



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 01:21 AM UTC

PDB ID : 8QNS / pdb_00008qns
Title : Crystal structure of murine AIF bound to N-terminal domain of CHCHD4
Authors : Fagnani, E.; Milani, M.
Deposited on : 2023-09-27
Resolution : 3.21 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) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

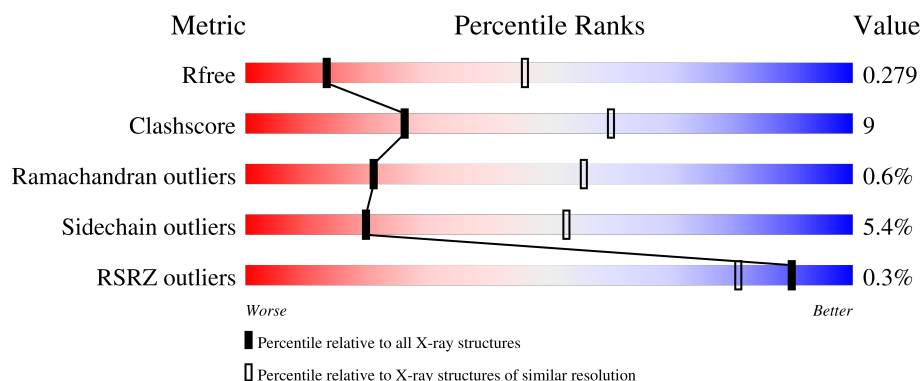
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.21 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))
R_{free}	180053	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

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 $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	524	
1	D	524	
1	G	524	
1	J	524	
2	M	27	

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Mol	Chain	Length	Quality of chain
2	N	27	37%19%15%.26%
2	O	27	33%22%11%33%
2	P	27	26%15%59%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 14701 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 Apoptosis-inducing factor 1, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	449	Total	C	N	O	S	0	0	0
			3472	2207	619	635	11			
1	D	439	Total	C	N	O	S	0	0	0
			3400	2162	605	622	11			
1	G	440	Total	C	N	O	S	0	0	0
			3400	2162	604	623	11			
1	J	437	Total	C	N	O	S	0	0	0
			3382	2152	601	618	11			

There are 52 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	101	MET	-	initiating methionine	UNP Q9Z0X1
A	613	LEU	-	expression tag	UNP Q9Z0X1
A	614	VAL	-	expression tag	UNP Q9Z0X1
A	615	PRO	-	expression tag	UNP Q9Z0X1
A	616	ARG	-	expression tag	UNP Q9Z0X1
A	617	GLY	-	expression tag	UNP Q9Z0X1
A	618	SER	-	expression tag	UNP Q9Z0X1
A	619	HIS	-	expression tag	UNP Q9Z0X1
A	620	HIS	-	expression tag	UNP Q9Z0X1
A	621	HIS	-	expression tag	UNP Q9Z0X1
A	622	HIS	-	expression tag	UNP Q9Z0X1
A	623	HIS	-	expression tag	UNP Q9Z0X1
A	624	HIS	-	expression tag	UNP Q9Z0X1
D	101	MET	-	initiating methionine	UNP Q9Z0X1
D	613	LEU	-	expression tag	UNP Q9Z0X1
D	614	VAL	-	expression tag	UNP Q9Z0X1
D	615	PRO	-	expression tag	UNP Q9Z0X1
D	616	ARG	-	expression tag	UNP Q9Z0X1
D	617	GLY	-	expression tag	UNP Q9Z0X1
D	618	SER	-	expression tag	UNP Q9Z0X1
D	619	HIS	-	expression tag	UNP Q9Z0X1

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Chain	Residue	Modelled	Actual	Comment	Reference
D	620	HIS	-	expression tag	UNP Q9Z0X1
D	621	HIS	-	expression tag	UNP Q9Z0X1
D	622	HIS	-	expression tag	UNP Q9Z0X1
D	623	HIS	-	expression tag	UNP Q9Z0X1
D	624	HIS	-	expression tag	UNP Q9Z0X1
G	101	MET	-	initiating methionine	UNP Q9Z0X1
G	613	LEU	-	expression tag	UNP Q9Z0X1
G	614	VAL	-	expression tag	UNP Q9Z0X1
G	615	PRO	-	expression tag	UNP Q9Z0X1
G	616	ARG	-	expression tag	UNP Q9Z0X1
G	617	GLY	-	expression tag	UNP Q9Z0X1
G	618	SER	-	expression tag	UNP Q9Z0X1
G	619	HIS	-	expression tag	UNP Q9Z0X1
G	620	HIS	-	expression tag	UNP Q9Z0X1
G	621	HIS	-	expression tag	UNP Q9Z0X1
G	622	HIS	-	expression tag	UNP Q9Z0X1
G	623	HIS	-	expression tag	UNP Q9Z0X1
G	624	HIS	-	expression tag	UNP Q9Z0X1
J	101	MET	-	initiating methionine	UNP Q9Z0X1
J	613	LEU	-	expression tag	UNP Q9Z0X1
J	614	VAL	-	expression tag	UNP Q9Z0X1
J	615	PRO	-	expression tag	UNP Q9Z0X1
J	616	ARG	-	expression tag	UNP Q9Z0X1
J	617	GLY	-	expression tag	UNP Q9Z0X1
J	618	SER	-	expression tag	UNP Q9Z0X1
J	619	HIS	-	expression tag	UNP Q9Z0X1
J	620	HIS	-	expression tag	UNP Q9Z0X1
J	621	HIS	-	expression tag	UNP Q9Z0X1
J	622	HIS	-	expression tag	UNP Q9Z0X1
J	623	HIS	-	expression tag	UNP Q9Z0X1
J	624	HIS	-	expression tag	UNP Q9Z0X1

- Molecule 2 is a protein called Mitochondrial intermembrane space import and assembly protein 40.

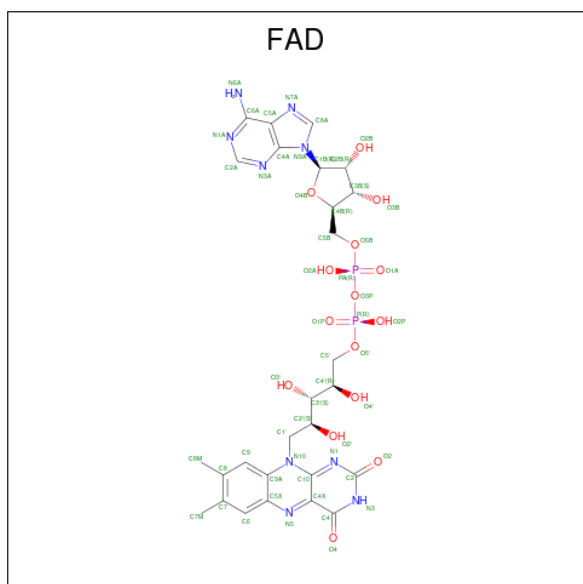
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	M	18	Total	C	N	O	S	0	0	0
			152	94	27	30	1			
2	N	20	Total	C	N	O	S	0	0	0
			171	105	31	34	1			
2	O	18	Total	C	N	O	S	0	0	0
			156	97	29	29	1			

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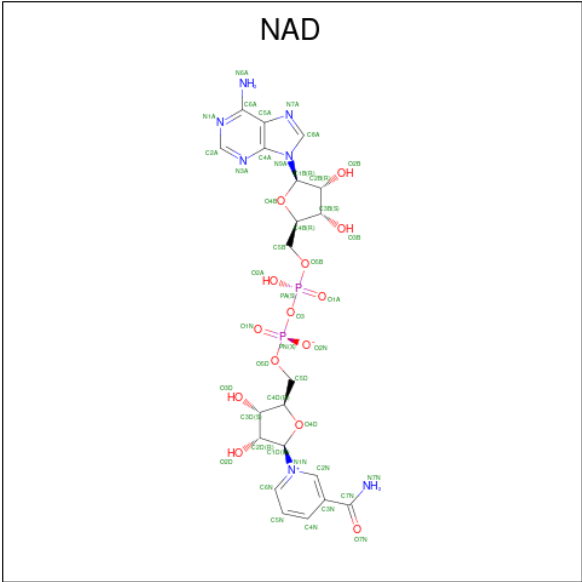
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	P	11	Total	C	N	O	S	0	0	0
			94	58	19	16	1			

- 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 53	C 27	N 9	O 15	P 2	0	0
3	D	1	Total 53	C 27	N 9	O 15	P 2	0	0
3	G	1	Total 53	C 27	N 9	O 15	P 2	0	0
3	J	1	Total 53	C 27	N 9	O 15	P 2	0	0

- Molecule 4 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (CCD ID: NAD) (formula: $\text{C}_{21}\text{H}_{27}\text{N}_7\text{O}_{14}\text{P}_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
4	D	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
4	G	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
4	J	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

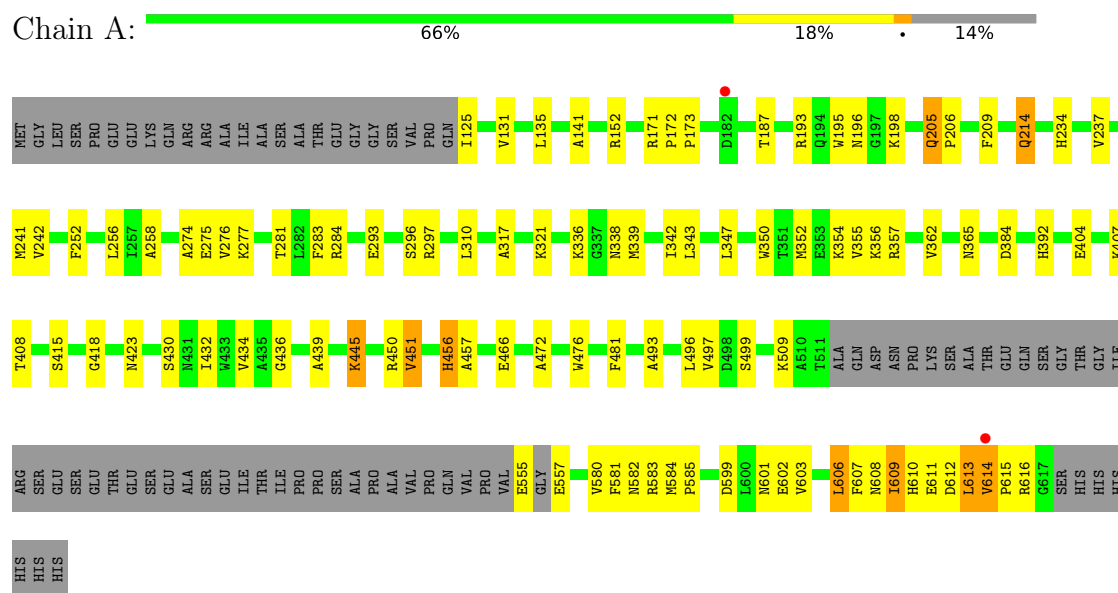
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	26	Total	O	0	0
			26	26		
5	D	20	Total	O	0	0
			20	20		
5	G	21	Total	O	0	0
			21	21		
5	J	14	Total	O	0	0
			14	14		
5	M	3	Total	O	0	0
			3	3		
5	N	2	Total	O	0	0
			2	2		

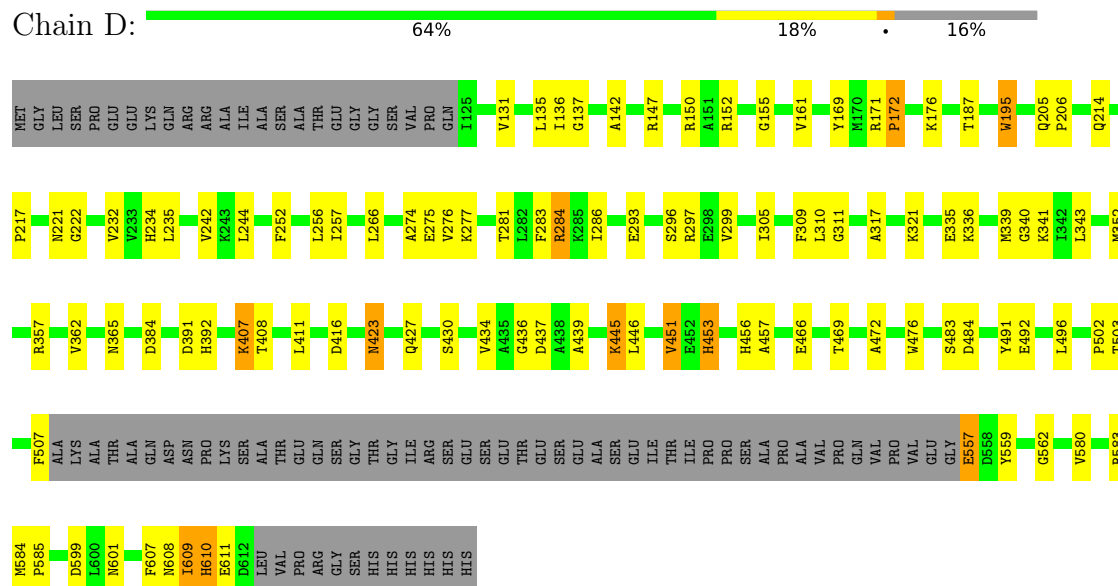
3 Residue-property plots

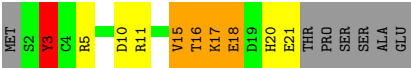
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: Apoptosis-inducing factor 1, mitochondrial

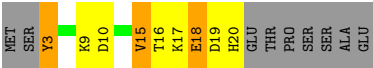
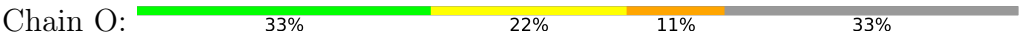


- Molecule 1: Apoptosis-inducing factor 1, mitochondrial

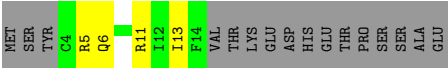




- Molecule 2: Mitochondrial intermembrane space import and assembly protein 40



- Molecule 2: Mitochondrial intermembrane space import and assembly protein 40



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	110.09Å 115.62Å 192.73Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	99.15 – 3.21 99.15 – 3.21	Depositor EDS
% Data completeness (in resolution range)	60.7 (99.15-3.21) 60.7 (99.15-3.21)	Depositor EDS
R_{merge}	0.30	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.43 (at 3.19Å)	Xtriage
Refinement program	REFMAC 5.8.0419	Depositor
R, R_{free}	0.233 , 0.284 0.230 , 0.279	Depositor DCC
R_{free} test set	1236 reflections (3.01%)	wwPDB-VP
Wilson B-factor (Å ²)	84.1	Xtriage
Anisotropy	0.158	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 56.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	0.029 for k,h,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	14701	wwPDB-VP
Average B, all atoms (Å ²)	77.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 18.30% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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

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	0.61	0/3541	1.02	3/4783 (0.1%)
1	D	0.56	0/3469	0.97	3/4687 (0.1%)
1	G	0.53	0/3469	0.93	1/4687 (0.0%)
1	J	0.50	0/3451	0.89	2/4661 (0.0%)
2	M	0.66	0/153	1.12	1/202 (0.5%)
2	N	0.63	0/173	1.16	1/229 (0.4%)
2	O	0.48	0/158	1.07	0/209
2	P	0.52	0/94	0.90	0/122
All	All	0.55	0/14508	0.96	11/19580 (0.1%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	G	0	1
All	All	0	2

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	456	HIS	CA-CB-CG	-5.96	107.84	113.80
1	D	172	PRO	N-CA-C	5.95	117.95	110.70
1	J	172	PRO	N-CA-C	5.68	117.63	110.70
1	G	172	PRO	N-CA-C	5.59	117.52	110.70
1	A	357	ARG	CB-CG-CD	5.54	124.05	111.30
1	D	484	ASP	CA-CB-CG	5.52	118.12	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	N	3	TYR	CA-CB-CG	5.37	123.56	113.90
1	J	262	THR	CB-CA-C	5.17	115.84	108.76
1	A	205	GLN	CB-CA-C	5.14	118.11	109.32
2	M	16	THR	CA-CB-OG1	-5.03	102.06	109.60
1	D	453	HIS	CA-CB-CG	-5.00	108.80	113.80

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	193	ARG	Sidechain
1	G	583	ARG	Sidechain

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	3472	0	3514	65	0
1	D	3400	0	3436	70	0
1	G	3400	0	3435	56	0
1	J	3382	0	3422	59	0
2	M	152	0	148	7	0
2	N	171	0	161	12	0
2	O	156	0	150	12	0
2	P	94	0	95	3	0
3	A	53	0	31	1	0
3	D	53	0	31	5	0
3	G	53	0	31	5	0
3	J	53	0	31	2	0
4	A	44	0	26	0	0
4	D	44	0	26	3	0
4	G	44	0	26	3	0
4	J	44	0	26	1	0
5	A	26	0	0	3	0
5	D	20	0	0	2	0
5	G	21	0	0	5	0
5	J	14	0	0	6	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	M	3	0	0	0	0
5	N	2	0	0	1	0
All	All	14701	0	14589	273	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 9.

All (273) 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:608:ASN:OD1	5:D:801:HOH:O	1.72	1.06
2:O:18:GLU:OE2	2:O:19:ASP:N	2.02	0.92
1:D:583:ARG:HH22	1:D:608:ASN:HB2	1.36	0.91
1:A:583:ARG:HH22	1:A:608:ASN:HB2	1.42	0.84
1:D:293:GLU:O	1:D:296:SER:OG	1.97	0.82
1:A:583:ARG:HE	1:A:613:LEU:HD11	1.44	0.81
1:G:612:ASP:O	5:G:801:HOH:O	2.00	0.80
1:D:195:TRP:CZ2	5:D:819:HOH:O	2.34	0.80
1:A:205:GLN:OE1	1:A:209:PHE:HB3	1.86	0.74
1:D:583:ARG:NH2	1:D:608:ASN:HB2	2.03	0.73
1:J:283:PHE:HB3	1:J:310:LEU:HD21	1.69	0.73
1:J:580:VAL:HG12	1:J:583:ARG:HG3	1.72	0.72
1:A:283:PHE:HB3	1:A:310:LEU:HD21	1.72	0.70
1:A:583:ARG:NH2	1:A:608:ASN:HB2	2.04	0.70
1:A:205:GLN:OE1	1:A:209:PHE:CB	2.39	0.70
1:G:283:PHE:HB3	1:G:310:LEU:HD21	1.73	0.70
2:O:18:GLU:OE2	2:O:19:ASP:OD1	2.10	0.69
1:D:283:PHE:HB3	1:D:310:LEU:HD21	1.75	0.69
1:G:580:VAL:HG12	1:G:583:ARG:HG3	1.75	0.69
1:D:580:VAL:HG12	1:D:583:ARG:HG3	1.74	0.68
2:N:20:HIS:O	2:N:21:GLU:C	2.35	0.68
1:A:499:SER:O	2:M:9:LYS:NZ	2.25	0.67
1:G:245:ASN:HB2	5:G:813:HOH:O	1.95	0.66
1:G:607:PHE:O	1:G:608:ASN:C	2.39	0.65
1:J:152:ARG:HD3	1:J:466:GLU:HA	1.81	0.62
1:G:583:ARG:NH2	1:G:608:ASN:HB2	2.14	0.62
1:A:352:MET:HG3	1:A:362:VAL:HG11	1.82	0.61
1:A:451:VAL:HG13	1:A:456:HIS:CE1	2.36	0.61
1:A:580:VAL:HG12	1:A:583:ARG:HG3	1.84	0.59
1:J:390:THR:HG22	5:J:802:HOH:O	2.01	0.59
1:G:483:SER:HB3	1:G:491:TYR:CE1	2.39	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:583:ARG:HH21	1:A:613:LEU:HG	1.68	0.58
1:D:451:VAL:HG23	1:D:453:HIS:HD2	1.69	0.58
1:A:606:LEU:HD23	1:G:586:ILE:HA	1.86	0.58
1:D:234:HIS:ND1	1:D:408:THR:HB	2.20	0.57
1:D:607:PHE:O	1:D:608:ASN:C	2.47	0.57
2:M:16:THR:OG1	2:M:18:GLU:OE1	2.17	0.57
1:A:581:PHE:HB2	1:A:614:VAL:O	2.05	0.57
1:D:281:THR:HG23	1:D:392:HIS:NE2	2.19	0.57
2:O:18:GLU:O	2:O:20:HIS:N	2.39	0.56
1:D:451:VAL:HG13	1:D:456:HIS:ND1	2.20	0.56
1:J:176:LYS:HE2	3:J:701:FAD:O4	2.05	0.56
1:G:451:VAL:HG23	1:G:453:HIS:HD2	1.71	0.56
1:J:162:SER:HB3	5:J:808:HOH:O	2.06	0.56
2:O:18:GLU:O	2:O:19:ASP:C	2.49	0.56
1:A:580:VAL:HG13	1:A:613:LEU:HD12	1.88	0.55
1:D:341:LYS:HD3	4:D:702:NAD:H3B	1.88	0.55
1:G:152:ARG:HD3	1:G:466:GLU:HA	1.87	0.55
1:J:172:PRO:HA	3:J:701:FAD:N5	2.20	0.55
1:A:607:PHE:O	1:A:608:ASN:C	2.49	0.55
1:J:317:ALA:O	1:J:321:LYS:HG2	2.06	0.55
1:G:281:THR:HG23	1:G:392:HIS:NE2	2.22	0.55
1:J:451:VAL:HG23	1:J:453:HIS:HD2	1.72	0.55
1:D:437:ASP:OD2	3:D:701:FAD:O3'	2.19	0.54
1:A:584:MET:N	1:A:585:PRO:CD	2.70	0.54
1:A:196:ASN:OD1	1:A:198:LYS:HD2	2.07	0.54
1:J:451:VAL:HG13	1:J:456:HIS:CE1	2.43	0.53
1:A:582:ASN:HA	5:A:814:HOH:O	2.08	0.53
1:D:176:LYS:HE2	3:D:701:FAD:O4	2.08	0.53
1:A:445:LYS:HE2	1:A:476:TRP:CE3	2.43	0.53
2:M:16:THR:O	2:M:18:GLU:N	2.40	0.53
1:A:274:ALA:HA	1:A:277:LYS:HB2	1.90	0.52
1:A:317:ALA:O	1:A:321:LYS:HG2	2.09	0.52
1:G:317:ALA:O	1:G:321:LYS:HG2	2.10	0.52
1:J:339:MET:HB2	1:J:343:LEU:HD12	1.91	0.52
2:O:16:THR:OG1	2:O:18:GLU:OE1	2.23	0.52
1:D:339:MET:HB2	1:D:343:LEU:HD12	1.91	0.52
1:D:451:VAL:HG13	1:D:456:HIS:CE1	2.45	0.52
1:D:341:LYS:CD	4:D:702:NAD:H3B	2.40	0.52
2:O:15:VAL:HG12	2:O:19:ASP:OD2	2.10	0.52
1:A:152:ARG:HD3	1:A:466:GLU:HA	1.90	0.52
1:J:281:THR:HG23	1:J:392:HIS:NE2	2.25	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:423:ASN:ND2	1:J:427:GLN:H	2.08	0.52
1:D:205:GLN:HB2	1:D:206:PRO:HD2	1.92	0.51
1:A:172:PRO:HA	3:A:701:FAD:N5	2.26	0.51
1:G:445:LYS:HE3	1:G:445:LYS:H	1.76	0.51
1:A:582:ASN:HD22	1:A:614:VAL:HB	1.75	0.51
2:N:3:TYR:CD1	2:N:3:TYR:C	2.89	0.51
1:D:257:ILE:HB	1:D:434:VAL:HG22	1.93	0.51
1:J:506:VAL:HA	2:P:13:ILE:O	2.10	0.51
1:A:338:ASN:HD22	1:A:352:MET:HA	1.76	0.51
2:P:6:GLN:OE1	2:P:11:ARG:HD2	2.12	0.51
1:D:152:ARG:HD3	1:D:466:GLU:HA	1.93	0.50
1:G:451:VAL:HG13	1:G:456:HIS:ND1	2.26	0.50
1:J:451:VAL:HG13	1:J:456:HIS:ND1	2.27	0.50
1:A:237:VAL:HG13	1:A:432:ILE:CD1	2.41	0.50
1:G:423:ASN:C	1:G:423:ASN:ND2	2.69	0.49
1:A:355:VAL:O	1:A:356:LYS:C	2.53	0.49
1:D:423:ASN:ND2	1:D:427:GLN:H	2.10	0.49
1:G:171:ARG:N	1:G:172:PRO:CD	2.76	0.49
1:J:226:VAL:HG12	5:J:808:HOH:O	2.12	0.49
1:J:584:MET:N	1:J:585:PRO:CD	2.76	0.49
1:G:156:ALA:O	1:G:222:GLY:HA3	2.13	0.49
1:J:212:SER:HA	5:J:801:HOH:O	2.12	0.49
1:D:483:SER:HB3	1:D:491:TYR:CE1	2.47	0.49
1:A:350:TRP:CZ2	1:A:354:LYS:HE3	2.48	0.49
1:G:583:ARG:HH21	1:G:608:ASN:HB2	1.77	0.49
1:A:472:ALA:HB1	1:D:430:SER:OG	2.13	0.48
1:G:165:PRO:HD3	5:G:820:HOH:O	2.13	0.48
1:G:509:LYS:N	2:N:15:VAL:O	2.43	0.48
3:G:701:FAD:C5X	4:G:702:NAD:C5N	2.91	0.48
1:J:156:ALA:HA	5:J:809:HOH:O	2.12	0.48
1:D:274:ALA:HA	1:D:277:LYS:HB2	1.94	0.48
1:G:483:SER:HB3	1:G:491:TYR:CZ	2.48	0.48
1:J:610:HIS:CG	1:J:611:GLU:N	2.81	0.48
1:G:451:VAL:HG13	1:G:456:HIS:CE1	2.48	0.48
1:D:608:ASN:O	1:D:610:HIS:N	2.46	0.48
1:J:138:GLY:HA3	1:J:162:SER:HB2	1.95	0.48
1:D:584:MET:N	1:D:585:PRO:CD	2.77	0.48
1:G:567:LEU:HD23	1:G:570:LYS:HA	1.95	0.48
1:J:451:VAL:HG13	1:J:456:HIS:CG	2.49	0.48
2:N:16:THR:O	2:N:17:LYS:C	2.57	0.48
1:A:339:MET:HB2	1:A:343:LEU:HD12	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:171:ARG:O	1:A:172:PRO:C	2.57	0.47
1:D:284:ARG:HD2	3:D:701:FAD:C8M	2.43	0.47
1:A:293:GLU:OE1	1:A:297:ARG:HD3	2.15	0.47
1:D:352:MET:HG3	1:D:362:VAL:HG11	1.94	0.47
1:D:436:GLY:HA2	1:D:457:ALA:HA	1.97	0.47
1:G:352:MET:HG3	1:G:362:VAL:HG11	1.96	0.47
1:D:451:VAL:HG13	1:D:456:HIS:CG	2.48	0.47
1:G:293:GLU:O	1:G:296:SER:OG	2.25	0.47
1:J:226:VAL:CG1	5:J:808:HOH:O	2.62	0.47
2:O:19:ASP:O	2:O:20:HIS:CG	2.67	0.47
1:J:595:GLU:HB3	1:J:597:HIS:CE1	2.49	0.47
1:G:343:LEU:HD22	1:G:347:LEU:HD23	1.96	0.47
1:J:235:LEU:HD23	1:J:411:LEU:HD11	1.97	0.47
1:A:509:LYS:HG2	2:M:16:THR:HG23	1.97	0.46
1:D:293:GLU:OE1	1:D:297:ARG:HD3	2.15	0.46
1:J:293:GLU:O	1:J:296:SER:OG	2.33	0.46
1:A:293:GLU:O	1:A:296:SER:OG	2.28	0.46
1:A:336:LYS:HG2	1:A:365:ASN:HD21	1.80	0.46
1:A:171:ARG:N	1:A:172:PRO:CD	2.79	0.46
1:G:423:ASN:C	1:G:423:ASN:HD22	2.23	0.46
1:A:436:GLY:HA2	1:A:457:ALA:HA	1.97	0.46
1:J:274:ALA:HA	1:J:277:LYS:HB2	1.96	0.46
2:N:16:THR:HB	2:N:17:LYS:HE3	1.97	0.46
1:A:205:GLN:HB2	1:A:206:PRO:HD2	1.97	0.46
1:A:445:LYS:HE3	1:A:445:LYS:H	1.81	0.46
1:D:434:VAL:HG12	1:D:439:ALA:HB2	1.98	0.46
1:G:431:ASN:C	1:G:432:ILE:HG13	2.41	0.46
1:J:266:LEU:HD13	4:J:702:NAD:H2A	1.98	0.46
1:A:418:GLY:O	1:A:450:ARG:HD3	2.15	0.46
1:G:355:VAL:O	1:G:356:LYS:C	2.59	0.46
1:G:577:LEU:HD13	1:G:583:ARG:HB2	1.98	0.46
1:J:446:LEU:HD13	1:J:496:LEU:HD13	1.97	0.46
1:J:336:LYS:HG2	1:J:365:ASN:HD21	1.80	0.46
1:D:446:LEU:HD13	1:D:496:LEU:HD13	1.97	0.45
1:G:281:THR:HG23	1:G:392:HIS:CE1	2.51	0.45
1:J:577:LEU:HD13	1:J:583:ARG:HB2	1.98	0.45
1:A:555:GLU:OE2	1:A:555:GLU:HA	2.16	0.45
2:N:3:TYR:C	2:N:3:TYR:HD1	2.24	0.45
1:G:397:VAL:HG12	4:G:702:NAD:N3A	2.31	0.45
1:J:559:TYR:OH	2:P:5:ARG:NH2	2.46	0.45
1:A:281:THR:HG23	1:A:392:HIS:NE2	2.32	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:595:GLU:HB3	1:G:597:HIS:CE1	2.52	0.45
1:D:136:ILE:HG12	1:D:232:VAL:HG21	1.98	0.45
1:D:284:ARG:HD2	3:D:701:FAD:HM83	1.99	0.45
1:A:237:VAL:HG13	1:A:432:ILE:HD11	1.99	0.45
1:A:430:SER:OG	1:D:472:ALA:HB1	2.16	0.45
1:D:135:LEU:HD22	1:D:256:LEU:HB3	1.99	0.45
1:A:599:ASP:OD2	1:A:601:ASN:HB3	2.17	0.44
1:A:583:ARG:HE	1:A:613:LEU:CD1	2.24	0.44
1:J:205:GLN:OE1	1:J:209:PHE:O	2.35	0.44
1:J:237:VAL:HG13	1:J:432:ILE:CD1	2.47	0.44
2:M:2:SER:O	2:M:4:CYS:N	2.51	0.44
2:M:16:THR:OG1	2:M:18:GLU:CD	2.60	0.44
2:N:11:ARG:NH2	5:N:101:HOH:O	2.49	0.44
1:D:335:GLU:HG2	1:D:340:GLY:HA3	2.00	0.44
1:G:557:GLU:HG3	1:G:559:TYR:CE2	2.52	0.44
1:J:234:HIS:ND1	1:J:408:THR:HB	2.33	0.44
1:A:141:ALA:HB3	1:A:258:ALA:HB1	1.99	0.44
1:D:557:GLU:HG3	1:D:559:TYR:CE2	2.53	0.44
1:G:456:HIS:HA	1:G:479:SER:CB	2.47	0.44
1:J:583:ARG:NH2	1:J:608:ASN:HB2	2.33	0.44
1:D:150:ARG:HH22	1:D:155:GLY:H	1.65	0.44
1:A:172:PRO:HB2	1:A:173:PRO:HD3	2.00	0.44
1:D:235:LEU:HD23	1:D:411:LEU:HD11	2.00	0.44
1:G:227:LEU:HD22	1:G:230:LYS:HG3	2.00	0.44
1:J:281:THR:HG23	1:J:392:HIS:CE1	2.53	0.44
1:J:583:ARG:HH22	1:J:608:ASN:HB2	1.83	0.44
2:N:17:LYS:HD2	2:N:18:GLU:HG3	2.00	0.44
1:A:131:VAL:O	1:A:252:PHE:HA	2.18	0.43
1:A:603:VAL:HA	1:A:606:LEU:HD12	1.99	0.43
1:J:174:LEU:HD23	1:J:178:LEU:HD12	2.00	0.43
1:J:299:VAL:HB	1:J:391:ASP:HB2	1.99	0.43
1:J:418:GLY:O	1:J:450:ARG:HD3	2.17	0.43
1:G:509:LYS:O	1:G:510:ALA:HB2	2.18	0.43
1:J:171:ARG:N	1:J:172:PRO:CD	2.82	0.43
1:D:309:PHE:HB3	4:D:702:NAD:C4N	2.49	0.43
1:G:139:GLY:HA3	3:G:701:FAD:O5B	2.18	0.43
1:A:451:VAL:HG13	1:A:456:HIS:ND1	2.32	0.43
1:G:174:LEU:HD23	1:G:178:LEU:HD12	2.00	0.43
1:G:304:VAL:O	1:G:331:GLN:HA	2.18	0.43
1:D:492:GLU:HB3	1:D:584:MET:HE2	2.00	0.43
1:J:141:ALA:HB3	1:J:258:ALA:HB1	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:483:SER:HB3	1:J:491:TYR:CE1	2.53	0.43
1:A:125:ILE:HG23	1:A:214:GLN:NE2	2.34	0.43
1:A:234:HIS:ND1	1:A:408:THR:HB	2.33	0.43
1:A:407:LYS:H	1:A:407:LYS:HG2	1.56	0.43
1:D:137:GLY:O	1:D:142:ALA:HB2	2.19	0.43
1:G:423:ASN:ND2	1:G:427:GLN:H	2.17	0.43
1:A:481:PHE:CZ	1:A:493:ALA:HB3	2.54	0.42
1:D:293:GLU:OE2	1:D:297:ARG:NH2	2.51	0.42
1:J:319:GLY:O	1:J:322:SER:OG	2.33	0.42
1:J:423:ASN:HD22	1:J:427:GLN:H	1.67	0.42
2:O:18:GLU:OE2	2:O:19:ASP:CA	2.67	0.42
1:A:434:VAL:HG12	1:A:439:ALA:HB2	2.01	0.42
1:G:275:GLU:HB2	5:G:818:HOH:O	2.19	0.42
1:J:352:MET:HG3	1:J:362:VAL:HG11	2.00	0.42
1:D:336:LYS:HG2	1:D:365:ASN:HD21	1.85	0.42
1:D:407:LYS:H	1:D:407:LYS:HG3	1.63	0.42
1:D:599:ASP:OD2	1:D:601:ASN:HB3	2.19	0.42
1:G:509:LYS:HB3	2:N:16:THR:HG23	2.00	0.42
1:G:257:ILE:HB	1:G:434:VAL:HG22	2.01	0.42
2:M:2:SER:O	2:M:3:TYR:C	2.62	0.42
1:D:310:LEU:O	1:D:311:GLY:C	2.63	0.42
1:D:445:LYS:HE2	1:D:476:TRP:CE3	2.55	0.42
1:J:205:GLN:HB2	1:J:206:PRO:HD2	2.01	0.42
1:A:135:LEU:HD22	1:A:256:LEU:HB3	2.02	0.42
1:D:221:ASN:O	1:D:222:GLY:C	2.62	0.42
1:D:317:ALA:O	1:D:321:LYS:HG2	2.20	0.42
1:D:503:THR:O	2:O:10:ASP:HA	2.20	0.42
1:J:136:ILE:HG12	1:J:232:VAL:HG21	2.02	0.42
1:G:503:THR:O	2:N:10:ASP:HA	2.20	0.42
1:G:208:SER:HB2	5:G:805:HOH:O	2.20	0.42
1:G:446:LEU:HD13	1:G:496:LEU:HD13	2.01	0.42
1:G:581:PHE:O	1:G:583:ARG:HG2	2.19	0.42
1:J:508:ALA:O	1:J:509:LYS:CB	2.68	0.42
1:G:172:PRO:HG3	3:G:701:FAD:C4X	2.50	0.42
2:N:17:LYS:H	2:N:17:LYS:HG3	1.55	0.42
1:D:232:VAL:HG13	1:D:242:VAL:HB	2.02	0.42
1:D:483:SER:HB3	1:D:491:TYR:CZ	2.55	0.42
1:A:343:LEU:HD22	1:A:347:LEU:HD23	2.02	0.41
1:D:136:ILE:HD11	1:D:242:VAL:HG11	2.02	0.41
1:D:171:ARG:N	1:D:172:PRO:CD	2.83	0.41
1:J:557:GLU:HG3	1:J:559:TYR:CE2	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:214:GLN:O	1:G:217:PRO:HD2	2.20	0.41
1:G:274:ALA:HA	1:G:277:LYS:HB2	2.01	0.41
1:J:599:ASP:OD2	1:J:601:ASN:HB3	2.20	0.41
1:G:472:ALA:HB1	1:J:430:SER:OG	2.19	0.41
1:J:140:THR:HG21	1:J:454:HIS:CE1	2.55	0.41
1:A:205:GLN:OE1	1:A:209:PHE:HB2	2.20	0.41
1:A:342:ILE:HB	1:A:497:VAL:HG12	2.01	0.41
1:A:609:ILE:O	1:A:610:HIS:C	2.64	0.41
1:D:502:PRO:HB3	2:O:9:LYS:HB2	2.03	0.41
1:D:284:ARG:HA	3:D:701:FAD:HM81	2.03	0.41
1:A:439:ALA:O	5:A:801:HOH:O	2.22	0.41
1:D:131:VAL:O	1:D:252:PHE:HA	2.20	0.41
1:D:214:GLN:O	1:D:217:PRO:HD2	2.20	0.41
1:D:445:LYS:H	1:D:445:LYS:HE3	1.85	0.41
1:J:492:GLU:HB3	1:J:584:MET:HE2	2.02	0.41
1:J:227:LEU:HD22	1:J:230:LYS:HG3	2.02	0.41
2:N:18:GLU:HG3	2:N:18:GLU:H	1.42	0.41
1:A:241:MET:HG2	1:A:242:VAL:N	2.35	0.41
3:G:701:FAD:HM83	3:G:701:FAD:HM71	1.91	0.41
1:J:443:ASP:HB3	1:J:447:GLY:O	2.21	0.41
1:D:266:LEU:HD22	1:D:305:ILE:HG21	2.02	0.41
1:D:281:THR:HG23	1:D:392:HIS:CE1	2.56	0.41
1:G:172:PRO:HA	3:G:701:FAD:N5	2.36	0.41
2:O:3:TYR:CD1	2:O:3:TYR:C	2.99	0.41
1:D:152:ARG:NE	1:D:469:THR:OG1	2.44	0.40
1:J:156:ALA:O	1:J:222:GLY:HA3	2.21	0.40
2:O:16:THR:OG1	2:O:18:GLU:CD	2.64	0.40
1:J:135:LEU:HD22	1:J:256:LEU:HB3	2.03	0.40
1:A:582:ASN:HD22	1:A:614:VAL:CG2	2.34	0.40
1:D:169:TYR:CE2	1:D:286:ILE:HG13	2.56	0.40
1:D:299:VAL:HB	1:D:391:ASP:HB2	2.03	0.40
1:G:397:VAL:CG1	4:G:702:NAD:N3A	2.85	0.40
1:G:445:LYS:HE2	1:G:476:TRP:CE3	2.57	0.40
1:A:456:HIS:HE1	5:A:801:HOH:O	2.04	0.40
1:D:161:VAL:HG11	1:D:244:LEU:HD21	2.03	0.40
1:D:507:PHE:HA	1:D:562:GLY:HA3	2.03	0.40
1:G:440:CYS:O	1:G:448:ARG:NH1	2.55	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all 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	444/524 (85%)	411 (93%)	31 (7%)	2 (0%)	24	59
1	D	435/524 (83%)	406 (93%)	26 (6%)	3 (1%)	18	52
1	G	436/524 (83%)	406 (93%)	28 (6%)	2 (0%)	24	59
1	J	433/524 (83%)	402 (93%)	28 (6%)	3 (1%)	18	52
2	M	16/27 (59%)	13 (81%)	2 (12%)	1 (6%)	1	8
2	N	18/27 (67%)	16 (89%)	2 (11%)	0	100	100
2	O	16/27 (59%)	14 (88%)	2 (12%)	0	100	100
2	P	9/27 (33%)	8 (89%)	1 (11%)	0	100	100
All	All	1807/2204 (82%)	1676 (93%)	120 (7%)	11 (1%)	21	56

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	284	ARG
1	J	608	ASN
2	M	3	TYR
1	A	284	ARG
1	D	610	HIS
1	G	284	ARG
1	A	615	PRO
1	D	609	ILE
1	J	284	ARG
1	G	609	ILE
1	J	609	ILE

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	368/429 (86%)	347 (94%)	21 (6%)	18	51
1	D	361/429 (84%)	346 (96%)	15 (4%)	26	60
1	G	360/429 (84%)	341 (95%)	19 (5%)	20	53
1	J	359/429 (84%)	346 (96%)	13 (4%)	31	63
2	M	17/25 (68%)	13 (76%)	4 (24%)	1	4
2	N	19/25 (76%)	13 (68%)	6 (32%)	0	1
2	O	17/25 (68%)	13 (76%)	4 (24%)	1	4
2	P	10/25 (40%)	10 (100%)	0	100	100
All	All	1511/1816 (83%)	1429 (95%)	82 (5%)	20	53

All (82) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	187	THR
1	A	195	TRP
1	A	214	GLN
1	A	275	GLU
1	A	276	VAL
1	A	384	ASP
1	A	404	GLU
1	A	415	SER
1	A	423	ASN
1	A	445	LYS
1	A	451	VAL
1	A	496	LEU
1	A	557	GLU
1	A	602	GLU
1	A	606	LEU
1	A	609	ILE
1	A	611	GLU
1	A	612	ASP
1	A	613	LEU
1	A	614	VAL
1	A	616	ARG
1	D	147	ARG
1	D	187	THR
1	D	195	TRP

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Mol	Chain	Res	Type
1	D	275	GLU
1	D	276	VAL
1	D	357	ARG
1	D	384	ASP
1	D	407	LYS
1	D	416	ASP
1	D	423	ASN
1	D	445	LYS
1	D	451	VAL
1	D	557	GLU
1	D	609	ILE
1	D	611	GLU
1	G	147	ARG
1	G	187	THR
1	G	195	TRP
1	G	214	GLN
1	G	219	ILE
1	G	275	GLU
1	G	276	VAL
1	G	345	GLN
1	G	357	ARG
1	G	384	ASP
1	G	404	GLU
1	G	407	LYS
1	G	423	ASN
1	G	445	LYS
1	G	451	VAL
1	G	496	LEU
1	G	557	GLU
1	G	609	ILE
1	G	612	ASP
1	J	187	THR
1	J	195	TRP
1	J	275	GLU
1	J	276	VAL
1	J	357	ARG
1	J	384	ASP
1	J	423	ASN
1	J	445	LYS
1	J	451	VAL
1	J	496	LEU
1	J	609	ILE

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Mol	Chain	Res	Type
1	J	610	HIS
1	J	611	GLU
2	M	15	VAL
2	M	17	LYS
2	M	18	GLU
2	M	19	ASP
2	N	3	TYR
2	N	5	ARG
2	N	15	VAL
2	N	16	THR
2	N	17	LYS
2	N	18	GLU
2	O	3	TYR
2	O	15	VAL
2	O	17	LYS
2	O	18	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (29) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	240	ASN
1	A	323	GLN
1	A	338	ASN
1	A	365	ASN
1	A	423	ASN
1	A	467	ASN
1	A	582	ASN
1	D	130	HIS
1	D	196	ASN
1	D	240	ASN
1	D	249	GLN
1	D	365	ASN
1	D	423	ASN
1	D	453	HIS
1	D	467	ASN
1	D	597	HIS
1	G	196	ASN
1	G	218	ASN
1	G	240	ASN
1	G	365	ASN
1	G	369	GLN
1	G	423	ASN

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Mol	Chain	Res	Type
1	G	454	HIS
1	G	597	HIS
1	J	218	ASN
1	J	240	ASN
1	J	365	ASN
1	J	423	ASN
1	J	597	HIS

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 ⓘ

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
4	NAD	J	702	-	46,48,48	0.71	2 (4%)	64,73,73	0.59	1 (1%)
3	FAD	D	701	-	58,58,58	0.51	0	85,89,89	0.63	0
3	FAD	J	701	-	58,58,58	0.56	0	85,89,89	0.72	1 (1%)
4	NAD	G	702	-	46,48,48	0.74	1 (2%)	64,73,73	0.73	1 (1%)
3	FAD	A	701	-	58,58,58	0.58	0	85,89,89	0.83	2 (2%)
3	FAD	G	701	-	58,58,58	0.53	0	85,89,89	0.85	5 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	NAD	A	702	-	46,48,48	0.76	2 (4%)	64,73,73	0.78	2 (3%)
4	NAD	D	702	-	46,48,48	0.86	4 (8%)	64,73,73	0.79	2 (3%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAD	J	702	-	-	10/30/62/62	0/5/5/5
3	FAD	D	701	-	-	5/34/50/50	0/6/6/6
3	FAD	J	701	-	-	5/34/50/50	0/6/6/6
4	NAD	G	702	-	-	7/30/62/62	0/5/5/5
3	FAD	A	701	-	-	5/34/50/50	0/6/6/6
3	FAD	G	701	-	-	6/34/50/50	0/6/6/6
4	NAD	A	702	-	-	8/30/62/62	0/5/5/5
4	NAD	D	702	-	-	2/30/62/62	0/5/5/5

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	702	NAD	C2N-N1N	3.23	1.38	1.35
4	D	702	NAD	C2N-N1N	3.09	1.38	1.35
4	G	702	NAD	C2N-N1N	2.83	1.38	1.35
4	J	702	NAD	C2N-N1N	2.73	1.38	1.35
4	D	702	NAD	PN-O3	2.69	1.62	1.59
4	D	702	NAD	PA-O3	2.50	1.62	1.59
4	A	702	NAD	O4D-C1D	2.26	1.43	1.40
4	D	702	NAD	O4D-C1D	2.07	1.43	1.40
4	J	702	NAD	O4D-C1D	2.00	1.43	1.40

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	701	FAD	O3'-C3'-C2'	-3.17	101.72	108.93
3	G	701	FAD	C3B-C2B-C1B	-2.93	95.91	101.46
3	J	701	FAD	C3B-C2B-C1B	-2.81	96.15	101.46
4	D	702	NAD	O2N-PN-O3	2.75	114.71	107.27
3	A	701	FAD	O2'-C2'-C3'	-2.59	103.18	109.25
4	D	702	NAD	C6N-N1N-C2N	-2.56	119.70	121.88

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	G	702	NAD	C6N-N1N-C2N	-2.45	119.79	121.88
3	G	701	FAD	O2P-P-O1P	2.43	123.73	112.44
4	A	702	NAD	C2N-C3N-C4N	2.23	120.85	118.26
3	G	701	FAD	O2B-C2B-C3B	2.13	118.64	111.82
3	G	701	FAD	C1B-N9A-C8A	2.10	131.76	127.09
4	J	702	NAD	C6N-N1N-C2N	-2.08	120.10	121.88
4	A	702	NAD	O2A-PA-O1A	2.08	122.10	112.44
3	G	701	FAD	C4A-N9A-C1B	-2.02	121.91	126.63

There are no chirality outliers.

All (48) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	701	FAD	N10-C1'-C2'-O2'
3	D	701	FAD	N10-C1'-C2'-O2'
3	D	701	FAD	N10-C1'-C2'-C3'
3	G	701	FAD	N10-C1'-C2'-O2'
3	J	701	FAD	N10-C1'-C2'-O2'
3	J	701	FAD	PA-O3P-P-O5'
4	A	702	NAD	C5D-O5D-PN-O3
4	A	702	NAD	C5D-O5D-PN-O1N
4	A	702	NAD	C5D-O5D-PN-O2N
4	G	702	NAD	C5D-O5D-PN-O3
4	G	702	NAD	C5D-O5D-PN-O1N
4	G	702	NAD	C5D-O5D-PN-O2N
4	J	702	NAD	C5B-O5B-PA-O2A
4	J	702	NAD	C5B-O5B-PA-O3
4	J	702	NAD	C5D-O5D-PN-O3
4	J	702	NAD	C5D-O5D-PN-O2N
4	A	702	NAD	C3B-C4B-C5B-O5B
4	A	702	NAD	O4B-C4B-C5B-O5B
4	G	702	NAD	O4B-C4B-C5B-O5B
4	G	702	NAD	C3B-C4B-C5B-O5B
4	G	702	NAD	O4D-C4D-C5D-O5D
4	J	702	NAD	C3B-C4B-C5B-O5B
3	A	701	FAD	O4B-C4B-C5B-O5B
3	A	701	FAD	C3B-C4B-C5B-O5B
3	G	701	FAD	C3B-C4B-C5B-O5B
4	G	702	NAD	C3D-C4D-C5D-O5D
3	G	701	FAD	O4B-C4B-C5B-O5B
3	A	701	FAD	PA-O3P-P-O5'
3	D	701	FAD	PA-O3P-P-O5'

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Mol	Chain	Res	Type	Atoms
3	D	701	FAD	C2B-C1B-N9A-C8A
3	G	701	FAD	C2B-C1B-N9A-C8A
3	J	701	FAD	C2B-C1B-N9A-C8A
4	A	702	NAD	PN-O3-PA-O2A
4	J	702	NAD	O4B-C4B-C5B-O5B
4	D	702	NAD	C5B-O5B-PA-O1A
4	J	702	NAD	C5D-O5D-PN-O1N
4	J	702	NAD	O4D-C4D-C5D-O5D
4	J	702	NAD	C3D-C4D-C5D-O5D
4	A	702	NAD	PN-O3-PA-O1A
3	D	701	FAD	O4B-C1B-N9A-C8A
3	G	701	FAD	O4B-C1B-N9A-C8A
4	D	702	NAD	PN-O3-PA-O2A
4	J	702	NAD	PA-O3-PN-O2N
3	J	701	FAD	O4B-C1B-N9A-C8A
3	J	701	FAD	O4B-C4B-C5B-O5B
3	G	701	FAD	C2B-C1B-N9A-C4A
3	A	701	FAD	N10-C1'-C2'-C3'
4	A	702	NAD	C3D-C4D-C5D-O5D

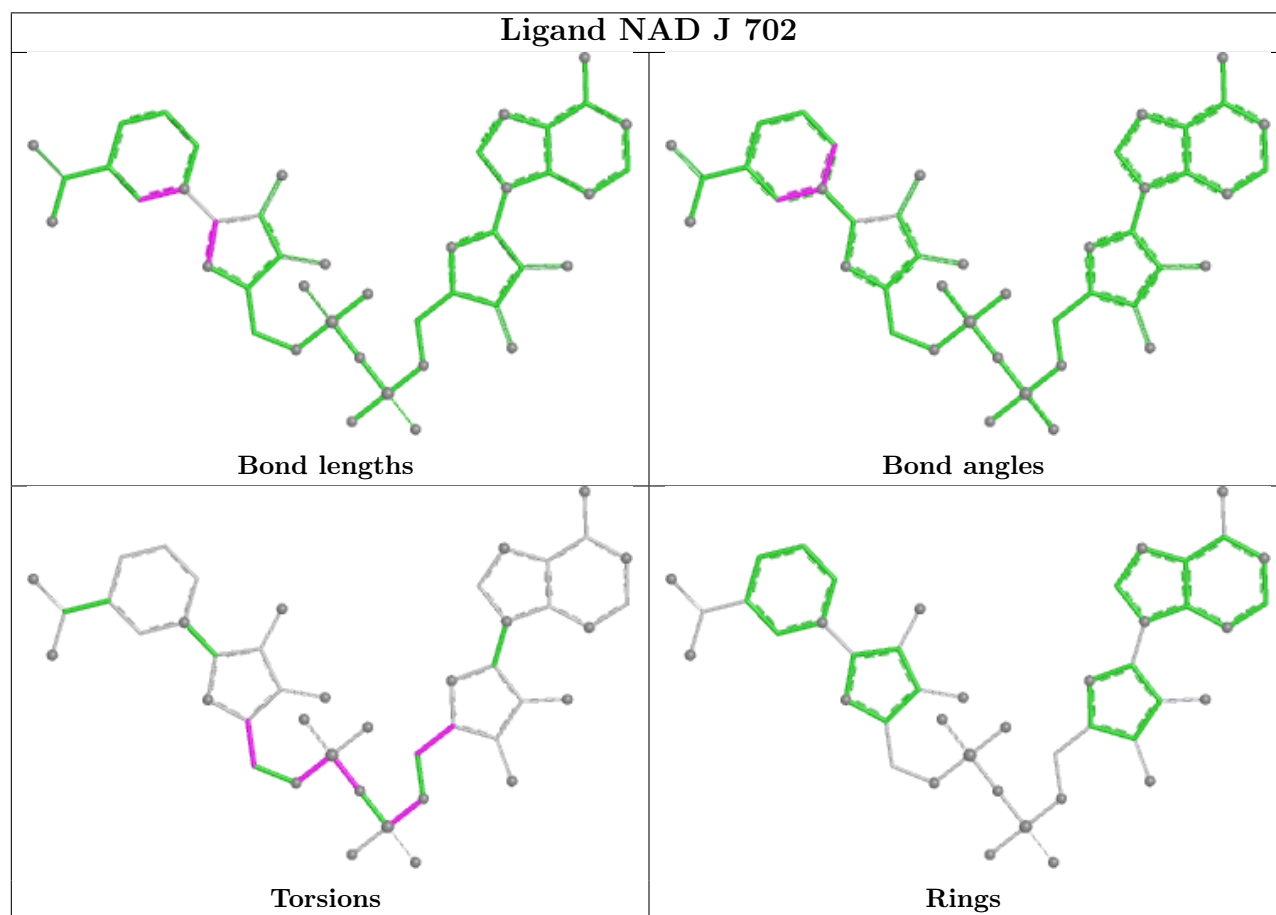
There are no ring outliers.

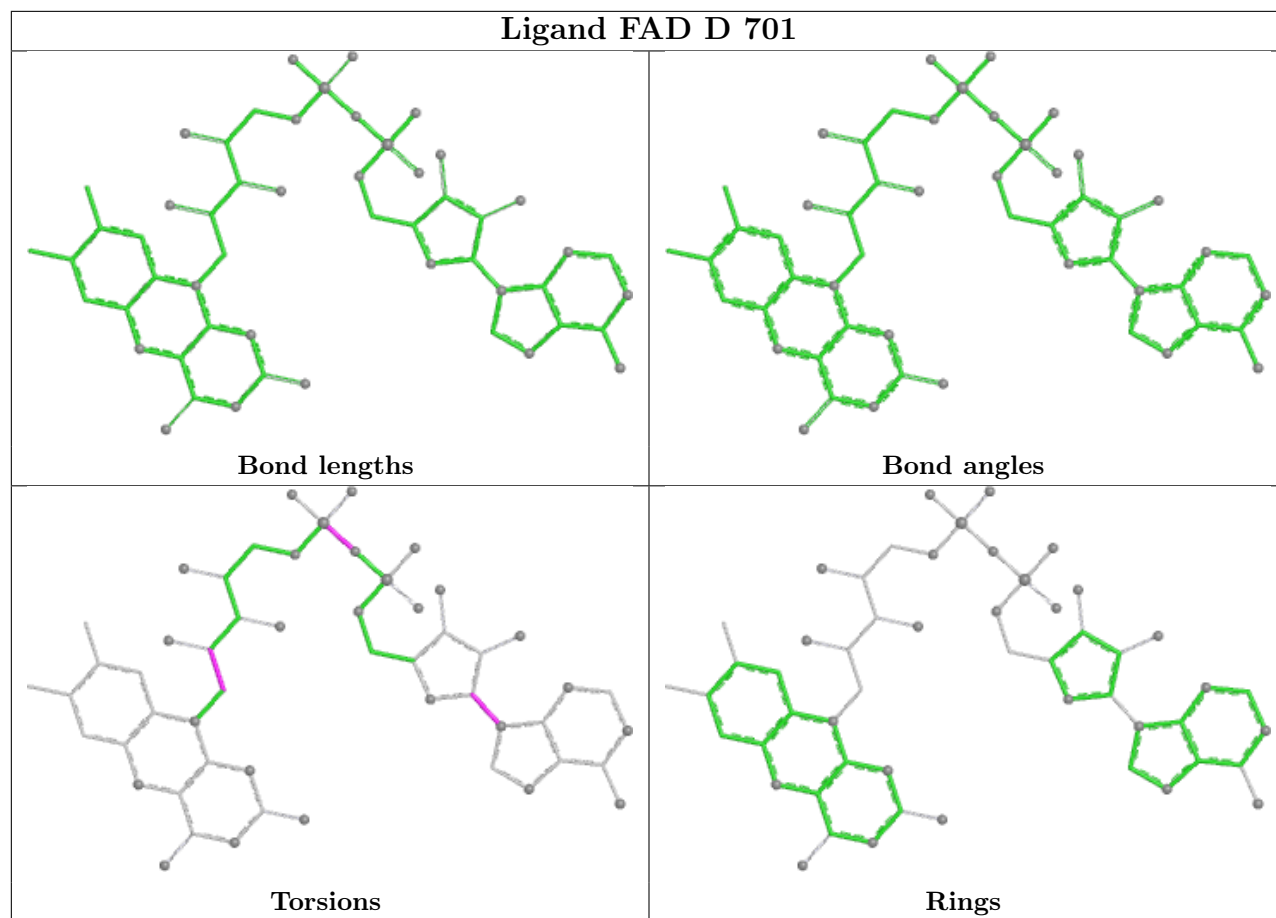
7 monomers are involved in 19 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	J	702	NAD	1	0
3	D	701	FAD	5	0
3	J	701	FAD	2	0
4	G	702	NAD	3	0
3	A	701	FAD	1	0
3	G	701	FAD	5	0
4	D	702	NAD	3	0

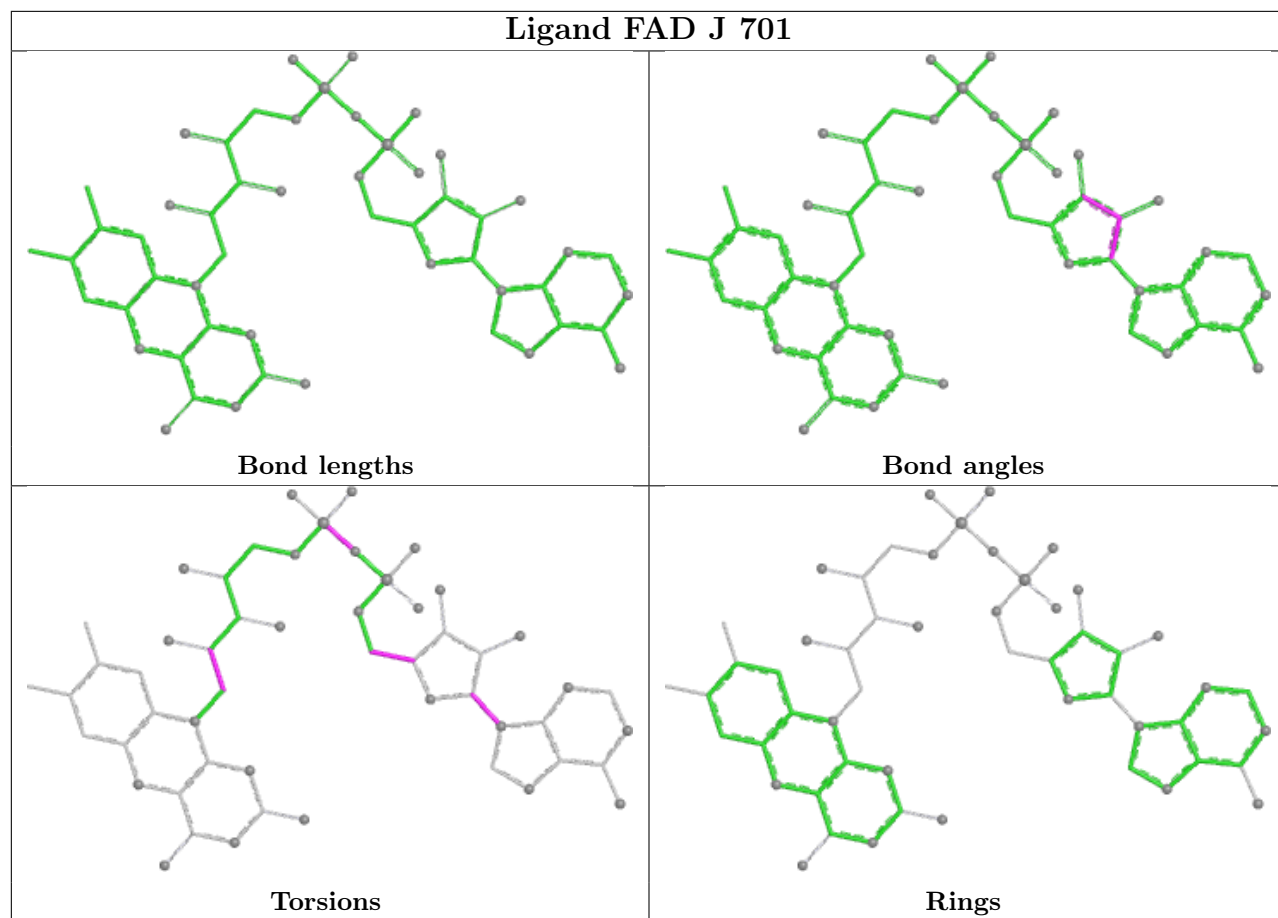
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

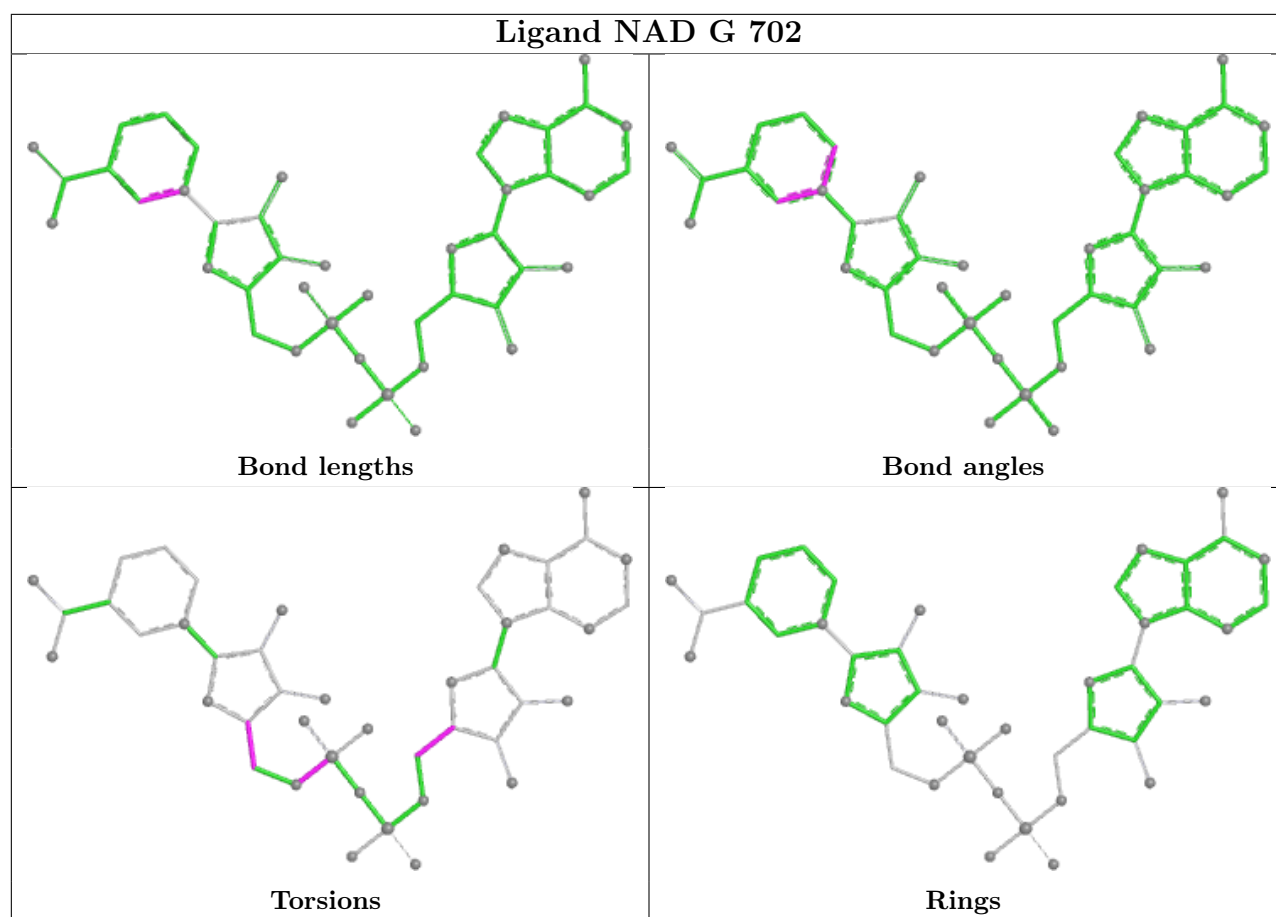
The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

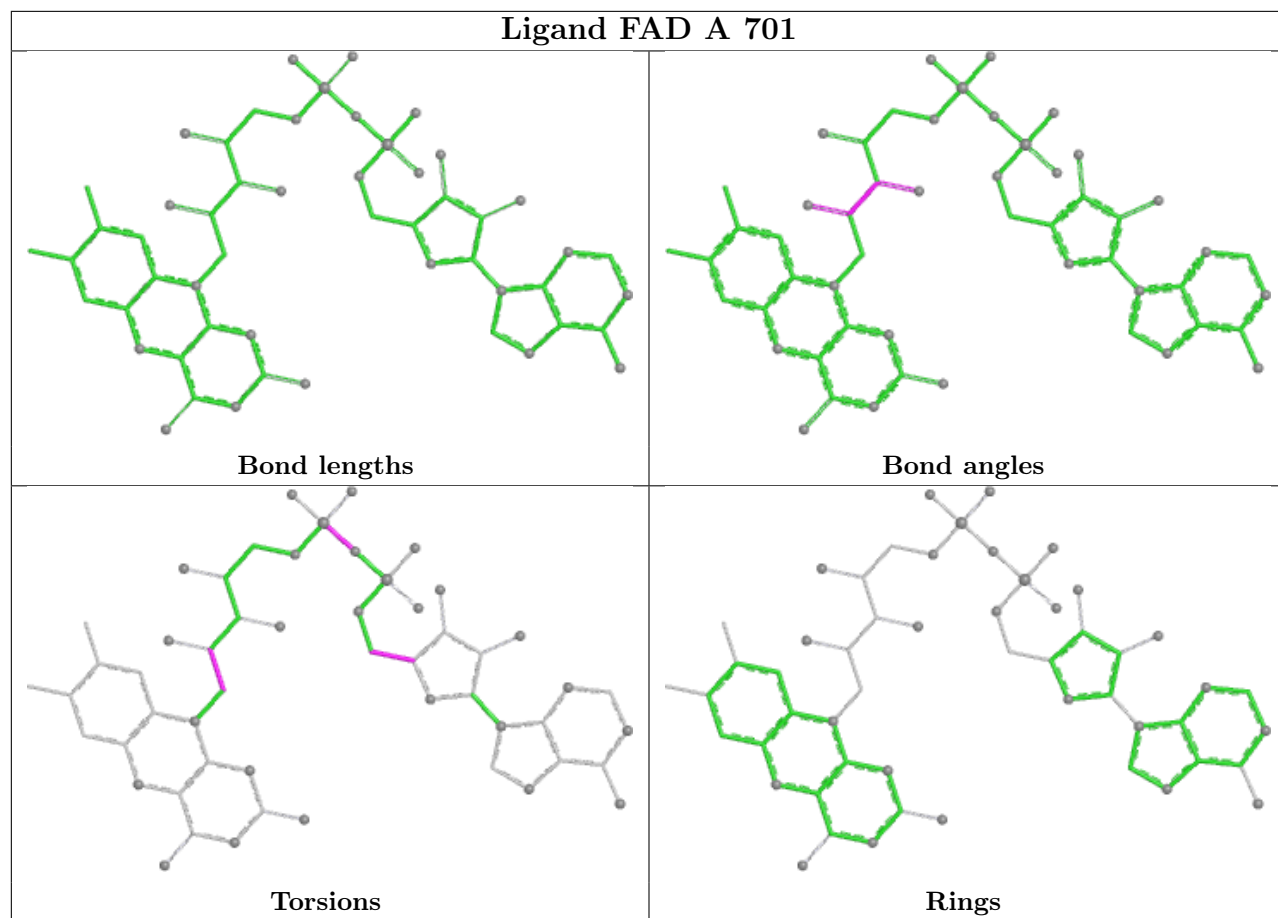


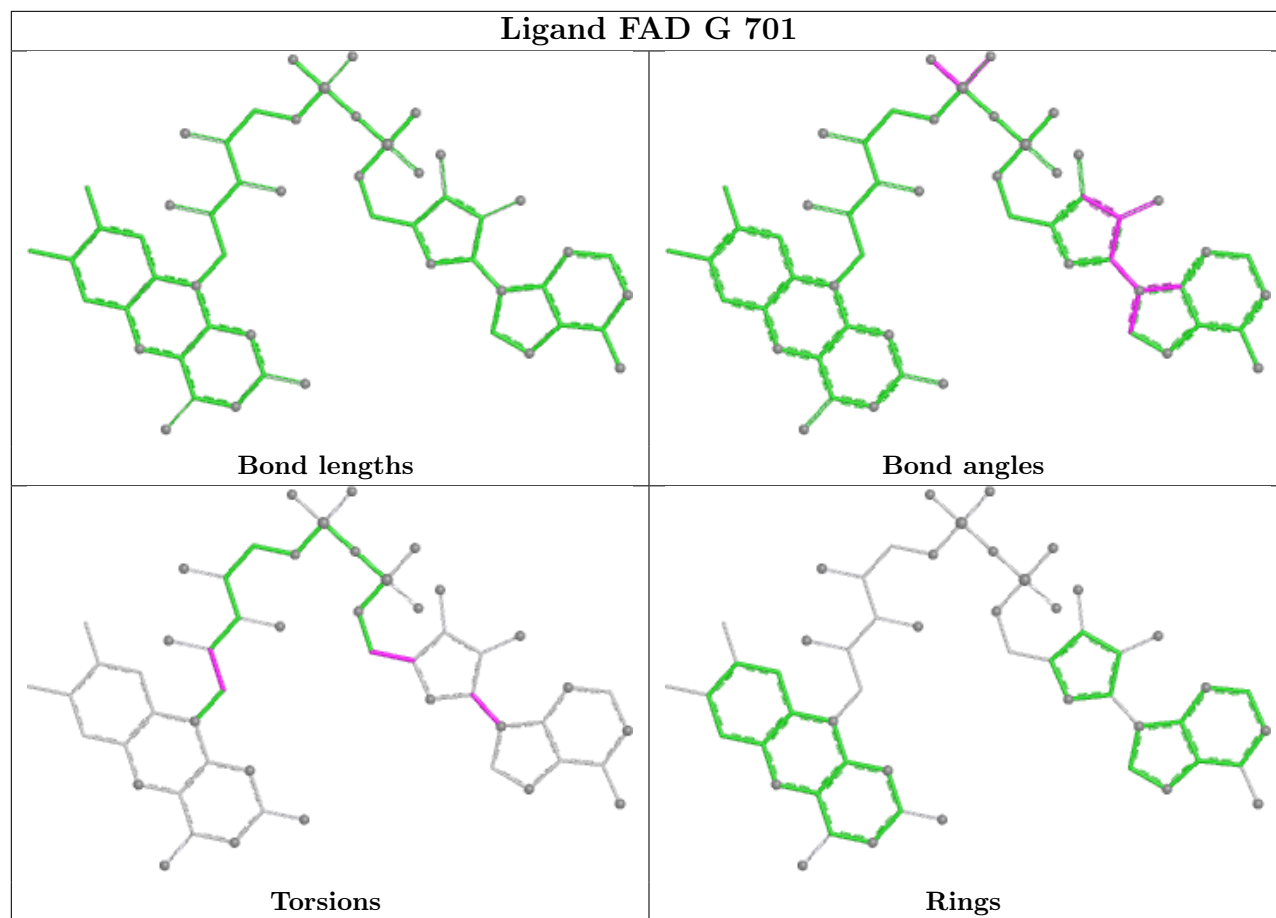


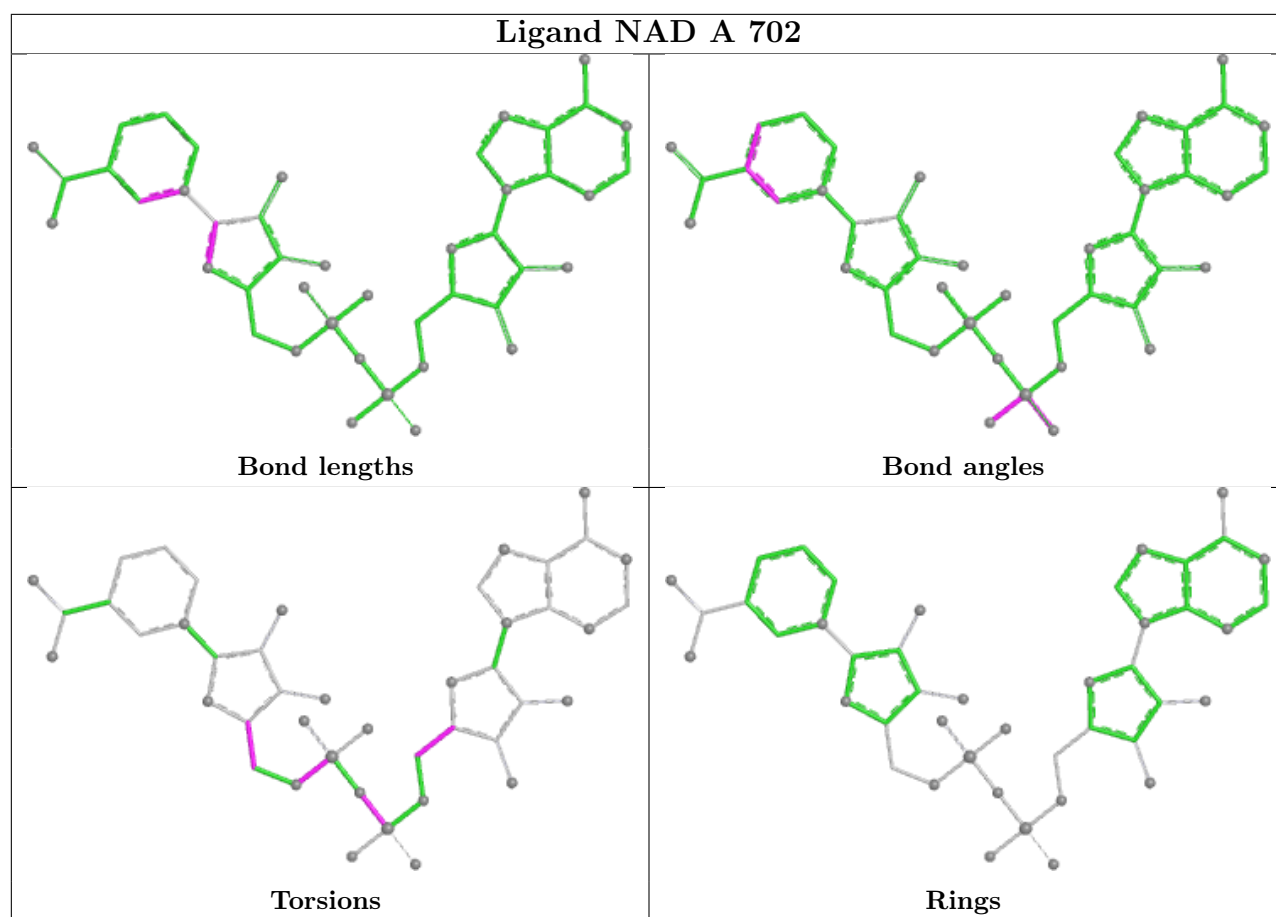
Ligand FAD J 701

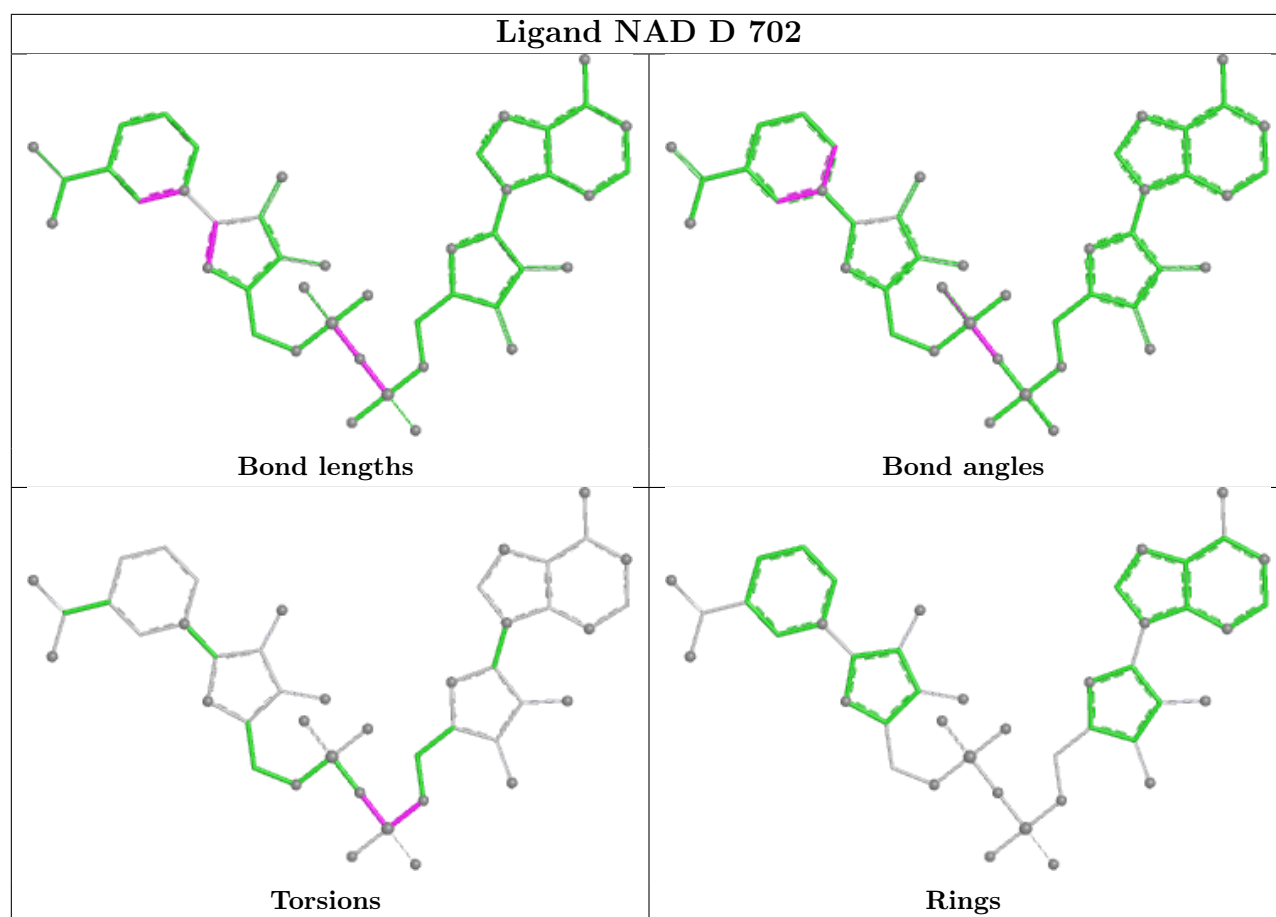












5.7 Other polymers [i](#)

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	449/524 (85%)	-0.46	2 (0%) 88 79	28, 52, 89, 117	0
1	D	439/524 (83%)	-0.54	0 100 100	33, 65, 108, 141	0
1	G	440/524 (83%)	-0.42	1 (0%) 91 85	52, 76, 107, 134	0
1	J	437/524 (83%)	-0.20	3 (0%) 84 69	58, 106, 176, 221	0
2	M	18/27 (66%)	-0.34	0 100 100	61, 71, 85, 86	0
2	N	20/27 (74%)	-0.05	0 100 100	74, 78, 82, 82	0
2	O	18/27 (66%)	-0.17	0 100 100	111, 114, 127, 129	0
2	P	11/27 (40%)	0.02	0 100 100	87, 88, 91, 92	0
All	All	1832/2204 (83%)	-0.40	6 (0%) 90 81	28, 73, 140, 221	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	614	VAL	2.6
1	J	190	LEU	2.5
1	G	510	ALA	2.4
1	J	366	ALA	2.2
1	A	182	ASP	2.2
1	J	318	LEU	2.2

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

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

6.3 Carbohydrates [i](#)

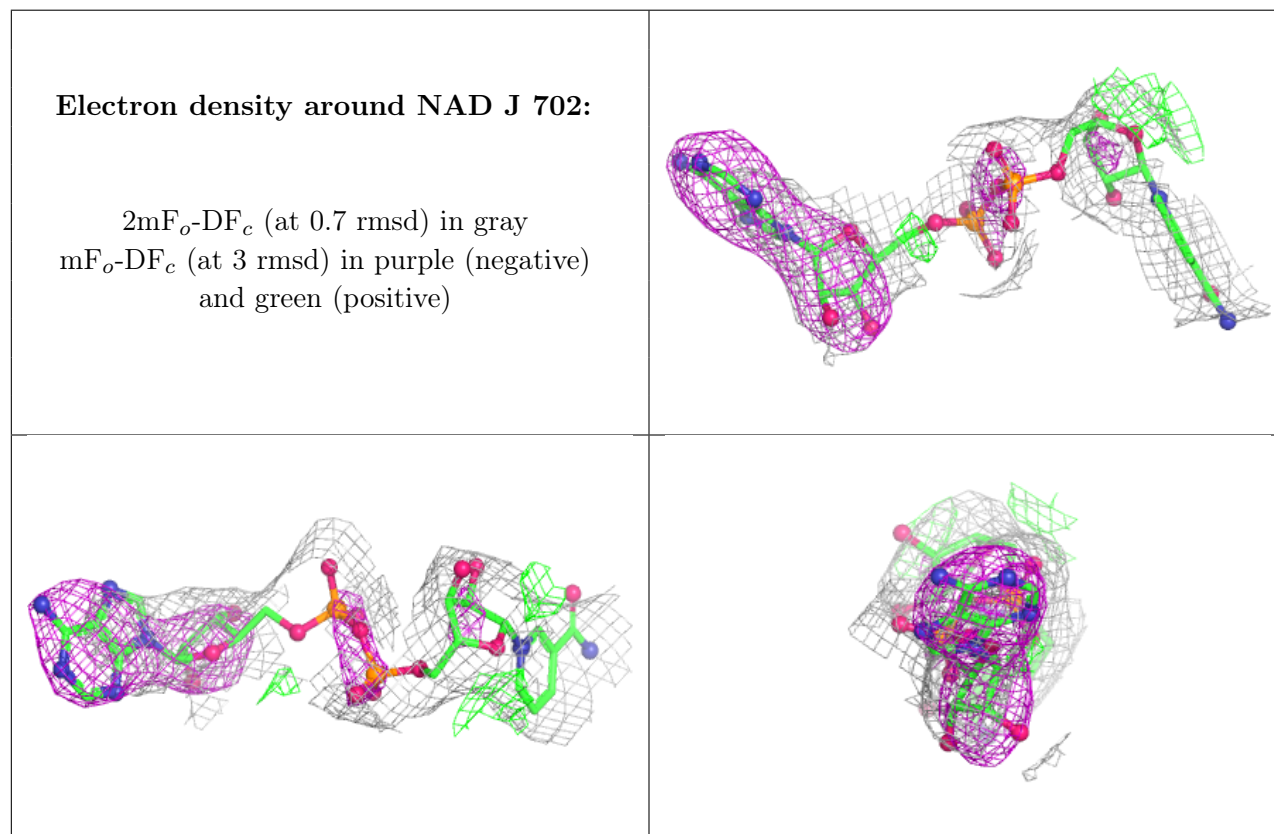
There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

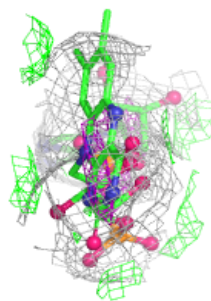
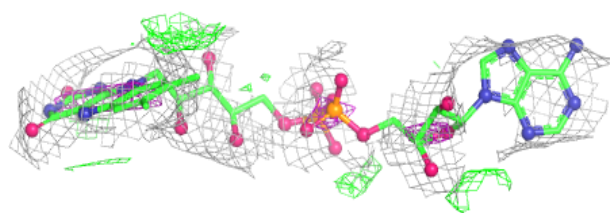
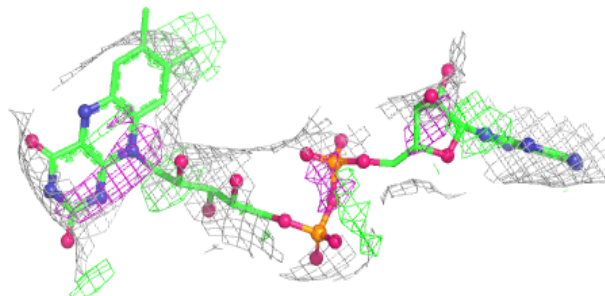
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	NAD	J	702	44/44	0.75	0.12	61,61,61,61	0
3	FAD	J	701	53/53	0.85	0.10	61,61,61,61	0
4	NAD	G	702	44/44	0.88	0.09	61,61,61,61	0
3	FAD	G	701	53/53	0.89	0.09	61,61,61,61	0
4	NAD	D	702	44/44	0.92	0.08	61,61,61,61	0
4	NAD	A	702	44/44	0.93	0.11	61,61,61,61	0
3	FAD	A	701	53/53	0.94	0.12	61,61,61,61	0
3	FAD	D	701	53/53	0.95	0.09	61,61,61,61	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

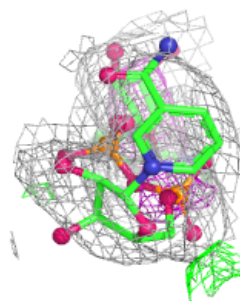
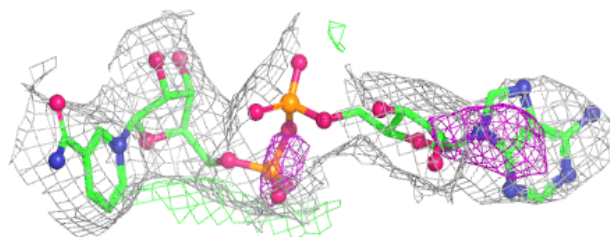
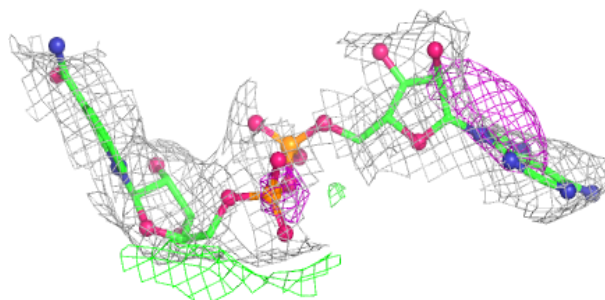


Electron density around FAD J 701:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

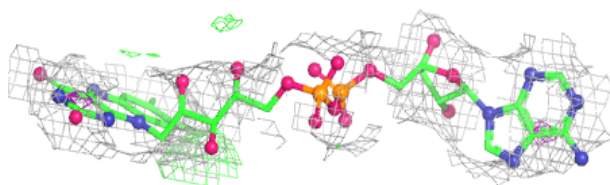
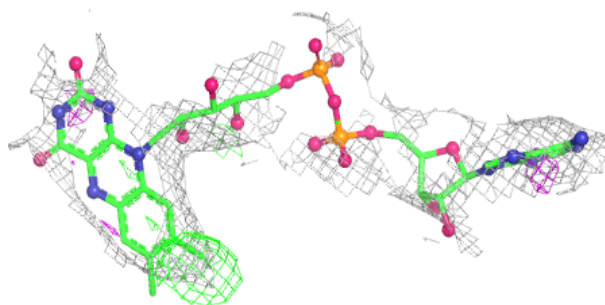
**Electron density around NAD G 702:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

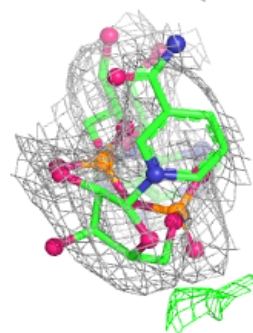
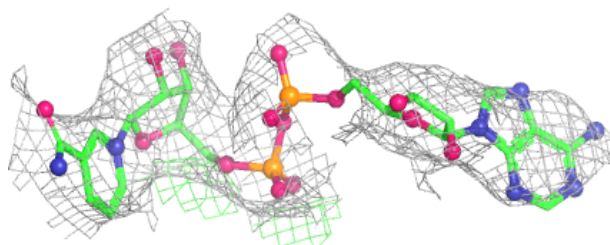
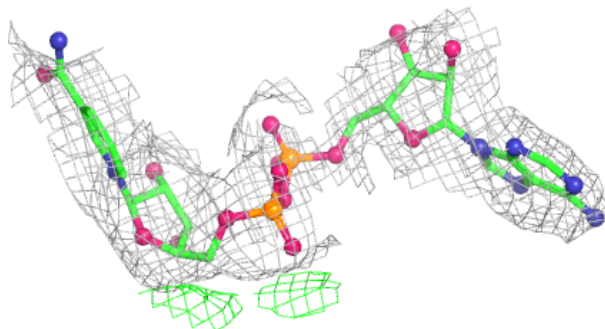


Electron density around FAD G 701:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

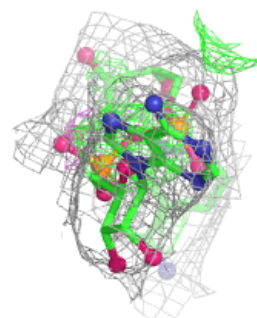
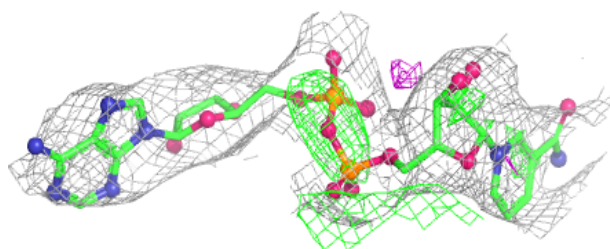
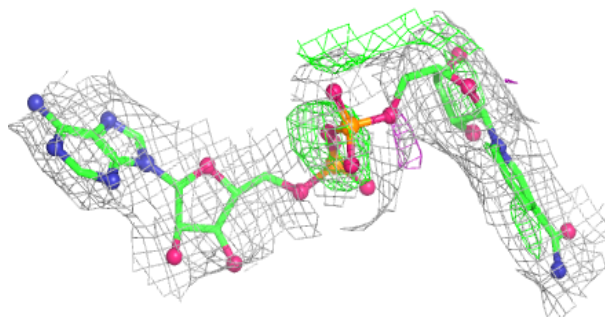
**Electron density around NAD D 702:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

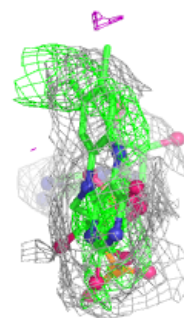
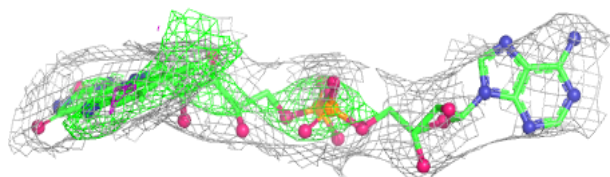
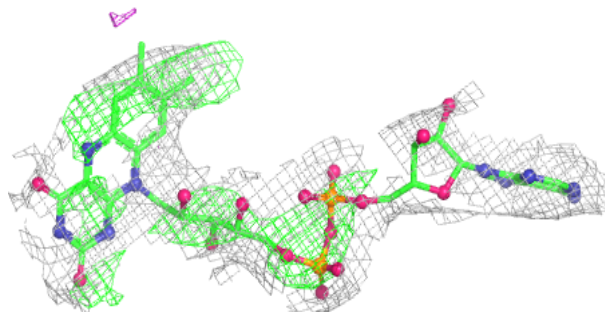


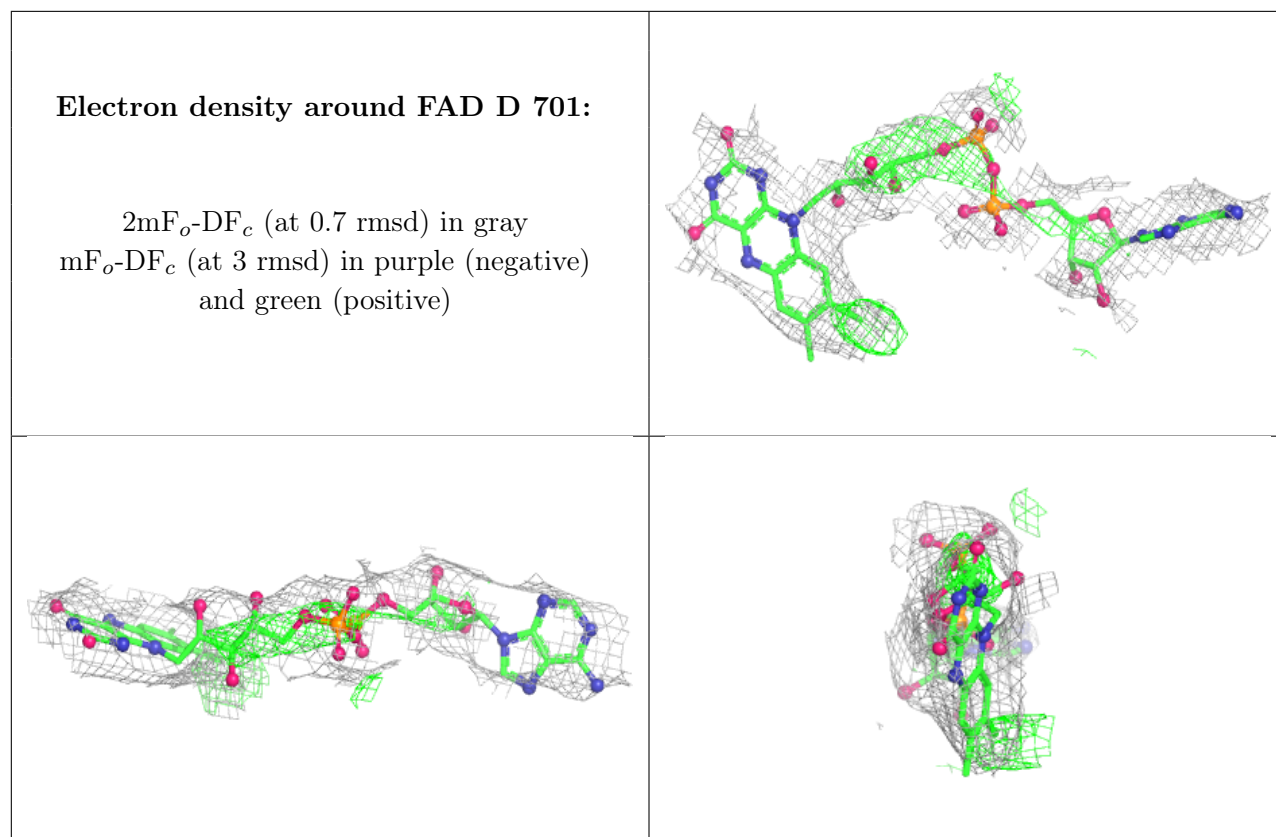
Electron density around NAD A 702:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around FAD A 701:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)





6.5 Other polymers [i](#)

There are no such residues in this entry.