



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 12:01 PM GMT

PDB ID : 3A9Q
Title : Crystal Structure Analysis of E173A variant of the soybean ferritin SFER4
Authors : Masuda, T.; Goto, F.; Yoshihara, T.; Mikami, B.
Deposited on : 2009-11-05
Resolution : 1.90 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

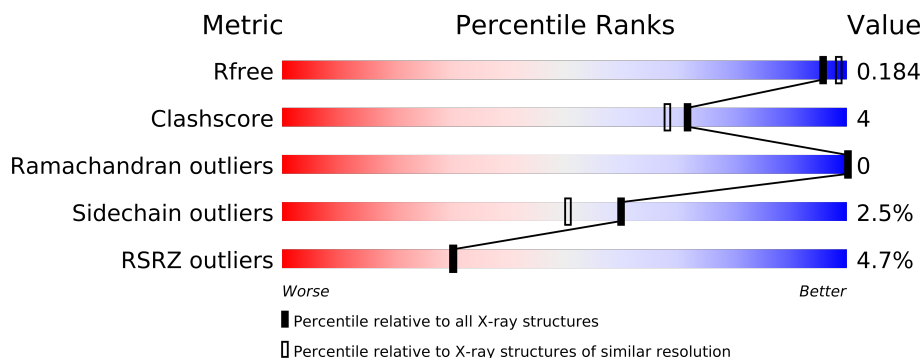
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 1.90 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.












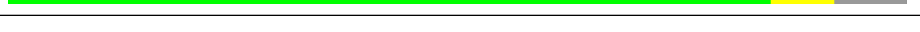
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	66092	3684 (1.90-1.90)
Clashscore	79885	4465 (1.90-1.90)
Ramachandran outliers	78287	4413 (1.90-1.90)
Sidechain outliers	78261	4414 (1.90-1.90)
RSRZ outliers	66119	3686 (1.90-1.90)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	212	
1	B	212	
1	C	212	
1	D	212	
1	E	212	
1	F	212	
1	G	212	
1	H	212	
1	I	212	
1	J	212	
1	K	212	
1	L	212	
1	M	212	
1	N	212	

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Mol	Chain	Length	Quality of chain
1	O	212	
1	P	212	
1	Q	212	
1	R	212	
1	S	212	
1	T	212	
1	U	212	
1	V	212	
1	W	212	
1	X	212	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
2	CA	B	216	-	X
2	CA	B	218[B]	-	X
2	CA	C	216[B]	-	X
2	CA	C	217	-	X
2	CA	D	214	-	X
2	CA	D	216[B]	-	X
2	CA	E	217[B]	-	X
2	CA	F	216	-	X
2	CA	F	218[B]	-	X
2	CA	G	216	-	X
2	CA	G	218[B]	-	X
2	CA	H	216[B]	-	X
2	CA	I	213	-	X
2	CA	I	216[A]	-	X
2	CA	I	216[B]	-	X
2	CA	J	215[A]	-	X
2	CA	J	215[B]	-	X
2	CA	J	216	-	X
2	CA	K	216[A]	-	X
2	CA	K	218	-	X
2	CA	L	218	-	X
2	CA	M	215	-	X
2	CA	M	217[B]	-	X
2	CA	N	218[B]	-	X
2	CA	O	215	-	X
2	CA	O	216	-	X
2	CA	P	217[A]	-	X

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Mol	Type	Chain	Res	Geometry	Electron density
2	CA	P	217[B]	-	X
2	CA	Q	216[B]	-	X
2	CA	R	217[B]	-	X
2	CA	R	218	-	X
2	CA	S	216[B]	-	X
2	CA	S	217	-	X
2	CA	T	215[A]	-	X
2	CA	T	217	-	X
2	CA	U	215[B]	-	X
2	CA	U	216	-	X
2	CA	V	216[A]	-	X
2	CA	W	215	-	X
2	CA	W	216	-	X
2	CA	X	213	-	X
2	CA	X	217[A]	-	X
3	ACY	G	219	-	X
3	ACY	V	221	-	X
3	ACY	X	241	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 42812 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Ferritin-4, chloroplastic.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	194	Total	C	N	O	S	0	9	0
			1621	1033	270	312	6			
1	B	194	Total	C	N	O	S	0	11	0
			1633	1039	271	316	7			
1	C	194	Total	C	N	O	S	0	16	0
			1669	1059	277	328	5			
1	D	194	Total	C	N	O	S	0	15	0
			1661	1056	273	326	6			
1	E	178	Total	C	N	O	S	0	11	0
			1492	945	252	289	6			
1	F	194	Total	C	N	O	S	0	12	0
			1636	1043	271	316	6			
1	G	193	Total	C	N	O	S	0	10	0
			1617	1027	269	315	6			
1	H	193	Total	C	N	O	S	0	12	0
			1631	1037	272	317	5			
1	I	193	Total	C	N	O	S	0	20	0
			1677	1069	274	328	6			
1	J	194	Total	C	N	O	S	0	11	0
			1637	1041	275	316	5			
1	K	193	Total	C	N	O	S	0	6	0
			1599	1013	270	311	5			
1	L	193	Total	C	N	O	S	0	9	0
			1618	1029	270	314	5			
1	M	193	Total	C	N	O	S	0	15	0
			1651	1050	275	320	6			
1	N	193	Total	C	N	O	S	0	13	0
			1631	1036	270	319	6			
1	O	193	Total	C	N	O	S	0	16	0
			1657	1051	274	327	5			
1	P	193	Total	C	N	O	S	0	12	0
			1631	1035	270	321	5			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	Q	193	Total	C	N	O	S	0	13	0
			1630	1040	271	312	7			
1	R	196	Total	C	N	O	S	0	11	0
			1643	1043	275	319	6			
1	S	193	Total	C	N	O	S	0	12	0
			1622	1033	269	315	5			
1	T	194	Total	C	N	O	S	0	13	0
			1650	1049	274	321	6			
1	U	193	Total	C	N	O	S	0	15	0
			1641	1045	271	318	7			
1	V	193	Total	C	N	O	S	0	13	0
			1633	1037	270	320	6			
1	W	194	Total	C	N	O	S	0	13	0
			1649	1048	275	320	6			
1	X	194	Total	C	N	O	S	0	15	0
			1652	1050	272	322	8			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	173	ALA	GLU	ENGINEERED	UNP Q948P5
B	173	ALA	GLU	ENGINEERED	UNP Q948P5
C	173	ALA	GLU	ENGINEERED	UNP Q948P5
D	173	ALA	GLU	ENGINEERED	UNP Q948P5
E	173	ALA	GLU	ENGINEERED	UNP Q948P5
F	173	ALA	GLU	ENGINEERED	UNP Q948P5
G	173	ALA	GLU	ENGINEERED	UNP Q948P5
H	173	ALA	GLU	ENGINEERED	UNP Q948P5
I	173	ALA	GLU	ENGINEERED	UNP Q948P5
J	173	ALA	GLU	ENGINEERED	UNP Q948P5
K	173	ALA	GLU	ENGINEERED	UNP Q948P5
L	173	ALA	GLU	ENGINEERED	UNP Q948P5
M	173	ALA	GLU	ENGINEERED	UNP Q948P5
N	173	ALA	GLU	ENGINEERED	UNP Q948P5
O	173	ALA	GLU	ENGINEERED	UNP Q948P5
P	173	ALA	GLU	ENGINEERED	UNP Q948P5
Q	173	ALA	GLU	ENGINEERED	UNP Q948P5
R	173	ALA	GLU	ENGINEERED	UNP Q948P5
S	173	ALA	GLU	ENGINEERED	UNP Q948P5
T	173	ALA	GLU	ENGINEERED	UNP Q948P5
U	173	ALA	GLU	ENGINEERED	UNP Q948P5
V	173	ALA	GLU	ENGINEERED	UNP Q948P5
W	173	ALA	GLU	ENGINEERED	UNP Q948P5

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Chain	Residue	Modelled	Actual	Comment	Reference
X	173	ALA	GLU	ENGINEERED	UNP Q948P5

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

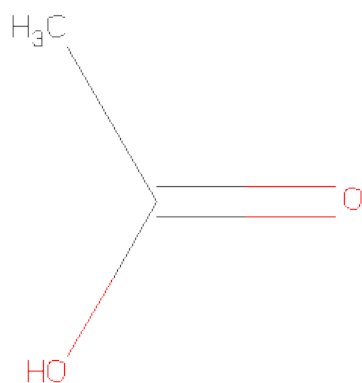
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	P	5	Total Ca 6 6	0	1
2	K	6	Total Ca 6 6	0	2
2	B	6	Total Ca 6 6	0	2
2	W	4	Total Ca 4 4	0	0
2	N	6	Total Ca 6 6	0	2
2	X	7	Total Ca 7 7	0	2
2	S	5	Total Ca 5 5	0	2
2	J	5	Total Ca 6 6	0	1
2	E	5	Total Ca 5 5	0	2
2	V	5	Total Ca 5 5	0	2
2	A	4	Total Ca 5 5	0	1
2	R	6	Total Ca 6 6	0	2
2	M	5	Total Ca 5 5	0	2
2	D	4	Total Ca 4 4	0	2
2	I	5	Total Ca 6 6	0	1
2	U	4	Total Ca 4 4	0	2
2	L	6	Total Ca 7 7	0	1
2	G	6	Total Ca 6 6	0	2
2	Q	4	Total Ca 4 4	0	2

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	H	5	Total	Ca	0	2
			5	5		
2	C	5	Total	Ca	0	1
			6	6		
2	T	5	Total	Ca	0	2
			5	5		
2	O	5	Total	Ca	0	0
			5	5		
2	F	6	Total	Ca	0	2
			6	6		

- Molecule 3 is ACETIC ACID (three-letter code: ACY) (formula: C₂H₄O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	2	2		
3	E	1	Total	C	O	0	0
			4	2	2		
3	F	1	Total	C	O	0	0
			4	2	2		
3	G	1	Total	C	O	0	0
			4	2	2		
3	H	1	Total	C	O	0	0
			4	2	2		
3	I	1	Total	C	O	0	0
			4	2	2		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	J	1	Total C O 4 2 2	0	0
3	L	1	Total C O 4 2 2	0	0
3	M	1	Total C O 4 2 2	0	0
3	O	1	Total C O 4 2 2	0	0
3	P	1	Total C O 4 2 2	0	0
3	Q	1	Total C O 4 2 2	0	0
3	R	1	Total C O 4 2 2	0	0
3	S	1	Total C O 4 2 2	0	0
3	T	1	Total C O 4 2 2	0	0
3	V	1	Total C O 4 2 2	0	0
3	W	1	Total C O 4 2 2	0	0
3	X	1	Total C O 4 2 2	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	158	Total O 158 158	0	0
4	B	149	Total O 149 149	0	0
4	C	143	Total O 143 143	0	0
4	D	130	Total O 130 130	0	0
4	E	150	Total O 150 150	0	0
4	F	146	Total O 146 146	0	0
4	G	155	Total O 155 155	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	H	143	Total 145	O 145	0	2
4	I	133	Total 133	O 133	0	0
4	J	142	Total 142	O 142	0	0
4	K	129	Total 129	O 129	0	0
4	L	136	Total 136	O 136	0	0
4	M	134	Total 134	O 134	0	0
4	N	144	Total 145	O 145	0	1
4	O	162	Total 162	O 162	0	0
4	P	136	Total 136	O 136	0	0
4	Q	147	Total 148	O 148	0	1
4	R	150	Total 150	O 150	0	0
4	S	137	Total 137	O 137	0	0
4	T	135	Total 135	O 135	0	0
4	U	135	Total 135	O 135	0	0
4	V	146	Total 146	O 146	0	0
4	W	137	Total 137	O 137	0	0
4	X	148	Total 148	O 148	0	0

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

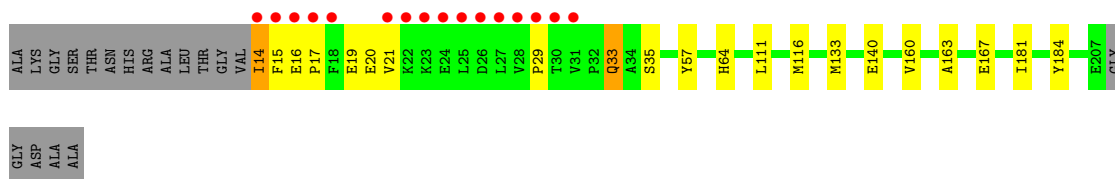
- Molecule 1: Ferritin-4, chloroplastic

Chain A: 



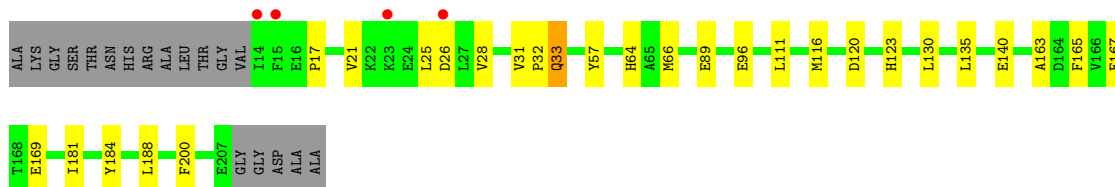
- Molecule 1: Ferritin-4, chloroplastic

Chain B: 



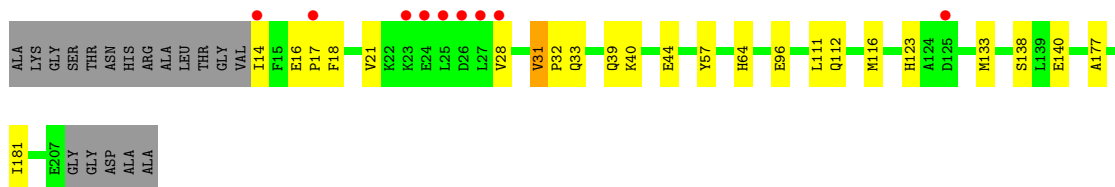
- Molecule 1: Ferritin-4, chloroplastic

Chain C: 



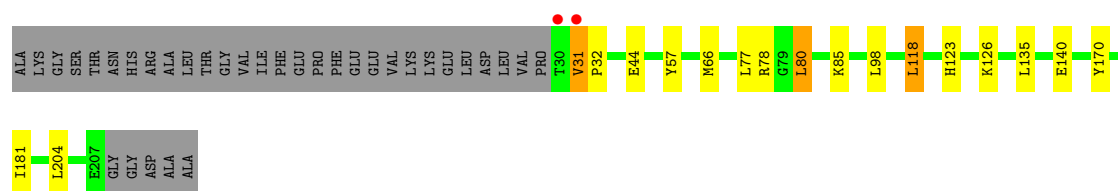
- Molecule 1: Ferritin-4, chloroplastic

Chain D: 



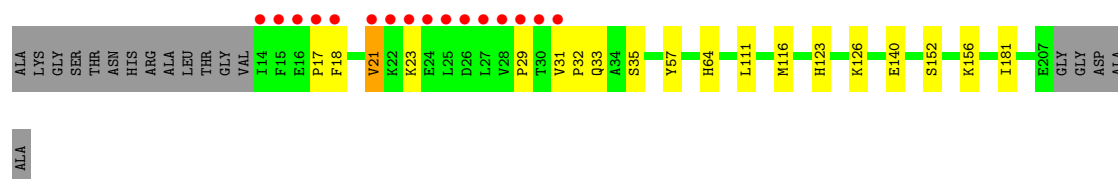
- Molecule 1: Ferritin-4, chloroplastic

Chain E: 



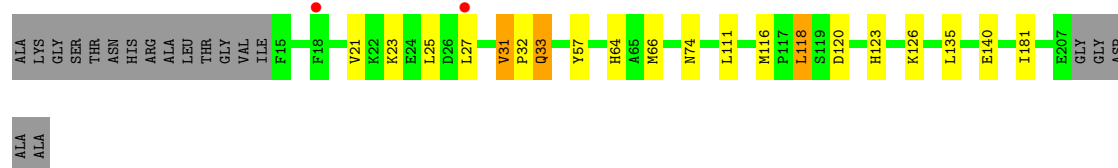
- Molecule 1: Ferritin-4, chloroplastic

Chain F:



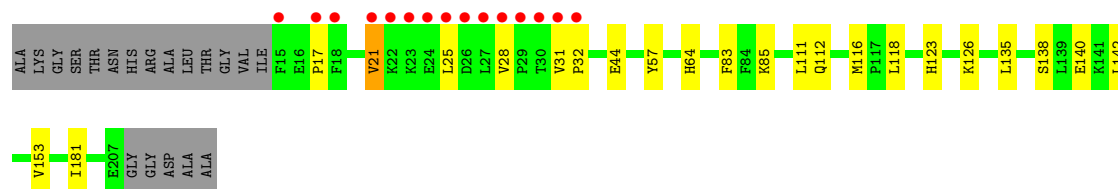
- Molecule 1: Ferritin-4, chloroplastic

Chain G:



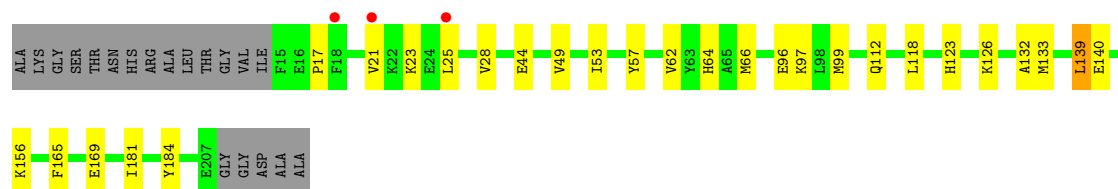
- Molecule 1: Ferritin-4, chloroplastic

Chain H:



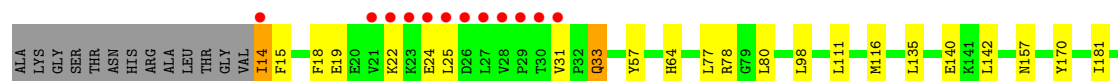
- Molecule 1: Ferritin-4, chloroplastic

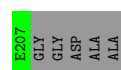
Chain I:



- Molecule 1: Ferritin-4, chloroplastic

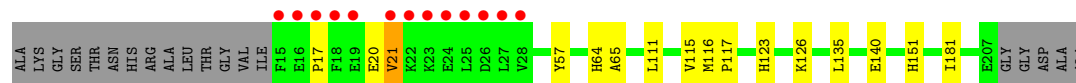
Chain J:





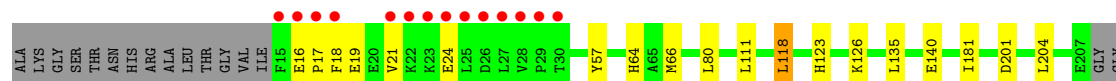
- Molecule 1: Ferritin-4, chloroplastic

Chain K:



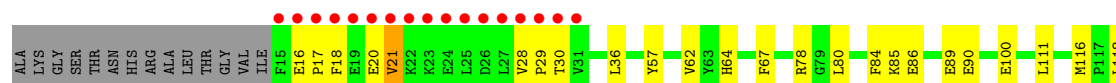
- Molecule 1: Ferritin-4, chloroplastic

Chain L:



- Molecule 1: Ferritin-4, chloroplastic

Chain M:



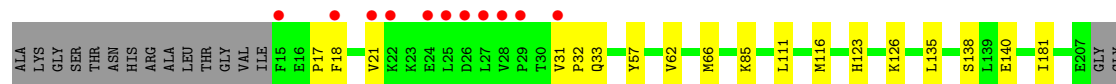
- Molecule 1: Ferritin-4, chloroplastic

Chain N:



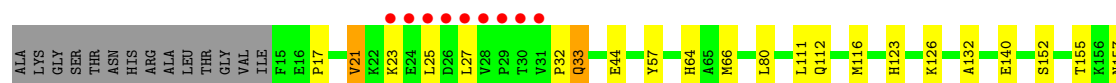
- Molecule 1: Ferritin-4, chloroplastic

Chain O:



- Molecule 1: Ferritin-4, chloroplastic

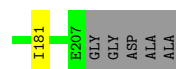
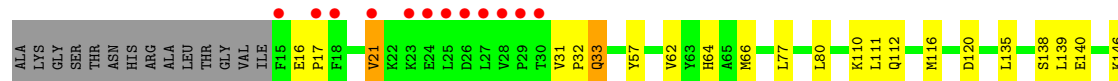
Chain P:





- Molecule 1: Ferritin-4, chloroplastic

Chain Q: 



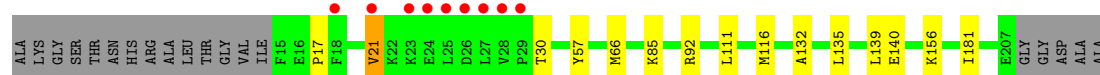
- Molecule 1: Ferritin-4, chloroplastic

Chain R: 



- Molecule 1: Ferritin-4, chloroplastic

Chain S: 



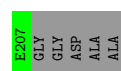
- Molecule 1: Ferritin-4, chloroplastic

Chain T: 



- Molecule 1: Ferritin-4, chloroplastic

Chain U: 

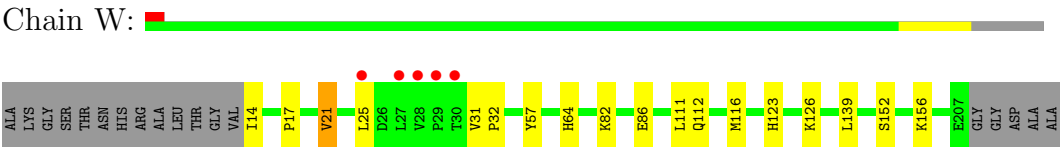


- Molecule 1: Ferritin-4, chloroplastic

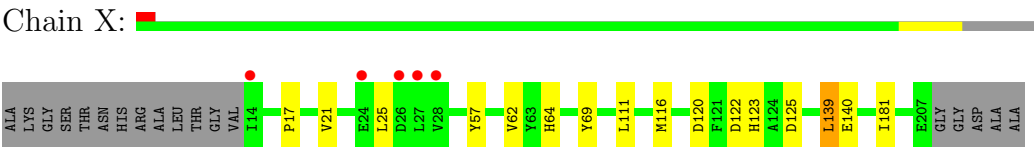
Chain V: 



● Molecule 1: Ferritin-4, chloroplastic



● Molecule 1: Ferritin-4, chloroplastic



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	223.50Å 221.82Å 122.72Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.67 – 1.90 49.90 – 1.90	Depositor EDS
% Data completeness (in resolution range)	96.8 (49.67-1.90) 99.0 (49.90-1.90)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.15 (at 1.90Å)	Xtriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, R_{free}	0.150 , 0.185 0.150 , 0.184	Depositor DCC
R_{free} test set	23767 reflections (5.02%)	DCC
Wilson B-factor (Å ²)	15.2	Xtriage
Anisotropy	0.483	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 32.8	EDS
Estimated twinning fraction	0.007 for k,h,-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 473652 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	42812	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.61% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

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

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	0.34	0/1672	0.47	0/2253
1	B	0.34	0/1684	0.47	0/2270
1	C	0.33	0/1729	0.46	0/2329
1	D	0.33	0/1721	0.44	0/2318
1	E	0.33	0/1547	0.46	0/2080
1	F	0.33	0/1696	0.45	0/2285
1	G	0.33	0/1671	0.46	0/2252
1	H	0.33	0/1688	0.45	0/2274
1	I	0.31	0/1752	0.45	0/2361
1	J	0.32	0/1691	0.46	0/2279
1	K	0.31	0/1641	0.45	0/2211
1	L	0.32	0/1666	0.45	0/2245
1	M	0.33	0/1717	0.46	0/2313
1	N	0.33	0/1697	0.47	0/2286
1	O	0.34	0/1726	0.47	0/2325
1	P	0.32	0/1687	0.44	0/2272
1	Q	0.33	0/1693	0.46	0/2279
1	R	0.33	0/1700	0.46	0/2291
1	S	0.34	0/1682	0.46	0/2266
1	T	0.32	0/1706	0.44	0/2298
1	U	0.34	0/1707	0.45	0/2298
1	V	0.34	0/1692	0.45	0/2278
1	W	0.31	0/1703	0.45	0/2294
1	X	0.32	0/1718	0.46	0/2313
All	All	0.33	0/40586	0.46	0/54670

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1621	0	1604	20	0
1	B	1633	0	1610	18	0
1	C	1669	0	1639	23	0
1	D	1661	0	1633	24	0
1	E	1492	0	1476	13	0
1	F	1636	0	1623	13	0
1	G	1617	0	1587	20	0
1	H	1631	0	1608	17	0
1	I	1677	0	1662	24	0
1	J	1637	0	1614	22	0
1	K	1599	0	1561	14	0
1	L	1618	0	1591	16	0
1	M	1651	0	1634	26	0
1	N	1631	0	1607	16	0
1	O	1657	0	1629	14	0
1	P	1631	0	1596	18	0
1	Q	1630	0	1628	22	0
1	R	1643	0	1617	17	0
1	S	1622	0	1608	11	0
1	T	1650	0	1623	15	0
1	U	1641	0	1628	22	0
1	V	1633	0	1606	13	0
1	W	1649	0	1627	16	0
1	X	1652	0	1631	16	0
2	A	5	0	0	0	0
2	B	6	0	0	0	0
2	C	6	0	0	0	0
2	D	4	0	0	0	0
2	E	5	0	0	0	0
2	F	6	0	0	0	0
2	G	6	0	0	0	0
2	H	5	0	0	0	0
2	I	6	0	0	0	0
2	J	6	0	0	0	0
2	K	6	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	L	7	0	0	0	0
2	M	5	0	0	0	0
2	N	6	0	0	0	0
2	O	5	0	0	0	0
2	P	6	0	0	0	0
2	Q	4	0	0	0	0
2	R	6	0	0	0	0
2	S	5	0	0	0	0
2	T	5	0	0	0	0
2	U	4	0	0	0	0
2	V	5	0	0	0	0
2	W	4	0	0	0	0
2	X	7	0	0	0	0
3	A	4	0	3	0	0
3	E	4	0	3	0	0
3	F	4	0	3	0	0
3	G	4	0	3	0	0
3	H	4	0	3	0	0
3	I	4	0	3	0	0
3	J	4	0	3	0	0
3	L	4	0	3	0	0
3	M	4	0	3	0	0
3	O	4	0	3	0	0
3	P	4	0	3	0	0
3	Q	4	0	3	0	0
3	R	4	0	3	0	0
3	S	4	0	3	0	0
3	T	4	0	3	0	0
3	V	4	0	3	0	0
3	W	4	0	3	0	0
3	X	4	0	3	0	0
4	A	158	0	0	0	0
4	B	149	0	0	1	0
4	C	143	0	0	0	0
4	D	130	0	0	1	0
4	E	150	0	0	1	0
4	F	146	0	0	0	0
4	G	155	0	0	0	0
4	H	145	0	0	1	0
4	I	133	0	0	2	0
4	J	142	0	0	2	0
4	K	129	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	L	136	0	0	0	0
4	M	134	0	0	1	0
4	N	145	0	0	1	0
4	O	162	0	0	0	0
4	P	136	0	0	2	0
4	Q	148	0	0	2	0
4	R	150	0	0	0	0
4	S	137	0	0	1	0
4	T	135	0	0	2	0
4	U	135	0	0	2	0
4	V	146	0	0	1	0
4	W	137	0	0	2	0
4	X	148	0	0	0	0
All	All	42812	0	38696	341	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 4.

All (341) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:T:123:HIS:HD2	1:T:126:LYS:H	1.22	0.86
1:I:21:VAL:HG11	1:W:139[A]:LEU:HD23	1.59	0.83
1:U:133[B]:MET:HE3	1:U:181:ILE:HG23	1.62	0.81
1:T:123:HIS:CD2	1:T:126:LYS:H	1.99	0.79
1:K:123:HIS:HD2	1:K:126:LYS:H	1.31	0.78
1:N:139:LEU:HD22	1:Q:21:VAL:HG21	1.64	0.77
1:E:123:HIS:HD2	1:E:126:LYS:H	1.32	0.77
1:P:123:HIS:HD2	1:P:126:LYS:H	1.33	0.76
1:D:14:ILE:HG21	1:I:112[A]:GLN:NE2	2.01	0.75
1:B:163:ALA:O	1:B:167:GLU:HG2	1.86	0.75
1:H:123:HIS:CD2	1:H:126:LYS:H	2.04	0.75
1:J:14:ILE:HG21	1:Q:112[B]:GLN:HE22	1.52	0.74
1:L:16:GLU:HG2	1:L:19:GLU:HG2	1.67	0.74
1:R:123:HIS:HD2	1:R:126:LYS:H	1.34	0.74
1:J:142[A]:LEU:HD11	1:N:21:VAL:HA	1.69	0.73
1:F:123:HIS:HD2	1:F:126:LYS:H	1.34	0.73
1:D:133[B]:MET:HE3	1:D:181[B]:ILE:HG23	1.70	0.73
1:G:31:VAL:HG22	1:G:33:GLN:HE22	1.54	0.73
1:M:123:HIS:HD2	1:M:126:LYS:H	1.36	0.73
1:R:123:HIS:CD2	1:R:126:LYS:H	2.07	0.73
1:C:31:VAL:HB	1:C:33:GLN:HE21	1.54	0.73
1:H:123:HIS:HD2	1:H:126:LYS:H	1.34	0.72

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:P:123:HIS:CD2	1:P:126:LYS:H	2.07	0.72
1:R:23:LYS:H	1:R:23:LYS:HD2	1.56	0.71
1:E:123:HIS:CD2	1:E:126:LYS:H	2.08	0.70
1:C:31:VAL:HB	1:C:33:GLN:NE2	2.06	0.70
1:F:123:HIS:CD2	1:F:126:LYS:H	2.09	0.69
1:U:133[B]:MET:CE	1:U:181:ILE:HG23	2.22	0.69
1:I:53[B]:ILE:HD12	1:I:99:MET:HG3	1.75	0.69
1:M:62:VAL:HG11	1:M:139:LEU:HD11	1.74	0.68
1:I:133[B]:MET:HE3	1:I:181:ILE:HG23	1.73	0.68
1:C:188[B]:LEU:HD11	1:C:200:PHE:CE2	2.27	0.68
1:K:17:PRO:O	1:K:21:VAL:HB	1.94	0.68
1:A:17:PRO:O	1:A:21:VAL:HB	1.94	0.68
1:M:123:HIS:CD2	1:M:126:LYS:H	2.12	0.68
1:W:123:HIS:HD2	1:W:126:LYS:H	1.42	0.67
1:K:123:HIS:CD2	1:K:126:LYS:H	2.11	0.67
1:I:17:PRO:O	1:I:21:VAL:HG13	1.95	0.67
1:C:130:LEU:HA	1:C:188[B]:LEU:HD23	1.75	0.67
1:X:62:VAL:HG11	1:X:139:LEU:HD21	1.78	0.66
1:N:123:HIS:HD2	1:N:126:LYS:H	1.41	0.66
1:I:62:VAL:HG11	1:I:139:LEU:HD21	1.78	0.66
1:P:17:PRO:O	1:P:21:VAL:HB	1.96	0.65
1:F:18:PHE:HE1	1:G:135[B]:LEU:HD21	1.62	0.65
1:G:123:HIS:CD2	1:G:126:LYS:H	2.15	0.65
1:W:123:HIS:CD2	1:W:126:LYS:H	2.14	0.64
1:N:163:ALA:O	1:N:167:GLU:HG2	1.97	0.64
1:N:123:HIS:CD2	1:N:126:LYS:H	2.16	0.64
1:J:14:ILE:HG21	1:Q:112[B]:GLN:NE2	2.12	0.64
1:C:188[B]:LEU:HD11	1:C:200:PHE:CD2	2.32	0.63
1:J:24:GLU:HG3	1:Q:138[B]:SER:HB2	1.80	0.63
1:J:31:VAL:HG12	1:J:33[B]:GLN:HG2	1.79	0.63
1:A:111:LEU:O	1:G:116[A]:MET:HG3	1.99	0.63
1:G:123:HIS:HD2	1:G:126:LYS:H	1.45	0.63
1:J:135[B]:LEU:HD21	1:N:18:PHE:HE1	1.63	0.62
1:P:25:LEU:HD21	1:U:135[A]:LEU:HD23	1.81	0.62
1:E:78:ARG:HD3	4:E:1888:HOH:O	2.00	0.62
1:O:123:HIS:CD2	1:O:126:LYS:H	2.18	0.62
1:G:25:LEU:HD21	1:O:135[B]:LEU:HD23	1.82	0.62
1:I:133[B]:MET:CE	1:I:181:ILE:HG23	2.29	0.61
1:H:44[A]:GLU:HG3	4:H:2001:HOH:O	1.99	0.61
1:R:31:VAL:HG22	1:R:32:PRO:HD2	1.83	0.61
1:O:123:HIS:HD2	1:O:126:LYS:H	1.48	0.61
1:J:25:LEU:HG	1:Q:135[A]:LEU:HD23	1.81	0.61

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:Q:17:PRO:O	1:Q:21:VAL:HB	2.01	0.61
1:B:14:ILE:HD13	1:B:15:PHE:H	1.66	0.60
1:E:31:VAL:HG22	1:E:32:PRO:HD2	1.82	0.60
1:U:133[B]:MET:HE2	1:U:184:TYR:HB2	1.82	0.60
1:S:140:GLU:HG3	1:S:181[B]:ILE:HD12	1.82	0.60
1:M:17:PRO:O	1:M:21:VAL:HB	2.02	0.60
1:W:152[B]:SER:O	1:W:156[B]:LYS:HG2	2.02	0.60
1:D:14:ILE:HG21	1:I:112[A]:GLN:HE22	1.67	0.60
1:R:23:LYS:HD2	1:R:23:LYS:N	2.17	0.59
1:A:135[A]:LEU:CD2	1:X:25:LEU:HD11	2.32	0.59
1:B:111:LEU:O	1:F:116[A]:MET:HG3	2.01	0.59
1:P:152:SER:HA	1:P:155[B]:THR:HG22	1.83	0.59
1:I:123:HIS:CD2	1:I:126:LYS:H	2.20	0.59
1:S:17:PRO:O	1:S:21:VAL:HB	2.02	0.59
1:B:20:GLU:C	1:V:142[B]:LEU:HD11	2.23	0.59
1:O:17:PRO:O	1:O:21:VAL:HB	2.03	0.59
1:C:66:MET:HE1	1:C:135[A]:LEU:HD23	1.84	0.59
1:F:17:PRO:O	1:F:21:VAL:HB	2.02	0.59
1:L:16:GLU:HG2	1:L:19:GLU:CG	2.33	0.58
1:O:62:VAL:HG12	1:O:66:MET:HE3	1.84	0.58
1:I:21:VAL:HG11	1:W:139[A]:LEU:CD2	2.30	0.58
1:M:140:GLU:HG3	1:M:181:ILE:HD12	1.86	0.58
1:N:17:PRO:O	1:N:21:VAL:HB	2.04	0.58
1:Q:116[A]:MET:HG3	1:U:111:LEU:O	2.03	0.58
1:I:21:VAL:CG1	1:W:139[A]:LEU:HD23	2.31	0.57
1:G:23:LYS:O	1:G:27:LEU:HD23	2.04	0.57
1:M:80[A]:LEU:HD12	1:M:84:PHE:CE2	2.39	0.57
1:H:17:PRO:O	1:H:21:VAL:HB	2.05	0.57
1:F:152:SER:O	1:F:156[A]:LYS:HB2	2.05	0.57
1:D:44[B]:GLU:HG3	4:D:1776:HOH:O	2.04	0.57
1:G:66:MET:HE1	1:G:135[B]:LEU:HD23	1.87	0.56
1:B:133[B]:MET:HE3	1:B:181:ILE:HG23	1.86	0.56
1:D:140:GLU:HG3	1:D:181[B]:ILE:HD12	1.88	0.56
1:I:123:HIS:HD2	1:I:126:LYS:H	1.53	0.56
1:L:123:HIS:HD2	1:L:126:LYS:H	1.54	0.56
1:K:135[A]:LEU:CD2	1:U:25:LEU:HD11	2.36	0.55
1:B:116[B]:MET:HG3	1:F:111:LEU:O	2.05	0.55
1:W:17:PRO:O	1:W:21:VAL:HB	2.06	0.55
1:B:140:GLU:HG3	1:B:181:ILE:HD12	1.88	0.54
1:V:140:GLU:HG3	1:V:181:ILE:HD12	1.89	0.54
1:U:140:GLU:HG3	1:U:181:ILE:HD12	1.90	0.54
1:D:18:PHE:CE1	1:I:118:LEU:HD22	2.42	0.54

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:R:25:LEU:HD21	1:S:135[A]:LEU:HD23	1.90	0.54
1:A:25:LEU:O	1:A:28:VAL:HG12	2.08	0.54
1:N:120[B]:ASP:OD2	1:N:122[B]:ASP:OD1	2.25	0.54
1:A:135[A]:LEU:HD21	1:X:25:LEU:HD11	1.89	0.53
1:P:66:MET:CE	1:P:132:ALA:HB1	2.38	0.53
1:C:111:LEU:O	1:O:116:MET:HG3	2.09	0.53
1:K:135[A]:LEU:HD21	1:U:25:LEU:HD11	1.90	0.53
1:T:66:MET:CE	1:T:132:ALA:HB1	2.38	0.53
1:E:135[B]:LEU:CD2	1:V:25:LEU:HD11	2.39	0.53
1:C:17:PRO:O	1:C:21:VAL:HB	2.09	0.53
1:O:140:GLU:HG3	1:O:181:ILE:HD12	1.91	0.52
1:A:21:VAL:HA	1:H:142[B]:LEU:HD11	1.90	0.52
1:V:44[A]:GLU:HG3	4:V:2636:HOH:O	2.10	0.52
1:U:133[B]:MET:HE1	1:U:184:TYR:HD2	1.75	0.52
1:M:18:PHE:HE1	1:T:135[B]:LEU:HD21	1.74	0.52
1:I:49:VAL:O	1:I:53[B]:ILE:HG12	2.09	0.51
1:H:44[A]:GLU:HG2	1:H:153:VAL:HG13	1.92	0.51
1:Q:62:VAL:O	1:Q:66[A]:MET:HG3	2.09	0.51
1:P:23:LYS:O	1:P:27:LEU:HG	2.11	0.51
1:A:14:ILE:HG21	1:H:112[B]:GLN:HE22	1.76	0.51
1:K:111:LEU:O	1:W:116[A]:MET:HG3	2.10	0.51
1:J:111:LEU:O	1:X:116[B]:MET:HG3	2.11	0.51
1:C:33:GLN:H	1:C:33:GLN:CD	2.14	0.51
1:E:135[B]:LEU:HD23	1:V:25:LEU:HD11	1.93	0.51
1:X:123:HIS:CE1	1:X:125[A]:ASP:HB3	2.45	0.51
1:A:23:LYS:HG2	1:A:23:LYS:O	2.11	0.51
1:A:142[A]:LEU:HD11	1:X:21:VAL:CA	2.40	0.51
1:L:17:PRO:O	1:L:21:VAL:HB	2.11	0.50
1:N:19:GLU:O	1:N:22:LYS:HG2	2.11	0.50
1:E:140:GLU:HG3	1:E:181:ILE:HD12	1.93	0.50
1:X:123:HIS:HE1	1:X:125[A]:ASP:HB3	1.77	0.50
1:A:142[A]:LEU:HD11	1:X:21:VAL:HA	1.92	0.50
1:Q:140:GLU:HG3	1:Q:181:ILE:HD12	1.93	0.50
1:V:66:MET:CE	1:V:132:ALA:HB1	2.42	0.50
1:Q:32:PRO:HG2	1:Q:33[A]:GLN:OE1	2.11	0.49
1:D:133[B]:MET:CE	1:D:181[B]:ILE:HG23	2.42	0.49
1:L:111:LEU:O	1:V:116[A]:MET:HG3	2.12	0.49
1:A:21:VAL:HG22	1:H:138:SER:HB3	1.95	0.49
1:C:66:MET:HE1	1:C:135[A]:LEU:CD2	2.42	0.49
1:X:140:GLU:HG3	1:X:181:ILE:HD12	1.95	0.49
1:X:69:TYR:OH	1:X:122[A]:ASP:HB2	2.13	0.49
4:K:2193:HOH:O	1:W:116[B]:MET:HE3	2.13	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:142[A]:LEU:HD11	1:X:21:VAL:N	2.28	0.49
1:K:140:GLU:HG3	1:K:181:ILE:HD12	1.95	0.49
1:D:116:MET:HG3	1:H:111:LEU:O	2.13	0.48
1:M:29:PRO:HD2	1:M:36[B]:LEU:HD13	1.93	0.48
1:I:133[B]:MET:HE2	1:I:184:TYR:HB2	1.94	0.48
1:L:80[B]:LEU:HD12	1:L:204:LEU:HD23	1.95	0.48
1:B:17:PRO:O	1:B:21:VAL:HG22	2.14	0.48
1:N:116:MET:SD	1:R:112[B]:GLN:OE1	2.72	0.48
1:F:18:PHE:CD1	1:G:118:LEU:HD22	2.48	0.48
1:A:25:LEU:HD13	1:H:135:LEU:HD23	1.95	0.48
1:D:138:SER:HB3	1:W:21:VAL:HG22	1.94	0.48
1:M:89[B]:GLU:HG3	1:S:92:ARG:NH1	2.28	0.48
1:L:80[A]:LEU:HD23	1:L:201:ASP:OD1	2.14	0.47
1:I:165:PHE:CE1	1:I:169:GLU:HG3	2.48	0.47
1:U:59:VAL:HG12	1:U:91:GLU:HG3	1.96	0.47
1:J:18:PHE:HE1	1:Q:135[B]:LEU:HD21	1.79	0.47
1:B:16:GLU:OE1	1:B:19:GLU:HG3	2.15	0.47
1:Q:77:LEU:HD13	1:Q:80[B]:LEU:HD12	1.95	0.47
1:C:25:LEU:O	1:C:28:VAL:HG12	2.13	0.47
1:J:31:VAL:CG1	1:J:33[B]:GLN:HG2	2.44	0.47
4:K:977:HOH:O	1:U:160[B]:VAL:HG13	2.14	0.47
1:P:116:MET:SD	1:T:112[B]:GLN:OE1	2.73	0.47
1:C:184:TYR:O	1:C:188[B]:LEU:HD13	2.14	0.47
1:P:111:LEU:O	1:T:116[B]:MET:HG3	2.14	0.47
1:E:118:LEU:HD22	1:V:18:PHE:CD1	2.49	0.47
1:D:28:VAL:HG22	4:I:2736:HOH:O	2.15	0.47
1:R:80[A]:LEU:HD12	1:R:204:LEU:HD23	1.97	0.47
1:M:116[A]:MET:HG3	1:S:111:LEU:O	2.14	0.47
1:G:31:VAL:HG22	1:G:32:PRO:HD2	1.97	0.47
1:Q:33[A]:GLN:NE2	1:Q:33[A]:GLN:H	2.12	0.47
1:G:140:GLU:HG3	1:G:181:ILE:HD12	1.97	0.47
1:I:53[B]:ILE:HD12	1:I:99:MET:CG	2.45	0.46
1:M:80[A]:LEU:HD11	1:M:133:MET:CE	2.46	0.46
1:P:32:PRO:HD2	1:P:33:GLN:OE1	2.15	0.46
1:D:18:PHE:CD1	1:I:118:LEU:HD22	2.50	0.46
1:J:140:GLU:HG3	1:J:181:ILE:HD12	1.97	0.46
1:M:86:GLU:O	1:M:90[B]:GLU:HG3	2.14	0.46
1:G:25:LEU:HD22	1:O:138:SER:CB	2.46	0.46
1:E:66[A]:MET:CE	1:E:135[A]:LEU:HD23	2.45	0.46
1:L:123:HIS:CD2	1:L:126:LYS:H	2.32	0.46
1:E:77:LEU:HD13	1:E:80[A]:LEU:HD23	1.96	0.46
1:I:44[A]:GLU:HG3	4:I:3135:HOH:O	2.16	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:112[B]:GLN:OE1	1:W:14:ILE:HG21	2.16	0.46
1:P:25:LEU:HD11	1:U:135[A]:LEU:CD2	2.46	0.46
1:C:165:PHE:CE1	1:C:169:GLU:HG3	2.51	0.46
1:A:116[A]:MET:HG3	1:G:111:LEU:O	2.16	0.46
1:M:78:ARG:HD3	4:M:745:HOH:O	2.16	0.46
1:D:31:VAL:HG12	1:D:33:GLN:OE1	2.16	0.46
1:R:17:PRO:O	1:R:21:VAL:HG13	2.16	0.46
1:R:31:VAL:HG22	1:R:33:GLN:OE1	2.15	0.46
1:X:17:PRO:O	1:X:21:VAL:HB	2.15	0.46
1:U:62:VAL:HG11	1:U:139:LEU:HD11	1.98	0.46
1:G:21:VAL:HG13	1:G:25:LEU:HD23	1.97	0.46
1:J:24:GLU:HG3	1:Q:138[B]:SER:CB	2.45	0.45
1:D:111:LEU:O	1:H:116:MET:HG3	2.16	0.45
1:K:116:MET:HG3	1:W:111:LEU:O	2.15	0.45
1:A:31:VAL:HA	1:A:32:PRO:HD3	1.85	0.45
1:G:31:VAL:HG22	1:G:33:GLN:NE2	2.27	0.45
1:Q:31:VAL:HG22	4:Q:3414:HOH:O	2.16	0.45
1:Q:110[A]:LYS:HE2	4:Q:2516:HOH:O	2.16	0.45
1:U:133[B]:MET:HE2	1:U:184:TYR:CB	2.45	0.45
1:U:39:GLN:HG2	1:U:40:LYS:HG2	1.99	0.45
1:T:19:GLU:O	1:T:22:LYS:HB2	2.17	0.45
1:E:85:LYS:HE2	1:I:96:GLU:CD	2.37	0.45
1:A:135[A]:LEU:HD23	1:X:25:LEU:HD11	1.98	0.45
1:H:140:GLU:HG3	1:H:181[A]:ILE:HD12	1.98	0.45
1:I:97:LYS:HE3	1:I:169:GLU:HB3	1.98	0.45
1:F:29:PRO:HG2	1:F:35:SER:O	2.16	0.45
1:J:24:GLU:HG3	1:Q:138[A]:SER:CB	2.47	0.44
1:V:44[A]:GLU:HG2	1:V:153:VAL:HG13	1.99	0.44
1:R:21:VAL:HG11	1:S:139:LEU:HD23	1.98	0.44
1:J:78:ARG:HD3	4:J:1050:HOH:O	2.17	0.44
1:N:112[B]:GLN:HA	1:R:116[B]:MET:HE3	1.99	0.44
4:K:977:HOH:O	1:U:160[A]:VAL:HG23	2.17	0.44
1:D:177:ALA:O	1:D:181[A]:ILE:HG12	2.17	0.44
1:B:29:PRO:HG2	1:B:35:SER:O	2.16	0.44
1:J:116:MET:HG3	1:X:111:LEU:O	2.17	0.44
1:L:66:MET:HE1	1:L:135[A]:LEU:HD23	1.99	0.44
1:L:18:PHE:CD2	1:M:118:LEU:HD23	2.52	0.44
1:L:18:PHE:CG	1:M:118:LEU:HD23	2.53	0.44
1:C:96:GLU:CD	1:O:85:LYS:HE2	2.38	0.44
1:H:83:PHE:HZ	1:H:181[B]:ILE:HD11	1.83	0.44
1:B:133[B]:MET:CE	1:B:181:ILE:HG23	2.45	0.43
1:D:32:PRO:HG2	1:D:33:GLN:NE2	2.33	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:140:GLU:HG3	1:C:181:ILE:HD12	1.98	0.43
1:G:74:ASN:H	1:G:74:ASN:ND2	2.16	0.43
1:K:151:HIS:NE2	1:U:160[B]:VAL:HG11	2.33	0.43
1:H:31:VAL:HA	1:H:32:PRO:HD3	1.82	0.43
1:B:160[A]:VAL:HG23	4:B:2658:HOH:O	2.17	0.43
1:P:207[B]:GLU:HA	1:P:207[B]:GLU:OE2	2.18	0.43
1:I:140:GLU:HG3	1:I:181:ILE:HD12	2.01	0.43
1:F:18:PHE:CE1	1:G:118:LEU:HD22	2.54	0.43
1:L:66:MET:HE1	1:L:135[A]:LEU:CD2	2.48	0.43
1:H:25:LEU:O	1:H:28:VAL:HG12	2.19	0.43
1:J:15:PHE:CE2	1:Q:146:LYS:HE3	2.54	0.43
1:R:33:GLN:H	1:R:33:GLN:CD	2.22	0.43
1:B:16:GLU:HA	1:B:17:PRO:HD2	1.85	0.43
1:O:31:VAL:HG13	1:O:33[A]:GLN:OE1	2.19	0.43
1:M:111:LEU:O	1:S:116:MET:HG3	2.19	0.43
1:P:112:GLN:HA	1:T:116[B]:MET:HE2	1.99	0.43
4:P:2489:HOH:O	1:T:89[B]:GLU:HG3	2.17	0.43
1:N:140:GLU:HG3	1:N:181:ILE:HD12	2.00	0.43
1:C:31:VAL:HA	1:C:32:PRO:HD3	1.87	0.43
1:K:65:ALA:HB2	1:W:111:LEU:HD12	2.01	0.43
1:D:17:PRO:O	1:D:21:VAL:HB	2.19	0.43
1:O:18:PHE:HA	1:O:21:VAL:HG12	2.01	0.43
1:K:116:MET:HA	1:K:117:PRO:HD3	1.93	0.43
1:J:19:GLU:O	1:J:22:LYS:HG3	2.19	0.43
1:R:152:SER:O	1:R:156:LYS:HG3	2.19	0.43
1:Q:111:LEU:O	1:U:116:MET:HG3	2.18	0.43
1:B:16:GLU:O	1:B:20:GLU:HG3	2.18	0.42
1:N:111:LEU:O	1:R:116[A]:MET:HG3	2.19	0.42
1:U:110:LYS:HE2	4:U:2980:HOH:O	2.19	0.42
1:G:66:MET:HE1	1:G:135[B]:LEU:CD2	2.48	0.42
1:A:140:GLU:HG3	1:A:181:ILE:HD12	2.02	0.42
1:D:39:GLN:HG2	1:D:40:LYS:HG3	2.02	0.42
1:D:16:GLU:HA	1:D:17:PRO:HD2	1.90	0.42
1:A:80[B]:LEU:HD23	1:A:201:ASP:OD1	2.19	0.42
1:M:80[A]:LEU:HD11	1:M:133:MET:HE2	2.01	0.42
1:P:167:GLU:OE1	4:P:2510:HOH:O	2.22	0.42
1:X:123:HIS:HE1	1:X:125[B]:ASP:HB2	1.85	0.42
1:H:28:VAL:HG13	1:H:28:VAL:O	2.20	0.42
1:M:20:GLU:C	1:T:142[B]:LEU:HD11	2.39	0.42
1:O:31:VAL:HA	1:O:32:PRO:HD3	1.85	0.42
1:T:31:VAL:HG13	1:T:33[A]:GLN:OE1	2.20	0.42
1:I:28:VAL:HG22	4:W:3518:HOH:O	2.20	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:J:77:LEU:HD13	1:J:80[B]:LEU:HD12	2.00	0.42
1:G:32:PRO:HG2	1:G:33:GLN:OE1	2.20	0.42
1:W:82:LYS:O	1:W:86:GLU:HG3	2.20	0.42
1:D:96:GLU:CD	1:H:85:LYS:HE3	2.40	0.41
1:B:33:GLN:O	1:B:33:GLN:HG2	2.20	0.41
1:C:116:MET:HG3	1:O:111:LEU:O	2.19	0.41
1:J:142[A]:LEU:HD11	1:N:21:VAL:CA	2.46	0.41
1:P:116:MET:HG3	1:T:111:LEU:O	2.19	0.41
1:M:100:GLU:OE1	1:S:85:LYS:NZ	2.47	0.41
1:B:133[B]:MET:HE2	1:B:184:TYR:HB2	2.01	0.41
1:F:140:GLU:HG3	1:F:181:ILE:HD12	2.02	0.41
1:P:44[A]:GLU:CD	1:P:157:ASN:HD21	2.23	0.41
1:J:24:GLU:HG3	1:Q:138[A]:SER:HB3	2.01	0.41
1:K:116:MET:SD	1:W:112[B]:GLN:NE2	2.93	0.41
4:N:3353:HOH:O	1:R:116[B]:MET:HE3	2.20	0.41
1:S:156:LYS:HE3	4:S:3449:HOH:O	2.20	0.41
1:Q:16:GLU:HA	1:Q:17:PRO:HD2	1.85	0.41
1:M:89[B]:GLU:HG3	1:S:92:ARG:HH12	1.84	0.41
1:I:66:MET:CE	1:I:132:ALA:HB1	2.50	0.41
1:R:140:GLU:HG3	1:R:181:ILE:HD12	2.01	0.41
1:J:98:LEU:HG	1:J:170:TYR:OH	2.21	0.41
1:A:24:GLU:HG3	1:A:24:GLU:O	2.21	0.41
1:L:24:GLU:OE1	1:M:142[A]:LEU:HD22	2.21	0.41
1:J:157[A]:ASN:ND2	4:J:3300:HOH:O	2.54	0.41
1:M:120[B]:ASP:OD1	1:M:122[B]:ASP:CG	2.59	0.41
1:U:42[B]:VAL:HG22	4:U:869:HOH:O	2.21	0.41
1:C:89[A]:GLU:OE2	1:C:89[A]:GLU:HA	2.21	0.41
1:W:31:VAL:HA	1:W:32:PRO:HD3	1.83	0.41
1:V:44[A]:GLU:HG2	1:V:153:VAL:CG1	2.51	0.41
1:B:160[A]:VAL:HG21	1:V:151:HIS:NE2	2.35	0.41
1:P:140:GLU:HG3	1:P:181:ILE:HD12	2.02	0.41
1:V:32:PRO:HG2	1:V:33:GLN:NE2	2.36	0.41
1:K:115:VAL:HG23	4:W:2194:HOH:O	2.20	0.41
1:F:23:LYS:HD3	1:F:23:LYS:O	2.21	0.41
1:B:21:VAL:HG13	1:V:142[B]:LEU:HD12	2.03	0.41
1:M:67:PHE:CE1	1:M:85:LYS:HE3	2.55	0.41
1:M:28:VAL:O	1:M:28:VAL:HG13	2.21	0.41
1:Q:139[B]:LEU:HD12	1:Q:139[B]:LEU:HA	1.89	0.41
1:M:80[A]:LEU:HD12	1:M:84:PHE:HE2	1.85	0.40
1:L:118:LEU:HD22	1:T:18:PHE:CD1	2.56	0.40
1:G:31:VAL:CG2	1:G:32:PRO:HD2	2.52	0.40
1:M:16:GLU:HA	1:M:17:PRO:HD2	1.82	0.40

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:28:VAL:O	1:A:28:VAL:HG13	2.21	0.40
1:X:123:HIS:CE1	1:X:125[B]:ASP:HB2	2.55	0.40
1:T:110:LYS:HE2	4:T:3460:HOH:O	2.20	0.40
1:P:80[B]:LEU:HD13	1:P:204:LEU:HD23	2.02	0.40
1:U:17:PRO:O	1:U:21:VAL:HB	2.21	0.40
1:E:98:LEU:HG	1:E:170:TYR:OH	2.21	0.40
1:T:206:HIS:HE1	4:T:3470:HOH:O	2.03	0.40
1:K:135[A]:LEU:HD23	1:U:25:LEU:HD11	2.03	0.40
1:E:80[A]:LEU:HD13	1:E:204:LEU:HD23	2.03	0.40
1:N:31:VAL:HA	1:N:32:PRO:HD3	1.86	0.40
1:S:66:MET:CE	1:S:132:ALA:HB1	2.52	0.40
1:O:62:VAL:HG12	1:O:66:MET:CE	2.52	0.40
1:N:22:LYS:HG2	1:N:22:LYS:H	1.73	0.40
1:L:66:MET:HE1	1:L:135[A]:LEU:HG	2.03	0.40
1:C:163:ALA:O	1:C:167:GLU:HG3	2.21	0.40
1:L:140:GLU:HG3	1:L:181:ILE:HD12	2.03	0.40
1:F:31:VAL:HA	1:F:32:PRO:HD3	1.87	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	201/212 (95%)	201 (100%)	0	0	100	100
1	B	203/212 (96%)	203 (100%)	0	0	100	100
1	C	208/212 (98%)	208 (100%)	0	0	100	100
1	D	207/212 (98%)	206 (100%)	1 (0%)	0	100	100
1	E	186/212 (88%)	186 (100%)	0	0	100	100
1	F	204/212 (96%)	204 (100%)	0	0	100	100
1	G	201/212 (95%)	201 (100%)	0	0	100	100
1	H	203/212 (96%)	203 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	I	211/212 (100%)	211 (100%)	0	0	100	100
1	J	203/212 (96%)	202 (100%)	1 (0%)	0	100	100
1	K	197/212 (93%)	197 (100%)	0	0	100	100
1	L	200/212 (94%)	200 (100%)	0	0	100	100
1	M	206/212 (97%)	206 (100%)	0	0	100	100
1	N	204/212 (96%)	204 (100%)	0	0	100	100
1	O	208/212 (98%)	208 (100%)	0	0	100	100
1	P	202/212 (95%)	202 (100%)	0	0	100	100
1	Q	204/212 (96%)	204 (100%)	0	0	100	100
1	R	205/212 (97%)	204 (100%)	1 (0%)	0	100	100
1	S	203/212 (96%)	203 (100%)	0	0	100	100
1	T	204/212 (96%)	204 (100%)	0	0	100	100
1	U	206/212 (97%)	205 (100%)	1 (0%)	0	100	100
1	V	203/212 (96%)	203 (100%)	0	0	100	100
1	W	205/212 (97%)	205 (100%)	0	0	100	100
1	X	207/212 (98%)	207 (100%)	0	0	100	100
All	All	4881/5088 (96%)	4877 (100%)	4 (0%)	0	100	100

There are no Ramachandran outliers to report.

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	177/178 (99%)	172 (97%)	5 (3%)	56	45
1	B	179/178 (101%)	175 (98%)	4 (2%)	64	57
1	C	184/178 (103%)	178 (97%)	6 (3%)	50	37
1	D	183/178 (103%)	180 (98%)	3 (2%)	75	70
1	E	163/178 (92%)	157 (96%)	6 (4%)	45	32
1	F	180/178 (101%)	176 (98%)	4 (2%)	64	57

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	G	177/178 (99%)	170 (96%)	7 (4%)	42	29
1	H	179/178 (101%)	175 (98%)	4 (2%)	64	57
1	I	187/178 (105%)	181 (97%)	6 (3%)	51	39
1	J	179/178 (101%)	174 (97%)	5 (3%)	56	45
1	K	173/178 (97%)	169 (98%)	4 (2%)	63	55
1	L	176/178 (99%)	173 (98%)	3 (2%)	73	68
1	M	182/178 (102%)	174 (96%)	8 (4%)	39	25
1	N	180/178 (101%)	178 (99%)	2 (1%)	84	82
1	O	184/178 (103%)	183 (100%)	1 (0%)	94	94
1	P	179/178 (101%)	175 (98%)	4 (2%)	64	57
1	Q	180/178 (101%)	174 (97%)	6 (3%)	50	37
1	R	180/178 (101%)	175 (97%)	5 (3%)	56	45
1	S	179/178 (101%)	176 (98%)	3 (2%)	73	68
1	T	181/178 (102%)	175 (97%)	6 (3%)	50	37
1	U	182/178 (102%)	176 (97%)	6 (3%)	50	37
1	V	180/178 (101%)	177 (98%)	3 (2%)	73	68
1	W	181/178 (102%)	177 (98%)	4 (2%)	64	57
1	X	183/178 (103%)	179 (98%)	4 (2%)	64	57
All	All	4308/4272 (101%)	4199 (98%)	109 (2%)	60	50

All (109) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	21	VAL
1	A	30	THR
1	A	31	VAL
1	A	57	TYR
1	A	64	HIS
1	B	14	ILE
1	B	33	GLN
1	B	57	TYR
1	B	64	HIS
1	C	26	ASP
1	C	33	GLN
1	C	57	TYR
1	C	64	HIS

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Mol	Chain	Res	Type
1	C	120[A]	ASP
1	C	120[B]	ASP
1	D	31	VAL
1	D	57	TYR
1	D	64	HIS
1	E	31	VAL
1	E	44	GLU
1	E	57	TYR
1	E	80[A]	LEU
1	E	80[B]	LEU
1	E	118	LEU
1	F	21	VAL
1	F	33	GLN
1	F	57	TYR
1	F	64	HIS
1	G	31	VAL
1	G	33	GLN
1	G	57	TYR
1	G	64	HIS
1	G	118	LEU
1	G	120[A]	ASP
1	G	120[B]	ASP
1	H	21	VAL
1	H	57	TYR
1	H	64	HIS
1	H	118	LEU
1	I	23	LYS
1	I	25	LEU
1	I	57	TYR
1	I	64	HIS
1	I	139	LEU
1	I	156	LYS
1	J	14	ILE
1	J	33[A]	GLN
1	J	33[B]	GLN
1	J	57	TYR
1	J	64	HIS
1	K	20	GLU
1	K	21	VAL
1	K	57	TYR
1	K	64	HIS
1	L	57	TYR

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Mol	Chain	Res	Type
1	L	64	HIS
1	L	118	LEU
1	M	21	VAL
1	M	30	THR
1	M	57	TYR
1	M	64	HIS
1	M	120[A]	ASP
1	M	120[B]	ASP
1	M	142[A]	LEU
1	M	142[B]	LEU
1	N	57	TYR
1	N	167	GLU
1	O	57	TYR
1	P	21	VAL
1	P	33	GLN
1	P	57	TYR
1	P	64	HIS
1	Q	21	VAL
1	Q	33[A]	GLN
1	Q	33[B]	GLN
1	Q	57	TYR
1	Q	64	HIS
1	Q	120	ASP
1	R	23	LYS
1	R	31	VAL
1	R	33	GLN
1	R	57	TYR
1	R	64	HIS
1	S	21	VAL
1	S	30	THR
1	S	57	TYR
1	T	57	TYR
1	T	64	HIS
1	T	120[A]	ASP
1	T	120[B]	ASP
1	T	160[A]	VAL
1	T	160[B]	VAL
1	U	21	VAL
1	U	31	VAL
1	U	57	TYR
1	U	64	HIS
1	U	160[A]	VAL

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Mol	Chain	Res	Type
1	U	160[B]	VAL
1	V	23	LYS
1	V	57	TYR
1	V	64	HIS
1	W	21	VAL
1	W	25	LEU
1	W	57	TYR
1	W	64	HIS
1	X	57	TYR
1	X	64	HIS
1	X	120	ASP
1	X	139	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (38) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	102	GLN
1	A	157	ASN
1	B	33	GLN
1	B	52	GLN
1	B	58	ASN
1	B	112	GLN
1	C	33	GLN
1	C	123	HIS
1	C	157	ASN
1	D	123	HIS
1	E	74	ASN
1	E	123	HIS
1	F	33	GLN
1	F	112	GLN
1	F	123	HIS
1	G	33	GLN
1	G	58	ASN
1	G	74	ASN
1	G	123	HIS
1	H	123	HIS
1	I	58	ASN
1	I	123	HIS
1	I	157	ASN
1	K	123	HIS
1	L	123	HIS
1	M	123	HIS

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Mol	Chain	Res	Type
1	N	123	HIS
1	O	123	HIS
1	O	206	HIS
1	P	112	GLN
1	P	123	HIS
1	P	157	ASN
1	R	123	HIS
1	R	206	HIS
1	T	123	HIS
1	T	206	HIS
1	V	206	HIS
1	W	123	HIS

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 148 ligands modelled in this entry, 130 are monoatomic - leaving 18 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	$\# Z > 2$
3	ACY	A	217	-	3,3,3	0.65	0	3,3,3	0.70	0
3	ACY	E	218	-	3,3,3	0.70	0	3,3,3	0.55	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	ACY	F	219	-	3,3,3	0.67	0	3,3,3	0.63	0
3	ACY	G	219	-	3,3,3	0.66	0	3,3,3	0.61	0
3	ACY	H	218	-	3,3,3	0.64	0	3,3,3	0.72	0
3	ACY	I	218	-	3,3,3	0.64	0	3,3,3	0.70	0
3	ACY	J	218	-	3,3,3	0.65	0	3,3,3	0.67	0
3	ACY	L	219	-	3,3,3	0.66	0	3,3,3	0.65	0
3	ACY	M	218	-	3,3,3	0.68	0	3,3,3	0.61	0
3	ACY	O	218	-	3,3,3	0.67	0	3,3,3	0.59	0
3	ACY	P	218	-	3,3,3	0.66	0	3,3,3	0.64	0
3	ACY	Q	217	-	3,3,3	0.72	0	3,3,3	0.51	0
3	ACY	R	219	-	3,3,3	0.63	0	3,3,3	0.79	0
3	ACY	S	218	-	3,3,3	0.66	0	3,3,3	0.66	0
3	ACY	T	218	-	3,3,3	0.67	0	3,3,3	0.59	0
3	ACY	V	221	-	3,3,3	0.66	0	3,3,3	0.63	0
3	ACY	W	231	-	3,3,3	0.70	0	3,3,3	0.59	0
3	ACY	X	241	-	3,3,3	0.68	0	3,3,3	0.59	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	ACY	A	217	-	-	0/0/0/0	0/0/0/0
3	ACY	E	218	-	-	0/0/0/0	0/0/0/0
3	ACY	F	219	-	-	0/0/0/0	0/0/0/0
3	ACY	G	219	-	-	0/0/0/0	0/0/0/0
3	ACY	H	218	-	-	0/0/0/0	0/0/0/0
3	ACY	I	218	-	-	0/0/0/0	0/0/0/0
3	ACY	J	218	-	-	0/0/0/0	0/0/0/0
3	ACY	L	219	-	-	0/0/0/0	0/0/0/0
3	ACY	M	218	-	-	0/0/0/0	0/0/0/0
3	ACY	O	218	-	-	0/0/0/0	0/0/0/0
3	ACY	P	218	-	-	0/0/0/0	0/0/0/0
3	ACY	Q	217	-	-	0/0/0/0	0/0/0/0
3	ACY	R	219	-	-	0/0/0/0	0/0/0/0
3	ACY	S	218	-	-	0/0/0/0	0/0/0/0
3	ACY	T	218	-	-	0/0/0/0	0/0/0/0
3	ACY	V	221	-	-	0/0/0/0	0/0/0/0
3	ACY	W	231	-	-	0/0/0/0	0/0/0/0
3	ACY	X	241	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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	194/212 (91%)	-0.27	13 (6%)	17 17	8, 13, 65, 106	0
1	B	194/212 (91%)	-0.22	16 (8%)	12 11	8, 13, 69, 96	0
1	C	194/212 (91%)	-0.49	4 (2%)	60 62	8, 13, 42, 81	0
1	D	194/212 (91%)	-0.33	9 (4%)	31 31	9, 14, 56, 89	0
1	E	178/212 (83%)	-0.59	2 (1%)	77 79	8, 13, 32, 73	0
1	F	194/212 (91%)	-0.28	16 (8%)	12 11	7, 13, 62, 94	0
1	G	193/212 (91%)	-0.53	2 (1%)	79 81	9, 13, 43, 75	0
1	H	193/212 (91%)	-0.28	15 (7%)	13 13	7, 14, 66, 93	0
1	I	193/212 (91%)	-0.52	3 (1%)	68 70	8, 15, 46, 70	0
1	J	194/212 (91%)	-0.26	12 (6%)	20 20	7, 14, 55, 100	0
1	K	193/212 (91%)	-0.25	13 (6%)	17 17	9, 16, 63, 91	0
1	L	193/212 (91%)	-0.24	14 (7%)	15 14	8, 15, 74, 100	0
1	M	193/212 (91%)	-0.29	18 (9%)	9 8	9, 13, 69, 92	0
1	N	193/212 (91%)	-0.41	7 (3%)	41 42	8, 14, 50, 82	0
1	O	193/212 (91%)	-0.36	11 (5%)	23 23	7, 12, 60, 86	0
1	P	193/212 (91%)	-0.33	9 (4%)	30 30	9, 15, 52, 93	0
1	Q	193/212 (91%)	-0.38	12 (6%)	20 20	7, 13, 65, 95	0
1	R	196/212 (92%)	-0.52	6 (3%)	47 48	7, 13, 45, 90	0
1	S	193/212 (91%)	-0.40	9 (4%)	30 30	7, 12, 53, 89	0
1	T	194/212 (91%)	-0.37	11 (5%)	23 23	9, 15, 55, 91	0
1	U	193/212 (91%)	-0.49	4 (2%)	60 62	8, 13, 50, 89	0
1	V	193/212 (91%)	-0.54	1 (0%)	88 90	8, 14, 41, 78	0
1	W	194/212 (91%)	-0.49	5 (2%)	53 55	10, 15, 55, 89	0
1	X	194/212 (91%)	-0.49	5 (2%)	53 55	8, 14, 46, 82	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å ²)	Q<0.9
All	All	4629/5088 (90%)	-0.39	217 (4%)	30	30	7, 14, 57, 106	0

All (217) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	J	28	VAL	9.7
1	H	28	VAL	8.5
1	J	25	LEU	8.4
1	B	25	LEU	8.3
1	A	27	LEU	8.2
1	M	27	LEU	7.9
1	A	14	ILE	7.8
1	A	25	LEU	7.6
1	L	27	LEU	7.6
1	H	27	LEU	7.5
1	B	21	VAL	7.4
1	J	24	GLU	7.3
1	P	27	LEU	7.1
1	J	26	ASP	7.0
1	B	14	ILE	7.0
1	J	27	LEU	6.7
1	L	25	LEU	6.6
1	L	21	VAL	6.5
1	H	25	LEU	6.4
1	P	25	LEU	6.2
1	K	21	VAL	6.1
1	B	27	LEU	6.0
1	P	28	VAL	5.9
1	A	21	VAL	5.7
1	H	29	PRO	5.5
1	D	14	ILE	5.5
1	J	30	THR	5.5
1	L	28	VAL	5.3
1	H	30	THR	5.1
1	S	25	LEU	5.0
1	M	26	ASP	5.0
1	A	26	ASP	5.0
1	M	28	VAL	4.9
1	F	25	LEU	4.9
1	M	30	THR	4.9
1	K	23	LYS	4.9
1	D	25	LEU	4.9

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Mol	Chain	Res	Type	RSRZ
1	D	27	LEU	4.9
1	C	14	ILE	4.9
1	A	24	GLU	4.9
1	D	26	ASP	4.9
1	P	26	ASP	4.8
1	L	26	ASP	4.8
1	S	27	LEU	4.8
1	M	23	LYS	4.7
1	E	30	THR	4.7
1	O	18	PHE	4.6
1	S	26	ASP	4.6
1	F	27	LEU	4.6
1	K	17	PRO	4.5
1	T	14	ILE	4.5
1	K	18	PHE	4.5
1	K	22	LYS	4.5
1	M	31	VAL	4.5
1	N	21	VAL	4.5
1	U	26	ASP	4.4
1	M	29	PRO	4.4
1	L	24	GLU	4.4
1	Q	25	LEU	4.4
1	Q	26	ASP	4.4
1	K	26	ASP	4.3
1	N	17	PRO	4.3
1	F	23	LYS	4.2
1	B	26	ASP	4.2
1	L	23	LYS	4.2
1	H	23	LYS	4.2
1	H	26	ASP	4.2
1	K	16	GLU	4.1
1	B	23	LYS	4.1
1	K	27	LEU	4.1
1	O	27	LEU	4.1
1	S	23	LYS	4.1
1	K	15	PHE	4.1
1	W	28	VAL	4.1
1	K	25	LEU	4.0
1	B	18	PHE	4.0
1	F	28	VAL	4.0
1	H	31	VAL	4.0
1	F	14	ILE	4.0

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Mol	Chain	Res	Type	RSRZ
1	P	23	LYS	4.0
1	B	28	VAL	3.9
1	M	22	LYS	3.9
1	N	18	PHE	3.9
1	T	18	PHE	3.9
1	L	18	PHE	3.8
1	M	21	VAL	3.7
1	O	21	VAL	3.7
1	F	17	PRO	3.7
1	H	18	PHE	3.7
1	T	23	LYS	3.7
1	K	24	GLU	3.6
1	L	29	PRO	3.6
1	B	30	THR	3.6
1	J	29	PRO	3.6
1	P	30	THR	3.6
1	A	29	PRO	3.6
1	D	28	VAL	3.5
1	X	27	LEU	3.5
1	F	30	THR	3.5
1	T	25	LEU	3.5
1	U	25	LEU	3.4
1	L	15	PHE	3.4
1	R	27	LEU	3.4
1	U	24	GLU	3.4
1	B	17	PRO	3.4
1	O	25	LEU	3.4
1	Q	23	LYS	3.3
1	B	16	GLU	3.3
1	T	17	PRO	3.3
1	J	23	LYS	3.3
1	F	21	VAL	3.3
1	H	22	LYS	3.2
1	N	15	PHE	3.2
1	Q	21	VAL	3.2
1	T	22	LYS	3.2
1	F	26	ASP	3.2
1	M	18	PHE	3.1
1	O	28	VAL	3.1
1	T	27	LEU	3.1
1	I	21	VAL	3.1
1	L	17	PRO	3.1

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Mol	Chain	Res	Type	RSRZ
1	E	31	VAL	3.1
1	R	25	LEU	3.1
1	R	28	VAL	3.0
1	X	24	GLU	3.0
1	W	25	LEU	3.0
1	C	26	ASP	3.0
1	B	22	LYS	3.0
1	B	24	GLU	3.0
1	P	29	PRO	3.0
1	A	30	THR	3.0
1	T	26	ASP	2.9
1	L	30	THR	2.9
1	H	21	VAL	2.9
1	K	28	VAL	2.9
1	Q	24	GLU	2.9
1	N	16	GLU	2.8
1	D	23	LYS	2.8
1	L	22	LYS	2.8
1	F	15	PHE	2.8
1	F	18	PHE	2.8
1	P	24	GLU	2.8
1	N	22	LYS	2.8
1	O	31	VAL	2.8
1	A	23	LYS	2.7
1	U	23[A]	LYS	2.7
1	A	28	VAL	2.7
1	T	24	GLU	2.7
1	X	28	VAL	2.7
1	Q	18	PHE	2.7
1	W	29	PRO	2.7
1	H	24	GLU	2.7
1	M	19	GLU	2.7
1	M	24	GLU	2.7
1	K	19	GLU	2.7
1	I	18	PHE	2.7
1	O	15	PHE	2.6
1	P	31	VAL	2.6
1	F	29	PRO	2.6
1	M	25	LEU	2.6
1	M	17	PRO	2.6
1	Q	29	PRO	2.6
1	T	21	VAL	2.6

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Mol	Chain	Res	Type	RSRZ
1	X	26	ASP	2.6
1	I	25	LEU	2.5
1	F	24	GLU	2.5
1	H	32	PRO	2.5
1	B	31	VAL	2.5
1	L	16	GLU	2.5
1	O	26	ASP	2.5
1	Q	28	VAL	2.5
1	A	18	PHE	2.5
1	W	30	THR	2.5
1	F	31	VAL	2.5
1	A	31	VAL	2.4
1	D	125[A]	ASP	2.4
1	Q	27	LEU	2.4
1	N	26	ASP	2.4
1	G	27	LEU	2.4
1	S	24	GLU	2.4
1	Q	30	THR	2.3
1	D	17	PRO	2.3
1	Q	17	PRO	2.3
1	M	20	GLU	2.3
1	R	29	PRO	2.3
1	H	15	PHE	2.2
1	O	22	LYS	2.2
1	O	29	PRO	2.2
1	S	21	VAL	2.2
1	R	26	ASP	2.2
1	W	27	LEU	2.2
1	B	29	PRO	2.2
1	S	28	VAL	2.2
1	O	24[A]	GLU	2.2
1	F	22	LYS	2.2
1	H	17	PRO	2.2
1	D	24	GLU	2.2
1	R	30	THR	2.2
1	F	16	GLU	2.1
1	T	28	VAL	2.1
1	Q	15	PHE	2.1
1	J	21	VAL	2.1
1	J	22	LYS	2.1
1	M	125[A]	ASP	2.1
1	M	16	GLU	2.1

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Mol	Chain	Res	Type	RSRZ
1	J	31	VAL	2.1
1	G	18	PHE	2.1
1	S	18	PHE	2.1
1	S	29	PRO	2.1
1	B	15	PHE	2.1
1	X	14	ILE	2.0
1	A	22	LYS	2.0
1	C	15	PHE	2.0
1	J	14	ILE	2.0
1	M	15	PHE	2.0
1	V	18	PHE	2.0
1	C	23	LYS	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	CA	T	217	1/1	0.14	47.67	59,59,59,59	0
2	CA	R	218	1/1	0.30	43.48	93,93,93,93	0
2	CA	R	217[B]	1/1	0.31	34.64	51,51,51,51	1
2	CA	Q	216[B]	1/1	0.27	28.75	52,52,52,52	1
2	CA	J	216	1/1	0.17	25.98	69,69,69,69	0
2	CA	B	218[B]	1/1	0.27	19.50	53,53,53,53	1
2	CA	L	218	1/1	0.10	17.67	48,48,48,48	0
2	CA	F	218[B]	1/1	0.34	16.37	57,57,57,57	1
2	CA	O	216	1/1	0.29	14.36	95,95,95,95	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	CA	N	218[B]	1/1	0.13	12.37	42,42,42,42	1
2	CA	S	216[B]	1/1	0.24	10.69	46,46,46,46	1
2	CA	K	216[A]	1/1	0.14	10.43	36,36,36,36	1
2	CA	F	216	1/1	0.10	10.20	52,52,52,52	0
2	CA	E	217[B]	1/1	0.18	9.46	40,40,40,40	1
2	CA	W	215	1/1	0.09	9.40	53,53,53,53	0
2	CA	M	217[B]	1/1	0.19	9.00	41,41,41,41	1
2	CA	C	217	1/1	0.12	9.00	58,58,58,58	0
2	CA	U	215[B]	1/1	0.18	8.70	50,50,50,50	1
2	CA	P	217[A]	1/1	0.17	8.67	49,49,49,49	1
2	CA	H	216[B]	1/1	0.17	6.92	45,45,45,45	1
2	CA	T	215[A]	1/1	0.16	6.59	37,37,37,37	1
2	CA	X	217[A]	1/1	0.14	5.82	35,35,35,35	1
3	ACY	G	219	4/4	0.23	5.77	21,22,50,98	0
2	CA	K	218	1/1	0.12	5.62	52,52,52,52	0
2	CA	D	216[B]	1/1	0.20	5.59	45,45,45,45	1
3	ACY	V	221	4/4	0.20	5.34	18,20,47,104	0
2	CA	J	215[B]	1/1	0.14	5.21	49,49,49,49	1
2	CA	D	214	1/1	0.12	5.17	61,61,61,61	0
2	CA	S	217	1/1	0.11	5.13	56,56,56,56	0
2	CA	V	216[A]	1/1	0.12	5.07	31,31,31,31	1
2	CA	G	216	1/1	0.09	5.00	42,42,42,42	0
2	CA	M	215	1/1	0.07	5.00	44,44,44,44	0
2	CA	J	215[A]	1/1	0.14	4.78	27,27,27,27	1
2	CA	G	218[B]	1/1	0.13	4.64	38,38,38,38	1
2	CA	P	217[B]	1/1	0.17	4.47	29,29,29,29	1
2	CA	B	216	1/1	0.12	4.12	54,54,54,54	0
2	CA	U	216	1/1	0.10	4.11	50,50,50,50	0
2	CA	X	213	1/1	0.15	3.92	52,52,52,52	0
3	ACY	X	241	4/4	0.21	2.60	23,24,35,99	0
2	CA	I	216[A]	1/1	0.09	2.30	25,25,25,25	1
2	CA	I	216[B]	1/1	0.09	2.27	35,35,35,35	1
2	CA	C	216[B]	1/1	0.11	2.26	40,40,40,40	1
2	CA	W	216	1/1	0.10	2.17	46,46,46,46	0
2	CA	O	215	1/1	0.16	2.07	65,65,65,65	0
2	CA	I	213	1/1	0.11	2.07	48,48,48,48	0
2	CA	A	216[A]	1/1	0.12	1.96	28,28,28,28	1
2	CA	D	215[A]	1/1	0.12	1.84	36,36,36,36	1
3	ACY	I	218	4/4	0.20	1.81	20,28,54,110	0
2	CA	C	216[A]	1/1	0.11	1.77	25,25,25,25	1
2	CA	A	216[B]	1/1	0.12	1.49	54,54,54,54	1

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	ACY	R	219	4/4	0.16	1.10	17,27,32,57	0
2	CA	N	213	1/1	0.11	1.03	46,46,46,46	0
2	CA	N	216	1/1	0.07	1.00	38,38,38,38	0
2	CA	B	214	1/1	0.09	0.89	52,52,52,52	0
3	ACY	W	231	4/4	0.17	0.84	18,23,36,48	0
2	CA	C	214	1/1	0.09	0.82	36,36,36,36	0
3	ACY	E	218	4/4	0.12	0.75	19,20,30,51	0
3	ACY	T	218	4/4	0.17	0.73	20,22,59,75	0
2	CA	V	217[B]	1/1	0.10	0.69	33,33,33,33	1
3	ACY	S	218	4/4	0.13	0.66	16,18,60,62	0
3	ACY	P	218	4/4	0.15	0.63	18,21,45,60	0
2	CA	G	215	1/1	0.09	0.59	33,33,33,33	0
3	ACY	Q	217	4/4	0.14	0.47	16,23,31,46	0
2	CA	L	217[B]	1/1	0.08	0.39	34,34,34,34	1
3	ACY	M	218	4/4	0.18	0.26	19,26,34,39	0
2	CA	L	217[A]	1/1	0.08	0.25	23,23,23,23	1
2	CA	G	217[A]	1/1	0.09	0.17	29,29,29,29	1
3	ACY	A	217	4/4	0.12	0.11	21,26,35,48	0
3	ACY	F	219	4/4	0.11	0.09	20,22,39,82	0
2	CA	H	215[A]	1/1	0.10	0.07	27,27,27,27	1
3	ACY	H	218	4/4	0.14	-0.06	19,19,33,72	0
3	ACY	L	219	4/4	0.11	-0.07	18,24,31,37	0
3	ACY	O	218	4/4	0.11	-0.18	16,20,38,63	0
2	CA	T	216[B]	1/1	0.08	-0.26	25,25,25,25	1
2	CA	X	219	1/1	0.07	-0.29	58,58,58,58	0
3	ACY	J	218	4/4	0.09	-0.39	19,24,31,68	0
2	CA	L	215	1/1	0.07	-0.40	26,26,26,26	0
2	CA	Q	215[A]	1/1	0.08	-0.45	23,23,23,23	1
2	CA	F	214	1/1	0.08	-0.72	36,36,36,36	0
2	CA	X	214	1/1	0.08	-0.95	52,52,52,52	0
2	CA	U	214[A]	1/1	0.07	-0.98	24,24,24,24	1
2	CA	L	216	1/1	0.06	-0.99	31,31,31,31	0
2	CA	P	213	1/1	0.07	-1.09	31,31,31,31	0
2	CA	S	215[A]	1/1	0.07	-1.14	27,27,27,27	1
2	CA	R	216[A]	1/1	0.08	-1.18	25,25,25,25	1
2	CA	X	218[B]	1/1	0.07	-1.28	27,27,27,27	1
2	CA	C	213	1/1	0.06	-1.33	17,17,17,17	0
2	CA	M	214	1/1	0.05	-1.43	17,17,17,17	0
2	CA	F	217[A]	1/1	0.07	-1.52	25,25,25,25	1
2	CA	K	213	1/1	0.05	-1.59	26,26,26,26	0
2	CA	E	216[A]	1/1	0.06	-1.68	23,23,23,23	1
2	CA	R	213	1/1	0.05	-1.80	36,36,36,36	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	CA	K	214	1/1	0.06	-1.81	21,21,21,21	0
2	CA	M	216[A]	1/1	0.07	-1.82	27,27,27,27	1
2	CA	N	217[A]	1/1	0.07	-2.00	26,26,26,26	1
2	CA	N	214	1/1	0.05	-2.02	29,29,29,29	0
2	CA	L	214	1/1	0.05	-2.14	21,21,21,21	0
2	CA	B	217[A]	1/1	0.05	-2.18	20,20,20,20	1
2	CA	J	217	1/1	0.06	-2.19	39,39,39,39	0
2	CA	R	215	1/1	0.05	-2.31	18,18,18,18	0
2	CA	E	214	1/1	0.05	-2.38	17,17,17,17	0
2	CA	K	217[B]	1/1	0.07	-2.60	32,32,32,32	1
2	CA	C	215	1/1	0.03	-2.61	27,27,27,27	0
2	CA	U	213	1/1	0.04	-2.78	18,18,18,18	0
2	CA	G	214	1/1	0.04	-2.84	18,18,18,18	0
2	CA	J	214	1/1	0.05	-2.88	19,19,19,19	0
2	CA	P	215	1/1	0.04	-2.94	22,22,22,22	0
2	CA	K	215	1/1	0.04	-3.11	33,33,33,33	0
2	CA	W	214	1/1	0.05	-3.14	20,20,20,20	0
2	CA	A	214	1/1	0.05	-3.20	17,17,17,17	0
2	CA	A	215	1/1	0.06	-3.29	49,49,49,49	0
2	CA	E	213	1/1	0.05	-3.32	28,28,28,28	0
2	CA	Q	214	1/1	0.04	-3.53	18,18,18,18	0
2	CA	D	213	1/1	0.03	-3.71	17,17,17,17	0
2	CA	S	214	1/1	0.05	-3.78	16,16,16,16	0
2	CA	Q	213	1/1	0.03	-3.96	22,22,22,22	0
2	CA	O	213	1/1	0.03	-3.99	19,19,19,19	0
2	CA	B	213	1/1	0.03	-4.08	14,14,14,14	0
2	CA	I	215	1/1	0.03	-4.08	19,19,19,19	0
2	CA	V	214	1/1	0.04	-4.09	18,18,18,18	0
2	CA	T	214	1/1	0.04	-4.34	20,20,20,20	0
2	CA	H	213	1/1	0.03	-4.39	22,22,22,22	0
2	CA	O	214	1/1	0.03	-4.92	14,14,14,14	0
2	CA	S	213	1/1	0.04	-5.16	26,26,26,26	0
2	CA	F	213	1/1	0.03	-5.54	22,22,22,22	0
2	CA	X	216	1/1	0.04	-5.68	18,18,18,18	0
2	CA	X	215	1/1	0.02	-5.77	21,21,21,21	0
2	CA	W	213	1/1	0.04	-5.84	25,25,25,25	0
2	CA	E	215	1/1	0.03	-5.88	19,19,19,19	0
2	CA	A	213	1/1	0.03	-5.90	20,20,20,20	0
2	CA	L	213	1/1	0.04	-6.26	21,21,21,21	0
2	CA	H	214	1/1	0.04	-6.44	16,16,16,16	0
2	CA	F	215	1/1	0.03	-6.48	21,21,21,21	0
2	CA	R	214	1/1	0.04	-6.48	22,22,22,22	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	CA	I	217	1/1	0.05	-6.71	48,48,48,48	0
2	CA	V	213	1/1	0.03	-7.00	25,25,25,25	0
2	CA	M	213	1/1	0.03	-7.49	23,23,23,23	0
2	CA	I	214	1/1	0.03	-9.10	25,25,25,25	0
2	CA	P	214	1/1	0.03	-9.35	23,23,23,23	0
2	CA	G	213	1/1	0.03	-9.81	27,27,27,27	0
2	CA	J	213	1/1	0.03	-10.81	23,23,23,23	0
2	CA	N	215	1/1	0.04	-15.00	19,19,19,19	0
2	CA	B	215	1/1	0.04	-29.39	24,24,24,24	0
2	CA	T	213	1/1	0.03	-72.17	25,25,25,25	0
2	CA	P	216	1/1	0.26	-	77,77,77,77	0
2	CA	V	215	1/1	0.12	-	71,71,71,71	0
2	CA	O	217	1/1	0.12	-	46,46,46,46	0
2	CA	H	217	1/1	0.15	-	66,66,66,66	0

6.5 Other polymers ⓘ

There are no such residues in this entry.