



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 10:39 AM UTC

PDB ID : 3DJJ / pdb_00003djj
Title : Catalytic cycle of human glutathione reductase near 1 Å resolution
Authors : Berkholz, D.S.; Faber, H.R.; Savvides, S.N.; Karplus, P.A.
Deposited on : 2008-06-23
Resolution : 1.10 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 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

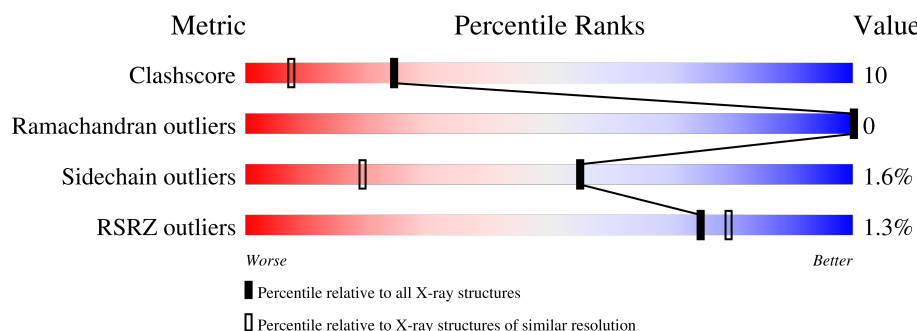
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.10 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	190562	1656 (1.12-1.08)
Ramachandran outliers	187476	1616 (1.12-1.08)
Sidechain outliers	187428	1612 (1.12-1.08)
RSRZ outliers	180081	1625 (1.12-1.08)

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	478	80% 16% ..

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	PO4	A	487	-	-	X	-
6	GOL	A	603	-	-	X	-
6	GOL	A	607	-	-	X	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
6	GOL	A	608	-	-	X	-

2 Entry composition [i](#)

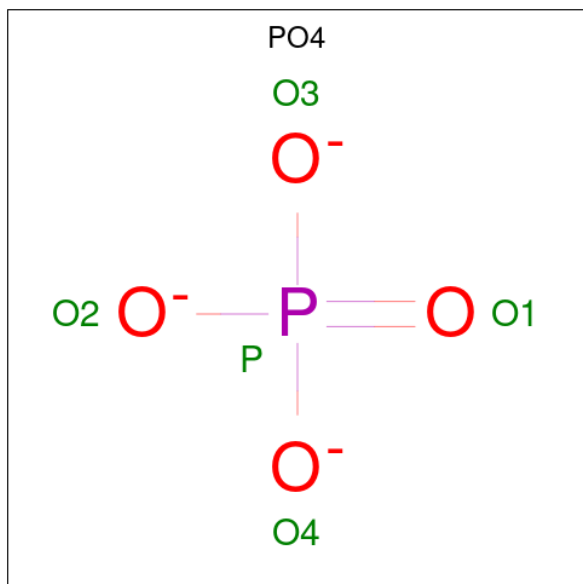
There are 7 unique types of molecules in this entry. The entry contains 4923 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Glutathione reductase, mitochondrial precursor.

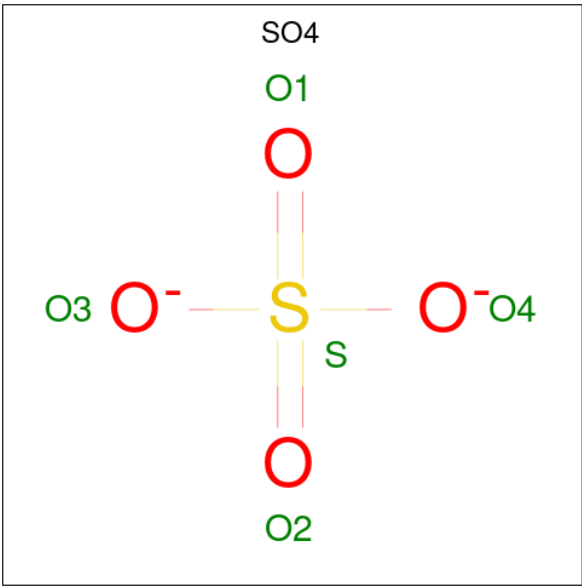
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	462	3842	2444	646	720	32	0	70	0

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



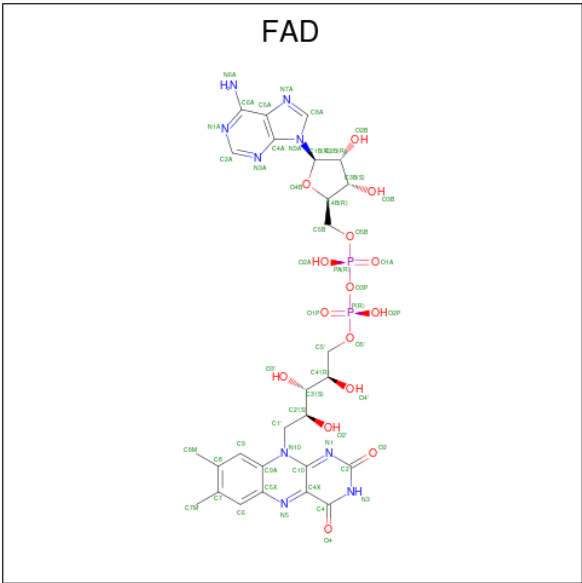
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	P	0	0
			5	4	1		
2	A	1	Total	O	P	0	0
			5	4	1		
2	A	1	Total	O	P	0	0
			5	4	1		
2	A	1	Total	O	P	0	0
			5	4	1		
2	A	1	Total	O	P	0	0
			5	4	1		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



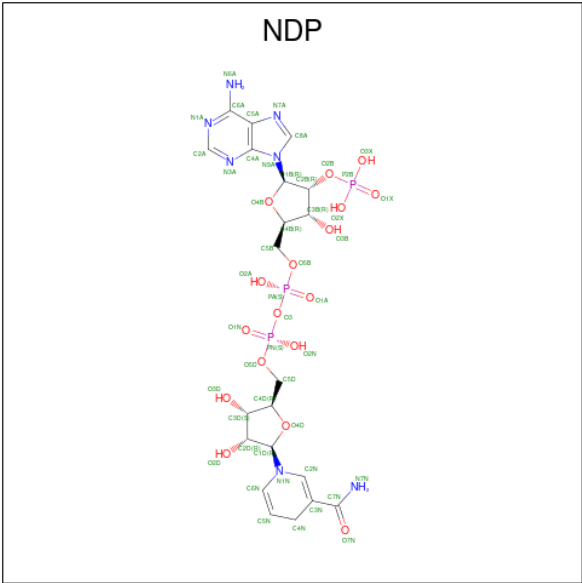
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O 1 1	0	0
3	A	1	Total O 1 1	0	0
3	A	1	Total O 1 1	0	0
3	A	1	Total O 1 1	0	0

- Molecule 4 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: C₂₇H₃₃N₉O₁₅P₂).



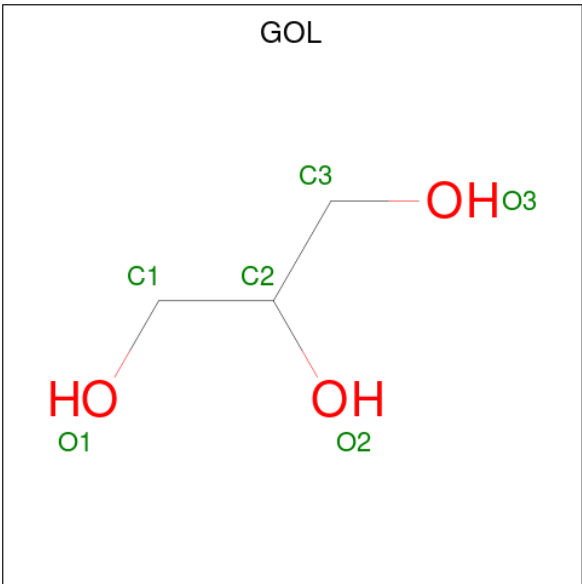
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 5 is NADPH DIHYDRO-NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (CCD ID: NDP) (formula: $C_{21}H_{30}N_7O_{17}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	P	0	0
			48	21	7	17	3		

- Molecule 6 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

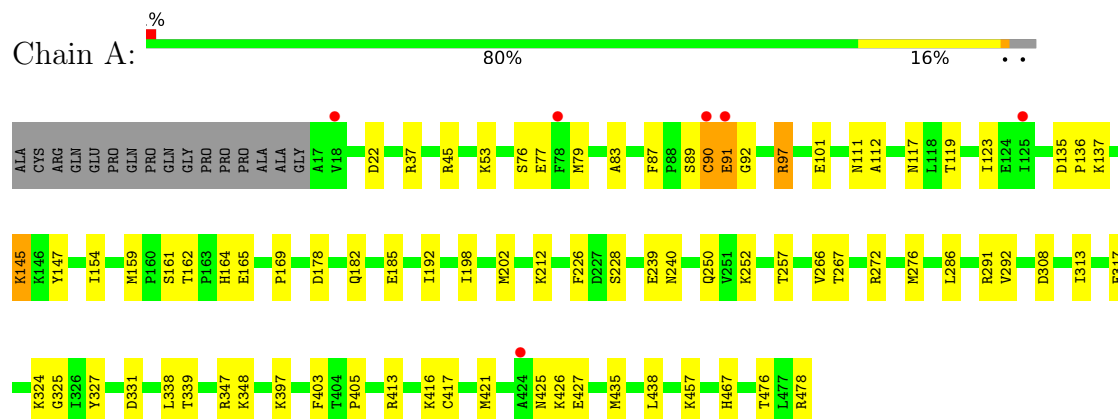
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	878	Total	O	0	35
			903	903		

3 Residue-property plots

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

- Molecule 1: Glutathione reductase, mitochondrial precursor



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	119.67Å 62.36Å 84.16Å 90.00° 122.29° 90.00°	Depositor
Resolution (Å)	100.00 – 1.10 100.00 – 1.10	Depositor EDS
% Data completeness (in resolution range)	90.9 (100.00-1.10) 39.1 (100.00-1.10)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.80 (at 1.10Å)	Xtriage
Refinement program	SHELX, SHELXL-97	Depositor
R, R_{free}	0.113 , 0.147 0.121 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	12.0	Xtriage
Anisotropy	0.090	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 66.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	4923	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.82% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: FAD, NDP, GOL, SO4, PO4

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	0.95	2/4126 (0.0%)	1.29	23/5570 (0.4%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	37	ARG	CD-NE	-6.31	1.37	1.46
1	A	435	MET	SD-CE	-5.49	1.65	1.79

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	37	ARG	CD-NE-CZ	17.02	148.23	124.40
1	A	117	ASN	OD1-CG-ND2	-7.24	115.36	122.60
1	A	308	ASP	CA-CB-CG	7.09	119.69	112.60
1	A	397	LYS	CB-CG-CD	6.75	126.82	111.30
1	A	425	ASN	OD1-CG-ND2	-6.75	115.85	122.60
1	A	331	ASP	CA-CB-CG	6.53	119.13	112.60
1	A	178	ASP	CA-CB-CG	6.45	119.05	112.60
1	A	239	GLU	CB-CG-CD	6.36	123.42	112.60
1	A	427	GLU	CB-CG-CD	5.89	122.61	112.60
1	A	111	ASN	CA-CB-CG	5.84	118.44	112.60
1	A	135	ASP	CA-CB-CG	5.75	118.35	112.60
1	A	291	ARG	CB-CG-CD	5.73	124.48	111.30
1	A	347	ARG	NE-CZ-NH1	-5.51	115.99	121.50
1	A	416	LYS	CA-CB-CG	5.46	125.03	114.10
1	A	291	ARG	NE-CZ-NH2	5.45	124.11	119.20
1	A	164	HIS	CA-CB-CG	-5.37	108.43	113.80
1	A	467	HIS	CA-CB-CG	5.24	119.04	113.80
1	A	425	ASN	CB-CG-ND2	5.23	124.24	116.40
1	A	435	MET	CG-SD-CE	-5.23	89.40	100.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	97	ARG	NE-CZ-NH2	5.11	123.80	119.20
1	A	317	GLU	CB-CG-CD	5.05	121.19	112.60
1	A	136	PRO	CA-C-N	-5.00	114.65	121.61
1	A	136	PRO	C-N-CA	-5.00	114.65	121.61

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	3842	0	3880	65	0
2	A	25	0	0	4	0
3	A	4	0	0	0	0
4	A	53	0	31	0	0
5	A	48	0	26	2	0
6	A	48	0	60	23	0
7	A	903	0	0	45	0
All	All	4923	0	3997	80	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 (80) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:91[A]:GLU:N	7:A:1807[A]:HOH:O	1.76	1.19
2:A:486:PO4:P	7:A:1622:HOH:O	2.38	0.81
1:A:257:THR:HA	7:A:1366:HOH:O	1.84	0.78
6:A:601:GOL:H32	7:A:1300:HOH:O	1.86	0.75
1:A:476[B]:THR:HG21	7:A:1303:HOH:O	1.86	0.74
1:A:338:LEU:HD21	7:A:1300:HOH:O	1.90	0.70
6:A:607:GOL:H31	7:A:1316:HOH:O	1.94	0.68
1:A:276[B]:MET:HE1	7:A:3149:HOH:O	1.94	0.67
1:A:266[B]:VAL:HG22	1:A:276[B]:MET:SD	2.35	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:A:604:GOL:H31	7:A:1695:HOH:O	1.96	0.65
1:A:161[A]:SER:OG	1:A:292[A]:VAL:HG13	1.97	0.65
1:A:137:LYS:HE3	7:A:1523:HOH:O	1.96	0.65
1:A:324[B]:LYS:HD2	7:A:1210:HOH:O	1.95	0.64
5:A:480:NDP:H2D	7:A:1300:HOH:O	1.97	0.64
1:A:324[B]:LYS:HE3	7:A:1258:HOH:O	1.97	0.63
1:A:90[A]:CYS:C	7:A:1807[A]:HOH:O	2.24	0.62
1:A:324[B]:LYS:HG3	7:A:3788:HOH:O	2.00	0.61
1:A:250:GLN:HG2	7:A:1257:HOH:O	1.99	0.61
1:A:92[B]:GLY:HA3	7:A:1613:HOH:O	2.01	0.60
6:A:607:GOL:H11	7:A:9004[B]:HOH:O	2.01	0.59
1:A:162[B]:THR:HG21	7:A:1797:HOH:O	2.05	0.57
1:A:438:LEU:HD21	7:A:1794:HOH:O	2.04	0.57
1:A:421:MET:SD	1:A:457[B]:LYS:HD3	2.45	0.56
1:A:101[A]:GLU:HG3	7:A:2792[A]:HOH:O	2.05	0.56
1:A:405:PRO:HG3	1:A:417[B]:CYS:SG	2.46	0.56
6:A:601:GOL:HO2	6:A:608:GOL:HO1	1.52	0.56
1:A:212[A]:LYS:HD3	7:A:3423:HOH:O	2.05	0.56
1:A:403:PHE:O	1:A:417[B]:CYS:HB2	2.06	0.55
1:A:97:ARG:HG3	7:A:1328:HOH:O	2.06	0.55
1:A:145[A]:LYS:HE2	1:A:147:TYR:OH	2.07	0.54
1:A:226:PHE:HA	6:A:608:GOL:H32	1.90	0.54
1:A:267:THR:HG1	6:A:606:GOL:HO2	1.56	0.52
1:A:212[A]:LYS:HD2	7:A:1164:HOH:O	2.08	0.52
1:A:325:GLY:HA2	7:A:1203:HOH:O	2.09	0.52
2:A:487:PO4:O4	6:A:607:GOL:O1	2.28	0.51
1:A:119:THR:HA	7:A:1576:HOH:O	2.11	0.51
1:A:240:ASN:HB3	7:A:1447:HOH:O	2.11	0.50
1:A:267:THR:OG1	6:A:606:GOL:O2	2.27	0.50
1:A:112:ALA:HA	6:A:607:GOL:H12	1.94	0.50
1:A:226:PHE:CE1	6:A:608:GOL:H11	2.46	0.50
1:A:324[A]:LYS:HG3	6:A:603:GOL:H31	1.93	0.49
6:A:606:GOL:H12	7:A:1571:HOH:O	2.13	0.49
1:A:198:ILE:CG2	1:A:202[B]:MET:HE2	2.43	0.48
1:A:76[A]:SER:OG	7:A:2730[A]:HOH:O	2.19	0.48
1:A:165[B]:GLU:CD	1:A:169:PRO:HA	2.39	0.48
2:A:487:PO4:P	6:A:607:GOL:O1	2.72	0.47
6:A:608:GOL:H31	7:A:1159:HOH:O	2.15	0.46
1:A:202[B]:MET:HE3	1:A:286:LEU:HD11	1.97	0.46
1:A:272[A]:ARG:HD3	7:A:1245:HOH:O	2.14	0.46
1:A:77[A]:GLU:HG3	7:A:3430[A]:HOH:O	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:426:LYS:NZ	7:A:1470:HOH:O	2.49	0.46
1:A:83:ALA:HA	1:A:87:PHE:O	2.16	0.45
1:A:77[B]:GLU:OE2	1:A:413:ARG:NH2	2.49	0.45
1:A:123:ILE:N	7:A:1576:HOH:O	2.50	0.45
1:A:154:ILE:HG21	1:A:313[B]:ILE:CD1	2.47	0.45
1:A:22[B]:ASP:CG	1:A:45:ARG:HE	2.25	0.44
1:A:182[A]:GLN:NE2	7:A:3192:HOH:O	2.49	0.44
1:A:97:ARG:NH1	1:A:185:GLU:OE2	2.50	0.44
1:A:327:TYR:OH	6:A:603:GOL:H2	2.18	0.44
6:A:608:GOL:H31	7:A:2763[A]:HOH:O	2.17	0.44
1:A:348[B]:LYS:NZ	7:A:1586:HOH:O	2.50	0.43
1:A:478:ARG:HG2	7:A:1358:HOH:O	2.19	0.43
1:A:22[B]:ASP:OD2	1:A:45:ARG:NH2	2.50	0.43
1:A:53:LYS:NZ	7:A:1596:HOH:O	2.50	0.43
1:A:89[B]:SER:O	1:A:90[B]:CYS:HB2	2.19	0.42
1:A:276[B]:MET:HE2	7:A:1685:HOH:O	2.17	0.42
1:A:228:SER:OG	6:A:602:GOL:H12	2.18	0.42
5:A:480:NDP:H2A	7:A:1257:HOH:O	2.18	0.42
1:A:90[A]:CYS:HA	7:A:1807[A]:HOH:O	2.20	0.42
1:A:226:PHE:CZ	6:A:608:GOL:H11	2.54	0.42
1:A:198:ILE:HG23	1:A:202[B]:MET:HE2	2.02	0.42
1:A:79[B]:MET:HG2	1:A:90[B]:CYS:O	2.20	0.41
1:A:325:GLY:H	6:A:603:GOL:H2	1.84	0.41
1:A:325:GLY:H	6:A:603:GOL:C2	2.34	0.41
1:A:272[B]:ARG:NH2	7:A:1718:HOH:O	2.50	0.41
1:A:438:LEU:HD11	7:A:1794:HOH:O	2.20	0.41
2:A:487:PO4:O2	6:A:607:GOL:O1	2.39	0.40
1:A:226:PHE:CD1	6:A:608:GOL:H11	2.57	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	530/478 (111%)	511 (96%)	19 (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	447/393 (114%)	436 (98%)	11 (2%)	42	8

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	90[A]	CYS
1	A	90[B]	CYS
1	A	91[A]	GLU
1	A	91[B]	GLU
1	A	145[A]	LYS
1	A	145[B]	LYS
1	A	159[A]	MET
1	A	159[B]	MET
1	A	252[A]	LYS
1	A	252[B]	LYS
1	A	339	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	233	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 19 ligands modelled in this entry, 4 are modelled with single atom - leaving 15 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
5	NDP	A	480	-	51,52,52	1.25	7 (13%)	71,80,80	1.04	5 (7%)
2	PO4	A	481	-	4,4,4	0.83	0	6,6,6	0.66	0
6	GOL	A	604	-	5,5,5	1.35	1 (20%)	5,5,5	1.26	0
6	GOL	A	605	-	5,5,5	1.01	0	5,5,5	1.80	1 (20%)
6	GOL	A	608	-	5,5,5	0.66	0	5,5,5	0.59	0
2	PO4	A	485	-	4,4,4	1.20	1 (25%)	6,6,6	0.76	0
6	GOL	A	601	-	5,5,5	0.69	0	5,5,5	1.51	1 (20%)
6	GOL	A	602	-	5,5,5	1.03	0	5,5,5	1.20	0
4	FAD	A	479	-	58,58,58	1.29	8 (13%)	85,89,89	1.09	5 (5%)
6	GOL	A	603	-	5,5,5	0.83	0	5,5,5	5.74	1 (20%)
2	PO4	A	487	-	4,4,4	0.81	0	6,6,6	1.26	1 (16%)
6	GOL	A	607	-	5,5,5	0.91	0	5,5,5	0.67	0
2	PO4	A	482	-	4,4,4	0.83	0	6,6,6	0.70	0
6	GOL	A	606	-	5,5,5	0.50	0	5,5,5	7.33	1 (20%)
2	PO4	A	486	-	4,4,4	1.21	0	6,6,6	1.58	2 (33%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns.

'-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NDP	A	480	-	-	4/34/77/77	0/5/5/5
6	GOL	A	604	-	-	1/4/4/4	-
6	GOL	A	605	-	-	1/4/4/4	-
6	GOL	A	608	-	-	4/4/4/4	-
6	GOL	A	601	-	-	1/4/4/4	-
6	GOL	A	602	-	-	4/4/4/4	-
6	GOL	A	603	-	-	2/4/4/4	-
4	FAD	A	479	-	-	3/34/50/50	0/6/6/6
6	GOL	A	607	-	-	2/4/4/4	-
6	GOL	A	606	-	-	4/4/4/4	-

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	479	FAD	C1'-C2'	3.35	1.57	1.52
5	A	480	NDP	C2D-C3D	-3.21	1.44	1.53
4	A	479	FAD	C4X-N5	3.16	1.37	1.30
5	A	480	NDP	PN-O3	3.07	1.62	1.59
5	A	480	NDP	P2B-O2B	2.90	1.64	1.59
4	A	479	FAD	O4B-C1B	2.89	1.48	1.42
4	A	479	FAD	O2-C2	2.63	1.29	1.24
4	A	479	FAD	C6-C5X	2.56	1.43	1.40
4	A	479	FAD	C8M-C8	2.31	1.55	1.51
2	A	485	PO4	P-O1	2.28	1.55	1.50
5	A	480	NDP	PN-O2N	-2.23	1.45	1.55
4	A	479	FAD	P-O3P	2.22	1.61	1.59
4	A	479	FAD	C10-N1	2.12	1.37	1.33
6	A	604	GOL	C3-C2	2.10	1.59	1.51
5	A	480	NDP	PA-O2A	-2.08	1.45	1.55
5	A	480	NDP	O4D-C4D	2.06	1.49	1.45
5	A	480	NDP	C4A-N9A	2.05	1.41	1.37

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	606	GOL	C3-C2-C1	-16.38	51.74	111.80
6	A	603	GOL	C3-C2-C1	-12.64	65.44	111.80
6	A	605	GOL	C3-C2-C1	-3.82	97.80	111.80
6	A	601	GOL	C3-C2-C1	-2.99	100.81	111.80
5	A	480	NDP	C1D-N1N-C2N	-2.96	116.27	121.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	480	NDP	P2B-O2B-C2B	-2.38	117.08	123.43
5	A	480	NDP	C2A-N1A-C6A	2.37	122.63	118.73
5	A	480	NDP	C3D-C2D-C1D	2.32	105.86	101.46
4	A	479	FAD	O5B-PA-O1A	2.27	117.94	108.94
4	A	479	FAD	C5A-C4A-N9A	2.25	108.27	105.81
4	A	479	FAD	C4X-C10-N10	2.14	119.55	116.48
5	A	480	NDP	O2N-PN-O1N	2.10	122.23	112.44
2	A	487	PO4	O2-P-O1	-2.10	103.54	110.95
4	A	479	FAD	C4-N3-C2	-2.03	122.04	125.64
4	A	479	FAD	O4B-C1B-N9A	-2.01	104.22	108.09
2	A	486	PO4	O2-P-O1	-2.01	103.84	110.95
2	A	486	PO4	O4-P-O3	2.00	114.14	107.91

There are no chirality outliers.

All (26) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	479	FAD	PA-O3P-P-O5'
6	A	602	GOL	O1-C1-C2-C3
6	A	606	GOL	O1-C1-C2-C3
6	A	606	GOL	C1-C2-C3-O3
6	A	608	GOL	O1-C1-C2-C3
6	A	608	GOL	C1-C2-C3-O3
5	A	480	NDP	O4D-C1D-N1N-C2N
6	A	603	GOL	O2-C2-C3-O3
6	A	603	GOL	C1-C2-C3-O3
6	A	607	GOL	C1-C2-C3-O3
6	A	602	GOL	O1-C1-C2-O2
6	A	608	GOL	O1-C1-C2-O2
6	A	608	GOL	O2-C2-C3-O3
5	A	480	NDP	C3B-C4B-C5B-O5B
5	A	480	NDP	O4B-C4B-C5B-O5B
6	A	602	GOL	C1-C2-C3-O3
4	A	479	FAD	P-O3P-PA-O1A
6	A	606	GOL	O1-C1-C2-O2
6	A	606	GOL	O2-C2-C3-O3
6	A	607	GOL	O2-C2-C3-O3
6	A	602	GOL	O2-C2-C3-O3
6	A	601	GOL	O1-C1-C2-C3
4	A	479	FAD	P-O3P-PA-O2A
6	A	604	GOL	C1-C2-C3-O3
5	A	480	NDP	C3D-C4D-C5D-O5D

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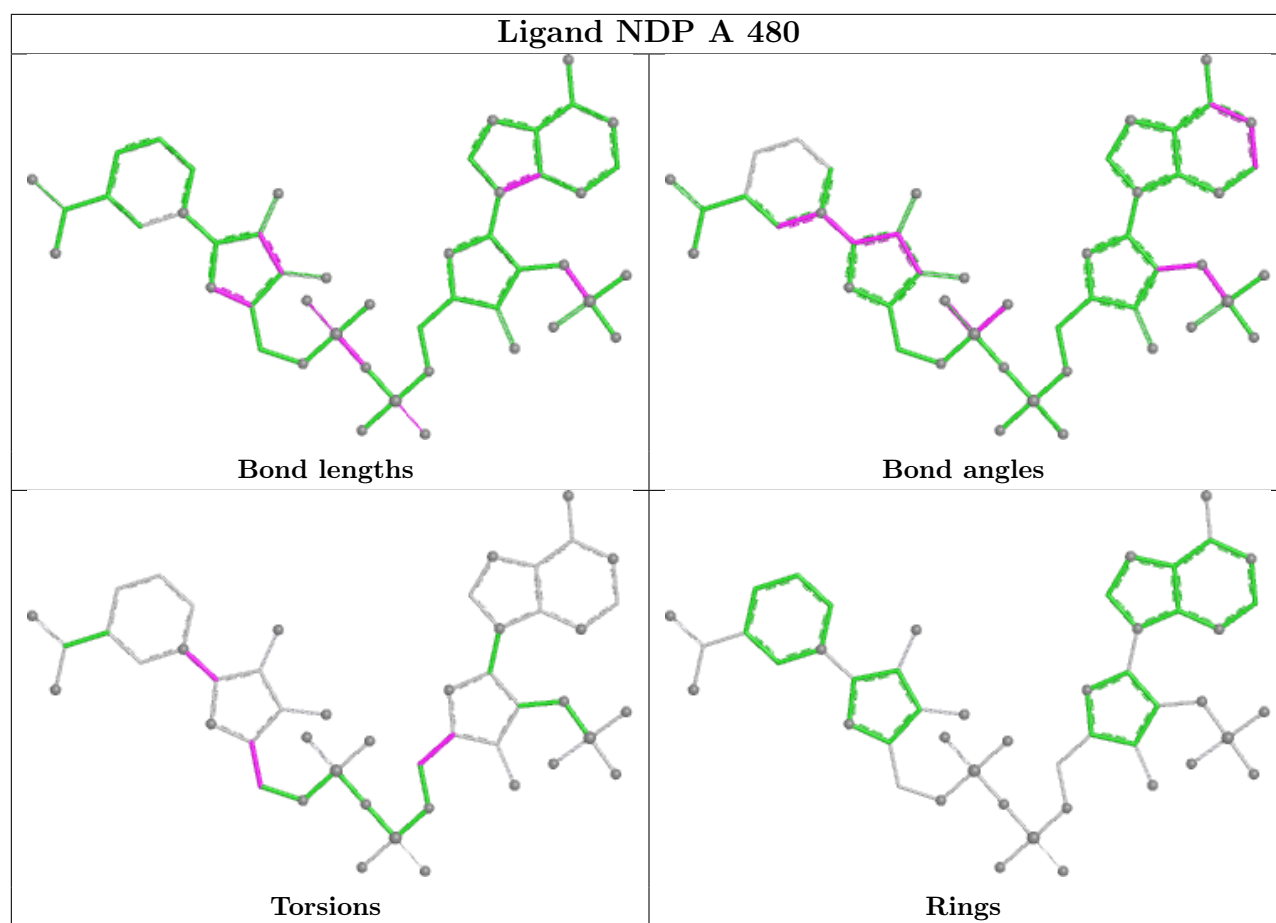
Mol	Chain	Res	Type	Atoms
6	A	605	GOL	O1-C1-C2-C3

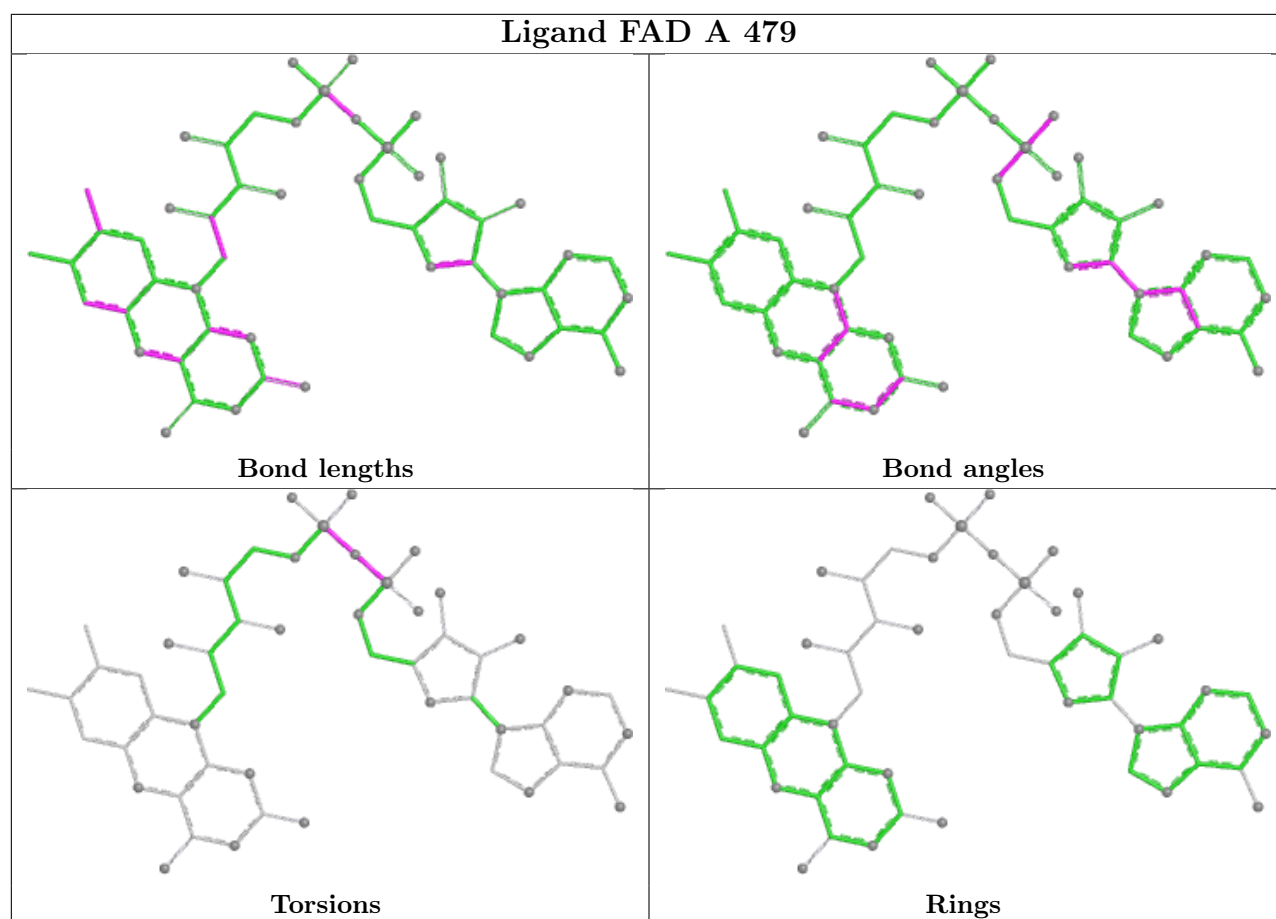
There are no ring outliers.

10 monomers are involved in 26 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	480	NDP	2	0
6	A	604	GOL	1	0
6	A	608	GOL	7	0
6	A	601	GOL	2	0
6	A	602	GOL	1	0
6	A	603	GOL	4	0
2	A	487	PO4	3	0
6	A	607	GOL	6	0
6	A	606	GOL	3	0
2	A	486	PO4	1	0

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





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	462/478 (96%)	0.25	6 (1%) 75 79	6, 14, 28, 58	70 (15%)

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	90[A]	CYS	4.1
1	A	125	ILE	2.8
1	A	78[A]	PHE	2.5
1	A	91[A]	GLU	2.2
1	A	18[A]	VAL	2.1
1	A	424	ALA	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(Å ²)	Q<0.9
2	PO4	A	487	5/5	0.83	0.13	29,45,63,67	0

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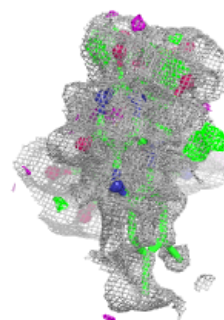
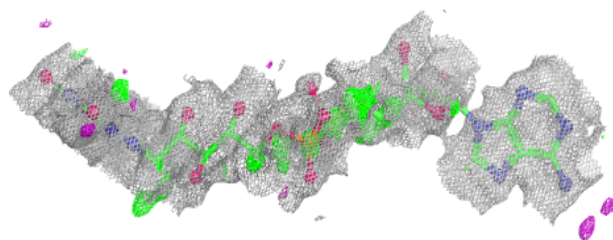
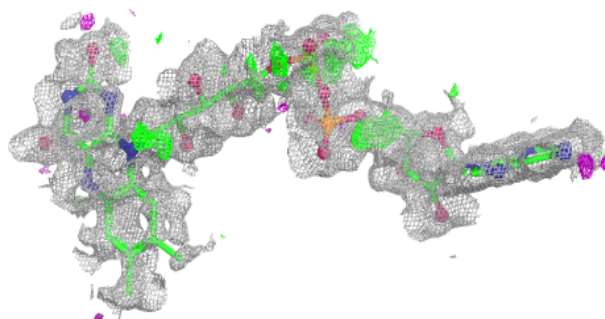
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	PO4	A	482	5/5	0.84	0.11	42,57,69,73	0
6	GOL	A	603	6/6	0.84	0.10	37,46,50,51	0
6	GOL	A	602	6/6	0.86	0.13	20,29,60,87	0
2	PO4	A	486	5/5	0.87	0.10	52,54,61,63	5
6	GOL	A	606	6/6	0.89	0.13	64,77,80,80	0
2	PO4	A	485	5/5	0.90	0.10	29,31,39,43	5
3	SO4	A	3585	1/5	0.91	0.13	54,54,54,54	0
6	GOL	A	607	6/6	0.92	0.09	29,49,59,87	0
3	SO4	A	3586	1/5	0.93	0.10	44,44,44,44	0
3	SO4	A	3587	1/5	0.93	0.11	58,58,58,58	0
2	PO4	A	481	5/5	0.93	0.10	21,29,38,44	0
6	GOL	A	608	6/6	0.94	0.10	30,40,52,57	0
6	GOL	A	601	6/6	0.95	0.07	18,20,22,23	0
6	GOL	A	605	6/6	0.96	0.08	28,29,30,35	0
3	SO4	A	3584	1/5	0.97	0.06	63,63,63,63	0
6	GOL	A	604	6/6	0.97	0.09	12,24,32,65	0
4	FAD	A	479	53/53	0.99	0.05	9,11,14,16	0
5	NDP	A	480	48/48	0.99	0.04	9,12,14,17	0

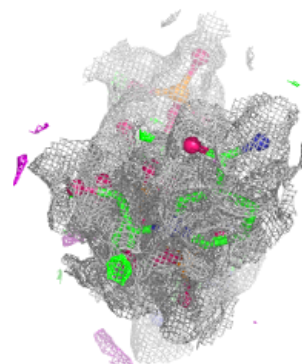
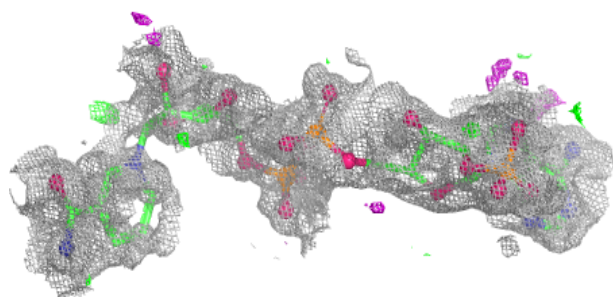
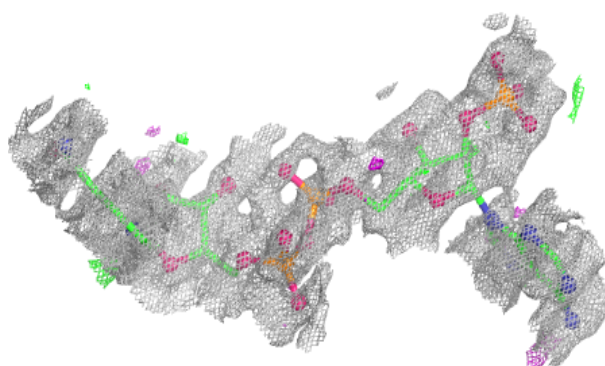
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around FAD A 479:

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

**Electron density around NDP A 480:**

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



6.5 Other polymers [i](#)

There are no such residues in this entry.