Full wwPDB X-ray Structure Validation Report ⓘ

Jan 6, 2025 – 02:57 pm GMT

PDB ID	:	8PIO
Title	:	Crystal structure of Ser33 in complex with PHP (3-phosphohydroxypyruvate)
Authors	:	Perrone, S.; Cifuente, J.O.; Marina, A.; Mastrella, L.; Trastoy, B.; Linster, C.L.; Guerin, M.E.
Deposited on	:	2023-06-22
Resolution	:	2.60 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:

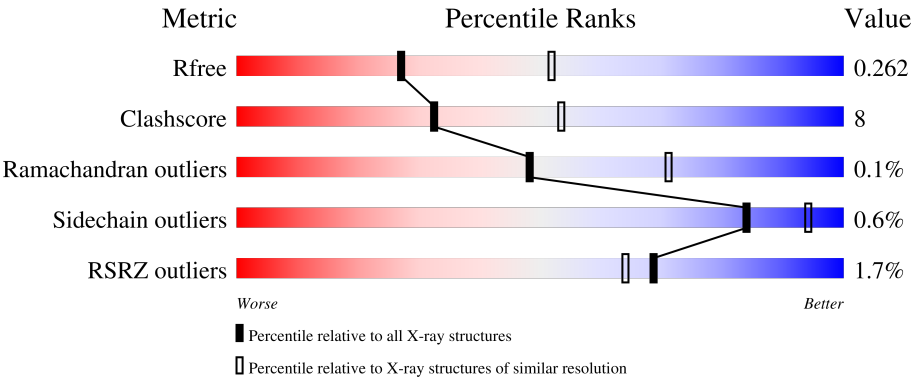
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	3.0
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.40

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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}	164625	3775 (2.60-2.60)
Clashscore	180529	4181 (2.60-2.60)
Ramachandran outliers	177936	4129 (2.60-2.60)
Sidechain outliers	177891	4129 (2.60-2.60)
RSRZ outliers	164620	3775 (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 <=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	469	2% 77% 17% 6%
1	B	469	77% 14% 9%
1	C	469	2% 76% 16% 8%
1	F	469	80% 15% 5%

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
3	HPV	C	503	-	X	-	-

2 Entry composition (i)

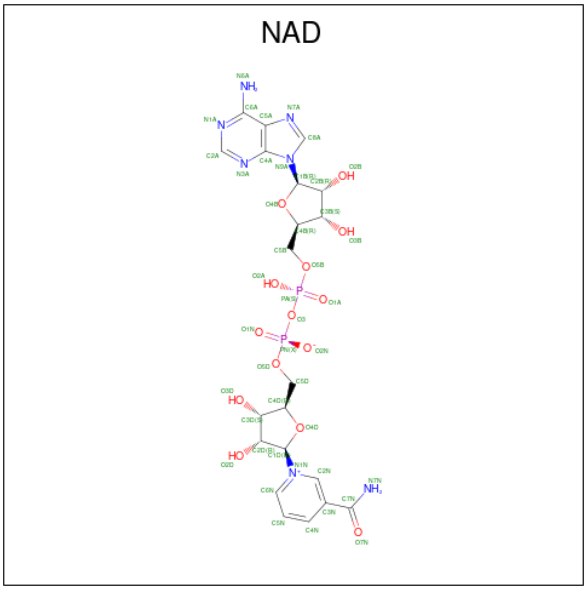
There are 5 unique types of molecules in this entry. The entry contains 13225 atoms, of which 0 are hydrogens and 0 are deuteriums.

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	445	Total	C	N	O	S	0	0	0
			3337	2111	577	639	10			
1	A	443	Total	C	N	O	S	0	0	0
			3273	2073	567	623	10			
1	B	429	Total	C	N	O	S	0	0	0
			3181	2007	545	619	10			
1	C	431	Total	C	N	O	S	0	0	0
			3155	1991	543	612	9			

- Molecule 2 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: $C_{21}H_{27}N_7O_{14}P_2$) (labeled 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		

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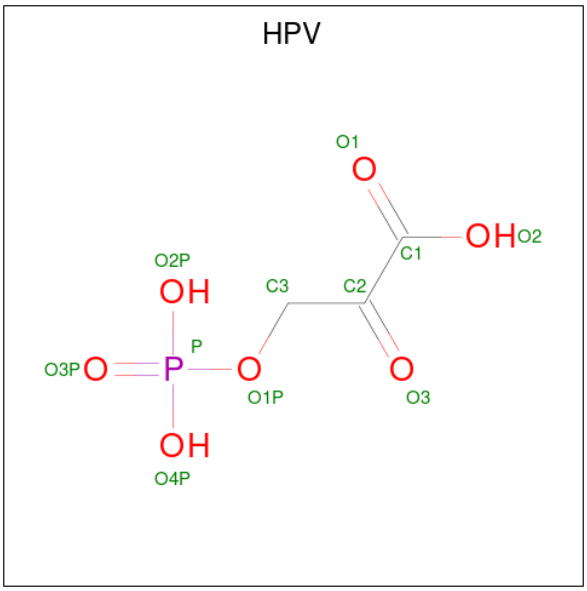
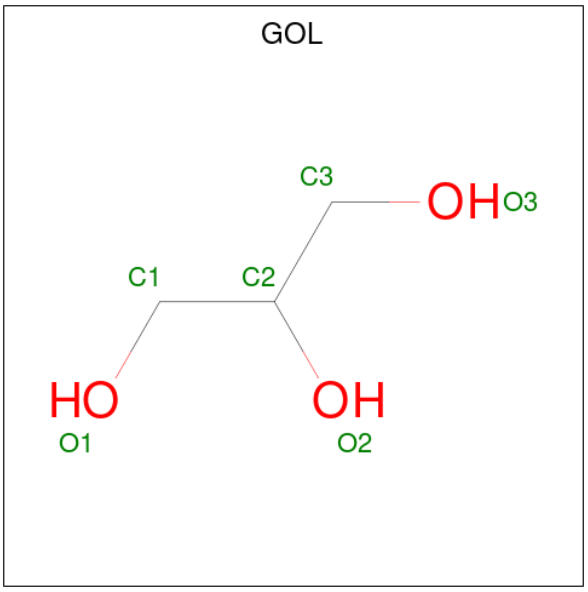
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total C N O P 44 21 7 14 2	0	0
2	C	1	Total C N O P 44 21 7 14 2	0	0

- Molecule 3 is 2-oxo-3-(phosphonooxy)propanoic acid (three-letter code: HPV) (formula: $C_3H_5O_7P$) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 5 is water.

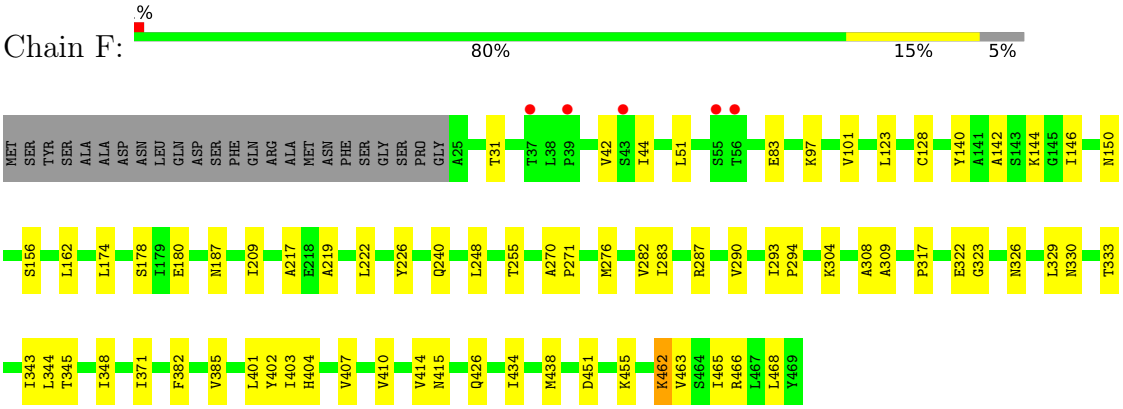
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	F	12	Total O 12 12	0	0
5	A	15	Total O 15 15	0	0
5	B	2	Total O 2 2	0	0
5	C	6	Total O 6 6	0	0

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.

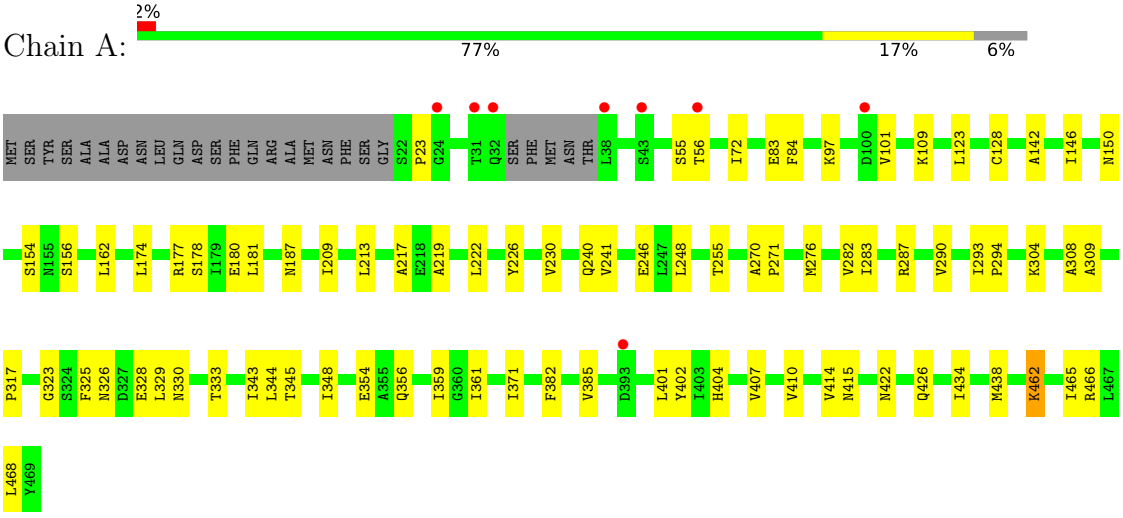
• Molecule 1: phosphoglycerate dehydrogenase

Chain F:



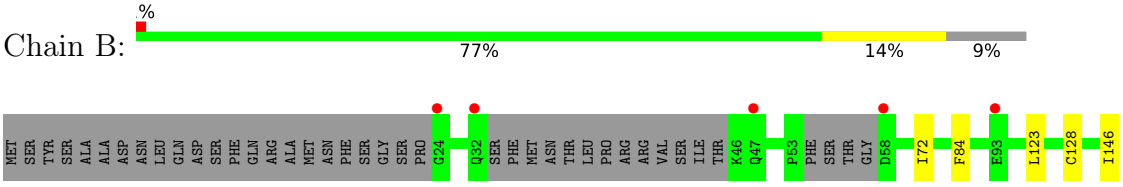
• Molecule 1: phosphoglycerate dehydrogenase

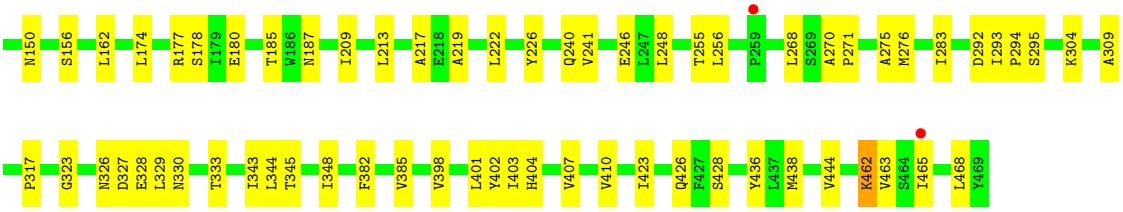
Chain A:



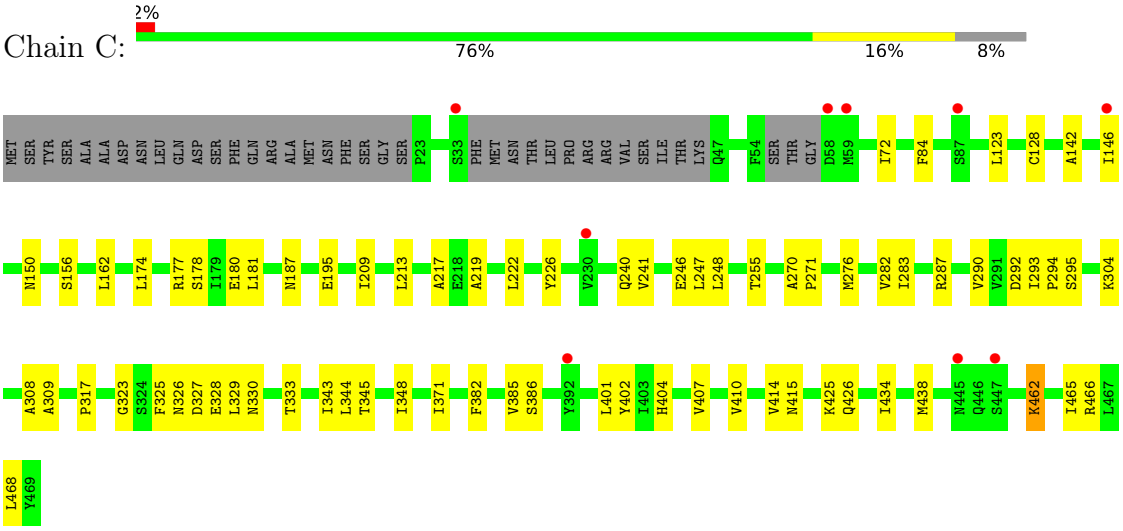
• Molecule 1: phosphoglycerate dehydrogenase

Chain B:





• Molecule 1: phosphoglycerate dehydrogenase



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α, β, γ	60.60Å 95.15Å 108.37Å 69.84° 88.72° 85.65°	Depositor
Resolution (Å)	52.04 – 2.60 52.04 – 2.60	Depositor EDS
% Data completeness (in resolution range)	97.9 (52.04-2.60) 97.9 (52.04-2.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.29 (at 2.61Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.222 , 0.262 0.223 , 0.262	Depositor DCC
R_{free} test set	3375 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å^2)	68.8	Xtriage
Anisotropy	0.130	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 40.8	EDS
L-test for twinning[2]	$< \|L\| > = 0.51$, $< L^2 > = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	13225	wwPDB-VP
Average B, all atoms (Å^2)	79.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.70% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: GOL, NAD, HPV

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	0/3328	0.69	0/4532
1	B	0.74	0/3233	0.69	0/4402
1	C	0.74	0/3205	0.69	0/4368
1	F	0.73	0/3393	0.69	0/4611
All	All	0.73	0/13159	0.69	0/17913

There are no bond length outliers.

There are no bond angle outliers.

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	3273	0	3203	65	0
1	B	3181	0	3070	56	0
1	C	3155	0	3037	53	0
1	F	3337	0	3293	56	0
2	A	44	0	26	0	0
2	B	44	0	26	0	0
2	C	44	0	26	0	0
2	F	44	0	26	0	0
3	A	11	0	2	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	11	0	2	0	0
3	C	11	0	2	0	0
3	F	11	0	2	0	0
4	A	12	0	16	1	0
4	B	6	0	8	0	0
4	C	6	0	8	3	0
5	A	15	0	0	2	0
5	B	2	0	0	0	0
5	C	6	0	0	1	0
5	F	12	0	0	0	0
All	All	13225	0	12747	201	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 (201) 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:F:219:ALA:HB1	1:B:219:ALA:HB1	1.53	0.91
1:F:426:GLN:HE21	1:A:426:GLN:HE21	1.26	0.84
1:B:398:VAL:HG23	1:B:444:VAL:HG21	1.67	0.76
1:A:466:ARG:HD3	5:A:614:HOH:O	1.85	0.75
1:B:426:GLN:HE21	1:C:426:GLN:HE21	1.35	0.74
1:B:398:VAL:HG23	1:B:444:VAL:CG2	2.18	0.73
1:B:398:VAL:CG2	1:B:444:VAL:HG21	2.19	0.72
1:B:323:GLY:O	1:B:326:ASN:ND2	2.26	0.68
1:F:323:GLY:O	1:F:326:ASN:ND2	2.25	0.68
1:A:415:ASN:HD21	1:A:426:GLN:HE22	1.42	0.67
1:C:323:GLY:O	1:C:326:ASN:ND2	2.28	0.67
1:B:403:ILE:HG22	1:B:463:VAL:HG21	1.78	0.66
1:A:323:GLY:O	1:A:326:ASN:ND2	2.29	0.65
1:A:219:ALA:HB1	1:C:219:ALA:HB1	1.79	0.64
1:F:403:ILE:HG22	1:F:463:VAL:HG21	1.79	0.64
1:F:403:ILE:C	1:F:463:VAL:HG22	2.19	0.64
1:B:403:ILE:C	1:B:463:VAL:HG22	2.18	0.63
1:C:415:ASN:HD21	1:C:426:GLN:HE22	1.48	0.62
1:F:403:ILE:HG22	1:F:463:VAL:CG2	2.31	0.61
1:F:426:GLN:NE2	1:A:426:GLN:HE21	1.96	0.60
1:C:270:ALA:HB3	1:C:271:PRO:HD3	1.83	0.60
1:F:178:SER:HA	1:B:344:LEU:O	2.01	0.60
1:B:403:ILE:HG22	1:B:463:VAL:CG2	2.31	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:403:ILE:O	1:F:463:VAL:HG22	2.02	0.59
1:B:270:ALA:HB3	1:B:271:PRO:HD3	1.84	0.59
1:F:270:ALA:HB3	1:F:271:PRO:HD3	1.84	0.59
1:B:403:ILE:O	1:B:463:VAL:HG22	2.02	0.58
1:A:97:LYS:O	1:A:101:VAL:HG23	2.04	0.58
1:A:270:ALA:HB3	1:A:271:PRO:HD3	1.84	0.58
1:A:344:LEU:O	1:C:178:SER:HA	2.04	0.58
1:F:344:LEU:O	1:B:178:SER:HA	2.04	0.58
1:F:97:LYS:O	1:F:101:VAL:HG23	2.04	0.57
1:A:328:GLU:N	1:A:328:GLU:OE1	2.38	0.57
1:B:423:ILE:H	4:C:502:GOL:C1	2.18	0.56
1:A:23:PRO:HG3	1:B:275:ALA:O	2.04	0.56
1:A:178:SER:HA	1:C:344:LEU:O	2.05	0.56
1:F:123:LEU:HD23	1:F:146:ILE:HD13	1.87	0.56
1:A:123:LEU:HD23	1:A:146:ILE:HD13	1.88	0.56
1:F:317:PRO:HD3	1:F:329:LEU:HD21	1.89	0.55
1:B:317:PRO:HD3	1:B:329:LEU:HD21	1.89	0.55
1:C:386:SER:HA	5:C:604:HOH:O	2.07	0.54
1:A:180:GLU:OE1	1:A:187:ASN:HB3	2.08	0.54
1:B:180:GLU:OE1	1:B:187:ASN:HB3	2.07	0.54
1:B:123:LEU:HD23	1:B:146:ILE:HD13	1.90	0.54
1:F:180:GLU:OE1	1:F:187:ASN:HB3	2.08	0.53
1:A:142:ALA:O	1:A:434:ILE:CD1	2.56	0.53
1:A:317:PRO:HD3	1:A:329:LEU:HD21	1.90	0.53
1:C:180:GLU:OE1	1:C:187:ASN:HB3	2.08	0.53
1:A:156:SER:HA	1:A:209:ILE:HG12	1.91	0.53
1:F:142:ALA:O	1:F:434:ILE:CD1	2.57	0.52
1:B:423:ILE:H	4:C:502:GOL:H11	1.74	0.52
1:F:462:LYS:HD2	1:F:462:LYS:C	2.30	0.52
1:F:226:TYR:CZ	1:F:240:GLN:HB2	2.44	0.52
1:A:401:LEU:HD11	1:A:468:LEU:HD11	1.92	0.51
1:B:156:SER:HA	1:B:209:ILE:HG12	1.92	0.51
1:C:226:TYR:CZ	1:C:240:GLN:HB2	2.46	0.51
1:F:255:THR:HA	1:F:283:ILE:O	2.11	0.51
1:C:402:TYR:CE2	1:C:404:HIS:HB3	2.46	0.51
1:F:415:ASN:OD1	1:A:415:ASN:ND2	2.43	0.51
1:B:226:TYR:CZ	1:B:240:GLN:HB2	2.45	0.51
1:C:317:PRO:HD3	1:C:329:LEU:HD21	1.92	0.51
1:F:426:GLN:HE21	1:A:426:GLN:NE2	2.03	0.51
1:A:255:THR:HA	1:A:283:ILE:O	2.10	0.51
1:A:462:LYS:HD2	1:A:462:LYS:C	2.31	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:248:LEU:O	1:A:276:MET:HA	2.11	0.51
1:B:462:LYS:HD2	1:B:462:LYS:C	2.31	0.51
1:F:451:ASP:OD2	1:F:455:LYS:NZ	2.45	0.51
1:A:226:TYR:CZ	1:A:240:GLN:HB2	2.46	0.50
1:B:255:THR:HA	1:B:283:ILE:O	2.12	0.50
1:C:123:LEU:HD23	1:C:146:ILE:HD13	1.93	0.50
1:C:142:ALA:O	1:C:434:ILE:CD1	2.59	0.50
1:F:248:LEU:O	1:F:276:MET:HA	2.11	0.50
1:C:255:THR:HA	1:C:283:ILE:O	2.11	0.50
1:F:156:SER:HA	1:F:209:ILE:HG12	1.93	0.50
1:C:462:LYS:HD2	1:C:462:LYS:C	2.31	0.50
1:B:248:LEU:O	1:B:276:MET:HA	2.12	0.49
1:A:309:ALA:HA	1:A:343:ILE:O	2.12	0.49
1:C:248:LEU:O	1:C:276:MET:HA	2.12	0.49
1:F:309:ALA:HA	1:F:343:ILE:O	2.13	0.49
1:F:402:TYR:HD1	1:F:465:ILE:HD13	1.77	0.49
1:B:402:TYR:CE2	1:B:404:HIS:HB3	2.47	0.49
1:C:309:ALA:HA	1:C:343:ILE:O	2.12	0.49
1:C:345:THR:HB	1:C:348:ILE:HG12	1.95	0.49
1:F:402:TYR:CE2	1:F:404:HIS:HB3	2.49	0.48
1:A:402:TYR:HD1	1:A:465:ILE:HD13	1.78	0.48
1:F:322:GLU:HA	1:B:185:THR:HG23	1.94	0.48
1:A:402:TYR:CE2	1:A:404:HIS:HB3	2.48	0.48
1:F:293:ILE:N	1:F:294:PRO:CD	2.77	0.48
1:F:401:LEU:HD11	1:F:468:LEU:HD11	1.95	0.48
1:B:402:TYR:HD1	1:B:465:ILE:HD13	1.79	0.48
1:F:42:VAL:O	1:A:230:VAL:HG13	2.14	0.48
1:A:142:ALA:O	1:A:434:ILE:HD13	2.13	0.48
1:B:309:ALA:HA	1:B:343:ILE:O	2.13	0.47
1:B:426:GLN:HE21	1:C:426:GLN:NE2	2.08	0.47
1:B:385:VAL:HG12	1:B:438:MET:SD	2.54	0.47
1:C:330:ASN:HB2	1:C:333:THR:HG23	1.97	0.47
1:C:385:VAL:HG12	1:C:438:MET:SD	2.54	0.47
1:B:345:THR:HB	1:B:348:ILE:HG12	1.96	0.47
1:B:401:LEU:HD11	1:B:468:LEU:HD11	1.97	0.47
1:A:276:MET:O	1:A:304:LYS:NZ	2.43	0.47
1:A:348:ILE:HD13	1:C:177:ARG:HG3	1.97	0.47
1:C:293:ILE:N	1:C:294:PRO:CD	2.77	0.47
1:A:385:VAL:HG12	1:A:438:MET:SD	2.55	0.46
1:B:293:ILE:N	1:B:294:PRO:CD	2.78	0.46
1:A:217:ALA:O	1:A:222:LEU:HB2	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:241:VAL:CG1	1:B:246:GLU:HB3	2.45	0.46
1:C:410:VAL:O	1:C:414:VAL:HG23	2.14	0.46
1:C:276:MET:O	1:C:304:LYS:NZ	2.45	0.46
1:F:31:THR:HG21	1:A:361:ILE:HD12	1.97	0.46
1:F:330:ASN:HB2	1:F:333:THR:HG23	1.97	0.46
1:F:385:VAL:HG12	1:F:438:MET:SD	2.55	0.46
1:F:226:TYR:CE1	1:F:240:GLN:HB2	2.51	0.46
1:A:293:ILE:N	1:A:294:PRO:CD	2.79	0.46
1:B:217:ALA:O	1:B:222:LEU:HB2	2.16	0.46
1:F:142:ALA:O	1:F:434:ILE:HD13	2.17	0.45
1:A:410:VAL:O	1:A:414:VAL:HG23	2.16	0.45
1:A:109:LYS:HA	1:A:109:LYS:HD3	1.80	0.45
1:B:330:ASN:HB2	1:B:333:THR:HG23	1.98	0.45
1:F:217:ALA:O	1:F:222:LEU:HB2	2.17	0.45
1:A:287:ARG:HB2	1:A:290:VAL:HG23	1.97	0.45
1:A:241:VAL:CG1	1:A:246:GLU:HB3	2.47	0.45
1:A:156:SER:HA	1:A:209:ILE:CG1	2.47	0.45
1:F:348:ILE:HD13	1:B:177:ARG:HG3	1.99	0.45
1:C:217:ALA:O	1:C:222:LEU:HB2	2.17	0.45
1:A:174:LEU:HD21	1:C:162:LEU:CD1	2.47	0.44
1:C:407:VAL:O	1:C:410:VAL:HG23	2.16	0.44
1:F:162:LEU:CD1	1:B:174:LEU:HD21	2.47	0.44
1:F:410:VAL:O	1:F:414:VAL:HG23	2.17	0.44
1:B:72:ILE:HD11	1:B:84:PHE:CD1	2.53	0.44
1:C:156:SER:HA	1:C:209:ILE:HG12	1.98	0.44
1:F:345:THR:HB	1:F:348:ILE:HG12	1.98	0.44
1:C:326:ASN:OD1	1:C:328:GLU:HB2	2.18	0.44
1:A:330:ASN:HB2	1:A:333:THR:HG23	1.98	0.44
1:C:226:TYR:CE1	1:C:240:GLN:HB2	2.52	0.44
1:F:407:VAL:O	1:F:410:VAL:HG23	2.18	0.44
1:A:72:ILE:HD11	1:A:84:PHE:CD1	2.53	0.44
1:B:407:VAL:O	1:B:410:VAL:HG23	2.16	0.44
1:F:156:SER:HA	1:F:209:ILE:CG1	2.48	0.44
1:C:401:LEU:O	1:C:465:ILE:HD12	2.18	0.44
1:F:426:GLN:NE2	1:A:426:GLN:NE2	2.65	0.43
1:A:241:VAL:HG13	1:A:246:GLU:OE1	2.18	0.43
1:B:156:SER:HA	1:B:209:ILE:CG1	2.48	0.43
1:B:292:ASP:OD2	1:B:295:SER:OG	2.32	0.43
1:C:402:TYR:HD1	1:C:465:ILE:HD13	1.82	0.43
1:C:241:VAL:HG11	1:C:247:LEU:HB2	2.00	0.43
1:A:128:CYS:O	1:A:150:ASN:HA	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:241:VAL:HG13	1:B:246:GLU:OE1	2.18	0.43
1:A:83:GLU:OE2	1:A:97:LYS:CE	2.67	0.43
1:B:401:LEU:O	1:B:465:ILE:HD12	2.19	0.43
1:C:156:SER:HA	1:C:209:ILE:CG1	2.48	0.42
1:F:401:LEU:O	1:F:465:ILE:HD12	2.19	0.42
1:A:181:LEU:HD22	1:C:325:PHE:CG	2.54	0.42
1:A:345:THR:HB	1:A:348:ILE:HG12	1.99	0.42
1:C:72:ILE:HD11	1:C:84:PHE:CD1	2.54	0.42
1:F:140:TYR:OH	1:F:144:LYS:HE3	2.19	0.42
1:F:174:LEU:HD21	1:B:162:LEU:CD1	2.49	0.42
1:A:83:GLU:OE2	1:A:97:LYS:HE3	2.18	0.42
1:F:371:ILE:O	1:F:466:ARG:NH2	2.53	0.42
1:C:213:LEU:HD21	1:C:255:THR:HG21	2.01	0.42
1:B:241:VAL:HG11	1:B:246:GLU:HB3	2.02	0.42
1:F:83:GLU:OE2	1:F:97:LYS:HE3	2.19	0.42
1:A:326:ASN:OD1	1:A:328:GLU:HB2	2.18	0.42
1:A:371:ILE:O	1:A:466:ARG:NH2	2.52	0.42
1:A:401:LEU:O	1:A:465:ILE:HD12	2.19	0.42
1:A:162:LEU:CD1	1:C:174:LEU:HD21	2.49	0.42
1:A:213:LEU:HD21	1:A:255:THR:HG21	2.01	0.42
1:A:422:ASN:OD1	4:A:501:GOL:O2	2.38	0.42
1:C:292:ASP:OD2	1:C:295:SER:OG	2.32	0.42
1:F:128:CYS:O	1:F:150:ASN:HA	2.19	0.42
1:A:154:SER:OG	1:A:359:ILE:HG12	2.20	0.42
1:C:371:ILE:O	1:C:466:ARG:NH2	2.53	0.42
1:C:282:VAL:O	1:C:308:ALA:HA	2.19	0.42
1:A:407:VAL:O	1:A:410:VAL:HG23	2.20	0.41
1:A:325:PHE:CG	1:C:181:LEU:HD22	2.55	0.41
1:B:276:MET:O	1:B:304:LYS:NZ	2.44	0.41
1:B:423:ILE:HB	4:C:502:GOL:H2	2.01	0.41
1:A:177:ARG:HG3	1:C:348:ILE:HD13	2.02	0.41
1:F:415:ASN:HD22	1:F:415:ASN:HA	1.70	0.41
1:A:356:GLN:NE2	5:A:602:HOH:O	2.52	0.41
1:B:128:CYS:O	1:B:150:ASN:HA	2.20	0.41
1:C:401:LEU:HD11	1:C:468:LEU:HD11	2.01	0.41
1:A:55:SER:O	1:A:56:THR:C	2.58	0.41
1:A:282:VAL:O	1:A:308:ALA:HA	2.20	0.41
1:B:326:ASN:OD1	1:B:328:GLU:HB2	2.20	0.41
1:F:282:VAL:O	1:F:308:ALA:HA	2.20	0.41
1:A:241:VAL:HG13	1:A:246:GLU:CD	2.41	0.41
1:F:51:LEU:HB3	1:F:468:LEU:HB3	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:276:MET:O	1:F:304:LYS:NZ	2.45	0.41
1:B:256:LEU:HD11	1:B:268:LEU:HD12	2.03	0.41
1:C:128:CYS:O	1:C:150:ASN:HA	2.20	0.41
1:F:287:ARG:HB2	1:F:290:VAL:HG23	2.03	0.41
1:A:354:GLU:HG2	1:C:195:GLU:HB2	2.03	0.41
1:B:428:SER:HA	1:B:436:TYR:O	2.21	0.41
1:B:426:GLN:NE2	1:C:426:GLN:HE21	2.09	0.40
1:C:241:VAL:CG2	1:C:246:GLU:CD	2.89	0.40
1:C:287:ARG:HB2	1:C:290:VAL:HG23	2.03	0.40
1:B:213:LEU:HD21	1:B:255:THR:HG21	2.03	0.40
1:F:83:GLU:OE2	1:F:97:LYS:CE	2.69	0.40
1:B:428:SER:O	1:C:425:LYS:HA	2.22	0.40
1:C:327:ASP:HA	1:C:330:ASN:O	2.22	0.40
1:B:327:ASP:HA	1:B:330:ASN:O	2.22	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	439/469 (94%)	416 (95%)	23 (5%)	0	100	100
1	B	423/469 (90%)	407 (96%)	16 (4%)	0	100	100
1	C	425/469 (91%)	409 (96%)	16 (4%)	0	100	100
1	F	443/469 (94%)	417 (94%)	25 (6%)	1 (0%)	44	66
All	All	1730/1876 (92%)	1649 (95%)	80 (5%)	1 (0%)	48	71

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	F	44	ILE

5.3.2 Protein sidechains (i)

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	338/401 (84%)	336 (99%)	2 (1%)	84 94
1	B	329/401 (82%)	327 (99%)	2 (1%)	84 94
1	C	322/401 (80%)	320 (99%)	2 (1%)	84 94
1	F	351/401 (88%)	349 (99%)	2 (1%)	84 94
All	All	1340/1604 (84%)	1332 (99%)	8 (1%)	84 94

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	F	382	PHE
1	F	462	LYS
1	A	382	PHE
1	A	462	LYS
1	B	382	PHE
1	B	462	LYS
1	C	382	PHE
1	C	462	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	F	78	GLN
1	F	212	GLN
1	F	415	ASN
1	F	426	GLN
1	A	372	ASN
1	A	394	GLN
1	A	415	ASN
1	B	415	ASN
1	B	426	GLN
1	C	32	GLN
1	C	372	ASN

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Mol	Chain	Res	Type
1	C	415	ASN

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

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	$\#\|Z\| > 2$
2	NAD	B	501	-	42,48,48	0.62	0	50,73,73	0.63	1 (2%)
4	GOL	A	501	-	5,5,5	0.15	0	5,5,5	0.28	0
4	GOL	A	503	-	5,5,5	0.17	0	5,5,5	0.32	0
2	NAD	A	502	-	42,48,48	0.70	0	50,73,73	0.63	1 (2%)
3	HPV	C	503	-	10,10,10	1.63	3 (30%)	13,14,14	1.79	3 (23%)
3	HPV	A	504	-	10,10,10	1.85	3 (30%)	13,14,14	1.67	3 (23%)
3	HPV	B	503	-	10,10,10	1.10	1 (10%)	13,14,14	1.10	2 (15%)
4	GOL	C	502	-	5,5,5	0.13	0	5,5,5	0.26	0
2	NAD	F	501	-	42,48,48	0.72	0	50,73,73	0.61	1 (2%)
3	HPV	F	502	-	10,10,10	1.99	3 (30%)	13,14,14	1.33	2 (15%)
4	GOL	B	502	-	5,5,5	0.21	0	5,5,5	0.27	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
2	NAD	C	501	-	42,48,48	0.74	0	50,73,73	0.58	1 (2%)

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	501	-	-	7/26/62/62	0/5/5/5
4	GOL	A	501	-	-	2/4/4/4	-
4	GOL	A	503	-	-	2/4/4/4	-
2	NAD	A	502	-	-	9/26/62/62	0/5/5/5
3	HPV	C	503	-	-	9/9/10/10	-
3	HPV	A	504	-	-	7/9/10/10	-
3	HPV	B	503	-	-	6/9/10/10	-
4	GOL	C	502	-	-	2/4/4/4	-
2	NAD	F	501	-	-	8/26/62/62	0/5/5/5
3	HPV	F	502	-	-	5/9/10/10	-
4	GOL	B	502	-	-	3/4/4/4	-
2	NAD	C	501	-	-	6/26/62/62	0/5/5/5

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	F	502	HPV	O1-C1	4.83	1.35	1.22
3	C	503	HPV	O1-C1	3.99	1.33	1.22
3	A	504	HPV	O1-C1	3.97	1.33	1.22
3	A	504	HPV	P-O1P	2.85	1.69	1.60
3	F	502	HPV	P-O1P	2.82	1.69	1.60
3	A	504	HPV	C2-C1	2.36	1.56	1.53
3	C	503	HPV	O2-C1	-2.32	1.23	1.30
3	B	503	HPV	C2-C1	2.24	1.56	1.53
3	C	503	HPV	P-O1P	2.19	1.67	1.60
3	F	502	HPV	O2-C1	-2.17	1.24	1.30

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	C	503	HPV	O1-C1-C2	-4.71	115.43	121.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	504	HPV	O1-C1-C2	-4.45	115.78	121.72
3	F	502	HPV	O1-C1-C2	-3.65	116.85	121.72
3	A	504	HPV	O2-C1-C2	2.62	121.13	113.97
3	C	503	HPV	O2-C1-C2	2.59	121.07	113.97
2	A	502	NAD	C5A-C6A-N6A	2.44	124.06	120.35
3	B	503	HPV	O3-C2-C1	2.41	122.92	119.43
2	B	501	NAD	C5A-C6A-N6A	2.39	123.99	120.35
2	C	501	NAD	C5A-C6A-N6A	2.33	123.89	120.35
3	C	503	HPV	O3-C2-C1	2.32	122.80	119.43
2	F	501	NAD	C5A-C6A-N6A	2.28	123.82	120.35
3	A	504	HPV	O3-C2-C1	2.18	122.59	119.43
3	B	503	HPV	O1P-C3-C2	-2.14	107.00	110.32
3	F	502	HPV	O2-C1-C2	2.13	119.79	113.97

There are no chirality outliers.

All (66) 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	502	NAD	O4D-C1D-N1N-C2N
2	A	502	NAD	O4D-C1D-N1N-C6N
2	A	502	NAD	C2D-C1D-N1N-C2N
2	A	502	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	HPV	O3-C2-C3-O1P
3	F	502	HPV	C1-C2-C3-O1P
3	F	502	HPV	C3-O1P-P-O4P
3	F	502	HPV	C3-O1P-P-O2P
3	A	504	HPV	O3-C2-C3-O1P
3	A	504	HPV	C1-C2-C3-O1P
3	A	504	HPV	C3-O1P-P-O4P
3	A	504	HPV	C3-O1P-P-O2P

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Mol	Chain	Res	Type	Atoms
3	B	503	HPV	O3-C2-C3-O1P
3	B	503	HPV	C1-C2-C3-O1P
3	B	503	HPV	O1-C1-C2-O3
3	B	503	HPV	C3-O1P-P-O4P
3	B	503	HPV	C3-O1P-P-O2P
3	C	503	HPV	O3-C2-C3-O1P
3	C	503	HPV	C1-C2-C3-O1P
3	C	503	HPV	C3-O1P-P-O4P
3	C	503	HPV	C3-O1P-P-O2P
3	C	503	HPV	C2-C3-O1P-P
4	A	501	GOL	O1-C1-C2-C3
4	B	502	GOL	C1-C2-C3-O3
4	A	501	GOL	O1-C1-C2-O2
4	C	502	GOL	C1-C2-C3-O3
4	B	502	GOL	O2-C2-C3-O3
4	C	502	GOL	O2-C2-C3-O3
2	A	502	NAD	O4D-C4D-C5D-O5D
2	F	501	NAD	O4B-C4B-C5B-O5B
2	A	502	NAD	C3D-C4D-C5D-O5D
3	A	504	HPV	C3-O1P-P-O3P
3	C	503	HPV	C3-O1P-P-O3P
4	A	503	GOL	O1-C1-C2-O2
2	B	501	NAD	PN-O3-PA-O1A
3	A	504	HPV	O1-C1-C2-C3
3	A	504	HPV	O1-C1-C2-O3
3	B	503	HPV	O1-C1-C2-C3
3	C	503	HPV	O1-C1-C2-C3
3	C	503	HPV	O2-C1-C2-O3
2	F	501	NAD	PN-O3-PA-O2A
2	A	502	NAD	PN-O3-PA-O1A
2	A	502	NAD	PN-O3-PA-O2A
2	B	501	NAD	PN-O3-PA-O2A
2	C	501	NAD	PN-O3-PA-O2A
4	B	502	GOL	O1-C1-C2-C3
2	A	502	NAD	O4B-C4B-C5B-O5B
2	B	501	NAD	O4B-C4B-C5B-O5B
2	F	501	NAD	PN-O3-PA-O1A
2	F	501	NAD	C3B-C4B-C5B-O5B
3	F	502	HPV	C3-O1P-P-O3P
2	C	501	NAD	O4B-C4B-C5B-O5B
4	A	503	GOL	O1-C1-C2-C3
3	C	503	HPV	O1-C1-C2-O3

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	501	GOL	1	0
4	C	502	GOL	3	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 then 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.

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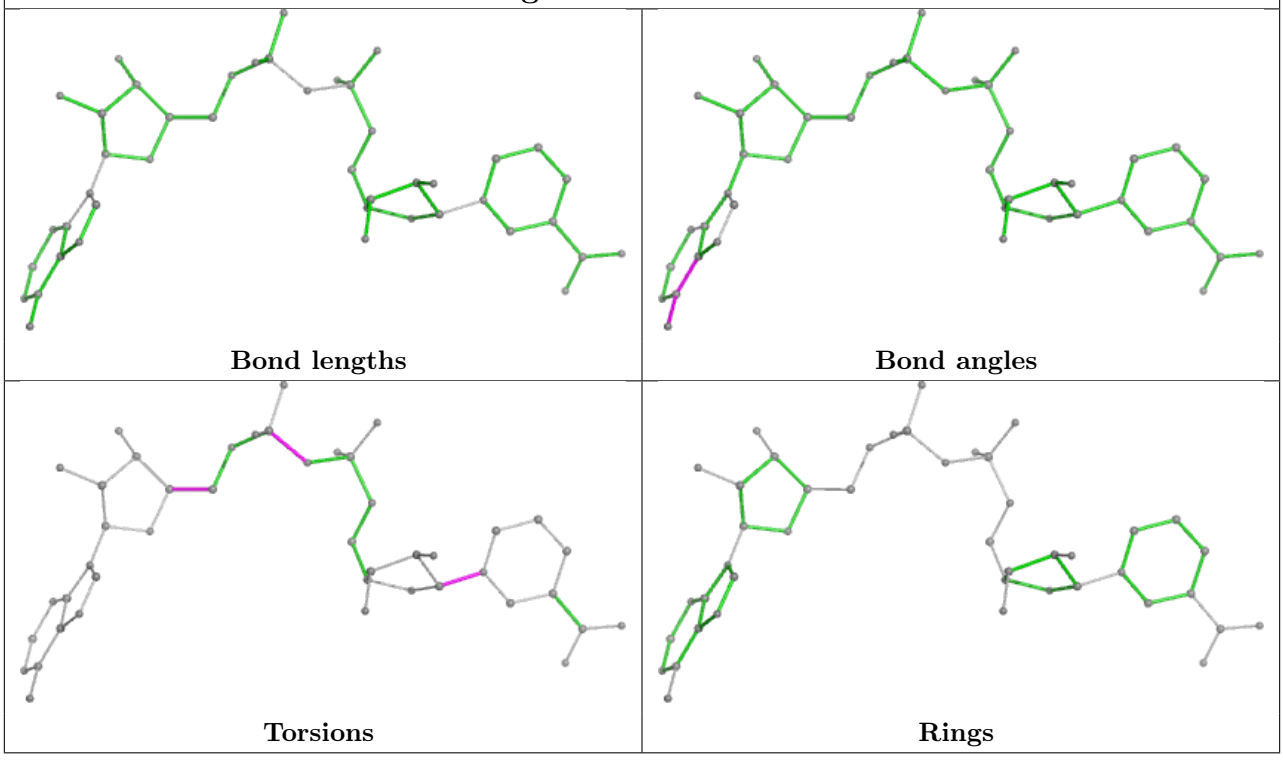
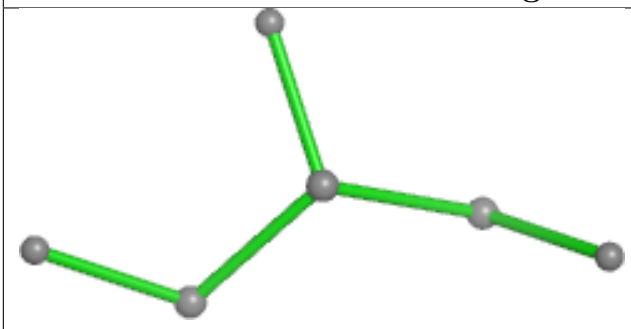


Ligand GOL A 501



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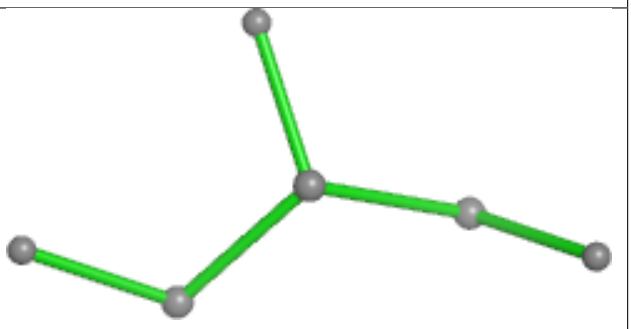
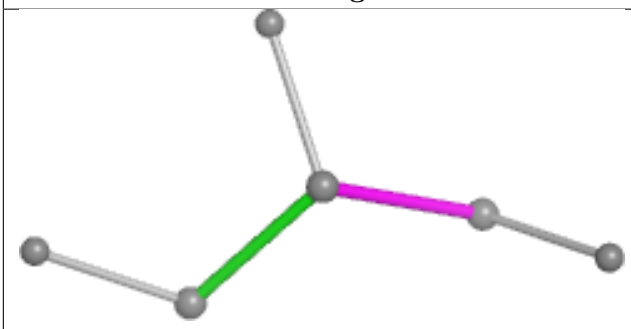
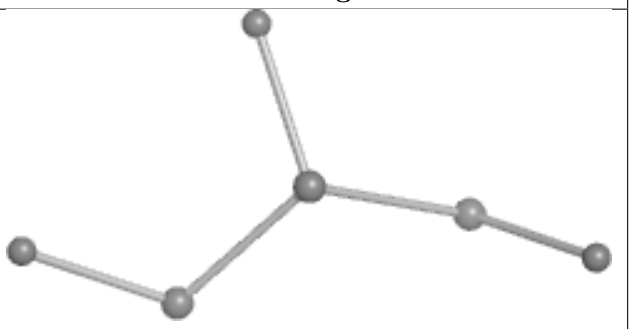


Bond angles

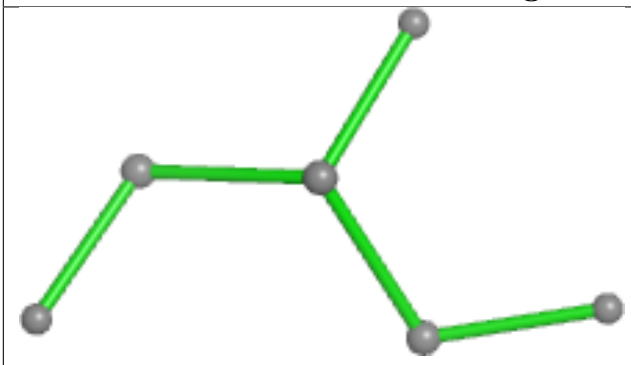


Torsions



Rings

Ligand GOL A 503



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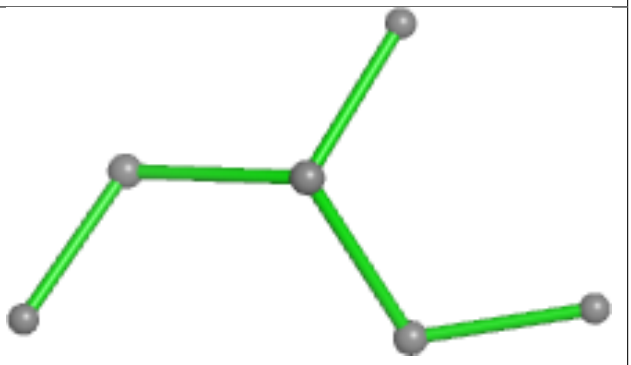

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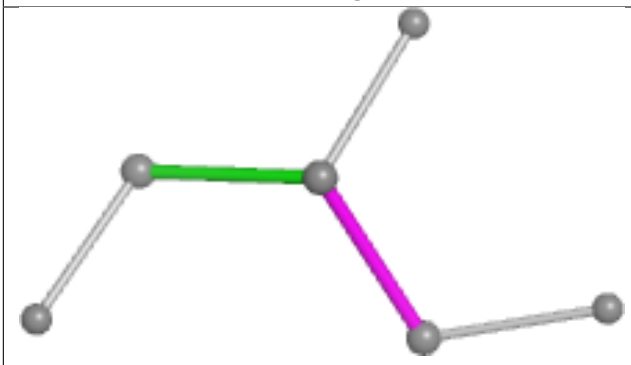

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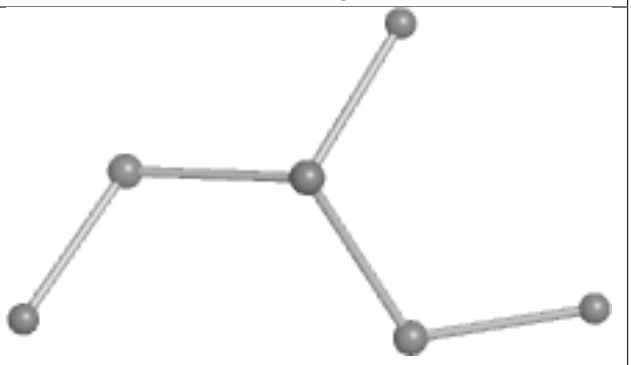
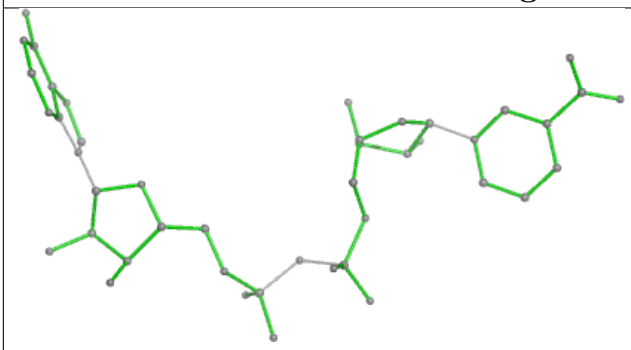
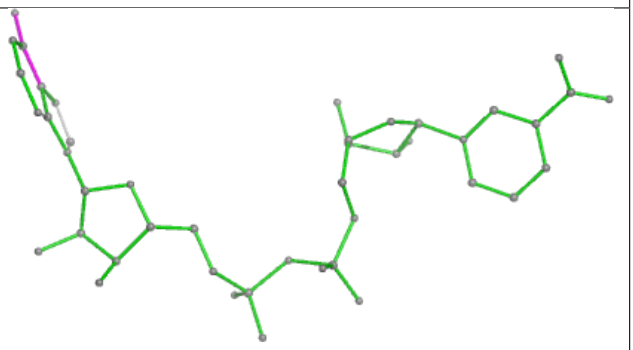


Rings

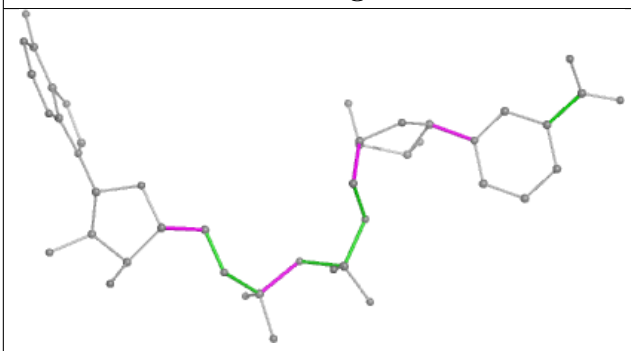
Ligand NAD A 502



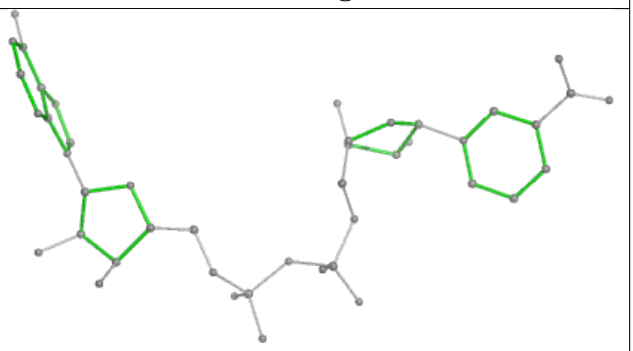
Bond lengths



Bond angles

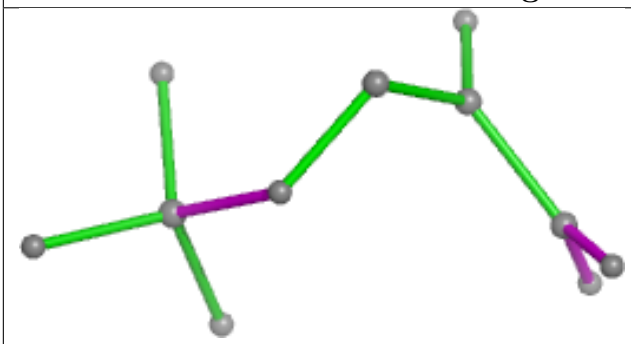


Torsions



Rings

Ligand HPV C 503



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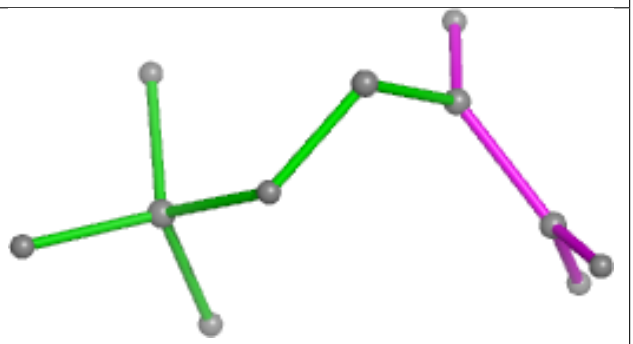
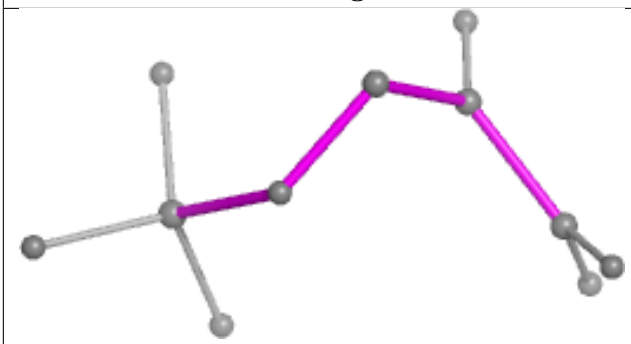
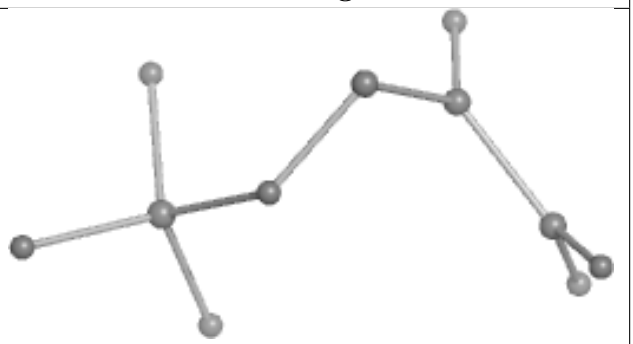


Bond angles

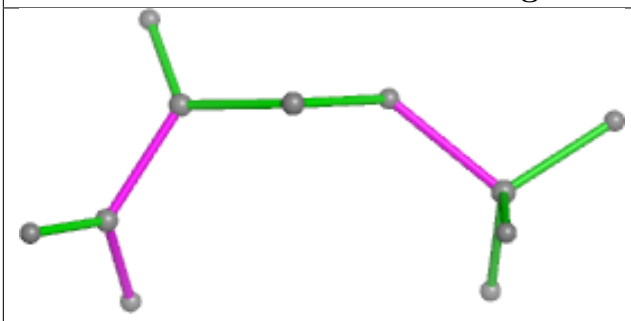


Torsions

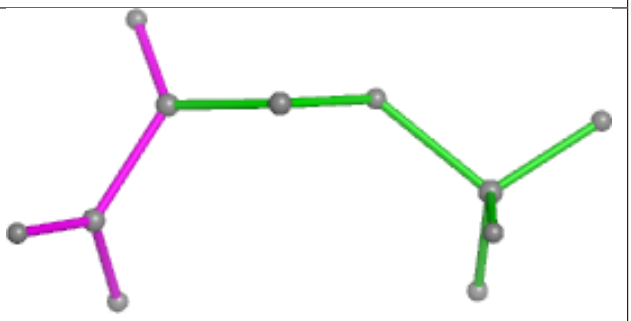


Rings

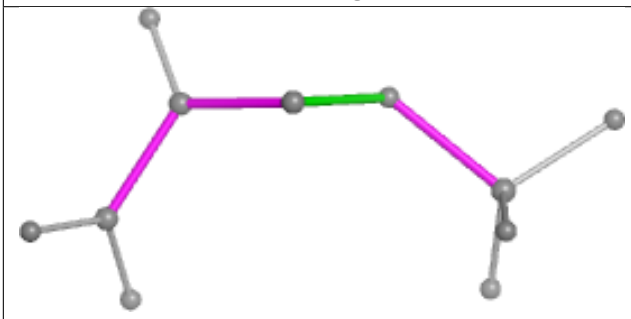
Ligand HPV A 504



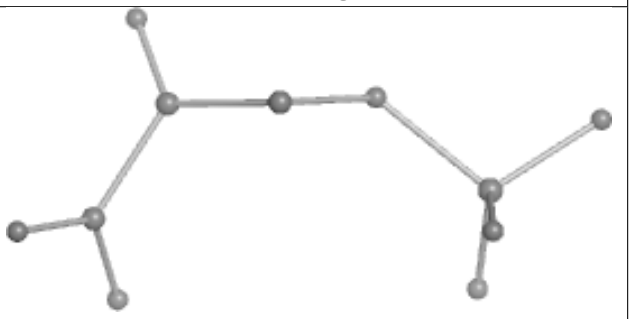
Bond lengths



Bond angles

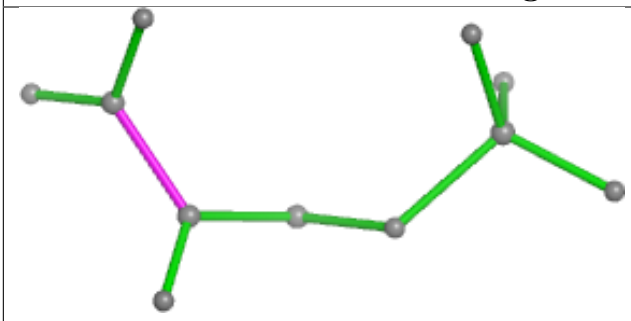


Torsions

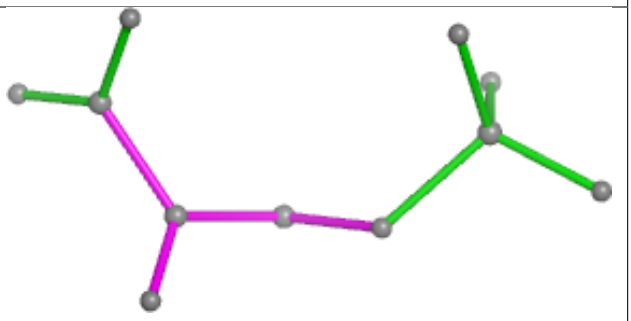


Rings

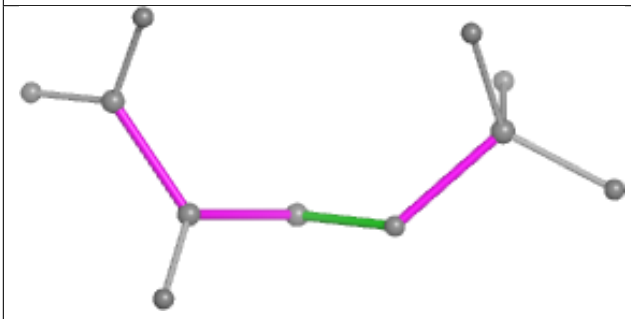
Ligand HPV B 503



Bond lengths



Bond angles



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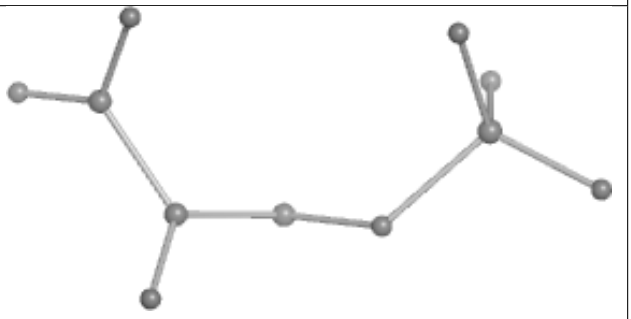
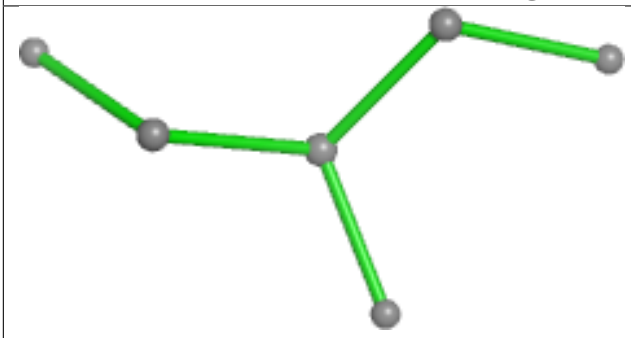


Rings

Ligand GOL C 502



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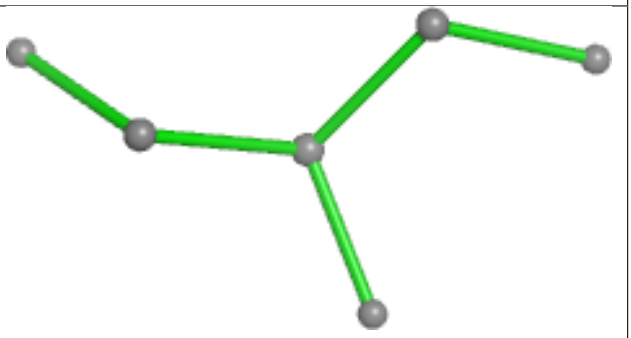
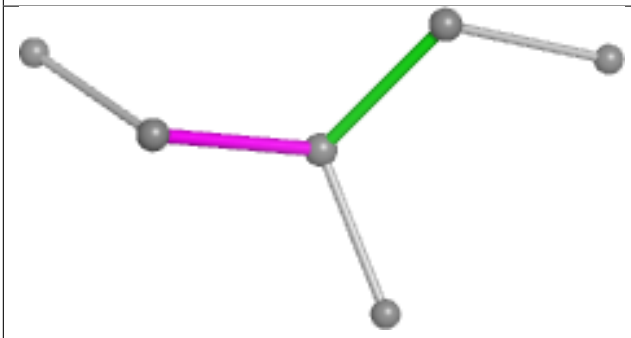


Bond angles



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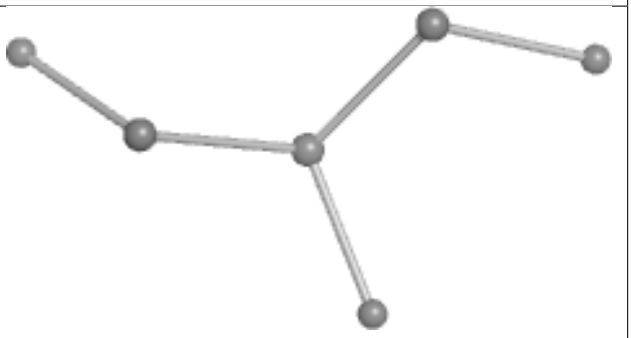
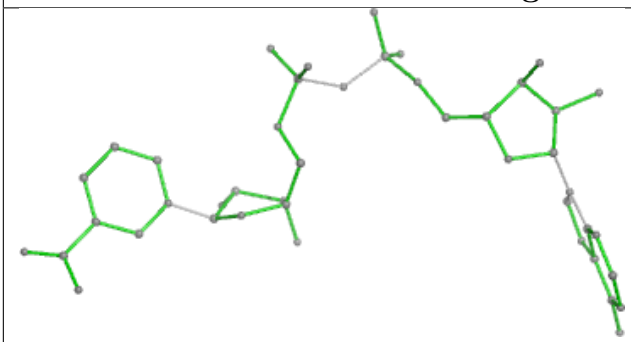
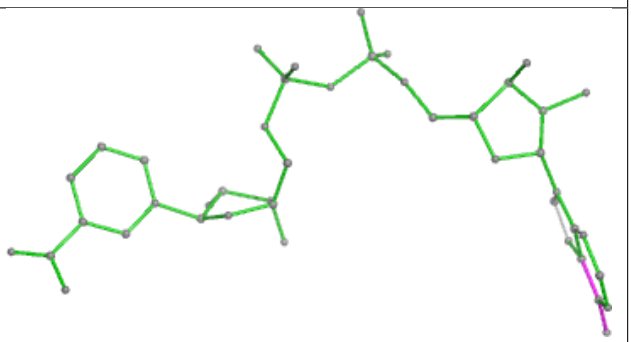


Rings

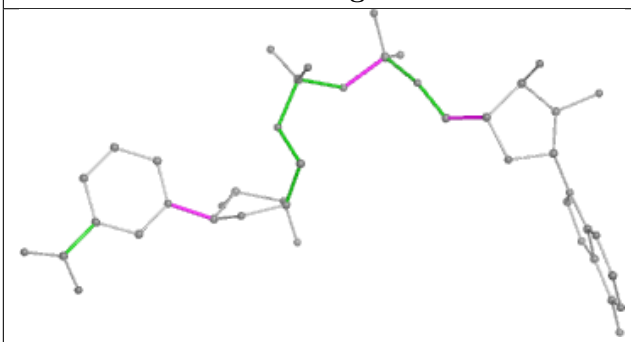
Ligand NAD F 501



Bond lengths



Bond angles



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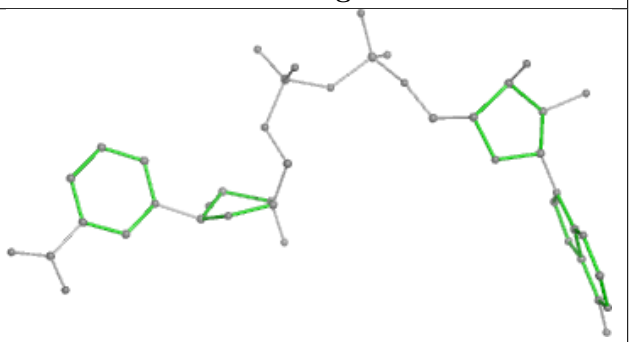
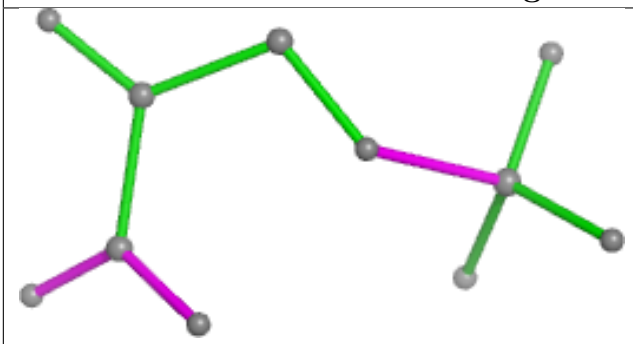


Rings

Ligand HPV F 502



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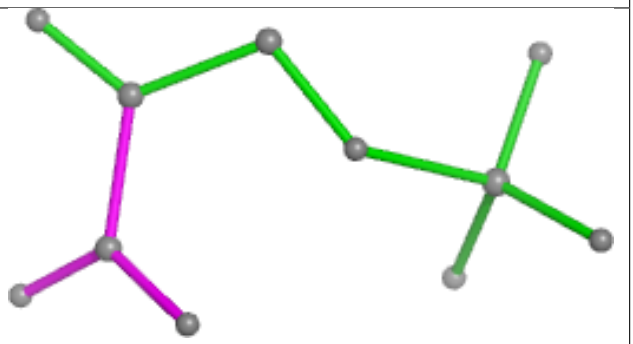
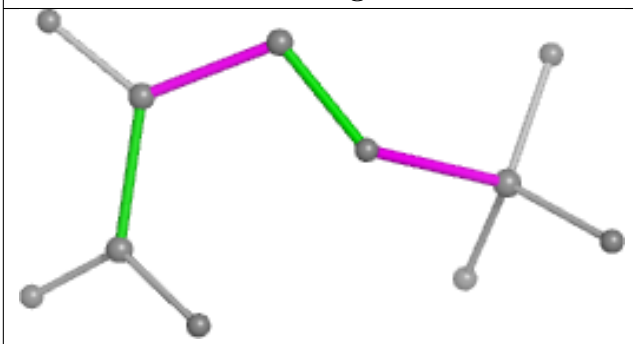


Bond angles



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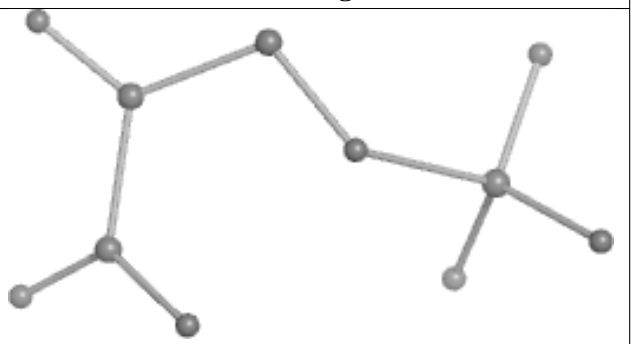
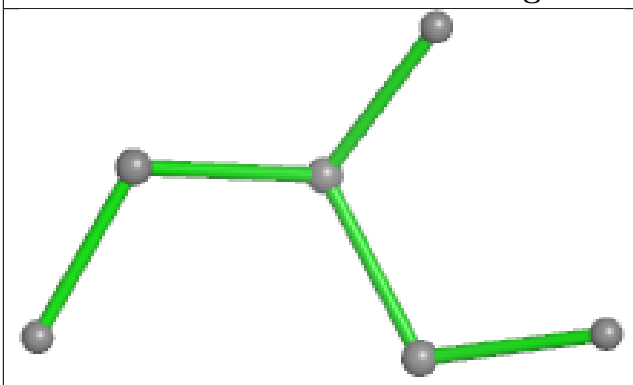
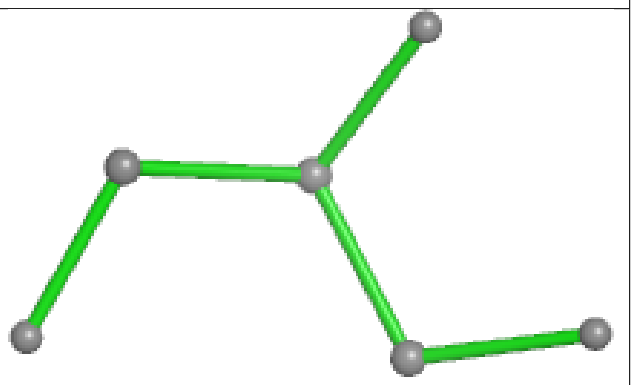


Rings

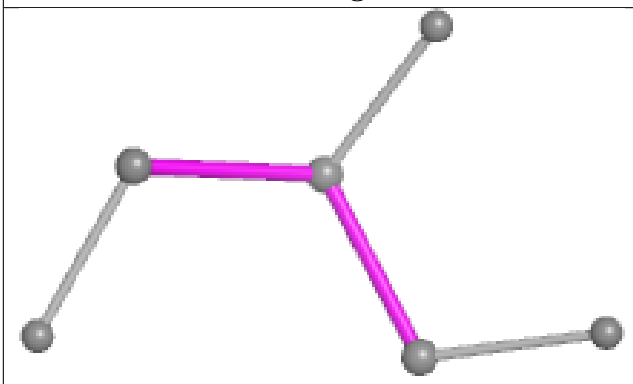
Ligand GOL B 502



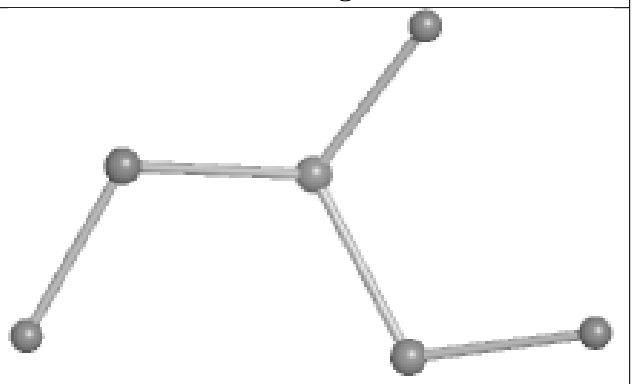
Bond lengths



Bond angles

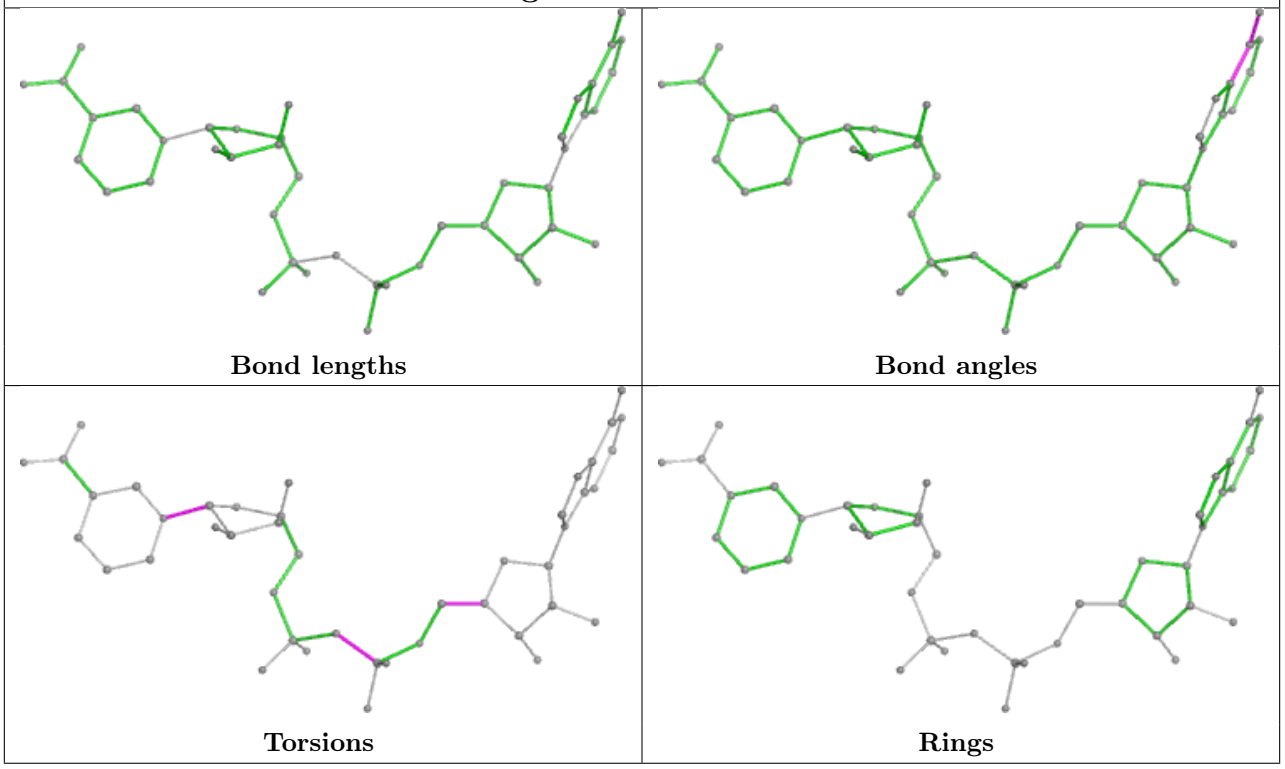


Torsions



Rings

Ligand NAD C 501



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, 95^{th} 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	443/469 (94%)	0.10	8 (1%)	67	62	45, 73, 117, 167	0
1	B	429/469 (91%)	0.35	7 (1%)	70	65	50, 85, 119, 148	0
1	C	431/469 (91%)	0.33	9 (2%)	63	58	48, 81, 119, 145	0
1	F	445/469 (94%)	-0.14	5 (1%)	77	74	45, 66, 105, 174	0
All	All	1748/1876 (93%)	0.16	29 (1%)	69	64	45, 76, 117, 174	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	32	GLN	3.4
1	A	31	THR	3.0
1	B	47	GLN	2.9
1	B	24	GLY	2.9
1	C	447	SER	2.7
1	F	39	PRO	2.7
1	C	58	ASP	2.6
1	F	56	THR	2.6
1	B	58	ASP	2.5
1	C	87	SER	2.5
1	C	392	TYR	2.4
1	A	100	ASP	2.4
1	A	393	ASP	2.4
1	B	93	GLU	2.4
1	A	43	SER	2.4
1	A	56	THR	2.4
1	C	146	ILE	2.4
1	A	38	LEU	2.3
1	C	230	VAL	2.3
1	F	43	SER	2.2
1	B	259	PRO	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	32	GLN	2.2
1	B	465	ILE	2.2
1	A	24	GLY	2.2
1	F	37	THR	2.2
1	C	445	ASN	2.2
1	C	59	MET	2.2
1	F	55	SER	2.1
1	C	33	SER	2.1

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 monosaccharides 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, 95^{th} 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.

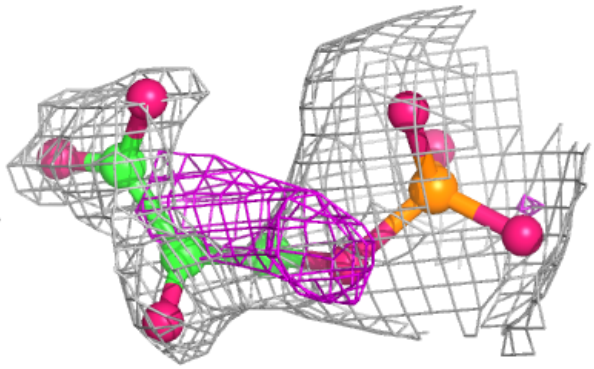
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	HPV	B	503	11/11	0.83	0.12	88,103,112,114	0
4	GOL	B	502	6/6	0.85	0.13	59,75,80,82	0
3	HPV	C	503	11/11	0.86	0.11	88,102,113,122	0
4	GOL	C	502	6/6	0.86	0.15	69,73,76,91	0
3	HPV	F	502	11/11	0.87	0.11	65,78,88,92	0
3	HPV	A	504	11/11	0.87	0.13	75,89,114,114	0
2	NAD	B	501	44/44	0.90	0.09	59,75,84,101	0
2	NAD	C	501	44/44	0.91	0.10	60,71,79,83	0
4	GOL	A	503	6/6	0.91	0.10	53,59,59,63	0
4	GOL	A	501	6/6	0.94	0.12	57,71,74,75	0
2	NAD	A	502	44/44	0.96	0.07	47,58,73,75	0
2	NAD	F	501	44/44	0.97	0.06	36,51,58,70	0

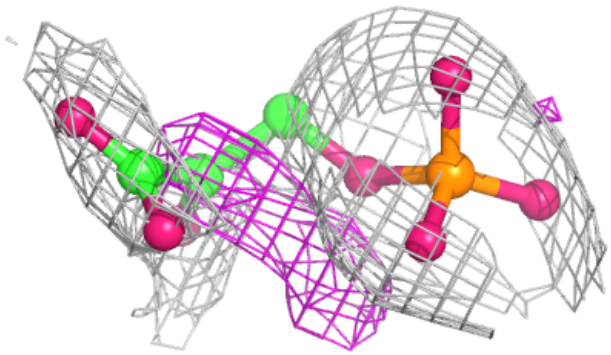
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers

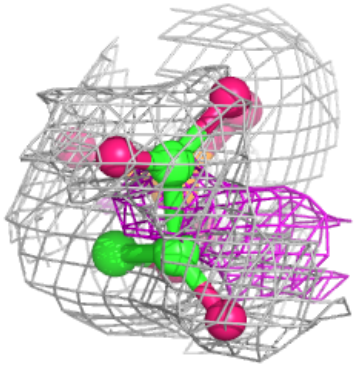
as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around HPV B 503:

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

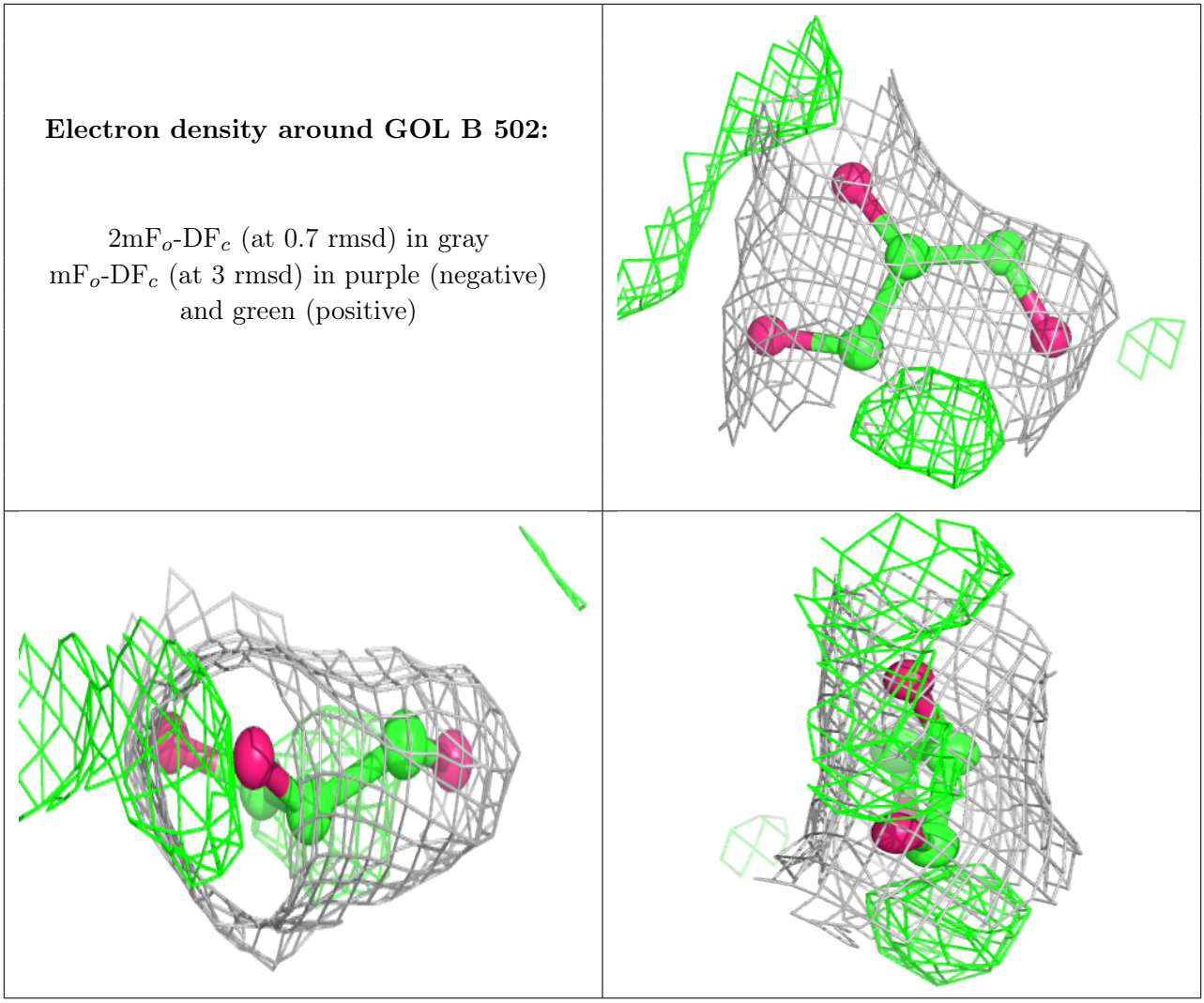






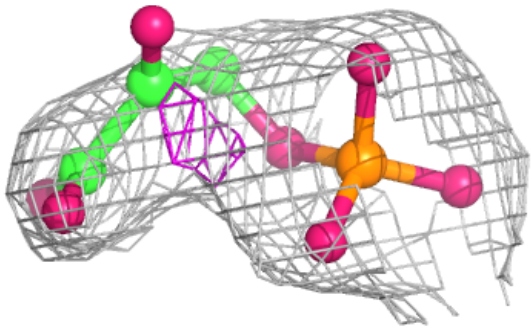
Electron density around GOL B 502:

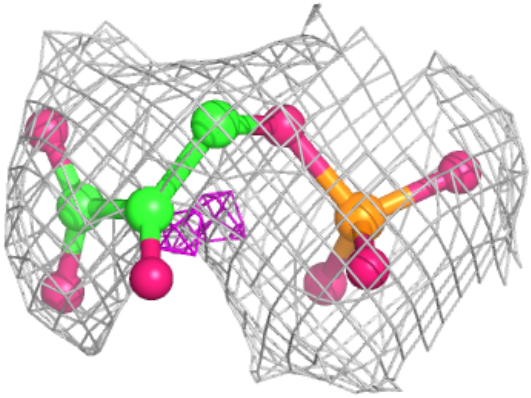
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

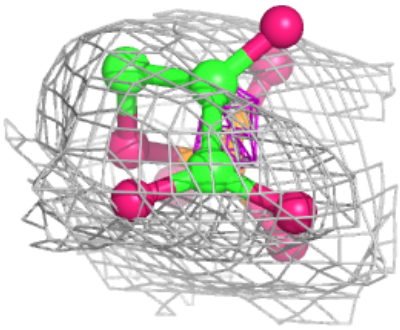


Electron density around HPV C 503:

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

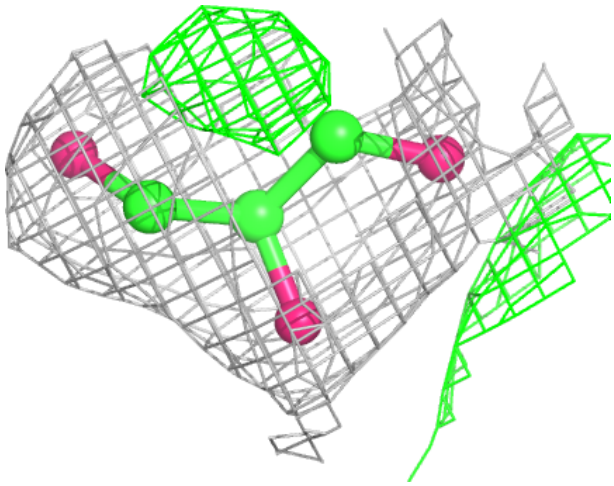


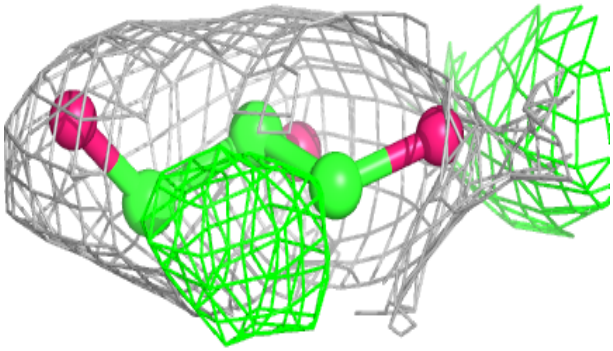


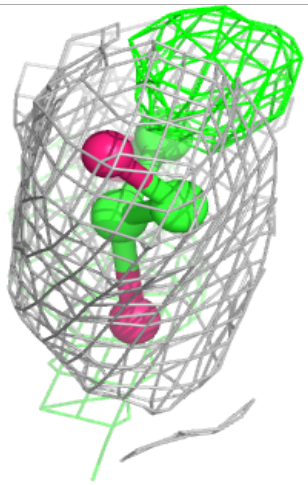


Electron density around GOL C 502:

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

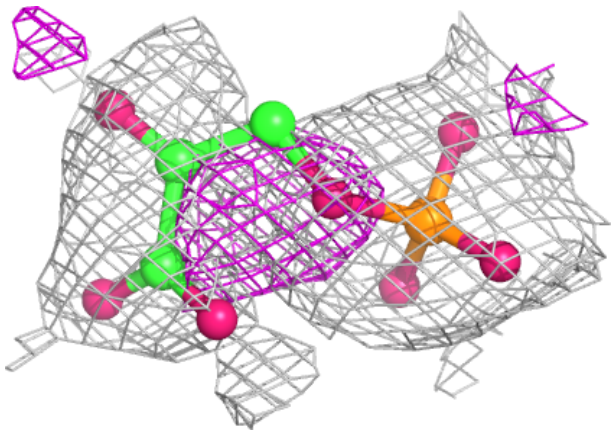


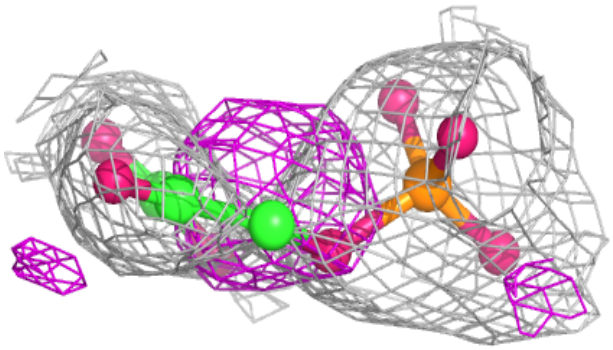


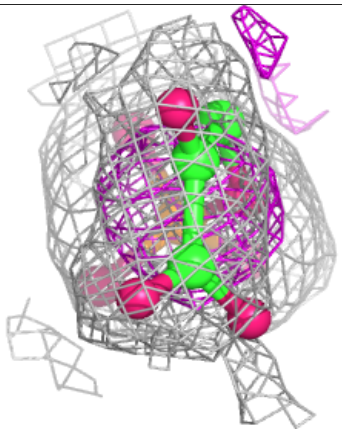


Electron density around HPV F 502:

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

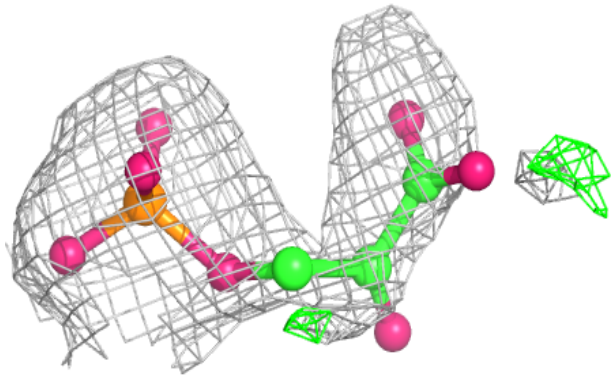


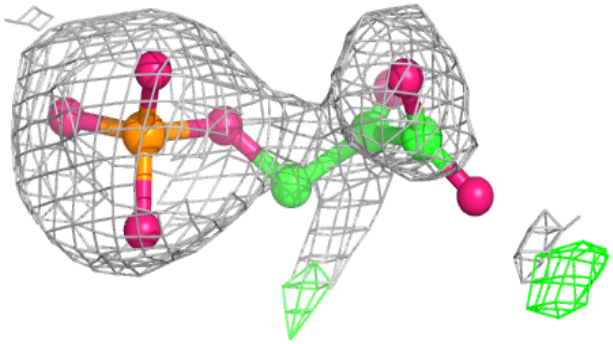


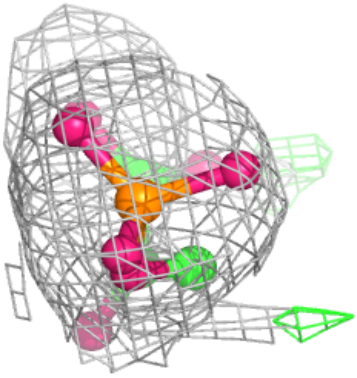


Electron density around HPV A 504:

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

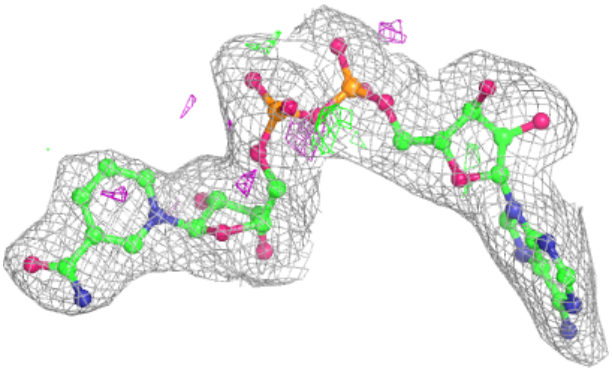


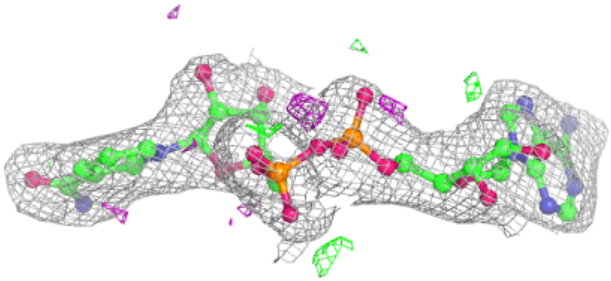


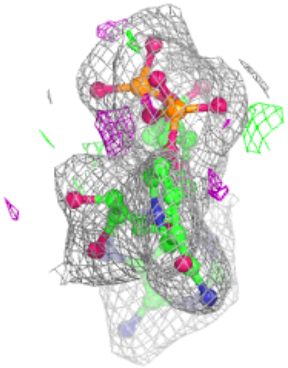


Electron density around NAD B 501:

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

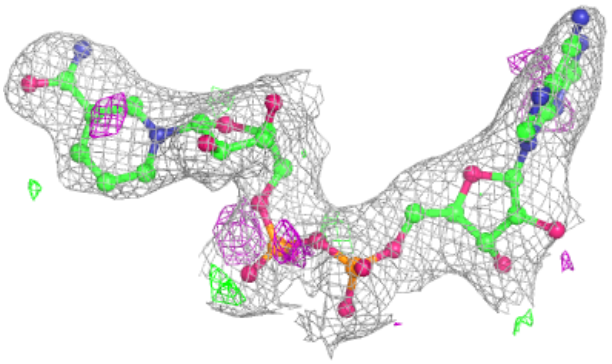


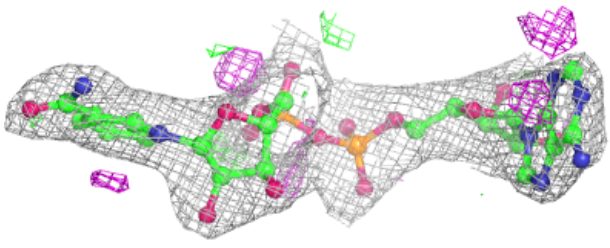


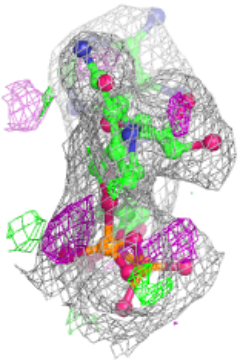


Electron density around NAD C 501:

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

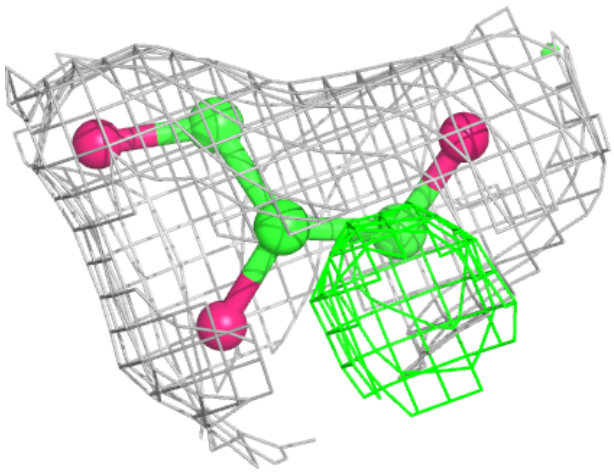


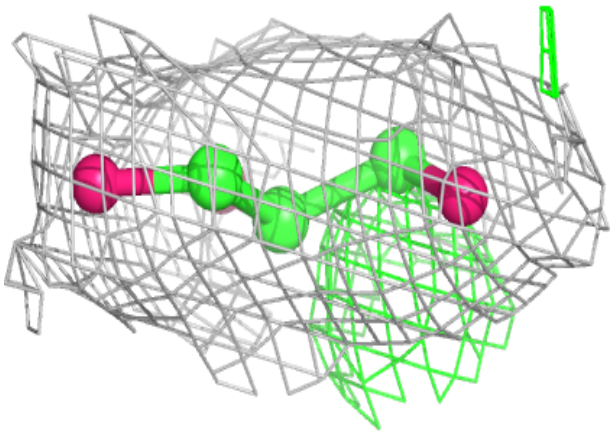


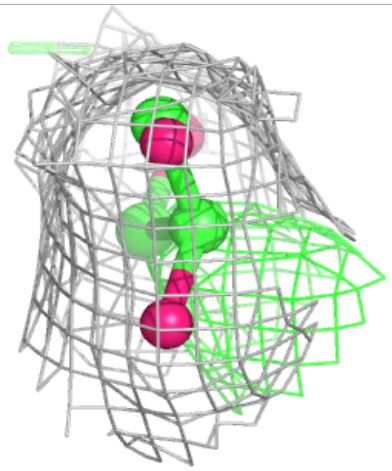


Electron density around GOL A 503:

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

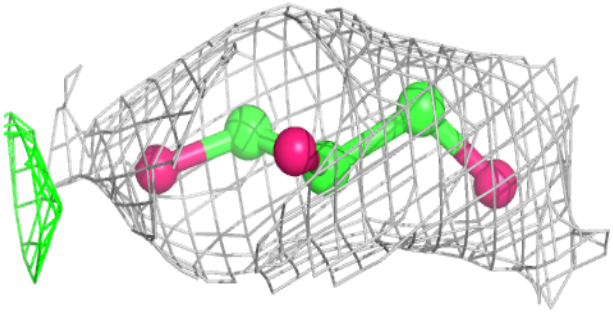


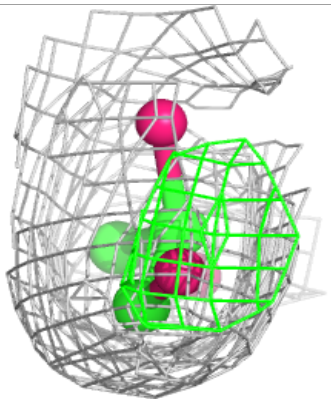




Electron density around GOL A 501:

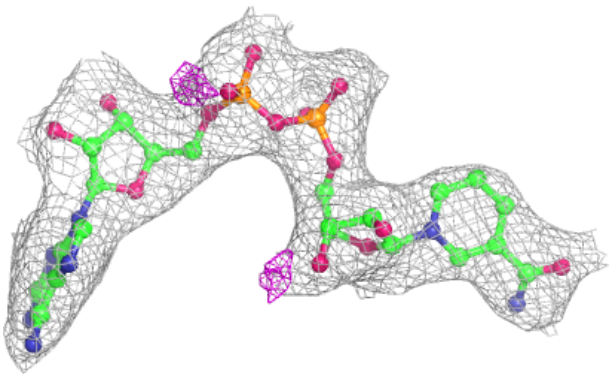
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

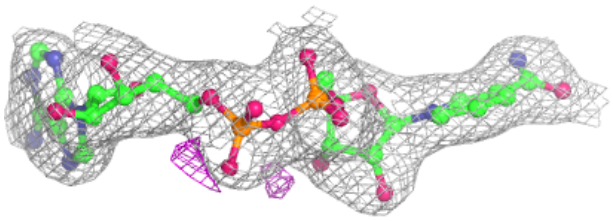


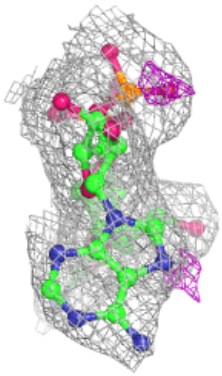


Electron density around NAD A 502:

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

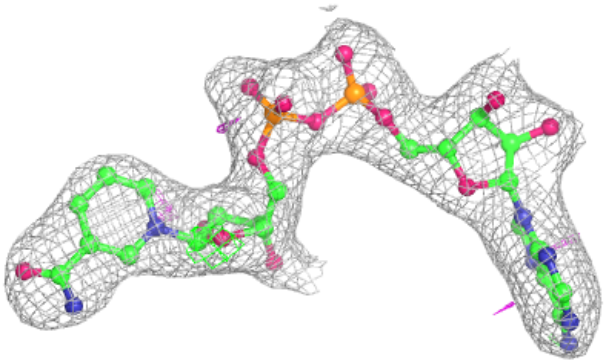


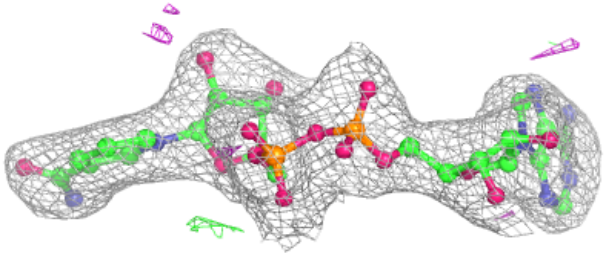


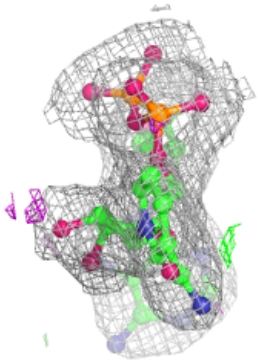


Electron density around NAD F 501:

$2mF_o$-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.