



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 07:09 AM UTC

PDB ID : 4WZ7 / pdb_00004wz7
Title : Crystal structure of mitochondrial NADH:ubiquinone oxidoreductase from Yarrowia lipolytica.
Authors : Wirth, C.; Zickermann, V.; Brandt, U.; Hunte, C.
Deposited on : 2014-11-18
Resolution : 3.60 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

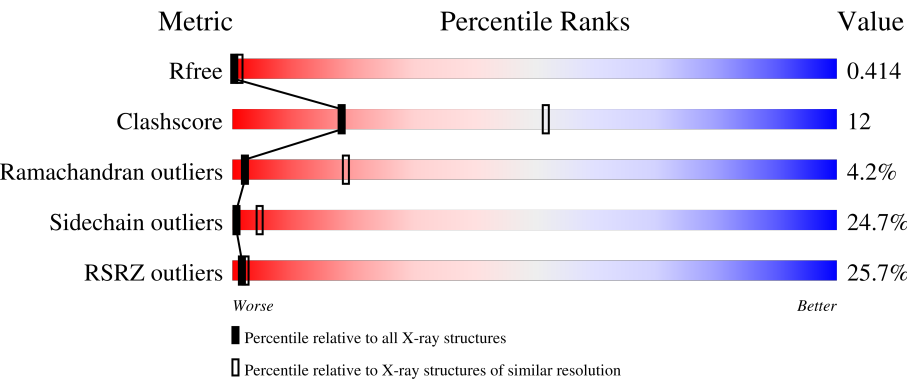
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.60 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	1747 (3.70-3.50)
Clashscore	190562	1827 (3.70-3.50)
Ramachandran outliers	187476	1773 (3.70-3.50)
Sidechain outliers	187428	1772 (3.70-3.50)
RSRZ outliers	180081	1745 (3.70-3.50)

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	1	327	9% 60%25%10%.
2	2	438	6% 54%32%12%.
3	3	89	11% 40%38%17%.
4	4	470	5% 63%27%9%.
5	5	619	13% 67%26%7%.

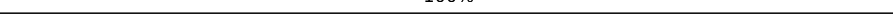
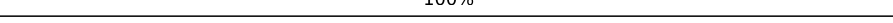
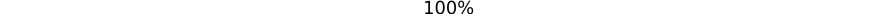
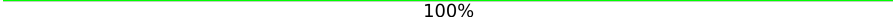
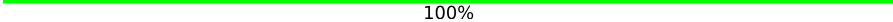
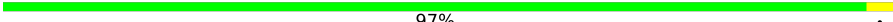
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Mol	Chain	Length	Quality of chain
6	6	185	
7	A	628	
8	B	370	
9	C	444	
10	E	195	
11	G	133	
12	H	154	
13	I	137	
14	K	183	
15	L	89	
16	D	57	
16	Z	57	
17	F	54	
18	J	63	
19	M	29	
20	N	50	
21	O	70	
22	P	46	
23	Q	51	
24	R	30	
25	S	69	
26	AH	15	
26	T	15	
27	U	26	
28	V	22	

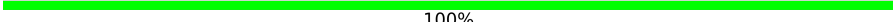
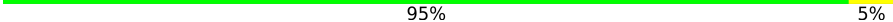
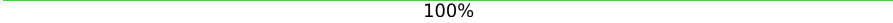
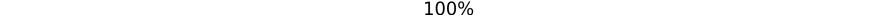
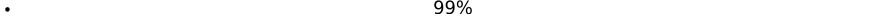
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Mol	Chain	Length	Quality of chain
29	AB	9	 100%
29	AY	9	 100%
29	BE	9	 100%
29	W	9	 78% 22%
30	AJ	16	 100%
30	AV	16	 100%
30	BH	16	 88% 12%
30	X	16	 100%
31	AR	13	 100%
31	AT	13	 100%
31	Y	13	 100%
32	AA	18	 100%
32	AW	18	 100%
32	BB	18	 100%
32	BG	18	 100%
33	AC	47	 85% 15%
33	AD	47	 100%
34	AE	48	 92% 8%
35	AF	35	 100%
36	AG	25	 100%
37	AI	36	 100%
37	AL	36	 100%
38	AK	76	 100%
38	AN	76	 97% .
39	AM	17	 100%

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Mol	Chain	Length	Quality of chain
40	AO	32	 100%
41	AP	11	 100%
41	AS	11	 100%
42	AQ	8	 100%
42	BA	8	 100%
43	AU	58	 100%
44	AX	39	 100%
45	AZ	40	 95% 5%
46	BC	20	 90% 10%
47	BD	19	 100%
47	BF	19	 100%
48	BI	905	 99%

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
50	SF4	K	500	-	-	X	-

2 Entry composition

There are 50 unique types of molecules in this entry. The entry contains 35169 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 NADH-ubiquinone oxidoreductase chain 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	1	318	Total	C	N	O	S	0	0	0
			2177	1431	352	389	5			

- Molecule 2 is a protein called NADH dehydrogenase subunit 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	2	438	Total	C	N	O	S	0	0	0
			3142	2092	482	556	12			

- Molecule 3 is a protein called NADH-ubiquinone oxidoreductase chain 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	3	89	Total	C	N	O	S	0	0	0
			641	444	91	104	2			

- Molecule 4 is a protein called NADH-ubiquinone oxidoreductase chain 4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	4	470	Total	C	N	O	S	0	0	0
			3017	1952	507	546	12			

- Molecule 5 is a protein called NADH-ubiquinone oxidoreductase chain 5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	5	619	Total	C	N	O	S	0	0	0
			4065	2636	677	727	25			

- Molecule 6 is a protein called NADH-ubiquinone oxidoreductase chain 6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	6	149	Total	C	N	O	S	0	0	0
			1078	735	156	180	7			

- Molecule 7 is a protein called NUAM protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	A	628	Total	C	N	O	S	0	0	0
			3226	1933	639	639	15			

- Molecule 8 is a protein called NUBM protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	B	370	Total	C	N	O	S	0	0	0
			2013	1214	391	399	9			

- Molecule 9 is a protein called NUCM protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
9	C	366	Total	C	N	O	S	0	0	0
			2647	1675	455	496	21			

- Molecule 10 is a protein called 39-kDa subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	E	195	Total	C	N	O	S	0	0	0
			975	585	195	195				

- Molecule 11 is a protein called NUGM protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
11	G	133	Total	C	N	O	S	0	0	0
			880	558	154	164	4			

- Molecule 12 is a protein called Subunit NUHM of protein NADH:Ubiquinone Oxidoreductase (Complex I).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
12	H	154	Total	C	N	O	S	0	0	0
			803	476	156	164	7			

- Molecule 13 is a protein called Subunit NUIM of protein NADH:Ubiquinone Oxidoreductase (Complex I).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
13	I	137	Total	C	N	O	S	0	0	0
			857	533	145	169	10			

- Molecule 14 is a protein called Subunit NUKM of protein NADH:Ubiquinone Oxidoreductase (Complex I).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
14	K	147	Total	C	N	O	S	0	0	0
			1023	642	182	187	12			

- Molecule 15 is a protein called NADH-ubiquinone oxidoreductase chain 4L.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
15	L	89	Total	C	N	O	S	0	0	0
			660	437	108	112	3			

- Molecule 16 is a protein called unknown subunits 1.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
16	Z	57	Total	C	N	O	0	0	0
			285	171	57	57			
16	D	57	Total	C	N	O	0	0	0
			285	171	57	57			

- Molecule 17 is a protein called unknown subunits 2.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
17	F	54	Total	C	N	O	0	0	0
			270	162	54	54			

- Molecule 18 is a protein called unknown subunits.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
18	J	63	Total	C	N	O	0	0	0
			315	189	63	63			

- Molecule 19 is a protein called unknown subunits.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
19	M	29	Total	C	N	O	0	0	0
			145	87	29	29			

- Molecule 20 is a protein called unknown subunits.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
20	N	50	Total	C	N	O	0	0	0
			250	150	50	50			

- Molecule 21 is a protein called unknown subunits.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
21	O	70	Total	C	N	O	0	0	0
			350	210	70	70			

- Molecule 22 is a protein called unknown subunits.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
22	P	46	Total	C	N	O	0	0	0
			230	138	46	46			

- Molecule 23 is a protein called unknown subunits.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
23	Q	51	Total	C	N	O	0	0	0
			255	153	51	51			

- Molecule 24 is a protein called unknown subunits.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
24	R	30	Total	C	N	O	0	0	0
			150	90	30	30			

- Molecule 25 is a protein called unknown subunits.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
25	S	69	Total	C	N	O	0	0	0
			345	207	69	69			

- Molecule 26 is a protein called unknown subunits.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
26	T	15	Total	C	N	O	0	0	0
			75	45	15	15			
26	AH	15	Total	C	N	O	0	0	0
			75	45	15	15			

- Molecule 27 is a protein called unknown subunits.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
27	U	26	Total	C	N	O	0	0	0
			130	78	26	26			

- Molecule 28 is a protein called unknown subunits.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
28	V	22	Total	C	N	O	0	0	0
			110	66	22	22			

- Molecule 29 is a protein called unknown subunits.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
29	W	9	Total	C	N	O	0	0	0
			45	27	9	9			
29	AB	9	Total	C	N	O	0	0	0
			45	27	9	9			
29	AY	9	Total	C	N	O	0	0	0
			45	27	9	9			
29	BE	9	Total	C	N	O	0	0	0
			45	27	9	9			

- Molecule 30 is a protein called unknown subunits.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
30	X	16	Total	C	N	O	0	0	0
			80	48	16	16			
30	AJ	16	Total	C	N	O	0	0	0
			80	48	16	16			
30	AV	16	Total	C	N	O	0	0	0
			80	48	16	16			
30	BH	16	Total	C	N	O	0	0	0
			80	48	16	16			

- Molecule 31 is a protein called unknown subunits.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
31	Y	13	Total	C	N	O	0	0	0
			65	39	13	13			
31	AR	13	Total	C	N	O	0	0	0
			65	39	13	13			
31	AT	13	Total	C	N	O	0	0	0
			65	39	13	13			

- Molecule 32 is a protein called unknown subunits.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
32	AA	18	Total	C	N	O	0	0	0
			90	54	18	18			
32	AW	18	Total	C	N	O	0	0	0
			90	54	18	18			
32	BB	18	Total	C	N	O	0	0	0
			90	54	18	18			
32	BG	18	Total	C	N	O	0	0	0
			90	54	18	18			

- Molecule 33 is a protein called unknown subunits.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
33	AC	47	Total	C	N	O	0	0	0
			235	141	47	47			
33	AD	47	Total	C	N	O	0	0	0
			235	141	47	47			

- Molecule 34 is a protein called unknown subunits.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
34	AE	48	Total	C	N	O	0	0	0
			240	144	48	48			

- Molecule 35 is a protein called unknown subunits.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
35	AF	35	Total	C	N	O	0	0	0
			175	105	35	35			

- Molecule 36 is a protein called unknown subunits.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
36	AG	25	Total	C	N	O	0	0	0
			125	75	25	25			

- Molecule 37 is a protein called unknown subunits.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
37	AI	36	Total	C	N	O	0	0	0
			180	108	36	36			

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
37	AL	36	Total	C	N	O	0	0	0
			180	108	36	36			

- Molecule 38 is a protein called unknown subunits.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
38	AK	76	Total	C	N	O	0	0	0
			380	228	76	76			
38	AN	76	Total	C	N	O	0	0	0
			380	228	76	76			

- Molecule 39 is a protein called unknown subunits.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
39	AM	17	Total	C	N	O	0	0	0
			85	51	17	17			

- Molecule 40 is a protein called unknown subunits.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
40	AO	32	Total	C	N	O	0	0	0
			160	96	32	32			

- Molecule 41 is a protein called unknown subunits.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
41	AP	11	Total	C	N	O	0	0	0
			55	33	11	11			
41	AS	11	Total	C	N	O	0	0	0
			54	32	11	11			

- Molecule 42 is a protein called unknown subunits.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
42	AQ	8	Total	C	N	O	0	0	0
			40	24	8	8			
42	BA	8	Total	C	N	O	0	0	0
			40	24	8	8			

- Molecule 43 is a protein called unknown subunits.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
43	AU	58	Total	C	N	O	0	0	0
			290	174	58	58			

- Molecule 44 is a protein called unknown subunits.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
44	AX	39	Total	C	N	O	0	0	0
			195	117	39	39			

- Molecule 45 is a protein called unknown subunits.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
45	AZ	40	Total	C	N	O	0	0	0
			200	120	40	40			

- Molecule 46 is a protein called unknown subunits.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
46	BC	20	Total	C	N	O	0	0	0
			100	60	20	20			

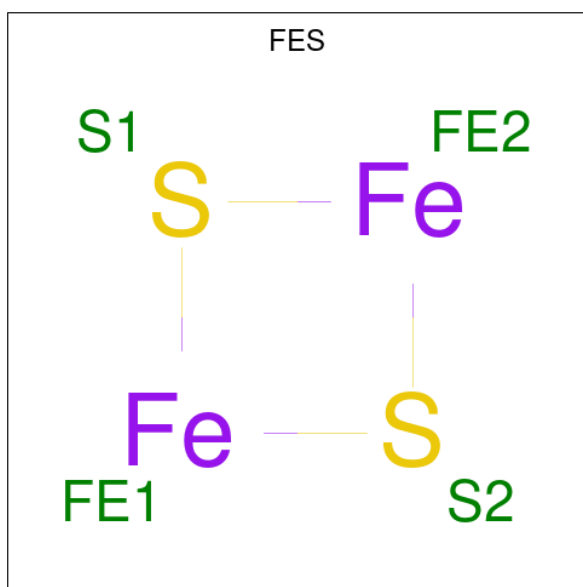
- Molecule 47 is a protein called unknown subunits.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
47	BD	19	Total	C	N	O	0	0	0
			95	57	19	19			
47	BF	19	Total	C	N	O	0	0	0
			95	57	19	19			

- Molecule 48 is a protein called unknown subunits.

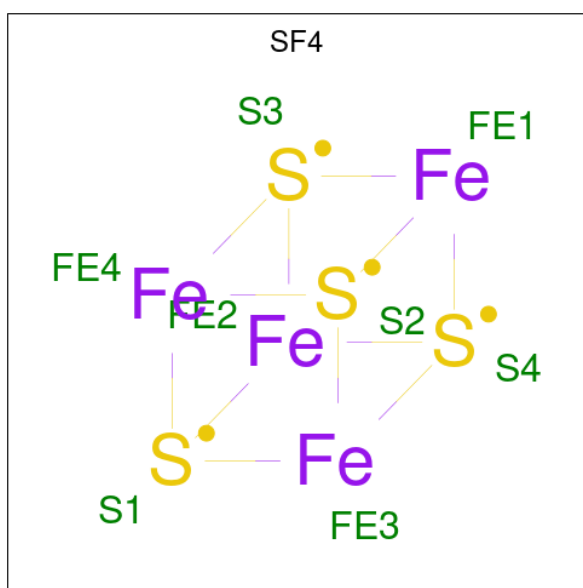
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
48	BI	9	Total	C	N	O	0	0	0
			45	27	9	9			

- Molecule 49 is FE2/S2 (INORGANIC) CLUSTER (CCD ID: FES) (formula: Fe₂S₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
49	A	1	Total	Fe	S	0	0
			4	2	2		
49	H	1	Total	Fe	S	0	0
			4	2	2		

- Molecule 50 is IRON/SULFUR CLUSTER (CCD ID: SF4) (formula: Fe_4S_4).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
50	A	1	Total	Fe	S	0	0
			8	4	4		
50	A	1	Total	Fe	S	0	0
			8	4	4		

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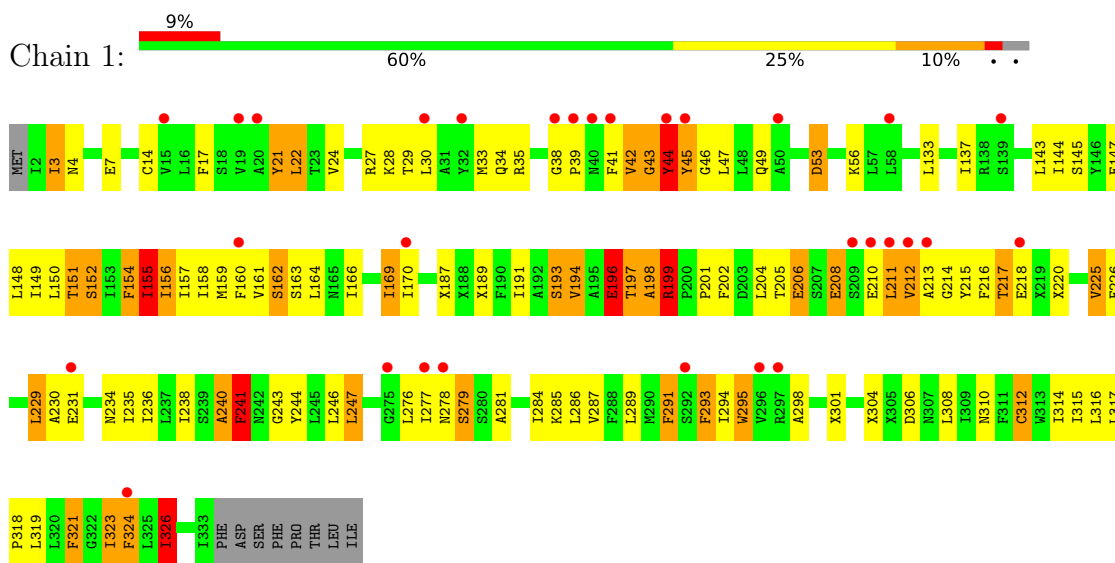
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
50	B	1	Total	Fe	S	0	0
			8	4	4		
50	I	1	Total	Fe	S	0	0
			8	4	4		
50	I	1	Total	Fe	S	0	0
			8	4	4		
50	K	1	Total	Fe	S	0	0
			8	4	4		

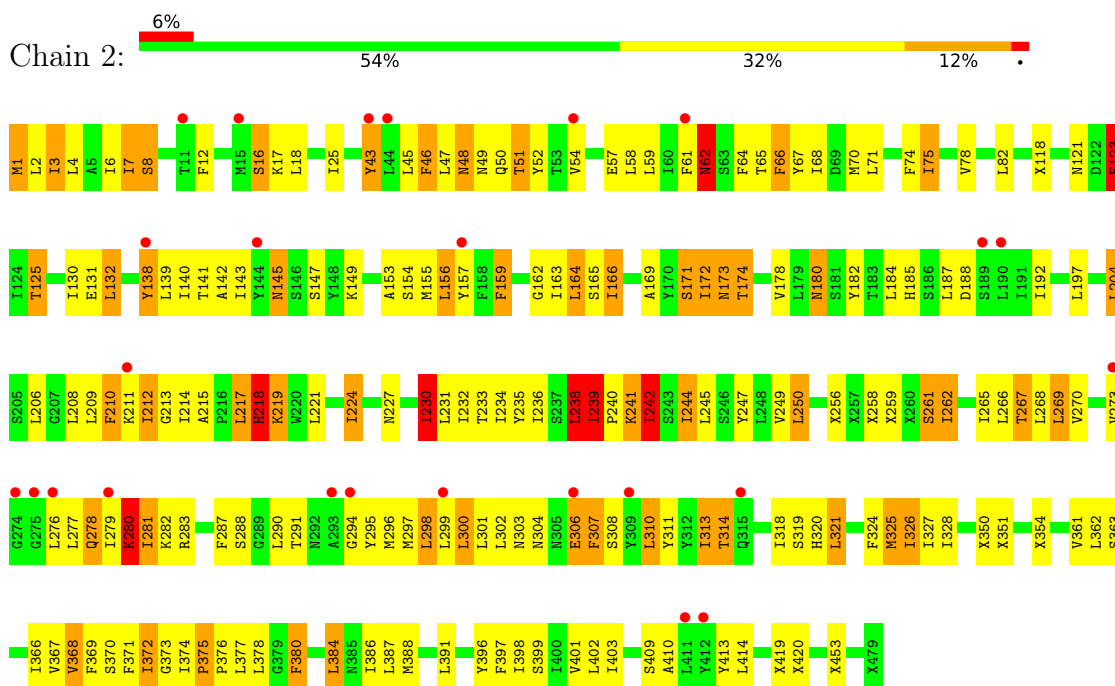
3 Residue-property plots

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.

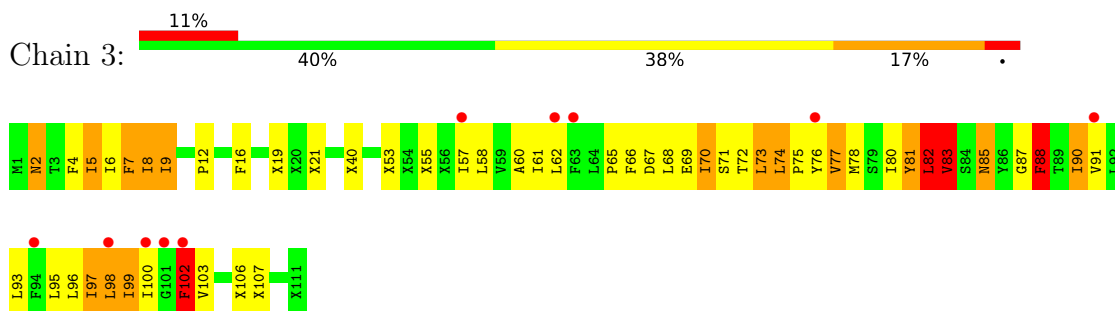
• Molecule 1: NADH-ubiquinone oxidoreductase chain 1



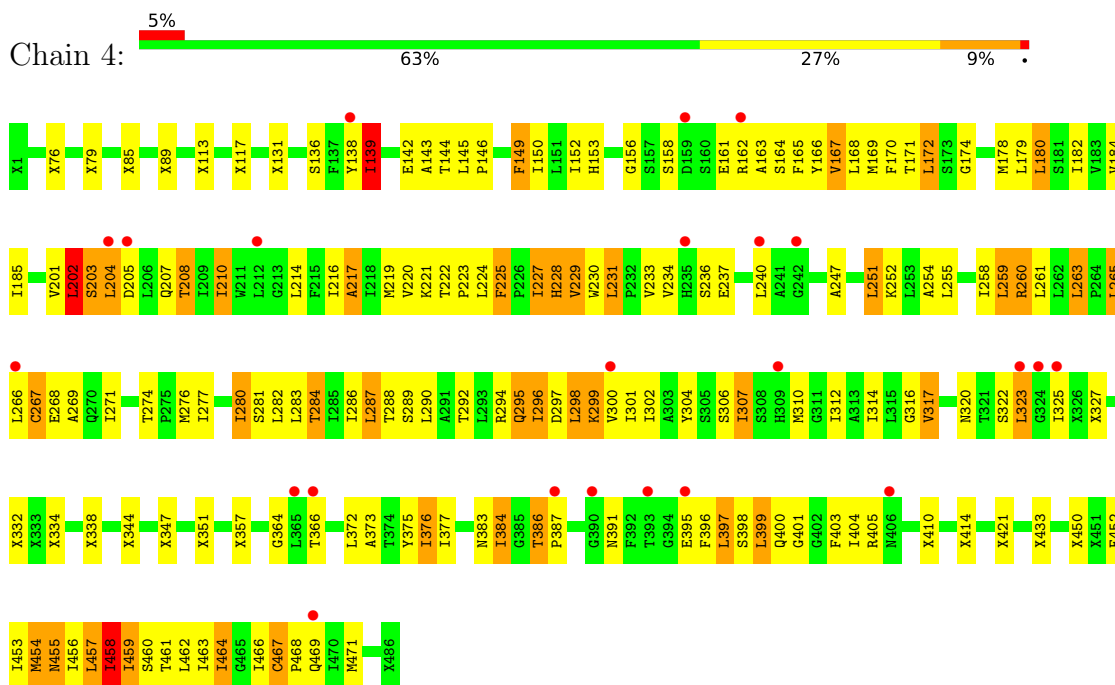
• Molecule 2: NADH dehydrogenase subunit 2



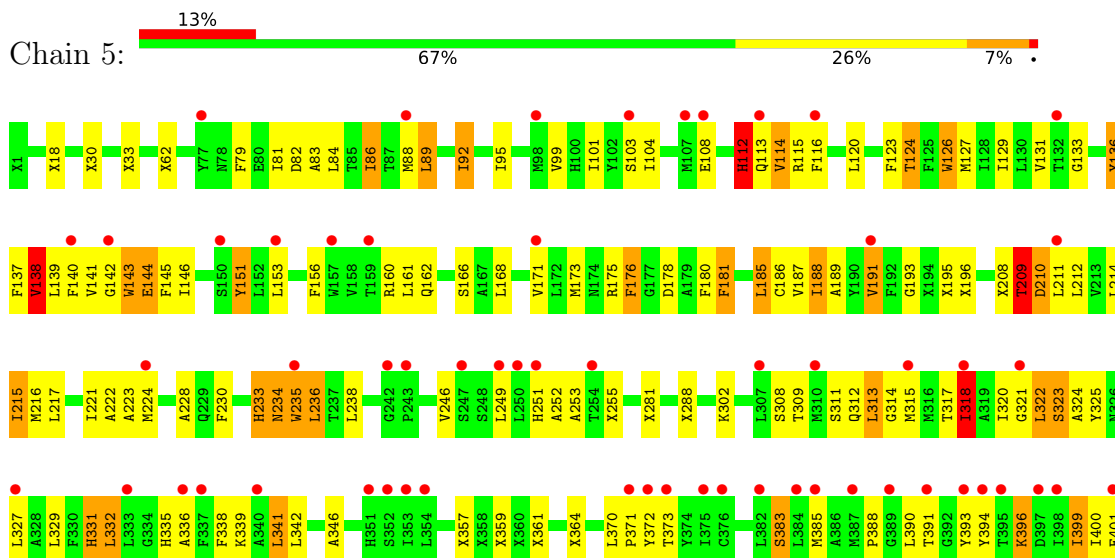
- Molecule 3: NADH-ubiquinone oxidoreductase chain 3

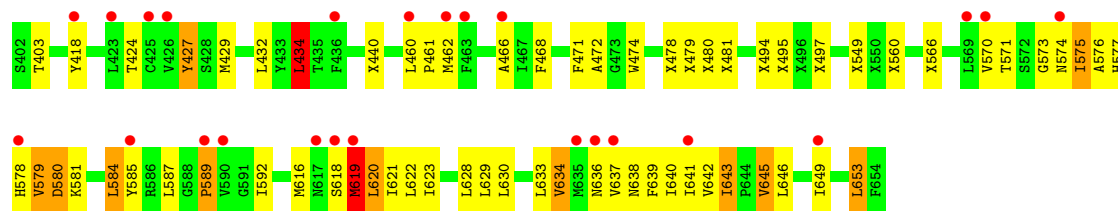


- Molecule 4: NADH-ubiquinone oxidoreductase chain 4

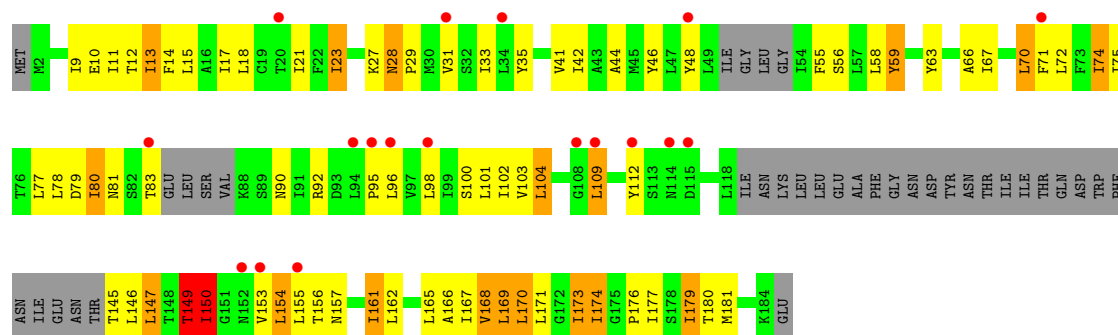


- Molecule 5: NADH-ubiquinone oxidoreductase chain 5

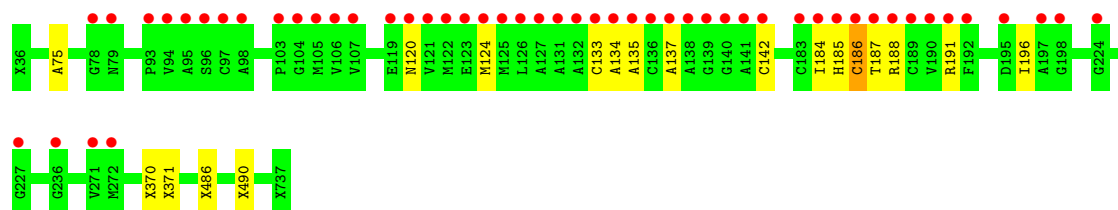




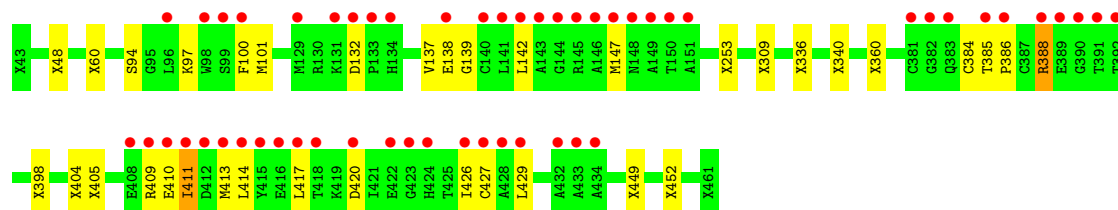
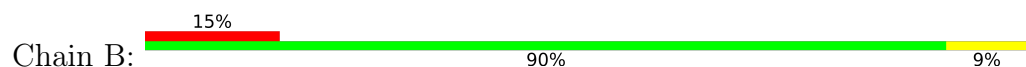
• Molecule 6: NADH-ubiquinone oxidoreductase chain 6



• Molecule 7: NUAM protein

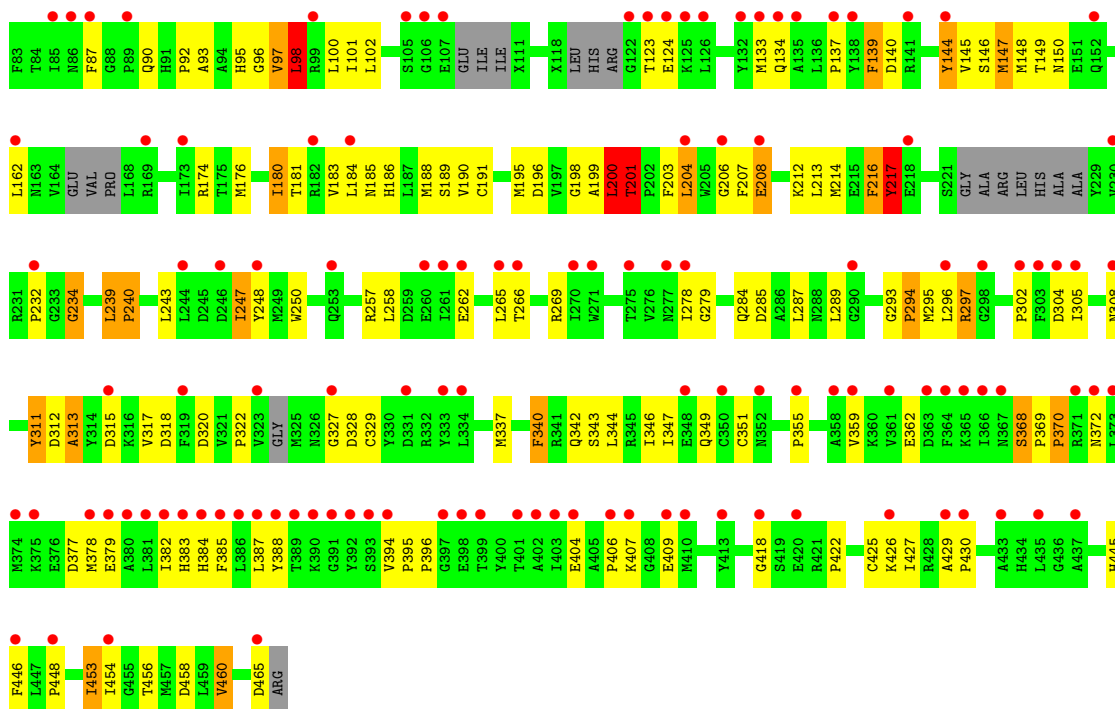


• Molecule 8: NUBM protein

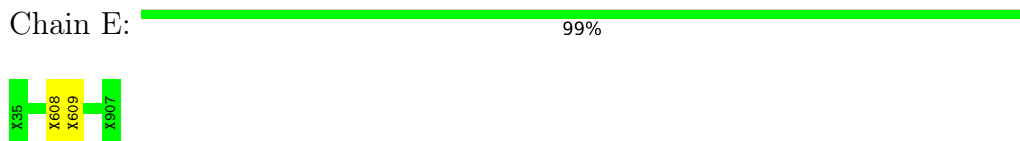


• Molecule 9: NUCM protein

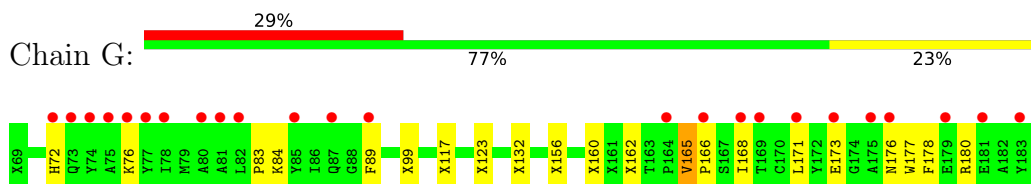




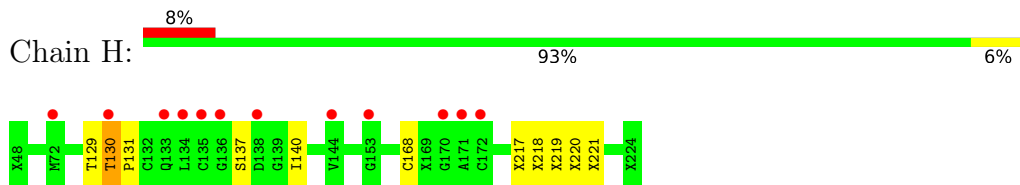
- Molecule 10: 39-kDa subunit



- Molecule 11: NUGM protein

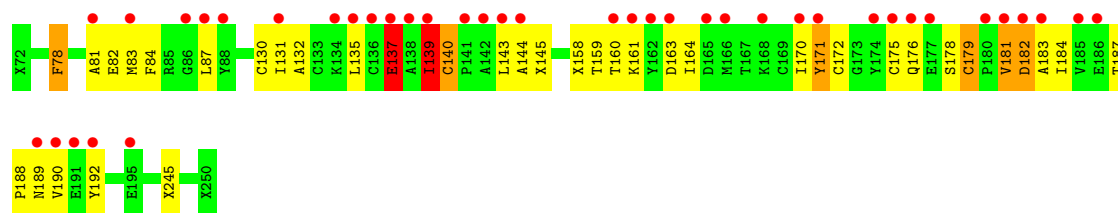


- Molecule 12: Subunit NUHM of protein NADH:Ubiquinone Oxidoreductase (Complex I)

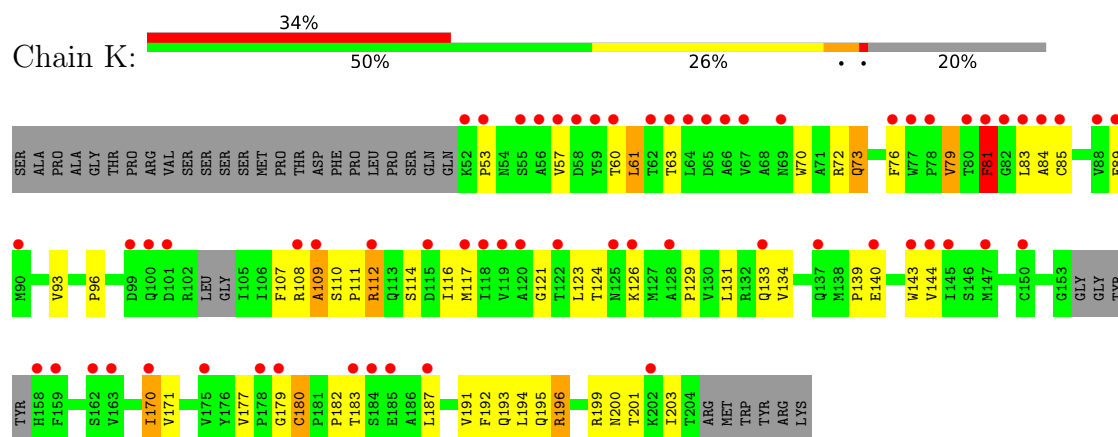


- Molecule 13: Subunit NUIM of protein NADH:Ubiquinone Oxidoreductase (Complex I)

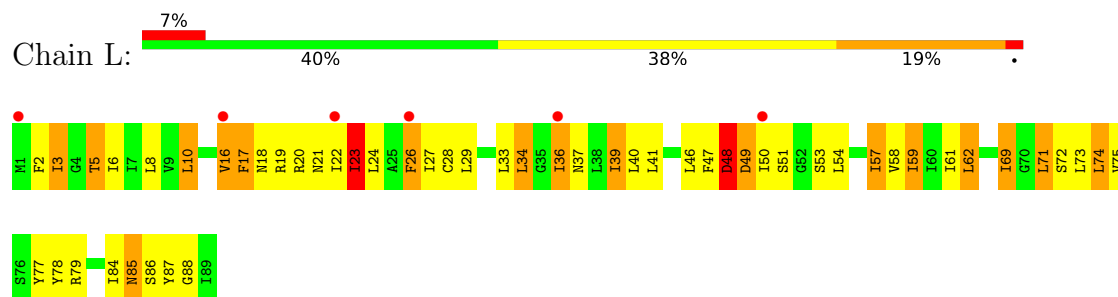




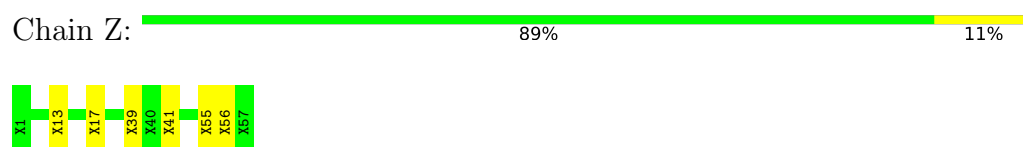
- Molecule 14: Subunit NUKM of protein NADH:Ubiquinone Oxidoreductase (Complex I)



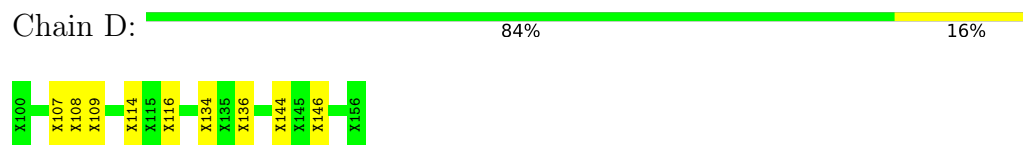
- Molecule 15: NADH-ubiquinone oxidoreductase chain 4L



- Molecule 16: unknown subunits 1



- Molecule 16: unknown subunits 1



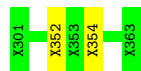
- Molecule 17: unknown subunits 2



There are no outlier residues recorded for this chain.

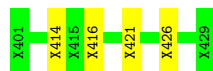
- Molecule 18: unknown subunits

Chain J:  97% .



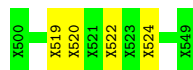
- Molecule 19: unknown subunits

Chain M:  86% 14%



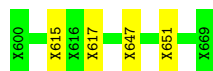
- Molecule 20: unknown subunits

Chain N:  92% 8%



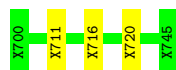
- Molecule 21: unknown subunits

Chain O:  94% 6%



- Molecule 22: unknown subunits

Chain P:  93% 7%



- Molecule 23: unknown subunits

Chain Q:  98% .



- Molecule 24: unknown subunits

Chain R:  100%

There are no outlier residues recorded for this chain.

- Molecule 25: unknown subunits

Chain S:  94% 6%



- Molecule 26: unknown subunits

Chain T: 100%

There are no outlier residues recorded for this chain.

- Molecule 26: unknown subunits

Chain AH: 100%

There are no outlier residues recorded for this chain.

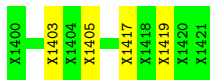
- Molecule 27: unknown subunits

Chain U: 100%

There are no outlier residues recorded for this chain.

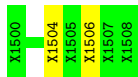
- Molecule 28: unknown subunits

Chain V: 82% 18%



- Molecule 29: unknown subunits

Chain W: 78% 22%



- Molecule 29: unknown subunits

Chain AB: 100%

There are no outlier residues recorded for this chain.

- Molecule 29: unknown subunits

Chain AY: 100%

There are no outlier residues recorded for this chain.

- Molecule 29: unknown subunits

Chain BE: 100%

There are no outlier residues recorded for this chain.

- Molecule 30: unknown subunits

Chain X:  100%

There are no outlier residues recorded for this chain.

- Molecule 30: unknown subunits

Chain AJ:  100%

There are no outlier residues recorded for this chain.

- Molecule 30: unknown subunits

Chain AV:  100%

There are no outlier residues recorded for this chain.

- Molecule 30: unknown subunits

Chain BH:  88% 12%



- Molecule 31: unknown subunits

Chain Y:  100%

There are no outlier residues recorded for this chain.

- Molecule 31: unknown subunits

Chain AR:  100%

There are no outlier residues recorded for this chain.

- Molecule 31: unknown subunits

Chain AT:  100%

There are no outlier residues recorded for this chain.

- Molecule 32: unknown subunits

Chain AA:  100%

There are no outlier residues recorded for this chain.

- Molecule 32: unknown subunits

Chain AW:  100%

There are no outlier residues recorded for this chain.

- Molecule 32: unknown subunits

Chain BB:  100%

There are no outlier residues recorded for this chain.

- Molecule 32: unknown subunits

Chain BG:  100%

There are no outlier residues recorded for this chain.

- Molecule 33: unknown subunits

Chain AC:  85% 15%



- Molecule 33: unknown subunits

Chain AD:  100%

There are no outlier residues recorded for this chain.

- Molecule 34: unknown subunits

Chain AE:  92% 8%



- Molecule 35: unknown subunits

Chain AF:  100%

There are no outlier residues recorded for this chain.

- Molecule 36: unknown subunits

Chain AG:  100%

There are no outlier residues recorded for this chain.

- Molecule 37: unknown subunits

Chain AI:  100%

There are no outlier residues recorded for this chain.

- Molecule 37: unknown subunits

Chain AL:  100%

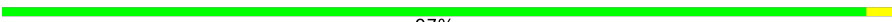
There are no outlier residues recorded for this chain.

- Molecule 38: unknown subunits

Chain AK:  100%

There are no outlier residues recorded for this chain.

- Molecule 38: unknown subunits

Chain AN:  97%



- Molecule 39: unknown subunits

Chain AM:  100%

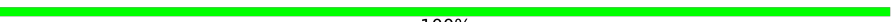
There are no outlier residues recorded for this chain.

- Molecule 40: unknown subunits

Chain AO:  100%

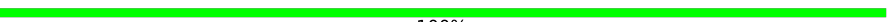
There are no outlier residues recorded for this chain.

- Molecule 41: unknown subunits

Chain AP:  100%

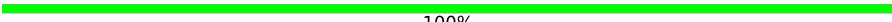
There are no outlier residues recorded for this chain.

- Molecule 41: unknown subunits

Chain AS:  100%

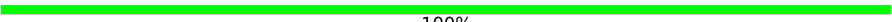
There are no outlier residues recorded for this chain.

- Molecule 42: unknown subunits

Chain AQ:  100%

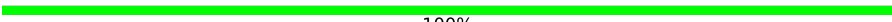
There are no outlier residues recorded for this chain.

- Molecule 42: unknown subunits

Chain BA:  100%

There are no outlier residues recorded for this chain.

- Molecule 43: unknown subunits

Chain AU:  100%

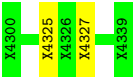
There are no outlier residues recorded for this chain.

- Molecule 44: unknown subunits

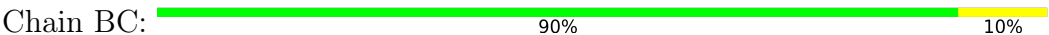


There are no outlier residues recorded for this chain.

- Molecule 45: unknown subunits



- Molecule 46: unknown subunits



- Molecule 47: unknown subunits



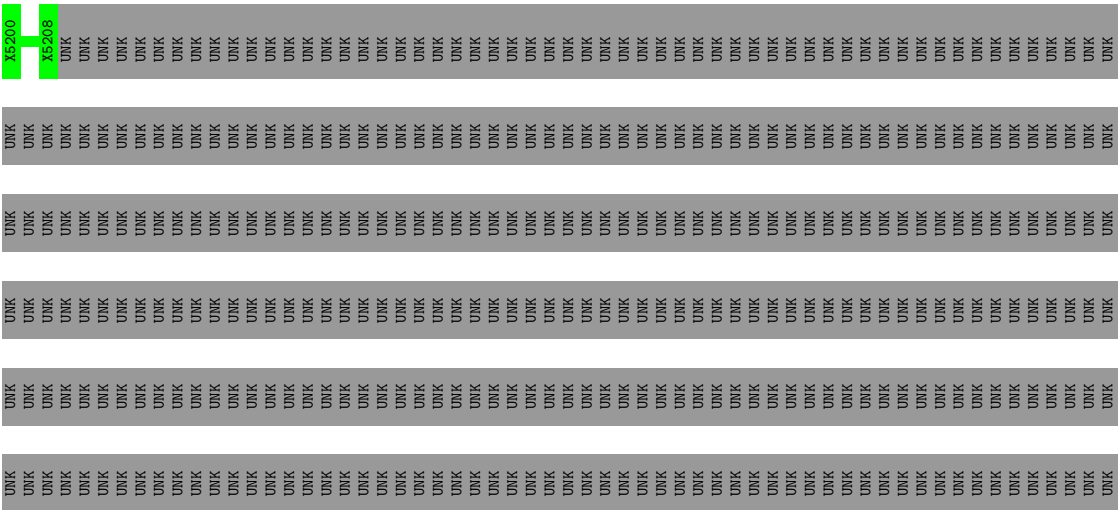
There are no outlier residues recorded for this chain.

- Molecule 47: unknown subunits



There are no outlier residues recorded for this chain.

- Molecule 48: unknown subunits



[illegible]

4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	317.74Å 317.74Å 818.97Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	25.00 – 3.60 25.00 – 3.60	Depositor EDS
% Data completeness (in resolution range)	84.8 (25.00-3.60) 84.6 (25.00-3.60)	Depositor EDS
R_{merge}	0.52	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.01 (at 3.57Å)	Xtriage
Refinement program	BUSTER 2.10.0	Depositor
R, R_{free}	0.316 , 0.341 0.379 , 0.414	Depositor DCC
R_{free} test set	2191 reflections (1.20%)	wwPDB-VP
Wilson B-factor (Å ²)	106.8	Xtriage
Anisotropy	0.020	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.24 , 345.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.41$, $\langle L^2 \rangle = 0.24$	Xtriage
Estimated twinning fraction	0.039 for $-1/3^*h+1/3^*k+1/3^*l,-k,8/3^*h+4/3^*k+1/3^*l$ 0.097 for $-2/3^*h-1/3^*k-1/3^*l,-1/3^*h-2/3^*k+1/3^*l,-4/3^*h+4/3^*k+1/3^*l$ 0.067 for $-h,1/3^*h-1/3^*k-1/3^*l,-4/3^*h-8/3^*k+1/3^*l$	Xtriage
F_o, F_c correlation	0.78	EDS
Total number of atoms	35169	wwPDB-VP
Average B, all atoms (Å ²)	109.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.62% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: FES, SF4

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	1	0.81	0/1638	1.63	30/2224 (1.3%)
2	2	0.91	1/2613 (0.0%)	1.64	57/3550 (1.6%)
3	3	0.86	1/533 (0.2%)	1.70	11/728 (1.5%)
4	4	0.91	0/1920	1.62	26/2615 (1.0%)
5	5	0.88	2/2813 (0.1%)	1.58	37/3820 (1.0%)
6	6	0.92	1/1090 (0.1%)	1.59	7/1491 (0.5%)
7	A	0.82	0/412	1.37	3/531 (0.6%)
8	B	0.77	1/527 (0.2%)	1.43	2/701 (0.3%)
9	C	0.79	0/2653	1.56	33/3592 (0.9%)
11	G	0.68	0/566	1.36	7/766 (0.9%)
12	H	0.76	0/178	1.55	0/220
13	I	0.70	0/499	1.43	2/675 (0.3%)
14	K	0.84	0/1042	1.57	14/1424 (1.0%)
15	L	0.95	0/666	1.69	12/902 (1.3%)
All	All	0.85	6/17150 (0.0%)	1.59	241/23239 (1.0%)

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	5	643	ILE	CA-CB	7.50	1.57	1.54
6	6	66	ALA	CA-C	-6.53	1.47	1.52
2	2	239	ILE	CA-C	5.55	1.58	1.52
8	B	411	ILE	CG1-CD1	-5.35	1.30	1.51
3	3	82	LEU	C-N	5.16	1.40	1.33
5	5	92	ILE	CG1-CD1	5.03	1.71	1.51

All (241) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	C	395	PRO	N-CA-CB	13.85	110.42	103.22

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	1	241	PHE	CA-CB-CG	12.97	126.77	113.80
4	4	384	ILE	N-CA-C	-11.88	101.08	112.96
2	2	46	PHE	CA-CB-CG	11.60	125.40	113.80
3	3	82	LEU	CA-C-N	10.66	140.88	121.70
3	3	82	LEU	C-N-CA	10.66	140.88	121.70
1	1	3	ILE	N-CA-C	-10.62	102.68	111.81
6	6	10	GLU	N-CA-C	-10.27	100.75	113.18
1	1	154	PHE	CA-CB-CG	9.91	123.71	113.80
3	3	88	PHE	CA-CB-CG	9.86	123.66	113.80
1	1	17	PHE	CA-CB-CG	9.70	123.50	113.80
5	5	210	ASP	CA-CB-CG	8.65	121.25	112.60
5	5	643	ILE	N-CA-CB	8.48	116.20	110.52
3	3	102	PHE	CA-CB-CG	8.16	121.96	113.80
13	I	78	PHE	CA-CB-CG	8.14	121.94	113.80
2	2	380	PHE	CA-CB-CG	8.12	121.92	113.80
3	3	9	ILE	N-CA-C	-8.08	105.33	112.12
2	2	375	PRO	N-CA-C	8.00	120.45	110.70
9	C	422	PRO	N-CA-CB	7.91	109.93	103.36
1	1	43	GLY	CA-C-N	7.89	135.90	121.70
1	1	43	GLY	C-N-CA	7.89	135.90	121.70
2	2	238	LEU	N-CA-C	7.74	119.62	111.03
5	5	176	PHE	CA-CB-CG	7.62	121.42	113.80
6	6	48	TYR	N-CA-C	-7.47	103.75	113.17
15	L	26	PHE	CA-CB-CG	7.43	121.23	113.80
15	L	77	TYR	N-CA-C	-7.34	101.38	110.41
9	C	240	PRO	N-CA-CB	7.33	110.94	103.25
9	C	265	LEU	N-CA-C	7.27	119.29	111.36
9	C	315	ASP	CA-CB-CG	7.27	119.87	112.60
14	K	61	LEU	CA-C-N	7.24	134.73	121.70
14	K	61	LEU	C-N-CA	7.24	134.73	121.70
2	2	306	GLU	CA-C-N	7.23	130.31	120.54
2	2	306	GLU	C-N-CA	7.23	130.31	120.54
15	L	59	ILE	N-CA-CB	7.22	119.50	110.47
4	4	217	ALA	CA-C-N	7.15	129.83	120.60
4	4	217	ALA	C-N-CA	7.15	129.83	120.60
9	C	340	PHE	CA-CB-CG	7.13	120.93	113.80
11	G	165	VAL	N-CA-C	7.12	116.72	108.96
1	1	324	PHE	CA-CB-CG	7.11	120.91	113.80
5	5	338	PHE	CA-CB-CG	7.09	120.89	113.80
5	5	311	SER	CA-C-N	7.04	129.71	120.28
5	5	311	SER	C-N-CA	7.04	129.71	120.28
14	K	53	PRO	N-CA-CB	7.04	109.87	103.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	2	413	TYR	N-CA-C	-6.98	103.75	112.90
6	6	14	PHE	CA-CB-CG	6.96	120.76	113.80
5	5	142	GLY	CA-C-N	6.95	133.30	121.14
5	5	142	GLY	C-N-CA	6.95	133.30	121.14
2	2	267	THR	N-CA-CB	6.92	120.89	110.22
3	3	73	LEU	N-CA-C	-6.89	102.87	112.45
2	2	326	ILE	N-CA-C	-6.89	103.59	110.62
5	5	646	LEU	N-CA-C	-6.87	104.86	113.18
2	2	54	VAL	N-CA-C	6.83	117.68	107.51
9	C	195	MET	CA-C-N	6.82	129.75	120.54
9	C	195	MET	C-N-CA	6.82	129.75	120.54
2	2	48	ASN	N-CA-C	6.81	120.89	111.90
1	1	4	ASN	CA-CB-CG	6.79	119.39	112.60
2	2	409	SER	CA-C-N	6.77	129.65	120.38
2	2	409	SER	C-N-CA	6.77	129.65	120.38
5	5	318	ILE	N-CA-CB	6.74	119.71	110.54
1	1	326	ILE	N-CA-CB	6.73	122.34	111.23
15	L	16	VAL	CA-C-N	6.70	133.76	121.70
15	L	16	VAL	C-N-CA	6.70	133.76	121.70
9	C	396	PRO	N-CA-CB	6.68	110.26	103.25
11	G	166	PRO	N-CA-C	6.64	121.50	111.14
9	C	239	LEU	CA-C-N	6.54	128.02	119.84
9	C	239	LEU	C-N-CA	6.54	128.02	119.84
9	C	200	LEU	N-CA-C	6.49	118.05	110.97
4	4	174	GLY	CA-C-N	6.49	128.97	120.28
4	4	174	GLY	C-N-CA	6.49	128.97	120.28
2	2	145	ASN	CA-CB-CG	6.48	119.08	112.60
4	4	271	ILE	N-CA-CB	6.47	118.12	110.55
14	K	107	PHE	N-CA-C	-6.47	105.25	113.01
9	C	139	PHE	N-CA-C	6.43	118.83	111.11
9	C	139	PHE	CA-CB-CG	-6.41	107.39	113.80
2	2	307	PHE	N-CA-C	6.41	118.80	111.11
2	2	313	ILE	N-CA-CB	6.38	120.16	110.58
5	5	252	ALA	N-CA-C	-6.38	105.51	113.55
9	C	123	THR	CA-C-N	6.37	128.81	120.28
9	C	123	THR	C-N-CA	6.37	128.81	120.28
11	G	178	PHE	CA-CB-CG	6.32	120.12	113.80
4	4	386	THR	CB-CA-C	6.32	120.09	109.67
2	2	184	LEU	CA-C-N	6.25	128.99	120.54
2	2	184	LEU	C-N-CA	6.25	128.99	120.54
2	2	306	GLU	N-CA-C	-6.25	104.39	111.14
5	5	341	LEU	N-CA-C	-6.17	104.45	112.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	5	138	VAL	N-CA-CB	6.16	119.82	110.58
9	C	429	ALA	CA-C-N	6.13	125.34	118.97
9	C	429	ALA	C-N-CA	6.13	125.34	118.97
15	L	23	ILE	CA-C-N	6.13	128.49	120.28
15	L	23	ILE	C-N-CA	6.13	128.49	120.28
9	C	232	PRO	N-CA-CB	6.13	110.02	103.52
2	2	325	MET	CA-C-N	6.12	128.85	120.46
2	2	325	MET	C-N-CA	6.12	128.85	120.46
5	5	645	VAL	N-CA-C	-6.11	107.90	113.71
2	2	185	HIS	CA-CB-CG	6.10	119.90	113.80
14	K	109	ALA	N-CA-C	6.09	118.00	111.36
2	2	242	ILE	N-CA-CB	6.07	117.24	110.62
5	5	144	GLU	N-CA-C	-6.04	105.73	113.23
2	2	165	SER	N-CA-C	-6.01	104.81	111.36
9	C	199	ALA	CA-C-N	5.98	128.54	120.65
9	C	199	ALA	C-N-CA	5.98	128.54	120.65
4	4	228	HIS	CA-C-N	5.96	132.70	121.97
4	4	228	HIS	C-N-CA	5.96	132.70	121.97
5	5	434	LEU	N-CA-C	-5.95	105.33	113.30
5	5	577	HIS	N-CA-C	-5.94	104.38	111.69
3	3	67	ASP	CA-CB-CG	5.93	118.53	112.60
2	2	66	PHE	CA-CB-CG	-5.93	107.87	113.80
5	5	236	LEU	N-CA-C	5.93	118.25	111.02
9	C	144	TYR	N-CA-C	5.92	117.74	111.28
2	2	48	ASN	CA-C-N	5.92	132.35	121.70
2	2	48	ASN	C-N-CA	5.92	132.35	121.70
5	5	581	LYS	CA-C-N	5.91	126.54	119.98
5	5	581	LYS	C-N-CA	5.91	126.54	119.98
9	C	430	PRO	N-CA-CB	5.91	109.55	103.23
9	C	196	ASP	CA-CB-CG	5.91	118.50	112.60
2	2	230	ILE	N-CA-C	5.88	121.57	109.34
1	1	198	ALA	N-CA-C	5.88	123.32	110.80
5	5	642	VAL	CA-C-N	5.86	126.06	120.43
5	5	642	VAL	C-N-CA	5.86	126.06	120.43
5	5	151	TYR	CA-C-N	5.84	128.91	120.79
5	5	151	TYR	C-N-CA	5.84	128.91	120.79
1	1	293	PHE	CA-CB-CG	-5.83	107.97	113.80
2	2	210	PHE	CA-CB-CG	5.82	119.62	113.80
4	4	383	ASN	CA-C-N	5.82	128.41	121.84
4	4	383	ASN	C-N-CA	5.82	128.41	121.84
2	2	280	LYS	N-CA-C	5.80	119.70	109.96
1	1	240	ALA	CA-C-N	5.78	127.95	120.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	1	240	ALA	C-N-CA	5.78	127.95	120.44
14	K	73	GLN	N-CA-C	5.77	118.31	111.33
14	K	110	SER	N-CA-C	5.75	120.32	112.55
2	2	218	HIS	CA-C-N	5.75	128.31	120.54
2	2	218	HIS	C-N-CA	5.75	128.31	120.54
2	2	396	TYR	N-CA-C	5.74	118.41	111.40
15	L	85	ASN	CA-C-N	5.74	132.51	121.54
15	L	85	ASN	C-N-CA	5.74	132.51	121.54
9	C	199	ALA	N-CA-C	5.74	119.31	110.42
9	C	372	ASN	N-CA-C	-5.73	106.84	113.88
6	6	150	ILE	CA-C-N	5.73	126.34	119.98
6	6	150	ILE	C-N-CA	5.73	126.34	119.98
3	3	7	PHE	CA-CB-CG	5.72	119.52	113.80
1	1	46	GLY	CA-C-N	5.71	128.51	120.28
1	1	46	GLY	C-N-CA	5.71	128.51	120.28
15	L	21	ASN	N-CA-C	5.71	122.96	110.80
2	2	270	VAL	CA-C-N	5.70	126.31	119.98
2	2	270	VAL	C-N-CA	5.70	126.31	119.98
9	C	328	ASP	CA-CB-CG	5.69	118.29	112.60
4	4	284	THR	CA-C-N	5.67	128.53	120.53
4	4	284	THR	C-N-CA	5.67	128.53	120.53
1	1	44	TYR	CA-C-N	5.67	132.36	121.54
1	1	44	TYR	C-N-CA	5.67	132.36	121.54
5	5	575	ILE	N-CA-CB	5.62	118.19	110.54
14	K	96	PRO	N-CA-CB	5.59	108.93	102.33
2	2	75	ILE	CA-C-N	5.59	128.04	120.44
2	2	75	ILE	C-N-CA	5.59	128.04	120.44
7	A	196	ILE	N-CA-C	5.58	116.23	110.82
1	1	155	ILE	N-CA-CB	5.57	118.12	110.54
1	1	154	PHE	N-CA-C	-5.57	104.42	111.11
9	C	384	HIS	CA-C-N	5.57	127.68	120.44
9	C	384	HIS	C-N-CA	5.57	127.68	120.44
2	2	169	ALA	CA-C-N	5.56	128.04	120.54
2	2	169	ALA	C-N-CA	5.56	128.04	120.54
11	G	165	VAL	CB-CA-C	5.55	117.07	110.68
2	2	43	TYR	CA-C-N	5.54	130.05	120.58
2	2	43	TYR	C-N-CA	5.54	130.05	120.58
4	4	322	SER	CA-C-N	5.52	131.26	120.99
4	4	322	SER	C-N-CA	5.52	131.26	120.99
9	C	98	LEU	N-CA-C	5.52	122.56	110.80
4	4	458	ILE	CA-C-N	5.46	128.18	120.42
4	4	458	ILE	C-N-CA	5.46	128.18	120.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
13	I	137	GLU	N-CA-C	5.46	116.92	110.97
14	K	81	PHE	CA-C-N	5.43	125.87	120.98
14	K	81	PHE	C-N-CA	5.43	125.87	120.98
8	B	137	VAL	CA-C-N	5.43	127.50	120.44
8	B	137	VAL	C-N-CA	5.43	127.50	120.44
1	1	216	PHE	CA-C-N	5.42	131.89	121.54
1	1	216	PHE	C-N-CA	5.42	131.89	121.54
14	K	179	GLY	CA-C-N	-5.42	116.89	122.85
14	K	179	GLY	C-N-CA	-5.42	116.89	122.85
2	2	159	PHE	CA-CB-CG	-5.42	108.39	113.80
3	3	87	GLY	CA-C-N	5.39	127.94	120.29
3	3	87	GLY	C-N-CA	5.39	127.94	120.29
1	1	196	GLU	CB-CG-CD	5.37	121.72	112.60
5	5	112	HIS	CA-C-N	5.36	131.78	121.54
5	5	112	HIS	C-N-CA	5.36	131.78	121.54
9	C	150	ASN	CA-CB-CG	5.36	117.96	112.60
5	5	371	PRO	CA-C-N	5.35	127.98	120.28
5	5	371	PRO	C-N-CA	5.35	127.98	120.28
4	4	464	ILE	CA-C-N	5.35	125.87	119.94
4	4	464	ILE	C-N-CA	5.35	125.87	119.94
2	2	62	ASN	CA-C-N	5.34	127.75	120.54
2	2	62	ASN	C-N-CA	5.34	127.75	120.54
2	2	147	SER	N-CA-C	5.31	117.93	109.96
1	1	160	PHE	CA-CB-CG	5.29	119.09	113.80
3	3	85	ASN	CA-CB-CG	5.29	117.89	112.60
5	5	427	TYR	CA-C-N	5.29	127.31	120.44
5	5	427	TYR	C-N-CA	5.29	127.31	120.44
1	1	310	ASN	CA-C-N	5.28	127.35	120.28
1	1	310	ASN	C-N-CA	5.28	127.35	120.28
2	2	48	ASN	CA-CB-CG	5.26	117.86	112.60
2	2	123	PHE	N-CA-C	5.24	121.95	110.80
2	2	308	SER	N-CA-C	5.22	116.65	111.07
1	1	21	TYR	CA-C-N	5.21	127.99	120.38
1	1	21	TYR	C-N-CA	5.21	127.99	120.38
2	2	204	LEU	N-CA-C	-5.21	105.50	111.07
14	K	70	TRP	N-CA-C	-5.21	105.51	111.14
1	1	206	GLU	CA-C-N	5.19	131.46	121.54
1	1	206	GLU	C-N-CA	5.19	131.46	121.54
4	4	149	PHE	CA-CB-CG	5.17	118.97	113.80
5	5	399	ILE	CA-C-N	5.17	127.76	120.42
5	5	399	ILE	C-N-CA	5.17	127.76	120.42
7	A	137	ALA	CA-C-N	5.16	130.99	121.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	A	137	ALA	C-N-CA	5.16	130.99	121.70
2	2	217	LEU	CA-C-N	5.16	131.40	121.54
2	2	217	LEU	C-N-CA	5.16	131.40	121.54
2	2	262	ILE	N-CA-C	5.16	116.50	110.62
9	C	304	ASP	CA-CB-CG	5.16	117.76	112.60
11	G	76	LYS	N-CA-C	-5.15	105.83	112.68
5	5	233	HIS	N-CA-C	5.14	118.82	112.54
4	4	202	LEU	N-CA-C	5.14	121.75	110.80
9	C	217	TYR	CA-CB-CG	5.14	123.15	113.90
2	2	318	ILE	CA-C-N	5.13	127.41	120.38
2	2	318	ILE	C-N-CA	5.13	127.41	120.38
6	6	10	GLU	CA-C-N	5.13	129.43	120.86
6	6	10	GLU	C-N-CA	5.13	129.43	120.86
2	2	244	ILE	N-CA-C	5.11	115.33	110.42
4	4	167	VAL	N-CA-CB	5.11	116.43	110.65
4	4	139	ILE	CA-C-N	5.10	127.11	120.28
4	4	139	ILE	C-N-CA	5.10	127.11	120.28
5	5	372	TYR	CA-C-N	5.08	127.40	120.54
5	5	372	TYR	C-N-CA	5.08	127.40	120.54
15	L	61	ILE	CA-C-N	5.06	127.06	120.28
15	L	61	ILE	C-N-CA	5.06	127.06	120.28
11	G	201	THR	CA-C-N	5.05	128.03	120.90
11	G	201	THR	C-N-CA	5.05	128.03	120.90
2	2	366	ILE	N-CA-CB	5.05	116.46	110.55
2	2	16	SER	N-CA-C	5.04	117.71	109.59
4	4	204	LEU	CA-C-N	5.03	131.14	121.54
4	4	204	LEU	C-N-CA	5.03	131.14	121.54
14	K	180	CYS	N-CA-C	5.02	117.28	108.75
5	5	234	ASN	CA-CB-CG	5.00	117.60	112.60
1	1	291	PHE	CA-CB-CG	5.00	118.80	113.80

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	1	2177	0	1809	76	0
2	2	3142	0	2868	137	0
3	3	641	0	608	45	0
4	4	3017	0	2276	109	0
5	5	4065	0	3176	94	0
6	6	1078	0	1125	55	0
7	A	3226	0	1030	10	0
8	B	2013	0	839	23	0
9	C	2647	0	2375	71	0
10	E	975	0	229	1	0
11	G	880	0	581	21	0
12	H	803	0	302	7	0
13	I	857	0	563	31	0
14	K	1023	0	899	24	0
15	L	660	0	701	35	0
16	D	285	0	64	5	0
16	Z	285	0	63	3	0
17	F	270	0	60	0	0
18	J	315	0	71	1	0
19	M	145	0	36	2	0
20	N	250	0	53	2	0
21	O	350	0	75	2	0
22	P	230	0	52	2	0
23	Q	255	0	55	2	0
24	R	150	0	32	0	0
25	S	345	0	77	2	0
26	AH	75	0	17	0	0
26	T	75	0	17	0	0
27	U	130	0	29	0	0
28	V	110	0	24	2	0
29	AB	45	0	14	0	0
29	AY	45	0	11	0	0
29	BE	45	0	12	0	0
29	W	45	0	11	1	0
30	AJ	80	0	19	0	0
30	AV	80	0	19	0	0
30	BH	80	0	18	1	0
30	X	80	0	19	0	0
31	AR	65	0	16	0	0
31	AT	65	0	15	0	0
31	Y	65	0	16	0	0
32	AA	90	0	20	0	0
32	AW	90	0	24	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
32	BB	90	0	21	0	0
32	BG	90	0	22	0	0
33	AC	235	0	49	4	0
33	AD	235	0	52	0	0
34	AE	240	0	54	2	0
35	AF	175	0	37	0	0
36	AG	125	0	29	0	0
37	AI	180	0	39	0	0
37	AL	180	0	44	0	0
38	AK	380	0	79	0	0
38	AN	380	0	84	1	0
39	AM	85	0	20	0	0
40	AO	160	0	34	0	0
41	AP	55	0	13	0	0
41	AS	54	0	13	0	0
42	AQ	40	0	10	0	0
42	BA	40	0	10	0	0
43	AU	290	0	64	0	0
44	AX	195	0	44	0	0
45	AZ	200	0	48	1	0
46	BC	100	0	22	1	0
47	BD	95	0	22	0	0
47	BF	95	0	21	0	0
48	BI	45	0	11	0	0
49	A	4	0	0	0	0
49	H	4	0	0	1	0
50	A	16	0	0	1	0
50	B	8	0	0	0	0
50	I	16	0	0	2	0
50	K	8	0	0	2	0
All	All	35169	0	21162	688	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 12.

All (688) 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)
5:5:585:TYR:O	5:5:589:PRO:HD2	1.36	1.25
3:3:98:LEU:HD11	6:6:170:LEU:HA	1.27	1.15
13:I:175:CYS:HB2	13:I:184:ILE:HD13	1.17	1.09

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:H:217:UNK:CB	12:H:220:UNK:O	2.01	1.08
6:6:28:ASN:HB3	6:6:29:PRO:HD3	1.34	1.06
5:5:95:ILE:HD11	5:5:336:ALA:HB1	1.41	1.02
4:4:225:PHE:H	4:4:284:THR:HG22	1.27	0.99
2:2:247:TYR:HA	2:2:250:LEU:HD23	1.44	0.96
1:1:208:GLU:HA	1:1:212:VAL:O	1.68	0.94
8:B:139:GLY:HA2	8:B:142:LEU:HD12	1.48	0.92
1:1:159:MET:HG2	3:3:77:VAL:HG21	1.49	0.92
5:5:585:TYR:O	5:5:589:PRO:CD	2.16	0.92
4:4:223:PRO:HG3	4:4:231:LEU:HG	1.52	0.91
9:C:385:PHE:HA	13:I:135:LEU:HD21	1.55	0.88
2:2:209:LEU:HD23	2:2:212:ILE:HD12	1.57	0.87
4:4:280:ILE:O	4:4:284:THR:HG23	1.75	0.87
8:B:410:GLU:O	8:B:414:LEU:HG	1.75	0.86
5:5:173:MET:HB3	5:5:235:TRP:HD1	1.40	0.85
9:C:293:GLY:HA2	9:C:295:MET:N	1.92	0.85
6:6:23:ILE:HG22	6:6:35:TYR:HB3	1.59	0.85
9:C:146:SER:HA	9:C:181:THR:HG22	1.58	0.85
2:2:1:MET:HA	2:2:4:LEU:HD13	1.57	0.85
5:5:151:TYR:HB2	5:5:171:VAL:HG21	1.57	0.85
13:I:175:CYS:CB	13:I:184:ILE:HD13	2.03	0.85
11:G:168:ILE:O	11:G:168:ILE:HG22	1.77	0.85
6:6:176:PRO:HA	6:6:179:ILE:HG22	1.59	0.84
9:C:388:TYR:O	13:I:139:ILE:HB	1.78	0.84
3:3:98:LEU:CD1	6:6:170:LEU:HA	2.06	0.83
2:2:78:VAL:HG11	2:2:321:LEU:HD21	1.60	0.83
3:3:61:ILE:O	3:3:65:PRO:HD3	1.78	0.83
13:I:139:ILE:HG23	13:I:140:CYS:H	1.42	0.82
2:2:230:ILE:O	2:2:234:ILE:HG13	1.80	0.82
5:5:576:ALA:HA	5:5:579:VAL:HB	1.61	0.82
34:AE:2212:UNK:CB	34:AE:2227:UNK:HA	2.10	0.80
1:1:194:VAL:HG12	1:1:293:PHE:HZ	1.47	0.78
12:H:217:UNK:CB	12:H:221:UNK:HA	2.13	0.78
5:5:126:TRP:CZ3	5:5:146:ILE:HG12	2.18	0.78
1:1:213:ALA:HB1	1:1:218:GLU:H	1.46	0.77
9:C:293:GLY:HA2	9:C:296:LEU:H	1.49	0.77
3:3:98:LEU:HD13	6:6:170:LEU:HD13	1.65	0.77
2:2:67:TYR:CE1	2:2:314:THR:HG21	2.21	0.76
7:A:135:ALA:HB2	13:I:131:ILE:HG22	1.66	0.76
1:1:45:TYR:HB3	1:1:47:LEU:HG	1.66	0.75
4:4:265:LEU:HG	4:4:266:LEU:N	2.01	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:2:138:TYR:OH	15:L:73:LEU:HB2	1.86	0.75
4:4:288:THR:HG21	4:4:307:ILE:HG12	1.67	0.75
13:I:176:GLN:HA	13:I:184:ILE:HG23	1.69	0.75
2:2:71:LEU:HD21	2:2:314:THR:HA	1.68	0.74
6:6:12:THR:HG21	6:6:46:TYR:HB2	1.69	0.74
2:2:125:THR:HA	6:6:161:ILE:HD11	1.68	0.74
9:C:337:MET:HA	9:C:340:PHE:HD2	1.52	0.74
13:I:181:VAL:O	13:I:182:ASP:HB2	1.87	0.73
9:C:329:CYS:HB3	9:C:456:THR:HG21	1.70	0.73
2:2:262:ILE:O	2:2:266:LEU:HG	1.88	0.73
2:2:155:MET:HA	15:L:73:LEU:HD21	1.68	0.73
1:1:240:ALA:HA	1:1:286:LEU:HD21	1.70	0.73
1:1:235:ILE:HA	1:1:238:ILE:HG22	1.71	0.73
4:4:327:UNK:CB	4:4:397:LEU:HB3	2.18	0.72
6:6:28:ASN:HB3	6:6:29:PRO:CD	2.17	0.72
34:AE:2208:UNK:CB	34:AE:2231:UNK:HA	2.19	0.72
6:6:74:ILE:HG13	6:6:78:LEU:HD13	1.71	0.72
8:B:60:UNK:HA	8:B:138:GLU:OE2	1.89	0.72
9:C:289:LEU:HD23	11:G:173:GLU:HB2	1.71	0.72
4:4:85:UNK:O	4:4:89:UNK:CB	2.37	0.72
5:5:173:MET:HB3	5:5:235:TRP:CD1	2.24	0.72
13:I:175:CYS:HB2	13:I:184:ILE:CD1	2.10	0.71
3:3:91:VAL:O	3:3:95:LEU:HG	1.90	0.71
2:2:180:ASN:HA	15:L:47:PHE:HE2	1.56	0.71
2:2:247:TYR:HA	2:2:250:LEU:CD2	2.18	0.71
6:6:28:ASN:CB	6:6:29:PRO:HD3	2.18	0.71
15:L:54:LEU:O	15:L:57:ILE:HG22	1.90	0.71
1:1:240:ALA:HA	1:1:286:LEU:CD2	2.21	0.71
5:5:401:GLU:HG2	5:5:495:UNK:O	1.90	0.71
4:4:163:ALA:O	4:4:167:VAL:HG23	1.91	0.71
5:5:313:LEU:O	5:5:317:THR:HG23	1.91	0.70
2:2:208:LEU:HB3	2:2:244:ILE:HD11	1.74	0.70
2:2:173:ASN:HD21	15:L:40:LEU:HD13	1.56	0.70
4:4:450:UNK:HA	4:4:453:ILE:HD12	1.71	0.70
4:4:138:TYR:CZ	4:4:179:LEU:HB2	2.25	0.70
1:1:154:PHE:CZ	1:1:241:PHE:HE1	2.10	0.69
9:C:134:GLN:O	9:C:137:PRO:HD2	1.93	0.69
3:3:61:ILE:O	3:3:65:PRO:CD	2.41	0.69
1:1:194:VAL:HG12	1:1:293:PHE:CZ	2.26	0.69
14:K:61:LEU:HA	14:K:63:THR:H	1.56	0.69
4:4:178:MET:HB2	4:4:217:ALA:HB2	1.73	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:4:254:ALA:O	4:4:258:ILE:HG13	1.93	0.69
9:C:289:LEU:CD2	11:G:173:GLU:HB2	2.24	0.68
14:K:81:PHE:HE2	14:K:134:VAL:HG11	1.59	0.68
1:1:241:PHE:HA	1:1:244:TYR:CE2	2.28	0.68
3:3:9:ILE:C	3:3:12:PRO:HD2	2.19	0.68
5:5:143:TRP:HA	5:5:146:ILE:HG13	1.75	0.67
5:5:126:TRP:CH2	5:5:146:ILE:HG12	2.29	0.67
5:5:619:MET:O	5:5:620:LEU:HB2	1.94	0.67
3:3:98:LEU:HD12	6:6:173:ILE:HG12	1.74	0.67
2:2:214:ILE:HG23	2:2:268:LEU:HD22	1.77	0.67
2:2:269:LEU:O	2:2:273:VAL:HB	1.94	0.67
2:2:354:UNK:HA	2:2:361:VAL:HG21	1.75	0.67
2:2:212:ILE:HG12	2:2:244:ILE:CG2	2.25	0.66
2:2:376:PRO:HD3	4:4:142:GLU:HB2	1.77	0.66
2:2:75:ILE:HG12	2:2:238:LEU:HD23	1.76	0.66
4:4:387:PRO:HB2	5:5:141:VAL:HG13	1.78	0.66
13:I:184:ILE:HG23	13:I:184:ILE:O	1.94	0.66
8:B:413:MET:O	8:B:417:LEU:HG	1.95	0.66
4:4:400:GLN:HA	5:5:186:CYS:SG	2.36	0.65
4:4:458:ILE:O	4:4:462:LEU:HG	1.97	0.65
9:C:293:GLY:CA	9:C:296:LEU:H	2.08	0.65
5:5:341:LEU:HD11	5:5:466:ALA:HA	1.78	0.65
2:2:265:ILE:HD12	2:2:298:LEU:HG	1.79	0.65
4:4:459:ILE:O	4:4:463:ILE:HG13	1.96	0.65
9:C:378:MET:HE3	9:C:382:ILE:HD11	1.79	0.65
13:I:181:VAL:O	13:I:182:ASP:CB	2.45	0.65
5:5:89:LEU:HB2	5:5:131:VAL:HG11	1.77	0.65
9:C:95:HIS:CG	9:C:96:GLY:H	2.15	0.65
9:C:293:GLY:HA2	9:C:295:MET:H	1.60	0.65
4:4:288:THR:HG21	4:4:307:ILE:CG1	2.26	0.64
11:G:168:ILE:O	11:G:168:ILE:CG2	2.45	0.64
1:1:149:ILE:HG12	1:1:316:LEU:HD23	1.80	0.64
2:2:213:GLY:HA3	2:2:221:LEU:HD12	1.78	0.64
6:6:70:LEU:HB3	15:L:71:LEU:HD23	1.80	0.64
5:5:364:UNK:HA	5:5:440:UNK:HA	1.80	0.64
11:G:168:ILE:HG23	11:G:171:LEU:HB3	1.79	0.64
4:4:247:ALA:HA	4:4:251:LEU:HD13	1.79	0.63
14:K:196:ARG:O	14:K:201:THR:CB	2.46	0.63
2:2:3:ILE:O	2:2:7:ILE:HG13	1.98	0.63
9:C:279:GLY:H	9:C:445:HIS:CE1	2.17	0.63
9:C:217:TYR:C	9:C:217:TYR:HD1	2.07	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:2:212:ILE:HG12	2:2:244:ILE:HG23	1.81	0.62
5:5:126:TRP:HA	5:5:129:ILE:HD12	1.81	0.62
2:2:414:LEU:HB3	4:4:165:PHE:HE2	1.64	0.62
5:5:288:UNK:CB	5:5:314:GLY:HA3	2.29	0.62
2:2:123:PHE:HB2	2:2:182:TYR:HB3	1.81	0.62
2:2:212:ILE:HG23	2:2:241:LYS:HG3	1.81	0.62
7:A:120:ASN:HD21	8:B:420:ASP:HB3	1.64	0.62
2:2:164:LEU:HB3	2:2:210:PHE:HB2	1.82	0.62
2:2:262:ILE:HG23	2:2:301:LEU:HD22	1.82	0.62
4:4:240:LEU:HD11	4:4:344:UNK:HA	1.81	0.62
6:6:101:LEU:HA	6:6:104:LEU:HD23	1.81	0.62
2:2:209:LEU:HD23	2:2:212:ILE:CD1	2.30	0.61
5:5:312:GLN:HB3	5:5:335:HIS:NE2	2.15	0.61
6:6:176:PRO:HA	6:6:179:ILE:CG2	2.29	0.61
33:AC:2012:UNK:HA	33:AC:2013:UNK:C	2.30	0.61
1:1:279:SER:HB2	22:P:711:UNK:HA	1.82	0.61
9:C:320:ASP:O	9:C:322:PRO:HD3	2.01	0.61
5:5:322:LEU:HD13	5:5:403:THR:HG22	1.83	0.61
5:5:123:PHE:HA	5:5:126:TRP:CD1	2.36	0.61
9:C:180:ILE:HD11	9:C:247:ILE:HD11	1.82	0.61
2:2:132:LEU:HG	6:6:168:VAL:HG23	1.83	0.61
2:2:139:LEU:HD11	15:L:69:ILE:HD11	1.82	0.61
4:4:152:ILE:O	4:4:156:GLY:HA3	2.01	0.61
5:5:30:UNK:HA	5:5:114:VAL:HG22	1.83	0.61
2:2:391:LEU:HD13	4:4:184:VAL:HG22	1.82	0.60
6:6:33:ILE:CD1	6:6:75:ILE:HG12	2.31	0.60
2:2:64:PHE:CZ	2:2:250:LEU:HD22	2.37	0.60
7:A:185:HIS:CD2	8:B:388:ARG:HD3	2.36	0.60
1:1:170:ILE:HG12	1:1:244:TYR:HB2	1.84	0.60
5:5:566:UNK:O	5:5:570:VAL:HG23	2.02	0.60
1:1:225:VAL:O	1:1:229:LEU:HB2	2.02	0.60
1:1:202:PHE:HA	1:1:205:THR:HB	1.83	0.60
2:2:262:ILE:O	2:2:265:ILE:HG12	2.01	0.60
4:4:204:LEU:HD22	4:4:268:GLU:HG2	1.84	0.60
5:5:92:ILE:HG22	5:5:124:THR:HG22	1.84	0.60
1:1:22:LEU:HG	1:1:236:ILE:HD11	1.83	0.59
33:AC:2020:UNK:HA	33:AC:2021:UNK:CB	2.32	0.59
5:5:618:SER:HB2	15:L:18:ASN:CB	2.33	0.59
14:K:117:MET:HB3	14:K:144:VAL:HG22	1.84	0.59
2:2:376:PRO:HA	4:4:139:ILE:HG22	1.85	0.59
8:B:411:ILE:HD13	8:B:449:UNK:HA	1.85	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:2:59:LEU:HB2	2:2:121:ASN:HB2	1.84	0.59
5:5:143:TRP:HH2	5:5:222:ALA:O	1.86	0.59
4:4:161:GLU:O	4:4:165:PHE:HD1	1.85	0.58
4:4:316:GLY:HA3	4:4:398:SER:HB2	1.85	0.58
5:5:370:LEU:O	5:5:373:THR:HB	2.04	0.58
14:K:111:PRO:HG3	14:K:134:VAL:HG13	1.86	0.58
2:2:138:TYR:CE2	15:L:69:ILE:HG23	2.39	0.58
5:5:633:LEU:HD11	6:6:109:LEU:HD22	1.85	0.58
2:2:164:LEU:HB3	2:2:210:PHE:CB	2.34	0.58
6:6:17:ILE:HG13	15:L:10:LEU:HD11	1.86	0.58
9:C:217:TYR:C	9:C:217:TYR:CD1	2.82	0.58
9:C:191:CYS:HB3	9:C:203:PHE:HA	1.86	0.58
4:4:143:ALA:O	4:4:146:PRO:HD2	2.04	0.58
4:4:228:HIS:HE1	4:4:287:LEU:HD13	1.69	0.58
1:1:152:SER:HB3	1:1:324:PHE:HZ	1.69	0.57
1:1:213:ALA:HB1	1:1:218:GLU:N	2.18	0.57
4:4:219:MET:HG2	4:4:224:LEU:HD12	1.85	0.57
6:6:176:PRO:HB3	15:L:72:SER:HB2	1.86	0.57
2:2:171:SER:HA	2:2:174:THR:OG1	2.04	0.57
2:2:259:UNK:O	2:2:262:ILE:HG12	2.04	0.57
14:K:199:ARG:O	14:K:200:ASN:HB2	2.03	0.57
5:5:191:VAL:HG23	5:5:211:LEU:HD23	1.85	0.57
8:B:309:UNK:HA	8:B:360:UNK:HA	1.85	0.57
1:1:314:ILE:HG12	3:3:103:VAL:HG12	1.86	0.57
2:2:231:LEU:HA	2:2:234:ILE:HD12	1.86	0.57
3:3:98:LEU:HD11	6:6:170:LEU:CA	2.17	0.57
13:I:187:THR:HB	13:I:188:PRO:CD	2.34	0.57
3:3:66:PHE:CE1	3:3:99:ILE:HG13	2.40	0.57
1:1:24:VAL:HG13	1:1:39:PRO:HG2	1.85	0.56
2:2:369:PHE:HD1	2:2:374:ILE:HG13	1.69	0.56
13:I:132:ALA:HB2	13:I:158:UNK:C	2.35	0.56
2:2:261:SER:O	2:2:265:ILE:HG23	2.05	0.56
4:4:364:GLY:N	4:4:433:UNK:O	2.33	0.56
5:5:308:SER:O	5:5:312:GLN:HG2	2.06	0.56
5:5:312:GLN:HB3	5:5:335:HIS:CE1	2.40	0.56
10:E:608:UNK:N	10:E:609:UNK:HA	2.19	0.56
4:4:455:ASN:O	4:4:459:ILE:HD12	2.05	0.56
1:1:201:PRO:O	1:1:205:THR:N	2.29	0.56
2:2:166:ILE:HD11	15:L:33:LEU:HG	1.87	0.56
4:4:399:LEU:HD21	4:4:414:UNK:HA	1.87	0.56
13:I:137:GLU:HG3	13:I:145:UNK:O	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:2:142:ALA:HB2	2:2:154:SER:HB2	1.88	0.56
14:K:79:VAL:HB	14:K:114:SER:OG	2.06	0.56
1:1:158:ILE:O	1:1:162:SER:N	2.39	0.56
2:2:51:THR:HB	2:2:61:PHE:HA	1.88	0.56
6:6:165:LEU:HD11	15:L:58:VAL:HG22	1.87	0.56
4:4:202:LEU:HD22	4:4:207:GLN:HG2	1.86	0.56
4:4:347:UNK:O	4:4:351:UNK:N	2.38	0.56
1:1:154:PHE:CZ	1:1:241:PHE:CE1	2.94	0.56
4:4:387:PRO:HD3	5:5:144:GLU:HB3	1.88	0.56
5:5:208:UNK:HA	5:5:209:THR:HG23	1.88	0.56
5:5:388:PRO:HA	5:5:393:TYR:HB2	1.88	0.56
2:2:299:LEU:HD11	2:2:386:ILE:HG12	1.88	0.55
1:1:28:LYS:HD3	1:1:38:GLY:H	1.70	0.55
2:2:256:UNK:O	2:2:258:UNK:N	2.39	0.55
2:2:375:PRO:HG2	4:4:138:TYR:HE2	1.71	0.55
5:5:143:TRP:CZ2	5:5:223:ALA:HA	2.40	0.55
4:4:220:VAL:HG22	4:4:227:ILE:HG21	1.87	0.55
4:4:295:GLN:HE21	4:4:300:VAL:HG11	1.72	0.55
5:5:92:ILE:HD12	5:5:127:MET:HG2	1.88	0.55
1:1:161:VAL:HG12	1:1:163:SER:H	1.71	0.55
13:I:145:UNK:HA	13:I:163:ASP:O	2.07	0.55
2:2:66:PHE:O	2:2:70:MET:HG2	2.06	0.55
2:2:149:LYS:O	2:2:227:ASN:HB3	2.07	0.55
2:2:269:LEU:HD21	2:2:298:LEU:HD11	1.89	0.55
4:4:384:ILE:HG12	4:4:421:UNK:HA	1.89	0.55
14:K:79:VAL:HG23	14:K:108:ARG:HA	1.88	0.55
14:K:192:PHE:HA	14:K:195:GLN:HB2	1.88	0.55
1:1:214:GLY:HA2	14:K:109:ALA:HB3	1.88	0.55
9:C:407:LYS:HE2	9:C:460:VAL:HG23	1.89	0.55
33:AC:2013:UNK:O	33:AC:2015:UNK:N	2.39	0.55
1:1:241:PHE:HA	1:1:244:TYR:CD2	2.42	0.54
2:2:211:LYS:HE3	2:2:244:ILE:HD13	1.89	0.54
2:2:295:TYR:HE1	2:2:402:LEU:HG	1.73	0.54
8:B:101:MET:HE2	8:B:147:MET:HB3	1.89	0.54
22:P:716:UNK:O	22:P:720:UNK:N	2.40	0.54
1:1:34:GLN:HE22	9:C:200:LEU:HD12	1.71	0.54
2:2:173:ASN:ND2	15:L:40:LEU:HD13	2.22	0.54
2:2:372:ILE:HB	2:2:410:ALA:CB	2.38	0.54
3:3:71:SER:O	3:3:74:LEU:HB3	2.06	0.54
4:4:366:THR:HG22	4:4:373:ALA:HB1	1.90	0.54
13:I:140:CYS:SG	13:I:144:ALA:HB2	2.47	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:43:GLY:HA3	1:1:44:TYR:C	2.32	0.54
9:C:140:ASP:HB2	9:C:147:MET:HB2	1.90	0.54
2:2:234:ILE:HG12	2:2:324:PHE:CG	2.43	0.54
6:6:153:VAL:HG13	6:6:157:ASN:HB3	1.89	0.54
14:K:121:GLY:HA2	50:K:500:SF4:S4	2.47	0.54
4:4:184:VAL:HB	4:4:210:ILE:HG21	1.90	0.54
8:B:142:LEU:HD11	8:B:253:UNK:CB	2.38	0.54
2:2:131:GLU:CB	15:L:62:LEU:HD21	2.38	0.54
4:4:219:MET:O	4:4:223:PRO:HA	2.07	0.54
5:5:549:UNK:O	5:5:560:UNK:N	2.41	0.54
12:H:137:SER:HA	12:H:140:ILE:HD12	1.89	0.54
3:3:65:PRO:CB	6:6:173:ILE:HG22	2.38	0.54
4:4:467:CYS:HB3	4:4:469:GLN:HG3	1.89	0.54
5:5:136:TYR:HB2	5:5:195:UNK:O	2.07	0.54
5:5:251:HIS:ND1	5:5:309:THR:HG21	2.23	0.54
5:5:424:THR:HA	5:5:427:TYR:CE2	2.43	0.53
4:4:255:LEU:HA	4:4:258:ILE:HD12	1.90	0.53
14:K:170:ILE:HG22	14:K:171:VAL:HG23	1.90	0.53
6:6:33:ILE:HD11	6:6:75:ILE:HG12	1.90	0.53
6:6:71:PHE:HE1	15:L:26:PHE:HB2	1.71	0.53
4:4:274:THR:HA	4:4:277:ILE:HD12	1.90	0.53
6:6:150:ILE:HD11	15:L:53:SER:HB2	1.90	0.53
13:I:189:ASN:HD21	13:I:245:UNK:CB	2.22	0.53
11:G:176:ASN:HA	11:G:193:HIS:HE1	1.73	0.53
4:4:292:THR:HA	4:4:295:GLN:HG2	1.90	0.53
9:C:149:THR:HG21	9:C:406:PRO:HD3	1.90	0.53
1:1:49:GLN:O	1:1:53:ASP:HB2	2.09	0.53
4:4:297:ASP:HB3	4:4:300:VAL:HB	1.90	0.53
9:C:147:MET:HG2	9:C:148:MET:N	2.24	0.53
3:3:62:LEU:O	3:3:65:PRO:HD2	2.09	0.53
5:5:331:HIS:CE1	5:5:396:LYS:HG2	2.44	0.53
6:6:173:ILE:HG12	6:6:174:ILE:HD12	1.91	0.53
9:C:87:PHE:HB3	9:C:100:LEU:HB3	1.91	0.53
6:6:55:PHE:O	6:6:59:TYR:HB2	2.09	0.52
1:1:187:UNK:HA	1:1:246:LEU:HD13	1.90	0.52
2:2:142:ALA:HB2	2:2:154:SER:CB	2.38	0.52
7:A:75:ALA:HB3	7:A:191:ARG:HD2	1.91	0.52
8:B:411:ILE:HD11	8:B:452:UNK:CB	2.39	0.52
2:2:397:PHE:O	2:2:401:VAL:HG23	2.08	0.52
4:4:251:LEU:HD23	4:4:310:MET:HG3	1.91	0.52
9:C:262:GLU:HG3	9:C:337:MET:HG2	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:C:313:ALA:O	9:C:317:VAL:HG23	2.09	0.52
16:D:144:UNK:C	16:D:146:UNK:H	2.22	0.52
3:3:69:GLU:HB3	3:3:95:LEU:HD13	1.92	0.52
4:4:375:TYR:CZ	4:4:454:MET:HG2	2.45	0.52
8:B:384:CYS:SG	8:B:426:ILE:HB	2.49	0.52
14:K:89:GLU:HG2	14:K:182:PRO:O	2.09	0.52
3:3:2:ASN:HB2	3:3:5:ILE:HG12	1.91	0.52
5:5:18:UNK:CB	5:5:115:ARG:HA	2.40	0.52
8:B:404:UNK:O	8:B:405:UNK:CB	2.57	0.52
1:1:159:MET:HG2	3:3:77:VAL:CG2	2.32	0.52
16:Z:39:UNK:C	16:Z:41:UNK:H	2.22	0.52
1:1:326:ILE:HG23	23:Q:933:UNK:CB	2.40	0.52
1:1:315:ILE:O	1:1:318:PRO:HD2	2.10	0.52
6:6:176:PRO:HB3	15:L:72:SER:CB	2.39	0.51
2:2:215:ALA:H	2:2:268:LEU:HD13	1.74	0.51
4:4:168:LEU:O	4:4:172:LEU:HB2	2.10	0.51
4:4:252:LYS:HG3	4:4:332:UNK:CB	2.41	0.51
12:H:218:UNK:C	12:H:220:UNK:N	2.72	0.51
2:2:64:PHE:CE1	2:2:250:LEU:HD22	2.45	0.51
14:K:116:ILE:HG13	14:K:143:TRP:HB2	1.90	0.51
9:C:95:HIS:CD2	9:C:96:GLY:H	2.28	0.51
9:C:124:GLU:HB3	9:C:426:LYS:NZ	2.25	0.51
2:2:217:LEU:O	2:2:219:LYS:N	2.44	0.51
3:3:93:LEU:O	3:3:97:ILE:HD12	2.10	0.51
4:4:334:UNK:O	4:4:338:UNK:N	2.44	0.51
4:4:396:PHE:CE1	5:5:185:LEU:HB3	2.46	0.51
5:5:571:THR:O	5:5:575:ILE:HG13	2.11	0.51
1:1:189:UNK:O	1:1:193:SER:OG	2.28	0.51
5:5:281:UNK:CB	5:5:321:GLY:HA3	2.41	0.51
1:1:34:GLN:HE21	9:C:201:THR:HA	1.76	0.51
15:L:16:VAL:HA	15:L:17:PHE:C	2.36	0.51
16:Z:13:UNK:O	16:Z:17:UNK:N	2.44	0.51
2:2:214:ILE:HD12	2:2:268:LEU:CD2	2.40	0.51
4:4:228:HIS:CE1	4:4:287:LEU:HD13	2.46	0.51
2:2:163:ILE:O	2:2:166:ILE:HG22	2.11	0.50
2:2:354:UNK:CA	2:2:361:VAL:HG21	2.40	0.50
1:1:198:ALA:HB3	1:1:235:ILE:HD11	1.93	0.50
3:3:72:THR:O	3:3:76:TYR:CD2	2.65	0.50
4:4:142:GLU:HA	4:4:145:LEU:HD13	1.93	0.50
13:I:179:CYS:HA	50:I:501:SF4:S2	2.51	0.50
2:2:57:GLU:O	2:2:121:ASN:HB3	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:5:112:HIS:O	5:5:114:VAL:N	2.45	0.50
9:C:180:ILE:HG13	9:C:347:ILE:HD12	1.93	0.50
3:3:69:GLU:HG2	3:3:95:LEU:HD22	1.94	0.50
4:4:225:PHE:CE1	4:4:287:LEU:HD12	2.46	0.50
9:C:388:TYR:HD1	13:I:139:ILE:HG21	1.77	0.50
5:5:136:TYR:H	5:5:196:UNK:HA	1.77	0.50
6:6:154:LEU:HD22	6:6:162:LEU:HB2	1.93	0.50
9:C:368:SER:C	9:C:370:PRO:HD3	2.37	0.50
15:L:47:PHE:O	15:L:48:ASP:HB2	2.11	0.50
9:C:394:VAL:O	9:C:418:GLY:HA2	2.11	0.50
4:4:391:ASN:O	4:4:395:GLU:HB2	2.12	0.50
14:K:76:PHE:HB2	14:K:116:ILE:HD13	1.94	0.50
4:4:373:ALA:O	4:4:376:ILE:HG22	2.11	0.49
1:1:243:GLY:HA2	1:1:247:LEU:HB2	1.93	0.49
2:2:155:MET:HE1	15:L:74:LEU:HG	1.94	0.49
4:4:297:ASP:OD1	4:4:299:LYS:HG3	2.11	0.49
8:B:410:GLU:HA	8:B:413:MET:HB2	1.93	0.49
1:1:154:PHE:HZ	1:1:241:PHE:HE1	1.56	0.49
2:2:294:GLY:O	2:2:297:MET:HB2	2.13	0.49
2:2:296:MET:SD	2:2:313:ILE:HG12	2.52	0.49
5:5:385:MET:HE1	5:5:429:MET:HE2	1.94	0.49
9:C:279:GLY:H	9:C:445:HIS:HE1	1.57	0.49
15:L:39:ILE:HG13	15:L:40:LEU:N	2.27	0.49
1:1:145:SER:HB3	1:1:312:CYS:SG	2.52	0.49
1:1:281:ALA:O	1:1:285:LYS:HG2	2.12	0.49
2:2:235:TYR:O	2:2:239:ILE:HG13	2.12	0.49
5:5:92:ILE:CD1	5:5:127:MET:HG2	2.42	0.49
9:C:379:GLU:O	9:C:383:HIS:HB2	2.11	0.49
1:1:191:ILE:HG23	1:1:293:PHE:HE1	1.77	0.49
2:2:234:ILE:HG12	2:2:324:PHE:CD1	2.48	0.49
2:2:350:UNK:O	2:2:420:UNK:HA	2.12	0.49
25:S:1133:UNK:O	25:S:1137:UNK:N	2.45	0.49
38:AN:3124:UNK:C	38:AN:3126:UNK:N	2.76	0.49
2:2:212:ILE:HG12	2:2:244:ILE:HG21	1.94	0.49
2:2:266:LEU:HD21	2:2:398:ILE:HG23	1.94	0.49
3:3:90:ILE:HA	3:3:93:LEU:HD12	1.93	0.49
14:K:126:LYS:O	14:K:129:PRO:HD2	2.13	0.49
1:1:151:THR:HA	1:1:154:PHE:HD2	1.78	0.49
2:2:118:UNK:HA	2:2:247:TYR:HB2	1.94	0.49
2:2:325:MET:HE3	2:2:453:UNK:HA	1.95	0.49
4:4:170:PHE:O	4:4:220:VAL:HG11	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:C:294:PRO:HA	9:C:297:ARG:HG3	1.95	0.49
1:1:143:LEU:HD23	1:1:197:THR:HG22	1.95	0.48
28:V:1417:UNK:C	28:V:1419:UNK:N	2.76	0.48
1:1:321:PHE:HA	1:1:324:PHE:CE2	2.47	0.48
3:3:96:LEU:O	3:3:100:ILE:HG13	2.13	0.48
4:4:290:LEU:HB3	5:5:573:GLY:HA2	1.95	0.48
5:5:173:MET:CB	5:5:235:TRP:HD1	2.18	0.48
9:C:446:PHE:HB3	9:C:448:PRO:HD2	1.95	0.48
1:1:321:PHE:HA	1:1:324:PHE:CD2	2.48	0.48
2:2:172:ILE:HD13	2:2:173:ASN:N	2.29	0.48
6:6:147:LEU:C	6:6:149:THR:H	2.21	0.48
11:G:160:UNK:C	11:G:162:UNK:H	2.24	0.48
11:G:180:ARG:HH11	11:G:197:ARG:HB3	1.77	0.48
4:4:403:PHE:HA	4:4:410:UNK:CB	2.44	0.48
15:L:37:ASN:HD21	15:L:59:ILE:HB	1.78	0.48
25:S:1161:UNK:HA	25:S:1162:UNK:C	2.43	0.48
5:5:115:ARG:HB3	5:5:156:PHE:HZ	1.78	0.48
2:2:375:PRO:O	2:2:380:PHE:HB2	2.13	0.48
4:4:292:THR:CG2	4:4:304:TYR:HB3	2.44	0.48
8:B:385:THR:N	8:B:386:PRO:HD2	2.28	0.48
1:1:152:SER:HB3	1:1:324:PHE:CZ	2.47	0.48
6:6:165:LEU:HA	6:6:168:VAL:HG12	1.95	0.48
20:N:522:UNK:HA	20:N:524:UNK:N	2.28	0.48
2:2:1:MET:HG2	2:2:4:LEU:HD22	1.94	0.47
5:5:187:VAL:HG21	5:5:214:LEU:HD22	1.96	0.47
2:2:138:TYR:HB2	2:2:157:TYR:HB3	1.95	0.47
2:2:153:ALA:HB1	2:2:224:ILE:HG12	1.95	0.47
1:1:323:ILE:O	1:1:326:ILE:HG13	2.15	0.47
3:3:71:SER:O	3:3:75:PRO:HD3	2.13	0.47
4:4:225:PHE:CD1	4:4:283:LEU:HB3	2.49	0.47
4:4:292:THR:HG22	4:4:304:TYR:HB3	1.95	0.47
7:A:486:UNK:O	7:A:490:UNK:N	2.47	0.47
2:2:46:PHE:HB3	2:2:62:ASN:HB3	1.96	0.47
2:2:47:LEU:HB2	2:2:50:GLN:HG3	1.95	0.47
3:3:62:LEU:C	3:3:65:PRO:HD2	2.38	0.47
4:4:255:LEU:HG	4:4:317:VAL:HG21	1.97	0.47
9:C:293:GLY:HA2	9:C:296:LEU:N	2.26	0.47
1:1:295:TRP:HE3	13:I:87:LEU:HD22	1.78	0.47
2:2:295:TYR:CE1	2:2:402:LEU:HG	2.50	0.47
4:4:323:LEU:HD11	4:4:401:GLY:HA3	1.96	0.47
13:I:137:GLU:HA	13:I:144:ALA:HB3	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:226:PHE:HE1	3:3:16:PHE:HB3	1.79	0.47
2:2:67:TYR:HE1	2:2:314:THR:HG21	1.77	0.47
16:D:108:UNK:HA	16:D:109:UNK:HA	1.75	0.47
2:2:138:TYR:HE1	15:L:73:LEU:HD13	1.78	0.47
5:5:79:PHE:CE1	5:5:138:VAL:HG11	2.50	0.47
2:2:131:GLU:HB3	15:L:62:LEU:HD21	1.97	0.47
3:3:76:TYR:O	3:3:80:ILE:HG23	2.15	0.47
3:3:99:ILE:HA	3:3:102:PHE:HD2	1.80	0.47
6:6:29:PRO:O	6:6:33:ILE:HG13	2.15	0.47
4:4:228:HIS:O	4:4:231:LEU:HB2	2.15	0.46
5:5:153:LEU:HB3	5:5:246:VAL:HG11	1.97	0.46
1:1:30:LEU:HD23	1:1:298:ALA:HB2	1.97	0.46
2:2:245:LEU:O	2:2:249:VAL:HG23	2.15	0.46
2:2:279:ILE:HB	2:2:280:LYS:HE2	1.95	0.46
3:3:97:ILE:HG22	6:6:170:LEU:HD11	1.96	0.46
4:4:298:LEU:HD21	4:4:357:UNK:O	2.14	0.46
5:5:468:PHE:HA	5:5:472:ALA:HB3	1.96	0.46
9:C:92:PRO:HG2	9:C:98:LEU:HD13	1.97	0.46
9:C:465:ASP:HA	11:G:200:MET:HE3	1.96	0.46
2:2:295:TYR:O	2:2:298:LEU:HB2	2.16	0.46
4:4:312:ILE:HG12	4:4:399:LEU:HD13	1.97	0.46
5:5:99:VAL:HG11	5:5:249:LEU:HD13	1.96	0.46
11:G:168:ILE:HG23	11:G:171:LEU:HD23	1.98	0.46
2:2:372:ILE:HB	2:2:410:ALA:HB2	1.97	0.46
4:4:222:THR:HB	4:4:314:ILE:HD11	1.96	0.46
4:4:229:VAL:HG13	5:5:584:LEU:HG	1.98	0.46
4:4:263:LEU:O	4:4:267:CYS:SG	2.69	0.46
5:5:221:ILE:HG12	5:5:224:MET:HE2	1.97	0.46
13:I:171:TYR:HA	50:I:500:SF4:S1	2.55	0.46
5:5:322:LEU:O	5:5:323:SER:HB2	2.14	0.46
5:5:589:PRO:HA	5:5:592:ILE:HB	1.97	0.46
13:I:179:CYS:SG	13:I:183:ALA:HB3	2.55	0.46
14:K:84:ALA:HB3	50:K:500:SF4:S4	2.56	0.46
1:1:28:LYS:HG2	1:1:39:PRO:HD2	1.97	0.46
1:1:196:GLU:HG3	1:1:308:LEU:HD11	1.97	0.46
1:1:301:UNK:C	9:C:198:GLY:HA2	2.46	0.46
2:2:239:ILE:HG13	2:2:239:ILE:H	1.48	0.46
4:4:178:MET:HB2	4:4:217:ALA:CB	2.43	0.46
4:4:294:ARG:O	4:4:296:ILE:N	2.49	0.46
4:4:456:ILE:HG23	4:4:457:LEU:HD23	1.98	0.46
5:5:221:ILE:HA	5:5:224:MET:HG2	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:C:190:VAL:HG21	9:C:340:PHE:HE2	1.81	0.46
14:K:93:VAL:HG21	14:K:187:LEU:HD23	1.98	0.46
3:3:98:LEU:HD22	6:6:170:LEU:HD22	1.97	0.46
5:5:33:UNK:CB	5:5:114:VAL:HG13	2.46	0.46
5:5:361:UNK:HA	5:5:434:LEU:HD23	1.96	0.46
9:C:200:LEU:HD23	9:C:200:LEU:H	1.80	0.46
2:2:235:TYR:CE1	2:2:239:ILE:HG12	2.51	0.46
2:2:8:SER:O	2:2:12:PHE:HD2	1.98	0.46
2:2:351:UNK:HA	2:2:419:UNK:O	2.16	0.46
4:4:295:GLN:HG3	4:4:300:VAL:HG12	1.97	0.46
2:2:149:LYS:HB3	2:2:227:ASN:HA	1.97	0.45
2:2:384:LEU:O	2:2:388:MET:HB2	2.15	0.45
5:5:180:PHE:HB3	5:5:222:ALA:HB2	1.98	0.45
7:A:370:UNK:HA	7:A:371:UNK:CB	2.46	0.45
9:C:139:PHE:CZ	9:C:427:ILE:HG12	2.52	0.45
1:1:196:GLU:HA	1:1:201:PRO:HG3	1.98	0.45
5:5:62:UNK:N	5:5:79:PHE:O	2.49	0.45
5:5:83:ALA:O	5:5:86:ILE:HG13	2.16	0.45
1:1:166:ILE:HA	1:1:169:ILE:HD12	1.98	0.45
1:1:315:ILE:C	1:1:318:PRO:HD2	2.41	0.45
3:3:80:ILE:HB	3:3:81:TYR:H	1.56	0.45
4:4:230:TRP:CE3	4:4:234:VAL:HG21	2.51	0.45
5:5:143:TRP:HZ2	5:5:223:ALA:HA	1.80	0.45
8:B:94:SER:HA	8:B:97:LYS:HB3	1.98	0.45
3:3:53:UNK:O	3:3:55:UNK:N	2.49	0.45
5:5:478:UNK:HA	5:5:479:UNK:C	2.46	0.45
14:K:112:ARG:HA	14:K:139:PRO:HD3	1.99	0.45
2:2:206:LEU:HA	2:2:209:LEU:HD12	1.98	0.45
3:3:19:UNK:O	3:3:21:UNK:N	2.50	0.45
3:3:60:ALA:O	6:6:70:LEU:HD21	2.17	0.45
4:4:224:LEU:HD23	4:4:284:THR:HG21	1.98	0.45
5:5:188:ILE:O	5:5:191:VAL:HG12	2.16	0.45
1:1:304:UNK:O	1:1:306:ASP:N	2.49	0.45
5:5:357:UNK:C	5:5:359:UNK:N	2.80	0.45
1:1:196:GLU:HA	1:1:201:PRO:CG	2.47	0.45
2:2:262:ILE:HG23	2:2:301:LEU:CD2	2.47	0.45
4:4:162:ARG:O	4:4:166:TYR:HB2	2.16	0.45
11:G:180:ARG:HB3	11:G:199:ILE:HD11	1.98	0.45
12:H:131:PRO:HD2	49:H:300:FES:S2	2.56	0.45
1:1:34:GLN:HB3	9:C:204:LEU:HD22	1.99	0.45
4:4:240:LEU:CD1	4:4:344:UNK:HA	2.46	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:6:44:ALA:HB1	6:6:56:SER:HB3	1.98	0.45
1:1:218:GLU:O	1:1:220:UNK:N	2.49	0.45
3:3:77:VAL:O	3:3:80:ILE:HG13	2.16	0.45
5:5:383:SER:HB3	5:5:391:THR:HG22	1.99	0.45
8:B:398:UNK:CB	8:B:414:LEU:HD22	2.47	0.45
2:2:298:LEU:HD21	2:2:398:ILE:HG22	1.99	0.45
4:4:323:LEU:CD1	4:4:401:GLY:HA3	2.47	0.45
5:5:189:ALA:O	5:5:193:GLY:N	2.48	0.45
9:C:212:LYS:HD3	9:C:250:TRP:HD1	1.82	0.45
3:3:106:UNK:O	3:3:107:UNK:O	2.35	0.44
9:C:279:GLY:HA2	9:C:453:ILE:HD11	1.99	0.44
1:1:230:ALA:O	1:1:234:ASN:HB2	2.17	0.44
2:2:64:PHE:CE2	2:2:250:LEU:HD13	2.52	0.44
5:5:228:ALA:HB2	5:5:235:TRP:HB3	1.99	0.44
9:C:183:VAL:HG11	9:C:250:TRP:HZ2	1.81	0.44
2:2:278:GLN:HG2	2:2:280:LYS:HE3	1.98	0.44
2:2:310:LEU:O	2:2:314:THR:HG23	2.16	0.44
4:4:161:GLU:O	4:4:165:PHE:CD1	2.69	0.44
4:4:265:LEU:CG	4:4:266:LEU:N	2.74	0.44
4:4:452:PHE:HA	4:4:455:ASN:HB2	1.98	0.44
5:5:101:ILE:HD12	5:5:461:PRO:HB3	1.99	0.44
11:G:99:UNK:CB	11:G:156:UNK:HA	2.47	0.44
5:5:123:PHE:CE1	5:5:253:ALA:HB2	2.53	0.44
1:1:29:THR:HB	1:1:294:ILE:HG21	1.99	0.44
1:1:210:GLU:HG2	1:1:211:LEU:HG	1.98	0.44
2:2:302:LEU:HD23	2:2:303:ASN:N	2.33	0.44
3:3:82:LEU:CB	3:3:83:VAL:HB	2.46	0.44
13:I:187:THR:HB	13:I:188:PRO:HD3	1.98	0.44
2:2:166:ILE:CD1	15:L:33:LEU:HG	2.48	0.44
5:5:166:SER:HB2	5:5:238:LEU:O	2.17	0.44
5:5:634:VAL:O	5:5:638:ASN:N	2.51	0.44
9:C:407:LYS:HE3	9:C:458:ASP:OD1	2.18	0.44
5:5:339:LYS:HD2	5:5:339:LYS:HA	1.73	0.44
5:5:480:UNK:HA	5:5:481:UNK:HA	1.79	0.44
15:L:33:LEU:O	15:L:36:ILE:HG22	2.18	0.44
4:4:259:LEU:HD12	4:4:260:ARG:HD2	1.99	0.43
5:5:318:ILE:O	5:5:322:LEU:HB2	2.18	0.43
6:6:79:ASP:HB3	15:L:78:TYR:OH	2.18	0.43
13:I:160:THR:HG23	13:I:161:LYS:HG2	2.01	0.43
28:V:1403:UNK:C	28:V:1405:UNK:N	2.81	0.43
2:2:159:PHE:O	2:2:163:ILE:HG13	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:C:185:ASN:HB3	9:C:406:PRO:HG2	1.99	0.43
16:D:134:UNK:O	16:D:136:UNK:N	2.51	0.43
1:1:194:VAL:HG22	1:1:238:ILE:HG12	2.00	0.43
4:4:76:UNK:H	4:4:139:ILE:HD11	1.83	0.43
6:6:96:LEU:HD22	6:6:96:LEU:H	1.84	0.43
9:C:388:TYR:CD1	13:I:178:SER:HB3	2.52	0.43
6:6:71:PHE:HA	6:6:74:ILE:HG22	1.99	0.43
8:B:48:UNK:HA	8:B:132:ASP:OD1	2.18	0.43
2:2:211:LYS:O	2:2:221:LEU:HD11	2.18	0.43
2:2:310:LEU:HD13	2:2:310:LEU:HA	1.92	0.43
4:4:468:PRO:HB2	4:4:471:MET:HG2	2.00	0.43
6:6:13:ILE:H	6:6:13:ILE:HG13	1.63	0.43
6:6:27:LYS:HA	6:6:28:ASN:HA	1.42	0.43
15:L:23:ILE:H	15:L:23:ILE:HG13	1.56	0.43
30:BH:5106:UNK:O	30:BH:5108:UNK:N	2.52	0.43
2:2:287:PHE:O	2:2:291:THR:HG23	2.18	0.43
4:4:460:SER:HA	4:4:463:ILE:HD12	2.01	0.43
5:5:151:TYR:HB2	5:5:171:VAL:CG2	2.39	0.43
6:6:95:PRO:HG2	6:6:96:LEU:HD22	2.00	0.43
9:C:191:CYS:SG	9:C:206:GLY:HA3	2.59	0.43
14:K:61:LEU:HA	14:K:63:THR:N	2.31	0.43
16:D:107:UNK:HA	16:D:108:UNK:HA	1.72	0.43
45:AZ:4325:UNK:C	45:AZ:4327:UNK:H	2.30	0.43
2:2:211:LYS:HB2	2:2:244:ILE:HG21	2.01	0.43
4:4:165:PHE:CD1	4:4:165:PHE:N	2.86	0.43
5:5:175:ARG:HA	5:5:178:ASP:HB2	2.01	0.43
5:5:255:UNK:O	5:5:332:LEU:HD11	2.19	0.43
1:1:41:PHE:O	1:1:43:GLY:N	2.52	0.43
2:2:213:GLY:O	2:2:218:HIS:HA	2.19	0.43
8:B:409:ARG:O	8:B:413:MET:N	2.52	0.43
9:C:174:ARG:HH21	9:C:234:GLY:HA2	1.84	0.43
3:3:99:ILE:O	3:3:103:VAL:HG22	2.19	0.42
4:4:260:ARG:NE	4:4:260:ARG:HA	2.33	0.42
8:B:336:UNK:O	8:B:340:UNK:N	2.51	0.42
11:G:202:ASP:HA	14:K:126:LYS:HD2	2.01	0.42
20:N:519:UNK:HA	20:N:520:UNK:HA	1.84	0.42
2:2:7:ILE:HG13	2:2:7:ILE:H	1.50	0.42
2:2:399:SER:O	2:2:402:LEU:HB3	2.19	0.42
3:3:95:LEU:O	3:3:98:LEU:HB2	2.18	0.42
5:5:136:TYR:O	5:5:140:PHE:HD2	2.01	0.42
7:A:184:ILE:HG22	7:A:184:ILE:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:2:238:LEU:HD22	2:2:320:HIS:CE1	2.55	0.42
4:4:113:UNK:O	4:4:117:UNK:CB	2.68	0.42
4:4:228:HIS:HE2	5:5:580:ASP:CG	2.26	0.42
11:G:165:VAL:O	11:G:190:PHE:HA	2.19	0.42
12:H:218:UNK:O	12:H:219:UNK:C	2.66	0.42
3:3:82:LEU:HB2	3:3:83:VAL:HB	2.01	0.42
4:4:180:LEU:O	4:4:184:VAL:HG23	2.19	0.42
8:B:385:THR:HG22	8:B:388:ARG:HH12	1.83	0.42
9:C:97:VAL:HG13	9:C:98:LEU:H	1.85	0.42
9:C:337:MET:O	9:C:340:PHE:HB2	2.19	0.42
1:1:21:TYR:HE2	1:1:47:LEU:HD22	1.84	0.42
1:1:326:ILE:CG2	23:Q:933:UNK:CB	2.97	0.42
2:2:368:VAL:O	2:2:372:ILE:HG23	2.19	0.42
2:2:373:GLY:O	2:2:375:PRO:HD3	2.20	0.42
4:4:289:SER:O	4:4:292:THR:OG1	2.37	0.42
9:C:305:ILE:HB	9:C:404:GLU:HB2	2.01	0.42
2:2:74:PHE:O	2:2:78:VAL:HG23	2.19	0.42
4:4:208:THR:HA	4:4:269:ALA:HB2	2.02	0.42
9:C:311:TYR:HB2	9:C:312:ASP:H	1.73	0.42
9:C:337:MET:HA	9:C:340:PHE:CD2	2.43	0.42
11:G:123:UNK:HA	11:G:132:UNK:CB	2.49	0.42
21:O:615:UNK:C	21:O:617:UNK:N	2.83	0.42
1:1:152:SER:O	1:1:156:ILE:HG23	2.19	0.42
3:3:77:VAL:HA	3:3:88:PHE:HZ	1.85	0.42
11:G:193:HIS:HA	11:G:194:PRO:HD3	1.97	0.42
33:AC:2035:UNK:C	33:AC:2037:UNK:N	2.83	0.42
12:H:129:THR:HG21	12:H:168:CYS:HB2	2.01	0.42
29:W:1504:UNK:O	29:W:1506:UNK:N	2.53	0.42
2:2:71:LEU:O	2:2:75:ILE:HD12	2.19	0.42
4:4:214:LEU:HD23	4:4:261:LEU:HB3	2.01	0.42
6:6:63:TYR:HE2	15:L:34:LEU:HB2	1.85	0.42
2:2:140:ILE:HD13	2:2:235:TYR:HE2	1.85	0.42
2:2:153:ALA:HA	2:2:156:LEU:HB2	2.01	0.42
2:2:236:ILE:O	2:2:240:PRO:HD2	2.19	0.42
2:2:238:LEU:HD22	2:2:320:HIS:ND1	2.35	0.42
6:6:13:ILE:HG12	15:L:3:ILE:HG12	2.01	0.42
2:2:281:ILE:HD12	2:2:326:ILE:HD11	2.02	0.41
4:4:153:HIS:HB2	4:4:164:SER:OG	2.20	0.41
16:Z:55:UNK:N	16:Z:56:UNK:HA	2.35	0.41
1:1:148:LEU:HD21	3:3:70:ILE:HD13	2.01	0.41
2:2:61:PHE:HB3	2:2:64:PHE:HE2	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:6:145:THR:HA	6:6:146:LEU:HA	1.89	0.41
7:A:185:HIS:HB2	8:B:388:ARG:NH1	2.34	0.41
9:C:184:LEU:HD12	9:C:213:LEU:HB2	2.02	0.41
9:C:208:GLU:H	9:C:208:GLU:HG3	1.75	0.41
5:5:370:LEU:HB2	5:5:373:THR:OG1	2.19	0.41
7:A:188:ARG:HB2	50:A:902:SF4:S3	2.59	0.41
11:G:168:ILE:CG2	11:G:171:LEU:HB3	2.50	0.41
11:G:190:PHE:HB2	11:G:196:LEU:HD21	2.02	0.41
13:I:140:CYS:CB	13:I:144:ALA:HB2	2.50	0.41
1:1:56:LYS:HG3	14:K:72:ARG:HG2	2.02	0.41
2:2:180:ASN:HA	15:L:47:PHE:CE2	2.45	0.41
4:4:165:PHE:HD1	4:4:165:PHE:H	1.68	0.41
9:C:258:LEU:CD1	9:C:340:PHE:HB3	2.50	0.41
9:C:318:ASP:HB2	9:C:349:GLN:HE22	1.84	0.41
13:I:176:GLN:CA	13:I:184:ILE:HG23	2.46	0.41
2:2:75:ILE:HD13	2:2:242:ILE:CD1	2.50	0.41
2:2:139:LEU:HD23	6:6:179:ILE:HD12	2.01	0.41
4:4:171:THR:HG22	4:4:221:LYS:HE3	2.02	0.41
7:A:186:CYS:SG	7:A:188:ARG:HG3	2.60	0.41
9:C:216:PHE:HE1	9:C:243:LEU:HD12	1.86	0.41
9:C:343:SER:O	9:C:347:ILE:HG12	2.20	0.41
11:G:72:HIS:HA	11:G:89:PHE:HE2	1.84	0.41
21:O:647:UNK:O	21:O:651:UNK:N	2.53	0.41
1:1:27:ARG:HD2	1:1:39:PRO:HG3	2.01	0.41
2:2:209:LEU:HB3	2:2:214:ILE:HG12	2.03	0.41
11:G:190:PHE:O	11:G:196:LEU:HD11	2.21	0.41
13:I:132:ALA:HA	13:I:159:THR:HG23	2.02	0.41
3:3:91:VAL:HG13	6:6:166:ALA:HB2	2.03	0.41
4:4:468:PRO:CB	4:4:471:MET:HG2	2.51	0.41
5:5:494:UNK:O	5:5:497:UNK:O	2.38	0.41
19:M:414:UNK:C	19:M:416:UNK:N	2.82	0.41
19:M:421:UNK:O	19:M:426:UNK:N	2.54	0.41
1:1:199:ARG:HH21	1:1:231:GLU:HG2	1.86	0.41
3:3:40:UNK:HA	14:K:133:GLN:HB3	2.03	0.41
4:4:145:LEU:HD23	4:4:168:LEU:HD12	2.02	0.41
16:D:114:UNK:C	16:D:116:UNK:N	2.84	0.41
1:1:152:SER:O	1:1:155:ILE:HG13	2.21	0.41
2:2:403:ILE:HD11	4:4:180:LEU:HB2	2.01	0.41
5:5:211:LEU:O	5:5:215:ILE:HD12	2.21	0.41
6:6:18:LEU:O	6:6:21:ILE:HG22	2.20	0.41
13:I:170:ILE:O	13:I:170:ILE:HG22	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
18:J:352:UNK:C	18:J:354:UNK:N	2.84	0.41
3:3:69:GLU:HG3	6:6:169:LEU:HD22	2.02	0.41
9:C:90:GLN:O	9:C:92:PRO:HD3	2.20	0.41
15:L:71:LEU:O	15:L:75:VAL:HG23	2.21	0.41
1:1:148:LEU:HD22	1:1:317:LEU:HD11	2.03	0.40
5:5:140:PHE:CD1	5:5:181:PHE:CD2	3.09	0.40
6:6:70:LEU:HB3	15:L:71:LEU:CD2	2.50	0.40
13:I:184:ILE:O	13:I:184:ILE:CG2	2.64	0.40
2:2:247:TYR:CA	2:2:250:LEU:HD23	2.32	0.40
8:B:97:LYS:HA	8:B:100:PHE:HD2	1.86	0.40
9:C:305:ILE:HD11	9:C:409:GLU:HG3	2.03	0.40
14:K:83:LEU:HB2	14:K:121:GLY:HA3	2.04	0.40
15:L:5:THR:HA	15:L:8:LEU:HD12	2.03	0.40
2:2:130:ILE:HG13	2:2:211:LYS:HE2	2.03	0.40
2:2:142:ALA:CB	6:6:179:ILE:HD11	2.50	0.40
4:4:149:PHE:O	4:4:164:SER:OG	2.38	0.40
6:6:165:LEU:HD23	6:6:168:VAL:HG11	2.02	0.40
9:C:148:MET:H	9:C:181:THR:HG21	1.85	0.40
9:C:287:LEU:HD22	11:G:117:UNK:CB	2.51	0.40
46:BC:4609:UNK:O	46:BC:4613:UNK:N	2.54	0.40
2:2:212:ILE:H	2:2:212:ILE:HG13	1.45	0.40
4:4:79:UNK:HA	4:4:131:UNK:HA	2.04	0.40
4:4:294:ARG:HB2	5:5:574:ASN:HA	2.02	0.40
6:6:92:ARG:O	6:6:95:PRO:HD2	2.22	0.40
9:C:342:GLN:O	9:C:346:ILE:HG12	2.22	0.40
2:2:266:LEU:HD22	2:2:401:VAL:HG11	2.02	0.40
4:4:236:SER:HB2	4:4:299:LYS:HD2	2.03	0.40
4:4:373:ALA:O	4:4:377:ILE:HG12	2.22	0.40
5:5:323:SER:O	5:5:325:TYR:N	2.55	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	1	202/327 (62%)	166 (82%)	30 (15%)	6 (3%)	3	25
2	2	320/438 (73%)	278 (87%)	31 (10%)	11 (3%)	3	23
3	3	64/89 (72%)	54 (84%)	7 (11%)	3 (5%)	2	17
4	4	243/470 (52%)	212 (87%)	24 (10%)	7 (3%)	3	26
5	5	356/619 (58%)	295 (83%)	46 (13%)	15 (4%)	2	19
6	6	141/185 (76%)	118 (84%)	14 (10%)	9 (6%)	1	12
7	A	63/628 (10%)	55 (87%)	5 (8%)	3 (5%)	2	16
8	B	72/370 (20%)	67 (93%)	5 (7%)	0	100	100
9	C	348/444 (78%)	303 (87%)	27 (8%)	18 (5%)	1	15
11	G	65/133 (49%)	55 (85%)	8 (12%)	2 (3%)	3	25
12	H	30/154 (20%)	28 (93%)	1 (3%)	1 (3%)	3	24
13	I	64/137 (47%)	47 (73%)	11 (17%)	6 (9%)	0	6
14	K	141/183 (77%)	126 (89%)	12 (8%)	3 (2%)	5	31
15	L	87/89 (98%)	72 (83%)	6 (7%)	9 (10%)	0	5
All	All	2196/4266 (52%)	1876 (85%)	227 (10%)	93 (4%)	2	19

All (93) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	1	42	VAL
1	1	44	TYR
1	1	45	TYR
1	1	199	ARG
1	1	212	VAL
1	1	217	THR
2	2	43	TYR
2	2	123	PHE
2	2	218	HIS
3	3	83	VAL
4	4	158	SER
4	4	203	SER
4	4	208	THR
4	4	295	GLN
5	5	113	GLN
5	5	620	LEU
6	6	9	ILE
6	6	28	ASN
6	6	80	ILE

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Mol	Chain	Res	Type
6	6	90	ASN
9	C	97	VAL
9	C	98	LEU
9	C	239	LEU
9	C	240	PRO
9	C	294	PRO
9	C	368	SER
11	G	191	GLU
13	I	139	ILE
13	I	140	CYS
13	I	179	CYS
13	I	182	ASP
14	K	57	VAL
14	K	60	THR
14	K	203	ILE
15	L	22	ILE
15	L	50	ILE
15	L	86	SER
2	2	51	THR
2	2	62	ASN
2	2	230	ILE
2	2	327	ILE
5	5	235	TRP
5	5	324	ALA
5	5	396	LYS
5	5	649	ILE
5	5	653	LEU
6	6	149	THR
6	6	150	ILE
7	A	133	CYS
7	A	187	THR
9	C	93	ALA
9	C	327	GLY
11	G	83	PRO
13	I	81	ALA
13	I	171	TYR
15	L	87	TYR
2	2	52	TYR
2	2	300	LEU
3	3	4	PHE
4	4	202	LEU
4	4	229	VAL

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Mol	Chain	Res	Type
5	5	209	THR
5	5	323	SER
5	5	619	MET
5	5	639	PHE
6	6	81	ASN
7	A	134	ALA
9	C	269	ARG
9	C	313	ALA
9	C	362	GLU
9	C	370	PRO
15	L	48	ASP
15	L	49	ASP
15	L	85	ASN
5	5	133	GLY
5	5	346	ALA
6	6	147	LEU
9	C	302	PRO
2	2	180	ASN
3	3	8	ILE
6	6	181	MET
9	C	101	ILE
9	C	201	THR
9	C	369	PRO
15	L	17	PHE
4	4	225	PHE
5	5	641	ILE
2	2	162	GLY
12	H	130	THR
9	C	234	GLY
9	C	355	PRO
5	5	634	VAL
15	L	88	GLY

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	1	175/186 (94%)	125 (71%)	50 (29%)	0	3
2	2	291/291 (100%)	210 (72%)	81 (28%)	0	3
3	3	60/60 (100%)	37 (62%)	23 (38%)	0	1
4	4	215/215 (100%)	158 (74%)	57 (26%)	0	3
5	5	301/305 (99%)	218 (72%)	83 (28%)	0	3
6	6	110/167 (66%)	71 (64%)	39 (36%)	0	1
7	A	39/40 (98%)	36 (92%)	3 (8%)	12	38
8	B	53/56 (95%)	50 (94%)	3 (6%)	18	46
9	C	238/371 (64%)	200 (84%)	38 (16%)	2	15
11	G	55/55 (100%)	53 (96%)	2 (4%)	31	56
12	H	19/19 (100%)	18 (95%)	1 (5%)	20	48
13	I	55/55 (100%)	42 (76%)	13 (24%)	1	5
14	K	91/157 (58%)	74 (81%)	17 (19%)	1	10
15	L	69/77 (90%)	42 (61%)	27 (39%)	0	1
All	All	1771/2054 (86%)	1334 (75%)	437 (25%)	1	4

All (437) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	1	3	ILE
1	1	7	GLU
1	1	14	CYS
1	1	22	LEU
1	1	33	MET
1	1	35	ARG
1	1	42	VAL
1	1	53	ASP
1	1	133	LEU
1	1	137	ILE
1	1	144	ILE
1	1	147	GLU
1	1	150	LEU
1	1	151	THR
1	1	152	SER
1	1	155	ILE
1	1	156	ILE
1	1	157	ILE
1	1	162	SER

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Mol	Chain	Res	Type
1	1	164	LEU
1	1	169	ILE
1	1	193	SER
1	1	194	VAL
1	1	196	GLU
1	1	197	THR
1	1	199	ARG
1	1	204	LEU
1	1	206	GLU
1	1	208	GLU
1	1	211	LEU
1	1	215	TYR
1	1	217	THR
1	1	225	VAL
1	1	229	LEU
1	1	241	PHE
1	1	247	LEU
1	1	276	LEU
1	1	277	ILE
1	1	278	ASN
1	1	279	SER
1	1	284	ILE
1	1	287	VAL
1	1	289	LEU
1	1	291	PHE
1	1	295	TRP
1	1	312	CYS
1	1	319	LEU
1	1	321	PHE
1	1	323	ILE
1	1	326	ILE
2	2	1	MET
2	2	2	LEU
2	2	3	ILE
2	2	6	ILE
2	2	7	ILE
2	2	8	SER
2	2	16	SER
2	2	17	LYS
2	2	18	LEU
2	2	25	ILE
2	2	45	LEU

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Mol	Chain	Res	Type
2	2	48	ASN
2	2	49	ASN
2	2	58	LEU
2	2	65	THR
2	2	68	ILE
2	2	82	LEU
2	2	125	THR
2	2	132	LEU
2	2	138	TYR
2	2	141	THR
2	2	143	ILE
2	2	145	ASN
2	2	156	LEU
2	2	164	LEU
2	2	166	ILE
2	2	171	SER
2	2	172	ILE
2	2	173	ASN
2	2	174	THR
2	2	178	VAL
2	2	187	LEU
2	2	188	ASP
2	2	192	ILE
2	2	197	LEU
2	2	204	LEU
2	2	212	ILE
2	2	219	LYS
2	2	224	ILE
2	2	230	ILE
2	2	232	ILE
2	2	233	THR
2	2	238	LEU
2	2	239	ILE
2	2	241	LYS
2	2	242	ILE
2	2	250	LEU
2	2	261	SER
2	2	267	THR
2	2	269	LEU
2	2	276	LEU
2	2	277	LEU
2	2	278	GLN

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Mol	Chain	Res	Type
2	2	280	LYS
2	2	281	ILE
2	2	282	LYS
2	2	283	ARG
2	2	288	SER
2	2	290	LEU
2	2	298	LEU
2	2	300	LEU
2	2	304	ASN
2	2	306	GLU
2	2	307	PHE
2	2	310	LEU
2	2	311	TYR
2	2	314	THR
2	2	319	SER
2	2	321	LEU
2	2	328	ILE
2	2	362	LEU
2	2	363	SER
2	2	367	VAL
2	2	368	VAL
2	2	370	SER
2	2	371	PHE
2	2	372	ILE
2	2	377	LEU
2	2	378	LEU
2	2	384	LEU
2	2	387	LEU
3	3	2	ASN
3	3	5	ILE
3	3	6	ILE
3	3	7	PHE
3	3	8	ILE
3	3	57	ILE
3	3	58	LEU
3	3	68	LEU
3	3	70	ILE
3	3	73	LEU
3	3	74	LEU
3	3	77	VAL
3	3	78	MET
3	3	81	TYR

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Mol	Chain	Res	Type
3	3	82	LEU
3	3	83	VAL
3	3	85	ASN
3	3	88	PHE
3	3	90	ILE
3	3	97	ILE
3	3	98	LEU
3	3	99	ILE
3	3	102	PHE
4	4	136	SER
4	4	139	ILE
4	4	144	THR
4	4	150	ILE
4	4	169	MET
4	4	172	LEU
4	4	180	LEU
4	4	182	ILE
4	4	185	ILE
4	4	201	VAL
4	4	203	SER
4	4	205	ASP
4	4	210	ILE
4	4	216	ILE
4	4	227	ILE
4	4	231	LEU
4	4	233	VAL
4	4	237	GLU
4	4	251	LEU
4	4	259	LEU
4	4	260	ARG
4	4	263	LEU
4	4	265	LEU
4	4	267	CYS
4	4	276	MET
4	4	280	ILE
4	4	281	SER
4	4	282	LEU
4	4	286	ILE
4	4	287	LEU
4	4	296	ILE
4	4	298	LEU
4	4	299	LYS

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Mol	Chain	Res	Type
4	4	301	ILE
4	4	302	ILE
4	4	306	SER
4	4	307	ILE
4	4	317	VAL
4	4	320	ASN
4	4	323	LEU
4	4	325	ILE
4	4	372	LEU
4	4	376	ILE
4	4	386	THR
4	4	397	LEU
4	4	399	LEU
4	4	404	ILE
4	4	405	ARG
4	4	454	MET
4	4	455	ASN
4	4	457	LEU
4	4	458	ILE
4	4	459	ILE
4	4	461	THR
4	4	464	ILE
4	4	466	ILE
4	4	467	CYS
5	5	81	ILE
5	5	82	ASP
5	5	84	LEU
5	5	86	ILE
5	5	88	MET
5	5	89	LEU
5	5	103	SER
5	5	104	ILE
5	5	108	GLU
5	5	112	HIS
5	5	114	VAL
5	5	116	PHE
5	5	120	LEU
5	5	124	THR
5	5	126	TRP
5	5	136	TYR
5	5	137	PHE
5	5	138	VAL

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Mol	Chain	Res	Type
5	5	139	LEU
5	5	143	TRP
5	5	145	PHE
5	5	160	ARG
5	5	161	LEU
5	5	162	GLN
5	5	168	LEU
5	5	176	PHE
5	5	181	PHE
5	5	185	LEU
5	5	188	ILE
5	5	191	VAL
5	5	209	THR
5	5	210	ASP
5	5	212	LEU
5	5	215	ILE
5	5	216	MET
5	5	217	LEU
5	5	230	PHE
5	5	233	HIS
5	5	234	ASN
5	5	236	LEU
5	5	302	LYS
5	5	313	LEU
5	5	315	MET
5	5	318	ILE
5	5	320	ILE
5	5	322	LEU
5	5	327	LEU
5	5	329	LEU
5	5	331	HIS
5	5	332	LEU
5	5	342	LEU
5	5	383	SER
5	5	390	LEU
5	5	394	TYR
5	5	399	ILE
5	5	400	ILE
5	5	418	TYR
5	5	432	LEU
5	5	434	LEU
5	5	460	LEU

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Mol	Chain	Res	Type
5	5	462	MET
5	5	471	PHE
5	5	474	TRP
5	5	578	HIS
5	5	579	VAL
5	5	580	ASP
5	5	584	LEU
5	5	587	LEU
5	5	589	PRO
5	5	616	MET
5	5	619	MET
5	5	621	ILE
5	5	622	LEU
5	5	623	ILE
5	5	628	LEU
5	5	629	LEU
5	5	630	LEU
5	5	636	ASN
5	5	637	VAL
5	5	640	ILE
5	5	643	ILE
5	5	645	VAL
5	5	653	LEU
6	6	11	ILE
6	6	13	ILE
6	6	15	LEU
6	6	23	ILE
6	6	31	VAL
6	6	41	VAL
6	6	42	ILE
6	6	58	LEU
6	6	59	TYR
6	6	67	ILE
6	6	70	LEU
6	6	72	LEU
6	6	74	ILE
6	6	77	LEU
6	6	80	ILE
6	6	83	THR
6	6	98	LEU
6	6	100	SER
6	6	102	ILE

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Mol	Chain	Res	Type
6	6	103	VAL
6	6	104	LEU
6	6	109	LEU
6	6	112	TYR
6	6	149	THR
6	6	150	ILE
6	6	154	LEU
6	6	155	LEU
6	6	156	THR
6	6	161	ILE
6	6	167	ILE
6	6	168	VAL
6	6	169	LEU
6	6	170	LEU
6	6	171	LEU
6	6	173	ILE
6	6	174	ILE
6	6	177	ILE
6	6	179	ILE
6	6	180	THR
7	A	124	MET
7	A	142	CYS
7	A	186	CYS
8	B	388	ARG
8	B	427	CYS
8	B	429	LEU
9	C	102	LEU
9	C	133	MET
9	C	144	TYR
9	C	145	VAL
9	C	147	MET
9	C	162	LEU
9	C	176	MET
9	C	180	ILE
9	C	186	HIS
9	C	188	MET
9	C	189	SER
9	C	200	LEU
9	C	201	THR
9	C	204	LEU
9	C	207	PHE
9	C	208	GLU

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Mol	Chain	Res	Type
9	C	214	MET
9	C	216	PHE
9	C	217	TYR
9	C	247	ILE
9	C	248	TYR
9	C	257	ARG
9	C	266	THR
9	C	278	ILE
9	C	284	GLN
9	C	285	ASP
9	C	297	ARG
9	C	308	ASN
9	C	311	TYR
9	C	344	LEU
9	C	351	CYS
9	C	359	VAL
9	C	377	ASP
9	C	387	LEU
9	C	425	CYS
9	C	453	ILE
9	C	454	ILE
9	C	460	VAL
11	G	84	LYS
11	G	177	TRP
12	H	130	THR
13	I	78	PHE
13	I	82	GLU
13	I	83	MET
13	I	84	PHE
13	I	130	CYS
13	I	137	GLU
13	I	139	ILE
13	I	143	LEU
13	I	164	ILE
13	I	172	CYS
13	I	181	VAL
13	I	190	VAL
13	I	192	TYR
14	K	73	GLN
14	K	79	VAL
14	K	81	PHE
14	K	85	CYS

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Mol	Chain	Res	Type
14	K	112	ARG
14	K	123	LEU
14	K	124	THR
14	K	131	LEU
14	K	140	GLU
14	K	170	ILE
14	K	177	VAL
14	K	180	CYS
14	K	183	THR
14	K	191	VAL
14	K	193	GLN
14	K	194	LEU
14	K	196	ARG
15	L	2	PHE
15	L	3	ILE
15	L	5	THR
15	L	6	ILE
15	L	10	LEU
15	L	19	ARG
15	L	20	ARG
15	L	23	ILE
15	L	24	LEU
15	L	27	ILE
15	L	28	CYS
15	L	29	LEU
15	L	34	LEU
15	L	36	ILE
15	L	39	ILE
15	L	41	LEU
15	L	46	LEU
15	L	48	ASP
15	L	49	ASP
15	L	51	SER
15	L	57	ILE
15	L	62	LEU
15	L	69	ILE
15	L	71	LEU
15	L	74	LEU
15	L	79	ARG
15	L	84	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	1	242	ASN
2	2	62	ASN
2	2	145	ASN
2	2	173	ASN
2	2	218	HIS
2	2	227	ASN
2	2	304	ASN
2	2	305	ASN
2	2	315	GLN
4	4	400	GLN
4	4	406	ASN
4	4	455	ASN
5	5	233	HIS
5	5	636	ASN
6	6	114	ASN
7	A	185	HIS
9	C	90	GLN
9	C	445	HIS
13	I	189	ASN
14	K	113	GLN
15	L	37	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

8 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The

Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
50	SF4	I	501	13	0,12,12	-	-	-		
50	SF4	I	500	13	0,12,12	-	-	-		
50	SF4	K	500	14	0,12,12	-	-	-		
49	FES	A	900	7	0,4,4	-	-	-		
50	SF4	B	500	8	0,12,12	-	-	-		
49	FES	H	300	12	0,4,4	-	-	-		
50	SF4	A	901	7	0,12,12	-	-	-		
50	SF4	A	902	7	0,12,12	-	-	-		

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
50	SF4	I	501	13	-	-	0/6/5/5
50	SF4	I	500	13	-	-	0/6/5/5
50	SF4	K	500	14	-	-	0/6/5/5
49	FES	A	900	7	-	-	0/1/1/1
50	SF4	B	500	8	-	-	0/6/5/5
49	FES	H	300	12	-	-	0/1/1/1
50	SF4	A	901	7	-	-	0/6/5/5
50	SF4	A	902	7	-	-	0/6/5/5

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 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
50	I	501	SF4	1	0
50	I	500	SF4	1	0
50	K	500	SF4	2	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
49	H	300	FES	1	0
50	A	902	SF4	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
7	A	30
8	B	11
10	E	10
5	5	8
13	I	5
12	H	5
2	2	4
4	4	4
1	1	3
3	3	2
11	G	2
22	P	1
16	Z	1
37	AL	1
21	O	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	E	422:UNK	C	499:UNK	N	48.62
1	B	270:UNK	C	279:UNK	N	26.71
1	E	513:UNK	C	604:UNK	N	26.57
1	A	148:UNK	C	152:UNK	N	24.97
1	I	92:UNK	C	103:UNK	N	24.67
1	3	21:UNK	C	38:UNK	N	22.87
1	E	325:UNK	C	404:UNK	N	22.86
1	2	422:UNK	C	446:UNK	N	21.69
1	A	358:UNK	C	360:UNK	N	17.83
1	E	637:UNK	C	785:UNK	N	17.21

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Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	H	121:UNK	C	127:CYS	N	16.50
1	E	65:UNK	C	83:UNK	N	16.40
1	E	109:UNK	C	161:UNK	N	16.24
1	B	109:UNK	C	115:UNK	N	16.14
1	A	560:UNK	C	565:UNK	N	16.08
1	E	180:UNK	C	235:UNK	N	15.93
1	I	212:UNK	C	233:UNK	N	15.00
1	B	195:UNK	C	201:UNK	N	14.93
1	E	795:UNK	C	887:UNK	N	14.77
1	H	174:UNK	C	181:UNK	N	14.49
1	A	510:UNK	C	515:UNK	N	12.92
1	A	81:UNK	C	84:UNK	N	12.79
1	A	363:UNK	C	367:UNK	N	12.10
1	3	44:UNK	C	51:UNK	N	12.08
1	B	343:UNK	C	356:UNK	N	11.74
1	A	642:UNK	C	647:UNK	N	11.45
1	H	109:UNK	C	117:UNK	N	11.18
1	A	708:UNK	C	712:UNK	N	10.79
1	H	90:UNK	C	94:UNK	N	10.68
1	4	52:UNK	C	57:UNK	N	10.67
1	G	138:UNK	C	145:UNK	N	10.57
1	2	84:GLY	C	97:UNK	N	10.44
1	2	255:UNK	C	256:UNK	N	10.18
1	G	125:UNK	C	128:UNK	N	9.87
1	B	155:UNK	C	159:UNK	N	9.83
1	4	67:UNK	C	76:UNK	N	9.82
1	B	120:UNK	C	125:UNK	N	9.82
1	A	414:UNK	C	416:UNK	N	9.79
1	B	245:UNK	C	251:UNK	N	9.67
1	P	727:UNK	C	728:UNK	N	9.66
1	I	153:UNK	C	157:UNK	N	9.61
1	5	297:UNK	C	301:UNK	N	9.57
1	H	162:UNK	C	165:UNK	N	9.53
1	A	530:UNK	C	534:UNK	N	9.36
1	A	427:UNK	C	432:UNK	N	9.27
1	A	541:UNK	C	543:UNK	N	9.20
1	I	120:UNK	C	128:UNK	N	9.06
1	A	127:ALA	C	131:ALA	N	8.89
1	A	457:UNK	C	464:UNK	N	8.88
1	A	435:UNK	C	438:UNK	N	8.86
1	B	87:UNK	C	92:UNK	N	8.74
1	4	18:UNK	C	22:UNK	N	8.73

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Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	E	242:UNK	C	316:UNK	N	8.00
1	1	120:UNK	C	129:UNK	N	7.81
1	A	704:UNK	C	707:UNK	N	7.59
1	1	62:UNK	C	66:UNK	N	7.54
1	A	87:UNK	C	89:UNK	N	7.53
1	5	24:UNK	C	27:UNK	N	7.47
1	Z	17:UNK	C	18:UNK	N	7.27
1	A	492:UNK	C	496:UNK	N	7.13
1	I	235:UNK	C	238:UNK	N	7.03
1	A	273:UNK	C	276:UNK	N	6.99
1	4	435:UNK	C	437:UNK	N	6.95
1	A	293:UNK	C	297:UNK	N	6.62
1	A	257:UNK	C	260:UNK	N	6.51
1	A	246:UNK	C	248:UNK	N	6.34
1	A	408:UNK	C	409:UNK	N	6.33
1	A	580:UNK	C	586:UNK	N	6.31
1	5	604:UNK	C	613:UNK	N	6.10
1	B	303:UNK	C	306:UNK	N	5.89
1	A	236:GLY	C	240:UNK	N	5.83
1	5	447:UNK	C	455:UNK	N	5.76
1	A	181:UNK	C	183:CYS	N	5.74
1	A	309:UNK	C	313:UNK	N	5.28
1	E	57:UNK	C	59:UNK	N	5.13
1	5	437:UNK	C	439:UNK	N	5.04
1	B	406:UNK	C	407:UNK	N	4.53
1	B	238:UNK	C	240:UNK	N	4.43
1	AL	2916:UNK	C	2917:UNK	N	4.25
1	2	332:UNK	C	339:UNK	N	4.23
1	5	412:UNK	C	416:UNK	N	4.21
1	1	173:UNK	C	177:UNK	N	4.18
1	A	443:UNK	C	445:UNK	N	3.87
1	5	205:UNK	C	208:UNK	N	3.82
1	5	550:UNK	C	560:UNK	N	3.79
1	O	639:UNK	C	640:UNK	N	3.76
1	A	397:UNK	C	400:UNK	N	3.56
1	A	334:UNK	C	336:UNK	N	3.40

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	1	204/327 (62%)	1.01	30 (14%) 6 5	17, 70, 108, 141	0
2	2	322/438 (73%)	0.55	25 (7%) 19 13	11, 31, 76, 95	0
3	3	65/89 (73%)	0.95	10 (15%) 5 5	28, 57, 128, 142	0
4	4	243/470 (51%)	0.76	23 (9%) 14 10	15, 48, 76, 96	0
5	5	357/619 (57%)	1.37	82 (22%) 2 2	54, 87, 120, 138	0
6	6	149/185 (80%)	0.53	18 (12%) 8 7	12, 60, 134, 202	0
7	A	67/628 (10%)	3.87	52 (77%) 0 0	105, 183, 300, 300	0
8	B	72/370 (19%)	3.95	54 (75%) 0 0	124, 249, 300, 300	0
9	C	358/444 (80%)	1.81	120 (33%) 1 1	70, 112, 228, 240	0
10	E	0/195	-	-	-	-
11	G	66/133 (49%)	2.64	39 (59%) 0 0	93, 126, 243, 297	0
12	H	31/154 (20%)	1.93	12 (38%) 1 1	133, 170, 246, 298	0
13	I	64/137 (46%)	2.77	40 (62%) 0 0	96, 128, 182, 201	0
14	K	147/183 (80%)	1.87	63 (42%) 0 1	78, 106, 255, 284	0
15	L	89/89 (100%)	0.56	6 (6%) 24 15	16, 45, 105, 118	0
16	D	0/57	-	-	-	-
16	Z	0/57	-	-	-	-
17	F	0/54	-	-	-	-
18	J	0/63	-	-	-	-
19	M	0/29	-	-	-	-
20	N	0/50	-	-	-	-
21	O	0/70	-	-	-	-
22	P	0/46	-	-	-	-
23	Q	0/51	-	-	-	-
24	R	0/30	-	-	-	-
25	S	0/69	-	-	-	-
26	AH	0/15	-	-	-	-

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
26	T	0/15	-	-	-	-
27	U	0/26	-	-	-	-
28	V	0/22	-	-	-	-
29	AB	0/9	-	-	-	-
29	AY	0/9	-	-	-	-
29	BE	0/9	-	-	-	-
29	W	0/9	-	-	-	-
30	AJ	0/16	-	-	-	-
30	AV	0/16	-	-	-	-
30	BH	0/16	-	-	-	-
30	X	0/16	-	-	-	-
31	AR	0/13	-	-	-	-
31	AT	0/13	-	-	-	-
31	Y	0/13	-	-	-	-
32	AA	0/18	-	-	-	-
32	AW	0/18	-	-	-	-
32	BB	0/18	-	-	-	-
32	BG	0/18	-	-	-	-
33	AC	0/47	-	-	-	-
33	AD	0/47	-	-	-	-
34	AE	0/48	-	-	-	-
35	AF	0/35	-	-	-	-
36	AG	0/25	-	-	-	-
37	AI	0/36	-	-	-	-
37	AL	0/36	-	-	-	-
38	AK	0/76	-	-	-	-
38	AN	0/76	-	-	-	-
39	AM	0/17	-	-	-	-
40	AO	0/32	-	-	-	-
41	AP	0/11	-	-	-	-
41	AS	0/11	-	-	-	-
42	AQ	0/8	-	-	-	-
42	BA	0/8	-	-	-	-
43	AU	0/58	-	-	-	-
44	AX	0/39	-	-	-	-
45	AZ	0/40	-	-	-	-
46	BC	0/20	-	-	-	-
47	BD	0/19	-	-	-	-
47	BF	0/19	-	-	-	-
48	BI	0/905	-	-	-	-
All	All	2234/6939 (32%)	1.40	574 (25%) 1 2	11, 84, 227, 300	0

All (574) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
8	B	389	GLU	11.7
7	A	135	ALA	11.5
7	A	131	ALA	11.3
7	A	122	MET	10.9
7	A	185	HIS	10.3
8	B	134	HIS	10.0
13	I	137	GLU	9.6
13	I	165	ASP	9.0
8	B	416	GLU	9.0
7	A	134	ALA	9.0
8	B	145	ARG	8.4
11	G	166	PRO	8.2
8	B	411	ILE	8.1
9	C	122	GLY	8.0
7	A	107	VAL	7.6
7	A	127	ALA	7.5
8	B	415	TYR	7.5
14	K	122	THR	7.4
5	5	585	TYR	7.3
11	G	77	TYR	7.2
7	A	125	MET	7.0
14	K	158	HIS	6.9
8	B	138	GLU	6.9
7	A	123	GLU	6.8
8	B	410	GLU	6.8
12	H	172	CYS	6.8
8	B	98	TRP	6.8
8	B	96	LEU	6.8
9	C	333	TYR	6.8
9	C	383	HIS	6.7
9	C	393	SER	6.6
8	B	382	GLY	6.5
9	C	384	HIS	6.5
8	B	414	LEU	6.4
5	5	191	VAL	6.3
8	B	132	ASP	6.1
7	A	94	VAL	6.0
2	2	44	LEU	5.9
11	G	203	TYR	5.9
11	G	179	GLU	5.9
8	B	383	GLN	5.9
5	5	460	LEU	5.9
13	I	160	THR	5.9

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Mol	Chain	Res	Type	RSRZ
1	1	45	TYR	5.8
5	5	353	ILE	5.7
8	B	381	CYS	5.7
11	G	72	HIS	5.6
9	C	375	LYS	5.5
13	I	191	GLU	5.5
7	A	191	ARG	5.5
7	A	188	ARG	5.5
7	A	197	ALA	5.4
7	A	98	ALA	5.4
9	C	413	TYR	5.4
8	B	386	PRO	5.3
5	5	397	ASP	5.3
8	B	392	THR	5.3
6	6	98	LEU	5.2
8	B	409	ARG	5.2
14	K	82	GLY	5.1
8	B	408	GLU	5.1
9	C	387	LEU	5.1
5	5	617	ASN	5.1
5	5	142	GLY	5.1
1	1	160	PHE	5.1
1	1	210	GLU	5.1
5	5	393	TYR	5.1
8	B	388	ARG	5.1
7	A	132	ALA	5.0
11	G	76	LYS	5.0
13	I	86	GLY	5.0
5	5	251	HIS	5.0
13	I	186	GLU	5.0
13	I	181	VAL	5.0
8	B	100	PHE	4.9
14	K	65	ASP	4.9
11	G	189	PHE	4.9
5	5	211	LEU	4.9
1	1	278	ASN	4.9
9	C	144	TYR	4.9
9	C	388	TYR	4.9
9	C	389	THR	4.9
9	C	302	PRO	4.9
2	2	315	GLN	4.9
11	G	173	GLU	4.9

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Mol	Chain	Res	Type	RSRZ
7	A	183	CYS	4.9
3	3	100	ILE	4.9
5	5	307	LEU	4.8
14	K	143	TRP	4.8
9	C	365	LYS	4.8
7	A	126	LEU	4.8
8	B	140	CYS	4.8
11	G	197	ARG	4.8
9	C	89	PRO	4.7
9	C	394	VAL	4.7
9	C	390	LYS	4.7
9	C	319	PHE	4.7
11	G	181	GLU	4.7
13	I	168	LYS	4.7
7	A	184	ILE	4.7
12	H	135	CYS	4.6
4	4	387	PRO	4.6
9	C	296	LEU	4.6
8	B	417	LEU	4.6
13	I	171	TYR	4.6
8	B	418	THR	4.5
5	5	140	PHE	4.5
7	A	106	VAL	4.5
11	G	87	GLN	4.5
14	K	62	THR	4.5
13	I	138	ALA	4.5
5	5	649	ILE	4.5
5	5	394	TYR	4.5
11	G	202	ASP	4.4
13	I	134	LYS	4.4
8	B	428	ALA	4.4
14	K	66	ALA	4.4
9	C	204	LEU	4.4
8	B	385	THR	4.4
14	K	60	THR	4.4
6	6	112	TYR	4.3
1	1	297	ARG	4.3
9	C	448	PRO	4.3
7	A	124	MET	4.3
8	B	149	ALA	4.3
8	B	131	LYS	4.3
11	G	183	TYR	4.3

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Mol	Chain	Res	Type	RSRZ
9	C	379	GLU	4.3
13	I	143	LEU	4.3
9	C	391	GLY	4.2
7	A	121	VAL	4.2
9	C	248	TYR	4.2
7	A	103	PRO	4.2
11	G	204	GLY	4.2
9	C	308	ASN	4.2
7	A	97	CYS	4.2
9	C	106	GLY	4.1
13	I	182	ASP	4.1
7	A	138	ALA	4.1
8	B	413	MET	4.1
8	B	151	ALA	4.1
9	C	373	LEU	4.1
13	I	162	TYR	4.1
9	C	315	ASP	4.1
5	5	107	MET	4.1
8	B	142	LEU	4.1
12	H	134	LEU	4.1
9	C	271	TRP	4.1
8	B	432	ALA	4.1
14	K	84	ALA	4.1
2	2	279	ILE	4.1
14	K	170	ILE	4.1
5	5	310	MET	4.0
9	C	382	ILE	4.0
9	C	392	TYR	4.0
4	4	300	VAL	4.0
14	K	85	CYS	4.0
7	A	227	GLY	4.0
5	5	250	LEU	4.0
5	5	321	GLY	4.0
7	A	139	GLY	4.0
7	A	142	CYS	4.0
2	2	293	ALA	4.0
15	L	22	ILE	4.0
5	5	395	THR	4.0
12	H	130	THR	4.0
2	2	211	LYS	3.9
9	C	374	MET	3.9
9	C	409	GLU	3.9

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Mol	Chain	Res	Type	RSRZ
9	C	363	ASP	3.9
5	5	243	PRO	3.9
8	B	150	THR	3.9
5	5	336	ALA	3.9
5	5	103	SER	3.9
14	K	55	SER	3.9
2	2	190	LEU	3.9
6	6	115	ASP	3.9
9	C	304	ASP	3.9
8	B	99	SER	3.9
9	C	402	ALA	3.9
9	C	386	LEU	3.8
13	I	176	GLN	3.8
3	3	76	TYR	3.8
5	5	618	SER	3.8
11	G	198	ARG	3.8
9	C	208	GLU	3.8
9	C	378	MET	3.8
8	B	141	LEU	3.8
11	G	168	ILE	3.8
4	4	162	ARG	3.8
7	A	96	SER	3.7
9	C	465	ASP	3.7
1	1	40	ASN	3.7
11	G	175	ALA	3.7
12	H	171	ALA	3.7
14	K	115	ASP	3.7
7	A	78	GLY	3.7
11	G	206	GLU	3.7
4	4	406	ASN	3.7
9	C	126	LEU	3.7
9	C	381	LEU	3.7
14	K	99	ASP	3.7
14	K	179	GLY	3.7
9	C	355	PRO	3.7
11	G	74	TYR	3.7
2	2	273	VAL	3.6
9	C	385	PHE	3.6
1	1	211	LEU	3.6
2	2	299	LEU	3.6
8	B	391	THR	3.6
14	K	159	PHE	3.6

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Mol	Chain	Res	Type	RSRZ
8	B	434	ALA	3.6
9	C	358	ALA	3.6
5	5	132	THR	3.6
9	C	125	LYS	3.6
7	A	136	CYS	3.6
2	2	412	TYR	3.6
14	K	64	LEU	3.6
8	B	147	MET	3.6
9	C	107	GLU	3.6
9	C	403	ILE	3.5
11	G	89	PHE	3.5
6	6	94	LEU	3.5
9	C	270	ILE	3.5
1	1	277	ILE	3.5
7	A	140	GLY	3.5
8	B	144	GLY	3.5
14	K	133	GLN	3.5
1	1	212	VAL	3.5
9	C	334	LEU	3.5
9	C	173	ILE	3.5
5	5	387	MET	3.5
11	G	196	LEU	3.5
7	A	133	CYS	3.5
14	K	77	TRP	3.4
9	C	404	GLU	3.4
1	1	209	SER	3.4
14	K	183	THR	3.4
13	I	141	PRO	3.4
11	G	193	HIS	3.4
9	C	105	SER	3.4
4	4	240	LEU	3.4
5	5	340	ALA	3.4
7	A	105	MET	3.4
14	K	52	LYS	3.4
8	B	433	ALA	3.4
1	1	19	VAL	3.4
3	3	57	ILE	3.4
4	4	205	ASP	3.4
2	2	309	TYR	3.4
7	A	104	GLY	3.3
7	A	187	THR	3.3
9	C	331	ASP	3.3

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Mol	Chain	Res	Type	RSRZ
1	1	20	ALA	3.3
2	2	15	MET	3.3
3	3	98	LEU	3.3
7	A	79	ASN	3.3
14	K	69	ASN	3.3
5	5	425	CYS	3.3
9	C	123	THR	3.3
9	C	290	GLY	3.3
9	C	265	LEU	3.3
9	C	305	ILE	3.3
11	G	209	PRO	3.3
9	C	367	ASN	3.3
14	K	137	GLN	3.3
5	5	423	LEU	3.2
8	B	424	HIS	3.2
9	C	141	ARG	3.2
14	K	67	VAL	3.2
5	5	372	TYR	3.2
5	5	116	PHE	3.2
8	B	129	MET	3.2
1	1	38	GLY	3.2
5	5	337	PHE	3.2
13	I	131	ILE	3.2
7	A	95	ALA	3.2
14	K	117	MET	3.2
2	2	274	GLY	3.2
4	4	324	GLY	3.2
8	B	420	ASP	3.1
1	1	30	LEU	3.1
14	K	162	SER	3.1
11	G	195	ASP	3.1
7	A	93	PRO	3.1
14	K	145	ILE	3.1
14	K	76	PHE	3.1
6	6	48	TYR	3.1
1	1	218	GLU	3.1
5	5	318	ILE	3.1
14	K	83	LEU	3.1
14	K	140	GLU	3.1
7	A	137	ALA	3.1
9	C	359	VAL	3.1
7	A	236	GLY	3.1

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Mol	Chain	Res	Type	RSRZ
13	I	136	CYS	3.1
14	K	81	PHE	3.0
9	C	406	PRO	3.0
7	A	195	ASP	3.0
5	5	352	SER	3.0
1	1	41	PHE	3.0
7	A	192	PHE	3.0
13	I	189	ASN	3.0
2	2	11	THR	3.0
9	C	253	GLN	3.0
11	G	176	ASN	3.0
11	G	205	PHE	3.0
5	5	354	LEU	3.0
8	B	143	ALA	3.0
13	I	81	ALA	3.0
13	I	174	TYR	3.0
9	C	85	ILE	3.0
5	5	159	THR	3.0
5	5	315	MET	3.0
7	A	271	VAL	3.0
9	C	230	VAL	3.0
5	5	466	ALA	3.0
8	B	148	ASN	2.9
13	I	195	GLU	2.9
4	4	366	THR	2.9
9	C	275	THR	2.9
9	C	361	VAL	2.9
14	K	56	ALA	2.9
5	5	574	ASN	2.9
9	C	262	GLU	2.9
5	5	382	LEU	2.9
1	1	275	GLY	2.9
9	C	366	ILE	2.9
1	1	32	TYR	2.9
4	4	365	LEU	2.9
5	5	384	LEU	2.9
13	I	192	TYR	2.9
9	C	348	GLU	2.9
9	C	99	ARG	2.9
11	G	169	THR	2.9
11	G	80	ALA	2.9
12	H	72	MET	2.9

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Mol	Chain	Res	Type	RSRZ
14	K	59	TYR	2.9
14	K	185	GLU	2.9
13	I	190	VAL	2.9
1	1	50	ALA	2.9
9	C	380	ALA	2.9
9	C	364	PHE	2.9
5	5	88	MET	2.9
5	5	418	TYR	2.9
9	C	134	GLN	2.8
12	H	153	GLY	2.8
13	I	170	ILE	2.8
5	5	327	LEU	2.8
1	1	44	TYR	2.8
9	C	352	ASN	2.8
5	5	389	GLY	2.8
8	B	390	GLY	2.8
14	K	120	ALA	2.8
9	C	446	PHE	2.8
14	K	53	PRO	2.8
4	4	325	ILE	2.8
11	G	78	ILE	2.8
4	4	469	GLN	2.8
9	C	184	LEU	2.8
13	I	139	ILE	2.8
14	K	80	THR	2.8
14	K	58	ASP	2.8
14	K	126	LYS	2.8
4	4	212	LEU	2.8
9	C	430	PRO	2.8
5	5	113	GLN	2.8
9	C	420	GLU	2.8
8	B	146	ALA	2.8
9	C	429	ALA	2.8
9	C	454	ILE	2.7
5	5	108	GLU	2.7
5	5	641	ILE	2.7
5	5	157	TRP	2.7
6	6	31	VAL	2.7
7	A	272	MET	2.7
2	2	157	TYR	2.7
9	C	206	GLY	2.7
14	K	63	THR	2.7

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Mol	Chain	Res	Type	RSRZ
13	I	161	LYS	2.7
9	C	232	PRO	2.7
9	C	162	LEU	2.7
11	G	85	TYR	2.7
8	B	133	PRO	2.7
12	H	133	GLN	2.7
9	C	137	PRO	2.7
7	A	198	GLY	2.7
9	C	298	GLY	2.7
9	C	418	GLY	2.7
5	5	254	THR	2.7
9	C	218	GLU	2.7
9	C	260	GLU	2.7
14	K	90	MET	2.7
5	5	570	VAL	2.6
8	B	422	GLU	2.6
9	C	124	GLU	2.6
14	K	125	ASN	2.6
5	5	376	CYS	2.6
4	4	309	HIS	2.6
9	C	138	TYR	2.6
14	K	89	GLU	2.6
5	5	242	GLY	2.6
11	G	75	ALA	2.6
5	5	635	MET	2.6
6	6	96	LEU	2.6
9	C	132	TYR	2.6
13	I	135	LEU	2.6
7	A	224	GLY	2.6
13	I	88	TYR	2.6
2	2	189	SER	2.6
15	L	50	ILE	2.6
4	4	235	HIS	2.6
5	5	463	PHE	2.6
7	A	189	CYS	2.6
6	6	152	ASN	2.6
5	5	462	MET	2.6
11	G	73	GLN	2.6
14	K	78	PRO	2.6
3	3	102	PHE	2.6
5	5	150	SER	2.5
5	5	637	VAL	2.5

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Mol	Chain	Res	Type	RSRZ
14	K	88	VAL	2.5
13	I	163	ASP	2.5
4	4	204	LEU	2.5
6	6	155	LEU	2.5
5	5	224	MET	2.5
1	1	213	ALA	2.5
14	K	100	GLN	2.5
1	1	139	SER	2.5
4	4	395	GLU	2.5
5	5	590	VAL	2.5
14	K	118	ILE	2.5
9	C	303	PHE	2.5
6	6	95	PRO	2.5
2	2	294	GLY	2.5
4	4	390	GLY	2.5
9	C	371	ARG	2.5
5	5	153	LEU	2.5
2	2	54	VAL	2.5
3	3	91	VAL	2.5
5	5	426	VAL	2.5
5	5	619	MET	2.5
13	I	83	MET	2.5
3	3	101	GLY	2.5
4	4	323	LEU	2.5
9	C	401	THR	2.5
2	2	138	TYR	2.5
9	C	87	PHE	2.5
9	C	407	LYS	2.5
14	K	202	LYS	2.5
7	A	141	ALA	2.5
14	K	128	ALA	2.5
3	3	62	LEU	2.5
13	I	183	ALA	2.4
13	I	87	LEU	2.4
9	C	246	ASP	2.4
6	6	34	LEU	2.4
4	4	242	GLY	2.4
12	H	170	GLY	2.4
5	5	171	VAL	2.4
14	K	57	VAL	2.4
9	C	169	ARG	2.4
14	K	109	ALA	2.4

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Mol	Chain	Res	Type	RSRZ
14	K	163	VAL	2.4
9	C	133	MET	2.4
7	A	120	ASN	2.4
6	6	153	VAL	2.4
15	L	26	PHE	2.4
15	L	1	MET	2.4
5	5	636	ASN	2.4
2	2	43	TYR	2.4
1	1	324	PHE	2.3
8	B	426	ILE	2.3
12	H	144	VAL	2.3
15	L	16	VAL	2.3
14	K	101	ASP	2.3
14	K	150	CYS	2.3
9	C	135	ALA	2.3
11	G	81	ALA	2.3
1	1	170	ILE	2.3
2	2	275	GLY	2.3
6	6	108	GLY	2.3
5	5	391	THR	2.3
6	6	20	THR	2.3
13	I	166	MET	2.3
6	6	109	LEU	2.3
9	C	244	LEU	2.3
4	4	159	ASP	2.3
7	A	119	GLU	2.3
14	K	184	SER	2.3
1	1	15	VAL	2.3
5	5	235	TRP	2.3
8	B	427	CYS	2.3
3	3	63	PHE	2.3
9	C	398	GLU	2.3
2	2	144	TYR	2.3
5	5	77	TYR	2.3
5	5	589	PRO	2.3
9	C	182	ARG	2.3
13	I	180	PRO	2.2
5	5	249	LEU	2.2
8	B	412	ASP	2.2
14	K	187	LEU	2.2
1	1	231	GLU	2.2
5	5	401	GLU	2.2

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Mol	Chain	Res	Type	RSRZ
1	1	58	LEU	2.2
2	2	61	PHE	2.2
9	C	426	LYS	2.2
5	5	385	MET	2.2
2	2	306	GLU	2.2
13	I	185	VAL	2.2
14	K	119	VAL	2.2
9	C	277	ASN	2.2
13	I	142	ALA	2.2
14	K	147	MET	2.2
4	4	138	TYR	2.2
14	K	144	VAL	2.2
5	5	371	PRO	2.2
6	6	83	THR	2.2
9	C	350	CYS	2.2
5	5	98	MET	2.2
5	5	351	HIS	2.2
9	C	261	ILE	2.2
12	H	136	GLY	2.2
11	G	190	PHE	2.2
5	5	333	LEU	2.2
3	3	94	PHE	2.1
14	K	178	PRO	2.1
11	G	185	MET	2.1
14	K	112	ARG	2.1
9	C	323	VAL	2.1
15	L	36	ILE	2.1
9	C	397	GLY	2.1
9	C	372	ASN	2.1
1	1	296	VAL	2.1
14	K	175	VAL	2.1
8	B	423	GLY	2.1
11	G	191	GLU	2.1
13	I	177	GLU	2.1
4	4	393	THR	2.1
14	K	108	ARG	2.1
9	C	433	ALA	2.1
6	6	114	ASN	2.1
5	5	373	THR	2.1
7	A	186	CYS	2.1
4	4	266	LEU	2.1
9	C	435	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
11	G	82	LEU	2.1
5	5	578	HIS	2.1
9	C	327	GLY	2.1
1	1	39	PRO	2.1
11	G	164	PRO	2.1
9	C	266	THR	2.1
5	5	398	ILE	2.1
13	I	175	CYS	2.1
9	C	437	ALA	2.1
9	C	86	ASN	2.1
9	C	152	GLN	2.1
9	C	399	THR	2.0
12	H	138	ASP	2.0
5	5	247	SER	2.0
13	I	144	ALA	2.0
2	2	276	LEU	2.0
2	2	411	LEU	2.0
5	5	569	LEU	2.0
11	G	171	LEU	2.0
1	1	292	SER	2.0
5	5	436	PHE	2.0
6	6	71	PHE	2.0
7	A	190	VAL	2.0
9	C	410	MET	2.0
5	5	375	ILE	2.0
9	C	278	ILE	2.0
8	B	429	LEU	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,

median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
49	FES	H	300	4/4	0.86	0.08	144,147,151,153	0
50	SF4	A	902	8/8	0.86	0.09	183,184,185,186	0
50	SF4	B	500	8/8	0.87	0.09	219,225,240,246	0
50	SF4	A	901	8/8	0.91	0.08	207,208,210,210	0
49	FES	A	900	4/4	0.93	0.06	153,153,153,153	0
50	SF4	I	501	8/8	0.96	0.06	116,119,121,122	0
50	SF4	I	500	8/8	0.98	0.04	107,109,111,111	0
50	SF4	K	500	8/8	0.98	0.04	89,93,97,100	0

6.5 Other polymers ⓘ

There are no such residues in this entry.