



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 02:13 PM UTC

PDB ID : 2MAS / pdb_00002mas
Title : PURINE NUCLEOSIDE HYDROLASE WITH A TRANSITION STATE INHIBITOR
Authors : Degano, M.; Schramm, V.L.; Sacchettini, J.C.
Deposited on : 1996-10-17
Resolution : 2.30 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

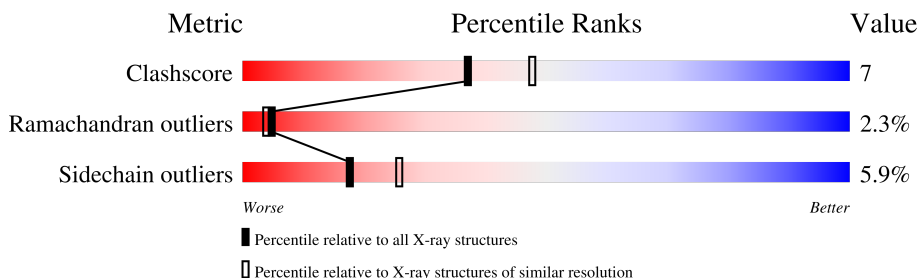
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)

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	314	79% 15% 5% .
1	B	314	76% 18% . .
1	C	314	80% 16% . .
1	D	314	77% 18% . .

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9774 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.

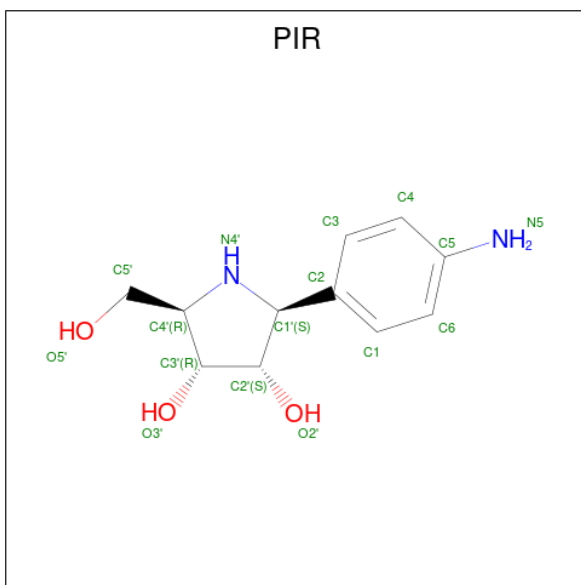
- Molecule 1 is a protein called INOSINE-URIDINE NUCLEOSIDE N-RIBOHYDROLASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	313	Total	C	N	O	S	0	0	0
			2387	1517	409	446	15			
1	B	313	Total	C	N	O	S	0	0	0
			2387	1517	409	446	15			
1	C	313	Total	C	N	O	S	0	0	0
			2387	1517	409	446	15			
1	D	313	Total	C	N	O	S	0	0	0
			2387	1517	409	446	15			

- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Ca	0	0
			1	1		
2	B	1	Total	Ca	0	0
			1	1		
2	C	1	Total	Ca	0	0
			1	1		
2	D	1	Total	Ca	0	0
			1	1		

- Molecule 3 is 2-(4-AMINO-PHENYL)-5-HYDROXYMETHYL-PYRROLIDINE-3,4-DIOL (CCD ID: PIR) (formula: C₁₁H₁₆N₂O₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			16	11	2	3		
3	B	1	Total	C	N	O	0	0
			16	11	2	3		
3	C	1	Total	C	N	O	0	0
			16	11	2	3		
3	D	1	Total	C	N	O	0	0
			16	11	2	3		

- Molecule 4 is water.

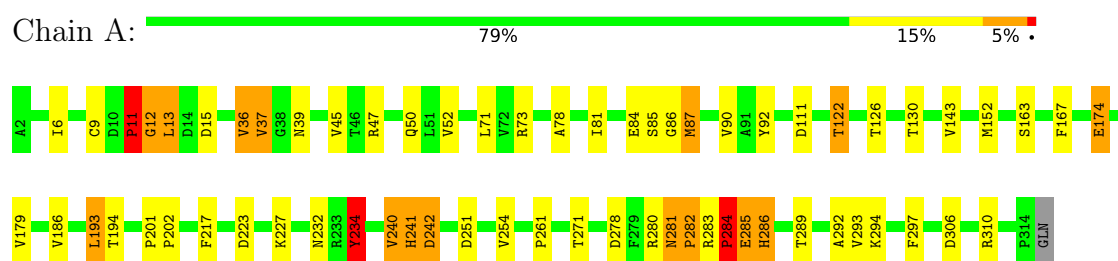
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	36	Total	O	0	0
			36	36		
4	B	30	Total	O	0	0
			30	30		
4	C	51	Total	O	0	0
			51	51		
4	D	41	Total	O	0	0
			41	41		

3 Residue-property plots [i](#)

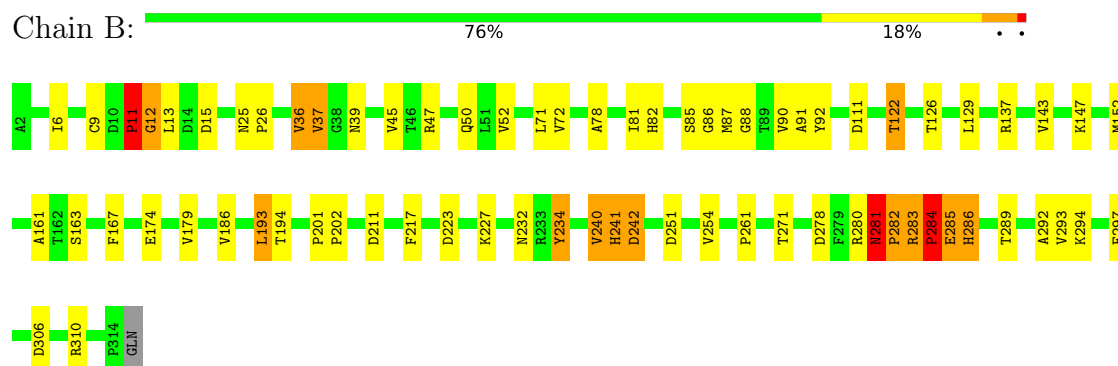
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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.

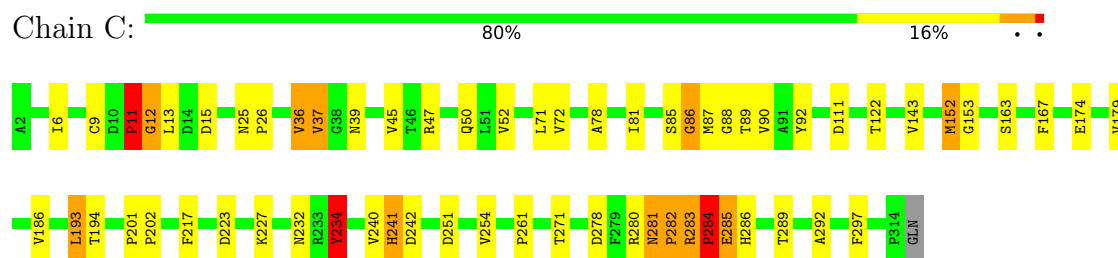
• Molecule 1: INOSINE-URIDINE NUCLEOSIDE N-RIBOHYDROLASE



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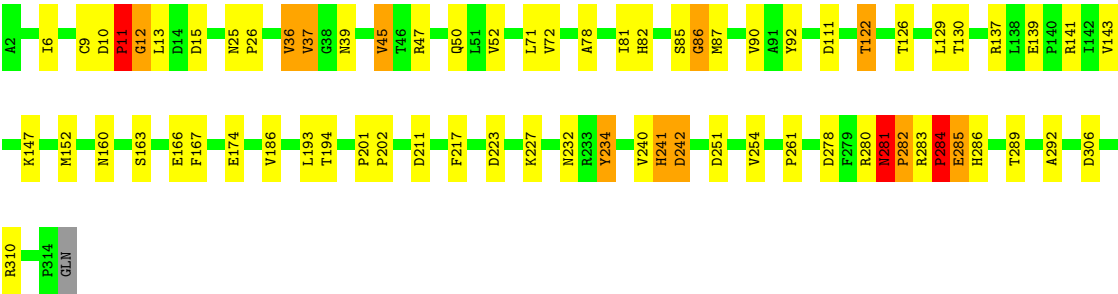


• Molecule 1: INOSINE-URIDINE NUCLEOSIDE N-RIBOHYDROLASE



• Molecule 1: INOSINE-URIDINE NUCLEOSIDE N-RIBOHYDROLASE





4 Data and refinement statistics

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

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	120.63Å 159.76Å 206.38Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	(Not available) – 2.30	Depositor
% Data completeness (in resolution range)	80.8 ((Not available)-2.30)	Depositor
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.205 , 0.234	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	9774	wwPDB-VP
Average B, all atoms (Å ²)	34.0	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: CA, PIR

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	0.79	1/2433 (0.0%)	1.56	19/3317 (0.6%)
1	B	0.87	0/2433	1.58	21/3317 (0.6%)
1	C	0.84	0/2433	1.55	20/3317 (0.6%)
1	D	0.78	0/2433	1.56	19/3317 (0.6%)
All	All	0.82	1/9732 (0.0%)	1.56	79/13268 (0.6%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	87	MET	SD-CE	-7.18	1.61	1.79

All (79) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	281	ASN	CA-C-N	42.82	173.37	119.84
1	A	281	ASN	C-N-CA	42.82	173.37	119.84
1	B	281	ASN	CA-C-N	42.40	172.84	119.84
1	B	281	ASN	C-N-CA	42.40	172.84	119.84
1	D	281	ASN	CA-C-N	42.38	172.81	119.84
1	D	281	ASN	C-N-CA	42.38	172.81	119.84
1	C	281	ASN	CA-C-N	42.15	172.52	119.84
1	C	281	ASN	C-N-CA	42.15	172.52	119.84
1	A	281	ASN	C-N-CD	-11.99	75.83	125.00
1	C	281	ASN	C-N-CD	-11.93	76.11	125.00
1	D	281	ASN	C-N-CD	-11.91	76.18	125.00
1	B	281	ASN	C-N-CD	-11.90	76.21	125.00
1	B	37	VAL	CB-CA-C	-11.11	99.01	111.45
1	A	37	VAL	CB-CA-C	-10.84	99.31	111.45
1	C	37	VAL	CB-CA-C	-10.31	99.90	111.45
1	B	241	HIS	N-CA-C	9.88	121.74	110.97

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	241	HIS	N-CA-C	9.66	121.50	110.97
1	C	241	HIS	N-CA-C	9.54	121.37	110.97
1	D	241	HIS	N-CA-C	9.11	120.89	110.97
1	D	37	VAL	CB-CA-C	-8.77	99.18	111.28
1	C	283	ARG	CA-C-N	8.50	130.46	119.84
1	C	283	ARG	C-N-CA	8.50	130.46	119.84
1	A	283	ARG	CA-C-N	8.41	130.35	119.84
1	A	283	ARG	C-N-CA	8.41	130.35	119.84
1	B	283	ARG	CA-C-N	8.36	130.29	119.84
1	B	283	ARG	C-N-CA	8.36	130.29	119.84
1	D	283	ARG	CA-C-N	8.14	130.02	119.84
1	D	283	ARG	C-N-CA	8.14	130.02	119.84
1	C	281	ASN	N-CA-C	7.28	120.18	109.62
1	D	36	VAL	CB-CA-C	-7.12	99.34	110.52
1	B	281	ASN	N-CA-C	6.98	119.75	109.62
1	D	12	GLY	N-CA-C	-6.82	97.02	113.18
1	B	12	GLY	N-CA-C	-6.71	97.27	113.18
1	A	281	ASN	N-CA-C	6.64	119.25	109.62
1	B	282	PRO	CA-N-CD	-6.63	102.72	112.00
1	D	281	ASN	N-CA-C	6.61	119.21	109.62
1	A	12	GLY	N-CA-C	-6.50	97.78	113.18
1	D	282	PRO	CA-N-CD	-6.47	102.94	112.00
1	D	242	ASP	N-CA-C	6.45	122.41	113.45
1	C	282	PRO	CA-N-CD	-6.39	103.05	112.00
1	A	282	PRO	CA-N-CD	-6.36	103.10	112.00
1	C	36	VAL	CB-CA-C	-6.30	99.65	110.30
1	D	234	TYR	N-CA-C	6.29	124.19	110.80
1	C	12	GLY	N-CA-C	-6.25	98.38	113.18
1	A	234	TYR	N-CA-C	6.21	124.04	110.80
1	A	36	VAL	CB-CA-C	-6.18	99.86	110.30
1	B	36	VAL	CB-CA-C	-6.17	99.88	110.30
1	C	271	THR	N-CA-C	6.09	119.66	112.72
1	C	234	TYR	N-CA-C	6.08	123.76	110.80
1	B	234	TYR	N-CA-C	5.99	123.55	110.80
1	B	271	THR	N-CA-C	5.95	120.24	112.92
1	D	129	LEU	N-CA-C	5.94	120.22	113.15
1	B	6	ILE	N-CA-C	-5.86	99.91	108.11
1	D	130	THR	N-CA-C	5.77	118.03	111.11
1	A	297	PHE	N-CA-C	5.71	118.24	111.33
1	B	129	LEU	N-CA-C	5.69	119.30	112.93
1	B	72	VAL	CB-CA-C	-5.62	105.41	110.91
1	C	242	ASP	N-CA-C	5.61	121.24	113.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	6	ILE	N-CA-C	-5.60	99.57	107.75
1	D	6	ILE	N-CA-C	-5.59	99.59	107.75
1	C	297	PHE	N-CA-C	5.56	117.78	111.11
1	B	242	ASP	N-CA-C	5.49	121.08	113.45
1	A	130	THR	N-CA-C	5.40	117.59	111.11
1	D	211	ASP	N-CA-C	5.40	118.64	111.30
1	A	6	ILE	N-CA-C	-5.38	99.89	107.75
1	A	242	ASP	N-CA-C	5.33	120.85	113.45
1	A	271	THR	N-CA-C	5.32	118.79	112.72
1	C	72	VAL	CB-CA-C	-5.26	105.75	110.91
1	C	179	VAL	N-CA-C	5.25	115.46	110.42
1	A	13	LEU	N-CA-C	5.16	117.32	111.02
1	D	86	GLY	N-CA-C	-5.15	108.95	115.08
1	C	86	GLY	N-CA-C	-5.14	108.96	115.08
1	B	179	VAL	N-CA-C	5.11	115.33	110.42
1	B	297	PHE	N-CA-C	5.09	117.49	111.33
1	B	161	ALA	N-CA-C	-5.07	105.83	111.36
1	D	72	VAL	N-CA-C	5.04	116.36	112.12
1	B	211	ASP	N-CA-C	5.04	118.27	111.17
1	C	72	VAL	N-CA-C	5.01	116.33	112.12
1	A	179	VAL	N-CA-C	5.00	115.22	110.42

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	2387	0	2424	33	0
1	B	2387	0	2424	40	1
1	C	2387	0	2424	31	1
1	D	2387	0	2424	37	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	16	0	14	3	0
3	B	16	0	15	3	0
3	C	16	0	15	3	0
3	D	16	0	14	3	0
4	A	36	0	0	2	0
4	B	30	0	0	0	0
4	C	51	0	0	0	0
4	D	41	0	0	1	0
All	All	9774	0	9754	137	2

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 7.

All (137) 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:B:87:MET:HE3	1:B:87:MET:HA	1.56	0.88
1:C:87:MET:HE3	1:C:87:MET:HA	1.56	0.85
1:A:87:MET:HA	1:A:87:MET:HE3	1.63	0.80
1:D:87:MET:HE3	1:D:87:MET:HA	1.66	0.78
1:C:194:THR:OG1	1:C:241:HIS:HD2	1.71	0.74
1:A:194:THR:OG1	1:A:241:HIS:HD2	1.74	0.69
1:B:194:THR:OG1	1:B:241:HIS:HD2	1.77	0.68
1:D:194:THR:OG1	1:D:241:HIS:HD2	1.76	0.68
1:B:12:GLY:HA2	1:B:86:GLY:O	1.96	0.65
1:D:9:CYS:SG	1:D:11:PRO:HD3	2.39	0.62
1:B:201:PRO:HB2	1:B:202:PRO:HD3	1.83	0.61
1:A:285:GLU:O	1:A:286:HIS:HB2	2.01	0.61
1:A:12:GLY:HA2	1:A:86:GLY:O	2.01	0.60
1:C:285:GLU:O	1:C:286:HIS:HB2	2.01	0.60
1:D:285:GLU:O	1:D:286:HIS:HB2	2.00	0.60
1:A:193:LEU:HD13	1:A:292:ALA:HB1	1.83	0.60
1:B:285:GLU:O	1:B:286:HIS:HB2	2.00	0.59
1:B:9:CYS:SG	1:B:11:PRO:HD3	2.42	0.59
1:B:85:SER:HB2	1:B:92:TYR:OH	2.02	0.59
1:C:201:PRO:HB2	1:C:202:PRO:HD3	1.85	0.59
1:B:11:PRO:HA	1:B:15:ASP:HB2	1.84	0.59
1:C:11:PRO:HA	1:C:15:ASP:HB2	1.85	0.59
1:D:11:PRO:HA	1:D:15:ASP:HB2	1.85	0.58
1:D:122:THR:HG23	4:D:416:HOH:O	2.03	0.58
1:C:12:GLY:HA2	1:C:86:GLY:O	2.02	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:193:LEU:HD13	1:D:292:ALA:HB1	1.86	0.58
1:D:284:PRO:O	1:D:285:GLU:HB3	2.04	0.58
1:A:84:GLU:HG3	4:A:434:HOH:O	2.05	0.57
1:C:193:LEU:HD13	1:C:292:ALA:HB1	1.85	0.57
1:C:284:PRO:O	1:C:285:GLU:HB3	2.05	0.57
1:D:251:ASP:O	1:D:254:VAL:HG22	2.04	0.57
1:B:193:LEU:HD13	1:B:292:ALA:HB1	1.87	0.56
1:B:278:ASP:OD1	1:B:280:ARG:HD3	2.05	0.56
1:A:284:PRO:O	1:A:285:GLU:HB3	2.04	0.56
1:D:12:GLY:HA2	1:D:86:GLY:O	2.06	0.55
1:A:280:ARG:HD2	1:D:163:SER:HA	1.88	0.55
1:A:78:ALA:HB2	1:A:167:PHE:CE2	2.42	0.55
1:B:281:ASN:O	1:C:234:TYR:HB2	2.07	0.54
1:A:278:ASP:OD1	1:A:280:ARG:HD3	2.08	0.54
1:C:278:ASP:OD1	1:C:280:ARG:HD3	2.06	0.54
1:B:81:ILE:HG13	3:B:400:PIR:N5	2.23	0.54
1:B:284:PRO:O	1:B:285:GLU:HB3	2.07	0.54
1:A:201:PRO:HB2	1:A:202:PRO:HD3	1.89	0.53
1:C:9:CYS:SG	1:C:11:PRO:HD3	2.49	0.53
1:D:261:PRO:CG	1:D:284:PRO:HG2	2.39	0.53
1:B:261:PRO:CG	1:B:284:PRO:HG2	2.38	0.53
1:A:11:PRO:HA	1:A:15:ASP:HB2	1.91	0.53
1:A:90:VAL:HG11	1:A:217:PHE:HE1	1.73	0.53
1:D:90:VAL:HG11	1:D:217:PHE:HE1	1.75	0.52
1:D:201:PRO:HB2	1:D:202:PRO:HD3	1.91	0.52
1:C:85:SER:HB2	1:C:92:TYR:OH	2.09	0.52
1:C:261:PRO:CG	1:C:284:PRO:HG2	2.40	0.52
1:B:47:ARG:HA	1:B:50:GLN:HE21	1.74	0.52
1:C:90:VAL:HG11	1:C:217:PHE:HE1	1.75	0.52
1:A:39:ASN:HD21	3:A:400:PIR:H1'	1.73	0.51
1:A:9:CYS:SG	1:A:11:PRO:HD3	2.51	0.51
1:B:251:ASP:O	1:B:254:VAL:HG22	2.10	0.51
1:A:261:PRO:CG	1:A:284:PRO:HG2	2.40	0.51
1:D:85:SER:HB2	1:D:92:TYR:OH	2.11	0.51
1:A:85:SER:HB2	1:A:92:TYR:OH	2.11	0.51
1:A:186:VAL:HB	1:A:289:THR:HG22	1.92	0.51
1:A:251:ASP:O	1:A:254:VAL:HG22	2.11	0.51
1:D:278:ASP:OD1	1:D:280:ARG:HD3	2.11	0.51
1:D:223:ASP:O	1:D:227:LYS:HD3	2.12	0.50
1:B:280:ARG:CD	1:C:163:SER:HA	2.42	0.50
1:D:47:ARG:HA	1:D:50:GLN:HE21	1.76	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:90:VAL:HG11	1:B:217:PHE:HE1	1.76	0.49
1:A:223:ASP:O	1:A:227:LYS:HD3	2.13	0.49
1:A:234:TYR:HB2	1:D:281:ASN:O	2.13	0.49
1:C:223:ASP:O	1:C:227:LYS:HD3	2.13	0.49
1:B:306:ASP:O	1:B:310:ARG:HG3	2.13	0.48
1:A:122:THR:HG21	4:A:401:HOH:O	2.13	0.48
1:C:78:ALA:HB2	1:C:167:PHE:CE2	2.48	0.48
1:C:81:ILE:HG13	3:C:400:PIR:N5	2.28	0.48
1:C:251:ASP:O	1:C:254:VAL:HG22	2.12	0.48
1:A:280:ARG:CD	1:D:163:SER:HA	2.43	0.48
1:B:280:ARG:HD2	1:C:163:SER:HA	1.95	0.48
1:B:39:ASN:HD21	3:B:400:PIR:H1'	1.79	0.47
1:B:223:ASP:O	1:B:227:LYS:HD3	2.14	0.47
1:B:25:ASN:HA	1:B:26:PRO:HD3	1.78	0.47
1:D:122:THR:HG22	1:D:147:LYS:HB3	1.96	0.47
1:A:306:ASP:O	1:A:310:ARG:HG3	2.14	0.47
1:C:39:ASN:HD21	3:C:400:PIR:H1'	1.80	0.47
1:D:39:ASN:HD21	3:D:400:PIR:H1'	1.80	0.47
1:D:78:ALA:HB2	1:D:167:PHE:CE2	2.50	0.47
1:C:47:ARG:HA	1:C:50:GLN:HE21	1.80	0.47
1:B:78:ALA:HB2	1:B:167:PHE:CE2	2.50	0.47
1:B:13:LEU:HD23	1:B:87:MET:CE	2.45	0.46
1:A:47:ARG:HA	1:A:50:GLN:HE21	1.80	0.46
1:B:13:LEU:HD23	1:B:87:MET:HE3	1.98	0.46
1:A:81:ILE:HG13	3:A:400:PIR:N5	2.30	0.46
1:D:306:ASP:O	1:D:310:ARG:HG3	2.15	0.46
1:B:82:HIS:HB3	1:B:86:GLY:O	2.16	0.46
1:D:139:GLU:OE2	1:D:141:ARG:HB2	2.16	0.46
1:D:25:ASN:HA	1:D:26:PRO:HD3	1.79	0.46
1:B:87:MET:HG2	1:B:90:VAL:HG21	1.98	0.45
1:B:13:LEU:HA	1:B:87:MET:HE1	1.99	0.45
1:C:152:MET:HE2	1:C:153:GLY:N	2.32	0.45
1:A:13:LEU:HD23	1:A:87:MET:CE	2.46	0.45
1:B:293:VAL:HG12	1:B:294:LYS:HD2	1.97	0.45
1:C:47:ARG:HH21	1:C:92:TYR:HB2	1.81	0.45
1:D:186:VAL:HB	1:D:289:THR:HG22	1.98	0.45
3:C:400:PIR:H1	3:C:400:PIR:H2'	1.75	0.44
1:D:10:ASP:HA	1:D:45:VAL:HG11	1.98	0.44
1:B:87:MET:HA	1:B:87:MET:CE	2.36	0.44
1:D:81:ILE:HG13	3:D:400:PIR:N5	2.32	0.44
1:C:186:VAL:HB	1:C:289:THR:HG22	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:9:CYS:C	1:D:11:PRO:HD3	2.43	0.43
1:B:186:VAL:HB	1:B:289:THR:HG22	2.00	0.43
1:B:227:LYS:HD2	1:B:227:LYS:N	2.33	0.43
1:B:240:VAL:O	1:B:240:VAL:CG2	2.67	0.43
1:C:25:ASN:HA	1:C:26:PRO:HD3	1.79	0.43
1:D:13:LEU:HD23	1:D:87:MET:CE	2.49	0.43
1:D:87:MET:HG2	1:D:90:VAL:HG21	2.00	0.43
1:B:47:ARG:HH21	1:B:92:TYR:HB2	1.84	0.43
1:C:13:LEU:HD23	1:C:87:MET:CE	2.48	0.43
1:C:283:ARG:HA	1:C:284:PRO:HD3	1.90	0.43
1:B:9:CYS:C	1:B:11:PRO:HD3	2.44	0.42
1:A:293:VAL:HG12	1:A:294:LYS:HD2	2.02	0.42
1:A:240:VAL:O	1:A:240:VAL:CG2	2.68	0.42
1:D:82:HIS:HB3	1:D:86:GLY:O	2.19	0.42
3:D:400:PIR:H1	3:D:400:PIR:H2'	1.73	0.42
1:A:73:ARG:HH11	1:A:174:GLU:CD	2.27	0.42
1:D:285:GLU:O	1:D:286:HIS:CB	2.68	0.42
1:A:163:SER:HA	1:D:280:ARG:HD2	2.02	0.41
1:C:13:LEU:HA	1:C:87:MET:HE1	2.01	0.41
1:B:283:ARG:HA	1:B:284:PRO:HD3	1.90	0.41
1:B:122:THR:HG22	1:B:147:LYS:HB3	2.03	0.41
1:D:160:ASN:OD1	1:D:166:GLU:HA	2.21	0.41
1:A:126:THR:HB	1:A:242:ASP:CG	2.46	0.41
3:A:400:PIR:H2'	3:A:400:PIR:H1	1.75	0.41
1:C:285:GLU:O	1:C:286:HIS:CB	2.69	0.41
1:B:163:SER:HA	1:C:280:ARG:HD2	2.02	0.40
3:B:400:PIR:H2'	3:B:400:PIR:H1	1.71	0.40
1:B:126:THR:HB	1:B:242:ASP:CG	2.45	0.40
1:A:227:LYS:N	1:A:227:LYS:HD2	2.35	0.40
1:D:126:THR:HB	1:D:242:ASP:CG	2.46	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:89:THR:O	1:C:89:THR:O[2_655]	2.11	0.09
1:B:91:ALA:N	1:B:91:ALA:N[2_755]	2.13	0.07

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	311/314 (99%)	292 (94%)	12 (4%)	7 (2%)	5	4
1	B	311/314 (99%)	290 (93%)	13 (4%)	8 (3%)	4	3
1	C	311/314 (99%)	291 (94%)	13 (4%)	7 (2%)	5	4
1	D	311/314 (99%)	294 (94%)	11 (4%)	6 (2%)	6	5
All	All	1244/1256 (99%)	1167 (94%)	49 (4%)	28 (2%)	5	4

All (28) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	282	PRO
1	B	282	PRO
1	C	282	PRO
1	D	282	PRO
1	A	232	ASN
1	B	232	ASN
1	C	232	ASN
1	D	232	ASN
1	A	11	PRO
1	B	11	PRO
1	D	11	PRO
1	B	285	GLU
1	C	11	PRO
1	C	285	GLU
1	A	234	TYR
1	A	284	PRO
1	A	285	GLU
1	A	286	HIS
1	B	88	GLY
1	B	234	TYR
1	B	284	PRO
1	B	286	HIS

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Mol	Chain	Res	Type
1	C	88	GLY
1	C	284	PRO
1	D	234	TYR
1	D	284	PRO
1	D	285	GLU
1	C	234	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	258/259 (100%)	243 (94%)	15 (6%)	18	26
1	B	258/259 (100%)	242 (94%)	16 (6%)	16	24
1	C	258/259 (100%)	243 (94%)	15 (6%)	18	26
1	D	258/259 (100%)	243 (94%)	15 (6%)	18	26
All	All	1032/1036 (100%)	971 (94%)	61 (6%)	18	26

All (61) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	PRO
1	A	36	VAL
1	A	37	VAL
1	A	45	VAL
1	A	52	VAL
1	A	71	LEU
1	A	111	ASP
1	A	122	THR
1	A	143	VAL
1	A	152	MET
1	A	174	GLU
1	A	193	LEU
1	A	240	VAL
1	A	281	ASN
1	A	284	PRO

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Mol	Chain	Res	Type
1	B	11	PRO
1	B	36	VAL
1	B	37	VAL
1	B	45	VAL
1	B	52	VAL
1	B	71	LEU
1	B	111	ASP
1	B	122	THR
1	B	137	ARG
1	B	143	VAL
1	B	152	MET
1	B	174	GLU
1	B	193	LEU
1	B	240	VAL
1	B	281	ASN
1	B	284	PRO
1	C	11	PRO
1	C	36	VAL
1	C	37	VAL
1	C	45	VAL
1	C	52	VAL
1	C	71	LEU
1	C	111	ASP
1	C	122	THR
1	C	143	VAL
1	C	152	MET
1	C	174	GLU
1	C	193	LEU
1	C	240	VAL
1	C	281	ASN
1	C	284	PRO
1	D	11	PRO
1	D	36	VAL
1	D	37	VAL
1	D	45	VAL
1	D	52	VAL
1	D	71	LEU
1	D	111	ASP
1	D	122	THR
1	D	137	ARG
1	D	143	VAL
1	D	152	MET

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Mol	Chain	Res	Type
1	D	174	GLU
1	D	240	VAL
1	D	281	ASN
1	D	284	PRO

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	23	HIS
1	A	39	ASN
1	A	50	GLN
1	A	157	HIS
1	A	181	ASN
1	A	241	HIS
1	B	23	HIS
1	B	39	ASN
1	B	50	GLN
1	B	157	HIS
1	B	181	ASN
1	B	241	HIS
1	C	23	HIS
1	C	39	ASN
1	C	50	GLN
1	C	157	HIS
1	C	181	ASN
1	C	241	HIS
1	D	23	HIS
1	D	39	ASN
1	D	50	GLN
1	D	157	HIS
1	D	181	ASN
1	D	241	HIS

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](#)

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 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$
3	PIR	D	400	2	17,17,17	1.04	1 (5%)	20,24,24	0.94	0
3	PIR	B	400	2	17,17,17	0.91	1 (5%)	20,24,24	0.99	1 (5%)
3	PIR	C	400	2	17,17,17	0.98	1 (5%)	20,24,24	0.96	1 (5%)
3	PIR	A	400	2	17,17,17	1.19	2 (11%)	20,24,24	0.99	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
3	PIR	D	400	2	-	5/6/22/22	0/2/2/2
3	PIR	B	400	2	-	6/6/22/22	0/2/2/2
3	PIR	C	400	2	-	5/6/22/22	0/2/2/2
3	PIR	A	400	2	-	6/6/22/22	0/2/2/2

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	400	PIR	C1'-N4'	-2.47	1.44	1.47
3	D	400	PIR	C1-C2	2.25	1.42	1.39
3	A	400	PIR	C3'-C4'	-2.23	1.51	1.53
3	B	400	PIR	C6-C1	2.16	1.42	1.38
3	A	400	PIR	C1-C2	2.00	1.42	1.39

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	400	PIR	C2-C1'-N4'	2.19	116.46	112.98
3	C	400	PIR	C4-C3-C2	-2.13	119.06	121.18

There are no chirality outliers.

All (22) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	400	PIR	C2'-C1'-C2-C1
3	A	400	PIR	C2'-C1'-C2-C3
3	B	400	PIR	C2'-C1'-C2-C1
3	B	400	PIR	C2'-C1'-C2-C3
3	C	400	PIR	C2'-C1'-C2-C1
3	C	400	PIR	C2'-C1'-C2-C3
3	D	400	PIR	C2'-C1'-C2-C1
3	D	400	PIR	C2'-C1'-C2-C3
3	C	400	PIR	N4'-C1'-C2-C1
3	B	400	PIR	N4'-C1'-C2-C1
3	A	400	PIR	N4'-C1'-C2-C1
3	D	400	PIR	N4'-C1'-C2-C1
3	B	400	PIR	N4'-C1'-C2-C3
3	A	400	PIR	N4'-C1'-C2-C3
3	C	400	PIR	N4'-C1'-C2-C3
3	D	400	PIR	N4'-C1'-C2-C3
3	A	400	PIR	N4'-C4'-C5'-O5'
3	B	400	PIR	N4'-C4'-C5'-O5'
3	A	400	PIR	C3'-C4'-C5'-O5'
3	B	400	PIR	C3'-C4'-C5'-O5'
3	C	400	PIR	N4'-C4'-C5'-O5'
3	D	400	PIR	N4'-C4'-C5'-O5'

There are no ring outliers.

4 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	400	PIR	3	0
3	B	400	PIR	3	0
3	C	400	PIR	3	0
3	A	400	PIR	3	0

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.