



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 10:46 AM UTC

PDB ID : 1CEB / pdb_00001ceb
Title : THE STRUCTURE OF THE NON-COVALENT COMPLEX OF RECOMBINANT KRINGLE 1 DOMAIN OF HUMAN PLASMINOGEN WITH AMCHA (TRANS-4-AMINOMETHYLCYCLOHEXANE-1-CARBOXYLIC ACID)
Authors : Tulinsky, A.; Mathews, I.I.
Deposited on : 1995-12-03
Resolution : 2.07 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

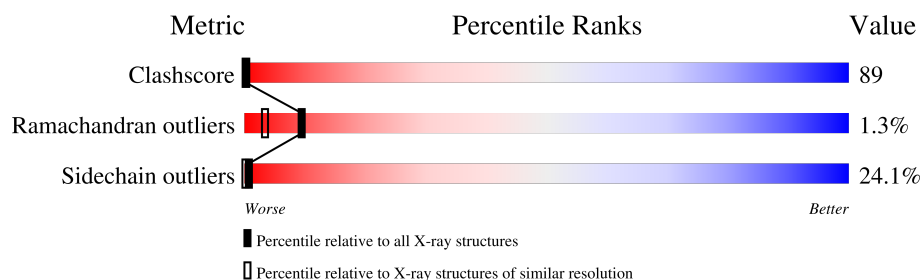
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.07 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	3883 (2.08-2.04)
Ramachandran outliers	187476	3860 (2.08-2.04)
Sidechain outliers	187428	3860 (2.08-2.04)

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	88	
1	B	88	

2 Entry composition [i](#)

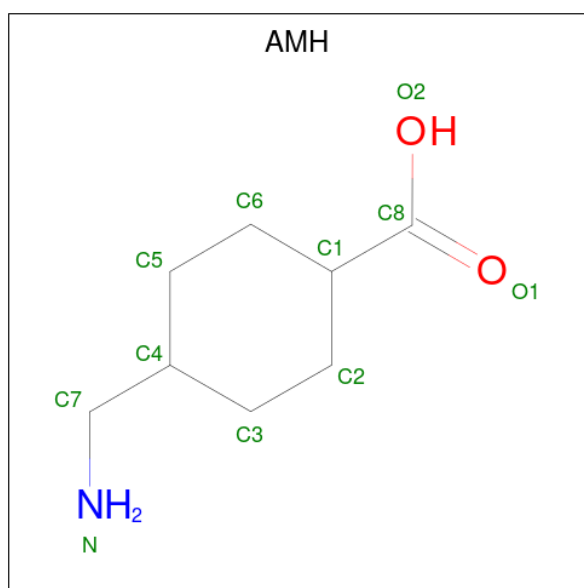
There are 3 unique types of molecules in this entry. The entry contains 1448 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 PLASMINOGEN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	80	Total	C	N	O	S	0	0	0
			641	390	115	129	7			
1	B	80	Total	C	N	O	S	0	0	0
			635	387	112	129	7			

- Molecule 2 is TRANS-4-AMINOMETHYLCYCLOHEXANE-1-CARBOXYLIC ACID (CCD ID: AMH) (formula: $C_8H_{15}NO_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			11	8	1	2		
2	B	1	Total	C	N	O	0	0
			11	8	1	2		

- Molecule 3 is water.

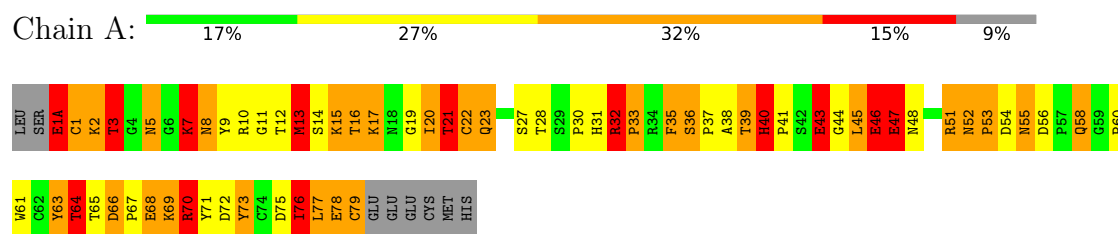
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	81	Total 81	O 81	0	0
3	B	69	Total 69	O 69	0	0

3 Residue-property plots

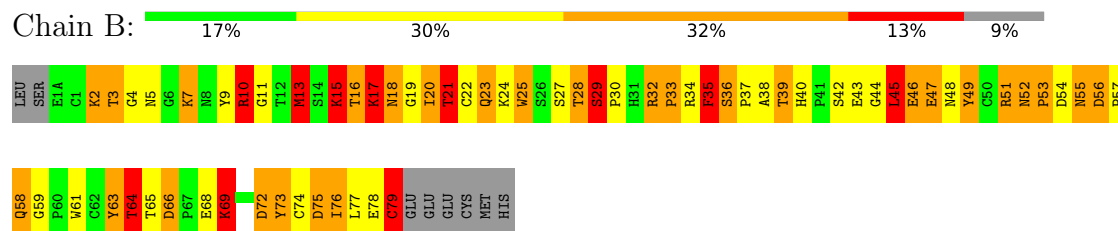
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded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



• Molecule 1: PLASMINOGEN



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	64.30Å 51.70Å 46.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	7.00 – 2.07	Depositor
% Data completeness (in resolution range)	57.0 (7.00-2.07)	Depositor
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	PROLSQ	Depositor
R, R_{free}	0.178 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	1448	wwPDB-VP
Average B, all atoms (Å ²)	17.0	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: AMH

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	1.37	2/660 (0.3%)	2.61	57/895 (6.4%)
1	B	1.49	1/654 (0.2%)	2.82	72/888 (8.1%)
All	All	1.43	3/1314 (0.2%)	2.72	129/1783 (7.2%)

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	0	3
1	B	0	1
All	All	0	4

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	13	MET	C-O	5.34	1.30	1.23
1	A	64	THR	CA-CB	5.18	1.61	1.53
1	B	58	GLN	C-O	5.07	1.30	1.24

All (129) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	72	ASP	CA-CB-CG	16.71	129.31	112.60
1	A	51	ARG	NE-CZ-NH2	16.70	134.23	119.20
1	B	46	GLU	CB-CG-CD	13.37	135.34	112.60
1	B	17	LYS	CA-CB-CG	10.01	134.12	114.10
1	A	78	GLU	CA-C-O	9.70	132.62	121.47
1	B	10	ARG	CA-C-N	9.30	134.70	121.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	10	ARG	C-N-CA	9.30	134.70	121.30
1	A	20	ILE	O-C-N	9.30	132.74	122.61
1	B	15	LYS	CA-C-O	-9.15	110.41	121.11
1	A	52	ASN	CA-CB-CG	-9.13	103.47	112.60
1	A	52	ASN	OD1-CG-ND2	8.88	131.48	122.60
1	B	53	PRO	O-C-N	8.61	132.99	122.23
1	B	46	GLU	CA-CB-CG	8.59	131.28	114.10
1	B	69	LYS	CG-CD-CE	8.58	131.03	111.30
1	B	69	LYS	CB-CG-CD	8.45	130.73	111.30
1	B	78	GLU	N-CA-CB	8.07	121.80	110.26
1	B	63	TYR	N-CA-C	-7.79	98.95	110.48
1	A	63	TYR	CA-C-N	7.71	135.75	122.64
1	A	63	TYR	C-N-CA	7.71	135.75	122.64
1	B	32	ARG	CA-C-O	-7.51	113.76	119.32
1	A	32	ARG	CB-CA-C	7.46	118.35	109.85
1	A	51	ARG	NH1-CZ-NH2	-7.45	109.61	119.30
1	A	55	ASN	OD1-CG-ND2	7.29	129.89	122.60
1	B	18	ASN	CA-CB-CG	-7.14	105.46	112.60
1	A	56	ASP	CA-CB-CG	7.13	119.73	112.60
1	B	75	ASP	CA-CB-CG	-7.09	105.51	112.60
1	A	38	ALA	CA-C-N	7.08	136.72	121.64
1	A	38	ALA	C-N-CA	7.08	136.72	121.64
1	B	78	GLU	CA-C-O	6.93	129.49	121.87
1	A	20	ILE	CB-CA-C	-6.79	100.94	111.26
1	A	3	THR	CB-CA-C	6.78	122.78	109.35
1	B	2	LYS	CA-C-N	-6.75	112.22	122.81
1	B	2	LYS	C-N-CA	-6.75	112.22	122.81
1	B	58	GLN	OE1-CD-NE2	-6.74	115.86	122.60
1	A	66	ASP	CB-CA-C	6.68	119.00	109.38
1	B	74	CYS	CA-C-N	6.63	132.07	122.85
1	B	74	CYS	C-N-CA	6.63	132.07	122.85
1	B	64	THR	OG1-CB-CG2	6.61	122.52	109.30
1	A	39	THR	CA-C-N	-6.52	115.59	122.89
1	A	39	THR	C-N-CA	-6.52	115.59	122.89
1	B	64	THR	N-CA-CB	-6.51	100.37	110.46
1	A	43	GLU	CA-C-O	-6.45	111.47	119.28
1	B	38	ALA	O-C-N	-6.41	114.38	122.27
1	A	23	GLN	N-CA-C	-6.41	100.21	110.14
1	B	36	SER	CA-C-O	6.34	126.22	120.71
1	B	32	ARG	O-C-N	6.29	127.13	121.35
1	A	77	LEU	CB-CA-C	6.27	120.38	109.65
1	A	69	LYS	N-CA-CB	-6.20	100.22	110.95

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	17	LYS	N-CA-C	-6.20	105.20	112.89
1	B	39	THR	CA-C-N	-6.20	116.04	122.85
1	B	39	THR	C-N-CA	-6.20	116.04	122.85
1	B	69	LYS	CB-CA-C	6.16	121.13	111.77
1	A	78	GLU	CA-C-N	6.14	132.76	121.70
1	A	78	GLU	C-N-CA	6.14	132.76	121.70
1	A	79	CYS	N-CA-CB	-6.12	100.10	110.50
1	B	35	PHE	CA-C-N	6.04	130.76	122.07
1	B	35	PHE	C-N-CA	6.04	130.76	122.07
1	A	55	ASN	CA-C-N	6.03	129.45	120.83
1	A	55	ASN	C-N-CA	6.03	129.45	120.83
1	B	10	ARG	CA-CB-CG	6.00	126.10	114.10
1	A	20	ILE	N-CA-CB	5.99	119.47	110.13
1	A	46	GLU	N-CA-C	5.93	119.92	109.96
1	A	16	THR	CA-CB-OG1	-5.92	100.71	109.60
1	B	34	ARG	O-C-N	5.89	130.23	122.33
1	B	51	ARG	NH1-CZ-NH2	-5.88	111.66	119.30
1	B	48	ASN	CA-C-O	5.87	126.41	119.48
1	B	23	GLN	OE1-CD-NE2	-5.87	116.73	122.60
1	A	35	PHE	CA-CB-CG	5.86	119.66	113.80
1	B	33	PRO	O-C-N	5.83	130.41	123.06
1	A	40	HIS	CA-C-O	5.81	127.89	121.15
1	B	58	GLN	CA-C-O	-5.79	112.24	120.51
1	A	51	ARG	NE-CZ-NH1	-5.78	115.72	121.50
1	A	76	ILE	O-C-N	5.78	129.45	123.03
1	B	51	ARG	O-C-N	-5.77	116.19	123.17
1	A	58	GLN	CA-C-O	-5.77	111.56	119.12
1	A	63	TYR	N-CA-C	-5.77	101.08	110.20
1	A	7	LYS	CA-C-O	5.74	126.64	120.55
1	B	76	ILE	CB-CA-C	5.73	119.15	110.62
1	B	56	ASP	CA-C-O	5.70	125.11	119.75
1	A	51	ARG	CA-C-N	5.66	130.64	123.34
1	A	51	ARG	C-N-CA	5.66	130.64	123.34
1	A	70	ARG	N-CA-CB	5.64	118.19	110.01
1	B	77	LEU	CB-CA-C	5.64	118.83	109.53
1	B	34	ARG	N-CA-CB	5.62	118.96	110.30
1	B	66	ASP	CB-CA-C	5.58	117.41	109.38
1	B	16	THR	CB-CA-C	-5.56	101.31	112.99
1	B	55	ASN	CA-C-N	5.54	128.18	120.65
1	B	55	ASN	C-N-CA	5.54	128.18	120.65
1	B	73	TYR	CA-C-N	5.52	131.76	122.87
1	B	73	TYR	C-N-CA	5.52	131.76	122.87

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1(A)	GLU	CB-CG-CD	-5.52	103.22	112.60
1	B	27	SER	CA-C-O	-5.52	114.65	121.11
1	A	52	ASN	CB-CG-OD1	-5.51	109.78	120.80
1	B	13	MET	CB-CA-C	5.50	118.85	109.72
1	A	33	PRO	N-CA-CB	5.48	108.40	103.19
1	B	21	THR	CA-CB-OG1	-5.48	101.38	109.60
1	B	47	GLU	CG-CD-OE1	5.47	130.99	118.40
1	B	16	THR	CA-C-O	-5.46	115.81	121.98
1	A	53	PRO	O-C-N	5.45	128.51	122.24
1	B	72	ASP	CB-CA-C	5.45	120.13	110.16
1	B	49	TYR	CA-C-O	-5.44	112.73	120.51
1	A	70	ARG	CA-C-O	-5.43	115.12	120.82
1	A	33	PRO	O-C-N	5.41	129.88	123.06
1	A	63	TYR	N-CA-CB	5.41	118.23	110.06
1	A	55	ASN	N-CA-CB	-5.37	103.84	111.84
1	B	34	ARG	N-CA-C	-5.36	105.48	112.23
1	B	72	ASP	O-C-N	-5.36	117.48	123.48
1	A	32	ARG	N-CA-CB	5.34	117.81	110.01
1	B	21	THR	O-C-N	5.33	129.49	122.93
1	A	76	ILE	CB-CA-C	5.28	118.03	110.84
1	B	51	ARG	NE-CZ-NH2	5.28	123.95	119.20
1	A	14	SER	CA-CB-OG	5.28	121.66	111.10
1	A	8	ASN	CB-CA-C	5.28	119.65	110.94
1	A	12	THR	CA-C-O	-5.27	114.31	120.53
1	B	47	GLU	CB-CG-CD	5.25	121.53	112.60
1	B	63	TYR	N-CA-CB	5.24	118.26	110.29
1	A	21	THR	O-C-N	5.21	129.34	122.93
1	B	29	SER	CA-C-O	5.21	124.33	119.13
1	B	79	CYS	N-CA-CB	-5.20	101.66	110.50
1	B	52	ASN	OD1-CG-ND2	5.18	127.78	122.60
1	B	42	SER	N-CA-C	5.15	119.25	112.92
1	B	3	THR	CA-CB-OG1	-5.14	101.90	109.60
1	B	76	ILE	O-C-N	5.12	128.71	123.18
1	B	68	GLU	CG-CD-OE2	5.12	130.17	118.40
1	A	35	PHE	N-CA-CB	5.11	118.92	110.69
1	A	47	GLU	CB-CA-C	-5.08	109.22	116.34
1	B	22	CYS	O-C-N	5.06	129.06	122.89
1	B	25	TRP	O-C-N	5.03	128.93	122.39
1	A	46	GLU	OE1-CD-OE2	5.02	134.94	122.90

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	22	CYS	Mainchain
1	A	68	GLU	Mainchain
1	A	70	ARG	Sidechain
1	B	58	GLN	Mainchain

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	641	0	583	147	0
1	B	635	0	572	82	0
2	A	11	0	13	1	0
2	B	11	0	14	2	0
3	A	81	0	0	14	0
3	B	69	0	0	3	0
All	All	1448	0	1182	221	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 89.

All (221) 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:64:THR:HG23	1:A:66:ASP:H	1.09	1.18
1:A:3:THR:HG22	1:A:79:CYS:CB	1.80	1.10
1:A:1(A):GLU:CA	1:A:1(A):GLU:OE1	1.92	1.09
1:A:28:THR:HB	1:A:32:ARG:HG3	1.22	1.09
1:A:28:THR:HB	1:A:32:ARG:CG	1.86	1.05
1:B:36:SER:HB2	1:B:37:PRO:HD2	1.42	1.02
1:A:22:CYS:O	1:A:65:THR:HG23	1.58	1.01
1:B:28:THR:HG22	1:B:32:ARG:HA	1.42	1.01
1:A:15:LYS:HG2	1:A:20:ILE:C	1.86	1.01
1:B:18:ASN:ND2	3:B:310:HOH:O	1.95	0.99
1:A:3:THR:CG2	1:A:79:CYS:HB3	1.93	0.97
1:A:3:THR:HG22	1:A:79:CYS:HB3	0.98	0.97
1:A:39:THR:HG22	1:A:40:HIS:CE1	2.00	0.96
1:A:64:THR:CG2	1:A:66:ASP:H	1.78	0.95

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:69:LYS:HE2	1:A:72:ASP:HB2	1.46	0.95
1:A:1(A):GLU:OE1	1:A:1(A):GLU:HA	1.13	0.94
1:A:15:LYS:HD2	1:A:19:GLY:O	1.66	0.94
1:A:13:MET:CE	1:A:15:LYS:O	2.15	0.94
1:B:28:THR:CG2	1:B:32:ARG:HG3	2.00	0.91
1:A:13:MET:HE3	1:A:15:LYS:O	1.70	0.91
1:A:64:THR:HG23	1:A:66:ASP:N	1.86	0.90
1:B:17:LYS:HB3	1:B:75:ASP:OD2	1.74	0.88
1:A:2:LYS:O	1:A:79:CYS:CB	2.20	0.88
1:B:17:LYS:HB2	1:B:17:LYS:HZ3	1.38	0.88
1:A:16:THR:HG21	1:A:72:ASP:HB3	1.54	0.87
1:B:28:THR:HG22	1:B:32:ARG:CA	2.05	0.85
1:A:39:THR:CG2	1:A:40:HIS:CE1	2.60	0.85
1:A:77:LEU:HD12	1:A:77:LEU:N	1.92	0.85
1:A:13:MET:HE2	3:A:376:HOH:O	1.76	0.85
1:A:40:HIS:CE1	1:B:39:THR:CG2	2.60	0.83
1:B:9:TYR:OH	1:B:13:MET:HB2	1.79	0.81
1:B:28:THR:HG22	1:B:32:ARG:CB	2.08	0.81
1:A:64:THR:HG21	1:A:69:LYS:HG3	1.62	0.80
1:B:28:THR:HG22	1:B:32:ARG:HG3	1.64	0.80
1:B:16:THR:OG1	1:B:20:ILE:HG23	1.82	0.79
1:A:2:LYS:O	1:A:79:CYS:HB3	1.82	0.79
1:B:44:GLY:HA2	1:B:46:GLU:OE1	1.84	0.78
1:B:10:ARG:NH1	1:B:43:GLU:O	2.18	0.77
1:A:40:HIS:NE2	1:B:39:THR:HG22	2.00	0.77
1:B:51:ARG:C	1:B:53:PRO:HD3	2.11	0.76
1:A:68:GLU:N	1:A:68:GLU:OE1	2.17	0.76
1:B:36:SER:HB2	1:B:37:PRO:CD	2.16	0.76
1:B:5:ASN:OD1	1:B:52:ASN:ND2	2.19	0.76
1:A:23:GLN:OE1	1:A:30:PRO:HD2	1.87	0.75
1:A:40:HIS:CE1	1:B:39:THR:HG22	2.22	0.75
1:B:28:THR:HG22	1:B:32:ARG:CG	2.17	0.75
1:A:55:ASN:OD1	1:A:55:ASN:O	2.04	0.74
1:B:33:PRO:HB3	1:B:63:TYR:CE2	2.23	0.74
1:B:20:ILE:HD12	1:B:21:THR:N	2.03	0.74
1:A:13:MET:HE1	1:A:15:LYS:O	1.86	0.74
1:A:39:THR:C	1:A:41:PRO:HD3	2.12	0.73
1:B:44:GLY:O	1:B:46:GLU:N	2.20	0.73
1:B:64:THR:HG23	1:B:66:ASP:H	1.54	0.73
1:A:69:LYS:CE	1:A:72:ASP:HB2	2.18	0.73
1:B:17:LYS:HB2	1:B:17:LYS:NZ	2.03	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:58:GLN:HB2	1:A:73:TYR:CE2	2.25	0.72
1:A:67:PRO:HG2	3:A:266:HOH:O	1.89	0.71
1:A:20:ILE:CG1	1:A:21:THR:N	2.54	0.71
1:A:44:GLY:O	1:A:46:GLU:HG2	1.90	0.70
1:A:16:THR:OG1	1:A:20:ILE:HG22	1.91	0.70
1:B:44:GLY:CA	1:B:46:GLU:OE1	2.40	0.70
1:A:69:LYS:HE2	1:A:71:TYR:O	1.91	0.69
1:B:75:ASP:C	1:B:76:ILE:HD13	2.18	0.69
1:A:20:ILE:HD13	1:A:64:THR:HG23	1.74	0.69
1:B:28:THR:CG2	1:B:32:ARG:CG	2.71	0.68
1:A:5:ASN:ND2	3:A:216:HOH:O	2.27	0.68
1:A:64:THR:CG2	1:A:66:ASP:HB3	2.24	0.67
1:B:64:THR:CG2	1:B:66:ASP:H	2.07	0.67
1:A:64:THR:CG2	1:A:66:ASP:N	2.52	0.67
1:A:55:ASN:O	1:A:55:ASN:CG	2.37	0.67
1:A:27:SER:O	3:A:277:HOH:O	2.13	0.67
1:A:2:LYS:HD3	1:A:78:GLU:OE2	1.94	0.66
1:A:15:LYS:HG3	1:A:21:THR:HA	1.76	0.66
1:A:2:LYS:O	1:A:79:CYS:N	2.28	0.66
1:A:69:LYS:CE	1:A:71:TYR:O	2.44	0.66
1:A:64:THR:HG22	1:A:66:ASP:O	1.96	0.66
1:B:54:ASP:O	1:B:55:ASN:HB3	1.95	0.65
1:A:28:THR:CB	1:A:32:ARG:HG3	2.14	0.65
1:A:75:ASP:C	1:A:76:ILE:HD12	2.22	0.65
1:A:77:LEU:N	1:A:77:LEU:CD1	2.60	0.64
1:A:2:LYS:HG2	1:A:3:THR:N	2.13	0.64
1:A:40:HIS:HE1	1:B:39:THR:CG2	2.06	0.64
1:A:16:THR:HB	1:A:73:TYR:O	1.98	0.63
1:B:73:TYR:HB3	3:B:333:HOH:O	1.99	0.63
1:A:2:LYS:O	1:A:79:CYS:HB2	1.99	0.63
1:A:31:HIS:CE1	1:A:70:ARG:HA	2.33	0.63
1:A:63:TYR:HE2	1:A:70:ARG:NH1	1.97	0.62
1:B:10:ARG:HH11	1:B:51:ARG:HD2	1.63	0.62
1:A:39:THR:HG23	1:B:40:HIS:CE1	2.35	0.62
1:A:77:LEU:HD12	1:A:77:LEU:H	1.63	0.62
1:A:2:LYS:C	1:A:79:CYS:HB3	2.24	0.62
1:B:9:TYR:CE1	1:B:11:GLY:HA3	2.35	0.61
1:B:59:GLY:O	1:B:61:TRP:HD1	1.82	0.61
1:B:28:THR:HG21	1:B:32:ARG:NE	2.15	0.61
1:A:54:ASP:OD2	2:A:90:AMH:H71	2.01	0.61
1:A:7:LYS:HG2	1:A:55:ASN:HD22	1.66	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:23:GLN:OE1	1:B:30:PRO:HD2	2.01	0.60
1:A:10:ARG:HE	1:A:51:ARG:HD2	1.65	0.60
1:A:35:PHE:CD2	1:A:53:PRO:HB2	2.37	0.60
1:B:64:THR:HB	1:B:69:LYS:O	2.01	0.60
1:A:64:THR:HG21	1:A:66:ASP:HB3	1.84	0.60
1:A:21:THR:HG22	1:A:22:CYS:O	2.02	0.59
1:A:39:THR:HG22	1:A:40:HIS:ND1	2.15	0.59
1:B:29:SER:HA	1:B:30:PRO:C	2.26	0.59
1:B:17:LYS:NZ	1:B:17:LYS:CB	2.64	0.59
1:A:77:LEU:CD1	1:A:77:LEU:H	2.14	0.58
1:A:10:ARG:HE	1:A:51:ARG:CD	2.16	0.57
1:B:61:TRP:CH2	2:B:90:AMH:H31	2.39	0.57
1:A:63:TYR:CE2	1:A:70:ARG:NH1	2.72	0.57
1:A:20:ILE:HG13	1:A:21:THR:N	2.18	0.57
1:A:10:ARG:HD3	1:A:51:ARG:NE	2.20	0.57
1:A:13:MET:CE	3:A:376:HOH:O	2.39	0.57
1:A:16:THR:OG1	1:A:20:ILE:CG2	2.52	0.57
1:A:22:CYS:O	1:A:65:THR:CG2	2.43	0.56
1:B:56:ASP:OD1	1:B:57:PRO:HD2	2.05	0.56
1:B:61:TRP:CZ2	2:B:90:AMH:H71	2.40	0.56
1:A:23:GLN:HG3	1:A:30:PRO:HD2	1.88	0.56
1:A:76:ILE:CD1	1:A:76:ILE:N	2.69	0.56
1:A:20:ILE:HG13	1:A:21:THR:H	1.71	0.55
1:B:35:PHE:CD1	1:B:35:PHE:N	2.74	0.55
1:A:39:THR:CG2	1:B:40:HIS:CE1	2.90	0.55
1:B:10:ARG:NH1	1:B:51:ARG:HD2	2.22	0.55
1:B:7:LYS:HD3	1:B:52:ASN:O	2.07	0.55
1:A:20:ILE:HG12	1:A:21:THR:N	2.21	0.54
1:B:28:THR:HG21	1:B:32:ARG:HG3	1.87	0.54
1:A:8:ASN:ND2	3:A:309:HOH:O	2.39	0.54
1:A:1(A):GLU:HG3	3:A:433:HOH:O	2.08	0.54
1:A:40:HIS:HB3	1:A:43:GLU:HG2	1.89	0.54
1:B:46:GLU:O	1:B:47:GLU:C	2.49	0.54
1:A:64:THR:HG21	1:A:69:LYS:CG	2.36	0.53
1:A:40:HIS:N	1:A:41:PRO:HD3	2.24	0.53
1:B:5:ASN:HD21	1:B:55:ASN:ND2	2.07	0.53
1:A:3:THR:HG22	1:A:79:CYS:SG	2.49	0.53
1:A:39:THR:HG21	1:A:40:HIS:CE1	2.43	0.53
1:A:5:ASN:OD1	1:A:52:ASN:ND2	2.41	0.52
1:B:24:LYS:HD3	1:B:47:GLU:O	2.09	0.52
1:A:15:LYS:CD	1:A:19:GLY:O	2.50	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:69:LYS:HE3	3:A:252:HOH:O	2.10	0.52
1:B:21:THR:CG2	1:B:65:THR:HG21	2.39	0.52
1:A:20:ILE:CD1	1:A:64:THR:HG23	2.40	0.52
1:B:20:ILE:HD12	1:B:21:THR:H	1.75	0.51
1:A:1(A):GLU:CG	3:A:433:HOH:O	2.58	0.51
1:B:2:LYS:O	1:B:79:CYS:HB2	2.10	0.51
1:B:28:THR:CG2	1:B:32:ARG:HA	2.28	0.51
1:A:75:ASP:HB3	3:A:271:HOH:O	2.10	0.50
1:A:20:ILE:HD12	1:A:66:ASP:HB2	1.93	0.50
1:A:1(A):GLU:HB3	1:A:9:TYR:CE1	2.47	0.49
1:A:9:TYR:CZ	1:A:11:GLY:HA3	2.47	0.49
1:B:5:ASN:ND2	3:B:315:HOH:O	2.37	0.49
1:A:5:ASN:OD1	1:A:55:ASN:HA	2.12	0.49
1:B:52:ASN:N	1:B:53:PRO:HD3	2.27	0.49
1:B:5:ASN:ND2	1:B:55:ASN:HD21	2.10	0.49
1:A:44:GLY:O	1:A:46:GLU:N	2.42	0.49
1:A:31:HIS:O	1:A:33:PRO:HD3	2.13	0.49
1:A:1(A):GLU:HB3	1:A:9:TYR:HE1	1.78	0.48
1:A:15:LYS:CG	1:A:21:THR:HA	2.42	0.48
1:A:69:LYS:HG3	1:A:69:LYS:O	2.14	0.48
1:A:10:ARG:HD3	1:A:51:ARG:HE	1.79	0.48
1:A:69:LYS:HD2	1:A:71:TYR:O	2.14	0.48
1:B:5:ASN:HD21	1:B:55:ASN:HD21	1.61	0.48
1:B:51:ARG:O	1:B:53:PRO:HD3	2.13	0.47
1:B:5:ASN:ND2	1:B:55:ASN:ND2	2.62	0.47
1:B:15:LYS:HG2	1:B:19:GLY:C	2.40	0.47
1:A:73:TYR:N	1:A:73:TYR:CD1	2.83	0.47
1:B:25:TRP:CE3	1:B:45:LEU:HG	2.50	0.46
1:A:40:HIS:CB	1:A:43:GLU:HG2	2.46	0.46
1:A:64:THR:HG23	1:A:66:ASP:HB3	1.98	0.46
1:A:33:PRO:HB3	1:A:63:TYR:CE1	2.50	0.46
1:A:36:SER:HB2	1:A:37:PRO:HD2	1.97	0.46
1:A:2:LYS:HG2	1:A:78:GLU:OE2	2.15	0.46
1:B:10:ARG:HH11	1:B:51:ARG:CD	2.28	0.46
1:A:40:HIS:N	1:A:41:PRO:CD	2.78	0.46
1:B:36:SER:CA	1:B:45:LEU:HD23	2.45	0.46
1:A:39:THR:C	1:A:41:PRO:CD	2.86	0.45
1:B:59:GLY:O	1:B:73:TYR:HD2	1.98	0.45
1:A:78:GLU:N	3:A:405:HOH:O	2.19	0.45
1:A:15:LYS:HG2	1:A:20:ILE:O	2.13	0.45
1:B:44:GLY:C	1:B:46:GLU:N	2.75	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:17:LYS:N	1:A:73:TYR:O	2.40	0.44
1:A:10:ARG:CD	1:A:51:ARG:HE	2.30	0.44
1:A:23:GLN:OE1	1:A:30:PRO:CD	2.60	0.44
1:B:9:TYR:HH	1:B:13:MET:HB2	1.81	0.44
1:A:40:HIS:NE2	1:B:39:THR:CG2	2.70	0.44
1:A:7:LYS:CD	1:A:55:ASN:HB2	2.48	0.43
1:A:64:THR:CG2	1:A:69:LYS:O	2.66	0.43
1:A:36:SER:CB	1:A:37:PRO:CD	2.96	0.43
1:A:76:ILE:HD12	1:A:76:ILE:N	2.33	0.43
1:B:36:SER:CB	1:B:37:PRO:HD2	2.28	0.43
1:A:1:CYS:HB2	1:A:79:CYS:HB2	1.71	0.43
1:A:15:LYS:HA	1:A:20:ILE:O	2.18	0.42
1:A:55:ASN:HB3	3:A:246:HOH:O	2.18	0.42
1:A:69:LYS:CD	1:A:71:TYR:O	2.66	0.42
1:B:21:THR:HG23	1:B:65:THR:HG21	2.01	0.42
1:A:20:ILE:HD11	1:A:66:ASP:N	2.34	0.42
1:B:28:THR:HG23	1:B:33:PRO:HD2	2.01	0.42
1:A:47:GLU:HB3	1:A:48:ASN:H	1.66	0.42
1:B:24:LYS:CD	1:B:47:GLU:O	2.67	0.42
1:B:61:TRP:CD1	1:B:61:TRP:N	2.87	0.42
1:A:7:LYS:HD3	1:A:55:ASN:HB2	2.01	0.42
1:A:51:ARG:C	1:A:53:PRO:HD3	2.45	0.42
1:B:13:MET:O	1:B:49:TYR:HA	2.20	0.42
1:B:15:LYS:HG3	1:B:20:ILE:C	2.45	0.42
1:B:49:TYR:HB3	1:B:51:ARG:HH12	1.84	0.41
1:A:39:THR:CG2	1:B:40:HIS:NE2	2.84	0.41
1:A:10:ARG:CD	1:A:51:ARG:NE	2.83	0.41
1:B:3:THR:O	1:B:4:GLY:C	2.63	0.41
1:A:15:LYS:HG2	1:A:21:THR:N	2.30	0.41
1:B:21:THR:HG22	1:B:65:THR:OG1	2.20	0.41
1:A:9:TYR:O	1:A:10:ARG:NE	2.54	0.41
1:A:28:THR:HB	1:A:32:ARG:HG2	1.89	0.41
1:A:8:ASN:HB2	3:A:367:HOH:O	2.20	0.41
1:A:20:ILE:HD13	1:A:64:THR:CG2	2.48	0.41
1:A:7:LYS:O	1:A:10:ARG:NH2	2.53	0.41
1:A:46:GLU:HG2	1:A:46:GLU:H	1.47	0.41
1:B:24:LYS:HD3	1:B:47:GLU:C	2.45	0.41
1:A:23:GLN:CG	1:A:30:PRO:HD2	2.50	0.41
1:A:61:TRP:HA	1:A:73:TYR:HA	2.03	0.41
1:A:27:SER:HB2	3:A:237:HOH:O	2.20	0.40
1:A:2:LYS:CD	1:A:78:GLU:OE2	2.66	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:64:THR:CG2	1:A:66:ASP:CB	2.98	0.40
1:A:36:SER:C	1:A:45:LEU:HD23	2.46	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	78/88 (89%)	71 (91%)	6 (8%)	1 (1%)	9	3
1	B	78/88 (89%)	71 (91%)	6 (8%)	1 (1%)	9	3
All	All	156/176 (89%)	142 (91%)	12 (8%)	2 (1%)	9	3

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	60	PRO
1	B	45	LEU

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	73/81 (90%)	53 (73%)	20 (27%)	0	0
1	B	72/81 (89%)	57 (79%)	15 (21%)	1	0

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	145/162 (90%)	110 (76%)	35 (24%)	1 0

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1(A)	GLU
1	A	1	CYS
1	A	2	LYS
1	A	3	THR
1	A	5	ASN
1	A	7	LYS
1	A	13	MET
1	A	15	LYS
1	A	17	LYS
1	A	21	THR
1	A	32	ARG
1	A	36	SER
1	A	40	HIS
1	A	43	GLU
1	A	45	LEU
1	A	46	GLU
1	A	47	GLU
1	A	64	THR
1	A	73	TYR
1	A	76	ILE
1	B	7	LYS
1	B	10	ARG
1	B	13	MET
1	B	15	LYS
1	B	17	LYS
1	B	20	ILE
1	B	21	THR
1	B	28	THR
1	B	29	SER
1	B	35	PHE
1	B	45	LEU
1	B	64	THR
1	B	69	LYS
1	B	72	ASP
1	B	79	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	8	ASN
1	A	31	HIS
1	A	55	ASN
1	A	58	GLN
1	B	18	ASN
1	B	31	HIS
1	B	55	ASN
1	B	58	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 [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	AMH	B	90	-	10,11,11	1.33	2 (20%)	13,14,14	2.52	6 (46%)
2	AMH	A	90	-	10,11,11	1.20	1 (10%)	13,14,14	2.42	4 (30%)

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	AMH	B	90	-	-	0/6/16/16	0/1/1/1
2	AMH	A	90	-	-	1/6/16/16	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	90	AMH	C3-C2	-2.75	1.46	1.52
2	A	90	AMH	C2-C1	-2.16	1.47	1.53
2	B	90	AMH	C5-C4	2.09	1.58	1.52

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	90	AMH	O2-C8-O1	-6.05	110.35	124.08
2	A	90	AMH	C6-C1-C2	5.60	121.81	110.00
2	A	90	AMH	C2-C1-C8	4.28	120.10	111.28
2	A	90	AMH	C5-C4-C3	3.18	117.06	109.29
2	B	90	AMH	C6-C1-C8	2.99	117.44	111.28
2	B	90	AMH	C5-C6-C1	-2.88	106.10	111.18
2	B	90	AMH	O1-C8-C1	2.67	129.93	122.86
2	A	90	AMH	O2-C8-O1	-2.27	118.94	124.08
2	B	90	AMH	O2-C8-C1	2.14	119.72	114.16
2	B	90	AMH	C6-C1-C2	2.03	114.28	110.00

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	90	AMH	C3-C4-C7-N

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	90	AMH	2	0
2	A	90	AMH	1	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.