



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 07:53 PM UTC

PDB ID : 1VFN / pdb_00001vfn
Title : PURINE NUCLEOSIDE PHOSPHORYLASE
Authors : Koellner, G.; Bzowska, A.
Deposited on : 1996-10-08
Resolution : 2.15 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

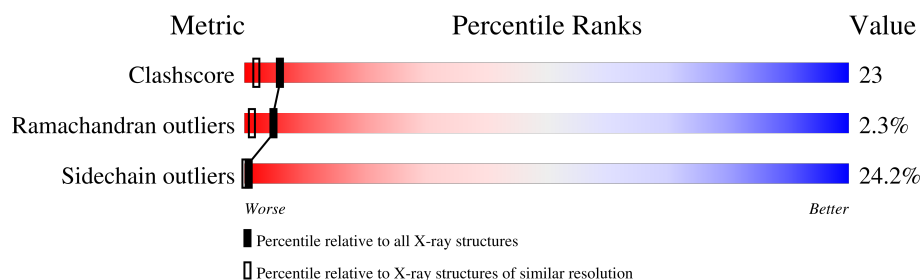
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.15 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	2159 (2.16-2.16)
Ramachandran outliers	187476	2134 (2.16-2.16)
Sidechain outliers	187428	2133 (2.16-2.16)

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	281	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 2243 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 PURINE-NUCLEOSIDE PHOSPHORYLASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	270	Total	C	N	O	S	0	0	0
			2102	1334	370	383	15			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	144	GLN	GLU	conflict	UNP P55859

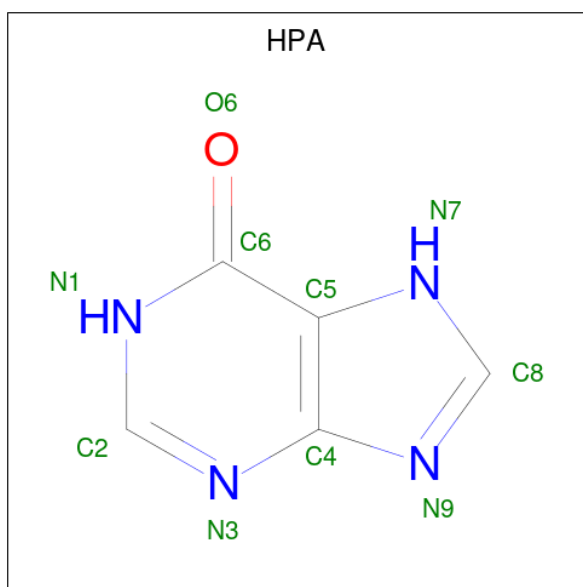
- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mg	0	0
			1	1		

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

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

- Molecule 4 is HYPOXANTHINE (CCD ID: HPA) (formula: C₅H₄N₄O).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			10	5	4	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	129	Total	O	0	0
			129	129		

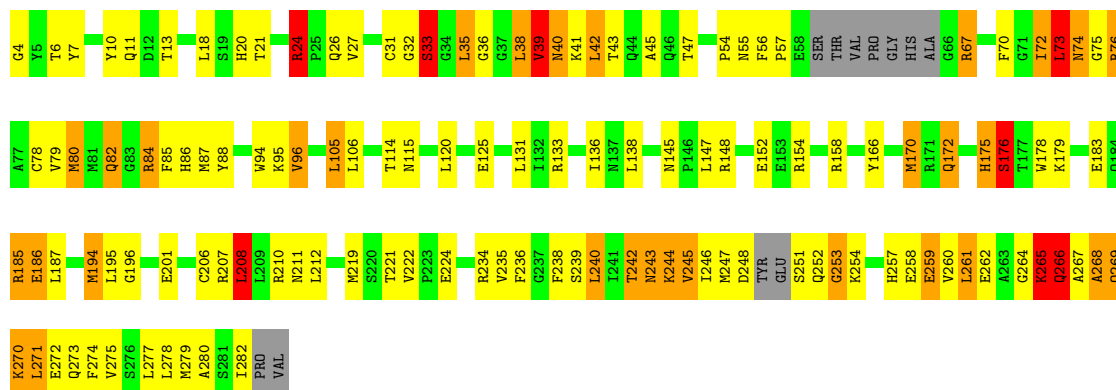
3 Residue-property plots

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: PURINE-NUCLEOSIDE PHOSPHORYLASE

Chain A: 



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 3	Depositor
Cell constants a, b, c, α , β , γ	94.11 Å 94.11 Å 94.11 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.15	Depositor
% Data completeness (in resolution range)	98.4 (20.00-2.15)	Depositor
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	TNT	Depositor
R, R_{free}	0.180 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2243	wwPDB-VP
Average B, all atoms (Å ²)	28.0	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: ZN, MG, HPA

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	1.54	15/2146 (0.7%)	1.88	54/2898 (1.9%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	1	0

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	243	ASN	CA-C	7.74	1.63	1.52
1	A	243	ASN	C-N	7.71	1.44	1.33
1	A	172	GLN	C-O	-6.16	1.17	1.24
1	A	242	THR	C-N	6.13	1.42	1.33
1	A	253	GLY	CA-C	6.09	1.57	1.52
1	A	74	ASN	CA-C	-5.52	1.45	1.53
1	A	176	SER	CA-C	-5.49	1.45	1.52
1	A	201	GLU	CA-C	-5.49	1.45	1.52
1	A	13	THR	C-O	-5.37	1.17	1.24
1	A	239	SER	CA-C	-5.20	1.46	1.52
1	A	158	ARG	CD-NE	-5.13	1.39	1.46
1	A	261	LEU	N-CA	-5.10	1.40	1.46
1	A	72	ILE	C-O	-5.09	1.18	1.24
1	A	166	TYR	N-CA	-5.09	1.40	1.46
1	A	86	HIS	C-N	-5.09	1.27	1.34

All (54) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	260	VAL	N-CA-C	-13.73	93.21	113.39
1	A	73	LEU	CA-C-N	-11.94	107.03	123.03
1	A	73	LEU	C-N-CA	-11.94	107.03	123.03
1	A	252	GLN	CA-C-N	10.84	130.92	121.82
1	A	252	GLN	C-N-CA	10.84	130.92	121.82
1	A	243	ASN	CA-C-N	10.00	139.69	121.70
1	A	243	ASN	C-N-CA	10.00	139.69	121.70
1	A	31	CYS	CA-C-N	-9.75	113.63	121.82
1	A	31	CYS	C-N-CA	-9.75	113.63	121.82
1	A	245	VAL	N-CA-C	8.90	127.86	109.34
1	A	24	ARG	CD-NE-CZ	7.97	135.55	124.40
1	A	196	GLY	N-CA-C	7.96	124.04	113.37
1	A	260	VAL	CA-C-O	7.51	126.41	119.20
1	A	74	ASN	CA-CB-CG	7.51	120.11	112.60
1	A	24	ARG	NE-CZ-NH1	7.49	128.99	121.50
1	A	259	GLU	CA-C-N	7.24	131.83	121.55
1	A	259	GLU	C-N-CA	7.24	131.83	121.55
1	A	243	ASN	O-C-N	-7.21	113.00	122.59
1	A	243	ASN	N-CA-C	7.07	125.85	110.80
1	A	40	ASN	N-CA-C	-6.63	101.98	110.53
1	A	47	THR	N-CA-CB	-6.48	99.99	111.08
1	A	24	ARG	NE-CZ-NH2	-6.45	113.39	119.20
1	A	20	HIS	CA-CB-CG	-6.35	107.45	113.80
1	A	175	HIS	N-CA-C	-6.33	96.85	107.99
1	A	280	ALA	CA-C-N	-6.29	112.75	122.49
1	A	280	ALA	C-N-CA	-6.29	112.75	122.49
1	A	85	PHE	CA-CB-CG	-6.25	107.55	113.80
1	A	242	THR	CA-C-O	-6.20	112.58	119.34
1	A	39	VAL	N-CA-C	-6.20	96.44	109.34
1	A	176	SER	N-CA-CB	6.09	120.79	110.49
1	A	57	PRO	N-CA-CB	5.99	108.90	103.33
1	A	96	VAL	N-CA-CB	5.96	118.64	110.54
1	A	158	ARG	NE-CZ-NH2	-5.84	113.94	119.20
1	A	175	HIS	CB-CA-C	-5.83	101.39	111.30
1	A	75	GLY	N-CA-C	-5.83	107.20	115.30
1	A	222	VAL	CA-C-O	5.79	123.11	118.71
1	A	243	ASN	CB-CA-C	5.76	121.89	110.42
1	A	78	CYS	CB-CA-C	-5.76	97.98	109.79
1	A	106	LEU	N-CA-C	-5.71	106.15	113.23
1	A	248	ASP	CA-CB-CG	5.71	118.31	112.60
1	A	131	LEU	CB-CA-C	-5.55	100.17	109.65
1	A	18	LEU	O-C-N	-5.49	115.80	122.22
1	A	26	GLN	N-CA-C	-5.41	107.33	114.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	257	HIS	CA-CB-CG	-5.26	108.54	113.80
1	A	176	SER	N-CA-C	-5.25	99.61	110.80
1	A	208	LEU	CA-C-N	5.18	127.22	120.28
1	A	208	LEU	C-N-CA	5.18	127.22	120.28
1	A	259	GLU	N-CA-C	-5.17	99.79	110.80
1	A	79	VAL	N-CA-CB	5.13	119.01	111.52
1	A	78	CYS	CA-CB-SG	-5.11	102.66	114.40
1	A	84	ARG	NE-CZ-NH2	-5.09	114.62	119.20
1	A	170	MET	CA-C-O	5.08	125.78	119.79
1	A	235	VAL	N-CA-C	5.05	115.75	108.48
1	A	84	ARG	CD-NE-CZ	5.01	131.41	124.40

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	243	ASN	CA

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	2102	0	2061	96	0
2	A	1	0	0	0	0
3	A	1	0	0	0	0
4	A	10	0	4	0	0
5	A	129	0	0	11	0
All	All	2243	0	2065	96	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 23.

All (96) 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:35:LEU:HD23	1:A:114:THR:HG23	1.33	1.08
1:A:145:ASN:HD22	1:A:147:LEU:H	1.02	0.98
1:A:38:LEU:HD12	1:A:38:LEU:H	1.34	0.92
1:A:261:LEU:HD11	1:A:265:LYS:NZ	1.85	0.91
1:A:35:LEU:HD23	1:A:114:THR:CG2	2.06	0.84
1:A:145:ASN:ND2	1:A:147:LEU:H	1.77	0.82
1:A:261:LEU:HD11	1:A:265:LYS:HZ2	1.44	0.81
1:A:244:LYS:HE2	1:A:244:LYS:C	2.06	0.80
1:A:88:TYR:CE2	1:A:219:MET:HE3	2.18	0.79
1:A:125:GLU:HG3	5:A:309:HOH:O	1.82	0.78
1:A:267:ALA:HB1	1:A:271:LEU:HD12	1.66	0.77
1:A:152:GLU:OE2	1:A:154:ARG:HB2	1.86	0.76
1:A:170:MET:HE2	1:A:236:PHE:CB	2.15	0.76
1:A:267:ALA:HB1	1:A:271:LEU:CD1	2.16	0.75
1:A:170:MET:HE2	1:A:236:PHE:HB2	1.70	0.73
1:A:35:LEU:C	5:A:333:HOH:O	2.31	0.72
1:A:244:LYS:HB3	1:A:244:LYS:NZ	1.91	0.71
1:A:206:CYS:SG	1:A:244:LYS:HE3	2.31	0.71
1:A:176:SER:N	5:A:427:HOH:O	2.24	0.70
1:A:246:ILE:CD1	1:A:247:MET:HE2	2.22	0.69
1:A:185:ARG:HG3	1:A:186:GLU:N	2.08	0.68
1:A:207:ARG:O	1:A:211:ASN:ND2	2.30	0.65
1:A:261:LEU:HD11	1:A:265:LYS:HZ3	1.63	0.64
1:A:38:LEU:HD12	1:A:38:LEU:N	2.09	0.64
1:A:264:GLY:O	1:A:266:GLN:N	2.31	0.63
1:A:269:GLN:O	1:A:273:GLN:HG2	1.99	0.62
1:A:33:SER:HB2	5:A:324:HOH:O	1.99	0.62
1:A:32:GLY:O	1:A:35:LEU:HD22	1.99	0.62
1:A:262:GLU:HA	1:A:265:LYS:HB2	1.81	0.61
1:A:238:PHE:HE1	1:A:271:LEU:HD21	1.66	0.60
1:A:40:ASN:C	1:A:41:LYS:HG3	2.27	0.60
1:A:39:VAL:HG12	1:A:40:ASN:N	2.17	0.59
1:A:84:ARG:NH2	1:A:114:THR:O	2.30	0.59
1:A:258:GLU:O	1:A:259:GLU:HG3	2.03	0.59
1:A:38:LEU:HD13	1:A:80:MET:HE1	1.84	0.58
1:A:246:ILE:HD11	1:A:247:MET:HE2	1.87	0.57
1:A:76:ARG:NH2	1:A:282:ILE:HB	2.20	0.56
1:A:282:ILE:HD12	1:A:282:ILE:N	2.19	0.56
1:A:145:ASN:HD22	1:A:147:LEU:N	1.87	0.56
1:A:32:GLY:O	1:A:35:LEU:HB2	2.06	0.56
1:A:240:LEU:HG	1:A:271:LEU:HD11	1.88	0.56
1:A:178:TRP:CZ2	1:A:183:GLU:HG3	2.42	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:172:GLN:O	1:A:175:HIS:O	2.25	0.55
1:A:40:ASN:C	1:A:42:LEU:H	2.15	0.54
1:A:88:TYR:HE2	1:A:219:MET:HE3	1.68	0.54
1:A:282:ILE:O	1:A:282:ILE:HG22	2.07	0.54
1:A:21:THR:O	1:A:24:ARG:NH2	2.42	0.53
1:A:73:LEU:HB3	1:A:279:MET:HG3	1.91	0.52
1:A:76:ARG:CG	1:A:76:ARG:HH11	2.22	0.52
1:A:210:ARG:O	1:A:210:ARG:HD2	2.10	0.52
1:A:7:TYR:HE2	1:A:11:GLN:OE1	1.93	0.52
1:A:238:PHE:HE1	1:A:271:LEU:CD2	2.22	0.52
1:A:244:LYS:HA	5:A:423:HOH:O	2.09	0.52
1:A:208:LEU:C	1:A:208:LEU:HD23	2.36	0.51
1:A:271:LEU:O	1:A:275:VAL:HG23	2.11	0.50
1:A:76:ARG:HH21	1:A:282:ILE:HB	1.76	0.50
1:A:265:LYS:O	1:A:266:GLN:HB3	2.11	0.50
1:A:32:GLY:HA3	1:A:84:ARG:HH21	1.78	0.49
1:A:138:LEU:HB2	1:A:194:MET:O	2.12	0.49
1:A:94:TRP:CZ3	1:A:95:LYS:HG3	2.49	0.48
1:A:221:THR:HA	1:A:224:GLU:OE1	2.14	0.48
1:A:282:ILE:N	1:A:282:ILE:CD1	2.77	0.47
1:A:246:ILE:HD12	1:A:246:ILE:C	2.39	0.47
1:A:67:ARG:HD3	1:A:82:GLN:HG2	1.96	0.47
1:A:170:MET:HE2	1:A:236:PHE:HB3	1.93	0.47
1:A:4:GLY:HA3	5:A:377:HOH:O	2.13	0.47
1:A:238:PHE:CE1	1:A:271:LEU:CD2	2.98	0.47
1:A:243:ASN:HB2	5:A:423:HOH:O	2.14	0.46
1:A:267:ALA:O	1:A:268:ALA:C	2.60	0.45
1:A:40:ASN:O	1:A:41:LYS:HG3	2.17	0.45
1:A:10:TYR:HB3	1:A:105:LEU:HD23	1.97	0.45
1:A:145:ASN:HB2	5:A:370:HOH:O	2.17	0.44
1:A:170:MET:SD	1:A:234:ARG:HD2	2.58	0.44
1:A:264:GLY:C	1:A:266:GLN:H	2.26	0.44
1:A:270:LYS:HA	1:A:273:GLN:HE21	1.83	0.44
1:A:80:MET:HB3	1:A:80:MET:HE3	1.63	0.43
1:A:211:ASN:ND2	5:A:338:HOH:O	2.50	0.43
1:A:274:PHE:O	1:A:278:LEU:HB2	2.17	0.43
1:A:136:ILE:HG22	1:A:138:LEU:HD13	2.00	0.43
1:A:76:ARG:CG	1:A:76:ARG:NH1	2.82	0.42
1:A:45:ALA:HA	1:A:70:PHE:O	2.19	0.42
1:A:244:LYS:HE2	1:A:245:VAL:N	2.35	0.42
1:A:154:ARG:HB2	1:A:154:ARG:HE	1.74	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:120:LEU:HD12	1:A:120:LEU:HA	1.90	0.42
1:A:148:ARG:NH1	5:A:420:HOH:O	2.39	0.41
1:A:35:LEU:N	1:A:35:LEU:CD1	2.82	0.41
1:A:73:LEU:HD23	1:A:73:LEU:HA	1.63	0.41
1:A:133:ARG:NE	5:A:413:HOH:O	2.53	0.41
1:A:259:GLU:HA	1:A:262:GLU:HG2	2.00	0.41
1:A:54:PRO:O	1:A:55:ASN:HB2	2.20	0.41
1:A:242:THR:O	1:A:243:ASN:HB3	2.19	0.41
1:A:245:VAL:HG21	1:A:253:GLY:O	2.20	0.41
1:A:114:THR:HG23	1:A:115:ASN:N	2.34	0.41
1:A:266:GLN:O	1:A:269:GLN:HG2	2.20	0.41
1:A:264:GLY:C	1:A:266:GLN:N	2.79	0.40
1:A:88:TYR:CE2	1:A:219:MET:CE	2.97	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	264/281 (94%)	244 (92%)	14 (5%)	6 (2%)	5 1

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	176	SER
1	A	265	LYS
1	A	266	GLN
1	A	33	SER
1	A	36	GLY
1	A	268	ALA

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	165/233 (71%)	125 (76%)	40 (24%)	1 0

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	THR
1	A	24	ARG
1	A	27	VAL
1	A	33	SER
1	A	35	LEU
1	A	38	LEU
1	A	39	VAL
1	A	42	LEU
1	A	43	THR
1	A	56	PHE
1	A	67	ARG
1	A	72	ILE
1	A	73	LEU
1	A	74	ASN
1	A	76	ARG
1	A	80	MET
1	A	82	GLN
1	A	87	MET
1	A	96	VAL
1	A	105	LEU
1	A	176	SER
1	A	179	LYS
1	A	185	ARG
1	A	186	GLU
1	A	187	LEU
1	A	194	MET
1	A	195	LEU
1	A	208	LEU
1	A	212	LEU
1	A	240	LEU

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Mol	Chain	Res	Type
1	A	244	LYS
1	A	251	SER
1	A	254	LYS
1	A	265	LYS
1	A	266	GLN
1	A	269	GLN
1	A	270	LYS
1	A	271	LEU
1	A	272	GLU
1	A	277	LEU

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

Mol	Chain	Res	Type
1	A	23	GLN
1	A	26	GLN
1	A	44	GLN
1	A	46	GLN
1	A	55	ASN
1	A	74	ASN
1	A	86	HIS
1	A	188	GLN
1	A	199	ASN
1	A	211	ASN
1	A	269	GLN
1	A	273	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 oligosaccharides in this entry.

5.6 Ligand geometry

Of 3 ligands modelled in this entry, 2 are monoatomic - leaving 1 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$
4	HPA	A	300	-	10,11,11	1.86	5 (50%)	11,15,15	1.28	1 (9%)

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	HPA	A	300	-	-	-	0/2/2/2

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	300	HPA	C8-N9	3.21	1.40	1.33
4	A	300	HPA	C2-N3	2.57	1.35	1.30
4	A	300	HPA	C8-N7	-2.45	1.30	1.35
4	A	300	HPA	C2-N1	2.29	1.39	1.35
4	A	300	HPA	C5-N7	2.07	1.41	1.38

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	300	HPA	C8-N7-C5	-2.74	103.51	106.59

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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.