



Full wwPDB EM Validation Report ⓘ

Mar 20, 2026 – 12:35 AM UTC

PDB ID : 8YLT / pdb_00008ylt
EMDB ID : EMD-39390
Title : The structure of DSR2 and NAD⁺ complex
Authors : Zheng, J.; Yang, X.
Deposited on : 2024-03-06
Resolution : 3.09 Å(reported)

This is a Full wwPDB EM Validation Report for a publicly released PDB entry.

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : **NOT EXECUTED**
MapQ : **FAILED**
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

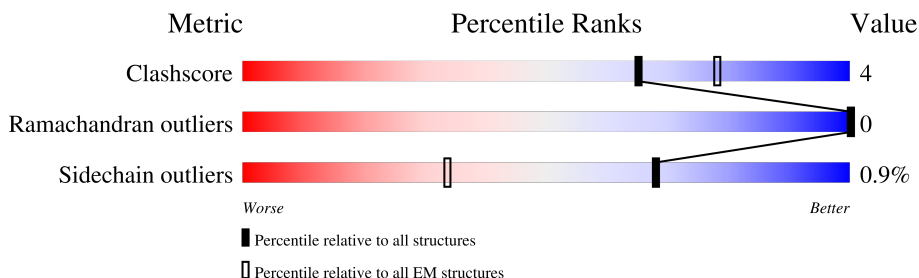
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.09 Å.

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



Metric	Whole archive (#Entries)	EM structures (#Entries)
Clashscore	229148	23984
Ramachandran outliers	224038	23583
Sidechain outliers	223484	23102

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

Mol	Chain	Length	Quality of chain
1	A	1005	
1	B	1005	
1	C	1005	
1	D	1005	

2 Entry composition [i](#)

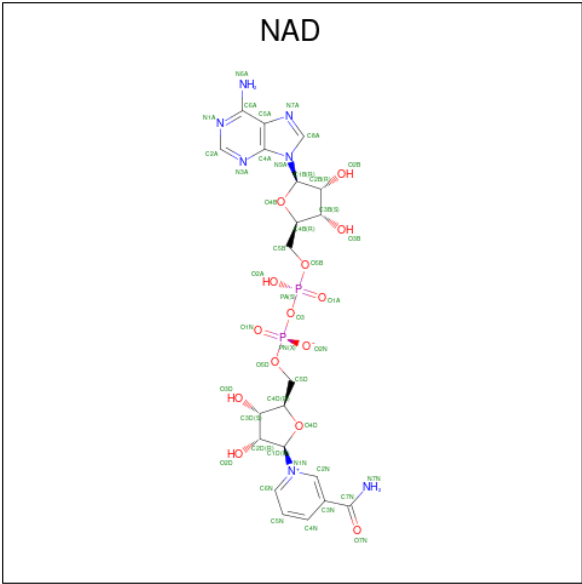
There are 2 unique types of molecules in this entry. The entry contains 65938 atoms, of which 32612 are hydrogens and 0 are deuteriums.

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

- Molecule 1 is a protein called SIR2-like domain-containing protein.

Mol	Chain	Residues	Atoms						AltConf	Trace
1	A	992	Total	C	H	N	O	S	0	0
			16370	5349	8105	1335	1549	32		
1	B	997	Total	C	H	N	O	S	0	0
			16459	5377	8149	1344	1557	32		
1	C	992	Total	C	H	N	O	S	0	0
			16370	5349	8105	1335	1549	32		
1	D	997	Total	C	H	N	O	S	0	0
			16459	5377	8149	1344	1557	32		

- Molecule 2 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (CCD ID: NAD) (formula: C₂₁H₂₇N₇O₁₄P₂) (labeled as "Ligand of Interest" by depositor).



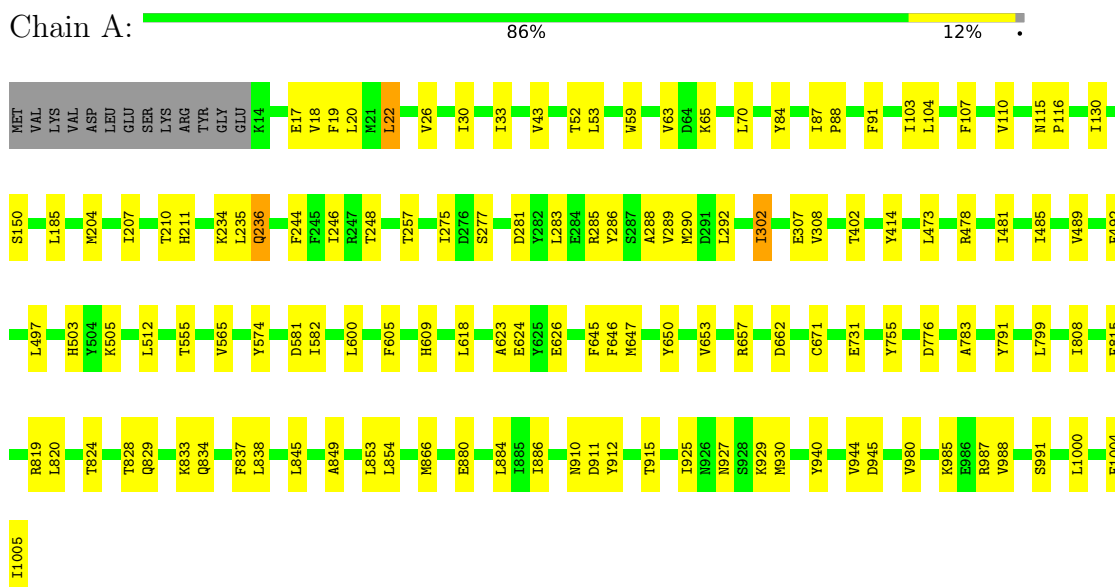
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Mol	Chain	Residues	Atoms						AltConf
2	C	1	Total	C	H	N	O	P	0
			70	21	26	7	14	2	
2	D	1	Total	C	H	N	O	P	0
			70	21	26	7	14	2	

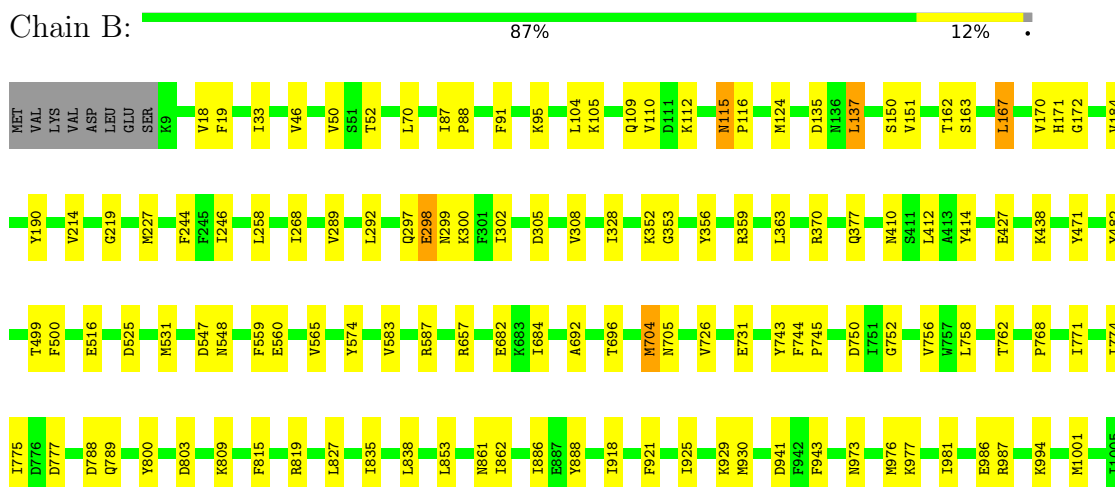
3 Residue-property plots

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. 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.

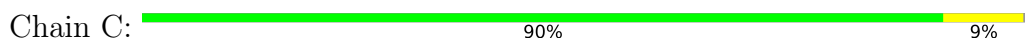
- Molecule 1: SIR2-like domain-containing protein

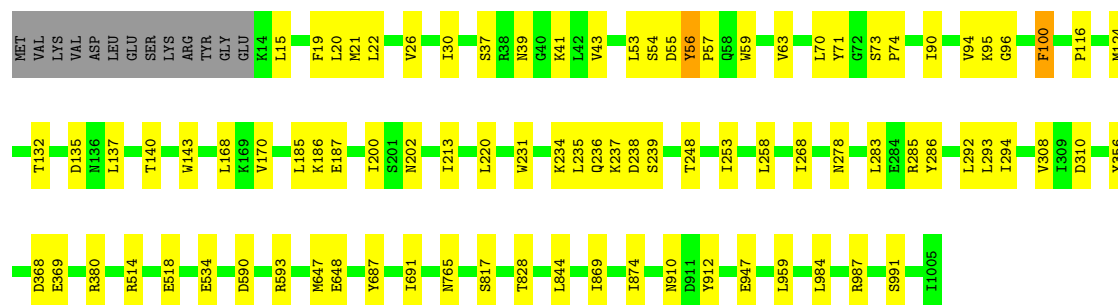


- Molecule 1: SIR2-like domain-containing protein



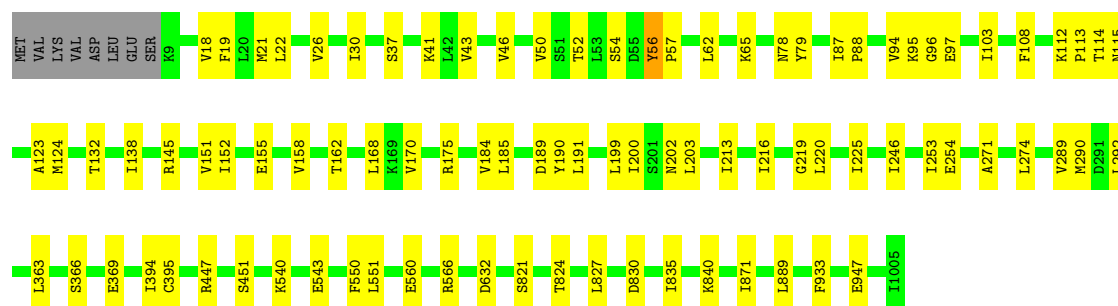
- Molecule 1: SIR2-like domain-containing protein





• Molecule 1: SIR2-like domain-containing protein

Chain D: 90% 9% .



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	259267	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	FEI TITAN	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	40	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	Not provided	
Image detector	FEI FALCON IV (4k x 4k)	Depositor

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: NAD

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	0.50	1/8458 (0.0%)	0.60	3/11391 (0.0%)
1	B	0.53	0/8504	0.60	8/11451 (0.1%)
1	C	0.43	0/8458	0.54	6/11391 (0.1%)
1	D	0.47	0/8504	0.54	6/11451 (0.1%)
All	All	0.49	1/33924 (0.0%)	0.57	23/45684 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	234	LYS	CA-C	-5.40	1.46	1.52

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	297	GLN	N-CA-C	-8.46	102.00	111.14
1	B	18	VAL	N-CA-C	-8.00	102.75	110.42
1	A	52	THR	N-CA-C	-7.14	104.38	113.23
1	B	219	GLY	N-CA-C	-6.82	100.86	112.51
1	B	19	PHE	N-CA-C	-6.56	101.25	110.50
1	D	138	ILE	N-CA-C	-6.54	104.12	110.72
1	B	292	LEU	N-CA-C	-6.50	104.28	111.82
1	C	63	VAL	N-CA-C	-6.46	104.03	110.62
1	D	96	GLY	N-CA-C	6.38	118.52	112.08
1	B	52	THR	N-CA-C	-6.36	104.78	112.54
1	C	124	MET	N-CA-C	-6.12	105.80	113.20
1	D	52	THR	N-CA-C	-6.03	105.02	112.38
1	D	219	GLY	N-CA-C	-6.00	103.10	112.58
1	A	302	ILE	N-CA-C	5.99	115.46	106.42
1	B	298	GLU	N-CA-C	-5.72	102.98	110.53
1	C	96	GLY	N-CA-C	5.50	119.88	112.17

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	53	LEU	N-CA-C	-5.30	105.62	111.71
1	C	100	PHE	N-CA-C	-5.26	105.44	111.07
1	A	286	TYR	N-CA-C	-5.21	105.69	111.36
1	B	172	GLY	N-CA-C	5.20	121.94	112.37
1	D	220	LEU	N-CA-C	-5.05	107.12	113.28
1	C	56	TYR	N-CA-C	-5.05	103.35	110.31
1	D	56	TYR	N-CA-C	-5.03	103.25	109.64

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	8265	8105	8104	87	0
1	B	8310	8149	8148	80	0
1	C	8265	8105	8104	55	0
1	D	8310	8149	8148	54	0
2	A	44	26	26	0	0
2	B	44	26	26	0	0
2	C	44	26	26	0	0
2	D	44	26	26	0	0
All	All	33326	32612	32608	266	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 4.

All (266) 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:489:VAL:HG21	1:A:512:LEU:HD21	1.56	0.87
1:C:59:TRP:CZ2	1:C:185:LEU:HD13	2.11	0.85
1:C:100:PHE:HZ	1:C:185:LEU:HD11	1.47	0.79
1:A:565:VAL:HG21	1:A:618:LEU:HD11	1.67	0.77
1:A:820:LEU:O	1:A:824:THR:HG23	1.88	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:427:GLU:O	1:B:438:LYS:NZ	2.22	0.71
1:C:987:ARG:NH2	1:D:632:ASP:OD2	2.25	0.70
1:A:824:THR:HG22	1:A:838:LEU:HD13	1.74	0.70
1:B:377:GLN:N	1:B:377:GLN:OE1	2.26	0.69
1:A:246:ILE:CG2	1:A:289:VAL:HG11	2.25	0.67
1:A:302:ILE:HD11	1:A:308:VAL:HA	1.76	0.66
1:C:185:LEU:C	1:C:185:LEU:HD12	2.21	0.65
1:A:565:VAL:CG2	1:A:618:LEU:HD11	2.27	0.64
1:C:100:PHE:CZ	1:C:185:LEU:HD11	2.32	0.64
1:B:758:LEU:O	1:B:762:THR:HG23	1.99	0.63
1:A:1000:LEU:HD21	1:A:1005:ILE:HD12	1.80	0.62
1:B:414:TYR:O	1:B:657:ARG:NH2	2.32	0.62
1:B:862:ILE:HG21	1:B:888:TYR:HB3	1.80	0.62
1:C:59:TRP:HZ2	1:C:185:LEU:HD13	1.63	0.62
1:A:277:SER:OG	1:A:281:ASP:HB2	1.99	0.62
1:A:911:ASP:OD1	1:A:915:THR:HG23	2.00	0.62
1:B:986:GLU:OE2	1:B:987:ARG:NH1	2.33	0.61
1:C:310:ASP:OD1	1:C:380:ARG:NH1	2.33	0.61
1:A:731:GLU:OE2	1:A:731:GLU:N	2.33	0.61
1:A:824:THR:HG21	1:A:845:LEU:HD22	1.81	0.61
1:C:30:ILE:HG23	1:C:294:ILE:HA	1.83	0.61
1:A:555:THR:HG22	1:B:559:PHE:CZ	2.37	0.60
1:D:158:VAL:HG11	1:D:199:LEU:HD13	1.84	0.60
1:D:246:ILE:HG23	1:D:289:VAL:HG11	1.83	0.60
1:D:132:THR:HG22	1:D:170:VAL:HG23	1.84	0.59
1:C:73:SER:HB2	1:C:74:PRO:HD2	1.84	0.59
1:C:59:TRP:CH2	1:C:185:LEU:HD13	2.36	0.59
1:A:414:TYR:O	1:A:657:ARG:NH2	2.35	0.59
1:D:19:PHE:HA	1:D:22:LEU:HG	1.85	0.59
1:C:116:PRO:HG2	1:C:283:LEU:HD21	1.85	0.59
1:C:947:GLU:N	1:C:947:GLU:OE1	2.34	0.58
1:C:234:LYS:O	1:C:235:LEU:C	2.46	0.58
1:D:26:VAL:HG21	1:D:274:LEU:HD21	1.85	0.58
1:D:947:GLU:N	1:D:947:GLU:OE1	2.37	0.58
1:A:246:ILE:HG23	1:A:289:VAL:HG11	1.86	0.58
1:C:22:LEU:HD21	1:C:293:LEU:HD21	1.87	0.57
1:C:185:LEU:HD12	1:C:186:LYS:N	2.19	0.57
1:B:33:ILE:HD11	1:B:244:PHE:CD1	2.39	0.57
1:B:921:PHE:CZ	1:B:976:MET:HE1	2.40	0.57
1:C:987:ARG:O	1:C:991:SER:N	2.36	0.57
1:A:574:TYR:OH	1:B:994:LYS:NZ	2.37	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:70:LEU:HD23	1:B:95:LYS:HE2	1.87	0.56
1:A:624:GLU:HG2	1:A:671:CYS:HB3	1.88	0.56
1:A:290:MET:HE2	1:A:290:MET:HA	1.88	0.56
1:B:731:GLU:N	1:B:731:GLU:OE2	2.39	0.56
1:A:497:LEU:HD21	1:A:503:HIS:O	2.05	0.56
1:A:581:ASP:OD1	1:A:582:ILE:N	2.39	0.55
1:A:33:ILE:HD11	1:A:244:PHE:CD1	2.42	0.55
1:A:854:LEU:HD12	1:A:854:LEU:C	2.31	0.55
1:D:540:LYS:NZ	1:D:543:GLU:OE1	2.40	0.55
1:B:471:TYR:HD1	1:B:531:MET:HE1	1.71	0.55
1:C:237:LYS:HD2	1:D:155:GLU:HG3	1.89	0.55
1:A:824:THR:CG2	1:A:845:LEU:HD22	2.37	0.55
1:D:560:GLU:O	1:D:566:ARG:NH1	2.35	0.55
1:B:777:ASP:OD1	1:B:819:ARG:NH2	2.40	0.55
1:B:918:ILE:HG23	1:B:943:PHE:CE2	2.42	0.54
1:A:783:ALA:HB2	1:A:837:PHE:HE2	1.73	0.54
1:C:687:TYR:CZ	1:C:691:ILE:HD11	2.43	0.54
1:B:471:TYR:CD1	1:B:531:MET:HE1	2.43	0.54
1:C:22:LEU:HD22	1:C:26:VAL:HG11	1.89	0.54
1:D:274:LEU:HD12	1:D:289:VAL:HG22	1.90	0.54
1:B:886:ILE:HD11	1:B:930:MET:HE3	1.89	0.54
1:A:246:ILE:HG21	1:A:289:VAL:HG11	1.90	0.54
1:C:22:LEU:HD11	1:C:293:LEU:HG	1.91	0.53
1:D:271:ALA:HB2	1:D:289:VAL:HG23	1.91	0.53
1:A:834:GLN:O	1:A:838:LEU:HG	2.09	0.53
1:A:940:TYR:CE2	1:A:944:VAL:HG21	2.45	0.52
1:A:107:PHE:O	1:A:110:VAL:HG12	2.08	0.52
1:A:402:THR:HG23	1:A:402:THR:O	2.08	0.52
1:D:363:LEU:HD23	1:D:369:GLU:HB3	1.90	0.52
1:C:817:SER:OG	1:C:844:LEU:O	2.26	0.52
1:A:988:VAL:CG1	1:B:1001:MET:HE1	2.39	0.52
1:B:827:LEU:HD21	1:B:835:ILE:HG12	1.91	0.51
1:C:647:MET:SD	1:C:648:GLU:N	2.83	0.51
1:B:684:ILE:HG21	1:B:726:VAL:HG21	1.91	0.51
1:D:184:VAL:HG12	1:D:190:TYR:CE1	2.45	0.51
1:A:70:LEU:HD21	1:A:91:PHE:HA	1.92	0.51
1:A:987:ARG:O	1:A:991:SER:N	2.43	0.51
1:A:489:VAL:HG22	1:A:505:LYS:HG2	1.93	0.51
1:A:22:LEU:HD21	1:A:30:ILE:HD12	1.93	0.51
1:A:985:LYS:HG3	1:B:1001:MET:HE2	1.93	0.51
1:B:482:TYR:OH	1:B:516:GLU:OE1	2.22	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:862:ILE:HG21	1:B:888:TYR:CB	2.40	0.51
1:A:828:THR:HG22	1:A:829:GLN:H	1.77	0.50
1:D:447:ARG:O	1:D:451:SER:OG	2.29	0.50
1:A:18:VAL:HG21	1:A:275:ILE:HD11	1.93	0.50
1:D:821:SER:O	1:D:824:THR:OG1	2.22	0.50
1:B:696:THR:HG22	1:B:743:TYR:CG	2.46	0.50
1:B:886:ILE:HD12	1:B:929:LYS:HG3	1.94	0.50
1:D:112:LYS:HB3	1:D:113:PRO:HD2	1.94	0.50
1:B:328:ILE:HD11	1:B:587:ARG:CD	2.41	0.50
1:B:258:LEU:HD13	1:B:268:ILE:HD11	1.93	0.49
1:B:838:LEU:HB3	1:B:853:LEU:HD21	1.94	0.49
1:B:328:ILE:HD11	1:B:587:ARG:HD3	1.94	0.49
1:B:827:LEU:HD21	1:B:835:ILE:CG1	2.42	0.49
1:D:22:LEU:CD2	1:D:274:LEU:HD22	2.42	0.49
1:B:800:TYR:N	1:B:803:ASP:OD2	2.39	0.49
1:B:930:MET:SD	1:B:930:MET:N	2.84	0.49
1:A:492:PHE:CG	1:A:497:LEU:HD22	2.48	0.49
1:A:605:PHE:O	1:A:609:HIS:ND1	2.43	0.49
1:D:78:ASN:O	1:D:79:TYR:C	2.55	0.49
1:A:485:ILE:O	1:A:489:VAL:HG23	2.12	0.49
1:B:692:ALA:O	1:B:696:THR:HG23	2.12	0.49
1:A:19:PHE:CE1	1:A:292:LEU:HD22	2.48	0.49
1:A:647:MET:N	1:A:647:MET:HE2	2.28	0.49
1:B:768:PRO:HD2	1:B:771:ILE:HD11	1.94	0.49
1:B:925:ILE:HG21	1:B:930:MET:HE2	1.95	0.49
1:C:19:PHE:HA	1:C:22:LEU:HG	1.94	0.49
1:A:911:ASP:OD1	1:A:911:ASP:O	2.31	0.49
1:B:305:ASP:OD2	1:B:359:ARG:NH1	2.46	0.49
1:B:771:ILE:HG22	1:B:774:ILE:HD12	1.95	0.49
1:C:94:VAL:HG23	1:C:95:LYS:HG2	1.95	0.48
1:C:248:THR:HG21	1:C:286:TYR:CZ	2.48	0.48
1:A:646:PHE:C	1:A:647:MET:HE2	2.37	0.48
1:D:132:THR:HG22	1:D:170:VAL:CG2	2.43	0.48
1:B:135:ASP:HB2	1:B:137:LEU:HD13	1.96	0.48
1:C:70:LEU:HD23	1:C:71:TYR:CE2	2.47	0.48
1:D:22:LEU:HD22	1:D:26:VAL:HG11	1.95	0.48
1:C:253:ILE:HG12	1:C:258:LEU:HG	1.95	0.48
1:D:62:LEU:CD1	1:D:108:PHE:CZ	2.97	0.48
1:D:184:VAL:HG13	1:D:189:ASP:HB3	1.96	0.48
1:B:33:ILE:HD12	1:B:214:VAL:CG2	2.44	0.47
1:C:237:LYS:O	1:C:238:ASP:C	2.56	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:46:VAL:HG13	1:B:50:VAL:HG21	1.95	0.47
1:C:90:ILE:O	1:C:94:VAL:HG22	2.14	0.47
1:D:123:ALA:HA	1:D:145:ARG:NH2	2.30	0.47
1:A:824:THR:OG1	1:A:849:ALA:HB1	2.15	0.47
1:A:910:ASN:HB3	1:A:912:TYR:CZ	2.49	0.47
1:B:300:LYS:HA	1:B:353:GLY:HA3	1.97	0.47
1:A:662:ASP:HB2	1:B:565:VAL:HG12	1.95	0.47
1:D:830:ASP:N	1:D:830:ASP:OD1	2.47	0.47
1:D:21:MET:O	1:D:22:LEU:C	2.56	0.47
1:A:791:TYR:O	1:A:833:LYS:NZ	2.37	0.47
1:A:277:SER:OG	1:A:281:ASP:CB	2.63	0.47
1:A:302:ILE:HD11	1:A:308:VAL:HG22	1.97	0.47
1:B:162:THR:HG23	1:B:163:SER:N	2.29	0.47
1:B:171:HIS:HE2	1:B:190:TYR:HE1	1.62	0.47
1:C:140:THR:HA	1:C:143:TRP:CE3	2.50	0.47
1:A:945:ASP:OD1	1:A:945:ASP:N	2.48	0.46
1:D:123:ALA:HA	1:D:145:ARG:HH21	1.80	0.46
1:D:366:SER:OG	1:D:369:GLU:OE1	2.32	0.46
1:B:744:PHE:CD1	1:B:745:PRO:HD2	2.50	0.46
1:C:368:ASP:OD1	1:C:369:GLU:N	2.48	0.46
1:D:46:VAL:HG13	1:D:50:VAL:HG21	1.98	0.46
1:A:980:VAL:HG11	1:A:1004:PHE:CZ	2.51	0.46
1:D:216:ILE:CD1	1:D:290:MET:HE3	2.46	0.46
1:B:171:HIS:CD2	1:B:184:VAL:HB	2.51	0.46
1:D:271:ALA:HB2	1:D:289:VAL:CG2	2.45	0.46
1:A:755:TYR:HD2	1:A:799:LEU:HD13	1.80	0.46
1:B:704:MET:O	1:B:705:ASN:C	2.58	0.46
1:C:70:LEU:HD11	1:C:94:VAL:HG21	1.97	0.46
1:B:246:ILE:HG23	1:B:289:VAL:HG11	1.97	0.45
1:C:308:VAL:HG13	1:C:356:TYR:HB2	1.97	0.45
1:D:87:ILE:HB	1:D:88:PRO:HD3	1.99	0.45
1:D:394:ILE:HG22	1:D:395:CYS:H	1.81	0.45
1:C:248:THR:HG21	1:C:286:TYR:CE1	2.51	0.45
1:A:248:THR:HB	1:A:285:ARG:HG3	1.99	0.45
1:C:135:ASP:HB2	1:C:137:LEU:HD13	1.99	0.45
1:C:534:GLU:OE1	1:D:162:THR:HG21	2.16	0.45
1:A:1000:LEU:CD2	1:A:1005:ILE:HD12	2.45	0.45
1:B:560:GLU:N	1:B:560:GLU:OE1	2.50	0.45
1:A:776:ASP:OD2	1:A:819:ARG:NH2	2.50	0.45
1:B:87:ILE:HB	1:B:88:PRO:HD3	1.99	0.45
1:D:550:PHE:CD2	1:D:551:LEU:HG	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:302:ILE:HG23	1:A:302:ILE:O	2.17	0.44
1:A:53:LEU:HD22	1:A:283:LEU:HG	1.98	0.44
1:A:275:ILE:HD13	1:A:288:ALA:HB1	1.99	0.44
1:A:43:VAL:HG21	1:A:207:ILE:CG2	2.47	0.44
1:A:489:VAL:HG21	1:A:512:LEU:CD2	2.38	0.44
1:A:837:PHE:CD2	1:A:838:LEU:HD23	2.53	0.44
1:B:299:ASN:HB3	1:B:302:ILE:HD12	2.00	0.44
1:B:788:ASP:OD1	1:B:789:GLN:N	2.51	0.44
1:C:168:LEU:HD11	1:C:200:ILE:HG23	1.97	0.44
1:C:590:ASP:OD1	1:C:593:ARG:NH2	2.51	0.44
1:D:216:ILE:HD12	1:D:290:MET:HE3	1.99	0.44
1:D:827:LEU:HD21	1:D:835:ILE:HD13	1.99	0.44
1:A:880:GLU:O	1:A:884:LEU:HG	2.17	0.44
1:C:21:MET:O	1:C:22:LEU:C	2.59	0.44
1:D:152:ILE:HB	1:D:168:LEU:HD12	1.98	0.44
1:A:22:LEU:HG	1:A:26:VAL:HB	2.00	0.43
1:B:110:VAL:HG21	1:B:112:LYS:HE2	1.99	0.43
1:B:151:VAL:HG12	1:B:167:LEU:HB3	1.99	0.43
1:D:94:VAL:HG23	1:D:95:LYS:HG2	2.00	0.43
1:B:758:LEU:HD21	1:B:775:ILE:HD11	1.99	0.43
1:A:65:LYS:HG2	1:A:103:ILE:HG23	2.00	0.43
1:B:115:ASN:HB2	1:B:116:PRO:CD	2.48	0.43
1:B:300:LYS:HE2	1:B:352:LYS:HB3	2.00	0.43
1:B:33:ILE:HD11	1:B:244:PHE:CG	2.53	0.43
1:A:302:ILE:HA	1:A:307:GLU:OE1	2.18	0.43
1:D:65:LYS:HG2	1:D:103:ILE:HG23	2.00	0.43
1:A:478:ARG:HA	1:A:481:ILE:HG22	2.01	0.43
1:A:925:ILE:HD12	1:A:930:MET:CE	2.49	0.43
1:B:574:TYR:CE1	1:B:583:VAL:HG11	2.54	0.43
1:A:988:VAL:HG13	1:B:1001:MET:HE1	2.00	0.43
1:C:235:LEU:O	1:C:236:GLN:C	2.61	0.43
1:A:87:ILE:HB	1:A:88:PRO:HD3	1.99	0.42
1:A:845:LEU:CD1	1:A:853:LEU:HD22	2.49	0.42
1:B:682:GLU:OE1	1:B:682:GLU:N	2.51	0.42
1:C:43:VAL:HB	1:C:213:ILE:HD13	2.01	0.42
1:D:22:LEU:HD21	1:D:292:LEU:HD13	2.01	0.42
1:D:22:LEU:HD23	1:D:274:LEU:HD22	2.00	0.42
1:B:412:LEU:HD22	1:B:412:LEU:H	1.83	0.42
1:B:574:TYR:HE1	1:B:583:VAL:HG11	1.83	0.42
1:C:54:SER:O	1:C:55:ASP:C	2.60	0.42
1:D:168:LEU:HD13	1:D:203:LEU:HD23	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:115:ASN:HB2	1:A:116:PRO:CD	2.50	0.42
1:B:308:VAL:HG13	1:B:356:TYR:HB2	2.01	0.42
1:C:15:LEU:HG	1:C:292:LEU:HB3	2.01	0.42
1:C:910:ASN:HB2	1:C:912:TYR:CD2	2.53	0.42
1:B:977:LYS:O	1:B:981:ILE:HG12	2.19	0.42
1:D:22:LEU:HD13	1:D:30:ILE:HD12	2.01	0.42
1:B:499:THR:HG23	1:B:500:PHE:CD1	2.55	0.42
1:B:809:LYS:NZ	1:B:815:PHE:O	2.45	0.42
1:D:889:LEU:HD22	1:D:933:PHE:CD2	2.54	0.42
1:B:973:ASN:CB	1:B:976:MET:HE2	2.50	0.42
1:C:56:TYR:O	1:C:57:PRO:C	2.62	0.42
1:A:17:GLU:HA	1:A:20:LEU:HD12	2.01	0.42
1:A:104:LEU:HD21	1:A:185:LEU:HD13	2.02	0.42
1:A:623:ALA:HA	1:A:645:PHE:HB3	2.01	0.42
1:A:63:VAL:HG21	1:A:84:TYR:HD1	1.85	0.42
1:B:752:GLY:O	1:B:756:VAL:HG13	2.20	0.42
1:D:54:SER:N	1:D:115:ASN:HD21	2.17	0.42
1:D:253:ILE:O	1:D:254:GLU:C	2.62	0.42
1:A:210:THR:HG23	1:A:211:HIS:ND1	2.34	0.42
1:C:70:LEU:HD21	1:C:90:ILE:CG2	2.50	0.42
1:C:869:ILE:HG13	1:C:874:ILE:HD11	2.02	0.42
1:C:514:ARG:NE	1:C:518:GLU:OE2	2.53	0.41
1:A:130:ILE:HG21	1:A:204:MET:HG3	2.02	0.41
1:A:555:THR:HG22	1:B:559:PHE:HZ	1.80	0.41
1:A:473:LEU:HD22	1:A:600:LEU:HD21	2.01	0.41
1:B:359:ARG:O	1:B:359:ARG:HD3	2.20	0.41
1:C:959:LEU:HD22	1:C:984:LEU:HD21	2.03	0.41
1:D:18:VAL:HG12	1:D:292:LEU:HD21	2.02	0.41
1:D:271:ALA:CB	1:D:289:VAL:HG23	2.50	0.41
1:A:886:ILE:HD12	1:A:929:LYS:CB	2.51	0.41
1:B:91:PHE:CE2	1:B:95:LYS:HG3	2.55	0.41
1:B:170:VAL:O	1:B:171:HIS:HB2	2.21	0.41
1:C:39:ASN:N	1:C:39:ASN:HD22	2.18	0.41
1:A:59:TRP:O	1:A:63:VAL:HG23	2.21	0.41
1:D:168:LEU:HD11	1:D:200:ILE:HG23	2.02	0.41
1:C:73:SER:HB2	1:C:74:PRO:CD	2.51	0.41
1:C:278:ASN:O	1:C:285:ARG:NH2	2.54	0.41
1:A:235:LEU:O	1:A:236:GLN:C	2.61	0.41
1:B:525:ASP:OD1	1:B:525:ASP:N	2.52	0.41
1:B:750:ASP:OD1	1:B:752:GLY:N	2.54	0.41
1:C:132:THR:HG22	1:C:170:VAL:CG2	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:220:LEU:HD21	1:C:268:ILE:HD12	2.02	0.41
1:D:43:VAL:HB	1:D:213:ILE:HD13	2.02	0.41
1:A:650:TYR:HA	1:A:653:VAL:HG12	2.03	0.41
1:B:547:ASP:OD1	1:B:548:ASN:N	2.54	0.41
1:D:56:TYR:O	1:D:57:PRO:C	2.62	0.41
1:B:363:LEU:HD23	1:B:370:ARG:HG2	2.03	0.40
1:B:941:ASP:N	1:B:941:ASP:OD1	2.53	0.40
1:B:105:LYS:HG3	1:B:109:GLN:HB2	2.02	0.40
1:A:808:ILE:HG21	1:A:815:PHE:CD2	2.57	0.40
1:C:231:TRP:HA	1:C:234:LYS:HD3	2.04	0.40
1:A:866:MET:HE1	1:A:912:TYR:CD1	2.56	0.40
1:A:927:ASN:HB3	1:A:930:MET:HG2	2.03	0.40
1:B:973:ASN:HB3	1:B:976:MET:HE2	2.03	0.40
1:D:151:VAL:HG21	1:D:175:ARG:NH2	2.37	0.40
1:D:840:LYS:HD2	1:D:871:ILE:HG23	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	990/1005 (98%)	962 (97%)	28 (3%)	0	100	100
1	B	995/1005 (99%)	963 (97%)	32 (3%)	0	100	100
1	C	990/1005 (98%)	964 (97%)	26 (3%)	0	100	100
1	D	995/1005 (99%)	972 (98%)	23 (2%)	0	100	100
All	All	3970/4020 (99%)	3861 (97%)	109 (3%)	0	100	100

There are no Ramachandran outliers to report.

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 PDB entries followed by that with respect to all EM entries.

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	911/923 (99%)	906 (100%)	5 (0%)	81	85
1	B	915/923 (99%)	904 (99%)	11 (1%)	63	78
1	C	911/923 (99%)	903 (99%)	8 (1%)	70	80
1	D	915/923 (99%)	906 (99%)	9 (1%)	68	79
All	All	3652/3692 (99%)	3619 (99%)	33 (1%)	68	80

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	LEU
1	A	150	SER
1	A	236	GLN
1	A	257	THR
1	A	626	GLU
1	B	104	LEU
1	B	115	ASN
1	B	124	MET
1	B	137	LEU
1	B	150	SER
1	B	167	LEU
1	B	227	MET
1	B	298	GLU
1	B	410	ASN
1	B	704	MET
1	B	861	ASN
1	C	20	LEU
1	C	37	SER
1	C	41	LYS
1	C	187	GLU
1	C	202	ASN
1	C	239	SER
1	C	765	ASN
1	C	828	THR

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Mol	Chain	Res	Type
1	D	37	SER
1	D	41	LYS
1	D	97	GLU
1	D	114	THR
1	D	124	MET
1	D	185	LEU
1	D	191	LEU
1	D	202	ASN
1	D	225	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (47) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	202	ASN
1	A	263	ASN
1	A	391	ASN
1	A	614	ASN
1	A	814	ASN
1	A	975	HIS
1	B	24	ASN
1	B	109	GLN
1	B	136	ASN
1	B	192	ASN
1	B	202	ASN
1	B	224	ASN
1	B	349	HIS
1	B	391	ASN
1	B	410	ASN
1	B	529	ASN
1	B	563	ASN
1	B	606	HIS
1	B	782	GLN
1	B	810	HIS
1	B	926	ASN
1	B	992	ASN
1	B	1002	ASN
1	C	182	ASN
1	C	202	ASN
1	C	297	GLN
1	C	339	HIS
1	C	477	ASN
1	C	549	GLN

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Mol	Chain	Res	Type
1	C	598	ASN
1	C	614	ASN
1	C	666	ASN
1	D	25	ASN
1	D	89	GLN
1	D	115	ASN
1	D	195	GLN
1	D	339	HIS
1	D	349	HIS
1	D	829	GLN
1	D	867	ASN
1	D	895	ASN
1	D	926	ASN
1	D	939	GLN
1	D	948	ASN
1	D	961	ASN
1	D	973	ASN
1	D	990	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](#)

4 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	NAD	C	1101	-	46,48,48	1.16	6 (13%)	64,73,73	1.57	11 (17%)
2	NAD	B	1101	-	46,48,48	1.13	6 (13%)	64,73,73	1.57	10 (15%)
2	NAD	D	1101	-	46,48,48	1.18	5 (10%)	64,73,73	1.60	11 (17%)
2	NAD	A	1101	-	46,48,48	1.16	5 (10%)	64,73,73	1.68	12 (18%)

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	NAD	C	1101	-	-	7/30/62/62	0/5/5/5
2	NAD	B	1101	-	-	8/30/62/62	0/5/5/5
2	NAD	D	1101	-	-	14/30/62/62	0/5/5/5
2	NAD	A	1101	-	-	11/30/62/62	0/5/5/5

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	1101	NAD	C5A-C4A	4.36	1.46	1.39
2	D	1101	NAD	C5A-C4A	4.22	1.46	1.39
2	A	1101	NAD	C5A-C4A	4.13	1.46	1.39
2	B	1101	NAD	C5A-C4A	3.92	1.46	1.39
2	B	1101	NAD	C5A-N7A	-2.75	1.34	1.39
2	D	1101	NAD	C5A-N7A	-2.66	1.34	1.39
2	C	1101	NAD	C5A-C6A	2.52	1.48	1.41
2	A	1101	NAD	C5A-C6A	2.44	1.47	1.41
2	D	1101	NAD	C5A-C6A	2.41	1.47	1.41
2	A	1101	NAD	C5A-N7A	-2.39	1.34	1.39
2	C	1101	NAD	C8A-N7A	2.38	1.36	1.31
2	B	1101	NAD	C5A-C6A	2.34	1.47	1.41
2	C	1101	NAD	C5A-N7A	-2.33	1.34	1.39
2	A	1101	NAD	C8A-N7A	2.29	1.36	1.31
2	A	1101	NAD	C4A-N9A	-2.26	1.33	1.37
2	B	1101	NAD	C4A-N9A	-2.25	1.33	1.37
2	C	1101	NAD	C4A-N9A	-2.22	1.33	1.37
2	D	1101	NAD	C8A-N7A	2.11	1.35	1.31
2	D	1101	NAD	C4A-N9A	-2.11	1.33	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	1101	NAD	O4D-C1D	2.10	1.43	1.40
2	B	1101	NAD	O4D-C1D	2.02	1.43	1.40
2	B	1101	NAD	C8A-N7A	2.01	1.35	1.31

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	1101	NAD	C5A-C4A-N3A	-5.69	118.88	126.72
2	B	1101	NAD	C5A-C4A-N3A	-5.57	119.05	126.72
2	C	1101	NAD	C5A-C4A-N3A	-5.45	119.21	126.72
2	A	1101	NAD	C5A-C4A-N3A	-5.42	119.25	126.72
2	D	1101	NAD	N3A-C4A-N9A	4.79	135.32	127.17
2	B	1101	NAD	N3A-C4A-N9A	4.41	134.66	127.17
2	A	1101	NAD	N3A-C4A-N9A	4.31	134.50	127.17
2	C	1101	NAD	N3A-C4A-N9A	4.22	134.34	127.17
2	A	1101	NAD	C2A-N3A-C4A	3.84	121.21	111.83
2	A	1101	NAD	N3A-C2A-N1A	-3.84	122.77	128.58
2	C	1101	NAD	C2A-N3A-C4A	3.81	121.14	111.83
2	B	1101	NAD	C2A-N3A-C4A	3.76	121.02	111.83
2	C	1101	NAD	N3A-C2A-N1A	-3.76	122.89	128.58
2	D	1101	NAD	C2A-N3A-C4A	3.74	120.95	111.83
2	A	1101	NAD	O4B-C1B-N9A	3.66	115.11	108.09
2	B	1101	NAD	N3A-C2A-N1A	-3.60	123.13	128.58
2	D	1101	NAD	N3A-C2A-N1A	-3.55	123.20	128.58
2	C	1101	NAD	C4A-C5A-N7A	-3.49	106.59	110.58
2	A	1101	NAD	C4A-C5A-N7A	-3.47	106.62	110.58
2	B	1101	NAD	C4A-C5A-N7A	-3.27	106.84	110.58
2	A	1101	NAD	C4A-N9A-C8A	3.23	109.13	105.74
2	D	1101	NAD	C4A-N9A-C8A	3.12	109.01	105.74
2	D	1101	NAD	C4A-C5A-N7A	-3.05	107.10	110.58
2	C	1101	NAD	C4A-N9A-C8A	2.98	108.87	105.74
2	B	1101	NAD	C4A-N9A-C8A	2.83	108.71	105.74
2	C	1101	NAD	C5A-N7A-C8A	2.66	107.64	103.45
2	B	1101	NAD	C5A-N7A-C8A	2.66	107.63	103.45
2	A	1101	NAD	C5A-N7A-C8A	2.63	107.59	103.45
2	D	1101	NAD	C5A-N7A-C8A	2.50	107.37	103.45
2	A	1101	NAD	N9A-C8A-N7A	-2.49	110.41	113.94
2	A	1101	NAD	C2B-C1B-N9A	-2.48	107.15	113.30
2	C	1101	NAD	N9A-C8A-N7A	-2.45	110.46	113.94
2	A	1101	NAD	C6A-C5A-N7A	2.45	136.81	132.09
2	B	1101	NAD	N9A-C8A-N7A	-2.40	110.53	113.94
2	B	1101	NAD	C3N-C7N-N7N	2.38	120.67	117.74

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1101	NAD	C6A-C5A-N7A	2.37	136.66	132.09
2	D	1101	NAD	N9A-C8A-N7A	-2.27	110.71	113.94
2	D	1101	NAD	C3B-C2B-C1B	2.24	105.70	101.46
2	A	1101	NAD	C2A-N1A-C6A	2.12	122.21	118.73
2	D	1101	NAD	O4B-C1B-N9A	2.07	112.06	108.09
2	C	1101	NAD	C2A-N1A-C6A	2.05	122.09	118.73
2	D	1101	NAD	C2B-C1B-N9A	-2.02	108.29	113.30
2	C	1101	NAD	C5N-C4N-C3N	-2.01	118.39	120.36
2	B	1101	NAD	C6A-C5A-N7A	2.00	135.95	132.09

There are no chirality outliers.

All (40) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1101	NAD	C5B-O5B-PA-O2A
2	A	1101	NAD	PN-O3-PA-O5B
2	A	1101	NAD	C5D-O5D-PN-O3
2	A	1101	NAD	C5D-O5D-PN-O1N
2	A	1101	NAD	C5D-O5D-PN-O2N
2	B	1101	NAD	C5B-O5B-PA-O3
2	C	1101	NAD	C5B-O5B-PA-O1A
2	D	1101	NAD	C5B-O5B-PA-O2A
2	D	1101	NAD	C5B-O5B-PA-O3
2	D	1101	NAD	O4B-C4B-C5B-O5B
2	D	1101	NAD	C5D-O5D-PN-O3
2	D	1101	NAD	C5D-O5D-PN-O1N
2	D	1101	NAD	C5D-O5D-PN-O2N
2	A	1101	NAD	C2N-C3N-C7N-O7N
2	A	1101	NAD	C2N-C3N-C7N-N7N
2	C	1101	NAD	C2N-C3N-C7N-O7N
2	C	1101	NAD	C2N-C3N-C7N-N7N
2	D	1101	NAD	C2N-C3N-C7N-O7N
2	A	1101	NAD	C4N-C3N-C7N-N7N
2	C	1101	NAD	C4N-C3N-C7N-N7N
2	A	1101	NAD	C4N-C3N-C7N-O7N
2	C	1101	NAD	C4N-C3N-C7N-O7N
2	A	1101	NAD	O4B-C4B-C5B-O5B
2	D	1101	NAD	C2N-C3N-C7N-N7N
2	D	1101	NAD	C4N-C3N-C7N-O7N
2	D	1101	NAD	C3B-C4B-C5B-O5B
2	D	1101	NAD	C4N-C3N-C7N-N7N
2	A	1101	NAD	C3B-C4B-C5B-O5B

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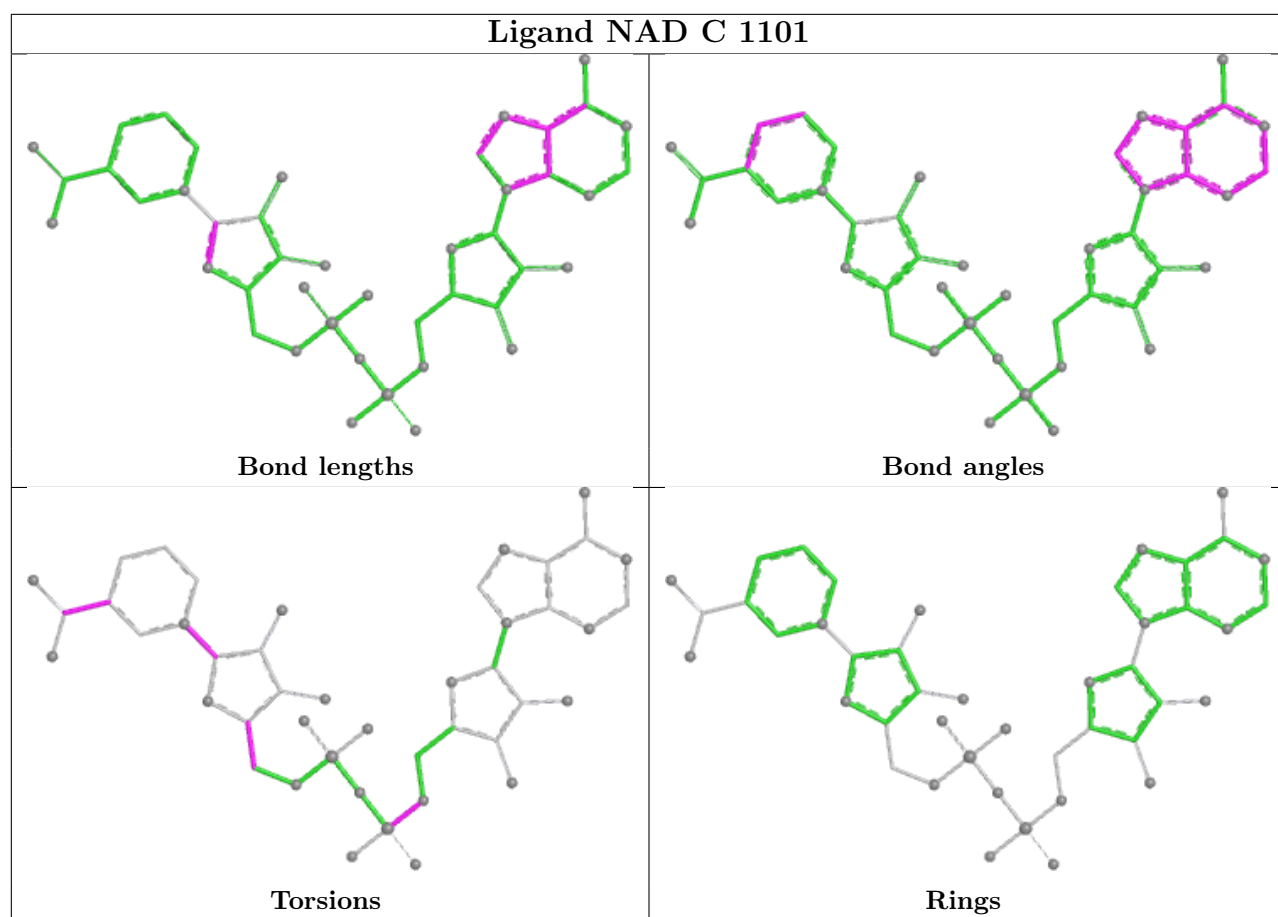
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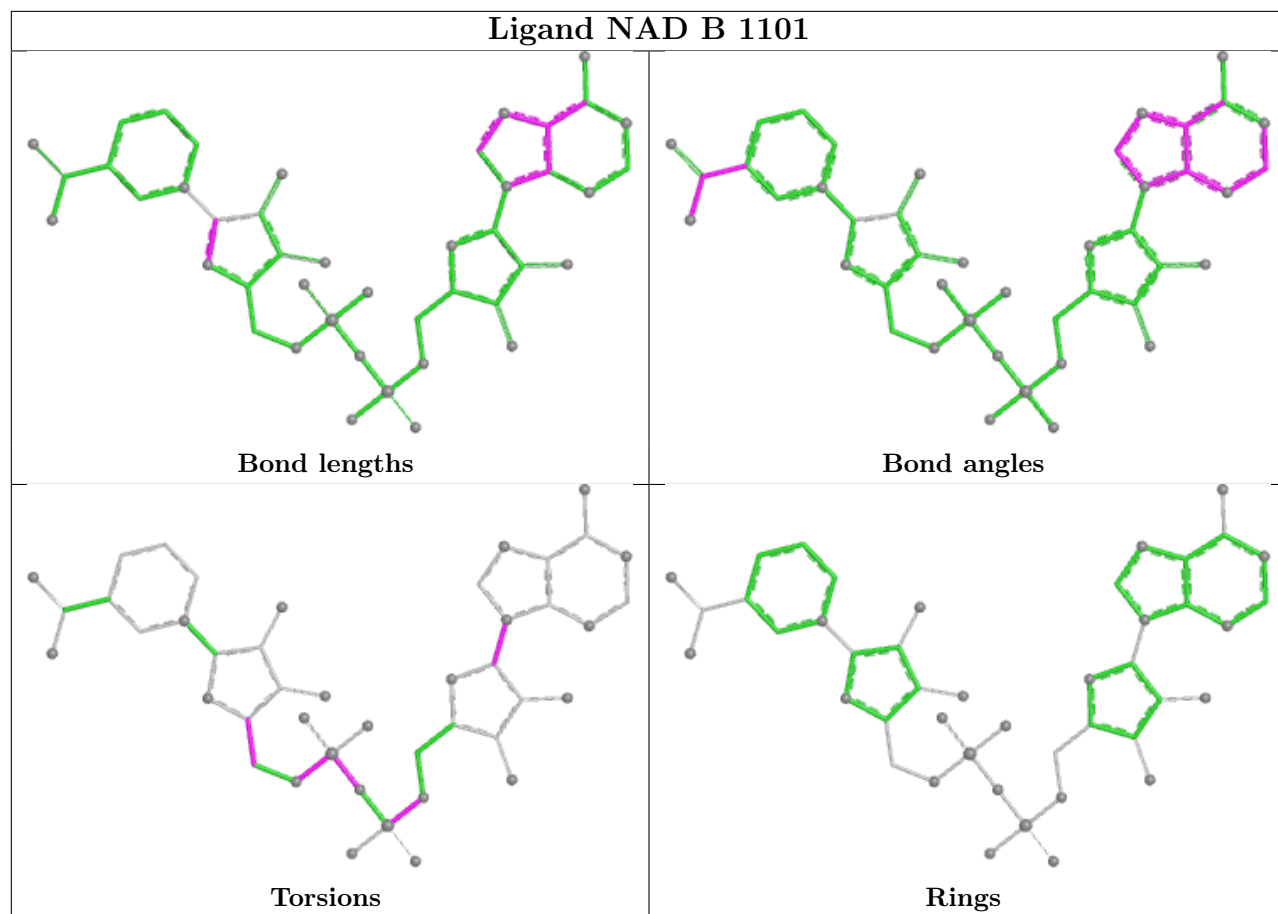
Mol	Chain	Res	Type	Atoms
2	D	1101	NAD	O4D-C4D-C5D-O5D
2	B	1101	NAD	C2B-C1B-N9A-C8A
2	B	1101	NAD	C5B-O5B-PA-O1A
2	B	1101	NAD	C5B-O5B-PA-O2A
2	B	1101	NAD	C5D-O5D-PN-O1N
2	D	1101	NAD	C5B-O5B-PA-O1A
2	C	1101	NAD	O4D-C4D-C5D-O5D
2	C	1101	NAD	O4D-C1D-N1N-C6N
2	B	1101	NAD	O4D-C4D-C5D-O5D
2	D	1101	NAD	C3D-C4D-C5D-O5D
2	B	1101	NAD	O4B-C1B-N9A-C8A
2	B	1101	NAD	PA-O3-PN-O2N

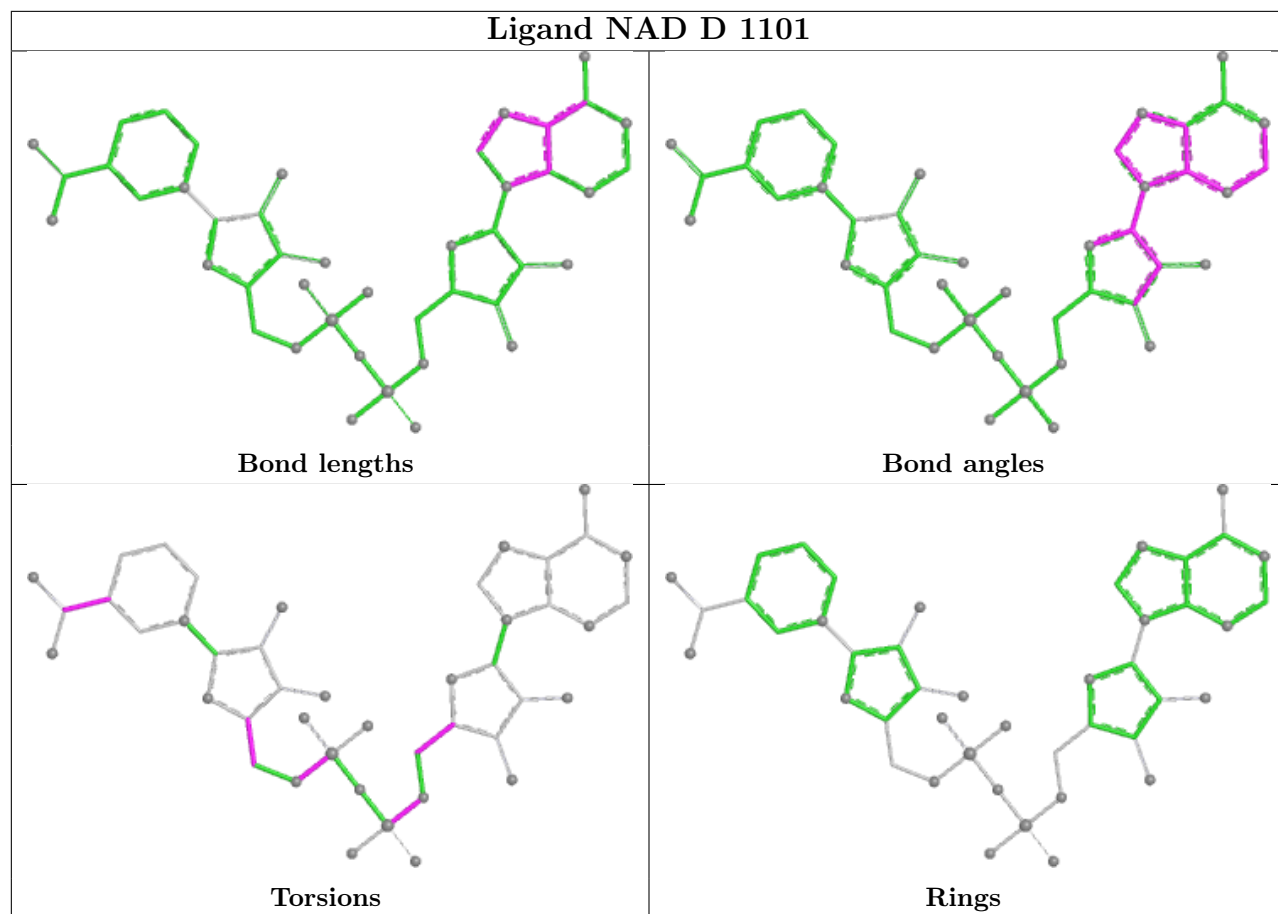
There are no ring outliers.

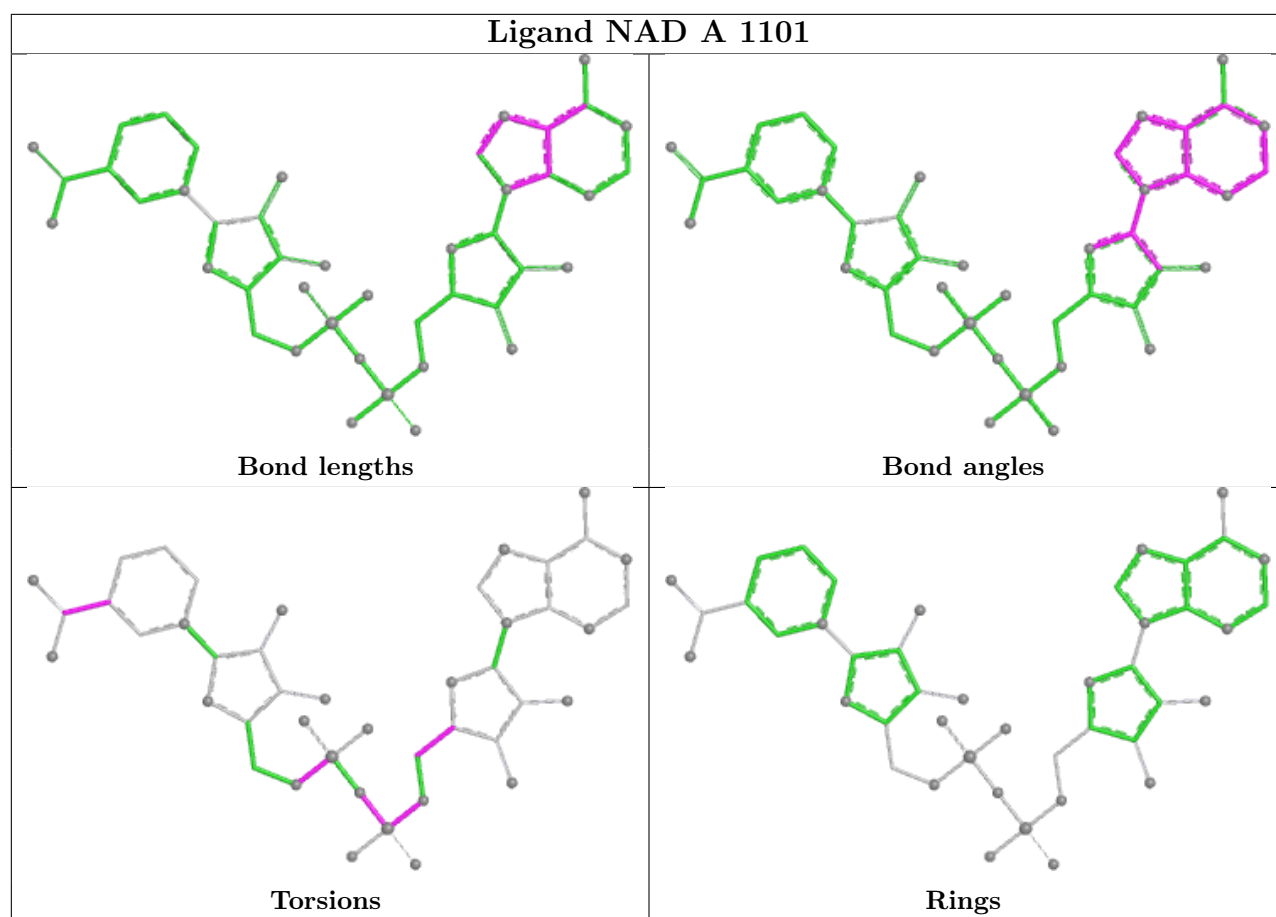
No monomer is involved in short contacts.

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









5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Map visualisation

This section contains visualisations of the EMDB entry EMD-39390. These allow visual inspection of the internal detail of the map and identification of artifacts.

Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections

This section was not generated.

6.2 Central slices

This section was not generated.

6.3 Largest variance slices

This section was not generated.

6.4 Orthogonal standard-deviation projections (False-color)

This section was not generated.

6.5 Orthogonal surface views

This section was not generated.

6.6 Mask visualisation

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

7 Map analysis ⓘ

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

7.1 Map-value distribution ⓘ

This section was not generated.

7.2 Volume estimate versus contour level ⓘ

This section was not generated.

7.3 Rotationally averaged power spectrum ⓘ

This section was not generated. The rotationally averaged power spectrum had issues being displayed.

8 Fourier-Shell correlation ⓘ

This section was not generated. No FSC curve or half-maps provided.

9 Map-model fit

This section was not generated.