Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2024 – 12:07 AM EST

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 X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

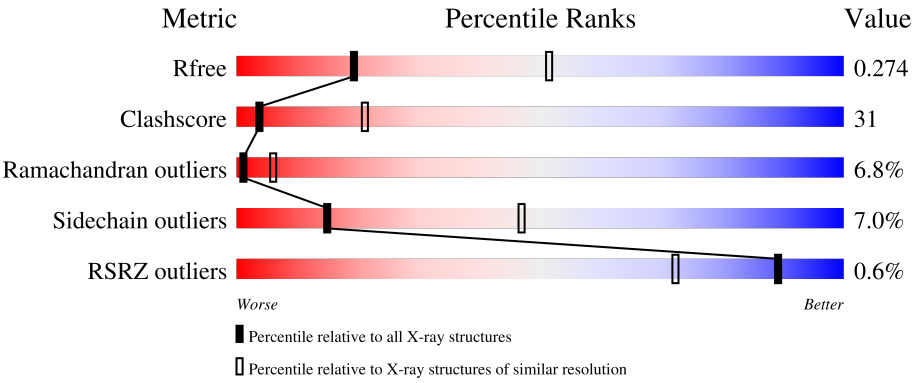
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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}	130704	2092 (3.00-3.00)
Clashscore	141614	2416 (3.00-3.00)
Ramachandran outliers	138981	2333 (3.00-3.00)
Sidechain outliers	138945	2336 (3.00-3.00)
RSRZ outliers	127900	1990 (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 of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=5% The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	336	% 37% 40% 8% • 15%
1	B	336	% 41% 38% 6% 15%
1	C	336	44% 34% 7% • 15%
2	G	5	20% 40% 40% 20%
2	H	5	20% 60% 20%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 7225 atoms, of which 0 are hydrogens and 0 are deuteriums.

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

- Molecule 1 is a protein called 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	C	N	O	S	0	0	0
			42	24	5	12	1			
2	H	5	Total	C	N	O	S	0	0	0
			42	24	5	12	1			

- Molecule 3 is water.

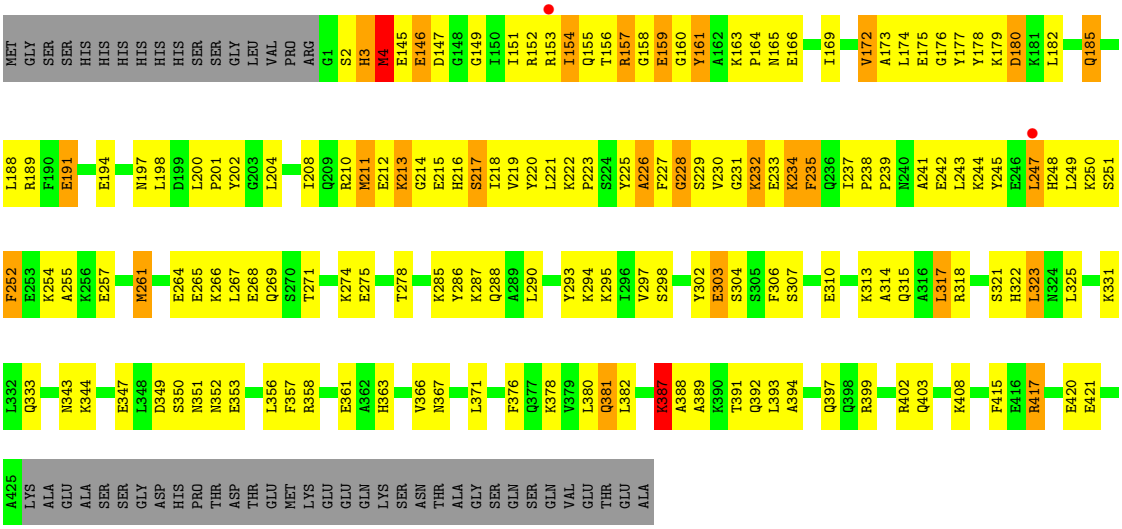
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	75	Total	O	0	0
			75	75		
3	B	64	Total	O	0	0
			64	64		
3	C	76	Total	O	0	0
			76	76		
3	G	4	Total	O	0	0
			4	4		
3	H	2	Total	O	0	0
			2	2		

3 Residue-property plots (i)

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

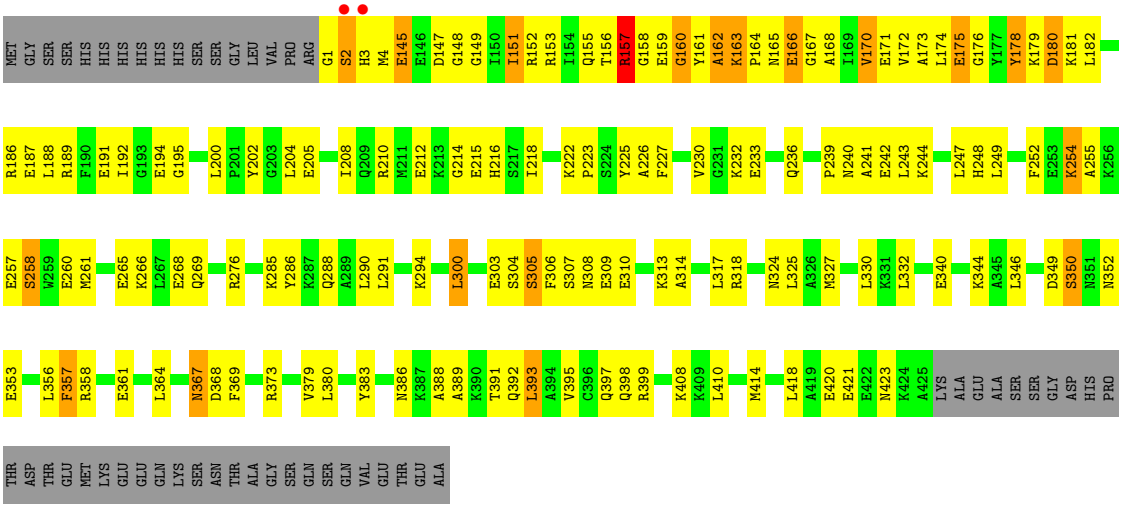
• Molecule 1: FK506-binding protein 4

Chain A:



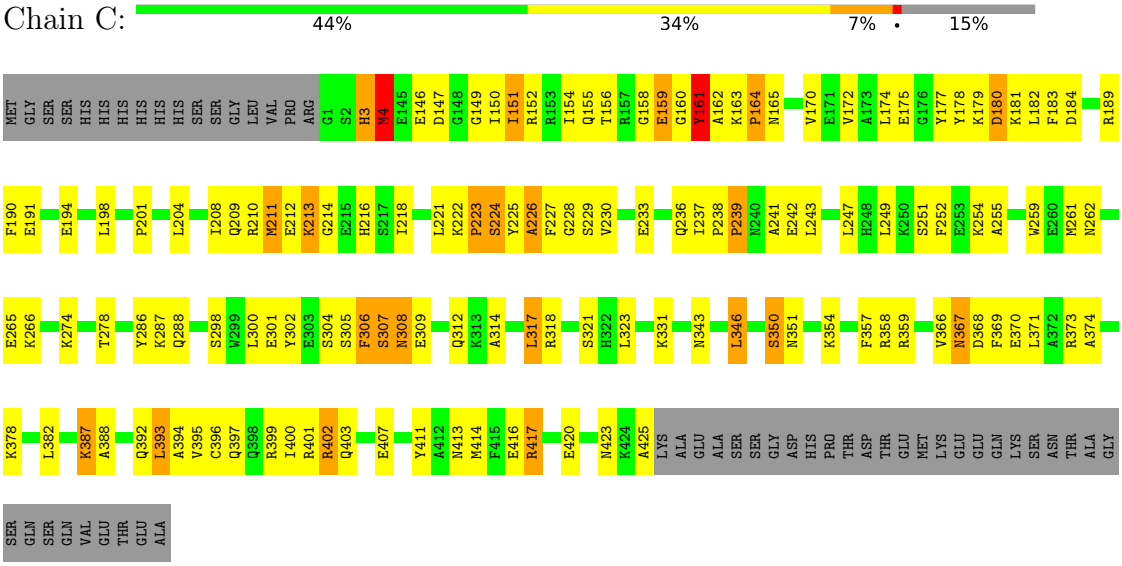
• Molecule 1: FK506-binding protein 4

Chain B:



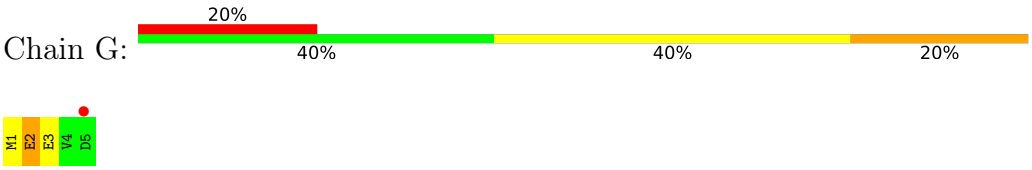
• Molecule 1: FK506-binding protein 4

Chain C:



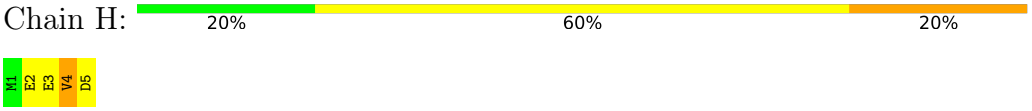
• Molecule 2: 5-mer peptide from Heat shock protein HSP 90

Chain G:



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

Chain H:



4 Data and refinement statistics (i)

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
$< I/\sigma(I) >$ [1]	2.15 (at 2.91Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.230 , 0.287 0.224 , 0.274	Depositor DCC
R_{free} test set	1344 reflections (4.35%)	wwPDB-VP
Wilson B-factor (Å^2)	51.5	Xtriage
Anisotropy	0.543	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 67.8	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	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.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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(^{o})	Ideal(^{o})
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 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2305	0	2296	171	0
1	B	2305	0	2296	143	0
1	C	2310	0	2301	135	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
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

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 31.

All (434) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:HB3	1:C:417:ARG:HH11	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:H	1:A:387:LYS:HD3	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:HD3	1:B:157:ARG:N	1.95	0.81
1:B:145:GLU:CD	1:B:145:GLU:H	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	Interatomic distance (Å)	Clash overlap (Å)
1:A:380:LEU:HD22	1:A:389:ALA:HB1	1.65	0.77
1:B:170:VAL:HG23	1:B:252:PHE:CB	2.14	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:HD12	1:A:154:ILE:H	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:O	1:B:410:LEU:HD12	1.89	0.71
1:A:227:PHE:HB3	1:A:231:GLY:O	1.91	0.71
1:C:182:LEU:HD23	1:C:182:LEU: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:194:GLU:HG3	1:C:288:GLN:HG2	1.72	0.70
1:C:201:PRO:HB3	1:C:226:ALA:HA	1.74	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:H	1:A:387:LYS:CD	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:226:ALA:C	1:A:228:GLY:H	2.00	0.65
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: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:204:LEU:HD11	1:C:247:LEU:HD11	1.79	0.65
1:A:251:SER:O	1:A:252:PHE:HB3	1.95	0.64
1:C:343:ASN:ND2	1:C:359:ARG:HH11	1.96	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:A:247:LEU:HD23	1:A:247:LEU:H	1.62	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: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:173:ALA:CB	1:B:248:HIS:HB3	2.28	0.62
1:B:174:LEU:HD11	1:B:200:LEU:HD21	1.82	0.62
1:C:223:PRO:O	1:C:224:SER:HB2	1.98	0.62
1:A:229:SER:O	1:A:239:PRO:HB3	2.00	0.61
1:B:261:MET:HB3	1:B:265:GLU:HB3	1.82	0.61
1:A:177:TYR:CE2	1:A:182:LEU:HB2	2.35	0.61
1:B:353:GLU:HB2	1:B:383:TYR:CE1	2.36	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:HH11	1:C:152:ARG:HG2	1.65	0.61
1:C:150:ILE:HD11	1:C:225:TYR:HD2	1.65	0.61
1:C:314:ALA:O	1:C:318:ARG:HG3	2.00	0.61
1:B:151:ILE:HD12	1:B:151:ILE:H	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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: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:A:318:ARG:HD3	3:A:476:HOH:O	2.02	0.60
1:B:174:LEU:HD12	1:B:186:ARG:HG3	1.83	0.60
1:B:1:GLY:O	1:B:2:SER:HB3	2.02	0.60
1:B:162:ALA:O	1:B:163:LYS:HB2	2.01	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:B:153:ARG:HB2	1:B:218:ILE:HG22	1.86	0.58
1:C:174:LEU:C	1:C:174:LEU:HD12	2.23	0.58
1:A:286:TYR:CE2	1:A:331:LYS:HG2	2.40	0.57
1:C:204:LEU:CD1	1:C:247:LEU:HD11	2.33	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:149:GLY:HA2	1:B:222:LYS:HD2	1.86	0.57
1:B:157:ARG:N	1:B:157:ARG:CD	2.68	0.57
1:B:174:LEU:HD23	1:B:247:LEU:CD2	2.35	0.57
1:B:165:ASN:HD21	1:B:255:ALA:HB3	1.68	0.56
1:B:286:TYR:CB	1:B:332:LEU:HD21	2.35	0.56
1:A:367:ASN:HB2	1:C:4:MET:HG2	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:145:GLU:CD	1:B:145:GLU:N	2.58	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
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:A:164:PRO:HG3	1:A:252:PHE:CD1	2.41	0.55
1:B:178:TYR:CG	1:B:179:LYS:N	2.73	0.55
1:B:392:GLN:OE1	1:B:392:GLN:HA	2.06	0.55
1:C:393:LEU:HD22	1:C:393:LEU:O	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:269:GLN:OE1	1:B:269:GLN:HA	2.08	0.54
1:B:357:PHE:CZ	1:B:386:ASN:OD1	2.60	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:B:361:GLU:OE1	1:B:392:GLN:HG3	2.09	0.53
1:C:407:GLU:HG2	1:C:411:TYR:CE1	2.42	0.53
1:C:425:ALA:HA	3:C:470:HOH:O	2.08	0.53
1:B:4:MET:HE1	1:B:210:ARG:CZ	2.39	0.53
1:C:177:TYR:HA	1:C:181:LYS:O	2.08	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:285:LYS:HB3	1:A:288:GLN:HE21	1.73	0.52
1:A:393:LEU:O	1:A:393:LEU:HD23	2.10	0.52
1:B:151:ILE:HD12	1:B:151:ILE:N	2.24	0.52
1:B:261:MET:HE3	1:B:266:LYS:HG2	1.91	0.52
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:A:307:SER:OG	1:A:310:GLU:HG3	2.10	0.52
1:C:366:VAL:HG12	1:C:366:VAL:O	2.10	0.52
1:A:172:VAL:CG2	1:A:249:LEU:HA	2.40	0.51
1:A:402:ARG:HD3	3:C:467:HOH:O	2.09	0.51
1:C:179:LYS:O	1:C:180:ASP:HB3	2.10	0.51
1:C:304:SER:O	1:C:306:PHE:N	2.40	0.51
1:A:177:TYR:HD1	1:A:244:LYS:HD3	1.74	0.51
1:A:351:ASN:OD1	1:A:382:LEU:HD21	2.10	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:A:313:LYS:HB2	1:A:313:LYS:NZ	2.24	0.51
1:B:4:MET:CE	1:B:210:ARG:CZ	2.88	0.51
1:C:307:SER:O	1:C:308:ASN:C	2.48	0.51
1:A:380:LEU:CD2	1:A:389:ALA:HB1	2.39	0.51
1:B:174:LEU:HD23	1:B:247:LEU:HD21	1.92	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: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:165:ASN:HD21	1:C:255:ALA:HB3	1.76	0.51
1:A:145:GLU:C	1:A:147:ASP:H	2.13	0.51
1:C:223:PRO:HB3	1:C:241:ALA:N	2.26	0.51
1:A:302:TYR:O	1:A:303:GLU:C	2.48	0.51
1:B:303:GLU:OE2	1:B:305:SER:OG	2.27	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:346:LEU:HD11	1:B:356:LEU:HD23	1.93	0.50
1:B:391:THR:O	1:B:395:VAL:HG23	2.12	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:161:TYR:N	1:A:161:TYR:CD1	2.79	0.49
1:B:239:PRO:O	1:B:240:ASN:HB3	2.12	0.49
1:A:169:ILE:HG21	1:A:189:ARG:NH2	2.26	0.49
1:B:410:LEU:HD12	1:B:410:LEU:C	2.30	0.49
1:B:181:LYS:HE3	3:B:500:HOH:O	2.13	0.49
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:A:267:LEU:HD21	1:A:314:ALA:HB2	1.95	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:349:ASP:CB	1:A:352:ASN:HB2	2.43	0.48
1:B:397:GLN:HA	1:B:397:GLN:OE1	2.13	0.48
1:C:308:ASN:O	1:C:312:GLN:HG2	2.13	0.48
1:A:177:TYR:HE2	1:A:182:LEU:HB2	1.77	0.48
1:A:191:GLU:O	1:A:194:GLU:HB3	2.12	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:153:ARG:HG3	1:A:153:ARG:HH11	1.78	0.48
1:A:166:GLU:O	1:A:255:ALA:HB3	2.14	0.48
1:B:156:THR:O	1:B:158:GLY:N	2.46	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:B:1:GLY:O	1:B:2:SER:CB	2.62	0.48
1:C:254:LYS:HD2	3:C:491:HOH:O	2.13	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:151:ILE:O	1:A:219:VAL:HA	2.15	0.47
1:A:261:MET:HB2	1:A:265:GLU:HB2	1.97	0.47
1:A:226:ALA:O	1:A:228:GLY:N	2.48	0.47
1:A:274:LYS:HG3	1:A:293:TYR:HE2	1.80	0.47
1:C:4:MET:SD	1:C:210:ARG:CZ	3.03	0.47
1:A:169:ILE:O	1:A:252:PHE:HA	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:264:GLU:HA	1:A:267:LEU:HD12	1.96	0.46
1:A:304:SER:C	1:A:306:PHE:H	2.18	0.46
1:B:157:ARG:HD3	1:B:157:ARG:H	1.79	0.46
1:C:221:LEU:HD13	1:C:226:ALA:HB2	1.98	0.46
1:C:262:ASN:OD1	1:C:265:GLU:HG3	2.15	0.46
1:A:215:GLU:HG2	1:A:216:HIS:N	2.30	0.46
1:A:315:GLN:HG3	3:A:468:HOH:O	2.14	0.46
1:B:226:ALA:HB1	1:B:243:LEU:CD1	2.40	0.46
1:A:285:LYS:HB3	1:A:288:GLN:NE2	2.30	0.46
1:C:177:TYR:CE2	1:C:182:LEU:HB2	2.51	0.46
1:A:178:TYR:O	1:A:180:ASP:N	2.44	0.46
1:A:221:LEU:HD13	1:A:226:ALA:HB2	1.96	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:274:LYS:O	1:C:278:THR:HG23	2.16	0.46
1:C:395:VAL:O	1:C:399:ARG:HG2	2.16	0.46
1:A:185:GLN:O	1:A:185:GLN:CD	2.54	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:B:380:LEU:HD11	1:B:389:ALA:HB1	1.97	0.46
1:C:177:TYR:HE2	1:C:182:LEU:HB2	1.81	0.46
1:A:198:LEU:O	1:A:200:LEU:HG	2.16	0.45
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:C:317:LEU:HD12	1:C:317:LEU:HA	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:A:408:LYS:HA	1:B:418:LEU:HD13	1.99	0.45
1:B:357:PHE:HZ	1:B:388:ALA:HB3	1.82	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:232:LYS:HE3	1:A:234:LYS:HE3	1.99	0.45
1:A:261:MET:CE	1:A:266:LYS:HA	2.46	0.45
1:A:317:LEU:HD12	1:A:317:LEU:HA	1.85	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:382:LEU:N	1:C:382:LEU:HD12	2.32	0.45
1:A:208:ILE:HA	1:A:211:MET:HG3	1.99	0.44
1:B:227:PHE:CD1	1:B:232:LYS:HB2	2.52	0.44
1:C:146:GLU:OE2	1:C:210:ARG:NH2	2.50	0.44
1:A:201:PRO:O	1:A:202:TYR:C	2.56	0.44
1:A:333:GLN:NE2	1:C:151:ILE:CD1	2.80	0.44
1:A:376:PHE:CE1	1:A:392:GLN:HB3	2.52	0.44
1:B:172:VAL:HG12	1:B:173:ALA:N	2.32	0.44
1:C:357:PHE:HZ	1:C:388:ALA:CB	2.20	0.44
1:B:344:LYS:HD3	1:B:344:LYS:HA	1.76	0.44
1:C:178:TYR:OH	1:C:179:LYS:HE3	2.17	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:367:ASN:ND2	1:B:369:PHE:CZ	2.85	0.44
1:B:373:ARG:HH21	1:C:304:SER:CB	2.31	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:O	1:B:156:THR:CG2	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:392:GLN:C	1:C:394:ALA:N	2.71	0.43
1:C:401:ARG:HH12	1:C:402:ARG:HH21	1.64	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:A:158:GLY:O	1:A:159:GLU:C	2.56	0.43
1:A:323:LEU:O	1:A:358:ARG:NH1	2.51	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:B:379:VAL:O	1:B:383:TYR:N	2.40	0.43
1:C:178:TYR:CE1	1:C:241:ALA:HB1	2.43	0.43
1:A:290:LEU:HD23	1:A:290:LEU:HA	1.85	0.43
1:B:4:MET:CE	1:B:210:ARG:NH2	2.82	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:155:GLN:O	1:B:156:THR:CB	2.67	0.43
1:B:172:VAL:O	1:B:187:GLU:HA	2.18	0.43
1:B:349:ASP:HB3	1:B:352:ASN:HB2	2.01	0.43
1:B:350:SER:H	1:C:350:SER:CB	2.20	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: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
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:A:321:SER:O	1:A:325:LEU:HD23	2.18	0.42
1:B:155:GLN:O	1:B:155:GLN:HG2	2.19	0.42
1:B:290:LEU:HD11	1:B:332:LEU:HD12	2.01	0.42
1:C:251:SER:O	1:C:252:PHE:HB3	2.19	0.42
1:C:318:ARG:O	1:C:321:SER:HB2	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:O	1:A:154:ILE:HG22	2.20	0.42
1:A:393:LEU:O	1:A:397:GLN:HG2	2.20	0.42
1:B:152:ARG:HA	1:B:218:ILE:O	2.20	0.42
1:B:176:GLY:HA2	1:B:244:LYS:O	2.19	0.42
1:B:392:GLN:O	1:B:393:LEU:C	2.58	0.42
1:A:185:GLN:HG3	3:A:518:HOH:O	2.19	0.42
1:A:4:MET:CE	1:A:4:MET:CG	2.94	0.42
1:A:211:MET:HB2	1:A:249:LEU:CD2	2.50	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:A:251:SER:O	1:A:252:PHE:CB	2.65	0.42
1:A:415:PHE:CE1	1:B:408:LYS:HG3	2.55	0.42
1:B:156:THR:O	1:B:157:ARG:C	2.58	0.42
1:B:232:LYS:HD2	1:B:232:LYS:HA	1.86	0.42
1:B:380:LEU:HD12	1:B:389:ALA:HB1	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:369:PHE:O	1:C:370:GLU:C	2.57	0.41
2:G:1:MET:HA	3:G:24:HOH:O	2.20	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:161:TYR:HB3	1:C:162:ALA:H	1.57	0.41
1:C:211:MET:HE3	1:C:249:LEU:HB2	2.01	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:N	1:C:180:ASP:OD1	2.54	0.41
1:A:149:GLY:HA2	1:A:222:LYS:CG	2.51	0.41
1:B:304:SER:HB2	1:C:374:ALA:HB2	2.01	0.41
1:B:330:LEU:HD23	1:B:330:LEU:HA	1.92	0.41
1:C:163:LYS:HB2	1:C:164:PRO:HD2	2.03	0.41
1:C:189:ARG:NH1	1:C:189:ARG:HG3	2.35	0.41
1:C:208:ILE:HG12	1:C:247:LEU:CD1	2.49	0.41
1:C:287:LYS:HD3	3:C:471:HOH:O	2.21	0.41
1:A:151:ILE:HA	3:A:463:HOH:O	2.21	0.41
1:B:162:ALA:HB1	1:B:252:PHE:CZ	2.56	0.41
1:B:174:LEU:HD23	1:B:247:LEU:HD22	2.03	0.41
1:B:421:GLU:OE1	1:B:421:GLU:HA	2.20	0.41
1:C:351:ASN:OD1	1:C:382:LEU:HD21	2.21	0.41
1:B:179:LYS:O	1:B:180:ASP:CB	2.64	0.40
1:C:189:ARG:HG3	1:C:189:ARG:HH11	1.86	0.40
1:C:183:PHE:HE2	1:C:243:LEU:HD22	1.87	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:C:223:PRO:HD3	1:C:242:GLU:HB3	2.02	0.40
1:A:188:LEU:HD13	1:A:198:LEU:HD13	2.03	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:290:LEU:CD2	1:A:325:LEU:HD12	2.52	0.40
1:B:172:VAL:CG1	1:B:173:ALA:N	2.85	0.40
1:B:233:GLU:C	1:B:236:GLN:H	2.25	0.40
1:C:183:PHE:CE2	1:C:243:LEU:HD22	2.56	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

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%)	1 6
1	B	283/336 (84%)	229 (81%)	34 (12%)	20 (7%)	1 5
1	C	283/336 (84%)	242 (86%)	23 (8%)	18 (6%)	1 7
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%)	1 6

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

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Mol	Chain	Res	Type
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
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 (i)

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%)	16 49
1	B	238/283 (84%)	220 (92%)	18 (8%)	13 43
1	C	239/283 (84%)	224 (94%)	15 (6%)	18 51
2	G	5/5 (100%)	3 (60%)	2 (40%)	0 0
2	H	5/5 (100%)	5 (100%)	0	100 100
All	All	725/859 (84%)	674 (93%)	51 (7%)	15 47

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

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Mol	Chain	Res	Type
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
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

Sometimes 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

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Mol	Chain	Res	Type
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
1	C	423	ASN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

There are no ligands in this entry.

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

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, 95^{th} 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%)	-0.19	2 (0%) 87 69	16, 51, 96, 130	0
1	B	285/336 (84%)	-0.31	2 (0%) 87 69	15, 50, 91, 121	0
1	C	285/336 (84%)	-0.27	0 100 100	10, 44, 92, 125	0
2	G	5/5 (100%)	1.22	1 (20%) 1 0	81, 86, 92, 93	0
2	H	5/5 (100%)	1.03	0 100 100	73, 82, 91, 95	0
All	All	865/1018 (84%)	-0.24	5 (0%) 89 72	10, 50, 93, 130	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	G	5	ASP	3.6
1	B	3	HIS	3.0
1	B	2	SER	2.7
1	A	153	ARG	2.5
1	A	247	LEU	2.4

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 monosaccharides in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.