



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 11:20 AM GMT

PDB ID : 3OQ5
Title : Crystal structure of the 3-MBT domain from human L3MBTL1 in complex with p53K382me1
Authors : Roy, S.; West, L.E.; Weiner, K.L.; Hayashi, R.; Shi, X.; Appella, E.; Gozani, O.; Kutateladze, T.
Deposited on : 2010-09-02
Resolution : 2.50 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

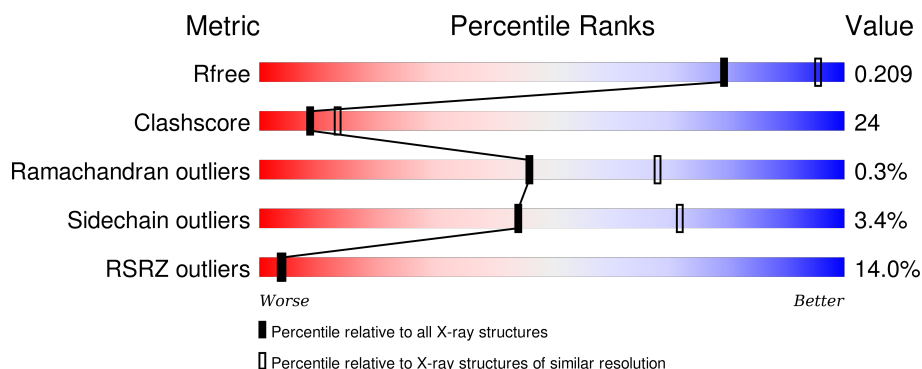
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)

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. 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 $\leq 5\%$. 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	345	<div> <div>6%</div> <div>53%</div> <div>35%</div> <div>10%</div> </div>
1	B	345	<div> <div>18%</div> <div>54%</div> <div>34%</div> <div>10%</div> </div>
1	C	345	<div> <div>14%</div> <div>51%</div> <div>37%</div> <div>10%</div> </div>
2	D	10	<div> <div>10%</div> <div>90%</div> </div>
2	E	10	<div> <div>10%</div> <div>90%</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8043 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 Lethal(3)malignant brain tumor-like protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	309	Total	C	N	O	S	0	0	0
			2509	1617	427	453	12			
1	B	309	Total	C	N	O	S	0	0	0
			2509	1617	427	453	12			
1	C	309	Total	C	N	O	S	0	0	0
			2509	1617	427	453	12			

There are 15 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	186	GLY	-	EXPRESSION TAG	UNP Q9Y468
A	187	PRO	-	EXPRESSION TAG	UNP Q9Y468
A	188	LEU	-	EXPRESSION TAG	UNP Q9Y468
A	189	GLY	-	EXPRESSION TAG	UNP Q9Y468
A	190	SER	-	EXPRESSION TAG	UNP Q9Y468
B	186	GLY	-	EXPRESSION TAG	UNP Q9Y468
B	187	PRO	-	EXPRESSION TAG	UNP Q9Y468
B	188	LEU	-	EXPRESSION TAG	UNP Q9Y468
B	189	GLY	-	EXPRESSION TAG	UNP Q9Y468
B	190	SER	-	EXPRESSION TAG	UNP Q9Y468
C	186	GLY	-	EXPRESSION TAG	UNP Q9Y468
C	187	PRO	-	EXPRESSION TAG	UNP Q9Y468
C	188	LEU	-	EXPRESSION TAG	UNP Q9Y468
C	189	GLY	-	EXPRESSION TAG	UNP Q9Y468
C	190	SER	-	EXPRESSION TAG	UNP Q9Y468

- Molecule 2 is a protein called Cellular tumor antigen p53.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	D	1	Total	C	N	O	0	0	0
			10	7	2	1			

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	E	1	Total	C	N	O	0	0	0
			10	7	2	1			

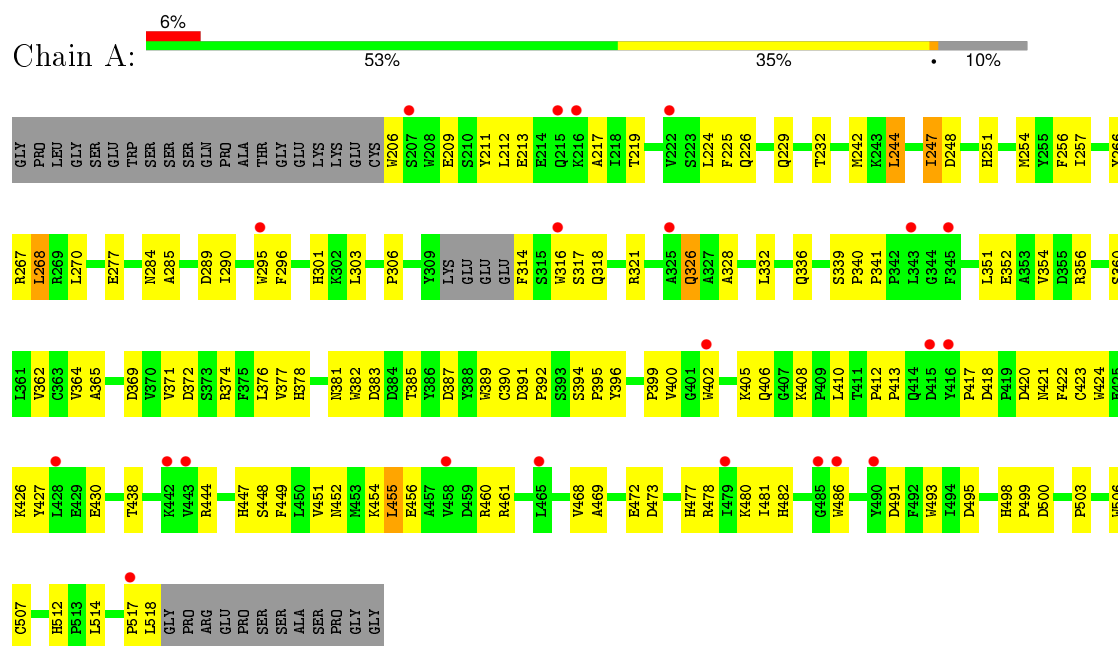
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	149	Total	O	0	0
			149	149		
3	B	155	Total	O	0	0
			155	155		
3	C	189	Total	O	0	0
			189	189		
3	D	3	Total	O	0	0
			3	3		

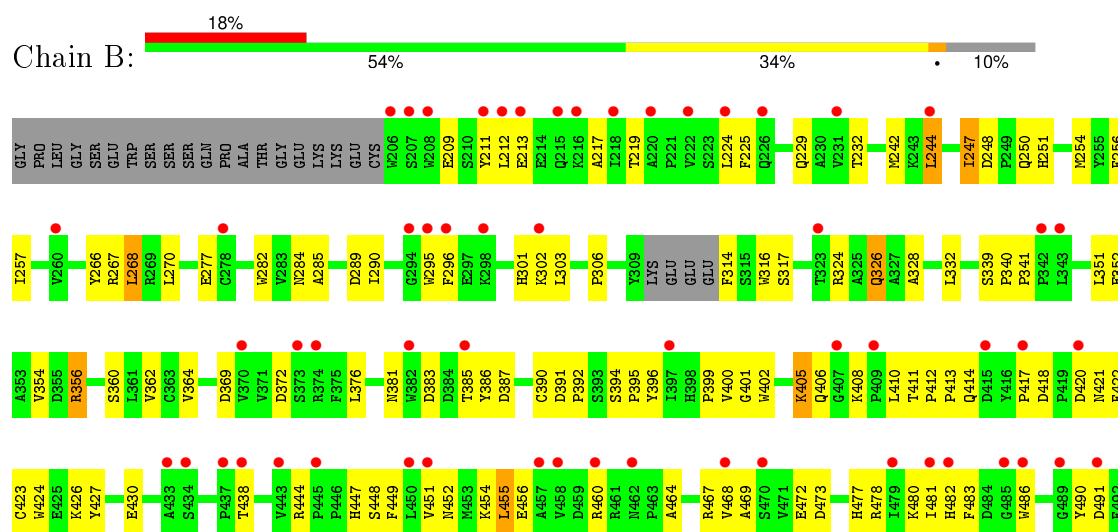
3 Residue-property plots

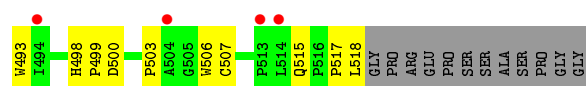
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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: Lethal(3)malignant brain tumor-like protein

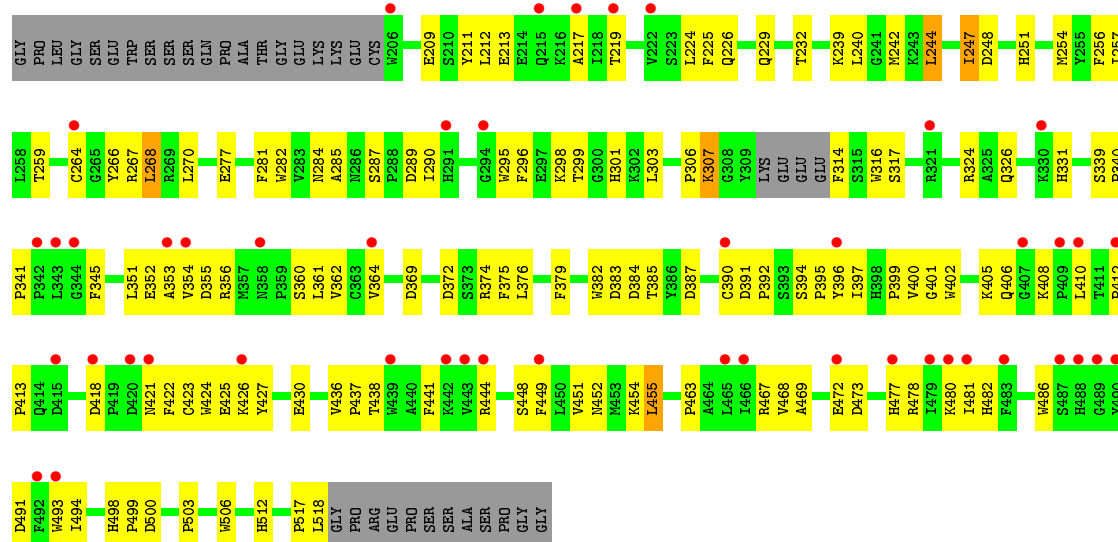


- Molecule 1: Lethal(3)malignant brain tumor-like protein

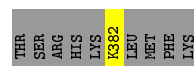




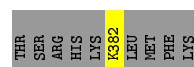
- Molecule 1: Lethal(3)malignant brain tumor-like protein



- Molecule 2: Cellular tumor antigen p53



- Molecule 2: Cellular tumor antigen p53



4 Data and refinement statistics

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, α , β , γ	109.08 Å 109.08 Å 90.05 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	46.65 – 2.50 46.65 – 2.50	Depositor EDS
% Data completeness (in resolution range)	100.0 (46.65-2.50) 100.0 (46.65-2.50)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.29 (at 2.51 Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.1_357)	Depositor
R, R_{free}	0.205 , 0.234 0.183 , 0.209	Depositor DCC
R_{free} test set	2086 reflections (5.31%)	DCC
Wilson B-factor (Å ²)	47.1	Xtriage
Anisotropy	0.380	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 48.7	EDS
Estimated twinning fraction	0.469 for -h,-k,l 0.406 for h,-h-k,-l 0.406 for -k,-h,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtriage
Outliers	0 of 41472 reflections	Xtriage
F_o, F_c correlation	0.84	EDS
Total number of atoms	8043	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.74% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MLZ

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	0.31	0/2607	0.43	0/3564
1	B	0.32	0/2607	0.46	2/3564 (0.1%)
1	C	0.32	0/2607	0.46	2/3564 (0.1%)
All	All	0.32	0/7821	0.45	4/10692 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	356	ARG	NE-CZ-NH1	5.93	123.27	120.30
1	B	356	ARG	NE-CZ-NH2	-5.87	117.36	120.30
1	C	356	ARG	NE-CZ-NH1	5.79	123.20	120.30
1	C	356	ARG	NE-CZ-NH2	-5.75	117.42	120.30

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	2509	0	2348	109	0
1	B	2509	0	2348	114	0
1	C	2509	0	2348	128	0
2	D	10	0	15	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	E	10	0	15	4	0
3	A	149	0	0	14	0
3	B	155	0	0	22	0
3	C	189	0	0	35	0
3	D	3	0	0	0	0
All	All	8043	0	7074	346	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 24.

All (346) 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:C:401:GLY:HA2	3:C:664:HOH:O	1.45	1.12
1:C:298:LYS:HB2	3:C:576:HOH:O	1.70	0.91
1:B:506:TRP:HB3	3:B:590:HOH:O	1.72	0.88
1:A:374:ARG:HD2	3:A:109:HOH:O	1.81	0.81
1:C:339:SER:HB2	1:C:340:PRO:HD2	1.63	0.80
1:C:493:TRP:HE3	3:C:638:HOH:O	1.65	0.79
1:B:339:SER:HB2	1:B:340:PRO:HD2	1.65	0.79
1:C:270:LEU:HD22	3:C:615:HOH:O	1.83	0.79
1:A:339:SER:HB2	1:A:340:PRO:HD2	1.64	0.78
1:C:240:LEU:HD23	3:C:607:HOH:O	1.82	0.78
1:C:382:TRP:CH2	2:E:382:MLZ:HB2	2.18	0.78
1:C:441:PHE:HA	3:C:594:HOH:O	1.85	0.77
1:C:390:CYS:SG	3:C:547:HOH:O	2.42	0.77
1:B:212:LEU:HD11	1:B:219:THR:HG23	1.66	0.76
1:C:212:LEU:HD11	1:C:219:THR:HG23	1.67	0.75
1:C:379:PHE:HB2	3:C:173:HOH:O	1.87	0.74
1:A:206:TRP:N	3:A:601:HOH:O	2.21	0.74
1:B:418:ASP:OD2	1:B:421:ASN:HB2	1.89	0.73
1:A:212:LEU:HD11	1:A:219:THR:HG23	1.70	0.72
1:C:418:ASP:OD2	1:C:421:ASN:HB2	1.90	0.71
1:C:266:TYR:CE1	1:C:285:ALA:HB3	2.26	0.71
1:A:365:ALA:HB1	3:A:133:HOH:O	1.91	0.71
1:A:418:ASP:OD2	1:A:421:ASN:HB2	1.91	0.70
1:C:473:ASP:H	1:C:480:LYS:HB2	1.56	0.70
1:B:473:ASP:H	1:B:480:LYS:HB2	1.57	0.69
1:C:354:VAL:HG22	1:C:362:VAL:HG22	1.74	0.69
1:C:240:LEU:HA	3:C:607:HOH:O	1.93	0.69
1:A:266:TYR:CE1	1:A:285:ALA:HB3	2.29	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:512:HIS:HB2	3:C:171:HOH:O	1.94	0.68
1:B:266:TYR:CE1	1:B:285:ALA:HB3	2.29	0.68
1:A:257:ILE:HD12	1:A:306:PRO:HD3	1.76	0.67
1:C:326:GLN:HB2	3:C:608:HOH:O	1.95	0.67
1:A:211:TYR:CE1	1:A:517:PRO:HG2	2.30	0.66
1:A:354:VAL:HG22	1:A:362:VAL:HG22	1.76	0.66
1:C:480:LYS:HZ1	1:C:491:ASP:HB3	1.60	0.66
1:A:473:ASP:H	1:A:480:LYS:HB2	1.59	0.66
1:A:364:VAL:HG13	1:A:424:TRP:CH2	2.30	0.66
1:C:211:TYR:CE1	1:C:517:PRO:HG2	2.30	0.66
1:A:267:ARG:HG2	1:A:284:ASN:HD22	1.60	0.66
1:B:364:VAL:HG13	1:B:424:TRP:CH2	2.31	0.66
1:B:354:VAL:HG22	1:B:362:VAL:HG22	1.76	0.66
1:A:461:ARG:HD3	3:A:606:HOH:O	1.95	0.66
1:C:251:HIS:CE1	3:C:174:HOH:O	2.48	0.65
1:C:480:LYS:NZ	1:C:491:ASP:HB3	2.11	0.65
1:C:267:ARG:HG2	1:C:284:ASN:HD22	1.61	0.65
1:C:467:ARG:NH2	3:C:116:HOH:O	2.30	0.64
1:B:257:ILE:HD12	1:B:306:PRO:HD3	1.78	0.64
1:C:257:ILE:HD12	1:C:306:PRO:HD3	1.79	0.64
1:B:480:LYS:NZ	1:B:491:ASP:HB3	2.12	0.64
1:C:364:VAL:HG13	1:C:424:TRP:CH2	2.33	0.64
1:A:480:LYS:NZ	1:A:491:ASP:HB3	2.13	0.64
1:B:211:TYR:CE1	1:B:517:PRO:HG2	2.32	0.64
1:C:240:LEU:N	3:C:607:HOH:O	2.29	0.64
1:A:224:LEU:HD22	1:A:448:SER:HB2	1.80	0.64
1:B:250:GLN:NE2	3:B:593:HOH:O	2.31	0.64
1:C:240:LEU:CA	3:C:607:HOH:O	2.46	0.63
1:B:267:ARG:HG2	1:B:284:ASN:HD22	1.63	0.63
1:C:268:LEU:HD12	1:C:270:LEU:HD21	1.81	0.63
1:B:480:LYS:HZ1	1:B:491:ASP:HB3	1.62	0.62
1:C:287:SER:HB2	3:C:172:HOH:O	1.99	0.62
1:C:299:THR:HG23	3:C:576:HOH:O	2.00	0.62
1:C:399:PRO:HG3	1:C:438:THR:HG23	1.82	0.62
1:B:224:LEU:HD22	1:B:448:SER:HB2	1.82	0.61
1:C:212:LEU:HD11	1:C:219:THR:CG2	2.31	0.61
1:B:212:LEU:HD11	1:B:219:THR:CG2	2.30	0.61
1:C:224:LEU:HD22	1:C:448:SER:HB2	1.83	0.61
1:B:395:PRO:HG2	1:B:396:TYR:CE2	2.36	0.61
1:B:399:PRO:HG3	1:B:438:THR:HG23	1.84	0.60
1:A:341:PRO:HA	1:A:372:ASP:O	2.02	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:395:PRO:HG2	1:A:396:TYR:CE2	2.37	0.59
1:A:268:LEU:HD12	1:A:270:LEU:HD21	1.84	0.59
1:C:472:GLU:HB3	1:C:480:LYS:HB3	1.85	0.59
1:A:212:LEU:HD11	1:A:219:THR:CG2	2.33	0.58
1:C:251:HIS:CG	3:C:174:HOH:O	2.56	0.58
1:C:282:TRP:HD1	3:C:604:HOH:O	1.84	0.58
1:B:268:LEU:HD12	1:B:270:LEU:HD21	1.84	0.58
1:A:472:GLU:HB3	1:A:480:LYS:HB3	1.85	0.58
1:C:341:PRO:HA	1:C:372:ASP:O	2.04	0.57
1:C:391:ASP:HB2	1:C:392:PRO:HD2	1.86	0.57
1:A:399:PRO:HG3	1:A:438:THR:HG23	1.85	0.57
1:B:456:GLU:CD	3:B:590:HOH:O	2.43	0.57
1:B:451:VAL:O	1:B:452:ASN:HB2	2.04	0.57
1:B:341:PRO:HA	1:B:372:ASP:O	2.04	0.57
1:A:480:LYS:HZ1	1:A:491:ASP:HB3	1.70	0.57
1:C:493:TRP:CE3	3:C:638:HOH:O	2.49	0.56
1:C:259:THR:C	3:C:615:HOH:O	2.44	0.56
1:A:382:TRP:CH2	2:D:382:MLZ:HE3	2.40	0.56
1:A:469:ALA:HB1	1:A:481:ILE:HG23	1.88	0.56
1:C:469:ALA:HB1	1:C:481:ILE:HG23	1.87	0.56
1:B:339:SER:CB	1:B:340:PRO:HD2	2.35	0.56
1:A:391:ASP:HB2	1:A:392:PRO:HD2	1.87	0.56
1:B:472:GLU:HB3	1:B:480:LYS:HB3	1.86	0.55
1:B:417:PRO:HD3	3:B:543:HOH:O	2.06	0.55
1:A:451:VAL:O	1:A:452:ASN:HB2	2.07	0.55
1:C:339:SER:CB	1:C:340:PRO:HD2	2.34	0.55
1:A:468:VAL:HG21	1:A:517:PRO:HD3	1.89	0.55
1:A:377:VAL:HG13	3:A:133:HOH:O	2.05	0.55
1:A:301:HIS:HE1	1:A:369:ASP:OD1	1.89	0.55
1:B:464:ALA:HB2	3:B:541:HOH:O	2.06	0.54
1:C:451:VAL:O	1:C:452:ASN:HB2	2.07	0.54
1:B:301:HIS:HE1	1:B:369:ASP:OD1	1.90	0.54
1:B:391:ASP:HB2	1:B:392:PRO:HD2	1.90	0.54
1:C:270:LEU:CD2	3:C:615:HOH:O	2.49	0.54
1:B:478:ARG:HB3	1:B:493:TRP:CE3	2.43	0.54
1:A:339:SER:CB	1:A:340:PRO:HD2	2.35	0.54
1:A:395:PRO:HG2	1:A:396:TYR:CD2	2.43	0.53
1:C:395:PRO:HG2	1:C:396:TYR:CE2	2.43	0.53
1:A:444:ARG:O	1:A:477:HIS:HE1	1.92	0.53
1:B:469:ALA:HB1	1:B:481:ILE:HG23	1.89	0.53
1:A:410:LEU:O	1:A:412:PRO:HD3	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:410:LEU:O	1:B:412:PRO:HD3	2.09	0.53
1:C:390:CYS:HB2	1:C:394:SER:HB2	1.90	0.53
1:C:468:VAL:HG21	1:C:517:PRO:HD3	1.91	0.53
1:B:395:PRO:HG2	1:B:396:TYR:CD2	2.43	0.53
1:C:444:ARG:O	1:C:477:HIS:HE1	1.91	0.53
1:A:224:LEU:HD13	1:A:449:PHE:HA	1.89	0.53
1:C:224:LEU:HD13	1:C:449:PHE:HA	1.90	0.53
1:A:356:ARG:NH2	1:A:495:ASP:OD2	2.42	0.53
1:C:375:PHE:CZ	3:C:547:HOH:O	2.54	0.52
1:A:257:ILE:HD12	1:A:306:PRO:HG3	1.91	0.52
1:B:444:ARG:O	1:B:477:HIS:HE1	1.92	0.52
1:B:225:PHE:CD2	1:B:503:PRO:HG3	2.44	0.52
1:A:478:ARG:HB3	1:A:493:TRP:CE3	2.45	0.52
1:C:410:LEU:O	1:C:412:PRO:HD3	2.10	0.52
1:A:224:LEU:CD2	1:A:448:SER:HB2	2.39	0.52
1:A:423:CYS:HB3	1:A:426:LYS:HB3	1.92	0.52
1:B:209:GLU:O	1:B:213:GLU:HB2	2.09	0.52
1:C:400:VAL:HG13	1:C:424:TRP:CG	2.45	0.52
1:C:423:CYS:HB3	1:C:426:LYS:HB3	1.92	0.52
1:C:478:ARG:HB3	1:C:493:TRP:CE3	2.45	0.52
1:A:382:TRP:CZ2	2:D:382:MLZ:HE3	2.45	0.52
1:B:390:CYS:HB2	1:B:394:SER:HB2	1.92	0.51
1:B:224:LEU:HD13	1:B:449:PHE:HA	1.91	0.51
1:B:302:LYS:HG3	3:B:154:HOH:O	2.10	0.51
1:B:423:CYS:HB3	1:B:426:LYS:HB3	1.93	0.51
1:C:301:HIS:HE1	1:C:369:ASP:OD1	1.93	0.51
1:B:455:LEU:HA	3:B:582:HOH:O	2.10	0.51
1:C:209:GLU:O	1:C:213:GLU:HB2	2.11	0.51
1:B:229:GLN:HA	3:B:542:HOH:O	2.10	0.51
1:C:486:TRP:CE2	1:C:518:LEU:HD13	2.46	0.51
1:B:468:VAL:HG21	1:B:517:PRO:HD3	1.93	0.51
1:A:225:PHE:CD2	1:A:503:PRO:HG3	2.46	0.50
1:B:490:TYR:HB2	3:B:577:HOH:O	2.10	0.50
1:C:296:PHE:CD1	1:C:303:LEU:HB2	2.47	0.50
1:A:209:GLU:O	1:A:213:GLU:HB2	2.11	0.50
1:A:390:CYS:HB2	1:A:394:SER:HB2	1.94	0.50
1:A:381:ASN:ND2	3:A:579:HOH:O	2.43	0.50
1:A:257:ILE:HD12	1:A:306:PRO:CD	2.40	0.50
1:B:224:LEU:CD2	1:B:448:SER:HB2	2.42	0.50
1:A:232:THR:HG22	1:A:506:TRP:CD1	2.46	0.50
1:C:467:ARG:NH1	3:C:548:HOH:O	2.30	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:480:LYS:HE3	1:B:421:ASN:HD21	1.77	0.49
1:B:478:ARG:HB3	1:B:493:TRP:HE3	1.77	0.49
1:C:494:ILE:N	3:C:638:HOH:O	2.44	0.49
1:B:400:VAL:HG13	1:B:424:TRP:CG	2.48	0.49
1:C:257:ILE:HD12	1:C:306:PRO:HG3	1.94	0.49
1:C:498:HIS:CE1	1:C:500:ASP:HB2	2.47	0.49
1:C:478:ARG:HB3	1:C:493:TRP:HE3	1.77	0.49
1:C:480:LYS:NZ	1:C:480:LYS:HB3	2.27	0.49
1:A:427:TYR:HA	1:A:430:GLU:HG2	1.95	0.49
1:C:486:TRP:CZ2	1:C:518:LEU:HD13	2.47	0.49
1:B:247:ILE:HG13	1:B:289:ASP:O	2.13	0.49
1:A:296:PHE:CD1	1:A:303:LEU:HB2	2.48	0.49
1:B:257:ILE:HD12	1:B:306:PRO:CD	2.43	0.49
1:A:486:TRP:CZ2	1:A:518:LEU:HD13	2.48	0.49
1:C:224:LEU:CD2	1:C:448:SER:HB2	2.41	0.49
1:B:232:THR:HG22	1:B:506:TRP:CD1	2.47	0.48
1:A:486:TRP:CE2	1:A:518:LEU:HD13	2.48	0.48
1:A:400:VAL:HG13	1:A:424:TRP:CG	2.47	0.48
1:A:277:GLU:CD	1:A:277:GLU:H	2.16	0.48
1:A:478:ARG:HB3	1:A:493:TRP:HE3	1.79	0.48
1:A:378:HIS:N	3:A:133:HOH:O	2.46	0.48
1:B:464:ALA:HB2	3:B:550:HOH:O	2.13	0.48
1:B:486:TRP:CE2	1:B:518:LEU:HD13	2.48	0.48
1:C:277:GLU:H	1:C:277:GLU:CD	2.17	0.48
1:C:480:LYS:HG3	1:C:493:TRP:CZ2	2.49	0.48
1:B:257:ILE:HD12	1:B:306:PRO:HG3	1.95	0.48
1:C:427:TYR:HA	1:C:430:GLU:HG2	1.95	0.48
1:B:426:LYS:HD3	3:B:589:HOH:O	2.13	0.48
1:B:277:GLU:H	1:B:277:GLU:CD	2.17	0.48
1:B:498:HIS:CE1	1:B:500:ASP:HB2	2.49	0.48
1:C:225:PHE:CD2	1:C:503:PRO:HG3	2.49	0.48
1:B:296:PHE:CD1	1:B:303:LEU:HB2	2.49	0.48
1:C:244:LEU:HD12	1:C:290:ILE:HB	1.96	0.48
1:A:257:ILE:HD12	1:A:306:PRO:CG	2.44	0.48
1:C:232:THR:HG22	1:C:506:TRP:CD1	2.49	0.48
1:B:217:ALA:CB	1:B:454:LYS:HE2	2.44	0.47
1:B:427:TYR:HA	1:B:430:GLU:HG2	1.95	0.47
1:C:361:LEU:CD1	2:E:382:MLZ:HE3	2.43	0.47
1:C:331:HIS:HD2	3:C:655:HOH:O	1.98	0.47
1:A:455:LEU:H	1:A:455:LEU:HD12	1.80	0.47
1:B:364:VAL:HG21	1:B:413:PRO:HD3	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:480:LYS:HG3	1:A:493:TRP:CZ2	2.50	0.47
1:B:480:LYS:NZ	1:B:480:LYS:HB3	2.29	0.47
1:A:285:ALA:O	3:A:120:HOH:O	2.20	0.47
1:A:267:ARG:HG2	1:A:284:ASN:ND2	2.28	0.47
1:C:217:ALA:CB	1:C:454:LYS:HE2	2.45	0.47
1:B:486:TRP:CZ2	1:B:518:LEU:HD13	2.50	0.47
1:C:251:HIS:ND1	3:C:174:HOH:O	2.34	0.46
1:A:217:ALA:CB	1:A:454:LYS:HE2	2.44	0.46
1:B:301:HIS:HA	3:B:154:HOH:O	2.14	0.46
1:A:257:ILE:CD1	1:A:306:PRO:HD3	2.44	0.46
1:C:395:PRO:HG2	1:C:396:TYR:CD2	2.50	0.46
1:B:455:LEU:H	1:B:455:LEU:HD12	1.81	0.46
1:B:356:ARG:HG2	3:B:105:HOH:O	2.15	0.46
1:A:421:ASN:HD21	1:B:480:LYS:HE3	1.80	0.46
1:A:257:ILE:HD13	1:A:314:PHE:CZ	2.50	0.46
1:B:364:VAL:HG21	1:B:422:PHE:CZ	2.50	0.46
1:C:257:ILE:HD12	1:C:306:PRO:CD	2.44	0.46
1:A:340:PRO:HA	1:A:341:PRO:HD3	1.75	0.46
1:B:212:LEU:CD1	1:B:219:THR:HG23	2.43	0.46
1:C:267:ARG:HG2	1:C:284:ASN:ND2	2.30	0.45
1:A:301:HIS:CE1	1:A:369:ASP:OD1	2.68	0.45
1:A:364:VAL:HG21	1:A:413:PRO:HD3	1.98	0.45
1:C:247:ILE:HG13	1:C:289:ASP:O	2.17	0.45
1:C:455:LEU:HD12	1:C:455:LEU:H	1.81	0.45
1:A:514:LEU:HD22	3:A:564:HOH:O	2.15	0.45
1:C:397:ILE:HD13	3:C:547:HOH:O	2.16	0.45
1:B:376:LEU:HD11	1:B:387:ASP:HB3	1.98	0.45
1:C:257:ILE:HG23	1:C:316:TRP:CH2	2.51	0.45
1:C:257:ILE:HD13	1:C:314:PHE:CZ	2.51	0.45
1:C:364:VAL:HG21	1:C:413:PRO:HD3	1.98	0.45
1:B:301:HIS:CE1	1:B:369:ASP:OD1	2.69	0.45
1:A:247:ILE:HG13	1:A:289:ASP:O	2.16	0.45
1:A:376:LEU:HD11	1:A:387:ASP:HB3	1.98	0.45
1:C:212:LEU:CD1	1:C:219:THR:HG23	2.43	0.45
1:A:364:VAL:HG21	1:A:422:PHE:CZ	2.52	0.45
1:C:498:HIS:CG	1:C:499:PRO:HD2	2.52	0.45
1:A:493:TRP:HD1	1:B:420:ASP:HB3	1.82	0.45
1:A:352:GLU:OE1	1:A:402:TRP:HB3	2.17	0.45
1:A:480:LYS:HB3	1:A:480:LYS:NZ	2.31	0.45
1:B:257:ILE:HD13	1:B:314:PHE:CZ	2.51	0.45
1:C:352:GLU:OE1	1:C:402:TRP:HB3	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:242:MET:HA	1:C:326:GLN:O	2.17	0.44
1:B:257:ILE:HD12	1:B:306:PRO:CG	2.47	0.44
1:A:498:HIS:CE1	1:A:500:ASP:HB2	2.52	0.44
1:B:352:GLU:OE1	1:B:402:TRP:HB3	2.17	0.44
1:A:244:LEU:HD12	1:A:290:ILE:HB	1.99	0.44
1:C:480:LYS:HG3	1:C:493:TRP:CE2	2.51	0.44
1:C:251:HIS:CD2	1:C:385:THR:HG22	2.52	0.44
1:C:364:VAL:HG21	1:C:422:PHE:CZ	2.52	0.44
1:B:386:TYR:O	3:B:162:HOH:O	2.20	0.44
1:B:248:ASP:HB2	1:B:256:PHE:CE2	2.52	0.44
1:C:239:LYS:HB2	1:C:242:MET:HE3	2.00	0.44
1:C:306:PRO:O	1:C:307:LYS:C	2.56	0.44
1:B:460:ARG:NH2	3:B:574:HOH:O	2.46	0.44
1:C:405:LYS:HB3	1:C:405:LYS:HE2	1.71	0.44
1:A:248:ASP:HB2	1:A:256:PHE:CE2	2.52	0.44
1:B:480:LYS:HG3	1:B:493:TRP:CZ2	2.52	0.44
1:A:257:ILE:HG23	1:A:316:TRP:CH2	2.52	0.44
1:C:301:HIS:CE1	1:C:369:ASP:OD1	2.70	0.44
1:B:498:HIS:CG	1:B:499:PRO:HD2	2.53	0.44
1:B:381:ASN:HB3	3:B:583:HOH:O	2.17	0.44
1:B:483:PHE:HB3	3:B:586:HOH:O	2.18	0.44
1:A:242:MET:HA	1:A:326:GLN:O	2.18	0.44
1:C:353:ALA:HB3	3:C:635:HOH:O	2.18	0.44
1:B:244:LEU:HD12	1:B:290:ILE:HB	1.99	0.44
1:C:376:LEU:HD11	1:C:387:ASP:HB3	2.00	0.44
1:A:405:LYS:NZ	3:A:571:HOH:O	2.28	0.44
1:A:417:PRO:HA	3:A:8:HOH:O	2.18	0.43
1:B:456:GLU:OE1	1:B:507:CYS:HB2	2.18	0.43
1:C:482:HIS:HD2	1:C:491:ASP:OD1	2.02	0.43
1:B:257:ILE:HG23	1:B:316:TRP:CH2	2.53	0.43
1:B:482:HIS:HD2	1:B:491:ASP:OD1	2.02	0.43
1:B:401:GLY:O	1:B:405:LYS:HB3	2.19	0.43
1:A:480:LYS:HG3	1:A:493:TRP:CE2	2.53	0.43
1:A:456:GLU:OE1	1:A:507:CYS:HB2	2.18	0.43
1:B:480:LYS:HG3	1:B:493:TRP:CE2	2.54	0.43
1:C:257:ILE:HD12	1:C:306:PRO:CG	2.48	0.43
1:A:295:TRP:CH2	1:A:301:HIS:CD2	3.07	0.43
1:A:512:HIS:HB2	3:A:123:HOH:O	2.18	0.43
1:A:498:HIS:CG	1:A:499:PRO:HD2	2.54	0.43
1:B:242:MET:HA	1:B:326:GLN:O	2.18	0.43
1:A:257:ILE:CD1	1:A:314:PHE:CZ	3.02	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:382:TRP:CH2	2:D:382:MLZ:HG3	2.53	0.43
1:B:295:TRP:CH2	1:B:301:HIS:CD2	3.06	0.43
1:C:480:LYS:CB	1:C:480:LYS:NZ	2.81	0.43
1:A:482:HIS:HD2	1:A:491:ASP:OD1	2.02	0.43
1:A:247:ILE:HA	1:A:254:MET:O	2.18	0.43
1:B:251:HIS:CD2	1:B:385:THR:HG22	2.54	0.43
1:C:355:ASP:OD2	2:E:382:MLZ:HD2	2.18	0.42
1:A:420:ASP:HB3	1:B:493:TRP:HD1	1.84	0.42
1:C:281:PHE:HA	3:C:604:HOH:O	2.18	0.42
1:C:257:ILE:CD1	1:C:314:PHE:CZ	3.02	0.42
1:A:460:ARG:HB2	3:A:605:HOH:O	2.18	0.42
1:A:336:GLN:OE1	3:A:551:HOH:O	2.21	0.42
1:C:361:LEU:HD13	2:E:382:MLZ:HE3	2.01	0.42
1:B:282:TRP:NE1	3:B:541:HOH:O	2.33	0.42
1:C:295:TRP:CH2	1:C:301:HIS:CD2	3.07	0.42
1:B:426:LYS:HE2	1:B:430:GLU:OE1	2.19	0.42
1:A:217:ALA:HB3	1:A:454:LYS:HE2	2.02	0.42
1:A:251:HIS:CD2	1:A:385:THR:HG22	2.54	0.42
1:C:247:ILE:HA	1:C:254:MET:O	2.19	0.42
1:C:340:PRO:HA	1:C:341:PRO:HD3	1.76	0.42
1:C:354:VAL:CG2	1:C:362:VAL:HG22	2.46	0.42
1:B:473:ASP:N	1:B:480:LYS:HB2	2.32	0.42
1:A:480:LYS:CE	1:B:421:ASN:HD21	2.33	0.42
1:B:354:VAL:CG2	1:B:362:VAL:HG22	2.49	0.42
1:B:257:ILE:CG1	1:B:303:LEU:HD21	2.50	0.41
1:C:463:PRO:HB2	3:C:604:HOH:O	2.20	0.41
1:C:345:PHE:HZ	3:C:594:HOH:O	2.01	0.41
1:C:264:CYS:SG	1:C:512:HIS:CD2	3.13	0.41
1:A:211:TYR:CZ	1:A:517:PRO:HG2	2.55	0.41
1:C:426:LYS:HE2	1:C:430:GLU:OE1	2.20	0.41
1:A:455:LEU:HD12	1:A:455:LEU:N	2.35	0.41
1:C:436:VAL:HA	1:C:437:PRO:HD3	1.93	0.41
1:C:425:GLU:CB	3:C:561:HOH:O	2.68	0.41
1:B:257:ILE:CD1	1:B:314:PHE:CZ	3.03	0.41
1:A:426:LYS:HE2	1:A:430:GLU:OE1	2.21	0.41
1:C:217:ALA:HB3	1:C:454:LYS:HE2	2.03	0.41
1:B:328:ALA:HB1	1:B:332:LEU:HD12	2.01	0.41
1:B:486:TRP:HB3	3:B:577:HOH:O	2.21	0.41
1:B:324:ARG:HA	1:B:324:ARG:HD3	1.84	0.41
1:C:324:ARG:HA	1:C:324:ARG:HD3	1.84	0.41
1:B:414:GLN:NE2	3:B:592:HOH:O	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:257:ILE:CD1	1:C:306:PRO:HD3	2.49	0.41
1:B:411:THR:HG21	3:B:160:HOH:O	2.19	0.41
1:B:480:LYS:NZ	1:B:480:LYS:CB	2.83	0.41
1:C:374:ARG:NE	3:C:663:HOH:O	2.54	0.41
1:B:217:ALA:HB3	1:B:454:LYS:HE2	2.03	0.41
1:A:406:GLN:HE21	1:A:406:GLN:HB3	1.71	0.41
1:B:406:GLN:HE21	1:B:406:GLN:HB3	1.71	0.41
1:A:318:GLN:HG3	1:A:321:ARG:HH22	1.85	0.41
1:B:467:ARG:HG2	1:B:515:GLN:HB2	2.02	0.41
1:A:257:ILE:CG1	1:A:303:LEU:HD21	2.51	0.41
1:A:354:VAL:CG2	1:A:362:VAL:HG22	2.48	0.41
1:B:455:LEU:HD12	1:B:455:LEU:N	2.35	0.41
1:A:328:ALA:HB1	1:A:332:LEU:HD12	2.03	0.41
1:A:226:GLN:HB2	1:A:229:GLN:HG3	2.03	0.41
1:A:480:LYS:CB	1:A:480:LYS:NZ	2.84	0.41
1:B:447:HIS:CE1	1:B:449:PHE:HB2	2.56	0.41
1:C:473:ASP:N	1:C:480:LYS:HB2	2.31	0.40
1:C:480:LYS:HZ2	1:C:480:LYS:HB3	1.86	0.40
1:B:340:PRO:HA	1:B:341:PRO:HD3	1.75	0.40
1:A:447:HIS:CE1	1:A:449:PHE:HB2	2.56	0.40
1:B:247:ILE:HA	1:B:254:MET:O	2.21	0.40
1:A:371:VAL:HG23	1:A:389:TRP:CH2	2.56	0.40
1:B:257:ILE:CD1	1:B:306:PRO:HD3	2.48	0.40
1:C:402:TRP:CE2	1:C:406:GLN:HG3	2.56	0.40
1:C:400:VAL:HG13	1:C:424:TRP:CD1	2.57	0.40
1:C:425:GLU:HB2	3:C:561:HOH:O	2.22	0.40
1:C:251:HIS:HB3	1:C:384:ASP:OD2	2.21	0.40
1:B:391:ASP:OD1	1:B:444:ARG:NH2	2.54	0.40
1:B:490:TYR:HB2	3:B:573:HOH:O	2.21	0.40
1:C:248:ASP:HB2	1:C:256:PHE:CE2	2.56	0.40
1:C:226:GLN:HB2	1:C:229:GLN:HG3	2.04	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	305/345 (88%)	289 (95%)	15 (5%)	1 (0%)	46	68
1	B	305/345 (88%)	292 (96%)	12 (4%)	1 (0%)	46	68
1	C	305/345 (88%)	291 (95%)	13 (4%)	1 (0%)	46	68
All	All	915/1035 (88%)	872 (95%)	40 (4%)	3 (0%)	46	68

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	383	ASP
1	A	383	ASP
1	B	383	ASP

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	273/301 (91%)	264 (97%)	9 (3%)	45	73
1	B	273/301 (91%)	263 (96%)	10 (4%)	41	68
1	C	273/301 (91%)	264 (97%)	9 (3%)	45	73
All	All	819/903 (91%)	791 (97%)	28 (3%)	44	72

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	244	LEU
1	A	247	ILE
1	A	268	LEU
1	A	317	SER
1	A	326	GLN
1	A	351	LEU
1	A	360	SER

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Mol	Chain	Res	Type
1	A	408	LYS
1	A	455	LEU
1	B	244	LEU
1	B	247	ILE
1	B	268	LEU
1	B	317	SER
1	B	326	GLN
1	B	351	LEU
1	B	360	SER
1	B	405	LYS
1	B	408	LYS
1	B	455	LEU
1	C	244	LEU
1	C	247	ILE
1	C	268	LEU
1	C	307	LYS
1	C	317	SER
1	C	351	LEU
1	C	360	SER
1	C	408	LYS
1	C	455	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (34) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	215	GLN
1	A	233	HIS
1	A	251	HIS
1	A	284	ASN
1	A	301	HIS
1	A	326	GLN
1	A	336	GLN
1	A	406	GLN
1	A	421	ASN
1	A	477	HIS
1	A	482	HIS
1	A	512	HIS
1	B	215	GLN
1	B	233	HIS
1	B	251	HIS
1	B	284	ASN
1	B	301	HIS

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Mol	Chain	Res	Type
1	B	326	GLN
1	B	406	GLN
1	B	421	ASN
1	B	477	HIS
1	B	482	HIS
1	B	512	HIS
1	C	215	GLN
1	C	284	ASN
1	C	301	HIS
1	C	304	GLN
1	C	326	GLN
1	C	331	HIS
1	C	358	ASN
1	C	406	GLN
1	C	421	ASN
1	C	482	HIS
1	C	512	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues are modelled in this entry.

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	MLZ	D	382	-	8,9,10	0.61	0	7,9,11	1.17	1 (14%)
2	MLZ	E	382	-	8,9,10	0.64	0	7,9,11	1.33	1 (14%)

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	MLZ	D	382	-	-	0/6/8/10	0/0/0/0
2	MLZ	E	382	-	-	0/6/8/10	0/0/0/0

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	D	382	MLZ	CM-NZ-CE	2.40	119.26	112.23
2	E	382	MLZ	CM-NZ-CE	2.70	120.14	112.23

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	382	MLZ	3	0
2	E	382	MLZ	4	0

5.5 Carbohydrates [i](#)

There are no carbohydrates 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, 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	309/345 (89%)	0.90	22 (7%) 19 21	27, 44, 62, 72	0
1	B	309/345 (89%)	1.28	61 (19%) 1 1	26, 44, 62, 71	0
1	C	309/345 (89%)	1.13	47 (15%) 3 3	27, 45, 63, 73	0
2	D	0/10	-	-	-	-
2	E	0/10	-	-	-	-
All	All	927/1055 (87%)	1.10	130 (14%) 4 3	26, 44, 62, 73	0

All (130) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	206	TRP	6.6
1	B	218	ILE	6.1
1	A	465	LEU	5.4
1	B	438	THR	5.2
1	B	468	VAL	4.8
1	C	407	GLY	4.7
1	C	420	ASP	4.5
1	C	483	PHE	4.4
1	B	343	LEU	4.4
1	B	222	VAL	4.3
1	B	514	LEU	4.3
1	C	410	LEU	4.3
1	B	278	CYS	4.1
1	A	343	LEU	4.0
1	B	458	VAL	3.9
1	B	212	LEU	3.8
1	A	486	TRP	3.7
1	C	358	ASN	3.6
1	B	504	ALA	3.6
1	C	479	ILE	3.5
1	C	480	LYS	3.5

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Mol	Chain	Res	Type	RSRZ
1	A	443	VAL	3.5
1	C	342	PRO	3.5
1	A	490	TYR	3.4
1	B	409	PRO	3.4
1	C	344	GLY	3.3
1	C	418	ASP	3.3
1	C	421	ASN	3.2
1	C	390	CYS	3.2
1	C	215	GLN	3.2
1	B	460	ARG	3.2
1	B	489	GLY	3.2
1	C	489	GLY	3.2
1	B	370	VAL	3.1
1	A	517	PRO	3.1
1	C	409	PRO	3.1
1	C	465	LEU	3.1
1	C	449	PHE	3.0
1	C	481	ILE	3.0
1	A	207	SER	3.0
1	B	231	VAL	3.0
1	A	215	GLN	3.0
1	B	415	ASP	3.0
1	C	472	GLU	3.0
1	B	485	GLY	3.0
1	C	217	ALA	3.0
1	B	433	ALA	2.9
1	A	479	ILE	2.9
1	B	207	SER	2.9
1	B	373	SER	2.9
1	B	295	TRP	2.9
1	B	482	HIS	2.9
1	B	407	GLY	2.8
1	C	442	LYS	2.8
1	B	481	ILE	2.8
1	C	488	HIS	2.8
1	A	216	LYS	2.8
1	B	437	PRO	2.8
1	C	490	TYR	2.7
1	C	206	TRP	2.7
1	B	342	PRO	2.7
1	A	325	ALA	2.7
1	B	443	VAL	2.7

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Mol	Chain	Res	Type	RSRZ
1	C	321	ARG	2.7
1	C	444	ARG	2.7
1	B	451	VAL	2.7
1	C	264	CYS	2.6
1	B	208	TRP	2.6
1	B	462	ASN	2.6
1	B	244	LEU	2.6
1	B	491	ASP	2.6
1	A	222	VAL	2.5
1	B	374	ARG	2.5
1	B	450	LEU	2.5
1	A	458	VAL	2.5
1	B	445	PRO	2.5
1	C	415	ASP	2.5
1	B	457	ALA	2.5
1	B	296	PHE	2.5
1	B	479	ILE	2.5
1	B	513	PRO	2.4
1	A	402	TRP	2.4
1	B	213	GLU	2.4
1	C	443	VAL	2.4
1	C	466	ILE	2.4
1	A	345	PHE	2.4
1	B	302	LYS	2.4
1	B	385	THR	2.4
1	C	354	VAL	2.4
1	B	224	LEU	2.4
1	C	219	THR	2.4
1	B	220	ALA	2.4
1	C	294	GLY	2.4
1	B	298	LYS	2.4
1	B	323	THR	2.4
1	C	493	TRP	2.3
1	B	216	LYS	2.3
1	C	426	LYS	2.3
1	C	222	VAL	2.3
1	C	439	TRP	2.3
1	B	382	TRP	2.3
1	B	486	TRP	2.3
1	C	412	PRO	2.2
1	C	364	VAL	2.2
1	C	477	HIS	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	260	VAL	2.2
1	C	353	ALA	2.2
1	B	470	SER	2.2
1	A	485	GLY	2.2
1	B	417	PRO	2.2
1	C	487	SER	2.1
1	A	442	LYS	2.1
1	B	397	ILE	2.1
1	B	226	GLN	2.1
1	C	291	HIS	2.1
1	A	416	TYR	2.1
1	B	211	TYR	2.1
1	C	396	TYR	2.1
1	A	415	ASP	2.1
1	B	420	ASP	2.1
1	B	294	GLY	2.1
1	A	316	TRP	2.1
1	C	330	LYS	2.1
1	C	343	LEU	2.1
1	B	434	SER	2.1
1	B	215	GLN	2.1
1	B	494	ILE	2.1
1	A	428	LEU	2.0
1	C	492	PHE	2.0
1	A	295	TRP	2.0

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

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	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	MLZ	D	382	10/11	0.77	0.24	-	27,36,49,53	0
2	MLZ	E	382	10/11	0.78	0.26	-	30,44,62,68	0

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.