



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 07:31 AM UTC

PDB ID : 8PIR / pdb_00008pir
Title : Crystal structure of Ser33 in complex with 3-PGA (3-phosphoglycerate)
Authors : Perrone, S.; Cifuentes, J.O.; Marina, A.; Mastrella, L.; Trastoy, B.; Linster, C.L.; Guerin, M.E.
Deposited on : 2023-06-22
Resolution : 2.70 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

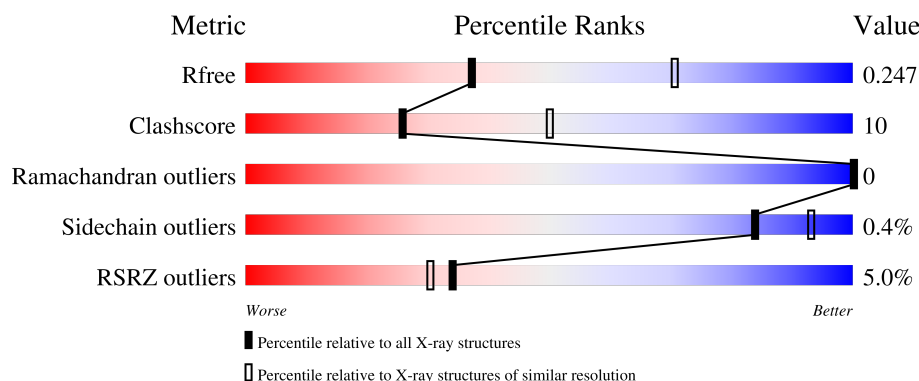
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	473	4% 75% 18% 7%
1	B	473	6% 77% 15% 8%
1	C	473	7% 78% 14% 8%
1	F	473	2% 75% 18% 7%

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	BU1	C	503	-	-	X	-
4	BU1	F	503	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 13319 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.

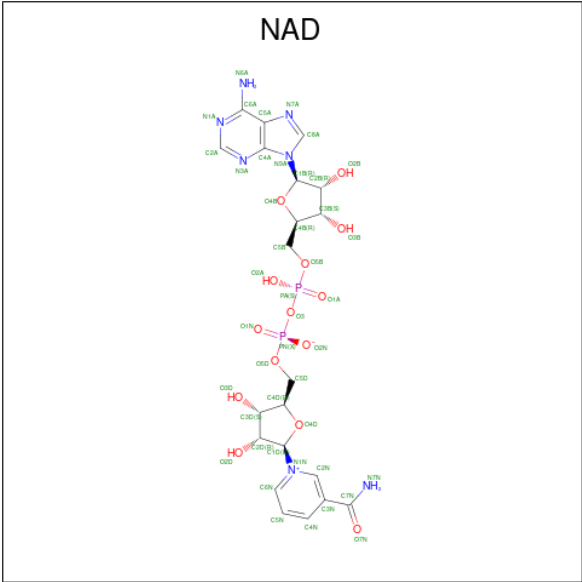
- Molecule 1 is a protein called phosphoglycerate dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	F	439	Total	C	N	O	S	0	0	0
			3318	2101	575	632	10			
1	A	440	Total	C	N	O	S	0	0	0
			3268	2072	561	625	10			
1	B	437	Total	C	N	O	S	0	0	0
			3240	2053	553	624	10			
1	C	436	Total	C	N	O	S	0	0	0
			3214	2036	553	615	10			

There are 16 discrepancies between the modelled and reference sequences:

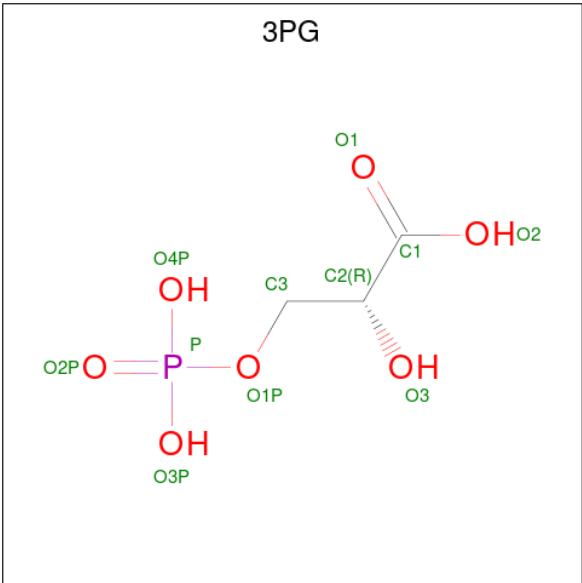
Chain	Residue	Modelled	Actual	Comment	Reference
F	-3	GLY	-	expression tag	UNP A0A8H4BZ61
F	-2	SER	-	expression tag	UNP A0A8H4BZ61
F	-1	GLY	-	expression tag	UNP A0A8H4BZ61
F	0	ALA	-	expression tag	UNP A0A8H4BZ61
A	-3	GLY	-	expression tag	UNP A0A8H4BZ61
A	-2	SER	-	expression tag	UNP A0A8H4BZ61
A	-1	GLY	-	expression tag	UNP A0A8H4BZ61
A	0	ALA	-	expression tag	UNP A0A8H4BZ61
B	-3	GLY	-	expression tag	UNP A0A8H4BZ61
B	-2	SER	-	expression tag	UNP A0A8H4BZ61
B	-1	GLY	-	expression tag	UNP A0A8H4BZ61
B	0	ALA	-	expression tag	UNP A0A8H4BZ61
C	-3	GLY	-	expression tag	UNP A0A8H4BZ61
C	-2	SER	-	expression tag	UNP A0A8H4BZ61
C	-1	GLY	-	expression tag	UNP A0A8H4BZ61
C	0	ALA	-	expression tag	UNP A0A8H4BZ61

- Molecule 2 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (CCD ID: NAD) (formula: $C_{21}H_{27}N_7O_{14}P_2$) (labelled as "Ligand of Interest" by depositor).



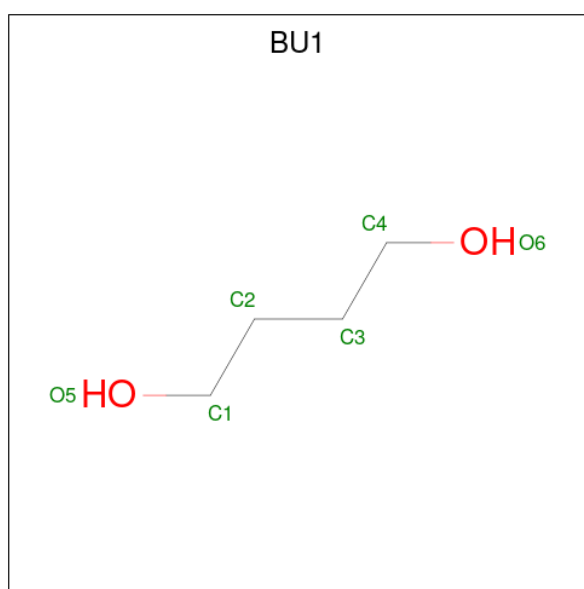
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	F	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	A	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	B	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	C	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

- Molecule 3 is 3-PHOSPHOGLYCERIC ACID (CCD ID: 3PG) (formula: C₃H₇O₇P) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	F	1	Total	C	O	P	0	0
			11	3	7	1		
3	A	1	Total	C	O	P	0	0
			11	3	7	1		
3	B	1	Total	C	O	P	0	0
			11	3	7	1		
3	C	1	Total	C	O	P	0	0
			11	3	7	1		

- Molecule 4 is 1,4-BUTANEDIOL (CCD ID: BU1) (formula: C₄H₁₀O₂) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	F	1	Total	C	O	0	0
			6	4	2		
4	A	1	Total	C	O	0	0
			6	4	2		
4	B	1	Total	C	O	0	0
			6	4	2		
4	C	1	Total	C	O	0	0
			6	4	2		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	F	22	Total	O	0	0
			22	22		

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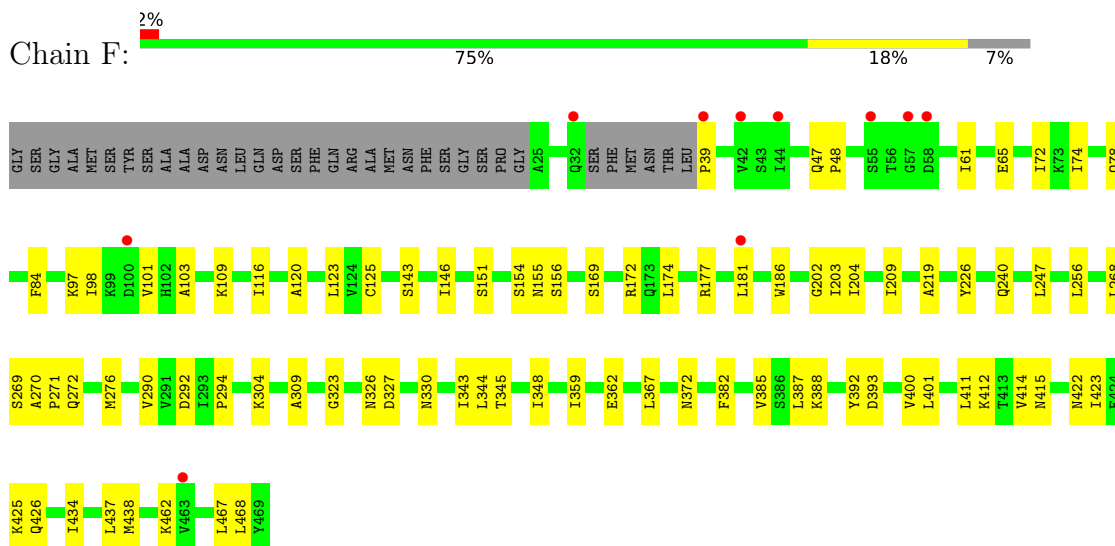
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	4	Total 4	O 4	0	0
5	B	7	Total 7	O 7	0	0
5	C	2	Total 2	O 2	0	0

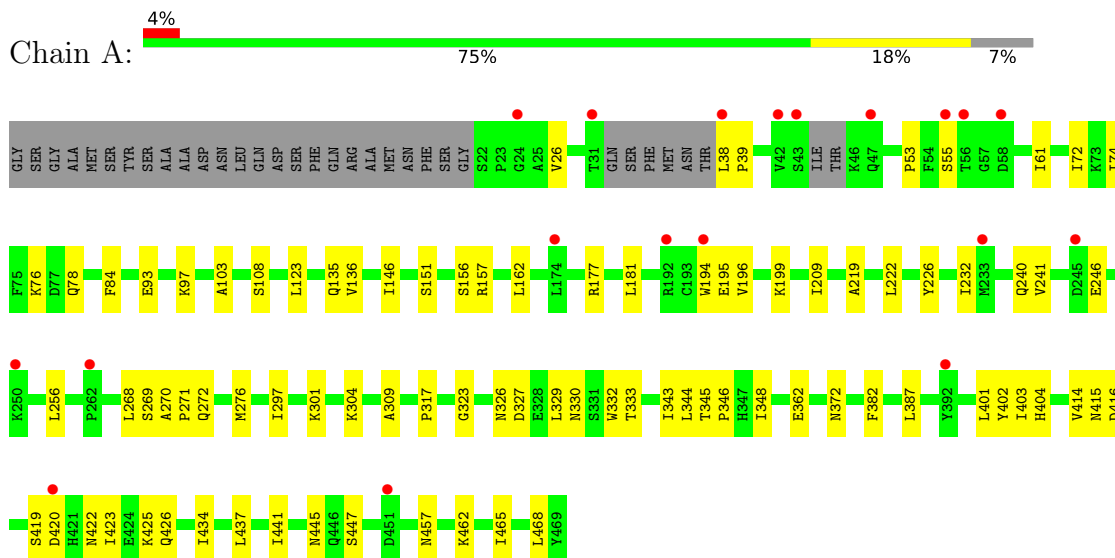
3 Residue-property plots [i](#)

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.

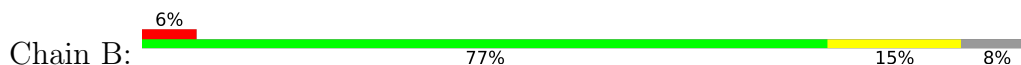
- Molecule 1: phosphoglycerate dehydrogenase

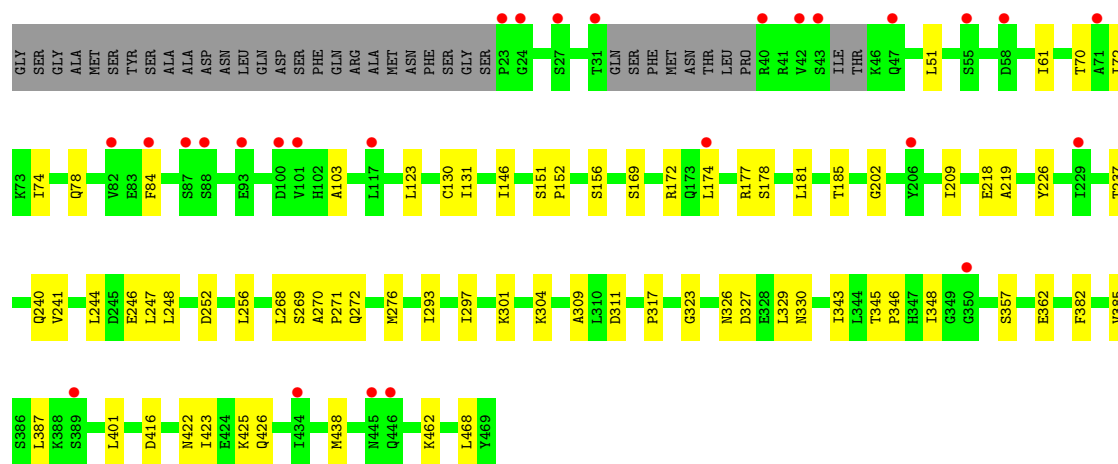


- Molecule 1: phosphoglycerate dehydrogenase

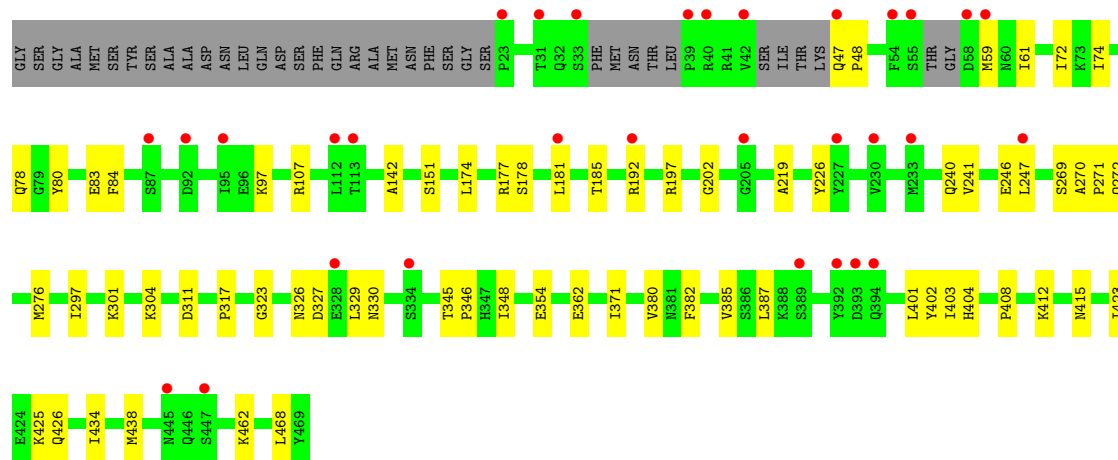
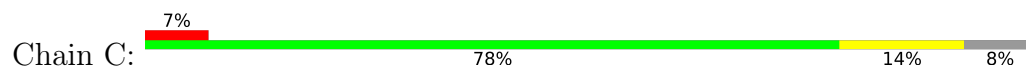


- Molecule 1: phosphoglycerate dehydrogenase





- Molecule 1: phosphoglycerate dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	60.60Å 95.73Å 109.25Å 69.92° 89.53° 86.06°	Depositor
Resolution (Å)	60.45 – 2.70 60.45 – 2.70	Depositor EDS
% Data completeness (in resolution range)	97.7 (60.45-2.70) 98.0 (60.45-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.44 (at 2.42Å)	Xtriage
Refinement program	PHENIX (1.19.2_4158: ???)	Depositor
R, R_{free}	0.215 , 0.245 0.216 , 0.247	Depositor DCC
R_{free} test set	4002 reflections (4.59%)	wwPDB-VP
Wilson B-factor (Å ²)	57.4	Xtriage
Anisotropy	0.258	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 39.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	13319	wwPDB-VP
Average B, all atoms (Å ²)	73.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.88% 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: 3PG, NAD, BU1

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	1.01	0/3323	1.33	1/4524 (0.0%)
1	B	1.02	0/3295	1.36	0/4485
1	C	1.02	0/3265	1.37	0/4445
1	F	1.01	0/3373	1.35	0/4579
All	All	1.02	0/13256	1.35	1/18033 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	420	ASP	N-CA-C	-5.49	106.55	113.20

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	3268	0	3193	76	0
1	B	3240	0	3147	65	0
1	C	3214	0	3118	70	0
1	F	3318	0	3307	68	0
2	A	44	0	26	0	0
2	B	44	0	26	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	44	0	26	0	0
2	F	44	0	26	0	0
3	A	11	0	4	1	0
3	B	11	0	4	1	0
3	C	11	0	4	1	0
3	F	11	0	4	0	0
4	A	6	0	10	3	0
4	B	6	0	10	3	0
4	C	6	0	10	8	0
4	F	6	0	10	5	0
5	A	4	0	0	2	0
5	B	7	0	0	2	0
5	C	2	0	0	0	0
5	F	22	0	0	4	0
All	All	13319	0	12925	249	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 10.

All (249) 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:256:LEU:HD11	1:B:268:LEU:HD12	1.36	1.07
1:A:256:LEU:HD11	1:A:268:LEU:HD12	1.33	1.05
1:C:61:ILE:CD1	1:C:80:TYR:HB3	1.97	0.95
1:A:38:LEU:HD12	1:A:39:PRO:HD2	1.45	0.95
1:B:256:LEU:HD11	1:B:268:LEU:CD1	2.02	0.90
1:A:256:LEU:HD11	1:A:268:LEU:CD1	2.07	0.84
1:B:202:GLY:CA	1:B:247:LEU:HD11	2.08	0.84
1:F:202:GLY:CA	1:F:247:LEU:HD11	2.07	0.84
1:F:172:ARG:HB2	1:F:174:LEU:CD1	2.10	0.81
1:B:172:ARG:HB2	1:B:174:LEU:CD1	2.13	0.79
1:B:244:LEU:HD21	1:B:248:LEU:HD11	1.64	0.78
1:C:202:GLY:CA	1:C:247:LEU:HD11	2.14	0.77
1:F:412:LYS:HE2	1:A:416:ASP:OD1	1.85	0.77
1:F:219:ALA:HB1	1:B:219:ALA:HB1	1.65	0.77
1:B:244:LEU:CD2	1:B:248:LEU:HD11	2.16	0.76
1:B:426:GLN:HE21	1:C:426:GLN:HE21	1.31	0.75
1:F:202:GLY:HA2	1:F:247:LEU:HD11	1.68	0.74
1:A:93:GLU:O	1:A:97:LYS:HG2	1.89	0.73
1:B:61:ILE:CD1	1:B:103:ALA:HB3	2.19	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:270:ALA:HB3	1:C:271:PRO:HD3	1.71	0.73
1:A:61:ILE:CD1	1:A:103:ALA:HB3	2.18	0.72
1:F:422:ASN:OD1	4:A:503:BU1:H31	1.89	0.72
1:A:415:ASN:HD21	1:A:426:GLN:HE22	1.37	0.72
1:B:202:GLY:HA2	1:B:247:LEU:HD11	1.70	0.72
1:B:130:CYS:HB2	3:B:502:3PG:H2	1.70	0.72
1:C:61:ILE:HD11	1:C:80:TYR:HB3	1.71	0.72
4:F:503:BU1:H41	5:A:603:HOH:O	1.90	0.72
1:C:83:GLU:OE2	1:C:97:LYS:HE3	1.90	0.71
1:F:426:GLN:HE21	1:A:426:GLN:HE21	1.37	0.71
1:F:204:ILE:HB	1:F:256:LEU:HD23	1.73	0.71
1:C:61:ILE:CD1	1:C:80:TYR:CB	2.69	0.70
1:F:270:ALA:HB3	1:F:271:PRO:HD3	1.74	0.69
1:B:151:SER:CB	1:B:362:GLU:OE1	2.41	0.69
1:B:270:ALA:HB3	1:B:271:PRO:HD3	1.74	0.69
1:A:151:SER:CB	1:A:362:GLU:OE1	2.41	0.68
1:C:151:SER:CB	1:C:362:GLU:OE1	2.41	0.68
1:C:415:ASN:HD21	1:C:426:GLN:HE22	1.39	0.68
1:A:270:ALA:HB3	1:A:271:PRO:HD3	1.74	0.68
1:A:241:VAL:HG13	1:A:246:GLU:CD	2.19	0.68
1:C:47:GLN:N	1:C:48:PRO:HD2	2.08	0.68
1:F:98:ILE:HD11	1:F:116:ILE:O	1.94	0.67
1:C:202:GLY:HA2	1:C:247:LEU:HD11	1.75	0.67
1:B:241:VAL:HG13	1:B:246:GLU:CD	2.19	0.67
1:C:61:ILE:HD11	1:C:80:TYR:CD2	2.30	0.67
1:A:177:ARG:HG3	1:C:348:ILE:HD13	1.77	0.67
1:F:348:ILE:HD13	1:B:177:ARG:HG3	1.76	0.66
1:F:202:GLY:HA3	1:F:247:LEU:HD11	1.78	0.65
4:F:503:BU1:C4	5:A:603:HOH:O	2.43	0.65
1:F:256:LEU:HD11	1:F:268:LEU:HD12	1.77	0.65
1:F:401:LEU:HD11	1:F:468:LEU:HD11	1.79	0.65
1:C:142:ALA:O	1:C:434:ILE:HD12	1.96	0.64
1:A:219:ALA:HB1	1:C:219:ALA:HB1	1.79	0.64
1:A:61:ILE:HD12	1:A:103:ALA:HB3	1.79	0.64
1:B:241:VAL:HG11	1:B:246:GLU:HB3	1.80	0.64
1:B:202:GLY:HA3	1:B:247:LEU:HD11	1.77	0.63
1:B:70:THR:HG21	1:B:357:SER:OG	1.98	0.63
1:A:346:PRO:N	1:C:181:LEU:HD11	2.14	0.63
1:C:297:ILE:O	1:C:301:LYS:HG3	1.98	0.63
1:F:151:SER:CB	1:F:362:GLU:OE1	2.47	0.62
1:A:108:SER:HB3	3:A:502:3PG:H32	1.82	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:F:503:BU1:H31	1:A:422:ASN:OD1	1.99	0.62
1:A:241:VAL:HG13	1:A:246:GLU:OE1	1.99	0.62
1:B:61:ILE:HD12	1:B:103:ALA:HB3	1.82	0.62
1:C:61:ILE:HD11	1:C:80:TYR:CB	2.30	0.62
1:B:241:VAL:CG1	1:B:246:GLU:HB3	2.30	0.62
1:A:401:LEU:HD11	1:A:468:LEU:HD11	1.82	0.61
1:A:297:ILE:O	1:A:301:LYS:HG3	2.00	0.61
1:A:345:THR:HB	1:A:348:ILE:HG12	1.83	0.61
1:F:411:LEU:HB2	5:F:619:HOH:O	2.01	0.61
1:F:155:ASN:HB2	5:F:617:HOH:O	2.01	0.61
1:A:177:ARG:O	1:A:181:LEU:HD13	2.01	0.61
1:A:53:PRO:C	1:A:55:SER:H	2.08	0.60
1:A:327:ASP:HA	1:A:330:ASN:O	2.01	0.60
1:B:131:ILE:HG23	1:B:152:PRO:HA	1.82	0.60
1:F:415:ASN:OD1	1:A:415:ASN:ND2	2.34	0.60
1:B:422:ASN:OD1	4:C:503:BU1:H31	2.02	0.60
1:A:241:VAL:CG1	1:A:246:GLU:HB3	2.31	0.60
1:A:162:LEU:HD11	1:C:174:LEU:HD21	1.84	0.59
1:B:297:ILE:O	1:B:301:LYS:HG3	2.01	0.59
1:B:387:LEU:HG	1:B:438:MET:HE2	1.85	0.59
1:B:401:LEU:HD11	1:B:468:LEU:HD11	1.83	0.59
1:C:202:GLY:HA3	1:C:247:LEU:HD11	1.83	0.58
1:A:196:VAL:O	1:A:222:LEU:HD23	2.04	0.58
1:C:401:LEU:HD11	1:C:468:LEU:HD11	1.86	0.58
1:F:387:LEU:HG	1:F:438:MET:HE2	1.85	0.58
1:A:38:LEU:CD1	1:A:39:PRO:HD2	2.28	0.58
1:C:181:LEU:HD23	1:C:185:THR:O	2.04	0.57
1:A:241:VAL:HG11	1:A:246:GLU:HB3	1.87	0.57
1:A:346:PRO:CD	1:C:181:LEU:HD11	2.35	0.57
1:F:256:LEU:HD13	1:F:290:VAL:HG12	1.85	0.57
1:A:346:PRO:HD3	1:C:181:LEU:HD11	1.87	0.57
1:C:107:ARG:NE	3:C:502:3PG:O3P	2.31	0.57
1:C:387:LEU:HG	1:C:438:MET:HE2	1.87	0.57
1:B:241:VAL:HG13	1:B:246:GLU:OE1	2.04	0.56
1:F:256:LEU:HD13	1:F:290:VAL:CG1	2.35	0.56
1:B:244:LEU:HD23	1:B:248:LEU:CD1	2.36	0.56
1:C:311:ASP:HA	1:C:345:THR:OG1	2.06	0.55
1:C:61:ILE:HD12	1:C:80:TYR:HB3	1.85	0.55
1:C:403:ILE:CG2	1:C:434:ILE:HD11	2.37	0.55
1:A:196:VAL:O	1:A:222:LEU:CD2	2.55	0.55
1:A:403:ILE:HG22	1:A:434:ILE:HD11	1.87	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:402:TYR:HD1	1:A:465:ILE:HD13	1.71	0.54
1:C:61:ILE:HD11	1:C:80:TYR:CG	2.43	0.53
1:F:177:ARG:O	1:F:181:LEU:HD13	2.09	0.53
1:A:345:THR:C	1:C:181:LEU:HD11	2.33	0.53
1:B:387:LEU:HB3	1:B:425:LYS:HE2	1.89	0.53
1:A:387:LEU:HB3	1:A:425:LYS:HE2	1.90	0.53
1:C:47:GLN:N	1:C:48:PRO:CD	2.72	0.53
1:A:72:ILE:HD11	1:A:84:PHE:CD1	2.44	0.53
4:B:503:BU1:H32	1:C:423:ILE:H	1.74	0.52
1:C:387:LEU:HB3	1:C:425:LYS:HE3	1.91	0.52
1:A:135:GLN:HG2	1:A:136:VAL:HG23	1.92	0.52
1:C:72:ILE:HD11	1:C:84:PHE:CD1	2.45	0.52
1:C:241:VAL:HG23	1:C:246:GLU:CD	2.35	0.52
1:F:72:ILE:HD11	1:F:84:PHE:CD1	2.46	0.51
1:F:344:LEU:C	1:B:181:LEU:HD12	2.36	0.51
1:C:276:MET:O	1:C:304:LYS:NZ	2.43	0.51
1:C:241:VAL:CG2	1:C:246:GLU:HB3	2.41	0.51
1:B:311:ASP:HA	1:B:345:THR:OG1	2.10	0.51
1:F:387:LEU:HB3	1:F:425:LYS:HE2	1.93	0.50
1:F:78:GLN:OE1	1:F:372:ASN:ND2	2.45	0.50
1:F:172:ARG:HB2	1:F:174:LEU:HD11	1.91	0.50
1:B:423:ILE:H	4:C:503:BU1:H32	1.76	0.50
1:F:385:VAL:HG12	1:F:438:MET:SD	2.51	0.50
1:A:445:ASN:OD1	1:A:447:SER:HB3	2.12	0.50
1:B:423:ILE:H	4:C:503:BU1:C3	2.24	0.50
1:A:457:ASN:HA	1:A:462:LYS:HD3	1.93	0.50
1:A:38:LEU:HG	1:A:39:PRO:O	2.12	0.50
1:C:317:PRO:HD3	1:C:329:LEU:HD21	1.94	0.50
1:F:74:ILE:O	1:F:78:GLN:HG2	2.11	0.50
1:B:276:MET:O	1:B:304:LYS:NZ	2.43	0.49
1:F:47:GLN:HG3	1:F:48:PRO:HD2	1.93	0.49
1:F:256:LEU:HD11	1:F:268:LEU:CD1	2.42	0.49
1:F:177:ARG:HG3	1:B:348:ILE:HD13	1.95	0.49
1:A:276:MET:O	1:A:304:LYS:NZ	2.43	0.49
1:A:53:PRO:C	1:A:55:SER:N	2.70	0.49
1:B:346:PRO:HG2	1:B:348:ILE:HG23	1.95	0.49
1:B:416:ASP:OD1	1:C:412:LYS:NZ	2.43	0.48
5:B:604:HOH:O	4:C:503:BU1:H41	2.13	0.48
1:F:169:SER:HA	1:F:174:LEU:HD13	1.95	0.48
1:F:344:LEU:O	1:B:178:SER:HA	2.13	0.48
1:A:195:GLU:HB2	1:C:354:GLU:HG2	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:72:ILE:HD11	1:B:84:PHE:CD1	2.47	0.48
1:A:332:TRP:O	1:A:333:THR:C	2.54	0.48
1:B:385:VAL:HG12	1:B:438:MET:SD	2.54	0.48
4:F:503:BU1:H32	1:A:423:ILE:N	2.29	0.48
1:A:162:LEU:CD1	1:C:174:LEU:HD21	2.43	0.48
1:A:177:ARG:HG3	1:C:348:ILE:CD1	2.43	0.48
1:B:423:ILE:N	4:C:503:BU1:H32	2.29	0.47
1:C:408:PRO:HA	4:C:503:BU1:H41	1.96	0.47
1:F:98:ILE:HD12	1:F:120:ALA:HB2	1.96	0.47
4:F:503:BU1:H32	1:A:423:ILE:H	1.80	0.47
1:F:61:ILE:CD1	1:F:103:ALA:HB3	2.44	0.47
1:F:123:LEU:HD23	1:F:146:ILE:HD13	1.95	0.47
1:B:169:SER:HA	1:B:174:LEU:HD13	1.97	0.47
1:A:194:TRP:HB3	1:A:199:LYS:HE2	1.97	0.47
1:B:61:ILE:HD13	1:B:103:ALA:HB3	1.96	0.47
1:C:380:VAL:HG12	1:C:380:VAL:O	2.14	0.47
1:F:345:THR:C	1:B:181:LEU:HD11	2.39	0.46
1:B:426:GLN:HE21	1:C:426:GLN:NE2	2.07	0.46
1:F:414:VAL:HG11	1:F:437:LEU:CD2	2.46	0.46
1:B:74:ILE:O	1:B:78:GLN:HG2	2.14	0.46
1:B:269:SER:N	1:B:272:GLN:OE1	2.47	0.46
1:F:276:MET:O	1:F:304:LYS:NZ	2.48	0.46
1:B:317:PRO:HD3	1:B:329:LEU:HD21	1.98	0.46
1:A:423:ILE:HG12	1:A:441:ILE:HG22	1.98	0.46
1:C:74:ILE:O	1:C:78:GLN:HG2	2.15	0.46
1:A:72:ILE:HG23	1:A:76:LYS:HE3	1.98	0.45
1:C:327:ASP:HA	1:C:330:ASN:O	2.16	0.45
1:C:346:PRO:HG2	1:C:348:ILE:CG2	2.47	0.45
4:B:503:BU1:H32	1:C:423:ILE:N	2.31	0.45
5:B:604:HOH:O	4:C:503:BU1:C4	2.64	0.45
1:F:65:GLU:OE1	1:F:109:LYS:HB2	2.17	0.45
1:A:269:SER:N	1:A:272:GLN:OE1	2.46	0.45
1:F:387:LEU:CG	1:F:438:MET:HE2	2.45	0.45
1:F:269:SER:N	1:F:272:GLN:OE1	2.47	0.45
1:B:181:LEU:HD23	1:B:185:THR:O	2.17	0.45
1:A:156:SER:HA	1:A:209:ILE:HG12	1.98	0.45
1:C:385:VAL:HG12	1:C:438:MET:SD	2.57	0.45
1:F:154:SER:OG	1:F:359:ILE:HG13	2.17	0.44
1:B:51:LEU:HB3	1:B:468:LEU:HB3	1.99	0.44
1:B:423:ILE:O	4:C:503:BU1:H11	2.17	0.44
1:F:388:LYS:HG3	5:F:620:HOH:O	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:269:SER:N	1:C:272:GLN:OE1	2.47	0.44
1:A:123:LEU:HD23	1:A:146:ILE:HD13	1.99	0.44
1:F:39:PRO:HB2	1:A:232:ILE:CG2	2.47	0.44
1:B:387:LEU:CG	1:B:438:MET:HE2	2.46	0.44
1:F:97:LYS:O	1:F:101:VAL:HG23	2.17	0.44
1:F:156:SER:HA	1:F:209:ILE:HG12	2.00	0.44
1:A:61:ILE:HD13	1:A:103:ALA:HB3	1.96	0.44
1:F:125:CYS:SG	1:F:367:LEU:HD22	2.58	0.43
1:F:143:SER:HA	1:F:434:ILE:HD11	2.00	0.43
1:A:344:LEU:O	1:C:178:SER:HA	2.18	0.43
1:B:226:TYR:CZ	1:B:240:GLN:HB2	2.53	0.43
1:B:327:ASP:HA	1:B:330:ASN:O	2.18	0.43
1:C:387:LEU:CG	1:C:438:MET:HE2	2.47	0.43
1:F:412:LYS:HD2	1:A:419:SER:OG	2.19	0.43
1:A:346:PRO:HB3	1:C:181:LEU:HD21	2.00	0.43
1:A:74:ILE:O	1:A:78:GLN:HG2	2.18	0.43
1:A:157:ARG:HG2	1:C:197:ARG:CZ	2.49	0.43
1:C:462:LYS:HD2	1:C:462:LYS:C	2.43	0.43
1:F:226:TYR:CZ	1:F:240:GLN:HB2	2.53	0.43
1:F:462:LYS:C	1:F:462:LYS:HD2	2.44	0.43
1:A:78:GLN:OE1	1:A:372:ASN:ND2	2.44	0.43
1:C:241:VAL:HG22	1:C:246:GLU:HB3	2.01	0.43
5:F:614:HOH:O	4:A:503:BU1:H41	2.19	0.42
1:B:156:SER:HA	1:B:209:ILE:HG12	2.00	0.42
1:F:400:VAL:HG22	1:F:467:LEU:CD2	2.49	0.42
1:F:172:ARG:HB2	1:F:174:LEU:HD12	1.97	0.42
1:B:172:ARG:HB2	1:B:174:LEU:HD11	1.96	0.42
1:F:323:GLY:O	1:F:326:ASN:ND2	2.52	0.42
1:A:323:GLY:O	1:A:326:ASN:ND2	2.53	0.42
1:F:204:ILE:HD12	1:F:256:LEU:CD2	2.49	0.42
1:F:204:ILE:HD12	1:F:256:LEU:HD23	2.01	0.42
1:B:387:LEU:CD1	1:B:438:MET:HE2	2.49	0.42
1:C:226:TYR:CZ	1:C:240:GLN:HB2	2.54	0.42
1:F:186:TRP:CG	1:B:317:PRO:HG2	2.54	0.42
1:A:344:LEU:C	1:C:181:LEU:HD12	2.45	0.42
1:B:462:LYS:HD2	1:B:462:LYS:C	2.45	0.42
1:F:423:ILE:O	4:A:503:BU1:H32	2.20	0.42
1:A:345:THR:C	1:C:181:LEU:CD1	2.93	0.42
1:F:327:ASP:HA	1:F:330:ASN:O	2.20	0.41
1:A:226:TYR:CZ	1:A:240:GLN:HB2	2.55	0.41
1:B:241:VAL:CG1	1:B:246:GLU:CD	2.91	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:142:ALA:O	1:C:434:ILE:CD1	2.67	0.41
1:C:323:GLY:O	1:C:326:ASN:ND2	2.53	0.41
1:F:392:TYR:CE1	1:F:393:ASP:HB3	2.55	0.41
1:F:411:LEU:HD12	1:A:423:ILE:HD12	2.02	0.41
1:A:309:ALA:HA	1:A:343:ILE:O	2.20	0.41
1:A:317:PRO:HD3	1:A:329:LEU:HD21	2.02	0.41
1:C:402:TYR:CE2	1:C:404:HIS:HB3	2.56	0.41
1:F:426:GLN:NE2	1:A:426:GLN:HE21	2.12	0.41
1:C:346:PRO:HG2	1:C:348:ILE:HG23	2.02	0.41
1:A:414:VAL:HG11	1:A:437:LEU:CD2	2.50	0.41
1:F:203:ILE:O	1:F:226:TYR:HA	2.20	0.41
1:A:26:VAL:HG22	1:B:252:ASP:OD1	2.21	0.41
1:A:402:TYR:CE2	1:A:404:HIS:HB3	2.56	0.41
1:B:323:GLY:O	1:B:326:ASN:ND2	2.54	0.41
1:C:83:GLU:OE2	1:C:97:LYS:CE	2.65	0.41
1:B:218:GLU:OE2	1:B:237:THR:OG1	2.29	0.40
1:B:426:GLN:NE2	1:C:426:GLN:HE21	2.10	0.40
1:F:292:ASP:OD1	1:F:294:PRO:HD2	2.22	0.40
1:F:309:ALA:HA	1:F:343:ILE:O	2.21	0.40
1:A:348:ILE:HD13	1:C:177:ARG:HG3	2.03	0.40
1:B:123:LEU:HD23	1:B:146:ILE:HD13	2.03	0.40
4:B:503:BU1:H11	1:C:423:ILE:O	2.20	0.40
1:B:309:ALA:HA	1:B:343:ILE:O	2.21	0.40
1:F:61:ILE:HD12	1:F:103:ALA:HB3	2.02	0.40
1:C:59:MET:HE2	1:C:371:ILE:HG23	2.02	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	434/473 (92%)	412 (95%)	22 (5%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	431/473 (91%)	413 (96%)	18 (4%)	0	100	100
1	C	428/473 (90%)	411 (96%)	17 (4%)	0	100	100
1	F	435/473 (92%)	416 (96%)	19 (4%)	0	100	100
All	All	1728/1892 (91%)	1652 (96%)	76 (4%)	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	339/402 (84%)	338 (100%)	1 (0%)	86	94
1	B	335/402 (83%)	333 (99%)	2 (1%)	78	91
1	C	328/402 (82%)	326 (99%)	2 (1%)	78	91
1	F	353/402 (88%)	352 (100%)	1 (0%)	86	94
All	All	1355/1608 (84%)	1349 (100%)	6 (0%)	84	93

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	F	382	PHE
1	A	382	PHE
1	B	293	ILE
1	B	382	PHE
1	C	192	ARG
1	C	382	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	F	330	ASN
1	F	415	ASN

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Mol	Chain	Res	Type
1	F	426	GLN
1	A	330	ASN
1	A	356	GLN
1	A	415	ASN
1	B	330	ASN
1	B	394	GLN
1	B	406	ASN
1	B	415	ASN
1	B	421	HIS
1	B	426	GLN
1	C	330	ASN
1	C	356	GLN
1	C	415	ASN
1	C	421	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](#)

12 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	3PG	C	502	-	9,10,10	0.16	0	11,14,14	0.47	0
3	3PG	F	502	-	9,10,10	1.52	2 (22%)	11,14,14	2.17	2 (18%)
3	3PG	A	502	-	9,10,10	1.53	2 (22%)	11,14,14	2.14	2 (18%)
2	NAD	C	501	-	46,48,48	0.40	0	64,73,73	0.43	0
4	BU1	B	503	-	5,5,5	0.15	0	4,4,4	0.24	0
2	NAD	B	501	-	46,48,48	0.36	0	64,73,73	0.44	0
2	NAD	A	501	-	46,48,48	0.41	0	64,73,73	0.43	0
4	BU1	C	503	-	5,5,5	0.09	0	4,4,4	0.17	0
3	3PG	B	502	-	9,10,10	0.16	0	11,14,14	0.41	0
4	BU1	A	503	-	5,5,5	0.15	0	4,4,4	0.14	0
2	NAD	F	501	-	46,48,48	0.40	0	64,73,73	0.54	0
4	BU1	F	503	-	5,5,5	0.08	0	4,4,4	0.15	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
3	3PG	C	502	-	-	2/10/10/10	-
3	3PG	F	502	-	-	4/10/10/10	-
3	3PG	A	502	-	-	4/10/10/10	-
2	NAD	C	501	-	-	8/30/62/62	0/5/5/5
4	BU1	B	503	-	-	2/3/3/3	-
2	NAD	B	501	-	-	4/30/62/62	0/5/5/5
2	NAD	A	501	-	-	9/30/62/62	0/5/5/5
4	BU1	C	503	-	-	1/3/3/3	-
3	3PG	B	502	-	-	4/10/10/10	-
4	BU1	A	503	-	-	2/3/3/3	-
2	NAD	F	501	-	-	6/30/62/62	0/5/5/5
4	BU1	F	503	-	-	1/3/3/3	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	502	3PG	O1-C1	3.32	1.31	1.22
3	F	502	3PG	O1-C1	3.28	1.31	1.22
3	F	502	3PG	O2-C1	-2.61	1.22	1.30
3	A	502	3PG	O2-C1	-2.60	1.22	1.30

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	502	3PG	O2-C1-C2	4.86	123.00	112.74
3	F	502	3PG	O1-C1-C2	-4.85	112.92	122.60
3	A	502	3PG	O2-C1-C2	4.72	122.72	112.74
3	A	502	3PG	O1-C1-C2	-4.70	113.21	122.60

There are no chirality outliers.

All (47) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	F	501	NAD	O4D-C1D-N1N-C2N
2	F	501	NAD	O4D-C1D-N1N-C6N
2	F	501	NAD	C2D-C1D-N1N-C2N
2	F	501	NAD	C2D-C1D-N1N-C6N
2	A	501	NAD	O4D-C1D-N1N-C2N
2	A	501	NAD	O4D-C1D-N1N-C6N
2	A	501	NAD	C2D-C1D-N1N-C2N
2	A	501	NAD	C2D-C1D-N1N-C6N
2	B	501	NAD	O4D-C1D-N1N-C2N
2	B	501	NAD	O4D-C1D-N1N-C6N
2	B	501	NAD	C2D-C1D-N1N-C2N
2	B	501	NAD	C2D-C1D-N1N-C6N
2	C	501	NAD	O4D-C1D-N1N-C2N
2	C	501	NAD	O4D-C1D-N1N-C6N
2	C	501	NAD	C2D-C1D-N1N-C2N
2	C	501	NAD	C2D-C1D-N1N-C6N
3	F	502	3PG	C1-C2-C3-O1P
3	F	502	3PG	C3-O1P-P-O3P
3	A	502	3PG	O1-C1-C2-O3
3	A	502	3PG	O2-C1-C2-O3
3	B	502	3PG	C1-C2-C3-O1P
3	B	502	3PG	O3-C2-C3-O1P
3	B	502	3PG	C3-O1P-P-O3P
3	B	502	3PG	C3-O1P-P-O4P
4	A	503	BU1	C2-C3-C4-O6
4	B	503	BU1	C2-C3-C4-O6
4	F	503	BU1	C2-C3-C4-O6
4	C	503	BU1	C2-C3-C4-O6
2	A	501	NAD	O4D-C4D-C5D-O5D
3	C	502	3PG	O1-C1-C2-O3
3	C	502	3PG	O2-C1-C2-O3
4	A	503	BU1	O5-C1-C2-C3

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Mol	Chain	Res	Type	Atoms
4	B	503	BU1	O5-C1-C2-C3
2	A	501	NAD	PN-O3-PA-O1A
3	F	502	3PG	O3-C2-C3-O1P
2	A	501	NAD	C3D-C4D-C5D-O5D
2	C	501	NAD	PN-O3-PA-O2A
2	C	501	NAD	C5B-O5B-PA-O1A
2	F	501	NAD	PN-O3-PA-O1A
2	F	501	NAD	PN-O3-PA-O2A
3	A	502	3PG	O2-C1-C2-C3
3	F	502	3PG	C3-O1P-P-O4P
2	C	501	NAD	O4B-C4B-C5B-O5B
2	A	501	NAD	PN-O3-PA-O2A
2	C	501	NAD	PN-O3-PA-O1A
2	A	501	NAD	C4N-C3N-C7N-N7N
3	A	502	3PG	O1-C1-C2-C3

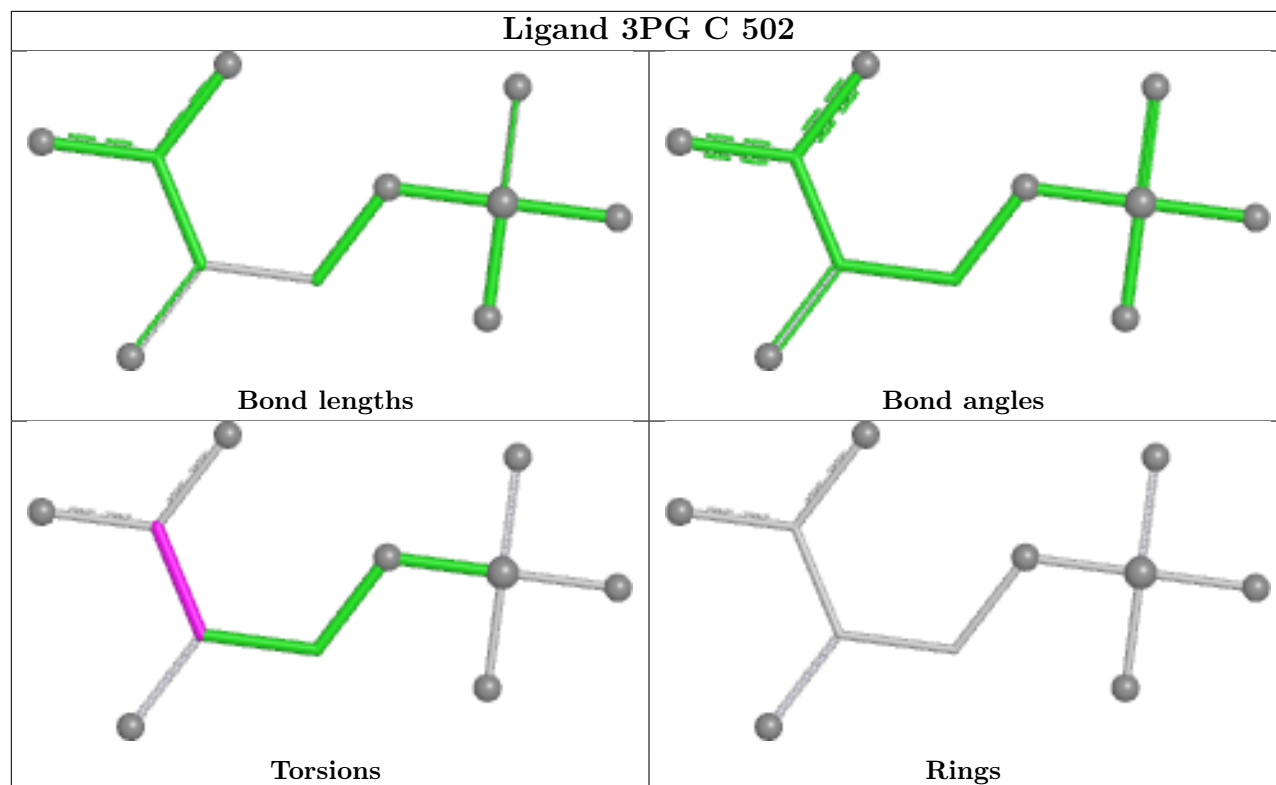
There are no ring outliers.

7 monomers are involved in 22 short contacts:

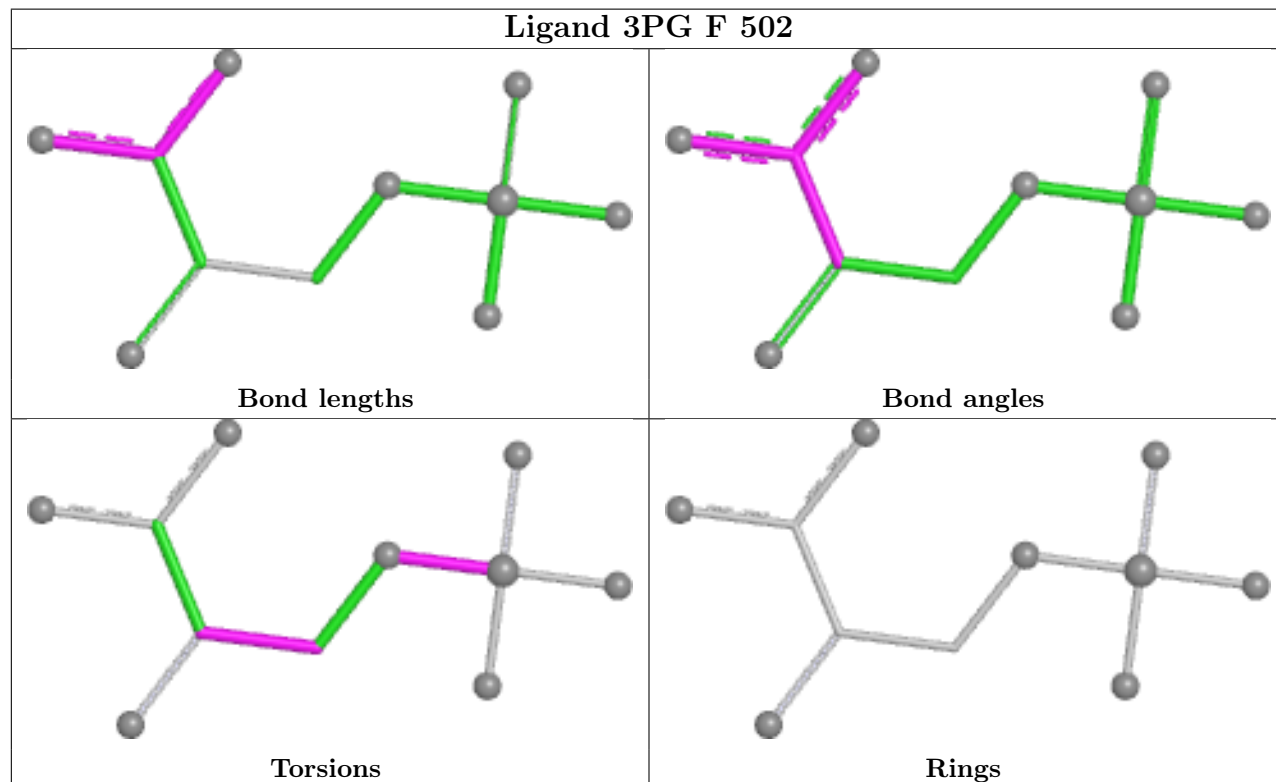
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	502	3PG	1	0
3	A	502	3PG	1	0
4	B	503	BU1	3	0
4	C	503	BU1	8	0
3	B	502	3PG	1	0
4	A	503	BU1	3	0
4	F	503	BU1	5	0

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

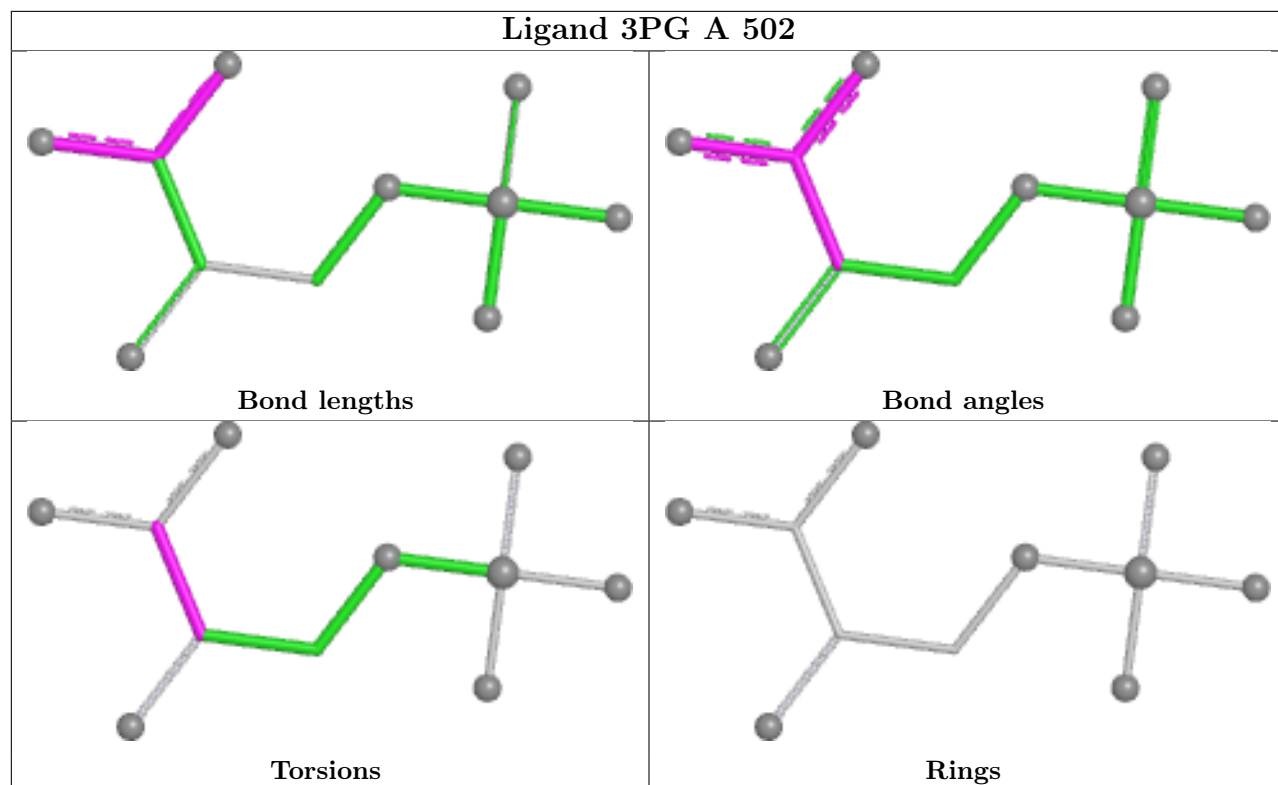
Ligand 3PG C 502



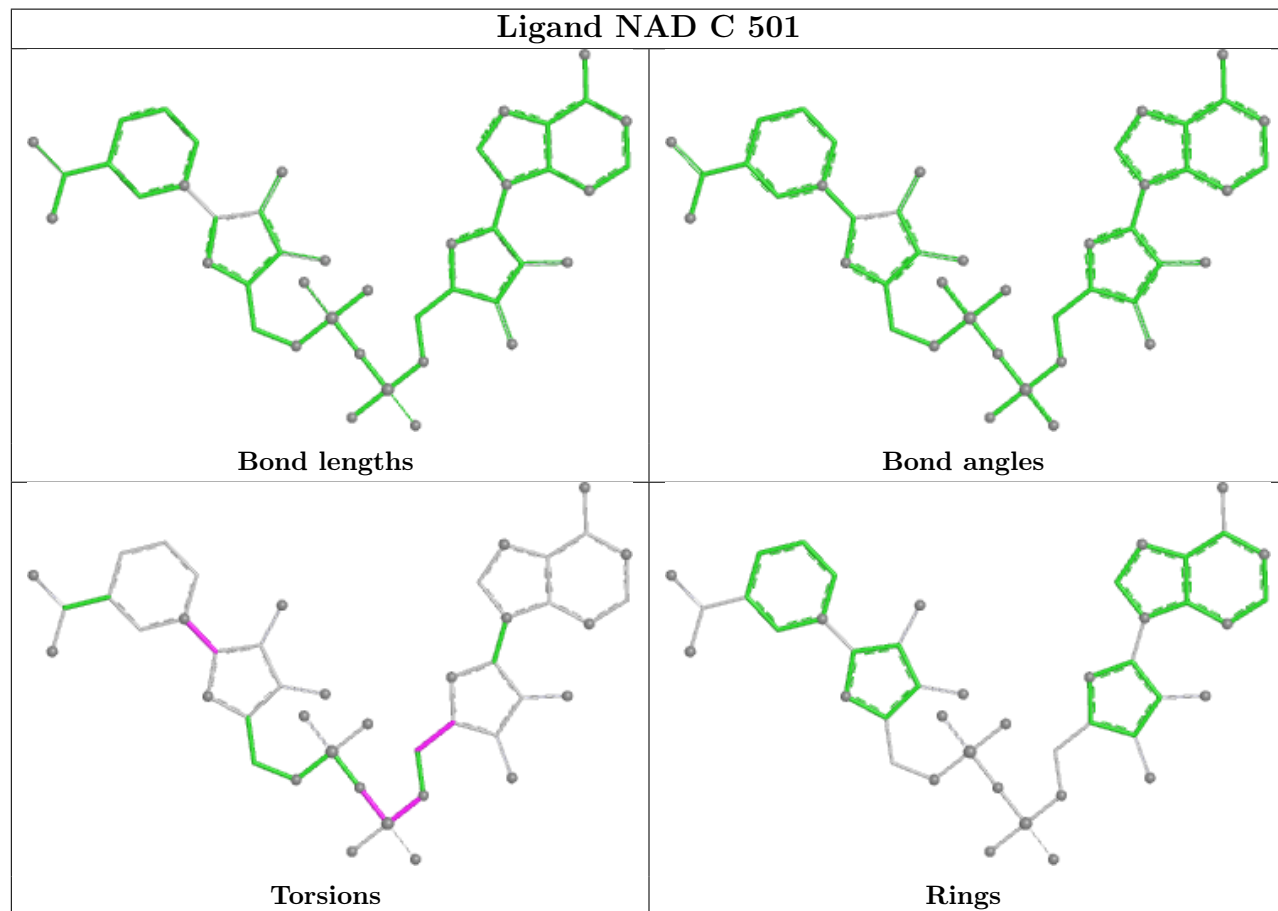
Ligand 3PG F 502

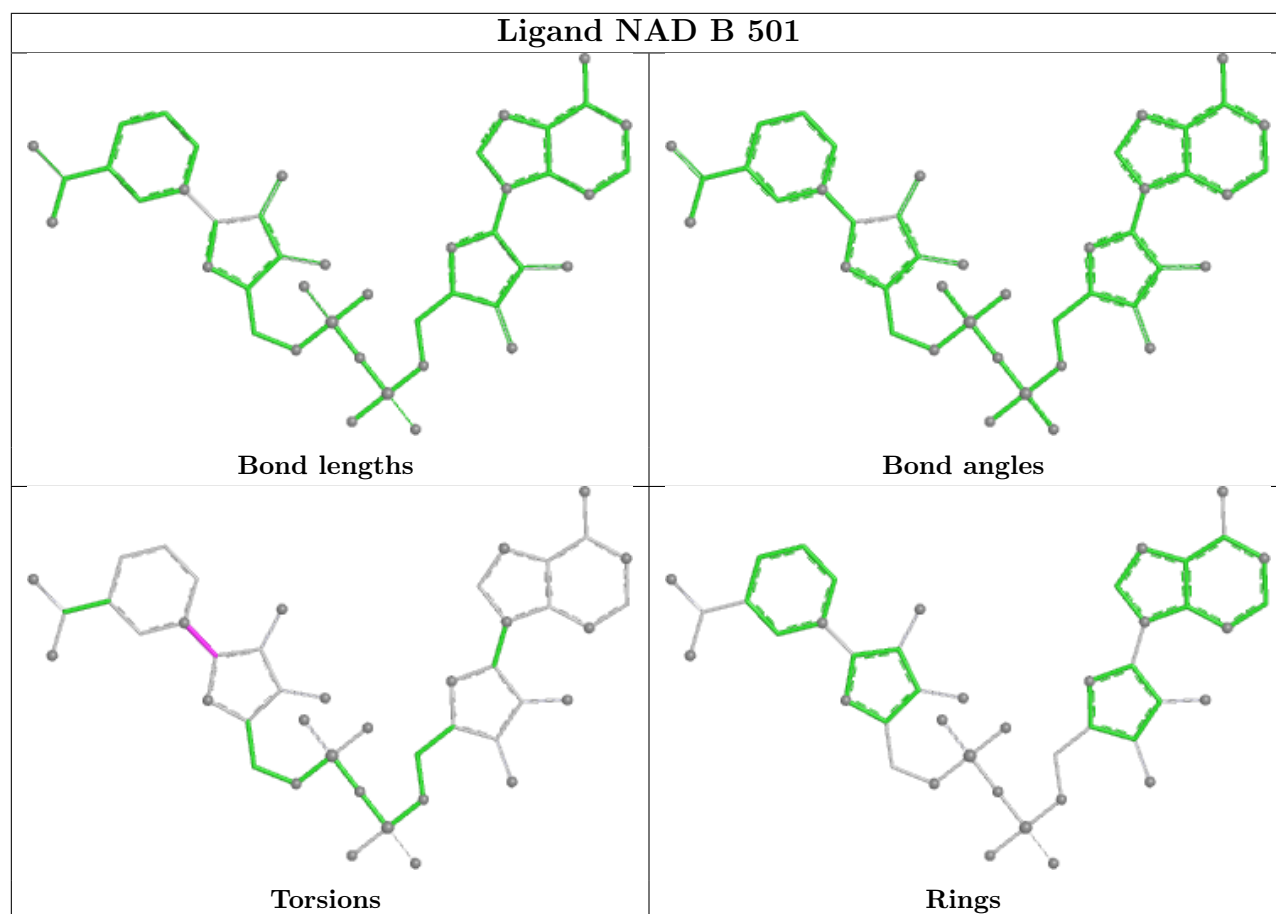
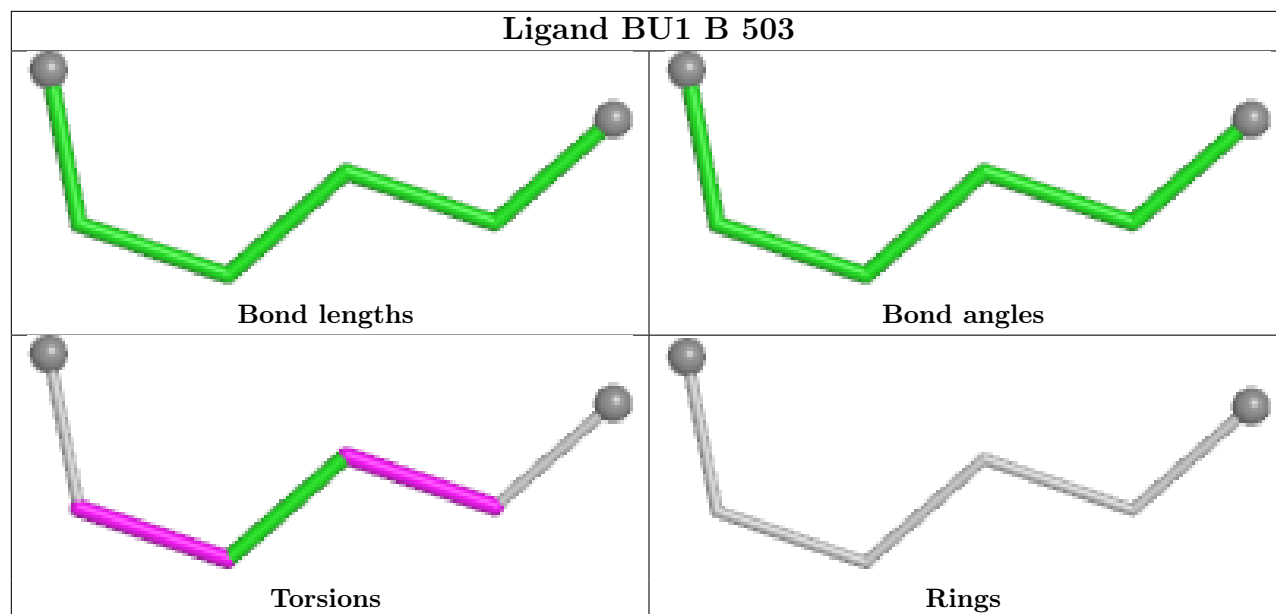


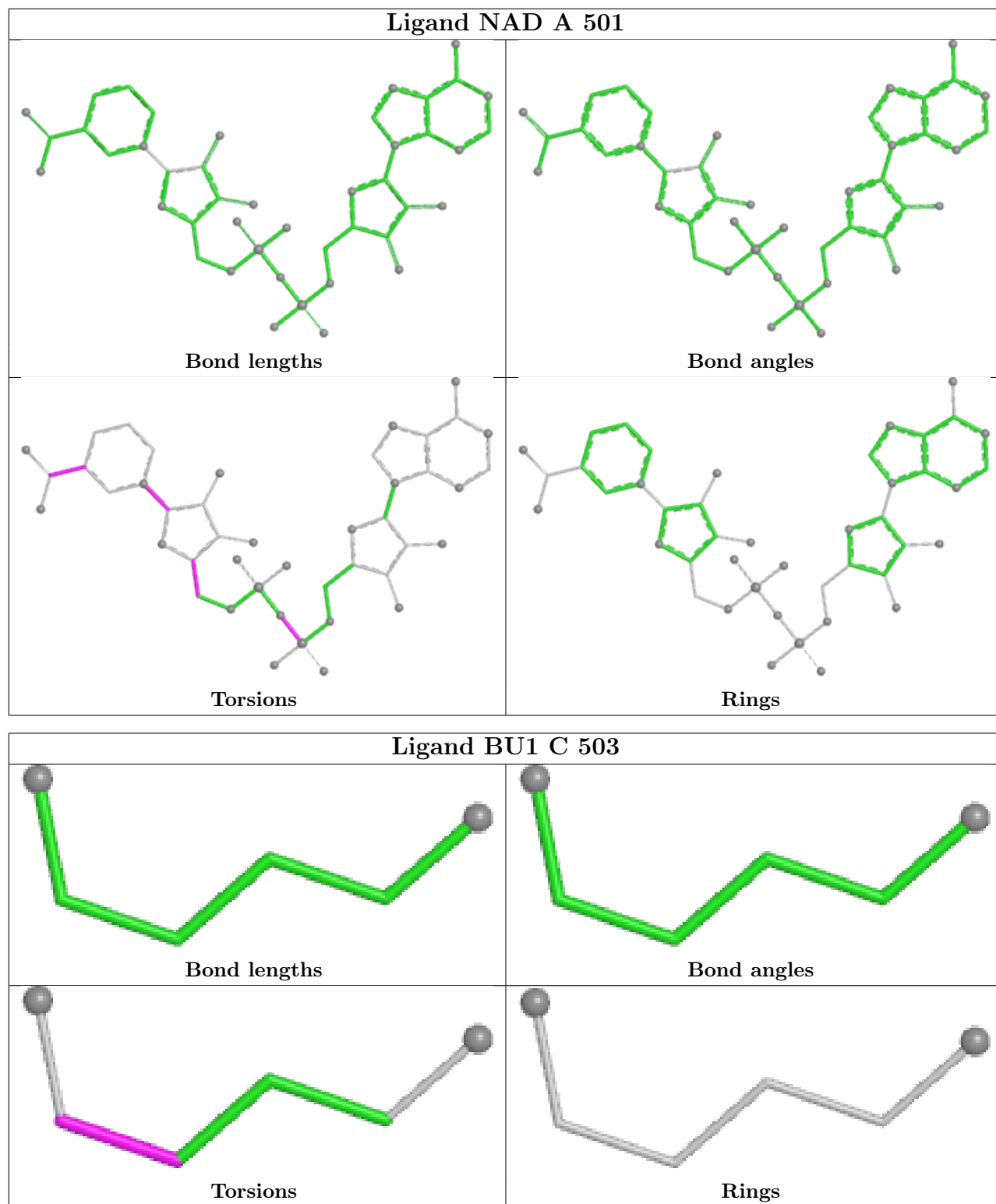
Ligand 3PG A 502

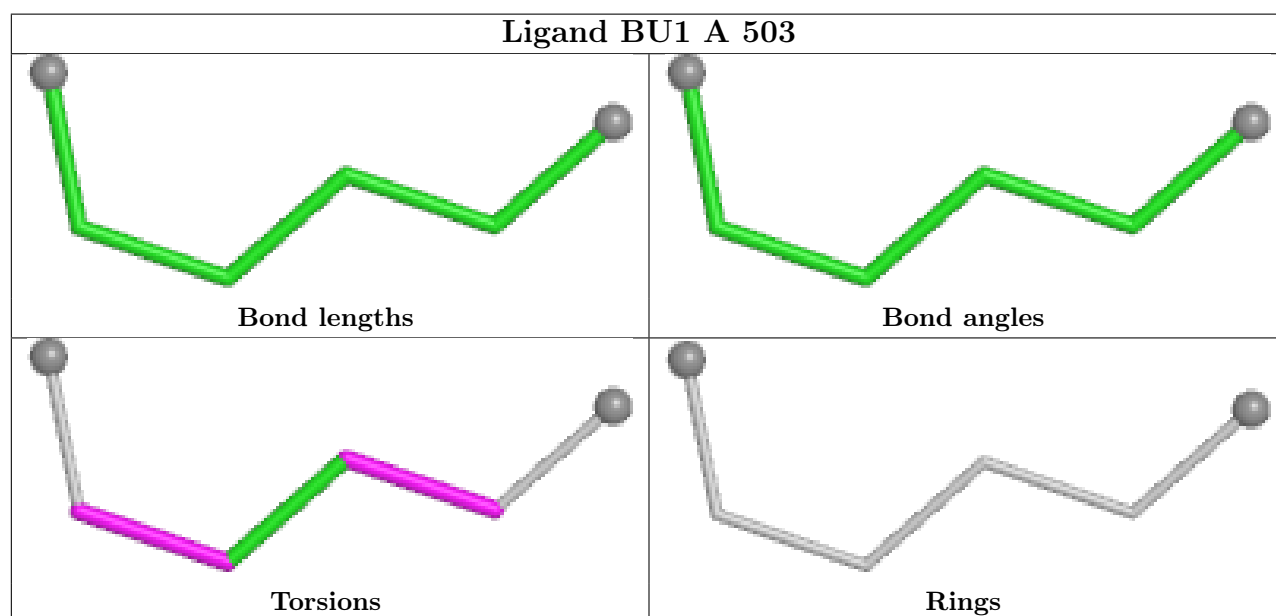
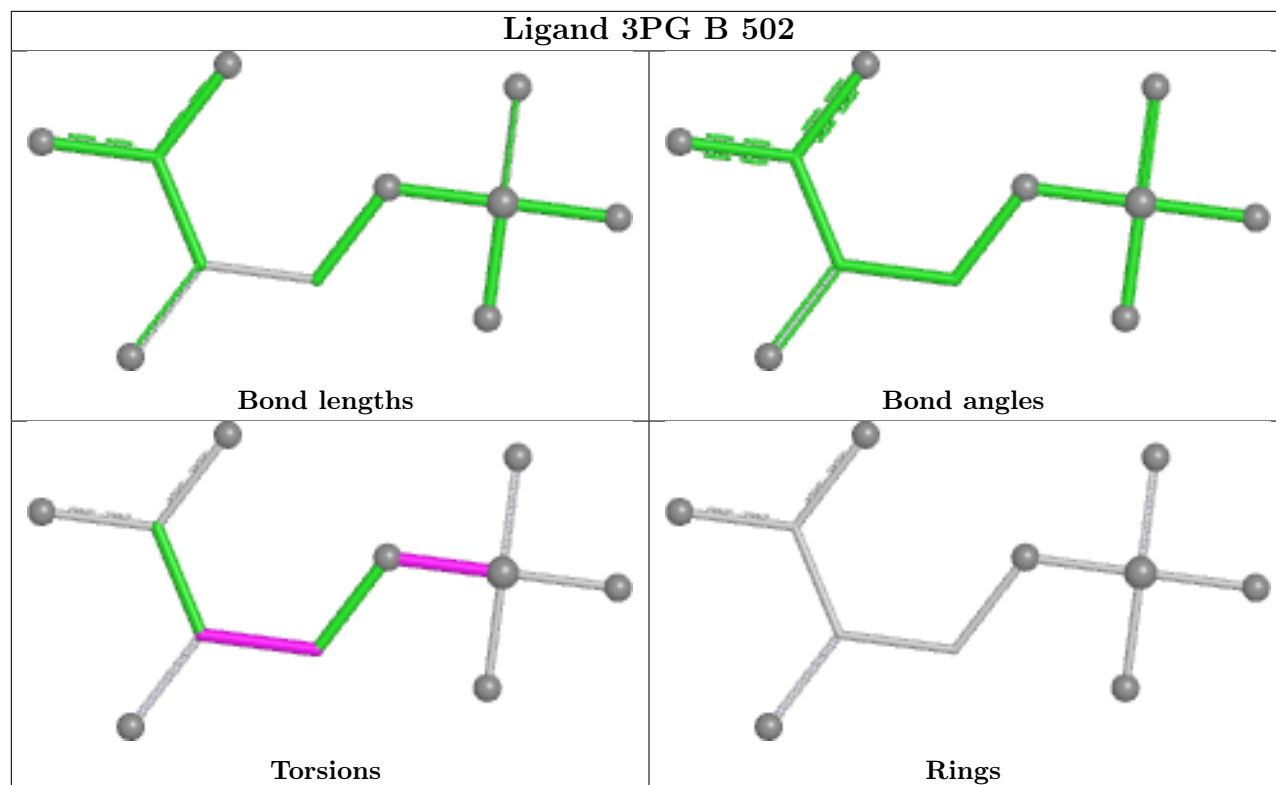


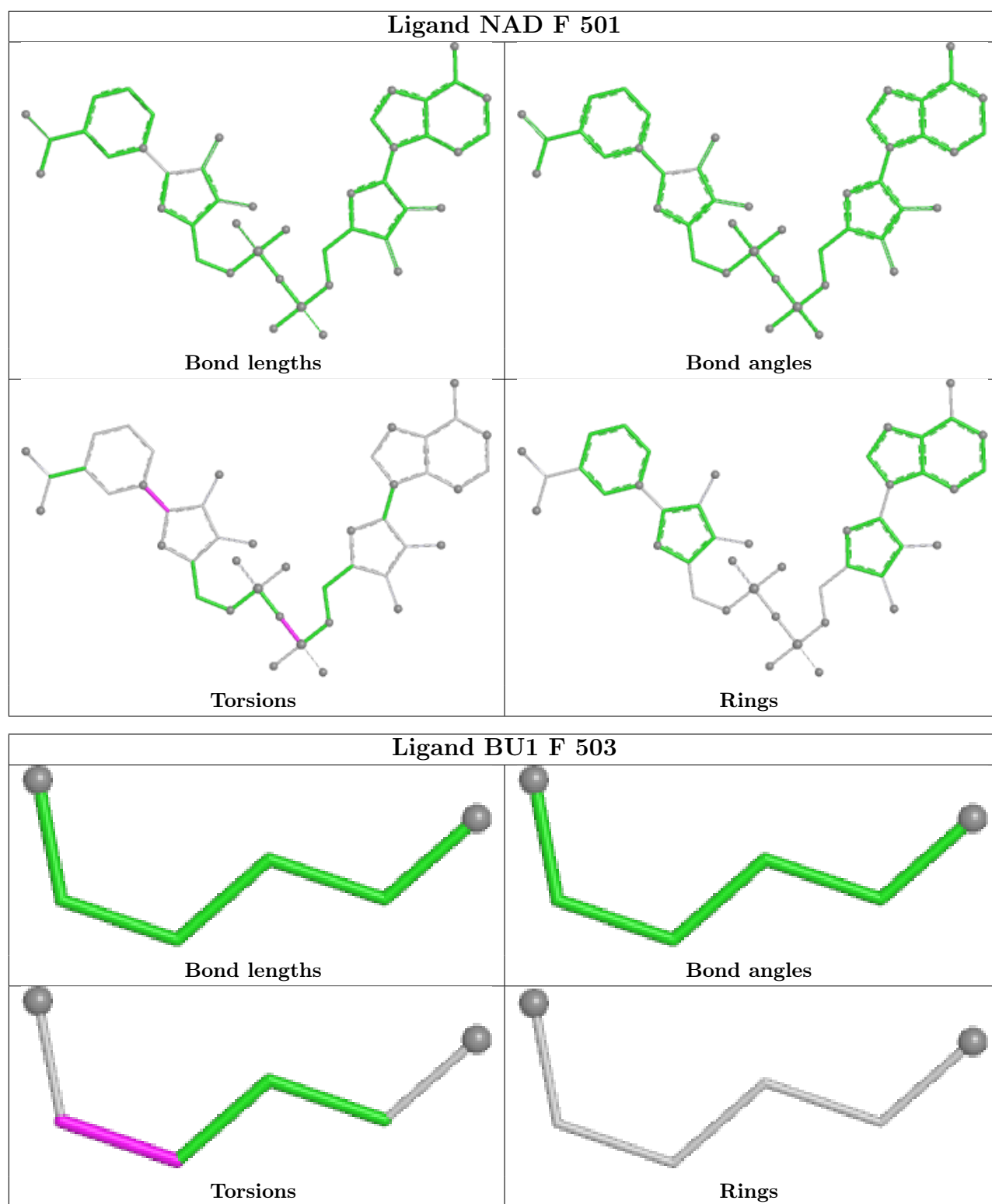
Ligand NAD C 501











5.7 Other polymers [i](#)

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	440/473 (93%)	0.57	19 (4%) 40 36	47, 72, 100, 123	0
1	B	437/473 (92%)	0.81	27 (6%) 26 23	54, 76, 103, 119	0
1	C	436/473 (92%)	0.77	31 (7%) 22 19	54, 75, 101, 112	0
1	F	439/473 (92%)	0.32	10 (2%) 61 58	46, 62, 88, 112	0
All	All	1752/1892 (92%)	0.62	87 (4%) 34 30	46, 71, 100, 123	0

All (87) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	43	SER	6.2
1	A	31	THR	5.5
1	B	23	PRO	5.4
1	B	31	THR	5.1
1	C	55	SER	4.2
1	B	42	VAL	4.2
1	A	43	SER	4.1
1	F	32	GLN	4.0
1	C	47	GLN	4.0
1	B	58	ASP	4.0
1	C	42	VAL	3.9
1	F	55	SER	3.7
1	C	392	TYR	3.5
1	F	57	GLY	3.4
1	A	38	LEU	3.3
1	C	40	ARG	3.2
1	A	24	GLY	3.1
1	C	23	PRO	3.0
1	A	56	THR	3.0
1	B	445	ASN	2.8
1	B	446	GLN	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	229	ILE	2.8
1	A	58	ASP	2.8
1	A	451	ASP	2.8
1	B	47	GLN	2.7
1	F	100	ASP	2.7
1	B	100	ASP	2.7
1	B	93	GLU	2.6
1	C	58	ASP	2.6
1	A	392	TYR	2.6
1	A	192	ARG	2.6
1	C	447	SER	2.6
1	C	92	ASP	2.6
1	B	434	ILE	2.5
1	B	101	VAL	2.5
1	C	33	SER	2.5
1	C	230	VAL	2.5
1	C	54	PHE	2.5
1	B	350	GLY	2.4
1	F	42	VAL	2.4
1	B	206	TYR	2.4
1	A	233	MET	2.4
1	C	205	GLY	2.4
1	C	95	ILE	2.4
1	A	245	ASP	2.4
1	C	39	PRO	2.4
1	C	394	GLN	2.4
1	C	445	ASN	2.3
1	F	58	ASP	2.3
1	B	117	LEU	2.3
1	B	71	ALA	2.2
1	C	393	ASP	2.2
1	B	24	GLY	2.2
1	A	55	SER	2.2
1	C	334	SER	2.2
1	B	84	PHE	2.2
1	A	174	LEU	2.2
1	A	250	LYS	2.2
1	B	27	SER	2.2
1	B	55	SER	2.2
1	A	194	TRP	2.2
1	C	328	GLU	2.2
1	B	87	SER	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	88	SER	2.2
1	F	39	PRO	2.1
1	A	262	PRO	2.1
1	F	463	VAL	2.1
1	C	113	THR	2.1
1	C	227	TYR	2.1
1	A	420	ASP	2.1
1	A	47	GLN	2.1
1	B	174	LEU	2.1
1	C	247	LEU	2.1
1	F	44	ILE	2.1
1	B	82	VAL	2.1
1	C	192	ARG	2.0
1	C	87	SER	2.0
1	C	389	SER	2.0
1	C	233	MET	2.0
1	F	181	LEU	2.0
1	B	40	ARG	2.0
1	B	389	SER	2.0
1	C	31	THR	2.0
1	C	59	MET	2.0
1	C	112	LEU	2.0
1	C	181	LEU	2.0
1	A	42	VAL	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](#)

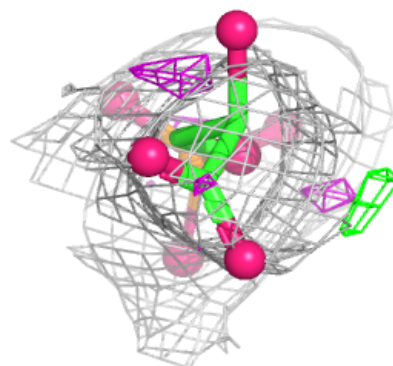
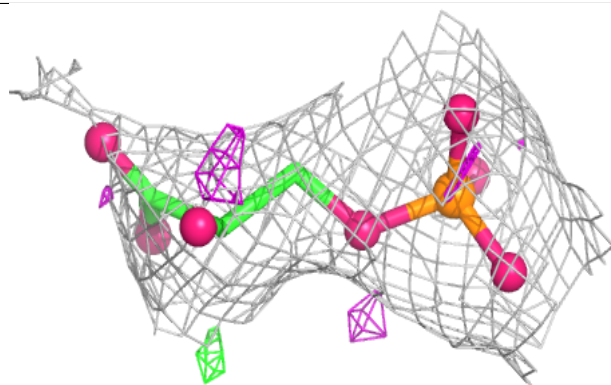
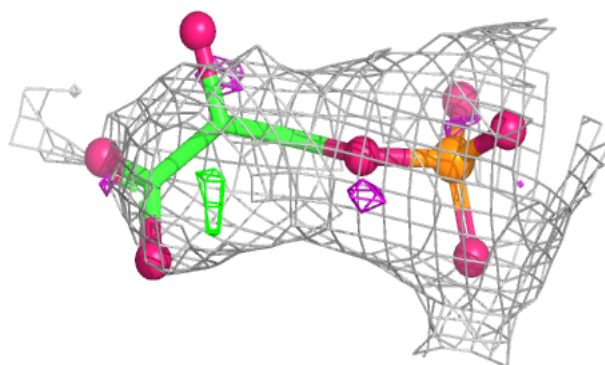
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
3	3PG	B	502	11/11	0.79	0.13	77,88,95,95	0
3	3PG	F	502	11/11	0.85	0.11	63,70,81,85	0
3	3PG	A	502	11/11	0.86	0.11	72,78,85,87	0
3	3PG	C	502	11/11	0.89	0.10	74,82,91,93	0
4	BU1	A	503	6/6	0.89	0.15	51,55,57,60	0
4	BU1	B	503	6/6	0.89	0.14	49,57,60,62	0
2	NAD	B	501	44/44	0.92	0.10	51,70,85,91	0
2	NAD	C	501	44/44	0.93	0.09	52,68,80,84	0
2	NAD	A	501	44/44	0.94	0.10	48,62,76,81	0
4	BU1	F	503	6/6	0.94	0.12	50,56,57,59	0
4	BU1	C	503	6/6	0.94	0.12	58,61,63,63	0
2	NAD	F	501	44/44	0.97	0.07	43,51,58,62	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.

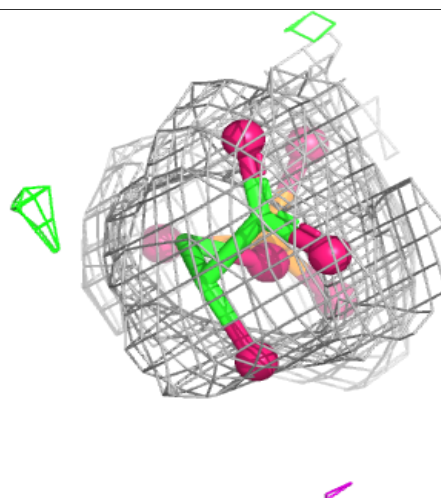
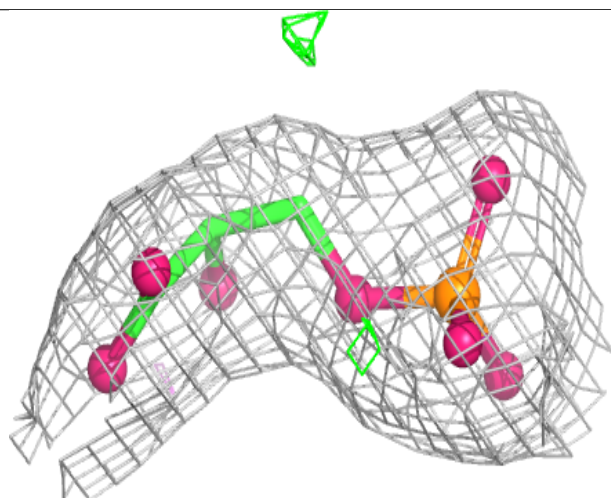
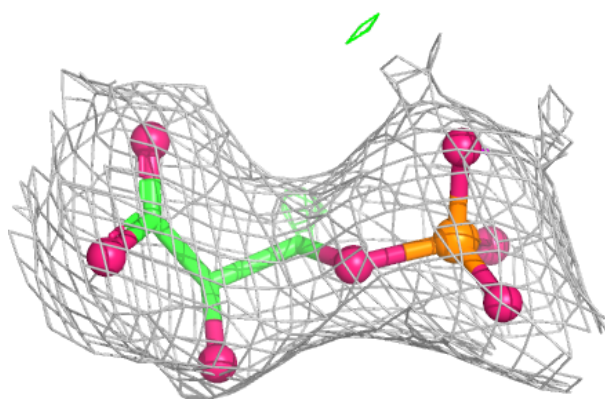
Electron density around 3PG B 502:

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



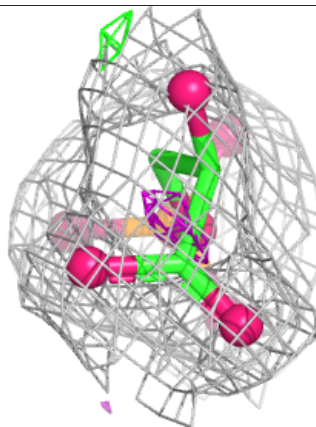
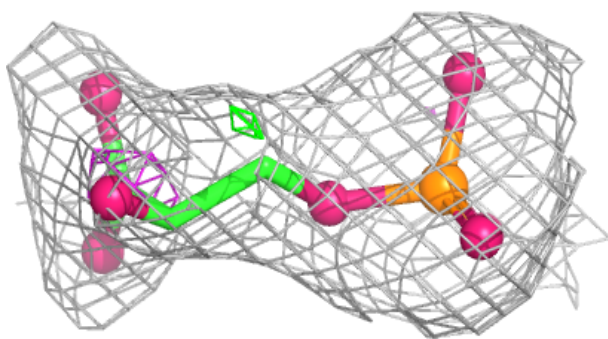
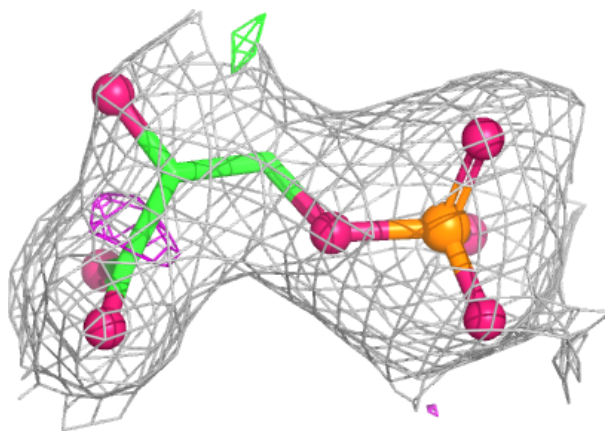
Electron density around 3PG F 502:

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

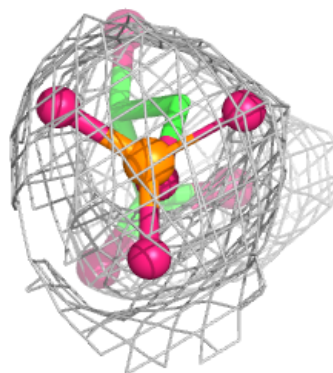
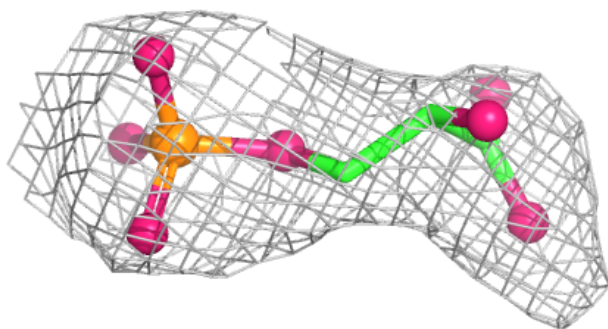
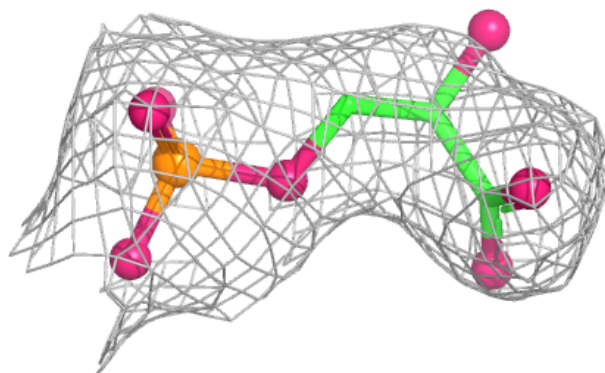


Electron density around 3PG A 502:

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

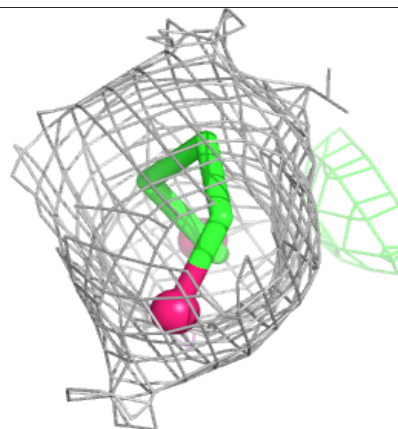
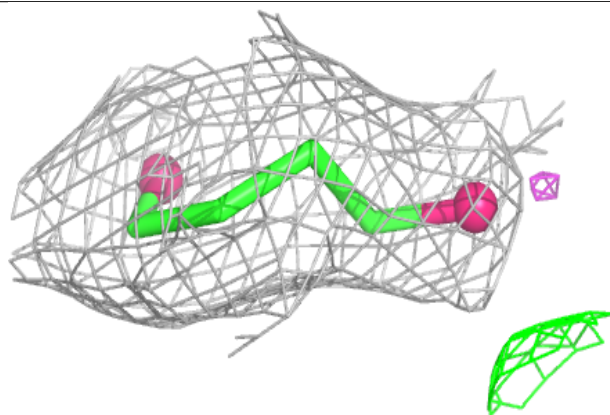
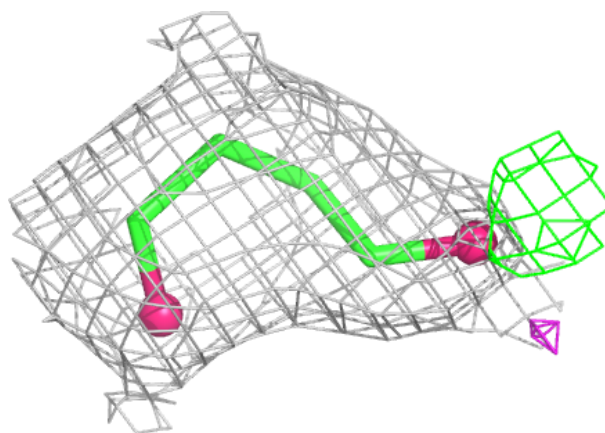
**Electron density around 3PG C 502:**

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



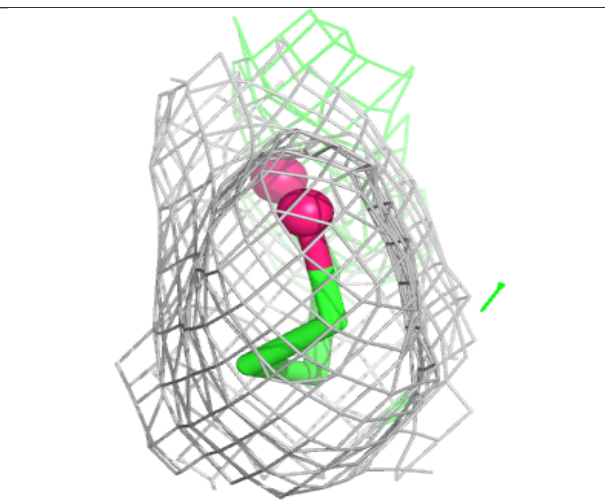
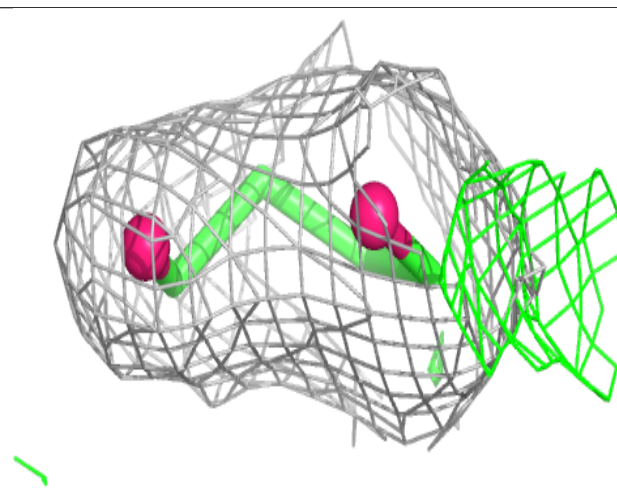
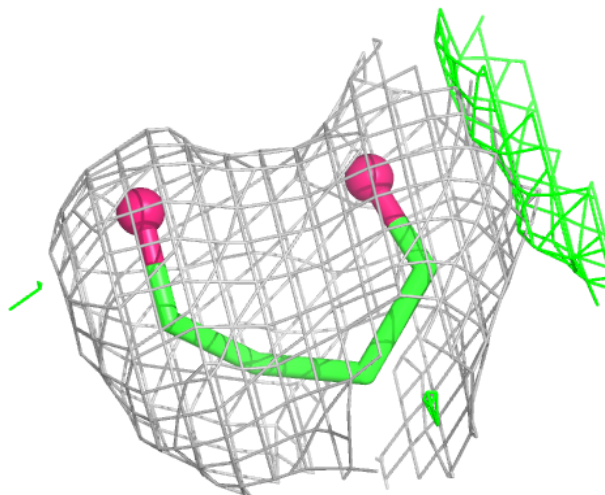
Electron density around BU1 A 503:

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



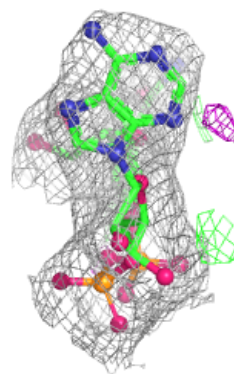
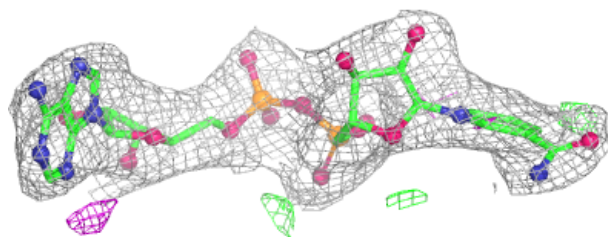
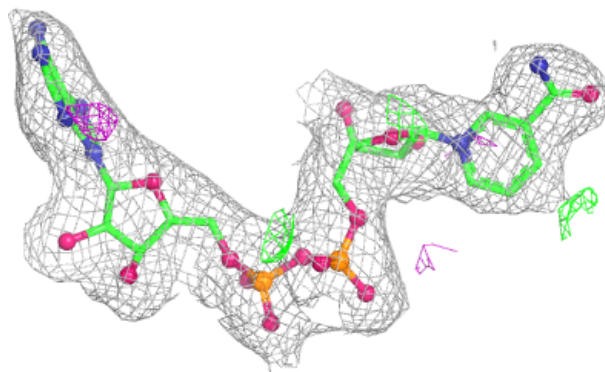
Electron density around BU1 B 503:

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

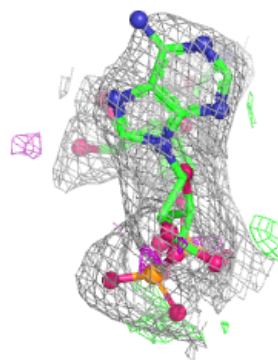
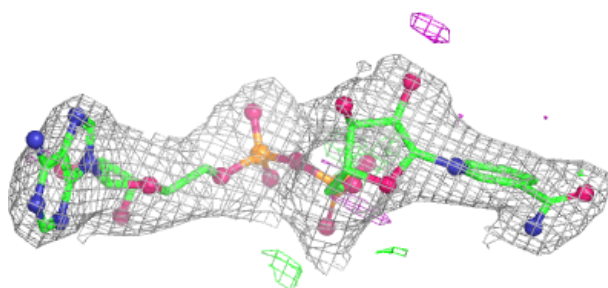
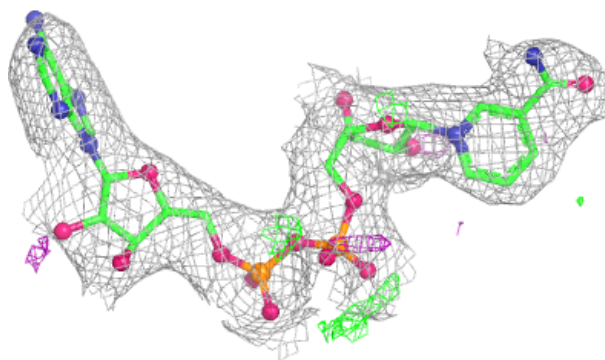


Electron density around NAD B 501:

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

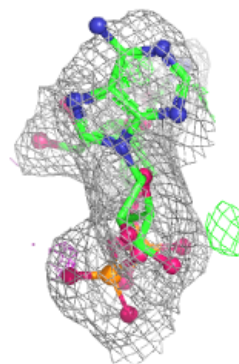
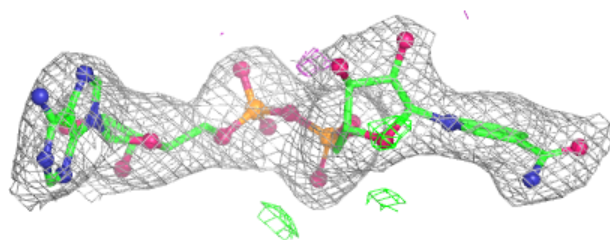
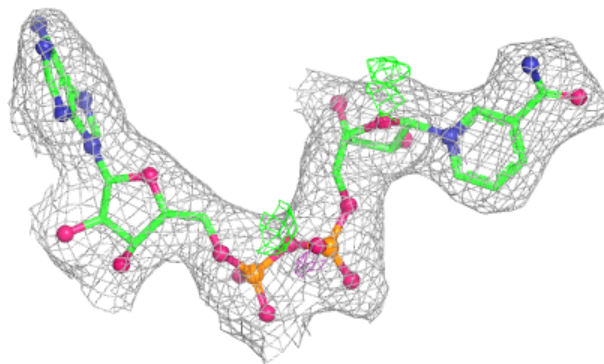
**Electron density around NAD C 501:**

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

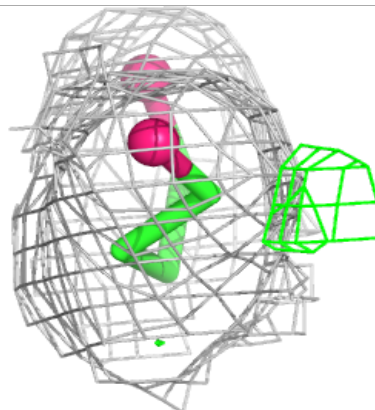
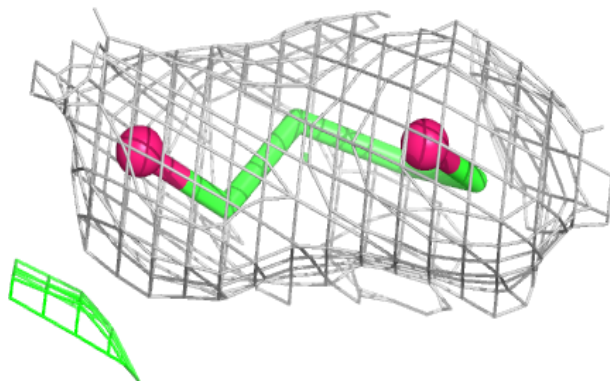
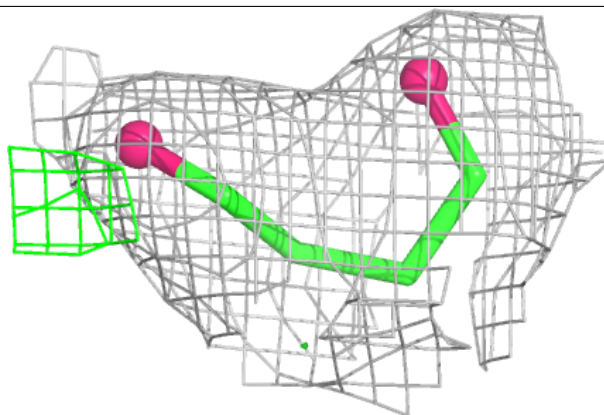


Electron density around NAD A 501:

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

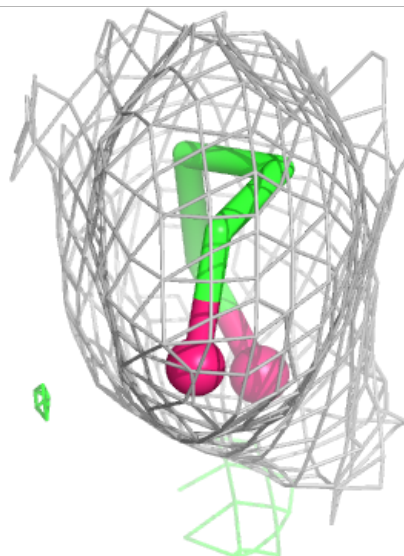
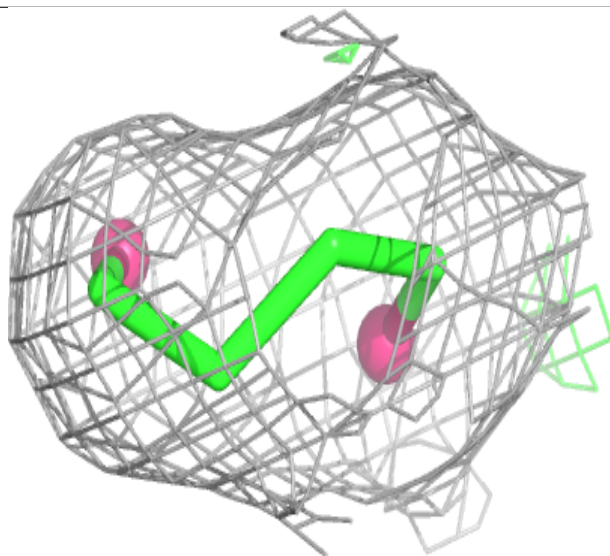
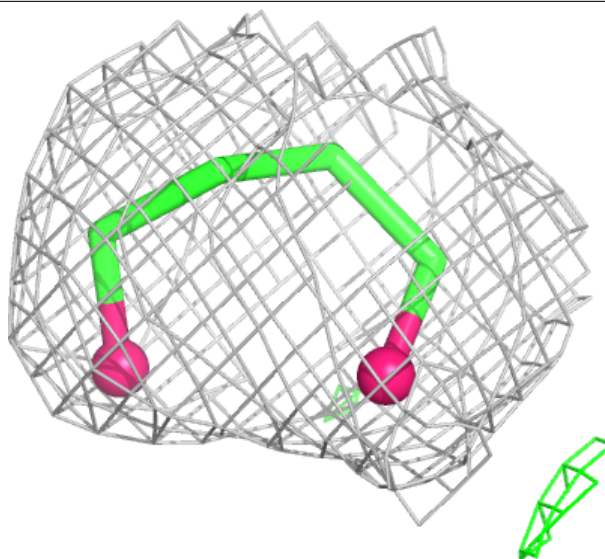
**Electron density around BU1 F 503:**

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



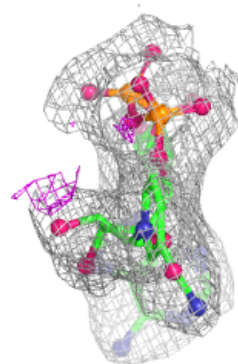
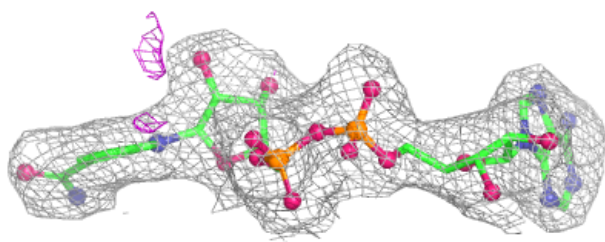
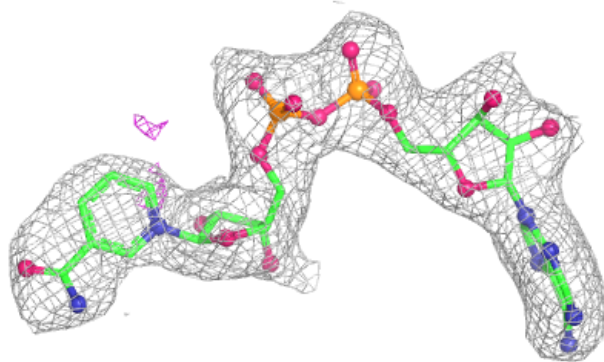
Electron density around BU1 C 503:

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



Electron density around NAD F 501:

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



6.5 Other polymers ⓘ

There are no such residues in this entry.