



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 10:56 PM UTC

PDB ID : 3ZOK / pdb_00003zok
Title : Structure of 3-Dehydroquinate Synthase from Actinidia chinensis in complex with NAD
Authors : Mittelstaedt, G.; Negron, L.; Schofield, L.R.; Marsh, K.; Parker, E.J.
Deposited on : 2013-02-22
Resolution : 2.40 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

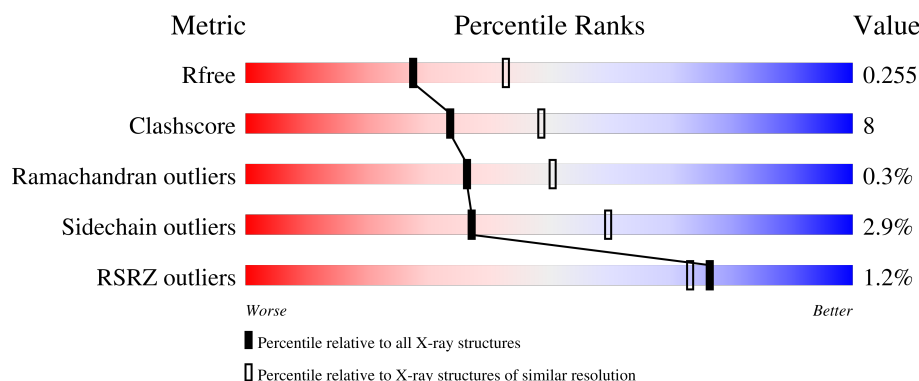
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	378	0% 78% 17% ..
1	B	378	0% 79% 17% ..
1	C	378	2% 78% 17% ..
1	D	378	84% 12% ..

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
2	PO4	B	1379	-	-	X	-
5	GLY	B	1385	-	X	-	-

2 Entry composition [i](#)

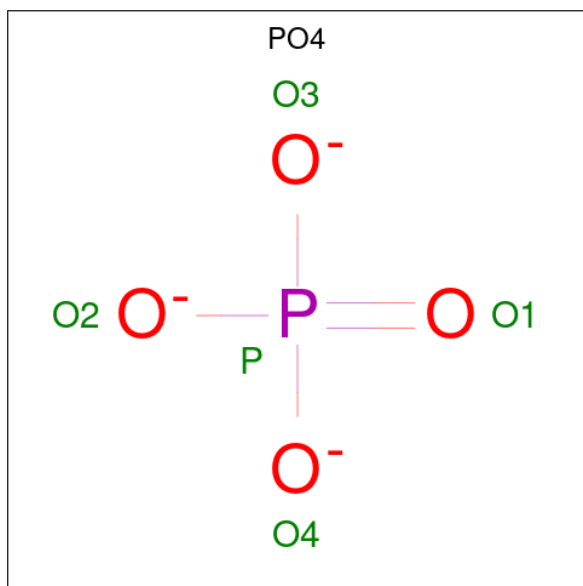
There are 9 unique types of molecules in this entry. The entry contains 11573 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 3-DEHYDROQUINATE SYNTHASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	365	Total	C	N	O	S	0	0	0
			2760	1752	474	516	18			
1	B	365	Total	C	N	O	S	0	0	0
			2758	1748	470	522	18			
1	C	365	Total	C	N	O	S	0	0	0
			2752	1744	474	516	18			
1	D	365	Total	C	N	O	S	0	0	0
			2768	1755	474	521	18			

- Molecule 2 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



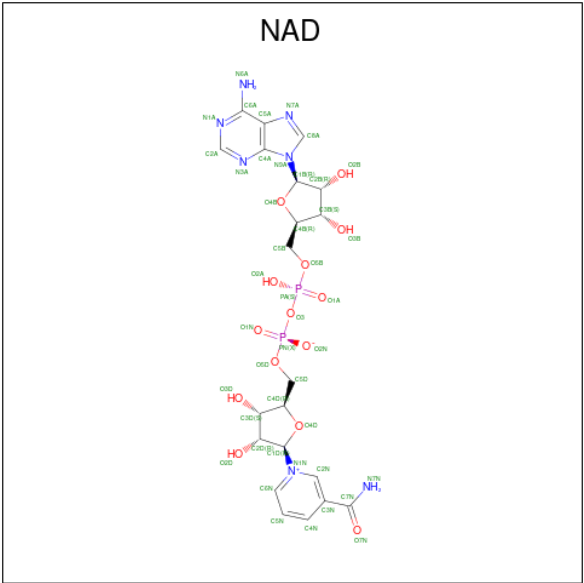
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	P	0	0
			5	4	1		
2	A	1	Total	O	P	0	0
			5	4	1		

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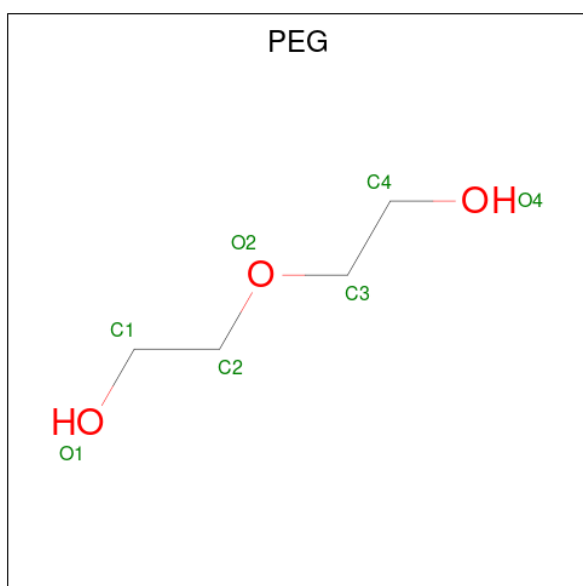
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	P		
2	A	1	5	4	1	0	0
2	B	1	5	4	1	0	0
2	B	1	5	4	1	0	0
2	B	1	5	4	1	0	0
2	B	1	5	4	1	0	0
2	C	1	5	4	1	0	0
2	C	1	5	4	1	0	0
2	C	1	5	4	1	0	0
2	D	1	5	4	1	0	0
2	D	1	5	4	1	0	0
2	D	1	5	4	1	0	0
2	D	1	5	4	1	0	0

- Molecule 3 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (CCD ID: NAD) (formula: C₂₁H₂₇N₇O₁₄P₂).



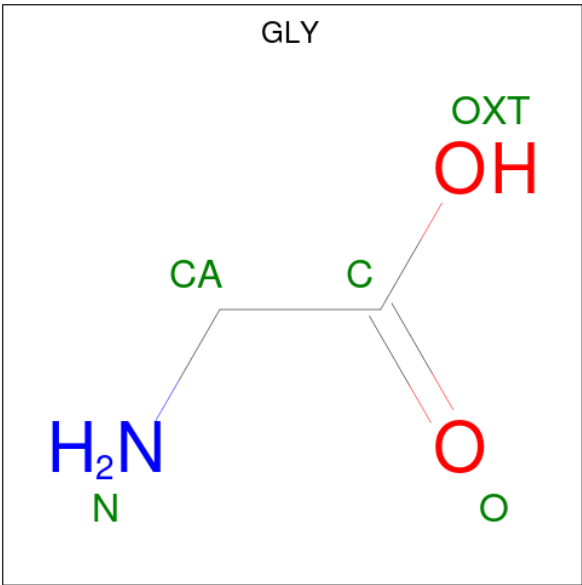
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
3	B	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
3	C	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
3	D	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

- Molecule 4 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



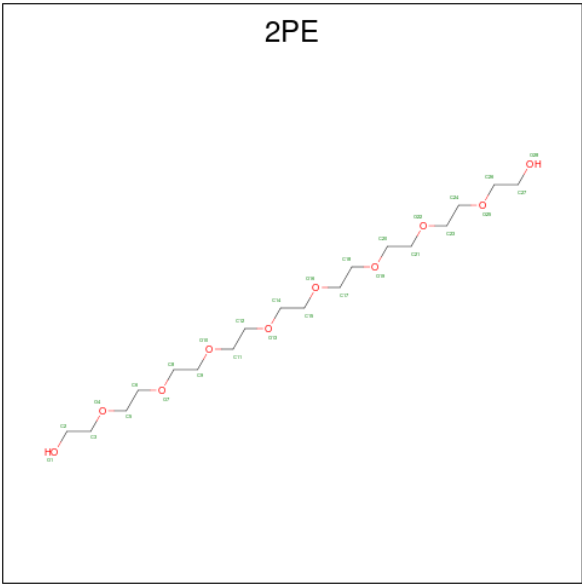
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			7	4	3		
4	A	1	Total	C	O	0	0
			7	4	3		
4	D	1	Total	C	O	0	0
			7	4	3		
4	D	1	Total	C	O	0	0
			7	4	3		
4	D	1	Total	C	O	0	0
			7	4	3		

- Molecule 5 is GLYCINE (CCD ID: GLY) (formula: $C_2H_5NO_2$).



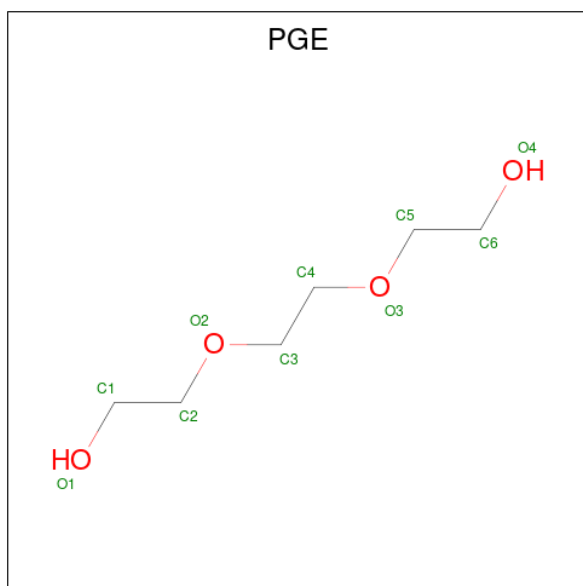
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			5	2	1	2		
5	B	1	Total	C	N	O	0	0
			5	2	1	2		
5	C	1	Total	C	N	O	0	0
			5	2	1	2		
5	D	1	Total	C	N	O	0	0
			5	2	1	2		
5	D	1	Total	C	N	O	0	0
			5	2	1	2		

- Molecule 6 is NONAETHYLENE GLYCOL (CCD ID: 2PE) (formula: C₁₈H₃₈O₁₀).



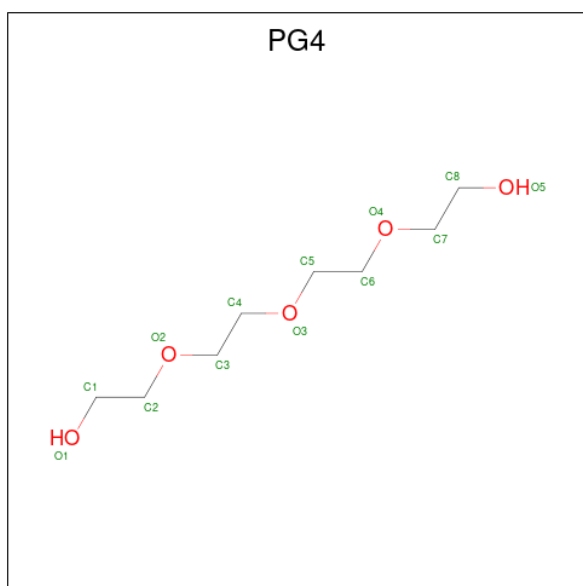
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	B	1	Total	C	O	0	0
			25	16	9		

- Molecule 7 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: $C_6H_{14}O_4$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	B	1	Total	C	O	0	0
			10	6	4		

- Molecule 8 is TETRAETHYLENE GLYCOL (CCD ID: PG4) (formula: $C_8H_{18}O_5$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	C	1	Total	C	O	0	0
			13	8	5		

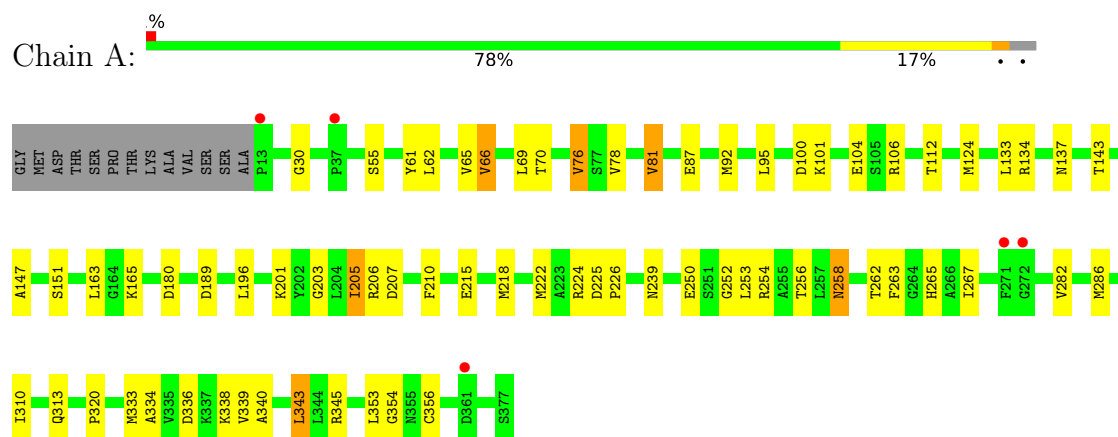
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	48	Total	O	0	0
			48	48		
9	B	40	Total	O	0	0
			40	40		
9	C	53	Total	O	0	0
			53	53		
9	D	40	Total	O	0	0
			40	40		

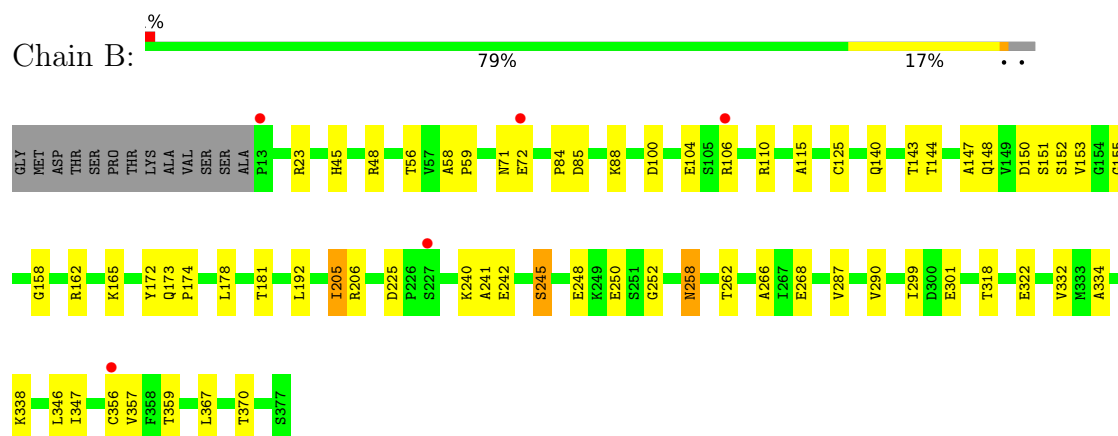
3 Residue-property plots [i](#)

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.

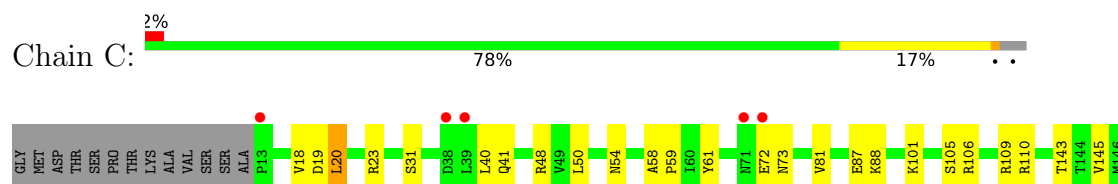
• Molecule 1: 3-DEHYDROQUINATE SYNTHASE

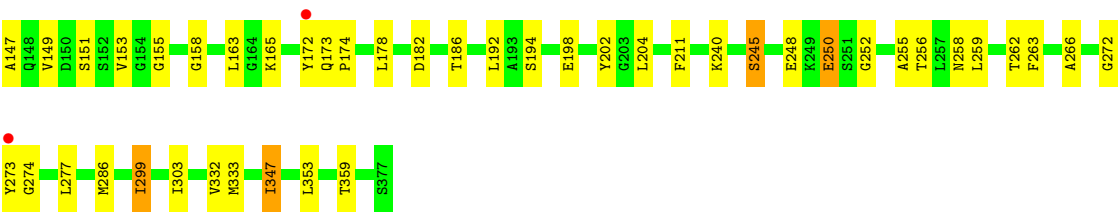


• Molecule 1: 3-DEHYDROQUINATE SYNTHASE

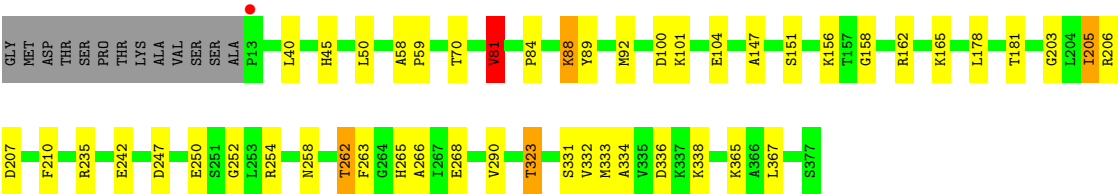
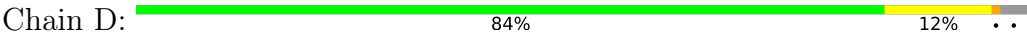


• Molecule 1: 3-DEHYDROQUINATE SYNTHASE





● Molecule 1: 3-DEHYDROQUINATE SYNTHASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	58.06Å 150.97Å 84.75Å 90.00° 101.25° 90.00°	Depositor
Resolution (Å)	45.50 – 2.40 45.50 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.5 (45.50-2.40) 99.5 (45.50-2.40)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.04 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.183 , 0.253 0.195 , 0.255	Depositor DCC
R_{free} test set	2834 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	29.9	Xtriage
Anisotropy	0.218	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 32.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	11573	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.42% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: PGE, 2PE, PG4, PO4, NAD, PEG

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	0.72	2/2811 (0.1%)	0.92	3/3814 (0.1%)
1	B	0.71	2/2808 (0.1%)	0.97	5/3812 (0.1%)
1	C	0.75	2/2802 (0.1%)	0.96	7/3802 (0.2%)
1	D	0.78	1/2819 (0.0%)	0.94	4/3825 (0.1%)
All	All	0.74	7/11240 (0.1%)	0.95	19/15253 (0.1%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	173	GLN	C-N	6.06	1.40	1.33
1	C	173	GLN	C-N	5.74	1.41	1.33
1	B	174	PRO	N-CD	5.46	1.55	1.47
1	C	174	PRO	N-CD	5.43	1.55	1.47
1	D	92	MET	C-O	-5.38	1.17	1.24
1	A	225	ASP	C-N	5.20	1.41	1.34
1	A	226	PRO	N-CD	5.18	1.55	1.47

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	272	GLY	N-CA-C	-9.96	98.83	112.18
1	B	205	ILE	N-CA-C	8.81	118.82	110.53
1	B	173	GLN	CA-C-N	-6.93	113.17	120.03
1	B	173	GLN	C-N-CA	-6.93	113.17	120.03
1	A	205	ILE	N-CA-C	6.86	116.98	110.53
1	C	173	GLN	CA-C-N	-6.70	113.08	119.85
1	C	173	GLN	C-N-CA	-6.70	113.08	119.85
1	D	81	VAL	CB-CA-C	-6.69	100.73	110.63
1	C	158	GLY	N-CA-C	6.25	119.26	110.38
1	D	205	ILE	N-CA-C	6.07	116.24	110.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	158	GLY	N-CA-C	5.88	118.37	110.43
1	D	290	VAL	N-CA-C	-5.49	105.37	110.53
1	B	250	GLU	N-CA-C	5.47	119.18	111.52
1	C	274	GLY	N-CA-C	5.32	121.03	114.69
1	D	158	GLY	N-CA-C	5.31	118.05	110.46
1	A	225	ASP	CA-C-N	-5.26	113.21	119.05
1	A	225	ASP	C-N-CA	-5.26	113.21	119.05
1	C	73	ASN	N-CA-C	5.13	117.49	108.82
1	C	250	GLU	N-CA-C	5.01	118.53	111.52

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	2760	0	2760	46	0
1	B	2758	0	2743	51	0
1	C	2752	0	2746	43	0
1	D	2768	0	2763	33	0
2	A	15	0	0	1	0
2	B	20	0	0	3	0
2	C	15	0	0	2	0
2	D	20	0	0	1	0
3	A	44	0	26	1	0
3	B	44	0	26	3	0
3	C	44	0	26	1	0
3	D	44	0	26	0	0
4	A	14	0	20	1	0
4	D	21	0	30	2	0
5	A	5	0	2	0	0
5	B	5	0	2	1	0
5	C	5	0	2	0	0
5	D	10	0	4	0	0
6	B	25	0	33	2	0
7	B	10	0	14	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
8	C	13	0	18	0	0
9	A	48	0	0	1	0
9	B	40	0	0	2	0
9	C	53	0	0	3	0
9	D	40	0	0	2	0
All	All	11573	0	11241	175	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 8.

All (175) 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:B:71:ASN:O	1:B:72:GLU:HG2	1.47	1.13
1:B:357:VAL:HG12	1:B:359:THR:HG23	1.40	1.02
1:B:206:ARG:NH2	1:B:242:GLU:CD	2.25	0.95
1:B:153:VAL:O	1:B:240:LYS:HE2	1.73	0.88
1:B:71:ASN:O	1:B:72:GLU:CG	2.23	0.87
1:A:147:ALA:HA	1:A:151:SER:OG	1.76	0.85
1:D:205:ILE:HG13	1:D:206:ARG:HG2	1.61	0.83
1:B:206:ARG:HH22	1:B:242:GLU:CD	1.87	0.81
1:A:92:MET:HE1	1:A:95:LEU:HD23	1.61	0.81
1:C:81:VAL:HG22	1:C:101:LYS:HG2	1.62	0.79
1:B:258:ASN:ND2	1:B:338:LYS:HE2	1.99	0.79
1:C:155:GLY:HA3	1:C:172:TYR:O	1.85	0.75
1:B:357:VAL:HG12	1:B:359:THR:CG2	2.18	0.74
1:C:23:ARG:NH2	1:C:248:GLU:OE1	2.21	0.73
1:A:206:ARG:NH1	1:A:239:ASN:OD1	2.20	0.73
1:B:266:ALA:HA	1:B:332:VAL:HG11	1.71	0.73
1:B:357:VAL:CG1	1:B:359:THR:HG23	2.17	0.73
1:B:45:HIS:O	1:B:110:ARG:NH2	2.24	0.71
1:D:262:THR:OG1	1:D:338:LYS:NZ	2.23	0.70
1:C:163:LEU:HD22	4:D:1384:PEG:H22	1.73	0.70
1:A:262:THR:OG1	1:A:338:LYS:NZ	2.21	0.70
1:C:155:GLY:CA	1:C:172:TYR:O	2.40	0.69
1:A:196:LEU:HD12	1:A:222:MET:HE3	1.72	0.69
1:B:258:ASN:HD22	1:B:338:LYS:CE	2.05	0.69
1:B:258:ASN:HD22	1:B:338:LYS:HE2	1.58	0.69
1:B:262:THR:OG1	1:B:338:LYS:NZ	2.25	0.69
1:C:147:ALA:HA	1:C:151:SER:OG	1.94	0.68
5:B:1385:GLY:N	9:B:2021:HOH:O	2.25	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:143:THR:HB	3:C:1381:NAD:H61A	1.58	0.68
1:C:87:GLU:HG3	1:C:273:TYR:HB3	1.76	0.67
1:D:206:ARG:HH22	1:D:242:GLU:CD	2.02	0.67
1:D:147:ALA:HA	1:D:151:SER:OG	1.96	0.66
1:A:263:PHE:CE1	1:A:333:MET:HE1	2.31	0.65
1:D:336:ASP:OD1	1:D:338:LYS:HE3	1.97	0.65
1:A:81:VAL:HG22	1:A:101:LYS:HG2	1.79	0.65
1:D:250:GLU:HG3	1:D:254:ARG:HB3	1.78	0.64
1:B:71:ASN:C	1:B:72:GLU:HG2	2.21	0.63
1:B:338:LYS:HD3	1:B:347:ILE:HD11	1.81	0.63
1:D:100:ASP:O	1:D:104:GLU:HG2	1.98	0.63
1:B:357:VAL:CG1	1:B:359:THR:CG2	2.77	0.62
1:C:87:GLU:OE2	1:C:165:LYS:NZ	2.31	0.62
1:B:206:ARG:NH2	1:B:242:GLU:OE1	2.26	0.62
1:C:194:SER:OG	1:C:277:LEU:HD12	2.00	0.62
1:D:263:PHE:CE1	1:D:333:MET:HE1	2.35	0.61
1:B:150:ASP:O	1:B:240:LYS:NZ	2.32	0.61
1:D:81:VAL:HG22	1:D:101:LYS:HG2	1.82	0.60
1:A:196:LEU:CD1	1:A:222:MET:HE3	2.31	0.59
1:B:150:ASP:OD1	1:B:240:LYS:NZ	2.23	0.59
1:B:153:VAL:O	1:B:240:LYS:CE	2.48	0.58
1:D:263:PHE:CD1	1:D:333:MET:HE1	2.39	0.58
1:C:48:ARG:NH2	1:C:105:SER:O	2.37	0.58
1:C:61:TYR:HH	1:C:186:THR:HG1	1.52	0.58
1:D:263:PHE:CD1	1:D:333:MET:CE	2.86	0.57
1:C:40:LEU:HD21	1:C:178:LEU:HD22	1.86	0.56
1:C:263:PHE:CE1	1:C:333:MET:HE1	2.40	0.56
1:A:267:ILE:HD11	1:A:320:PRO:HB3	1.88	0.56
1:C:18:VAL:HG12	1:C:20:LEU:HD22	1.87	0.56
1:C:50:LEU:HD11	1:C:81:VAL:HG23	1.86	0.56
1:A:207:ASP:CG	1:A:210:PHE:HB2	2.30	0.56
1:A:258:ASN:HD22	1:A:338:LYS:HE2	1.71	0.55
1:D:45:HIS:HE1	9:D:2003:HOH:O	1.89	0.55
1:B:338:LYS:HD3	1:B:347:ILE:CD1	2.37	0.55
1:A:92:MET:HE1	1:A:95:LEU:CD2	2.34	0.55
1:A:265:HIS:HB3	9:A:2029:HOH:O	2.06	0.54
2:C:1378:PO4:O3	2:C:1379:PO4:O3	2.26	0.54
1:B:241:ALA:O	1:B:245:SER:OG	2.24	0.54
1:C:255:ALA:HB1	1:C:347:ILE:HD11	1.89	0.54
1:A:334:ALA:HB1	1:B:106:ARG:HD3	1.89	0.54
1:B:165:LYS:HE2	1:B:268:GLU:OE2	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:145:VAL:HG21	1:C:192:LEU:HD12	1.88	0.53
1:A:205:ILE:HG13	1:A:206:ARG:N	2.24	0.53
1:D:165:LYS:HG2	1:D:265:HIS:CD2	2.44	0.53
2:B:1378:PO4:O2	2:B:1379:PO4:O2	2.28	0.52
1:D:40:LEU:HD21	1:D:178:LEU:HD22	1.90	0.52
1:B:258:ASN:ND2	1:B:338:LYS:CE	2.68	0.51
1:B:205:ILE:HG13	1:B:206:ARG:HG2	1.91	0.51
1:C:266:ALA:HA	1:C:332:VAL:HG11	1.90	0.51
1:A:336:ASP:OD1	1:A:338:LYS:HE3	2.11	0.51
1:C:259:LEU:HD23	1:C:286:MET:HE3	1.92	0.51
1:C:41:GLN:OE1	1:C:72:GLU:OE1	2.29	0.51
1:A:61:TYR:O	1:A:65:VAL:HG23	2.10	0.50
1:C:145:VAL:O	1:C:149:VAL:HG22	2.11	0.50
1:A:87:GLU:OE2	1:A:165:LYS:NZ	2.39	0.50
1:A:253:LEU:O	1:A:256:THR:OG1	2.20	0.50
1:C:204:LEU:HD12	1:C:286:MET:CE	2.42	0.50
1:D:165:LYS:HE2	1:D:268:GLU:OE1	2.12	0.50
1:B:84:PRO:HB3	6:B:1383:2PE:C9	2.41	0.50
1:C:110:ARG:HD2	9:C:2003:HOH:O	2.10	0.49
1:B:88:LYS:HB2	9:B:2007:HOH:O	2.12	0.49
1:A:262:THR:HG22	1:A:333:MET:HG2	1.93	0.49
1:C:18:VAL:HG12	1:C:20:LEU:CD2	2.43	0.49
1:A:203:GLY:HA2	1:A:207:ASP:O	2.12	0.49
1:B:100:ASP:O	1:B:104:GLU:HG2	2.12	0.48
1:A:62:LEU:O	1:A:66:VAL:HG12	2.13	0.48
1:A:282:VAL:HG12	1:A:286:MET:HE2	1.96	0.48
1:A:70:THR:HG22	1:A:76:VAL:HG12	1.95	0.48
1:A:106:ARG:HE	7:B:1384:PGE:H62	1.79	0.48
1:B:56:THR:OG1	1:B:85:ASP:OD2	2.27	0.48
1:C:256:THR:HA	1:C:347:ILE:HG12	1.96	0.48
1:B:205:ILE:HG13	1:B:206:ARG:N	2.29	0.48
1:A:215:GLU:HA	1:A:310:ILE:HD13	1.95	0.48
1:A:100:ASP:O	1:A:104:GLU:HG2	2.13	0.48
1:A:133:LEU:O	1:A:134:ARG:HB2	2.14	0.47
1:A:201:LYS:O	1:A:205:ILE:HG12	2.15	0.47
1:B:258:ASN:HD22	1:B:338:LYS:NZ	2.13	0.47
1:C:40:LEU:CD2	1:C:178:LEU:HD22	2.45	0.47
1:D:58:ALA:HB3	1:D:59:PRO:HD3	1.96	0.46
1:D:207:ASP:HB3	1:D:210:PHE:HB3	1.96	0.46
1:A:250:GLU:HG3	1:A:254:ARG:HB3	1.98	0.46
1:B:266:ALA:HA	1:B:332:VAL:CG1	2.44	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:198:GLU:O	1:C:202:TYR:CD2	2.68	0.46
1:B:287:VAL:HG21	1:B:318:THR:HG22	1.98	0.46
1:C:262:THR:HG22	1:C:333:MET:HG2	1.98	0.46
1:C:299:ILE:HG13	1:C:303:ILE:HD12	1.98	0.45
1:A:189:ASP:OD1	1:A:224:ARG:NH1	2.49	0.45
2:A:1378:PO4:O2	2:A:1379:PO4:O4	2.34	0.45
1:A:339:VAL:HA	1:A:343:LEU:O	2.16	0.45
1:D:247:ASP:OD2	1:D:250:GLU:HA	2.17	0.45
1:D:266:ALA:HA	1:D:332:VAL:HG11	1.98	0.45
1:C:263:PHE:CE2	2:C:1380:PO4:O3	2.70	0.45
1:D:165:LYS:HE3	1:D:268:GLU:OE2	2.17	0.44
1:B:155:GLY:CA	1:B:172:TYR:O	2.66	0.44
1:B:58:ALA:HB3	1:B:59:PRO:HD3	1.99	0.44
1:D:205:ILE:HG13	1:D:206:ARG:N	2.33	0.44
1:B:162:ARG:CD	1:D:162:ARG:HD2	2.48	0.44
1:D:88:LYS:HD3	1:D:89:TYR:CZ	2.53	0.44
1:B:115:ALA:HB2	1:B:125:CYS:SG	2.58	0.44
1:A:256:THR:HG22	1:A:356:CYS:SG	2.58	0.44
1:B:140:GLN:HG2	1:B:152:SER:HB2	1.99	0.44
1:A:124:MET:HE2	1:A:124:MET:HB3	1.91	0.43
1:C:54:ASN:C	1:C:54:ASN:OD1	2.60	0.43
1:D:50:LEU:HD11	1:D:81:VAL:HG23	2.00	0.43
1:A:30:GLY:O	1:A:180:ASP:HA	2.18	0.43
1:A:340:ALA:HB2	1:A:345:ARG:CZ	2.49	0.43
2:B:1379:PO4:O1	3:B:1382:NAD:O2D	2.31	0.43
1:D:235:ARG:HD3	2:D:1381:PO4:O2	2.18	0.43
1:C:353:LEU:HD12	9:C:2042:HOH:O	2.18	0.43
1:D:50:LEU:HD11	1:D:81:VAL:CG2	2.49	0.43
1:D:323:THR:O	1:D:323:THR:CG2	2.66	0.43
1:B:147:ALA:HA	1:B:151:SER:OG	2.19	0.43
1:C:106:ARG:NH2	9:C:2012:HOH:O	2.51	0.43
1:A:206:ARG:HA	1:A:354:GLY:O	2.19	0.42
1:C:31:SER:OG	1:C:182:ASP:OD1	2.28	0.42
1:A:163:LEU:HB3	4:A:1383:PEG:H32	2.01	0.42
1:B:144:THR:O	1:B:148:GLN:HG3	2.18	0.42
1:C:109:ARG:HH12	1:D:250:GLU:HB3	1.85	0.42
1:A:143:THR:HB	3:A:1381:NAD:H61A	1.83	0.42
1:B:143:THR:HB	3:B:1382:NAD:H61A	1.84	0.42
1:C:19:ASP:O	1:C:245:SER:HB3	2.19	0.42
1:D:203:GLY:HA2	1:D:207:ASP:HB3	2.02	0.42
1:B:258:ASN:CG	1:B:338:LYS:HE2	2.45	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:218:MET:HG2	1:A:313:GLN:HB3	2.00	0.42
1:B:225:ASP:C	1:B:225:ASP:OD1	2.62	0.42
1:C:23:ARG:HH22	1:C:248:GLU:CD	2.23	0.42
1:C:263:PHE:CD1	1:C:333:MET:HE1	2.54	0.42
1:D:156:LYS:HB2	9:D:2028:HOH:O	2.20	0.42
1:A:112:THR:HA	1:A:137:ASN:O	2.19	0.41
1:B:23:ARG:NH2	1:B:248:GLU:OE1	2.37	0.41
1:C:153:VAL:O	1:C:240:LYS:HE2	2.20	0.41
1:D:334:ALA:HA	4:D:1385:PEG:H32	2.02	0.41
1:B:165:LYS:HE2	1:B:268:GLU:CD	2.45	0.41
1:D:367:LEU:C	1:D:367:LEU:HD13	2.46	0.41
1:B:346:LEU:HD22	2:B:1380:PO4:O4	2.21	0.41
1:A:205:ILE:HG13	1:A:206:ARG:HG2	2.03	0.41
1:B:151:SER:HA	3:B:1382:NAD:C4N	2.51	0.41
1:B:334:ALA:HA	7:B:1384:PGE:H22	2.03	0.41
1:A:92:MET:CE	1:A:95:LEU:HD23	2.41	0.41
1:C:58:ALA:HB3	1:C:59:PRO:HD3	2.02	0.41
1:A:65:VAL:O	1:A:69:LEU:HD13	2.21	0.41
1:A:258:ASN:ND2	1:A:338:LYS:HE2	2.35	0.41
1:B:367:LEU:O	1:B:370:THR:HG22	2.21	0.41
1:C:250:GLU:O	1:C:250:GLU:HG2	2.20	0.41
1:A:70:THR:CG2	1:A:78:VAL:HG23	2.51	0.40
6:B:1383:2PE:H151	1:D:84:PRO:HG3	2.03	0.40
1:C:211:PHE:CD2	1:C:353:LEU:HD21	2.56	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	363/378 (96%)	352 (97%)	10 (3%)	1 (0%)	36 50

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	363/378 (96%)	355 (98%)	7 (2%)	1 (0%)	36	50
1	C	363/378 (96%)	350 (96%)	12 (3%)	1 (0%)	36	50
1	D	363/378 (96%)	353 (97%)	9 (2%)	1 (0%)	36	50
All	All	1452/1512 (96%)	1410 (97%)	38 (3%)	4 (0%)	36	50

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	252	GLY
1	B	252	GLY
1	C	252	GLY
1	D	252	GLY

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	293/312 (94%)	286 (98%)	7 (2%)	43	65
1	B	293/312 (94%)	282 (96%)	11 (4%)	29	49
1	C	291/312 (93%)	284 (98%)	7 (2%)	43	65
1	D	295/312 (95%)	286 (97%)	9 (3%)	35	57
All	All	1172/1248 (94%)	1138 (97%)	34 (3%)	37	60

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	55	SER
1	A	66	VAL
1	A	76	VAL
1	A	81	VAL
1	A	258	ASN
1	A	343	LEU
1	A	353	LEU

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Mol	Chain	Res	Type
1	B	48	ARG
1	B	178	LEU
1	B	181	THR
1	B	192	LEU
1	B	245	SER
1	B	258	ASN
1	B	290	VAL
1	B	299	ILE
1	B	301	GLU
1	B	322	GLU
1	B	356	CYS
1	C	20	LEU
1	C	88	LYS
1	C	245	SER
1	C	258	ASN
1	C	299	ILE
1	C	347	ILE
1	C	359	THR
1	D	70	THR
1	D	81	VAL
1	D	88	LYS
1	D	181	THR
1	D	258	ASN
1	D	262	THR
1	D	323	THR
1	D	331	SER
1	D	365	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	258	ASN
1	A	313	GLN
1	A	372	HIS
1	B	217	ASN
1	B	258	ASN
1	B	275	GLN
1	C	175	GLN
1	C	217	ASN
1	C	258	ASN
1	D	45	HIS
1	D	209	ASN

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Mol	Chain	Res	Type
1	D	258	ASN
1	D	275	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

31 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
3	NAD	B	1382	-	46,48,48	1.26	6 (13%)	64,73,73	1.58	10 (15%)
8	PG4	C	1382	-	12,12,12	0.59	0	11,11,11	0.30	0
4	PEG	A	1383	-	6,6,6	0.58	0	5,5,5	0.68	0
2	PO4	D	1380	-	4,4,4	1.06	0	6,6,6	0.89	0
2	PO4	B	1381	-	4,4,4	1.06	0	6,6,6	0.52	0
2	PO4	D	1381	-	4,4,4	1.07	0	6,6,6	0.69	0
5	GLY	C	1383	-	4,4,4	1.18	1 (25%)	3,4,4	1.50	0
2	PO4	C	1378	-	4,4,4	0.77	0	6,6,6	0.78	0
5	GLY	D	1387	-	4,4,4	1.12	1 (25%)	3,4,4	1.37	0
4	PEG	D	1385	-	6,6,6	0.45	0	5,5,5	0.42	0
3	NAD	C	1381	-	46,48,48	1.20	5 (10%)	64,73,73	1.65	10 (15%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
7	PGE	B	1384	-	9,9,9	0.70	0	8,8,8	0.43	0
5	GLY	D	1386	-	4,4,4	1.22	1 (25%)	3,4,4	1.92	2 (66%)
4	PEG	D	1383	-	6,6,6	0.50	0	5,5,5	0.12	0
4	PEG	A	1382	-	6,6,6	0.59	0	5,5,5	0.41	0
2	PO4	C	1379	-	4,4,4	1.04	0	6,6,6	0.52	0
3	NAD	A	1381	-	46,48,48	1.28	6 (13%)	64,73,73	1.64	13 (20%)
5	GLY	B	1385	-	4,4,4	1.20	1 (25%)	3,4,4	1.80	2 (66%)
2	PO4	C	1380	-	4,4,4	0.97	0	6,6,6	0.76	0
2	PO4	D	1379	-	4,4,4	1.20	0	6,6,6	0.93	0
2	PO4	B	1378	-	4,4,4	1.05	0	6,6,6	0.82	0
2	PO4	A	1378	-	4,4,4	1.46	0	6,6,6	1.14	1 (16%)
2	PO4	D	1378	-	4,4,4	1.36	0	6,6,6	1.16	1 (16%)
2	PO4	B	1379	-	4,4,4	1.03	0	6,6,6	0.46	0
5	GLY	A	1384	-	4,4,4	1.00	0	3,4,4	1.93	2 (66%)
6	2PE	B	1383	-	24,24,27	0.48	0	23,23,26	0.50	0
2	PO4	A	1380	-	4,4,4	0.69	0	6,6,6	1.13	0
4	PEG	D	1384	-	6,6,6	0.63	0	5,5,5	0.67	0
3	NAD	D	1382	-	46,48,48	1.31	7 (15%)	64,73,73	1.58	10 (15%)
2	PO4	A	1379	-	4,4,4	1.07	0	6,6,6	0.29	0
2	PO4	B	1380	-	4,4,4	1.09	0	6,6,6	0.86	0

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
5	GLY	D	1387	-	-	0/2/2/2	-
3	NAD	B	1382	-	-	14/30/62/62	0/5/5/5
4	PEG	D	1383	-	-	3/4/4/4	-
4	PEG	A	1382	-	-	1/4/4/4	-
4	PEG	A	1383	-	-	2/4/4/4	-
5	GLY	A	1384	-	-	0/2/2/2	-
6	2PE	B	1383	-	-	11/22/22/25	-
3	NAD	A	1381	-	-	11/30/62/62	0/5/5/5
4	PEG	D	1384	-	-	2/4/4/4	-
5	GLY	B	1385	-	-	2/2/2/2	-
8	PG4	C	1382	-	-	6/10/10/10	-
3	NAD	D	1382	-	-	5/30/62/62	0/5/5/5

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	PEG	D	1385	-	-	2/4/4/4	-
5	GLY	C	1383	-	-	0/2/2/2	-
3	NAD	C	1381	-	-	2/30/62/62	0/5/5/5
7	PGE	B	1384	-	-	4/7/7/7	-
5	GLY	D	1386	-	-	0/2/2/2	-

All (28) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	1382	NAD	C5A-C4A	3.93	1.46	1.39
3	B	1382	NAD	C5A-C4A	3.86	1.46	1.39
3	A	1381	NAD	C5A-C4A	3.82	1.45	1.39
3	C	1381	NAD	C5A-C4A	3.59	1.45	1.39
3	C	1381	NAD	C5A-N7A	-3.00	1.33	1.39
3	D	1382	NAD	C5A-N7A	-2.96	1.33	1.39
3	A	1381	NAD	C4A-N9A	-2.88	1.31	1.37
3	A	1381	NAD	C5A-N7A	-2.86	1.33	1.39
3	B	1382	NAD	C4A-N9A	-2.81	1.31	1.37
3	A	1381	NAD	C2N-N1N	-2.77	1.31	1.35
3	B	1382	NAD	C5A-N7A	-2.70	1.34	1.39
3	A	1381	NAD	C8A-N9A	-2.70	1.33	1.37
3	C	1381	NAD	C8A-N9A	-2.68	1.33	1.37
3	C	1381	NAD	C4A-N9A	-2.65	1.32	1.37
3	D	1382	NAD	C2N-N1N	-2.60	1.32	1.35
3	B	1382	NAD	C8A-N9A	-2.50	1.33	1.37
3	D	1382	NAD	C8A-N9A	-2.50	1.33	1.37
3	D	1382	NAD	C4A-N9A	-2.49	1.32	1.37
3	B	1382	NAD	C5A-C6A	2.49	1.47	1.41
3	D	1382	NAD	C5A-C6A	2.48	1.47	1.41
3	D	1382	NAD	O7N-C7N	-2.43	1.19	1.24
3	A	1381	NAD	C5A-C6A	2.40	1.47	1.41
5	D	1386	GLY	OXT-C	-2.39	1.22	1.30
3	C	1381	NAD	C5A-C6A	2.39	1.47	1.41
3	B	1382	NAD	C2N-N1N	-2.38	1.32	1.35
5	B	1385	GLY	OXT-C	-2.35	1.23	1.30
5	C	1383	GLY	OXT-C	-2.08	1.23	1.30
5	D	1387	GLY	OXT-C	-2.04	1.24	1.30

All (51) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	1382	NAD	C5A-C4A-N3A	-5.83	118.70	126.72
3	C	1381	NAD	C5A-C4A-N3A	-5.55	119.08	126.72
3	B	1382	NAD	C5A-C4A-N3A	-5.46	119.20	126.72
3	D	1382	NAD	N3A-C4A-N9A	5.08	135.81	127.17
3	A	1381	NAD	C5A-C4A-N3A	-5.07	119.74	126.72
3	C	1381	NAD	N3A-C4A-N9A	4.65	135.08	127.17
3	B	1382	NAD	N3A-C4A-N9A	4.36	134.58	127.17
3	A	1381	NAD	N3A-C4A-N9A	4.23	134.37	127.17
3	C	1381	NAD	C2A-N3A-C4A	3.76	121.01	111.83
3	B	1382	NAD	C4A-C5A-N7A	-3.73	106.31	110.58
3	A	1381	NAD	N3A-C2A-N1A	-3.70	122.97	128.58
3	A	1381	NAD	C2A-N3A-C4A	3.66	120.77	111.83
3	B	1382	NAD	C2A-N3A-C4A	3.64	120.73	111.83
3	D	1382	NAD	C2A-N3A-C4A	3.62	120.68	111.83
3	C	1381	NAD	C4A-N9A-C8A	3.53	109.45	105.74
3	C	1381	NAD	C4A-C5A-N7A	-3.51	106.57	110.58
3	C	1381	NAD	N3A-C2A-N1A	-3.51	123.27	128.58
3	A	1381	NAD	C4A-C5A-N7A	-3.46	106.62	110.58
3	A	1381	NAD	C4A-N9A-C8A	3.44	109.35	105.74
3	D	1382	NAD	C4A-C5A-N7A	-3.34	106.76	110.58
3	B	1382	NAD	N3A-C2A-N1A	-3.32	123.56	128.58
3	D	1382	NAD	C4A-N9A-C8A	3.25	109.15	105.74
3	B	1382	NAD	C4A-N9A-C8A	3.17	109.06	105.74
3	D	1382	NAD	N3A-C2A-N1A	-3.11	123.87	128.58
3	A	1381	NAD	O2N-PN-O3	2.86	115.01	107.27
3	C	1381	NAD	C5A-N7A-C8A	2.74	107.76	103.45
3	D	1382	NAD	C5A-N7A-C8A	2.64	107.59	103.45
5	D	1386	GLY	OXT-C-O	-2.62	116.60	123.33
3	A	1381	NAD	C6A-C5A-N7A	2.61	137.12	132.09
3	B	1382	NAD	C5A-N7A-C8A	2.57	107.49	103.45
3	B	1382	NAD	C6A-C5A-N7A	2.57	137.04	132.09
3	C	1381	NAD	N9A-C8A-N7A	-2.52	110.37	113.94
5	A	1384	GLY	OXT-C-O	-2.48	116.95	123.33
3	A	1381	NAD	C5N-C4N-C3N	-2.46	117.94	120.36
3	A	1381	NAD	C5A-N7A-C8A	2.45	107.29	103.45
2	D	1378	PO4	O4-P-O3	2.37	115.29	107.91
3	C	1381	NAD	C6A-C5A-N7A	2.35	136.63	132.09
3	D	1382	NAD	C3N-C7N-N7N	2.33	120.61	117.74
5	B	1385	GLY	OXT-C-O	-2.27	117.50	123.33
3	A	1381	NAD	C2A-N1A-C6A	2.26	122.45	118.73
5	A	1384	GLY	OXT-C-CA	2.21	122.16	113.38
3	A	1381	NAD	N9A-C8A-N7A	-2.19	110.83	113.94
3	B	1382	NAD	N9A-C8A-N7A	-2.15	110.89	113.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1378	PO4	O3-P-O1	-2.14	103.39	110.95
5	B	1385	GLY	OXT-C-CA	2.11	121.76	113.38
3	B	1382	NAD	C2A-N1A-C6A	2.08	122.14	118.73
3	D	1382	NAD	N9A-C8A-N7A	-2.07	111.00	113.94
3	A	1381	NAD	O4B-C1B-N9A	-2.06	104.13	108.09
3	C	1381	NAD	O3-PN-O1N	-2.06	104.51	110.70
5	D	1386	GLY	OXT-C-CA	2.05	121.54	113.38
3	D	1382	NAD	C6A-C5A-N7A	2.01	135.97	132.09

There are no chirality outliers.

All (65) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1381	NAD	C5D-O5D-PN-O1N
3	A	1381	NAD	O4D-C1D-N1N-C2N
3	A	1381	NAD	O4D-C1D-N1N-C6N
3	A	1381	NAD	C2D-C1D-N1N-C2N
3	B	1382	NAD	C5D-O5D-PN-O3
3	B	1382	NAD	O4D-C1D-N1N-C2N
3	B	1382	NAD	O4D-C1D-N1N-C6N
3	C	1381	NAD	O4D-C1D-N1N-C6N
3	D	1382	NAD	O4D-C1D-N1N-C6N
5	B	1385	GLY	O-C-CA-N
5	B	1385	GLY	OXT-C-CA-N
3	B	1382	NAD	C3D-C4D-C5D-O5D
7	B	1384	PGE	O2-C3-C4-O3
8	C	1382	PG4	O2-C3-C4-O3
6	B	1383	2PE	O16-C17-C18-O19
6	B	1383	2PE	O4-C5-C6-O7
6	B	1383	2PE	O13-C14-C15-O16
3	B	1382	NAD	C3B-C4B-C5B-O5B
3	B	1382	NAD	O4D-C4D-C5D-O5D
7	B	1384	PGE	O3-C5-C6-O4
6	B	1383	2PE	O10-C11-C12-O13
4	D	1384	PEG	O1-C1-C2-O2
8	C	1382	PG4	O4-C7-C8-O5
6	B	1383	2PE	O1-C2-C3-O4
3	A	1381	NAD	O4D-C4D-C5D-O5D
3	A	1381	NAD	C3D-C4D-C5D-O5D
4	A	1383	PEG	O1-C1-C2-O2
4	D	1384	PEG	O2-C3-C4-O4
6	B	1383	2PE	O22-C23-C24-O25

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Mol	Chain	Res	Type	Atoms
4	D	1383	PEG	O2-C3-C4-O4
3	B	1382	NAD	O4B-C4B-C5B-O5B
3	D	1382	NAD	O4B-C4B-C5B-O5B
3	D	1382	NAD	C3B-C4B-C5B-O5B
8	C	1382	PG4	O3-C5-C6-O4
6	B	1383	2PE	C9-C8-O7-C6
8	C	1382	PG4	C6-C5-O3-C4
7	B	1384	PGE	C4-C3-O2-C2
8	C	1382	PG4	C4-C3-O2-C2
6	B	1383	2PE	C18-C17-O16-C15
4	D	1385	PEG	C1-C2-O2-C3
4	D	1385	PEG	O2-C3-C4-O4
3	A	1381	NAD	PA-O3-PN-O2N
6	B	1383	2PE	C8-C9-O10-C11
7	B	1384	PGE	C6-C5-O3-C4
3	B	1382	NAD	C5B-O5B-PA-O3
3	B	1382	NAD	C5D-O5D-PN-O1N
3	A	1381	NAD	PA-O3-PN-O1N
4	A	1382	PEG	C1-C2-O2-C3
4	A	1383	PEG	C4-C3-O2-C2
3	A	1381	NAD	C2D-C1D-N1N-C6N
3	B	1382	NAD	C2D-C1D-N1N-C2N
3	B	1382	NAD	C2D-C1D-N1N-C6N
8	C	1382	PG4	C1-C2-O2-C3
4	D	1383	PEG	C1-C2-O2-C3
3	A	1381	NAD	O4B-C4B-C5B-O5B
3	A	1381	NAD	C3B-C4B-C5B-O5B
3	C	1381	NAD	O4D-C1D-N1N-C2N
3	D	1382	NAD	O4D-C1D-N1N-C2N
6	B	1383	2PE	C15-C14-O13-C12
3	B	1382	NAD	PN-O3-PA-O1A
3	D	1382	NAD	C2B-C1B-N9A-C8A
4	D	1383	PEG	O1-C1-C2-O2
6	B	1383	2PE	O7-C8-C9-O10
3	B	1382	NAD	C2B-C1B-N9A-C8A
3	B	1382	NAD	PN-O3-PA-O2A

There are no ring outliers.

18 monomers are involved in 19 short contacts:

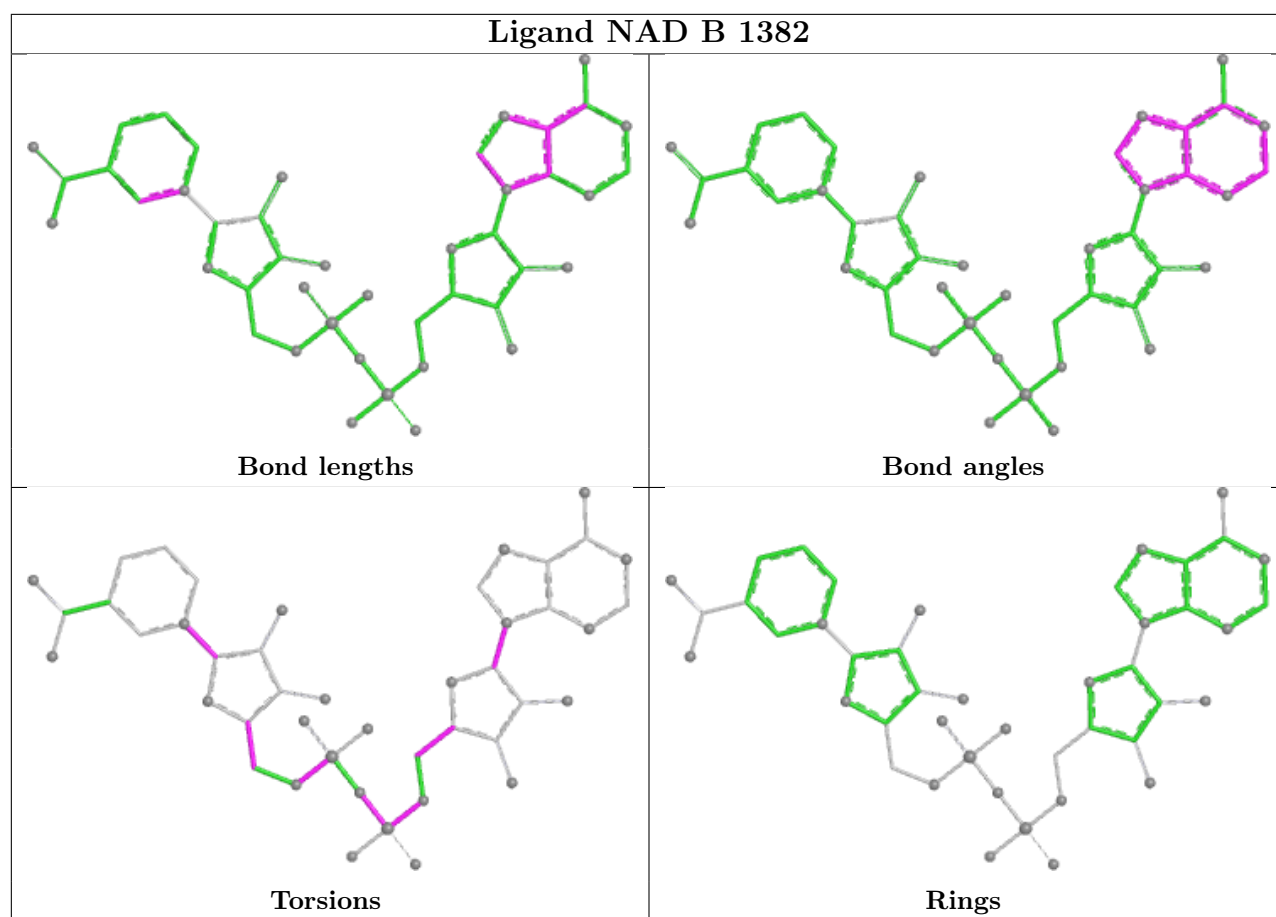
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1382	NAD	3	0

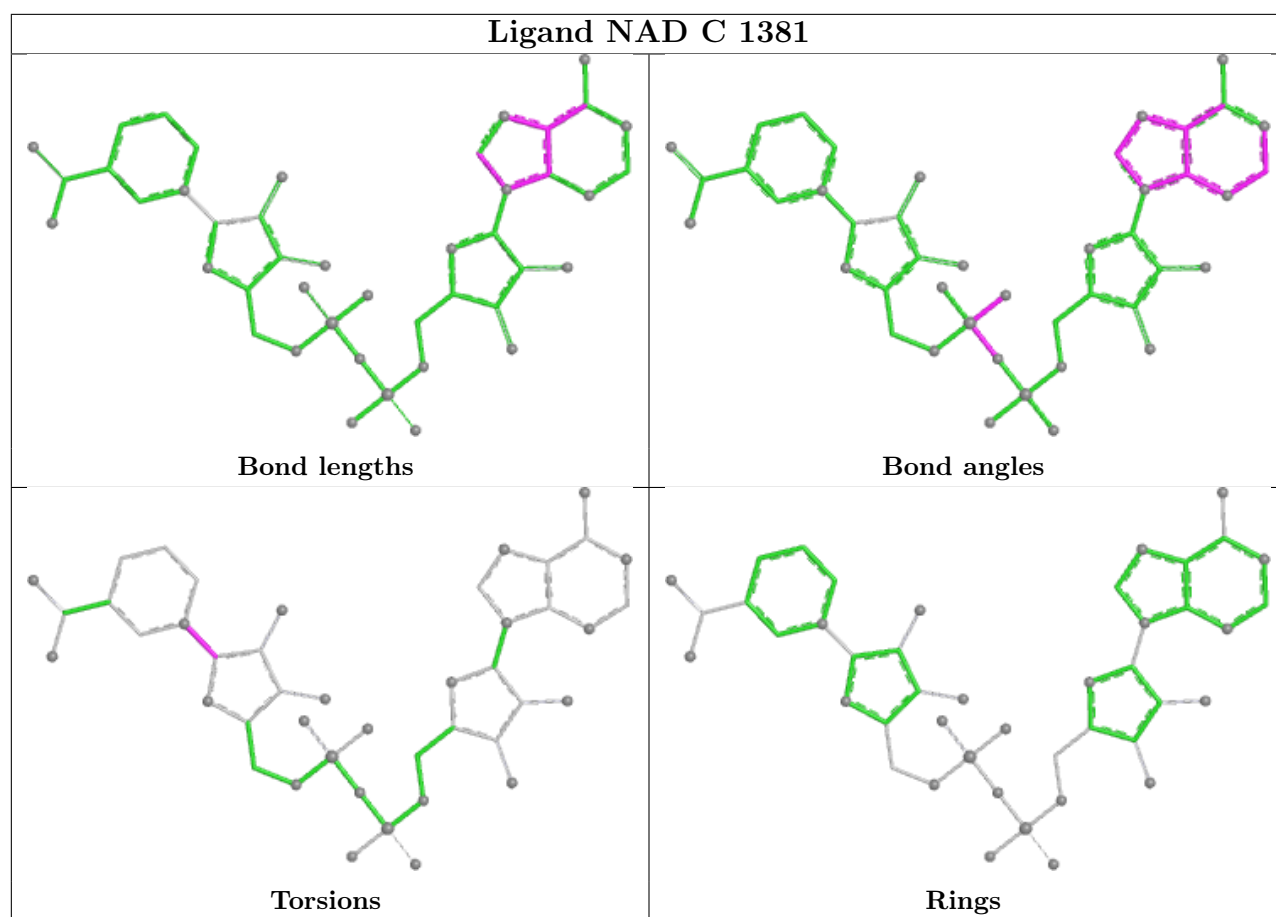
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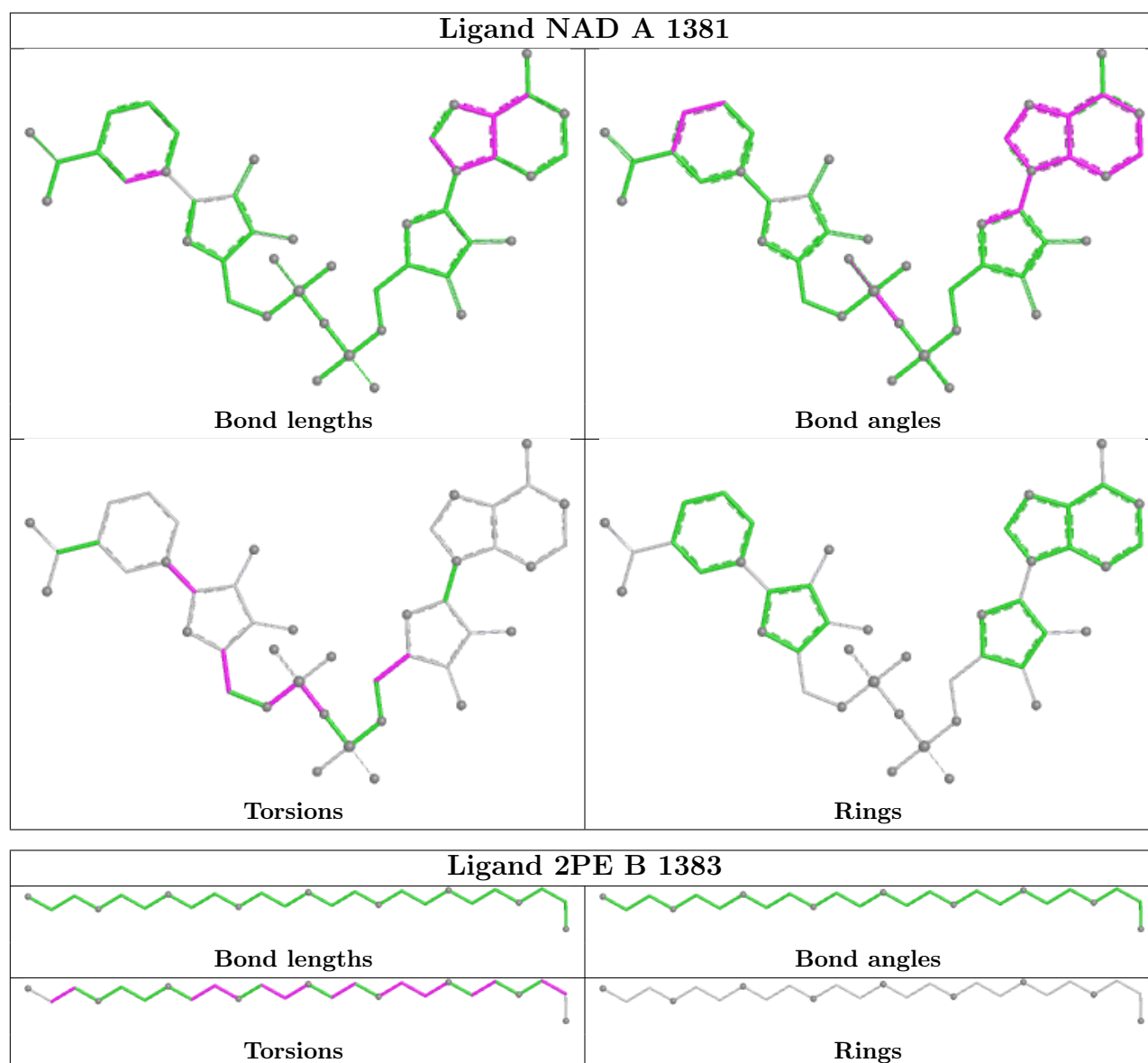
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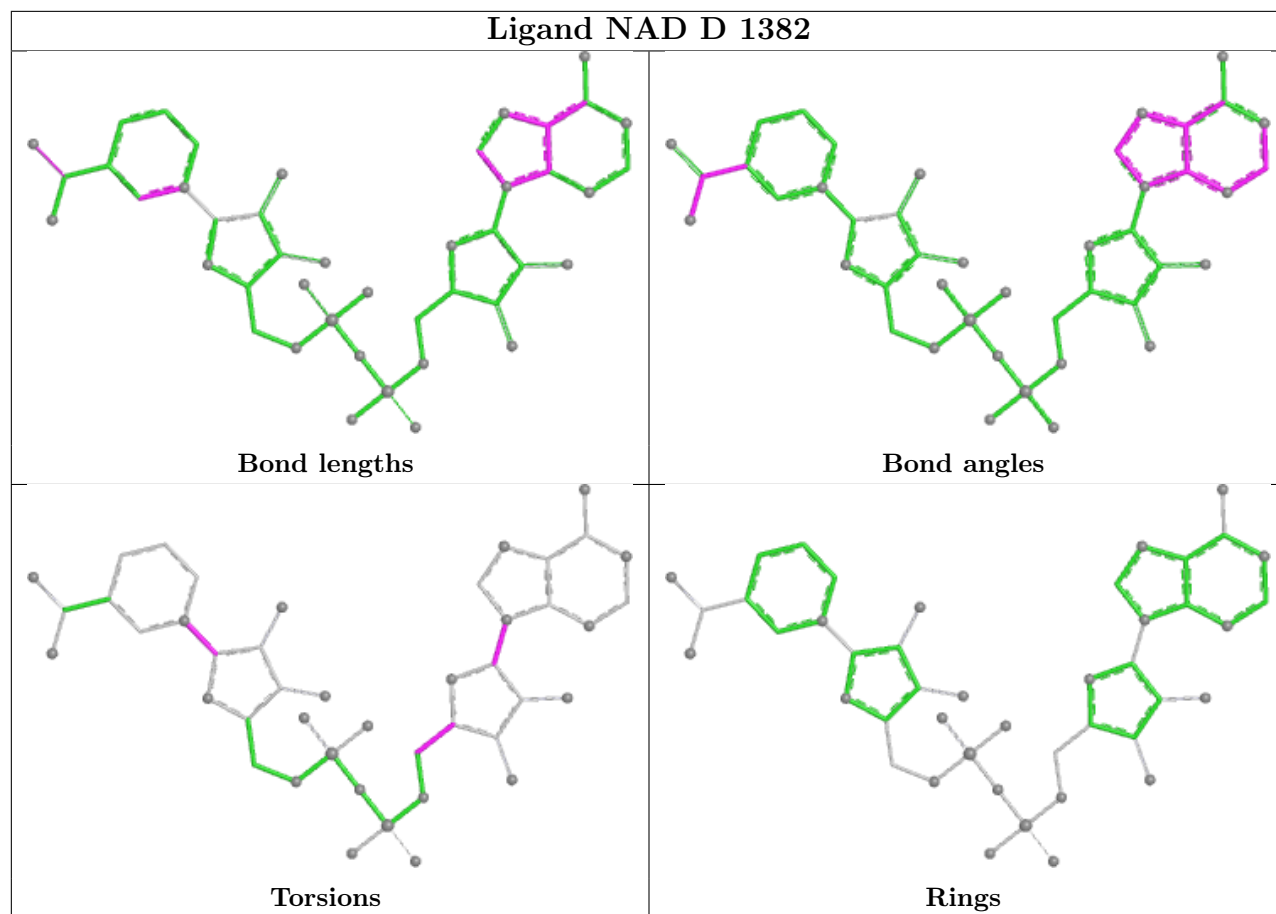
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1383	PEG	1	0
2	D	1381	PO4	1	0
2	C	1378	PO4	1	0
4	D	1385	PEG	1	0
3	C	1381	NAD	1	0
7	B	1384	PGE	2	0
2	C	1379	PO4	1	0
3	A	1381	NAD	1	0
5	B	1385	GLY	1	0
2	C	1380	PO4	1	0
2	B	1378	PO4	1	0
2	A	1378	PO4	1	0
2	B	1379	PO4	2	0
6	B	1383	2PE	2	0
4	D	1384	PEG	1	0
2	A	1379	PO4	1	0
2	B	1380	PO4	1	0

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









5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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	365/378 (96%)	-0.06	5 (1%) 73 69	20, 31, 48, 68	0
1	B	365/378 (96%)	0.00	5 (1%) 73 69	19, 34, 52, 64	0
1	C	365/378 (96%)	0.06	7 (1%) 66 62	19, 36, 56, 84	0
1	D	365/378 (96%)	-0.30	1 (0%) 90 88	18, 28, 41, 53	0
All	All	1460/1512 (96%)	-0.07	18 (1%) 76 73	18, 32, 51, 84	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	13	PRO	4.0
1	C	71	ASN	4.0
1	A	361	ASP	3.2
1	B	13	PRO	3.1
1	C	273	TYR	2.7
1	B	356	CYS	2.6
1	C	72	GLU	2.4
1	D	13	PRO	2.4
1	A	272	GLY	2.4
1	C	38	ASP	2.4
1	A	13	PRO	2.3
1	C	39	LEU	2.3
1	C	172	TYR	2.1
1	B	227	SER	2.1
1	B	106	ARG	2.1
1	A	271	PHE	2.0
1	B	72	GLU	2.0
1	A	37	PRO	2.0

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 [i](#)

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($\text{\AA}^2$)	Q<0.9
5	GLY	D	1386	5/5	0.84	0.13	45,45,46,48	0
4	PEG	A	1383	7/7	0.86	0.12	37,38,42,42	0
2	PO4	B	1381	5/5	0.87	0.12	56,59,62,66	0
4	PEG	D	1385	7/7	0.88	0.12	37,38,44,47	0
5	GLY	D	1387	5/5	0.88	0.15	47,49,52,55	0
8	PG4	C	1382	13/13	0.88	0.11	35,42,46,47	0
2	PO4	D	1381	5/5	0.89	0.13	64,68,70,72	0
2	PO4	B	1380	5/5	0.89	0.17	60,61,65,69	0
4	PEG	D	1383	7/7	0.89	0.10	40,43,45,46	0
4	PEG	D	1384	7/7	0.89	0.10	40,41,42,43	0
6	2PE	B	1383	25/28	0.91	0.10	31,34,41,43	0
7	PGE	B	1384	10/10	0.92	0.09	37,40,41,42	0
2	PO4	A	1380	5/5	0.92	0.15	32,33,39,42	0
2	PO4	D	1380	5/5	0.94	0.15	40,40,46,50	0
4	PEG	A	1382	7/7	0.95	0.09	39,40,42,42	0
3	NAD	A	1381	44/44	0.95	0.07	19,28,34,34	0
3	NAD	B	1382	44/44	0.96	0.07	21,24,28,30	0
3	NAD	C	1381	44/44	0.96	0.06	20,27,34,38	0
3	NAD	D	1382	44/44	0.96	0.07	17,21,23,24	0
2	PO4	C	1380	5/5	0.96	0.11	37,37,45,46	0
5	GLY	B	1385	5/5	0.96	0.07	25,27,31,35	0
5	GLY	C	1383	5/5	0.97	0.05	26,29,31,33	0
5	GLY	A	1384	5/5	0.97	0.06	19,20,21,21	0
2	PO4	A	1378	5/5	0.97	0.06	27,30,32,32	0
2	PO4	B	1378	5/5	0.98	0.04	28,28,30,31	0
2	PO4	C	1378	5/5	0.98	0.06	32,33,36,39	0
2	PO4	B	1379	5/5	0.98	0.04	32,32,35,35	0

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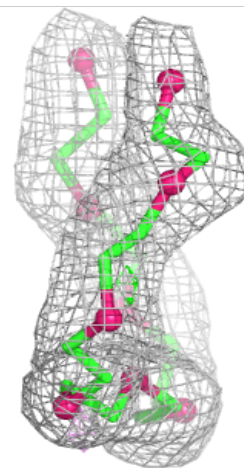
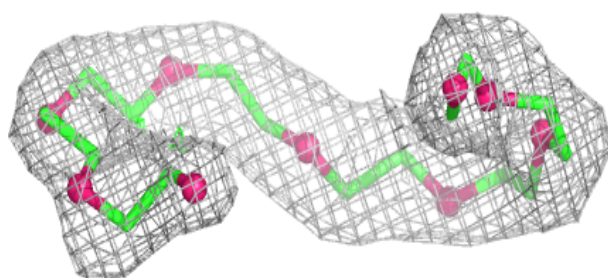
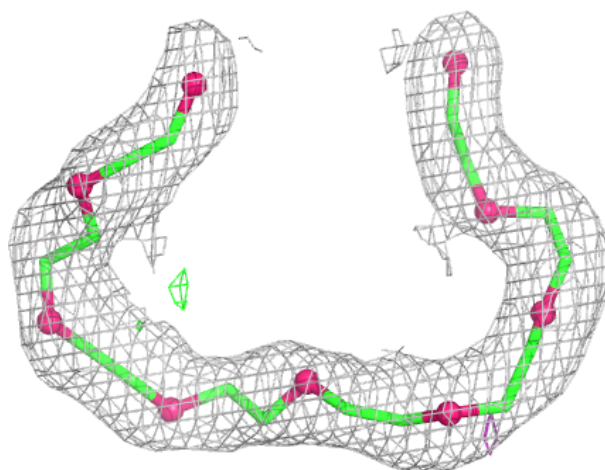
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	PO4	D	1378	5/5	0.98	0.06	27,28,30,31	0
2	PO4	D	1379	5/5	0.98	0.05	25,26,28,28	0
2	PO4	A	1379	5/5	0.98	0.05	28,28,32,33	0
2	PO4	C	1379	5/5	0.99	0.04	24,24,25,26	0

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.

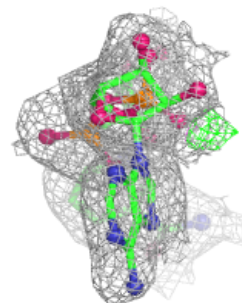
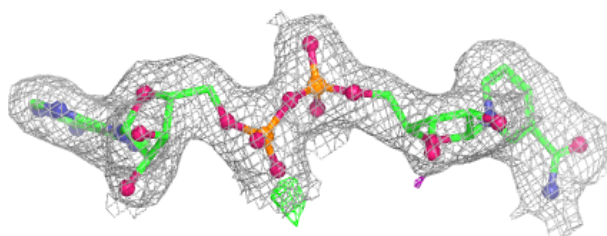
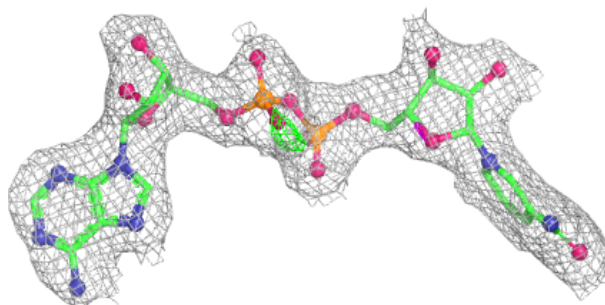
Electron density around 2PE B 1383:

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

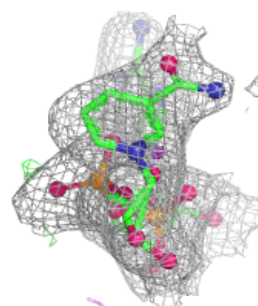
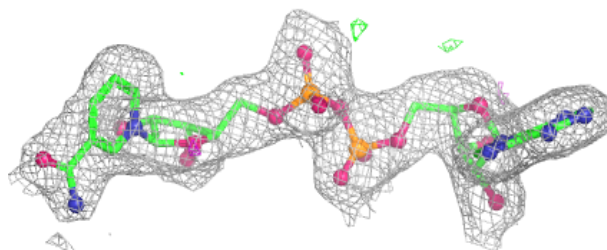
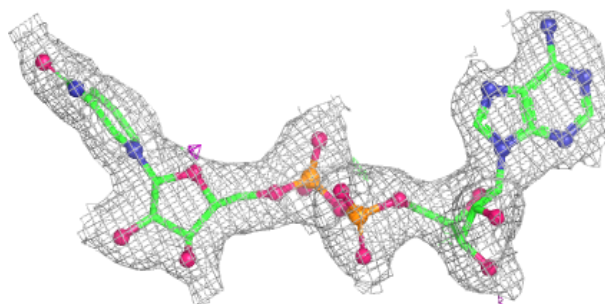


Electron density around NAD A 1381:

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

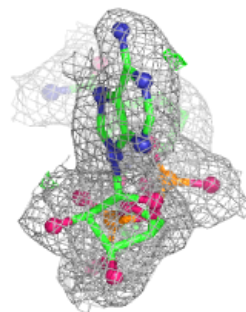
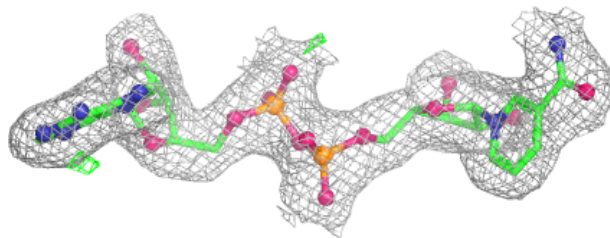
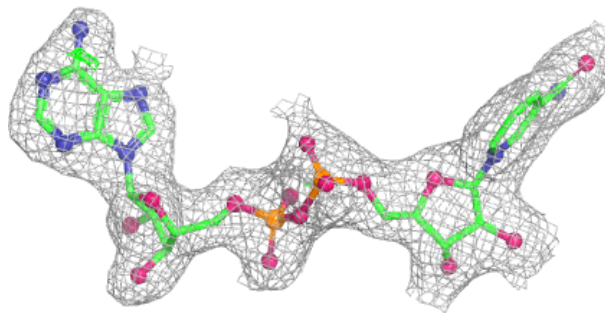
**Electron density around NAD B 1382:**

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

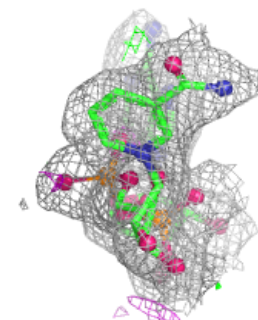
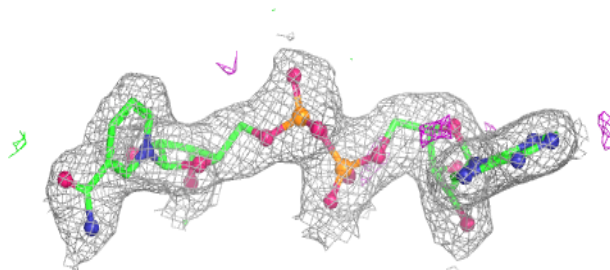
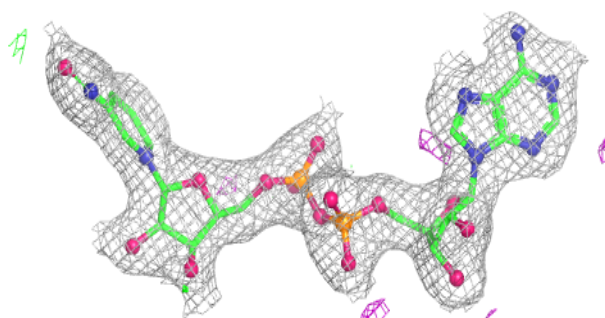


Electron density around NAD C 1381:

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

**Electron density around NAD D 1382:**

$2mF_o-DF_c$ (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.