



Full wwPDB X-ray Structure Validation Report ⓘ

Jul 26, 2023 – 12:16 AM EDT

PDB ID : 1AE5
Title : HUMAN HEPARIN BINDING PROTEIN
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Deposited on : 1997-03-05
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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.34
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) : 2.34

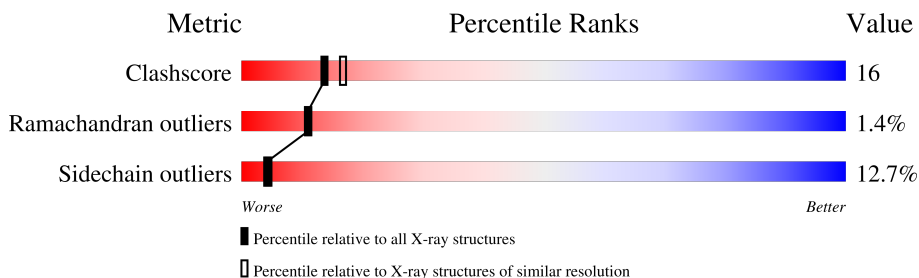
1 Overall quality at a glance

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	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (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 $\leq 5\%$

Mol	Chain	Length	Quality of chain
1	A	225	 60% 30% 8% ..

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 1778 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 HEPARIN BINDING PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	223	1687	1043	323	310	11	0	0	0

- Molecule 2 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
2	A	1	14	8	1	5	0	0
2	A	1	14	8	1	5	0	0

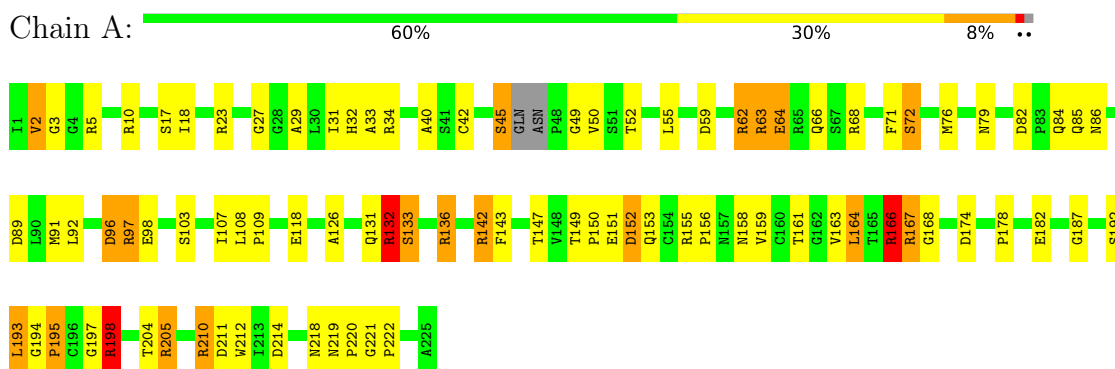
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	63	Total 63 O 63	0	0

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.

- Molecule 1: HEPARIN BINDING PROTEIN



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	39.19Å 66.12Å 101.36Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.30 19.38 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.2 (20.00-2.30) 99.3 (19.38-2.30)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.55 (at 2.30Å)	Xtrriage
Refinement program	TNT	Depositor
R, R_{free}	(Not available) , (Not available) 0.180 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	31.3	Xtrriage
Anisotropy	0.479	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 84.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	1778	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.99% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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: NAG

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	1.04	4/1721 (0.2%)	1.45	25/2329 (1.1%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	151	GLU	CD-OE2	7.09	1.33	1.25
1	A	182	GLU	CD-OE1	6.84	1.33	1.25
1	A	98	GLU	CD-OE2	6.40	1.32	1.25
1	A	118	GLU	CD-OE1	5.13	1.31	1.25

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	221	GLY	C-N-CD	-10.53	97.44	120.60
1	A	59	ASP	CB-CG-OD2	-9.80	109.48	118.30
1	A	59	ASP	CB-CG-OD1	9.13	126.52	118.30
1	A	166	ARG	NE-CZ-NH1	8.82	124.71	120.30
1	A	142	ARG	NE-CZ-NH2	-8.67	115.96	120.30
1	A	210	ARG	NE-CZ-NH1	8.51	124.56	120.30
1	A	89	ASP	CB-CG-OD2	-7.92	111.18	118.30
1	A	210	ARG	NE-CZ-NH2	-7.26	116.67	120.30
1	A	211	ASP	CB-CG-OD1	6.56	124.20	118.30
1	A	96	ASP	CB-CG-OD2	6.50	124.15	118.30
1	A	132	ARG	NE-CZ-NH1	6.33	123.47	120.30
1	A	205	ARG	NE-CZ-NH1	6.25	123.42	120.30
1	A	89	ASP	CB-CG-OD1	6.05	123.75	118.30
1	A	96	ASP	CB-CG-OD1	-6.03	112.87	118.30
1	A	152	ASP	CB-CG-OD2	-6.03	112.88	118.30
1	A	174	ASP	CB-CG-OD2	-5.96	112.94	118.30
1	A	211	ASP	CB-CG-OD2	-5.78	113.10	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	5	ARG	NE-CZ-NH2	-5.76	117.42	120.30
1	A	198	ARG	NE-CZ-NH2	-5.73	117.43	120.30
1	A	5	ARG	NE-CZ-NH1	5.73	123.17	120.30
1	A	152	ASP	CB-CG-OD1	5.55	123.29	118.30
1	A	142	ARG	NE-CZ-NH1	5.52	123.06	120.30
1	A	42	CYS	CA-CB-SG	-5.28	104.50	114.00
1	A	45	SER	N-CA-CB	-5.25	102.63	110.50
1	A	214	ASP	CB-CG-OD2	-5.16	113.66	118.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	1687	0	1655	54	0
2	A	28	0	26	1	0
3	A	63	0	0	3	0
All	All	1778	0	1681	54	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 16.

All (54) 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:193:LEU:H	1:A:193:LEU:HD12	1.36	0.89
1:A:32:HIS:CD2	1:A:34:ARG:H	1.96	0.83
1:A:96:ASP:OD1	1:A:97:ARG:HD3	1.84	0.77
1:A:32:HIS:HD2	1:A:34:ARG:H	1.33	0.74
1:A:18:ILE:HD13	1:A:18:ILE:N	2.06	0.71
1:A:91:MET:HG3	1:A:92:LEU:N	2.06	0.68
1:A:108:LEU:HD12	1:A:109:PRO:HD2	1.79	0.65
1:A:156:PRO:HD2	3:A:440:HOH:O	1.98	0.62
1:A:62:ARG:HB3	1:A:64:GLU:HG2	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:193:LEU:HD12	1:A:193:LEU:N	2.06	0.60
1:A:55:LEU:HD12	1:A:71:PHE:CE1	2.38	0.59
1:A:219:ASN:N	1:A:220:PRO:HD3	2.20	0.57
1:A:27:GLY:HA3	3:A:423:HOH:O	2.04	0.56
1:A:29:ALA:HB2	1:A:178:PRO:HG3	1.88	0.55
1:A:62:ARG:HG3	3:A:421:HOH:O	2.07	0.54
1:A:2:VAL:O	1:A:168:GLY:HA2	2.07	0.53
1:A:82:ASP:O	1:A:86:ASN:N	2.41	0.53
1:A:55:LEU:HD12	1:A:71:PHE:HE1	1.73	0.52
1:A:64:GLU:OE2	1:A:68:ARG:NE	2.39	0.51
1:A:72:SER:N	1:A:96:ASP:OD2	2.29	0.51
1:A:79:ASN:HB2	1:A:212:TRP:CE2	2.46	0.51
1:A:63:ARG:O	1:A:64:GLU:C	2.48	0.51
1:A:166:ARG:HG3	1:A:166:ARG:NH1	2.26	0.50
1:A:152:ASP:OD2	1:A:153:GLN:OE1	2.29	0.50
1:A:40:ALA:HA	1:A:91:MET:HE1	1.93	0.49
1:A:143:PHE:CE2	2:A:411:NAG:H82	2.47	0.49
1:A:155:ARG:HE	1:A:158:ASN:ND2	2.10	0.49
1:A:126:ALA:HA	1:A:142:ARG:O	2.12	0.49
1:A:198:ARG:HG2	1:A:198:ARG:NH1	2.28	0.48
1:A:149:THR:HB	1:A:150:PRO:HD2	1.95	0.47
1:A:187:GLY:HA2	1:A:204:THR:O	2.15	0.47
1:A:218:ASN:C	1:A:220:PRO:HD3	2.36	0.47
1:A:31:ILE:C	1:A:107:ILE:HG23	2.36	0.46
1:A:198:ARG:HG2	1:A:198:ARG:HH11	1.79	0.46
1:A:17:SER:C	1:A:18:ILE:HD13	2.36	0.46
1:A:163:VAL:HG11	1:A:168:GLY:HA3	1.97	0.45
1:A:31:ILE:HA	1:A:107:ILE:HG23	1.98	0.45
1:A:167:ARG:HD2	1:A:197:GLY:C	2.37	0.44
1:A:2:VAL:HG23	1:A:3:GLY:N	2.33	0.43
1:A:32:HIS:HD2	1:A:34:ARG:N	2.10	0.43
1:A:49:GLY:CA	1:A:76:MET:HE3	2.48	0.43
1:A:167:ARG:HD2	1:A:167:ARG:HA	1.66	0.43
1:A:163:VAL:O	1:A:163:VAL:HG23	2.20	0.42
1:A:132:ARG:O	1:A:133:SER:C	2.57	0.42
1:A:194:GLY:O	1:A:195:PRO:C	2.57	0.42
1:A:85:GLN:O	1:A:86:ASN:HB2	2.20	0.41
1:A:147:THR:O	1:A:161:THR:HA	2.19	0.41
1:A:164:LEU:HA	1:A:164:LEU:HD12	1.43	0.41
1:A:167:ARG:HD2	1:A:197:GLY:HA3	2.02	0.41
1:A:40:ALA:HA	1:A:91:MET:CE	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:136:ARG:HH11	1:A:136:ARG:HD2	1.68	0.41
1:A:159:VAL:CG2	1:A:205:ARG:HG3	2.51	0.41
1:A:10:ARG:HD2	1:A:103:SER:O	2.20	0.40
1:A:168:GLY:H	1:A:197:GLY:HA2	1.87	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	219/225 (97%)	203 (93%)	13 (6%)	3 (1%)	11 11

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	222	PRO
1	A	33	ALA
1	A	195	PRO

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	181/183 (99%)	158 (87%)	23 (13%)	4 4

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	VAL
1	A	23	ARG
1	A	45	SER
1	A	50	VAL
1	A	52	THR
1	A	62	ARG
1	A	63	ARG
1	A	64	GLU
1	A	66	GLN
1	A	72	SER
1	A	84	GLN
1	A	97	ARG
1	A	131	GLN
1	A	132	ARG
1	A	133	SER
1	A	136	ARG
1	A	164	LEU
1	A	166	ARG
1	A	167	ARG
1	A	192	SER
1	A	193	LEU
1	A	198	ARG
1	A	210	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	19	GLN
1	A	32	HIS
1	A	66	GLN
1	A	94	GLN
1	A	153	GLN
1	A	158	ASN
1	A	218	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 monosaccharides 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
2	NAG	A	411	1	14,14,15	1.24	1 (7%)	17,19,21	2.68	7 (41%)
2	NAG	A	401	1	14,14,15	1.31	2 (14%)	17,19,21	2.97	8 (47%)

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	NAG	A	411	1	-	0/6/23/26	0/1/1/1
2	NAG	A	401	1	-	2/6/23/26	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	411	NAG	O4-C4	2.49	1.48	1.43
2	A	401	NAG	C8-C7	2.35	1.55	1.50
2	A	401	NAG	O4-C4	2.31	1.48	1.43

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	NAG	C1-O5-C5	8.71	123.99	112.19
2	A	411	NAG	C1-O5-C5	6.37	120.83	112.19
2	A	401	NAG	C8-C7-N2	-4.78	108.01	116.10
2	A	411	NAG	O4-C4-C3	4.68	121.17	110.35
2	A	411	NAG	C2-N2-C7	-4.52	116.46	122.90
2	A	401	NAG	O5-C1-C2	-3.67	105.49	111.29
2	A	411	NAG	C8-C7-N2	-3.11	110.84	116.10
2	A	401	NAG	C2-N2-C7	-2.88	118.80	122.90
2	A	411	NAG	O3-C3-C2	-2.81	103.65	109.47
2	A	411	NAG	O5-C5-C4	-2.61	104.47	110.83
2	A	401	NAG	O4-C4-C3	2.57	116.30	110.35
2	A	401	NAG	O5-C5-C6	2.26	110.75	107.20
2	A	411	NAG	O5-C1-C2	-2.24	107.75	111.29
2	A	401	NAG	O7-C7-C8	2.03	125.83	122.06
2	A	401	NAG	O4-C4-C5	2.00	114.28	109.30

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401	NAG	O5-C5-C6-O6
2	A	401	NAG	C4-C5-C6-O6

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	411	NAG	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

Unable to reproduce the depositors R factor - this section is therefore empty.

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

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates

Unable to reproduce the depositors R factor - this section is therefore empty.

6.4 Ligands

Unable to reproduce the depositors R factor - this section is therefore empty.

6.5 Other polymers

Unable to reproduce the depositors R factor - this section is therefore empty.