



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2014 – 01:57 AM GMT

PDB ID : 3E20
Title : Crystal structure of S.pombe eRF1/eRF3 complex
Authors : Cheng, Z.; Lim, M.; Kong, C.; Song, H.
Deposited on : 2008-08-05
Resolution : 3.50 Å(reported)

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

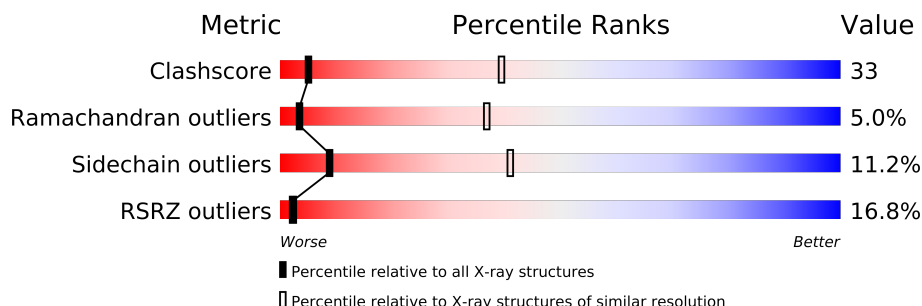
The following versions of software and data (see [references](#)) were used in the production of this report:

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

1 Overall quality at a glance

The reported resolution of this entry is 3.50 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	79885	1039 (3.66-3.34)
Ramachandran outliers	78287	1000 (3.66-3.34)
Sidechain outliers	78261	1000 (3.66-3.34)
RSRZ outliers	66119	1243 (3.70-3.30)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	201	
1	D	201	
1	E	201	
1	J	201	
2	B	441	
2	C	441	
2	H	441	
2	K	441	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 13096 atoms, of which 0 are hydrogen and 0 are deuterium.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Eukaryotic peptide chain release factor GTP-binding subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	196	Total	C	N	O	S	0	0	0
			1524	965	256	293	10			
1	D	196	Total	C	N	O	S	0	0	0
			1524	965	256	293	10			
1	E	196	Total	C	N	O	S	0	0	0
			1524	965	256	293	10			
1	J	196	Total	C	N	O	S	0	0	0
			1524	965	256	293	10			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	462	GLY	-	EXPRESSION TAG	UNP O74718
A	463	PRO	-	EXPRESSION TAG	UNP O74718
A	464	LEU	-	EXPRESSION TAG	UNP O74718
A	465	GLY	-	EXPRESSION TAG	UNP O74718
A	466	SER	-	EXPRESSION TAG	UNP O74718
D	462	GLY	-	EXPRESSION TAG	UNP O74718
D	463	PRO	-	EXPRESSION TAG	UNP O74718
D	464	LEU	-	EXPRESSION TAG	UNP O74718
D	465	GLY	-	EXPRESSION TAG	UNP O74718
D	466	SER	-	EXPRESSION TAG	UNP O74718
E	462	GLY	-	EXPRESSION TAG	UNP O74718
E	463	PRO	-	EXPRESSION TAG	UNP O74718
E	464	LEU	-	EXPRESSION TAG	UNP O74718
E	465	GLY	-	EXPRESSION TAG	UNP O74718
E	466	SER	-	EXPRESSION TAG	UNP O74718
J	462	GLY	-	EXPRESSION TAG	UNP O74718
J	463	PRO	-	EXPRESSION TAG	UNP O74718
J	464	LEU	-	EXPRESSION TAG	UNP O74718
J	465	GLY	-	EXPRESSION TAG	UNP O74718
J	466	SER	-	EXPRESSION TAG	UNP O74718

- Molecule 2 is a protein called Eukaryotic peptide chain release factor subunit 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	261	Total 2088	C 1321	N 347	O 405	S 15	0	0	0
2	B	261	Total 2088	C 1321	N 347	O 405	S 15	0	0	0
2	H	175	Total 1412	C 900	N 225	O 277	S 10	0	0	0
2	K	175	Total 1412	C 900	N 225	O 277	S 10	0	0	0

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-7	MET	-	EXPRESSION TAG	UNP P79063
C	-6	HIS	-	EXPRESSION TAG	UNP P79063
C	-5	HIS	-	EXPRESSION TAG	UNP P79063
C	-4	HIS	-	EXPRESSION TAG	UNP P79063
C	-3	HIS	-	EXPRESSION TAG	UNP P79063
C	-2	HIS	-	EXPRESSION TAG	UNP P79063
C	-1	HIS	-	EXPRESSION TAG	UNP P79063
C	0	MET	-	EXPRESSION TAG	UNP P79063
B	-7	MET	-	EXPRESSION TAG	UNP P79063
B	-6	HIS	-	EXPRESSION TAG	UNP P79063
B	-5	HIS	-	EXPRESSION TAG	UNP P79063
B	-4	HIS	-	EXPRESSION TAG	UNP P79063
B	-3	HIS	-	EXPRESSION TAG	UNP P79063
B	-2	HIS	-	EXPRESSION TAG	UNP P79063
B	-1	HIS	-	EXPRESSION TAG	UNP P79063
B	0	MET	-	EXPRESSION TAG	UNP P79063
H	-7	MET	-	EXPRESSION TAG	UNP P79063
H	-6	HIS	-	EXPRESSION TAG	UNP P79063
H	-5	HIS	-	EXPRESSION TAG	UNP P79063
H	-4	HIS	-	EXPRESSION TAG	UNP P79063
H	-3	HIS	-	EXPRESSION TAG	UNP P79063
H	-2	HIS	-	EXPRESSION TAG	UNP P79063
H	-1	HIS	-	EXPRESSION TAG	UNP P79063
H	0	MET	-	EXPRESSION TAG	UNP P79063
K	-7	MET	-	EXPRESSION TAG	UNP P79063
K	-6	HIS	-	EXPRESSION TAG	UNP P79063
K	-5	HIS	-	EXPRESSION TAG	UNP P79063
K	-4	HIS	-	EXPRESSION TAG	UNP P79063
K	-3	HIS	-	EXPRESSION TAG	UNP P79063
K	-2	HIS	-	EXPRESSION TAG	UNP P79063

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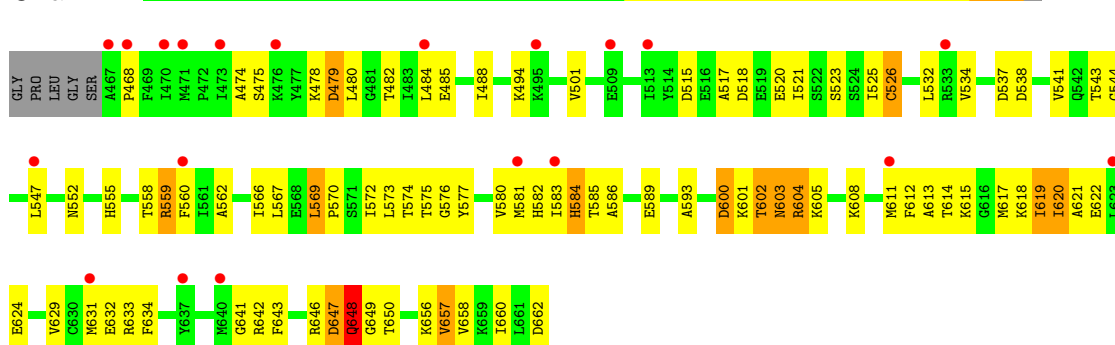
Chain	Residue	Modelled	Actual	Comment	Reference
K	-1	HIS	-	EXPRESSION TAG	UNP P79063
K	0	MET	-	EXPRESSION TAG	UNP P79063

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

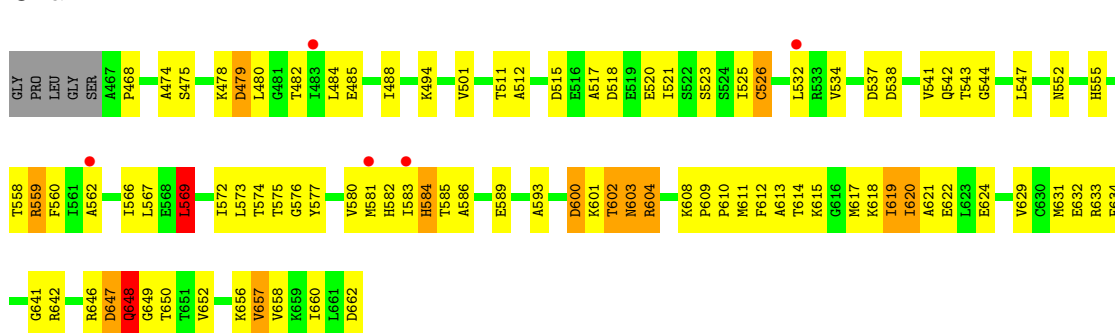
- Molecule 1: Eukaryotic peptide chain release factor GTP-binding subunit

Chain A:



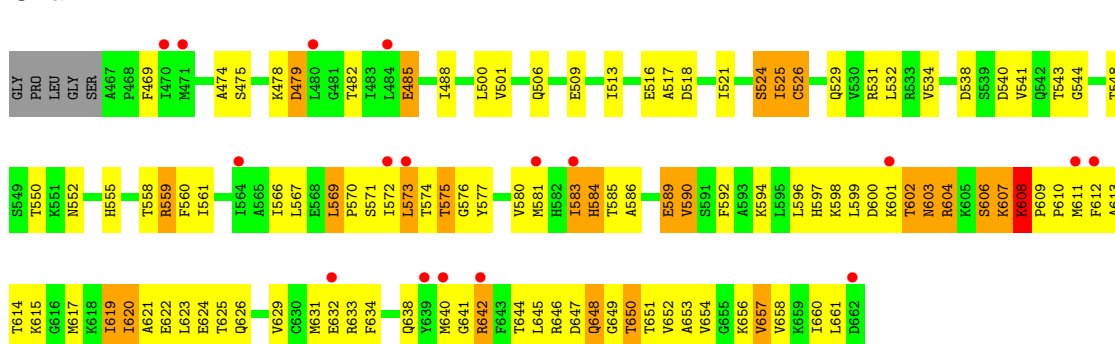
- Molecule 1: Eukaryotic peptide chain release factor GTP-binding subunit

Chain D:



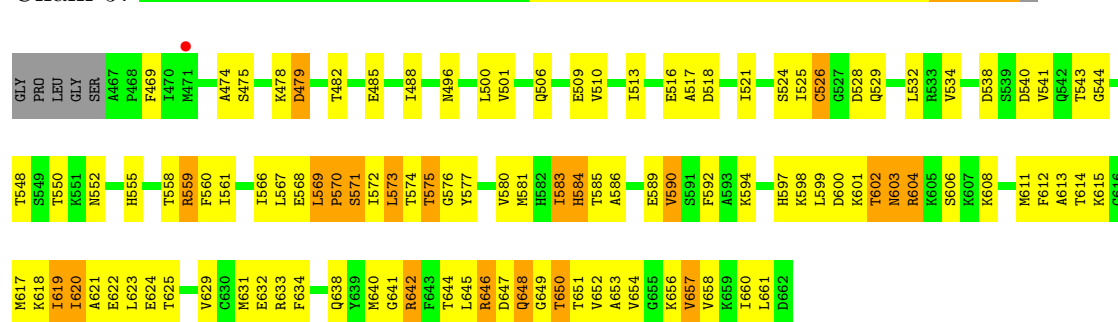
- Molecule 1: Eukaryotic peptide chain release factor GTP-binding subunit

Chain E:



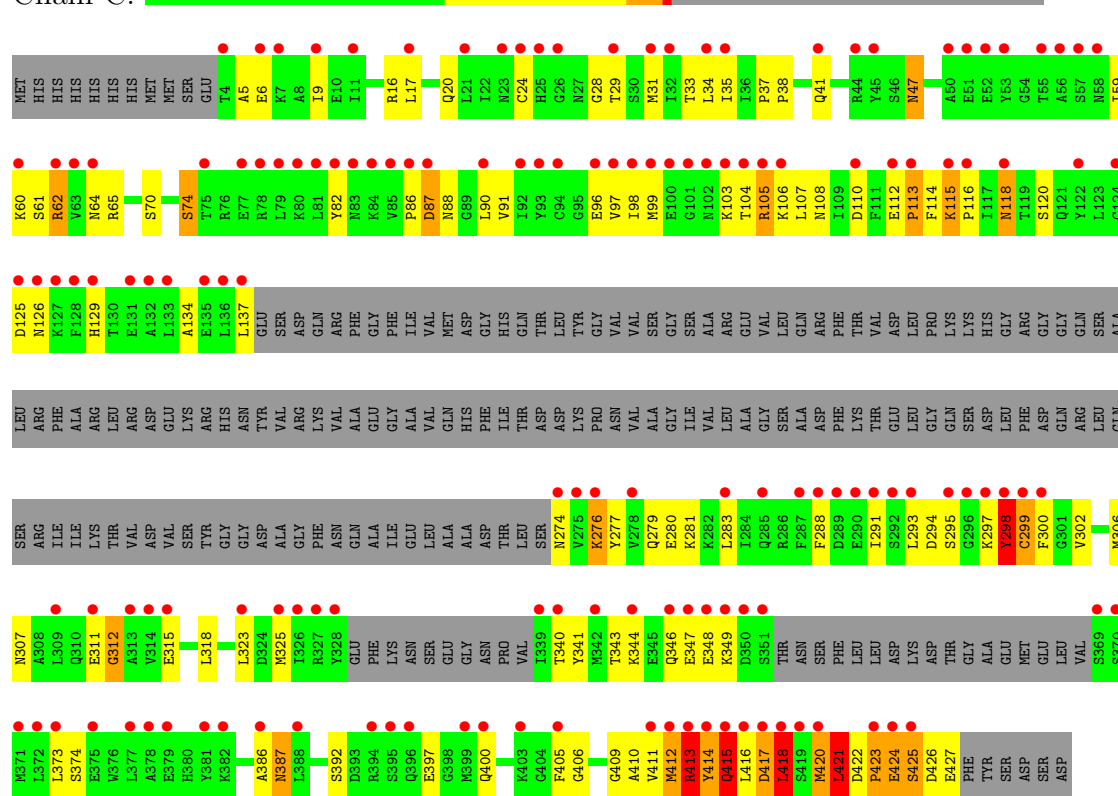
• Molecule 1: Eukaryotic peptide chain release factor GTP-binding subunit

Chain J:



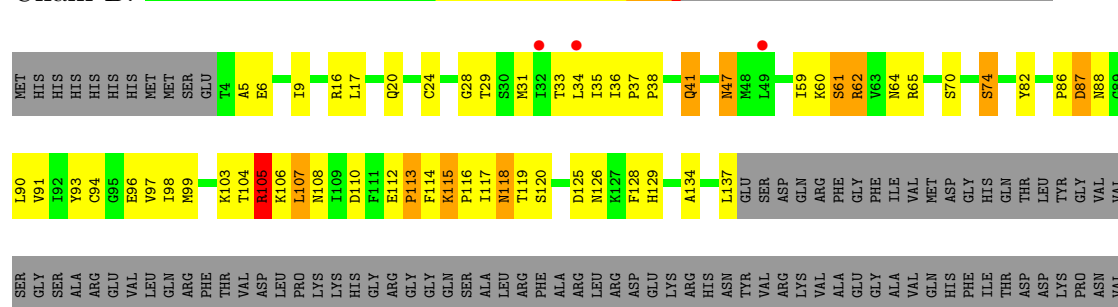
• Molecule 2: Eukaryotic peptide chain release factor subunit 1

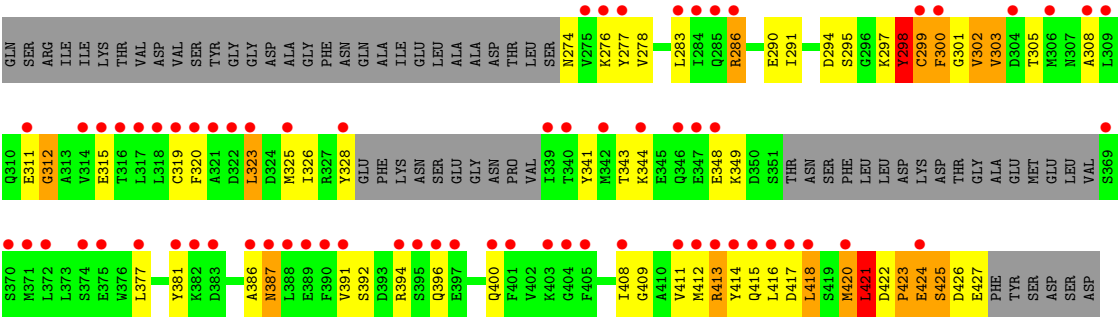
Chain C:



• Molecule 2: Eukaryotic peptide chain release factor subunit 1

Chain B:





4 Data and refinement statistics

Property	Value	Source
Space group	P 43	Depositor
Cell constants a, b, c, α , β , γ	129.85Å 129.85Å 332.64Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 3.50 88.51 – 3.50	Depositor EDS
% Data completeness (in resolution range)	99.7 (30.00-3.50) 99.7 (88.51-3.50)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.43 (at 3.49Å)	Xtriage
Refinement program	REFMAC 5.4.0077	Depositor
R, R_{free}	0.258 , 0.280 0.273 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	123.7	Xtriage
Anisotropy	0.046	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 156.5	EDS
Estimated twinning fraction	0.439 for h,-k,-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 68782 reflections	Xtriage
F_o, F_c correlation	0.81	EDS
Total number of atoms	13096	wwPDB-VP
Average B, all atoms (Å ²)	75.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.25% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	0.36	0/1547	0.57	1/2089 (0.0%)
1	D	0.36	0/1547	0.55	0/2089
1	E	0.49	0/1547	0.63	1/2089 (0.0%)
1	J	0.48	1/1547 (0.1%)	0.65	1/2089 (0.0%)
2	B	0.34	0/2119	0.61	6/2848 (0.2%)
2	C	0.34	0/2119	0.63	7/2848 (0.2%)
2	H	0.34	0/1425	0.55	1/1901 (0.1%)
2	K	0.34	0/1425	0.56	1/1901 (0.1%)
All	All	0.38	1/13276 (0.0%)	0.60	18/17854 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
1	D	0	2
1	E	0	3
1	J	0	1
2	B	0	1
2	C	0	1
2	H	0	1
2	K	0	1
All	All	0	12

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	J	646	ARG	CB-CG	-5.26	1.38	1.52

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	K	413	ARG	NE-CZ-NH2	9.76	125.18	120.30
2	B	105	ARG	CG-CD-NE	8.50	129.65	111.80
2	C	413	ARG	NE-CZ-NH2	-8.02	116.29	120.30
2	C	413	ARG	NE-CZ-NH1	7.61	124.11	120.30
2	B	413	ARG	NE-CZ-NH1	-7.29	116.66	120.30
2	C	65	ARG	NE-CZ-NH2	-6.79	116.90	120.30
2	B	413	ARG	NE-CZ-NH2	6.53	123.57	120.30
2	H	413	ARG	NE-CZ-NH2	6.46	123.53	120.30
2	B	414	TYR	N-CA-CB	-5.91	99.96	110.60
2	C	105	ARG	CG-CD-NE	5.68	123.73	111.80
2	B	413	ARG	N-CA-C	-5.68	95.66	111.00
2	C	413	ARG	N-CA-C	-5.67	95.70	111.00
1	A	570	PRO	N-CA-C	-5.66	97.38	112.10
2	C	414	TYR	N-CA-CB	-5.61	100.50	110.60
2	C	65	ARG	CB-CG-CD	-5.51	97.28	111.60
2	B	105	ARG	CB-CG-CD	-5.39	97.58	111.60
1	J	571	SER	N-CA-CB	-5.21	102.68	110.50
1	E	608	LYS	CB-CG-CD	-5.08	98.38	111.60

There are no chirality outliers.

All (12) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	569	LEU	Peptide
1	A	647	ASP	Peptide
2	B	413	ARG	Peptide
2	C	413	ARG	Peptide
1	D	569	LEU	Peptide
1	D	647	ASP	Peptide
1	E	570	PRO	Peptide
1	E	607	LYS	Peptide
1	E	608	LYS	Peptide
2	H	298	TYR	Peptide
1	J	570	PRO	Peptide
2	K	298	TYR	Peptide

5.2 Close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the

chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1524	0	1569	89	0
1	D	1524	0	1569	92	0
1	E	1524	0	1569	117	0
1	J	1524	0	1569	104	0
2	B	2088	0	2077	144	0
2	C	2088	0	2077	133	0
2	H	1412	0	1388	117	0
2	K	1412	0	1388	86	0
All	All	13096	0	13206	878	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 33.

All (878) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:425:SER:CB	2:B:426:ASP:HA	1.52	1.38
2:H:414:TYR:CE2	2:H:416:LEU:HD11	1.60	1.36
2:C:425:SER:CB	2:C:426:ASP:HA	1.52	1.34
2:H:425:SER:CB	2:H:426:ASP:HA	1.51	1.29
2:K:425:SER:CB	2:K:426:ASP:HA	1.52	1.23
2:K:425:SER:OG	2:K:426:ASP:HA	1.38	1.22
2:C:420:MET:HG2	2:C:421:LEU:HD12	1.24	1.20
2:H:425:SER:OG	2:H:426:ASP:HA	1.37	1.19
2:C:425:SER:OG	2:C:426:ASP:HA	1.40	1.18
2:B:425:SER:OG	2:B:426:ASP:HA	1.40	1.17
2:C:386:ALA:HA	2:C:387:ASN:HB2	1.18	1.17
2:H:418:LEU:HD22	2:H:418:LEU:N	1.54	1.17
2:B:420:MET:HA	2:B:421:LEU:CB	1.73	1.17
2:B:386:ALA:HA	2:B:387:ASN:HB2	1.18	1.16
2:B:420:MET:HG2	2:B:421:LEU:HD12	1.24	1.15
2:K:418:LEU:H	2:K:418:LEU:CD2	1.54	1.15
2:C:420:MET:HA	2:C:421:LEU:CB	1.74	1.14
2:H:421:LEU:HD22	2:H:421:LEU:O	1.47	1.14
2:H:420:MET:HA	2:H:421:LEU:CB	1.76	1.13
2:H:325:MET:HE1	2:H:341:TYR:HB3	1.28	1.12
1:D:572:ILE:HD11	1:D:612:PHE:CB	1.79	1.11
2:K:421:LEU:HD22	2:K:421:LEU:O	1.49	1.11
2:B:414:TYR:CE2	2:B:416:LEU:HD11	1.84	1.11
1:A:572:ILE:HD11	1:A:612:PHE:CB	1.79	1.11
2:B:420:MET:CA	2:B:421:LEU:HB3	1.82	1.10

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:K:418:LEU:H	2:K:418:LEU:HD22	1.06	1.10
2:C:414:TYR:CE2	2:C:416:LEU:HD11	1.84	1.10
2:H:414:TYR:CD2	2:H:416:LEU:HD11	1.87	1.10
2:C:38:PRO:HB3	2:C:88:ASN:HA	1.21	1.10
2:C:420:MET:CA	2:C:421:LEU:HB3	1.82	1.09
2:B:421:LEU:HD22	2:B:421:LEU:O	1.51	1.09
2:K:420:MET:HA	2:K:421:LEU:CB	1.76	1.09
2:B:38:PRO:HB3	2:B:88:ASN:HA	1.20	1.09
2:H:425:SER:CB	2:H:426:ASP:CA	2.31	1.08
2:C:423:PRO:HA	2:C:424:GLU:C	1.73	1.08
2:B:425:SER:HB2	2:B:426:ASP:CA	1.84	1.08
2:B:425:SER:CB	2:B:426:ASP:CA	2.32	1.08
2:C:421:LEU:O	2:C:421:LEU:HD22	1.51	1.08
1:E:646:ARG:CD	1:E:648:GLN:HG2	1.82	1.08
2:C:425:SER:CB	2:C:426:ASP:CA	2.32	1.07
2:C:425:SER:HB2	2:C:426:ASP:CA	1.85	1.07
1:E:646:ARG:HG2	1:E:648:GLN:HA	1.36	1.07
1:J:572:ILE:HD11	1:J:612:PHE:CB	1.84	1.07
2:K:425:SER:CB	2:K:426:ASP:CA	2.32	1.06
2:K:420:MET:CA	2:K:421:LEU:HB3	1.84	1.06
2:K:423:PRO:HA	2:K:424:GLU:C	1.74	1.06
1:E:646:ARG:HD3	1:E:648:GLN:CG	1.84	1.06
1:E:646:ARG:HD3	1:E:648:GLN:HG2	1.08	1.06
2:H:420:MET:CG	2:H:421:LEU:HD12	1.86	1.06
2:K:420:MET:CG	2:K:421:LEU:HD12	1.85	1.06
2:B:423:PRO:HA	2:B:424:GLU:C	1.72	1.06
2:K:418:LEU:HD22	2:K:418:LEU:N	1.53	1.06
2:K:325:MET:HE1	2:K:341:TYR:HB3	1.35	1.05
1:E:559:ARG:CG	1:E:559:ARG:HH11	1.69	1.05
2:H:425:SER:HB2	2:H:426:ASP:CA	1.86	1.05
2:H:420:MET:CA	2:H:421:LEU:HB3	1.84	1.05
2:K:425:SER:HB2	2:K:426:ASP:HA	1.39	1.05
2:H:297:LYS:O	2:H:298:TYR:HB3	1.54	1.05
1:J:559:ARG:CG	1:J:559:ARG:HH11	1.71	1.04
2:K:386:ALA:HA	2:K:387:ASN:HB2	1.06	1.04
2:H:425:SER:HB2	2:H:426:ASP:HA	1.39	1.03
2:K:425:SER:HB2	2:K:426:ASP:CA	1.87	1.03
2:H:420:MET:HG2	2:H:421:LEU:HD12	1.06	1.03
2:H:423:PRO:HA	2:H:424:GLU:C	1.74	1.03
1:A:559:ARG:HH11	1:A:559:ARG:CG	1.71	1.03
1:D:559:ARG:HH11	1:D:559:ARG:CG	1.72	1.03
2:H:386:ALA:HA	2:H:387:ASN:HB2	1.06	1.03

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:572:ILE:HD11	1:D:612:PHE:CG	1.95	1.02
2:C:420:MET:CG	2:C:421:LEU:HD12	1.90	1.02
1:A:572:ILE:HD11	1:A:612:PHE:CG	1.94	1.02
2:B:425:SER:HB2	2:B:426:ASP:HA	1.37	1.01
2:K:420:MET:HG2	2:K:421:LEU:HD12	1.05	1.01
2:H:418:LEU:CD2	2:H:418:LEU:N	2.21	1.01
2:B:420:MET:CG	2:B:421:LEU:HD12	1.90	1.00
1:D:559:ARG:HG3	1:D:559:ARG:NH1	1.61	1.00
1:A:559:ARG:NH1	1:A:559:ARG:HG3	1.61	1.00
2:H:414:TYR:O	2:H:416:LEU:HD12	1.60	1.00
2:C:425:SER:HB2	2:C:426:ASP:HA	1.37	1.00
1:E:572:ILE:HD11	1:E:612:PHE:CB	1.91	0.99
1:J:559:ARG:HG3	1:J:559:ARG:NH1	1.61	0.99
2:H:414:TYR:CE2	2:H:416:LEU:CD1	2.46	0.98
2:K:297:LYS:O	2:K:298:TYR:HB3	1.59	0.97
1:D:572:ILE:HD11	1:D:612:PHE:HB2	1.43	0.97
1:A:572:ILE:HD11	1:A:612:PHE:HB2	1.44	0.97
2:B:24:CYS:HB2	2:B:97:VAL:HG21	1.47	0.96
1:A:559:ARG:HH11	1:A:559:ARG:HG3	0.79	0.96
2:C:24:CYS:HB2	2:C:97:VAL:HG21	1.48	0.96
2:B:418:LEU:N	2:B:418:LEU:HD22	1.80	0.96
2:H:386:ALA:CA	2:H:387:ASN:HB2	1.95	0.95
1:E:647:ASP:O	1:E:649:GLY:N	1.99	0.95
2:K:386:ALA:CA	2:K:387:ASN:HB2	1.95	0.95
2:C:413:ARG:O	2:C:414:TYR:CD1	2.20	0.94
2:K:420:MET:HG2	2:K:421:LEU:CD1	1.98	0.94
1:D:559:ARG:HG3	1:D:559:ARG:HH11	0.80	0.94
2:C:418:LEU:N	2:C:418:LEU:HD22	1.81	0.94
1:E:589:GLU:OE2	1:E:626:GLN:HG3	1.68	0.94
1:E:589:GLU:OE2	1:E:626:GLN:CG	2.16	0.93
1:J:572:ILE:HD11	1:J:612:PHE:HB2	1.49	0.93
1:E:559:ARG:HG3	1:E:559:ARG:HH11	0.78	0.93
2:K:325:MET:CE	2:K:341:TYR:HB3	1.98	0.93
2:B:297:LYS:O	2:B:298:TYR:HB3	1.69	0.93
2:C:297:LYS:O	2:C:298:TYR:HB3	1.67	0.93
2:H:420:MET:HG2	2:H:421:LEU:CD1	1.99	0.92
1:A:647:ASP:O	1:A:649:GLY:N	2.03	0.92
2:B:105:ARG:HH12	2:B:107:LEU:HD12	1.34	0.92
1:J:559:ARG:HG3	1:J:559:ARG:HH11	0.80	0.92
1:J:544:GLY:HA2	1:J:585:THR:HG23	1.52	0.92
1:E:613:ALA:HA	1:E:617:MET:HE1	1.52	0.92
2:H:325:MET:CE	2:H:341:TYR:HB3	2.00	0.92

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:415:GLN:O	2:B:416:LEU:HD12	1.71	0.92
1:J:647:ASP:O	1:J:649:GLY:N	2.03	0.91
2:C:415:GLN:O	2:C:416:LEU:HD12	1.70	0.91
2:B:86:PRO:HB2	2:B:112:GLU:OE1	1.70	0.91
1:E:544:GLY:HA2	1:E:585:THR:HG23	1.50	0.91
1:E:559:ARG:NH1	1:E:559:ARG:HG3	1.59	0.91
2:H:418:LEU:CD2	2:H:418:LEU:H	1.82	0.91
2:B:413:ARG:O	2:B:414:TYR:CD1	2.24	0.90
1:A:600:ASP:O	1:A:618:LYS:NZ	2.05	0.90
2:H:315:GLU:HB2	2:H:413:ARG:HG3	1.54	0.90
1:D:647:ASP:O	1:D:649:GLY:N	2.04	0.90
1:D:600:ASP:O	1:D:618:LYS:NZ	2.05	0.89
2:H:386:ALA:HA	2:H:387:ASN:CB	1.97	0.89
2:C:400:GLN:HA	1:D:576:GLY:HA3	1.54	0.89
2:C:86:PRO:HB2	2:C:112:GLU:OE1	1.72	0.89
2:C:298:TYR:H	2:C:412:MET:CE	1.86	0.88
1:A:576:GLY:HA3	2:B:400:GLN:HA	1.56	0.87
2:H:414:TYR:CD2	2:H:416:LEU:CD1	2.57	0.86
1:E:602:THR:HG23	1:E:604:ARG:HD2	1.55	0.85
1:A:572:ILE:HG12	1:A:573:LEU:H	1.39	0.85
2:B:298:TYR:H	2:B:412:MET:HE2	1.41	0.85
2:B:105:ARG:NH1	2:B:107:LEU:HD12	1.91	0.85
2:C:386:ALA:HA	2:C:387:ASN:CB	2.06	0.84
2:C:386:ALA:CA	2:C:387:ASN:HB2	2.04	0.84
2:C:298:TYR:N	2:C:412:MET:HE2	1.92	0.84
2:C:414:TYR:CD2	2:C:416:LEU:HD11	2.13	0.84
1:E:572:ILE:HD11	1:E:612:PHE:HB2	1.57	0.84
1:D:572:ILE:HG12	1:D:573:LEU:H	1.40	0.83
1:E:646:ARG:HG2	1:E:648:GLN:CA	2.08	0.83
2:B:386:ALA:CA	2:B:387:ASN:HB2	2.04	0.83
2:B:425:SER:HG	2:B:426:ASP:HA	1.44	0.82
2:B:386:ALA:HA	2:B:387:ASN:CB	2.06	0.82
2:B:414:TYR:CD2	2:B:416:LEU:HD11	2.14	0.82
2:B:59:ILE:HB	2:B:65:ARG:HG3	1.62	0.81
2:K:420:MET:HA	2:K:421:LEU:HB3	0.89	0.81
2:H:420:MET:HA	2:H:421:LEU:HB3	0.89	0.81
1:A:573:LEU:HD22	1:A:577:TYR:CG	2.16	0.81
1:D:573:LEU:HD22	1:D:577:TYR:CG	2.16	0.80
2:C:414:TYR:HE2	2:C:416:LEU:HD11	1.46	0.80
2:C:300:PHE:H	2:C:409:GLY:HA2	1.45	0.80
2:C:298:TYR:H	2:C:412:MET:HE2	1.44	0.80
2:B:325:MET:HE1	2:B:341:TYR:HB3	1.63	0.80

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:J:573:LEU:HD22	1:J:577:TYR:CG	2.17	0.79
2:H:18:VAL:HG23	2:H:19:LYS:H	1.46	0.79
2:C:47:ASN:C	2:C:47:ASN:HD22	1.86	0.79
2:B:298:TYR:N	2:B:412:MET:HE2	1.95	0.79
2:B:414:TYR:HE2	2:B:416:LEU:HD11	1.46	0.79
2:C:114:PHE:O	2:C:115:LYS:HB2	1.82	0.79
1:E:573:LEU:HD22	1:E:577:TYR:CG	2.18	0.79
2:B:47:ASN:HD22	2:B:47:ASN:C	1.86	0.78
2:B:298:TYR:H	2:B:412:MET:CE	1.95	0.78
2:B:300:PHE:H	2:B:409:GLY:CA	1.97	0.78
2:K:386:ALA:HA	2:K:387:ASN:CB	1.97	0.78
2:B:300:PHE:H	2:B:409:GLY:HA2	1.46	0.78
2:H:300:PHE:H	2:H:409:GLY:HA3	1.48	0.78
2:B:114:PHE:O	2:B:115:LYS:HB2	1.82	0.78
2:B:420:MET:HA	2:B:421:LEU:HB3	0.86	0.78
2:B:414:TYR:O	2:B:416:LEU:HD13	1.83	0.77
2:C:386:ALA:HB1	2:C:387:ASN:O	1.85	0.77
2:C:414:TYR:O	2:C:416:LEU:HD13	1.83	0.77
2:B:386:ALA:HB1	2:B:387:ASN:O	1.85	0.77
2:K:18:VAL:HG23	2:K:19:LYS:H	1.50	0.77
1:J:572:ILE:HD11	1:J:612:PHE:CG	2.19	0.77
1:J:574:THR:O	1:J:576:GLY:N	2.18	0.77
2:C:300:PHE:H	2:C:409:GLY:CA	1.96	0.77
2:K:425:SER:HG	2:K:426:ASP:HA	1.50	0.77
1:A:574:THR:O	1:A:576:GLY:N	2.18	0.77
1:A:494:LYS:HG2	1:A:523:SER:HB3	1.67	0.77
1:E:574:THR:O	1:E:576:GLY:N	2.18	0.76
1:D:560:PHE:HZ	1:D:581:MET:HE3	1.51	0.76
1:E:589:GLU:OE2	1:E:626:GLN:HG2	1.86	0.76
2:B:414:TYR:CD2	2:B:414:TYR:O	2.39	0.76
2:C:325:MET:CE	2:C:341:TYR:HB3	2.15	0.76
1:J:602:THR:HG23	1:J:604:ARG:HD2	1.68	0.76
1:A:478:LYS:O	1:A:479:ASP:HB2	1.86	0.76
1:D:494:LYS:HG2	1:D:523:SER:HB3	1.67	0.76
1:D:574:THR:O	1:D:576:GLY:N	2.18	0.76
1:E:646:ARG:CG	1:E:648:GLN:HG2	2.15	0.76
2:B:325:MET:CE	2:B:341:TYR:HB3	2.16	0.75
2:C:33:THR:HG21	2:C:129:HIS:H	1.52	0.75
2:B:33:THR:HG21	2:B:129:HIS:H	1.51	0.75
2:C:420:MET:HA	2:C:421:LEU:HB3	0.86	0.75
2:C:325:MET:HE1	2:C:341:TYR:HB3	1.67	0.75
2:K:411:VAL:O	2:K:411:VAL:HG23	1.87	0.75

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:J:580:VAL:HG21	1:J:648:GLN:HG3	1.69	0.75
2:C:414:TYR:O	2:C:414:TYR:CD2	2.40	0.74
1:A:572:ILE:HD11	1:A:612:PHE:CD1	2.22	0.74
2:B:414:TYR:O	2:B:416:LEU:CD1	2.36	0.74
1:D:478:LYS:O	1:D:479:ASP:HB2	1.87	0.74
2:B:425:SER:HB2	2:B:426:ASP:C	2.09	0.73
1:E:572:ILE:HD11	1:E:612:PHE:CG	2.23	0.73
2:H:91:VAL:O	2:H:110:ASP:HA	1.89	0.73
1:D:572:ILE:CD1	1:D:612:PHE:HB2	2.19	0.73
2:C:425:SER:HB2	2:C:426:ASP:C	2.09	0.73
2:H:411:VAL:HG23	2:H:411:VAL:O	1.88	0.73
1:J:613:ALA:HA	1:J:617:MET:HE1	1.69	0.73
2:C:413:ARG:O	2:C:414:TYR:HD1	1.72	0.73
1:E:607:LYS:O	1:E:608:LYS:HB2	1.88	0.73
2:C:417:ASP:C	2:C:418:LEU:HD22	2.09	0.72
2:H:414:TYR:O	2:H:416:LEU:CD1	2.35	0.72
2:H:415:GLN:O	2:H:416:LEU:HD12	1.90	0.72
2:H:301:GLY:O	2:H:305:THR:OG1	2.02	0.72
2:H:425:SER:HB2	2:H:427:GLU:OE1	1.89	0.72
2:C:425:SER:HB2	2:C:427:GLU:OE1	1.90	0.72
2:H:425:SER:HG	2:H:426:ASP:HA	1.52	0.72
2:K:423:PRO:CA	2:K:424:GLU:C	2.57	0.72
2:C:414:TYR:O	2:C:416:LEU:CD1	2.37	0.72
1:A:572:ILE:CD1	1:A:612:PHE:HB2	2.19	0.72
1:J:572:ILE:CD1	1:J:612:PHE:HB2	2.20	0.71
1:D:572:ILE:HD11	1:D:612:PHE:CD1	2.24	0.71
2:B:417:ASP:C	2:B:418:LEU:HD22	2.10	0.71
1:D:501:VAL:HG11	1:D:541:VAL:HG22	1.72	0.71
1:J:478:LYS:O	1:J:479:ASP:HB2	1.89	0.71
2:C:425:SER:HG	2:C:426:ASP:HA	1.55	0.71
1:D:475:SER:HB2	1:D:485:GLU:HG3	1.73	0.71
1:A:572:ILE:CD1	1:A:612:PHE:CD1	2.74	0.71
2:B:59:ILE:HG22	2:B:61:SER:H	1.56	0.71
1:J:555:HIS:HD2	1:J:632:GLU:OE2	1.74	0.71
1:J:634:PHE:CZ	1:J:638:GLN:HG2	2.26	0.71
1:A:501:VAL:HG11	1:A:541:VAL:HG22	1.72	0.71
2:B:420:MET:HG2	2:B:421:LEU:CD1	2.13	0.70
1:A:475:SER:HB2	1:A:485:GLU:HG3	1.73	0.70
1:A:560:PHE:HZ	1:A:581:MET:HE3	1.55	0.70
2:C:59:ILE:HG22	2:C:61:SER:H	1.56	0.70
2:H:425:SER:HB2	2:H:426:ASP:C	2.11	0.70
1:J:576:GLY:HA3	2:K:400:GLN:HA	1.73	0.70

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:602:THR:HG23	1:A:604:ARG:HD2	1.73	0.70
1:E:560:PHE:HZ	1:E:581:MET:HE3	1.54	0.70
1:J:572:ILE:HG13	1:J:612:PHE:HD1	1.55	0.70
2:H:417:ASP:C	2:H:418:LEU:HD22	2.12	0.70
1:E:517:ALA:O	1:E:518:ASP:HB2	1.90	0.70
1:A:544:GLY:HA2	1:A:585:THR:HG23	1.73	0.69
2:K:425:SER:HB2	2:K:426:ASP:C	2.12	0.69
1:E:576:GLY:HA3	2:H:400:GLN:HA	1.75	0.69
1:D:544:GLY:HA2	1:D:585:THR:HG23	1.75	0.69
2:C:423:PRO:CA	2:C:424:GLU:C	2.58	0.69
1:J:517:ALA:O	1:J:518:ASP:HB2	1.93	0.69
2:K:425:SER:HB2	2:K:427:GLU:OE1	1.92	0.69
1:D:572:ILE:HG12	1:D:573:LEU:N	2.08	0.69
1:D:572:ILE:CD1	1:D:612:PHE:CD1	2.76	0.68
2:H:423:PRO:CA	2:H:424:GLU:C	2.57	0.68
1:A:572:ILE:HG12	1:A:573:LEU:N	2.08	0.68
1:E:572:ILE:HG13	1:E:612:PHE:HD1	1.58	0.68
1:A:614:THR:HG22	1:A:617:MET:SD	2.34	0.68
2:B:425:SER:HB2	2:B:427:GLU:OE1	1.94	0.68
1:J:646:ARG:HB3	1:J:648:GLN:HG2	1.75	0.68
1:D:614:THR:HG22	1:D:617:MET:SD	2.34	0.68
1:E:572:ILE:HG13	1:E:612:PHE:CD1	2.29	0.68
2:C:420:MET:HG2	2:C:421:LEU:CD1	2.14	0.68
1:E:646:ARG:CG	1:E:648:GLN:HA	2.19	0.68
1:E:634:PHE:CZ	1:E:638:GLN:HG2	2.29	0.67
2:B:423:PRO:HA	2:B:424:GLU:O	1.95	0.67
2:H:377:LEU:O	2:H:381:TYR:HB3	1.94	0.67
2:B:293:LEU:O	2:B:294:ASP:HB3	1.94	0.67
2:H:297:LYS:O	2:H:298:TYR:CB	2.38	0.67
1:E:478:LYS:O	1:E:479:ASP:HB2	1.93	0.67
2:H:413:ARG:O	2:H:414:TYR:HB3	1.94	0.66
2:B:411:VAL:O	2:B:411:VAL:HG23	1.96	0.66
2:B:415:GLN:C	2:B:416:LEU:HD12	2.16	0.66
2:C:415:GLN:C	2:C:416:LEU:HD12	2.15	0.66
1:J:572:ILE:HG13	1:J:612:PHE:CD1	2.30	0.66
1:J:488:ILE:HB	1:J:526:CYS:HA	1.77	0.66
1:D:602:THR:HG23	1:D:604:ARG:HD2	1.75	0.66
2:C:423:PRO:HA	2:C:424:GLU:O	1.94	0.66
2:C:315:GLU:HB2	2:C:413:ARG:HG2	1.77	0.65
1:E:559:ARG:CG	1:E:559:ARG:NH1	2.39	0.65
1:A:559:ARG:HD3	1:A:622:GLU:OE2	1.96	0.65
2:C:411:VAL:O	2:C:411:VAL:HG23	1.97	0.65

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:61:SER:HB2	2:B:64:ASN:HB3	1.79	0.65
1:E:488:ILE:HD12	1:E:525:ILE:O	1.95	0.65
1:E:575:THR:O	1:E:592:PHE:O	2.14	0.65
2:K:377:LEU:O	2:K:381:TYR:HB3	1.97	0.65
2:B:283:LEU:HD21	2:B:318:LEU:HD22	1.79	0.65
2:C:344:LYS:HA	2:C:346:GLN:HE21	1.63	0.64
2:C:283:LEU:HD21	2:C:318:LEU:HD22	1.79	0.64
1:J:516:GLU:OE2	1:J:529:GLN:HB3	1.97	0.64
2:H:421:LEU:HD22	2:H:421:LEU:C	2.17	0.64
1:J:559:ARG:HD3	1:J:622:GLU:OE2	1.97	0.64
1:E:572:ILE:CD1	1:E:612:PHE:HB2	2.26	0.64
1:D:559:ARG:HD3	1:D:622:GLU:OE2	1.98	0.64
2:C:61:SER:HB2	2:C:64:ASN:HB3	1.80	0.64
2:K:421:LEU:C	2:K:421:LEU:HD22	2.17	0.64
1:J:573:LEU:CD2	1:J:577:TYR:CG	2.81	0.63
1:J:478:LYS:HE3	1:J:482:THR:OG1	1.97	0.63
2:B:422:ASP:HB3	2:B:423:PRO:HD2	1.80	0.63
2:B:413:ARG:O	2:B:414:TYR:HD1	1.76	0.63
1:E:555:HIS:HD2	1:E:632:GLU:OE2	1.80	0.63
2:B:414:TYR:CD2	2:B:416:LEU:CD1	2.82	0.63
2:B:344:LYS:HA	2:B:346:GLN:HE21	1.63	0.63
2:K:418:LEU:CD2	2:K:418:LEU:N	2.21	0.63
2:C:422:ASP:HB3	2:C:423:PRO:HD2	1.80	0.63
1:D:567:LEU:O	1:D:615:LYS:HG3	1.99	0.63
1:E:488:ILE:HB	1:E:526:CYS:HA	1.79	0.62
1:D:559:ARG:NH1	1:D:559:ARG:CG	2.42	0.62
1:J:501:VAL:HG11	1:J:541:VAL:HG22	1.80	0.62
1:J:559:ARG:CG	1:J:559:ARG:NH1	2.40	0.62
1:E:559:ARG:HD3	1:E:622:GLU:OE2	1.99	0.62
2:H:348:GLU:HG3	2:H:349:LYS:H	1.65	0.62
2:B:423:PRO:CA	2:B:424:GLU:C	2.58	0.62
1:E:599:LEU:HD11	1:E:620:ILE:HG22	1.80	0.62
1:D:475:SER:HB2	1:D:485:GLU:CG	2.30	0.62
1:A:478:LYS:O	1:A:479:ASP:CB	2.48	0.62
1:A:567:LEU:O	1:A:615:LYS:HG3	2.00	0.62
2:K:348:GLU:HG3	2:K:349:LYS:H	1.65	0.61
1:D:573:LEU:CD2	1:D:577:TYR:CG	2.83	0.61
2:B:418:LEU:N	2:B:418:LEU:CD2	2.54	0.61
1:A:573:LEU:CD2	1:A:577:TYR:CG	2.82	0.61
2:C:418:LEU:CD2	2:C:418:LEU:N	2.55	0.61
1:E:613:ALA:HA	1:E:617:MET:CE	2.28	0.61
1:A:475:SER:HB2	1:A:485:GLU:CG	2.30	0.61

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:421:LEU:C	2:B:421:LEU:HD22	2.21	0.61
2:C:414:TYR:CD2	2:C:416:LEU:CD1	2.82	0.61
1:E:516:GLU:OE2	1:E:529:GLN:HB3	2.01	0.61
1:E:573:LEU:CD2	1:E:577:TYR:CG	2.84	0.61
1:J:575:THR:O	1:J:592:PHE:O	2.18	0.61
2:H:414:TYR:HE2	2:H:416:LEU:HD21	1.65	0.60
2:K:300:PHE:H	2:K:409:GLY:HA3	1.66	0.60
1:A:560:PHE:HZ	1:A:581:MET:CE	2.14	0.60
1:J:560:PHE:HZ	1:J:581:MET:HE3	1.66	0.60
1:A:631:MET:HG3	1:A:660:ILE:HD11	1.84	0.60
1:A:559:ARG:CG	1:A:559:ARG:NH1	2.41	0.59
1:D:560:PHE:HZ	1:D:581:MET:CE	2.15	0.59
2:C:300:PHE:N	2:C:409:GLY:HA2	2.16	0.59
1:J:569:LEU:HD22	1:J:652:VAL:HB	1.83	0.59
1:E:580:VAL:HG21	1:E:648:GLN:HG3	1.84	0.59
2:H:415:GLN:O	2:H:416:LEU:CD1	2.50	0.59
2:C:295:SER:OG	2:C:297:LYS:HG2	2.02	0.59
2:C:421:LEU:C	2:C:421:LEU:HD22	2.21	0.59
1:E:478:LYS:HE3	1:E:482:THR:OG1	2.02	0.59
2:H:414:TYR:HE2	2:H:416:LEU:CG	2.15	0.59
2:B:414:TYR:O	2:B:415:GLN:C	2.41	0.59
1:D:572:ILE:CG1	1:D:573:LEU:H	2.12	0.59
1:J:633:ARG:HA	1:J:657:VAL:HG22	1.85	0.59
1:J:613:ALA:HA	1:J:617:MET:CE	2.32	0.59
1:J:634:PHE:CE1	1:J:638:GLN:HG2	2.38	0.59
2:B:281:LYS:HE3	2:B:397:GLU:OE1	2.03	0.59
2:B:74:SER:OG	2:B:106:LYS:HE2	2.02	0.59
1:E:646:ARG:HD3	1:E:648:GLN:CD	2.22	0.58
1:E:633:ARG:HA	1:E:657:VAL:HG22	1.86	0.58
1:D:631:MET:HG3	1:D:660:ILE:HD11	1.85	0.58
2:C:281:LYS:HE3	2:C:397:GLU:OE1	2.03	0.58
2:B:422:ASP:CB	2:B:423:PRO:HD2	2.34	0.58
2:C:414:TYR:O	2:C:415:GLN:C	2.42	0.58
1:J:488:ILE:HD12	1:J:525:ILE:O	2.03	0.58
2:K:411:VAL:CG2	2:K:411:VAL:O	2.50	0.58
2:H:423:PRO:HA	2:H:424:GLU:O	2.03	0.58
2:B:300:PHE:N	2:B:409:GLY:HA2	2.16	0.58
2:C:422:ASP:CB	2:C:423:PRO:HD2	2.34	0.57
1:E:558:THR:O	1:E:624:GLU:HA	2.03	0.57
2:K:319:CYS:SG	2:K:323:LEU:HD12	2.43	0.57
2:H:386:ALA:HB1	2:H:387:ASN:O	2.04	0.57
2:K:75:THR:HA	2:K:78:ARG:HG3	1.85	0.57

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:572:ILE:CG1	1:A:573:LEU:H	2.12	0.57
1:A:584:HIS:HD2	1:A:586:ALA:H	1.51	0.57
2:B:423:PRO:CA	2:B:424:GLU:O	2.52	0.57
1:D:584:HIS:HD2	1:D:586:ALA:H	1.50	0.57
1:E:560:PHE:HZ	1:E:581:MET:CE	2.18	0.57
2:B:70:SER:O	2:B:74:SER:HB2	2.05	0.57
2:K:300:PHE:H	2:K:409:GLY:CA	2.18	0.57
2:C:423:PRO:CA	2:C:424:GLU:O	2.53	0.57
1:J:573:LEU:C	1:J:573:LEU:HD13	2.25	0.57
1:D:573:LEU:HD13	1:D:573:LEU:C	2.26	0.56
1:A:573:LEU:HD13	1:A:573:LEU:C	2.26	0.56
2:K:308:ALA:CB	2:K:412:MET:HE3	2.35	0.56
2:H:290:GLU:OE1	2:H:295:SER:OG	2.20	0.56
2:C:134:ALA:HA	2:C:137:LEU:HD12	1.87	0.56
1:D:478:LYS:O	1:D:479:ASP:CB	2.49	0.56
2:C:426:ASP:C	2:C:427:GLU:OE1	2.44	0.56
2:H:415:GLN:C	2:H:416:LEU:HD12	2.25	0.56
1:E:572:ILE:CG1	1:E:612:PHE:HB2	2.36	0.56
1:E:573:LEU:HD13	1:E:573:LEU:C	2.25	0.56
1:E:634:PHE:CE1	1:E:638:GLN:HG2	2.41	0.56
2:H:415:GLN:O	2:H:416:LEU:CG	2.54	0.56
2:H:418:LEU:HD23	2:H:418:LEU:H	1.70	0.56
2:K:423:PRO:HA	2:K:424:GLU:O	2.03	0.56
2:H:415:GLN:O	2:H:416:LEU:HG	2.06	0.56
2:C:416:LEU:CD1	2:C:416:LEU:N	2.69	0.56
2:C:417:ASP:OD1	2:C:417:ASP:N	2.37	0.56
2:C:416:LEU:C	2:C:417:ASP:OD1	2.45	0.56
2:B:113:PRO:HG2	2:B:116:PRO:HG3	1.88	0.56
2:H:411:VAL:O	2:H:411:VAL:CG2	2.53	0.56
2:H:326:ILE:HD12	2:H:328:TYR:CE2	2.41	0.56
1:E:602:THR:CG2	1:E:604:ARG:HD2	2.33	0.56
1:D:560:PHE:CZ	1:D:581:MET:HE3	2.37	0.56
2:C:6:GLU:O	2:C:9:ILE:HG22	2.06	0.56
1:E:607:LYS:O	1:E:608:LYS:CB	2.53	0.55
2:B:294:ASP:O	2:B:295:SER:C	2.44	0.55
2:H:300:PHE:H	2:H:409:GLY:CA	2.17	0.55
1:D:613:ALA:HA	1:D:617:MET:HE1	1.87	0.55
2:B:6:GLU:O	2:B:9:ILE:HG22	2.06	0.55
1:E:631:MET:HG3	1:E:660:ILE:HD11	1.88	0.55
2:K:326:ILE:HD12	2:K:328:TYR:CE2	2.41	0.55
2:B:426:ASP:C	2:B:427:GLU:OE1	2.44	0.55
1:J:560:PHE:HZ	1:J:581:MET:CE	2.19	0.55

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:134:ALA:HA	2:B:137:LEU:HD12	1.86	0.55
2:B:24:CYS:CB	2:B:97:VAL:HG21	2.30	0.55
1:J:634:PHE:HB2	1:J:656:LYS:HB2	1.89	0.55
1:J:572:ILE:HD11	1:J:612:PHE:HB3	1.83	0.55
1:J:567:LEU:CD1	1:J:654:VAL:HG13	2.37	0.55
1:J:572:ILE:CG1	1:J:612:PHE:HB2	2.36	0.55
1:A:525:ILE:HG13	1:A:526:CYS:O	2.06	0.55
2:C:113:PRO:HG2	2:C:116:PRO:HG3	1.89	0.55
2:K:290:GLU:OE1	2:K:297:LYS:HG3	2.07	0.55
1:E:501:VAL:HG11	1:E:541:VAL:HG22	1.89	0.55
1:D:525:ILE:HG13	1:D:526:CYS:O	2.07	0.55
1:D:572:ILE:CD1	1:D:612:PHE:CG	2.81	0.55
1:E:559:ARG:HG2	1:E:624:GLU:HG2	1.89	0.55
1:J:558:THR:O	1:J:624:GLU:HA	2.07	0.55
2:K:386:ALA:HB1	2:K:387:ASN:O	2.07	0.55
2:H:348:GLU:HG3	2:H:349:LYS:N	2.22	0.55
2:H:308:ALA:CB	2:H:412:MET:HE3	2.37	0.55
2:H:290:GLU:OE1	2:H:297:LYS:HG3	2.07	0.54
2:H:414:TYR:CE2	2:H:416:LEU:CG	2.90	0.54
2:H:414:TYR:HD2	2:H:415:GLN:O	1.90	0.54
2:B:417:ASP:N	2:B:417:ASP:OD1	2.37	0.54
2:K:297:LYS:O	2:K:298:TYR:CB	2.39	0.54
1:J:566:ILE:HG22	1:J:615:LYS:HA	1.90	0.54
2:K:301:GLY:HA2	2:K:408:ILE:O	2.07	0.54
1:J:521:ILE:HD11	1:J:524:SER:OG	2.08	0.54
2:H:423:PRO:CA	2:H:424:GLU:O	2.56	0.54
1:J:572:ILE:HG12	1:J:573:LEU:N	2.21	0.54
1:J:567:LEU:O	1:J:615:LYS:HG3	2.07	0.54
2:K:348:GLU:HG3	2:K:349:LYS:N	2.22	0.54
2:K:423:PRO:CA	2:K:424:GLU:O	2.56	0.54
2:K:418:LEU:H	2:K:418:LEU:HD23	1.60	0.54
2:B:416:LEU:C	2:B:417:ASP:OD1	2.46	0.54
2:K:301:GLY:O	2:K:305:THR:OG1	2.15	0.54
2:B:315:GLU:HB2	2:B:413:ARG:HG2	1.89	0.54
2:B:82:TYR:CE1	2:B:90:LEU:HD21	2.42	0.54
1:E:634:PHE:HB2	1:E:656:LYS:HB2	1.90	0.54
1:J:569:LEU:CD2	1:J:652:VAL:HB	2.37	0.54
1:J:559:ARG:HG2	1:J:624:GLU:HG2	1.91	0.53
2:K:291:ILE:HG12	2:K:298:TYR:CE1	2.43	0.53
1:E:602:THR:HG23	1:E:604:ARG:CD	2.33	0.53
1:E:567:LEU:O	1:E:615:LYS:HG3	2.08	0.53
1:A:619:ILE:HG12	1:A:620:ILE:N	2.23	0.53

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:295:SER:OG	2:B:297:LYS:HG2	2.08	0.53
2:C:420:MET:HG3	2:C:421:LEU:HD12	1.87	0.53
2:C:82:TYR:CE1	2:C:90:LEU:HD21	2.42	0.53
1:J:572:ILE:CD1	1:J:612:PHE:CG	2.92	0.53
1:E:647:ASP:C	1:E:649:GLY:H	2.04	0.53
1:J:590:VAL:HG23	1:J:625:THR:HG22	1.91	0.53
1:J:469:PHE:O	1:J:548:THR:HA	2.08	0.53
2:C:37:PRO:HD3	2:C:120:SER:HA	1.91	0.53
1:J:513:ILE:HG13	1:J:532:LEU:CD2	2.39	0.53
1:A:633:ARG:HA	1:A:657:VAL:HG22	1.91	0.53
1:A:573:LEU:CD2	1:A:577:TYR:CD1	2.92	0.53
2:H:416:LEU:O	2:H:417:ASP:C	2.47	0.53
2:K:422:ASP:CB	2:K:423:PRO:HD2	2.39	0.53
1:E:597:HIS:HB2	1:E:620:ILE:HG23	1.91	0.53
2:H:426:ASP:C	2:H:427:GLU:OE1	2.46	0.52
2:C:415:GLN:C	2:C:416:LEU:CD1	2.77	0.52
2:B:299:CYS:HA	2:B:300:PHE:HB3	1.91	0.52
2:K:426:ASP:C	2:K:427:GLU:OE1	2.47	0.52
2:B:38:PRO:HB3	2:B:88:ASN:CA	2.14	0.52
2:C:24:CYS:CB	2:C:97:VAL:HG21	2.31	0.52
1:J:597:HIS:HB2	1:J:620:ILE:HG23	1.92	0.52
1:A:634:PHE:HE1	1:A:642:ARG:HH11	1.57	0.52
1:E:513:ILE:HG13	1:E:532:LEU:CD2	2.39	0.52
1:D:573:LEU:CD2	1:D:577:TYR:CD1	2.93	0.52
2:K:417:ASP:C	2:K:417:ASP:OD1	2.47	0.52
2:B:416:LEU:CD1	2:B:416:LEU:N	2.71	0.52
1:D:634:PHE:HE1	1:D:642:ARG:HH11	1.57	0.52
2:H:422:ASP:CB	2:H:423:PRO:HD2	2.40	0.52
2:C:47:ASN:C	2:C:47:ASN:ND2	2.59	0.52
1:A:560:PHE:CZ	1:A:581:MET:CE	2.92	0.52
2:H:422:ASP:O	2:H:423:PRO:O	2.28	0.52
2:K:295:SER:O	2:K:415:GLN:NE2	2.43	0.52
2:B:37:PRO:HD3	2:B:120:SER:HA	1.91	0.52
2:B:415:GLN:C	2:B:416:LEU:CD1	2.78	0.52
1:D:520:GLU:HG3	1:D:521:ILE:N	2.25	0.52
2:C:311:GLU:O	2:C:312:GLY:C	2.48	0.52
2:B:288:PHE:HE1	2:B:405:PHE:HE2	1.57	0.51
2:K:425:SER:OG	2:K:426:ASP:CA	2.33	0.51
2:K:290:GLU:OE1	2:K:295:SER:OG	2.25	0.51
2:C:299:CYS:HA	2:C:300:PHE:HB3	1.92	0.51
2:H:414:TYR:HE2	2:H:416:LEU:CD2	2.22	0.51
2:C:422:ASP:O	2:C:423:PRO:O	2.28	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:J:474:ALA:HA	1:J:585:THR:OG1	2.09	0.51
1:J:538:ASP:C	1:J:540:ASP:H	2.13	0.51
2:H:15:ARG:HG2	2:H:137:LEU:HD13	1.93	0.51
1:J:567:LEU:HD11	1:J:654:VAL:HG13	1.92	0.51
1:D:633:ARG:HA	1:D:657:VAL:HG22	1.91	0.51
2:K:422:ASP:O	2:K:423:PRO:O	2.29	0.51
1:E:631:MET:HG3	1:E:660:ILE:CD1	2.40	0.51
2:C:288:PHE:HE1	2:C:405:PHE:HE2	1.57	0.51
2:B:422:ASP:O	2:B:423:PRO:O	2.29	0.51
2:B:414:TYR:HD2	2:B:414:TYR:O	1.92	0.51
2:C:38:PRO:HB3	2:C:88:ASN:CA	2.14	0.51
2:C:59:ILE:O	2:C:60:LYS:HB2	2.11	0.51
2:H:414:TYR:CD2	2:H:414:TYR:O	2.64	0.51
1:E:614:THR:HG22	1:E:617:MET:SD	2.51	0.51
2:B:47:ASN:ND2	2:B:47:ASN:C	2.59	0.51
1:D:560:PHE:CZ	1:D:581:MET:CE	2.93	0.51
2:B:91:VAL:O	2:B:110:ASP:HA	2.11	0.51
1:A:602:THR:HG23	1:A:604:ARG:CD	2.41	0.51
2:K:299:CYS:HA	2:K:300:PHE:O	2.11	0.51
1:D:619:ILE:HG12	1:D:620:ILE:N	2.25	0.51
2:C:99:MET:HE3	2:C:105:ARG:HB3	1.93	0.51
2:H:295:SER:O	2:H:415:GLN:NE2	2.43	0.50
2:H:414:TYR:HE2	2:H:416:LEU:CD1	2.17	0.50
2:H:421:LEU:CD2	2:H:421:LEU:O	2.40	0.50
2:H:92:ILE:HG12	2:H:110:ASP:HB3	1.93	0.50
1:E:566:ILE:HG22	1:E:615:LYS:HA	1.93	0.50
2:H:319:CYS:SG	2:H:323:LEU:HD12	2.50	0.50
2:H:414:TYR:CE2	2:H:416:LEU:HD21	2.45	0.50
2:C:306:MET:HA	2:C:306:MET:CE	2.41	0.50
1:A:520:GLU:HG3	1:A:521:ILE:N	2.26	0.50
2:C:91:VAL:O	2:C:110:ASP:HA	2.12	0.50
1:D:569:LEU:HD23	1:D:652:VAL:CG1	2.42	0.50
1:E:573:LEU:HD22	1:E:577:TYR:CD2	2.47	0.50
2:C:414:TYR:O	2:C:414:TYR:HD2	1.93	0.50
2:B:306:MET:CE	2:B:306:MET:HA	2.41	0.50
1:E:469:PHE:O	1:E:548:THR:HA	2.12	0.50
1:E:474:ALA:HA	1:E:585:THR:OG1	2.11	0.50
1:E:596:LEU:O	1:E:609:PRO:HB3	2.11	0.50
1:E:572:ILE:HG12	1:E:573:LEU:N	2.26	0.49
2:B:59:ILE:O	2:B:60:LYS:HB2	2.10	0.49
2:B:311:GLU:O	2:B:312:GLY:C	2.48	0.49
2:B:300:PHE:H	2:B:409:GLY:HA3	1.73	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:J:580:VAL:CG2	1:J:648:GLN:HG3	2.38	0.49
2:C:294:ASP:CG	2:C:294:ASP:O	2.50	0.49
1:E:538:ASP:C	1:E:540:ASP:H	2.14	0.49
1:J:573:LEU:HD22	1:J:577:TYR:CD2	2.47	0.49
1:A:613:ALA:HA	1:A:617:MET:CE	2.43	0.49
1:D:613:ALA:HA	1:D:617:MET:CE	2.43	0.49
1:J:631:MET:HG3	1:J:660:ILE:CD1	2.42	0.49
2:B:315:GLU:HB3	2:B:411:VAL:O	2.13	0.49
1:D:584:HIS:CD2	1:D:586:ALA:H	2.30	0.49
1:E:590:VAL:HG23	1:E:625:THR:HG22	1.95	0.49
1:E:521:ILE:HD11	1:E:524:SER:OG	2.12	0.49
2:H:425:SER:OG	2:H:426:ASP:CA	2.32	0.49
1:A:572:ILE:CD1	1:A:612:PHE:CG	2.80	0.49
1:J:647:ASP:C	1:J:649:GLY:H	2.08	0.49
1:E:646:ARG:HD3	1:E:648:GLN:OE1	2.13	0.49
2:K:311:GLU:O	2:K:312:GLY:C	2.51	0.49
2:B:86:PRO:O	2:B:87:ASP:C	2.50	0.49
2:B:125:ASP:CG	2:B:126:ASN:H	2.16	0.49
2:B:99:MET:HE3	2:B:105:ARG:HB3	1.94	0.48
2:C:62:ARG:HB3	2:C:62:ARG:NH1	2.28	0.48
1:J:631:MET:HG3	1:J:660:ILE:HD11	1.94	0.48
2:K:5:ALA:C	2:K:7:LYS:H	2.16	0.48
2:H:10:GLU:O	2:H:14:ILE:HG12	2.13	0.48
2:K:274:ASN:HB3	2:K:277:TYR:HB2	1.94	0.48
2:C:422:ASP:O	2:C:423:PRO:C	2.51	0.48
1:D:574:THR:HG22	1:D:612:PHE:HA	1.95	0.48
1:A:572:ILE:HG13	1:A:612:PHE:HD1	1.78	0.48
1:A:468:PRO:HB3	1:A:552:ASN:HB2	1.95	0.48
1:D:468:PRO:HB3	1:D:552:ASN:HB2	1.95	0.48
2:C:125:ASP:CG	2:C:126:ASN:H	2.17	0.48
2:H:315:GLU:HA	2:H:413:ARG:HE	1.78	0.48
1:E:572:ILE:HD11	1:E:612:PHE:HB3	1.87	0.48
2:C:300:PHE:H	2:C:409:GLY:HA3	1.73	0.48
1:D:566:ILE:HG22	1:D:615:LYS:HA	1.96	0.48
2:H:315:GLU:CB	2:H:413:ARG:HE	2.26	0.48
2:C:86:PRO:O	2:C:87:ASP:C	2.51	0.48
1:E:632:GLU:HB2	1:E:640:MET:HB2	1.96	0.48
1:E:650:THR:OG1	1:E:651:THR:N	2.47	0.48
2:B:274:ASN:HB3	2:B:277:TYR:HB2	1.96	0.48
2:H:421:LEU:O	2:H:422:ASP:CG	2.52	0.48
1:J:574:THR:HG22	1:J:612:PHE:HA	1.94	0.48
2:K:294:ASP:O	2:K:294:ASP:OD2	2.31	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:H:422:ASP:O	2:H:423:PRO:C	2.51	0.48
1:E:589:GLU:HG3	1:E:589:GLU:O	2.12	0.48
2:C:274:ASN:HB3	2:C:277:TYR:HB2	1.95	0.48
2:K:421:LEU:O	2:K:422:ASP:CG	2.52	0.48
2:B:62:ARG:HB3	2:B:62:ARG:NH1	2.28	0.48
2:C:315:GLU:HB3	2:C:411:VAL:O	2.13	0.48
2:H:34:LEU:HD23	2:H:92:ILE:HD12	1.96	0.48
1:A:555:HIS:HD2	1:A:632:GLU:OE2	1.97	0.48
2:H:414:TYR:CD2	2:H:414:TYR:C	2.87	0.48
2:C:413:ARG:C	2:C:414:TYR:CD1	2.87	0.48
2:C:17:LEU:O	2:C:20:GLN:HG2	2.14	0.48
2:K:422:ASP:O	2:K:423:PRO:C	2.51	0.47
2:B:417:ASP:O	2:B:418:LEU:O	2.32	0.47
1:A:569:LEU:CD2	1:A:573:LEU:HB2	2.44	0.47
1:A:574:THR:HG22	1:A:612:PHE:HA	1.95	0.47
1:E:574:THR:HG22	1:E:612:PHE:HA	1.96	0.47
1:E:598:LYS:HB2	1:E:610:PRO:HG3	1.95	0.47
2:H:302:VAL:HG23	2:H:303:VAL:H	1.79	0.47
2:C:70:SER:O	2:C:74:SER:HB2	2.14	0.47
1:D:569:LEU:HD11	1:D:573:LEU:HB2	1.96	0.47
1:J:568:GLU:O	1:J:652:VAL:HG12	2.14	0.47
2:C:293:LEU:O	2:C:294:ASP:C	2.53	0.47
2:B:422:ASP:O	2:B:423:PRO:C	2.52	0.47
1:E:572:ILE:CD1	1:E:612:PHE:CG	2.95	0.47
2:B:31:MET:HG3	2:B:126:ASN:C	2.35	0.47
2:H:75:THR:O	2:H:75:THR:HG22	2.14	0.47
2:C:347:GLU:H	2:C:347:GLU:CD	2.18	0.47
2:B:291:ILE:HG12	2:B:298:TYR:CE1	2.50	0.47
2:C:291:ILE:HG12	2:C:298:TYR:CE1	2.50	0.47
1:J:602:THR:OG1	1:J:602:THR:O	2.29	0.47
1:J:646:ARG:HD3	1:J:648:GLN:HG2	1.97	0.47
1:J:632:GLU:HB2	1:J:640:MET:HB2	1.96	0.47
1:A:560:PHE:CZ	1:A:581:MET:HE3	2.43	0.47
1:D:602:THR:HG23	1:D:604:ARG:CD	2.42	0.47
2:C:96:GLU:HB3	2:C:106:LYS:HG3	1.95	0.47
1:D:569:LEU:HD23	1:D:652:VAL:HB	1.97	0.47
1:A:478:LYS:HE2	1:A:482:THR:OG1	2.15	0.47
1:E:569:LEU:HD22	1:E:652:VAL:HB	1.95	0.47
1:J:619:ILE:HD11	1:J:621:ALA:HB2	1.97	0.47
1:D:609:PRO:HA	1:D:610:PRO:HD3	1.82	0.47
1:D:646:ARG:HB3	1:D:648:GLN:HG2	1.97	0.47
2:K:416:LEU:O	2:K:418:LEU:HD22	2.15	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:646:ARG:HG2	1:E:648:GLN:HG2	1.95	0.47
2:K:308:ALA:HB2	2:K:412:MET:HE3	1.96	0.47
1:D:517:ALA:O	1:D:518:ASP:HB2	2.15	0.47
2:C:16:ARG:HA	2:C:16:ARG:HD2	1.71	0.47
2:B:413:ARG:C	2:B:414:TYR:CD1	2.88	0.46
1:J:555:HIS:CD2	1:J:632:GLU:OE2	2.62	0.46
2:K:283:LEU:O	2:K:286:ARG:HB3	2.15	0.46
2:C:411:VAL:CG2	2:C:411:VAL:O	2.63	0.46
1:E:580:VAL:CG2	1:E:648:GLN:HG3	2.45	0.46
1:J:572:ILE:CD1	1:J:612:PHE:CD1	2.98	0.46
2:K:308:ALA:HB1	2:K:412:MET:HE3	1.97	0.46
2:B:17:LEU:O	2:B:20:GLN:HG2	2.15	0.46
1:D:572:ILE:HG13	1:D:612:PHE:HD1	1.79	0.46
1:J:573:LEU:CD2	1:J:577:TYR:CD1	2.99	0.46
2:K:291:ILE:HG12	2:K:298:TYR:CD1	2.51	0.46
2:C:417:ASP:O	2:C:418:LEU:O	2.32	0.46
1:J:478:LYS:O	1:J:479:ASP:CB	2.59	0.46
1:J:500:LEU:HB2	1:J:550:THR:HG22	1.96	0.46
1:E:500:LEU:HB2	1:E:550:THR:HG22	1.97	0.46
1:D:555:HIS:HD2	1:D:632:GLU:OE2	1.98	0.46
1:D:562:ALA:HA	1:D:658:VAL:HG12	1.97	0.46
1:A:517:ALA:O	1:A:518:ASP:HB2	2.15	0.46
1:J:581:MET:SD	1:J:583:ILE:HG13	2.56	0.46
1:J:566:ILE:N	1:J:566:ILE:HD12	2.30	0.46
2:H:311:GLU:O	2:H:312:GLY:C	2.54	0.46
1:E:600:ASP:HB3	1:E:601:LYS:H	1.62	0.46
2:B:16:ARG:HA	2:B:16:ARG:HD2	1.72	0.46
1:E:573:LEU:CD1	1:E:573:LEU:C	2.84	0.46
2:B:62:ARG:CZ	2:B:62:ARG:HB3	2.46	0.46
1:E:641:GLY:HA2	1:E:657:VAL:HG13	1.98	0.46
1:A:562:ALA:HA	1:A:658:VAL:HG12	1.97	0.46
1:D:569:LEU:HD23	1:D:652:VAL:HG11	1.97	0.46
1:A:566:ILE:HG22	1:A:615:LYS:HA	1.98	0.46
2:K:91:VAL:O	2:K:110:ASP:HA	2.15	0.46
2:C:348:GLU:HG3	2:C:349:LYS:N	2.31	0.46
2:B:347:GLU:CD	2:B:347:GLU:H	2.18	0.46
1:J:650:THR:OG1	1:J:651:THR:N	2.49	0.46
1:D:478:LYS:HE2	1:D:482:THR:OG1	2.16	0.46
1:A:641:GLY:HA2	1:A:657:VAL:CG1	2.46	0.46
1:J:561:ILE:O	1:J:658:VAL:HG12	2.16	0.46
1:A:558:THR:O	1:A:624:GLU:HA	2.16	0.45
1:J:602:THR:HG23	1:J:604:ARG:CD	2.44	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:C:31:MET:HG3	2:C:126:ASN:C	2.36	0.45
2:K:392:SER:OG	2:K:394:ARG:HG2	2.16	0.45
2:B:421:LEU:CD2	2:B:421:LEU:O	2.43	0.45
2:B:417:ASP:CA	2:B:418:LEU:HD22	2.46	0.45
2:B:276:LYS:HD3	2:B:276:LYS:N	2.32	0.45
2:C:421:LEU:O	2:C:422:ASP:CG	2.55	0.45
2:B:411:VAL:O	2:B:411:VAL:CG2	2.62	0.45
1:D:641:GLY:HA2	1:D:657:VAL:CG1	2.46	0.45
2:H:315:GLU:CA	2:H:413:ARG:HE	2.30	0.45
1:A:584:HIS:CD2	1:A:586:ALA:H	2.31	0.45
1:E:634:PHE:HA	1:E:641:GLY:HA3	1.98	0.45
1:D:631:MET:HG3	1:D:660:ILE:CD1	2.46	0.45
2:B:99:MET:HB3	2:B:103:LYS:O	2.17	0.45
1:A:613:ALA:HA	1:A:617:MET:HE1	1.97	0.45
1:A:631:MET:HG3	1:A:660:ILE:CD1	2.46	0.45
1:J:538:ASP:C	1:J:540:ASP:N	2.68	0.45
2:C:99:MET:HB3	2:C:103:LYS:O	2.17	0.45
2:H:417:ASP:OD1	2:H:417:ASP:C	2.55	0.45
1:D:573:LEU:CD1	1:D:573:LEU:C	2.85	0.45
2:K:315:GLU:O	2:K:387:ASN:CB	2.65	0.45
1:D:474:ALA:HA	1:D:585:THR:OG1	2.16	0.45
1:E:538:ASP:C	1:E:540:ASP:N	2.70	0.45
2:K:302:VAL:HG23	2:K:303:VAL:H	1.82	0.45
2:H:283:LEU:O	2:H:286:ARG:HB3	2.16	0.45
1:D:515:ASP:OD1	1:D:515:ASP:C	2.55	0.45
2:H:131:GLU:HA	2:H:134:ALA:HB3	1.99	0.45
2:H:315:GLU:O	2:H:387:ASN:CB	2.64	0.45
2:C:417:ASP:O	2:C:418:LEU:C	2.55	0.45
1:A:474:ALA:HA	1:A:585:THR:OG1	2.16	0.45
2:B:348:GLU:HG3	2:B:349:LYS:N	2.32	0.45
1:D:582:HIS:CD2	1:D:582:HIS:N	2.85	0.45
1:A:582:HIS:CD2	1:A:582:HIS:N	2.84	0.45
1:J:573:LEU:C	1:J:573:LEU:CD1	2.85	0.45
1:A:602:THR:CG2	1:A:604:ARG:HD2	2.42	0.45
2:C:134:ALA:HA	2:C:137:LEU:CD1	2.47	0.45
1:J:634:PHE:CB	1:J:656:LYS:HB2	2.47	0.45
2:B:421:LEU:O	2:B:422:ASP:CG	2.55	0.45
1:A:573:LEU:C	1:A:573:LEU:CD1	2.85	0.45
2:H:308:ALA:HB2	2:H:412:MET:HE3	1.97	0.45
1:D:558:THR:O	1:D:624:GLU:HA	2.16	0.44
1:D:559:ARG:HG2	1:D:624:GLU:HG2	2.00	0.44
2:C:295:SER:OG	2:C:297:LYS:CG	2.66	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:602:THR:CG2	1:D:604:ARG:HD2	2.45	0.44
1:E:567:LEU:CD1	1:E:654:VAL:HG13	2.47	0.44
1:J:538:ASP:O	1:J:540:ASP:N	2.50	0.44
1:E:485:GLU:HG3	1:E:485:GLU:O	2.17	0.44
2:C:62:ARG:CZ	2:C:62:ARG:HB3	2.47	0.44
2:B:134:ALA:HA	2:B:137:LEU:CD1	2.47	0.44
2:H:291:ILE:HG12	2:H:298:TYR:CE1	2.51	0.44
2:C:421:LEU:CD2	2:C:421:LEU:O	2.43	0.44
1:A:559:ARG:HG2	1:A:624:GLU:HG2	1.98	0.44
1:E:555:HIS:CD2	1:E:632:GLU:OE2	2.67	0.44
2:H:274:ASN:HB3	2:H:277:TYR:HB2	1.98	0.44
2:C:417:ASP:CA	2:C:418:LEU:HD22	2.47	0.44
1:D:593:ALA:HB2	1:D:624:GLU:HG3	1.99	0.44
1:E:581:MET:SD	1:E:583:ILE:HG13	2.58	0.44
1:D:488:ILE:CG2	1:D:526:CYS:HA	2.48	0.44
1:J:584:HIS:C	1:J:586:ALA:H	2.21	0.44
2:K:421:LEU:O	2:K:422:ASP:CB	2.65	0.44
1:E:560:PHE:CZ	1:E:581:MET:CE	3.00	0.44
2:H:308:ALA:O	2:H:311:GLU:O	2.35	0.44
1:J:594:LYS:O	1:J:621:ALA:HB1	2.18	0.44
1:E:561:ILE:O	1:E:658:VAL:HG12	2.17	0.44
2:B:416:LEU:N	2:B:416:LEU:HD12	2.32	0.44
1:E:559:ARG:HB3	1:E:661:LEU:HD12	2.00	0.44
2:B:113:PRO:HG2	2:B:116:PRO:CG	2.46	0.44
2:H:18:VAL:HG23	2:H:19:LYS:N	2.23	0.44
1:E:634:PHE:CB	1:E:656:LYS:HB2	2.48	0.44
2:H:308:ALA:HB1	2:H:412:MET:HE3	1.99	0.44
1:D:484:LEU:HD12	1:D:534:VAL:HG11	1.99	0.44
2:B:41:GLN:HE21	2:B:41:GLN:HB2	1.68	0.44
2:B:291:ILE:HG12	2:B:298:TYR:CD1	2.53	0.44
1:A:584:HIS:HD2	1:A:586:ALA:N	2.16	0.44
1:D:580:VAL:HG21	1:D:648:GLN:HG3	1.99	0.44
2:H:343:THR:OG1	2:H:344:LYS:N	2.51	0.44
1:A:484:LEU:HD12	1:A:534:VAL:HG11	2.00	0.44
1:A:515:ASP:C	1:A:515:ASP:OD1	2.56	0.44
1:J:618:LYS:HE2	1:J:618:LYS:HB2	1.76	0.44
1:E:619:ILE:HD11	1:E:621:ALA:HB2	1.99	0.44
2:B:420:MET:HG3	2:B:421:LEU:HD12	1.87	0.43
1:J:602:THR:O	1:J:603:ASN:C	2.56	0.43
1:E:613:ALA:HB1	1:E:617:MET:HE2	2.00	0.43
2:B:306:MET:HE2	2:B:306:MET:HA	1.99	0.43
1:A:580:VAL:HG21	1:A:648:GLN:HG3	2.00	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:C:276:LYS:HD3	2:C:276:LYS:N	2.32	0.43
1:J:572:ILE:HG12	1:J:573:LEU:H	1.83	0.43
2:B:288:PHE:HE1	2:B:405:PHE:CE2	2.35	0.43
2:B:417:ASP:O	2:B:418:LEU:C	2.55	0.43
2:C:288:PHE:HE1	2:C:405:PHE:CE2	2.35	0.43
1:D:532:LEU:HD12	1:D:547:LEU:CD1	2.48	0.43
1:D:572:ILE:CG1	1:D:573:LEU:N	2.75	0.43
2:K:421:LEU:CD2	2:K:421:LEU:O	2.42	0.43
2:C:113:PRO:HG2	2:C:116:PRO:CG	2.47	0.43
2:H:274:ASN:HB2	2:H:278:VAL:HG13	2.01	0.43
1:A:532:LEU:HD12	1:A:547:LEU:CD1	2.49	0.43
1:J:598:LYS:HB3	1:J:598:LYS:HE2	1.78	0.43
2:B:96:GLU:HA	2:B:106:LYS:HA	2.01	0.43
1:J:569:LEU:HB3	1:J:571:SER:O	2.18	0.43
1:A:619:ILE:HG12	1:A:620:ILE:H	1.83	0.43
1:E:594:LYS:O	1:E:621:ALA:HB1	2.19	0.43
2:H:422:ASP:HB3	2:H:423:PRO:HD2	2.01	0.43
2:B:294:ASP:O	2:B:296:GLY:N	2.52	0.43
2:H:349:LYS:NZ	2:H:350:ASP:O	2.52	0.43
2:K:326:ILE:HD12	2:K:328:TYR:HE2	1.82	0.43
1:J:600:ASP:HB3	1:J:601:LYS:H	1.63	0.43
2:H:294:ASP:O	2:H:294:ASP:OD2	2.36	0.43
2:H:417:ASP:OD1	2:H:418:LEU:N	2.52	0.43
2:C:291:ILE:HG12	2:C:298:TYR:CD1	2.53	0.43
2:K:294:ASP:O	2:K:294:ASP:CG	2.56	0.43
1:E:569:LEU:HB3	1:E:571:SER:O	2.19	0.43
2:H:320:PHE:HA	2:H:391:VAL:O	2.19	0.43
1:A:488:ILE:HB	1:A:526:CYS:HA	2.00	0.42
1:E:598:LYS:HE2	1:E:598:LYS:HB3	1.76	0.42
2:H:421:LEU:CD2	2:H:421:LEU:C	2.86	0.42
1:J:634:PHE:HE1	1:J:642:ARG:HH11	1.66	0.42
1:A:488:ILE:CG2	1:A:526:CYS:HA	2.49	0.42
1:A:619:ILE:HD11	1:A:621:ALA:HB2	2.00	0.42
2:C:61:SER:O	2:C:62:ARG:C	2.58	0.42
1:E:634:PHE:HE1	1:E:642:ARG:HH11	1.65	0.42
1:D:619:ILE:HD11	1:D:621:ALA:HB2	2.01	0.42
2:C:306:MET:HA	2:C:306:MET:HE2	2.00	0.42
1:D:537:ASP:OD2	1:D:538:ASP:N	2.53	0.42
2:K:421:LEU:C	2:K:421:LEU:CD2	2.86	0.42
1:J:560:PHE:CZ	1:J:581:MET:CE	3.01	0.42
1:A:537:ASP:OD2	1:A:538:ASP:N	2.53	0.42
2:K:320:PHE:HA	2:K:391:VAL:O	2.20	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:479:ASP:O	1:D:480:LEU:C	2.58	0.42
2:K:422:ASP:HB3	2:K:423:PRO:HD2	2.01	0.42
2:C:416:LEU:HD12	2:C:416:LEU:N	2.30	0.42
1:A:593:ALA:HB2	1:A:624:GLU:HG3	2.00	0.42
1:J:634:PHE:HA	1:J:641:GLY:HA3	2.02	0.42
1:J:567:LEU:HD12	1:J:653:ALA:HA	2.02	0.42
2:H:421:LEU:O	2:H:422:ASP:CB	2.66	0.42
2:C:298:TYR:HA	2:C:410:ALA:O	2.20	0.42
2:B:61:SER:O	2:B:62:ARG:C	2.58	0.42
2:C:34:LEU:C	2:C:35:ILE:HD12	2.39	0.42
2:C:300:PHE:CD1	2:C:300:PHE:C	2.92	0.42
1:J:614:THR:HG22	1:J:617:MET:SD	2.59	0.42
2:C:98:ILE:HA	2:C:104:THR:HG22	2.02	0.42
1:E:573:LEU:CD2	1:E:577:TYR:CD1	3.03	0.42
2:B:112:GLU:HA	2:B:113:PRO:HD3	1.75	0.42
1:A:560:PHE:CZ	1:A:581:MET:HE1	2.55	0.42
2:H:326:ILE:HD12	2:H:328:TYR:HE2	1.83	0.42
2:K:315:GLU:O	2:K:387:ASN:HB3	2.20	0.41
2:B:280:GLU:CD	2:B:392:SER:H	2.23	0.41
2:K:10:GLU:O	2:K:14:ILE:HG12	2.20	0.41
2:H:315:GLU:O	2:H:387:ASN:HB3	2.20	0.41
1:A:569:LEU:HD21	1:A:573:LEU:HB2	2.02	0.41
1:E:647:ASP:C	1:E:649:GLY:N	2.67	0.41
1:A:618:LYS:HB2	1:A:618:LYS:HE2	1.84	0.41
1:E:517:ALA:O	1:E:518:ASP:CB	2.64	0.41
1:D:488:ILE:HB	1:D:526:CYS:HA	2.01	0.41
1:E:567:LEU:HD11	1:E:654:VAL:HG13	2.01	0.41
2:B:34:LEU:C	2:B:35:ILE:HD12	2.40	0.41
2:K:325:MET:HE2	2:K:341:TYR:HB3	1.94	0.41
1:J:647:ASP:C	1:J:649:GLY:N	2.71	0.41
1:A:600:ASP:O	1:A:618:LYS:CE	2.67	0.41
1:E:602:THR:O	1:E:603:ASN:C	2.59	0.41
1:A:479:ASP:O	1:A:480:LEU:C	2.58	0.41
1:E:518:ASP:OD1	1:E:531:ARG:NH2	2.52	0.41
1:E:567:LEU:HD12	1:E:653:ALA:HA	2.01	0.41
1:A:605:LYS:HB2	1:A:605:LYS:HE2	1.78	0.41
1:J:559:ARG:HB3	1:J:661:LEU:HD12	2.01	0.41
2:C:325:MET:HE3	2:C:341:TYR:HB3	2.00	0.41
1:A:602:THR:O	1:A:603:ASN:C	2.59	0.41
2:B:307:ASN:ND2	2:B:420:MET:O	2.54	0.41
1:D:647:ASP:C	1:D:649:GLY:N	2.73	0.41
2:H:325:MET:HB3	2:H:325:MET:HE3	1.89	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:573:LEU:HD22	1:A:577:TYR:CD2	2.56	0.41
1:E:646:ARG:HG2	1:E:648:GLN:N	2.35	0.41
2:B:300:PHE:C	2:B:300:PHE:CD1	2.93	0.41
1:D:584:HIS:CD2	1:D:585:THR:HG22	2.56	0.41
2:C:343:THR:OG1	2:C:344:LYS:N	2.53	0.41
1:A:537:ASP:OD2	1:A:537:ASP:C	2.59	0.41
2:B:421:LEU:CD2	2:B:421:LEU:C	2.87	0.41
1:A:600:ASP:HB3	1:A:601:LYS:H	1.60	0.41
1:J:526:CYS:C	1:J:528:ASP:H	2.24	0.41
2:B:36:ILE:HA	2:B:37:PRO:HD2	1.95	0.41
1:D:619:ILE:HG12	1:D:620:ILE:H	1.85	0.41
2:C:280:GLU:CD	2:C:392:SER:H	2.23	0.41
2:B:93:TYR:HD1	2:B:128:PHE:CD1	2.38	0.41
1:D:602:THR:O	1:D:603:ASN:C	2.59	0.41
1:E:509:GLU:O	1:E:534:VAL:HA	2.21	0.41
2:H:5:ALA:C	2:H:7:LYS:H	2.23	0.41
2:H:386:ALA:CA	2:H:387:ASN:CB	2.75	0.41
2:C:307:ASN:ND2	2:C:420:MET:O	2.54	0.41
1:E:572:ILE:CD1	1:E:612:PHE:CB	2.80	0.41
1:D:600:ASP:O	1:D:618:LYS:CE	2.67	0.41
2:C:96:GLU:HA	2:C:106:LYS:HA	2.03	0.41
1:D:537:ASP:OD2	1:D:537:ASP:C	2.59	0.41
2:B:38:PRO:CB	2:B:88:ASN:HA	2.15	0.41
2:B:98:ILE:HA	2:B:104:THR:HG22	2.03	0.41
1:E:572:ILE:CG1	1:E:612:PHE:CD1	3.03	0.40
2:B:325:MET:HE1	2:B:341:TYR:CB	2.43	0.40
1:J:613:ALA:HB1	1:J:617:MET:HE2	2.03	0.40
1:D:584:HIS:HD2	1:D:586:ALA:N	2.16	0.40
1:E:538:ASP:O	1:E:540:ASP:N	2.54	0.40
2:K:343:THR:OG1	2:K:344:LYS:N	2.52	0.40
1:D:511:THR:O	1:D:512:ALA:HB2	2.22	0.40
1:J:544:GLY:HA2	1:J:585:THR:CG2	2.36	0.40
1:A:581:MET:HE2	1:A:643:PHE:CD1	2.56	0.40
2:B:343:THR:OG1	2:B:344:LYS:N	2.53	0.40
1:A:646:ARG:CB	1:A:648:GLN:HG2	2.51	0.40
1:J:496:ASN:HA	1:J:510:VAL:O	2.21	0.40
2:B:117:ILE:O	2:B:119:THR:N	2.54	0.40
2:H:424:GLU:OE1	2:H:424:GLU:N	2.35	0.40
2:B:414:TYR:CD1	2:B:414:TYR:N	2.81	0.40
1:D:641:GLY:HA2	1:D:657:VAL:HG13	2.02	0.40
1:E:584:HIS:C	1:E:586:ALA:H	2.25	0.40
1:J:509:GLU:O	1:J:534:VAL:HA	2.21	0.40

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:C:421:LEU:O	2:C:422:ASP:CB	2.70	0.40
1:E:572:ILE:CD1	1:E:612:PHE:CD1	3.04	0.40
2:H:300:PHE:C	2:H:300:PHE:CD1	2.94	0.40
1:J:641:GLY:HA2	1:J:657:VAL:HG13	2.03	0.40
1:J:599:LEU:HD11	1:J:620:ILE:HG22	2.02	0.40
2:B:372:LEU:O	2:B:373:LEU:C	2.59	0.40
2:H:392:SER:OG	2:H:394:ARG:HG2	2.22	0.40
2:K:414:TYR:CE2	2:K:416:LEU:HD11	2.57	0.40
2:B:94:CYS:HA	2:B:107:LEU:O	2.21	0.40
1:D:600:ASP:HB3	1:D:601:LYS:H	1.61	0.40
1:A:584:HIS:CD2	1:A:585:THR:HG22	2.56	0.40
2:K:274:ASN:HB2	2:K:278:VAL:HG13	2.04	0.40
1:D:646:ARG:CB	1:D:648:GLN:HG2	2.51	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	194/201 (96%)	170 (88%)	20 (10%)	4 (2%)	11	62
1	D	194/201 (96%)	170 (88%)	20 (10%)	4 (2%)	11	62
1	E	194/201 (96%)	164 (84%)	23 (12%)	7 (4%)	5	49
1	J	194/201 (96%)	166 (86%)	23 (12%)	5 (3%)	8	57
2	B	253/441 (57%)	201 (79%)	33 (13%)	19 (8%)	2	24
2	C	253/441 (57%)	203 (80%)	31 (12%)	19 (8%)	2	24
2	H	155/441 (35%)	113 (73%)	30 (19%)	12 (8%)	1	22
2	K	155/441 (35%)	114 (74%)	31 (20%)	10 (6%)	2	29
All	All	1592/2568 (62%)	1301 (82%)	211 (13%)	80 (5%)	3	37

All (80) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	479	ASP
1	A	575	THR
1	A	648	GLN
2	C	87	ASP
2	C	299	CYS
2	C	413	ARG
2	C	423	PRO
1	D	479	ASP
1	D	575	THR
1	D	648	GLN
2	B	87	ASP
2	B	299	CYS
2	B	413	ARG
2	B	423	PRO
1	E	479	ASP
1	E	575	THR
1	E	648	GLN
2	H	298	TYR
2	H	423	PRO
1	J	479	ASP
1	J	575	THR
1	J	648	GLN
2	K	299	CYS
2	K	423	PRO
2	C	29	THR
2	C	113	PRO
2	C	312	GLY
2	C	418	LEU
2	B	5	ALA
2	B	29	THR
2	B	113	PRO
2	B	312	GLY
2	B	418	LEU
2	H	7	LYS
2	H	76	ARG
2	H	420	MET
2	K	298	TYR
2	K	420	MET
2	C	5	ALA
2	C	28	GLY
2	C	298	TYR
2	C	420	MET
2	C	424	GLU

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Mol	Chain	Res	Type
2	B	28	GLY
2	B	298	TYR
2	B	420	MET
2	B	424	GLU
1	E	606	SER
1	E	608	LYS
2	H	312	GLY
2	H	387	ASN
2	H	424	GLU
1	J	570	PRO
1	J	603	ASN
2	K	47	ASN
2	K	300	PHE
2	K	312	GLY
2	K	387	ASN
2	K	424	GLU
1	A	603	ASN
2	C	387	ASN
2	C	415	GLN
1	D	603	ASN
2	B	387	ASN
2	B	415	GLN
1	E	603	ASN
2	H	46	SER
2	H	421	LEU
2	K	421	LEU
2	C	118	ASN
2	C	412	MET
2	C	421	LEU
2	B	118	ASN
2	B	412	MET
2	B	421	LEU
2	H	414	TYR
2	B	406	GLY
2	H	47	ASN
2	C	406	GLY
1	E	525	ILE

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	172/175 (98%)	153 (89%)	19 (11%)	9	43
1	D	172/175 (98%)	151 (88%)	21 (12%)	7	36
1	E	172/175 (98%)	144 (84%)	28 (16%)	3	21
1	J	172/175 (98%)	145 (84%)	27 (16%)	4	23
2	B	232/383 (61%)	208 (90%)	24 (10%)	10	47
2	C	232/383 (61%)	210 (90%)	22 (10%)	12	51
2	H	154/383 (40%)	144 (94%)	10 (6%)	24	71
2	K	154/383 (40%)	141 (92%)	13 (8%)	16	59
All	All	1460/2232 (65%)	1296 (89%)	164 (11%)	9	41

All (164) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	526	CYS
1	A	543	THR
1	A	559	ARG
1	A	583	ILE
1	A	584	HIS
1	A	589	GLU
1	A	600	ASP
1	A	602	THR
1	A	604	ARG
1	A	608	LYS
1	A	611	MET
1	A	619	ILE
1	A	620	ILE
1	A	629	VAL
1	A	648	GLN
1	A	650	THR
1	A	656	LYS
1	A	657	VAL
1	A	662	ASP
2	C	41	GLN
2	C	47	ASN
2	C	62	ARG
2	C	74	SER
2	C	107	LEU

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Mol	Chain	Res	Type
2	C	108	ASN
2	C	115	LYS
2	C	118	ASN
2	C	276	LYS
2	C	279	GLN
2	C	298	TYR
2	C	302	VAL
2	C	323	LEU
2	C	340	THR
2	C	373	LEU
2	C	374	SER
2	C	413	ARG
2	C	415	GLN
2	C	417	ASP
2	C	418	LEU
2	C	421	LEU
2	C	425	SER
1	D	526	CYS
1	D	542	GLN
1	D	543	THR
1	D	559	ARG
1	D	569	LEU
1	D	583	ILE
1	D	584	HIS
1	D	589	GLU
1	D	600	ASP
1	D	602	THR
1	D	604	ARG
1	D	608	LYS
1	D	611	MET
1	D	619	ILE
1	D	620	ILE
1	D	629	VAL
1	D	648	GLN
1	D	650	THR
1	D	656	LYS
1	D	657	VAL
1	D	662	ASP
2	B	41	GLN
2	B	47	ASN
2	B	61	SER
2	B	62	ARG

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Mol	Chain	Res	Type
2	B	74	SER
2	B	105	ARG
2	B	107	LEU
2	B	108	ASN
2	B	115	LYS
2	B	118	ASN
2	B	276	LYS
2	B	279	GLN
2	B	298	TYR
2	B	302	VAL
2	B	323	LEU
2	B	340	THR
2	B	373	LEU
2	B	374	SER
2	B	413	ARG
2	B	415	GLN
2	B	417	ASP
2	B	418	LEU
2	B	421	LEU
2	B	425	SER
1	E	475	SER
1	E	485	GLU
1	E	506	GLN
1	E	524	SER
1	E	526	CYS
1	E	543	THR
1	E	552	ASN
1	E	559	ARG
1	E	569	LEU
1	E	573	LEU
1	E	583	ILE
1	E	584	HIS
1	E	589	GLU
1	E	590	VAL
1	E	602	THR
1	E	604	ARG
1	E	606	SER
1	E	608	LYS
1	E	611	MET
1	E	619	ILE
1	E	620	ILE
1	E	623	LEU

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Mol	Chain	Res	Type
1	E	629	VAL
1	E	642	ARG
1	E	644	THR
1	E	645	LEU
1	E	650	THR
1	E	657	VAL
2	H	276	LYS
2	H	286	ARG
2	H	298	TYR
2	H	302	VAL
2	H	303	VAL
2	H	323	LEU
2	H	396	GLN
2	H	418	LEU
2	H	421	LEU
2	H	425	SER
1	J	475	SER
1	J	485	GLU
1	J	506	GLN
1	J	526	CYS
1	J	543	THR
1	J	552	ASN
1	J	559	ARG
1	J	569	LEU
1	J	573	LEU
1	J	583	ILE
1	J	584	HIS
1	J	589	GLU
1	J	590	VAL
1	J	602	THR
1	J	604	ARG
1	J	606	SER
1	J	608	LYS
1	J	611	MET
1	J	619	ILE
1	J	620	ILE
1	J	623	LEU
1	J	629	VAL
1	J	642	ARG
1	J	644	THR
1	J	645	LEU
1	J	650	THR

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Mol	Chain	Res	Type
1	J	657	VAL
2	K	78	ARG
2	K	131	GLU
2	K	276	LYS
2	K	286	ARG
2	K	298	TYR
2	K	302	VAL
2	K	303	VAL
2	K	323	LEU
2	K	396	GLN
2	K	413	ARG
2	K	418	LEU
2	K	421	LEU
2	K	425	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (43) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	496	ASN
1	A	552	ASN
1	A	555	HIS
1	A	563	GLN
1	A	584	HIS
2	C	25	HIS
2	C	41	GLN
2	C	47	ASN
2	C	88	ASN
2	C	108	ASN
2	C	118	ASN
2	C	285	GLN
2	C	310	GLN
2	C	346	GLN
2	C	396	GLN
2	C	400	GLN
2	C	415	GLN
1	D	496	ASN
1	D	552	ASN
1	D	555	HIS
1	D	563	GLN
1	D	584	HIS
2	B	41	GLN
2	B	47	ASN

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Mol	Chain	Res	Type
2	B	88	ASN
2	B	108	ASN
2	B	118	ASN
2	B	285	GLN
2	B	310	GLN
2	B	346	GLN
2	B	396	GLN
2	B	400	GLN
2	B	415	GLN
1	E	496	ASN
1	E	555	HIS
1	E	563	GLN
1	E	584	HIS
2	H	310	GLN
1	J	542	GLN
1	J	555	HIS
1	J	563	GLN
1	J	584	HIS
2	K	310	GLN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	196/201 (97%)	0.99	20 (10%) 7 5	69, 77, 80, 82	0
1	D	196/201 (97%)	0.68	5 (2%) 53 24	69, 77, 80, 82	0
1	E	196/201 (97%)	0.88	17 (8%) 10 6	49, 58, 67, 72	0
1	J	196/201 (97%)	0.57	1 (0%) 88 64	49, 58, 67, 72	0
2	B	261/441 (59%)	0.44	4 (1%) 70 36	66, 79, 83, 83	0
2	C	261/441 (59%)	3.13	149 (57%) 0 1	66, 79, 83, 83	0
2	H	175/441 (39%)	0.46	9 (5%) 27 11	66, 83, 116, 122	0
2	K	175/441 (39%)	1.96	74 (42%) 1 1	66, 83, 118, 138	0
All	All	1656/2568 (64%)	1.19	279 (16%) 2 2	49, 77, 98, 138	0

All (279) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	C	99	MET	25.0
2	C	326	ILE	12.4
2	C	371	MET	11.6
2	C	328	TYR	11.0
2	C	104	THR	10.6
2	C	98	ILE	10.4
2	C	105	ARG	10.1
2	C	346	GLN	10.0
2	C	102	ASN	9.7
2	K	328	TYR	9.3
2	K	388	LEU	9.2
2	C	97	VAL	9.2
2	C	412	MET	8.8
2	C	276	LYS	8.7
2	C	290	GLU	8.4
2	C	297	LYS	8.4

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Mol	Chain	Res	Type	RSRZ
2	C	82	TYR	8.3
2	C	127	LYS	8.1
2	C	100	GLU	8.1
2	C	275	VAL	8.1
2	C	129	HIS	8.0
2	K	396	GLN	7.7
2	K	321	ALA	7.4
2	K	412	MET	7.4
2	K	397	GLU	7.3
2	C	347	GLU	7.2
2	C	386	ALA	7.2
2	K	300	PHE	7.0
2	K	413	ARG	7.0
2	C	25	HIS	6.9
2	C	293	LEU	6.9
2	C	340	THR	6.6
2	C	133	LEU	6.5
2	C	348	GLU	6.5
2	C	349	LYS	6.4
2	C	101	GLY	6.3
2	C	292	SER	6.3
2	C	370	SER	6.2
2	C	414	TYR	6.1
2	K	389	GLU	6.1
2	C	26	GLY	6.0
2	C	84	LYS	5.9
2	C	415	GLN	5.9
2	C	424	GLU	5.8
2	K	275	VAL	5.8
2	K	414	TYR	5.8
2	K	276	LYS	5.7
2	K	390	PHE	5.7
2	C	400	GLN	5.7
2	C	399	MET	5.5
2	C	103	LYS	5.4
2	C	41	GLN	5.4
2	C	413	ARG	5.4
2	C	375	GLU	5.3
2	C	287	PHE	5.3
2	C	125	ASP	5.3
2	K	137	LEU	5.2
2	C	291	ILE	5.2

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Mol	Chain	Res	Type	RSRZ
2	C	136	LEU	5.1
2	C	128	PHE	5.1
2	K	370	SER	5.1
2	K	369	SER	5.1
2	C	420	MET	5.0
2	K	315	GLU	5.0
2	C	372	LEU	4.9
2	C	369	SER	4.9
2	C	416	LEU	4.9
2	K	416	LEU	4.9
2	C	351	SER	4.8
2	C	122	TYR	4.8
2	C	411	VAL	4.7
2	K	411	VAL	4.7
2	C	63	VAL	4.6
2	C	295	SER	4.6
2	C	396	GLN	4.5
2	C	78	ARG	4.4
2	C	86	PRO	4.3
2	C	62	ARG	4.3
2	C	405	PHE	4.3
2	C	377	LEU	4.3
2	C	339	ILE	4.3
2	C	32	ILE	4.2
2	K	317	LEU	4.2
2	C	378	ALA	4.2
2	C	350	ASP	4.2
2	K	340	THR	4.2
2	C	327	ARG	4.1
2	K	403	LYS	4.1
2	K	391	VAL	4.1
2	K	308	ALA	4.1
2	C	137	LEU	4.1
2	C	29	THR	4.1
2	C	394	ARG	4.1
2	K	405	PHE	4.0
2	C	325	MET	4.0
2	K	371	MET	3.9
1	A	468	PRO	3.9
2	K	346	GLN	3.9
2	C	112	GLU	3.9
2	K	135	GLU	3.8

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Mol	Chain	Res	Type	RSRZ
2	K	395	SER	3.8
2	C	296	GLY	3.8
1	A	631	MET	3.7
2	C	313	ALA	3.7
2	C	21	LEU	3.7
2	C	6	GLU	3.7
2	C	7	LYS	3.7
1	A	611	MET	3.6
2	C	31	MET	3.6
2	C	395	SER	3.6
2	C	115	LYS	3.6
2	C	106	LYS	3.6
2	C	85	VAL	3.6
1	E	583	ILE	3.6
2	K	342	MET	3.6
2	K	401	PHE	3.5
2	K	400	GLN	3.5
2	C	381	TYR	3.5
1	E	572	ILE	3.5
1	A	467	ALA	3.5
2	K	372	LEU	3.4
2	C	403	LYS	3.4
2	K	418	LEU	3.4
2	C	96	GLU	3.4
2	C	34	LEU	3.4
2	C	342	MET	3.3
2	K	286	ARG	3.3
1	E	642	ARG	3.3
2	K	311	GLU	3.3
2	C	24	CYS	3.3
2	K	375	GLU	3.3
2	C	382	LYS	3.3
1	A	470	ILE	3.2
2	C	298	TYR	3.2
2	C	113	PRO	3.2
2	B	411	VAL	3.2
2	C	373	LEU	3.2
2	C	309	LEU	3.2
2	C	87	ASP	3.2
2	K	381	TYR	3.2
2	C	417	ASP	3.2
2	C	77	GLU	3.1

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Mol	Chain	Res	Type	RSRZ
2	C	274	ASN	3.1
2	K	325	MET	3.1
2	K	383	ASP	3.1
2	K	322	ASP	3.1
1	E	640	MET	3.1
2	C	299	CYS	3.1
2	C	288	PHE	3.1
2	C	344	LYS	3.0
2	C	17	LEU	3.0
2	C	56	ALA	3.0
2	C	92	ILE	3.0
2	K	111	PHE	3.0
2	K	323	LEU	3.0
2	C	315	GLU	3.0
1	A	583	ILE	3.0
2	K	299	CYS	3.0
2	K	347	GLU	3.0
2	C	135	GLU	2.9
2	C	289	ASP	2.9
2	K	386	ALA	2.9
2	C	379	GLU	2.9
1	A	560	PHE	2.9
2	K	283	LEU	2.9
2	K	377	LEU	2.9
2	C	311	GLU	2.9
2	C	126	ASN	2.9
2	K	304	ASP	2.8
2	C	423	PRO	2.8
1	E	573	LEU	2.8
2	C	323	LEU	2.8
2	C	94	CYS	2.8
2	C	132	ALA	2.8
2	K	339	ILE	2.7
2	C	79	LEU	2.7
2	C	60	LYS	2.7
2	C	52	GLU	2.7
2	C	110	ASP	2.7
2	C	64	ASN	2.7
2	K	309	LEU	2.7
1	A	640	MET	2.6
2	C	81	LEU	2.6
2	C	93	TYR	2.6

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Mol	Chain	Res	Type	RSRZ
2	C	80	LYS	2.6
2	C	388	LEU	2.6
2	C	116	PRO	2.6
2	K	424	GLU	2.6
1	A	495	LYS	2.6
2	C	45	TYR	2.6
1	A	623	LEU	2.6
2	C	57	SER	2.5
2	K	387	ASN	2.5
1	A	581	MET	2.5
2	C	50	ALA	2.5
2	C	118	ASN	2.5
1	E	611	MET	2.5
2	H	45	TYR	2.5
2	K	382	LYS	2.5
2	C	58	ASN	2.5
2	K	133	LEU	2.5
2	K	284	ILE	2.5
2	C	23	ASN	2.5
2	C	418	LEU	2.5
1	A	533	ARG	2.4
2	C	124	CYS	2.4
2	C	9	ILE	2.4
2	C	419	SER	2.4
1	E	470	ILE	2.4
2	H	319	CYS	2.4
2	C	285	GLN	2.4
2	H	306	MET	2.4
2	K	320	PHE	2.4
1	E	612	PHE	2.4
2	C	314	VAL	2.3
2	C	283	LEU	2.3
2	B	32	ILE	2.3
2	K	314	VAL	2.3
1	E	639	TYR	2.3
2	C	55	THR	2.3
1	J	471	MET	2.3
2	K	344	LYS	2.3
1	E	581	MET	2.3
1	D	483	ILE	2.3
2	K	420	MET	2.3
1	E	632	GLU	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	484	LEU	2.3
1	A	473	ILE	2.3
2	H	328	TYR	2.2
2	B	49	LEU	2.2
2	K	316	THR	2.2
2	H	92	ILE	2.2
1	A	509	GLU	2.2
2	C	90	LEU	2.2
2	K	285	GLN	2.2
2	K	394	ARG	2.2
2	K	374	SER	2.2
2	C	131	GLU	2.2
1	D	583	ILE	2.2
2	C	35	ILE	2.2
2	H	373	LEU	2.2
2	K	417	ASP	2.2
1	E	471	MET	2.2
2	K	348	GLU	2.2
1	D	562	ALA	2.2
1	A	471	MET	2.2
2	C	75	THR	2.2
2	K	408	ILE	2.2
2	C	300	PHE	2.2
2	C	4	THR	2.2
1	A	547	LEU	2.2
2	C	44	ARG	2.1
2	C	53	TYR	2.1
2	K	277	TYR	2.1
1	D	532	LEU	2.1
2	K	306	MET	2.1
2	K	319	CYS	2.1
1	E	564	ILE	2.1
2	K	404	GLY	2.1
2	H	323	LEU	2.1
1	E	601	LYS	2.1
2	H	325	MET	2.1
1	A	513	ILE	2.1
2	K	415	GLN	2.1
2	C	83	ASN	2.1
1	E	662	ASP	2.1
1	A	637	TYR	2.1
1	A	476	LYS	2.1

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Mol	Chain	Res	Type	RSRZ
2	C	425	SER	2.1
2	C	278	VAL	2.1
2	H	276	LYS	2.0
2	K	136	LEU	2.0
2	C	51	GLU	2.0
1	D	581	MET	2.0
2	B	34	LEU	2.0
2	K	318	LEU	2.0
2	C	11	ILE	2.0
1	E	480	LEU	2.0
1	E	484	LEU	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.