



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:18 AM UTC

PDB ID : 6I8S / pdb_00006i8s
Title : Discovery and characterisation of an antibody that selectively modulates the inhibitory activity of plasminogen activator inhibitor-1
Authors : Vousden, K.A.; Lundqvist, T.; Popovic, B.
Deposited on : 2018-11-21
Resolution : 2.90 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:

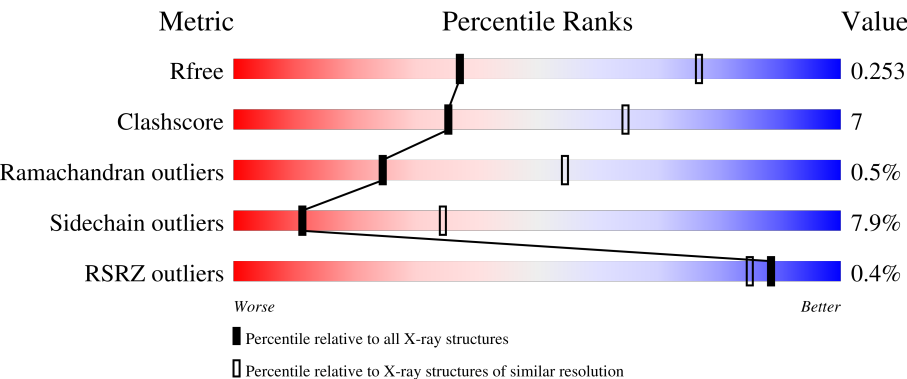
MolProbity : 4-5-2 with Phenix2.0
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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	379	80%14% . .
1	B	379	78%16% . . .
1	C	379	% 78%16% . .
1	D	379	% 75%16%5% .
2	E	223	% 73%19% . .

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Mol	Chain	Length	Quality of chain
2	F	223	% 61%30%5% .
2	G	223	83%10% . .
2	H	223	82%12% . .
3	I	214	84%11% . .
3	J	214	75%21% . .
3	K	214	85%13% ..
3	L	214	89%9% .

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 24618 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 activator inhibitor 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	363	Total	C	N	O	S	0	0	0
			2901	1862	495	529	15			
1	B	363	Total	C	N	O	S	0	0	0
			2901	1862	495	529	15			
1	C	363	Total	C	N	O	S	0	0	0
			2901	1862	495	529	15			
1	D	363	Total	C	N	O	S	0	0	0
			2901	1862	495	529	15			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	150	HIS	ASN	engineered mutation	UNP P05121
A	154	THR	LYS	engineered mutation	UNP P05121
A	319	LEU	GLN	engineered mutation	UNP P05121
A	354	ILE	MET	engineered mutation	UNP P05121
B	150	HIS	ASN	engineered mutation	UNP P05121
B	154	THR	LYS	engineered mutation	UNP P05121
B	319	LEU	GLN	engineered mutation	UNP P05121
B	354	ILE	MET	engineered mutation	UNP P05121
C	150	HIS	ASN	engineered mutation	UNP P05121
C	154	THR	LYS	engineered mutation	UNP P05121
C	319	LEU	GLN	engineered mutation	UNP P05121
C	354	ILE	MET	engineered mutation	UNP P05121
D	150	HIS	ASN	engineered mutation	UNP P05121
D	154	THR	LYS	engineered mutation	UNP P05121
D	319	LEU	GLN	engineered mutation	UNP P05121
D	354	ILE	MET	engineered mutation	UNP P05121

- Molecule 2 is a protein called HEAVY CHAIN OF FAB FRAGMENT FROM AN PAI-1 ANTIBODY.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	214	Total	C	N	O	S	0	0	0
			1614	1021	274	313	6			
2	F	214	Total	C	N	O	S	0	0	0
			1614	1021	274	313	6			
2	G	214	Total	C	N	O	S	0	0	0
			1614	1021	274	313	6			
2	H	214	Total	C	N	O	S	0	0	0
			1614	1021	274	313	6			

- Molecule 3 is a protein called LIGHT CHAIN OF FAB FRAGMENT FROM AN PAI-1 ANTIBODY.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	I	211	Total	C	N	O	S	0	0	0
			1628	1022	274	327	5			
3	J	211	Total	C	N	O	S	0	0	0
			1628	1022	274	327	5			
3	K	211	Total	C	N	O	S	0	0	0
			1628	1022	274	327	5			
3	L	211	Total	C	N	O	S	0	0	0
			1628	1022	274	327	5			

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	6	Total	O	0	0
			6	6		
4	B	5	Total	O	0	0
			5	5		
4	C	7	Total	O	0	0
			7	7		
4	D	5	Total	O	0	0
			5	5		
4	F	2	Total	O	0	0
			2	2		
4	H	3	Total	O	0	0
			3	3		
4	I	4	Total	O	0	0
			4	4		
4	J	3	Total	O	0	0
			3	3		
4	K	7	Total	O	0	0
			7	7		

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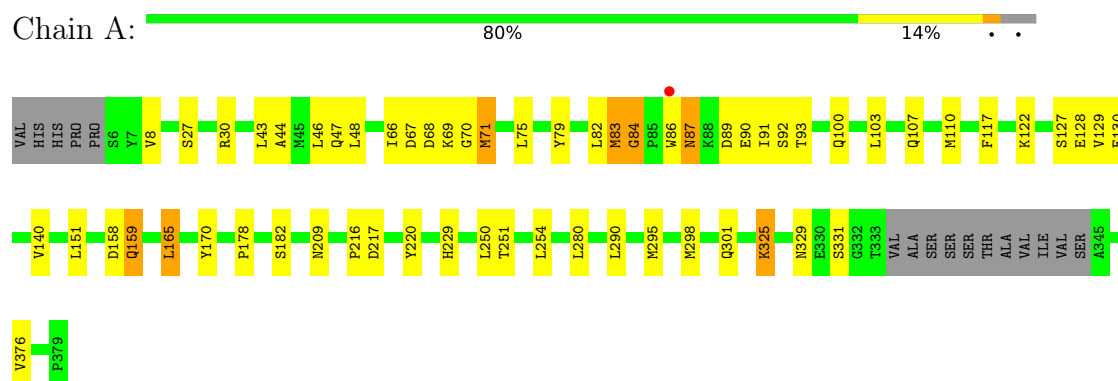
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	L	4	Total	O	0	0
			4	4		

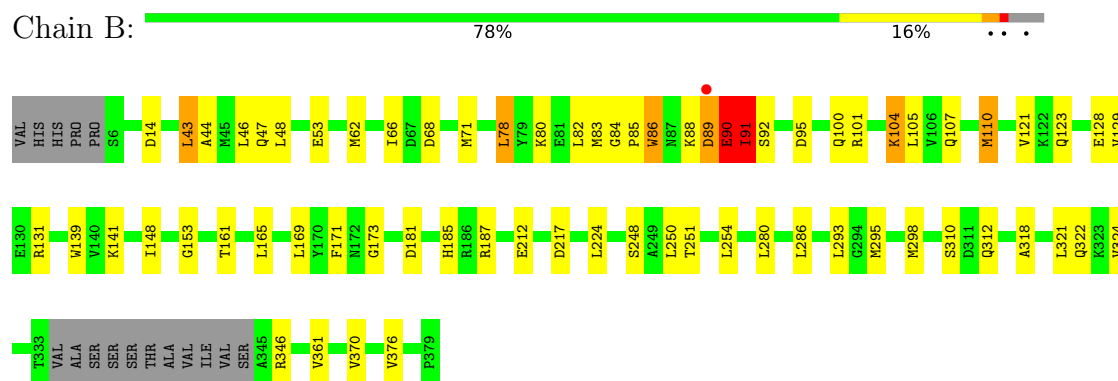
3 Residue-property plots

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.

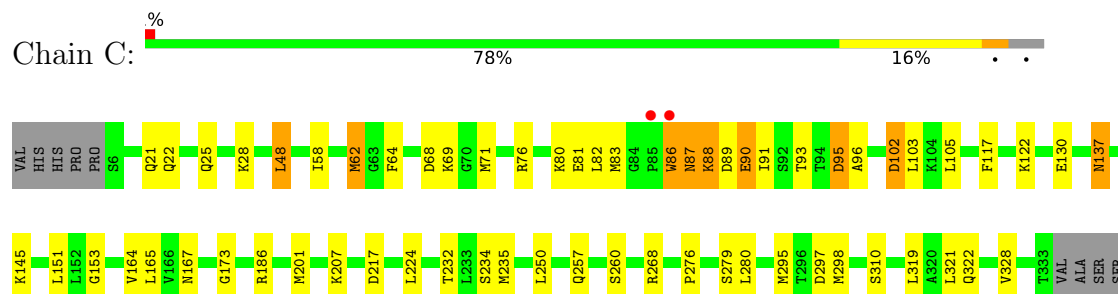
• Molecule 1: Plasminogen activator inhibitor 1



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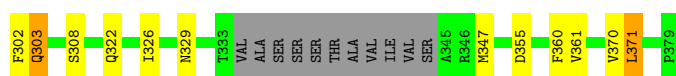
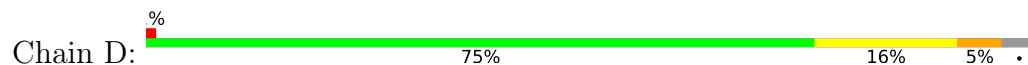


• Molecule 1: Plasminogen activator inhibitor 1

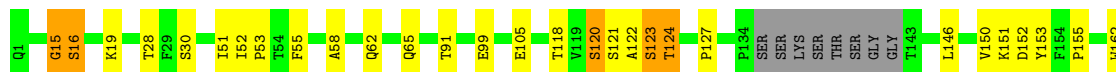
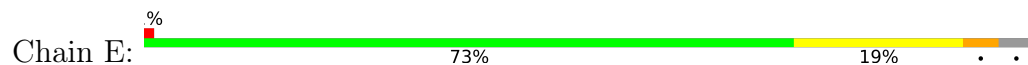




- Molecule 1: Plasminogen activator inhibitor 1



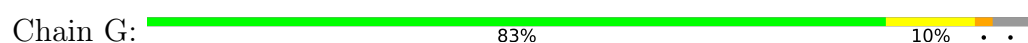
- Molecule 2: HEAVY CHAIN OF FAB FRAGMENT FROM AN PAI-1 ANTIBODY



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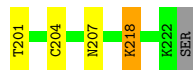
- Molecule 2: HEAVY CHAIN OF FAB FRAGMENT FROM AN PAI-1 ANTIBODY





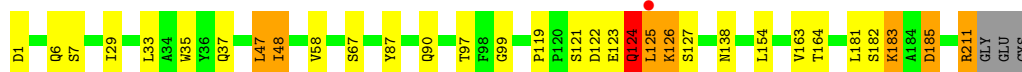
• Molecule 2: HEAVY CHAIN OF FAB FRAGMENT FROM AN PAI-1 ANTIBODY

Chain H: 82% 12% ..



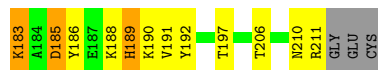
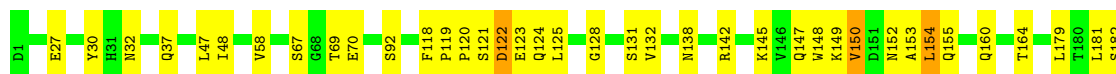
• Molecule 3: LIGHT CHAIN OF FAB FRAGMENT FROM AN PAI-1 ANTIBODY

Chain I: 84% 11% ..



• Molecule 3: LIGHT CHAIN OF FAB FRAGMENT FROM AN PAI-1 ANTIBODY

Chain J: 75% 21% ..



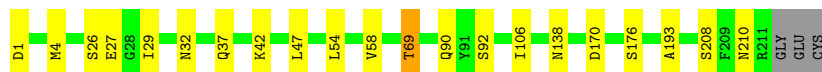
• Molecule 3: LIGHT CHAIN OF FAB FRAGMENT FROM AN PAI-1 ANTIBODY

Chain K: 85% 13% ..



• Molecule 3: LIGHT CHAIN OF FAB FRAGMENT FROM AN PAI-1 ANTIBODY

Chain L: 89% 9% ..



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	90.22Å 89.66Å 249.77Å 90.00° 99.35° 90.00°	Depositor
Resolution (Å)	14.99 – 2.90 14.99 – 2.90	Depositor EDS
% Data completeness (in resolution range)	99.2 (14.99-2.90) 98.4 (14.99-2.90)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.72 (at 2.91Å)	Xtriage
Refinement program	BUSTER 2.10.0	Depositor
R, R_{free}	0.219 , 0.243 0.228 , 0.253	Depositor DCC
R_{free} test set	4317 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	65.0	Xtriage
Anisotropy	0.068	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 34.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.028 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	24618	wwPDB-VP
Average B, all atoms (Å ²)	63.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.93% 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	0.80	1/2971 (0.0%)	1.28	9/4025 (0.2%)
1	B	0.80	1/2971 (0.0%)	1.25	6/4025 (0.1%)
1	C	0.74	0/2971	1.24	3/4025 (0.1%)
1	D	0.77	0/2971	1.33	9/4025 (0.2%)
2	E	0.88	0/1653	1.29	17/2250 (0.8%)
2	F	1.08	0/1653	1.29	9/2250 (0.4%)
2	G	0.75	0/1653	1.16	3/2250 (0.1%)
2	H	0.66	0/1653	1.14	2/2250 (0.1%)
3	I	0.78	1/1664 (0.1%)	1.11	4/2259 (0.2%)
3	J	0.77	1/1664 (0.1%)	1.17	5/2259 (0.2%)
3	K	0.83	0/1664	1.10	2/2259 (0.1%)
3	L	0.72	0/1664	1.10	4/2259 (0.2%)
All	All	0.80	4/25152 (0.0%)	1.22	73/34136 (0.2%)

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	D	0	1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	91	ILE	C-O	-5.96	1.17	1.24
3	J	118	PHE	C-N	5.70	1.38	1.33
3	I	124	GLN	CA-C	-5.43	1.46	1.52
1	A	220	TYR	C-O	-5.39	1.17	1.23

All (73) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	90	GLU	N-CA-C	13.09	125.27	111.14
1	D	26	ALA	N-CA-C	10.03	122.17	111.03
1	A	117	PHE	N-CA-C	8.90	124.76	112.93
1	D	26	ALA	CA-C-N	8.59	137.17	121.70
1	D	26	ALA	C-N-CA	8.59	137.17	121.70
1	C	88	LYS	N-CA-C	8.15	122.89	113.19
2	F	154	PHE	C-N-CD	7.64	137.42	120.60
2	G	30	SER	N-CA-C	7.64	119.24	111.07
1	A	84	GLY	C-N-CD	7.63	137.39	120.60
2	E	120	SER	CA-C-N	7.62	135.42	121.70
2	E	120	SER	C-N-CA	7.62	135.42	121.70
2	F	153	TYR	N-CA-C	7.49	121.06	110.50
2	F	156	GLN	C-N-CD	7.18	136.39	120.60
2	E	15	GLY	CA-C-N	7.17	134.60	121.70
2	E	15	GLY	C-N-CA	7.17	134.60	121.70
3	J	119	PRO	O-C-N	7.17	124.61	121.31
3	L	69	THR	CB-CA-C	6.87	120.61	109.07
2	E	123	SER	N-CA-C	6.84	118.39	111.07
3	J	118	PHE	CA-C-N	-6.38	115.00	119.66
3	J	118	PHE	C-N-CA	-6.38	115.00	119.66
3	J	189	HIS	N-CA-C	6.32	119.95	109.96
1	B	90	GLU	N-CA-C	-6.08	104.94	112.72
1	B	129	VAL	N-CA-C	5.92	116.57	110.36
2	E	151	LYS	N-CA-C	5.80	118.67	109.50
2	E	123	SER	CA-C-N	5.80	132.61	121.54
2	E	123	SER	C-N-CA	5.80	132.61	121.54
2	F	28	THR	CB-CA-C	5.78	117.74	109.71
3	I	124	GLN	N-CA-C	-5.77	104.91	111.14
3	I	119	PRO	CA-C-N	-5.69	114.11	119.85
3	I	119	PRO	C-N-CA	-5.69	114.11	119.85
2	F	52	ILE	N-CA-C	-5.58	102.03	107.55
1	A	129	VAL	N-CA-C	5.56	116.02	110.23
1	C	90	GLU	N-CA-C	-5.50	106.62	113.55
3	L	210	ASN	CA-C-N	5.48	131.56	121.70
3	L	210	ASN	C-N-CA	5.48	131.56	121.70
2	F	184	TYR	N-CA-C	5.47	117.88	109.07
2	H	52	ILE	N-CA-C	-5.45	102.16	107.55
1	B	128	GLU	CA-C-N	5.44	127.62	120.60
1	B	128	GLU	C-N-CA	5.44	127.62	120.60
2	E	192	VAL	CA-C-N	-5.42	114.34	119.76
2	E	192	VAL	C-N-CA	-5.42	114.34	119.76
3	J	122	ASP	N-CA-C	-5.42	104.92	112.45
2	H	30	SER	N-CA-C	5.41	117.61	111.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	117	PHE	CA-C-N	5.40	131.42	121.70
1	A	117	PHE	C-N-CA	5.40	131.42	121.70
2	E	55	PHE	CA-CB-CG	5.39	119.19	113.80
2	G	52	ILE	N-CA-C	-5.39	102.22	107.55
1	A	128	GLU	CA-C-N	5.36	127.79	120.77
1	A	128	GLU	C-N-CA	5.36	127.79	120.77
1	D	296	THR	N-CA-C	5.33	117.17	111.36
2	E	52	ILE	N-CA-C	-5.27	102.33	107.55
1	D	26	ALA	CA-C-O	-5.26	115.48	120.90
2	E	122	ALA	CA-C-N	5.25	127.26	120.44
2	E	122	ALA	C-N-CA	5.25	127.26	120.44
3	K	198	HIS	CA-C-N	5.21	127.52	120.38
3	K	198	HIS	C-N-CA	5.21	127.52	120.38
2	E	28	THR	CB-CA-C	5.17	118.01	110.26
2	E	164	SER	N-CA-C	5.15	118.69	111.74
3	L	170	ASP	CA-CB-CG	5.14	117.74	112.60
1	C	297	ASP	N-CA-C	5.11	116.53	111.07
1	D	243	LYS	CA-C-N	5.10	127.62	120.28
1	D	243	LYS	C-N-CA	5.10	127.62	120.28
2	F	206	VAL	N-CA-C	5.10	115.31	107.77
1	D	81	GLU	N-CA-C	-5.08	105.74	111.28
3	I	123	GLU	CB-CA-C	-5.08	102.68	110.81
2	E	186	LEU	N-CA-C	-5.08	101.58	109.24
1	A	89	ASP	CB-CA-C	5.06	119.61	111.66
1	B	139	TRP	CA-C-N	5.03	126.90	120.56
1	B	139	TRP	C-N-CA	5.03	126.90	120.56
2	F	192	VAL	CA-C-N	-5.03	114.58	119.76
2	F	192	VAL	C-N-CA	-5.03	114.58	119.76
2	G	55	PHE	CA-CB-CG	5.00	118.81	113.80
1	A	301	GLN	N-CA-C	5.00	117.12	111.02

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	88	LYS	Peptide

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	2901	0	2891	29	0
1	B	2901	0	2891	42	0
1	C	2901	0	2891	35	0
1	D	2901	0	2891	63	0
2	E	1614	0	1589	22	0
2	F	1614	0	1589	76	0
2	G	1614	0	1589	10	0
2	H	1614	0	1589	10	0
3	I	1628	0	1591	14	0
3	J	1628	0	1591	45	0
3	K	1628	0	1591	21	0
3	L	1628	0	1591	5	0
4	A	6	0	0	0	0
4	B	5	0	0	0	0
4	C	7	0	0	0	0
4	D	5	0	0	0	0
4	F	2	0	0	0	0
4	H	3	0	0	0	0
4	I	4	0	0	0	0
4	J	3	0	0	0	0
4	K	7	0	0	0	0
4	L	4	0	0	0	0
All	All	24618	0	24284	362	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 7.

All (362) 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)
2:F:203:ILE:HG23	2:F:217:LYS:O	1.44	1.18
2:E:200:GLN:NE2	2:E:201:THR:O	1.88	1.06
1:B:88:LYS:H	1:B:89:ASP:HA	1.28	0.98
1:D:83:MET:HE2	1:D:86:TRP:CD1	2.00	0.96
3:J:125:LEU:HB3	3:J:183:LYS:HE2	1.46	0.96
1:A:84:GLY:H	1:A:87:ASN:HD21	1.08	0.96
1:D:83:MET:HE2	1:D:86:TRP:HD1	1.32	0.95
2:F:203:ILE:CG2	2:F:217:LYS:O	2.15	0.95
2:F:179:GLN:OE1	2:F:185:SER:OG	1.85	0.93

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:87:ASN:HD22	1:A:87:ASN:H	1.12	0.92
1:C:90:GLU:HG2	1:C:173:GLY:HA2	1.50	0.91
2:F:153:TYR:CD1	2:F:184:TYR:O	2.25	0.89
1:D:82:LEU:HD11	1:D:371:LEU:HD21	1.52	0.89
3:J:150:VAL:HG13	3:J:191:VAL:O	1.73	0.88
1:C:83:MET:CE	1:C:93:THR:HG22	2.03	0.88
2:F:209:LYS:HG2	2:F:210:PRO:HD3	1.54	0.88
2:F:151:LYS:HG2	2:F:152:ASP:OD1	1.74	0.87
2:F:212:ASN:O	2:F:212:ASN:ND2	2.08	0.87
2:F:151:LYS:HA	2:F:185:SER:HB3	1.57	0.86
2:E:163:ASN:HB3	2:E:166:ALA:HB3	1.58	0.85
3:J:150:VAL:O	3:J:153:ALA:N	2.09	0.85
2:F:211:SER:OG	2:F:213:THR:OG1	1.94	0.85
1:B:90:GLU:HG2	1:B:173:GLY:HA2	1.58	0.84
1:B:88:LYS:HB3	1:B:89:ASP:O	1.76	0.83
2:F:153:TYR:CE2	2:F:158:VAL:HG23	2.13	0.83
1:D:83:MET:HB3	1:D:87:ASN:HB2	1.60	0.83
3:J:150:VAL:HG23	3:J:155:GLN:NE2	1.94	0.82
1:D:83:MET:HG2	1:D:86:TRP:HE1	1.45	0.82
2:F:67:ARG:HH12	2:F:90:ASP:CG	1.86	0.82
3:K:121:SER:O	3:K:125:LEU:HD12	1.79	0.82
1:C:87:ASN:C	1:C:89:ASP:HA	2.04	0.81
1:D:301:GLN:HG2	1:D:302:PHE:H	1.44	0.81
2:F:168:THR:O	2:F:171:VAL:HG23	1.80	0.81
3:J:125:LEU:HA	3:J:183:LYS:CE	2.11	0.80
1:D:83:MET:HG2	1:D:86:TRP:NE1	1.96	0.80
2:F:154:PHE:CD1	2:F:155:PRO:HB3	2.17	0.80
2:F:194:SER:O	2:F:197:LEU:HB2	1.82	0.79
2:E:194:SER:O	2:E:197:LEU:HB2	1.83	0.79
3:J:148:TRP:O	3:J:155:GLN:HG2	1.81	0.79
2:F:125:LYS:HD2	2:F:183:LEU:HD21	1.65	0.79
2:F:217:LYS:NZ	3:J:123:GLU:OE2	2.13	0.79
3:J:125:LEU:CB	3:J:183:LYS:HE2	2.12	0.77
2:E:200:GLN:HE21	2:E:201:THR:N	1.81	0.77
2:F:154:PHE:CE1	2:F:155:PRO:HB3	2.20	0.77
1:C:83:MET:HE1	1:C:93:THR:HG22	1.67	0.77
1:A:83:MET:HE2	1:A:93:THR:HG22	1.67	0.76
1:B:78:LEU:O	1:B:82:LEU:HG	1.87	0.75
3:K:149:LYS:HG2	3:K:154:LEU:HD23	1.67	0.74
1:C:83:MET:HE2	1:C:93:THR:HG22	1.69	0.74
1:D:301:GLN:HG2	1:D:302:PHE:N	1.99	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:J:150:VAL:HG23	3:J:155:GLN:HE22	1.52	0.73
1:A:84:GLY:H	1:A:87:ASN:ND2	1.85	0.72
2:F:182:GLY:O	2:F:183:LEU:HD12	1.88	0.72
3:I:121:SER:OG	3:I:124:GLN:HB3	1.90	0.72
2:F:207:ASN:N	2:F:207:ASN:HD22	1.88	0.71
2:F:209:LYS:CG	2:F:210:PRO:HD3	2.21	0.71
1:A:79:TYR:O	1:A:83:MET:HB2	1.90	0.71
2:H:30:SER:HA	2:H:53:PRO:HG2	1.72	0.71
2:G:64:PHE:HB2	2:G:68:VAL:HG12	1.71	0.71
3:J:125:LEU:O	3:J:183:LYS:HE3	1.91	0.71
1:D:83:MET:HG2	1:D:86:TRP:CD1	2.25	0.71
2:F:67:ARG:NH1	2:F:90:ASP:OD2	2.23	0.71
1:C:76:ARG:O	1:C:80:LYS:HG3	1.91	0.71
3:J:185:ASP:OD1	3:J:185:ASP:N	2.20	0.71
2:F:203:ILE:HG23	2:F:217:LYS:C	2.16	0.70
1:B:104:LYS:H	1:B:312:GLN:HE21	1.37	0.70
2:F:163:ASN:CB	2:F:166:ALA:HB3	2.22	0.69
3:J:125:LEU:HA	3:J:183:LYS:HE3	1.74	0.69
1:B:85:PRO:HB3	1:B:86:TRP:CD1	2.28	0.69
1:B:43:LEU:HB3	1:B:62:MET:HE1	1.75	0.69
1:B:100:GLN:OE1	1:B:101:ARG:N	2.25	0.68
2:E:163:ASN:CB	2:E:166:ALA:HB3	2.24	0.68
3:J:121:SER:OG	3:J:124:GLN:HG3	1.94	0.68
1:B:88:LYS:HB3	1:B:89:ASP:C	2.18	0.67
2:F:182:GLY:C	2:F:183:LEU:HD12	2.19	0.67
3:I:6:GLN:HE22	3:I:87:TYR:HA	1.60	0.67
1:A:84:GLY:N	1:A:87:ASN:HD21	1.89	0.67
1:B:86:TRP:HA	1:B:86:TRP:CE3	2.30	0.66
1:A:87:ASN:HD22	1:A:87:ASN:N	1.89	0.66
1:D:83:MET:CE	1:D:86:TRP:HD1	2.06	0.66
1:D:57:GLN:OE1	1:D:297:ASP:HB2	1.96	0.66
2:F:178:LEU:HB2	2:F:184:TYR:CE1	2.31	0.66
1:A:295:MET:HB3	1:A:298:MET:HE2	1.77	0.65
3:J:181:LEU:HD22	3:J:185:ASP:HB2	1.78	0.65
1:D:88:LYS:HA	1:D:90:GLU:N	2.12	0.65
1:C:81:GLU:O	1:C:86:TRP:HZ2	1.80	0.65
2:G:64:PHE:CB	2:G:68:VAL:HG12	2.27	0.65
3:J:125:LEU:CA	3:J:183:LYS:CE	2.75	0.65
3:K:149:LYS:HG2	3:K:154:LEU:CD2	2.27	0.65
1:D:88:LYS:HZ1	1:D:145:LYS:HD2	1.61	0.64
1:B:88:LYS:H	1:B:89:ASP:CA	2.06	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:163:ASN:HB3	2:F:166:ALA:HB3	1.80	0.64
1:B:185:HIS:HE1	1:B:187:ARG:HE	1.45	0.64
2:F:173:THR:HG23	2:F:188:SER:HB2	1.78	0.64
1:D:83:MET:HB3	1:D:87:ASN:CB	2.28	0.64
1:C:87:ASN:O	1:C:89:ASP:HA	1.98	0.64
3:J:150:VAL:CG1	3:J:191:VAL:O	2.45	0.63
2:G:64:PHE:HB2	2:G:68:VAL:CG1	2.27	0.63
1:B:104:LYS:HB2	1:B:312:GLN:HG2	1.80	0.63
2:F:153:TYR:HD1	2:F:184:TYR:O	1.77	0.62
3:J:125:LEU:CA	3:J:183:LYS:HE2	2.29	0.62
2:H:28:THR:HG23	2:H:32:TYR:HE2	1.64	0.62
1:D:33:VAL:HG12	1:D:360:PHE:HZ	1.64	0.62
1:B:85:PRO:HA	1:B:86:TRP:CD2	2.34	0.62
3:K:149:LYS:CG	3:K:154:LEU:HD23	2.28	0.62
2:F:196:SER:OG	2:F:202:TYR:OH	2.04	0.62
3:K:124:GLN:O	3:K:129:THR:O	2.18	0.61
1:B:110:MET:HE1	1:B:121:VAL:H	1.65	0.61
1:B:89:ASP:OD2	1:B:92:SER:OG	2.19	0.61
2:F:211:SER:CB	2:F:213:THR:HG1	2.13	0.61
1:D:91:ILE:HD12	1:D:92:SER:N	2.16	0.61
3:K:150:VAL:HG23	3:K:155:GLN:OE1	2.01	0.60
2:E:124:THR:HG23	2:E:155:PRO:HD3	1.83	0.60
3:K:182:SER:OG	3:K:185:ASP:HB2	2.02	0.60
1:A:83:MET:CE	1:A:93:THR:HG22	2.30	0.60
3:J:128:GLY:O	3:J:183:LYS:HG3	2.02	0.60
1:A:87:ASN:H	1:A:87:ASN:ND2	1.90	0.59
1:B:88:LYS:N	1:B:89:ASP:HA	2.07	0.59
1:D:88:LYS:HA	1:D:90:GLU:H	1.67	0.59
2:E:196:SER:HG	2:E:202:TYR:HH	1.50	0.59
2:E:120:SER:HA	2:E:121:SER:HB3	1.84	0.59
3:K:182:SER:O	3:K:185:ASP:N	2.34	0.59
2:E:127:PRO:HB3	2:E:153:TYR:HB3	1.84	0.59
2:F:154:PHE:CD1	2:F:155:PRO:CB	2.85	0.58
2:F:153:TYR:CE1	2:F:184:TYR:O	2.57	0.58
2:F:154:PHE:CD1	2:F:155:PRO:CA	2.86	0.58
1:D:61:ALA:HB3	1:D:295:MET:HE3	1.86	0.58
1:D:88:LYS:CE	1:D:145:LYS:HD2	2.34	0.58
2:F:213:THR:HG22	2:F:214:LYS:N	2.19	0.58
1:C:86:TRP:H	1:C:86:TRP:CD1	2.21	0.57
3:J:182:SER:OG	3:J:183:LYS:N	2.35	0.57
1:D:87:ASN:HD22	1:D:88:LYS:N	2.02	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:86:TRP:CD1	1:C:86:TRP:N	2.73	0.57
2:F:214:LYS:C	2:F:215:VAL:HG23	2.30	0.56
1:D:26:ALA:HB3	1:D:27:SER:HB2	1.87	0.56
3:I:125:LEU:HD11	3:I:211:ARG:HH22	1.70	0.56
1:D:43:LEU:HB3	1:D:62:MET:HE1	1.88	0.56
2:F:155:PRO:HD2	2:F:210:PRO:HG2	1.88	0.56
1:C:58:ILE:HA	1:C:295:MET:HE2	1.88	0.55
2:F:163:ASN:HB2	2:F:166:ALA:HB3	1.88	0.55
1:C:153:GLY:HA3	1:C:321:LEU:HD11	1.89	0.55
3:I:37:GLN:HB2	3:I:47:LEU:HD11	1.89	0.55
3:L:4:MET:HE3	3:L:90:GLN:HG2	1.87	0.55
2:F:147:GLY:HA3	2:F:189:VAL:HG12	1.88	0.54
2:F:209:LYS:CG	2:F:210:PRO:CD	2.86	0.54
3:K:183:LYS:O	3:K:186:TYR:HB3	2.07	0.54
2:F:174:PHE:CE1	3:J:164:THR:HG23	2.42	0.54
3:K:149:LYS:CG	3:K:154:LEU:CD2	2.84	0.54
2:F:154:PHE:HD1	2:F:155:PRO:CB	2.20	0.54
1:B:68:ASP:HB2	1:B:71:MET:HG3	1.89	0.54
1:A:251:THR:HA	1:A:254:LEU:HD12	1.89	0.54
2:F:153:TYR:CE2	2:F:158:VAL:CG2	2.87	0.54
3:L:37:GLN:HB2	3:L:47:LEU:HD11	1.89	0.54
1:C:83:MET:HE2	1:C:93:THR:CG2	2.37	0.53
2:F:209:LYS:HG3	2:F:210:PRO:N	2.22	0.53
1:A:280:LEU:HD11	1:A:376:VAL:HG22	1.90	0.53
3:J:120:PRO:HD3	3:J:132:VAL:HG22	1.90	0.53
3:K:125:LEU:O	3:K:183:LYS:NZ	2.41	0.53
2:F:154:PHE:HD1	2:F:155:PRO:CA	2.21	0.53
3:K:182:SER:O	3:K:186:TYR:N	2.37	0.53
1:D:76:ARG:HH12	1:D:116:LEU:HA	1.74	0.53
3:I:90:GLN:HE21	3:I:97:THR:H	1.57	0.53
1:D:286:LEU:HD11	1:D:322:GLN:HB2	1.91	0.52
1:C:48:LEU:HD13	1:C:117:PHE:HE2	1.75	0.52
1:C:280:LEU:HD11	1:C:376:VAL:HG22	1.91	0.52
1:D:88:LYS:NZ	1:D:145:LYS:HD2	2.24	0.52
1:A:8:VAL:HG22	1:A:71:MET:HE3	1.92	0.52
1:C:88:LYS:N	1:C:89:ASP:CA	2.72	0.52
1:D:65:LYS:HB2	1:D:68:ASP:HB2	1.91	0.52
1:D:87:ASN:ND2	1:D:91:ILE:CG2	2.72	0.52
2:F:153:TYR:CD2	2:F:158:VAL:HG23	2.44	0.52
1:A:170:TYR:HD1	1:A:325:LYS:HB3	1.75	0.52
1:B:280:LEU:HD11	1:B:376:VAL:HG22	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:K:37:GLN:HB2	3:K:47:LEU:HD11	1.91	0.52
1:A:46:LEU:HD13	1:A:165:LEU:HD21	1.91	0.51
2:F:162:TRP:CZ3	2:F:204:CYS:HB3	2.45	0.51
3:I:185:ASP:N	3:I:185:ASP:OD1	2.39	0.51
1:C:295:MET:HB3	1:C:298:MET:HE3	1.92	0.51
2:F:127:PRO:HB3	2:F:153:TYR:HB3	1.93	0.51
1:D:87:ASN:ND2	1:D:88:LYS:N	2.58	0.51
2:H:4:LEU:HD11	2:H:98:ARG:HB2	1.93	0.51
3:I:35:TRP:HB2	3:I:48:ILE:HB	1.93	0.51
3:K:155:GLN:HA	3:K:155:GLN:HE21	1.75	0.51
2:F:161:SER:HB3	2:F:205:ASN:HB2	1.93	0.51
2:E:120:SER:HA	2:E:121:SER:CB	2.40	0.51
1:D:58:ILE:HA	1:D:295:MET:HE2	1.93	0.51
1:A:82:LEU:C	1:A:86:TRP:CZ2	2.89	0.50
1:D:88:LYS:HZ1	1:D:145:LYS:CD	2.25	0.50
2:F:179:GLN:HG3	2:F:183:LEU:O	2.11	0.50
3:L:32:ASN:HD22	3:L:92:SER:HA	1.76	0.50
3:L:47:LEU:HA	3:L:58:VAL:HG21	1.93	0.50
2:F:174:PHE:O	2:F:186:LEU:HD13	2.12	0.50
1:D:302:PHE:O	1:D:303:GLN:HB2	2.12	0.50
3:K:35:TRP:HB2	3:K:48:ILE:HB	1.93	0.50
3:J:37:GLN:HB2	3:J:47:LEU:HD11	1.93	0.50
1:C:105:LEU:HD23	1:C:310:SER:HB2	1.94	0.49
3:I:125:LEU:HD11	3:I:211:ARG:NH2	2.26	0.49
3:J:125:LEU:CA	3:J:183:LYS:HE3	2.36	0.49
2:G:1:GLN:HE22	2:G:110:TYR:HE1	1.59	0.49
1:B:88:LYS:N	1:B:89:ASP:CA	2.72	0.49
1:D:51:GLY:HA2	1:D:55:GLN:HE21	1.77	0.49
1:D:87:ASN:ND2	1:D:91:ILE:HG23	2.27	0.49
1:D:183:SER:HB3	1:D:203:ALA:HB3	1.95	0.49
2:F:153:TYR:OH	2:F:186:LEU:HD23	2.12	0.49
1:A:178:PRO:HB3	1:A:331:SER:HB3	1.94	0.49
1:C:137:ASN:HD21	1:C:151:LEU:H	1.59	0.49
2:E:171:VAL:HG22	2:E:190:VAL:HG23	1.94	0.49
1:D:82:LEU:HG	1:D:82:LEU:O	2.13	0.49
3:J:125:LEU:HA	3:J:183:LYS:HE2	1.88	0.49
2:E:162:TRP:CE2	2:E:204:CYS:HB3	2.48	0.49
1:B:153:GLY:HA3	1:B:321:LEU:HD11	1.94	0.49
1:C:82:LEU:HD22	1:C:371:LEU:HD21	1.93	0.49
2:F:182:GLY:C	2:F:183:LEU:CD1	2.85	0.48
2:F:148:CYS:HB2	2:F:162:TRP:CH2	2.48	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:I:122:ASP:OD1	3:I:126:LYS:HE3	2.12	0.48
1:A:127:SER:HA	1:A:159:GLN:HG2	1.95	0.48
1:D:79:TYR:HA	1:D:82:LEU:O	2.13	0.48
1:D:137:ASN:HD21	1:D:151:LEU:H	1.60	0.48
2:E:174:PHE:CD1	3:I:164:THR:HG23	2.48	0.48
2:F:154:PHE:CD1	2:F:155:PRO:HA	2.48	0.48
1:B:212:GLU:HB2	3:J:30:TYR:HE1	1.78	0.48
2:F:174:PHE:CD1	3:J:164:THR:HG23	2.49	0.48
3:J:149:LYS:HG2	3:J:154:LEU:HD23	1.94	0.48
3:J:210:ASN:O	3:J:211:ARG:HB2	2.14	0.48
3:J:188:LYS:HD3	3:J:189:HIS:CE1	2.49	0.48
2:F:91:THR:HG23	2:F:118:THR:HA	1.96	0.48
1:C:207:LYS:HB3	1:C:268:ARG:HG2	1.96	0.48
1:D:88:LYS:CA	1:D:90:GLU:H	2.26	0.48
1:A:69:LYS:HA	1:A:70:GLY:HA2	1.71	0.48
1:D:88:LYS:HZ1	1:D:145:LYS:CE	2.26	0.48
2:F:193:PRO:HB2	2:F:196:SER:HB3	1.96	0.48
1:C:88:LYS:N	1:C:89:ASP:C	2.71	0.47
1:C:88:LYS:N	1:C:89:ASP:HA	2.30	0.47
3:J:47:LEU:HA	3:J:58:VAL:HG21	1.95	0.47
1:B:212:GLU:HB2	3:J:30:TYR:CE1	2.49	0.47
1:C:164:VAL:HG22	1:C:319:LEU:HB3	1.97	0.47
1:D:172:ASN:HD21	1:D:329:ASN:HD21	1.62	0.47
3:K:149:LYS:HE2	3:K:154:LEU:HD21	1.95	0.47
2:G:150:VAL:HB	2:G:186:LEU:HB3	1.97	0.47
1:D:361:VAL:HG13	1:D:370:VAL:HG13	1.96	0.47
2:E:62:GLN:HA	2:E:65:GLN:HG3	1.97	0.47
1:A:82:LEU:O	1:A:86:TRP:CZ2	2.68	0.47
2:G:91:THR:HG23	2:G:118:THR:HA	1.96	0.46
1:B:161:THR:HG23	1:B:318:ALA:HB3	1.97	0.46
1:D:190:HIS:N	1:D:355:ASP:O	2.45	0.46
2:F:167:LEU:HD13	2:F:190:VAL:HG21	1.98	0.46
1:D:88:LYS:HE3	1:D:145:LYS:HD2	1.98	0.46
3:L:193:ALA:HB2	3:L:208:SER:HB3	1.97	0.46
1:D:89:ASP:HB3	1:D:90:GLU:HG3	1.97	0.46
2:F:208:HIS:HD2	2:F:211:SER:OG	1.99	0.46
1:B:85:PRO:HA	1:B:86:TRP:CE3	2.51	0.45
2:F:162:TRP:CH2	2:F:204:CYS:HB3	2.50	0.45
1:C:95:ASP:HB3	1:C:167:ASN:HD22	1.80	0.45
1:D:91:ILE:HD12	1:D:92:SER:H	1.82	0.45
3:K:181:LEU:HD22	3:K:185:ASP:HB3	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:295:MET:HB3	1:B:298:MET:HE3	1.98	0.45
1:C:103:LEU:HD12	1:C:103:LEU:HA	1.78	0.45
1:D:79:TYR:O	1:D:79:TYR:CD1	2.69	0.45
2:E:15:GLY:CA	2:E:16:SER:HB2	2.46	0.45
1:A:100:GLN:HB3	1:A:103:LEU:HB2	1.97	0.45
1:C:201:MET:HA	1:C:276:PRO:HA	1.99	0.45
1:B:104:LYS:H	1:B:312:GLN:NE2	2.11	0.45
1:A:30:ARG:HH21	1:B:185:HIS:HB2	1.81	0.45
3:J:124:GLN:HE22	3:J:131:SER:HB2	1.81	0.45
2:E:152:ASP:HB3	2:E:183:LEU:HD13	1.98	0.45
1:A:27:SER:HB3	1:A:30:ARG:HB2	1.99	0.44
1:A:82:LEU:O	1:A:86:TRP:CH2	2.70	0.44
1:B:251:THR:HA	1:B:254:LEU:HD12	1.99	0.44
1:C:235:MET:HE1	1:C:328:VAL:HG21	1.99	0.44
1:D:78:LEU:O	1:D:82:LEU:HD23	2.17	0.44
2:H:51:ILE:HD12	2:H:58:ALA:HB2	1.99	0.44
3:J:150:VAL:C	3:J:152:ASN:N	2.73	0.44
1:B:84:GLY:O	1:B:86:TRP:CE3	2.70	0.44
1:C:28:LYS:HA	1:C:375:GLN:HE22	1.82	0.44
1:D:347:MET:HE3	2:H:59:ASN:HB2	1.99	0.44
2:F:98:ARG:NH1	2:F:99:GLU:O	2.50	0.44
3:K:155:GLN:HE21	3:K:155:GLN:CA	2.30	0.44
1:B:361:VAL:HG13	1:B:370:VAL:HG13	2.00	0.44
2:G:150:VAL:HG11	2:G:158:VAL:HG11	1.99	0.44
1:C:64:PHE:HB2	1:C:71:MET:HE2	2.00	0.44
1:C:167:ASN:O	1:C:322:GLN:HA	2.18	0.44
1:D:58:ILE:HA	1:D:295:MET:CE	2.48	0.44
1:A:140:VAL:HG21	1:A:151:LEU:HD11	1.99	0.43
1:D:87:ASN:ND2	1:D:91:ILE:HG22	2.33	0.43
2:E:30:SER:HA	2:E:53:PRO:HB2	2.00	0.43
2:F:48:MET:HE1	2:F:94:TYR:HD2	1.84	0.43
2:F:177:VAL:HG11	3:J:160:GLN:HB3	2.00	0.43
2:F:97:ALA:HB1	2:F:108:PHE:HB3	2.01	0.43
1:B:91:ILE:HB	1:B:171:PHE:HD1	1.82	0.43
2:H:201:THR:HG23	2:H:218:LYS:HE2	2.00	0.43
2:E:150:VAL:HB	2:E:186:LEU:HB3	2.01	0.43
1:C:96:ALA:HB1	1:C:122:LYS:HD3	2.01	0.43
2:F:209:LYS:O	2:F:212:ASN:N	2.52	0.43
3:I:47:LEU:HA	3:I:58:VAL:HG21	2.00	0.43
3:J:186:TYR:CD1	3:J:192:TYR:CZ	3.07	0.43
1:A:90:GLU:OE2	1:A:229:HIS:NE2	2.43	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:83:MET:O	1:B:83:MET:HG3	2.19	0.43
1:B:85:PRO:HA	1:B:86:TRP:HA	1.66	0.43
1:D:295:MET:HB3	1:D:298:MET:HE3	2.01	0.43
1:A:290:LEU:HD13	1:A:298:MET:HE1	2.01	0.43
1:D:251:THR:HA	1:D:254:LEU:HD12	1.99	0.43
2:E:91:THR:HG23	2:E:118:THR:HA	2.00	0.42
3:J:32:ASN:HD22	3:J:92:SER:HA	1.84	0.42
2:F:215:VAL:HG12	2:F:216:ASP:N	2.32	0.42
1:B:85:PRO:HA	1:B:86:TRP:CG	2.54	0.42
2:F:197:LEU:HD23	2:F:197:LEU:HA	1.90	0.42
3:J:125:LEU:C	3:J:183:LYS:HE3	2.43	0.42
2:H:48:MET:HE1	2:H:94:TYR:HD2	1.84	0.42
1:B:286:LEU:HD11	1:B:322:GLN:HB2	2.01	0.42
3:J:128:GLY:O	3:J:183:LYS:CG	2.65	0.42
1:D:188:LEU:HD22	1:D:196:THR:HB	2.01	0.42
2:F:60:TYR:HB2	2:F:65:GLN:HG3	2.01	0.42
2:F:213:THR:CG2	2:F:214:LYS:N	2.82	0.42
1:B:83:MET:N	1:B:84:GLY:HA3	2.35	0.42
1:C:58:ILE:O	1:C:62:MET:HB2	2.20	0.42
2:E:51:ILE:HD12	2:E:58:ALA:HB2	2.02	0.42
1:C:362:VAL:HB	1:C:372:PHE:HB2	2.01	0.41
2:F:209:LYS:HG3	2:F:210:PRO:CD	2.49	0.41
2:F:214:LYS:C	2:F:215:VAL:CG2	2.93	0.41
3:J:147:GLN:HE21	3:J:154:LEU:HD22	1.85	0.41
1:B:85:PRO:HB3	1:B:86:TRP:NE1	2.33	0.41
1:B:169:LEU:HB2	1:B:324:VAL:HG22	2.02	0.41
1:D:110:MET:HE2	1:D:110:MET:HA	2.01	0.41
1:D:236:PHE:O	1:D:360:PHE:HA	2.21	0.41
2:H:154:PHE:HB2	2:H:183:LEU:HD23	2.01	0.41
3:J:190:LYS:HE2	3:J:210:ASN:CG	2.45	0.41
1:B:46:LEU:HD13	1:B:165:LEU:HD11	2.01	0.41
1:C:347:MET:HE1	2:G:51:ILE:HA	2.02	0.41
1:D:33:VAL:HG11	1:D:326:ILE:HG22	2.01	0.41
3:J:188:LYS:HD3	3:J:189:HIS:HE1	1.85	0.41
2:H:176:ALA:HA	2:H:186:LEU:HB3	2.01	0.41
2:F:209:LYS:CG	2:F:210:PRO:N	2.83	0.41
3:I:6:GLN:HE21	3:I:99:GLY:HA3	1.85	0.41
2:F:154:PHE:HA	2:F:155:PRO:HA	1.79	0.41
3:I:125:LEU:O	3:I:183:LYS:HD2	2.21	0.41
1:D:82:LEU:HA	1:D:83:MET:HA	1.86	0.41
1:D:179:PHE:CD2	1:D:202:MET:HB3	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:J:183:LYS:HB3	3:J:183:LYS:NZ	2.35	0.41
2:F:163:ASN:O	2:F:164:SER:HB2	2.20	0.41
2:G:64:PHE:HB2	2:G:65:GLN:H	1.69	0.41
2:E:197:LEU:HD23	2:E:197:LEU:HA	1.77	0.41
2:F:203:ILE:HG22	2:F:204:CYS:N	2.34	0.41
2:G:30:SER:HA	2:G:53:PRO:HB2	2.03	0.41
3:K:183:LYS:HA	3:K:186:TYR:HB3	2.02	0.41
1:A:82:LEU:C	1:A:86:TRP:HZ2	2.29	0.41
1:B:44:ALA:HB1	1:B:66:ILE:HD13	2.02	0.41
1:D:89:ASP:CB	1:D:174:GLN:NE2	2.84	0.41
1:D:87:ASN:ND2	1:D:87:ASN:C	2.79	0.40
2:H:99:GLU:HG2	2:H:107:HIS:O	2.21	0.40
1:B:104:LYS:O	1:B:310:SER:OG	2.26	0.40
1:D:88:LYS:C	1:D:90:GLU:H	2.29	0.40
2:F:162:TRP:CD1	2:F:171:VAL:CG1	3.05	0.40
3:J:148:TRP:CE2	3:J:179:LEU:HB2	2.57	0.40
2:E:167:LEU:HD13	2:E:190:VAL:HG21	2.04	0.40
3:J:150:VAL:O	3:J:153:ALA:HB3	2.21	0.40
3:K:181:LEU:HD23	3:K:181:LEU:HA	1.76	0.40
1:A:44:ALA:HB1	1:A:66:ILE:HD13	2.02	0.40
1:D:100:GLN:HE22	1:D:125:ASP:HA	1.87	0.40
3:I:182:SER:OG	3:I:185:ASP:OD1	2.38	0.40
1:D:302:PHE:O	1:D:303:GLN:CB	2.69	0.40
3:J:145:LYS:HB3	3:J:197:THR:HB	2.03	0.40
3:K:155:GLN:H	3:K:155:GLN:HG2	1.71	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	359/379 (95%)	345 (96%)	14 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	359/379 (95%)	345 (96%)	14 (4%)	0	100	100
1	C	359/379 (95%)	347 (97%)	11 (3%)	1 (0%)	36	65
1	D	359/379 (95%)	345 (96%)	10 (3%)	4 (1%)	11	36
2	E	210/223 (94%)	200 (95%)	8 (4%)	2 (1%)	12	39
2	F	210/223 (94%)	199 (95%)	11 (5%)	0	100	100
2	G	210/223 (94%)	197 (94%)	11 (5%)	2 (1%)	12	39
2	H	210/223 (94%)	200 (95%)	9 (4%)	1 (0%)	24	54
3	I	209/214 (98%)	198 (95%)	10 (5%)	1 (0%)	24	54
3	J	209/214 (98%)	193 (92%)	15 (7%)	1 (0%)	24	54
3	K	209/214 (98%)	193 (92%)	13 (6%)	3 (1%)	9	30
3	L	209/214 (98%)	199 (95%)	9 (4%)	1 (0%)	24	54
All	All	3112/3264 (95%)	2961 (95%)	135 (4%)	16 (0%)	24	54

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	E	124	THR
1	D	88	LYS
1	D	303	GLN
2	G	28	THR
3	J	138	ASN
3	K	138	ASN
2	G	152	ASP
3	I	138	ASN
3	K	151	ASP
3	L	138	ASN
2	E	16	SER
1	C	102	ASP
1	D	27	SER
3	K	143	GLU
1	D	129	VAL
2	H	155	PRO

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	319/333 (96%)	294 (92%)	25 (8%)	11	35
1	B	319/333 (96%)	292 (92%)	27 (8%)	10	31
1	C	319/333 (96%)	292 (92%)	27 (8%)	10	31
1	D	319/333 (96%)	288 (90%)	31 (10%)	8	25
2	E	179/186 (96%)	166 (93%)	13 (7%)	13	38
2	F	179/186 (96%)	161 (90%)	18 (10%)	7	24
2	G	179/186 (96%)	164 (92%)	15 (8%)	10	31
2	H	179/186 (96%)	166 (93%)	13 (7%)	13	38
3	I	185/187 (99%)	168 (91%)	17 (9%)	8	27
3	J	185/187 (99%)	173 (94%)	12 (6%)	15	44
3	K	185/187 (99%)	175 (95%)	10 (5%)	20	51
3	L	185/187 (99%)	176 (95%)	9 (5%)	22	54
All	All	2732/2824 (97%)	2515 (92%)	217 (8%)	11	34

All (217) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	43	LEU
1	A	47	GLN
1	A	48	LEU
1	A	67	ASP
1	A	68	ASP
1	A	71	MET
1	A	75	LEU
1	A	83	MET
1	A	87	ASN
1	A	91	ILE
1	A	92	SER
1	A	107	GLN
1	A	110	MET
1	A	122	LYS
1	A	130	GLU
1	A	158	ASP
1	A	159	GLN
1	A	165	LEU
1	A	182	SER

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Mol	Chain	Res	Type
1	A	209	ASN
1	A	216	PRO
1	A	217	ASP
1	A	250	LEU
1	A	325	LYS
1	A	329	ASN
1	B	14	ASP
1	B	43	LEU
1	B	47	GLN
1	B	48	LEU
1	B	53	GLU
1	B	78	LEU
1	B	80	LYS
1	B	86	TRP
1	B	89	ASP
1	B	90	GLU
1	B	91	ILE
1	B	95	ASP
1	B	104	LYS
1	B	105	LEU
1	B	107	GLN
1	B	110	MET
1	B	123	GLN
1	B	131	ARG
1	B	141	LYS
1	B	148	ILE
1	B	181	ASP
1	B	217	ASP
1	B	224	LEU
1	B	248	SER
1	B	250	LEU
1	B	293	LEU
1	B	346	ARG
1	C	21	GLN
1	C	22	GLN
1	C	25	GLN
1	C	48	LEU
1	C	62	MET
1	C	68	ASP
1	C	69	LYS
1	C	86	TRP
1	C	87	ASN

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Mol	Chain	Res	Type
1	C	91	ILE
1	C	95	ASP
1	C	102	ASP
1	C	130	GLU
1	C	137	ASN
1	C	145	LYS
1	C	165	LEU
1	C	186	ARG
1	C	217	ASP
1	C	224	LEU
1	C	232	THR
1	C	234	SER
1	C	250	LEU
1	C	257	GLN
1	C	260	SER
1	C	279	SER
1	C	346	ARG
1	C	347	MET
1	D	14	ASP
1	D	42	VAL
1	D	43	LEU
1	D	48	LEU
1	D	55	GLN
1	D	62	MET
1	D	69	LYS
1	D	75	LEU
1	D	78	LEU
1	D	80	LYS
1	D	81	GLU
1	D	82	LEU
1	D	83	MET
1	D	87	ASN
1	D	88	LYS
1	D	91	ILE
1	D	99	VAL
1	D	102	ASP
1	D	110	MET
1	D	137	ASN
1	D	140	VAL
1	D	158	ASP
1	D	202	MET
1	D	217	ASP

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Mol	Chain	Res	Type
1	D	224	LEU
1	D	250	LEU
1	D	263	LYS
1	D	268	ARG
1	D	298	MET
1	D	308	SER
1	D	371	LEU
2	E	19	LYS
2	E	99	GLU
2	E	105	GLU
2	E	123	SER
2	E	146	LEU
2	E	168	THR
2	E	194	SER
2	E	196	SER
2	E	200	GLN
2	E	213	THR
2	E	217	LYS
2	E	220	GLU
2	E	222	LYS
2	F	10	GLU
2	F	34	ILE
2	F	43	GLN
2	F	105	GLU
2	F	120	SER
2	F	171	VAL
2	F	178	LEU
2	F	181	SER
2	F	185	SER
2	F	186	LEU
2	F	190	VAL
2	F	194	SER
2	F	195	SER
2	F	199	THR
2	F	201	THR
2	F	207	ASN
2	F	214	LYS
2	F	220	GLU
2	G	1	GLN
2	G	10	GLU
2	G	43	GLN
2	G	64	PHE

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Mol	Chain	Res	Type
2	G	67	ARG
2	G	68	VAL
2	G	81	MET
2	G	99	GLU
2	G	105	GLU
2	G	148	CYS
2	G	156	GLN
2	G	158	VAL
2	G	167	LEU
2	G	217	LYS
2	G	218	LYS
2	H	10	GLU
2	H	81	MET
2	H	98	ARG
2	H	99	GLU
2	H	104	LEU
2	H	105	GLU
2	H	148	CYS
2	H	156	GLN
2	H	167	LEU
2	H	199	THR
2	H	204	CYS
2	H	207	ASN
2	H	218	LYS
3	I	1	ASP
3	I	7	SER
3	I	29	ILE
3	I	33	LEU
3	I	47	LEU
3	I	48	ILE
3	I	67	SER
3	I	124	GLN
3	I	125	LEU
3	I	126	LYS
3	I	127	SER
3	I	154	LEU
3	I	163	VAL
3	I	181	LEU
3	I	183	LYS
3	I	185	ASP
3	I	211	ARG
3	J	27	GLU

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Mol	Chain	Res	Type
3	J	48	ILE
3	J	67	SER
3	J	69	THR
3	J	70	GLU
3	J	122	ASP
3	J	142	ARG
3	J	150	VAL
3	J	154	LEU
3	J	183	LYS
3	J	185	ASP
3	J	206	THR
3	K	3	GLN
3	K	7	SER
3	K	18	ARG
3	K	54	LEU
3	K	79	GLN
3	K	105	GLU
3	K	155	GLN
3	K	183	LYS
3	K	187	GLU
3	K	211	ARG
3	L	1	ASP
3	L	26	SER
3	L	27	GLU
3	L	29	ILE
3	L	42	LYS
3	L	54	LEU
3	L	69	THR
3	L	106	ILE
3	L	176	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (68) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	22	GLN
1	A	87	ASN
1	A	100	GLN
1	A	172	ASN
1	A	261	HIS
1	A	265	ASN
1	A	303	GLN
1	A	316	HIS

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Mol	Chain	Res	Type
1	A	365	ASN
1	A	375	GLN
1	B	21	GLN
1	B	22	GLN
1	B	55	GLN
1	B	56	GLN
1	B	150	HIS
1	B	167	ASN
1	B	185	HIS
1	B	261	HIS
1	B	292	ASN
1	B	301	GLN
1	B	312	GLN
1	C	25	GLN
1	C	137	ASN
1	C	159	GLN
1	C	252	ASN
1	C	257	GLN
1	C	265	ASN
1	C	301	GLN
1	D	22	GLN
1	D	55	GLN
1	D	59	GLN
1	D	87	ASN
1	D	137	ASN
1	D	150	HIS
1	D	159	GLN
1	D	167	ASN
1	D	174	GLN
1	D	261	HIS
1	D	292	ASN
1	D	329	ASN
2	E	62	GLN
2	E	65	GLN
2	E	172	HIS
2	E	179	GLN
2	E	200	GLN
2	F	43	GLN
2	F	102	GLN
2	F	172	HIS
2	F	207	ASN
2	F	208	HIS

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Mol	Chain	Res	Type
2	G	1	GLN
2	G	43	GLN
2	G	200	GLN
2	G	208	HIS
2	H	172	HIS
2	H	200	GLN
3	I	6	GLN
3	I	31	HIS
3	I	79	GLN
3	I	90	GLN
3	I	137	ASN
3	J	147	GLN
3	J	155	GLN
3	K	155	GLN
3	L	32	ASN
3	L	137	ASN
3	L	138	ASN
3	L	189	HIS

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](#)

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 ⓘ

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	363/379 (95%)	-0.22	1 (0%) 90 87	34, 49, 72, 83	0
1	B	363/379 (95%)	-0.09	1 (0%) 90 87	37, 53, 85, 99	0
1	C	363/379 (95%)	-0.11	3 (0%) 82 77	44, 62, 88, 98	0
1	D	363/379 (95%)	0.15	2 (0%) 85 81	54, 72, 113, 127	0
2	E	214/223 (95%)	0.16	2 (0%) 81 75	52, 77, 109, 122	0
2	F	214/223 (95%)	0.18	3 (1%) 73 65	49, 75, 104, 121	0
2	G	214/223 (95%)	-0.16	1 (0%) 87 83	44, 61, 83, 97	0
2	H	214/223 (95%)	-0.05	0 100 100	44, 63, 80, 90	0
3	I	211/214 (98%)	-0.13	1 (0%) 87 83	37, 56, 82, 92	0
3	J	211/214 (98%)	-0.06	0 100 100	41, 57, 116, 130	0
3	K	211/214 (98%)	-0.18	0 100 100	32, 47, 97, 110	0
3	L	211/214 (98%)	-0.31	0 100 100	35, 44, 74, 86	0
All	All	3152/3264 (96%)	-0.07	14 (0%) 88 85	32, 61, 99, 130	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	E	203	ILE	3.2
1	C	85	PRO	2.7
3	I	125	LEU	2.7
2	F	199	THR	2.6
2	F	127	PRO	2.6
1	D	150	HIS	2.5
1	C	86	TRP	2.4
1	B	89	ASP	2.2
2	G	65	GLN	2.1
1	A	86	TRP	2.1
1	C	345	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
2	E	199	THR	2.1
2	F	212	ASN	2.0
1	D	119	SER	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.