



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 12:51 PM UTC

PDB ID : 3Q6I / pdb_00003q6i
Title : Crystal structure of FabG4 and coenzyme binary complex
Authors : Dutta, D.; Bhattacharyya, S.; Das, A.K.
Deposited on : 2011-01-01
Resolution : 2.59 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

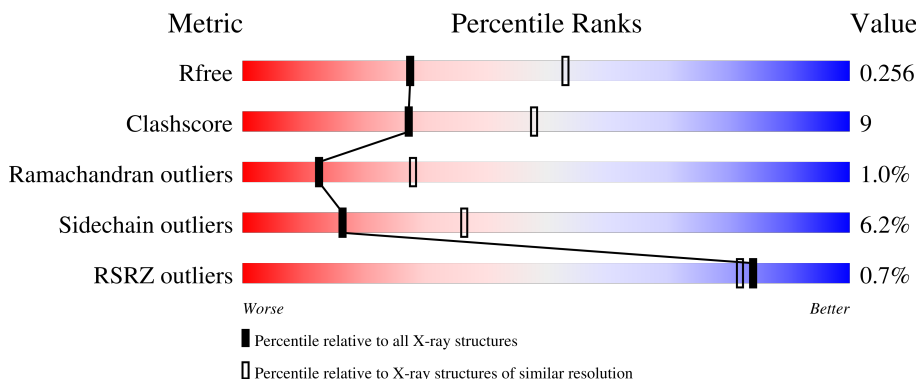
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.59 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	446	 74% 18% • •
1	B	446	 74% 20% • 5%
1	C	446	 74% 17% • 6%
1	D	446	 2% 68% 22% • 8%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 12462 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-oxoacyl-(Acyl-carrier-protein) reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	427	Total	C	N	O	S	0	0	0
			3080	1925	553	595	7			
1	B	425	Total	C	N	O	S	0	0	0
			3059	1916	543	593	7			
1	C	418	Total	C	N	O	S	0	0	0
			2991	1871	530	584	6			
1	D	409	Total	C	N	O	S	0	0	0
			2927	1836	514	571	6			

There are 32 discrepancies between the modelled and reference sequences:

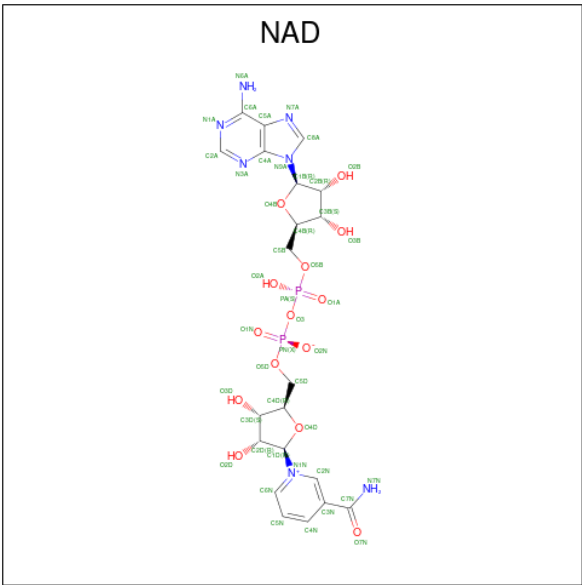
Chain	Residue	Modelled	Actual	Comment	Reference
A	9	HIS	-	expression tag	UNP O53665
A	10	HIS	-	expression tag	UNP O53665
A	11	HIS	-	expression tag	UNP O53665
A	12	HIS	-	expression tag	UNP O53665
A	13	HIS	-	expression tag	UNP O53665
A	14	HIS	-	expression tag	UNP O53665
A	15	GLY	-	expression tag	UNP O53665
A	16	SER	-	expression tag	UNP O53665
B	9	HIS	-	expression tag	UNP O53665
B	10	HIS	-	expression tag	UNP O53665
B	11	HIS	-	expression tag	UNP O53665
B	12	HIS	-	expression tag	UNP O53665
B	13	HIS	-	expression tag	UNP O53665
B	14	HIS	-	expression tag	UNP O53665
B	15	GLY	-	expression tag	UNP O53665
B	16	SER	-	expression tag	UNP O53665
C	9	HIS	-	expression tag	UNP O53665
C	10	HIS	-	expression tag	UNP O53665
C	11	HIS	-	expression tag	UNP O53665
C	12	HIS	-	expression tag	UNP O53665
C	13	HIS	-	expression tag	UNP O53665

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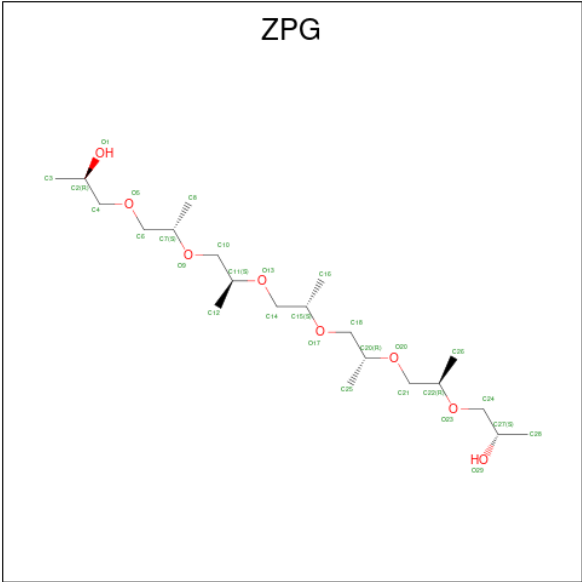
Chain	Residue	Modelled	Actual	Comment	Reference
C	14	HIS	-	expression tag	UNP O53665
C	15	GLY	-	expression tag	UNP O53665
C	16	SER	-	expression tag	UNP O53665
D	9	HIS	-	expression tag	UNP O53665
D	10	HIS	-	expression tag	UNP O53665
D	11	HIS	-	expression tag	UNP O53665
D	12	HIS	-	expression tag	UNP O53665
D	13	HIS	-	expression tag	UNP O53665
D	14	HIS	-	expression tag	UNP O53665
D	15	GLY	-	expression tag	UNP O53665
D	16	SER	-	expression tag	UNP O53665

- Molecule 2 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (CCD ID: NAD) (formula: C₂₁H₂₇N₇O₁₄P₂).



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

- Molecule 3 is (2S,5R,8R,11S,14S,17S,21R)-5,8,11,14,17-PENTAMETHYL-4,7,10,13,16,19-H EXAOXADOCOSANE-2,21-DIOL (CCD ID: ZPG) (formula: C₂₁H₄₄O₈).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			13	9	4		

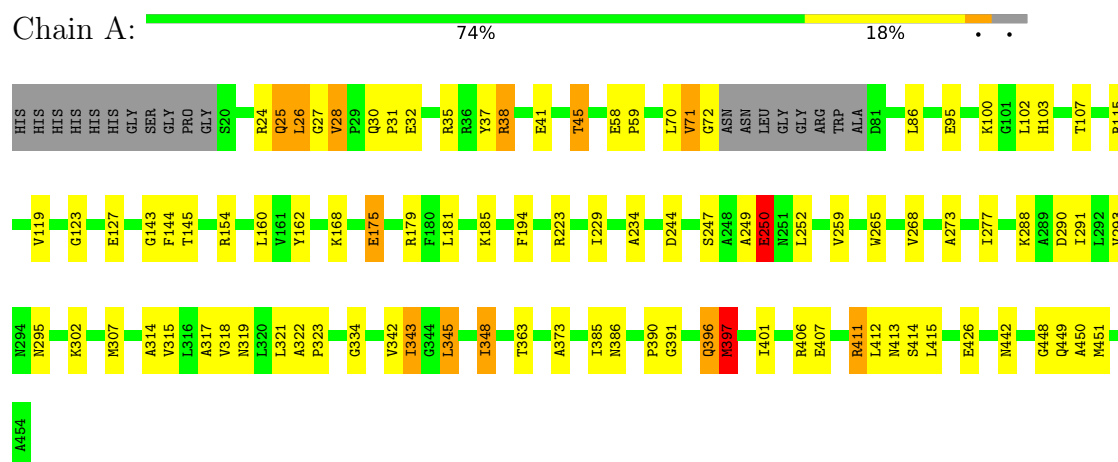
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	78	Total	O	0	0
			78	78		
4	B	69	Total	O	0	0
			69	69		
4	C	39	Total	O	0	0
			39	39		
4	D	30	Total	O	0	0
			30	30		

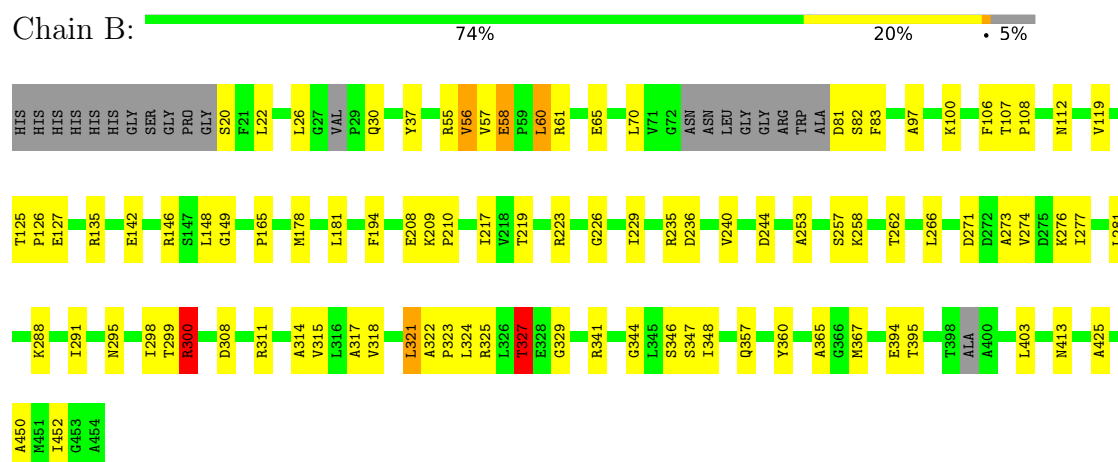
3 Residue-property plots

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

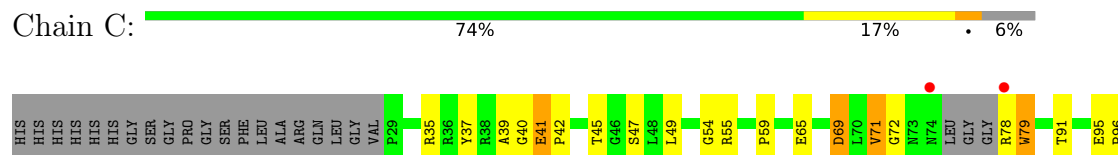
- Molecule 1: 3-oxoacyl-(Acyl-carrier-protein) reductase

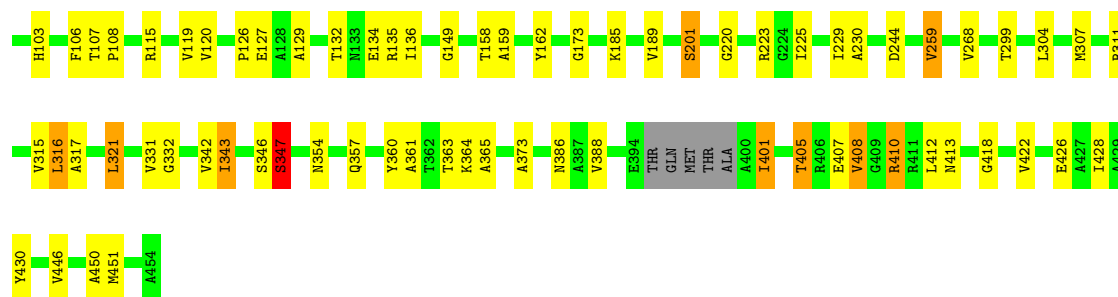


- Molecule 1: 3-oxoacyl-(Acyl-carrier-protein) reductase

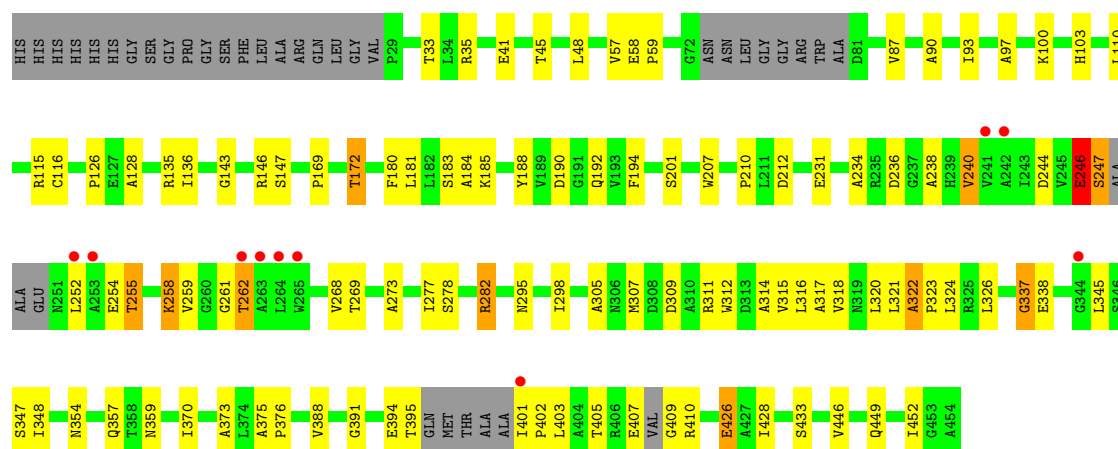


- Molecule 1: 3-oxoacyl-(Acyl-carrier-protein) reductase





- Molecule 1: 3-oxoacyl-(Acyl-carrier-protein) reductase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	63.07Å 71.03Å 92.92Å 105.02° 97.06° 93.66°	Depositor
Resolution (Å)	19.78 – 2.59 19.78 – 2.59	Depositor EDS
% Data completeness (in resolution range)	95.5 (19.78-2.59) 95.5 (19.78-2.59)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.15	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.79 (at 2.59Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.168 , 0.242 0.195 , 0.256	Depositor DCC
R_{free} test set	2322 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	40.3	Xtriage
Anisotropy	0.129	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 41.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	12462	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.81% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: ZPG, 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	0.92	0/3125	1.11	7/4248 (0.2%)
1	B	0.91	0/3102	1.09	9/4212 (0.2%)
1	C	0.87	0/3036	1.06	3/4132 (0.1%)
1	D	0.80	0/2968	1.05	9/4032 (0.2%)
All	All	0.88	0/12231	1.08	28/16624 (0.2%)

There are no bond length outliers.

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	125	THR	CA-C-N	-8.03	110.62	118.97
1	B	125	THR	C-N-CA	-8.03	110.62	118.97
1	A	175	GLU	N-CA-C	6.84	118.82	111.36
1	C	365	ALA	N-CA-C	-6.65	104.11	111.36
1	D	192	GLN	N-CA-C	6.14	117.97	110.41
1	A	396	GLN	N-CA-C	-6.14	100.57	109.59
1	D	143	GLY	N-CA-C	-5.95	105.60	112.50
1	B	365	ALA	N-CA-C	-5.82	105.02	111.36
1	C	347	SER	N-CA-C	5.77	123.08	110.80
1	D	246	GLU	N-CA-C	-5.64	105.22	111.36
1	B	327	THR	CB-CA-C	5.63	119.64	110.96
1	A	143	GLY	N-CA-C	-5.58	105.89	113.37
1	A	70	LEU	N-CA-C	5.57	117.17	109.15
1	D	322	ALA	CA-C-N	-5.55	112.89	119.05
1	D	322	ALA	C-N-CA	-5.55	112.89	119.05
1	D	240	VAL	N-CA-C	5.36	115.67	108.17
1	A	250	GLU	N-CA-C	-5.32	99.46	110.80
1	B	209	LYS	CA-C-N	5.30	124.93	119.05
1	B	209	LYS	C-N-CA	5.30	124.93	119.05
1	D	410	ARG	N-CA-C	-5.29	106.77	114.12

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	65	GLU	N-CA-C	5.19	116.94	111.28
1	A	107	THR	CA-C-N	-5.13	115.05	121.00
1	A	107	THR	C-N-CA	-5.13	115.05	121.00
1	B	240	VAL	N-CA-C	5.13	115.48	107.99
1	D	136	ILE	N-CA-C	-5.12	105.55	110.72
1	D	403	LEU	N-CA-C	5.04	117.16	111.02
1	C	115	ARG	N-CA-C	-5.01	102.86	110.28
1	B	83	PHE	N-CA-C	5.00	116.91	109.25

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	3080	0	3100	64	0
1	B	3059	0	3073	59	0
1	C	2991	0	2963	55	0
1	D	2927	0	2916	57	0
2	A	44	0	26	4	0
2	B	44	0	26	3	0
2	C	44	0	26	1	0
2	D	44	0	26	1	0
3	A	13	0	19	0	0
4	A	78	0	0	1	0
4	B	69	0	0	4	0
4	C	39	0	0	1	0
4	D	30	0	0	0	0
All	All	12462	0	12175	222	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 9.

All (222) 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:38:ARG:HG2	1:A:38:ARG:HH11	0.96	1.10
1:D:337:GLY:HA2	1:D:338:GLU:HB2	1.51	0.92
1:A:38:ARG:HH11	1:A:38:ARG:CG	1.82	0.92
1:B:299:THR:O	1:B:300:ARG:HB2	1.72	0.89
1:C:173:GLY:HA2	1:C:201:SER:OG	1.72	0.89
1:A:38:ARG:HG2	1:A:38:ARG:NH1	1.75	0.88
1:D:407:GLU:O	1:D:409:GLY:N	2.07	0.88
1:A:315:VAL:O	1:A:319:ASN:HB2	1.79	0.82
1:B:395:THR:HG21	2:B:3:NAD:H72N	1.46	0.81
1:B:298:ILE:HG12	1:B:300:ARG:HH12	1.47	0.79
1:D:337:GLY:CA	1:D:338:GLU:HB2	2.13	0.78
1:A:396:GLN:O	1:A:397:MET:HB3	1.86	0.76
1:A:249:ALA:O	1:A:250:GLU:HB2	1.91	0.71
1:A:343:ILE:HD12	1:A:343:ILE:N	2.05	0.71
1:D:246:GLU:O	1:D:247:SER:C	2.32	0.70
1:B:82:SER:HA	1:B:112:ASN:ND2	2.06	0.70
1:B:82:SER:HA	1:B:112:ASN:HD21	1.55	0.70
1:C:311:ARG:NE	4:C:476:HOH:O	2.25	0.70
1:B:298:ILE:CG1	1:B:300:ARG:HH12	2.05	0.69
1:C:129:ALA:HB1	1:C:134:GLU:HB3	1.75	0.68
1:B:308:ASP:OD1	1:B:311:ARG:HD2	1.94	0.67
1:C:316:LEU:HD13	1:C:363:THR:HG22	1.77	0.66
1:D:255:THR:O	1:D:259:VAL:HG22	1.95	0.66
1:C:49:LEU:HD13	1:C:71:VAL:CG2	2.26	0.66
1:B:322:ALA:HB3	1:B:323:PRO:HD3	1.77	0.65
1:D:322:ALA:O	1:D:326:LEU:HG	1.97	0.65
1:D:278:SER:HB3	1:D:282:ARG:HH21	1.63	0.64
1:A:401:ILE:O	1:A:406:ARG:NH1	2.31	0.63
1:C:41:GLU:HB2	1:C:42:PRO:HD2	1.80	0.63
1:A:268:VAL:HG22	2:A:1:NAD:N1A	2.13	0.63
1:A:223:ARG:HD2	4:A:500:HOH:O	1.98	0.62
1:A:426:GLU:OE1	1:A:426:GLU:HA	1.98	0.62
1:A:295:ASN:HA	1:A:345:LEU:HD13	1.80	0.62
1:B:323:PRO:O	1:B:327:THR:HG22	1.99	0.62
1:D:259:VAL:HG23	1:D:261:GLY:H	1.64	0.61
1:B:219:THR:O	1:B:295:ASN:HB3	2.00	0.61
1:B:323:PRO:O	1:B:327:THR:CG2	2.50	0.60
1:A:25:GLN:HB3	1:B:22:LEU:HD11	1.83	0.60
1:A:396:GLN:O	1:A:396:GLN:HG3	2.00	0.60
1:B:344:GLY:HA3	4:B:520:HOH:O	2.01	0.58
1:A:414:SER:HB2	1:A:449:GLN:HB3	1.85	0.57
1:B:60:LEU:HD11	1:B:178:MET:HG3	1.85	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:322:ALA:HB3	1:A:323:PRO:HD3	1.86	0.57
1:C:132:THR:O	1:C:136:ILE:HG13	2.04	0.57
1:D:337:GLY:HA2	1:D:338:GLU:CB	2.32	0.57
1:B:298:ILE:CG1	1:B:300:ARG:NH1	2.67	0.57
1:C:103:HIS:HE1	1:D:307:MET:O	1.88	0.57
1:A:415:LEU:HD12	1:A:448:GLY:HA2	1.87	0.56
1:A:32:GLU:OE2	1:A:115:ARG:HG2	2.06	0.56
1:B:298:ILE:HG13	1:B:300:ARG:NH1	2.19	0.56
1:A:100:LYS:O	1:A:103:HIS:HB3	2.06	0.56
1:C:230:ALA:HB1	1:C:259:VAL:HG11	1.88	0.56
1:D:354:ASN:HB3	1:D:357:GLN:HG3	1.87	0.55
1:A:38:ARG:CG	1:A:38:ARG:NH1	2.52	0.55
1:B:299:THR:O	1:B:300:ARG:CB	2.48	0.55
1:D:185:LYS:HB3	1:D:426:GLU:HG2	1.88	0.55
1:C:126:PRO:O	1:C:135:ARG:HG3	2.07	0.55
1:A:373:ALA:CB	1:B:135:ARG:HD3	2.37	0.54
1:C:413:ASN:HB3	1:C:450:ALA:HA	1.89	0.54
1:B:413:ASN:HB3	1:B:450:ALA:HA	1.89	0.54
1:D:172:THR:HB	1:D:201:SER:H	1.73	0.54
1:A:86:LEU:O	1:A:119:VAL:HA	2.07	0.54
1:C:49:LEU:CD1	1:C:71:VAL:HG22	2.37	0.54
1:C:49:LEU:HD13	1:C:71:VAL:HG22	1.89	0.54
1:C:189:VAL:HG22	1:C:446:VAL:HG22	1.90	0.54
1:B:314:ALA:O	1:B:318:VAL:HG23	2.08	0.53
1:D:126:PRO:C	1:D:128:ALA:H	2.16	0.53
1:D:273:ALA:O	1:D:277:ILE:HG13	2.08	0.53
1:D:181:LEU:HD21	1:D:194:PHE:CZ	2.44	0.53
1:A:26:LEU:HD23	1:B:22:LEU:HD13	1.91	0.53
1:D:295:ASN:HA	1:D:345:LEU:HD13	1.90	0.52
1:B:107:THR:HB	1:B:108:PRO:HD3	1.91	0.52
1:D:48:LEU:HD11	1:D:87:VAL:HG23	1.91	0.52
1:B:226:GLY:HA2	1:B:229:ILE:HD12	1.90	0.52
1:C:135:ARG:HD3	1:D:373:ALA:CB	2.40	0.52
1:D:348:ILE:HD12	1:D:391:GLY:HA2	1.90	0.52
1:B:311:ARG:O	1:B:315:VAL:HG23	2.10	0.52
1:A:115:ARG:NH2	1:A:154:ARG:HD2	2.24	0.52
1:C:149:GLY:HA2	1:C:158:THR:HG23	1.91	0.51
1:A:373:ALA:HB2	1:B:135:ARG:HD3	1.92	0.51
1:D:317:ALA:HA	1:D:321:LEU:HB2	1.91	0.51
1:A:234:ALA:HB2	1:A:259:VAL:HB	1.92	0.51
1:D:188:TYR:HD2	1:D:446:VAL:HG12	1.76	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:295:ASN:HB2	1:A:345:LEU:HD13	1.91	0.51
1:B:56:VAL:O	1:B:60:LEU:HB2	2.11	0.51
1:C:47:SER:HA	1:C:69:ASP:HB3	1.91	0.51
1:C:315:VAL:HG12	1:C:363:THR:HG21	1.91	0.51
2:B:3:NAD:N7N	2:B:3:NAD:O1N	2.43	0.50
1:B:273:ALA:O	1:B:277:ILE:HG13	2.11	0.50
1:C:346:SER:O	1:C:347:SER:O	2.29	0.50
1:A:26:LEU:O	1:A:28:VAL:N	2.45	0.50
1:D:115:ARG:NH1	1:D:115:ARG:HB3	2.26	0.50
1:D:269:THR:HG21	1:D:317:ALA:HB1	1.94	0.50
1:A:102:LEU:HD13	1:A:144:PHE:CG	2.46	0.50
1:B:181:LEU:HD21	1:B:194:PHE:CZ	2.47	0.50
1:C:304:LEU:HD23	1:D:110:LEU:HD11	1.92	0.50
1:D:190:ASP:HB3	1:D:449:GLN:N	2.27	0.50
1:D:231:GLU:OE2	1:D:258:LYS:HD3	2.11	0.50
1:A:268:VAL:HG22	2:A:1:NAD:C6A	2.41	0.50
1:B:274:VAL:HG21	1:B:329:GLY:HA3	1.94	0.50
1:A:30:GLN:NE2	1:A:31:PRO:HD2	2.27	0.50
1:A:244:ASP:HB2	1:A:252:LEU:HD13	1.94	0.50
1:D:298:ILE:HD11	1:D:314:ALA:O	2.11	0.50
1:D:246:GLU:O	1:D:247:SER:O	2.30	0.49
1:C:107:THR:HB	1:C:108:PRO:HD3	1.93	0.49
1:D:347:SER:HB3	2:D:4:NAD:H6N	1.94	0.49
1:D:180:PHE:O	1:D:183:SER:OG	2.27	0.49
1:A:342:VAL:O	1:A:385:ILE:HA	2.13	0.49
1:D:35:ARG:NH1	1:D:184:ALA:HB2	2.28	0.49
1:B:208:GLU:C	1:B:210:PRO:HD3	2.38	0.48
1:A:451:MET:O	1:B:452:ILE:HA	2.14	0.48
1:B:55:ARG:CZ	1:B:165:PRO:HA	2.44	0.48
1:B:58:GLU:HB2	4:B:505:HOH:O	2.14	0.47
1:C:225:ILE:HG22	1:C:229:ILE:HD11	1.96	0.47
1:A:86:LEU:HB2	1:A:119:VAL:HG22	1.95	0.47
1:C:106:PHE:O	1:C:107:THR:C	2.56	0.47
1:D:320:LEU:C	1:D:323:PRO:HD2	2.40	0.47
1:A:319:ASN:HB3	1:A:363:THR:CB	2.45	0.47
1:A:319:ASN:HB3	1:A:363:THR:HB	1.96	0.47
1:D:402:PRO:O	1:D:405:THR:HB	2.14	0.47
1:A:145:THR:HG21	1:A:160:LEU:HB2	1.97	0.47
1:D:115:ARG:HB3	1:D:115:ARG:CZ	2.46	0.46
1:D:244:ASP:HB2	1:D:252:LEU:HD13	1.97	0.46
1:A:181:LEU:HD21	1:A:194:PHE:CZ	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:311:ARG:O	1:D:315:VAL:HG23	2.15	0.46
1:A:413:ASN:HB3	1:A:450:ALA:HA	1.97	0.46
1:D:259:VAL:HG23	1:D:261:GLY:N	2.31	0.46
1:C:299:THR:HG22	1:C:360:TYR:CD1	2.50	0.46
1:C:401:ILE:HD13	1:C:401:ILE:H	1.81	0.46
1:A:295:ASN:CB	1:A:345:LEU:HD13	2.45	0.46
1:A:342:VAL:HG12	1:A:385:ILE:HG13	1.96	0.46
1:C:408:VAL:O	1:C:412:LEU:HB2	2.15	0.46
1:D:146:ARG:HG2	1:D:146:ARG:HH11	1.81	0.46
1:C:119:VAL:O	1:C:158:THR:HA	2.16	0.45
1:B:210:PRO:HD2	1:B:236:ASP:HB3	1.97	0.45
1:A:123:GLY:O	1:A:162:TYR:HA	2.16	0.45
1:C:35:ARG:NH2	1:C:42:PRO:O	2.50	0.45
1:C:422:VAL:O	1:C:426:GLU:HG2	2.16	0.45
1:B:291:ILE:HA	1:B:341:ARG:O	2.16	0.45
1:B:258:LYS:HE2	4:B:500:HOH:O	2.16	0.45
1:B:298:ILE:HD12	1:B:298:ILE:HA	1.89	0.45
1:D:210:PRO:HD2	1:D:236:ASP:HB3	1.99	0.45
1:A:229:ILE:HD13	1:A:293:VAL:HG11	1.99	0.45
1:C:54:GLY:HA3	1:C:91:THR:OG1	2.17	0.44
1:C:410:ARG:HG3	1:C:418:GLY:H	1.82	0.44
1:B:324:LEU:O	1:B:325:ARG:C	2.60	0.44
1:D:126:PRO:C	1:D:128:ALA:N	2.75	0.44
1:A:415:LEU:HD12	1:A:448:GLY:CA	2.46	0.44
1:C:373:ALA:HB2	1:D:135:ARG:HG2	1.99	0.44
1:A:175:GLU:HG2	1:A:179:ARG:HD2	1.99	0.44
1:A:407:GLU:OE2	1:A:411:ARG:HD3	2.16	0.44
1:B:253:ALA:O	1:B:257:SER:HB2	2.17	0.44
1:B:148:LEU:O	1:B:149:GLY:C	2.60	0.44
1:C:405:THR:O	1:C:405:THR:OG1	2.34	0.44
1:C:220:GLY:HA2	1:C:244:ASP:OD1	2.18	0.44
1:C:361:ALA:CB	1:D:147:SER:OG	2.66	0.44
1:C:95:GLU:N	1:C:95:GLU:OE1	2.50	0.43
1:C:120:VAL:HG22	1:C:159:ALA:HB3	1.99	0.43
1:C:346:SER:HA	1:C:364:LYS:HE3	2.00	0.43
1:A:244:ASP:OD2	2:A:1:NAD:H1B	2.17	0.43
1:A:288:LYS:NZ	1:A:334:GLY:O	2.51	0.43
1:A:390:PRO:HB2	2:A:1:NAD:C5N	2.49	0.43
1:B:37:TYR:CE2	1:B:235:ARG:HD2	2.54	0.43
1:D:388:VAL:HG21	1:D:428:ILE:HG13	1.98	0.43
1:A:386:ASN:OD1	1:A:442:ASN:HB2	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:318:VAL:O	1:D:323:PRO:HD3	2.18	0.43
1:C:55:ARG:O	1:C:59:PRO:HG2	2.17	0.43
1:B:57:VAL:O	1:B:61:ARG:HG3	2.17	0.43
1:D:312:TRP:CE3	1:D:359:ASN:HB3	2.54	0.43
1:A:314:ALA:O	1:A:318:VAL:HG23	2.18	0.43
1:C:354:ASN:O	1:D:147:SER:HB3	2.19	0.43
1:C:331:VAL:O	1:C:332:GLY:C	2.62	0.43
1:A:317:ALA:HA	1:A:321:LEU:HB2	2.02	0.42
1:B:452:ILE:HD13	1:B:452:ILE:HG21	1.78	0.42
1:C:307:MET:O	1:D:103:HIS:HE1	2.02	0.42
1:A:273:ALA:O	1:A:277:ILE:HG13	2.20	0.42
1:C:127:GLU:H	1:C:127:GLU:CD	2.27	0.42
1:B:223:ARG:HD3	2:B:3:NAD:O2A	2.19	0.42
1:D:57:VAL:HG13	1:D:58:GLU:HG2	2.00	0.42
1:D:234:ALA:HA	1:D:238:ALA:O	2.19	0.42
1:C:346:SER:O	1:C:347:SER:C	2.63	0.42
1:D:59:PRO:HG3	1:D:169:PRO:O	2.19	0.42
1:B:127:GLU:CD	1:B:127:GLU:H	2.27	0.42
1:C:342:VAL:C	1:C:343:ILE:HG13	2.44	0.42
1:B:97:ALA:O	1:B:100:LYS:HG2	2.19	0.42
1:B:26:LEU:HD23	1:B:26:LEU:HA	1.91	0.42
1:B:357:GLN:HB3	1:B:360:TYR:HB3	2.02	0.42
1:B:323:PRO:HG2	1:B:367:MET:HE1	2.01	0.41
1:C:268:VAL:HG22	2:C:2:NAD:N1A	2.34	0.41
1:D:207:TRP:O	1:D:210:PRO:HG3	2.20	0.41
1:A:295:ASN:CA	1:A:345:LEU:HD13	2.48	0.41
1:C:357:GLN:HB3	1:C:360:TYR:HB3	2.02	0.41
1:A:290:ASP:C	1:A:291:ILE:HG13	2.44	0.41
1:D:90:ALA:HB1	1:D:93:ILE:HD12	2.01	0.41
1:A:26:LEU:CD2	1:B:22:LEU:HD13	2.50	0.41
1:B:142:GLU:OE2	1:B:146:ARG:NE	2.40	0.41
1:C:126:PRO:HD3	1:C:162:TYR:CE1	2.55	0.41
1:A:26:LEU:HD23	1:B:22:LEU:CD1	2.50	0.41
1:C:317:ALA:HA	1:C:321:LEU:HB2	2.02	0.41
1:A:37:TYR:CD1	1:A:185:LYS:HE2	2.54	0.41
1:C:37:TYR:CD1	1:C:185:LYS:HE3	2.56	0.41
1:C:78:ARG:O	1:C:79:TRP:CB	2.67	0.41
1:B:106:PHE:CE1	1:B:119:VAL:HG13	2.56	0.41
1:B:394:GLU:H	1:B:394:GLU:HG2	1.66	0.41
1:C:95:GLU:HB2	1:C:96:PRO:HD2	2.02	0.41
1:C:343:ILE:HA	1:C:386:ASN:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:268:VAL:HB	1:D:326:LEU:HD11	2.02	0.41
1:A:249:ALA:HB2	1:A:265:TRP:CD1	2.56	0.41
1:B:217:ILE:HD11	1:B:281:LEU:HD11	2.03	0.41
1:B:229:ILE:HA	1:B:425:ALA:HB1	2.03	0.41
1:D:324:LEU:HD13	1:D:370:ILE:HG12	2.01	0.41
1:A:58:GLU:HB3	1:A:59:PRO:HD3	2.02	0.41
1:A:348:ILE:HD12	1:A:391:GLY:HA2	2.03	0.41
1:B:317:ALA:HA	1:B:321:LEU:HB2	2.02	0.41
1:D:240:VAL:O	1:D:262:THR:N	2.52	0.41
1:A:71:VAL:HA	1:A:72:GLY:HA2	1.75	0.40
1:C:388:VAL:HG21	1:C:428:ILE:CG1	2.52	0.40
1:B:81:ASP:CB	4:B:514:HOH:O	2.68	0.40
1:D:375:ALA:N	1:D:376:PRO:HD2	2.37	0.40
1:A:35:ARG:CZ	1:A:45:THR:HG23	2.51	0.40
1:B:126:PRO:O	1:B:135:ARG:HG3	2.21	0.40
1:C:127:GLU:OE1	1:C:127:GLU:N	2.45	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	423/446 (95%)	398 (94%)	22 (5%)	3 (1%)	18	38
1	B	417/446 (94%)	393 (94%)	21 (5%)	3 (1%)	18	38
1	C	412/446 (92%)	384 (93%)	22 (5%)	6 (2%)	8	18
1	D	399/446 (90%)	364 (91%)	30 (8%)	5 (1%)	9	21
All	All	1651/1784 (92%)	1539 (93%)	95 (6%)	17 (1%)	12	28

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	250	GLU
1	B	300	ARG
1	C	347	SER
1	A	397	MET
1	C	39	ALA
1	C	40	GLY
1	C	45	THR
1	D	337	GLY
1	B	346	SER
1	B	347	SER
1	D	116	CYS
1	C	79	TRP
1	D	97	ALA
1	D	100	LYS
1	D	305	ALA
1	A	27	GLY
1	C	72	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	306/323 (95%)	286 (94%)	20 (6%)	15	35
1	B	303/323 (94%)	286 (94%)	17 (6%)	19	40
1	C	292/323 (90%)	275 (94%)	17 (6%)	18	39
1	D	287/323 (89%)	267 (93%)	20 (7%)	14	31
All	All	1188/1292 (92%)	1114 (94%)	74 (6%)	16	36

All (74) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	ARG
1	A	25	GLN
1	A	26	LEU
1	A	28	VAL
1	A	38	ARG

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Mol	Chain	Res	Type
1	A	41	GLU
1	A	45	THR
1	A	71	VAL
1	A	95	GLU
1	A	127	GLU
1	A	168	LYS
1	A	247	SER
1	A	302	LYS
1	A	307	MET
1	A	343	ILE
1	A	345	LEU
1	A	348	ILE
1	A	397	MET
1	A	411	ARG
1	A	412	LEU
1	B	20	SER
1	B	30	GLN
1	B	56	VAL
1	B	58	GLU
1	B	60	LEU
1	B	70	LEU
1	B	244	ASP
1	B	262	THR
1	B	266	LEU
1	B	271	ASP
1	B	276	LYS
1	B	288	LYS
1	B	300	ARG
1	B	321	LEU
1	B	327	THR
1	B	348	ILE
1	B	403	LEU
1	C	41	GLU
1	C	65	GLU
1	C	69	ASP
1	C	71	VAL
1	C	201	SER
1	C	223	ARG
1	C	259	VAL
1	C	316	LEU
1	C	321	LEU
1	C	343	ILE

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Mol	Chain	Res	Type
1	C	401	ILE
1	C	405	THR
1	C	407	GLU
1	C	408	VAL
1	C	410	ARG
1	C	430	TYR
1	C	451	MET
1	D	33	THR
1	D	41	GLU
1	D	45	THR
1	D	172	THR
1	D	212	ASP
1	D	246	GLU
1	D	247	SER
1	D	254	GLU
1	D	255	THR
1	D	258	LYS
1	D	262	THR
1	D	282	ARG
1	D	309	ASP
1	D	316	LEU
1	D	394	GLU
1	D	395	THR
1	D	401	ILE
1	D	426	GLU
1	D	433	SER
1	D	452	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	103	HIS
1	A	251	ASN
1	A	294	ASN
1	B	25	GLN
1	B	30	GLN
1	B	112	ASN
1	B	192	GLN
1	B	239	HIS
1	B	284	HIS
1	C	103	HIS
1	C	192	GLN

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Mol	Chain	Res	Type
1	C	251	ASN
1	C	285	HIS
1	C	420	GLN
1	D	192	GLN
1	D	251	ASN
1	D	294	ASN
1	D	306	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](#)

5 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	B	3	-	46,48,48	1.92	6 (13%)	64,73,73	1.76	10 (15%)
2	NAD	D	4	-	46,48,48	1.93	8 (17%)	64,73,73	1.69	11 (17%)
2	NAD	A	1	-	46,48,48	1.97	7 (15%)	64,73,73	1.78	12 (18%)
2	NAD	C	2	-	46,48,48	2.01	7 (15%)	64,73,73	1.77	12 (18%)
3	ZPG	A	455	-	12,12,28	0.72	0	14,14,34	1.20	2 (14%)

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	B	3	-	-	9/30/62/62	0/5/5/5
2	NAD	D	4	-	-	5/30/62/62	0/5/5/5
2	NAD	A	1	-	-	4/30/62/62	0/5/5/5
2	NAD	C	2	-	-	8/30/62/62	0/5/5/5
3	ZPG	A	455	-	-	2/12/12/31	-

All (28) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	3	NAD	O7N-C7N	10.14	1.43	1.24
2	C	2	NAD	O7N-C7N	9.93	1.42	1.24
2	D	4	NAD	O7N-C7N	9.72	1.42	1.24
2	A	1	NAD	O7N-C7N	9.48	1.41	1.24
2	A	1	NAD	C2N-N1N	4.15	1.39	1.35
2	C	2	NAD	PN-O3	3.94	1.63	1.59
2	C	2	NAD	PA-O3	3.83	1.63	1.59
2	A	1	NAD	PA-O3	3.69	1.63	1.59
2	C	2	NAD	C2N-N1N	3.42	1.38	1.35
2	B	3	NAD	PA-O3	3.33	1.63	1.59
2	D	4	NAD	PA-O3	3.29	1.63	1.59
2	D	4	NAD	PN-O3	3.17	1.62	1.59
2	B	3	NAD	C2N-N1N	2.83	1.38	1.35
2	A	1	NAD	C8A-N7A	2.76	1.37	1.31
2	D	4	NAD	C2A-N3A	2.74	1.38	1.33
2	D	4	NAD	C2N-N1N	2.72	1.38	1.35
2	D	4	NAD	C2A-N1A	2.68	1.38	1.33
2	B	3	NAD	C2A-N1A	2.64	1.38	1.33
2	C	2	NAD	C2A-N3A	2.62	1.38	1.33
2	B	3	NAD	C2A-N3A	2.62	1.38	1.33
2	B	3	NAD	PN-O3	2.54	1.62	1.59
2	A	1	NAD	C2A-N1A	2.49	1.38	1.33
2	C	2	NAD	C2A-N1A	2.49	1.38	1.33
2	A	1	NAD	C5A-N7A	-2.45	1.34	1.39
2	D	4	NAD	C5A-N7A	-2.41	1.34	1.39
2	A	1	NAD	O4D-C4D	-2.28	1.39	1.45
2	D	4	NAD	C8A-N7A	2.18	1.35	1.31
2	C	2	NAD	C8A-N7A	2.12	1.35	1.31

All (47) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1	NAD	N3A-C2A-N1A	-7.14	117.78	128.58
2	B	3	NAD	N3A-C2A-N1A	-6.28	119.08	128.58
2	C	2	NAD	N3A-C2A-N1A	-6.13	119.30	128.58
2	D	4	NAD	N3A-C2A-N1A	-5.21	120.70	128.58
2	C	2	NAD	N9A-C8A-N7A	-5.03	106.80	113.94
2	D	4	NAD	C5A-C4A-N3A	-4.84	120.06	126.72
2	B	3	NAD	N9A-C8A-N7A	-4.76	107.18	113.94
2	A	1	NAD	O4B-C1B-N9A	4.50	116.73	108.09
2	D	4	NAD	C6N-N1N-C2N	-4.27	118.24	121.88
2	A	1	NAD	N9A-C8A-N7A	-4.06	108.18	113.94
2	B	3	NAD	C5A-C4A-N3A	-4.02	121.18	126.72
2	B	3	NAD	C6N-N1N-C2N	-3.94	118.52	121.88
2	C	2	NAD	C5A-N7A-C8A	3.92	109.61	103.45
2	B	3	NAD	C5A-N7A-C8A	3.91	109.59	103.45
2	B	3	NAD	C2A-N3A-C4A	3.70	120.86	111.83
2	C	2	NAD	O2A-PA-O3	3.58	116.95	107.27
2	C	2	NAD	C5A-C4A-N3A	-3.56	121.82	126.72
2	A	1	NAD	C2A-N3A-C4A	3.46	120.27	111.83
2	C	2	NAD	C2A-N3A-C4A	3.34	119.98	111.83
2	D	4	NAD	C2A-N3A-C4A	3.32	119.95	111.83
2	D	4	NAD	C4D-O4D-C1D	3.21	112.87	109.92
2	B	3	NAD	C4A-N9A-C8A	3.18	109.08	105.74
2	C	2	NAD	C4A-N9A-C8A	3.12	109.02	105.74
2	D	4	NAD	N3A-C4A-N9A	3.09	132.41	127.17
2	B	3	NAD	N3A-C4A-N9A	3.01	132.28	127.17
2	A	1	NAD	C5A-C4A-N3A	-3.01	122.57	126.72
2	A	1	NAD	C5A-N7A-C8A	2.85	107.93	103.45
2	D	4	NAD	C5A-N7A-C8A	2.80	107.85	103.45
2	D	4	NAD	N9A-C8A-N7A	-2.72	110.07	113.94
2	B	3	NAD	C4D-O4D-C1D	2.61	112.31	109.92
2	B	3	NAD	C4A-C5A-N7A	-2.60	107.61	110.58
2	C	2	NAD	N3A-C4A-N9A	2.56	131.53	127.17
2	D	4	NAD	C6A-C5A-C4A	2.56	120.67	117.18
2	D	4	NAD	C2N-N1N-C1D	2.46	124.57	119.13
2	A	1	NAD	C2A-N1A-C6A	2.44	122.74	118.73
2	C	2	NAD	C6N-N1N-C2N	-2.35	119.88	121.88
3	A	455	ZPG	O20-C20-C18	2.35	114.89	108.64
2	A	1	NAD	O2N-PN-O3	2.33	113.58	107.27
2	C	2	NAD	C4A-C5A-N7A	-2.32	107.93	110.58
2	C	2	NAD	C4A-N9A-C1B	-2.27	121.32	126.63
2	C	2	NAD	C4D-O4D-C1D	-2.26	107.85	109.92
2	A	1	NAD	C4D-O4D-C1D	2.24	111.97	109.92

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	455	ZPG	C21-O20-C20	2.19	118.27	115.05
2	A	1	NAD	O7N-C7N-N7N	-2.14	119.52	122.62
2	A	1	NAD	C4A-N9A-C8A	2.11	107.96	105.74
2	A	1	NAD	C4A-N9A-C1B	-2.11	121.71	126.63
2	D	4	NAD	C4A-C5A-N7A	-2.02	108.28	110.58

There are no chirality outliers.

All (28) torsion outliers are listed below:

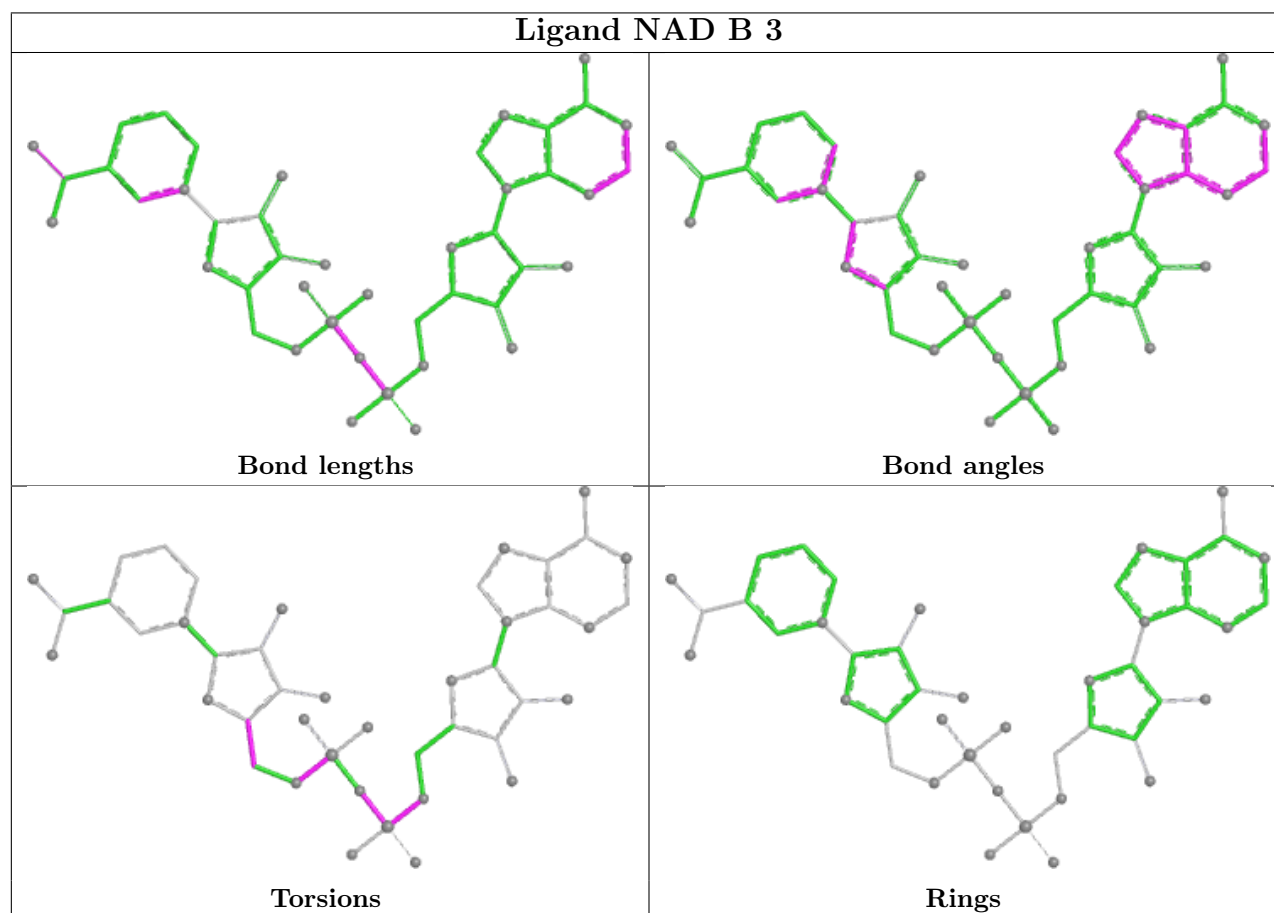
Mol	Chain	Res	Type	Atoms
2	A	1	NAD	O4D-C1D-N1N-C2N
2	B	3	NAD	C5B-O5B-PA-O1A
2	B	3	NAD	C5B-O5B-PA-O2A
2	B	3	NAD	C5B-O5B-PA-O3
2	B	3	NAD	C5D-O5D-PN-O3
2	C	2	NAD	C5D-O5D-PN-O2N
2	C	2	NAD	O4D-C4D-C5D-O5D
2	D	4	NAD	C4B-C5B-O5B-PA
2	D	4	NAD	C5D-O5D-PN-O3
2	C	2	NAD	C3D-C4D-C5D-O5D
2	D	4	NAD	O4B-C4B-C5B-O5B
2	D	4	NAD	C3B-C4B-C5B-O5B
3	A	455	ZPG	C27-C24-O23-C22
2	B	3	NAD	O4D-C4D-C5D-O5D
2	B	3	NAD	PN-O3-PA-O5B
2	A	1	NAD	PN-O3-PA-O2A
2	B	3	NAD	C5D-O5D-PN-O1N
2	C	2	NAD	C5D-O5D-PN-O3
2	C	2	NAD	C5D-O5D-PN-O1N
2	D	4	NAD	C5D-O5D-PN-O1N
2	A	1	NAD	O4D-C1D-N1N-C6N
2	C	2	NAD	O4D-C1D-N1N-C2N
2	C	2	NAD	O4D-C1D-N1N-C6N
3	A	455	ZPG	C25-C20-O20-C21
2	B	3	NAD	PN-O3-PA-O1A
2	B	3	NAD	C3D-C4D-C5D-O5D
2	C	2	NAD	PA-O3-PN-O1N
2	A	1	NAD	O4B-C1B-N9A-C8A

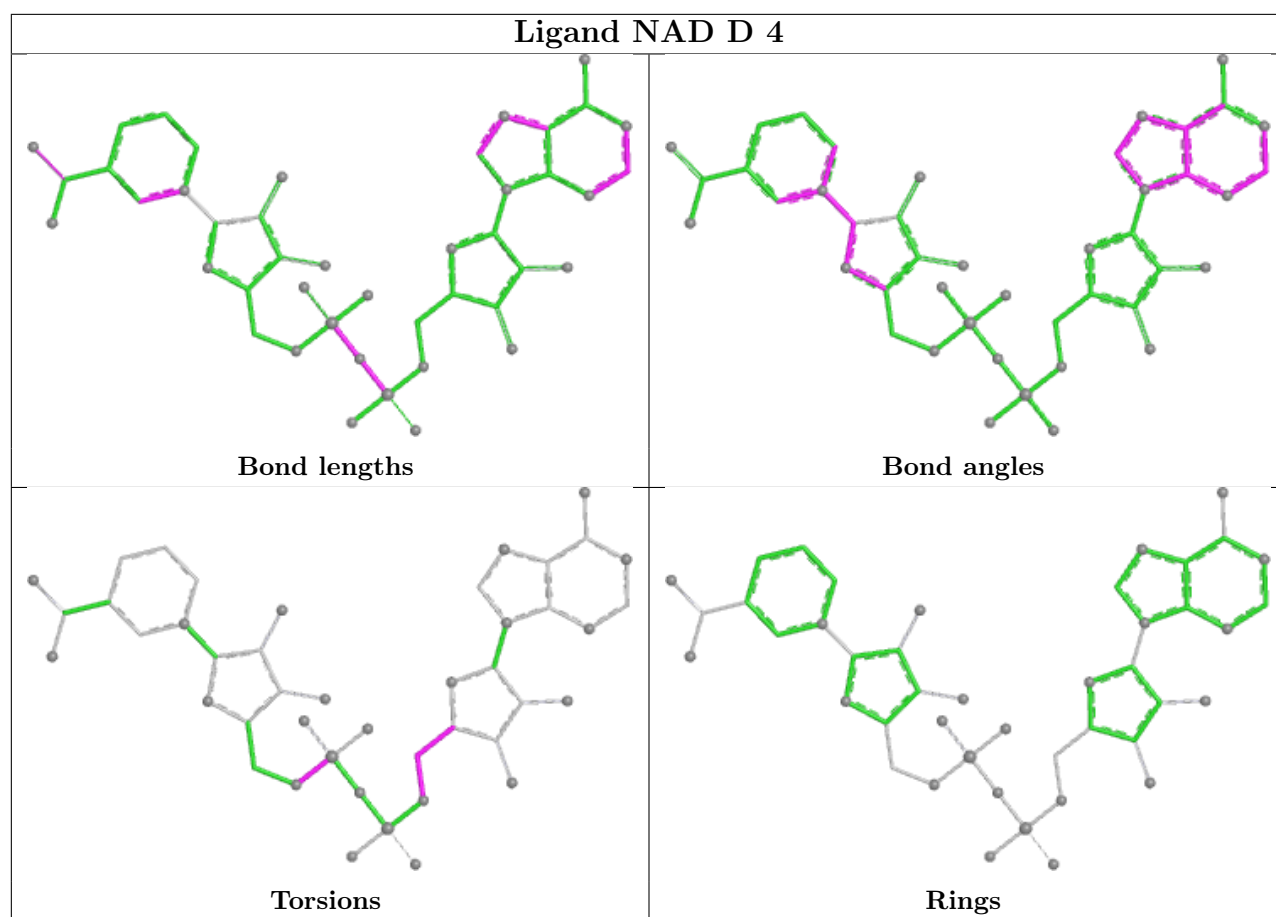
There are no ring outliers.

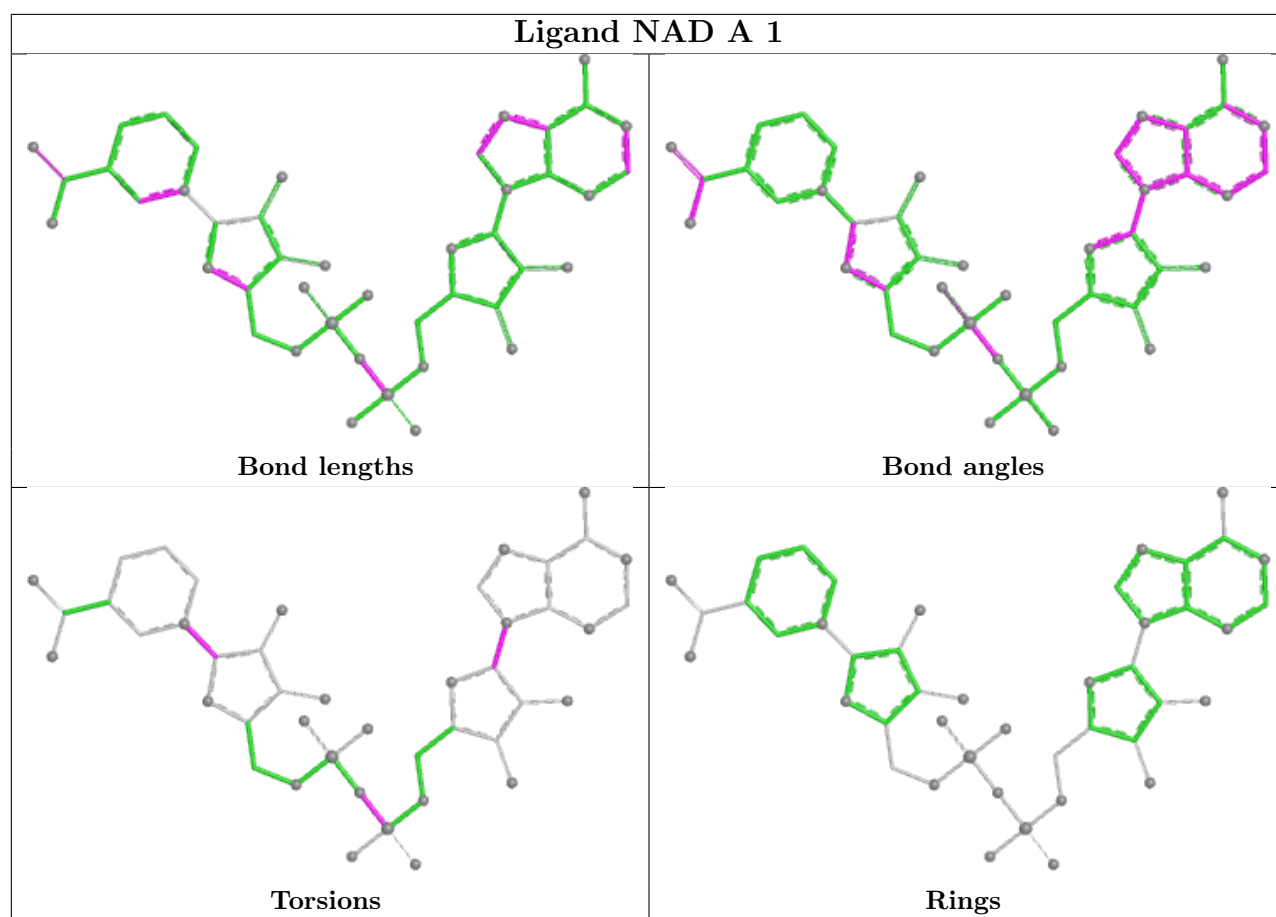
4 monomers are involved in 9 short contacts:

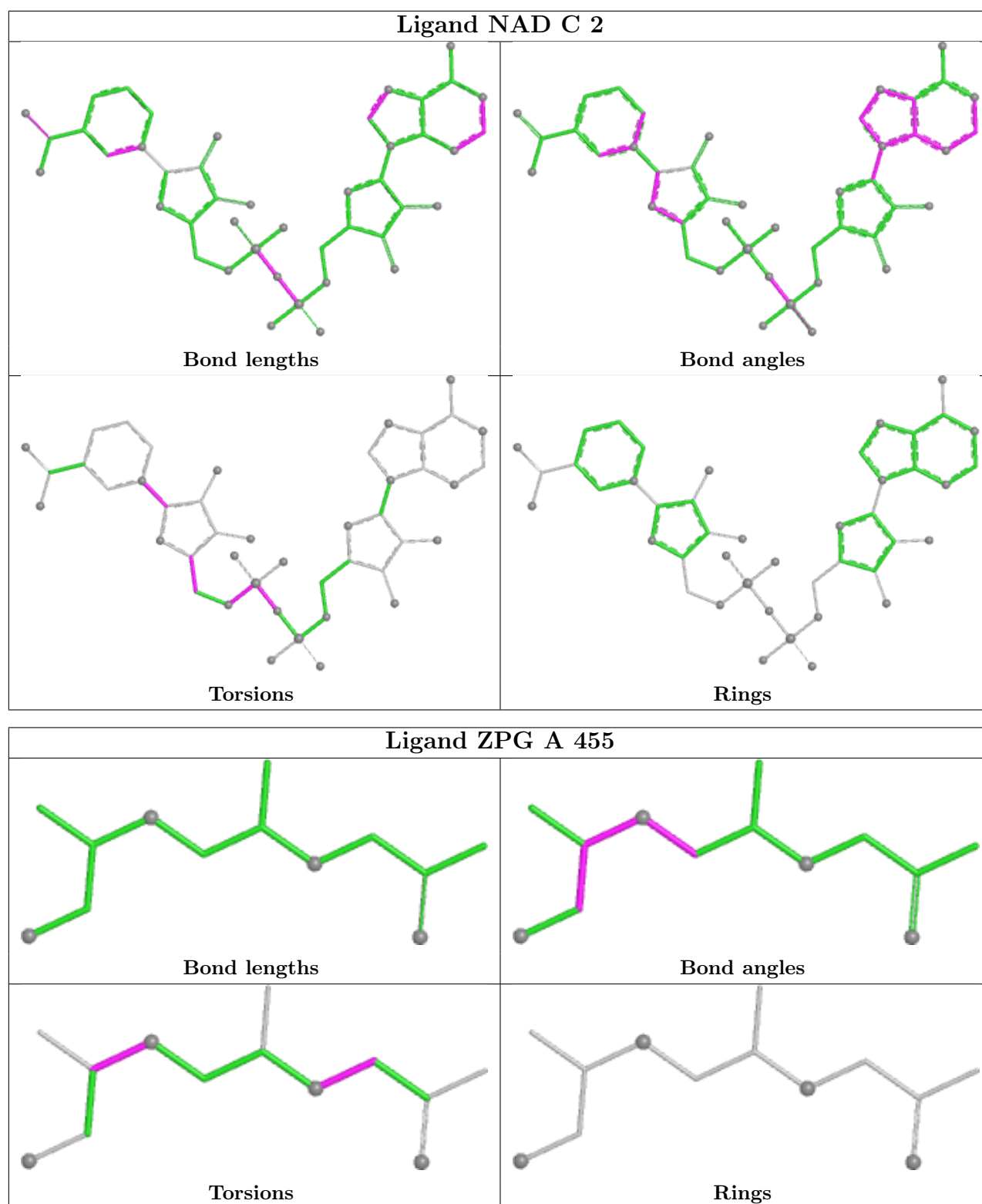
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	3	NAD	3	0
2	D	4	NAD	1	0
2	A	1	NAD	4	0
2	C	2	NAD	1	0

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 ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	427/446 (95%)	-0.44	0 100 100	17, 33, 54, 67	0
1	B	425/446 (95%)	-0.30	0 100 100	14, 36, 71, 83	0
1	C	418/446 (93%)	-0.30	2 (0%) 87 85	19, 40, 66, 78	0
1	D	409/446 (91%)	0.21	10 (2%) 59 54	27, 52, 88, 103	0
All	All	1679/1784 (94%)	-0.21	12 (0%) 84 82	14, 40, 75, 103	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	242	ALA	3.3
1	D	401	ILE	2.9
1	D	241	VAL	2.8
1	D	265	TRP	2.5
1	D	263	ALA	2.4
1	D	252	LEU	2.3
1	D	253	ALA	2.3
1	C	74	ASN	2.2
1	D	344	GLY	2.1
1	D	264	LEU	2.1
1	C	78	ARG	2.0
1	D	262	THR	2.0

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

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

6.3 Carbohydrates ⓘ

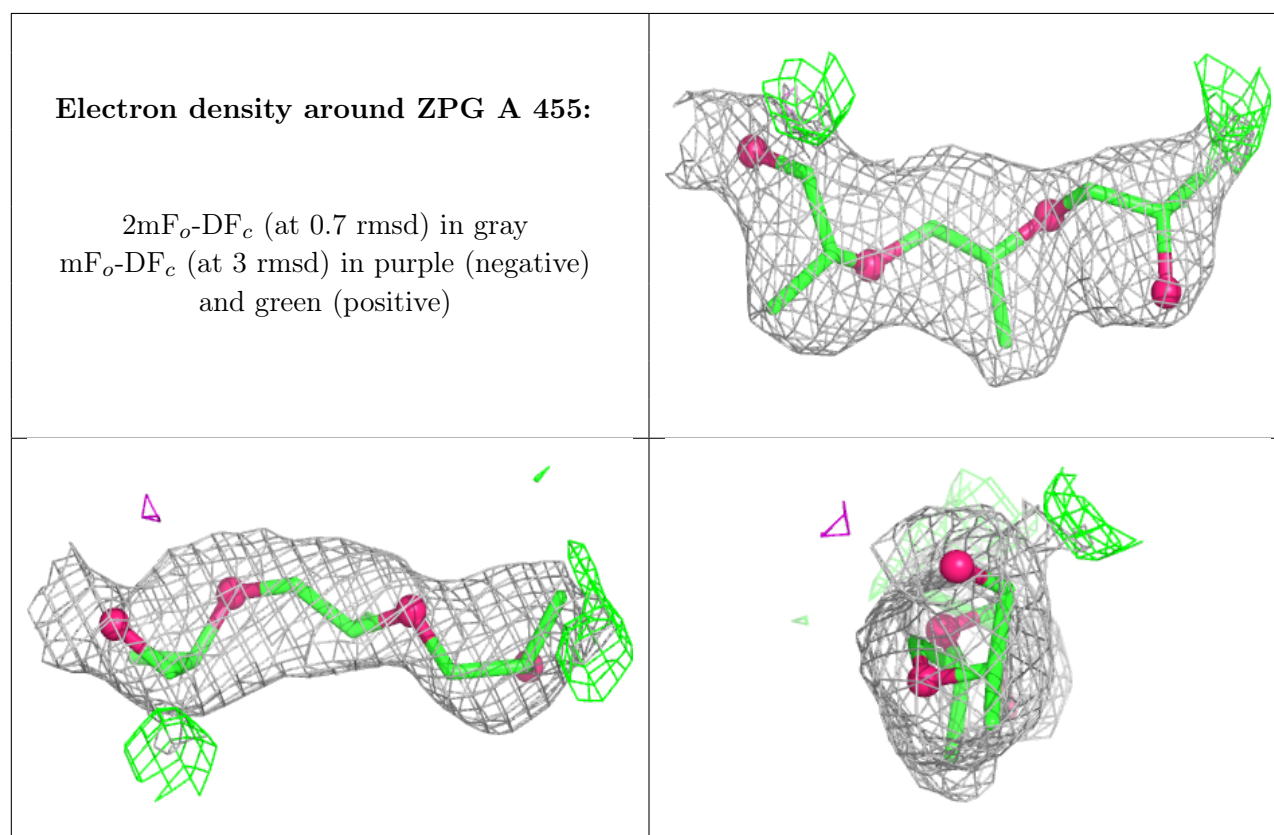
There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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

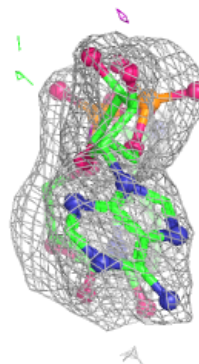
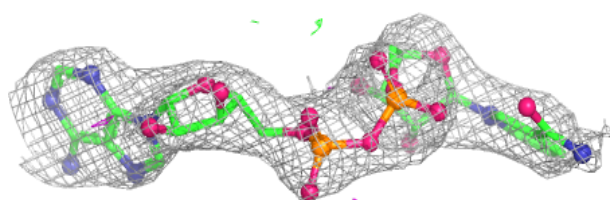
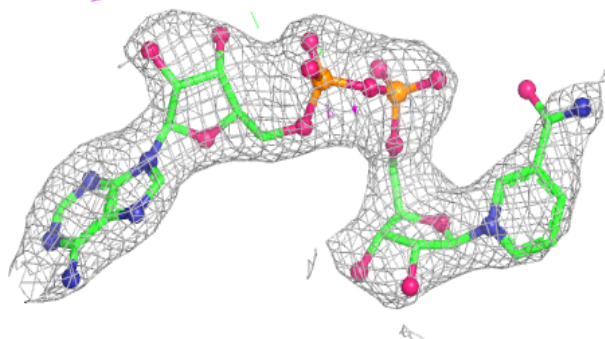
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	ZPG	A	455	13/29	0.80	0.12	60,63,66,67	0
2	NAD	D	4	44/44	0.89	0.09	54,66,71,74	0
2	NAD	B	3	44/44	0.90	0.09	55,62,70,71	0
2	NAD	C	2	44/44	0.94	0.07	39,46,53,55	0
2	NAD	A	1	44/44	0.95	0.07	30,36,46,50	0

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.

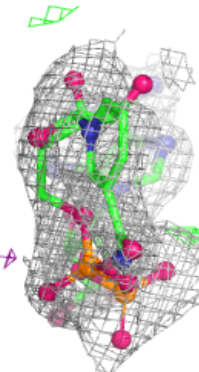
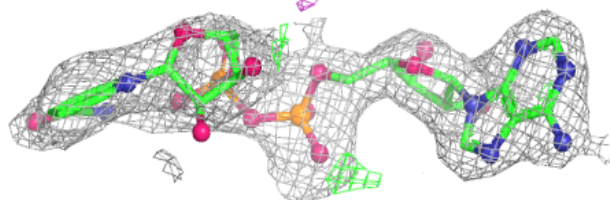
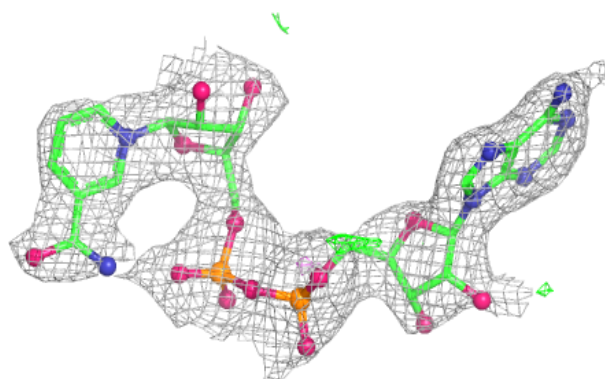


Electron density around NAD D 4:

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

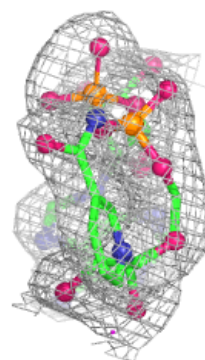
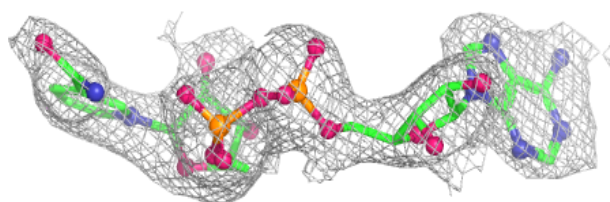
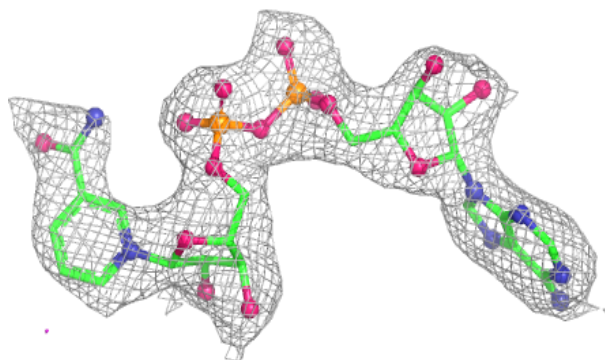
**Electron density around NAD B 3:**

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

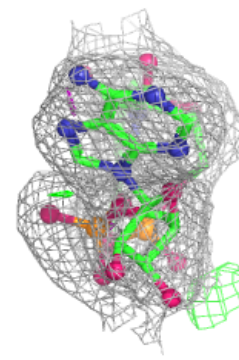
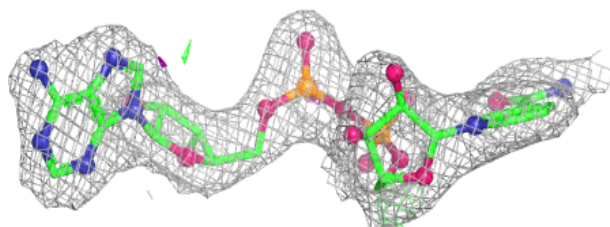
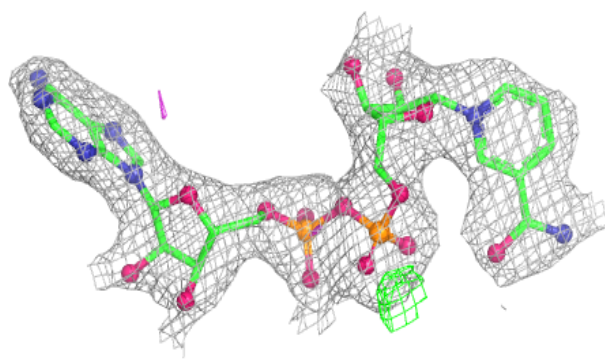


Electron density around NAD C 2:

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

**Electron density around NAD A 1:**

$2mF_o-DF_c$ (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.