



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 23, 2026 – 08:32 AM EDT

PDB ID : 1FNC / pdb_00001fnc
Title : REFINED CRYSTAL STRUCTURE OF SPINACH FERREDOXIN REDUCTASE AT 1.7 ANGSTROMS RESOLUTION: OXIDIZED, REDUCED, AND 2'-PHOSPHO-5'-AMP BOUND STATES
Authors : Bruns, C.M.; Karplus, P.A.
Deposited on : 1995-01-05
Resolution : 2.00 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

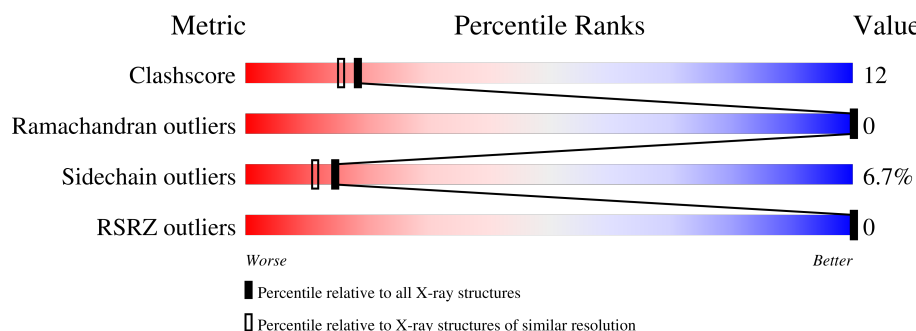
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	314	

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	SO4	A	317	-	-	X	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 2680 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 FERREDONIN-NADP+ REDUCTASE.

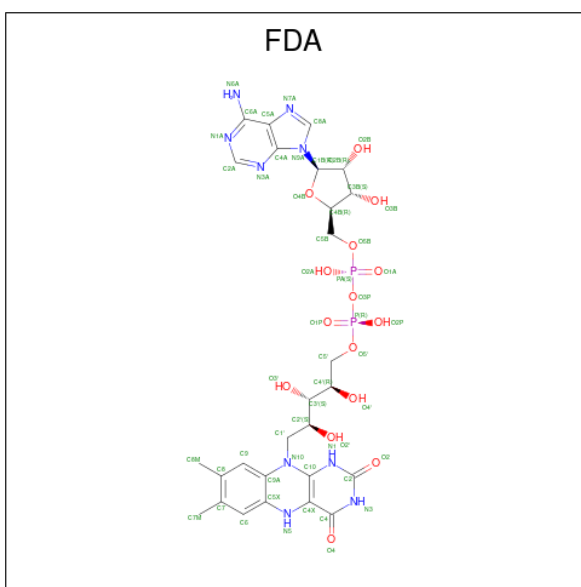
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	296	2352	1506	388	440	18	0	0	0

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



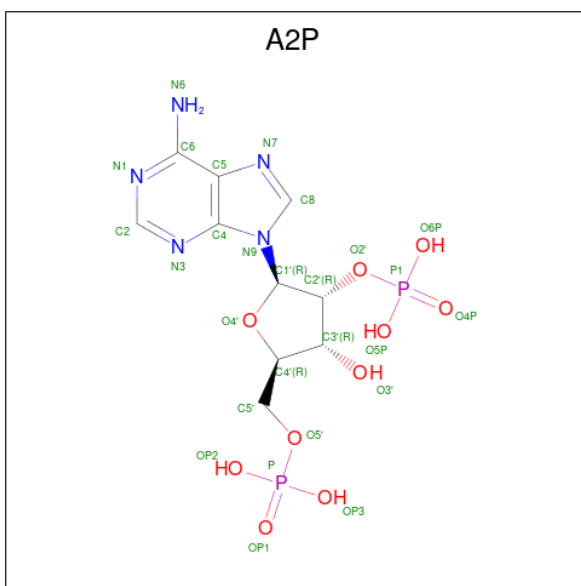
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	S		
2	A	1	5	4	1	0	0

- Molecule 3 is DIHYDROFLAVINE-ADENINE DINUCLEOTIDE (CCD ID: FDA) (formula: C₂₇H₃₅N₉O₁₅P₂).



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

- Molecule 4 is ADENOSINE-2'-5'-DIPHOSPHATE (CCD ID: A2P) (formula: $\text{C}_{10}\text{H}_{15}\text{N}_5\text{O}_{10}\text{P}_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total 54	C 20	N 10	O 20	P 4	22	1

- Molecule 5 is water.

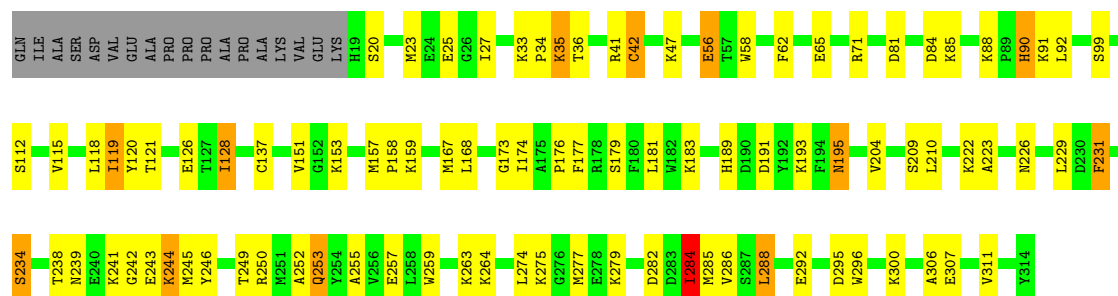
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	216	Total 216	O 216	0	0

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: FERREDOXIN-NADP+ REDUCTASE

Chain A: 



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	90.70Å 57.70Å 68.10Å 90.00° 100.00° 90.00°	Depositor
Resolution (Å)	(Not available) – 2.00 67.06 – 2.00	Depositor EDS
% Data completeness (in resolution range)	(Not available) ((Not available)-2.00) 90.2 (67.06-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$	-	Xtriage
Refinement program	TNT 5A	Depositor
R, R_{free}	0.149 , (Not available) 0.140 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	16.7	Xtriage
Anisotropy	0.412	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 131.0	EDS
L-test for twinning ¹	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	2680	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.27% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Theoretical 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: FDA, SO4, A2P

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	1.39	3/2406 (0.1%)	1.62	19/3236 (0.6%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	42	CYS	N-CA	-5.60	1.39	1.46
1	A	295	ASP	CA-C	5.53	1.59	1.52
1	A	231	PHE	C-O	-5.12	1.17	1.23

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	177	PHE	CA-CB-CG	-6.91	106.89	113.80
1	A	311	VAL	N-CA-C	6.74	118.50	108.46
1	A	222	LYS	CB-CA-C	-6.66	99.53	110.85
1	A	189	HIS	CA-CB-CG	-6.16	107.64	113.80
1	A	253	GLN	N-CA-CB	-5.87	100.12	110.39
1	A	35	LYS	CA-C-N	-5.72	117.05	122.37
1	A	35	LYS	C-N-CA	-5.72	117.05	122.37
1	A	284	ILE	CB-CA-C	-5.67	103.08	112.26
1	A	115	VAL	N-CA-C	5.62	116.22	108.12
1	A	229	LEU	N-CA-CB	-5.59	101.42	110.81
1	A	234	SER	N-CA-C	5.56	118.05	111.33
1	A	36	THR	N-CA-CB	-5.35	104.66	110.71
1	A	253	GLN	CB-CG-CD	-5.34	103.52	112.60
1	A	151	VAL	CB-CA-C	-5.31	102.19	110.69
1	A	81	ASP	CA-CB-CG	5.30	117.90	112.60
1	A	168	LEU	CA-C-N	-5.29	117.53	121.61
1	A	168	LEU	C-N-CA	-5.29	117.53	121.61
1	A	90	HIS	CA-CB-CG	-5.13	108.67	113.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	173	GLY	N-CA-C	-5.05	109.04	115.31

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	2352	0	2342	58	0
2	A	5	0	0	2	0
3	A	53	0	33	0	0
4	A	54	0	22	0	0
5	A	216	0	0	6	0
All	All	2680	0	2397	58	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 12.

All (58) 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:255:ALA:HB1	1:A:288:LEU:HD21	1.48	0.92
1:A:157:MET:HE1	1:A:183:LYS:HD3	1.54	0.87
1:A:285:MET:HA	1:A:285:MET:HE3	1.56	0.86
1:A:33:LYS:HB3	1:A:34:PRO:HD2	1.65	0.78
1:A:35:LYS:HG3	5:A:480:HOH:O	1.84	0.77
1:A:157:MET:HE1	1:A:183:LYS:CD	2.20	0.71
1:A:274:LEU:O	1:A:277:MET:HG2	1.94	0.66
1:A:257:GLU:HB2	5:A:719:HOH:O	1.95	0.65
1:A:121:THR:HA	1:A:126:GLU:O	1.97	0.65
1:A:239:ASN:O	1:A:242:GLY:N	2.30	0.65
1:A:157:MET:HE2	5:A:427:HOH:O	1.96	0.65
1:A:285:MET:HE3	1:A:285:MET:CA	2.26	0.63
1:A:167:MET:HE1	1:A:181:LEU:HD21	1.81	0.62
1:A:238:THR:HA	1:A:243:GLU:O	2.00	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:90:HIS:HA	2:A:317:SO4:O1	2.00	0.61
1:A:20:SER:O	1:A:159:LYS:NZ	2.28	0.60
1:A:239:ASN:ND2	1:A:243:GLU:OE1	2.31	0.59
1:A:47:LYS:HD2	1:A:56:GLU:HB3	1.85	0.59
1:A:239:ASN:OD1	1:A:243:GLU:N	2.29	0.58
1:A:58:TRP:CD1	1:A:137:CYS:HG	2.23	0.57
1:A:246:TYR:O	1:A:249:THR:HB	2.04	0.57
1:A:259:TRP:NE1	1:A:263:LYS:HE3	2.19	0.57
1:A:84:ASP:N	1:A:88:LYS:O	2.32	0.57
1:A:27:ILE:HG22	1:A:71:ARG:HD2	1.87	0.55
1:A:90:HIS:ND1	2:A:317:SO4:O1	2.29	0.54
1:A:296:TRP:CZ2	1:A:300:LYS:HG3	2.43	0.54
1:A:252:ALA:HB2	1:A:284:ILE:HD12	1.89	0.54
1:A:284:ILE:HG22	1:A:288:LEU:HD12	1.88	0.53
1:A:274:LEU:N	1:A:274:LEU:HD12	2.23	0.53
1:A:27:ILE:HD11	1:A:153:LYS:CG	2.39	0.53
1:A:195:ASN:HB2	5:A:522:HOH:O	2.08	0.53
1:A:118:LEU:C	1:A:119:ILE:HG12	2.36	0.51
1:A:120:TYR:O	1:A:128:ILE:N	2.26	0.51
1:A:234:SER:O	1:A:244:LYS:HG2	2.11	0.50
1:A:99:SER:HB3	1:A:112:SER:HB2	1.93	0.50
1:A:250:ARG:O	1:A:253:GLN:HB3	2.11	0.50
1:A:223:ALA:HB1	1:A:226:ASN:HB2	1.96	0.47
1:A:157:MET:HE1	1:A:183:LYS:CE	2.44	0.47
1:A:264:LYS:NZ	5:A:741:HOH:O	2.39	0.47
1:A:241:LYS:HG2	1:A:241:LYS:O	2.15	0.46
1:A:245:MET:SD	1:A:249:THR:HG22	2.56	0.45
1:A:282:ASP:O	1:A:286:VAL:HG23	2.18	0.44
1:A:204:VAL:HG21	1:A:210:LEU:HD23	1.98	0.44
1:A:241:LYS:HB3	1:A:243:GLU:HG3	1.99	0.44
1:A:246:TYR:HD1	1:A:246:TYR:HA	1.51	0.43
1:A:274:LEU:HD12	1:A:274:LEU:H	1.84	0.42
1:A:306:ALA:O	1:A:307:GLU:HB2	2.18	0.42
1:A:210:LEU:HD22	1:A:231:PHE:CD1	2.55	0.42
1:A:23:MET:HE3	1:A:158:PRO:O	2.19	0.42
1:A:25:GLU:OE2	1:A:159:LYS:HE2	2.20	0.42
1:A:33:LYS:HB3	1:A:34:PRO:CD	2.44	0.42
1:A:176:PRO:O	1:A:179:SER:HB3	2.20	0.42
1:A:65:GLU:HA	5:A:516:HOH:O	2.19	0.42
1:A:92:LEU:C	1:A:92:LEU:HD12	2.45	0.42
1:A:274:LEU:H	1:A:274:LEU:CD1	2.32	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:259:TRP:CD1	1:A:263:LYS:HE3	2.55	0.41
1:A:42:CYS:HA	1:A:62:PHE:HD1	1.85	0.41
1:A:27:ILE:CG2	1:A:71:ARG:HD2	2.50	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	294/314 (94%)	284 (97%)	10 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	253/267 (95%)	236 (93%)	17 (7%)	15	11

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	41	ARG
1	A	56	GLU
1	A	85	LYS

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Mol	Chain	Res	Type
1	A	91	LYS
1	A	119	ILE
1	A	128	ILE
1	A	174	ILE
1	A	191	ASP
1	A	193	LYS
1	A	195	ASN
1	A	209	SER
1	A	244	LYS
1	A	275	LYS
1	A	279	LYS
1	A	284	ILE
1	A	288	LEU
1	A	292	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	30	ASN
1	A	162	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

4 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
4	A2P	A	316[A]	-	29,29,29	1.41	4 (13%)	44,45,45	1.19	4 (9%)
4	A2P	A	316[B]	-	29,29,29	2.22	6 (20%)	44,45,45	1.26	5 (11%)
3	FDA	A	315	-	57,58,58	1.98	13 (22%)	78,89,89	1.39	13 (16%)
2	SO4	A	317	-	4,4,4	0.80	0	6,6,6	0.24	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
4	A2P	A	316[A]	-	-	9/15/31/31	0/3/3/3
4	A2P	A	316[B]	-	-	3/15/31/31	0/3/3/3
3	FDA	A	315	-	-	2/34/50/50	0/6/6/6

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	316[B]	A2P	P1-O2'	9.60	1.76	1.59
3	A	315	FDA	PA-O3P	6.09	1.66	1.59
3	A	315	FDA	P-O3P	5.90	1.65	1.59
3	A	315	FDA	C1'-C2'	5.22	1.59	1.52
3	A	315	FDA	C10-N1	-4.66	1.29	1.37
4	A	316[A]	A2P	P-OP2	-3.72	1.41	1.54
3	A	315	FDA	C5'-C4'	3.08	1.56	1.51
3	A	315	FDA	O4-C4	-3.04	1.17	1.23
3	A	315	FDA	C9A-N10	-2.97	1.36	1.41
4	A	316[B]	A2P	O5'-C5'	2.71	1.54	1.44
3	A	315	FDA	C8A-N9A	2.61	1.42	1.37
3	A	315	FDA	O4B-C1B	2.53	1.47	1.42
3	A	315	FDA	C2-N3	2.45	1.41	1.37
4	A	316[A]	A2P	O4'-C1'	2.43	1.47	1.42
3	A	315	FDA	C4X-N5	-2.38	1.31	1.35
3	A	315	FDA	C5X-C9A	2.35	1.43	1.40
4	A	316[B]	A2P	C5-N7	-2.33	1.34	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	316[B]	A2P	O4'-C1'	2.30	1.47	1.42
3	A	315	FDA	C4-N3	-2.26	1.34	1.38
4	A	316[A]	A2P	P1-O2'	2.12	1.63	1.59
4	A	316[B]	A2P	P-OP2	-2.08	1.47	1.54
4	A	316[B]	A2P	P-OP1	-2.06	1.44	1.50
4	A	316[A]	A2P	P-OP3	-2.05	1.47	1.54

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	315	FDA	O2A-PA-O1A	2.96	126.20	112.44
3	A	315	FDA	C9A-C5X-N5	2.87	122.87	119.37
3	A	315	FDA	C2A-N1A-C6A	2.67	123.12	118.73
3	A	315	FDA	C4B-O4B-C1B	-2.59	103.74	109.47
4	A	316[A]	A2P	C2-N1-C6	2.54	122.90	118.73
4	A	316[B]	A2P	C3'-C2'-C1'	-2.52	97.98	102.81
3	A	315	FDA	C5X-N5-C4X	-2.49	115.34	121.08
4	A	316[B]	A2P	O2'-P1-O4P	-2.41	100.74	109.33
4	A	316[B]	A2P	OP3-P-O5'	2.40	112.92	106.67
3	A	315	FDA	N3A-C2A-N1A	-2.38	124.97	128.58
3	A	315	FDA	O2P-P-O1P	2.35	123.38	112.44
3	A	315	FDA	C4A-C5A-N7A	2.33	113.24	110.58
3	A	315	FDA	O4-C4-C4X	-2.32	121.66	127.26
4	A	316[A]	A2P	P1-O2'-C2'	-2.31	117.27	123.43
3	A	315	FDA	C4A-N9A-C8A	2.29	108.14	105.74
4	A	316[B]	A2P	C2-N1-C6	2.28	122.47	118.73
4	A	316[A]	A2P	O5'-P-OP1	2.24	112.49	106.44
4	A	316[B]	A2P	O4'-C1'-C2'	-2.15	102.88	106.59
3	A	315	FDA	C5A-C4A-N9A	-2.12	103.50	105.81
4	A	316[A]	A2P	OP2-P-O5'	-2.07	101.28	106.67
3	A	315	FDA	N9A-C8A-N7A	-2.06	111.02	113.94
3	A	315	FDA	C4A-N9A-C1B	-2.05	121.83	126.63

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	316[A]	A2P	C5'-O5'-P-OP3
4	A	316[A]	A2P	C5'-O5'-P-OP2
4	A	316[A]	A2P	C3'-C4'-C5'-O5'
4	A	316[A]	A2P	O4'-C4'-C5'-O5'
4	A	316[A]	A2P	C5'-O5'-P-OP1

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Mol	Chain	Res	Type	Atoms
4	A	316[A]	A2P	C2'-C1'-N9-C8
4	A	316[B]	A2P	C2'-C1'-N9-C8
4	A	316[A]	A2P	C2'-C1'-N9-C4
4	A	316[B]	A2P	C2'-C1'-N9-C4
4	A	316[A]	A2P	O4'-C1'-N9-C8
3	A	315	FDA	PA-O3P-P-O1P
3	A	315	FDA	PA-O3P-P-O2P
4	A	316[B]	A2P	O4'-C1'-N9-C8
4	A	316[A]	A2P	C2'-O2'-P1-O4P

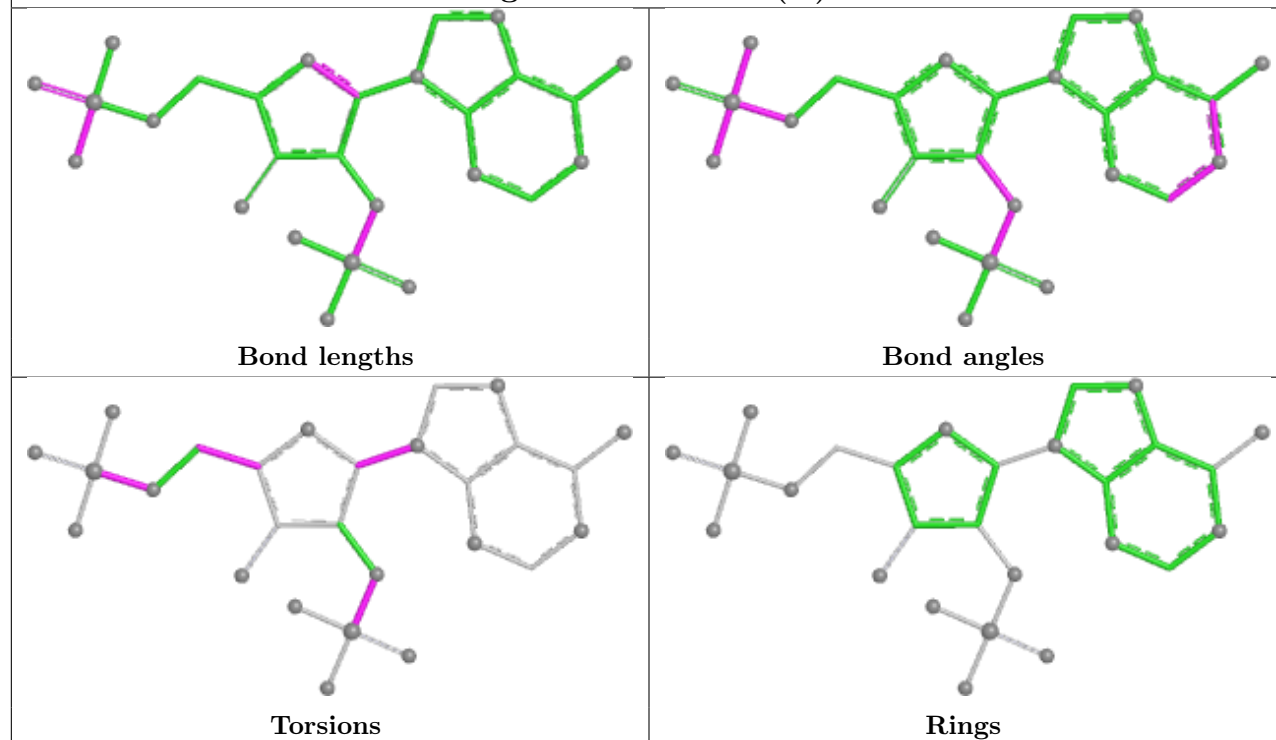
There are no ring outliers.

1 monomer is involved in 2 short contacts:

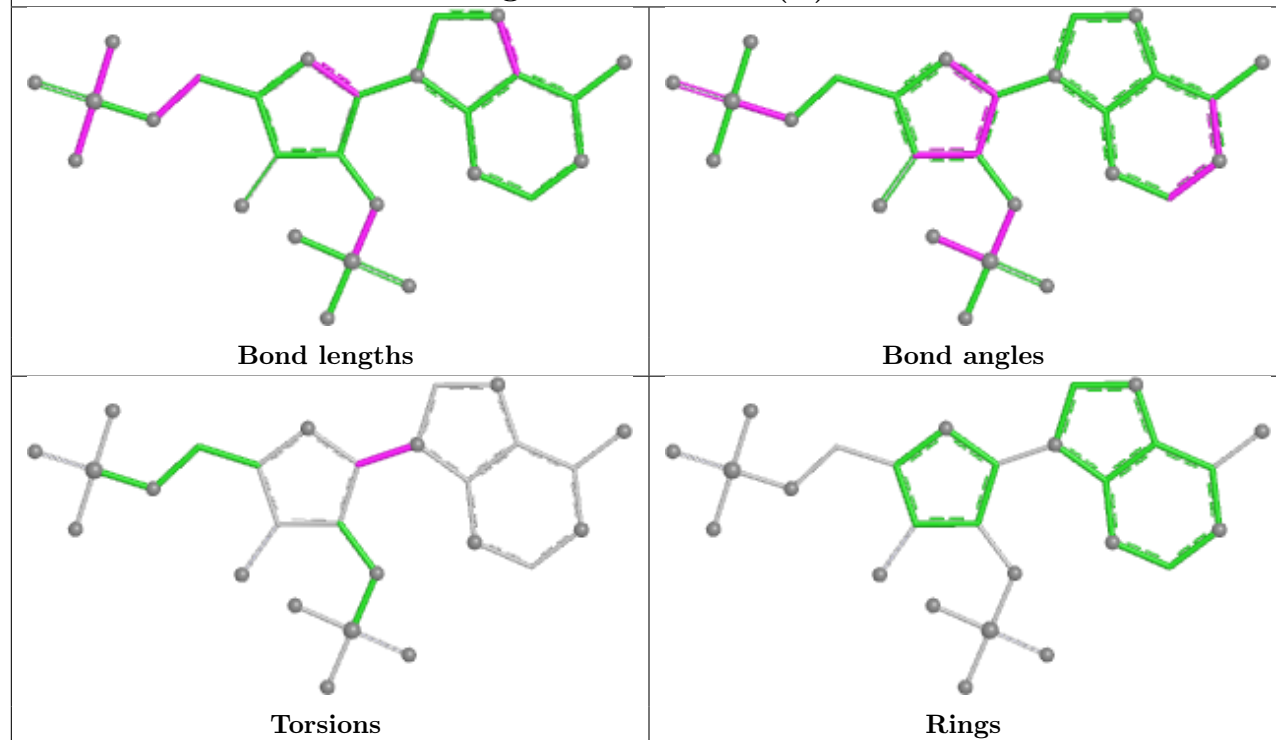
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	317	SO4	2	0

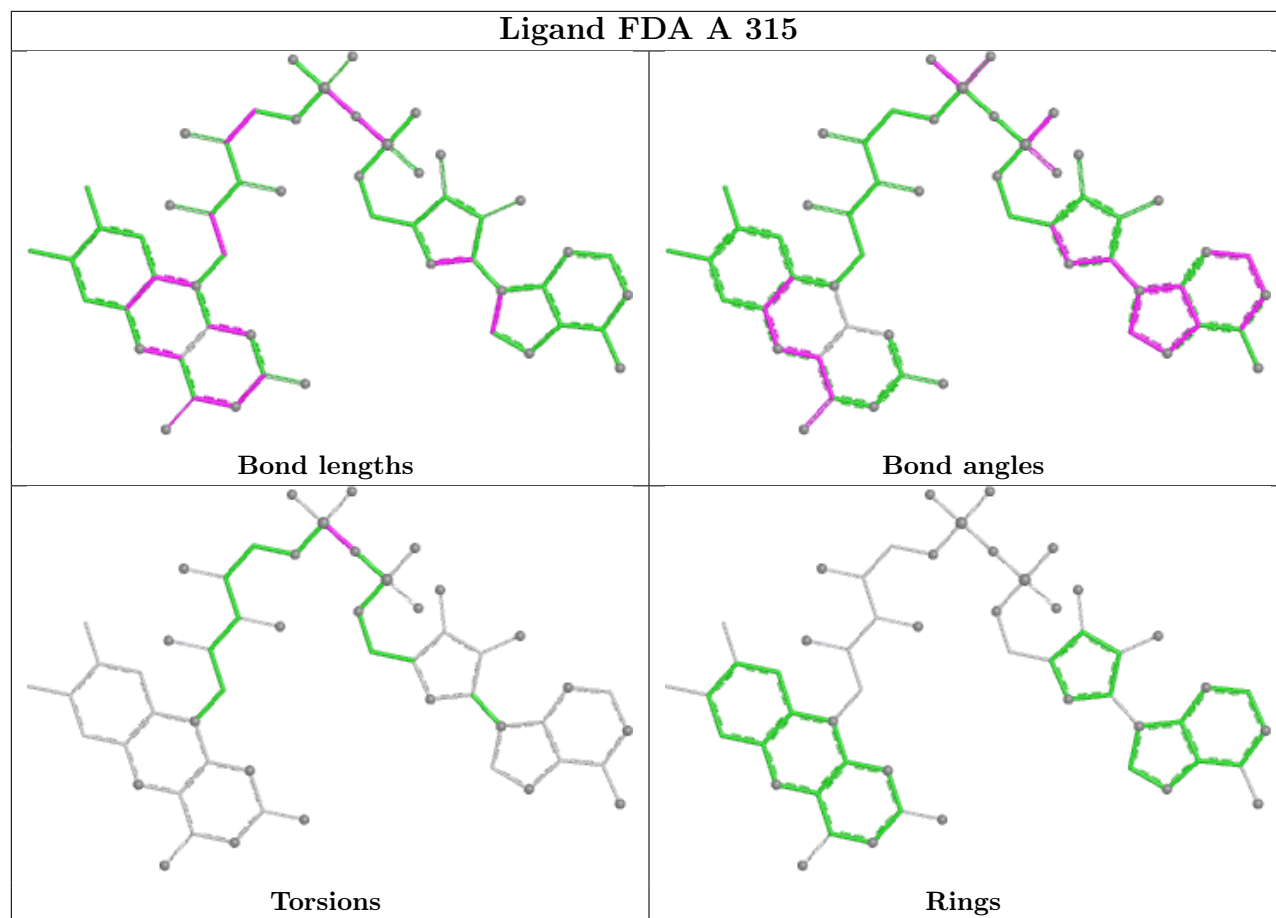
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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 A2P A 316 (A)



Ligand A2P A 316 (B)





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	296/314 (94%)	-0.80	0 100 100	3, 19, 52, 95	0

There are no RSRZ outliers to report.

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](#)

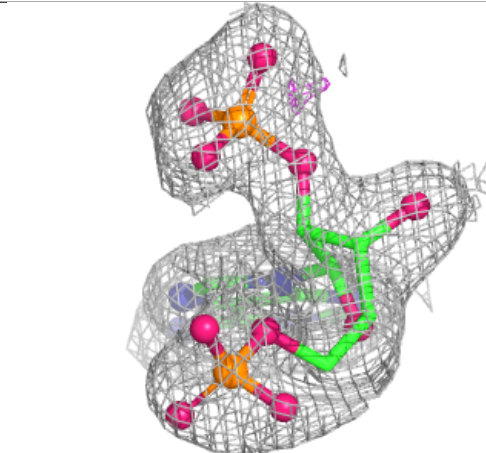
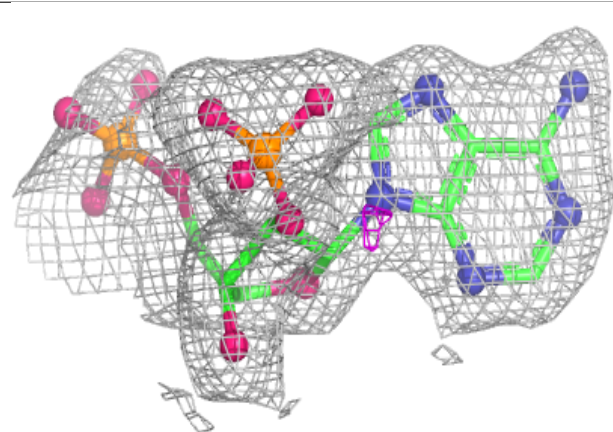
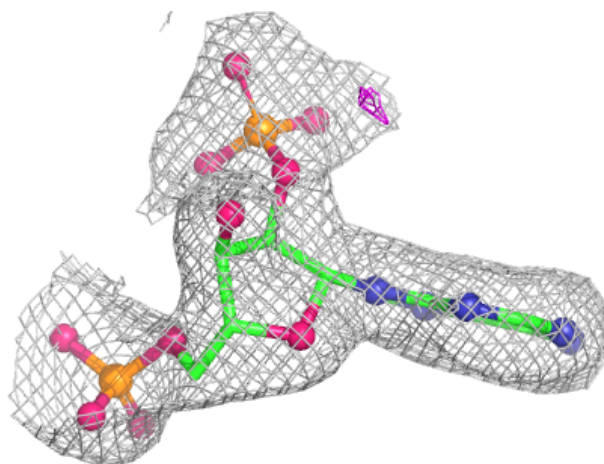
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q < 0.9’ lists 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	SO4	A	317	5/5	0.97	0.09	26,37,100,100	0
4	A2P	A	316[A]	27/27	0.98	0.09	10,25,38,100	27
4	A2P	A	316[B]	27/27	0.98	0.09	9,26,36,100	5
3	FDA	A	315	53/53	0.99	0.06	1,15,100,100	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.

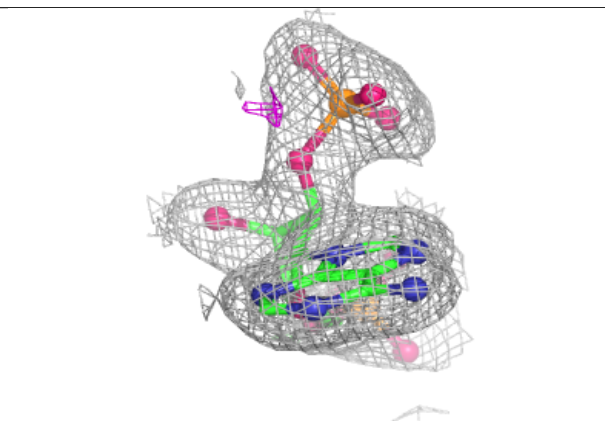
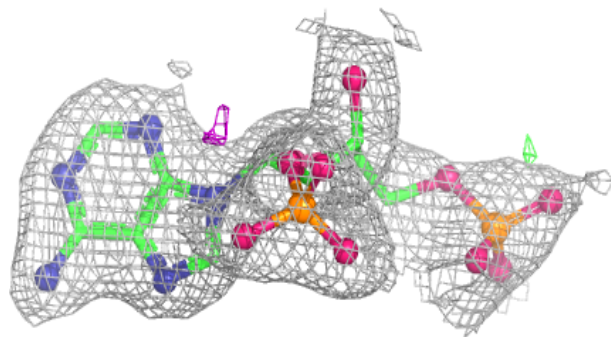
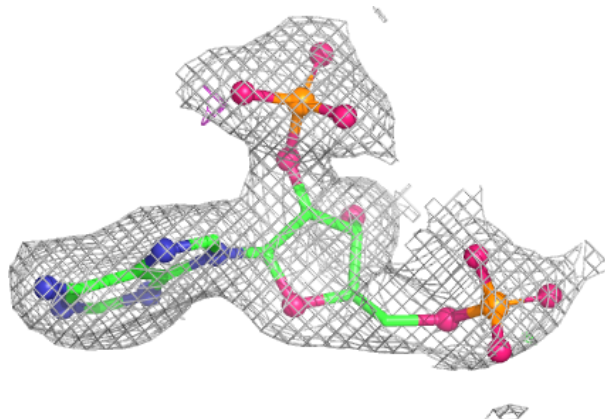
Electron density around A2P A 316 (A):

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

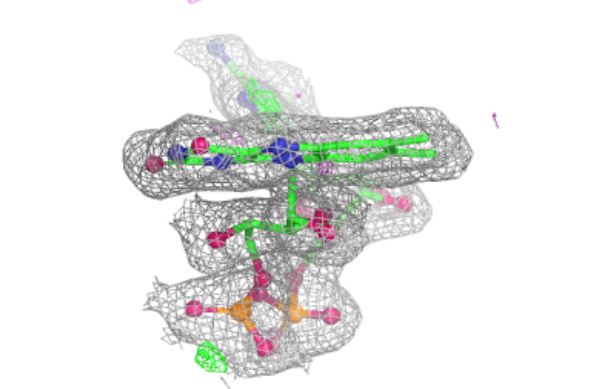
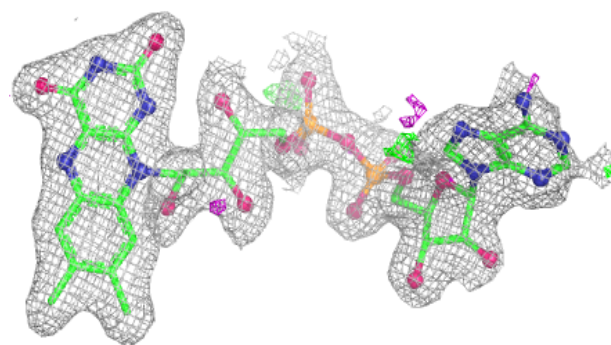
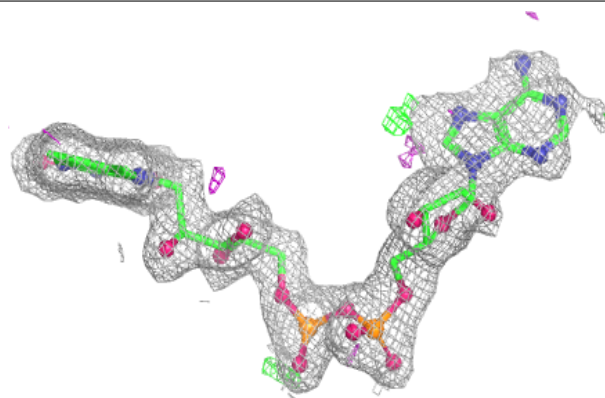


Electron density around A2P A 316 (B):

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

**Electron density around FDA A 315:**

$2mF_o-DF_c$ (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.