



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:34 PM UTC

PDB ID : 1A5Z / pdb_00001a5z
Title : LACTATE DEHYDROGENASE FROM THERMOTOGA MARITIMA (TMLDH)
Authors : Auerbach, G.; Ostendorp, R.; Prade, L.; Korndorfer, I.; Dams, T.; Huber, R.; Jaenicke, R.
Deposited on : 1998-02-18
Resolution : 2.10 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

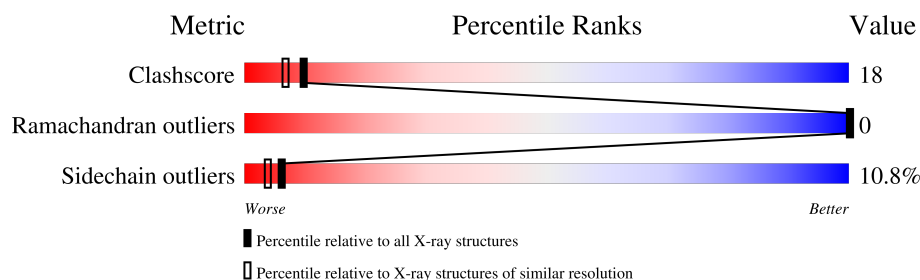
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)

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	319	

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
4	OXM	A	355	-	X	-	-

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 2852 atoms, of which 12 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.

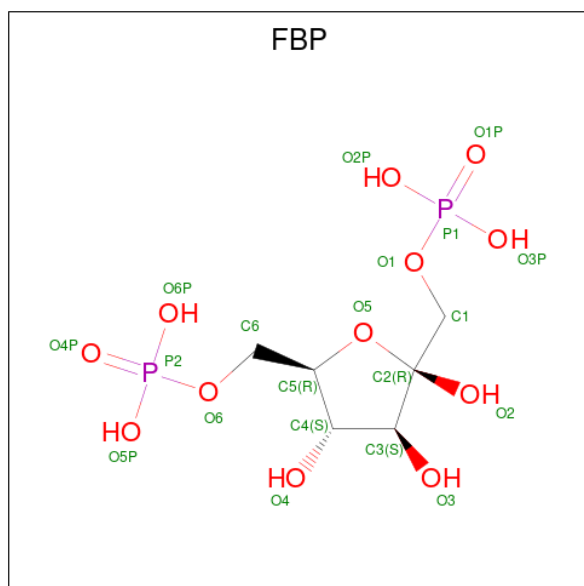
- Molecule 1 is a protein called L-LACTATE DEHYDROGENASE.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	312	2415	1526	12	416	449	12	79	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	215	VAL	ILE	conflict	UNP P16115

- Molecule 2 is 1,6-di-O-phosphono-beta-D-fructofuranose (CCD ID: FBP) (formula: $C_6H_{14}O_{12}P_2$).

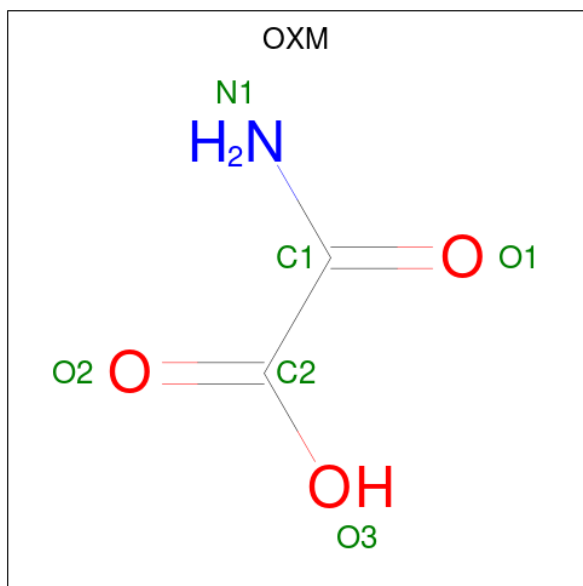


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	O	P		
2	A	1	20	6	12	2	0	0
2	A	1	20	6	12	2	0	0

- Molecule 3 is CADMIUM ION (CCD ID: CD) (formula: Cd).

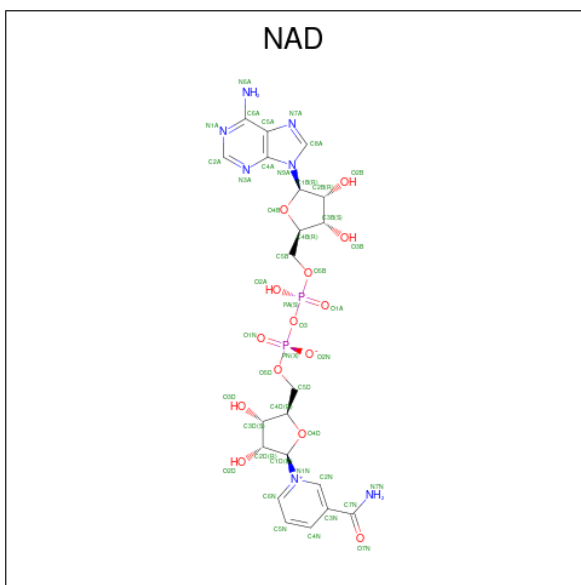
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	3	Total	Cd		0	0
			3	3			

- Molecule 4 is OXAMIC ACID (CCD ID: OXM) (formula: C₂H₃NO₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			6	2	1	3		

- Molecule 5 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (CCD ID: NAD) (formula: C₂₁H₂₇N₇O₁₄P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total 44	C 21	N 7	O 14	P 2	0	0

- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	344	Total O 344 344	0	0

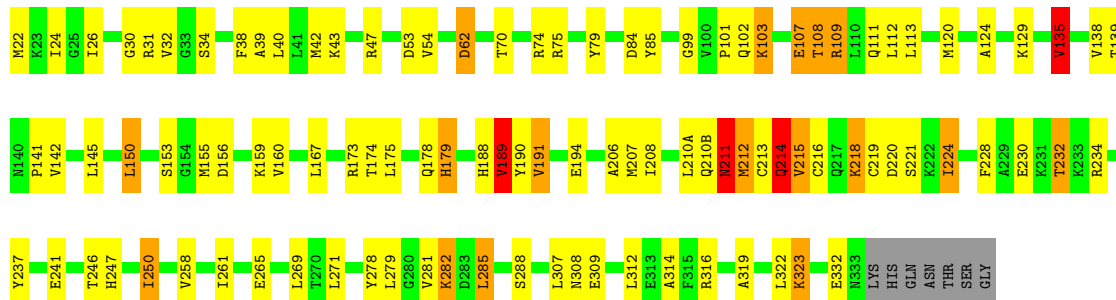
3 Residue-property plots

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: L-LACTATE DEHYDROGENASE

Chain A: 



4 Data and refinement statistics

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

Property	Value	Source
Space group	I 41 2 2	Depositor
Cell constants a, b, c, α , β , γ	105.48Å 105.48Å 187.70Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 2.10	Depositor
% Data completeness (in resolution range)	97.7 (8.00-2.10)	Depositor
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.8	Depositor
R, R_{free}	0.200 , 0.286	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2852	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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

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	0.93	1/2440 (0.0%)	1.29	18/3296 (0.5%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	26	ILE	CA-CB	5.23	1.60	1.54

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	215	VAL	N-CA-C	-21.12	93.40	111.56
1	A	142	VAL	N-CA-C	8.80	119.60	110.62
1	A	70	THR	CA-C-N	-7.59	111.82	119.56
1	A	70	THR	C-N-CA	-7.59	111.82	119.56
1	A	211	ASN	N-CA-C	-6.99	103.27	111.03
1	A	47	ARG	N-CA-C	-6.97	102.76	112.45
1	A	129	LYS	N-CA-C	6.96	118.52	111.07
1	A	189	VAL	CB-CA-C	-6.86	100.04	111.29
1	A	191	VAL	N-CA-CB	6.17	119.64	111.25
1	A	214	GLN	N-CA-C	6.03	119.38	112.57
1	A	85	TYR	N-CA-C	5.99	119.42	111.75
1	A	156	ASP	CA-C-N	5.28	124.99	119.82
1	A	156	ASP	C-N-CA	5.28	124.99	119.82
1	A	261	ILE	N-CA-C	-5.25	105.59	110.53
1	A	265	GLU	N-CA-C	5.22	118.81	112.23
1	A	135	VAL	N-CA-C	5.18	115.42	108.17
1	A	221	SER	N-CA-C	5.15	117.29	111.11
1	A	124	ALA	N-CA-C	5.12	116.55	111.07

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	2403	12	2461	84	1
2	A	40	0	20	4	1
3	A	3	0	0	0	0
4	A	6	0	2	0	0
5	A	44	0	26	8	0
6	A	344	0	0	15	0
All	All	2840	12	2509	84	1

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 18.

All (84) 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:79:TYR:HB2	2:A:354:FBP:O2	1.74	0.87
1:A:30:GLY:HA3	5:A:352:NAD:O1A	1.78	0.82
1:A:246:THR:H	5:A:352:NAD:H72N	1.25	0.81
1:A:103:LYS:O	1:A:107:GLU:HB2	1.86	0.75
1:A:174:THR:O	1:A:178:GLN:HG2	1.87	0.74
1:A:228:PHE:O	1:A:232:THR:HG23	1.86	0.74
1:A:194:GLU:HG3	1:A:322:LEU:HD21	1.73	0.70
1:A:175:LEU:HD23	6:A:557:HOH:O	1.91	0.70
1:A:109:ARG:NH2	1:A:141:PRO:HG3	2.08	0.68
1:A:210(A):LEU:HD12	1:A:210(A):LEU:O	1.94	0.68
1:A:210(A):LEU:O	1:A:213:CYS:HB2	1.95	0.66
1:A:230:GLU:O	1:A:234:ARG:HG3	1.95	0.65
1:A:103:LYS:HB2	1:A:107:GLU:HG3	1.78	0.65
1:A:138:VAL:HG11	1:A:250:ILE:HD11	1.78	0.64
1:A:218:LYS:HZ3	1:A:224:ILE:HD12	1.63	0.63
1:A:282:LYS:HE2	1:A:282:LYS:O	1.98	0.63
1:A:250:ILE:HG12	6:A:638:HOH:O	2.00	0.62
1:A:175:LEU:HA	6:A:557:HOH:O	1.98	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:32:VAL:HG13	1:A:250:ILE:HD12	1.82	0.61
1:A:150:LEU:HB2	6:A:667:HOH:O	1.99	0.61
1:A:237:TYR:O	1:A:241:GLU:HG3	2.01	0.61
1:A:135:VAL:HG13	1:A:160:VAL:HG22	1.84	0.60
1:A:188:HIS:ND1	2:A:353:FBP:O5P	2.31	0.59
1:A:173:ARG:HG2	1:A:189:VAL:HG23	1.84	0.59
1:A:250:ILE:C	1:A:250:ILE:HD13	2.28	0.59
1:A:108:THR:HB	1:A:111:GLN:HG3	1.84	0.59
1:A:211:ASN:O	1:A:215:VAL:HG23	2.02	0.59
1:A:74:ARG:HB2	6:A:656:HOH:O	2.02	0.59
1:A:74:ARG:HG2	1:A:75:ARG:N	2.17	0.58
1:A:53:ASP:OD2	5:A:352:NAD:H1B	2.03	0.58
1:A:99:GLY:O	5:A:352:NAD:H51A	2.05	0.56
1:A:208:ILE:HB	1:A:212:MET:HG2	1.87	0.56
1:A:308:ASN:OD1	1:A:308:ASN:C	2.50	0.55
1:A:234:ARG:NH2	6:A:588:HOH:O	2.34	0.54
1:A:32:VAL:CG1	1:A:250:ILE:HD12	2.38	0.54
1:A:109:ARG:HH22	1:A:141:PRO:HG3	1.72	0.53
1:A:38:PHE:CZ	1:A:42:MET:HE2	2.44	0.53
1:A:210(B):GLN:HB2	6:A:642:HOH:O	2.08	0.53
1:A:282:LYS:H	1:A:282:LYS:CE	2.23	0.52
1:A:38:PHE:CE1	1:A:42:MET:HE2	2.45	0.52
1:A:79:TYR:CD1	2:A:354:FBP:H61	2.46	0.51
1:A:34:SER:HB3	1:A:62:ASP:OD2	2.12	0.50
1:A:153:SER:HB2	1:A:155:MET:HE2	1.94	0.50
1:A:74:ARG:HG2	1:A:75:ARG:O	2.12	0.49
1:A:214:GLN:OE1	1:A:219:CYS:O	2.32	0.48
1:A:22:MET:HE1	1:A:258:VAL:HG13	1.95	0.48
1:A:108:THR:HB	1:A:111:GLN:CG	2.44	0.48
1:A:179:HIS:NE2	1:A:218:LYS:NZ	2.62	0.48
1:A:234:ARG:NH1	6:A:588:HOH:O	2.40	0.48
1:A:314:ALA:HA	6:A:457:HOH:O	2.14	0.47
1:A:282:LYS:CD	1:A:282:LYS:N	2.77	0.47
1:A:101:PRO:HD3	6:A:467:HOH:O	2.13	0.47
1:A:319:ALA:O	1:A:323:LYS:HG3	2.14	0.47
1:A:323:LYS:NZ	1:A:323:LYS:HB3	2.30	0.46
1:A:312:LEU:HG	1:A:316:ARG:NH1	2.32	0.45
1:A:206:ALA:C	1:A:207:MET:HG3	2.42	0.45
1:A:79:TYR:HB2	2:A:354:FBP:HO2	1.75	0.45
1:A:30:GLY:CA	5:A:352:NAD:O1A	2.60	0.45
1:A:102:GLN:HA	1:A:112:LEU:HD13	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:246:THR:N	5:A:352:NAD:H72N	2.03	0.44
1:A:211:ASN:OD1	1:A:211:ASN:N	2.50	0.44
1:A:31:ARG:HH21	1:A:247:HIS:HE1	1.64	0.44
1:A:159:LYS:HE2	1:A:159:LYS:HB3	1.68	0.44
1:A:31:ARG:HH21	1:A:247:HIS:CE1	2.36	0.43
1:A:84:ASP:HB2	6:A:593:HOH:O	2.18	0.43
1:A:323:LYS:HB3	1:A:323:LYS:HZ2	1.83	0.43
1:A:224:ILE:HG22	6:A:681:HOH:O	2.17	0.43
1:A:108:THR:HG22	1:A:111:GLN:H	1.83	0.43
1:A:250:ILE:CG1	6:A:638:HOH:O	2.63	0.42
1:A:278:TYR:O	1:A:281:VAL:HG22	2.18	0.42
1:A:282:LYS:N	1:A:282:LYS:HD3	2.34	0.42
1:A:285:LEU:HD13	1:A:323:LYS:CG	2.48	0.42
1:A:271:LEU:O	1:A:288:SER:HA	2.19	0.42
1:A:22:MET:HA	6:A:492:HOH:O	2.20	0.42
1:A:139:THR:HG23	5:A:352:NAD:C4D	2.50	0.42
1:A:216:CYS:SG	1:A:219:CYS:HB2	2.59	0.42
1:A:39:ALA:O	1:A:43:LYS:HG2	2.20	0.41
1:A:188:HIS:CE1	1:A:190:TYR:CE2	3.09	0.41
1:A:323:LYS:NZ	6:A:544:HOH:O	2.52	0.41
1:A:22:MET:HE3	1:A:24:ILE:HD11	2.02	0.41
1:A:120:MET:HG3	1:A:145:LEU:HB3	2.02	0.41
1:A:139:THR:HG23	5:A:352:NAD:H4D	2.03	0.40
1:A:279:LEU:HD21	1:A:307:LEU:HG	2.03	0.40
1:A:282:LYS:NZ	1:A:323:LYS:HD2	2.36	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:188:HIS:ND1	2:A:353:FBP:C3[10_765]	2.16	0.04

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	310/319 (97%)	291 (94%)	19 (6%)	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 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	259/265 (98%)	231 (89%)	28 (11%)	6	4

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

Mol	Chain	Res	Type
1	A	40	LEU
1	A	54	VAL
1	A	62	ASP
1	A	103	LYS
1	A	107	GLU
1	A	108	THR
1	A	109	ARG
1	A	113	LEU
1	A	135	VAL
1	A	150	LEU
1	A	167	LEU
1	A	179	HIS
1	A	189	VAL
1	A	191	VAL
1	A	211	ASN
1	A	212	MET
1	A	214	GLN
1	A	218	LYS
1	A	220	ASP
1	A	224	ILE
1	A	232	THR

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Mol	Chain	Res	Type
1	A	250	ILE
1	A	269	LEU
1	A	282	LYS
1	A	285	LEU
1	A	309	GLU
1	A	323	LYS
1	A	332	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	77	ASN
1	A	140	ASN
1	A	178	GLN
1	A	195	HIS
1	A	247	HIS

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](#)

Of 7 ligands modelled in this entry, 3 are monoatomic - leaving 4 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	OXM	A	355	-	5,5,5	2.61	1 (20%)	2,6,6	2.11	2 (100%)
2	FBP	A	354	-	18,20,20	1.77	6 (33%)	21,32,32	1.50	5 (23%)
5	NAD	A	352	-	46,48,48	1.62	5 (10%)	64,73,73	1.88	9 (14%)
2	FBP	A	353	-	18,20,20	1.44	3 (16%)	21,32,32	1.58	5 (23%)

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	OXM	A	355	-	-	3/4/4/4	-
2	FBP	A	354	-	-	4/13/32/32	0/1/1/1
5	NAD	A	352	-	-	3/30/62/62	0/5/5/5
2	FBP	A	353	-	-	7/13/32/32	0/1/1/1

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	352	NAD	O4D-C1D	5.57	1.48	1.40
5	A	352	NAD	C2N-N1N	5.30	1.40	1.35
4	A	355	OXM	C1-C2	-5.18	1.49	1.55
5	A	352	NAD	C5A-N7A	-3.41	1.32	1.39
2	A	354	FBP	O5-C5	3.08	1.50	1.43
2	A	354	FBP	P1-O1	-2.67	1.51	1.60
2	A	354	FBP	C4-C5	2.64	1.59	1.53
2	A	354	FBP	P2-O6	-2.54	1.52	1.60
2	A	354	FBP	P1-O2P	-2.45	1.45	1.54
2	A	353	FBP	P2-O6	-2.42	1.52	1.60
5	A	352	NAD	C6N-N1N	2.39	1.40	1.35
2	A	353	FBP	P1-O1	-2.31	1.53	1.60
2	A	354	FBP	O5-C2	2.24	1.47	1.43
5	A	352	NAD	C3N-C7N	2.19	1.53	1.50
2	A	353	FBP	P1-O2P	-2.16	1.46	1.54

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	352	NAD	C5A-C4A-N3A	-7.24	116.74	126.72
5	A	352	NAD	N3A-C4A-N9A	6.83	138.78	127.17

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	352	NAD	C2A-N3A-C4A	3.82	121.15	111.83
5	A	352	NAD	C6A-C5A-C4A	3.65	122.16	117.18
5	A	352	NAD	N3A-C2A-N1A	-3.60	123.13	128.58
5	A	352	NAD	C6A-C5A-N7A	-3.44	125.46	132.09
2	A	353	FBP	C6-C5-C4	-3.31	103.28	115.21
5	A	352	NAD	O5B-C5B-C4B	3.05	119.37	108.99
2	A	354	FBP	O3P-P1-O2P	2.90	118.67	107.80
5	A	352	NAD	O4B-C4B-C5B	-2.85	100.21	109.33
2	A	353	FBP	O3P-P1-O2P	2.75	118.12	107.80
2	A	354	FBP	O5P-P2-O6	-2.73	99.56	106.67
2	A	353	FBP	O6-C6-C5	2.71	118.24	108.99
5	A	352	NAD	O4B-C1B-N9A	2.63	113.15	108.09
2	A	354	FBP	O6P-P2-O6	2.51	113.21	106.67
2	A	354	FBP	O6-C6-C5	2.36	117.03	108.99
2	A	353	FBP	O5-C5-C4	-2.26	99.73	105.42
2	A	353	FBP	O3P-P1-O1	2.24	112.51	106.67
4	A	355	OXM	O3-C2-O2	-2.20	118.66	123.90
2	A	354	FBP	O4-C4-C5	2.02	116.89	111.08
4	A	355	OXM	O1-C1-N1	-2.01	118.39	122.89

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	353	FBP	C1-O1-P1-O2P
2	A	353	FBP	C1-O1-P1-O3P
2	A	353	FBP	C5-C6-O6-P2
2	A	353	FBP	C6-O6-P2-O4P
2	A	353	FBP	C6-O6-P2-O5P
2	A	354	FBP	C6-O6-P2-O4P
2	A	354	FBP	C6-O6-P2-O5P
2	A	354	FBP	C6-O6-P2-O6P
4	A	355	OXM	N1-C1-C2-O2
4	A	355	OXM	N1-C1-C2-O3
5	A	352	NAD	C4B-C5B-O5B-PA
2	A	354	FBP	C5-C6-O6-P2
2	A	353	FBP	C1-O1-P1-O1P
4	A	355	OXM	O1-C1-C2-O2
5	A	352	NAD	O4D-C4D-C5D-O5D
2	A	353	FBP	C6-O6-P2-O6P
5	A	352	NAD	C3D-C4D-C5D-O5D

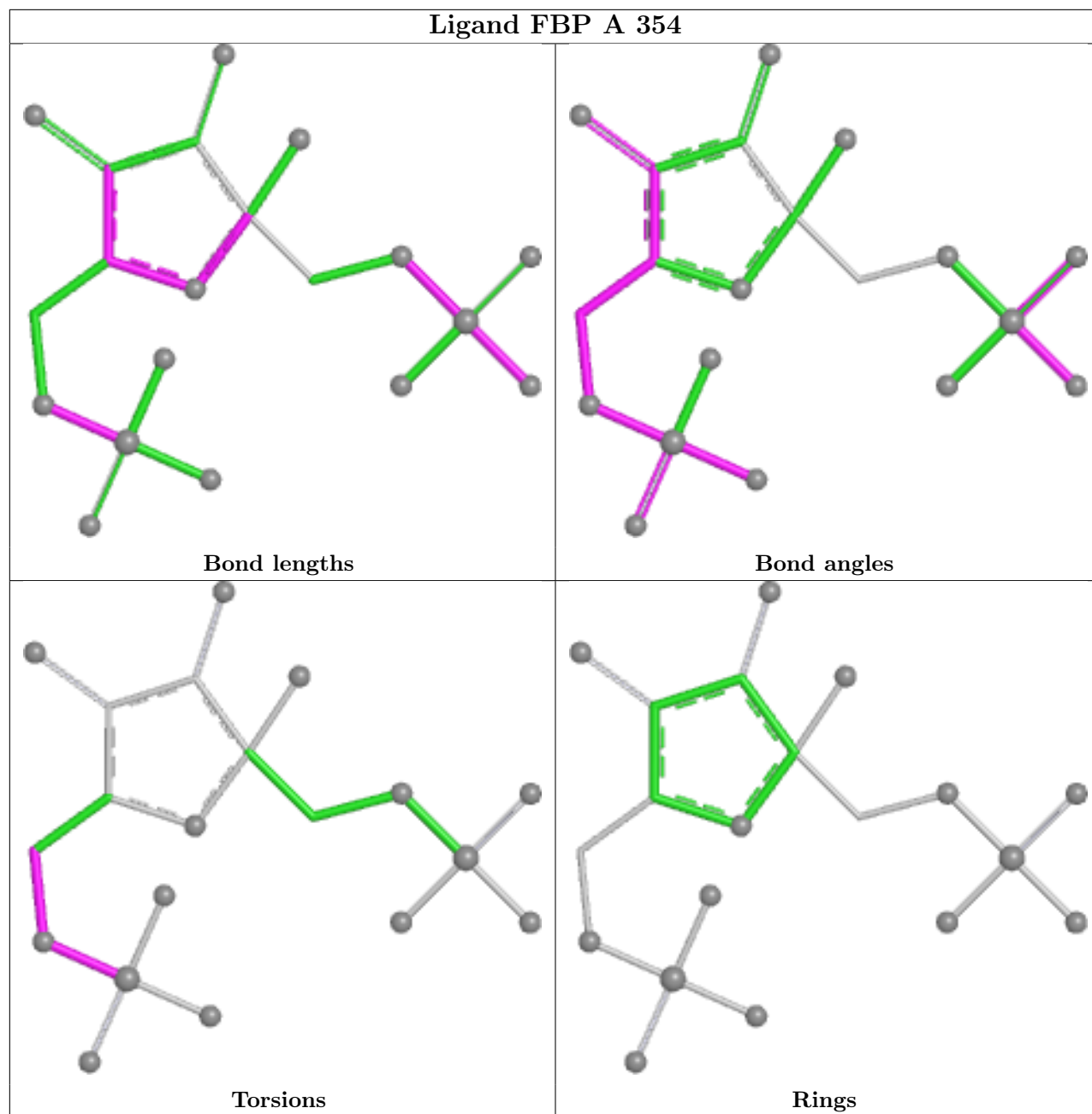
There are no ring outliers.

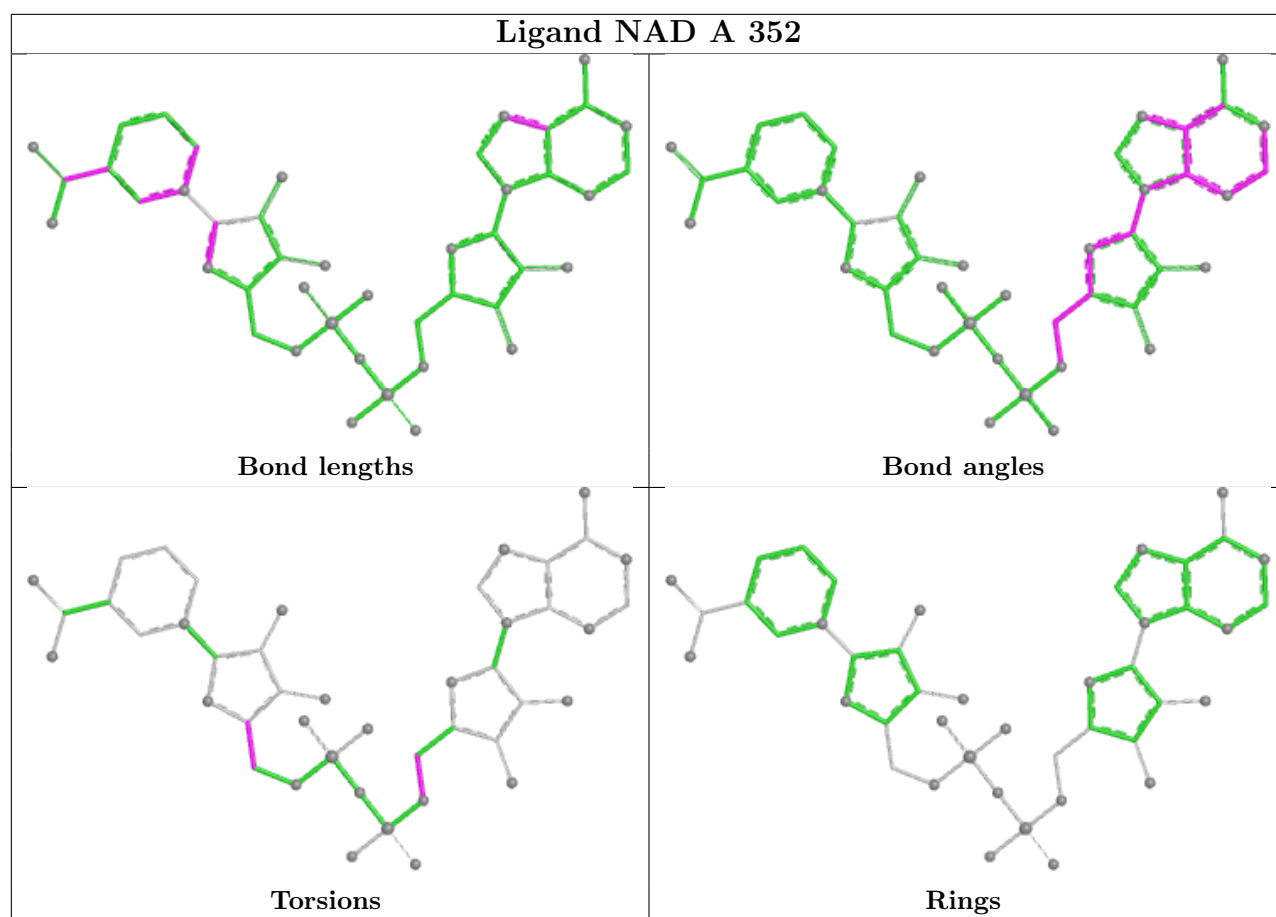
3 monomers are involved in 13 short contacts:

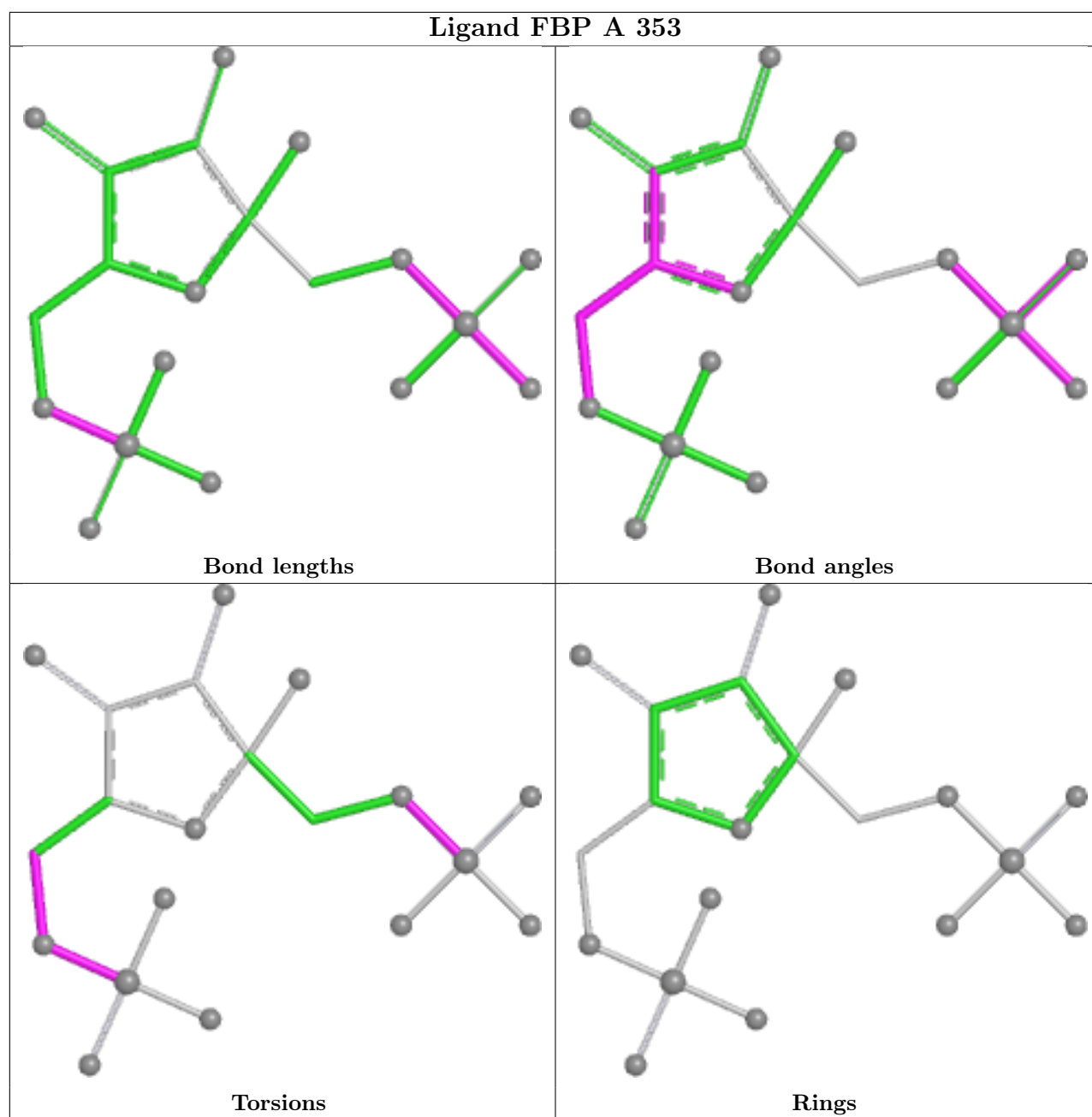
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	354	FBP	3	0
5	A	352	NAD	8	0
2	A	353	FBP	1	1

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 FBP A 354







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.