



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:41 PM UTC

PDB ID : 4AST / pdb_00004ast
Title : The apo structure of a bacterial aldo-keto reductase AKR14A1
Authors : Zhu, X.; Ellis, E.M.; Lapthorn, A.
Deposited on : 2012-05-03
Resolution : 2.38 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
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

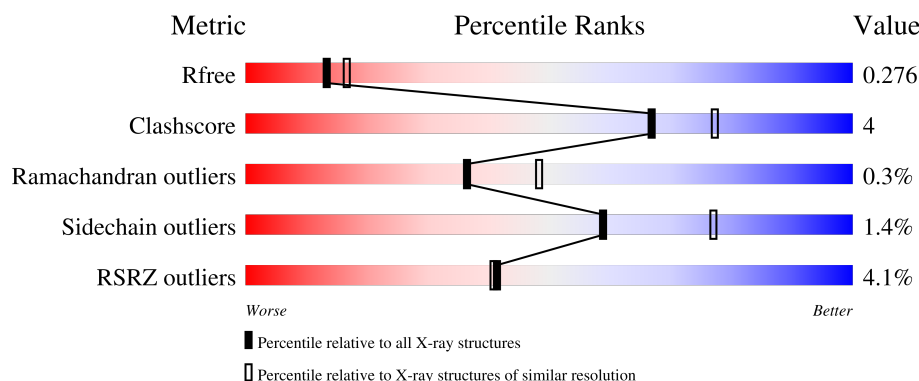
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.38 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))
R_{free}	180053	7164 (2.40-2.36)
Clashscore	190562	7722 (2.40-2.36)
Ramachandran outliers	187476	7626 (2.40-2.36)
Sidechain outliers	187428	7627 (2.40-2.36)
RSRZ outliers	180081	7170 (2.40-2.36)

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 $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	346	
1	B	346	
1	C	346	
1	D	346	
1	E	346	

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Mol	Chain	Length	Quality of chain
1	F	346	4% 79% 7% 12%
1	G	346	6% 79% 8% 12%
1	H	346	8% 78% 7% 14%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 20048 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 ALDO-KETO REDUCTASE AKR14A1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	307	Total	C	N	O	S	0	1	0
			2429	1533	428	458	10			
1	B	311	Total	C	N	O	S	0	0	0
			2449	1546	431	462	10			
1	C	310	Total	C	N	O	S	0	1	0
			2453	1548	431	464	10			
1	D	308	Total	C	N	O	S	0	2	0
			2447	1546	431	460	10			
1	E	307	Total	C	N	O	S	0	1	0
			2428	1533	427	458	10			
1	F	303	Total	C	N	O	S	0	1	0
			2403	1518	423	452	10			
1	G	303	Total	C	N	O	S	0	1	0
			2400	1516	422	452	10			
1	H	297	Total	C	N	O	S	0	1	0
			2358	1492	415	441	10			

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	113	Total	O	0	0
			113	113		
2	B	114	Total	O	0	0
			114	114		
2	C	93	Total	O	0	0
			93	93		
2	D	96	Total	O	0	0
			96	96		
2	E	86	Total	O	0	0
			86	86		
2	F	63	Total	O	0	0
			63	63		

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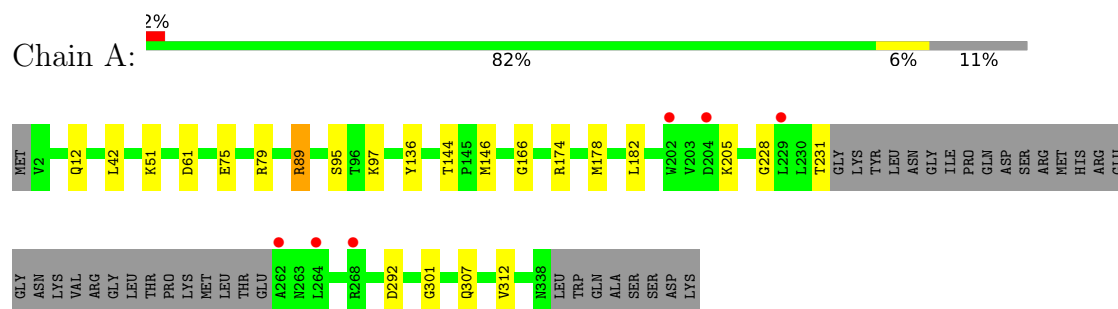
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	G	54	Total	O	0	0
			54	54		
2	H	62	Total	O	0	0
			62	62		

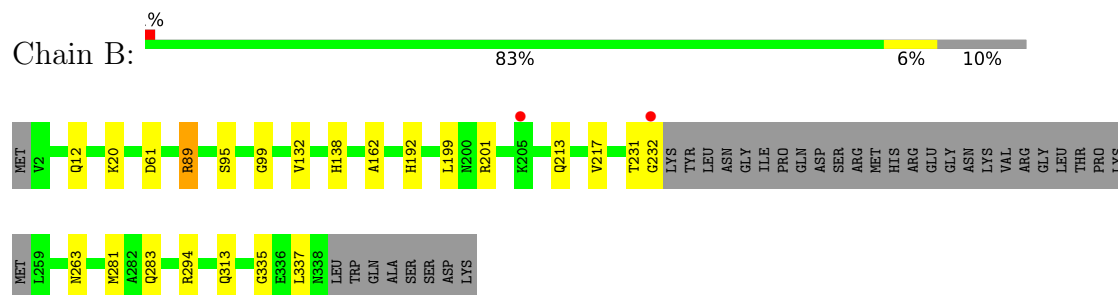
3 Residue-property plots [i](#)

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 and electron density. 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. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). 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.

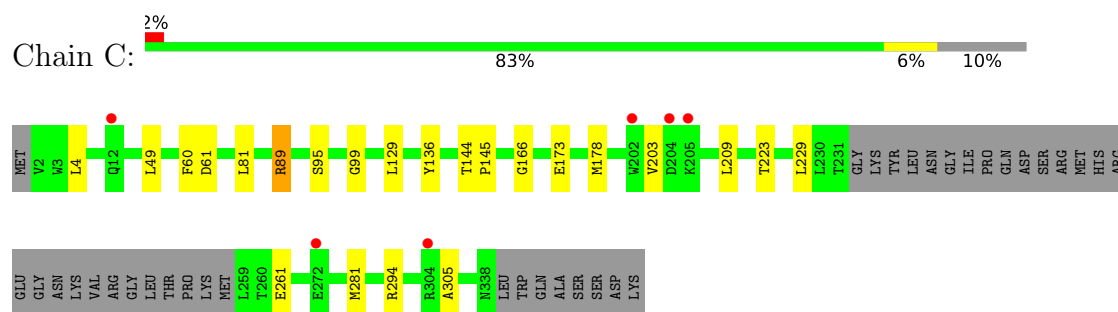
• Molecule 1: ALDO-KETO REDUCTASE AKR14A1



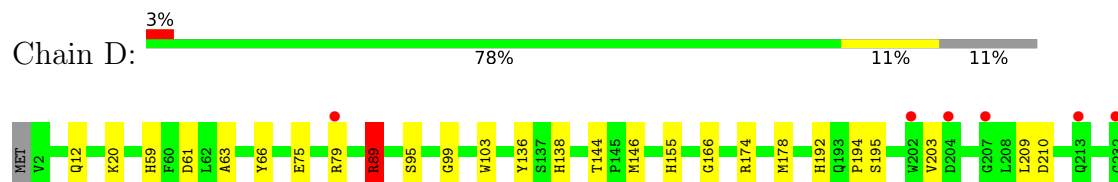
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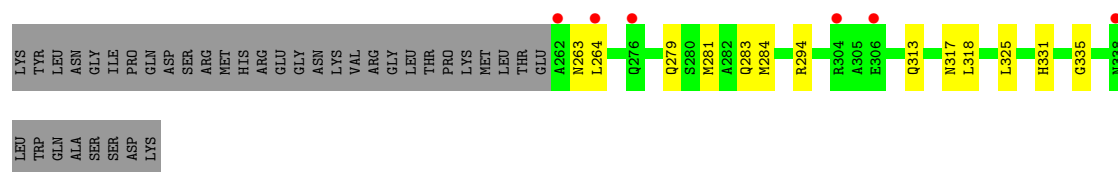


• Molecule 1: ALDO-KETO REDUCTASE AKR14A1

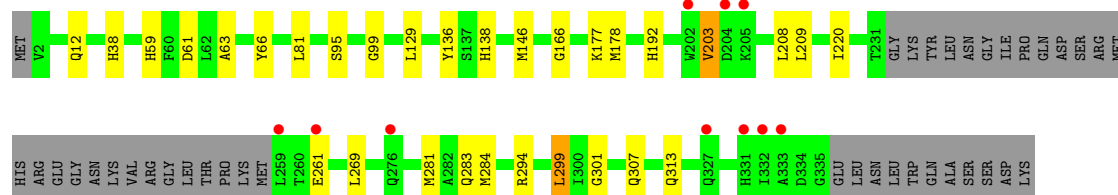
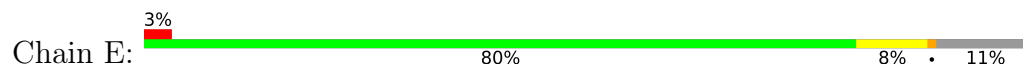


• Molecule 1: ALDO-KETO REDUCTASE AKR14A1

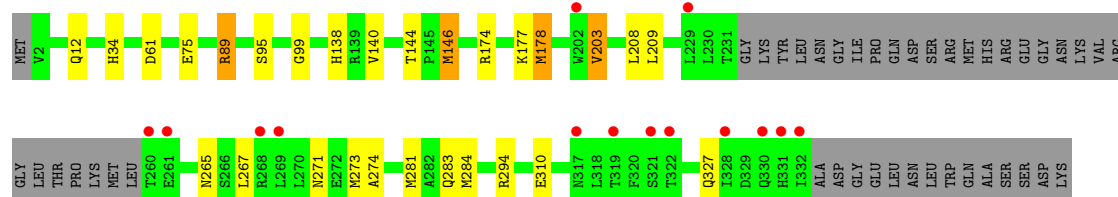
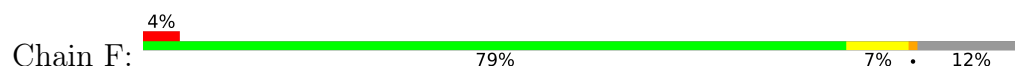




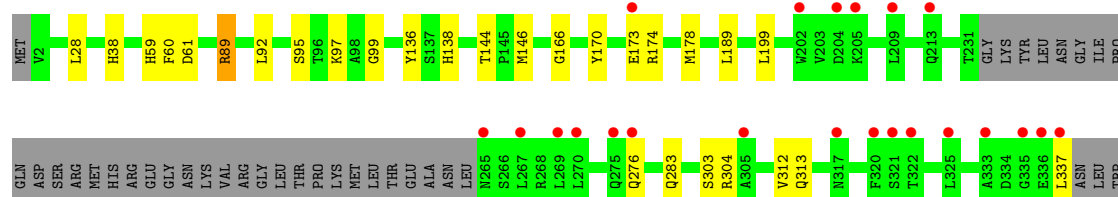
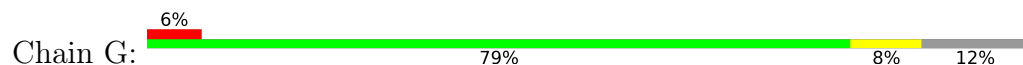
• Molecule 1: ALDO-KETO REDUCTASE AKR14A1



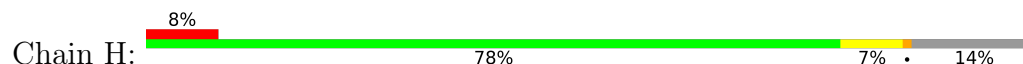
• Molecule 1: ALDO-KETO REDUCTASE AKR14A1

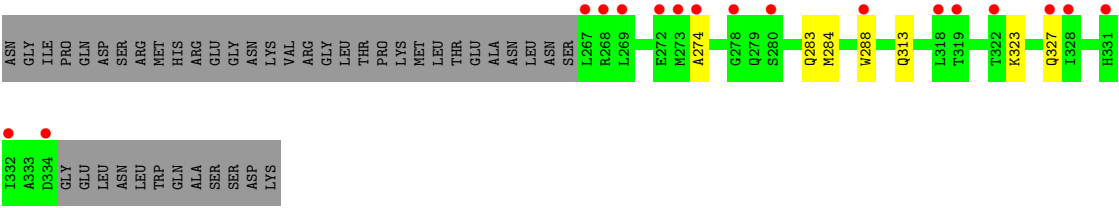


• Molecule 1: ALDO-KETO REDUCTASE AKR14A1



• Molecule 1: ALDO-KETO REDUCTASE AKR14A1





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	98.81Å 144.32Å 113.90Å 90.00° 96.51° 90.00°	Depositor
Resolution (Å)	113.16 – 2.38 113.16 – 2.38	Depositor EDS
% Data completeness (in resolution range)	93.8 (113.16-2.38) 93.9 (113.16-2.38)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.76 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.239 , 0.279 0.237 , 0.276	Depositor DCC
R_{free} test set	6339 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	36.8	Xtriage
Anisotropy	0.031	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 32.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	20048	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.04% 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 ⓘ

5.1 Standard geometry ⓘ

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	1.18	1/2479 (0.0%)	1.01	5/3357 (0.1%)
1	B	1.21	5/2499 (0.2%)	1.02	5/3384 (0.1%)
1	C	1.14	1/2503 (0.0%)	0.99	6/3390 (0.2%)
1	D	1.16	8/2499 (0.3%)	0.97	4/3385 (0.1%)
1	E	1.13	6/2478 (0.2%)	0.98	2/3356 (0.1%)
1	F	1.15	4/2453 (0.2%)	1.01	9/3322 (0.3%)
1	G	1.08	3/2450 (0.1%)	0.95	3/3317 (0.1%)
1	H	1.07	4/2408 (0.2%)	0.95	5/3260 (0.2%)
All	All	1.14	32/19769 (0.2%)	0.99	39/26771 (0.1%)

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	D	0	1

All (32) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	59	HIS	ND1-CE1	7.42	1.40	1.32
1	D	138	HIS	CG-CD2	7.41	1.44	1.35
1	E	138	HIS	CG-CD2	6.81	1.43	1.35
1	H	38	HIS	CG-CD2	6.71	1.43	1.35
1	E	38	HIS	CG-CD2	6.69	1.43	1.35
1	F	138	HIS	CG-CD2	6.40	1.42	1.35
1	H	155	HIS	CG-CD2	6.37	1.42	1.35
1	C	223	THR	CA-CB	5.97	1.56	1.52
1	G	138	HIS	CG-CD2	5.93	1.42	1.35
1	B	138	HIS	CG-CD2	5.71	1.42	1.35
1	D	138	HIS	CE1-NE2	5.60	1.38	1.32
1	D	155	HIS	CG-CD2	5.52	1.42	1.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	231	THR	C-O	-5.49	1.17	1.24
1	D	103	TRP	NE1-CE2	-5.45	1.31	1.37
1	G	59	HIS	CG-CD2	5.42	1.41	1.35
1	B	232	GLY	N-CA	5.41	1.54	1.45
1	F	34	HIS	CG-CD2	5.37	1.41	1.35
1	E	59	HIS	ND1-CE1	5.37	1.38	1.32
1	F	34	HIS	ND1-CE1	5.28	1.37	1.32
1	D	331	HIS	CG-ND1	-5.26	1.32	1.38
1	E	192	HIS	CG-CD2	5.26	1.41	1.35
1	H	138	HIS	CG-CD2	5.25	1.41	1.35
1	F	146	MET	N-CA	-5.24	1.39	1.46
1	H	38	HIS	CE1-NE2	5.19	1.37	1.32
1	B	192	HIS	CG-CD2	5.19	1.41	1.35
1	D	59	HIS	CG-CD2	5.18	1.41	1.35
1	E	138	HIS	CB-CG	5.13	1.57	1.50
1	A	301	GLY	N-CA	5.10	1.50	1.44
1	G	38	HIS	CG-CD2	5.06	1.41	1.35
1	D	192	HIS	CE1-NE2	5.00	1.37	1.32
1	E	301	GLY	N-CA	5.00	1.50	1.44
1	B	217	VAL	N-CA	-5.00	1.40	1.46

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	146	MET	CB-CA-C	7.76	126.53	110.31
1	H	146	MET	CB-CA-C	7.57	126.14	110.31
1	E	307	GLN	CB-CG-CD	7.47	125.29	112.60
1	C	294	ARG	CG-CD-NE	7.30	128.06	112.00
1	F	146	MET	N-CA-CB	-6.87	98.98	110.39
1	G	89	ARG	CG-CD-NE	6.51	126.31	112.00
1	H	146	MET	N-CA-CB	-6.35	99.86	110.39
1	H	89	ARG	CG-CD-NE	6.28	125.81	112.00
1	B	89	ARG	CG-CD-NE	6.27	125.80	112.00
1	C	89	ARG	CG-CD-NE	6.12	125.45	112.00
1	D	89	ARG	CG-CD-NE	6.12	125.45	112.00
1	A	89	ARG	CG-CD-NE	6.10	125.43	112.00
1	F	89	ARG	CG-CD-NE	6.10	125.41	112.00
1	G	144	THR	CA-C-N	-6.02	113.74	119.76
1	G	144	THR	C-N-CA	-6.02	113.74	119.76
1	A	144	THR	CA-C-N	-5.71	113.88	119.76
1	A	144	THR	C-N-CA	-5.71	113.88	119.76
1	B	281	MET	CG-SD-CE	-5.69	88.38	100.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	174	ARG	CG-CD-NE	-5.69	99.49	112.00
1	F	144	THR	CA-C-N	-5.63	113.96	119.76
1	F	144	THR	C-N-CA	-5.63	113.96	119.76
1	F	294	ARG	CG-CD-NE	5.63	124.39	112.00
1	H	146	MET	CA-CB-CG	5.61	125.32	114.10
1	B	213	GLN	CA-CB-CG	5.58	125.26	114.10
1	F	294	ARG	CD-NE-CZ	5.58	132.21	124.40
1	B	294	ARG	CG-CD-NE	5.49	124.08	112.00
1	F	146	MET	CA-CB-CG	5.45	125.00	114.10
1	D	195	SER	CB-CA-C	5.39	118.62	109.84
1	C	281	MET	CG-SD-CE	-5.23	89.39	100.90
1	H	203	VAL	N-CA-C	5.21	115.42	110.42
1	E	294	ARG	CG-CD-NE	5.17	123.37	112.00
1	C	144	THR	CA-C-N	-5.16	114.46	119.78
1	C	144	THR	C-N-CA	-5.16	114.46	119.78
1	C	305	ALA	N-CA-C	5.15	117.56	111.33
1	F	294	ARG	NE-CZ-NH1	5.12	126.62	121.50
1	B	201	ARG	CB-CA-C	-5.07	102.96	111.13
1	D	144	THR	CA-C-N	-5.07	114.69	119.76
1	D	144	THR	C-N-CA	-5.07	114.69	119.76
1	A	307	GLN	CB-CG-CD	-5.06	104.00	112.60

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	89	ARG	Sidechain

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	2429	0	2390	12	0
1	B	2449	0	2414	6	0
1	C	2453	0	2414	13	1
1	D	2447	0	2402	25	1
1	E	2428	0	2391	16	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	F	2403	0	2368	28	0
1	G	2400	0	2362	17	1
1	H	2358	0	2324	30	0
2	A	113	0	0	1	0
2	B	114	0	0	0	1
2	C	93	0	0	0	0
2	D	96	0	0	2	0
2	E	86	0	0	0	0
2	F	63	0	0	3	0
2	G	54	0	0	0	0
2	H	62	0	0	1	0
All	All	20048	0	19065	141	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 4.

All (141) 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:G:146:MET:CE	1:G:178:MET:HE2	1.62	1.29
1:G:146:MET:HE3	1:G:178:MET:CE	1.64	1.27
1:D:146:MET:CE	1:D:178:MET:HE2	1.71	1.19
1:E:146:MET:HE3	1:E:178:MET:CE	1.72	1.16
1:F:146:MET:CE	1:F:177:LYS:HB3	1.78	1.14
1:F:146:MET:HG3	1:F:178:MET:CE	1.79	1.12
1:D:146:MET:HE2	1:D:178:MET:HE2	1.32	1.11
1:F:146:MET:CG	1:F:178:MET:HE1	1.84	1.07
1:E:146:MET:CE	1:E:178:MET:HE2	1.88	1.03
1:C:203:VAL:CG1	1:C:209:LEU:HD21	1.96	0.94
1:H:146:MET:HE1	1:H:174:ARG:HH11	1.32	0.93
1:F:146:MET:HG3	1:F:178:MET:HE1	0.94	0.93
1:D:146:MET:HE3	1:D:178:MET:HE2	1.50	0.92
1:E:146:MET:HE3	1:E:178:MET:HE2	0.92	0.88
1:D:209:LEU:HB2	2:D:2075:HOH:O	1.73	0.87
1:C:203:VAL:CG1	1:C:209:LEU:CD2	2.54	0.85
1:D:146:MET:HE3	1:D:178:MET:CE	2.09	0.82
1:F:146:MET:HE2	1:F:177:LYS:HB3	1.59	0.82
1:H:170:TYR:CE2	1:H:178:MET:HE3	2.15	0.80
1:D:146:MET:CE	1:D:178:MET:CE	2.58	0.80
1:F:146:MET:HE2	1:F:177:LYS:CG	2.12	0.80
1:C:203:VAL:HG11	1:C:209:LEU:HD21	1.63	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:146:MET:CG	1:F:178:MET:CE	2.55	0.75
1:H:198:LEU:HD21	1:H:288:TRP:CD2	2.24	0.73
1:F:146:MET:HE2	1:F:177:LYS:CB	2.18	0.73
1:H:170:TYR:HE2	1:H:178:MET:HE3	1.53	0.72
1:H:198:LEU:HD21	1:H:288:TRP:CG	2.24	0.72
1:H:198:LEU:CD2	1:H:288:TRP:CD2	2.74	0.71
1:H:146:MET:HE1	1:H:174:ARG:NH1	2.08	0.68
1:F:146:MET:HE1	1:F:177:LYS:HB3	1.73	0.67
1:D:283:GLN:NE2	1:D:313:GLN:HB2	2.09	0.66
1:G:146:MET:HE3	1:G:178:MET:HE2	0.76	0.62
1:A:146:MET:HE3	1:A:178:MET:HE2	1.80	0.62
1:F:12:GLN:HG3	2:F:2008:HOH:O	2.00	0.62
1:G:146:MET:HE1	1:G:174:ARG:O	2.00	0.61
1:D:281:MET:CE	1:D:284:MET:HE2	2.30	0.61
1:H:170:TYR:HE2	1:H:178:MET:CE	2.14	0.61
1:C:203:VAL:HG13	1:C:209:LEU:CD2	2.31	0.61
1:F:146:MET:CE	1:F:177:LYS:CB	2.64	0.60
1:B:283:GLN:NE2	1:B:313:GLN:HB2	2.16	0.60
1:D:281:MET:HE1	1:D:284:MET:CE	2.32	0.60
1:F:146:MET:HE1	1:F:174:ARG:HA	1.81	0.60
1:D:281:MET:HE1	1:D:284:MET:HE2	1.83	0.59
1:F:283:GLN:NE2	1:F:310:GLU:HG3	2.18	0.59
1:C:203:VAL:HG13	1:C:209:LEU:HD23	1.85	0.59
1:D:209:LEU:HD13	1:D:294:ARG:HD3	1.85	0.58
1:F:75:GLU:HB3	1:H:79:ARG:NH2	2.19	0.57
1:H:198:LEU:HD23	1:H:288:TRP:CD2	2.40	0.56
1:G:170:TYR:HE2	1:G:178:MET:HE3	1.71	0.55
1:H:203:VAL:CG1	1:H:209:LEU:HG	2.36	0.55
1:H:274:ALA:HB2	1:H:284:MET:HE2	1.89	0.55
1:E:220:ILE:HG12	1:E:299:LEU:HD22	1.90	0.54
1:H:198:LEU:CD2	1:H:288:TRP:CE2	2.90	0.54
1:H:323:LYS:HE2	1:H:327:GLN:HE22	1.72	0.54
1:E:146:MET:CE	1:E:178:MET:CE	2.66	0.54
1:D:263:ASN:ND2	1:D:335:GLY:O	2.41	0.53
1:E:220:ILE:HD11	1:E:299:LEU:HD13	1.90	0.53
1:E:61:ASP:HA	1:E:95:SER:OG	2.10	0.52
1:G:283:GLN:NE2	1:G:313:GLN:HB2	2.25	0.51
1:E:146:MET:HE1	1:E:177:LYS:HB3	1.91	0.51
1:A:61:ASP:HA	1:A:95:SER:OG	2.11	0.51
1:H:198:LEU:HD21	1:H:288:TRP:CD1	2.47	0.50
1:E:203:VAL:HG13	1:E:209:LEU:HG	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:61:ASP:HA	1:C:95:SER:OG	2.10	0.50
1:F:274:ALA:HB2	1:F:284:MET:HE2	1.92	0.50
1:G:170:TYR:CE2	1:G:178:MET:HE3	2.47	0.50
1:F:61:ASP:HA	1:F:95:SER:OG	2.12	0.50
1:G:61:ASP:OD2	1:G:97:LYS:NZ	2.44	0.50
1:H:146:MET:HG3	1:H:178:MET:HE1	1.94	0.50
1:D:61:ASP:HA	1:D:95:SER:OG	2.12	0.49
1:H:61:ASP:HA	1:H:95:SER:OG	2.11	0.49
1:H:203:VAL:HG13	1:H:209:LEU:HG	1.95	0.49
1:E:203:VAL:CG1	1:E:209:LEU:HG	2.43	0.49
1:E:281:MET:HE1	1:E:284:MET:HE2	1.95	0.49
1:F:283:GLN:HE21	1:F:310:GLU:HG3	1.76	0.49
1:H:146:MET:HG3	1:H:178:MET:CE	2.43	0.49
1:G:61:ASP:HA	1:G:95:SER:OG	2.13	0.48
1:H:283:GLN:NE2	1:H:313:GLN:HB2	2.29	0.48
1:F:146:MET:HE3	1:F:177:LYS:HB3	1.83	0.48
1:E:283:GLN:NE2	1:E:313:GLN:HB2	2.29	0.47
1:E:81:LEU:HD21	1:E:129:LEU:HD21	1.96	0.47
1:G:60:PHE:HE2	1:G:92:LEU:HD22	1.79	0.47
1:A:61:ASP:OD2	1:A:97:LYS:NZ	2.46	0.47
1:D:63:ALA:HB3	1:D:66:TYR:CD1	2.50	0.47
1:F:203:VAL:CG1	1:F:209:LEU:HG	2.45	0.47
1:F:267:LEU:O	1:F:281:MET:HE1	2.15	0.47
1:H:146:MET:HE3	1:H:146:MET:HB2	1.63	0.47
1:B:20:LYS:O	1:C:145:PRO:HG3	2.14	0.46
1:H:146:MET:CE	1:H:174:ARG:HH11	2.18	0.46
1:B:61:ASP:HA	1:B:95:SER:OG	2.14	0.46
1:D:281:MET:HE3	1:D:281:MET:HA	1.96	0.46
1:F:146:MET:HE2	1:F:177:LYS:HG2	1.92	0.46
1:F:75:GLU:HB3	1:H:79:ARG:HH22	1.80	0.46
1:G:303:SER:O	1:G:304:ARG:HG3	2.16	0.46
1:C:81:LEU:HD21	1:C:129:LEU:HD21	1.98	0.46
1:C:203:VAL:HG12	1:C:209:LEU:HD21	1.90	0.46
1:D:283:GLN:HE21	1:D:313:GLN:HB2	1.78	0.46
1:A:292:ASP:HB2	2:A:2101:HOH:O	2.16	0.45
1:A:79:ARG:NH2	1:D:75:GLU:HB3	2.31	0.45
1:H:198:LEU:HD21	1:H:288:TRP:CE2	2.49	0.45
1:C:49:LEU:HD22	1:C:60:PHE:CE2	2.51	0.45
1:F:265:ASN:HB3	2:F:2055:HOH:O	2.17	0.45
1:B:199:LEU:HD22	1:B:337:LEU:HB3	1.98	0.44
1:H:146:MET:CE	1:H:174:ARG:NH1	2.79	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:271:ASN:HB2	1:F:281:MET:HE2	2.00	0.44
1:D:146:MET:HE1	1:D:178:MET:HB2	2.00	0.44
1:D:146:MET:CE	1:D:178:MET:HB2	2.47	0.44
1:D:146:MET:HE1	1:D:174:ARG:O	2.17	0.44
1:H:213:GLN:HG2	2:H:2050:HOH:O	2.18	0.43
1:A:146:MET:HE1	1:A:178:MET:HB2	2.01	0.43
1:E:203:VAL:HG22	1:E:208:LEU:HD23	2.01	0.43
1:F:140:VAL:HG21	2:F:2047:HOH:O	2.19	0.43
1:G:146:MET:HG2	1:G:178:MET:HE1	2.00	0.43
1:H:178:MET:HE2	1:H:178:MET:HB2	1.75	0.43
1:C:203:VAL:HG12	1:C:209:LEU:CD2	2.46	0.42
1:A:75:GLU:HB3	1:D:79:ARG:NH2	2.34	0.42
1:A:136:TYR:HA	1:A:166:GLY:O	2.19	0.42
1:F:146:MET:SD	1:F:178:MET:CE	3.07	0.42
1:G:166:GLY:HA2	1:G:189:LEU:HD11	2.02	0.42
1:B:132:VAL:O	1:B:162:ALA:HA	2.20	0.42
1:D:279:GLN:NE2	1:D:317:ASN:HB3	2.35	0.42
1:E:136:TYR:HA	1:E:166:GLY:O	2.19	0.42
1:F:273:MET:CE	1:F:327:GLN:HB2	2.50	0.42
1:D:20:LYS:O	1:H:145:PRO:HG3	2.20	0.42
1:H:136:TYR:HA	1:H:166:GLY:O	2.19	0.42
1:B:263:ASN:ND2	1:B:335:GLY:O	2.53	0.41
1:C:178:MET:CA	1:C:178:MET:HE3	2.50	0.41
1:G:146:MET:HE3	1:G:178:MET:HB2	2.02	0.41
1:A:42:LEU:HD21	1:A:79:ARG:HD3	2.01	0.41
1:A:228:GLY:O	1:A:231:THR:HG23	2.20	0.41
1:E:63:ALA:HB3	1:E:66:TYR:CD1	2.56	0.41
1:A:182:LEU:HD23	1:A:182:LEU:HA	1.91	0.41
1:D:209:LEU:CB	2:D:2075:HOH:O	2.51	0.41
1:F:203:VAL:HG22	1:F:208:LEU:HD23	2.03	0.41
1:G:28:LEU:HD13	1:G:312:VAL:HA	2.01	0.41
1:A:51:LYS:HA	1:A:51:LYS:HD3	1.95	0.40
1:D:136:TYR:HA	1:D:166:GLY:O	2.20	0.40
1:G:136:TYR:HA	1:G:166:GLY:O	2.21	0.40
1:H:51:LYS:HA	1:H:51:LYS:HD3	1.95	0.40
1:G:199:LEU:HD22	1:G:337:LEU:HB3	2.03	0.40
1:C:136:TYR:HA	1:C:166:GLY:O	2.22	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:D:210:ASP:OD2	2:B:2108:HOH:O[1_556]	2.15	0.05
1:C:173:GLU:OE2	1:G:276:GLN:OE1[2_545]	2.19	0.01

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	304/346 (88%)	298 (98%)	6 (2%)	0	100	100
1	B	307/346 (89%)	301 (98%)	5 (2%)	1 (0%)	36	48
1	C	307/346 (89%)	301 (98%)	5 (2%)	1 (0%)	36	48
1	D	306/346 (88%)	301 (98%)	3 (1%)	2 (1%)	18	26
1	E	304/346 (88%)	297 (98%)	6 (2%)	1 (0%)	36	48
1	F	300/346 (87%)	293 (98%)	6 (2%)	1 (0%)	36	48
1	G	300/346 (87%)	293 (98%)	6 (2%)	1 (0%)	36	48
1	H	294/346 (85%)	290 (99%)	3 (1%)	1 (0%)	36	48
All	All	2422/2768 (88%)	2374 (98%)	40 (2%)	8 (0%)	36	48

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	194	PRO
1	E	99	GLY
1	G	99	GLY
1	B	99	GLY
1	F	99	GLY
1	H	99	GLY
1	C	99	GLY
1	D	99	GLY

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	257/290 (89%)	253 (98%)	4 (2%)	55	74
1	B	259/290 (89%)	257 (99%)	2 (1%)	73	85
1	C	260/290 (90%)	256 (98%)	4 (2%)	57	75
1	D	258/290 (89%)	252 (98%)	6 (2%)	44	64
1	E	257/290 (89%)	252 (98%)	5 (2%)	50	69
1	F	255/290 (88%)	252 (99%)	3 (1%)	63	79
1	G	254/290 (88%)	252 (99%)	2 (1%)	73	85
1	H	249/290 (86%)	246 (99%)	3 (1%)	63	79
All	All	2049/2320 (88%)	2020 (99%)	29 (1%)	59	77

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	GLN
1	A	89	ARG
1	A	205	LYS
1	A	312	VAL
1	B	12	GLN
1	B	89	ARG
1	C	4	LEU
1	C	89	ARG
1	C	229	LEU
1	C	261	GLU
1	D	12	GLN
1	D	89	ARG
1	D	203	VAL
1	D	264	LEU
1	D	318	LEU
1	D	325	LEU
1	E	12	GLN
1	E	203	VAL
1	E	261	GLU

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Mol	Chain	Res	Type
1	E	269	LEU
1	E	299	LEU
1	F	89	ARG
1	F	178	MET
1	F	203	VAL
1	G	89	ARG
1	G	173	GLU
1	H	89	ARG
1	H	203	VAL
1	H	229	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	12	GLN
1	A	215	ASN
1	A	263	ASN
1	C	35	ASN
1	C	215	ASN
1	C	338	ASN
1	D	143	ASN
1	D	331	HIS
1	E	176	GLN
1	E	215	ASN
1	E	271	ASN
1	F	35	ASN
1	F	215	ASN
1	G	76	ASN
1	G	215	ASN
1	G	331	HIS
1	H	35	ASN
1	H	143	ASN
1	H	215	ASN
1	H	327	GLN

5.3.3 RNA ⓘ

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](#)

There are no ligands in this entry.

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 ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	307/346 (88%)	-0.19	6 (1%) 65 63	16, 25, 48, 71	2 (0%)
1	B	311/346 (89%)	-0.25	2 (0%) 85 85	16, 25, 48, 69	1 (0%)
1	C	310/346 (89%)	-0.06	6 (1%) 66 65	18, 30, 52, 79	2 (0%)
1	D	308/346 (89%)	0.04	12 (3%) 43 43	18, 29, 56, 78	3 (0%)
1	E	307/346 (88%)	0.14	10 (3%) 49 49	20, 32, 64, 94	2 (0%)
1	F	303/346 (87%)	0.29	14 (4%) 37 37	20, 34, 71, 92	2 (0%)
1	G	303/346 (87%)	0.57	22 (7%) 21 20	24, 39, 80, 99	2 (0%)
1	H	297/346 (85%)	0.73	28 (9%) 14 13	20, 40, 75, 95	2 (0%)
All	All	2446/2768 (88%)	0.16	100 (4%) 41 41	16, 32, 67, 99	16 (0%)

All (100) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	202	TRP	5.0
1	E	204[A]	ASP	5.0
1	H	267	LEU	4.9
1	D	262	ALA	4.6
1	G	204[A]	ASP	4.4
1	H	204[A]	ASP	4.2
1	G	276	GLN	4.1
1	E	202	TRP	4.0
1	B	232	GLY	3.7
1	G	337	LEU	3.7
1	D	207	GLY	3.5
1	G	265	ASN	3.5
1	A	262	ALA	3.4
1	E	276	GLN	3.4
1	H	318	LEU	3.2
1	G	335	GLY	3.2

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Mol	Chain	Res	Type	RSRZ
1	H	199	LEU	3.1
1	H	269	LEU	3.1
1	F	268	ARG	3.1
1	D	202[A]	TRP	3.0
1	H	288	TRP	3.0
1	H	229	LEU	3.0
1	G	275	GLN	3.0
1	A	264	LEU	3.0
1	H	274	ALA	2.9
1	H	230	LEU	2.9
1	G	205	LYS	2.9
1	H	196	TYR	2.9
1	G	305	ALA	2.9
1	F	269	LEU	2.9
1	H	198	LEU	2.9
1	G	202	TRP	2.8
1	H	278	GLY	2.8
1	F	330	GLN	2.7
1	H	20	LYS	2.7
1	G	320	PHE	2.7
1	F	332	ILE	2.7
1	E	333	ALA	2.7
1	F	328	ILE	2.7
1	D	232	GLY	2.6
1	F	202	TRP	2.6
1	H	272	GLU	2.6
1	D	264	LEU	2.6
1	F	229	LEU	2.6
1	H	319	THR	2.6
1	G	267	LEU	2.5
1	G	269	LEU	2.5
1	G	325	LEU	2.5
1	C	12	GLN	2.5
1	G	209	LEU	2.5
1	C	202	TRP	2.5
1	F	322	THR	2.4
1	C	204[A]	ASP	2.4
1	D	79	ARG	2.4
1	D	306	GLU	2.4
1	H	328	ILE	2.4
1	D	276	GLN	2.4
1	H	322	THR	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	272	GLU	2.4
1	D	338	ASN	2.3
1	F	260	THR	2.3
1	F	319	THR	2.3
1	H	280	SER	2.3
1	H	332	ILE	2.3
1	G	213	GLN	2.3
1	F	317	ASN	2.3
1	H	11	GLY	2.3
1	G	333	ALA	2.3
1	G	321	SER	2.3
1	E	332	ILE	2.3
1	E	205	LYS	2.3
1	D	204[A]	ASP	2.3
1	H	334	ASP	2.3
1	A	202	TRP	2.3
1	G	322	THR	2.3
1	C	205	LYS	2.3
1	D	213	GLN	2.2
1	H	268	ARG	2.2
1	F	261	GLU	2.2
1	F	331	HIS	2.2
1	G	173	GLU	2.2
1	H	33	TRP	2.2
1	E	327	GLN	2.2
1	E	331	HIS	2.2
1	H	327	GLN	2.1
1	F	321	SER	2.1
1	B	205	LYS	2.1
1	E	259	LEU	2.1
1	H	209	LEU	2.1
1	E	261	GLU	2.1
1	A	268	ARG	2.1
1	C	304	ARG	2.1
1	H	273	MET	2.1
1	A	204[A]	ASP	2.1
1	G	317	ASN	2.1
1	H	331	HIS	2.0
1	A	229	LEU	2.0
1	D	304	ARG	2.0
1	G	336	GLU	2.0
1	G	270	LEU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.