



Full wwPDB X-ray Structure Validation Report

Apr 9, 2014 – 08:46 PM EDT

PDB ID : 2ED6
Title : Crystal Structure of Envelope Protein VP28 from White Spot Syndrome Virus (WSSV)
Authors : Hew, C.L.; Sivaraman, J.; Tang, X.H.
Deposited on : 2007-02-14
Resolution : 2.00 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>

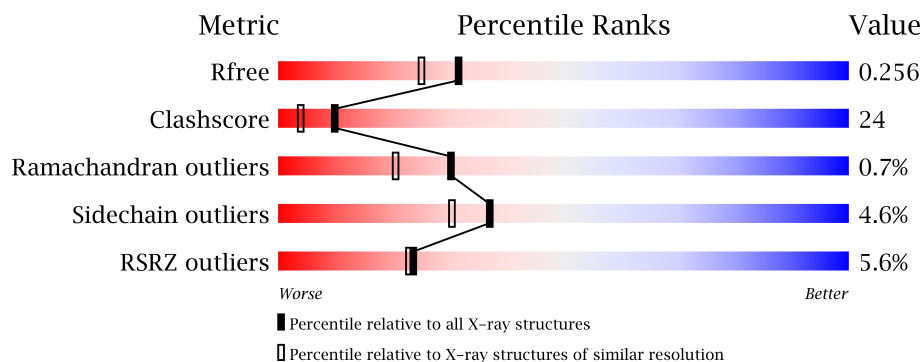
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1439
EDS : stable22978
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) : stable22978

1 Overall quality at a glance

The reported resolution of this entry is 2.00 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	66092	4888 (2.00-2.00)
Clashscore	79885	6188 (2.00-2.00)
Ramachandran outliers	78287	6102 (2.00-2.00)
Sidechain outliers	78261	6100 (2.00-2.00)
RSRZ outliers	66119	4890 (2.00-2.00)

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	170	
1	B	170	
1	C	170	
1	D	170	
1	E	170	
1	F	170	
1	G	170	
1	H	170	
1	I	170	
1	J	170	
1	K	170	
1	L	170	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 16843 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 25kDa structural protein VP25.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	170	Total	C	N	O	S	0	0	0
			1288	803	211	267	7			
1	B	170	Total	C	N	O	S	0	0	0
			1288	803	211	267	7			
1	C	170	Total	C	N	O	S	0	0	0
			1288	803	211	267	7			
1	D	170	Total	C	N	O	S	0	0	0
			1288	803	211	267	7			
1	E	170	Total	C	N	O	S	0	0	0
			1288	803	211	267	7			
1	F	170	Total	C	N	O	S	0	0	0
			1288	803	211	267	7			
1	G	170	Total	C	N	O	S	0	0	0
			1288	803	211	267	7			
1	H	170	Total	C	N	O	S	0	0	0
			1288	803	211	267	7			
1	I	170	Total	C	N	O	S	0	0	0
			1288	803	211	267	7			
1	J	170	Total	C	N	O	S	0	0	0
			1288	803	211	267	7			
1	K	170	Total	C	N	O	S	0	0	0
			1288	803	211	267	7			
1	L	170	Total	C	N	O	S	0	0	0
			1288	803	211	267	7			

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	107	Total	O	0	0
			107	107		
2	B	133	Total	O	0	0
			133	133		

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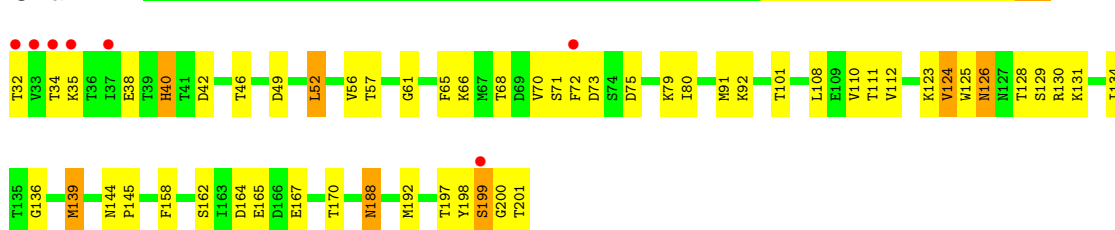
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	C	124	Total 124	O 124	0	0
2	D	112	Total 112	O 112	0	0
2	E	138	Total 138	O 138	0	0
2	F	154	Total 154	O 154	0	0
2	G	136	Total 136	O 136	0	0
2	H	122	Total 122	O 122	0	0
2	I	125	Total 125	O 125	0	0
2	J	102	Total 102	O 102	0	0
2	K	53	Total 53	O 53	0	0
2	L	81	Total 81	O 81	0	0

3 Residue-property plots

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

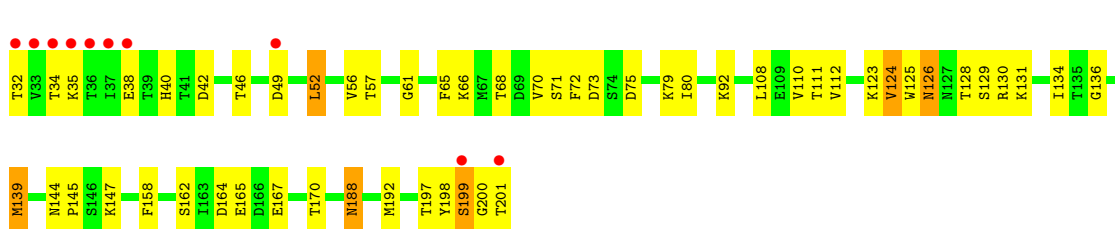
• Molecule 1: 25kDa structural protein VP25

Chain A:



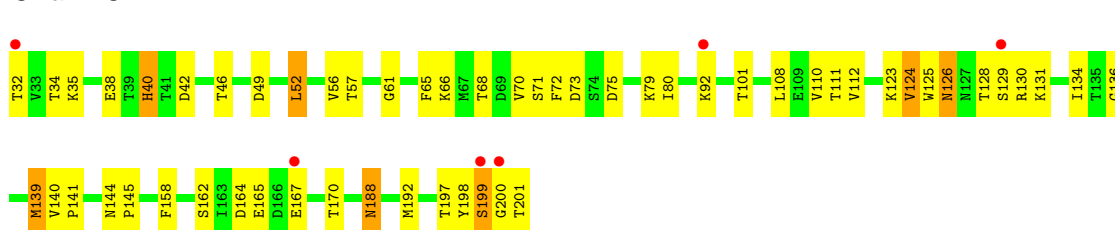
• Molecule 1: 25kDa structural protein VP25

Chain B:



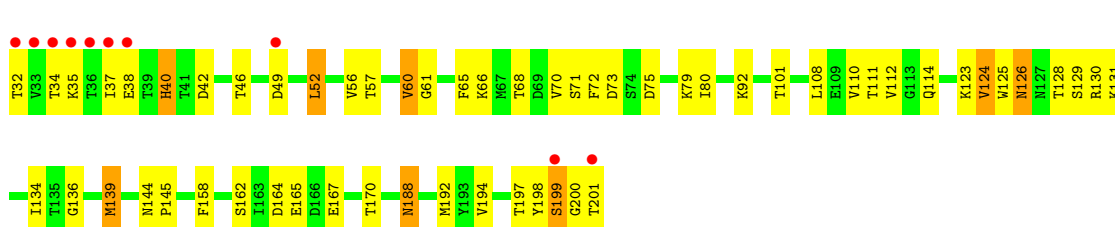
• Molecule 1: 25kDa structural protein VP25

Chain C:

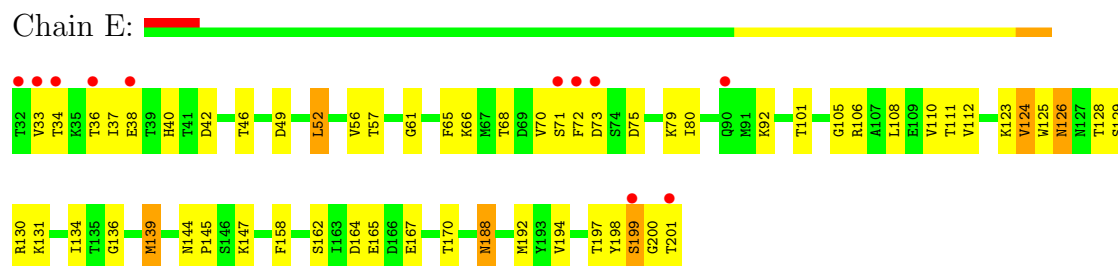


• Molecule 1: 25kDa structural protein VP25

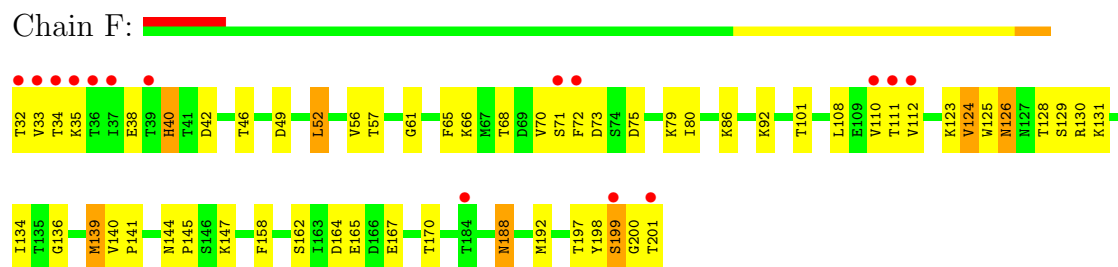
Chain D:



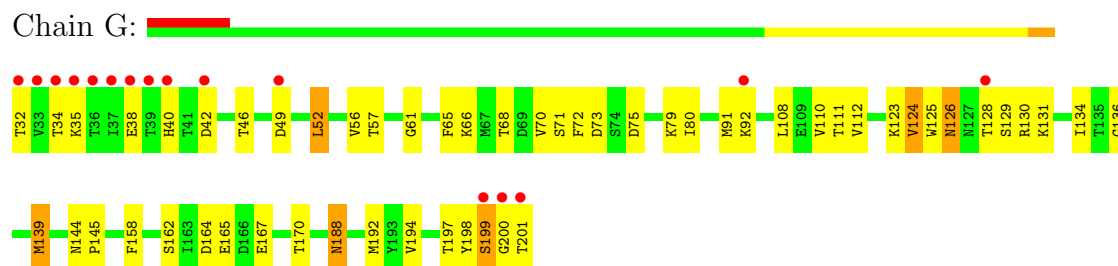
- Molecule 1: 25kDa structural protein VP25



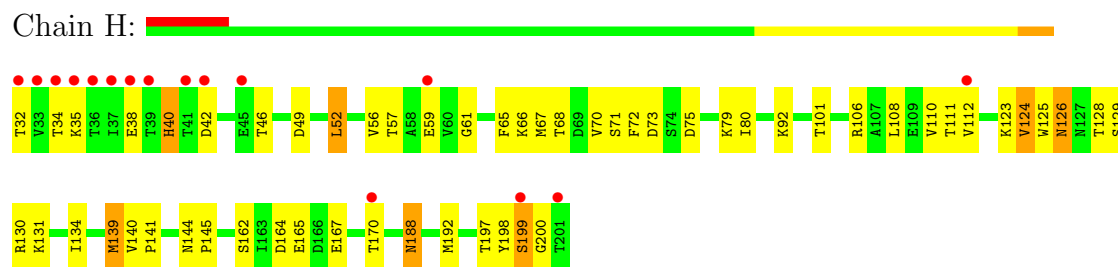
- Molecule 1: 25kDa structural protein VP25



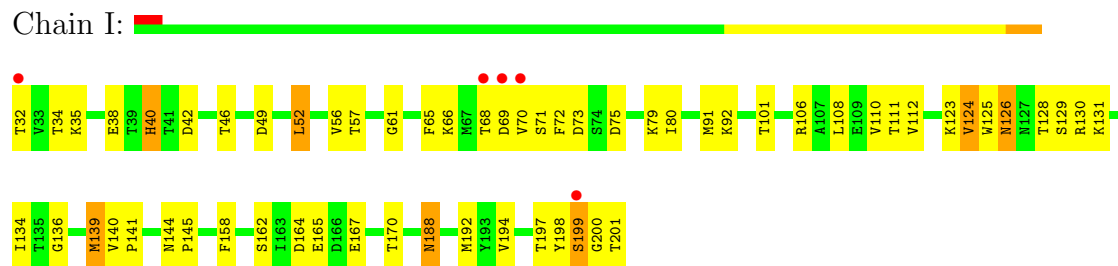
- Molecule 1: 25kDa structural protein VP25



- Molecule 1: 25kDa structural protein VP25

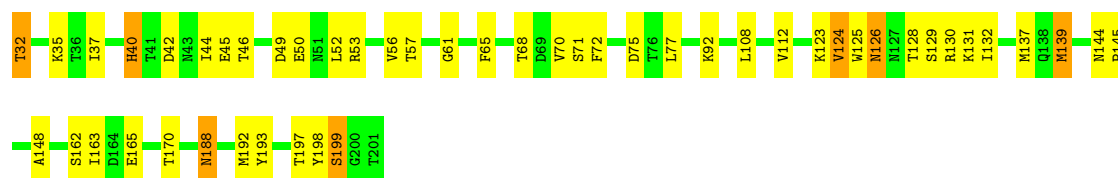


- Molecule 1: 25kDa structural protein VP25



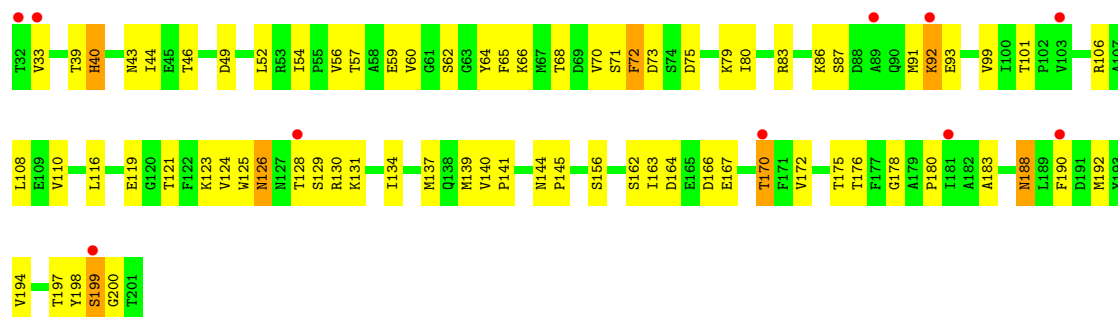
- Molecule 1: 25kDa structural protein VP25

Chain J: 



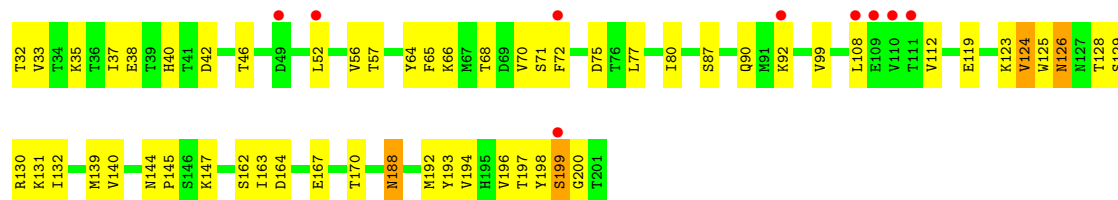
- Molecule 1: 25kDa structural protein VP25

Chain K: 



- Molecule 1: 25kDa structural protein VP25

Chain L: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	105.33Å 106.71Å 200.37Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.00 – 2.00 49.87 – 2.00	Depositor EDS
% Data completeness (in resolution range)	(Not available) (45.00-2.00) 99.6 (49.87-2.00)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	8.63 (at 2.00Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.248 , 0.281 0.252 , 0.256	Depositor DCC
R_{free} test set	7648 reflections (5.02%)	DCC
Wilson B-factor (Å ²)	15.3	Xtriage
Anisotropy	0.271	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 37.0	EDS
Estimated twinning fraction	0.020 for k,h,-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Outliers	0 of 152458 reflections	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	16843	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.28% 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	0.33	0/1309	0.61	0/1775
1	B	0.35	0/1309	0.61	0/1775
1	C	0.34	0/1309	0.61	0/1775
1	D	0.41	1/1309 (0.1%)	0.62	0/1775
1	E	0.35	0/1309	0.61	0/1775
1	F	0.37	0/1309	0.62	0/1775
1	G	0.37	0/1309	0.62	0/1775
1	H	0.37	0/1309	0.61	0/1775
1	I	0.34	0/1309	0.61	0/1775
1	J	0.37	0/1309	0.64	0/1775
1	K	0.32	0/1309	0.57	0/1775
1	L	0.33	0/1309	0.58	0/1775
All	All	0.36	1/15708 (0.0%)	0.61	0/21300

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	60	VAL	CB-CG1	-7.91	1.36	1.52

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1288	0	1256	59	0
1	B	1288	0	1256	63	0
1	C	1288	0	1256	62	0
1	D	1288	0	1256	64	1
1	E	1288	0	1256	71	0
1	F	1288	0	1256	62	0
1	G	1288	0	1256	62	0
1	H	1288	0	1256	63	1
1	I	1288	0	1256	73	0
1	J	1288	0	1256	65	0
1	K	1288	0	1256	71	0
1	L	1288	0	1256	54	0
2	A	107	0	0	5	0
2	B	133	0	0	5	0
2	C	124	0	0	8	0
2	D	112	0	0	8	0
2	E	138	0	0	5	0
2	F	154	0	0	5	0
2	G	136	0	0	3	0
2	H	122	0	0	7	0
2	I	125	0	0	6	0
2	J	102	0	0	5	0
2	K	53	0	0	2	0
2	L	81	0	0	2	0
All	All	16843	0	15072	732	1

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 24.

All (732) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:F:56:VAL:HG11	1:F:65:PHE:HB3	1.28	1.14
1:B:56:VAL:HG11	1:B:65:PHE:HB3	1.29	1.13
1:D:56:VAL:HG11	1:D:65:PHE:HB3	1.30	1.13
1:A:56:VAL:HG11	1:A:65:PHE:HB3	1.29	1.12
1:G:56:VAL:HG11	1:G:65:PHE:HB3	1.27	1.10
1:E:56:VAL:HG11	1:E:65:PHE:HB3	1.29	1.09
1:C:56:VAL:HG11	1:C:65:PHE:HB3	1.30	1.09
1:H:56:VAL:HG11	1:H:65:PHE:HB3	1.29	1.08
1:I:56:VAL:HG11	1:I:65:PHE:HB3	1.31	1.07
1:H:46:THR:HG22	1:H:130:ARG:HH22	1.19	1.06
1:G:46:THR:HG22	1:G:130:ARG:HH22	1.23	1.04
1:D:197:THR:HG22	1:D:199:SER:H	1.22	1.04

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:L:46:THR:HG22	1:L:130:ARG:HH22	1.23	1.03
1:A:197:THR:HG22	1:A:199:SER:H	1.23	1.03
1:L:56:VAL:HG11	1:L:65:PHE:HB3	1.38	1.03
1:I:46:THR:HG22	1:I:130:ARG:HH22	1.21	1.03
1:F:197:THR:HG22	1:F:199:SER:H	1.22	1.03
1:C:46:THR:HG22	1:C:130:ARG:HH22	1.23	1.03
1:B:46:THR:HG22	1:B:130:ARG:HH22	1.22	1.02
1:B:197:THR:HG22	1:B:199:SER:H	1.22	1.02
1:E:46:THR:HG22	1:E:130:ARG:HH22	1.24	1.02
1:K:197:THR:HG22	1:K:199:SER:H	1.17	1.02
1:C:197:THR:HG22	1:C:199:SER:H	1.24	1.01
1:A:46:THR:HG22	1:A:130:ARG:HH22	1.22	1.01
1:D:46:THR:HG22	1:D:130:ARG:HH22	1.21	1.01
1:H:197:THR:HG22	1:H:199:SER:H	1.23	1.00
1:E:197:THR:HG22	1:E:199:SER:H	1.24	1.00
1:F:46:THR:HG22	1:F:130:ARG:HH22	1.21	0.99
1:I:197:THR:HG22	1:I:199:SER:H	1.23	0.99
1:G:197:THR:HG22	1:G:199:SER:H	1.22	0.99
1:K:139:MET:HE1	1:K:190:PHE:HB2	1.44	0.98
1:J:197:THR:HG22	1:J:199:SER:H	1.28	0.97
1:D:126:ASN:HD21	1:D:128:THR:HB	1.34	0.91
1:A:126:ASN:HD21	1:A:128:THR:HB	1.36	0.91
1:I:126:ASN:HD21	1:I:128:THR:HB	1.36	0.91
1:E:126:ASN:HD21	1:E:128:THR:HB	1.36	0.90
1:H:126:ASN:HD21	1:H:128:THR:HB	1.36	0.90
1:C:126:ASN:HD21	1:C:128:THR:HB	1.36	0.90
1:J:126:ASN:HD21	1:J:128:THR:HB	1.36	0.89
1:J:46:THR:HG22	1:J:130:ARG:HH22	1.38	0.89
1:B:126:ASN:HD21	1:B:128:THR:HB	1.37	0.89
1:F:126:ASN:HD21	1:F:128:THR:HB	1.35	0.89
1:J:56:VAL:HG11	1:J:65:PHE:HB3	1.54	0.89
1:E:49:ASP:HB2	2:F:1516:HOH:O	1.73	0.89
1:G:126:ASN:HD21	1:G:128:THR:HB	1.35	0.88
1:K:39:THR:HG22	1:K:43:ASN:HD21	1.38	0.88
1:K:56:VAL:HG11	1:K:65:PHE:HB3	1.55	0.88
1:H:46:THR:HG22	1:H:130:ARG:NH2	1.90	0.86
1:L:126:ASN:HD21	1:L:128:THR:HB	1.41	0.86
1:F:46:THR:HG22	1:F:130:ARG:NH2	1.90	0.86
1:I:46:THR:HG22	1:I:130:ARG:NH2	1.91	0.86
1:L:197:THR:HG22	1:L:199:SER:H	1.41	0.85
1:B:46:THR:HG22	1:B:130:ARG:NH2	1.91	0.85
1:G:46:THR:HG22	1:G:130:ARG:NH2	1.91	0.85

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:K:46:THR:HG22	1:K:130:ARG:HH22	1.39	0.85
1:A:46:THR:HG22	1:A:130:ARG:NH2	1.91	0.84
1:C:46:THR:HG22	1:C:130:ARG:NH2	1.92	0.84
1:D:46:THR:HG22	1:D:130:ARG:NH2	1.90	0.84
1:E:46:THR:HG22	1:E:130:ARG:NH2	1.92	0.83
1:L:46:THR:HG22	1:L:130:ARG:NH2	1.93	0.82
1:B:49:ASP:HB2	2:C:1518:HOH:O	1.79	0.82
1:J:188:ASN:H	1:J:188:ASN:HD22	1.27	0.81
1:D:49:ASP:HB2	2:D:1519:HOH:O	1.81	0.81
1:L:188:ASN:H	1:L:188:ASN:HD22	1.29	0.81
1:J:46:THR:HG22	1:J:130:ARG:NH2	1.96	0.80
1:K:141:PRO:HA	1:K:190:PHE:HD1	1.46	0.80
1:G:49:ASP:HB2	2:I:1515:HOH:O	1.80	0.79
1:F:131:LYS:HE2	1:F:162:SER:HB2	1.63	0.79
1:D:131:LYS:HE2	1:D:162:SER:HB2	1.64	0.79
1:K:124:VAL:HG11	1:K:134:ILE:HD11	1.65	0.78
1:H:131:LYS:HE2	1:H:162:SER:HB2	1.63	0.78
1:I:131:LYS:HE2	1:I:162:SER:HB2	1.64	0.78
1:E:147:LYS:HE3	2:F:2214:HOH:O	1.82	0.78
1:A:131:LYS:HE2	1:A:162:SER:HB2	1.65	0.77
1:E:131:LYS:HE2	1:E:162:SER:HB2	1.65	0.77
1:G:131:LYS:HE2	1:G:162:SER:HB2	1.66	0.76
1:B:131:LYS:HE2	1:B:162:SER:HB2	1.66	0.76
1:F:188:ASN:HD22	1:F:188:ASN:H	1.30	0.76
1:F:139:MET:HE3	1:F:192:MET:HA	1.68	0.75
1:I:188:ASN:H	1:I:188:ASN:HD22	1.34	0.75
1:C:131:LYS:HE2	1:C:162:SER:HB2	1.67	0.75
2:H:2339:HOH:O	1:I:49:ASP:HB2	1.87	0.75
1:E:139:MET:HE3	1:E:192:MET:HA	1.69	0.75
1:H:139:MET:HE3	1:H:192:MET:HA	1.67	0.75
1:B:188:ASN:HD22	1:B:188:ASN:H	1.35	0.74
1:A:188:ASN:HD22	1:A:188:ASN:H	1.35	0.74
1:L:128:THR:HG21	1:L:198:TYR:OH	1.88	0.74
1:J:68:THR:O	1:J:70:VAL:HG23	1.88	0.74
1:H:188:ASN:H	1:H:188:ASN:HD22	1.32	0.74
1:K:126:ASN:HD21	1:K:128:THR:HB	1.52	0.73
1:C:139:MET:HE3	1:C:192:MET:HA	1.68	0.73
1:A:139:MET:HE3	1:A:192:MET:HA	1.68	0.73
1:C:188:ASN:H	1:C:188:ASN:HD22	1.35	0.73
1:G:56:VAL:CG1	1:G:65:PHE:HB3	2.15	0.73
1:I:139:MET:HE3	1:I:192:MET:HA	1.68	0.73
1:G:188:ASN:HD22	1:G:188:ASN:H	1.35	0.73

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:L:68:THR:O	1:L:70:VAL:HG23	1.89	0.73
2:G:1517:HOH:O	1:H:49:ASP:HB2	1.87	0.73
1:J:144:ASN:HD21	1:J:148:ALA:HB3	1.54	0.72
1:E:188:ASN:H	1:E:188:ASN:HD22	1.37	0.72
1:D:188:ASN:H	1:D:188:ASN:HD22	1.36	0.72
1:K:164:ASP:HB2	1:K:167:GLU:OE1	1.90	0.72
1:B:139:MET:HE3	1:B:192:MET:HA	1.70	0.71
1:B:139:MET:HA	1:B:139:MET:HE3	1.72	0.71
1:D:139:MET:HE3	1:D:139:MET:HA	1.71	0.71
1:D:139:MET:HE3	1:D:192:MET:HA	1.73	0.71
1:J:52:LEU:HD11	1:J:72:PHE:CE1	2.25	0.71
1:G:139:MET:HE3	1:G:192:MET:HA	1.71	0.70
1:E:106:ARG:HB3	1:I:68:THR:CG2	2.21	0.70
1:K:139:MET:CE	1:K:190:PHE:HB2	2.20	0.70
1:L:139:MET:HE2	1:L:192:MET:HB2	1.74	0.70
1:B:139:MET:CE	1:B:139:MET:HA	2.22	0.70
1:H:199:SER:O	1:I:75:ASP:OD2	2.09	0.70
1:F:56:VAL:CG1	1:F:65:PHE:HB3	2.16	0.69
1:J:75:ASP:OD2	1:K:199:SER:O	2.10	0.69
1:L:52:LEU:HD11	1:L:77:LEU:HD12	1.73	0.69
1:G:199:SER:O	1:H:75:ASP:OD2	2.11	0.69
1:D:139:MET:CE	1:D:139:MET:HA	2.22	0.69
1:I:139:MET:CE	1:I:139:MET:HA	2.23	0.69
1:J:139:MET:HE3	1:J:192:MET:HA	1.74	0.69
1:F:139:MET:HA	1:F:139:MET:CE	2.23	0.69
1:K:197:THR:HG22	1:K:199:SER:N	2.00	0.68
1:J:52:LEU:HD11	1:J:72:PHE:HE1	1.59	0.68
1:A:139:MET:CE	1:A:139:MET:HA	2.24	0.68
1:C:139:MET:HA	1:C:139:MET:CE	2.23	0.68
1:K:197:THR:CG2	1:K:199:SER:H	1.99	0.68
1:J:199:SER:O	1:L:75:ASP:OD2	2.12	0.68
1:G:128:THR:HG21	1:G:198:TYR:OH	1.94	0.68
1:H:139:MET:HA	1:H:139:MET:CE	2.24	0.68
1:L:131:LYS:HE2	1:L:162:SER:HB2	1.76	0.67
1:B:92:LYS:HG3	1:B:125:TRP:CZ2	2.29	0.67
1:G:92:LYS:HG3	1:G:125:TRP:CZ2	2.29	0.67
1:J:131:LYS:HE2	1:J:162:SER:HB2	1.76	0.67
1:L:92:LYS:HG3	1:L:125:TRP:CZ2	2.27	0.67
1:A:56:VAL:CG1	1:A:65:PHE:HB3	2.17	0.67
1:F:197:THR:HG21	2:F:1083:HOH:O	1.93	0.67
1:G:197:THR:HG21	2:G:1088:HOH:O	1.94	0.67
1:G:139:MET:HA	1:G:139:MET:CE	2.24	0.67

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:K:188:ASN:HD22	1:K:188:ASN:H	1.41	0.67
1:D:49:ASP:OD2	1:E:49:ASP:OD2	2.13	0.67
1:E:106:ARG:HG2	1:I:69:ASP:O	1.94	0.67
1:G:68:THR:O	1:G:70:VAL:HG23	1.95	0.67
1:E:139:MET:CE	1:E:139:MET:HA	2.24	0.67
1:I:92:LYS:HG3	1:I:125:TRP:CZ2	2.30	0.67
1:I:139:MET:HE3	1:I:139:MET:HA	1.77	0.67
1:E:56:VAL:CG1	1:E:65:PHE:HB3	2.17	0.66
1:F:92:LYS:HG3	1:F:125:TRP:CZ2	2.30	0.66
1:D:56:VAL:CG1	1:D:65:PHE:HB3	2.18	0.66
1:H:92:LYS:HG3	1:H:125:TRP:CZ2	2.30	0.66
1:K:39:THR:HG22	1:K:43:ASN:ND2	2.09	0.66
1:G:139:MET:HA	1:G:139:MET:HE3	1.77	0.66
1:K:46:THR:HG22	1:K:130:ARG:NH2	2.09	0.66
1:C:92:LYS:HG3	1:C:125:TRP:CZ2	2.31	0.66
1:A:199:SER:O	1:C:75:ASP:OD2	2.14	0.65
1:A:75:ASP:OD2	1:B:199:SER:O	2.13	0.65
1:L:42:ASP:O	1:L:46:THR:HG23	1.95	0.65
1:L:87:SER:OG	1:L:90:GLN:HG3	1.96	0.65
1:C:128:THR:HG21	1:C:198:TYR:OH	1.97	0.65
1:D:49:ASP:OD2	1:F:49:ASP:OD2	2.15	0.65
1:J:124:VAL:HG22	1:J:163:ILE:CD1	2.25	0.65
1:A:92:LYS:HG3	1:A:125:TRP:CZ2	2.31	0.65
1:C:68:THR:O	1:C:70:VAL:HG23	1.96	0.64
1:D:92:LYS:HG3	1:D:125:TRP:CZ2	2.32	0.64
1:E:92:LYS:HG3	1:E:125:TRP:CZ2	2.31	0.64
1:D:75:ASP:OD2	1:E:199:SER:O	2.15	0.64
1:L:52:LEU:HD11	1:L:77:LEU:CD1	2.27	0.64
1:A:139:MET:HA	1:A:139:MET:HE3	1.80	0.64
1:A:128:THR:HG21	1:A:198:TYR:OH	1.98	0.64
1:I:128:THR:HG21	1:I:198:TYR:OH	1.98	0.64
1:J:139:MET:CE	1:J:139:MET:HA	2.28	0.64
1:F:128:THR:HG21	1:F:198:TYR:OH	1.97	0.64
1:I:68:THR:O	1:I:70:VAL:HG23	1.98	0.64
1:K:131:LYS:HE2	1:K:162:SER:HB2	1.80	0.63
1:F:170:THR:HG23	2:F:1304:HOH:O	1.98	0.63
1:F:126:ASN:ND2	1:F:128:THR:HB	2.12	0.63
1:D:128:THR:HG21	1:D:198:TYR:OH	1.99	0.63
1:J:126:ASN:ND2	1:J:128:THR:HB	2.10	0.63
1:B:56:VAL:CG1	1:B:65:PHE:HB3	2.18	0.63
1:C:197:THR:HG21	2:C:1084:HOH:O	1.98	0.63
1:E:128:THR:HG21	1:E:198:TYR:OH	1.99	0.63

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:128:THR:HG21	1:B:198:TYR:OH	1.99	0.62
1:D:197:THR:HG21	2:D:1918:HOH:O	1.97	0.62
1:K:75:ASP:OD2	1:L:199:SER:O	2.15	0.62
1:G:170:THR:HG23	2:G:2286:HOH:O	1.99	0.62
1:H:67:MET:HE3	2:H:2129:HOH:O	1.99	0.62
1:K:72:PHE:CE1	1:K:80:ILE:HB	2.34	0.62
1:H:56:VAL:CG1	1:H:65:PHE:HB3	2.17	0.62
1:L:124:VAL:HG22	1:L:163:ILE:CD1	2.29	0.62
1:B:68:THR:O	1:B:70:VAL:HG23	1.98	0.62
1:C:32:THR:HG23	2:C:2063:HOH:O	1.99	0.62
1:D:126:ASN:ND2	1:D:128:THR:HB	2.12	0.62
1:E:139:MET:HE3	1:E:139:MET:HA	1.80	0.62
1:A:40:HIS:HD2	2:A:1254:HOH:O	1.84	0.61
1:E:49:ASP:OD2	1:F:49:ASP:OD2	2.19	0.61
1:G:126:ASN:ND2	1:G:128:THR:HB	2.12	0.61
1:F:139:MET:HA	1:F:139:MET:HE3	1.81	0.61
1:C:126:ASN:ND2	1:C:128:THR:HB	2.13	0.61
1:C:56:VAL:CG1	1:C:65:PHE:HB3	2.18	0.61
1:B:70:VAL:HG21	2:B:1802:HOH:O	1.99	0.61
1:D:68:THR:O	1:D:70:VAL:HG23	2.01	0.61
1:J:37:ILE:CD1	1:K:33:VAL:HG13	2.31	0.61
1:A:126:ASN:ND2	1:A:128:THR:HB	2.13	0.61
1:C:139:MET:HA	1:C:139:MET:HE3	1.81	0.60
1:A:49:ASP:OD2	1:B:49:ASP:OD2	2.20	0.60
1:A:68:THR:O	1:A:70:VAL:HG23	2.01	0.60
1:H:128:THR:HG21	1:H:198:TYR:OH	2.00	0.60
1:J:128:THR:HG21	1:J:198:TYR:OH	2.02	0.60
1:K:123:LYS:HA	1:K:170:THR:HA	1.83	0.60
1:J:188:ASN:ND2	1:J:188:ASN:H	1.99	0.59
1:C:56:VAL:HG12	1:C:57:THR:N	2.17	0.59
1:I:56:VAL:CG1	1:I:65:PHE:HB3	2.19	0.59
1:L:144:ASN:HB2	1:L:145:PRO:CD	2.32	0.59
1:B:126:ASN:ND2	1:B:128:THR:HB	2.14	0.59
1:G:56:VAL:HG12	1:G:57:THR:N	2.18	0.59
1:H:40:HIS:HD2	2:H:1104:HOH:O	1.86	0.59
1:H:68:THR:O	1:H:70:VAL:HG23	2.02	0.59
1:J:70:VAL:HG12	1:J:71:SER:N	2.17	0.59
1:E:56:VAL:HG11	1:E:65:PHE:CB	2.20	0.59
1:H:139:MET:HA	1:H:139:MET:HE3	1.83	0.59
1:J:56:VAL:CG1	1:J:65:PHE:HB3	2.31	0.59
1:A:56:VAL:HG12	1:A:57:THR:N	2.18	0.59
1:E:68:THR:O	1:E:70:VAL:HG23	2.03	0.58

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:I:56:VAL:HG12	1:I:57:THR:N	2.18	0.58
1:J:45:GLU:HG3	1:J:130:ARG:HH11	1.68	0.58
1:K:137:MET:HG3	1:K:192:MET:HE2	1.85	0.58
1:I:126:ASN:ND2	1:I:128:THR:HB	2.14	0.58
1:D:37:ILE:CD1	1:E:33:VAL:HG13	2.34	0.58
1:F:56:VAL:HG12	1:F:57:THR:N	2.19	0.58
1:H:56:VAL:HG12	1:H:57:THR:N	2.18	0.58
1:J:45:GLU:HG3	1:J:130:ARG:NH1	2.19	0.58
1:B:56:VAL:HG12	1:B:57:THR:N	2.19	0.58
1:K:49:ASP:OD1	1:K:197:THR:HG23	2.03	0.58
1:E:75:ASP:OD2	1:F:199:SER:O	2.22	0.58
1:K:70:VAL:HG12	1:K:71:SER:N	2.17	0.58
1:E:126:ASN:ND2	1:E:128:THR:HB	2.13	0.57
1:G:188:ASN:HD22	1:G:188:ASN:N	2.01	0.57
1:E:92:LYS:HB2	2:E:1525:HOH:O	2.04	0.57
1:A:49:ASP:OD2	1:C:49:ASP:OD2	2.22	0.57
1:B:123:LYS:HB3	1:B:170:THR:HG22	1.86	0.57
1:H:126:ASN:ND2	1:H:128:THR:HB	2.14	0.57
1:K:124:VAL:HG13	1:K:163:ILE:CD1	2.34	0.57
1:L:188:ASN:H	1:L:188:ASN:ND2	2.00	0.57
1:A:197:THR:HG21	2:C:2012:HOH:O	2.03	0.57
1:E:46:THR:CG2	1:E:130:ARG:HH12	2.18	0.57
1:C:52:LEU:HD21	1:C:80:ILE:HD13	1.87	0.57
1:J:37:ILE:HD11	1:K:33:VAL:HG13	1.87	0.57
1:E:56:VAL:HG12	1:E:57:THR:N	2.20	0.57
1:I:123:LYS:HB3	1:I:170:THR:HG22	1.87	0.56
1:J:126:ASN:ND2	1:J:128:THR:H	2.02	0.56
1:H:56:VAL:HG11	1:H:65:PHE:CB	2.20	0.56
1:K:73:ASP:OD1	1:K:79:LYS:HG3	2.04	0.56
1:B:52:LEU:HD21	1:B:80:ILE:HD13	1.87	0.56
1:C:123:LYS:HB3	1:C:170:THR:HG22	1.87	0.56
1:F:56:VAL:HG11	1:F:65:PHE:CB	2.20	0.56
1:B:56:VAL:HG11	1:B:65:PHE:CB	2.19	0.56
1:D:199:SER:O	1:F:75:ASP:OD2	2.24	0.56
1:I:32:THR:HG23	2:I:1412:HOH:O	2.05	0.56
1:A:123:LYS:HB3	1:A:170:THR:HG22	1.87	0.56
1:B:170:THR:HG23	2:B:2148:HOH:O	2.05	0.56
1:E:170:THR:HG23	2:E:2491:HOH:O	2.04	0.56
1:E:52:LEU:HD21	1:E:80:ILE:HD13	1.87	0.56
1:E:70:VAL:HG12	1:E:71:SER:N	2.21	0.56
1:H:123:LYS:HB3	1:H:170:THR:HG22	1.88	0.56
1:G:123:LYS:HB3	1:G:170:THR:HG22	1.87	0.56

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:I:68:THR:HG23	2:I:1306:HOH:O	2.06	0.56
1:L:52:LEU:HB2	1:L:194:VAL:CG2	2.36	0.56
1:C:46:THR:HG21	2:C:1470:HOH:O	2.05	0.56
1:D:56:VAL:HG12	1:D:57:THR:N	2.20	0.56
1:E:46:THR:HG21	2:E:1404:HOH:O	2.05	0.56
1:F:123:LYS:HB3	1:F:170:THR:HG22	1.87	0.56
1:H:110:VAL:HG11	2:H:1380:HOH:O	2.05	0.55
1:J:129:SER:HA	1:J:165:GLU:OE2	2.06	0.55
1:A:56:VAL:HG11	1:A:65:PHE:CB	2.20	0.55
1:D:123:LYS:HB3	1:D:170:THR:HG22	1.87	0.55
1:D:52:LEU:HD21	1:D:80:ILE:HD13	1.87	0.55
1:G:124:VAL:HG11	1:G:134:ILE:HD11	1.87	0.55
1:E:123:LYS:HB3	1:E:170:THR:HG22	1.87	0.55
1:B:46:THR:CG2	1:B:130:ARG:HH12	2.19	0.55
1:A:52:LEU:HD21	1:A:80:ILE:HD13	1.89	0.55
1:H:52:LEU:HD21	1:H:80:ILE:HD13	1.88	0.55
2:D:1588:HOH:O	1:F:147:LYS:HE3	2.07	0.55
1:F:70:VAL:HG12	1:F:71:SER:N	2.22	0.55
1:B:75:ASP:OD2	1:C:199:SER:O	2.24	0.55
1:F:52:LEU:HD21	1:F:80:ILE:HD13	1.88	0.55
1:I:124:VAL:HG11	1:I:134:ILE:HD11	1.88	0.55
1:I:188:ASN:HD22	1:I:188:ASN:N	2.00	0.55
1:J:46:THR:CG2	1:J:130:ARG:HH12	2.20	0.55
1:D:46:THR:CG2	1:D:130:ARG:HH12	2.20	0.54
1:D:73:ASP:OD1	1:D:79:LYS:HG2	2.08	0.54
1:K:197:THR:HG21	2:K:1095:HOH:O	2.06	0.54
1:B:70:VAL:HG12	1:B:71:SER:N	2.23	0.54
1:D:124:VAL:HG11	1:D:134:ILE:HD11	1.89	0.54
1:H:70:VAL:HG12	1:H:71:SER:N	2.22	0.54
1:G:46:THR:CG2	1:G:130:ARG:HH12	2.21	0.54
1:I:73:ASP:OD1	1:I:79:LYS:HG2	2.08	0.54
1:J:188:ASN:N	1:J:188:ASN:HD22	1.92	0.54
1:F:46:THR:CG2	1:F:130:ARG:HH12	2.20	0.54
1:J:92:LYS:HG3	1:J:125:TRP:CZ2	2.43	0.54
1:K:108:LEU:O	1:K:110:VAL:HG23	2.06	0.54
1:J:123:LYS:HB3	1:J:170:THR:HG22	1.90	0.54
1:L:139:MET:CE	1:L:192:MET:HB2	2.37	0.54
1:G:73:ASP:OD1	1:G:79:LYS:HG2	2.08	0.54
1:I:52:LEU:HD21	1:I:80:ILE:HD13	1.90	0.54
1:K:64:TYR:CD2	1:K:119:GLU:HB3	2.43	0.54
1:A:46:THR:CG2	1:A:130:ARG:HH12	2.21	0.54
1:E:106:ARG:HB3	1:I:68:THR:HG22	1.88	0.54

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:H:73:ASP:OD1	1:H:79:LYS:HG2	2.08	0.54
1:K:56:VAL:HG12	1:K:57:THR:N	2.23	0.54
1:K:92:LYS:HE3	1:K:166:ASP:OD2	2.08	0.53
1:B:73:ASP:OD1	1:B:79:LYS:HG2	2.09	0.53
1:D:70:VAL:HG12	1:D:71:SER:N	2.23	0.53
1:G:49:ASP:OD2	1:H:49:ASP:OD2	2.25	0.53
1:G:52:LEU:HD21	1:G:80:ILE:HD13	1.90	0.53
1:G:70:VAL:HG12	1:G:71:SER:N	2.23	0.53
1:A:73:ASP:OD1	1:A:79:LYS:HG2	2.08	0.53
1:C:70:VAL:HG12	1:C:71:SER:N	2.23	0.53
1:G:56:VAL:HG11	1:G:65:PHE:CB	2.18	0.53
1:A:70:VAL:HG12	1:A:71:SER:N	2.23	0.53
1:H:124:VAL:HG11	1:H:134:ILE:HD11	1.91	0.53
1:J:170:THR:HG23	2:J:1246:HOH:O	2.09	0.53
1:K:99:VAL:O	1:K:124:VAL:HA	2.08	0.53
1:C:73:ASP:OD1	1:C:79:LYS:HG2	2.08	0.53
1:F:73:ASP:OD1	1:F:79:LYS:HG2	2.08	0.53
1:I:46:THR:CG2	1:I:130:ARG:HH12	2.22	0.53
1:I:70:VAL:HG12	1:I:71:SER:N	2.24	0.52
1:K:101:THR:CG2	1:K:123:LYS:HZ2	2.23	0.52
1:K:83:ARG:HB3	1:K:86:LYS:HD2	1.91	0.52
1:E:105:GLY:O	1:I:68:THR:HG23	2.10	0.52
1:F:124:VAL:HG11	1:F:134:ILE:HD11	1.91	0.52
1:C:164:ASP:HB2	1:C:167:GLU:CD	2.30	0.52
1:L:123:LYS:HB3	1:L:170:THR:HG22	1.92	0.52
1:E:73:ASP:OD1	1:E:79:LYS:HG2	2.08	0.52
1:D:164:ASP:HB2	1:D:167:GLU:CD	2.30	0.52
1:B:164:ASP:HB2	1:B:167:GLU:CD	2.30	0.52
1:L:126:ASN:ND2	1:L:128:THR:HB	2.18	0.52
1:L:56:VAL:HG12	1:L:57:THR:N	2.25	0.52
1:A:124:VAL:HG11	1:A:134:ILE:HD11	1.91	0.52
1:G:75:ASP:OD2	1:I:199:SER:O	2.28	0.52
1:J:131:LYS:CE	1:J:162:SER:HB2	2.40	0.52
1:J:70:VAL:HG12	1:J:71:SER:H	1.75	0.52
1:C:56:VAL:HG11	1:C:65:PHE:CB	2.21	0.52
1:L:46:THR:CG2	1:L:130:ARG:HH12	2.23	0.52
1:K:139:MET:CE	1:K:190:PHE:CB	2.88	0.52
1:L:70:VAL:HG12	1:L:71:SER:N	2.25	0.52
1:C:46:THR:CG2	1:C:130:ARG:HH12	2.22	0.51
1:C:124:VAL:HG11	1:C:134:ILE:HD11	1.91	0.51
1:E:124:VAL:HG11	1:E:134:ILE:HD11	1.91	0.51
1:F:68:THR:O	1:F:70:VAL:HG23	2.09	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:J:46:THR:HG22	1:J:130:ARG:CZ	2.40	0.51
1:J:42:ASP:O	1:J:46:THR:HG23	2.11	0.51
1:I:164:ASP:HB2	1:I:167:GLU:CD	2.30	0.51
1:D:56:VAL:HG13	1:D:66:LYS:O	2.11	0.51
1:H:56:VAL:HG13	1:H:66:LYS:O	2.10	0.51
1:K:46:THR:CG2	1:K:130:ARG:HH12	2.24	0.51
1:A:164:ASP:HB2	1:A:167:GLU:CD	2.31	0.51
1:A:46:THR:HG21	2:A:1673:HOH:O	2.10	0.51
1:L:46:THR:HG21	2:L:2137:HOH:O	2.11	0.51
1:D:56:VAL:HG11	1:D:65:PHE:CB	2.22	0.51
1:K:106:ARG:HA	1:K:106:ARG:NE	2.26	0.51
1:K:54:ILE:O	1:K:56:VAL:HG23	2.10	0.51
1:K:60:VAL:HG21	1:K:183:ALA:H	1.74	0.51
1:B:124:VAL:HG11	1:B:134:ILE:HD11	1.93	0.51
1:H:188:ASN:HD22	1:H:188:ASN:N	2.00	0.51
1:E:164:ASP:HB2	1:E:167:GLU:CD	2.30	0.51
1:F:86:LYS:HE2	2:F:2344:HOH:O	2.09	0.51
1:L:46:THR:HG22	1:L:130:ARG:HH12	1.75	0.51
1:B:46:THR:HG21	2:B:1396:HOH:O	2.11	0.50
1:C:56:VAL:HG13	1:C:66:LYS:O	2.12	0.50
1:K:56:VAL:HG13	1:K:66:LYS:O	2.12	0.50
1:K:178:GLY:HA3	2:K:2471:HOH:O	2.12	0.50
1:E:92:LYS:HD2	2:E:1525:HOH:O	2.11	0.50
1:K:144:ASN:HB2	1:K:145:PRO:CD	2.42	0.50
1:L:144:ASN:HB2	1:L:145:PRO:HD2	1.92	0.50
1:C:56:VAL:CG1	1:C:57:THR:N	2.74	0.50
2:J:1397:HOH:O	1:L:147:LYS:HE3	2.12	0.50
1:D:46:THR:HG21	2:D:1566:HOH:O	2.11	0.50
2:D:1346:HOH:O	1:F:72:PHE:HB2	2.11	0.50
1:G:128:THR:HG22	1:G:129:SER:N	2.26	0.50
1:J:45:GLU:CG	1:J:130:ARG:HH11	2.25	0.50
1:L:132:ILE:HD11	1:L:196:VAL:HG11	1.94	0.50
1:E:188:ASN:HD22	1:E:188:ASN:N	2.03	0.49
1:J:128:THR:HG22	1:J:129:SER:N	2.26	0.49
1:H:128:THR:HG22	1:H:129:SER:N	2.27	0.49
1:H:56:VAL:CG1	1:H:57:THR:N	2.75	0.49
1:A:56:VAL:CG1	1:A:57:THR:N	2.75	0.49
1:G:56:VAL:CG1	1:G:57:THR:N	2.75	0.49
1:K:137:MET:HE3	1:K:156:SER:HB3	1.94	0.49
1:H:131:LYS:HE2	1:H:162:SER:CB	2.40	0.49
2:H:1104:HOH:O	1:I:40:HIS:HD2	1.95	0.49
1:I:56:VAL:CG1	1:I:57:THR:N	2.75	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:G:164:ASP:HB2	1:G:167:GLU:CD	2.32	0.49
1:B:56:VAL:HG13	1:B:66:LYS:O	2.13	0.49
1:I:128:THR:HG22	1:I:129:SER:N	2.26	0.49
1:B:56:VAL:CG1	1:B:57:THR:N	2.76	0.49
1:E:37:ILE:CD1	1:F:33:VAL:HG13	2.42	0.49
1:H:164:ASP:HB2	1:H:167:GLU:CD	2.32	0.49
1:I:144:ASN:HB2	1:I:145:PRO:CD	2.43	0.49
1:K:199:SER:OG	1:K:200:GLY:N	2.43	0.49
1:D:128:THR:HG22	1:D:129:SER:N	2.28	0.49
1:E:128:THR:HG22	1:E:129:SER:N	2.28	0.49
1:F:128:THR:HG22	1:F:129:SER:N	2.28	0.49
1:F:56:VAL:CG1	1:F:57:THR:N	2.76	0.49
1:I:188:ASN:ND2	1:I:188:ASN:H	2.08	0.49
1:J:139:MET:CE	1:J:192:MET:HA	2.42	0.49
1:L:128:THR:HG22	1:L:129:SER:N	2.27	0.49
1:E:56:VAL:CG1	1:E:57:THR:N	2.76	0.48
1:A:56:VAL:HG13	1:A:66:LYS:O	2.13	0.48
1:C:34:THR:O	1:C:38:GLU:HG3	2.13	0.48
1:D:56:VAL:CG1	1:D:57:THR:N	2.76	0.48
1:H:46:THR:CG2	1:H:130:ARG:HH12	2.25	0.48
1:J:139:MET:HE3	1:J:192:MET:CA	2.43	0.48
1:F:34:THR:O	1:F:38:GLU:HG3	2.14	0.48
1:L:164:ASP:HB2	1:L:167:GLU:OE1	2.13	0.48
1:L:132:ILE:HD11	1:L:196:VAL:CG1	2.43	0.48
1:L:46:THR:HG22	1:L:130:ARG:CZ	2.43	0.48
1:B:128:THR:HG22	1:B:129:SER:N	2.27	0.48
1:C:128:THR:HG22	1:C:129:SER:N	2.28	0.48
1:D:34:THR:O	1:D:38:GLU:HG3	2.14	0.48
1:E:56:VAL:HG13	1:E:66:LYS:O	2.13	0.48
1:K:128:THR:HG22	1:K:129:SER:N	2.28	0.48
1:K:188:ASN:HD22	1:K:188:ASN:N	2.07	0.48
1:F:46:THR:HG23	1:F:130:ARG:HH12	1.77	0.48
1:F:164:ASP:HB2	1:F:167:GLU:CD	2.33	0.48
1:G:49:ASP:OD2	1:I:49:ASP:OD2	2.32	0.48
1:C:188:ASN:N	1:C:188:ASN:HD22	2.01	0.48
1:G:46:THR:HG22	1:G:130:ARG:CZ	2.44	0.48
1:E:101:THR:HB	1:E:123:LYS:HZ2	1.78	0.48
1:G:197:THR:HG22	1:G:199:SER:N	2.07	0.48
1:H:34:THR:O	1:H:38:GLU:HG3	2.13	0.48
1:K:52:LEU:HD11	1:K:72:PHE:CE1	2.48	0.48
1:F:56:VAL:HG13	1:F:66:LYS:O	2.14	0.48
1:D:201:THR:HG23	1:F:73:ASP:O	2.14	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:37:ILE:HD11	1:E:33:VAL:HG13	1.96	0.47
1:E:46:THR:HG22	1:E:130:ARG:CZ	2.43	0.47
1:K:49:ASP:OD1	1:K:197:THR:CG2	2.62	0.47
1:L:32:THR:HG22	1:L:35:LYS:H	1.79	0.47
1:A:110:VAL:HG11	2:A:1543:HOH:O	2.14	0.47
1:A:128:THR:HG22	1:A:129:SER:N	2.28	0.47
1:D:188:ASN:N	1:D:188:ASN:HD22	2.02	0.47
1:E:34:THR:O	1:E:38:GLU:HG3	2.12	0.47
1:J:49:ASP:HB2	1:K:49:ASP:OD2	2.14	0.47
1:A:34:THR:O	1:A:38:GLU:HG3	2.14	0.47
1:B:46:THR:HG22	1:B:130:ARG:CZ	2.44	0.47
1:D:46:THR:HG22	1:D:130:ARG:CZ	2.44	0.47
1:G:61:GLY:O	1:G:112:VAL:CG1	2.62	0.47
1:D:42:ASP:O	1:D:46:THR:HG23	2.15	0.47
1:F:139:MET:HE3	1:F:192:MET:CA	2.41	0.47
1:F:101:THR:HB	1:F:123:LYS:HZ2	1.79	0.47
1:B:34:THR:O	1:B:38:GLU:HG3	2.15	0.47
2:E:1568:HOH:O	1:I:70:VAL:HG22	2.14	0.47
1:C:110:VAL:HG11	2:C:1316:HOH:O	2.15	0.47
1:C:139:MET:HE3	1:C:192:MET:CA	2.41	0.47
1:B:49:ASP:OD2	1:C:49:ASP:OD2	2.33	0.47
1:I:34:THR:O	1:I:38:GLU:HG3	2.14	0.47
1:J:50:GLU:HG2	1:J:77:LEU:CD1	2.45	0.47
2:A:1254:HOH:O	1:C:40:HIS:HD2	1.98	0.47
1:K:140:VAL:O	1:K:190:PHE:HB3	2.15	0.47
1:F:131:LYS:HE2	1:F:162:SER:CB	2.41	0.47
1:H:61:GLY:O	1:H:112:VAL:CG1	2.62	0.47
1:J:61:GLY:O	1:J:112:VAL:CG1	2.63	0.47
1:J:32:THR:HG23	1:J:35:LYS:H	1.80	0.47
1:L:99:VAL:O	1:L:124:VAL:HA	2.15	0.47
1:D:114:GLN:HG2	2:D:1203:HOH:O	2.15	0.46
1:D:61:GLY:O	1:D:112:VAL:CG1	2.64	0.46
1:G:34:THR:O	1:G:38:GLU:HG3	2.14	0.46
1:A:79:LYS:HB3	2:A:2075:HOH:O	2.14	0.46
1:E:131:LYS:HE2	1:E:162:SER:CB	2.42	0.46
1:B:164:ASP:HB2	1:B:167:GLU:OE1	2.15	0.46
1:G:56:VAL:HG13	1:G:66:LYS:O	2.15	0.46
1:J:139:MET:HE2	1:J:139:MET:HA	1.96	0.46
1:K:124:VAL:HG11	1:K:134:ILE:CD1	2.41	0.46
1:L:46:THR:HG22	1:L:130:ARG:NH1	2.30	0.46
1:I:46:THR:HG23	1:I:130:ARG:HH12	1.81	0.46
1:E:106:ARG:CB	1:I:68:THR:CG2	2.91	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:L:64:TYR:CD2	1:L:119:GLU:HB3	2.50	0.46
1:A:46:THR:HG22	1:A:130:ARG:CZ	2.45	0.46
1:H:101:THR:HB	1:H:123:LYS:HZ2	1.81	0.46
1:H:197:THR:HG23	2:H:1253:HOH:O	2.15	0.46
1:I:56:VAL:HG13	1:I:66:LYS:O	2.15	0.46
1:C:188:ASN:ND2	1:C:188:ASN:H	2.11	0.46
1:C:42:ASP:O	1:C:46:THR:HG23	2.15	0.46
1:F:72:PHE:CE1	1:F:80:ILE:HD12	2.51	0.46
1:G:144:ASN:HB2	1:G:145:PRO:CD	2.45	0.46
1:I:164:ASP:HB2	1:I:167:GLU:OE1	2.16	0.46
1:J:197:THR:HG23	2:J:1534:HOH:O	2.15	0.46
1:K:59:GLU:HB3	1:K:62:SER:OG	2.16	0.46
1:L:198:TYR:O	1:L:199:SER:O	2.33	0.46
1:C:46:THR:HG22	1:C:130:ARG:CZ	2.45	0.46
1:C:61:GLY:O	1:C:112:VAL:CG1	2.64	0.46
1:F:46:THR:HG22	1:F:130:ARG:CZ	2.44	0.46
1:J:46:THR:HG22	1:J:130:ARG:NH1	2.30	0.46
1:J:128:THR:HG22	1:J:130:ARG:H	1.80	0.46
1:A:46:THR:HG23	1:A:130:ARG:HH12	1.80	0.46
1:B:72:PHE:CE1	1:B:80:ILE:HD12	2.51	0.46
1:C:101:THR:HB	1:C:123:LYS:HZ2	1.80	0.46
1:E:46:THR:HG23	1:E:130:ARG:HH12	1.80	0.46
1:D:144:ASN:HB2	1:D:145:PRO:CD	2.46	0.46
1:C:170:THR:HG23	2:C:2423:HOH:O	2.16	0.45
1:J:137:MET:HA	1:J:193:TYR:O	2.15	0.45
1:C:32:THR:HG22	1:C:35:LYS:H	1.81	0.45
1:E:188:ASN:ND2	1:E:188:ASN:H	2.11	0.45
1:G:61:GLY:O	1:G:112:VAL:HG13	2.15	0.45
1:I:139:MET:HE3	1:I:192:MET:CA	2.43	0.45
1:J:128:THR:CG2	1:J:129:SER:N	2.78	0.45
1:C:131:LYS:HE2	1:C:162:SER:CB	2.43	0.45
1:D:40:HIS:HD2	2:D:1558:HOH:O	1.99	0.45
1:G:201:THR:CG2	1:H:72:PHE:CD1	2.99	0.45
1:A:61:GLY:O	1:A:112:VAL:CG1	2.64	0.45
1:B:199:SER:OG	1:B:200:GLY:N	2.49	0.45
1:E:46:THR:HG22	1:E:130:ARG:HH12	1.81	0.45
1:E:144:ASN:HB2	1:E:145:PRO:CD	2.47	0.45
1:G:128:THR:CG2	1:G:129:SER:N	2.80	0.45
1:H:46:THR:HG23	1:H:130:ARG:HH12	1.82	0.45
1:J:124:VAL:HG22	1:J:163:ILE:HD12	1.94	0.45
1:J:139:MET:HE3	1:J:139:MET:HA	1.98	0.45
1:L:197:THR:HG22	1:L:199:SER:N	2.21	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:128:THR:O	1:D:165:GLU:HG2	2.16	0.45
1:E:42:ASP:O	1:E:46:THR:HG23	2.17	0.45
1:L:188:ASN:HD22	1:L:188:ASN:N	1.95	0.45
1:A:139:MET:HE3	1:A:192:MET:CA	2.42	0.45
1:G:91:MET:SD	1:G:123:LYS:NZ	2.81	0.45
1:I:128:THR:CG2	1:I:129:SER:N	2.79	0.45
1:L:56:VAL:CG1	1:L:57:THR:N	2.79	0.45
1:A:128:THR:O	1:A:165:GLU:HG2	2.17	0.45
1:B:46:THR:HG23	1:B:130:ARG:HH12	1.79	0.45
1:G:136:GLY:HA2	1:G:158:PHE:CZ	2.52	0.45
1:H:197:THR:HG21	2:I:1092:HOH:O	2.17	0.45
1:B:144:ASN:HB2	1:B:145:PRO:CD	2.47	0.45
1:C:128:THR:O	1:C:165:GLU:HG2	2.16	0.45
1:L:139:MET:HE2	1:L:192:MET:CB	2.46	0.45
1:B:188:ASN:N	1:B:188:ASN:HD22	2.01	0.45
1:E:128:THR:CG2	1:E:129:SER:N	2.80	0.45
1:F:140:VAL:HA	1:F:141:PRO:HD3	1.89	0.45
1:H:139:MET:HE3	1:H:192:MET:CA	2.40	0.45
1:I:56:VAL:HG11	1:I:65:PHE:CB	2.23	0.45
1:B:42:ASP:O	1:B:46:THR:HG23	2.17	0.44
1:E:139:MET:HE3	1:E:192:MET:CA	2.43	0.44
1:G:46:THR:HG23	1:G:130:ARG:HH12	1.82	0.44
1:H:61:GLY:O	1:H:112:VAL:HG13	2.17	0.44
1:H:128:THR:CG2	1:H:129:SER:N	2.80	0.44
1:L:52:LEU:HB2	1:L:194:VAL:HG22	1.98	0.44
1:L:56:VAL:HG13	1:L:66:LYS:O	2.16	0.44
1:A:144:ASN:HB2	1:A:145:PRO:CD	2.47	0.44
1:D:199:SER:OG	1:D:200:GLY:N	2.50	0.44
1:G:188:ASN:ND2	1:G:188:ASN:H	2.10	0.44
1:J:53:ARG:HG3	1:J:193:TYR:CE1	2.52	0.44
1:K:46:THR:HG23	1:K:130:ARG:HH12	1.82	0.44
1:D:46:THR:HG23	1:D:130:ARG:HH12	1.81	0.44
1:E:164:ASP:HB2	1:E:167:GLU:OE1	2.17	0.44
1:I:42:ASP:O	1:I:46:THR:HG23	2.18	0.44
1:I:61:GLY:O	1:I:112:VAL:CG1	2.66	0.44
1:K:121:THR:HG23	1:K:172:VAL:HA	2.00	0.44
1:A:42:ASP:O	1:A:46:THR:HG23	2.18	0.44
1:D:128:THR:CG2	1:D:129:SER:N	2.80	0.44
1:D:61:GLY:O	1:D:112:VAL:HG13	2.18	0.44
1:F:188:ASN:HD22	1:F:188:ASN:N	1.98	0.44
1:G:201:THR:HG22	1:G:201:THR:O	2.17	0.44
1:I:110:VAL:HG11	2:I:1192:HOH:O	2.17	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:I:32:THR:HG22	1:I:35:LYS:H	1.82	0.44
1:A:128:THR:CG2	1:A:129:SER:N	2.80	0.44
1:A:188:ASN:HD22	1:A:188:ASN:N	2.02	0.44
1:A:72:PHE:CE1	1:A:80:ILE:HD12	2.52	0.44
1:B:128:THR:CG2	1:B:129:SER:N	2.80	0.44
1:B:92:LYS:HD2	2:B:2227:HOH:O	2.17	0.44
1:G:128:THR:O	1:G:165:GLU:HG2	2.17	0.44
1:H:72:PHE:CE1	1:H:80:ILE:HD12	2.52	0.44
1:I:144:ASN:HB2	1:I:145:PRO:HD2	2.00	0.44
1:J:46:THR:HG22	1:J:130:ARG:HH12	1.83	0.44
1:K:87:SER:O	1:K:91:MET:HG3	2.17	0.44
1:J:46:THR:HG23	1:J:130:ARG:HH12	1.83	0.44
1:H:140:VAL:HA	1:H:141:PRO:HD3	1.89	0.44
1:I:46:THR:HG22	1:I:130:ARG:CZ	2.46	0.44
1:B:128:THR:O	1:B:165:GLU:HG2	2.18	0.44
1:E:199:SER:OG	1:E:200:GLY:N	2.50	0.44
1:H:144:ASN:HB2	1:H:145:PRO:CD	2.48	0.44
1:A:164:ASP:HB2	1:A:167:GLU:OE1	2.18	0.43
1:B:61:GLY:O	1:B:112:VAL:CG1	2.66	0.43
1:I:131:LYS:HE2	1:I:162:SER:CB	2.41	0.43
1:J:144:ASN:HB2	1:J:145:PRO:CD	2.48	0.43
1:D:72:PHE:CE1	1:D:80:ILE:HD12	2.53	0.43
1:E:61:GLY:O	1:E:112:VAL:CG1	2.66	0.43
1:E:46:THR:HG22	1:E:130:ARG:NH1	2.33	0.43
1:F:128:THR:CG2	1:F:129:SER:N	2.81	0.43
1:H:199:SER:OG	1:H:200:GLY:N	2.50	0.43
1:I:91:MET:SD	1:I:123:LYS:NZ	2.85	0.43
1:K:175:THR:HG22	1:K:176:THR:N	2.34	0.43
1:C:128:THR:CG2	1:C:129:SER:N	2.81	0.43
1:C:164:ASP:HB2	1:C:167:GLU:OE1	2.17	0.43
1:D:37:ILE:HG21	1:E:36:THR:HG21	2.00	0.43
1:E:194:VAL:HG23	1:E:194:VAL:O	2.17	0.43
1:G:164:ASP:HB2	1:G:167:GLU:OE1	2.18	0.43
1:G:199:SER:OG	1:G:200:GLY:N	2.50	0.43
1:J:45:GLU:OE1	1:J:130:ARG:HD2	2.18	0.43
1:J:32:THR:CG2	1:J:35:LYS:H	2.31	0.43
1:K:116:LEU:HD23	1:K:180:PRO:HA	2.01	0.43
1:A:61:GLY:O	1:A:112:VAL:HG13	2.18	0.43
1:I:199:SER:OG	1:I:200:GLY:N	2.49	0.43
1:E:72:PHE:CE1	1:E:80:ILE:HD12	2.53	0.43
1:F:128:THR:O	1:F:165:GLU:HG2	2.19	0.43
1:I:197:THR:HG21	2:I:1086:HOH:O	2.18	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:164:ASP:HB2	1:D:167:GLU:OE1	2.18	0.43
1:J:56:VAL:HG11	1:J:65:PHE:CB	2.38	0.43
1:C:46:THR:HG23	1:C:130:ARG:HH12	1.84	0.43
1:C:144:ASN:HB2	1:C:145:PRO:CD	2.48	0.43
1:C:61:GLY:O	1:C:112:VAL:HG13	2.19	0.43
1:G:42:ASP:O	1:G:46:THR:HG23	2.19	0.43
1:A:101:THR:HB	1:A:123:LYS:HZ2	1.84	0.43
1:I:128:THR:O	1:I:165:GLU:HG2	2.18	0.43
1:A:32:THR:HG22	1:A:35:LYS:H	1.84	0.43
1:E:105:GLY:O	1:I:68:THR:CG2	2.67	0.43
1:E:128:THR:O	1:E:165:GLU:HG2	2.19	0.43
1:K:40:HIS:O	1:K:44:ILE:HG13	2.18	0.43
1:L:33:VAL:HG12	1:L:37:ILE:HG13	2.01	0.43
1:F:144:ASN:HB2	1:F:145:PRO:CD	2.48	0.43
1:G:72:PHE:CE1	1:G:80:ILE:HD12	2.53	0.43
1:I:194:VAL:HG23	1:I:194:VAL:O	2.19	0.43
1:D:124:VAL:HG11	1:D:134:ILE:CD1	2.49	0.42
1:J:132:ILE:HD12	1:J:198:TYR:CZ	2.54	0.42
1:K:68:THR:O	1:K:70:VAL:HG23	2.19	0.42
1:G:46:THR:HG22	1:G:130:ARG:HH12	1.83	0.42
1:J:40:HIS:O	1:J:44:ILE:HG13	2.20	0.42
1:B:46:THR:HG22	1:B:130:ARG:NH1	2.34	0.42
1:C:136:GLY:HA2	1:C:158:PHE:CZ	2.54	0.42
1:F:136:GLY:HA2	1:F:158:PHE:CZ	2.54	0.42
1:I:72:PHE:CE1	1:I:80:ILE:HD12	2.54	0.42
1:D:194:VAL:O	1:D:194:VAL:HG23	2.19	0.42
1:D:201:THR:HG22	1:D:201:THR:O	2.20	0.42
1:F:32:THR:HG22	1:F:35:LYS:H	1.84	0.42
1:G:144:ASN:HB2	1:G:145:PRO:HD2	2.00	0.42
1:I:140:VAL:HA	1:I:141:PRO:HD3	1.90	0.42
1:K:52:LEU:HB3	1:K:194:VAL:HG22	2.02	0.42
1:L:72:PHE:CE1	1:L:80:ILE:HB	2.54	0.42
1:A:199:SER:OG	1:A:200:GLY:N	2.50	0.42
1:B:139:MET:CE	1:B:139:MET:CA	2.96	0.42
1:B:147:LYS:HE3	2:C:1887:HOH:O	2.19	0.42
1:D:110:VAL:HG12	1:D:111:THR:N	2.34	0.42
1:D:46:THR:HG22	1:D:130:ARG:HH12	1.84	0.42
1:G:124:VAL:HG11	1:G:134:ILE:CD1	2.50	0.42
1:A:91:MET:SD	1:A:123:LYS:NZ	2.88	0.42
1:B:32:THR:HG22	1:B:35:LYS:H	1.84	0.42
1:H:197:THR:HG22	1:H:199:SER:N	2.09	0.42
1:E:144:ASN:HB2	1:E:145:PRO:HD2	2.02	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:F:72:PHE:HE1	1:F:80:ILE:HD12	1.85	0.42
1:F:42:ASP:O	1:F:46:THR:HG23	2.20	0.42
1:H:42:ASP:O	1:H:46:THR:HG23	2.19	0.42
1:I:136:GLY:HA2	1:I:158:PHE:CZ	2.55	0.42
1:J:53:ARG:HG3	1:J:193:TYR:CZ	2.55	0.42
1:K:56:VAL:CG1	1:K:57:THR:N	2.83	0.42
1:F:199:SER:OG	1:F:200:GLY:N	2.52	0.42
1:H:128:THR:O	1:H:165:GLU:HG2	2.19	0.42
1:K:141:PRO:HA	1:K:190:PHE:CD1	2.38	0.42
1:K:70:VAL:CG1	1:K:71:SER:N	2.82	0.42
1:A:110:VAL:HG12	1:A:111:THR:N	2.35	0.41
1:A:136:GLY:HA2	1:A:158:PHE:CZ	2.55	0.41
1:B:46:THR:HG22	1:B:130:ARG:HH12	1.83	0.41
1:D:136:GLY:HA2	1:D:158:PHE:CZ	2.54	0.41
1:E:201:THR:O	1:E:201:THR:HG22	2.20	0.41
2:D:1558:HOH:O	1:F:40:HIS:HD2	2.03	0.41
1:G:110:VAL:HG12	1:G:111:THR:N	2.35	0.41
1:H:106:ARG:HA	1:H:106:ARG:NE	2.34	0.41
1:L:128:THR:CG2	1:L:129:SER:N	2.83	0.41
1:A:131:LYS:HE2	1:A:162:SER:CB	2.42	0.41
1:A:201:THR:O	1:A:201:THR:HG22	2.20	0.41
1:B:131:LYS:HE2	1:B:162:SER:CB	2.43	0.41
1:C:110:VAL:HG12	1:C:111:THR:N	2.35	0.41
1:C:201:THR:O	1:C:201:THR:HG22	2.20	0.41
1:H:61:GLY:C	1:H:112:VAL:HG13	2.40	0.41
1:C:140:VAL:HA	1:C:141:PRO:HD3	1.88	0.41
1:K:92:LYS:HG3	1:K:125:TRP:CZ2	2.55	0.41
2:J:2246:HOH:O	1:L:140:VAL:HG22	2.20	0.41
1:B:197:THR:HG22	1:B:199:SER:N	2.07	0.41
1:B:61:GLY:O	1:B:112:VAL:HG13	2.20	0.41
1:C:46:THR:HG22	1:C:130:ARG:HH12	1.84	0.41
1:D:144:ASN:HB2	1:D:145:PRO:HD2	2.02	0.41
1:F:110:VAL:HG12	1:F:111:THR:N	2.34	0.41
1:G:194:VAL:O	1:G:194:VAL:HG23	2.21	0.41
1:G:32:THR:HG22	1:G:35:LYS:H	1.86	0.41
1:J:56:VAL:CG1	1:J:57:THR:N	2.82	0.41
1:D:101:THR:HB	1:D:123:LYS:HZ2	1.85	0.41
1:F:61:GLY:O	1:F:112:VAL:CG1	2.68	0.41
1:F:139:MET:HE2	1:F:139:MET:HA	2.01	0.41
1:G:131:LYS:HE2	1:G:162:SER:CB	2.43	0.41
1:I:124:VAL:HG11	1:I:134:ILE:CD1	2.50	0.41
1:K:144:ASN:HB2	1:K:145:PRO:HD2	2.01	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:K:59:GLU:OE1	1:K:66:LYS:HE3	2.21	0.41
1:D:61:GLY:C	1:D:112:VAL:HG13	2.41	0.41
1:D:197:THR:HG22	1:D:198:TYR:N	2.36	0.41
1:G:46:THR:HG22	1:G:130:ARG:NH1	2.35	0.41
1:H:164:ASP:HB2	1:H:167:GLU:OE1	2.19	0.41
1:H:32:THR:HG22	1:H:35:LYS:H	1.86	0.41
1:L:52:LEU:O	1:L:193:TYR:HA	2.20	0.41
1:C:199:SER:OG	1:C:200:GLY:N	2.50	0.41
1:C:72:PHE:CE1	1:C:80:ILE:HD12	2.55	0.41
1:D:32:THR:HG22	1:D:35:LYS:H	1.85	0.41
1:F:201:THR:HG22	1:F:201:THR:O	2.20	0.41
1:L:197:THR:HG23	2:L:1350:HOH:O	2.20	0.41
1:L:199:SER:OG	1:L:200:GLY:N	2.50	0.41
1:E:37:ILE:HD11	1:F:33:VAL:HG13	2.03	0.41
1:H:139:MET:HA	1:H:139:MET:HE2	2.01	0.41
1:I:106:ARG:NE	1:I:106:ARG:HA	2.36	0.41
1:I:110:VAL:HG12	1:I:111:THR:N	2.34	0.41
1:K:91:MET:C	1:K:93:GLU:H	2.22	0.41
1:I:61:GLY:C	1:I:112:VAL:HG13	2.41	0.41
1:A:144:ASN:HB2	1:A:145:PRO:HD2	2.02	0.41
1:J:92:LYS:HG2	2:J:1873:HOH:O	2.21	0.41
1:K:124:VAL:HG13	1:K:163:ILE:HD12	2.03	0.41
1:F:188:ASN:H	1:F:188:ASN:ND2	2.07	0.41
1:B:110:VAL:HG12	1:B:111:THR:N	2.35	0.40
1:B:72:PHE:HE1	1:B:80:ILE:HD12	1.86	0.40
1:A:46:THR:HG22	1:A:130:ARG:HH12	1.86	0.40
1:B:201:THR:HG22	1:B:201:THR:O	2.20	0.40
1:B:70:VAL:HG22	2:B:1302:HOH:O	2.20	0.40
1:G:139:MET:HE3	1:G:192:MET:CA	2.46	0.40
1:H:144:ASN:HB2	1:H:145:PRO:HD2	2.04	0.40
1:H:46:THR:HG22	1:H:130:ARG:CZ	2.48	0.40
1:I:61:GLY:O	1:I:112:VAL:HG13	2.21	0.40
1:H:49:ASP:OD2	1:I:49:ASP:OD2	2.40	0.40
1:B:139:MET:HE3	1:B:192:MET:CA	2.45	0.40
1:B:144:ASN:HB2	1:B:145:PRO:HD2	2.02	0.40
1:C:46:THR:HG22	1:C:130:ARG:NH1	2.36	0.40
1:I:201:THR:O	1:I:201:THR:HG22	2.21	0.40
1:B:136:GLY:HA2	1:B:158:PHE:CZ	2.56	0.40
1:E:136:GLY:HA2	1:E:158:PHE:CZ	2.56	0.40
1:H:110:VAL:HG12	1:H:111:THR:N	2.35	0.40
1:H:141:PRO:HG3	2:H:1283:HOH:O	2.21	0.40
1:I:139:MET:CA	1:I:139:MET:CE	2.97	0.40

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:K:92:LYS:O	1:K:92:LYS:HG2	2.21	0.40
1:C:61:GLY:C	1:C:112:VAL:HG13	2.42	0.40
1:E:110:VAL:HG12	1:E:111:THR:N	2.36	0.40
1:I:101:THR:HB	1:I:123:LYS:HZ2	1.86	0.40
1:J:126:ASN:HD22	1:J:126:ASN:C	2.25	0.40
1:K:198:TYR:O	1:K:199:SER:O	2.39	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:60:VAL:CG1	1:H:59:GLU:OE2[3_656]	1.92	0.28

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	168/170 (99%)	163 (97%)	4 (2%)	1 (1%)	33	24
1	B	168/170 (99%)	163 (97%)	4 (2%)	1 (1%)	33	24
1	C	168/170 (99%)	163 (97%)	4 (2%)	1 (1%)	33	24
1	D	168/170 (99%)	163 (97%)	4 (2%)	1 (1%)	33	24
1	E	168/170 (99%)	162 (96%)	5 (3%)	1 (1%)	33	24
1	F	168/170 (99%)	163 (97%)	4 (2%)	1 (1%)	33	24
1	G	168/170 (99%)	162 (96%)	5 (3%)	1 (1%)	33	24
1	H	168/170 (99%)	163 (97%)	4 (2%)	1 (1%)	33	24
1	I	168/170 (99%)	162 (96%)	5 (3%)	1 (1%)	33	24
1	J	168/170 (99%)	159 (95%)	8 (5%)	1 (1%)	33	24
1	K	168/170 (99%)	158 (94%)	8 (5%)	2 (1%)	19	9
1	L	168/170 (99%)	163 (97%)	3 (2%)	2 (1%)	19	9
All	All	2016/2040 (99%)	1944 (96%)	58 (3%)	14 (1%)	30	20

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	199	SER
1	B	199	SER
1	C	199	SER
1	D	199	SER
1	E	199	SER
1	F	199	SER
1	G	199	SER
1	H	199	SER
1	I	199	SER
1	K	199	SER
1	L	199	SER
1	J	199	SER
1	L	112	VAL
1	K	92	LYS

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	146/146 (100%)	139 (95%)	7 (5%)	35	28
1	B	146/146 (100%)	139 (95%)	7 (5%)	35	28
1	C	146/146 (100%)	139 (95%)	7 (5%)	35	28
1	D	146/146 (100%)	139 (95%)	7 (5%)	35	28
1	E	146/146 (100%)	139 (95%)	7 (5%)	35	28
1	F	146/146 (100%)	139 (95%)	7 (5%)	35	28
1	G	146/146 (100%)	139 (95%)	7 (5%)	35	28
1	H	146/146 (100%)	139 (95%)	7 (5%)	35	28
1	I	146/146 (100%)	139 (95%)	7 (5%)	35	28
1	J	146/146 (100%)	139 (95%)	7 (5%)	35	28
1	K	146/146 (100%)	141 (97%)	5 (3%)	49	45
1	L	146/146 (100%)	140 (96%)	6 (4%)	41	35
All	All	1752/1752 (100%)	1671 (95%)	81 (5%)	37	30

All (81) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	40	HIS
1	A	52	LEU
1	A	108	LEU
1	A	124	VAL
1	A	126	ASN
1	A	139	MET
1	A	188	ASN
1	B	40	HIS
1	B	52	LEU
1	B	108	LEU
1	B	124	VAL
1	B	126	ASN
1	B	139	MET
1	B	188	ASN
1	C	40	HIS
1	C	52	LEU
1	C	108	LEU
1	C	124	VAL
1	C	126	ASN
1	C	139	MET
1	C	188	ASN
1	D	40	HIS
1	D	52	LEU
1	D	108	LEU
1	D	124	VAL
1	D	126	ASN
1	D	139	MET
1	D	188	ASN
1	E	40	HIS
1	E	52	LEU
1	E	108	LEU
1	E	124	VAL
1	E	126	ASN
1	E	139	MET
1	E	188	ASN
1	F	40	HIS
1	F	52	LEU
1	F	108	LEU
1	F	124	VAL
1	F	126	ASN
1	F	139	MET
1	F	188	ASN

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Mol	Chain	Res	Type
1	G	40	HIS
1	G	52	LEU
1	G	108	LEU
1	G	124	VAL
1	G	126	ASN
1	G	139	MET
1	G	188	ASN
1	H	40	HIS
1	H	52	LEU
1	H	108	LEU
1	H	124	VAL
1	H	126	ASN
1	H	139	MET
1	H	188	ASN
1	I	40	HIS
1	I	52	LEU
1	I	108	LEU
1	I	124	VAL
1	I	126	ASN
1	I	139	MET
1	I	188	ASN
1	J	32	THR
1	J	40	HIS
1	J	108	LEU
1	J	124	VAL
1	J	126	ASN
1	J	139	MET
1	J	188	ASN
1	K	40	HIS
1	K	72	PHE
1	K	126	ASN
1	K	170	THR
1	K	188	ASN
1	L	38	GLU
1	L	40	HIS
1	L	108	LEU
1	L	124	VAL
1	L	126	ASN
1	L	188	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (48) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	40	HIS
1	A	126	ASN
1	A	127	ASN
1	A	188	ASN
1	B	126	ASN
1	B	127	ASN
1	B	188	ASN
1	C	40	HIS
1	C	126	ASN
1	C	127	ASN
1	C	188	ASN
1	D	40	HIS
1	D	126	ASN
1	D	127	ASN
1	D	188	ASN
1	E	40	HIS
1	E	126	ASN
1	E	127	ASN
1	E	188	ASN
1	F	40	HIS
1	F	126	ASN
1	F	127	ASN
1	F	188	ASN
1	G	40	HIS
1	G	126	ASN
1	G	127	ASN
1	G	188	ASN
1	H	40	HIS
1	H	126	ASN
1	H	127	ASN
1	H	188	ASN
1	I	40	HIS
1	I	126	ASN
1	I	127	ASN
1	I	188	ASN
1	J	40	HIS
1	J	126	ASN
1	J	127	ASN
1	J	188	ASN
1	K	40	HIS
1	K	43	ASN
1	K	126	ASN
1	K	127	ASN

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Mol	Chain	Res	Type
1	K	188	ASN
1	L	126	ASN
1	L	127	ASN
1	L	138	GLN
1	L	188	ASN

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	170/170 (100%)	0.38	7 (4%) 35 35	8, 17, 47, 74	0
1	B	170/170 (100%)	0.37	10 (5%) 22 21	6, 14, 40, 110	0
1	C	170/170 (100%)	0.24	6 (3%) 42 41	7, 15, 38, 67	0
1	D	170/170 (100%)	0.41	10 (5%) 22 21	6, 16, 39, 87	0
1	E	170/170 (100%)	0.35	11 (6%) 18 18	7, 13, 39, 84	0
1	F	170/170 (100%)	0.48	15 (8%) 10 9	4, 13, 38, 66	0
1	G	170/170 (100%)	0.68	16 (9%) 9 8	5, 13, 48, 124	0
1	H	170/170 (100%)	0.67	16 (9%) 9 8	6, 17, 53, 89	0
1	I	170/170 (100%)	0.35	5 (2%) 49 49	6, 16, 37, 60	0
1	J	170/170 (100%)	0.18	0 100 100	7, 17, 38, 56	0
1	K	170/170 (100%)	0.82	10 (5%) 22 21	12, 29, 51, 68	0
1	L	170/170 (100%)	0.42	9 (5%) 25 25	10, 21, 41, 66	0
All	All	2040/2040 (100%)	0.45	115 (5%) 24 23	4, 17, 45, 124	0

All (115) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	32	THR	20.3
1	G	34	THR	18.3
1	B	32	THR	13.1
1	G	36	THR	9.7
1	G	33	VAL	9.1
1	F	33	VAL	8.8
1	G	37	ILE	8.5
1	H	37	ILE	8.3
1	H	33	VAL	8.3
1	G	35	LYS	7.8
1	H	34	THR	7.4

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Mol	Chain	Res	Type	RSRZ
1	G	38	GLU	6.9
1	D	34	THR	6.3
1	E	32	THR	5.6
1	K	32	THR	5.6
1	D	33	VAL	5.5
1	A	34	THR	5.4
1	D	35	LYS	5.1
1	F	32	THR	5.1
1	E	33	VAL	5.1
1	H	32	THR	4.7
1	D	201	THR	4.7
1	B	35	LYS	4.7
1	D	32	THR	4.6
1	F	36	THR	4.6
1	H	36	THR	4.6
1	H	35	LYS	4.6
1	E	201	THR	4.4
1	B	33	VAL	4.3
1	G	39	THR	4.2
1	F	72	PHE	4.2
1	F	34	THR	4.0
1	A	199	SER	4.0
1	K	103	VAL	4.0
1	L	199	SER	3.9
1	D	37	ILE	3.8
1	D	199	SER	3.8
1	H	38	GLU	3.8
1	B	34	THR	3.7
1	E	36	THR	3.7
1	B	199	SER	3.7
1	F	39	THR	3.7
1	H	199	SER	3.7
1	E	199	SER	3.6
1	F	110	VAL	3.6
1	A	35	LYS	3.6
1	F	199	SER	3.5
1	I	199	SER	3.5
1	B	38	GLU	3.3
1	A	37	ILE	3.3
1	A	33	VAL	3.3
1	F	112	VAL	3.3
1	H	42	ASP	3.2

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Mol	Chain	Res	Type	RSRZ
1	L	110	VAL	3.2
1	F	71	SER	3.1
1	E	72	PHE	3.1
1	A	32	THR	3.0
1	C	199	SER	3.0
1	I	32	THR	3.0
1	G	199	SER	3.0
1	B	37	ILE	3.0
1	B	201	THR	3.0
1	G	92	LYS	3.0
1	L	72	PHE	3.0
1	F	37	ILE	2.9
1	D	36	THR	2.9
1	K	199	SER	2.9
1	F	201	THR	2.9
1	I	69	ASP	2.8
1	K	190	PHE	2.8
1	B	36	THR	2.8
1	K	181	ILE	2.8
1	A	72	PHE	2.8
1	F	35	LYS	2.7
1	C	92	LYS	2.7
1	K	33	VAL	2.7
1	D	38	GLU	2.6
1	K	89	ALA	2.6
1	H	201	THR	2.6
1	C	129	SER	2.6
1	H	112	VAL	2.6
1	E	34	THR	2.6
1	H	59	GLU	2.6
1	K	92	LYS	2.5
1	E	38	GLU	2.5
1	F	184	THR	2.4
1	I	70	VAL	2.4
1	H	41	THR	2.4
1	H	39	THR	2.4
1	I	68	THR	2.3
1	K	128	THR	2.3
1	G	200	GLY	2.3
1	L	92	LYS	2.3
1	L	49	ASP	2.3
1	C	32	THR	2.3

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Mol	Chain	Res	Type	RSRZ
1	K	170	THR	2.3
1	G	40	HIS	2.2
1	G	128	THR	2.2
1	B	49	ASP	2.2
1	F	111	THR	2.2
1	L	109	GLU	2.2
1	E	90	GLN	2.2
1	G	201	THR	2.2
1	D	49	ASP	2.2
1	C	167	GLU	2.1
1	L	108	LEU	2.1
1	E	73	ASP	2.1
1	L	111	THR	2.1
1	C	200	GLY	2.1
1	L	52	LEU	2.1
1	H	170	THR	2.0
1	E	71	SER	2.0
1	G	42	ASP	2.0
1	G	49	ASP	2.0
1	H	45	GLU	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.