



Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 01:26 am BST

PDB ID : 1PAG
Title : THE 2.5 ANGSTROMS STRUCTURE OF POKEWEED ANTIVIRAL PROTEIN
Authors : Monzingo, A.F.; Collins, E.J.; Ernst, S.R.; Irvin, J.D.; Robertus, J.D.
Deposited on : 1992-10-19
Resolution : 2.80 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

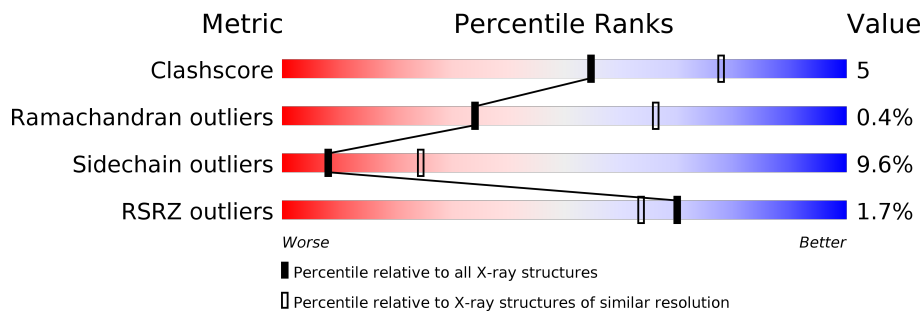
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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 on 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	262	<div style="display: flex; align-items: center;"> <div style="width: 2%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 73%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 23%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 5%; height: 10px; background-color: orange;"></div> </div> <p style="margin-left: 20px;">2% 73% 23% 5%</p>
1	B	262	<div style="display: flex; align-items: center;"> <div style="width: 1%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 71%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 24%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: orange;"></div> </div> <p style="margin-left: 20px;">% 71% 24% ..</p>

2 Entry composition [i](#)

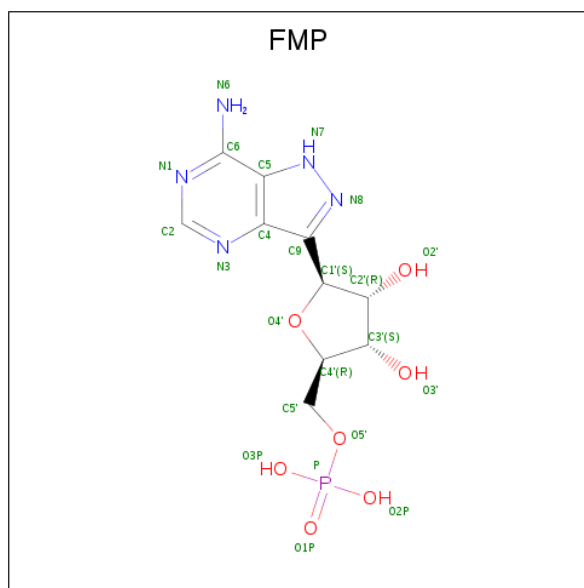
There are 3 unique types of molecules in this entry. The entry contains 4244 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 POKWEED ANTIVIRAL PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	262	Total 2060	C 1300	N 355	O 396	S 9	0	0	0
1	B	262	Total 2060	C 1300	N 355	O 396	S 9	0	0	0

- Molecule 2 is FORMYCIN-5'-MONOPHOSPHATE (three-letter code: FMP) (formula: $C_{10}H_{14}N_5O_7P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	A	1	Total 23	C 10	N 5	O 7	P 1	0	0
2	B	1	Total 23	C 10	N 5	O 7	P 1	0	0

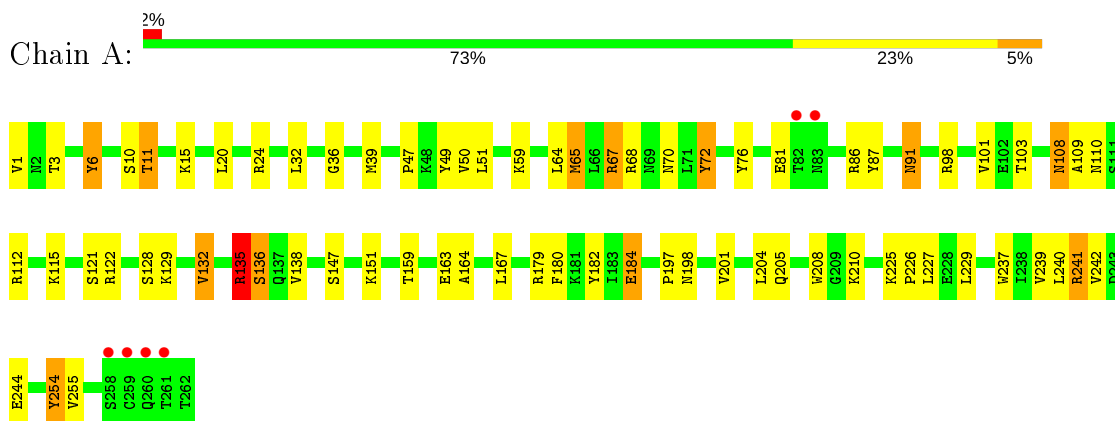
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	37	Total 37	O 37	0	0
3	B	41	Total 41	O 41	0	0

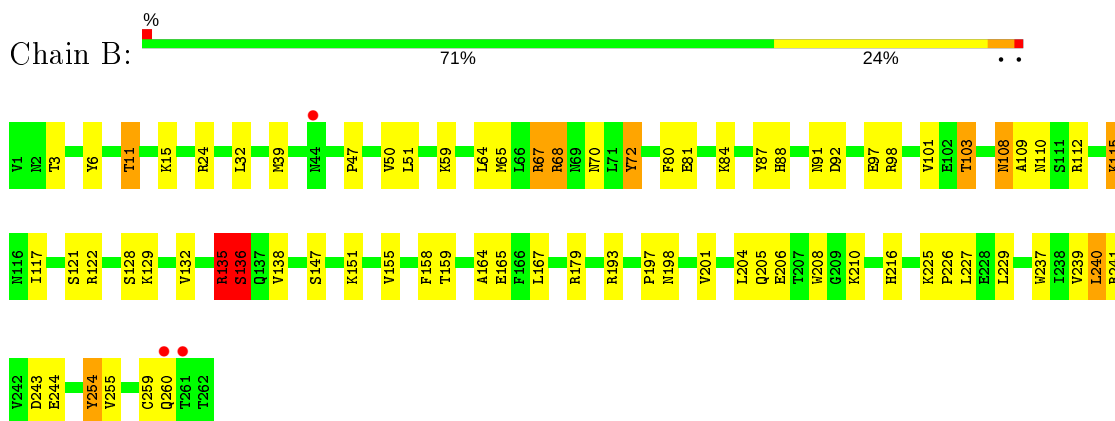
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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: POKEWEED ANTIVIRAL PROTEIN



- Molecule 1: POKEWEED ANTIVIRAL PROTEIN



4 Data and refinement statistics i

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	49.40Å 50.10Å 65.20Å 80.00° 113.20° 116.50°	Depositor
Resolution (Å)	(Not available) – 2.80 19.98 – 2.80	Depositor EDS
% Data completeness (in resolution range)	(Not available) ((Not available)-2.80) 94.4 (19.98-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.30 (at 2.79Å)	Xtrriage
Refinement program	X-PLOR	Depositor
R, R_{free}	0.200 , (Not available) 0.193 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	26.3	Xtrriage
Anisotropy	0.421	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 66.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.36$	Xtrriage
Estimated twinning fraction	0.000 for -h,-k,h+1	Xtrriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	4244	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.41% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: FMP

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	1.34	6/2096 (0.3%)	1.71	37/2840 (1.3%)
1	B	1.32	4/2096 (0.2%)	1.73	40/2840 (1.4%)
All	All	1.33	10/4192 (0.2%)	1.72	77/5680 (1.4%)

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	184	GLU	CD-OE2	-7.93	1.17	1.25
1	B	208	TRP	NE1-CE2	-6.70	1.28	1.37
1	A	10	SER	CA-CB	-6.40	1.43	1.52
1	B	165	GLU	CD-OE1	-6.20	1.18	1.25
1	A	163	GLU	CD-OE2	-5.60	1.19	1.25
1	A	6	TYR	CE1-CZ	-5.55	1.31	1.38
1	B	237	TRP	NE1-CE2	-5.47	1.30	1.37
1	A	241	ARG	NE-CZ	5.43	1.40	1.33
1	A	208	TRP	CD1-NE1	-5.12	1.29	1.38
1	B	216	HIS	CB-CG	5.10	1.59	1.50

All (77) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	112	ARG	NE-CZ-NH2	-14.33	113.13	120.30
1	B	112	ARG	NE-CZ-NH1	13.45	127.02	120.30
1	A	112	ARG	NE-CZ-NH2	-12.51	114.05	120.30
1	B	208	TRP	CE2-CD2-CG	-11.30	98.26	107.30
1	A	122	ARG	NE-CZ-NH2	-11.22	114.69	120.30
1	B	135	ARG	NE-CZ-NH2	-11.08	114.76	120.30
1	B	68	ARG	NE-CZ-NH2	-10.68	114.96	120.30
1	B	179	ARG	NE-CZ-NH1	9.91	125.26	120.30
1	A	122	ARG	NE-CZ-NH1	9.90	125.25	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	135	ARG	NE-CZ-NH1	9.85	125.22	120.30
1	A	112	ARG	NE-CZ-NH1	9.55	125.08	120.30
1	B	67	ARG	NE-CZ-NH1	9.50	125.05	120.30
1	A	67	ARG	NE-CZ-NH1	9.43	125.01	120.30
1	B	98	ARG	NE-CZ-NH1	8.92	124.76	120.30
1	A	135	ARG	NE-CZ-NH2	-8.88	115.86	120.30
1	B	208	TRP	CD1-CG-CD2	8.69	113.25	106.30
1	B	68	ARG	NE-CZ-NH1	8.69	124.64	120.30
1	A	179	ARG	NE-CZ-NH1	8.47	124.54	120.30
1	B	254	TYR	CB-CG-CD1	-8.01	116.19	121.00
1	A	76	TYR	CB-CG-CD2	-7.81	116.31	121.00
1	A	49	TYR	CB-CG-CD1	-7.74	116.35	121.00
1	A	98	ARG	NE-CZ-NH2	-7.50	116.55	120.30
1	B	92	ASP	CB-CG-OD1	7.49	125.04	118.30
1	A	24	ARG	NE-CZ-NH2	-7.46	116.57	120.30
1	B	24	ARG	NE-CZ-NH1	7.42	124.01	120.30
1	A	208	TRP	CD1-CG-CD2	6.94	111.85	106.30
1	A	72	TYR	CB-CG-CD2	-6.92	116.85	121.00
1	B	67	ARG	NE-CZ-NH2	-6.85	116.87	120.30
1	A	208	TRP	CE2-CD2-CG	-6.77	101.89	107.30
1	A	237	TRP	CG-CD2-CE3	6.64	139.88	133.90
1	A	67	ARG	NE-CZ-NH2	-6.56	117.02	120.30
1	A	98	ARG	NE-CZ-NH1	6.53	123.56	120.30
1	B	122	ARG	NE-CZ-NH1	6.51	123.56	120.30
1	B	237	TRP	CG-CD1-NE1	-6.48	103.62	110.10
1	B	208	TRP	NE1-CE2-CD2	6.45	113.75	107.30
1	B	108	ASN	CB-CG-ND2	6.39	132.05	116.70
1	B	254	TYR	CG-CD1-CE1	-6.39	116.19	121.30
1	A	254	TYR	CB-CG-CD1	-6.23	117.26	121.00
1	A	86	ARG	NE-CZ-NH1	6.20	123.40	120.30
1	B	237	TRP	CE2-CD2-CG	-6.10	102.42	107.30
1	B	158	PHE	CB-CG-CD2	-6.06	116.56	120.80
1	A	208	TRP	CG-CD1-NE1	-5.98	104.12	110.10
1	B	122	ARG	NE-CZ-NH2	-5.90	117.35	120.30
1	B	103	THR	N-CA-CB	-5.90	99.09	110.30
1	A	237	TRP	CD1-CG-CD2	5.89	111.02	106.30
1	A	242	VAL	CA-CB-CG1	-5.88	102.08	110.90
1	A	81	GLU	CA-CB-CG	-5.83	100.56	113.40
1	B	243	ASP	CB-CG-OD1	5.70	123.43	118.30
1	A	108	ASN	CB-CG-ND2	5.57	130.07	116.70
1	A	132	VAL	N-CA-CB	-5.57	99.24	111.50
1	B	237	TRP	CD1-CG-CD2	5.55	110.74	106.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	86	ARG	NE-CZ-NH2	-5.53	117.53	120.30
1	A	87	TYR	CG-CD2-CE2	-5.49	116.91	121.30
1	B	136	SER	N-CA-CB	5.45	118.68	110.50
1	A	91	ASN	CA-CB-CG	5.44	125.37	113.40
1	B	240	LEU	CA-CB-CG	5.41	127.73	115.30
1	B	208	TRP	CD2-CE2-CZ2	-5.40	115.82	122.30
1	B	260	GLN	CA-CB-CG	-5.34	101.65	113.40
1	A	182	TYR	CG-CD1-CE1	-5.33	117.04	121.30
1	A	180	PHE	CB-CG-CD2	-5.30	117.09	120.80
1	A	182	TYR	CB-CG-CD2	-5.29	117.83	121.00
1	B	72	TYR	CB-CG-CD2	-5.25	117.85	121.00
1	B	81	GLU	CA-CB-CG	-5.23	101.89	113.40
1	A	65	MET	CA-CB-CG	-5.18	104.48	113.30
1	A	180	PHE	CD1-CG-CD2	5.16	125.01	118.30
1	B	237	TRP	CB-CG-CD1	-5.15	120.30	127.00
1	B	179	ARG	NE-CZ-NH2	-5.12	117.74	120.30
1	B	115	LYS	N-CA-C	-5.12	97.17	111.00
1	B	72	TYR	CG-CD1-CE1	-5.11	117.21	121.30
1	B	87	TYR	CB-CG-CD2	-5.10	117.94	121.00
1	A	237	TRP	CE2-CD2-CG	-5.09	103.23	107.30
1	A	115	LYS	N-CA-C	-5.08	97.30	111.00
1	B	193	ARG	NE-CZ-NH2	-5.05	117.78	120.30
1	A	36	GLY	O-C-N	-5.04	114.64	122.70
1	B	206	GLU	O-C-N	-5.04	114.64	122.70
1	B	117	ILE	N-CA-C	-5.02	97.45	111.00
1	A	1	VAL	N-CA-C	-5.02	97.46	111.00

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	2060	0	2083	20	0
1	B	2060	0	2083	23	0
2	A	23	0	12	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	23	0	12	2	0
3	A	37	0	0	0	0
3	B	41	0	0	1	0
All	All	4244	0	4190	43	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 5.

All (43) 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:121:SER:HB2	2:A:263:FMP:HN61	1.49	0.77
1:A:136:SER:HA	1:A:197:PRO:HD2	1.65	0.76
1:B:121:SER:HB2	2:B:263:FMP:HN61	1.55	0.72
1:B:136:SER:HA	1:B:197:PRO:HD2	1.72	0.71
1:B:227:LEU:HB3	1:B:239:VAL:HG23	1.74	0.69
1:A:227:LEU:HB3	1:A:239:VAL:HG23	1.75	0.66
1:A:67:ARG:HH21	1:A:70:ASN:HD21	1.50	0.60
1:B:67:ARG:HH21	1:B:70:ASN:HD21	1.51	0.59
1:B:39:MET:HE3	1:B:255:VAL:H	1.68	0.57
1:A:39:MET:HE3	1:A:255:VAL:H	1.70	0.56
1:B:135:ARG:HH11	1:B:205:GLN:HE21	1.55	0.53
1:B:135:ARG:HH11	1:B:205:GLN:NE2	2.07	0.53
1:A:198:ASN:O	1:A:201:VAL:HG22	2.09	0.52
1:A:11:THR:HG22	1:A:15:LYS:HB2	1.91	0.52
1:A:135:ARG:HH11	1:A:205:GLN:HE21	1.59	0.51
1:A:201:VAL:O	1:A:205:GLN:HG3	2.11	0.51
1:A:135:ARG:HH11	1:A:205:GLN:NE2	2.08	0.51
1:B:11:THR:HG22	1:B:15:LYS:HB2	1.93	0.50
1:A:39:MET:HE2	1:A:254:TYR:HA	1.95	0.49
1:A:47:PRO:O	1:A:68:ARG:HD3	2.12	0.49
1:B:72:TYR:OH	2:B:263:FMP:H2'	2.14	0.48
1:B:259:CYS:HA	3:B:809:HOH:O	2.14	0.47
1:A:72:TYR:OH	2:A:263:FMP:H2'	2.14	0.47
1:B:198:ASN:O	1:B:201:VAL:HG22	2.15	0.46
1:B:129:LYS:HB2	1:B:164:ALA:HB2	1.99	0.45
1:A:6:TYR:OH	1:A:11:THR:HG21	2.16	0.45
1:B:109:ALA:O	1:B:110:ASN:HB2	2.17	0.44
1:B:201:VAL:O	1:B:205:GLN:HG3	2.17	0.44
1:A:241:ARG:HD3	1:A:244:GLU:OE2	2.17	0.44
1:A:129:LYS:HB2	1:A:164:ALA:HB2	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:65:MET:SD	1:A:101:VAL:HA	2.58	0.44
1:B:241:ARG:HD3	1:B:244:GLU:OE2	2.18	0.44
1:A:20:LEU:HD12	1:A:184:GLU:HG3	2.00	0.43
1:B:47:PRO:O	1:B:68:ARG:HD3	2.18	0.43
1:A:226:PRO:HB3	1:A:240:LEU:HD23	2.00	0.43
1:B:6:TYR:OH	1:B:11:THR:HG21	2.18	0.43
1:B:88:HIS:HA	1:B:115:LYS:O	2.19	0.41
1:B:65:MET:SD	1:B:101:VAL:HA	2.60	0.41
1:B:80:PHE:O	1:B:84:LYS:HB2	2.21	0.41
1:B:226:PRO:HB3	1:B:240:LEU:HD23	2.03	0.41
1:A:109:ALA:O	1:A:110:ASN:HB2	2.21	0.40
1:B:97:GLU:O	1:B:101:VAL:HG23	2.21	0.40
1:B:39:MET:HE2	1:B:254:TYR:CD1	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	260/262 (99%)	246 (95%)	13 (5%)	1 (0%)	34	66
1	B	260/262 (99%)	246 (95%)	13 (5%)	1 (0%)	34	66
All	All	520/524 (99%)	492 (95%)	26 (5%)	2 (0%)	34	66

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	136	SER
1	B	136	SER

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	234/234 (100%)	212 (91%)	22 (9%)	8	26
1	B	234/234 (100%)	211 (90%)	23 (10%)	8	24
All	All	468/468 (100%)	423 (90%)	45 (10%)	8	24

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	THR
1	A	11	THR
1	A	32	LEU
1	A	50	VAL
1	A	51	LEU
1	A	59	LYS
1	A	64	LEU
1	A	91	ASN
1	A	103	THR
1	A	108	ASN
1	A	128	SER
1	A	132	VAL
1	A	135	ARG
1	A	138	VAL
1	A	147	SER
1	A	151	LYS
1	A	159	THR
1	A	167	LEU
1	A	204	LEU
1	A	210	LYS
1	A	225	LYS
1	A	229	LEU
1	B	3	THR
1	B	11	THR
1	B	32	LEU
1	B	50	VAL
1	B	51	LEU

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Mol	Chain	Res	Type
1	B	59	LYS
1	B	64	LEU
1	B	91	ASN
1	B	103	THR
1	B	108	ASN
1	B	128	SER
1	B	132	VAL
1	B	135	ARG
1	B	138	VAL
1	B	147	SER
1	B	151	LYS
1	B	155	VAL
1	B	159	THR
1	B	167	LEU
1	B	204	LEU
1	B	210	LYS
1	B	225	LYS
1	B	229	LEU

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

Mol	Chain	Res	Type
1	A	205	GLN
1	B	205	GLN

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 carbohydrates in this entry.

5.6 Ligand geometry

2 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	FMP	B	263	-	24,25,25	2.03	7 (29%)	26,38,38	2.06	12 (46%)
2	FMP	A	263	-	24,25,25	2.26	6 (25%)	26,38,38	2.45	10 (38%)

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	FMP	B	263	-	-	2/6/26/26	0/3/3/3
2	FMP	A	263	-	-	2/6/26/26	0/3/3/3

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	263	FMP	C9-C4	-5.75	1.33	1.43
2	A	263	FMP	O4'-C1'	-4.54	1.37	1.44
2	A	263	FMP	C2'-C1'	-4.14	1.49	1.54
2	B	263	FMP	C2'-C1'	-4.14	1.49	1.54
2	B	263	FMP	C4-N3	3.91	1.42	1.37
2	A	263	FMP	N7-N8	-3.68	1.30	1.37
2	B	263	FMP	O4'-C1'	-3.55	1.39	1.44
2	B	263	FMP	N7-N8	-3.53	1.30	1.37
2	A	263	FMP	P-O3P	3.50	1.68	1.54
2	B	263	FMP	C9-C4	-3.47	1.37	1.43
2	B	263	FMP	C4-C5	-2.83	1.33	1.43
2	A	263	FMP	C3'-C4'	-2.70	1.46	1.53
2	B	263	FMP	P-O3P	2.47	1.64	1.54

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	263	FMP	C5-C4-N3	-6.13	117.80	124.92
2	A	263	FMP	C4'-O4'-C1'	-5.39	102.77	109.42
2	B	263	FMP	C5-C4-N3	-5.10	118.99	124.92
2	A	263	FMP	O4'-C4'-C3'	4.74	114.50	105.11
2	A	263	FMP	O5'-C5'-C4'	4.18	123.36	108.99
2	B	263	FMP	O5'-C5'-C4'	3.48	120.98	108.99
2	B	263	FMP	C4'-O4'-C1'	-3.04	105.68	109.42
2	B	263	FMP	C2-N3-C4	3.01	117.15	114.81
2	A	263	FMP	C2-N3-C4	2.97	117.12	114.81
2	A	263	FMP	O3'-C3'-C2'	-2.78	102.82	111.82
2	A	263	FMP	C2'-C3'-C4'	-2.76	97.29	102.64
2	B	263	FMP	O2'-C2'-C3'	-2.66	103.22	111.82
2	B	263	FMP	O3P-P-O5'	2.60	113.66	106.73
2	B	263	FMP	C2'-C3'-C4'	-2.47	97.85	102.64
2	B	263	FMP	C3'-C2'-C1'	2.27	104.54	101.93
2	A	263	FMP	O2P-P-O1P	2.26	119.54	110.68
2	B	263	FMP	O4'-C1'-C2'	-2.24	101.03	104.66
2	B	263	FMP	N3-C2-N1	-2.16	125.31	128.68
2	A	263	FMP	C2-N1-C6	-2.13	115.11	118.75
2	A	263	FMP	C3'-C2'-C1'	2.07	104.32	101.93
2	B	263	FMP	O2P-P-O1P	2.03	118.62	110.68
2	B	263	FMP	O4'-C4'-C3'	2.02	109.10	105.11

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	263	FMP	C3'-C4'-C5'-O5'
2	B	263	FMP	C3'-C4'-C5'-O5'
2	A	263	FMP	O4'-C4'-C5'-O5'
2	B	263	FMP	O4'-C4'-C5'-O5'

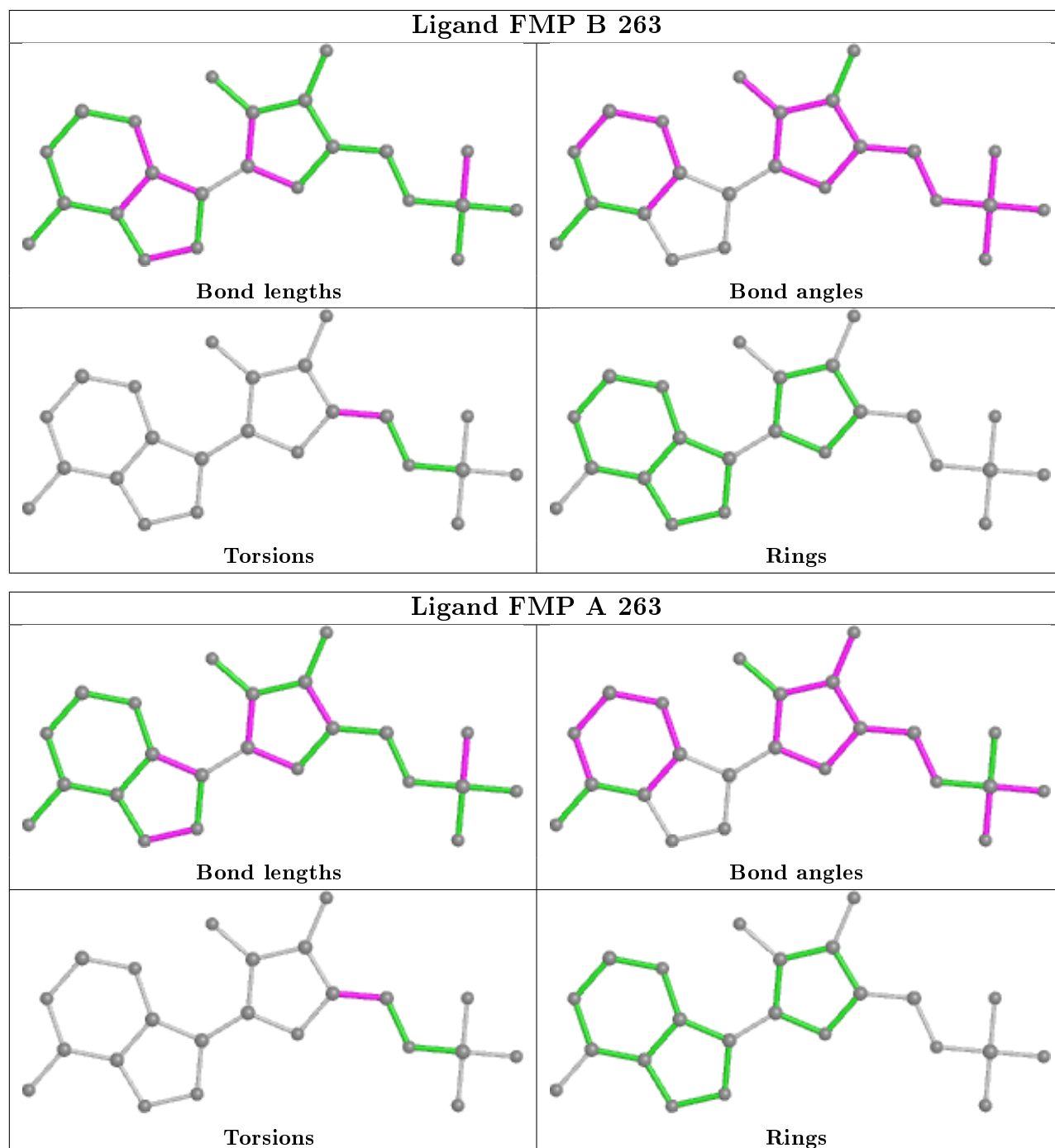
There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	263	FMP	2	0
2	A	263	FMP	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.



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	262/262 (100%)	-0.42	6 (2%) 60 51	6, 19, 34, 43	0
1	B	262/262 (100%)	-0.43	3 (1%) 80 75	5, 19, 33, 42	0
All	All	524/524 (100%)	-0.43	9 (1%) 70 63	5, 19, 34, 43	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	259	CYS	4.2
1	A	258	SER	3.7
1	A	261	THR	3.2
1	B	44	ASN	3.2
1	A	260	GLN	3.0
1	B	260	GLN	2.9
1	A	82	THR	2.8
1	A	83	ASN	2.3
1	B	261	THR	2.2

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 carbohydrates in this entry.

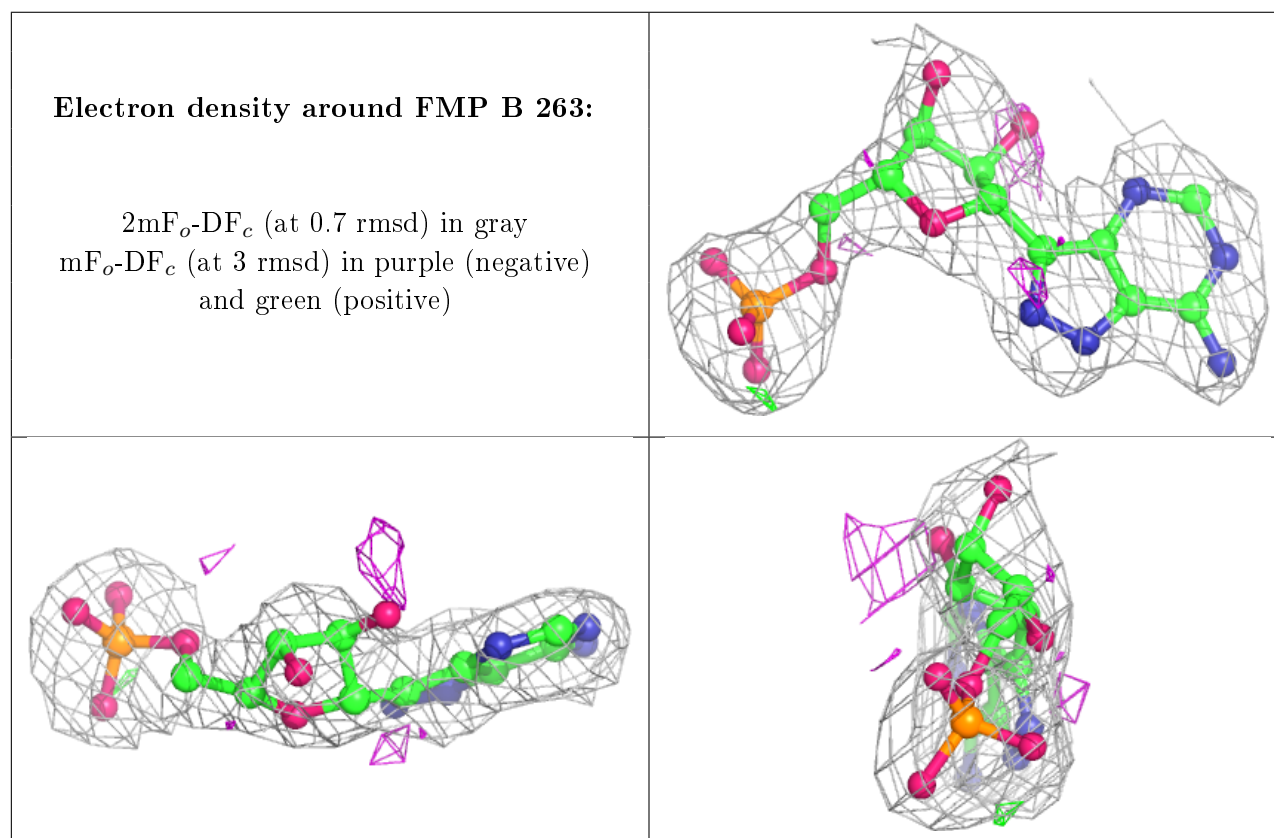
6.4 Ligands [i](#)

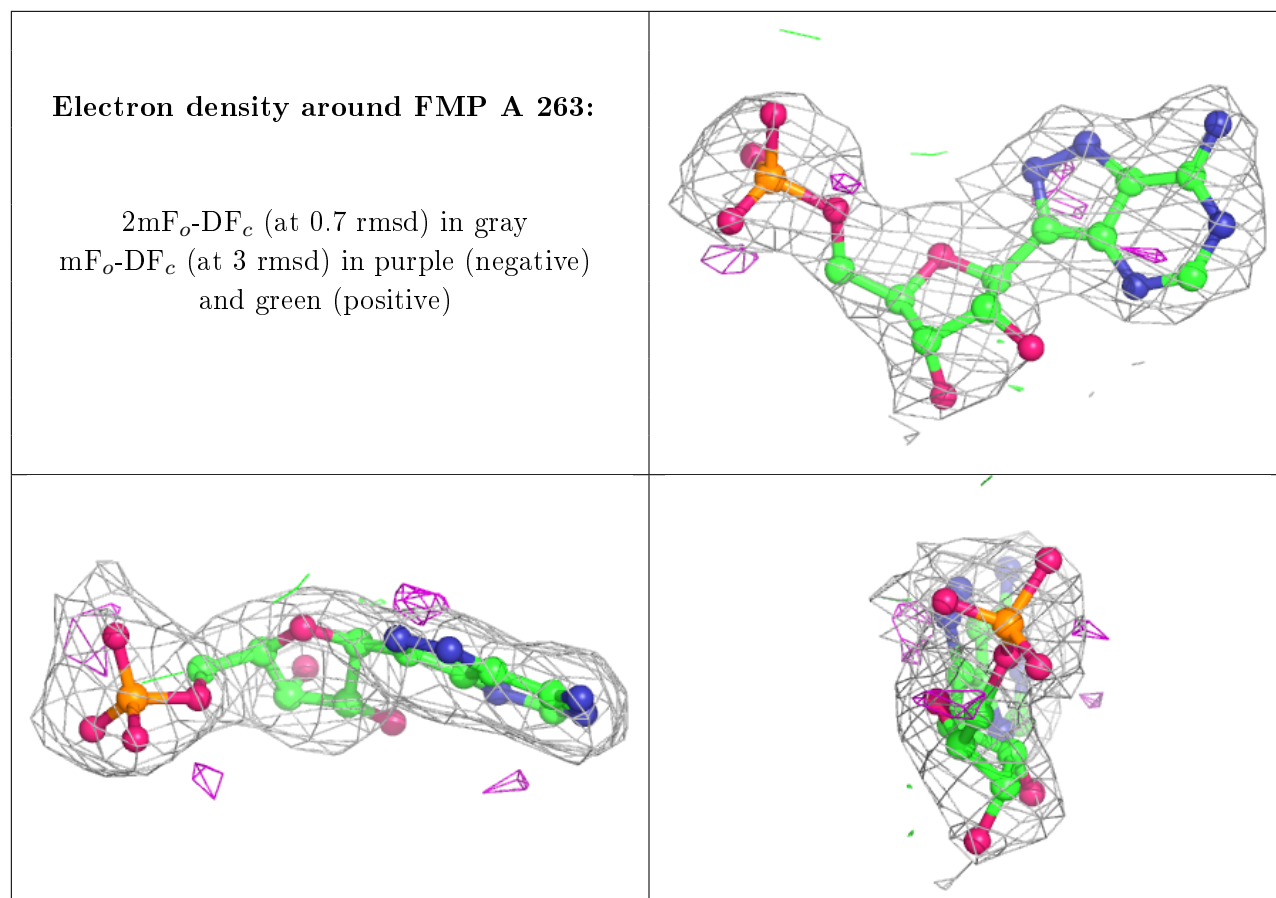
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,

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	FMP	B	263	23/23	0.92	0.19	14,22,27,30	0
2	FMP	A	263	23/23	0.94	0.18	14,21,26,29	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.





6.5 Other polymers [i](#)

There are no such residues in this entry.