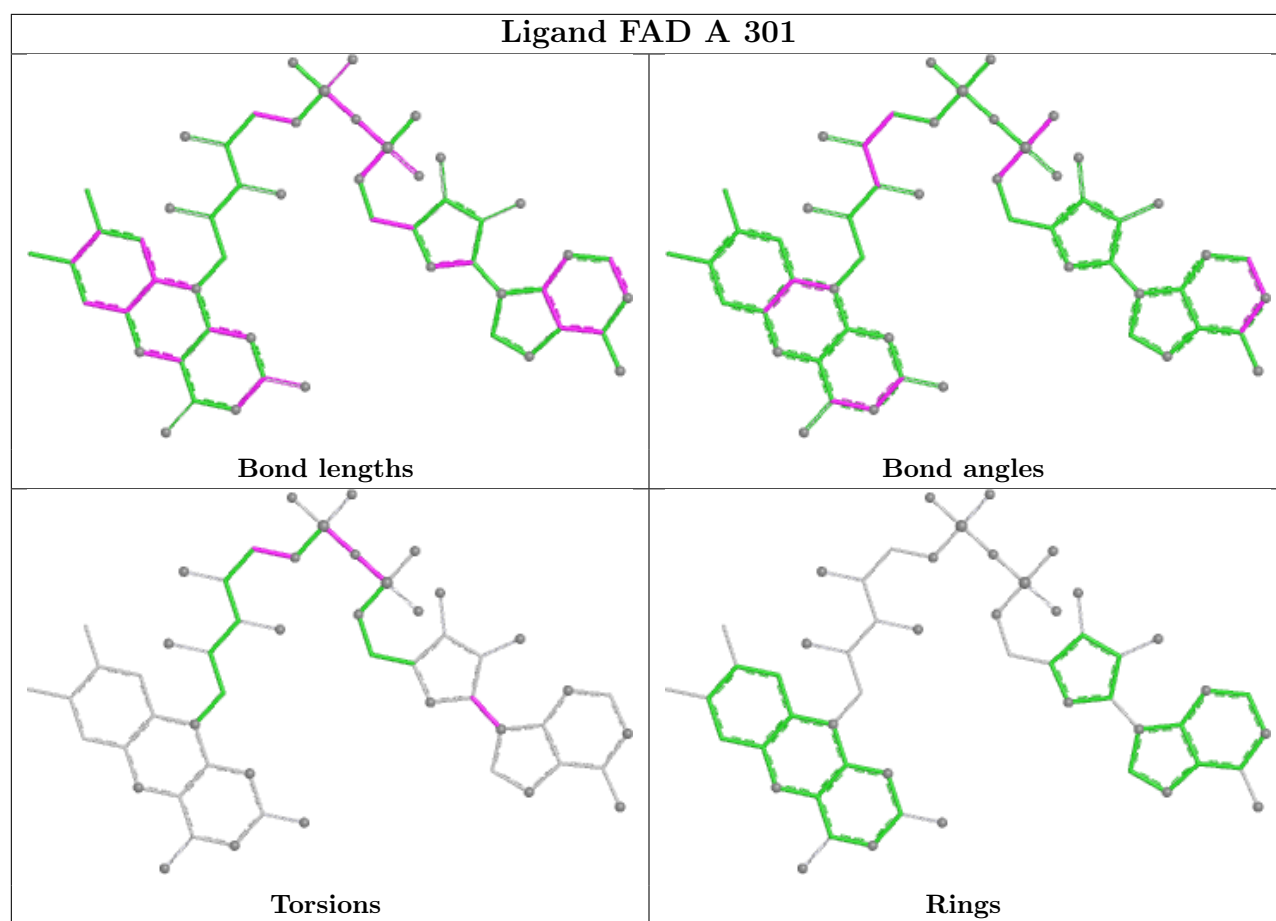
7 monomers are involved in 24 short contacts:

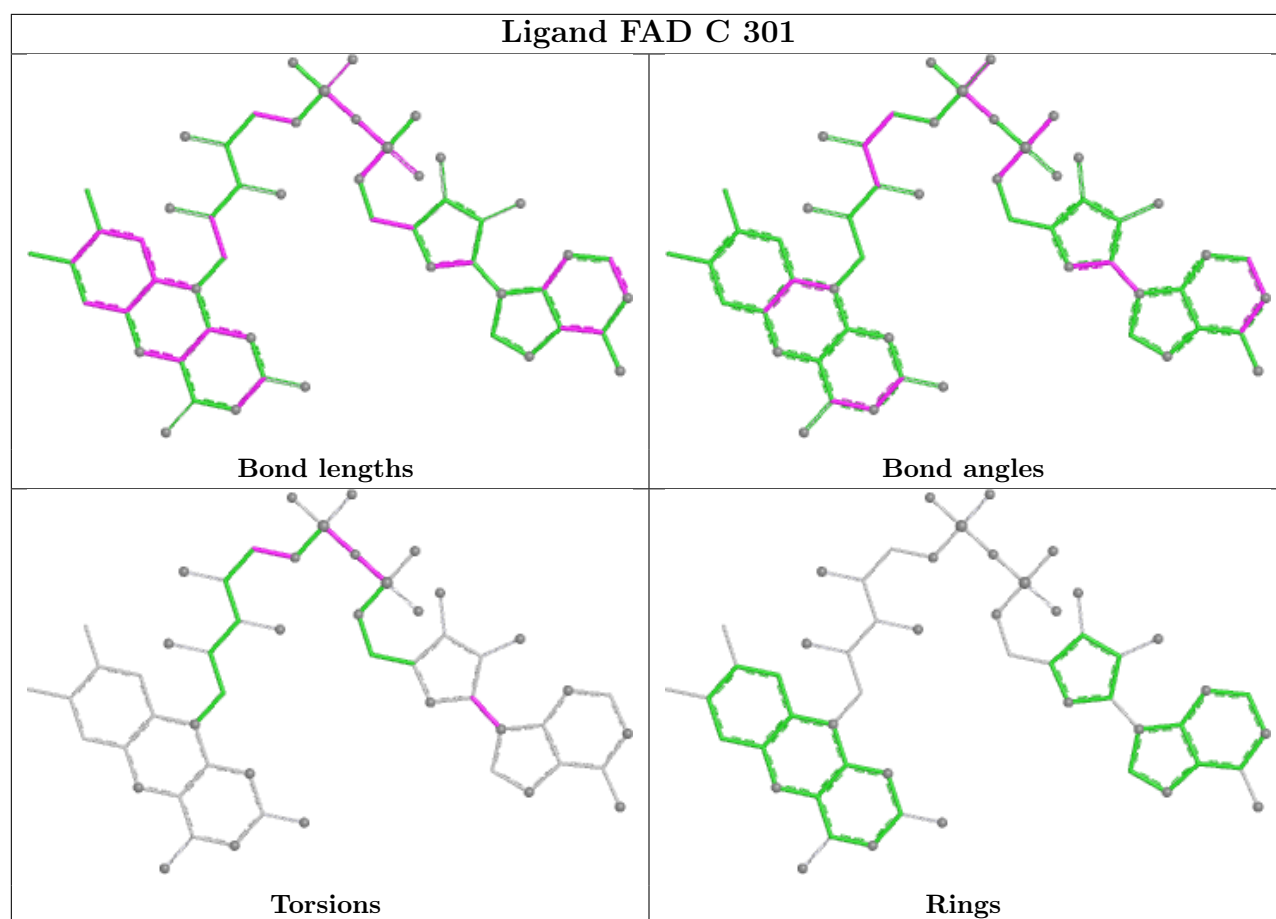
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	301	FAD	5	0
2	D	301	FAD	3	0
3	B	302	DQN	7	0
2	A	301	FAD	4	0
3	D	302	DQN	3	0
3	C	302	DQN	1	0
2	C	301	FAD	4	0

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.