



Full wwPDB X-ray Structure Validation Report

Mar 1, 2014 – 02:33 AM GMT

PDB ID : 1QZ2
Title : Crystal Structure of FKBP52 C-terminal Domain complex with the C-terminal peptide MEEVD of Hsp90
Authors : Wu, B.; Li, P.; Lou, Z.; Liu, Y.; Ding, Y.; Shu, C.; Shen, B.; Rao, Z.
Deposited on : 2003-09-15
Resolution : 3.00 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>

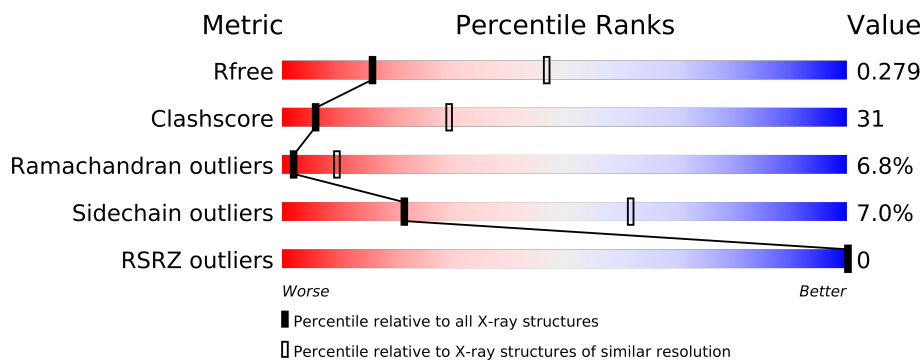
The following versions of software and data (see [references](#)) were used in the production of this report:

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

1 Overall quality at a glance

The reported resolution of this entry is 3.00 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	1216 (3.00-3.00)
Clashscore	79885	1594 (3.00-3.00)
Ramachandran outliers	78287	1537 (3.00-3.00)
Sidechain outliers	78261	1540 (3.00-3.00)
RSRZ outliers	66119	1217 (3.00-3.00)

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	336	
1	B	336	
1	C	336	
2	G	5	
2	H	5	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7225 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 FK506-binding protein 4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	285	Total	C	N	O	S	0	0	0
			2305	1455	408	434	8			
1	B	285	Total	C	N	O	S	0	0	0
			2305	1455	408	434	8			
1	C	285	Total	C	N	O	S	0	0	0
			2310	1458	408	436	8			

There are 63 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-16	MET	-	CLONING ARTIFACT	UNP Q02790
A	-15	GLY	-	CLONING ARTIFACT	UNP Q02790
A	-14	SER	-	CLONING ARTIFACT	UNP Q02790
A	-13	SER	-	CLONING ARTIFACT	UNP Q02790
A	-12	HIS	-	CLONING ARTIFACT	UNP Q02790
A	-11	HIS	-	CLONING ARTIFACT	UNP Q02790
A	-10	HIS	-	CLONING ARTIFACT	UNP Q02790
A	-9	HIS	-	CLONING ARTIFACT	UNP Q02790
A	-8	HIS	-	CLONING ARTIFACT	UNP Q02790
A	-7	HIS	-	CLONING ARTIFACT	UNP Q02790
A	-6	SER	-	CLONING ARTIFACT	UNP Q02790
A	-5	SER	-	CLONING ARTIFACT	UNP Q02790
A	-4	GLY	-	CLONING ARTIFACT	UNP Q02790
A	-3	LEU	-	CLONING ARTIFACT	UNP Q02790
A	-2	VAL	-	CLONING ARTIFACT	UNP Q02790
A	-1	PRO	-	CLONING ARTIFACT	UNP Q02790
A	0	ARG	-	CLONING ARTIFACT	UNP Q02790
A	1	GLY	-	CLONING ARTIFACT	UNP Q02790
A	2	SER	-	CLONING ARTIFACT	UNP Q02790
A	3	HIS	-	CLONING ARTIFACT	UNP Q02790
A	4	MET	-	CLONING ARTIFACT	UNP Q02790
B	-16	MET	-	CLONING ARTIFACT	UNP Q02790
B	-15	GLY	-	CLONING ARTIFACT	UNP Q02790

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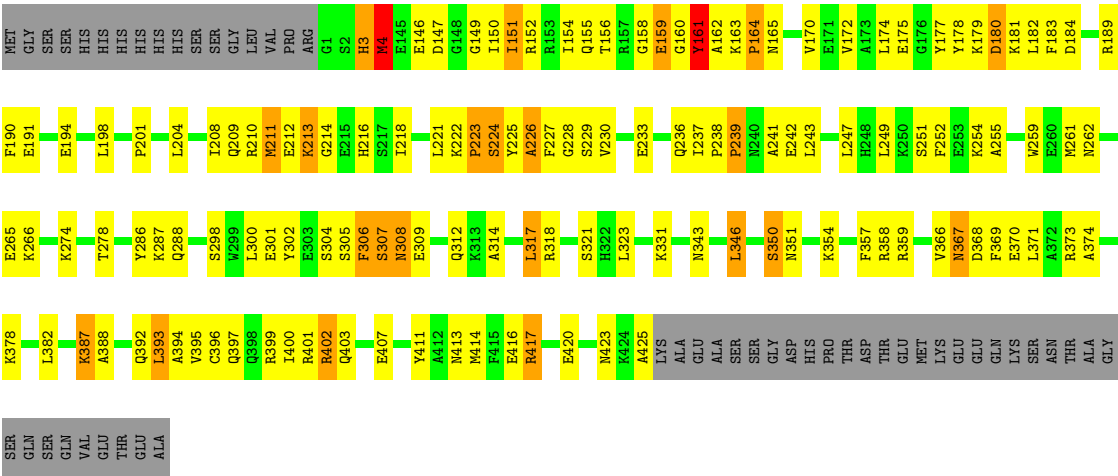
Chain	Residue	Modelled	Actual	Comment	Reference
B	-14	SER	-	CLONING ARTIFACT	UNP Q02790
B	-13	SER	-	CLONING ARTIFACT	UNP Q02790
B	-12	HIS	-	CLONING ARTIFACT	UNP Q02790
B	-11	HIS	-	CLONING ARTIFACT	UNP Q02790
B	-10	HIS	-	CLONING ARTIFACT	UNP Q02790
B	-9	HIS	-	CLONING ARTIFACT	UNP Q02790
B	-8	HIS	-	CLONING ARTIFACT	UNP Q02790
B	-7	HIS	-	CLONING ARTIFACT	UNP Q02790
B	-6	SER	-	CLONING ARTIFACT	UNP Q02790
B	-5	SER	-	CLONING ARTIFACT	UNP Q02790
B	-4	GLY	-	CLONING ARTIFACT	UNP Q02790
B	-3	LEU	-	CLONING ARTIFACT	UNP Q02790
B	-2	VAL	-	CLONING ARTIFACT	UNP Q02790
B	-1	PRO	-	CLONING ARTIFACT	UNP Q02790
B	0	ARG	-	CLONING ARTIFACT	UNP Q02790
B	1	GLY	-	CLONING ARTIFACT	UNP Q02790
B	2	SER	-	CLONING ARTIFACT	UNP Q02790
B	3	HIS	-	CLONING ARTIFACT	UNP Q02790
B	4	MET	-	CLONING ARTIFACT	UNP Q02790
C	-16	MET	-	CLONING ARTIFACT	UNP Q02790
C	-15	GLY	-	CLONING ARTIFACT	UNP Q02790
C	-14	SER	-	CLONING ARTIFACT	UNP Q02790
C	-13	SER	-	CLONING ARTIFACT	UNP Q02790
C	-12	HIS	-	CLONING ARTIFACT	UNP Q02790
C	-11	HIS	-	CLONING ARTIFACT	UNP Q02790
C	-10	HIS	-	CLONING ARTIFACT	UNP Q02790
C	-9	HIS	-	CLONING ARTIFACT	UNP Q02790
C	-8	HIS	-	CLONING ARTIFACT	UNP Q02790
C	-7	HIS	-	CLONING ARTIFACT	UNP Q02790
C	-6	SER	-	CLONING ARTIFACT	UNP Q02790
C	-5	SER	-	CLONING ARTIFACT	UNP Q02790
C	-4	GLY	-	CLONING ARTIFACT	UNP Q02790
C	-3	LEU	-	CLONING ARTIFACT	UNP Q02790
C	-2	VAL	-	CLONING ARTIFACT	UNP Q02790
C	-1	PRO	-	CLONING ARTIFACT	UNP Q02790
C	0	ARG	-	CLONING ARTIFACT	UNP Q02790
C	1	GLY	-	CLONING ARTIFACT	UNP Q02790
C	2	SER	-	CLONING ARTIFACT	UNP Q02790
C	3	HIS	-	CLONING ARTIFACT	UNP Q02790
C	4	MET	-	CLONING ARTIFACT	UNP Q02790

- Molecule 2 is a protein called 5-mer peptide from Heat shock protein HSP 90.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	G	5	Total 42	C 24	N 5	O 12	S 1	0	0	0
2	H	5	Total 42	C 24	N 5	O 12	S 1	0	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	75	Total 75	O 75	0	0
3	B	64	Total 64	O 64	0	0
3	C	76	Total 76	O 76	0	0
3	G	4	Total 4	O 4	0	0
3	H	2	Total 2	O 2	0	0



- Molecule 2: 5-mer peptide from Heat shock protein HSP 90

Chain G:



- Molecule 2: 5-mer peptide from Heat shock protein HSP 90

Chain H:



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	111.61Å 144.42Å 170.81Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 3.00 47.85 – 2.88	Depositor EDS
% Data completeness (in resolution range)	(Not available) (50.00-3.00) 91.5 (47.85-2.88)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.15 (at 2.91Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.230 , 0.287 0.228 , 0.279	Depositor DCC
R_{free} test set	1270 reflections (4.87%)	DCC
Wilson B-factor (Å ²)	51.5	Xtriage
Anisotropy	0.543	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 35.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 30914 reflections	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	7225	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.48% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality i

5.1 Standard geometry i

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	0.39	1/2345 (0.0%)	0.63	0/3138
1	B	0.38	0/2345	0.62	0/3138
1	C	0.40	0/2350	0.65	1/3145 (0.0%)
2	G	0.57	0/41	0.63	0/52
2	H	0.64	0/41	1.16	0/52
All	All	0.40	1/7122 (0.0%)	0.64	1/9525 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	4	MET	SD-CE	5.21	2.07	1.77

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	3	HIS	N-CA-C	6.21	127.78	111.00

There are no chirality outliers.

There are no planarity outliers.

5.2 Close contacts i

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	2305	0	2296	171	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	2305	0	2296	143	0
1	C	2310	0	2301	135	0
2	G	42	0	36	2	0
2	H	42	0	36	2	0
3	A	75	0	0	10	0
3	B	64	0	0	4	0
3	C	76	0	0	5	0
3	G	4	0	0	1	0
3	H	2	0	0	0	0
All	All	7225	0	6965	434	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 31.

All (434) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:4:MET:SD	1:A:4:MET:CE	2.07	1.42
1:B:204:LEU:HD11	1:B:247:LEU:HD11	1.25	1.10
1:C:417:ARG:HH11	1:C:417:ARG:HB3	1.24	1.01
1:B:223:PRO:HD3	1:B:242:GLU:HG2	1.43	1.00
1:A:212:GLU:HB2	1:A:215:GLU:HB2	1.44	0.97
1:C:387:LYS:HG3	1:C:388:ALA:H	1.33	0.93
1:A:387:LYS:HD3	1:A:387:LYS:H	1.34	0.92
1:A:261:MET:HE3	1:A:266:LYS:HG2	1.50	0.92
1:C:417:ARG:HB3	1:C:417:ARG:NH1	1.85	0.91
1:C:301:GLU:HG3	1:C:302:TYR:HD1	1.37	0.90
1:A:223:PRO:HD3	1:A:242:GLU:HG2	1.52	0.90
1:B:350:SER:H	1:C:350:SER:HB2	1.35	0.89
1:A:417:ARG:HD3	1:C:259:TRP:O	1.73	0.87
1:C:170:VAL:HG12	1:C:172:VAL:HG23	1.56	0.87
1:B:186:ARG:HD2	1:B:188:LEU:HD21	1.57	0.86
1:B:353:GLU:O	1:B:357:PHE:HB2	1.75	0.86
1:C:357:PHE:HZ	1:C:388:ALA:HB3	1.40	0.84
1:A:356:LEU:HD21	1:A:378:LYS:HD3	1.60	0.83
1:B:173:ALA:HB3	1:B:248:HIS:HB3	1.58	0.83
1:B:157:ARG:N	1:B:157:ARG:HD3	1.95	0.81
1:B:145:GLU:H	1:B:145:GLU:CD	1.83	0.81
1:C:208:ILE:HA	1:C:211:MET:HG3	1.60	0.80
1:A:4:MET:HG2	1:A:210:ARG:NH1	1.98	0.79
1:C:4:MET:H	1:C:4:MET:CE	1.95	0.79
1:C:223:PRO:HG3	1:C:242:GLU:N	1.97	0.78
1:A:156:THR:HB	1:A:216:HIS:HB2	1.64	0.78

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:170:VAL:HG23	1:B:252:PHE:CB	2.14	0.77
1:A:380:LEU:HD22	1:A:389:ALA:HB1	1.65	0.77
1:C:307:SER:O	1:C:309:GLU:N	2.17	0.77
1:A:349:ASP:O	1:A:351:ASN:N	2.17	0.77
1:C:226:ALA:HB1	1:C:243:LEU:HD12	1.66	0.76
1:B:166:GLU:O	1:B:168:ALA:N	2.19	0.75
1:B:226:ALA:HB1	1:B:243:LEU:HD12	1.68	0.74
1:A:154:ILE:H	1:A:154:ILE:HD12	1.51	0.74
1:A:154:ILE:HD12	1:A:154:ILE:N	2.04	0.73
1:C:178:TYR:HE1	1:C:241:ALA:HB1	1.54	0.73
1:B:191:GLU:O	1:B:194:GLU:HG2	1.89	0.73
1:B:160:GLY:HA2	1:B:212:GLU:OE1	1.89	0.72
1:B:166:GLU:O	1:B:192:ILE:HB	1.89	0.72
1:B:170:VAL:HG23	1:B:252:PHE:HB3	1.72	0.72
1:A:219:VAL:HG23	1:A:247:LEU:HD21	1.71	0.72
1:B:410:LEU:HD12	1:B:410:LEU:O	1.89	0.71
1:C:182:LEU:O	1:C:182:LEU:HD23	1.91	0.71
1:A:227:PHE:HB3	1:A:231:GLY:O	1.91	0.71
1:A:366:VAL:HB	1:C:4:MET:O	1.91	0.71
1:B:266:LYS:CD	1:B:303:GLU:HG2	2.21	0.71
1:C:357:PHE:CZ	1:C:388:ALA:HB3	2.26	0.70
1:B:179:LYS:O	1:B:180:ASP:HB2	1.90	0.70
1:C:201:PRO:HB3	1:C:226:ALA:HA	1.74	0.70
1:C:194:GLU:HG3	1:C:288:GLN:HG2	1.72	0.70
1:A:226:ALA:HB1	1:A:243:LEU:HD12	1.73	0.69
1:A:152:ARG:HA	1:A:218:ILE:O	1.93	0.69
1:A:349:ASP:HB3	1:A:352:ASN:HB2	1.74	0.69
1:C:165:ASN:OD1	1:C:254:LYS:HD3	1.92	0.69
1:C:4:MET:H	1:C:4:MET:HE3	1.57	0.69
1:C:368:ASP:HB3	1:C:371:LEU:HD12	1.74	0.68
1:A:218:ILE:HA	1:A:245:TYR:O	1.93	0.68
1:B:149:GLY:HA3	1:B:225:TYR:CE1	2.29	0.68
1:A:212:GLU:O	1:A:249:LEU:HD23	1.94	0.68
1:A:387:LYS:CD	1:A:387:LYS:H	2.05	0.67
1:C:223:PRO:HG3	1:C:241:ALA:C	2.15	0.67
1:C:298:SER:O	1:C:301:GLU:HG2	1.95	0.67
1:B:156:THR:HG22	1:B:156:THR:O	1.95	0.67
1:C:354:LYS:HD3	1:C:358:ARG:NH2	2.10	0.66
1:C:156:THR:HB	1:C:216:HIS:HB2	1.77	0.66
1:B:286:TYR:HB2	1:B:332:LEU:HD21	1.78	0.66
1:A:173:ALA:HB3	1:A:248:HIS:HB3	1.77	0.65
1:A:226:ALA:C	1:A:228:GLY:H	2.00	0.65

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:163:LYS:HD3	1:A:210:ARG:HA	1.79	0.65
1:A:174:LEU:HD11	1:A:245:TYR:CD2	2.32	0.65
1:B:266:LYS:HD2	1:B:303:GLU:HG2	1.79	0.65
1:C:204:LEU:HD11	1:C:247:LEU:HD11	1.79	0.65
1:B:148:GLY:O	1:B:222:LYS:HE3	1.96	0.65
1:B:257:GLU:HG3	1:B:260:GLU:OE2	1.97	0.65
1:C:343:ASN:ND2	1:C:359:ARG:HH11	1.96	0.64
1:A:251:SER:O	1:A:252:PHE:HB3	1.95	0.64
1:A:204:LEU:O	1:A:208:ILE:HG13	1.97	0.64
1:A:211:MET:HE3	1:A:249:LEU:HB2	1.80	0.64
1:C:301:GLU:HG3	1:C:302:TYR:CD1	2.26	0.64
1:B:373:ARG:HE	1:C:304:SER:HB2	1.62	0.64
1:A:247:LEU:HD23	1:A:247:LEU:N	2.13	0.64
1:B:204:LEU:HD11	1:B:247:LEU:CD1	2.15	0.63
1:A:146:GLU:HB2	1:A:225:TYR:OH	1.98	0.63
1:C:208:ILE:HG12	1:C:247:LEU:HD13	1.80	0.63
1:B:170:VAL:HG23	1:B:252:PHE:HB2	1.79	0.63
1:B:156:THR:HB	1:B:216:HIS:HB2	1.80	0.63
1:C:211:MET:HB2	1:C:249:LEU:HD22	1.79	0.63
1:A:247:LEU:H	1:A:247:LEU:HD23	1.62	0.63
1:C:204:LEU:O	1:C:208:ILE:HG13	1.99	0.62
1:A:176:GLY:HA2	1:A:244:LYS:O	1.99	0.62
1:B:170:VAL:HG22	1:B:249:LEU:HD11	1.80	0.62
1:B:174:LEU:HD11	1:B:200:LEU:HD21	1.82	0.62
1:B:173:ALA:CB	1:B:248:HIS:HB3	2.28	0.62
1:C:223:PRO:O	1:C:224:SER:HB2	1.98	0.62
1:B:261:MET:HB3	1:B:265:GLU:HB3	1.82	0.61
1:A:229:SER:O	1:A:239:PRO:HB3	2.00	0.61
1:B:353:GLU:HB2	1:B:383:TYR:CE1	2.36	0.61
1:A:177:TYR:CE2	1:A:182:LEU:HB2	2.35	0.61
1:A:155:GLN:HG2	1:A:217:SER:HA	1.82	0.61
1:B:149:GLY:HA3	1:B:225:TYR:HE1	1.65	0.61
1:C:152:ARG:HG2	1:C:152:ARG:HH11	1.65	0.61
1:C:314:ALA:O	1:C:318:ARG:HG3	2.00	0.61
1:C:150:ILE:HD11	1:C:225:TYR:HD2	1.65	0.61
1:B:151:ILE:H	1:B:151:ILE:HD12	1.66	0.61
1:A:174:LEU:HD11	1:A:245:TYR:HD2	1.65	0.61
1:A:294:LYS:O	1:A:297:VAL:HG22	2.01	0.61
1:B:174:LEU:O	1:B:175:GLU:HG3	2.00	0.61
1:B:305:SER:HA	1:C:373:ARG:NH1	2.16	0.61
1:C:223:PRO:HB3	1:C:241:ALA:H	1.64	0.61
1:A:357:PHE:CZ	1:A:388:ALA:HB3	2.36	0.60

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:170:VAL:CG2	1:B:249:LEU:HD11	2.30	0.60
1:C:343:ASN:HD21	1:C:359:ARG:HH11	1.49	0.60
1:B:174:LEU:HD12	1:B:186:ARG:HG3	1.83	0.60
1:A:318:ARG:HD3	3:A:476:HOH:O	2.02	0.60
1:A:233:GLU:C	1:A:235:PHE:H	2.01	0.60
1:A:233:GLU:O	1:A:235:PHE:N	2.33	0.60
1:B:162:ALA:O	1:B:163:LYS:HB2	2.01	0.60
1:B:1:GLY:O	1:B:2:SER:HB3	2.02	0.60
1:C:191:GLU:O	1:C:194:GLU:HB3	2.01	0.60
1:A:357:PHE:HZ	1:A:388:ALA:HB3	1.67	0.59
1:C:407:GLU:HG2	1:C:411:TYR:HE1	1.66	0.59
1:A:233:GLU:HG3	1:C:238:PRO:CG	2.32	0.59
1:A:154:ILE:H	1:A:154:ILE:CD1	2.15	0.59
1:A:177:TYR:CD1	1:A:244:LYS:HD3	2.37	0.59
1:A:399:ARG:HD2	3:A:462:HOH:O	2.02	0.59
1:C:387:LYS:HG3	1:C:388:ALA:N	2.13	0.59
1:B:327:MET:HB2	1:B:358:ARG:NH1	2.18	0.59
1:B:285:LYS:HB3	1:B:288:GLN:OE1	2.03	0.58
1:A:172:VAL:HG23	1:A:249:LEU:HD12	1.86	0.58
1:B:300:LEU:O	1:B:318:ARG:NH1	2.33	0.58
1:B:3:HIS:O	1:B:4:MET:HG2	2.03	0.58
1:C:174:LEU:C	1:C:174:LEU:HD12	2.23	0.58
1:B:153:ARG:HB2	1:B:218:ILE:HG22	1.86	0.58
1:C:204:LEU:CD1	1:C:247:LEU:HD11	2.33	0.57
1:A:286:TYR:CE2	1:A:331:LYS:HG2	2.40	0.57
1:A:197:ASN:HD22	1:A:288:GLN:HG2	1.69	0.57
1:A:2:SER:O	1:A:3:HIS:C	2.43	0.57
2:G:1:MET:HG2	2:G:2:GLU:N	2.19	0.57
1:A:233:GLU:C	1:A:235:PHE:N	2.58	0.57
1:B:307:SER:O	1:B:308:ASN:HB2	2.05	0.57
1:A:286:TYR:CZ	1:A:331:LYS:HG2	2.40	0.57
1:A:349:ASP:C	1:A:351:ASN:H	2.08	0.57
1:C:387:LYS:CG	1:C:388:ALA:H	2.09	0.57
1:B:174:LEU:HD23	1:B:247:LEU:CD2	2.35	0.57
1:B:157:ARG:N	1:B:157:ARG:CD	2.68	0.57
1:B:149:GLY:HA2	1:B:222:LYS:HD2	1.86	0.57
1:B:286:TYR:CB	1:B:332:LEU:HD21	2.35	0.56
1:B:165:ASN:HD21	1:B:255:ALA:HB3	1.68	0.56
1:B:145:GLU:N	1:B:145:GLU:CD	2.58	0.56
1:A:367:ASN:HB2	1:C:4:MET:HG2	1.87	0.56
1:A:197:ASN:ND2	1:A:288:GLN:HG2	2.19	0.56
1:C:149:GLY:HA2	1:C:222:LYS:HD2	1.87	0.56

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:175:GLU:HA	1:C:184:ASP:O	2.06	0.56
1:A:223:PRO:HA	1:A:241:ALA:O	2.06	0.56
1:A:343:ASN:O	1:A:347:GLU:HG3	2.05	0.56
1:A:268:GLU:O	1:A:271:THR:HB	2.06	0.56
1:A:391:THR:O	1:A:394:ALA:HB3	2.06	0.56
1:B:261:MET:CE	1:B:266:LYS:HG2	2.36	0.55
1:A:178:TYR:O	1:A:179:LYS:HB2	2.07	0.55
1:B:178:TYR:CG	1:B:179:LYS:N	2.73	0.55
1:A:164:PRO:HG3	1:A:252:PHE:CD1	2.41	0.55
1:B:392:GLN:HA	1:B:392:GLN:OE1	2.06	0.55
1:C:393:LEU:O	1:C:393:LEU:HD22	2.06	0.55
1:B:223:PRO:HG3	1:B:241:ALA:C	2.27	0.55
1:C:223:PRO:HG3	1:C:242:GLU:CA	2.35	0.55
1:C:286:TYR:CZ	1:C:331:LYS:HG2	2.41	0.55
1:A:211:MET:CE	1:A:249:LEU:HB2	2.37	0.55
1:A:237:ILE:HD13	1:A:243:LEU:HD21	1.90	0.54
1:B:171:GLU:CD	1:B:189:ARG:HE	2.10	0.54
1:B:380:LEU:CD1	1:B:389:ALA:HB1	2.37	0.54
1:A:218:ILE:HB	3:A:466:HOH:O	2.08	0.54
1:B:357:PHE:CZ	1:B:386:ASN:OD1	2.60	0.54
1:B:269:GLN:HA	1:B:269:GLN:OE1	2.08	0.54
1:C:160:GLY:O	1:C:161:TYR:HB2	2.08	0.54
1:C:304:SER:C	1:C:306:PHE:H	2.10	0.54
1:A:163:LYS:HG2	1:A:211:MET:O	2.08	0.54
1:B:309:GLU:O	1:B:313:LYS:HD2	2.08	0.53
1:B:357:PHE:CZ	1:B:388:ALA:HB3	2.43	0.53
1:B:157:ARG:HD2	3:B:491:HOH:O	2.08	0.53
1:C:378:LYS:O	1:C:382:LEU:HD13	2.09	0.53
1:B:164:PRO:HB3	1:B:252:PHE:CD1	2.44	0.53
1:C:407:GLU:HG2	1:C:411:TYR:CE1	2.42	0.53
1:B:361:GLU:OE1	1:B:392:GLN:HG3	2.09	0.53
1:C:425:ALA:HA	3:C:470:HOH:O	2.08	0.53
1:C:177:TYR:HA	1:C:181:LYS:O	2.08	0.53
1:B:4:MET:HE1	1:B:210:ARG:CZ	2.39	0.53
1:B:155:GLN:O	1:B:156:THR:HB	2.08	0.53
1:A:172:VAL:HG22	1:A:249:LEU:HA	1.91	0.53
1:C:213:LYS:HG3	1:C:214:GLY:N	2.24	0.53
1:C:170:VAL:CG1	1:C:172:VAL:HG23	2.34	0.52
1:B:261:MET:HE3	1:B:266:LYS:HG2	1.91	0.52
1:B:151:ILE:HD12	1:B:151:ILE:N	2.24	0.52
1:A:285:LYS:HB3	1:A:288:GLN:HE21	1.73	0.52
1:A:393:LEU:HD23	1:A:393:LEU:O	2.10	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:160:GLY:O	1:C:161:TYR:CB	2.58	0.52
1:B:276:ARG:HD3	3:B:499:HOH:O	2.08	0.52
1:A:287:LYS:O	1:A:290:LEU:HB2	2.09	0.52
1:A:387:LYS:HD3	1:A:387:LYS:N	2.15	0.52
1:C:286:TYR:CE2	1:C:331:LYS:HG2	2.45	0.52
1:C:366:VAL:HG12	1:C:366:VAL:O	2.10	0.52
1:A:307:SER:OG	1:A:310:GLU:HG3	2.10	0.52
1:A:172:VAL:CG2	1:A:249:LEU:HA	2.40	0.51
1:C:179:LYS:O	1:C:180:ASP:HB3	2.10	0.51
1:A:402:ARG:HD3	3:C:467:HOH:O	2.09	0.51
1:C:304:SER:O	1:C:306:PHE:N	2.40	0.51
1:A:351:ASN:OD1	1:A:382:LEU:HD21	2.10	0.51
1:A:177:TYR:HD1	1:A:244:LYS:HD3	1.74	0.51
1:A:261:MET:HE3	1:A:266:LYS:HA	1.92	0.51
1:B:307:SER:OG	1:B:310:GLU:HG3	2.10	0.51
1:C:307:SER:O	1:C:308:ASN:C	2.48	0.51
1:B:4:MET:CE	1:B:210:ARG:CZ	2.88	0.51
1:A:313:LYS:HB2	1:A:313:LYS:NZ	2.24	0.51
1:B:174:LEU:HD23	1:B:247:LEU:HD21	1.92	0.51
1:A:380:LEU:CD2	1:A:389:ALA:HB1	2.39	0.51
1:B:165:ASN:HB3	1:B:254:LYS:NZ	2.26	0.51
1:A:211:MET:HB2	1:A:249:LEU:HD22	1.92	0.51
1:C:165:ASN:HD21	1:C:255:ALA:HB3	1.76	0.51
1:A:353:GLU:O	1:A:357:PHE:HB2	2.11	0.51
1:C:160:GLY:HA2	1:C:212:GLU:HB3	1.93	0.51
1:C:223:PRO:HB3	1:C:241:ALA:N	2.26	0.51
1:A:145:GLU:C	1:A:147:ASP:H	2.13	0.51
1:B:303:GLU:OE2	1:B:305:SER:OG	2.27	0.51
1:A:302:TYR:O	1:A:303:GLU:C	2.48	0.51
1:B:304:SER:O	1:C:373:ARG:NH1	2.43	0.50
1:B:223:PRO:HG3	1:B:242:GLU:N	2.26	0.50
1:A:226:ALA:HB1	1:A:243:LEU:CD1	2.39	0.50
1:B:391:THR:O	1:B:395:VAL:HG23	2.12	0.50
1:B:346:LEU:HD11	1:B:356:LEU:HD23	1.93	0.50
1:A:178:TYR:CE1	1:A:241:ALA:HB1	2.47	0.50
1:C:396:CYS:O	1:C:400:ILE:HG13	2.11	0.49
1:A:392:GLN:HA	1:A:392:GLN:OE1	2.10	0.49
1:B:239:PRO:O	1:B:240:ASN:HB3	2.12	0.49
1:A:161:TYR:CD1	1:A:161:TYR:N	2.79	0.49
1:A:169:ILE:HG21	1:A:189:ARG:NH2	2.26	0.49
1:B:410:LEU:C	1:B:410:LEU:HD12	2.30	0.49
1:B:181:LYS:HE3	3:B:500:HOH:O	2.13	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:395:VAL:O	1:B:399:ARG:HG2	2.13	0.49
1:A:201:PRO:HG2	1:A:204:LEU:HB2	1.94	0.49
1:B:223:PRO:HB3	1:B:241:ALA:N	2.27	0.49
1:C:399:ARG:O	1:C:403:GLN:HG3	2.12	0.49
1:A:267:LEU:HD21	1:A:314:ALA:HB2	1.95	0.49
1:A:349:ASP:CB	1:A:352:ASN:HB2	2.43	0.48
1:C:308:ASN:O	1:C:312:GLN:HG2	2.13	0.48
1:B:397:GLN:HA	1:B:397:GLN:OE1	2.13	0.48
1:A:177:TYR:HE2	1:A:182:LEU:HB2	1.77	0.48
1:B:364:LEU:O	1:B:364:LEU:HD12	2.13	0.48
1:C:401:ARG:HH12	1:C:402:ARG:NH2	2.11	0.48
1:A:191:GLU:O	1:A:194:GLU:HB3	2.12	0.48
1:B:156:THR:O	1:B:158:GLY:N	2.46	0.48
1:A:153:ARG:HH11	1:A:153:ARG:HG3	1.78	0.48
1:A:166:GLU:O	1:A:255:ALA:HB3	2.14	0.48
1:A:349:ASP:C	1:A:351:ASN:N	2.67	0.48
1:B:3:HIS:C	1:B:4:MET:HG2	2.33	0.48
1:C:254:LYS:HD2	3:C:491:HOH:O	2.13	0.48
1:B:1:GLY:O	1:B:2:SER:CB	2.62	0.48
1:A:179:LYS:O	1:A:180:ASP:HB2	2.14	0.48
1:C:228:GLY:O	1:C:229:SER:HB2	2.14	0.48
1:C:397:GLN:HG3	3:C:498:HOH:O	2.13	0.47
1:A:356:LEU:HA	1:A:356:LEU:HD12	1.70	0.47
1:B:290:LEU:O	1:B:294:LYS:HG3	2.14	0.47
1:A:4:MET:HG2	1:A:210:ARG:CZ	2.44	0.47
1:A:322:HIS:CE1	1:A:344:LYS:HB3	2.49	0.47
1:A:363:HIS:ND1	1:A:371:LEU:HB3	2.29	0.47
1:A:233:GLU:HG3	1:C:238:PRO:HG2	1.94	0.47
1:A:397:GLN:OE1	1:A:397:GLN:HA	2.14	0.47
1:A:357:PHE:HZ	1:A:388:ALA:CB	2.28	0.47
1:A:261:MET:HB2	1:A:265:GLU:HB2	1.97	0.47
1:A:151:ILE:O	1:A:219:VAL:HA	2.15	0.47
1:A:226:ALA:O	1:A:228:GLY:N	2.48	0.47
1:C:4:MET:SD	1:C:210:ARG:CZ	3.03	0.47
1:A:274:LYS:HG3	1:A:293:TYR:HE2	1.80	0.47
1:A:169:ILE:O	1:A:252:PHE:HA	2.16	0.46
1:B:157:ARG:H	1:B:157:ARG:HD3	1.79	0.46
1:C:221:LEU:HD13	1:C:226:ALA:HB2	1.98	0.46
1:A:264:GLU:HA	1:A:267:LEU:HD12	1.96	0.46
1:C:262:ASN:OD1	1:C:265:GLU:HG3	2.15	0.46
1:A:304:SER:C	1:A:306:PHE:H	2.18	0.46
1:A:215:GLU:HG2	1:A:216:HIS:N	2.30	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:226:ALA:HB1	1:B:243:LEU:CD1	2.40	0.46
1:A:315:GLN:HG3	3:A:468:HOH:O	2.14	0.46
1:C:177:TYR:CE2	1:C:182:LEU:HB2	2.51	0.46
1:A:285:LYS:HB3	1:A:288:GLN:NE2	2.30	0.46
1:A:221:LEU:HD13	1:A:226:ALA:HB2	1.96	0.46
1:A:178:TYR:O	1:A:180:ASP:N	2.44	0.46
1:C:399:ARG:HB3	1:C:399:ARG:NH1	2.31	0.46
1:B:170:VAL:HG22	1:B:249:LEU:CD1	2.45	0.46
1:B:257:GLU:O	1:B:258:SER:C	2.53	0.46
1:B:208:ILE:C	1:B:210:ARG:H	2.18	0.46
1:C:395:VAL:O	1:C:399:ARG:HG2	2.16	0.46
1:C:274:LYS:O	1:C:278:THR:HG23	2.16	0.46
1:A:265:GLU:HB3	3:A:509:HOH:O	2.15	0.46
2:H:2:GLU:HG2	2:H:3:GLU:OE1	2.15	0.46
1:A:185:GLN:CD	1:A:185:GLN:O	2.54	0.46
1:C:177:TYR:HE2	1:C:182:LEU:HB2	1.81	0.46
1:B:380:LEU:HD11	1:B:389:ALA:HB1	1.97	0.46
1:A:226:ALA:C	1:A:228:GLY:N	2.67	0.45
1:A:285:LYS:HD3	1:A:288:GLN:NE2	2.31	0.45
1:A:198:LEU:O	1:A:200:LEU:HG	2.16	0.45
1:C:317:LEU:HA	1:C:317:LEU:HD12	1.85	0.45
1:A:165:ASN:HB3	1:A:254:LYS:HE2	1.97	0.45
1:C:420:GLU:HA	1:C:423:ASN:HD22	1.82	0.45
1:A:4:MET:HE3	3:A:471:HOH:O	2.16	0.45
1:C:237:ILE:HD13	1:C:243:LEU:HD21	1.99	0.45
1:A:175:GLU:OE1	1:A:185:GLN:HA	2.16	0.45
1:B:223:PRO:HB3	1:B:240:ASN:HA	1.97	0.45
1:B:324:ASN:O	1:B:327:MET:HB3	2.16	0.45
1:B:357:PHE:HZ	1:B:388:ALA:HB3	1.82	0.45
1:A:408:LYS:HA	1:B:418:LEU:HD13	1.99	0.45
1:A:352:ASN:O	1:A:356:LEU:HB2	2.16	0.45
1:B:4:MET:HE1	1:B:210:ARG:NH2	2.32	0.45
1:A:261:MET:CE	1:A:266:LYS:HA	2.46	0.45
1:A:232:LYS:HE3	1:A:234:LYS:HE3	1.99	0.45
1:C:382:LEU:N	1:C:382:LEU:HD12	2.32	0.45
1:A:317:LEU:HA	1:A:317:LEU:HD12	1.85	0.45
1:A:208:ILE:HA	1:A:211:MET:HG3	1.99	0.44
1:C:146:GLU:OE2	1:C:210:ARG:NH2	2.50	0.44
1:B:227:PHE:CD1	1:B:232:LYS:HB2	2.52	0.44
1:A:201:PRO:O	1:A:202:TYR:C	2.56	0.44
1:C:357:PHE:HZ	1:C:388:ALA:CB	2.20	0.44
1:B:172:VAL:HG12	1:B:173:ALA:N	2.32	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:376:PHE:CE1	1:A:392:GLN:HB3	2.52	0.44
1:A:333:GLN:NE2	1:C:151:ILE:CD1	2.80	0.44
1:C:178:TYR:OH	1:C:179:LYS:HE3	2.17	0.44
1:B:344:LYS:HD3	1:B:344:LYS:HA	1.76	0.44
1:C:152:ARG:HG2	1:C:152:ARG:NH1	2.30	0.44
1:C:190:PHE:HB2	1:C:198:LEU:CD1	2.48	0.44
1:A:146:GLU:HA	3:A:473:HOH:O	2.17	0.44
1:A:233:GLU:O	1:A:234:LYS:HG2	2.18	0.44
1:B:306:PHE:CE1	1:B:314:ALA:HB2	2.53	0.44
1:B:373:ARG:HH21	1:C:304:SER:CB	2.31	0.44
1:B:367:ASN:ND2	1:B:369:PHE:CZ	2.85	0.44
1:C:366:VAL:O	1:C:367:ASN:HB2	2.17	0.44
1:A:201:PRO:CB	1:A:226:ALA:HA	2.48	0.43
1:B:156:THR:CG2	1:B:156:THR:O	2.66	0.43
1:C:226:ALA:O	1:C:227:PHE:HB2	2.17	0.43
1:C:351:ASN:HB3	1:C:382:LEU:HD23	1.99	0.43
1:C:401:ARG:HH12	1:C:402:ARG:HH21	1.64	0.43
1:C:392:GLN:C	1:C:394:ALA:N	2.71	0.43
1:A:178:TYR:CZ	1:A:179:LYS:HE2	2.53	0.43
1:A:230:VAL:HG12	1:A:230:VAL:O	2.19	0.43
1:A:233:GLU:C	1:A:234:LYS:HG2	2.38	0.43
1:C:261:MET:O	1:C:266:LYS:HE3	2.18	0.43
1:A:420:GLU:HG3	1:A:421:GLU:N	2.33	0.43
1:A:208:ILE:C	1:A:210:ARG:H	2.21	0.43
1:B:166:GLU:C	1:B:168:ALA:H	2.19	0.43
1:C:233:GLU:O	1:C:236:GLN:N	2.52	0.43
1:B:379:VAL:O	1:B:383:TYR:N	2.40	0.43
1:A:381:GLN:HG2	1:A:382:LEU:HD12	2.01	0.43
1:B:156:THR:HG21	3:B:516:HOH:O	2.18	0.43
1:C:178:TYR:CE1	1:C:241:ALA:HB1	2.43	0.43
1:A:323:LEU:O	1:A:358:ARG:NH1	2.51	0.43
1:A:158:GLY:O	1:A:159:GLU:C	2.56	0.43
1:B:350:SER:H	1:C:350:SER:CB	2.20	0.43
1:B:172:VAL:O	1:B:187:GLU:HA	2.18	0.43
1:B:155:GLN:O	1:B:156:THR:CB	2.67	0.43
1:C:165:ASN:CG	1:C:254:LYS:HD3	2.39	0.43
1:C:368:ASP:CB	1:C:371:LEU:HD12	2.46	0.43
1:B:4:MET:CE	1:B:210:ARG:NH2	2.82	0.43
1:A:290:LEU:HD23	1:A:290:LEU:HA	1.85	0.43
1:B:349:ASP:HB3	1:B:352:ASN:HB2	2.01	0.43
1:A:156:THR:HG22	1:A:157:ARG:O	2.19	0.43
1:B:223:PRO:HB3	1:B:241:ALA:H	1.84	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:350:SER:N	1:C:350:SER:HB2	2.19	0.43
1:C:3:HIS:HA	1:C:4:MET:HE3	2.00	0.43
1:A:216:HIS:O	1:A:217:SER:HB3	2.18	0.43
1:A:220:TYR:HE1	3:A:466:HOH:O	2.01	0.43
1:A:152:ARG:NH1	1:A:215:GLU:OE2	2.50	0.42
1:C:164:PRO:HB2	1:C:209:GLN:HA	2.00	0.42
1:A:403:GLN:NE2	1:C:3:HIS:NE2	2.67	0.42
1:B:178:TYR:C	1:B:180:ASP:N	2.72	0.42
1:C:346:LEU:HD12	1:C:346:LEU:HA	1.74	0.42
1:C:413:ASN:O	1:C:416:GLU:HB3	2.19	0.42
1:A:275:GLU:O	1:A:278:THR:HB	2.19	0.42
1:B:155:GLN:HG2	1:B:155:GLN:O	2.19	0.42
1:B:290:LEU:HD11	1:B:332:LEU:HD12	2.01	0.42
1:C:318:ARG:O	1:C:321:SER:HB2	2.19	0.42
1:A:321:SER:O	1:A:325:LEU:HD23	2.18	0.42
1:C:251:SER:O	1:C:252:PHE:HB3	2.19	0.42
1:A:2:SER:O	1:A:3:HIS:O	2.38	0.42
1:A:213:LYS:HB3	1:A:214:GLY:H	1.63	0.42
1:A:154:ILE:HG22	1:A:154:ILE:O	2.20	0.42
1:B:152:ARG:HA	1:B:218:ILE:O	2.20	0.42
1:B:392:GLN:O	1:B:393:LEU:C	2.58	0.42
1:A:393:LEU:O	1:A:397:GLN:HG2	2.20	0.42
1:B:176:GLY:HA2	1:B:244:LYS:O	2.19	0.42
1:A:185:GLN:HG3	3:A:518:HOH:O	2.19	0.42
1:A:211:MET:HB2	1:A:249:LEU:CD2	2.50	0.42
1:A:4:MET:CG	1:A:4:MET:CE	2.94	0.42
1:C:155:GLN:HG2	1:C:218:ILE:HG13	2.02	0.42
1:C:158:GLY:O	1:C:159:GLU:O	2.37	0.42
1:B:156:THR:O	1:B:157:ARG:C	2.58	0.42
1:A:251:SER:O	1:A:252:PHE:CB	2.65	0.42
1:B:380:LEU:HD12	1:B:389:ALA:HB1	2.01	0.42
1:B:232:LYS:HD2	1:B:232:LYS:HA	1.86	0.42
1:A:415:PHE:CE1	1:B:408:LYS:HG3	2.55	0.42
2:G:1:MET:HA	3:G:24:HOH:O	2.20	0.41
1:C:369:PHE:O	1:C:370:GLU:C	2.57	0.41
2:H:4:VAL:O	2:H:5:ASP:CB	2.68	0.41
1:B:420:GLU:HA	1:B:423:ASN:HD22	1.86	0.41
1:B:163:LYS:HA	1:B:164:PRO:HD2	1.96	0.41
1:B:195:GLY:N	1:B:205:GLU:OE2	2.50	0.41
1:A:3:HIS:O	1:A:4:MET:C	2.58	0.41
1:B:170:VAL:HG13	1:B:172:VAL:HG23	2.02	0.41
1:C:211:MET:HE3	1:C:249:LEU:HB2	2.01	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:161:TYR:HB3	1:C:162:ALA:H	1.57	0.41
1:B:373:ARG:NE	1:C:304:SER:HB2	2.33	0.41
1:C:163:LYS:HA	1:C:252:PHE:CZ	2.56	0.41
1:C:180:ASP:OD1	1:C:180:ASP:N	2.54	0.41
1:C:208:ILE:HG12	1:C:247:LEU:CD1	2.49	0.41
1:B:304:SER:HB2	1:C:374:ALA:HB2	2.01	0.41
1:C:163:LYS:HB2	1:C:164:PRO:HD2	2.03	0.41
1:B:330:LEU:HD23	1:B:330:LEU:HA	1.92	0.41
1:C:189:ARG:NH1	1:C:189:ARG:HG3	2.35	0.41
1:A:149:GLY:HA2	1:A:222:LYS:CG	2.51	0.41
1:C:287:LYS:HD3	3:C:471:HOH:O	2.21	0.41
1:B:174:LEU:HD23	1:B:247:LEU:HD22	2.03	0.41
1:B:162:ALA:HB1	1:B:252:PHE:CZ	2.56	0.41
1:A:151:ILE:HA	3:A:463:HOH:O	2.21	0.41
1:C:351:ASN:OD1	1:C:382:LEU:HD21	2.21	0.41
1:B:421:GLU:OE1	1:B:421:GLU:HA	2.20	0.41
1:B:179:LYS:O	1:B:180:ASP:CB	2.64	0.40
1:C:189:ARG:HH11	1:C:189:ARG:HG3	1.86	0.40
1:C:183:PHE:HE2	1:C:243:LEU:HD22	1.87	0.40
1:C:223:PRO:HD3	1:C:242:GLU:HB3	2.02	0.40
1:A:361:GLU:OE2	1:A:392:GLN:HG3	2.21	0.40
1:A:415:PHE:HE1	1:B:408:LYS:HG2	1.86	0.40
1:A:225:TYR:O	1:A:226:ALA:HB2	2.22	0.40
1:A:238:PRO:HA	1:A:239:PRO:HD3	1.87	0.40
1:A:294:LYS:O	1:A:295:LYS:C	2.60	0.40
1:A:188:LEU:HD13	1:A:198:LEU:HD13	2.03	0.40
1:B:172:VAL:CG1	1:B:173:ALA:N	2.85	0.40
1:C:183:PHE:CE2	1:C:243:LEU:HD22	2.56	0.40
1:A:290:LEU:CD2	1:A:325:LEU:HD12	2.52	0.40
1:B:233:GLU:C	1:B:236:GLN:H	2.25	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	283/336 (84%)	225 (80%)	39 (14%)	19 (7%)	2	10
1	B	283/336 (84%)	229 (81%)	34 (12%)	20 (7%)	2	9
1	C	283/336 (84%)	242 (86%)	23 (8%)	18 (6%)	2	11
2	G	3/5 (60%)	1 (33%)	2 (67%)	0	100	100
2	H	3/5 (60%)	1 (33%)	1 (33%)	1 (33%)	0	0
All	All	855/1018 (84%)	698 (82%)	99 (12%)	58 (7%)	2	10

All (58) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	3	HIS
1	A	180	ASP
1	A	213	LYS
1	A	234	LYS
1	A	252	PHE
1	A	350	SER
1	B	2	SER
1	B	157	ARG
1	B	162	ALA
1	B	166	GLU
1	B	167	GLY
1	B	175	GLU
1	B	178	TYR
1	B	180	ASP
1	C	4	MET
1	C	159	GLU
1	C	161	TYR
1	C	180	ASP
1	C	213	LYS
1	C	224	SER
1	C	308	ASN
1	C	387	LYS
1	A	4	MET
1	A	387	LYS
1	B	147	ASP
1	B	159	GLU
1	B	160	GLY
1	B	258	SER
1	C	305	SER
1	C	306	PHE
1	C	307	SER

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Mol	Chain	Res	Type
2	H	4	VAL
1	A	157	ARG
1	A	159	GLU
1	A	250	LYS
1	A	298	SER
1	A	303	GLU
1	B	161	TYR
1	B	254	LYS
1	A	160	GLY
1	A	226	ALA
1	A	232	LYS
1	B	163	LYS
1	B	182	LEU
1	B	350	SER
1	C	154	ILE
1	C	226	ALA
1	C	350	SER
1	B	230	VAL
1	B	305	SER
1	A	217	SER
1	C	239	PRO
1	A	228	GLY
1	C	164	PRO
1	A	154	ILE
1	B	214	GLY
1	C	223	PRO
1	C	230	VAL

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	238/283 (84%)	222 (93%)	16 (7%)	23	64
1	B	238/283 (84%)	220 (92%)	18 (8%)	19	57
1	C	239/283 (84%)	224 (94%)	15 (6%)	25	66
2	G	5/5 (100%)	3 (60%)	2 (40%)	0	0

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	H	5/5 (100%)	5 (100%)	0	100	100
All	All	725/859 (84%)	674 (93%)	51 (7%)	21	62

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	146	GLU
1	A	161	TYR
1	A	172	VAL
1	A	185	GLN
1	A	191	GLU
1	A	211	MET
1	A	235	PHE
1	A	247	LEU
1	A	257	GLU
1	A	261	MET
1	A	269	GLN
1	A	317	LEU
1	A	323	LEU
1	A	381	GLN
1	A	387	LYS
1	A	417	ARG
1	B	145	GLU
1	B	151	ILE
1	B	157	ARG
1	B	170	VAL
1	B	202	TYR
1	B	215	GLU
1	B	268	GLU
1	B	291	LEU
1	B	300	LEU
1	B	317	LEU
1	B	325	LEU
1	B	340	GLU
1	B	357	PHE
1	B	367	ASN
1	B	368	ASP
1	B	393	LEU
1	B	398	GLN
1	B	414	MET
1	C	4	MET
1	C	147	ASP

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Mol	Chain	Res	Type
1	C	151	ILE
1	C	161	TYR
1	C	211	MET
1	C	239	PRO
1	C	300	LEU
1	C	317	LEU
1	C	323	LEU
1	C	346	LEU
1	C	367	ASN
1	C	393	LEU
1	C	402	ARG
1	C	414	MET
1	C	417	ARG
2	G	2	GLU
2	G	3	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	185	GLN
1	A	197	ASN
1	A	248	HIS
1	A	269	GLN
1	A	288	GLN
1	A	324	ASN
1	A	333	GLN
1	A	343	ASN
1	A	367	ASN
1	A	381	GLN
1	A	403	GLN
1	B	315	GLN
1	B	367	ASN
1	B	398	GLN
1	B	423	ASN
1	C	209	GLN
1	C	248	HIS
1	C	308	ASN
1	C	315	GLN
1	C	343	ASN
1	C	381	GLN
1	C	392	GLN
1	C	403	GLN

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Mol	Chain	Res	Type
1	C	423	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	285/336 (84%)	-0.18	0 100 100	16, 51, 96, 130	0
1	B	285/336 (84%)	-0.24	0 100 100	15, 50, 91, 121	0
1	C	285/336 (84%)	-0.21	0 100 100	10, 44, 92, 125	0
2	G	5/5 (100%)	0.37	0 100 100	81, 86, 92, 93	0
2	H	5/5 (100%)	0.56	0 100 100	73, 82, 91, 95	0
All	All	865/1018 (84%)	-0.21	0 100 100	10, 50, 93, 130	0

There are no RSRZ outliers to report.

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.