



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 07:40 AM UTC

PDB ID : 1AMR / pdb_00001amr
Title : X-RAY CRYSTALLOGRAPHIC STUDY OF PYRIDOXAMINE 5'-
PHOSPHATE-TYPE ASPARTATE AMINOTRANSFERASES FROM
ESCHERICHIA COLI IN THREE FORMS
Authors : Miyahara, I.; Hirotsu, K.; Hayashi, H.; Kagamiyama, H.
Deposited on : 1994-07-01
Resolution : 2.10 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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

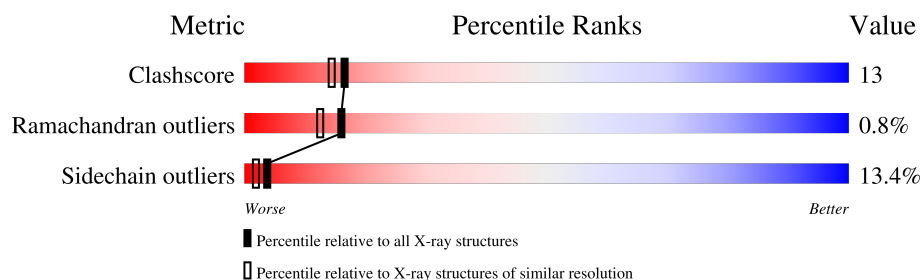
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	396	 67% 26% 6%

2 Entry composition [i](#)

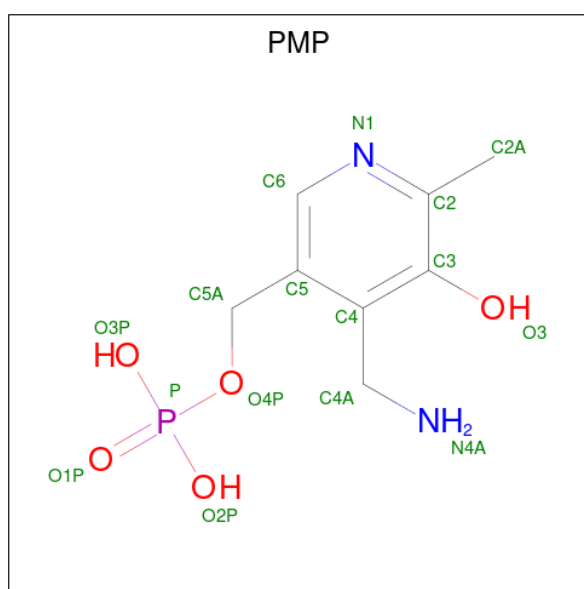
There are 4 unique types of molecules in this entry. The entry contains 3357 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 ASPARTATE AMINOTRANSFERASE.

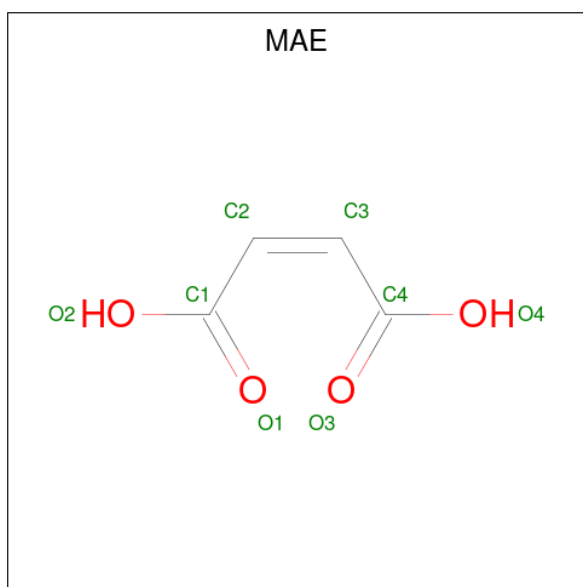
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	396	3069	1936	536	584	13	0	0	0

- Molecule 2 is 4'-DEOXY-4'-AMINOPYRIDOXAL-5'-PHOSPHATE (CCD ID: PMP) (formula: $C_8H_{13}N_2O_5P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	A	1	16	8	2	5	1	0	0

- Molecule 3 is MALEIC ACID (CCD ID: MAE) (formula: $C_4H_4O_4$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			8	4	4		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	264	Total	O	0	0
			264	264		

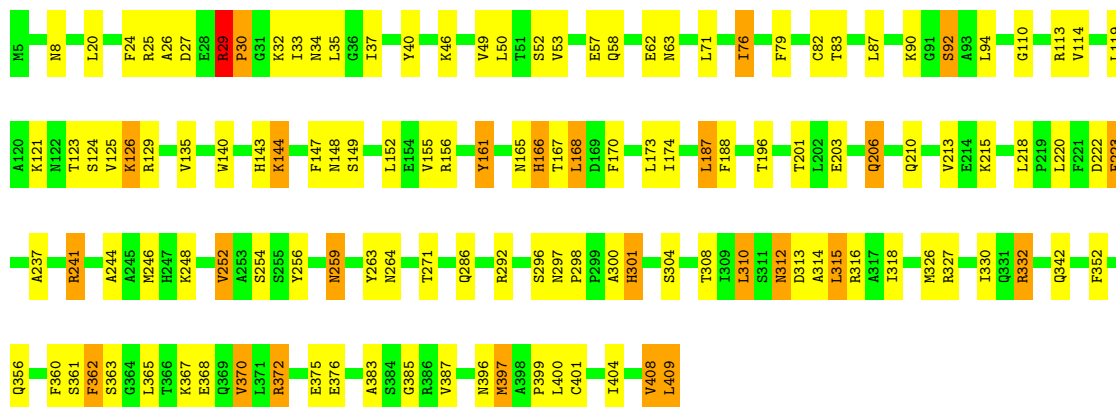
3 Residue-property plots

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. 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. 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.

Note EDS was not executed.

• Molecule 1: ASPARTATE AMINOTRANSFERASE

Chain A: 



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	157.60Å 85.70Å 78.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 2.10	Depositor
% Data completeness (in resolution range)	(Not available) (10.00-2.10)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR, TOM, TOM/FRODO	Depositor
R, R_{free}	0.194 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3357	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

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

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	0.54	0/3130	1.04	21/4240 (0.5%)

There are no bond length outliers.

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	63	ASN	N-CA-C	8.52	123.97	113.50
1	A	298	PRO	N-CA-C	7.67	120.06	110.70
1	A	300	ALA	N-CA-C	7.57	119.31	111.14
1	A	29	ARG	CA-C-N	7.15	128.78	119.84
1	A	29	ARG	C-N-CA	7.15	128.78	119.84
1	A	301	HIS	N-CA-C	6.66	122.29	111.37
1	A	360	PHE	N-CA-C	6.47	120.07	109.72
1	A	196	THR	N-CA-C	5.85	118.60	111.82
1	A	92	SER	N-CA-C	5.75	123.05	110.80
1	A	76	ILE	CA-C-N	5.56	125.03	119.24
1	A	76	ILE	C-N-CA	5.56	125.03	119.24
1	A	363	SER	N-CA-C	5.45	117.30	111.36
1	A	161	TYR	N-CA-C	5.38	118.01	109.24
1	A	26	ALA	N-CA-C	-5.31	107.34	113.88
1	A	201	THR	N-CA-C	-5.20	103.82	110.53
1	A	264	ASN	N-CA-C	5.16	119.44	113.20
1	A	71	LEU	N-CA-C	5.14	119.40	113.18
1	A	362	PHE	N-CA-C	-5.07	99.51	108.23
1	A	166	HIS	N-CA-C	-5.06	104.44	111.52
1	A	140	TRP	CA-C-N	5.00	125.27	119.47
1	A	140	TRP	C-N-CA	5.00	125.27	119.47

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	3069	0	3018	78	0
2	A	16	0	11	1	0
3	A	8	0	2	0	0
4	A	264	0	0	22	0
All	All	3357	0	3031	79	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 13.

All (79) 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:252:VAL:HG13	1:A:271:THR:HB	1.59	0.85
1:A:397:MET:HE3	1:A:397:MET:HA	1.61	0.81
1:A:318:ILE:HG22	4:A:642:HOH:O	1.83	0.78
1:A:286:GLN:HG2	4:A:510:HOH:O	1.84	0.76
1:A:135:VAL:HG11	1:A:144:LYS:HE2	1.71	0.71
1:A:52:SER:HB3	4:A:642:HOH:O	1.93	0.69
1:A:94:LEU:HD11	1:A:244:ALA:HB1	1.76	0.68
1:A:356:GLN:NE2	1:A:361:SER:HA	2.09	0.67
1:A:301:HIS:HD2	4:A:481:HOH:O	1.77	0.66
1:A:206:GLN:HG3	4:A:503:HOH:O	1.96	0.65
1:A:121:LYS:HE3	4:A:428:HOH:O	1.98	0.64
1:A:29:ARG:O	1:A:32:LYS:HG3	1.98	0.62
1:A:313:ASP:HB2	4:A:588:HOH:O	1.98	0.62
1:A:210:GLN:O	1:A:213:VAL:HG22	2.01	0.61
1:A:123:THR:HG22	1:A:124:SER:N	2.16	0.60
1:A:35:LEU:HD21	1:A:400:LEU:HD11	1.83	0.59
2:A:413:PMP:N4A	2:A:413:PMP:O3	2.35	0.59
1:A:312:ASN:ND2	1:A:315:LEU:H	1.99	0.59
1:A:126:LYS:HE3	4:A:559:HOH:O	2.02	0.58
1:A:58:GLN:O	1:A:62:GLU:HG2	2.03	0.57
1:A:33:ILE:HG22	1:A:35:LEU:HD13	1.86	0.57
1:A:187:LEU:HD11	1:A:222:ASP:HB2	1.86	0.57
1:A:79:PHE:O	1:A:83:THR:HG23	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:327:ARG:HG3	4:A:623:HOH:O	2.06	0.56
1:A:123:THR:CG2	1:A:125:VAL:HG23	2.36	0.55
1:A:312:ASN:HD21	1:A:314:ALA:HB3	1.71	0.55
1:A:82:CYS:HB3	1:A:310:LEU:HD23	1.89	0.55
1:A:27:ASP:O	1:A:32:LYS:HD3	2.06	0.55
1:A:143:HIS:HB2	1:A:144:LYS:HE3	1.88	0.55
1:A:57:GLU:OE2	1:A:301:HIS:HE1	1.91	0.54
1:A:123:THR:HG21	1:A:125:VAL:HG23	1.90	0.53
1:A:29:ARG:O	1:A:29:ARG:HG3	2.09	0.53
1:A:385:GLY:HA3	4:A:523:HOH:O	2.07	0.53
1:A:304:SER:O	1:A:308:THR:HG23	2.09	0.53
1:A:312:ASN:HD22	1:A:315:LEU:H	1.57	0.51
1:A:37:ILE:HG13	4:A:658:HOH:O	2.09	0.51
1:A:83:THR:HB	4:A:532:HOH:O	2.09	0.51
1:A:155:VAL:HG23	4:A:643:HOH:O	2.09	0.51
1:A:372:ARG:HG2	1:A:408:VAL:HG22	1.93	0.50
1:A:223:PHE:O	1:A:254:SER:HA	2.11	0.50
1:A:404:ILE:O	1:A:409:LEU:HD13	2.12	0.49
1:A:148:ASN:N	4:A:643:HOH:O	2.46	0.48
1:A:119:LEU:O	1:A:123:THR:HB	2.14	0.48
1:A:123:THR:CG2	1:A:124:SER:N	2.77	0.47
1:A:370:VAL:HG21	1:A:383:ALA:HA	1.96	0.47
1:A:83:THR:HG22	1:A:256:TYR:OH	2.15	0.46
1:A:368:GLU:HB2	4:A:564:HOH:O	2.15	0.46
1:A:170:PHE:O	1:A:174:ILE:HG12	2.16	0.46
1:A:396:ASN:O	1:A:399:PRO:HD2	2.16	0.44
1:A:24:PHE:CZ	1:A:34:ASN:HB2	2.52	0.44
1:A:256:TYR:HA	1:A:259:ASN:HD21	1.82	0.44
1:A:8:ASN:HB2	4:A:528:HOH:O	2.17	0.44
1:A:397:MET:HA	1:A:397:MET:CE	2.42	0.44
1:A:330:ILE:HD12	4:A:665:HOH:O	2.18	0.44
1:A:147:PHE:HB2	4:A:643:HOH:O	2.16	0.43
1:A:161:TYR:CE1	1:A:166:HIS:HA	2.53	0.43
1:A:187:LEU:HD12	1:A:188:PHE:N	2.34	0.43
1:A:40:TYR:CD2	1:A:326:MET:HE2	2.54	0.43
1:A:110:GLY:O	1:A:114:VAL:HG13	2.18	0.43
1:A:167:THR:CG2	1:A:168:LEU:N	2.81	0.43
1:A:76:ILE:O	1:A:79:PHE:HB3	2.19	0.43
1:A:396:ASN:C	1:A:399:PRO:HD2	2.44	0.43
1:A:332:ARG:NH1	4:A:678:HOH:O	2.51	0.43
1:A:332:ARG:NH1	4:A:488:HOH:O	2.51	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:292:ARG:HA	1:A:296:SER:HA	2.00	0.42
1:A:29:ARG:HA	1:A:30:PRO:HD2	1.60	0.42
1:A:352:PHE:HB2	4:A:624:HOH:O	2.18	0.42
1:A:123:THR:HG22	1:A:124:SER:H	1.83	0.42
1:A:168:LEU:HD21	1:A:173:LEU:HD12	2.01	0.42
1:A:223:PHE:HD1	1:A:223:PHE:HA	1.77	0.42
1:A:332:ARG:HA	1:A:332:ARG:HD2	1.77	0.41
1:A:113:ARG:NH2	1:A:149:SER:HB3	2.35	0.41
1:A:87:LEU:O	1:A:241:ARG:HD3	2.21	0.41
1:A:50:LEU:HB2	1:A:53:VAL:HG23	2.03	0.41
1:A:237:ALA:O	1:A:241:ARG:HG3	2.20	0.41
1:A:330:ILE:CD1	4:A:665:HOH:O	2.69	0.41
1:A:123:THR:CG2	1:A:124:SER:H	2.35	0.40
1:A:356:GLN:HE22	1:A:361:SER:HA	1.85	0.40
1:A:397:MET:HE2	1:A:401:CYS:SG	2.61	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	394/396 (100%)	375 (95%)	16 (4%)	3 (1%)	16 12

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	30	PRO
1	A	92	SER
1	A	263	TYR

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	320/320 (100%)	277 (87%)	43 (13%)	4 2

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	LEU
1	A	25	ARG
1	A	29	ARG
1	A	46	LYS
1	A	49	VAL
1	A	90	LYS
1	A	126	LYS
1	A	129	ARG
1	A	144	LYS
1	A	152	LEU
1	A	156	ARG
1	A	165	ASN
1	A	168	LEU
1	A	187	LEU
1	A	203	GLU
1	A	206	GLN
1	A	215	LYS
1	A	218	LEU
1	A	220	LEU
1	A	223	PHE
1	A	241	ARG
1	A	246	MET
1	A	248	LYS
1	A	252	VAL
1	A	259	ASN
1	A	297	ASN
1	A	310	LEU
1	A	312	ASN
1	A	315	LEU
1	A	316	ARG

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Mol	Chain	Res	Type
1	A	332	ARG
1	A	342	GLN
1	A	362	PHE
1	A	365	LEU
1	A	367	LYS
1	A	370	VAL
1	A	372	ARG
1	A	375	GLU
1	A	376	GLU
1	A	387	VAL
1	A	397	MET
1	A	408	VAL
1	A	409	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	69	ASN
1	A	84	GLN
1	A	96	ASN
1	A	166	HIS
1	A	210	GLN
1	A	259	ASN
1	A	286	GLN
1	A	301	HIS
1	A	312	ASN
1	A	321	GLN
1	A	328	GLN
1	A	342	GLN
1	A	356	GLN
1	A	357	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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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	$\# Z > 2$
3	MAE	A	414	-	7,7,7	1.43	1 (14%)	8,8,8	1.09	0
2	PMP	A	413	-	16,16,16	1.68	3 (18%)	22,23,23	1.67	5 (22%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	MAE	A	414	-	-	2/5/5/5	-
2	PMP	A	413	-	-	5/8/8/8	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	413	PMP	C3-C2	-4.12	1.36	1.41
2	A	413	PMP	C2A-C2	2.65	1.54	1.50
3	A	414	MAE	O2-C1	-2.21	1.24	1.30
2	A	413	PMP	P-O2P	-2.20	1.46	1.54

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	413	PMP	O4P-P-O1P	4.22	117.85	106.44
2	A	413	PMP	O3P-P-O4P	-2.69	99.66	106.67

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	413	PMP	O2P-P-O4P	-2.51	100.13	106.67
2	A	413	PMP	C5-C6-N1	-2.25	120.17	123.83
2	A	413	PMP	O3P-P-O2P	2.09	115.64	107.80

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	413	PMP	C5-C4-C4A-N4A
2	A	413	PMP	C5A-O4P-P-O1P
2	A	413	PMP	C5A-O4P-P-O2P
2	A	413	PMP	C5A-O4P-P-O3P
2	A	413	PMP	C3-C4-C4A-N4A
3	A	414	MAE	C2-C3-C4-O3
3	A	414	MAE	C2-C3-C4-O4

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	413	PMP	1	0

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

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.