



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 03:31 AM UTC

PDB ID : 1G9U / pdb_00001g9u
Title : CRYSTAL STRUCTURE OF YOPM-LEUCINE RICH EFFECTOR PROTEIN FROM YERSINIA PESTIS
Authors : Evdokimov, A.G.; Anderson, D.E.; Routzahn, K.M.; Waugh, D.S.
Deposited on : 2000-11-28
Resolution : 2.35 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)	:	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

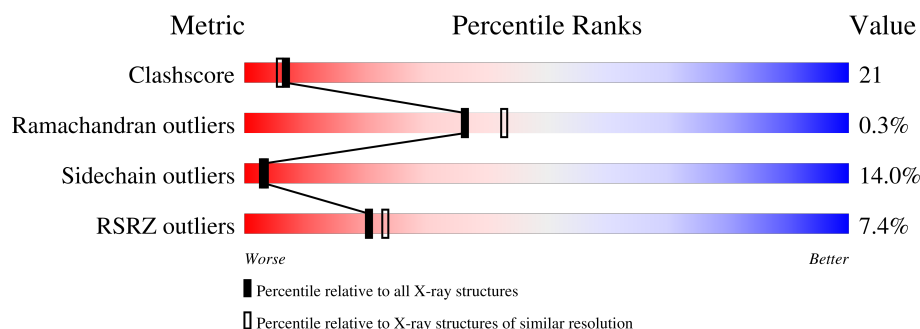
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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	1663 (2.36-2.36)
Ramachandran outliers	187476	1646 (2.36-2.36)
Sidechain outliers	187428	1646 (2.36-2.36)
RSRZ outliers	180081	1598 (2.36-2.36)

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	454	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	ACT	A	1001	-	-	X	-
2	ACT	A	1004	-	-	X	-
2	ACT	A	1005	-	-	X	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	ACT	A	1006	-	-	X	-
2	ACT	A	1008	-	-	X	-
4	HG	A	2006	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 3075 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 OUTER PROTEIN YOPM.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	353	Total	C	N	O	S	0	0	0
			2793	1773	457	558	5			

There are 45 discrepancies between the modelled and reference sequences:

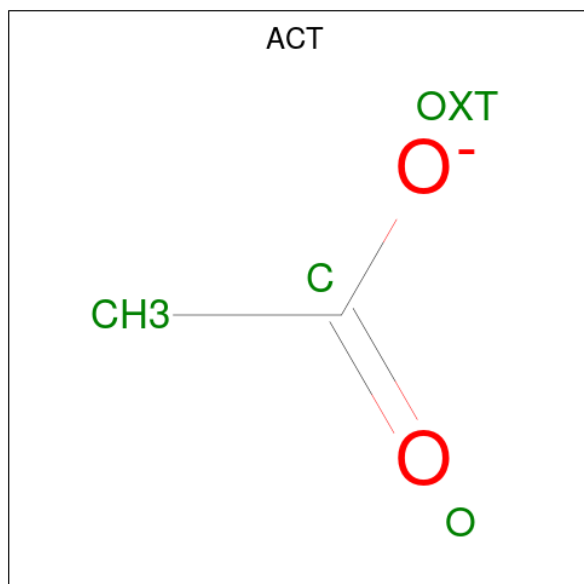
Chain	Residue	Modelled	Actual	Comment	Reference
A	385	ASN	-	SEE REMARK 999	UNP P17778
A	386	SER	-	SEE REMARK 999	UNP P17778
A	1387	HIS	-	SEE REMARK 999	UNP P17778
A	1388	LEU	-	SEE REMARK 999	UNP P17778
A	1389	ALA	-	SEE REMARK 999	UNP P17778
A	1390	GLU	-	SEE REMARK 999	UNP P17778
A	1391	VAL	-	SEE REMARK 999	UNP P17778
A	1392	PRO	-	SEE REMARK 999	UNP P17778
A	1393	GLU	-	SEE REMARK 999	UNP P17778
A	1394	LEU	-	SEE REMARK 999	UNP P17778
A	1395	PRO	-	SEE REMARK 999	UNP P17778
A	1396	GLN	-	SEE REMARK 999	UNP P17778
A	1397	ASN	-	SEE REMARK 999	UNP P17778
A	1398	LEU	-	SEE REMARK 999	UNP P17778
A	1399	LYS	-	SEE REMARK 999	UNP P17778
A	1400	GLN	-	SEE REMARK 999	UNP P17778
A	1401	LEU	-	SEE REMARK 999	UNP P17778
A	1402	HIS	-	SEE REMARK 999	UNP P17778
A	1403	VAL	-	SEE REMARK 999	UNP P17778
A	1404	GLU	-	SEE REMARK 999	UNP P17778
A	1405	THR	-	SEE REMARK 999	UNP P17778
A	1406	ASN	-	SEE REMARK 999	UNP P17778
A	1407	PRO	-	SEE REMARK 999	UNP P17778
A	1408	LEU	-	SEE REMARK 999	UNP P17778
A	1409	ARG	-	SEE REMARK 999	UNP P17778
A	1410	GLU	-	SEE REMARK 999	UNP P17778
A	1411	PHE	-	SEE REMARK 999	UNP P17778

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Chain	Residue	Modelled	Actual	Comment	Reference
A	1412	PRO	-	SEE REMARK 999	UNP P17778
A	1413	ASP	-	SEE REMARK 999	UNP P17778
A	1414	ILE	-	SEE REMARK 999	UNP P17778
A	1415	PRO	-	SEE REMARK 999	UNP P17778
A	1416	GLU	-	SEE REMARK 999	UNP P17778
A	1417	SER	-	SEE REMARK 999	UNP P17778
A	1418	VAL	-	SEE REMARK 999	UNP P17778
A	1419	GLU	-	SEE REMARK 999	UNP P17778
A	1420	ASP	-	SEE REMARK 999	UNP P17778
A	1421	LEU	-	SEE REMARK 999	UNP P17778
A	1422	ARG	-	SEE REMARK 999	UNP P17778
A	1423	MET	-	SEE REMARK 999	UNP P17778
A	1449	HIS	-	expression tag	UNP P17778
A	1450	HIS	-	expression tag	UNP P17778
A	1451	HIS	-	expression tag	UNP P17778
A	1452	HIS	-	expression tag	UNP P17778
A	1453	HIS	-	expression tag	UNP P17778
A	1454	HIS	-	expression tag	UNP P17778

- Molecule 2 is ACETATE ION (CCD ID: ACT) (formula: $C_2H_3O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			4	2	2		
2	A	1	Total	C	O	0	0
			4	2	2		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 4 2 2	0	0
2	A	1	Total C O 4 2 2	0	0
2	A	1	Total C O 4 2 2	0	0
2	A	1	Total C O 4 2 2	0	0
2	A	1	Total C O 4 2 2	0	0
2	A	1	Total C O 4 2 2	0	0
2	A	1	Total C O 4 2 2	0	0

- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	2	Total Ca 2 2	0	0

- Molecule 4 is MERCURY (II) ION (CCD ID: HG) (formula: Hg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	8	Total Hg 8 8	0	0

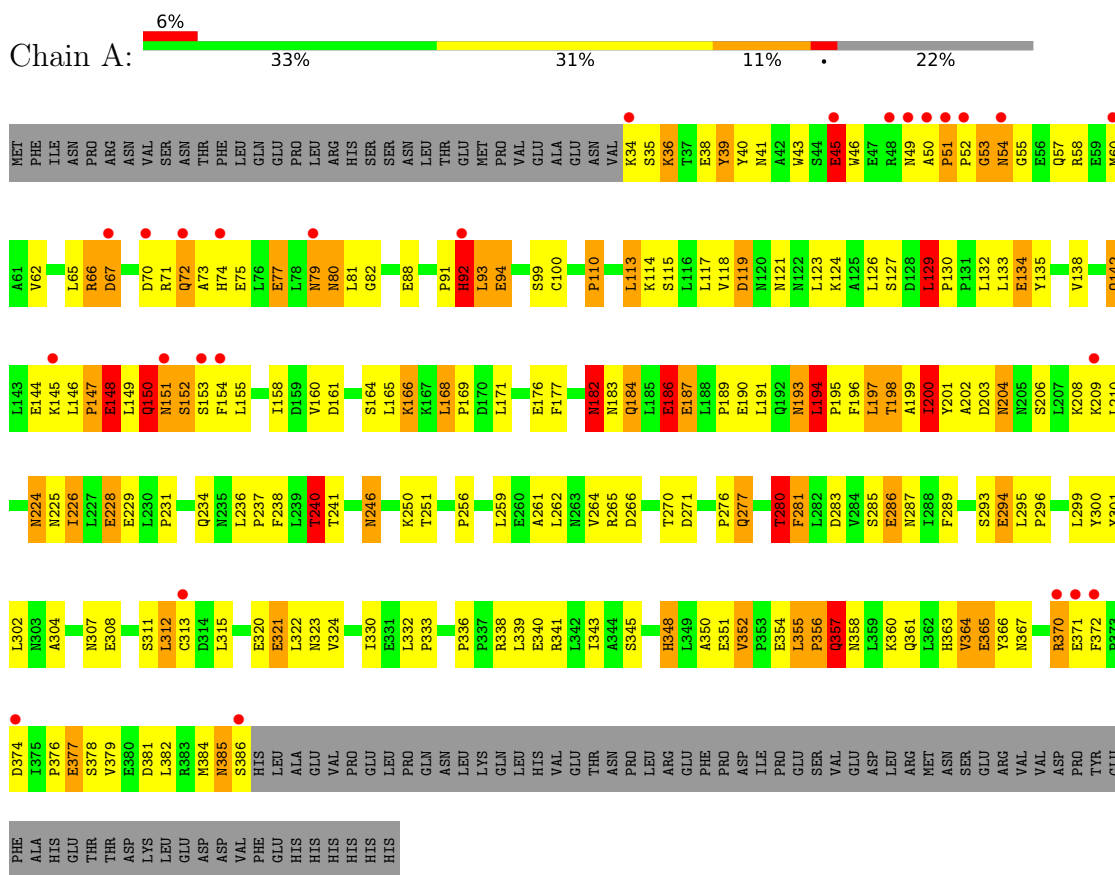
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	236	Total O 236 236	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 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: OUTER PROTEIN YOPM



4 Data and refinement statistics

Property	Value	Source
Space group	P 42 2 2	Depositor
Cell constants a, b, c, α , β , γ	109.36Å 109.36Å 101.50Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	100.00 – 2.35 100.00 – 2.37	Depositor EDS
% Data completeness (in resolution range)	97.0 (100.00-2.35) 92.0 (100.00-2.37)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.70 (at 2.37Å)	Xtriage
Refinement program	SHELXL-97	Depositor
R, R_{free}	0.200 , 0.230 0.186 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	21.6	Xtriage
Anisotropy	0.121	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 112.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3075	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.20% 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

Bond lengths and bond angles in the following residue types are not validated in this section: ACT, HG, CA

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	0.66	0/2853	2.25	142/3903 (3.6%)

There are no bond length outliers.

All (142) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	66	ARG	CD-NE-CZ	18.16	149.83	124.40
1	A	372	PHE	CA-CB-CG	14.14	127.94	113.80
1	A	73	ALA	CA-C-N	13.88	139.32	120.44
1	A	73	ALA	C-N-CA	13.88	139.32	120.44
1	A	45	GLU	CB-CG-CD	10.56	130.55	112.60
1	A	236	LEU	CA-C-O	9.76	127.61	119.71
1	A	364	VAL	CA-C-N	9.54	134.42	120.87
1	A	364	VAL	C-N-CA	9.54	134.42	120.87
1	A	182	ASN	CA-CB-CG	-9.47	103.13	112.60
1	A	240	THR	CA-C-O	9.39	129.20	118.77
1	A	161	ASP	CA-CB-CG	8.75	121.35	112.60
1	A	228	GLU	CA-C-O	8.55	129.27	119.27
1	A	113	LEU	CA-C-O	8.16	130.08	120.70
1	A	53	GLY	CA-C-N	8.15	137.11	121.54
1	A	53	GLY	C-N-CA	8.15	137.11	121.54
1	A	238	PHE	CA-CB-CG	-8.06	105.74	113.80
1	A	280	THR	N-CA-CB	-8.06	100.34	110.98
1	A	200	ILE	CB-CA-C	-7.93	98.75	110.33
1	A	312	LEU	CA-C-N	-7.81	106.61	121.54
1	A	312	LEU	C-N-CA	-7.81	106.61	121.54
1	A	311	SER	O-C-N	7.75	132.09	123.33
1	A	199	ALA	CA-C-N	7.70	133.69	122.99
1	A	199	ALA	C-N-CA	7.70	133.69	122.99
1	A	72	GLN	CB-CG-CD	7.67	125.64	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	165	LEU	O-C-N	7.62	131.89	123.13
1	A	256	PRO	O-C-N	7.59	130.42	121.46
1	A	177	PHE	CA-C-N	-7.50	111.91	122.71
1	A	177	PHE	C-N-CA	-7.50	111.91	122.71
1	A	80	ASN	N-CA-CB	-7.32	100.78	111.91
1	A	358	ASN	CA-C-O	-7.28	111.36	119.95
1	A	287	ASN	CA-CB-CG	-7.27	105.33	112.60
1	A	148	GLU	CA-C-N	7.19	130.63	120.28
1	A	148	GLU	C-N-CA	7.19	130.63	120.28
1	A	264	VAL	CA-C-O	7.12	129.68	120.78
1	A	357	GLN	CA-C-N	-7.08	111.42	122.67
1	A	357	GLN	C-N-CA	-7.08	111.42	122.67
1	A	385	ASN	N-CA-C	7.04	117.81	107.88
1	A	313	CYS	CB-CA-C	6.97	124.29	110.42
1	A	336	PRO	CA-C-O	6.95	130.64	120.56
1	A	250	LYS	CA-C-O	6.93	126.77	119.14
1	A	92	HIS	CA-C-N	-6.85	106.93	121.32
1	A	92	HIS	C-N-CA	-6.85	106.93	121.32
1	A	134	GLU	CA-C-N	6.77	132.31	122.77
1	A	134	GLU	C-N-CA	6.77	132.31	122.77
1	A	280	THR	CB-CA-C	6.69	120.25	109.27
1	A	332	LEU	CA-C-O	-6.52	113.91	120.63
1	A	304	ALA	N-CA-C	-6.44	104.03	112.41
1	A	371	GLU	O-C-N	6.44	131.52	123.15
1	A	228	GLU	N-CA-C	-6.41	105.29	113.23
1	A	176	GLU	CA-C-O	6.38	126.44	119.24
1	A	280	THR	CA-C-N	6.33	131.76	122.39
1	A	280	THR	C-N-CA	6.33	131.76	122.39
1	A	153	SER	CA-C-N	-6.33	112.02	122.08
1	A	153	SER	C-N-CA	-6.33	112.02	122.08
1	A	193	ASN	CA-C-O	6.31	126.74	119.35
1	A	240	THR	N-CA-CB	-6.30	102.52	110.90
1	A	202	ALA	CA-C-O	6.24	126.00	119.51
1	A	225	ASN	CA-C-N	-6.21	110.80	121.97
1	A	225	ASN	C-N-CA	-6.21	110.80	121.97
1	A	147	PRO	CA-C-N	6.20	132.19	122.26
1	A	147	PRO	C-N-CA	6.20	132.19	122.26
1	A	348	HIS	CA-C-N	-6.20	112.55	121.42
1	A	348	HIS	C-N-CA	-6.20	112.55	121.42
1	A	39	TYR	O-C-N	6.19	128.44	122.07
1	A	150	GLN	OE1-CD-NE2	6.13	128.73	122.60
1	A	300	TYR	CA-C-N	-6.10	112.09	122.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	300	TYR	C-N-CA	-6.10	112.09	122.33
1	A	194	LEU	CA-C-O	6.06	125.12	119.24
1	A	358	ASN	CA-CB-CG	-6.05	106.55	112.60
1	A	256	PRO	N-CA-C	-6.05	103.32	110.70
1	A	54	ASN	CA-C-O	-6.00	111.94	120.51
1	A	197	LEU	CA-C-N	5.99	132.67	122.36
1	A	197	LEU	C-N-CA	5.99	132.67	122.36
1	A	229	GLU	O-C-N	5.98	129.82	123.42
1	A	352	VAL	O-C-N	-5.97	114.30	121.10
1	A	294	GLU	CA-C-N	5.96	133.64	121.48
1	A	294	GLU	C-N-CA	5.96	133.64	121.48
1	A	153	SER	O-C-N	-5.95	114.68	122.59
1	A	187	GLU	CA-C-O	-5.93	114.92	121.33
1	A	168	LEU	CA-C-O	-5.90	114.70	120.19
1	A	338	ARG	CA-C-N	-5.88	112.72	120.95
1	A	338	ARG	C-N-CA	-5.88	112.72	120.95
1	A	372	PHE	O-C-N	-5.88	116.19	121.31
1	A	82	GLY	CA-C-N	-5.83	112.78	120.95
1	A	82	GLY	C-N-CA	-5.83	112.78	120.95
1	A	287	ASN	CA-C-N	-5.83	111.47	121.97
1	A	287	ASN	C-N-CA	-5.83	111.47	121.97
1	A	129	LEU	CA-C-O	-5.82	112.19	120.16
1	A	135	TYR	CA-C-O	5.77	126.36	120.24
1	A	355	LEU	CA-C-O	5.76	125.03	120.19
1	A	184	GLN	O-C-N	5.70	129.41	121.83
1	A	374	ASP	CA-CB-CG	5.65	118.25	112.60
1	A	123	LEU	CA-C-O	-5.65	114.83	121.16
1	A	53	GLY	O-C-N	5.64	127.61	122.19
1	A	367	ASN	O-C-N	5.59	126.61	121.80
1	A	286	GLU	CA-C-O	5.59	128.19	121.65
1	A	311	SER	CA-C-O	-5.55	114.78	120.99
1	A	203	ASP	CA-CB-CG	-5.54	107.06	112.60
1	A	281	PHE	CA-CB-CG	-5.50	108.30	113.80
1	A	164	SER	CA-C-N	-5.44	114.96	122.30
1	A	164	SER	C-N-CA	-5.44	114.96	122.30
1	A	308	GLU	N-CA-C	-5.42	103.94	110.88
1	A	296	PRO	O-C-N	-5.41	118.61	121.15
1	A	81	LEU	CA-C-O	5.39	126.47	119.95
1	A	234	GLN	CA-C-N	-5.37	113.73	122.65
1	A	234	GLN	C-N-CA	-5.37	113.73	122.65
1	A	261	ALA	O-C-N	5.36	129.58	123.31
1	A	142	GLN	CA-C-O	5.35	127.65	121.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	151	ASN	CA-C-O	5.35	125.55	119.18
1	A	240	THR	CA-CB-OG1	-5.33	101.60	109.60
1	A	283	ASP	CA-C-O	-5.30	114.83	120.40
1	A	321	GLU	CB-CG-CD	5.29	121.60	112.60
1	A	41	ASN	OD1-CG-ND2	-5.29	117.31	122.60
1	A	190	GLU	CA-C-N	-5.28	113.96	122.24
1	A	190	GLU	C-N-CA	-5.28	113.96	122.24
1	A	356	PRO	CA-C-O	5.27	127.35	121.34
1	A	186	GLU	CA-C-O	5.26	124.92	119.14
1	A	385	ASN	N-CA-CB	-5.24	102.27	111.49
1	A	226	ILE	CA-C-N	-5.24	113.62	120.95
1	A	226	ILE	C-N-CA	-5.24	113.62	120.95
1	A	307	ASN	CA-CB-CG	-5.22	107.38	112.60
1	A	289	PHE	O-C-N	5.19	129.14	123.01
1	A	121	ASN	OD1-CG-ND2	-5.19	117.41	122.60
1	A	199	ALA	O-C-N	5.16	129.36	123.48
1	A	246	ASN	OD1-CG-ND2	-5.16	117.44	122.60
1	A	372	PHE	CA-C-N	-5.13	114.95	120.03
1	A	372	PHE	C-N-CA	-5.13	114.95	120.03
1	A	377	GLU	CA-C-O	-5.13	112.64	119.10
1	A	237	PRO	O-C-N	5.12	129.49	122.38
1	A	171	LEU	N-CA-CB	-5.10	101.63	110.10
1	A	119	ASP	CA-CB-CG	-5.10	107.50	112.60
1	A	228	GLU	O-C-N	-5.10	115.48	122.46
1	A	176	GLU	O-C-N	-5.08	115.55	122.36
1	A	261	ALA	CA-C-O	-5.08	114.68	120.32
1	A	198	THR	N-CA-CB	-5.06	103.86	110.59
1	A	204	ASN	OD1-CG-ND2	5.06	127.66	122.60
1	A	323	ASN	CA-C-O	5.06	125.71	120.40
1	A	51	PRO	N-CA-C	-5.06	104.53	110.70
1	A	183	ASN	CA-CB-CG	-5.06	107.54	112.60
1	A	201	TYR	O-C-N	5.05	129.13	123.27
1	A	152	SER	CA-C-N	5.01	131.11	121.54
1	A	152	SER	C-N-CA	5.01	131.11	121.54

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	2793	0	2781	115	0
2	A	36	0	27	12	0
3	A	2	0	0	0	0
4	A	8	0	0	1	0
5	A	236	0	0	13	0
All	All	3075	0	2808	119	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 21.

All (119) 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:92:HIS:CE1	4:A:2004:HG:HG	1.68	1.06
1:A:35:SER:OG	1:A:38:GLU:HG3	1.84	0.78
1:A:62:VAL:O	1:A:66:ARG:HG3	1.84	0.77
1:A:93:LEU:HD12	1:A:110:PRO:HG2	1.68	0.76
1:A:77:GLU:HG2	1:A:79:ASN:OD1	1.86	0.75
1:A:271:ASP:HB3	2:A:1003:ACT:H3	1.73	0.70
1:A:182:ASN:HD22	1:A:204:ASN:HB3	1.58	0.69
1:A:144:GLU:O	1:A:145:LYS:HD2	1.94	0.68
1:A:315:LEU:HD23	2:A:1008:ACT:H3	1.75	0.67
1:A:155:LEU:HD13	1:A:158:ILE:HD11	1.75	0.66
1:A:94:GLU:HG2	1:A:114:LYS:HD2	1.75	0.66
1:A:130:PRO:HB2	5:A:3144:HOH:O	1.97	0.64
1:A:266:ASP:HA	1:A:286:GLU:O	1.98	0.63
1:A:376:PRO:HG2	1:A:379:VAL:HG23	1.78	0.63
1:A:80:ASN:HA	1:A:100:CYS:O	1.99	0.63
1:A:67:ASP:O	1:A:71:ARG:HG3	1.98	0.63
1:A:366:TYR:H	1:A:385:ASN:HB3	1.63	0.63
1:A:382:LEU:HD11	1:A:384:MET:HE2	1.80	0.62
1:A:312:LEU:HB2	1:A:333:PRO:HD3	1.83	0.61
1:A:45:GLU:HG2	1:A:46:TRP:N	2.16	0.60
1:A:138:VAL:HG23	1:A:160:VAL:HG12	1.82	0.60
1:A:382:LEU:HD11	1:A:384:MET:CE	2.32	0.60
1:A:138:VAL:CG2	1:A:160:VAL:HG12	2.33	0.59
1:A:118:VAL:HG23	1:A:138:VAL:HG12	1.84	0.59
1:A:124:LYS:HE2	1:A:144:GLU:OE2	2.03	0.59
1:A:66:ARG:HD3	5:A:3130:HOH:O	2.02	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:147:PRO:HB2	2:A:1005:ACT:O	2.03	0.58
1:A:321:GLU:HG2	1:A:341:ARG:HB2	1.84	0.58
1:A:210:LEU:HB2	1:A:231:PRO:HD3	1.85	0.58
1:A:126:LEU:HD13	1:A:129:LEU:HD21	1.85	0.58
1:A:195:PRO:HG2	1:A:196:PHE:CD1	2.39	0.57
1:A:350:ALA:O	1:A:351:GLU:HG3	2.05	0.57
1:A:150:GLN:H	1:A:150:GLN:HE21	1.51	0.56
2:A:1005:ACT:H1	5:A:3218:HOH:O	2.06	0.56
1:A:193:ASN:HB2	5:A:3080:HOH:O	2.03	0.56
1:A:149:LEU:HD13	1:A:158:ILE:HD13	1.87	0.56
1:A:39:TYR:OH	1:A:91:PRO:HG3	2.06	0.56
1:A:52:PRO:HD2	5:A:3230:HOH:O	2.07	0.55
1:A:330:ILE:HG22	1:A:348:HIS:HB3	1.89	0.55
1:A:194:LEU:HD23	1:A:197:LEU:HD22	1.90	0.54
1:A:345:SER:OG	1:A:365:GLU:HG3	2.08	0.54
1:A:350:ALA:HB1	1:A:370:ARG:HG3	1.89	0.54
1:A:113:LEU:HB3	1:A:130:PRO:CG	2.38	0.53
1:A:352:VAL:CG1	1:A:355:LEU:HD21	2.38	0.53
1:A:49:ASN:HB3	2:A:1004:ACT:OXT	2.08	0.52
1:A:182:ASN:HD22	1:A:204:ASN:CB	2.21	0.52
2:A:1004:ACT:H1	5:A:3127:HOH:O	2.08	0.52
1:A:186:GLU:OE1	1:A:208:LYS:HE2	2.09	0.52
1:A:50:ALA:HB2	1:A:58:ARG:HD2	1.92	0.51
1:A:197:LEU:HD21	1:A:200:ILE:HD13	1.93	0.51
1:A:341:ARG:HD2	5:A:3195:HOH:O	2.11	0.51
1:A:351:GLU:HG3	1:A:370:ARG:HB2	1.93	0.50
1:A:382:LEU:CD1	1:A:384:MET:HG3	2.41	0.50
1:A:149:LEU:HD12	1:A:169:PRO:HG2	1.93	0.50
1:A:354:GLU:HA	5:A:3183:HOH:O	2.10	0.50
1:A:382:LEU:HD11	1:A:384:MET:HG3	1.94	0.50
1:A:113:LEU:HB3	1:A:130:PRO:HG2	1.94	0.50
1:A:186:GLU:OE2	1:A:206:SER:O	2.30	0.50
1:A:151:ASN:HB2	2:A:1005:ACT:H2	1.93	0.49
1:A:67:ASP:OD2	1:A:71:ARG:HD2	2.12	0.49
1:A:281:PHE:CD1	1:A:301:TYR:HB3	2.47	0.49
1:A:118:VAL:CG2	1:A:138:VAL:HG12	2.43	0.49
1:A:295:LEU:HG	2:A:1008:ACT:H2	1.94	0.49
1:A:345:SER:HG	1:A:365:GLU:HG3	1.78	0.48
1:A:132:LEU:HB2	5:A:3144:HOH:O	2.14	0.48
1:A:343:ILE:HD11	1:A:361:GLN:OE1	2.13	0.48
1:A:364:VAL:O	1:A:364:VAL:HG22	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:99:SER:HB3	1:A:119:ASP:OD1	2.14	0.48
1:A:295:LEU:HD13	1:A:302:LEU:HD22	1.95	0.47
1:A:277:GLN:HG2	5:A:3207:HOH:O	2.14	0.47
1:A:240:THR:HG22	1:A:241:THR:OG1	2.14	0.47
1:A:182:ASN:ND2	1:A:204:ASN:HB3	2.29	0.47
1:A:40:TYR:HH	1:A:70:ASP:CG	2.19	0.47
1:A:191:LEU:HD13	1:A:200:ILE:CD1	2.45	0.47
1:A:352:VAL:HG11	1:A:355:LEU:HD21	1.97	0.47
1:A:146:LEU:HB3	1:A:147:PRO:HD2	1.96	0.47
1:A:166:LYS:HE2	1:A:184:GLN:OE1	2.15	0.47
1:A:94:GLU:CG	1:A:114:LYS:HD2	2.43	0.46
1:A:145:LYS:HE3	5:A:3056:HOH:O	2.14	0.46
1:A:280:THR:HA	1:A:299:LEU:HA	1.96	0.46
1:A:126:LEU:HD23	1:A:126:LEU:HA	1.79	0.46
1:A:384:MET:O	2:A:1001:ACT:H2	2.15	0.45
2:A:1006:ACT:H3	5:A:3047:HOH:O	2.16	0.45
1:A:265:ARG:HH11	1:A:265:ARG:HD2	1.66	0.45
1:A:320:GLU:C	1:A:339:LEU:HD12	2.42	0.45
1:A:365:GLU:HG3	1:A:365:GLU:H	1.37	0.45
1:A:129:LEU:HB3	1:A:130:PRO:HD2	1.98	0.45
1:A:134:GLU:HA	1:A:154:PHE:O	2.17	0.45
1:A:57:GLN:HB3	1:A:60:MET:SD	2.56	0.45
1:A:148:GLU:OE1	1:A:150:GLN:NE2	2.50	0.45
1:A:385:ASN:ND2	2:A:1001:ACT:O	2.49	0.45
1:A:312:LEU:HD23	1:A:312:LEU:HA	1.66	0.44
1:A:376:PRO:HG2	1:A:379:VAL:CG2	2.47	0.44
1:A:99:SER:O	1:A:100:CYS:HB2	2.18	0.44
1:A:280:THR:HG23	1:A:281:PHE:CE2	2.52	0.44
1:A:355:LEU:HA	1:A:356:PRO:HD3	1.89	0.44
1:A:187:GLU:N	5:A:3209:HOH:O	2.50	0.44
1:A:322:LEU:HD11	1:A:324:VAL:HG13	1.99	0.44
2:A:1006:ACT:O	2:A:1009:ACT:H1	2.17	0.44
1:A:36:LYS:HD3	1:A:40:TYR:OH	2.18	0.43
1:A:197:LEU:HD21	1:A:200:ILE:CD1	2.48	0.43
1:A:74:HIS:CD2	1:A:75:GLU:H	2.36	0.43
1:A:53:GLY:O	1:A:55:GLY:N	2.49	0.43
1:A:168:LEU:HB2	1:A:189:PRO:HD3	2.01	0.42
1:A:312:LEU:HD11	1:A:324:VAL:HG11	2.01	0.42
1:A:350:ALA:O	1:A:370:ARG:HB2	2.19	0.42
1:A:51:PRO:HA	1:A:52:PRO:HD3	1.58	0.42
1:A:265:ARG:HD2	1:A:285:SER:OG	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:224:ASN:HD22	1:A:246:ASN:HB3	1.85	0.41
1:A:166:LYS:HE2	1:A:184:GLN:CD	2.45	0.41
1:A:364:VAL:O	1:A:364:VAL:HG13	2.21	0.41
1:A:114:LYS:C	1:A:133:LEU:HD12	2.46	0.41
1:A:259:LEU:HD21	1:A:262:LEU:HB2	2.02	0.41
1:A:43:TRP:CZ3	1:A:65:LEU:HB3	2.55	0.40
1:A:357:GLN:OE1	1:A:357:GLN:HA	2.08	0.40
1:A:363:HIS:HE2	1:A:381:ASP:CG	2.29	0.40
1:A:110:PRO:O	1:A:113:LEU:HB2	2.21	0.40
1:A:119:ASP:OD1	1:A:119:ASP:N	2.49	0.40
1:A:58:ARG:O	1:A:62:VAL:N	2.48	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	351/454 (77%)	314 (90%)	36 (10%)	1 (0%)	36	43

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	54	ASN

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	329/427 (77%)	283 (86%)	46 (14%)	3 3

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	34	LYS
1	A	36	LYS
1	A	45	GLU
1	A	67	ASP
1	A	72	GLN
1	A	77	GLU
1	A	79	ASN
1	A	88	GLU
1	A	92	HIS
1	A	93	LEU
1	A	94	GLU
1	A	110	PRO
1	A	115	SER
1	A	117	LEU
1	A	127	SER
1	A	129	LEU
1	A	142	GLN
1	A	148	GLU
1	A	150	GLN
1	A	152	SER
1	A	166	LYS
1	A	182	ASN
1	A	186	GLU
1	A	194	LEU
1	A	198	THR
1	A	200	ILE
1	A	209	LYS
1	A	224	ASN
1	A	226	ILE
1	A	228	GLU
1	A	240	THR
1	A	251	THR
1	A	270	THR
1	A	276	PRO
1	A	277	GLN
1	A	280	THR
1	A	293	SER
1	A	294	GLU

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Mol	Chain	Res	Type
1	A	340	GLU
1	A	357	GLN
1	A	360	LYS
1	A	365	GLU
1	A	370	ARG
1	A	377	GLU
1	A	378	SER
1	A	386	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	74	HIS
1	A	80	ASN
1	A	122	ASN
1	A	150	GLN
1	A	182	ASN
1	A	192	GLN
1	A	224	ASN
1	A	357	GLN
1	A	358	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 19 ligands modelled in this entry, 10 are monoatomic - leaving 9 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$
2	ACT	A	1005	-	3,3,3	1.29	1 (33%)	3,3,3	0.31	0
2	ACT	A	1002	-	3,3,3	1.43	1 (33%)	3,3,3	1.89	1 (33%)
2	ACT	A	1008	-	3,3,3	1.42	1 (33%)	3,3,3	1.94	1 (33%)
2	ACT	A	1001	-	3,3,3	1.32	1 (33%)	3,3,3	0.69	0
2	ACT	A	1009	-	3,3,3	1.32	1 (33%)	3,3,3	0.48	0
2	ACT	A	1003	-	3,3,3	1.43	1 (33%)	3,3,3	0.93	0
2	ACT	A	1007	-	3,3,3	1.25	1 (33%)	3,3,3	0.58	0
2	ACT	A	1004	-	3,3,3	1.37	1 (33%)	3,3,3	1.52	1 (33%)
2	ACT	A	1006	-	3,3,3	1.31	1 (33%)	3,3,3	1.20	0

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1008	ACT	O-C	2.44	1.32	1.22
2	A	1002	ACT	O-C	2.43	1.32	1.22
2	A	1003	ACT	O-C	2.41	1.32	1.22
2	A	1004	ACT	O-C	2.29	1.32	1.22
2	A	1009	ACT	O-C	2.21	1.31	1.22
2	A	1006	ACT	O-C	2.20	1.31	1.22
2	A	1005	ACT	O-C	2.14	1.31	1.22
2	A	1001	ACT	O-C	2.12	1.31	1.22
2	A	1007	ACT	O-C	2.12	1.31	1.22

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1002	ACT	OXT-C-O	2.79	132.38	122.03
2	A	1008	ACT	OXT-C-O	2.78	132.35	122.03
2	A	1004	ACT	OXT-C-O	2.18	130.12	122.03

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

7 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1005	ACT	3	0
2	A	1008	ACT	2	0
2	A	1001	ACT	2	0
2	A	1009	ACT	1	0
2	A	1003	ACT	1	0
2	A	1004	ACT	2	0
2	A	1006	ACT	2	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 ⓘ

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	353/454 (77%)	0.36	26 (7%) 20 23	12, 28, 67, 111	0

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	372	PHE	6.0
1	A	386	SER	4.4
1	A	92	HIS	3.9
1	A	72	GLN	3.8
1	A	145	LYS	3.7
1	A	313	CYS	3.6
1	A	74	HIS	3.2
1	A	60	MET	3.2
1	A	45	GLU	3.1
1	A	49	ASN	3.1
1	A	370	ARG	3.0
1	A	374	ASP	2.9
1	A	151	ASN	2.8
1	A	48	ARG	2.8
1	A	67	ASP	2.6
1	A	79	ASN	2.6
1	A	154	PHE	2.5
1	A	34	LYS	2.5
1	A	371	GLU	2.5
1	A	153	SER	2.4
1	A	70	ASP	2.3
1	A	51	PRO	2.3
1	A	52	PRO	2.2
1	A	209	LYS	2.2
1	A	50	ALA	2.1
1	A	54	ASN	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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	HG	A	2006	1/1	0.47	0.50	112,112,112,112	1
2	ACT	A	1009	4/4	0.61	0.33	80,84,87,88	0
4	HG	A	2009	1/1	0.68	0.38	96,96,96,96	1
4	HG	A	2010	1/1	0.71	0.32	157,157,157,157	1
2	ACT	A	1005	4/4	0.79	0.20	56,74,78,83	0
2	ACT	A	1004	4/4	0.87	0.24	51,70,72,75	0
2	ACT	A	1001	4/4	0.87	0.26	66,69,80,85	0
4	HG	A	2008	1/1	0.89	0.27	92,92,92,92	1
4	HG	A	2007	1/1	0.92	0.18	46,46,46,46	1
2	ACT	A	1006	4/4	0.92	0.14	37,43,53,68	0
2	ACT	A	1002	4/4	0.93	0.16	16,51,55,61	0
2	ACT	A	1008	4/4	0.93	0.12	36,45,51,56	0
2	ACT	A	1003	4/4	0.94	0.12	39,44,49,49	0
4	HG	A	2005	1/1	0.95	0.06	34,34,34,34	1
2	ACT	A	1007	4/4	0.96	0.13	23,51,55,92	0
4	HG	A	2004	1/1	0.98	0.07	29,29,29,29	1
3	CA	A	2002	1/1	0.99	0.04	22,22,22,22	0
4	HG	A	2003	1/1	0.99	0.03	47,47,47,47	0
3	CA	A	2001	1/1	0.99	0.03	20,20,20,20	0

6.5 Other polymers [i](#)

There are no such residues in this entry.