



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 12:40 PM GMT

PDB ID : 3HM2
Title : Crystal Structure of Putative Precorrin-6Y C5,15-Methyltransferase Targeted Domain from *Corynebacterium diphtheriae*
Authors : Kim, Y.; Bigelow, L.; Keigher, L.; Joachimiak, A.; Midwest Center for Structural Genomics (MCSG)
Deposited on : 2009-05-28
Resolution : 2.21 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>

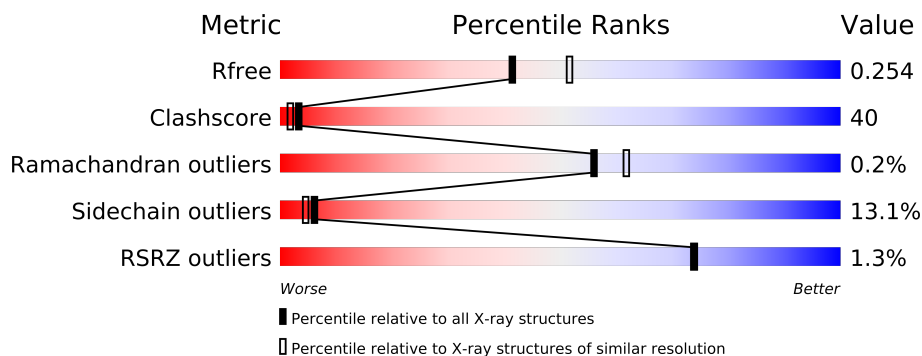
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.21 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	2938 (2.20-2.20)
Clashscore	79885	3751 (2.20-2.20)
Ramachandran outliers	78287	3681 (2.20-2.20)
Sidechain outliers	78261	3682 (2.20-2.20)
RSRZ outliers	66119	2939 (2.20-2.20)

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	178	
1	B	178	
1	C	178	
1	D	178	
1	E	178	
1	F	178	
1	G	178	
1	H	178	

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

Mol	Type	Chain	Res	Geometry	Electron density
2	ACY	F	176	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 10700 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 precorrin-6Y C5,15-methyltransferase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	171	Total	C	N	O	S	Se	0	2	0
			1322	842	239	238	1	2			
1	B	171	Total	C	N	O	S	Se	0	0	0
			1304	831	234	236	1	2			
1	C	169	Total	C	N	O	S	Se	0	0	0
			1294	826	232	233	1	2			
1	D	167	Total	C	N	O	S	Se	0	0	0
			1279	815	230	231	1	2			
1	E	170	Total	C	N	O	S	Se	0	0	0
			1298	828	233	234	1	2			
1	F	167	Total	C	N	O	S	Se	0	1	0
			1294	826	231	234	1	2			
1	G	171	Total	C	N	O	S	Se	0	0	0
			1295	824	232	236	1	2			
1	H	167	Total	C	N	O	S	Se	0	1	0
			1282	816	232	231	1	2			

There are 24 discrepancies between the modelled and reference sequences:

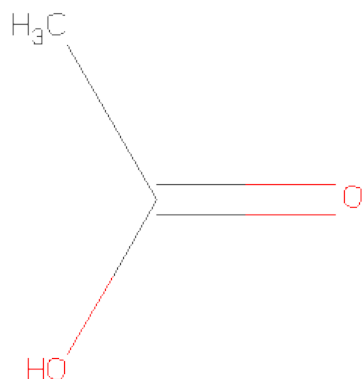
Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	expression tag	UNP Q6NHA4
A	-1	ASN	-	expression tag	UNP Q6NHA4
A	0	ALA	-	expression tag	UNP Q6NHA4
B	-2	SER	-	expression tag	UNP Q6NHA4
B	-1	ASN	-	expression tag	UNP Q6NHA4
B	0	ALA	-	expression tag	UNP Q6NHA4
C	-2	SER	-	expression tag	UNP Q6NHA4
C	-1	ASN	-	expression tag	UNP Q6NHA4
C	0	ALA	-	expression tag	UNP Q6NHA4
D	-2	SER	-	expression tag	UNP Q6NHA4
D	-1	ASN	-	expression tag	UNP Q6NHA4
D	0	ALA	-	expression tag	UNP Q6NHA4
E	-2	SER	-	expression tag	UNP Q6NHA4

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Chain	Residue	Modelled	Actual	Comment	Reference
E	-1	ASN	-	expression tag	UNP Q6NHA4
E	0	ALA	-	expression tag	UNP Q6NHA4
F	-2	SER	-	expression tag	UNP Q6NHA4
F	-1	ASN	-	expression tag	UNP Q6NHA4
F	0	ALA	-	expression tag	UNP Q6NHA4
G	-2	SER	-	expression tag	UNP Q6NHA4
G	-1	ASN	-	expression tag	UNP Q6NHA4
G	0	ALA	-	expression tag	UNP Q6NHA4
H	-2	SER	-	expression tag	UNP Q6NHA4
H	-1	ASN	-	expression tag	UNP Q6NHA4
H	0	ALA	-	expression tag	UNP Q6NHA4

- Molecule 2 is ACETIC ACID (three-letter code: ACY) (formula: C₂H₄O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			4	2	2		
2	D	1	Total	C	O	0	0
			4	2	2		
2	F	1	Total	C	O	0	0
			4	2	2		

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	G	1	Total Mg 1 1	0	0
3	B	1	Total Mg 1 1	0	0

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	G	1	Total Ca 1 1	0	0
4	D	1	Total Ca 1 1	0	0

- Molecule 5 is water.

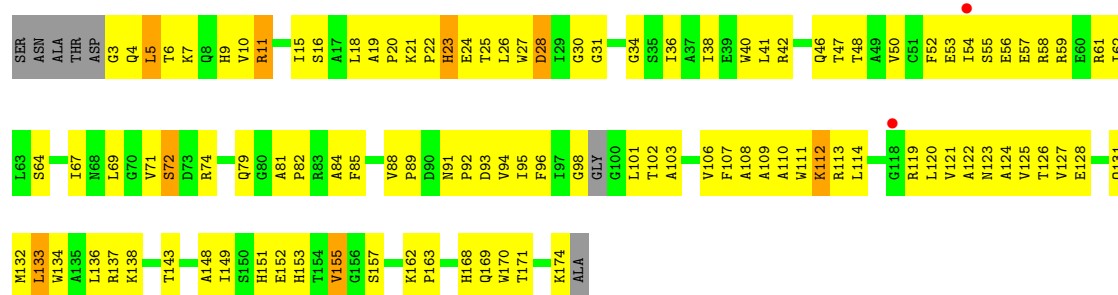
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	38	Total O 38 38	0	0
5	B	42	Total O 42 42	0	0
5	C	30	Total O 30 30	0	0
5	D	46	Total O 46 46	0	0
5	E	50	Total O 50 50	0	0
5	F	33	Total O 33 33	0	0
5	G	39	Total O 39 39	0	0
5	H	38	Total O 38 38	0	0

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.

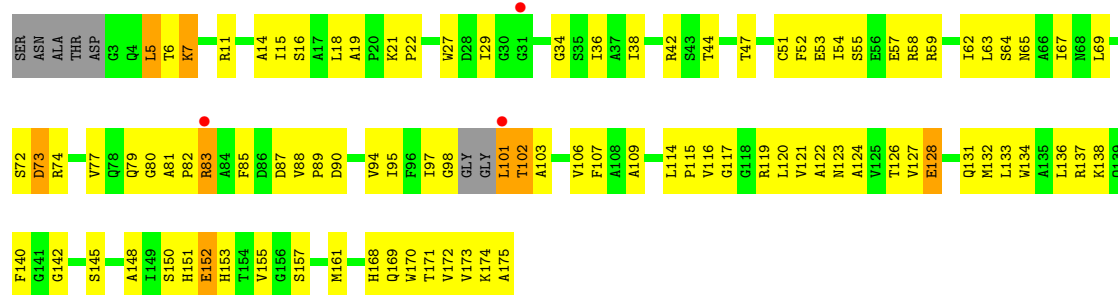
- Molecule 1: precorrin-6Y C5,15-methyltransferase

Chain A: 



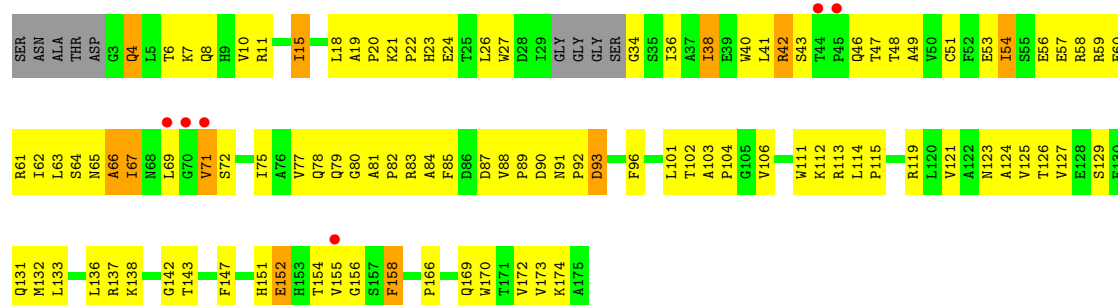
- Molecule 1: precorrin-6Y C5,15-methyltransferase

Chain B: 



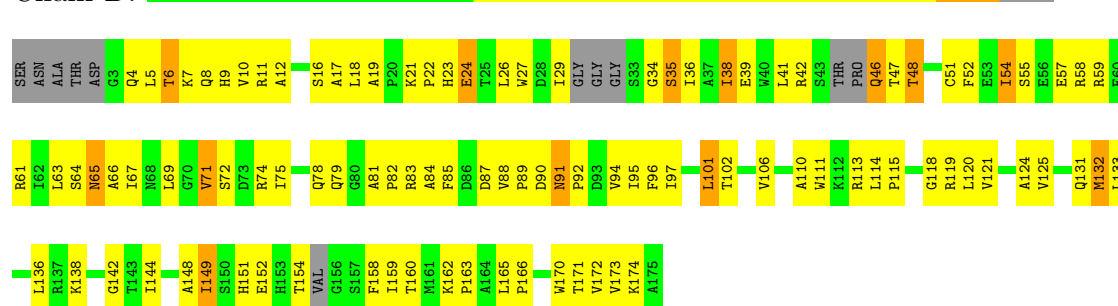
- Molecule 1: precorrin-6Y C5,15-methyltransferase

Chain C: 



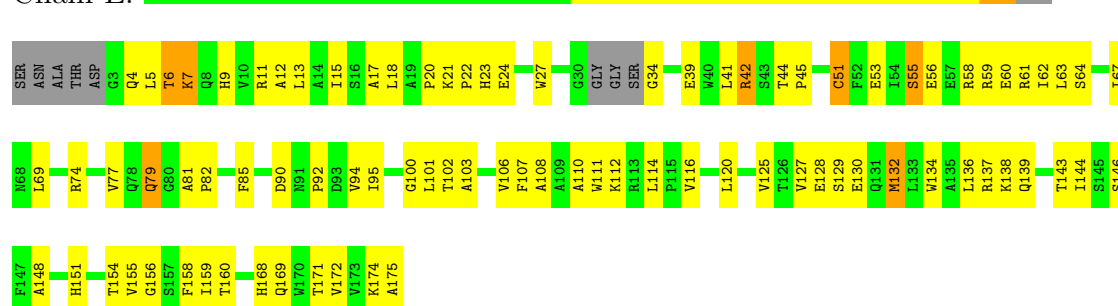
- Molecule 1: precorrin-6Y C5,15-methyltransferase

Chain D:



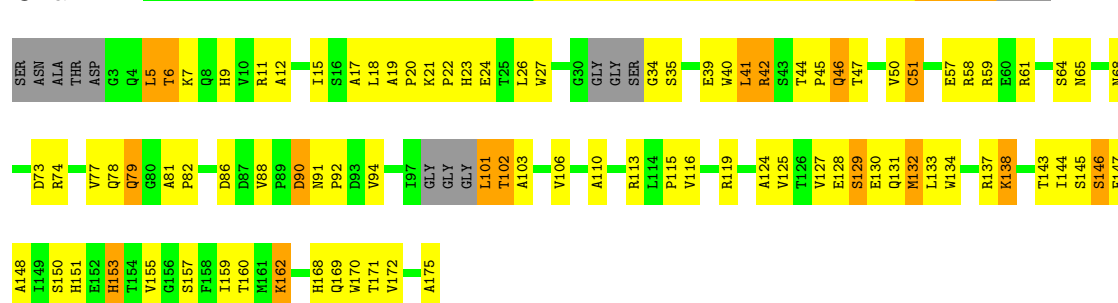
- Molecule 1: precorrin-6Y C5,15-methyltransferase

Chain E:



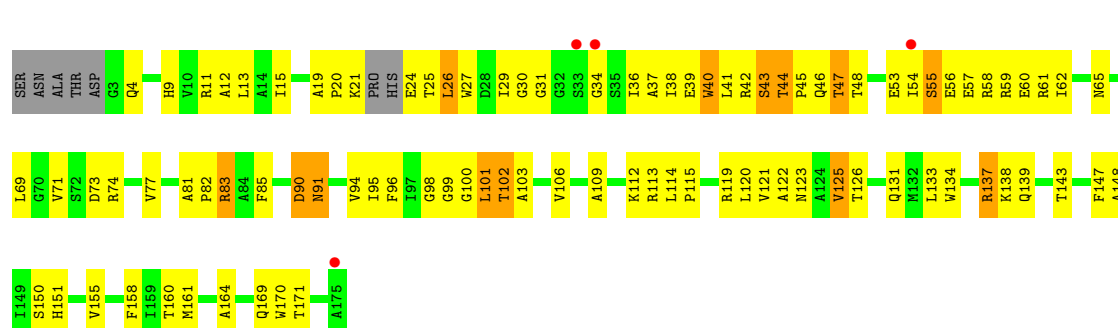
- Molecule 1: precorrin-6Y C5,15-methyltransferase

Chain F:



- Molecule 1: precorrin-6Y C5,15-methyltransferase

Chain G:



- Molecule 1: precorrin-6Y C5,15-methyltransferase

A148	A149	S150	H151	E152	H153	THR	VAL	G156	S157	F158	T159	M160	M161	K162	P163	W170	K174	A175	A176																																	
167	N68	V71	R74	G79	G80	A81	P82	R83	D86	D87	D90	N91	P92	D93	V94	P90	I95	F96	I97	G98	G99	G100	L101	V106	R113	L114	P115	V116	G117	G118	R119	L120	V121	A122	N123	A124	V125	T126	V127	E128	S129	E130	Q131	M132	L133	W134	A135	L136	R137	G138	S145	
ASN	THR	ASP	G3	Q4	L5	Q8	H9	V10	R11	A12	I15	S16	A17	L18	A19	P20	K21	P80	HIS	E24	T25	L26	W27	G31	GLY	SER	G34	S35	T36	A37	I38	E39	W40	L41	R42	S43	T44	P45	Q46	T47	T48	A49	V50	C51	I54	S55	E56	E57	R58	R59	E60	S64

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	77.78Å 113.85Å 84.65Å 90.00° 90.12° 90.00°	Depositor
Resolution (Å)	47.24 – 2.21 47.24 – 2.20	Depositor EDS
% Data completeness (in resolution range)	88.9 (47.24-2.21) 95.9 (47.24-2.20)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.17 (at 2.20Å)	Xtriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, R_{free}	0.210 , 0.260 0.214 , 0.254	Depositor DCC
R_{free} test set	3621 reflections (5.32%)	DCC
Wilson B-factor (Å ²)	36.2	Xtriage
Anisotropy	0.205	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 40.0	EDS
Estimated twinning fraction	0.420 for h,-k,-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	1 of 71721 reflections (0.001%)	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	10700	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 57.44 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 2.3673e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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

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	0.48	0/1350	0.64	1/1834 (0.1%)
1	B	0.52	0/1332	0.67	1/1810 (0.1%)
1	C	0.50	0/1322	0.63	0/1797
1	D	0.53	0/1304	0.69	0/1767
1	E	0.58	0/1326	0.69	0/1802
1	F	0.56	0/1321	0.71	1/1795 (0.1%)
1	G	0.57	0/1321	0.71	0/1793
1	H	0.53	1/1306 (0.1%)	0.66	0/1768
All	All	0.53	1/10582 (0.0%)	0.68	3/14366 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	H	57	GLU	C-N	5.86	1.47	1.34

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	119	ARG	NE-CZ-NH1	-6.47	117.06	120.30
1	B	133	LEU	CA-CB-CG	6.00	129.10	115.30
1	A	133	LEU	CA-CB-CG	5.28	127.45	115.30

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	1322	0	1334	132	0
1	B	1304	0	1313	103	0
1	C	1294	0	1305	106	0
1	D	1279	0	1285	116	0
1	E	1298	0	1308	90	0
1	F	1294	0	1301	100	0
1	G	1295	0	1305	114	0
1	H	1282	0	1291	124	0
2	A	4	0	3	0	0
2	D	4	0	3	0	0
2	F	4	0	3	1	0
3	B	1	0	0	0	0
3	G	1	0	0	0	0
4	D	1	0	0	0	0
4	G	1	0	0	0	0
5	A	38	0	0	10	0
5	B	42	0	0	1	0
5	C	30	0	0	4	0
5	D	46	0	0	8	0
5	E	50	0	0	2	0
5	F	33	0	0	2	0
5	G	39	0	0	4	0
5	H	38	0	0	2	0
All	All	10700	0	10451	838	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 40.

All (838) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:81:ALA:HB2	1:A:85:PHE:CE1	1.70	1.26
1:C:27:TRP:CE2	1:C:92:PRO:HG3	1.80	1.17
1:C:54:ILE:HD12	1:C:80:GLY:HA3	1.29	1.15
1:C:41:LEU:HB3	1:C:71:VAL:HG11	1.26	1.12
1:G:56:GLU:O	1:G:60:GLU:HG3	1.48	1.12
1:F:44:THR:CG2	1:F:47:THR:HG23	1.81	1.10
1:A:81:ALA:CB	1:A:85:PHE:CE1	2.37	1.08
1:H:55:SER:HB2	1:H:58[A]:ARG:HB2	1.35	1.07

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:H:57:GLU:HA	1:H:60:GLU:HG3	1.11	1.06
1:A:81:ALA:CB	1:A:85:PHE:CZ	2.40	1.05
1:H:55:SER:HB2	1:H:58[B]:ARG:HB2	1.35	1.04
1:H:57:GLU:HA	1:H:60:GLU:CG	1.87	1.03
1:H:57:GLU:CA	1:H:60:GLU:HG3	1.88	1.03
1:A:153:HIS:CD2	1:A:155:VAL:HG22	1.92	1.02
1:B:27:TRP:HB3	1:B:29:ILE:HD11	1.40	1.02
1:E:154:THR:HG22	1:E:156:GLY:H	1.21	1.01
1:F:44:THR:CG2	1:F:47:THR:CG2	2.39	1.00
1:G:20:PRO:HG3	1:G:40:TRP:CE3	1.96	0.99
1:A:81:ALA:HB1	1:A:85:PHE:CZ	1.95	0.99
1:F:7:LYS:HD2	1:F:11:ARG:HE	1.28	0.97
1:A:41:LEU:HD13	1:A:71:VAL:HG12	1.44	0.96
1:G:54:ILE:HD12	1:G:58:ARG:NH2	1.80	0.95
1:A:41:LEU:HB3	1:A:71:VAL:HG11	1.48	0.95
1:B:137:ARG:HG2	1:B:137:ARG:HH11	1.33	0.94
1:C:27:TRP:CD2	1:C:92:PRO:HG3	2.05	0.91
1:D:101:LEU:HD12	1:D:132:MSE:HG2	1.51	0.91
1:D:26:LEU:HD12	1:D:94:VAL:HG12	1.53	0.90
1:H:11:ARG:HD2	1:H:39:GLU:HG3	1.52	0.90
1:F:147:PHE:HD1	1:H:149:ILE:HG12	1.37	0.89
1:F:21:LYS:HG2	1:F:22:PRO:HD2	1.55	0.89
1:E:6:THR:HG23	1:E:151:HIS:NE2	1.88	0.88
1:D:87:ASP:O	1:D:89:PRO:HD3	1.73	0.88
1:D:24:GLU:H	1:D:47:THR:HG22	1.39	0.87
1:A:81:ALA:HB1	1:A:82:PRO:HA	1.57	0.86
1:H:9:HIS:NE2	1:H:151:HIS:HE1	1.73	0.85
1:H:91:ASN:HD21	1:H:113:ARG:HE	1.21	0.85
1:D:119:ARG:HG2	1:D:173:VAL:HG22	1.59	0.85
1:E:5:LEU:HD23	1:E:5:LEU:H	1.41	0.85
1:A:41:LEU:CD1	1:A:71:VAL:HG12	2.05	0.85
1:C:53:GLU:O	1:C:59:ARG:HD2	1.77	0.85
1:E:7:LYS:HD2	1:E:11:ARG:HH12	1.40	0.85
1:F:6:THR:HG23	1:F:151:HIS:NE2	1.92	0.84
1:A:153:HIS:HD2	1:A:155:VAL:HG22	1.38	0.84
1:C:59:ARG:NH2	1:C:79:GLN:HE21	1.75	0.84
1:B:101:LEU:O	1:B:101:LEU:HD12	1.77	0.84
1:D:82:PRO:HG3	1:D:106:VAL:HA	1.60	0.83
1:H:11:ARG:CD	1:H:39:GLU:HG3	2.09	0.82
1:B:148:ALA:HB3	1:D:148:ALA:HB3	1.60	0.82
1:A:24:GLU:CD	1:A:119:ARG:HH22	1.83	0.82
1:A:41:LEU:CB	1:A:71:VAL:HG11	2.10	0.81

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:116:VAL:HG22	1:B:175:ALA:HA	1.62	0.81
1:H:55:SER:HB2	1:H:58[A]:ARG:CB	2.10	0.81
1:E:17:ALA:HB2	1:E:169:GLN:HE22	1.45	0.81
1:G:101:LEU:HD11	1:G:122:ALA:HB1	1.63	0.80
1:D:6:THR:HA	1:D:9:HIS:HD2	1.45	0.80
1:G:47:THR:N	1:G:74:ARG:NH1	2.29	0.80
1:B:137:ARG:HG2	1:B:137:ARG:NH1	1.95	0.80
1:A:89:PRO:HD2	5:A:278:HOH:O	1.80	0.80
1:C:4:GLN:HA	1:C:7:LYS:HD2	1.64	0.79
1:A:41:LEU:HD13	1:A:71:VAL:CG1	2.13	0.78
1:D:63:LEU:O	1:D:67:ILE:HG13	1.84	0.78
1:A:132:MSE:HE3	1:A:136:LEU:HD11	1.65	0.78
1:E:158:PHE:CD2	1:H:126:THR:HG21	2.17	0.78
1:F:44:THR:HG22	1:F:47:THR:CG2	2.13	0.78
1:G:164:ALA:HB1	5:G:229:HOH:O	1.84	0.78
1:F:44:THR:HG21	1:F:47:THR:CG2	2.14	0.78
1:G:15:ILE:HD11	1:G:39:GLU:HB3	1.66	0.78
1:G:91:ASN:HB3	1:G:113:ARG:O	1.84	0.77
1:H:20:PRO:HG3	1:H:40:TRP:CE3	2.19	0.77
1:A:134:TRP:O	1:A:138:LYS:HG2	1.85	0.77
1:C:4:GLN:HG3	1:C:7:LYS:HD2	1.65	0.77
1:F:41:LEU:HD21	1:F:74:ARG:HB3	1.67	0.76
1:A:56:GLU:HG3	1:A:57:GLU:OE2	1.85	0.76
1:G:91:ASN:ND2	1:G:113:ARG:HE	1.84	0.76
1:A:41:LEU:HB3	1:A:71:VAL:CG1	2.15	0.75
1:C:133:LEU:HD22	1:C:170:TRP:HB2	1.69	0.75
1:F:44:THR:HG21	1:F:47:THR:HG23	1.66	0.75
1:E:155:VAL:HG12	1:E:155:VAL:O	1.86	0.75
1:B:81:ALA:HB1	1:B:85:PHE:CZ	2.21	0.75
1:G:47:THR:N	1:G:74:ARG:HH11	1.85	0.75
1:G:21:LYS:CE	1:G:119:ARG:HH21	2.00	0.75
1:C:111:TRP:CZ2	1:C:174:LYS:HG3	2.22	0.75
1:D:38:ILE:HG22	1:D:39:GLU:N	2.03	0.74
1:C:152:GLU:HG3	1:C:152:GLU:O	1.87	0.74
1:G:25:THR:HA	1:G:48:THR:OG1	1.86	0.74
1:C:59:ARG:HH22	1:C:79:GLN:HE21	1.34	0.74
1:F:44:THR:HG23	1:F:47:THR:HG23	1.68	0.74
1:C:57:GLU:O	1:C:61:ARG:HG3	1.88	0.73
1:A:153:HIS:ND1	1:A:162:LYS:HD2	2.03	0.73
1:A:41:LEU:CD1	1:A:71:VAL:CG1	2.67	0.73
1:H:11:ARG:HG3	1:H:36:ILE:HD13	1.71	0.73
1:B:53:GLU:HB3	1:B:59:ARG:HG2	1.69	0.73

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:22:PRO:HA	1:C:47:THR:HG23	1.71	0.72
1:C:41:LEU:HB3	1:C:71:VAL:CG1	2.13	0.72
1:H:82:PRO:HG3	1:H:106:VAL:HA	1.70	0.72
1:A:95:ILE:CD1	1:A:114[B]:LEU:HD11	2.19	0.72
1:C:82:PRO:HG3	1:C:106:VAL:HG22	1.71	0.72
1:E:6:THR:HG23	1:E:151:HIS:CD2	2.24	0.72
1:D:22:PRO:HB2	1:D:23:HIS:HD2	1.53	0.72
1:A:38:ILE:HG21	1:A:69:LEU:HD12	1.72	0.72
1:G:91:ASN:HD22	1:G:91:ASN:N	1.87	0.72
1:F:44:THR:CG2	1:F:47:THR:HG21	2.19	0.71
1:G:21:LYS:HE3	1:G:119:ARG:HH21	1.55	0.71
1:G:53:GLU:HB2	1:G:62:ILE:HD11	1.73	0.71
1:F:146:SER:OG	1:H:161:MSE:HE2	1.89	0.71
1:A:6:THR:HG23	1:A:151:HIS:NE2	2.05	0.71
1:D:23:HIS:HB3	1:G:155:VAL:HG13	1.72	0.71
1:B:11:ARG:O	1:B:15:ILE:HG12	1.91	0.71
1:C:34:GLY:HA3	1:C:62:ILE:HG23	1.71	0.71
1:F:44:THR:HG22	1:F:47:THR:HG21	1.72	0.71
1:H:9:HIS:NE2	1:H:151:HIS:CE1	2.58	0.71
1:E:61:ARG:O	1:E:64:SER:N	2.23	0.71
1:D:26:LEU:CD1	1:D:94:VAL:HG12	2.20	0.70
1:H:95:ILE:O	1:H:120:LEU:HA	1.91	0.70
1:F:5:LEU:HD13	1:F:9:HIS:CE1	2.26	0.70
1:G:69:LEU:O	1:G:71:VAL:HG23	1.91	0.70
1:E:116:VAL:HG22	1:E:175:ALA:HB3	1.71	0.70
1:E:134:TRP:CZ3	1:E:137:ARG:HD3	2.27	0.70
1:H:91:ASN:HD21	1:H:113:ARG:NE	1.90	0.70
1:G:90:ASP:O	1:G:113:ARG:NH2	2.25	0.70
1:A:111:TRP:HA	1:A:114[B]:LEU:HD13	1.74	0.70
1:H:150:SER:CB	1:H:162:LYS:O	2.40	0.69
1:H:56:GLU:HG3	1:H:57:GLU:N	2.07	0.69
1:C:129:SER:O	1:C:132:MSE:HB2	1.92	0.69
1:D:11:ARG:HD3	1:D:35:SER:HB3	1.73	0.69
1:C:124:ALA:HB1	1:C:129:SER:HB3	1.75	0.69
1:D:36:ILE:HG13	1:D:96:PHE:CE2	2.27	0.69
1:E:81:ALA:HB1	1:E:82:PRO:HA	1.73	0.69
1:H:116:VAL:HG13	1:H:175:ALA:HA	1.75	0.69
1:F:23:HIS:N	1:F:46:GLN:O	2.21	0.69
1:B:82:PRO:HG3	1:B:106:VAL:HA	1.75	0.69
1:H:150:SER:HB3	1:H:163:PRO:HA	1.73	0.69
1:B:124:ALA:HB3	1:B:168:HIS:HB2	1.74	0.68
1:C:4:GLN:HA	1:C:7:LYS:CD	2.23	0.68

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:F:61:ARG:O	1:F:64:SER:HB3	1.93	0.68
1:F:102:THR:HA	1:F:132:MSE:SE	2.43	0.68
1:G:21:LYS:HE3	1:G:119:ARG:NH2	2.07	0.68
1:D:46:GLN:HA	1:D:74:ARG:NH1	2.09	0.68
1:D:22:PRO:HB2	1:D:23:HIS:CD2	2.29	0.68
1:H:82:PRO:HD3	1:H:106:VAL:HG22	1.76	0.68
1:F:134:TRP:HZ3	1:F:137:ARG:HH11	1.42	0.68
1:G:90:ASP:C	1:G:91:ASN:HD22	1.97	0.67
1:H:96:PHE:CE2	1:H:98:GLY:HA2	2.29	0.67
1:G:26:LEU:HG	1:G:27:TRP:N	2.08	0.67
1:G:96:PHE:HZ	1:G:123:ASN:ND2	1.93	0.67
1:G:83:ARG:HD2	1:G:83:ARG:O	1.93	0.67
1:H:100:GLY:HA2	5:H:309:HOH:O	1.93	0.67
1:H:150:SER:HB2	1:H:162:LYS:O	1.94	0.67
1:E:103:ALA:HB3	1:E:106:VAL:HG21	1.75	0.67
1:A:126:THR:HA	1:D:160:THR:HG22	1.75	0.67
1:C:92:PRO:HD2	1:C:113:ARG:O	1.94	0.67
1:F:21:LYS:O	1:F:24:GLU:HB2	1.95	0.67
1:A:137:ARG:NE	1:C:152:GLU:OE1	2.28	0.67
1:C:151:HIS:NE2	5:C:177:HOH:O	2.29	0.66
1:H:9:HIS:CD2	1:H:151:HIS:CE1	2.83	0.66
1:B:73:ASP:HB3	1:B:74:ARG:HG3	1.76	0.66
1:G:59:ARG:HB3	1:G:77:VAL:HG11	1.78	0.66
1:B:59:ARG:HD2	1:B:77:VAL:HG12	1.78	0.66
1:F:18:LEU:HD22	1:F:94:VAL:CG1	2.26	0.66
1:A:38:ILE:O	1:A:42:ARG:HB2	1.95	0.66
1:A:18:LEU:HD23	1:A:121:VAL:HG23	1.78	0.66
1:F:138:LYS:NZ	5:F:184:HOH:O	2.29	0.66
1:A:126:THR:HG21	1:D:158:PHE:CD2	2.31	0.66
1:A:56:GLU:HG2	5:A:277:HOH:O	1.95	0.65
1:A:134:TRP:NE1	5:A:189:HOH:O	2.25	0.65
1:D:132:MSE:HE3	1:D:136:LEU:HD21	1.78	0.65
1:E:103:ALA:CB	1:E:106:VAL:HG21	2.25	0.65
1:B:142:GLY:HA3	1:B:172:VAL:HB	1.78	0.65
1:D:102:THR:HG23	5:D:186:HOH:O	1.97	0.65
1:H:91:ASN:ND2	1:H:113:ARG:HE	1.93	0.65
1:B:128:GLU:HG3	1:C:158:PHE:CD1	2.31	0.65
1:C:88:VAL:O	1:C:113:ARG:NH2	2.30	0.65
1:C:27:TRP:CD2	1:C:92:PRO:CG	2.79	0.65
1:H:46:GLN:HA	1:H:74:ARG:NE	2.12	0.65
1:A:27:TRP:HB2	1:A:95:ILE:HG12	1.77	0.65
1:A:125:VAL:HG12	5:A:280:HOH:O	1.96	0.65

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:6:THR:O	1:C:10:VAL:HG23	1.96	0.65
1:F:90[A]:ASP:N	1:F:90[A]:ASP:OD1	2.28	0.65
1:A:94:VAL:HG13	1:A:119:ARG:HB2	1.79	0.65
1:E:17:ALA:HB2	1:E:169:GLN:NE2	2.10	0.65
1:B:19:ALA:O	1:B:119:ARG:NH2	2.28	0.65
1:E:128:GLU:HB2	5:E:298:HOH:O	1.97	0.65
1:G:133:LEU:HD22	1:G:170:TRP:CB	2.27	0.64
1:B:6:THR:CG2	1:B:151:HIS:NE2	2.59	0.64
1:D:71:VAL:HG12	1:D:74:ARG:HB2	1.79	0.64
1:C:114:LEU:HD12	1:C:115:PRO:HD2	1.78	0.64
1:B:27:TRP:HB2	1:B:95:ILE:HG12	1.79	0.64
1:F:147:PHE:CD1	1:H:149:ILE:HG12	2.27	0.64
1:A:21:LYS:HB3	1:A:22:PRO:CD	2.27	0.64
1:D:91:ASN:HD21	1:D:113:ARG:HE	1.46	0.64
1:E:103:ALA:HB3	1:E:106:VAL:CG2	2.27	0.64
1:F:44:THR:HG23	1:F:45:PRO:O	1.98	0.64
1:G:9:HIS:CE1	1:G:151:HIS:CE1	2.86	0.64
1:D:55:SER:O	1:D:59:ARG:HB2	1.97	0.64
1:G:53:GLU:CG	1:G:58:ARG:HB3	2.28	0.64
1:A:109:ALA:O	1:A:112:LYS:HE3	1.98	0.64
1:D:21:LYS:HB2	1:D:24:GLU:OE1	1.97	0.63
1:G:91:ASN:ND2	1:G:113:ARG:HH21	1.96	0.63
1:D:57:GLU:O	1:D:61:ARG:HG3	1.98	0.63
1:D:89:PRO:HB3	5:D:178:HOH:O	1.99	0.63
1:D:26:LEU:HG	1:D:27:TRP:N	2.12	0.63
1:D:6:THR:O	1:D:10:VAL:HG23	1.98	0.63
1:D:95:ILE:HB	1:D:120:LEU:HD13	1.81	0.62
1:H:15:ILE:CD1	1:H:40:TRP:HB2	2.29	0.62
1:B:82:PRO:HB3	1:B:109:ALA:CB	2.29	0.62
1:F:26:LEU:HG	1:F:27:TRP:N	2.14	0.62
1:G:150:SER:HB3	1:G:161:MSE:HE3	1.80	0.62
1:F:128:GLU:OE1	1:G:158:PHE:HD1	1.81	0.62
1:B:97:ILE:HG22	1:B:101:LEU:HB2	1.80	0.62
1:H:120:LEU:HD23	1:H:121:VAL:N	2.14	0.62
1:E:18:LEU:HD22	1:E:94:VAL:CG1	2.29	0.62
1:B:6:THR:HG23	1:B:151:HIS:NE2	2.14	0.62
1:F:21:LYS:HG2	1:F:22:PRO:CD	2.28	0.62
1:A:121:VAL:HA	1:A:170:TRP:O	2.00	0.62
1:B:172:VAL:HG22	1:B:173:VAL:N	2.15	0.62
1:B:27:TRP:HB3	1:B:29:ILE:CD1	2.25	0.62
1:D:46:GLN:HA	1:D:74:ARG:HH12	1.65	0.62
1:G:99:GLY:HA2	5:G:315:HOH:O	1.99	0.62

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:85:PHE:CZ	1:D:110:ALA:HB2	2.34	0.62
1:B:107:PHE:CZ	1:B:140:PHE:CE2	2.88	0.62
1:A:103:ALA:HB3	1:A:106:VAL:CG2	2.30	0.62
1:D:149:ILE:CD1	1:D:149:ILE:N	2.63	0.62
1:F:129:SER:O	1:F:132:MSE:HB3	1.99	0.62
1:G:150:SER:HB3	1:G:161:MSE:CE	2.30	0.62
1:A:3:GLY:N	1:A:5:LEU:HD12	2.15	0.62
1:A:53:GLU:O	1:A:79:GLN:HA	1.99	0.62
1:D:26:LEU:HD12	1:D:94:VAL:O	2.00	0.61
1:E:53:GLU:HG2	1:E:58:ARG:HB2	1.81	0.61
1:B:55:SER:CB	1:B:58:ARG:HG3	2.30	0.61
1:C:131:GLN:HG2	1:D:131:GLN:OE1	1.99	0.61
1:D:21:LYS:O	1:D:47:THR:CG2	2.48	0.61
1:G:20:PRO:HG3	1:G:40:TRP:CZ3	2.35	0.61
1:A:19:ALA:N	1:A:20:PRO:HD3	2.15	0.61
1:A:95:ILE:HD12	1:A:114[B]:LEU:HD21	1.81	0.61
1:G:47:THR:H	1:G:74:ARG:NH1	1.98	0.61
1:D:124:ALA:HB3	1:D:133:LEU:HD22	1.82	0.61
1:B:116:VAL:HG13	1:B:175:ALA:HB2	1.82	0.61
1:A:18:LEU:HD23	1:A:121:VAL:CG2	2.31	0.61
1:G:54:ILE:HD12	1:G:58:ARG:HH22	1.63	0.61
1:G:133:LEU:HD22	1:G:170:TRP:HB3	1.81	0.61
1:F:143:THR:N	1:F:171:THR:O	2.26	0.61
1:C:91:ASN:OD1	1:C:113:ARG:NE	2.20	0.60
1:A:111:TRP:CZ2	1:A:174:LYS:HE3	2.36	0.60
1:A:47:THR:OG1	1:A:74:ARG:NH1	2.33	0.60
1:H:96:PHE:CZ	1:H:98:GLY:HA2	2.36	0.60
1:F:102:THR:HG22	1:F:132:MSE:SE	2.52	0.60
1:F:18:LEU:HD22	1:F:94:VAL:HG11	1.82	0.60
1:B:80:GLY:O	1:B:83:ARG:HG2	2.01	0.60
1:D:67:ILE:HD12	1:F:61:ARG:NE	2.16	0.60
1:G:27:TRP:HB2	1:G:95:ILE:HG12	1.82	0.60
1:H:27:TRP:CD1	1:H:92:PRO:HB3	2.37	0.60
1:B:55:SER:HB3	1:B:58:ARG:HG3	1.84	0.60
1:E:42:ARG:NH2	1:E:69:LEU:HD22	2.16	0.60
1:F:81:ALA:HB1	1:F:82:PRO:HA	1.83	0.60
1:G:59:ARG:HD2	1:G:77:VAL:HG12	1.83	0.60
1:A:101:LEU:HD22	1:A:123:ASN:O	2.02	0.60
1:C:127:VAL:O	1:C:131:GLN:HG3	2.00	0.60
1:E:160:THR:HG22	1:H:126:THR:HG22	1.83	0.60
1:F:41:LEU:CD2	1:F:74:ARG:HB3	2.32	0.60
1:F:134:TRP:CZ3	1:F:137:ARG:HD3	2.37	0.60

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:11:ARG:O	1:C:15:ILE:HG12	2.01	0.60
1:H:126:THR:O	1:H:130:GLU:HB2	2.01	0.59
1:H:44:THR:HG22	1:H:45:PRO:O	2.02	0.59
1:B:127:VAL:O	1:B:131:GLN:HG3	2.01	0.59
1:B:47:THR:OG1	1:B:74:ARG:NH1	2.35	0.59
1:F:88:VAL:O	1:F:113:ARG:NH2	2.29	0.59
1:C:27:TRP:CE2	1:C:92:PRO:CG	2.73	0.59
1:E:106:VAL:HG12	1:E:107:PHE:N	2.18	0.59
1:D:16:SER:O	1:D:19:ALA:HB2	2.03	0.59
1:A:25:THR:HA	1:A:48:THR:O	2.01	0.59
1:E:63:LEU:O	1:E:67:ILE:HG13	2.03	0.59
1:B:67:ILE:HA	1:B:72:SER:HB3	1.84	0.59
1:C:60:GLU:O	1:C:64:SER:HB2	2.01	0.59
1:H:15:ILE:HD12	1:H:40:TRP:HB2	1.85	0.59
1:B:123:ASN:HA	1:B:168:HIS:O	2.02	0.59
1:H:9:HIS:CD2	1:H:151:HIS:HE1	2.20	0.58
1:G:38:ILE:HD13	1:G:65:ASN:HD22	1.68	0.58
1:F:113:ARG:O	1:F:115:PRO:HD3	2.02	0.58
1:F:128:GLU:OE1	1:G:158:PHE:HA	2.03	0.58
1:G:45:PRO:C	1:G:46:GLN:HG3	2.23	0.58
1:A:103:ALA:HB3	1:A:106:VAL:HG23	1.85	0.58
1:D:91:ASN:ND2	1:D:113:ARG:HE	2.01	0.58
1:D:24:GLU:N	1:D:47:THR:HG22	2.14	0.58
1:H:90:ASP:O	1:H:113:ARG:NH2	2.36	0.58
1:A:128:GLU:OE2	1:D:159:ILE:HG12	2.04	0.58
1:E:59:ARG:HG2	1:E:77:VAL:HG12	1.85	0.58
1:E:13:LEU:HD11	1:G:13:LEU:HD21	1.85	0.58
1:H:11:ARG:HA	1:H:36:ILE:HD11	1.85	0.58
1:E:56:GLU:HB2	1:E:59:ARG:NH2	2.19	0.58
1:C:54:ILE:CD1	1:C:80:GLY:HA3	2.18	0.58
1:G:83:ARG:NH1	5:G:313:HOH:O	2.36	0.58
1:A:148:ALA:HA	5:A:199:HOH:O	2.04	0.58
1:D:6:THR:HA	1:D:9:HIS:CD2	2.35	0.58
1:A:169:GLN:NE2	5:A:273:HOH:O	2.36	0.58
1:A:67:ILE:HG12	1:A:72:SER:HB3	1.86	0.58
1:D:21:LYS:O	1:D:47:THR:HG21	2.03	0.58
1:D:9:HIS:CD2	1:D:151:HIS:HE1	2.22	0.57
1:B:134:TRP:O	1:B:138:LYS:HG3	2.03	0.57
1:H:94:VAL:HG22	1:H:119:ARG:HB2	1.86	0.57
1:A:58:ARG:O	1:A:62:ILE:HG13	2.04	0.57
1:F:45:PRO:C	1:F:46:GLN:HG3	2.24	0.57
1:F:7:LYS:CD	1:F:11:ARG:HE	2.10	0.57

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:H:9:HIS:O	1:H:12:ALA:HB3	2.05	0.57
1:A:133:LEU:O	1:A:137:ARG:HB2	2.04	0.57
1:A:127:VAL:O	1:A:131:GLN:HG3	2.04	0.57
1:B:21:LYS:HB3	1:B:22:PRO:HD2	1.85	0.57
1:D:154:THR:HG23	1:D:159:ILE:HD13	1.86	0.57
1:D:88:VAL:O	1:D:88:VAL:HG12	2.05	0.57
1:H:15:ILE:HD13	1:H:18:LEU:HD13	1.86	0.57
1:C:11:ARG:HG3	1:C:36:ILE:CD1	2.35	0.57
1:E:59:ARG:HG2	1:E:77:VAL:CG1	2.35	0.57
1:B:97:ILE:CG2	1:B:101:LEU:HB2	2.34	0.56
1:D:72:SER:O	1:F:58:ARG:NH2	2.37	0.56
1:H:117:GLY:N	1:H:174:LYS:O	2.33	0.56
1:C:155:VAL:HG12	1:C:155:VAL:O	2.05	0.56
1:E:44:THR:HG22	1:E:45:PRO:O	2.04	0.56
1:H:98:GLY:O	1:H:123:ASN:HB2	2.04	0.56
1:G:150:SER:CB	1:G:161:MSE:HE2	2.35	0.56
1:E:34:GLY:HA2	1:E:51:CYS:SG	2.45	0.56
1:A:149:ILE:HG12	1:C:147:PHE:HD1	1.69	0.56
1:C:88:VAL:O	1:C:113:ARG:NH1	2.36	0.56
1:A:38:ILE:CG2	1:A:69:LEU:HD12	2.34	0.56
1:H:40:TRP:HA	1:H:43:SER:HB3	1.87	0.56
1:A:111:TRP:CE2	1:A:174:LYS:HE3	2.39	0.56
1:A:137:ARG:CZ	1:C:152:GLU:OE1	2.53	0.56
1:D:38:ILE:CG2	1:D:39:GLU:N	2.69	0.56
1:C:101:LEU:HD22	1:C:136:LEU:HD11	1.88	0.56
1:A:28:ASP:C	1:A:28:ASP:OD1	2.43	0.56
1:D:82:PRO:O	1:D:85:PHE:HB2	2.06	0.56
1:B:128:GLU:HG3	1:C:158:PHE:HD1	1.69	0.56
1:B:107:PHE:CE2	1:B:140:PHE:HE2	2.23	0.56
1:B:107:PHE:CZ	1:B:140:PHE:HE2	2.23	0.56
1:A:109:ALA:O	1:A:110:ALA:C	2.43	0.56
1:B:64:SER:HA	1:B:67:ILE:HD12	1.87	0.56
1:D:5:LEU:HG	1:D:9:HIS:NE2	2.21	0.56
1:F:127:VAL:O	1:F:131:GLN:HB2	2.06	0.56
1:G:53:GLU:CB	1:G:62:ILE:HD11	2.36	0.56
1:G:82:PRO:HG3	1:G:106:VAL:HA	1.86	0.56
1:C:6:THR:HG1	1:C:151:HIS:CD2	2.21	0.55
1:C:21:LYS:HB3	1:C:22:PRO:HD2	1.87	0.55
1:F:58:ARG:NH1	5:F:183:HOH:O	2.38	0.55
1:H:129:SER:O	1:H:132:MSE:HB3	2.06	0.55
1:H:44:THR:HB	1:H:47:THR:OG1	2.05	0.55
1:B:95:ILE:HB	1:B:120:LEU:HD12	1.89	0.55

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:25:THR:H	1:A:93:ASP:HB2	1.71	0.55
1:D:18:LEU:HD23	1:D:121:VAL:HG23	1.87	0.55
1:E:103:ALA:CB	1:E:106:VAL:CG2	2.84	0.55
1:G:55:SER:OG	1:G:58:ARG:HG3	2.06	0.55
1:F:110:ALA:O	1:F:113:ARG:N	2.38	0.55
1:G:9:HIS:HE1	1:G:151:HIS:CE1	2.24	0.55
1:B:82:PRO:HB3	1:B:109:ALA:HB1	1.89	0.55
1:B:126:THR:HG21	1:C:158:PHE:CG	2.42	0.55
1:D:5:LEU:O	1:D:8:GLN:N	2.40	0.54
1:F:125:VAL:O	1:F:125:VAL:HG22	2.07	0.54
1:C:34:GLY:HA3	1:C:62:ILE:HG12	1.88	0.54
1:D:101:LEU:CD1	1:D:132:MSE:HG2	2.31	0.54
1:C:34:GLY:CA	1:C:62:ILE:HG12	2.37	0.54
1:B:132:MSE:HE3	1:B:136:LEU:HD11	1.88	0.54
1:D:171:THR:HG21	5:D:179:HOH:O	2.07	0.54
1:B:5:LEU:HB3	1:B:151:HIS:CE1	2.43	0.54
1:E:129:SER:O	1:E:132:MSE:HB3	2.08	0.54
1:E:21:LYS:HB2	1:E:24:GLU:CD	2.28	0.54
1:D:51:CYS:O	1:D:78:GLN:HB2	2.07	0.54
1:E:154:THR:HG22	1:E:156:GLY:N	2.06	0.54
1:A:21:LYS:O	1:A:24:GLU:HB2	2.07	0.54
1:B:174:LYS:O	1:B:175:ALA:CB	2.56	0.54
1:F:101:LEU:N	1:F:101:LEU:HD12	2.23	0.54
1:B:172:VAL:HG22	1:B:173:VAL:O	2.08	0.54
1:H:55:SER:HB2	1:H:58[B]:ARG:CB	2.12	0.54
1:C:59:ARG:NH2	1:C:79:GLN:NE2	2.53	0.54
1:D:149:ILE:HD12	1:D:149:ILE:N	2.23	0.54
1:E:134:TRP:HZ3	1:E:137:ARG:NH1	2.06	0.54
1:G:11:ARG:HD2	1:G:39:GLU:HB2	1.90	0.53
1:G:53:GLU:HG2	1:G:58:ARG:HB3	1.88	0.53
1:B:101:LEU:CD2	1:B:123:ASN:O	2.56	0.53
1:B:67:ILE:HA	1:B:72:SER:CB	2.38	0.53
1:A:153:HIS:CD2	1:A:155:VAL:CG2	2.80	0.53
1:B:81:ALA:CB	1:B:85:PHE:CZ	2.92	0.53
1:B:142:GLY:CA	1:B:172:VAL:HB	2.38	0.53
1:B:150:SER:HB3	1:B:161:MSE:HE3	1.91	0.53
1:E:111:TRP:CE3	1:E:120:LEU:HD13	2.43	0.53
1:G:29:ILE:HD12	1:G:95:ILE:HG23	1.90	0.53
1:G:47:THR:O	1:G:74:ARG:HD2	2.08	0.53
1:F:148:ALA:N	1:H:148:ALA:O	2.36	0.53
1:B:27:TRP:CB	1:B:29:ILE:HD11	2.26	0.53
1:C:58:ARG:O	1:C:61:ARG:HB2	2.08	0.53

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:63:LEU:HG	1:E:63:LEU:O	2.08	0.53
1:H:132:MSE:O	1:H:135:ALA:HB3	2.08	0.53
1:G:38:ILE:HG21	1:G:69:LEU:HD12	1.91	0.53
1:A:21:LYS:NZ	5:A:186:HOH:O	2.38	0.53
1:H:81:ALA:HB1	1:H:82:PRO:HA	1.90	0.53
1:A:22:PRO:HB2	1:A:23:HIS:CD2	2.44	0.53
1:B:29:ILE:HD13	1:B:95:ILE:HG23	1.91	0.53
1:F:11:ARG:HD3	1:F:35:SER:OG	2.09	0.53
1:D:23:HIS:HA	1:D:46:GLN:O	2.09	0.53
1:C:82:PRO:O	1:C:85:PHE:CD2	2.62	0.53
1:B:6:THR:HG22	1:B:151:HIS:NE2	2.23	0.53
1:B:136:LEU:O	1:B:140:PHE:HB2	2.09	0.53
1:B:122:ALA:O	1:B:169:GLN:HA	2.09	0.53
1:B:16:SER:HB2	1:D:12:ALA:HB2	1.90	0.53
1:H:133:LEU:HD22	1:H:170:TRP:CB	2.39	0.53
1:G:131:GLN:NE2	5:G:311:HOH:O	2.42	0.53
1:B:16:SER:HB2	1:D:12:ALA:CB	2.39	0.52
1:H:96:PHE:HA	1:H:121:VAL:O	2.09	0.52
1:C:24:GLU:HG2	1:C:93:ASP:CB	2.39	0.52
1:F:50:VAL:CG1	1:F:78:GLN:HG3	2.39	0.52
1:F:7:LYS:O	1:F:11:ARG:HG3	2.09	0.52
1:G:53:GLU:HB3	1:G:59:ARG:HG2	1.90	0.52
1:D:149:ILE:HG22	1:D:151:HIS:NE2	2.24	0.52
1:H:150:SER:HB3	1:H:162:LYS:O	2.09	0.52
1:D:84:ALA:O	1:D:87:ASP:HB2	2.09	0.52
1:C:26:LEU:HB2	1:C:40:TRP:CE2	2.45	0.52
1:D:138:LYS:HB2	5:D:192:HOH:O	2.09	0.52
1:G:59:ARG:HA	1:G:62:ILE:HD12	1.92	0.52
1:D:18:LEU:HD23	1:D:121:VAL:CG2	2.38	0.52
1:A:55:SER:HB3	1:A:58:ARG:HB2	1.91	0.52
1:D:83:ARG:HB3	5:D:182:HOH:O	2.09	0.52
1:D:4:GLN:HA	1:D:7:LYS:HD2	1.91	0.52
1:F:11:ARG:HB3	1:F:39:GLU:OE1	2.09	0.52
1:B:121:VAL:HA	1:B:170:TRP:O	2.10	0.52
1:A:101:LEU:O	1:A:101:LEU:HD12	2.09	0.52
1:A:137:ARG:HH11	1:A:138:LYS:HE2	1.74	0.52
1:C:133:LEU:HD22	1:C:170:TRP:CB	2.37	0.52
1:E:11:ARG:HD2	1:E:39:GLU:OE1	2.10	0.52
1:F:6:THR:HG23	1:F:151:HIS:CE1	2.45	0.52
1:H:44:THR:HG21	1:H:47:THR:HG23	1.92	0.52
1:B:54:ILE:HG23	1:B:55:SER:H	1.75	0.52
1:D:120:LEU:HD12	1:D:121:VAL:N	2.25	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:64:SER:O	1:C:67:ILE:HB	2.10	0.52
1:A:95:ILE:HG22	1:A:96:PHE:N	2.24	0.51
1:A:149:ILE:HG12	1:C:147:PHE:CD1	2.45	0.51
1:E:114:LEU:HD23	1:E:174:LYS:HB2	1.91	0.51
1:G:150:SER:CB	1:G:161:MSE:CE	2.89	0.51
1:B:174:LYS:O	1:B:175:ALA:HB2	2.09	0.51
1:A:59:ARG:CZ	1:A:79:GLN:HG2	2.40	0.51
1:C:112:LYS:HG3	5:C:197:HOH:O	2.11	0.51
1:H:39:GLU:OE2	1:H:42:ARG:NH1	2.43	0.51
1:H:56:GLU:O	1:H:60:GLU:HG3	2.11	0.51
1:E:63:LEU:O	1:E:67:ILE:CG1	2.59	0.51
1:F:17:ALA:HB2	1:F:169:GLN:HE22	1.75	0.51
1:F:19:ALA:N	1:F:20:PRO:HD3	2.26	0.51
1:B:87:ASP:O	1:B:89:PRO:HD3	2.11	0.51
1:E:134:TRP:CZ2	1:F:127:VAL:HG21	2.45	0.51
1:E:107:PHE:CE2	1:E:136:LEU:HD22	2.46	0.51
1:H:124:ALA:HB1	1:H:129:SER:HB2	1.92	0.51
1:F:130:GLU:HG3	1:F:168:HIS:CD2	2.45	0.51
1:D:111:TRP:NE1	1:D:174:LYS:HE3	2.25	0.51
1:B:101:LEU:HD22	1:B:123:ASN:O	2.09	0.51
1:A:22:PRO:HD2	5:A:198:HOH:O	2.10	0.51
1:D:5:LEU:O	1:D:8:GLN:HB3	2.11	0.51
1:E:7:LYS:O	1:E:7:LYS:HG2	2.10	0.50
1:H:15:ILE:HD13	1:H:18:LEU:CD1	2.40	0.50
1:B:52:PHE:HD2	1:B:79:GLN:O	1.94	0.50
1:B:53:GLU:O	1:B:79:GLN:HA	2.11	0.50
1:F:27:TRP:CZ2	1:F:88:VAL:HG11	2.46	0.50
1:B:21:LYS:HB3	1:B:22:PRO:CD	2.41	0.50
1:A:81:ALA:HB1	1:A:85:PHE:CE2	2.41	0.50
1:H:11:ARG:HD3	1:H:35:SER:OG	2.11	0.50
1:A:24:GLU:OE1	1:A:119:ARG:NH2	2.43	0.50
1:B:117:GLY:N	1:B:174:LYS:O	2.44	0.50
1:F:15:ILE:HG22	1:F:15:ILE:O	2.10	0.50
1:C:102:THR:HA	1:C:132:MSE:HE3	1.92	0.50
1:A:91:ASN:ND2	1:A:113:ARG:HE	2.09	0.50
1:F:103:ALA:HB3	1:F:106:VAL:HG21	1.93	0.50
1:H:40:TRP:CD1	1:H:40:TRP:C	2.84	0.50
1:B:107:PHE:HZ	1:B:140:PHE:CE2	2.28	0.50
1:F:153:HIS:CD2	1:F:162:LYS:HD2	2.46	0.50
1:D:17:ALA:O	5:D:179:HOH:O	2.20	0.50
1:E:11:ARG:HD2	1:E:39:GLU:CD	2.31	0.50
1:E:9:HIS:O	1:E:12:ALA:HB3	2.12	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:125:VAL:O	1:D:125:VAL:HG22	2.10	0.50
1:D:81:ALA:HB1	1:D:82:PRO:HA	1.94	0.50
1:A:137:ARG:NH2	1:C:152:GLU:OE1	2.45	0.50
1:C:82:PRO:HG3	1:C:106:VAL:CG2	2.42	0.50
1:B:7:LYS:CE	1:B:11:ARG:HD3	2.41	0.50
1:F:78:GLN:NE2	2:F:176:ACY:O	2.44	0.50
1:D:4:GLN:HA	1:D:7:LYS:CD	2.42	0.50
1:G:150:SER:HB2	1:G:161:MSE:HE2	1.93	0.50
1:A:67:ILE:CG1	1:A:72:SER:HB3	2.42	0.50
1:G:94:VAL:HG12	1:G:95:ILE:H	1.77	0.49
1:A:121:VAL:HG13	1:A:171:THR:HB	1.94	0.49
1:A:69:LEU:O	1:A:71:VAL:HG23	2.12	0.49
1:A:22:PRO:C	1:A:23:HIS:CD2	2.86	0.49
1:G:134:TRP:O	1:G:138:LYS:HG3	2.11	0.49
1:A:24:GLU:CD	1:A:119:ARG:NH2	2.61	0.49
1:A:18:LEU:HD21	1:A:96:PHE:HB2	1.94	0.49
1:C:82:PRO:O	1:C:85:PHE:HD2	1.96	0.49
1:E:103:ALA:HB3	1:E:106:VAL:CB	2.42	0.49
1:H:71:VAL:HG13	1:H:71:VAL:O	2.10	0.49
1:A:109:ALA:HA	1:A:112:LYS:HE3	1.93	0.49
1:H:55:SER:HB2	1:H:58[A]:ARG:CG	2.42	0.49
1:G:19:ALA:N	1:G:20:PRO:HD3	2.28	0.49
1:C:111:TRP:CE2	1:C:174:LYS:HE3	2.48	0.49
1:D:83:ARG:CB	5:D:182:HOH:O	2.60	0.49
1:B:63:LEU:O	1:B:67:ILE:HG13	2.12	0.49
1:C:56:GLU:HG2	1:C:59:ARG:NH1	2.27	0.49
1:A:19:ALA:O	1:A:119:ARG:NH1	2.46	0.49
1:D:11:ARG:HB3	1:D:39:GLU:OE1	2.12	0.49
1:E:56:GLU:O	1:E:60:GLU:HG3	2.13	0.49
1:A:122:ALA:O	1:A:169:GLN:HA	2.12	0.49
1:H:57:GLU:OE1	1:H:60:GLU:OE1	2.31	0.49
1:G:94:VAL:HG13	1:G:119:ARG:HB2	1.95	0.49
1:C:59:ARG:C	1:C:61:ARG:N	2.65	0.49
1:E:18:LEU:HD22	1:E:94:VAL:HG12	1.94	0.49
1:D:65:ASN:O	1:D:69:LEU:HD23	2.12	0.49
1:B:29:ILE:CD1	1:B:95:ILE:HG23	2.43	0.49
1:G:91:ASN:CG	1:G:113:ARG:HE	2.16	0.49
1:G:134:TRP:CZ2	1:H:127:VAL:HG21	2.48	0.49
1:H:96:PHE:CE2	1:H:98:GLY:CA	2.95	0.48
1:F:5:LEU:CD1	1:F:9:HIS:CE1	2.94	0.48
1:F:116:VAL:HG13	1:F:175:ALA:HA	1.95	0.48
1:A:52:PHE:HD2	1:A:84:ALA:HB3	1.77	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:120:LEU:C	1:A:120:LEU:HD23	2.33	0.48
1:B:11:ARG:CZ	1:B:11:ARG:HB2	2.42	0.48
1:G:81:ALA:HB1	1:G:85:PHE:CZ	2.48	0.48
1:A:114[B]:LEU:N	1:A:114[B]:LEU:HD12	2.29	0.48
1:A:133:LEU:HD22	1:A:170:TRP:HB2	1.94	0.48
1:H:150:SER:HB2	1:H:161:MSE:HE3	1.96	0.48
1:D:152:GLU:HG3	1:D:152:GLU:O	2.11	0.48
1:G:47:THR:H	1:G:74:ARG:HH12	1.61	0.48
1:F:26:LEU:HB2	1:F:40:TRP:CE2	2.49	0.48
1:G:30:GLY:O	1:G:31:GLY:C	2.51	0.48
1:E:85:PHE:CZ	1:E:110:ALA:HB2	2.48	0.48
1:D:120:LEU:O	1:D:171:THR:HA	2.14	0.48
1:B:103:ALA:HB3	1:B:106:VAL:CG2	2.44	0.48
1:G:96:PHE:CE2	1:G:98:GLY:HA2	2.49	0.48
1:B:59:ARG:CZ	1:B:79:GLN:HG2	2.44	0.48
1:H:44:THR:HB	1:H:47:THR:CG2	2.44	0.48
1:E:95:ILE:O	1:E:120:LEU:HA	2.14	0.48
1:C:87:ASP:O	1:C:89:PRO:HD3	2.13	0.48
1:D:23:HIS:N	1:D:47:THR:HG22	2.28	0.48
1:C:6:THR:HA	1:C:151:HIS:CE1	2.49	0.48
1:H:122:ALA:HB3	1:H:133:LEU:HD21	1.96	0.48
1:D:111:TRP:CZ2	1:D:174:LYS:HG3	2.48	0.48
1:F:59:ARG:HG2	1:F:77:VAL:HG12	1.95	0.48
1:G:95:ILE:O	1:G:120:LEU:HA	2.14	0.48
1:D:27:TRP:CD1	1:D:92:PRO:HB3	2.49	0.48
1:E:103:ALA:HB3	1:E:106:VAL:HB	1.95	0.48
1:C:11:ARG:HG3	1:C:36:ILE:HD12	1.96	0.48
1:F:65:ASN:O	1:F:68:ASN:HB2	2.14	0.48
1:F:102:THR:CG2	1:F:132:MSE:SE	3.12	0.48
1:G:147:PHE:HE2	1:G:169:GLN:HB2	1.77	0.48
1:C:41:LEU:HD21	1:C:49:ALA:N	2.29	0.47
1:A:125:VAL:HG13	1:A:126:THR:HG23	1.95	0.47
1:B:172:VAL:CG2	1:B:173:VAL:N	2.78	0.47
1:H:133:LEU:HD22	1:H:170:TRP:HB3	1.94	0.47
1:F:59:ARG:HG2	1:F:77:VAL:CG1	2.44	0.47
1:A:11[A]:ARG:O	1:A:15:ILE:HG12	2.14	0.47
1:H:96:PHE:HD1	1:H:121:VAL:HG12	1.79	0.47
1:A:95:ILE:CG2	1:A:96:PHE:N	2.76	0.47
1:E:120:LEU:HB3	1:E:172:VAL:HG12	1.95	0.47
1:G:20:PRO:CG	1:G:40:TRP:CE3	2.85	0.47
1:H:116:VAL:HA	1:H:174:LYS:O	2.15	0.47
1:B:34:GLY:HA2	1:B:51:CYS:SG	2.55	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:27:TRP:CZ2	1:C:92:PRO:HG3	2.43	0.47
1:G:59:ARG:CB	1:G:77:VAL:HG11	2.43	0.47
1:D:97:ILE:HG21	1:D:101:LEU:HB3	1.97	0.47
1:H:11:ARG:CG	1:H:39:GLU:HG3	2.44	0.47
1:H:113:ARG:HA	1:H:113:ARG:HD2	1.55	0.47
1:C:22:PRO:O	1:C:46:GLN:O	2.33	0.47
1:B:7:LYS:HE2	1:B:11:ARG:HD3	1.96	0.47
1:E:21:LYS:HB2	1:E:24:GLU:OE1	2.14	0.47
1:H:82:PRO:CD	1:H:106:VAL:HG22	2.44	0.47
1:B:132:MSE:HE3	1:B:136:LEU:HD21	1.97	0.47
1:A:81:ALA:HB2	1:A:85:PHE:HE1	1.61	0.47
1:C:111:TRP:CH2	1:C:174:LYS:HG3	2.50	0.47
1:F:103:ALA:O	1:F:106:VAL:HB	2.15	0.47
1:H:93:ASP:OD1	1:H:93:ASP:N	2.48	0.47
1:H:19:ALA:N	1:H:20:PRO:HD3	2.29	0.47
1:D:59:ARG:NH2	1:D:79:GLN:HG2	2.29	0.47
1:B:18:LEU:HD23	1:B:121:VAL:CG2	2.45	0.47
1:D:92:PRO:HD2	1:D:113:ARG:O	2.15	0.46
1:B:94:VAL:HA	1:B:119:ARG:O	2.15	0.46
1:E:100:GLY:O	1:E:102:THR:N	2.48	0.46
1:B:16:SER:OG	1:D:8:GLN:NE2	2.48	0.46
1:C:26:LEU:HD22	1:C:40:TRP:CG	2.50	0.46
1:C:113:ARG:HD2	1:C:113:ARG:HA	1.72	0.46
1:H:27:TRP:CD2	1:H:92:PRO:HG3	2.50	0.46
1:E:111:TRP:CZ3	1:E:172:VAL:HG11	2.50	0.46
1:A:30:GLY:O	1:A:31:GLY:C	2.53	0.46
1:E:41:LEU:HD22	1:E:74:ARG:HB3	1.98	0.46
1:B:137:ARG:CG	1:B:137:ARG:HH11	2.11	0.46
1:A:27:TRP:CD1	1:A:50:VAL:HB	2.50	0.46
1:E:101:LEU:HG	1:E:132:MSE:HG2	1.96	0.46
1:C:24:GLU:HG2	1:C:93:ASP:HB2	1.96	0.46
1:A:137:ARG:NH2	1:C:152:GLU:OE2	2.47	0.46
1:D:34:GLY:O	1:D:38:ILE:HB	2.16	0.46
1:H:54:ILE:HD12	1:H:55:SER:H	1.81	0.46
1:H:56:GLU:CG	1:H:57:GLU:N	2.74	0.46
1:D:132:MSE:O	1:D:136:LEU:HG	2.16	0.46
1:H:18:LEU:O	1:H:119:ARG:NH1	2.48	0.46
1:A:27:TRP:CZ2	1:A:92:PRO:HD3	2.51	0.46
1:A:125:VAL:HG22	1:A:125:VAL:O	2.16	0.46
1:A:103:ALA:HB3	1:A:106:VAL:HG21	1.98	0.46
1:D:29:ILE:HD11	1:D:85:PHE:HZ	1.81	0.46
1:G:96:PHE:CE1	1:G:121:VAL:HG12	2.51	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:H:15:ILE:HD11	1:H:40:TRP:HB2	1.97	0.46
1:G:34:GLY:O	1:G:38:ILE:HD12	2.16	0.46
1:G:82:PRO:HB3	1:G:109:ALA:HB3	1.97	0.46
1:B:38:ILE:O	1:B:42:ARG:HG3	2.16	0.46
1:B:114:LEU:HD12	1:B:115:PRO:HD2	1.98	0.46
1:G:25:THR:CA	1:G:48:THR:OG1	2.62	0.46
1:C:124:ALA:CB	1:C:129:SER:HB3	2.45	0.46
1:C:142:GLY:N	1:C:172:VAL:HG23	2.31	0.46
1:C:18:LEU:HD21	1:C:121:VAL:HB	1.98	0.46
1:G:133:LEU:HD22	1:G:170:TRP:HB2	1.98	0.45
1:G:39:GLU:OE2	1:G:42:ARG:HD2	2.16	0.45
1:A:98:GLY:HA2	1:A:123:ASN:HB2	1.97	0.45
1:E:102:THR:HG22	1:E:132:MSE:SE	2.66	0.45
1:E:111:TRP:HE3	1:E:120:LEU:HD13	1.82	0.45
1:A:26:LEU:HB2	1:A:40:TRP:CE2	2.51	0.45
1:B:107:PHE:CE2	1:B:140:PHE:CE2	3.04	0.45
1:B:171:THR:HG23	5:B:284:HOH:O	2.16	0.45
1:A:153:HIS:HD2	1:A:155:VAL:CG2	2.21	0.45
1:C:123:ASN:OD1	1:C:169:GLN:HG3	2.16	0.45
1:D:41:LEU:HD21	1:D:48:THR:HA	1.97	0.45
1:H:54:ILE:H	1:H:54:ILE:HG13	1.34	0.45
1:G:137:ARG:C	1:G:137:ARG:HD3	2.37	0.45
1:G:96:PHE:HA	1:G:121:VAL:O	2.15	0.45
1:H:86:ASP:OD1	1:H:86:ASP:O	2.34	0.45
1:B:98:GLY:O	1:B:101:LEU:HB3	2.16	0.45
1:D:58:ARG:HA	1:D:61:ARG:HB2	1.99	0.45
1:B:88:VAL:HA	1:B:89:PRO:HD3	1.70	0.45
1:D:142:GLY:N	1:D:172:VAL:HG23	2.32	0.45
1:G:40:TRP:HH2	1:G:94:VAL:HG21	1.81	0.45
1:G:94:VAL:HG12	1:G:95:ILE:N	2.31	0.45
1:A:88:VAL:HA	1:A:89:PRO:HD3	1.63	0.45
1:E:137:ARG:O	1:E:138:LYS:C	2.55	0.45
1:D:114:LEU:HA	1:D:115:PRO:HD2	1.71	0.45
1:E:127:VAL:HG12	1:E:127:VAL:O	2.15	0.45
1:G:103:ALA:HB3	1:G:106:VAL:HG21	1.98	0.45
1:C:81:ALA:HB1	1:C:82:PRO:HA	1.98	0.45
1:F:101:LEU:CD1	1:F:132:MSE:HG2	2.47	0.45
1:D:58:ARG:O	1:D:61:ARG:HB2	2.17	0.45
1:G:114:LEU:HA	1:G:115:PRO:HD3	1.77	0.45
1:C:103:ALA:HA	1:C:104:PRO:HD3	1.71	0.45
1:E:159:ILE:HD12	1:H:128:GLU:OE2	2.17	0.45
1:H:57:GLU:HA	1:H:60:GLU:CD	2.37	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:154:THR:CG2	1:E:156:GLY:H	2.10	0.45
1:F:41:LEU:CD2	1:F:74:ARG:CB	2.95	0.45
1:G:9:HIS:O	1:G:12:ALA:HB3	2.17	0.45
1:A:41:LEU:HD12	1:A:71:VAL:CG1	2.45	0.45
1:B:11:ARG:O	1:B:14:ALA:HB3	2.17	0.45
1:H:116:VAL:CG1	1:H:175:ALA:HA	2.45	0.45
1:H:82:PRO:HG3	1:H:106:VAL:HG22	1.99	0.44
1:H:114:LEU:HD12	1:H:115:PRO:HD2	1.98	0.44
1:D:67:ILE:HD12	1:F:61:ARG:CZ	2.46	0.44
1:E:81:ALA:CB	1:E:82:PRO:HA	2.37	0.44
1:E:127:VAL:HG11	1:F:134:TRP:CE2	2.52	0.44
1:F:15:ILE:HA	1:F:18:LEU:HB2	1.98	0.44
1:E:53:GLU:HG3	1:E:55:SER:H	1.82	0.44
1:E:42:ARG:HH22	1:E:69:LEU:HD22	1.81	0.44
1:D:71:VAL:HG12	1:D:74:ARG:HD2	1.98	0.44
1:F:124:ALA:HB1	1:F:129:SER:HB3	1.99	0.44
1:H:41:LEU:O	1:H:74:ARG:NH1	2.49	0.44
1:G:37:ALA:O	1:G:41:LEU:HG	2.18	0.44
1:D:165:LEU:HA	1:D:166:PRO:HD3	1.83	0.44
1:A:41:LEU:CD1	1:A:71:VAL:HG11	2.47	0.44
1:A:6:THR:O	1:A:10:VAL:HG23	2.17	0.44
1:C:6:THR:HA	5:C:177:HOH:O	2.18	0.44
1:F:12:ALA:CB	1:H:16:SER:HB2	2.48	0.44
1:A:82:PRO:HG3	1:A:106:VAL:HA	1.99	0.44
1:C:91:ASN:HA	1:C:92:PRO:HD3	1.72	0.44
1:H:41:LEU:HD21	1:H:49:ALA:HB2	2.00	0.44
1:C:125:VAL:HG23	1:C:166:PRO:O	2.17	0.44
1:D:66:ALA:CB	1:D:75:ILE:HG21	2.47	0.44
1:E:11:ARG:HG3	1:E:39:GLU:HG2	1.99	0.44
1:E:155:VAL:CG1	1:E:155:VAL:O	2.58	0.44
1:E:56:GLU:HA	1:E:59:ARG:HB2	1.99	0.44
1:C:88:VAL:HA	5:C:192:HOH:O	2.16	0.44
1:G:21:LYS:N	1:G:24:GLU:OE1	2.51	0.44
1:H:38:ILE:O	1:H:42:ARG:CD	2.66	0.44
1:E:158:PHE:HD2	1:H:126:THR:HG21	1.76	0.44
1:C:21:LYS:O	1:C:47:THR:CG2	2.66	0.44
1:F:26:LEU:O	1:F:27:TRP:HD1	2.01	0.44
1:E:100:GLY:O	1:E:101:LEU:C	2.57	0.44
1:G:139:GLN:O	1:G:139:GLN:HG2	2.17	0.44
1:H:36:ILE:HD13	1:H:36:ILE:HA	1.72	0.43
1:E:125:VAL:HG22	1:E:125:VAL:O	2.18	0.43
1:E:59:ARG:NH1	1:E:79:GLN:HA	2.33	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:27:TRP:CD1	1:E:92:PRO:HB3	2.53	0.43
1:C:66:ALA:HB1	1:C:72:SER:HA	1.99	0.43
1:G:121:VAL:HA	1:G:170:TRP:O	2.18	0.43
1:A:5:LEU:O	1:A:9:HIS:CD2	2.71	0.43
1:B:58:ARG:O	1:B:62:ILE:HG13	2.18	0.43
1:F:150:SER:HA	1:F:162:LYS:O	2.18	0.43
1:D:54:ILE:H	1:D:54:ILE:HG13	1.49	0.43
1:A:81:ALA:HA	1:A:82:PRO:C	2.39	0.43
1:D:55:SER:OG	1:D:58:ARG:HG3	2.18	0.43
1:F:34:GLY:HA2	1:F:51:CYS:SG	2.59	0.43
1:C:51:CYS:O	1:C:78:GLN:HB2	2.17	0.43
1:G:54:ILE:HG13	1:G:54:ILE:H	1.46	0.43
1:D:82:PRO:HG3	1:D:106:VAL:CA	2.39	0.43
1:H:81:ALA:CB	1:H:82:PRO:HA	2.48	0.43
1:G:131:GLN:HE22	1:H:134:TRP:CB	2.31	0.43
1:A:52:PHE:CD2	1:A:84:ALA:HB3	2.53	0.43
1:E:130:GLU:HG3	1:E:168:HIS:CG	2.54	0.43
1:G:100:GLY:C	1:G:102:THR:H	2.22	0.43
1:C:38:ILE:O	1:C:41:LEU:HB2	2.19	0.43
1:E:55:SER:OG	1:E:58:ARG:HG2	2.19	0.43
1:C:19:ALA:N	1:C:20:PRO:HD3	2.34	0.43
1:C:142:GLY:CA	1:C:172:VAL:HG23	2.48	0.43
1:G:125:VAL:HG12	1:G:126:THR:HG23	2.00	0.43
1:A:21:LYS:HB3	1:A:22:PRO:HD2	2.01	0.43
1:B:7:LYS:O	1:B:11:ARG:HG3	2.18	0.43
1:A:112:LYS:HG2	1:A:113:ARG:HG2	1.99	0.43
1:H:91:ASN:ND2	1:H:113:ARG:NE	2.59	0.43
1:D:18:LEU:O	1:D:119:ARG:NH1	2.32	0.43
1:E:146:SER:OG	1:G:150:SER:HB2	2.19	0.43
1:D:52:PHE:CD1	1:D:52:PHE:N	2.87	0.43
1:E:15:ILE:O	1:E:20:PRO:HD3	2.19	0.43
1:E:61:ARG:O	1:E:62:ILE:C	2.58	0.43
1:E:103:ALA:O	1:E:106:VAL:HB	2.19	0.43
1:A:46:GLN:O	1:A:47:THR:CG2	2.67	0.43
1:D:7:LYS:NZ	5:D:180:HOH:O	2.51	0.43
1:G:122:ALA:HB3	1:G:133:LEU:HD21	2.00	0.42
1:A:52:PHE:HE2	1:A:84:ALA:O	2.02	0.42
1:C:124:ALA:HB1	1:C:129:SER:CB	2.46	0.42
1:C:65:ASN:C	1:C:67:ILE:N	2.72	0.42
1:G:113:ARG:HG3	1:G:113:ARG:NH1	2.35	0.42
1:D:61:ARG:O	1:D:64:SER:HB3	2.20	0.42
1:G:150:SER:C	1:G:161:MSE:CE	2.88	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:63:LEU:O	1:C:67:ILE:HD12	2.19	0.42
1:F:125:VAL:O	1:G:160:THR:HG22	2.19	0.42
1:C:84:ALA:O	1:C:87:ASP:HB2	2.19	0.42
1:B:38:ILE:HG23	1:B:69:LEU:CD1	2.50	0.42
1:H:114:LEU:HA	1:H:115:PRO:HD2	1.74	0.42
1:H:96:PHE:CD1	1:H:121:VAL:HG12	2.54	0.42
1:F:27:TRP:CE3	1:F:92:PRO:HG3	2.54	0.42
1:H:44:THR:CG2	1:H:47:THR:HG23	2.49	0.42
1:A:4:GLN:HA	1:A:7:LYS:HB2	2.01	0.42
1:E:15:ILE:HG13	1:E:39:GLU:HG2	2.00	0.42
1:A:107:PHE:O	1:A:111:TRP:N	2.48	0.42
1:F:92:PRO:HD2	1:F:113:ARG:O	2.19	0.42
1:D:114:LEU:HD11	1:D:118:GLY:HA3	2.02	0.42
1:E:148:ALA:HB3	1:G:148:ALA:HB3	2.02	0.42
1:G:91:ASN:ND2	1:G:113:ARG:NE	2.60	0.42
1:F:137:ARG:HD2	1:F:144:ILE:HG12	2.02	0.42
1:H:25:THR:HA	1:H:48:THR:HB	2.01	0.42
1:F:172:VAL:HG13	1:F:172:VAL:O	2.20	0.42
1:A:101:LEU:HD21	1:A:133:LEU:HD21	2.02	0.42
1:E:144:ILE:HD12	1:E:144:ILE:HG23	1.70	0.42
1:G:20:PRO:HG3	1:G:40:TRP:HE3	1.73	0.42
1:H:91:ASN:ND2	1:H:113:ARG:O	2.53	0.42
1:F:26:LEU:CG	1:F:27:TRP:N	2.82	0.42
1:B:38:ILE:HG23	1:B:69:LEU:HD13	2.01	0.42
1:H:17:ALA:CB	5:H:178:HOH:O	2.67	0.42
1:E:108:ALA:O	1:E:112:LYS:HG3	2.20	0.42
1:C:119:ARG:HG2	1:C:173:VAL:HG22	2.01	0.42
1:A:61:ARG:O	1:A:61:ARG:HG3	2.20	0.42
1:E:171:THR:HG22	5:E:199:HOH:O	2.19	0.42
1:D:63:LEU:HD11	1:F:58:ARG:NH2	2.35	0.42
1:A:107:PHE:O	1:A:108:ALA:C	2.58	0.42
1:A:95:ILE:HD12	1:A:114[B]:LEU:CD2	2.50	0.42
1:E:56:GLU:HB2	1:E:59:ARG:HH21	1.85	0.42
1:B:128:GLU:HG3	1:C:158:PHE:CE1	2.54	0.41
1:C:20:PRO:HG3	1:C:40:TRP:CE3	2.55	0.41
1:D:87:ASP:O	1:D:89:PRO:CD	2.55	0.41
1:G:113:ARG:HA	1:G:113:ARG:HD2	1.33	0.41
1:H:15:ILE:HA	1:H:18:LEU:HB2	2.02	0.41
1:H:27:TRP:HB2	1:H:95:ILE:HG12	2.02	0.41
1:G:71:VAL:HG12	1:G:71:VAL:O	2.20	0.41
1:C:126:THR:H	1:C:129:SER:HB2	1.85	0.41
1:B:128:GLU:HG2	1:B:128:GLU:H	1.47	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:F:133:LEU:HD13	1:F:170:TRP:HB2	2.02	0.41
1:A:46:GLN:O	1:A:47:THR:HG23	2.19	0.41
1:B:67:ILE:HA	1:B:72:SER:OG	2.20	0.41
1:E:59:ARG:HH11	1:E:79:GLN:HA	1.85	0.41
1:G:43:SER:OG	1:G:44:THR:N	2.53	0.41
1:F:79:GLN:HB3	1:F:79:GLN:HE21	1.50	0.41
1:H:8:GLN:O	1:H:12:ALA:N	2.51	0.41
1:B:102:THR:O	1:B:103:ALA:C	2.58	0.41
1:F:15:ILE:HD12	1:F:40:TRP:HA	2.03	0.41
1:H:87:ASP:N	1:H:87:ASP:OD1	2.54	0.41
1:H:37:ALA:O	1:H:41:LEU:HG	2.21	0.41
1:E:53:GLU:HG2	1:E:58:ARG:CB	2.49	0.41
1:A:131:GLN:HB3	5:A:188:HOH:O	2.19	0.41
1:C:18:LEU:CD2	1:C:121:VAL:HB	2.51	0.41
1:A:26:LEU:HB2	1:A:40:TRP:CZ2	2.56	0.41
1:C:90:ASP:C	1:C:90:ASP:OD2	2.58	0.41
1:F:45:PRO:HB2	1:F:46:GLN:HG3	2.02	0.41
1:H:97:ILE:HG13	1:H:120:LEU:HD21	2.02	0.41
1:A:124:ALA:HA	5:A:280:HOH:O	2.20	0.41
1:E:85:PHE:CE2	1:E:110:ALA:HB2	2.55	0.41
1:C:42:ARG:CB	1:C:42:ARG:HH21	2.33	0.41
1:H:55:SER:HB2	1:H:58[A]:ARG:HG3	2.02	0.41
1:H:174:LYS:O	1:H:174:LYS:HG2	2.20	0.41
1:C:36:ILE:HG13	1:C:96:PHE:CE2	2.55	0.41
1:F:162:LYS:O	1:F:162:LYS:CG	2.69	0.41
1:C:59:ARG:C	1:C:61:ARG:H	2.23	0.41
1:A:133:LEU:HD22	1:A:170:TRP:CB	2.51	0.41
1:D:4:GLN:HG2	1:D:7:LYS:NZ	2.36	0.41
1:A:34:GLY:O	1:A:38:ILE:HD12	2.21	0.41
1:G:133:LEU:O	1:G:137:ARG:CB	2.69	0.41
1:H:98:GLY:O	1:H:123:ASN:CB	2.69	0.41
1:A:95:ILE:HD13	1:A:114[B]:LEU:HD11	2.02	0.41
1:A:123:ASN:HA	1:A:168:HIS:O	2.21	0.41
1:A:109:ALA:HA	1:A:112:LYS:CE	2.50	0.41
1:D:133:LEU:HG	1:D:170:TRP:HB2	2.02	0.41
1:B:152:GLU:HG2	1:D:144:ILE:O	2.20	0.41
1:B:153:HIS:HE1	1:B:155:VAL:HG23	1.85	0.41
1:H:67:ILE:HG22	1:H:68:ASN:N	2.36	0.41
1:F:42:ARG:NH2	1:F:42:ARG:HG2	2.36	0.41
1:D:162:LYS:HA	1:D:163:PRO:HD3	1.97	0.41
1:H:47:THR:H	1:H:74:ARG:HH11	1.67	0.41
1:B:38:ILE:CG2	1:B:69:LEU:CD1	2.99	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:F:7:LYS:HD2	1:F:11:ARG:NE	2.12	0.40
1:G:53:GLU:OE1	1:G:53:GLU:HA	2.21	0.40
1:D:71:VAL:CG1	1:D:74:ARG:HD2	2.51	0.40
1:H:133:LEU:HD22	1:H:170:TRP:HB2	2.02	0.40
1:F:160:THR:HG22	1:G:126:THR:HG22	2.04	0.40
1:G:36:ILE:O	1:G:40:TRP:HB3	2.21	0.40
1:F:144:ILE:HG22	1:F:145:SER:N	2.36	0.40
1:G:131:GLN:NE2	1:H:134:TRP:CB	2.84	0.40
1:C:41:LEU:CD1	1:C:75:ILE:HB	2.52	0.40
1:G:53:GLU:CD	1:G:54:ILE:N	2.74	0.40
1:C:82:PRO:HA	1:C:85:PHE:CE2	2.56	0.40
1:B:54:ILE:HG23	1:B:55:SER:N	2.36	0.40
1:D:111:TRP:CE2	1:D:174:LYS:HE3	2.56	0.40
1:H:55:SER:HB2	1:H:58[B]:ARG:CG	2.47	0.40
1:D:27:TRP:CZ3	1:D:85:PHE:CD1	3.10	0.40
1:F:146:SER:OG	1:H:161:MSE:CE	2.65	0.40
1:G:81:ALA:CB	1:G:85:PHE:CZ	3.04	0.40
1:A:163:PRO:CG	1:D:163:PRO:HB2	2.52	0.40
1:H:153:HIS:H	1:H:159:ILE:HG22	1.87	0.40
1:G:57:GLU:O	1:G:61:ARG:HG3	2.22	0.40
1:F:155:VAL:HG12	1:F:155:VAL:O	2.20	0.40
1:E:22:PRO:O	1:E:23:HIS:HB2	2.22	0.40
1:C:59:ARG:HG2	1:C:77:VAL:HG12	2.04	0.40
1:B:117:GLY:HA2	1:B:174:LYS:O	2.22	0.40
1:D:67:ILE:HG12	1:D:72:SER:HB2	2.03	0.40
1:G:15:ILE:HD12	1:G:39:GLU:HG3	2.04	0.40
1:H:94:VAL:CG1	1:H:95:ILE:N	2.83	0.40
1:A:137:ARG:NH1	1:A:138:LYS:HE2	2.36	0.40
1:F:101:LEU:HD11	1:F:132:MSE:HG2	2.04	0.40
1:F:20:PRO:HG3	1:F:40:TRP:CE3	2.57	0.40
1:G:82:PRO:HG3	1:G:106:VAL:CA	2.51	0.40
1:B:153:HIS:CE1	1:B:155:VAL:HG23	2.56	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	169/178 (95%)	149 (88%)	20 (12%)	0	100	100
1	B	167/178 (94%)	148 (89%)	19 (11%)	0	100	100
1	C	165/178 (93%)	151 (92%)	12 (7%)	2 (1%)	19	14
1	D	159/178 (89%)	147 (92%)	12 (8%)	0	100	100
1	E	166/178 (93%)	152 (92%)	14 (8%)	0	100	100
1	F	162/178 (91%)	156 (96%)	6 (4%)	0	100	100
1	G	167/178 (94%)	151 (90%)	16 (10%)	0	100	100
1	H	160/178 (90%)	143 (89%)	17 (11%)	0	100	100
All	All	1315/1424 (92%)	1197 (91%)	116 (9%)	2 (0%)	56	62

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	66	ALA
1	C	156	GLY

5.3.2 Protein sidechains ⓘ

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	139/139 (100%)	123 (88%)	16 (12%)	8	7
1	B	137/139 (99%)	122 (89%)	15 (11%)	9	8
1	C	136/139 (98%)	116 (85%)	20 (15%)	4	3
1	D	134/139 (96%)	119 (89%)	15 (11%)	9	7
1	E	136/139 (98%)	125 (92%)	11 (8%)	17	16
1	F	137/139 (99%)	114 (83%)	23 (17%)	3	2
1	G	135/139 (97%)	117 (87%)	18 (13%)	6	4
1	H	133/139 (96%)	107 (80%)	26 (20%)	2	1
All	All	1087/1112 (98%)	943 (87%)	144 (13%)	6	4

All (144) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	LEU
1	A	11[A]	ARG
1	A	11[B]	ARG
1	A	16	SER
1	A	23	HIS
1	A	28	ASP
1	A	36	ILE
1	A	54	ILE
1	A	64	SER
1	A	72	SER
1	A	102	THR
1	A	112	LYS
1	A	143	THR
1	A	152	GLU
1	A	155	VAL
1	A	157	SER
1	B	5	LEU
1	B	7	LYS
1	B	36	ILE
1	B	44	THR
1	B	57	GLU
1	B	65	ASN
1	B	73	ASP
1	B	83	ARG
1	B	90	ASP
1	B	101	LEU
1	B	102	THR
1	B	128	GLU
1	B	145	SER
1	B	152	GLU
1	B	157	SER
1	C	4	GLN
1	C	8	GLN
1	C	15	ILE
1	C	23	HIS
1	C	38	ILE
1	C	42	ARG
1	C	43	SER
1	C	48	THR
1	C	54	ILE
1	C	67	ILE
1	C	69	LEU

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Mol	Chain	Res	Type
1	C	71	VAL
1	C	83	ARG
1	C	93	ASP
1	C	137	ARG
1	C	138	LYS
1	C	143	THR
1	C	152	GLU
1	C	154	THR
1	C	158	PHE
1	D	6	THR
1	D	24	GLU
1	D	35	SER
1	D	38	ILE
1	D	42	ARG
1	D	46	GLN
1	D	48	THR
1	D	54	ILE
1	D	65	ASN
1	D	71	VAL
1	D	90	ASP
1	D	91	ASN
1	D	101	LEU
1	D	132	MSE
1	D	149	ILE
1	E	4	GLN
1	E	6	THR
1	E	7	LYS
1	E	42	ARG
1	E	51	CYS
1	E	55	SER
1	E	79	GLN
1	E	90	ASP
1	E	132	MSE
1	E	139	GLN
1	E	143	THR
1	F	5	LEU
1	F	6	THR
1	F	41	LEU
1	F	42	ARG
1	F	46	GLN
1	F	51	CYS
1	F	57	GLU

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Mol	Chain	Res	Type
1	F	73	ASP
1	F	79	GLN
1	F	86	ASP
1	F	90[A]	ASP
1	F	90[B]	ASP
1	F	91	ASN
1	F	101	LEU
1	F	102	THR
1	F	129	SER
1	F	132	MSE
1	F	138	LYS
1	F	146	SER
1	F	153	HIS
1	F	157	SER
1	F	159	ILE
1	F	162	LYS
1	G	4	GLN
1	G	26	LEU
1	G	40	TRP
1	G	43	SER
1	G	44	THR
1	G	47	THR
1	G	55	SER
1	G	73	ASP
1	G	83	ARG
1	G	90	ASP
1	G	91	ASN
1	G	101	LEU
1	G	102	THR
1	G	112	LYS
1	G	125	VAL
1	G	137	ARG
1	G	143	THR
1	G	171	THR
1	H	4	GLN
1	H	5	LEU
1	H	38	ILE
1	H	40	TRP
1	H	42	ARG
1	H	43	SER
1	H	47	THR
1	H	48	THR

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Mol	Chain	Res	Type
1	H	51	CYS
1	H	54	ILE
1	H	56	GLU
1	H	57	GLU
1	H	64	SER
1	H	71	VAL
1	H	79	GLN
1	H	83	ARG
1	H	86	ASP
1	H	87	ASP
1	H	93	ASP
1	H	101	LEU
1	H	123	ASN
1	H	137	ARG
1	H	138	LYS
1	H	145	SER
1	H	150	SER
1	H	159	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (30) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	9	HIS
1	A	23	HIS
1	A	91	ASN
1	A	168	HIS
1	B	169	GLN
1	C	78	GLN
1	C	79	GLN
1	C	153	HIS
1	D	8	GLN
1	D	9	HIS
1	D	23	HIS
1	D	65	ASN
1	D	91	ASN
1	E	65	ASN
1	E	91	ASN
1	E	169	GLN
1	F	9	HIS
1	F	23	HIS
1	F	65	ASN
1	F	79	GLN

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Mol	Chain	Res	Type
1	F	91	ASN
1	F	169	GLN
1	G	9	HIS
1	G	65	ASN
1	G	91	ASN
1	G	131	GLN
1	H	79	GLN
1	H	91	ASN
1	H	123	ASN
1	H	151	HIS

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 7 ligands modelled in this entry, 4 are monoatomic - leaving 3 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	ACY	A	176	-	3,3,3	0.62	0	3,3,3	1.02	0
2	ACY	D	177	-	3,3,3	0.60	0	3,3,3	1.04	0
2	ACY	F	176	-	3,3,3	0.65	0	3,3,3	0.60	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	ACY	A	176	-	-	0/0/0/0	0/0/0/0
2	ACY	D	177	-	-	0/0/0/0	0/0/0/0
2	ACY	F	176	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	171/178 (96%)	-0.03	2 (1%) 75 76	25, 42, 67, 90	0
1	B	171/178 (96%)	-0.08	3 (1%) 65 66	25, 43, 62, 90	0
1	C	169/178 (94%)	0.20	6 (3%) 41 41	29, 47, 81, 105	0
1	D	167/178 (93%)	-0.06	0 100 100	24, 44, 68, 97	0
1	E	170/178 (95%)	-0.17	0 100 100	22, 39, 62, 94	0
1	F	167/178 (93%)	-0.21	0 100 100	23, 38, 59, 74	0
1	G	171/178 (96%)	-0.01	4 (2%) 57 58	27, 44, 66, 83	0
1	H	167/178 (93%)	-0.02	3 (1%) 65 66	28, 46, 69, 89	0
All	All	1353/1424 (95%)	-0.05	18 (1%) 74 74	22, 43, 69, 105	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	70	GLY	8.7
1	C	69	LEU	6.4
1	C	44	THR	6.4
1	G	34	GLY	4.7
1	C	155	VAL	4.2
1	H	175	ALA	3.7
1	C	71	VAL	3.6
1	B	83	ARG	2.9
1	C	45	PRO	2.9
1	B	31	GLY	2.8
1	B	101	LEU	2.7
1	G	33	SER	2.5
1	H	99	GLY	2.2
1	A	54	ILE	2.2
1	G	54	ILE	2.1
1	A	118	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
1	G	175	ALA	2.1
1	H	157	SER	2.0

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

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

6.3 Carbohydrates

There are no carbohydrates in this entry.

6.4 Ligands

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	ACY	F	176	4/4	0.25	2.50	42,54,57,71	0
2	ACY	A	176	4/4	0.13	1.35	38,43,49,55	0
2	ACY	D	177	4/4	0.14	0.34	35,45,47,49	0
4	CA	G	177	1/1	0.15	0.15	56,56,56,56	0
3	MG	B	176	1/1	0.13	-0.65	39,39,39,39	0
3	MG	G	176	1/1	0.06	-2.26	43,43,43,43	0
4	CA	D	176	1/1	0.10	-3.57	64,64,64,64	0

6.5 Other polymers

There are no such residues in this entry.